



**Groundwater Sample Results, Level 2 Laboratory
Report, Level 4 Laboratory Report, Electronic Data
Deliverable, Data Validation Report, Sample Location
Report, SDG J237831**

*Bay Head Road Annex
NSWC Annapolis
Maryland*

December 2020

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

ANALYTICAL REPORT

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TestAmerica Job ID: 320-23783-1

Client Project/Site: Former Bay Head Road Annex (RFP JU06-01)

For:
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Authorized for release by:
12/9/2016 4:30:43 PM

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The test results in this report meet all 2003 NELAC and 2009 TNI requirements for accredited parameters, exceptions are noted in this report. This report may not be reproduced except in full, and with written approval from the laboratory. For questions please contact the Project Manager at the e-mail address or telephone number listed on this page.

This report has been electronically signed and authorized by the signatory. Electronic signature is intended to be the legally binding equivalent of a traditionally handwritten signature.

Results relate only to the items tested and the sample(s) as received by the laboratory.

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Definitions/Glossary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Qualifiers

LCMS

Qualifier	Qualifier Description
J	Estimated: The analyte was positively identified; the quantitation is an estimation
M	Manual integrated compound.
U	Undetected at the Limit of Detection.
Q	One or more quality control criteria failed.
E	Result exceeded calibration range.
D	The reported value is from a dilution.

Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
α	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
CFL	Contains Free Liquid
CNF	Contains no Free Liquid
DER	Duplicate error ratio (normalized absolute difference)
Dil Fac	Dilution Factor
DL, RA, RE, IN	Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample
DLC	Decision level concentration
MDA	Minimum detectable activity
EDL	Estimated Detection Limit
MDC	Minimum detectable concentration
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
NC	Not Calculated
ND	Not detected at the reporting limit (or MDL or EDL if shown)
PQL	Practical Quantitation Limit
QC	Quality Control
RER	Relative error ratio
RL	Reporting Limit or Requested Limit (Radiochemistry)
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)

Case Narrative

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Job ID: 320-23783-1

Laboratory: TestAmerica Sacramento

Narrative

CASE NARRATIVE

Client: AECOM Technical Services Inc.

Project: Former Bay Head Road Annex (RFP JU06-01)

Report Number: 320-23783-1

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica West Sacramento attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the applicable methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

TestAmerica utilizes USEPA approved methods and DOD QSM, where applicable, in all analytical work. The samples presented in this report were analyzed for the parameter(s) listed on the analytical methods summary page in accordance with the method(s) indicated. A summary of QC data for these analyses is included at the back of the report.

All parameters for which TestAmerica West Sacramento has certification were evaluated to the QSM specified reporting convention or to the client specified format if different from QSM. Parameters not certified under QSM, if any, were evaluated to the detection limit (DL) and include qualified results where applicable.

The sample(s) that contain constituents flagged with U are undetected. The result associated with this flag is the limit of detection (LOD).

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

RECEIPT

The samples were received on 11/23/2016; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 1.5 C.

Receipt Exceptions

The container label for the following sample did not match the information listed on the Chain-of-Custody (COC): DPT-16-28-SO-14-15 (320-23783-10). The container label lists DPT-16-22-SO-14-15, while the COC lists DPT-16-28-SO-14-15. The sample was logged in per the COC.

PFAS

The first level standard from the initial calibration curve is used to evaluate the tune criteria. The instrument mass windows are set at +/- 0.5amu; therefore, detection of the analyte serves as verification that the assigned mass is within +/- 0.5amu of the true value, which meets the DoD/DOE QSM tune criterion.

Case Narrative

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Job ID: 320-23783-1 (Continued)

Laboratory: TestAmerica Sacramento (Continued)

The concentration of one or more analytes associated with the following samples exceeded the instrument calibration range: DPT-16-35-GW-21-25 (320-23783-3), DPT-16-35-GW-31-35 (320-23783-4), DPT-16-19-SO-00-01 (320-23783-5), DPT-16-15-GW-18-22 (320-23783-12) and DPT-16-15-GW-31-35 (320-23783-13). These samples have been run at dilution and both sets of data were reported.

The Isotope Dilution Analyte (IDA) recoveries for $^{18}\text{O}_2$ PFHxS and $^{13}\text{C}_4$ PFOA associated with the following sample is below the method recommended limit: DPT-16-15-GW-18-22 (320-23783-12). Generally, data quality is not considered affected if the IDA signal-to-noise ratio is greater than 10:1, which is achieved for all IDA in the sample.

Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with preparation batch 320-139749.

Due to excessive amount of sediment in the sample bottles, the aqueous portion of these samples was decanted to new bottles prior to spiking and the extraction. DPT-16-35-GW-21-25 (320-23783-3), DPT-16-35-GW-31-35 (320-23783-4), DPT-16-15-GW-18-22 (320-23783-12) and DPT-16-15-GW-31-35 (320-23783-13)

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

PERCENT SOLIDS

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

Detection Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Lab Sample ID: 320-23783-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	12	M	0.58	0.12	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	28		0.58	0.15	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.21	J	0.46	0.12	ug/Kg	1	☒	537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-SO-19-20

Lab Sample ID: 320-23783-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.56	J	0.60	0.12	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	4.0	M	0.60	0.15	ug/Kg	1	☒	537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-GW-21-25

Lab Sample ID: 320-23783-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.60	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	1.9	E M	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.28		0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	0.76	D	0.025	0.0075	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	2.4	D	0.040	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.19	D	0.025	0.0092	ug/L	10		537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-GW-31-35

Lab Sample ID: 320-23783-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	1.1	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	0.72	E	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.58	E	0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	1.4	D	0.025	0.0075	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	0.80	D	0.040	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.38	D	0.025	0.0092	ug/L	10		537 (Modified)	Total/NA

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	3.8		0.57	0.12	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	130	M E	0.57	0.14	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.18	J	0.46	0.12	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	3.5	J D M	5.7	1.2	ug/Kg	10	☒	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	170	D	5.7	1.4	ug/Kg	10	☒	537 (Modified)	Total/NA

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.37	J	0.62	0.13	ug/Kg	1	☒	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	6.5	M	0.62	0.16	ug/Kg	1	☒	537 (Modified)	Total/NA

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

Detection Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-32-SO-14-15 (Continued)

Lab Sample ID: 320-23783-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	1.1		0.60	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	46	M	0.60	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.27	J	0.60	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	11	M	0.60	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	5.5		0.59	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	57	M	0.59	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-28-SO-14-15

Lab Sample ID: 320-23783-10

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.45	J	0.58	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	10	M	0.58	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-SO-14-15

Lab Sample ID: 320-23783-11

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.51	J	0.61	0.13	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	11	M	0.61	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-GW-18-22

Lab Sample ID: 320-23783-12

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	5.8	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	6.5	E M	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	1.2	E	0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	15	D	0.25	0.075	ug/L	100		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	12	D	0.40	0.13	ug/L	100		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.54	D	0.25	0.092	ug/L	100		537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-GW-31-35

Lab Sample ID: 320-23783-13

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.27	M	0.0025	0.00076	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	0.46	E	0.0041	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.048		0.0025	0.00093	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	0.27	D	0.025	0.0076	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	0.45	D	0.041	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.035	D	0.025	0.0093	ug/L	10		537 (Modified)	Total/NA

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24
 Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid
 Percent Solids: 86.4

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	12	M	0.58	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Perfluorooctane Sulfonate (PFOS)	28		0.58	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Perfluorobutanesulfonic acid (PFBS)	0.21	J	0.46	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	93		25 - 150				11/28/16 11:03	12/02/16 15:52	1
13C4 PFOS	73		25 - 150				11/28/16 11:03	12/02/16 15:52	1
18O2 PFHxS	86		25 - 150				11/28/16 11:03	12/02/16 15:52	1

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36
 Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid
 Percent Solids: 82.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.56	J	0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Perfluorooctane Sulfonate (PFOS)	4.0	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	112		25 - 150				11/28/16 11:03	12/02/16 15:59	1
13C4 PFOS	101		25 - 150				11/28/16 11:03	12/02/16 15:59	1
18O2 PFHxS	108		25 - 150				11/28/16 11:03	12/02/16 15:59	1

Client Sample ID: DPT-16-35-GW-21-25

Date Collected: 11/22/16 10:10
 Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-3

Matrix: Water

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.60	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:26	1
Perfluorooctane Sulfonate (PFOS)	1.9	E M	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:26	1
Perfluorobutanesulfonic acid (PFBS)	0.28		0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:26	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	70		25 - 150				11/28/16 16:55	12/03/16 16:26	1
13C4 PFOS	58		25 - 150				11/28/16 16:55	12/03/16 16:26	1
18O2 PFHxS	43		25 - 150				11/28/16 16:55	12/03/16 16:26	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.76	D	0.025	0.0075	ug/L		11/28/16 16:55	12/07/16 17:10	10
Perfluorooctane Sulfonate (PFOS)	2.4	D	0.040	0.013	ug/L		11/28/16 16:55	12/07/16 17:10	10
Perfluorobutanesulfonic acid (PFBS)	0.19	D	0.025	0.0092	ug/L		11/28/16 16:55	12/07/16 17:10	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	91		25 - 150				11/28/16 16:55	12/07/16 17:10	10
13C4 PFOS	87		25 - 150				11/28/16 16:55	12/07/16 17:10	10
18O2 PFHxS	83		25 - 150				11/28/16 16:55	12/07/16 17:10	10

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-GW-31-35

Lab Sample ID: 320-23783-4

Date Collected: 11/22/16 10:30

Matrix: Water

Date Received: 11/23/16 08:46

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.1	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:33	1
Perfluorooctane Sulfonate (PFOS)	0.72	E	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:33	1
Perfluorobutanesulfonic acid (PFBS)	0.58	E	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:33	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	52		25 - 150				11/28/16 16:55	12/03/16 16:33	1
13C4 PFOS	76		25 - 150				11/28/16 16:55	12/03/16 16:33	1
18O2 PFHxS	30		25 - 150				11/28/16 16:55	12/03/16 16:33	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.4	D	0.025	0.0075	ug/L		11/28/16 16:55	12/07/16 17:18	10
Perfluorooctane Sulfonate (PFOS)	0.80	D	0.040	0.013	ug/L		11/28/16 16:55	12/07/16 17:18	10
Perfluorobutanesulfonic acid (PFBS)	0.38	D	0.025	0.0092	ug/L		11/28/16 16:55	12/07/16 17:18	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	78		25 - 150				11/28/16 16:55	12/07/16 17:18	10
13C4 PFOS	91		25 - 150				11/28/16 16:55	12/07/16 17:18	10
18O2 PFHxS	73		25 - 150				11/28/16 16:55	12/07/16 17:18	10

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Date Collected: 11/22/16 11:20

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 86.7

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	3.8		0.57	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Perfluorooctane Sulfonate (PFOS)	130	M E	0.57	0.14	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Perfluorobutanesulfonic acid (PFBS)	0.18	J	0.46	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	96		25 - 150				11/28/16 11:03	12/02/16 16:07	1
13C4 PFOS	54		25 - 150				11/28/16 11:03	12/02/16 16:07	1
18O2 PFHxS	85		25 - 150				11/28/16 11:03	12/02/16 16:07	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	3.5	J D M	5.7	1.2	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Perfluorooctane Sulfonate (PFOS)	170	D	5.7	1.4	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Perfluorobutanesulfonic acid (PFBS)	3.4	U	4.6	1.2	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	136		25 - 150				11/28/16 11:03	12/07/16 17:33	10
13C4 PFOS	90		25 - 150				11/28/16 11:03	12/07/16 17:33	10
18O2 PFHxS	106		25 - 150				11/28/16 11:03	12/07/16 17:33	10

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Date Collected: 11/22/16 11:30

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 79.7

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.37	J	0.62	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Perfluorooctane Sulfonate (PFOS)	6.5	M	0.62	0.16	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.50	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	69		25 - 150				11/28/16 11:03	12/02/16 16:44	1
13C4 PFOS	64		25 - 150				11/28/16 11:03	12/02/16 16:44	1
18O2 PFHxS	73		25 - 150				11/28/16 11:03	12/02/16 16:44	1

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

Date Collected: 11/21/16 09:36

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 82.3

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.1		0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Perfluorooctane Sulfonate (PFOS)	46	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	96		25 - 150				11/28/16 11:03	12/02/16 16:52	1
13C4 PFOS	62		25 - 150				11/28/16 11:03	12/02/16 16:52	1
18O2 PFHxS	70		25 - 150				11/28/16 11:03	12/02/16 16:52	1

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Date Collected: 11/21/16 10:12

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	J	0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Perfluorooctane Sulfonate (PFOS)	11	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	108		25 - 150				11/28/16 11:03	12/02/16 16:59	1
13C4 PFOS	82		25 - 150				11/28/16 11:03	12/02/16 16:59	1
18O2 PFHxS	93		25 - 150				11/28/16 11:03	12/02/16 16:59	1

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Date Collected: 11/21/16 10:46

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	5.5		0.59	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Perfluorooctane Sulfonate (PFOS)	57	M	0.59	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	101		25 - 150				11/28/16 11:03	12/02/16 17:07	1
13C4 PFOS	63		25 - 150				11/28/16 11:03	12/02/16 17:07	1
18O2 PFHxS	90		25 - 150				11/28/16 11:03	12/02/16 17:07	1

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-28-SO-14-15

Date Collected: 11/21/16 11:54

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-10

Matrix: Solid

Percent Solids: 85.2

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.45	J	0.58	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Perfluorooctane Sulfonate (PFOS)	10	M	0.58	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Perfluorobutanesulfonic acid (PFBS)	0.35	U	0.47	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	105		25 - 150				11/28/16 11:03	12/02/16 17:14	1
13C4 PFOS	61		25 - 150				11/28/16 11:03	12/02/16 17:14	1
18O2 PFHxS	89		25 - 150				11/28/16 11:03	12/02/16 17:14	1

Client Sample ID: DPT-16-15-SO-14-15

Date Collected: 11/21/16 12:50

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-11

Matrix: Solid

Percent Solids: 80.6

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.51	J	0.61	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Perfluorooctane Sulfonate (PFOS)	11	M	0.61	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.49	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	101		25 - 150				11/28/16 11:03	12/02/16 17:22	1
13C4 PFOS	66		25 - 150				11/28/16 11:03	12/02/16 17:22	1
18O2 PFHxS	89		25 - 150				11/28/16 11:03	12/02/16 17:22	1

Client Sample ID: DPT-16-15-GW-18-22

Date Collected: 11/21/16 13:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-12

Matrix: Water

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	5.8	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:41	1
Perfluorooctane Sulfonate (PFOS)	6.5	E M	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:41	1
Perfluorobutanesulfonic acid (PFBS)	1.2	E	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:41	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	23	Q	25 - 150				11/28/16 16:55	12/03/16 16:41	1
13C4 PFOS	29		25 - 150				11/28/16 16:55	12/03/16 16:41	1
18O2 PFHxS	21	Q	25 - 150				11/28/16 16:55	12/03/16 16:41	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	15	D	0.25	0.075	ug/L		11/28/16 16:55	12/07/16 17:03	100
Perfluorooctane Sulfonate (PFOS)	12	D	0.40	0.13	ug/L		11/28/16 16:55	12/07/16 17:03	100
Perfluorobutanesulfonic acid (PFBS)	0.54	D	0.25	0.092	ug/L		11/28/16 16:55	12/07/16 17:03	100
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	83		25 - 150				11/28/16 16:55	12/07/16 17:03	100
13C4 PFOS	94		25 - 150				11/28/16 16:55	12/07/16 17:03	100
18O2 PFHxS	104		25 - 150				11/28/16 16:55	12/07/16 17:03	100

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-15-GW-31-35

Lab Sample ID: 320-23783-13

Date Collected: 11/21/16 14:00

Matrix: Water

Date Received: 11/23/16 08:46

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	M	0.0025	0.00076	ug/L		11/28/16 16:55	12/03/16 16:48	1
Perfluorooctane Sulfonate (PFOS)	0.46	E	0.0041	0.0013	ug/L		11/28/16 16:55	12/03/16 16:48	1
Perfluorobutanesulfonic acid (PFBS)	0.048		0.0025	0.00093	ug/L		11/28/16 16:55	12/03/16 16:48	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	22	Q	25 - 150				11/28/16 16:55	12/03/16 16:48	1
13C4 PFOS	82		25 - 150				11/28/16 16:55	12/03/16 16:48	1
18O2 PFHxS	66		25 - 150				11/28/16 16:55	12/03/16 16:48	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	D	0.025	0.0076	ug/L		11/28/16 16:55	12/07/16 17:25	10
Perfluorooctane Sulfonate (PFOS)	0.45	D	0.041	0.013	ug/L		11/28/16 16:55	12/07/16 17:25	10
Perfluorobutanesulfonic acid (PFBS)	0.035	D	0.025	0.0093	ug/L		11/28/16 16:55	12/07/16 17:25	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	33		25 - 150				11/28/16 16:55	12/07/16 17:25	10
13C4 PFOS	95		25 - 150				11/28/16 16:55	12/07/16 17:25	10
18O2 PFHxS	101		25 - 150				11/28/16 16:55	12/07/16 17:25	10

Isotope Dilution Summary

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Matrix: Solid

Prep Type: Total/NA

Percent Isotope Dilution Recovery (Acceptance Limits)

Lab Sample ID	Client Sample ID	Percent Isotope Dilution Recovery (Acceptance Limits)		
		3C4 PFO/ (25-150)	3C4 PFO: (25-150)	3O2 PFHx (25-150)
320-23783-1	DPT-16-35-SO-00-01	93	73	86
320-23783-2	DPT-16-35-SO-19-20	112	101	108
320-23783-5	DPT-16-19-SO-00-01	96	54	85
320-23783-5 - DL	DPT-16-19-SO-00-01	136	90	106
320-23783-6	DPT-16-19-SO-17-18	69	64	73
320-23783-7	DPT-16-32-SO-14-15	96	62	70
320-23783-8	DPT-16-31-SO-14-15	108	82	93
320-23783-9	DPT-16-30-SO-14-15	101	63	90
320-23783-10	DPT-16-28-SO-14-15	105	61	89
320-23783-11	DPT-16-15-SO-14-15	101	66	89
LCS 320-139627/2-A	Lab Control Sample	119	112	120
MB 320-139627/1-A	Method Blank	108	94	104

Surrogate Legend

13C4 PFOA = 13C4 PFOA
 13C4 PFOS = 13C4 PFOS
 18O2 PFHxS = 18O2 PFHxS

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Matrix: Water

Prep Type: Total/NA

Percent Isotope Dilution Recovery (Acceptance Limits)

Lab Sample ID	Client Sample ID	Percent Isotope Dilution Recovery (Acceptance Limits)		
		3C4 PFO/ (25-150)	3C4 PFO: (25-150)	3O2 PFHx (25-150)
320-23783-3	DPT-16-35-GW-21-25	70	58	43
320-23783-3 - DL	DPT-16-35-GW-21-25	91	87	83
320-23783-4	DPT-16-35-GW-31-35	52	76	30
320-23783-4 - DL	DPT-16-35-GW-31-35	78	91	73
320-23783-12	DPT-16-15-GW-18-22	23 Q	29	21 Q
320-23783-12 - DL	DPT-16-15-GW-18-22	83	94	104
320-23783-13	DPT-16-15-GW-31-35	22 Q	82	66
320-23783-13 - DL	DPT-16-15-GW-31-35	33	95	101
LCS 320-139749/2-A	Lab Control Sample	108	104	105
LCSD 320-139749/3-A	Lab Control Sample Dup	116	104	104
MB 320-139749/1-A	Method Blank	93	85	86

Surrogate Legend

13C4 PFOA = 13C4 PFOA
 13C4 PFOS = 13C4 PFOS
 18O2 PFHxS = 18O2 PFHxS

QC Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Lab Sample ID: MB 320-139627/1-A
Matrix: Solid
Analysis Batch: 140429

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 139627

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.30	U	0.50	0.10	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Perfluorooctane Sulfonate (PFOS)	0.30	U	0.50	0.13	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Perfluorobutanesulfonic acid (PFBS)	0.30	U	0.40	0.10	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Isotope Dilution	MB %Recovery	MB Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	108		25 - 150				11/28/16 11:03	12/02/16 14:59	1
13C4 PFOS	94		25 - 150				11/28/16 11:03	12/02/16 14:59	1
18O2 PFHxS	104		25 - 150				11/28/16 11:03	12/02/16 14:59	1

Lab Sample ID: LCS 320-139627/2-A
Matrix: Solid
Analysis Batch: 140429

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 139627

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	Limits
Perfluorooctanoic acid (PFOA)	4.00	4.25		ug/Kg		106	60 - 140
Perfluorooctane Sulfonate (PFOS)	3.71	3.66		ug/Kg		99	60 - 140
Perfluorobutanesulfonic acid (PFBS)	3.54	4.03		ug/Kg		114	50 - 150
Isotope Dilution	LCS %Recovery	LCS Qualifier	Limits				
13C4 PFOA	119		25 - 150				
13C4 PFOS	112		25 - 150				
18O2 PFHxS	120		25 - 150				

Lab Sample ID: MB 320-139749/1-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 139749

Analyte	MB Result	MB Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:03	1
Perfluorooctane Sulfonate (PFOS)	0.0030	U	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:03	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:03	1
Isotope Dilution	MB %Recovery	MB Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	93		25 - 150				11/28/16 16:55	12/03/16 16:03	1
13C4 PFOS	85		25 - 150				11/28/16 16:55	12/03/16 16:03	1
18O2 PFHxS	86		25 - 150				11/28/16 16:55	12/03/16 16:03	1

Lab Sample ID: LCS 320-139749/2-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 139749

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	Limits
Perfluorooctanoic acid (PFOA)	0.0400	0.0296		ug/L		74	60 - 140
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0261		ug/L		70	60 - 140
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0289		ug/L		82	50 - 150

TestAmerica Sacramento

QC Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

<i>Isotope Dilution</i>	<i>%Recovery</i>	<i>Qualifier</i>	<i>Limits</i>
13C4 PFOA	108		25 - 150
13C4 PFOS	104		25 - 150
18O2 PFHxS	105		25 - 150

Lab Sample ID: LCSD 320-139749/3-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 139749

<i>Analyte</i>	<i>Spike Added</i>	<i>LCSD Result</i>	<i>LCSD Qualifier</i>	<i>Unit</i>	<i>D</i>	<i>%Rec</i>	<i>%Rec. Limits</i>	<i>RPD</i>	<i>RPD Limit</i>
Perfluorooctanoic acid (PFOA)	0.0400	0.0314		ug/L		79	60 - 140	6	30
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0289		ug/L		78	60 - 140	10	30
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0325		ug/L		92	50 - 150	12	30

<i>Isotope Dilution</i>	<i>%Recovery</i>	<i>Qualifier</i>	<i>Limits</i>
13C4 PFOA	116		25 - 150
13C4 PFOS	104		25 - 150
18O2 PFHxS	104		25 - 150



QC Association Summary

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

LCMS

Prep Batch: 139627

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	SHAKE	
320-23783-5 - DL	DPT-16-19-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	SHAKE	
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	SHAKE	
MB 320-139627/1-A	Method Blank	Total/NA	Solid	SHAKE	
LCS 320-139627/2-A	Lab Control Sample	Total/NA	Solid	SHAKE	

Prep Batch: 139749

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3 - DL	DPT-16-35-GW-21-25	Total/NA	Water	3535	
320-23783-3	DPT-16-35-GW-21-25	Total/NA	Water	3535	
320-23783-4 - DL	DPT-16-35-GW-31-35	Total/NA	Water	3535	
320-23783-4	DPT-16-35-GW-31-35	Total/NA	Water	3535	
320-23783-12	DPT-16-15-GW-18-22	Total/NA	Water	3535	
320-23783-12 - DL	DPT-16-15-GW-18-22	Total/NA	Water	3535	
320-23783-13	DPT-16-15-GW-31-35	Total/NA	Water	3535	
320-23783-13 - DL	DPT-16-15-GW-31-35	Total/NA	Water	3535	
MB 320-139749/1-A	Method Blank	Total/NA	Water	3535	
LCS 320-139749/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-139749/3-A	Lab Control Sample Dup	Total/NA	Water	3535	

Analysis Batch: 140429

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	537 (Modified)	139627
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	537 (Modified)	139627
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	537 (Modified)	139627
MB 320-139627/1-A	Method Blank	Total/NA	Solid	537 (Modified)	139627
LCS 320-139627/2-A	Lab Control Sample	Total/NA	Solid	537 (Modified)	139627

Analysis Batch: 140564

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3	DPT-16-35-GW-21-25	Total/NA	Water	537 (Modified)	139749
320-23783-4	DPT-16-35-GW-31-35	Total/NA	Water	537 (Modified)	139749
320-23783-12	DPT-16-15-GW-18-22	Total/NA	Water	537 (Modified)	139749
320-23783-13	DPT-16-15-GW-31-35	Total/NA	Water	537 (Modified)	139749
MB 320-139749/1-A	Method Blank	Total/NA	Water	537 (Modified)	139749
LCS 320-139749/2-A	Lab Control Sample	Total/NA	Water	537 (Modified)	139749
LCSD 320-139749/3-A	Lab Control Sample Dup	Total/NA	Water	537 (Modified)	139749

TestAmerica Sacramento

QC Association Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

LCMS (Continued)

Analysis Batch: 141062

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3 - DL	DPT-16-35-GW-21-25	Total/NA	Water	537 (Modified)	139749
320-23783-4 - DL	DPT-16-35-GW-31-35	Total/NA	Water	537 (Modified)	139749
320-23783-5 - DL	DPT-16-19-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-12 - DL	DPT-16-15-GW-18-22	Total/NA	Water	537 (Modified)	139749
320-23783-13 - DL	DPT-16-15-GW-31-35	Total/NA	Water	537 (Modified)	139749

General Chemistry

Analysis Batch: 140129

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	D 2216	
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	D 2216	
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	D 2216	
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	D 2216	

Analysis Batch: 140181

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	D 2216	
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	D 2216	
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	D 2216	
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	D 2216	
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	D 2216	

Lab Chronicle

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid

Percent Solids: 86.4

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.01 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 15:52	SBC	TAL SAC

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid

Percent Solids: 82.8

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.07 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 15:59	SBC	TAL SAC

Client Sample ID: DPT-16-35-GW-21-25

Date Collected: 11/22/16 10:10

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-3

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			250.2 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140564	12/03/16 16:26	SBC	TAL SAC
Total/NA	Prep	3535	DL		250.2 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10			141062	12/07/16 17:10	CBW	TAL SAC

Client Sample ID: DPT-16-35-GW-31-35

Date Collected: 11/22/16 10:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-4

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			249.6 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140564	12/03/16 16:33	SBC	TAL SAC

TestAmerica Sacramento

Lab Chronicle

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-GW-31-35

Lab Sample ID: 320-23783-4

Date Collected: 11/22/16 10:30

Matrix: Water

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535	DL		249.6 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10			141062	12/07/16 17:18	CBW	TAL SAC

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Date Collected: 11/22/16 11:20

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Date Collected: 11/22/16 11:20

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 86.7

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.02 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 16:07	SBC	TAL SAC
Total/NA	Prep	SHAKE	DL		5.02 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10			141062	12/07/16 17:33	CBW	TAL SAC

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Date Collected: 11/22/16 11:30

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Date Collected: 11/22/16 11:30

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 79.7

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.06 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 16:44	SBC	TAL SAC

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

Date Collected: 11/21/16 09:36

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140181	12/01/16 11:23	KSL	TAL SAC

TestAmerica Sacramento

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

Date Collected: 11/21/16 09:36

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 82.3

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.03 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 16:52	SBC	TAL SAC

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Date Collected: 11/21/16 10:12

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Date Collected: 11/21/16 10:12

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.00 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 16:59	SBC	TAL SAC

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Date Collected: 11/21/16 10:46

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Date Collected: 11/21/16 10:46

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.02 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 17:07	SBC	TAL SAC

Client Sample ID: DPT-16-28-SO-14-15

Lab Sample ID: 320-23783-10

Date Collected: 11/21/16 11:54

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140181	12/01/16 11:23	KSL	TAL SAC

TestAmerica Sacramento

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-28-SO-14-15

Lab Sample ID: 320-23783-10

Date Collected: 11/21/16 11:54

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 85.2

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.03 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 17:14	SBC	TAL SAC

Client Sample ID: DPT-16-15-SO-14-15

Lab Sample ID: 320-23783-11

Date Collected: 11/21/16 12:50

Matrix: Solid

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1			140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-15-SO-14-15

Lab Sample ID: 320-23783-11

Date Collected: 11/21/16 12:50

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 80.6

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			5.05 g	1.00 mL	139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140429	12/02/16 17:22	SBC	TAL SAC

Client Sample ID: DPT-16-15-GW-18-22

Lab Sample ID: 320-23783-12

Date Collected: 11/21/16 13:30

Matrix: Water

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			249.7 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140564	12/03/16 16:41	SBC	TAL SAC
Total/NA	Prep	3535	DL		249.7 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	100			141062	12/07/16 17:03	CBW	TAL SAC

Client Sample ID: DPT-16-15-GW-31-35

Lab Sample ID: 320-23783-13

Date Collected: 11/21/16 14:00

Matrix: Water

Date Received: 11/23/16 08:46

Prep Type	Batch Type	Batch Method	Run	Dil Factor	Initial Amount	Final Amount	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			245.9 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1			140564	12/03/16 16:48	SBC	TAL SAC
Total/NA	Prep	3535	DL		245.9 mL	0.5 mL	139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10			141062	12/07/16 17:25	CBW	TAL SAC

Laboratory References:

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600

Certification Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Laboratory: TestAmerica Sacramento

Unless otherwise noted, all analytes for this laboratory were covered under each certification below.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2928-01	01-31-17

The following analytes are included in this report, but certification is not offered by the governing authority:

Analysis Method	Prep Method	Matrix	Analyte
537 (Modified)	SHAKE	Solid	Perfluorooctane Sulfonate (PFOS)
D 2216		Solid	Percent Moisture
D 2216		Solid	Percent Solids



Method Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method	Method Description	Protocol	Laboratory
537 (Modified)	Perfluorinated Hydrocarbons	EPA	TAL SAC
D 2216	Percent Moisture	ASTM	TAL SAC

Protocol References:

ASTM = ASTM International

EPA = US Environmental Protection Agency

Laboratory References:

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600



Sample Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-23783-1	DPT-16-35-SO-00-01	Solid	11/22/16 09:24	11/23/16 08:46
320-23783-2	DPT-16-35-SO-19-20	Solid	11/22/16 09:36	11/23/16 08:46
320-23783-3	DPT-16-35-GW-21-25	Water	11/22/16 10:10	11/23/16 08:46
320-23783-4	DPT-16-35-GW-31-35	Water	11/22/16 10:30	11/23/16 08:46
320-23783-5	DPT-16-19-SO-00-01	Solid	11/22/16 11:20	11/23/16 08:46
320-23783-6	DPT-16-19-SO-17-18	Solid	11/22/16 11:30	11/23/16 08:46
320-23783-7	DPT-16-32-SO-14-15	Solid	11/21/16 09:36	11/23/16 08:46
320-23783-8	DPT-16-31-SO-14-15	Solid	11/21/16 10:12	11/23/16 08:46
320-23783-9	DPT-16-30-SO-14-15	Solid	11/21/16 10:46	11/23/16 08:46
320-23783-10	DPT-16-28-SO-14-15	Solid	11/21/16 11:54	11/23/16 08:46
320-23783-11	DPT-16-15-SO-14-15	Solid	11/21/16 12:50	11/23/16 08:46
320-23783-12	DPT-16-15-GW-18-22	Water	11/21/16 13:30	11/23/16 08:46
320-23783-13	DPT-16-15-GW-31-35	Water	11/21/16 14:00	11/23/16 08:46

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Temperature on Receipt _____

Drinking Water? Yes No

Chain of Custody Record

TAL-4124 (1007)

Client **AECOM** Chain of Custody Number **299019**
 Address **3101 WILSON BLVD** Lab Number **11/22/16**
 City **ARLINGTON** State **VA** Zip Code **22201** Page **1** of **2**

Project Name and Location (State) **Former Bay Head Rd Annex - Annapolis**
 Contract/Purchase Order/Quote No. **32007449 / 60444465**
 Project Manager **Kurt VanGelder** Lab Contact **Jill Kellman**
 Telephone Number (Area Code)/Fax Number **(703) 682-9041** Carrier/Waybill Number **Fedex**

Special Instructions/
Conditions of Receipt

Analysis (Attach list if
more space is needed)

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix				Containers & Preservatives					Analysis (Attach list if more space is needed)	
			Air	Aqueous	Sed	Soil	Unpres	H2SO4	HNO3	HCl	NaOH		ZnAc/NaOH
DPT-16-35-50-00-01	11/22/16	0924	✓				1					PFDA	
DPT-16-35-50-19-20		0936	✓				1					PFOS	
DPT-16-35-50-GW-21-25		1010	✓				2					PFBS	
DPT-16-35-50-GW-31-35		1030	✓				2						
DPT-16-19-50-00-01		1120	✓				1						
DPT-16-19-50-17-18		1130	✓				1						
DPT-16-32-50-14-15	11/21/16	0936	✓				1						
DPT-16-31-50-14-15		1012	✓				1						
DPT-16-30-50-14-15		1046	✓				1						
DPT-16-28-50-14-15		1154	✓				1						
DPT-16-15-50-14-15		1250	✓				1						
DPT-16-15-GW-18-22		1330	✓				2						



320-23783 Chain of Custody

Possible Hazard Identification
 Non-Hazard Flammable Skin Irritant Poison B Unknown
 24 Hours 48 Hours 7 Days 14 Days 21 Days Other _____
 Sample Disposal: Return To Client Disposal By Lab Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)

1. Relinquished By	Date	Time
JMK/Kellman	11/22/16	1400
2. Relinquished By	Date	Time
3. Relinquished By	Date	Time

1. Received By **TRANSFER TO FEDEX** Date **11/22/16** Time **1400**
 2. Received By **Jerry G. Turpen THS** Date **11/25/16** Time **09:30**
 3. Received By

Comments **1.5' ice * Sample label states "DPT-16-22-50-14-15" TAT 11/25/16**

DISTRIBUTION: WHITE - Returned to Client with Report; CANARY - Stays with the Sample; PINK - Field Copy



TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Chain of Custody Record

TAL-4124 (1007)

Temperature on Receipt _____
 Drinking Water? Yes No

Client: **AECOM** Chain of Custody Number: **295762**
 Address: **3101 WILSON BLVD** Date: **11/22/16**
 City: **ARLINGTON VA** Zip Code: **22201** Lab Number: **2** of **2**
 Project Name and Location (State): **Former BxM Henry Rd Annex**
 Contract/Purchase Order/Quote No.: **32007449 / 60444465**

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Containers & Preservatives					Special Instructions/ Conditions of Receipt
			Matrix	U/Pres	H2SO4	HNO3	HCl	
DPY-16-15-GW-31-35	11/21/16	1400	Air					
			soenrthy	✓				✓ PFOA
			Sed					✓ PFOS
			Soil					✓ PFBS

Project Manager: **Kurt Van Gelder**
 Telephone Number (Area Code)/Fax Number: _____
 Site Contact: **MIKE GUMSKI** Lab Contact: **JILL K.**
 Carrier/Waybill Number: **FEDEX**
 Sample Disposal: Return To Client Unknown Poison B Skin Irritant Flammable Non-Hazard Possible Hazard Identification
 Disposal By Lab Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)
 OC Requirements (Specify):
 1. Relinquished By: **11/22/16** Time: **1400**
 2. Relinquished By: **11/25/16** Time: **09:30**
 3. Relinquished By: _____ Time: _____
 Comments: **1.5" x 1.5"**
 DISTRIBUTION: WHITE - Returned to Client with Report; CANARY - Stays with the Sample; PINK - Field Copy

Login Sample Receipt Checklist

Client: AECOM Technical Services Inc.

Job Number: 320-23783-1

Login Number: 23783

List Source: TestAmerica Sacramento

List Number: 1

Creator: Nelson, Kym D

Question	Answer	Comment
Radioactivity wasn't checked or is <=/ background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	CUSTODY SEALS
Sample custody seals, if present, are intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	N/A	
There are no discrepancies between the containers received and the COC.	False	IDs on containers do not match the COC. Logged in per COC.
Samples are received within Holding Time (excluding tests with immediate HTs)	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	N/A	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	



ANALYTICAL REPORT

Job Number: 320-23783-1

Job Description: Former Bay Head Road Annex (RFP JU06-01)

For:
AECOM Technical Services Inc.
3101 Wilson Blvd.
Suite 900
Arlington, VA 22201
Attention: Kurt Vangelder



Approved for release.
Jill Kellmann
Manager of Project Management
12/9/2016 4:36 PM

Jill Kellmann, Manager of Project Management
880 Riverside Parkway, West Sacramento, CA, 95605
(916)374-4402
jill.kellmann@testamericainc.com
12/09/2016

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Definitions/Glossary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Qualifiers

LCMS

Qualifier	Qualifier Description
J	Estimated: The analyte was positively identified; the quantitation is an estimation
M	Manual integrated compound.
U	Undetected at the Limit of Detection.
Q	One or more quality control criteria failed.
E	Result exceeded calibration range.
D	The reported value is from a dilution.

Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
α	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
CFL	Contains Free Liquid
CNF	Contains no Free Liquid
DER	Duplicate error ratio (normalized absolute difference)
Dil Fac	Dilution Factor
DL, RA, RE, IN	Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample
DLC	Decision level concentration
MDA	Minimum detectable activity
EDL	Estimated Detection Limit
MDC	Minimum detectable concentration
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
NC	Not Calculated
ND	Not detected at the reporting limit (or MDL or EDL if shown)
PQL	Practical Quantitation Limit
QC	Quality Control
RER	Relative error ratio
RL	Reporting Limit or Requested Limit (Radiochemistry)
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)

CASE NARRATIVE

Client: AECOM Technical Services Inc.

Project: Former Bay Head Road Annex (RFP JU06-01)

Report Number: 320-23783-1

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

TestAmerica West Sacramento attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the applicable methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

TestAmerica utilizes USEPA approved methods and DOD QSM, where applicable, in all analytical work. The samples presented in this report were analyzed for the parameter(s) listed on the analytical methods summary page in accordance with the method(s) indicated. A summary of QC data for these analyses is included at the back of the report.

All parameters for which TestAmerica West Sacramento has certification were evaluated to the QSM specified reporting convention or to the client specified format if different from QSM. Parameters not certified under QSM, if any, were evaluated to the detection limit (DL) and include qualified results where applicable.

The sample(s) that contain constituents flagged with U are undetected. The result associated with this flag is the limit of detection (LOD).

Calculations are performed before rounding to avoid round-off errors in calculated results.

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

All solid sample results are reported on an "as received" basis unless otherwise indicated by the presence of a % solids value in the method header.

This laboratory report is confidential and is intended for the sole use of TestAmerica and its client.

RECEIPT

The samples were received on 11/23/2016; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 1.5 C.

Receipt Exceptions

The container label for the following sample did not match the information listed on the Chain-of-Custody (COC): DPT-16-28-SO-14-15 (320-23783-10). The container label lists DPT-16-22-SO-14-15, while the COC lists DPT-16-28-SO-14-15. The sample was logged in per the COC.

PFAS

The first level standard from the initial calibration curve is used to evaluate the tune criteria. The instrument mass windows are set at +/- 0.5amu; therefore, detection of the analyte serves as verification that the assigned mass is within +/- 0.5amu of the true value, which meets the DoD/DOE QSM tune criterion.

The concentration of one or more analytes associated with the following samples exceeded the instrument calibration range: DPT-16-35-GW-21-25 (320-23783-3), DPT-16-35-GW-31-35 (320-23783-4), DPT-16-19-SO-00-01 (320-23783-5), DPT-16-15-GW-18-22 (320-23783-12) and DPT-16-15-GW-31-35 (320-23783-13). These samples have been run at dilution and both sets of data were reported.

The Isotope Dilution Analyte (IDA) recoveries for 18O2 PFHxS and 13C4 PFOA associated with the following sample is below the method recommended limit: DPT-16-15-GW-18-22 (320-23783-12). Generally, data quality is not considered affected if the IDA signal-to-noise ratio is greater than 10:1, which is achieved for all IDA in the sample.

Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with preparation batch 320-139749.

Due to excessive amount of sediment in the sample bottles, the aqueous portion of these samples was decanted to new bottles prior to spiking and the extraction. DPT-16-35-GW-21-25 (320-23783-3), DPT-16-35-GW-31-35 (320-23783-4), DPT-16-15-GW-18-22 (320-23783-12) and DPT-16-15-GW-31-35 (320-23783-13)

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

PERCENT SOLIDS

No analytical or quality issues were noted, other than those described above or in the Definitions/Glossary page.

Detection Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Lab Sample ID: 320-23783-1

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	12	M	0.58	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	28		0.58	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.21	J	0.46	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-SO-19-20

Lab Sample ID: 320-23783-2

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.56	J	0.60	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	4.0	M	0.60	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-GW-21-25

Lab Sample ID: 320-23783-3

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.60	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	1.9	E M	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.28		0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	0.76	D	0.025	0.0075	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	2.4	D	0.040	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.19	D	0.025	0.0092	ug/L	10		537 (Modified)	Total/NA

Client Sample ID: DPT-16-35-GW-31-35

Lab Sample ID: 320-23783-4

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	1.1	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	0.72	E	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.58	E	0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	1.4	D	0.025	0.0075	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	0.80	D	0.040	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.38	D	0.025	0.0092	ug/L	10		537 (Modified)	Total/NA

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	3.8		0.57	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	130	M E	0.57	0.14	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.18	J	0.46	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	3.5	J D M	5.7	1.2	ug/Kg	10	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	170	D	5.7	1.4	ug/Kg	10	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.37	J	0.62	0.13	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	6.5	M	0.62	0.16	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

Detection Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-32-SO-14-15 (Continued)

Lab Sample ID: 320-23783-7

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	1.1		0.60	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	46	M	0.60	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.27	J	0.60	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	11	M	0.60	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	5.5		0.59	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	57	M	0.59	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-28-SO-14-15

Lab Sample ID: 320-23783-10

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.45	J	0.58	0.12	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	10	M	0.58	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-SO-14-15

Lab Sample ID: 320-23783-11

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.51	J	0.61	0.13	ug/Kg	1	☼	537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	11	M	0.61	0.15	ug/Kg	1	☼	537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-GW-18-22

Lab Sample ID: 320-23783-12

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	5.8	E M	0.0025	0.00075	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	6.5	E M	0.0040	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	1.2	E	0.0025	0.00092	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	15	D	0.25	0.075	ug/L	100		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	12	D	0.40	0.13	ug/L	100		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.54	D	0.25	0.092	ug/L	100		537 (Modified)	Total/NA

Client Sample ID: DPT-16-15-GW-31-35

Lab Sample ID: 320-23783-13

Analyte	Result	Qualifier	RL	MDL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	0.27	M	0.0025	0.00076	ug/L	1		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS)	0.46	E	0.0041	0.0013	ug/L	1		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS)	0.048		0.0025	0.00093	ug/L	1		537 (Modified)	Total/NA
Perfluorooctanoic acid (PFOA) - DL	0.27	D	0.025	0.0076	ug/L	10		537 (Modified)	Total/NA
Perfluorooctane Sulfonate (PFOS) - DL	0.45	D	0.041	0.013	ug/L	10		537 (Modified)	Total/NA
Perfluorobutanesulfonic acid (PFBS) - DL	0.035	D	0.025	0.0093	ug/L	10		537 (Modified)	Total/NA

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24
Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid
Percent Solids: 86.4

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	12	M	0.58	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Perfluorooctane Sulfonate (PFOS)	28		0.58	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Perfluorobutanesulfonic acid (PFBS)	0.21	J	0.46	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:52	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	93		25 - 150				11/28/16 11:03	12/02/16 15:52	1
13C4 PFOS	73		25 - 150				11/28/16 11:03	12/02/16 15:52	1
18O2 PFHxS	86		25 - 150				11/28/16 11:03	12/02/16 15:52	1

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36
Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid
Percent Solids: 82.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.56	J	0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Perfluorooctane Sulfonate (PFOS)	4.0	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 15:59	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	112		25 - 150				11/28/16 11:03	12/02/16 15:59	1
13C4 PFOS	101		25 - 150				11/28/16 11:03	12/02/16 15:59	1
18O2 PFHxS	108		25 - 150				11/28/16 11:03	12/02/16 15:59	1

Client Sample ID: DPT-16-35-GW-21-25

Date Collected: 11/22/16 10:10
Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-3

Matrix: Water

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.60	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:26	1
Perfluorooctane Sulfonate (PFOS)	1.9	E M	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:26	1
Perfluorobutanesulfonic acid (PFBS)	0.28		0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:26	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	70		25 - 150				11/28/16 16:55	12/03/16 16:26	1
13C4 PFOS	58		25 - 150				11/28/16 16:55	12/03/16 16:26	1
18O2 PFHxS	43		25 - 150				11/28/16 16:55	12/03/16 16:26	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.76	D	0.025	0.0075	ug/L		11/28/16 16:55	12/07/16 17:10	10
Perfluorooctane Sulfonate (PFOS)	2.4	D	0.040	0.013	ug/L		11/28/16 16:55	12/07/16 17:10	10
Perfluorobutanesulfonic acid (PFBS)	0.19	D	0.025	0.0092	ug/L		11/28/16 16:55	12/07/16 17:10	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	91		25 - 150				11/28/16 16:55	12/07/16 17:10	10
13C4 PFOS	87		25 - 150				11/28/16 16:55	12/07/16 17:10	10
18O2 PFHxS	83		25 - 150				11/28/16 16:55	12/07/16 17:10	10

Client Sample Results

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-GW-31-35

Lab Sample ID: 320-23783-4

Date Collected: 11/22/16 10:30

Matrix: Water

Date Received: 11/23/16 08:46

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.1	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:33	1
Perfluorooctane Sulfonate (PFOS)	0.72	E	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:33	1
Perfluorobutanesulfonic acid (PFBS)	0.58	E	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:33	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	52		25 - 150				11/28/16 16:55	12/03/16 16:33	1
13C4 PFOS	76		25 - 150				11/28/16 16:55	12/03/16 16:33	1
18O2 PFHxS	30		25 - 150				11/28/16 16:55	12/03/16 16:33	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.4	D	0.025	0.0075	ug/L		11/28/16 16:55	12/07/16 17:18	10
Perfluorooctane Sulfonate (PFOS)	0.80	D	0.040	0.013	ug/L		11/28/16 16:55	12/07/16 17:18	10
Perfluorobutanesulfonic acid (PFBS)	0.38	D	0.025	0.0092	ug/L		11/28/16 16:55	12/07/16 17:18	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	78		25 - 150				11/28/16 16:55	12/07/16 17:18	10
13C4 PFOS	91		25 - 150				11/28/16 16:55	12/07/16 17:18	10
18O2 PFHxS	73		25 - 150				11/28/16 16:55	12/07/16 17:18	10

Client Sample ID: DPT-16-19-SO-00-01

Lab Sample ID: 320-23783-5

Date Collected: 11/22/16 11:20

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 86.7

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	3.8		0.57	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Perfluorooctane Sulfonate (PFOS)	130	M E	0.57	0.14	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Perfluorobutanesulfonic acid (PFBS)	0.18	J	0.46	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:07	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	96		25 - 150				11/28/16 11:03	12/02/16 16:07	1
13C4 PFOS	54		25 - 150				11/28/16 11:03	12/02/16 16:07	1
18O2 PFHxS	85		25 - 150				11/28/16 11:03	12/02/16 16:07	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	3.5	J D M	5.7	1.2	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Perfluorooctane Sulfonate (PFOS)	170	D	5.7	1.4	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Perfluorobutanesulfonic acid (PFBS)	3.4	U	4.6	1.2	ug/Kg	☼	11/28/16 11:03	12/07/16 17:33	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	136		25 - 150				11/28/16 11:03	12/07/16 17:33	10
13C4 PFOS	90		25 - 150				11/28/16 11:03	12/07/16 17:33	10
18O2 PFHxS	106		25 - 150				11/28/16 11:03	12/07/16 17:33	10

Client Sample Results

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-19-SO-17-18

Lab Sample ID: 320-23783-6

Date Collected: 11/22/16 11:30

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 79.7

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.37	J	0.62	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Perfluorooctane Sulfonate (PFOS)	6.5	M	0.62	0.16	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.50	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 16:44	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	69		25 - 150				11/28/16 11:03	12/02/16 16:44	1
13C4 PFOS	64		25 - 150				11/28/16 11:03	12/02/16 16:44	1
18O2 PFHxS	73		25 - 150				11/28/16 11:03	12/02/16 16:44	1

Client Sample ID: DPT-16-32-SO-14-15

Lab Sample ID: 320-23783-7

Date Collected: 11/21/16 09:36

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 82.3

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.1		0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Perfluorooctane Sulfonate (PFOS)	46	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:52	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	96		25 - 150				11/28/16 11:03	12/02/16 16:52	1
13C4 PFOS	62		25 - 150				11/28/16 11:03	12/02/16 16:52	1
18O2 PFHxS	70		25 - 150				11/28/16 11:03	12/02/16 16:52	1

Client Sample ID: DPT-16-31-SO-14-15

Lab Sample ID: 320-23783-8

Date Collected: 11/21/16 10:12

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	J	0.60	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Perfluorooctane Sulfonate (PFOS)	11	M	0.60	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 16:59	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	108		25 - 150				11/28/16 11:03	12/02/16 16:59	1
13C4 PFOS	82		25 - 150				11/28/16 11:03	12/02/16 16:59	1
18O2 PFHxS	93		25 - 150				11/28/16 11:03	12/02/16 16:59	1

Client Sample ID: DPT-16-30-SO-14-15

Lab Sample ID: 320-23783-9

Date Collected: 11/21/16 10:46

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 83.8

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	5.5		0.59	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Perfluorooctane Sulfonate (PFOS)	57	M	0.59	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:07	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	101		25 - 150				11/28/16 11:03	12/02/16 17:07	1
13C4 PFOS	63		25 - 150				11/28/16 11:03	12/02/16 17:07	1
18O2 PFHxS	90		25 - 150				11/28/16 11:03	12/02/16 17:07	1

TestAmerica Sacramento

Client Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-28-SO-14-15

Lab Sample ID: 320-23783-10

Date Collected: 11/21/16 11:54

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 85.2

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.45	J	0.58	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Perfluorooctane Sulfonate (PFOS)	10	M	0.58	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Perfluorobutanesulfonic acid (PFBS)	0.35	U	0.47	0.12	ug/Kg	☼	11/28/16 11:03	12/02/16 17:14	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	105		25 - 150				11/28/16 11:03	12/02/16 17:14	1
13C4 PFOS	61		25 - 150				11/28/16 11:03	12/02/16 17:14	1
18O2 PFHxS	89		25 - 150				11/28/16 11:03	12/02/16 17:14	1

Client Sample ID: DPT-16-15-SO-14-15

Lab Sample ID: 320-23783-11

Date Collected: 11/21/16 12:50

Matrix: Solid

Date Received: 11/23/16 08:46

Percent Solids: 80.6

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.51	J	0.61	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Perfluorooctane Sulfonate (PFOS)	11	M	0.61	0.15	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.49	0.13	ug/Kg	☼	11/28/16 11:03	12/02/16 17:22	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	101		25 - 150				11/28/16 11:03	12/02/16 17:22	1
13C4 PFOS	66		25 - 150				11/28/16 11:03	12/02/16 17:22	1
18O2 PFHxS	89		25 - 150				11/28/16 11:03	12/02/16 17:22	1

Client Sample ID: DPT-16-15-GW-18-22

Lab Sample ID: 320-23783-12

Date Collected: 11/21/16 13:30

Matrix: Water

Date Received: 11/23/16 08:46

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	5.8	E M	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:41	1
Perfluorooctane Sulfonate (PFOS)	6.5	E M	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:41	1
Perfluorobutanesulfonic acid (PFBS)	1.2	E	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:41	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	23	Q	25 - 150				11/28/16 16:55	12/03/16 16:41	1
13C4 PFOS	29		25 - 150				11/28/16 16:55	12/03/16 16:41	1
18O2 PFHxS	21	Q	25 - 150				11/28/16 16:55	12/03/16 16:41	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	15	D	0.25	0.075	ug/L		11/28/16 16:55	12/07/16 17:03	100
Perfluorooctane Sulfonate (PFOS)	12	D	0.40	0.13	ug/L		11/28/16 16:55	12/07/16 17:03	100
Perfluorobutanesulfonic acid (PFBS)	0.54	D	0.25	0.092	ug/L		11/28/16 16:55	12/07/16 17:03	100
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	83		25 - 150				11/28/16 16:55	12/07/16 17:03	100
13C4 PFOS	94		25 - 150				11/28/16 16:55	12/07/16 17:03	100
18O2 PFHxS	104		25 - 150				11/28/16 16:55	12/07/16 17:03	100

Client Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-15-GW-31-35

Lab Sample ID: 320-23783-13

Date Collected: 11/21/16 14:00

Matrix: Water

Date Received: 11/23/16 08:46

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	M	0.0025	0.00076	ug/L		11/28/16 16:55	12/03/16 16:48	1
Perfluorooctane Sulfonate (PFOS)	0.46	E	0.0041	0.0013	ug/L		11/28/16 16:55	12/03/16 16:48	1
Perfluorobutanesulfonic acid (PFBS)	0.048		0.0025	0.00093	ug/L		11/28/16 16:55	12/03/16 16:48	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	22	Q	25 - 150				11/28/16 16:55	12/03/16 16:48	1
13C4 PFOS	82		25 - 150				11/28/16 16:55	12/03/16 16:48	1
18O2 PFHxS	66		25 - 150				11/28/16 16:55	12/03/16 16:48	1

Method: 537 (Modified) - Perfluorinated Hydrocarbons - DL

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	0.27	D	0.025	0.0076	ug/L		11/28/16 16:55	12/07/16 17:25	10
Perfluorooctane Sulfonate (PFOS)	0.45	D	0.041	0.013	ug/L		11/28/16 16:55	12/07/16 17:25	10
Perfluorobutanesulfonic acid (PFBS)	0.035	D	0.025	0.0093	ug/L		11/28/16 16:55	12/07/16 17:25	10
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOA	33		25 - 150				11/28/16 16:55	12/07/16 17:25	10
13C4 PFOS	95		25 - 150				11/28/16 16:55	12/07/16 17:25	10
18O2 PFHxS	101		25 - 150				11/28/16 16:55	12/07/16 17:25	10

Default Detection Limits

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Prep: 3535

Analyte	RL	MDL	Units	Method
Perfluorobutanesulfonic acid (PFBS)	0.0025	0.00092	ug/L	537 (Modified)
Perfluorooctane Sulfonate (PFOS)	0.0040	0.0013	ug/L	537 (Modified)
Perfluorooctanoic acid (PFOA)	0.0025	0.00075	ug/L	537 (Modified)

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Prep: SHAKE

Analyte	RL	MDL	Units	Method
Perfluorobutanesulfonic acid (PFBS)	0.40	0.10	ug/Kg	537 (Modified)
Perfluorooctane Sulfonate (PFOS)	0.50	0.13	ug/Kg	537 (Modified)
Perfluorooctanoic acid (PFOA)	0.50	0.10	ug/Kg	537 (Modified)

Isotope Dilution Summary

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Matrix: Solid

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Isotope Dilution Recovery (Acceptance Limits)		
		3C4 PFO/ (25-150)	3C4 PFO/ (25-150)	3O2 PFHx (25-150)
320-23783-1	DPT-16-35-SO-00-01	93	73	86
320-23783-2	DPT-16-35-SO-19-20	112	101	108
320-23783-5	DPT-16-19-SO-00-01	96	54	85
320-23783-5 - DL	DPT-16-19-SO-00-01	136	90	106
320-23783-6	DPT-16-19-SO-17-18	69	64	73
320-23783-7	DPT-16-32-SO-14-15	96	62	70
320-23783-8	DPT-16-31-SO-14-15	108	82	93
320-23783-9	DPT-16-30-SO-14-15	101	63	90
320-23783-10	DPT-16-28-SO-14-15	105	61	89
320-23783-11	DPT-16-15-SO-14-15	101	66	89
LCS 320-139627/2-A	Lab Control Sample	119	112	120
MB 320-139627/1-A	Method Blank	108	94	104

Surrogate Legend

- 13C4 PFOA = 13C4 PFOA
- 13C4 PFOS = 13C4 PFOS
- 18O2 PFHxS = 18O2 PFHxS

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Isotope Dilution Recovery (Acceptance Limits)		
		3C4 PFO/ (25-150)	3C4 PFO/ (25-150)	3O2 PFHx (25-150)
320-23783-3	DPT-16-35-GW-21-25	70	58	43
320-23783-3 - DL	DPT-16-35-GW-21-25	91	87	83
320-23783-4	DPT-16-35-GW-31-35	52	76	30
320-23783-4 - DL	DPT-16-35-GW-31-35	78	91	73
320-23783-12	DPT-16-15-GW-18-22	23 Q	29	21 Q
320-23783-12 - DL	DPT-16-15-GW-18-22	83	94	104
320-23783-13	DPT-16-15-GW-31-35	22 Q	82	66
320-23783-13 - DL	DPT-16-15-GW-31-35	33	95	101
LCS 320-139749/2-A	Lab Control Sample	108	104	105
LCSD 320-139749/3-A	Lab Control Sample Dup	116	104	104
MB 320-139749/1-A	Method Blank	93	85	86

Surrogate Legend

- 13C4 PFOA = 13C4 PFOA
- 13C4 PFOS = 13C4 PFOS
- 18O2 PFHxS = 18O2 PFHxS

QC Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method: 537 (Modified) - Perfluorinated Hydrocarbons

Lab Sample ID: MB 320-139627/1-A
Matrix: Solid
Analysis Batch: 140429

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 139627

Analyte	MB	MB	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier							
Perfluorooctanoic acid (PFOA)	0.30	U	0.50	0.10	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Perfluorooctane Sulfonate (PFOS)	0.30	U	0.50	0.13	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Perfluorobutanesulfonic acid (PFBS)	0.30	U	0.40	0.10	ug/Kg		11/28/16 11:03	12/02/16 14:59	1
Isotope Dilution	MB	MB	Limits				Prepared	Analyzed	Dil Fac
	%Recovery	Qualifier							
13C4 PFOA	108		25 - 150				11/28/16 11:03	12/02/16 14:59	1
13C4 PFOS	94		25 - 150				11/28/16 11:03	12/02/16 14:59	1
18O2 PFHxS	104		25 - 150				11/28/16 11:03	12/02/16 14:59	1

Lab Sample ID: LCS 320-139627/2-A
Matrix: Solid
Analysis Batch: 140429

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 139627

Analyte	Spike Added	LCS	LCS	Unit	D	%Rec	Limits
		Result	Qualifier				
Perfluorooctanoic acid (PFOA)	4.00	4.25		ug/Kg		106	60 - 140
Perfluorooctane Sulfonate (PFOS)	3.71	3.66		ug/Kg		99	60 - 140
Perfluorobutanesulfonic acid (PFBS)	3.54	4.03		ug/Kg		114	50 - 150
Isotope Dilution	LCS	LCS	Limits				
	%Recovery	Qualifier					
13C4 PFOA	119		25 - 150				
13C4 PFOS	112		25 - 150				
18O2 PFHxS	120		25 - 150				

Lab Sample ID: MB 320-139749/1-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Method Blank
Prep Type: Total/NA
Prep Batch: 139749

Analyte	MB	MB	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier							
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		11/28/16 16:55	12/03/16 16:03	1
Perfluorooctane Sulfonate (PFOS)	0.0030	U	0.0040	0.0013	ug/L		11/28/16 16:55	12/03/16 16:03	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		11/28/16 16:55	12/03/16 16:03	1
Isotope Dilution	MB	MB	Limits				Prepared	Analyzed	Dil Fac
	%Recovery	Qualifier							
13C4 PFOA	93		25 - 150				11/28/16 16:55	12/03/16 16:03	1
13C4 PFOS	85		25 - 150				11/28/16 16:55	12/03/16 16:03	1
18O2 PFHxS	86		25 - 150				11/28/16 16:55	12/03/16 16:03	1

Lab Sample ID: LCS 320-139749/2-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Lab Control Sample
Prep Type: Total/NA
Prep Batch: 139749

Analyte	Spike Added	LCS	LCS	Unit	D	%Rec	Limits
		Result	Qualifier				
Perfluorooctanoic acid (PFOA)	0.0400	0.0296		ug/L		74	60 - 140
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0261		ug/L		70	60 - 140
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0289		ug/L		82	50 - 150

TestAmerica Sacramento

QC Sample Results

Client: AECOM Technical Services Inc.
 Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

<i>Isotope Dilution</i>	<i>LCS LCS</i>		<i>Limits</i>
	<i>%Recovery</i>	<i>Qualifier</i>	
13C4 PFOA	108		25 - 150
13C4 PFOS	104		25 - 150
18O2 PFHxS	105		25 - 150

Lab Sample ID: LCSD 320-139749/3-A
Matrix: Water
Analysis Batch: 140564

Client Sample ID: Lab Control Sample Dup
Prep Type: Total/NA
Prep Batch: 139749

<i>Analyte</i>	<i>Spike Added</i>	<i>LCSD Result</i>	<i>LCSD Qualifier</i>	<i>Unit</i>	<i>D</i>	<i>%Rec</i>	<i>Limits</i>	<i>RPD</i>	<i>RPD</i>
									<i>Limit</i>
Perfluorooctanoic acid (PFOA)	0.0400	0.0314		ug/L		79	60 - 140	6	30
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0289		ug/L		78	60 - 140	10	30
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0325		ug/L		92	50 - 150	12	30

<i>Isotope Dilution</i>	<i>LCSD LCSD</i>		<i>Limits</i>
	<i>%Recovery</i>	<i>Qualifier</i>	
13C4 PFOA	116		25 - 150
13C4 PFOS	104		25 - 150
18O2 PFHxS	104		25 - 150

QC Association Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

LCMS

Prep Batch: 139627

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	SHAKE	
320-23783-5 - DL	DPT-16-19-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	SHAKE	
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	SHAKE	
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	SHAKE	
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	SHAKE	
MB 320-139627/1-A	Method Blank	Total/NA	Solid	SHAKE	
LCS 320-139627/2-A	Lab Control Sample	Total/NA	Solid	SHAKE	

Prep Batch: 139749

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3	DPT-16-35-GW-21-25	Total/NA	Water	3535	
320-23783-3 - DL	DPT-16-35-GW-21-25	Total/NA	Water	3535	
320-23783-4	DPT-16-35-GW-31-35	Total/NA	Water	3535	
320-23783-4 - DL	DPT-16-35-GW-31-35	Total/NA	Water	3535	
320-23783-12	DPT-16-15-GW-18-22	Total/NA	Water	3535	
320-23783-12 - DL	DPT-16-15-GW-18-22	Total/NA	Water	3535	
320-23783-13	DPT-16-15-GW-31-35	Total/NA	Water	3535	
320-23783-13 - DL	DPT-16-15-GW-31-35	Total/NA	Water	3535	
MB 320-139749/1-A	Method Blank	Total/NA	Water	3535	
LCS 320-139749/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-139749/3-A	Lab Control Sample Dup	Total/NA	Water	3535	

Analysis Batch: 140429

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	537 (Modified)	139627
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	537 (Modified)	139627
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	537 (Modified)	139627
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	537 (Modified)	139627
MB 320-139627/1-A	Method Blank	Total/NA	Solid	537 (Modified)	139627
LCS 320-139627/2-A	Lab Control Sample	Total/NA	Solid	537 (Modified)	139627

Analysis Batch: 140564

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3	DPT-16-35-GW-21-25	Total/NA	Water	537 (Modified)	139749
320-23783-4	DPT-16-35-GW-31-35	Total/NA	Water	537 (Modified)	139749
320-23783-12	DPT-16-15-GW-18-22	Total/NA	Water	537 (Modified)	139749
320-23783-13	DPT-16-15-GW-31-35	Total/NA	Water	537 (Modified)	139749
MB 320-139749/1-A	Method Blank	Total/NA	Water	537 (Modified)	139749
LCS 320-139749/2-A	Lab Control Sample	Total/NA	Water	537 (Modified)	139749
LCSD 320-139749/3-A	Lab Control Sample Dup	Total/NA	Water	537 (Modified)	139749

TestAmerica Sacramento

QC Association Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

LCMS (Continued)

Analysis Batch: 141062

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-3 - DL	DPT-16-35-GW-21-25	Total/NA	Water	537 (Modified)	139749
320-23783-4 - DL	DPT-16-35-GW-31-35	Total/NA	Water	537 (Modified)	139749
320-23783-5 - DL	DPT-16-19-SO-00-01	Total/NA	Solid	537 (Modified)	139627
320-23783-12 - DL	DPT-16-15-GW-18-22	Total/NA	Water	537 (Modified)	139749
320-23783-13 - DL	DPT-16-15-GW-31-35	Total/NA	Water	537 (Modified)	139749

General Chemistry

Analysis Batch: 140129

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-1	DPT-16-35-SO-00-01	Total/NA	Solid	D 2216	
320-23783-2	DPT-16-35-SO-19-20	Total/NA	Solid	D 2216	
320-23783-5	DPT-16-19-SO-00-01	Total/NA	Solid	D 2216	
320-23783-6	DPT-16-19-SO-17-18	Total/NA	Solid	D 2216	

Analysis Batch: 140181

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-23783-7	DPT-16-32-SO-14-15	Total/NA	Solid	D 2216	
320-23783-8	DPT-16-31-SO-14-15	Total/NA	Solid	D 2216	
320-23783-9	DPT-16-30-SO-14-15	Total/NA	Solid	D 2216	
320-23783-10	DPT-16-28-SO-14-15	Total/NA	Solid	D 2216	
320-23783-11	DPT-16-15-SO-14-15	Total/NA	Solid	D 2216	

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-35-SO-00-01

Date Collected: 11/22/16 09:24

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-1

Matrix: Solid

Percent Solids: 86.4

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 15:52	SBC	TAL SAC

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-35-SO-19-20

Date Collected: 11/22/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-2

Matrix: Solid

Percent Solids: 82.8

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 15:59	SBC	TAL SAC

Client Sample ID: DPT-16-35-GW-21-25

Date Collected: 11/22/16 10:10

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-3

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140564	12/03/16 16:26	SBC	TAL SAC
Total/NA	Prep	3535	DL		139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10	141062	12/07/16 17:10	CBW	TAL SAC

Client Sample ID: DPT-16-35-GW-31-35

Date Collected: 11/22/16 10:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-4

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140564	12/03/16 16:33	SBC	TAL SAC

TestAmerica Sacramento

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-35-GW-31-35

Date Collected: 11/22/16 10:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-4

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535	DL		139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10	141062	12/07/16 17:18	CBW	TAL SAC

Client Sample ID: DPT-16-19-SO-00-01

Date Collected: 11/22/16 11:20

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-5

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-19-SO-00-01

Date Collected: 11/22/16 11:20

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-5

Matrix: Solid

Percent Solids: 86.7

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 16:07	SBC	TAL SAC
Total/NA	Prep	SHAKE	DL		139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10	141062	12/07/16 17:33	CBW	TAL SAC

Client Sample ID: DPT-16-19-SO-17-18

Date Collected: 11/22/16 11:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-6

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140129	12/01/16 09:31	CRM	TAL SAC

Client Sample ID: DPT-16-19-SO-17-18

Date Collected: 11/22/16 11:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-6

Matrix: Solid

Percent Solids: 79.7

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 16:44	SBC	TAL SAC

Client Sample ID: DPT-16-32-SO-14-15

Date Collected: 11/21/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-7

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140181	12/01/16 11:23	KSL	TAL SAC

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-32-SO-14-15

Date Collected: 11/21/16 09:36

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-7

Matrix: Solid

Percent Solids: 82.3

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 16:52	SBC	TAL SAC

Client Sample ID: DPT-16-31-SO-14-15

Date Collected: 11/21/16 10:12

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-8

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-31-SO-14-15

Date Collected: 11/21/16 10:12

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-8

Matrix: Solid

Percent Solids: 83.8

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 16:59	SBC	TAL SAC

Client Sample ID: DPT-16-30-SO-14-15

Date Collected: 11/21/16 10:46

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-9

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-30-SO-14-15

Date Collected: 11/21/16 10:46

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-9

Matrix: Solid

Percent Solids: 83.8

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 17:07	SBC	TAL SAC

Client Sample ID: DPT-16-28-SO-14-15

Date Collected: 11/21/16 11:54

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-10

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140181	12/01/16 11:23	KSL	TAL SAC

Lab Chronicle

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Client Sample ID: DPT-16-28-SO-14-15

Date Collected: 11/21/16 11:54

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-10

Matrix: Solid

Percent Solids: 85.2

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 17:14	SBC	TAL SAC

Client Sample ID: DPT-16-15-SO-14-15

Date Collected: 11/21/16 12:50

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-11

Matrix: Solid

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Analysis	D 2216		1	140181	12/01/16 11:23	KSL	TAL SAC

Client Sample ID: DPT-16-15-SO-14-15

Date Collected: 11/21/16 12:50

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-11

Matrix: Solid

Percent Solids: 80.6

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	SHAKE			139627	11/28/16 11:03	HJA	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140429	12/02/16 17:22	SBC	TAL SAC

Client Sample ID: DPT-16-15-GW-18-22

Date Collected: 11/21/16 13:30

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-12

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140564	12/03/16 16:41	SBC	TAL SAC
Total/NA	Prep	3535	DL		139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	100	141062	12/07/16 17:03	CBW	TAL SAC

Client Sample ID: DPT-16-15-GW-31-35

Date Collected: 11/21/16 14:00

Date Received: 11/23/16 08:46

Lab Sample ID: 320-23783-13

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)		1	140564	12/03/16 16:48	SBC	TAL SAC
Total/NA	Prep	3535	DL		139749	11/28/16 16:55	VPM	TAL SAC
Total/NA	Analysis	537 (Modified)	DL	10	141062	12/07/16 17:25	CBW	TAL SAC

Laboratory References:

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600

Certification Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Laboratory: TestAmerica Sacramento

Unless otherwise noted, all analytes for this laboratory were covered under each certification below.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2928-01	01-31-17

The following analytes are included in this report, but certification is not offered by the governing authority:

Analysis Method	Prep Method	Matrix	Analyte
537 (Modified)	SHAKE	Solid	Perfluorooctane Sulfonate (PFOS)
D 2216		Solid	Percent Moisture
D 2216		Solid	Percent Solids

Method Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Method	Method Description	Protocol	Laboratory
537 (Modified)	Perfluorinated Hydrocarbons	EPA	TAL SAC
D 2216	Percent Moisture	ASTM	TAL SAC

Protocol References:

ASTM = ASTM International
EPA = US Environmental Protection Agency

Laboratory References:

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600

Sample Summary

Client: AECOM Technical Services Inc.
Project/Site: Former Bay Head Road Annex (RFP JU06-01)

TestAmerica Job ID: 320-23783-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-23783-1	DPT-16-35-SO-00-01	Solid	11/22/16 09:24	11/23/16 08:46
320-23783-2	DPT-16-35-SO-19-20	Solid	11/22/16 09:36	11/23/16 08:46
320-23783-3	DPT-16-35-GW-21-25	Water	11/22/16 10:10	11/23/16 08:46
320-23783-4	DPT-16-35-GW-31-35	Water	11/22/16 10:30	11/23/16 08:46
320-23783-5	DPT-16-19-SO-00-01	Solid	11/22/16 11:20	11/23/16 08:46
320-23783-6	DPT-16-19-SO-17-18	Solid	11/22/16 11:30	11/23/16 08:46
320-23783-7	DPT-16-32-SO-14-15	Solid	11/21/16 09:36	11/23/16 08:46
320-23783-8	DPT-16-31-SO-14-15	Solid	11/21/16 10:12	11/23/16 08:46
320-23783-9	DPT-16-30-SO-14-15	Solid	11/21/16 10:46	11/23/16 08:46
320-23783-10	DPT-16-28-SO-14-15	Solid	11/21/16 11:54	11/23/16 08:46
320-23783-11	DPT-16-15-SO-14-15	Solid	11/21/16 12:50	11/23/16 08:46
320-23783-12	DPT-16-15-GW-18-22	Water	11/21/16 13:30	11/23/16 08:46
320-23783-13	DPT-16-15-GW-31-35	Water	11/21/16 14:00	11/23/16 08:46

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140382

Lab Sample ID: IC 320-140382/4 Client Sample ID: _____

Date Analyzed: 12/02/16 10:29 Lab File ID: 02DEC2016A_004.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorotetradecanoic acid (PFTeA)	4.94	Incomplete Integration	chandrase nas	12/02/16 13:37

Lab Sample ID: IC 320-140382/5 Client Sample ID: _____

Date Analyzed: 12/02/16 10:37 Lab File ID: 02DEC2016A_005.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.38	Isomers	chandrase nas	12/02/16 13:38

Lab Sample ID: IC 320-140382/10 Client Sample ID: _____

Date Analyzed: 12/02/16 11:14 Lab File ID: 02DEC2016A_010.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorotridecanoic Acid (PFTriA)	4.61	Incomplete Integration	chandrase nas	12/02/16 13:42

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140429

Lab Sample ID: 320-23783-1 Client Sample ID: DPT-16-35-SO-00-01

Date Analyzed: 12/02/16 15:52 Lab File ID: 02DEC2016C_008.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.89	Baseline	chandrase nas	12/04/16 18:22

Lab Sample ID: 320-23783-2 Client Sample ID: DPT-16-35-SO-19-20

Date Analyzed: 12/02/16 15:59 Lab File ID: 02DEC2016C_009.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.27	Baseline	chandrase nas	12/04/16 18:23

Lab Sample ID: 320-23783-5 Client Sample ID: DPT-16-19-SO-00-01

Date Analyzed: 12/02/16 16:07 Lab File ID: 02DEC2016C_010.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.27	Baseline	chandrase nas	12/04/16 18:25

Lab Sample ID: 320-23783-6 Client Sample ID: DPT-16-19-SO-17-18

Date Analyzed: 12/02/16 16:44 Lab File ID: 02DEC2016C_015.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.26	Baseline	chandrase nas	12/04/16 18:27

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140429

Lab Sample ID: 320-23783-7 Client Sample ID: DPT-16-32-SO-14-15

Date Analyzed: 12/02/16 16:52 Lab File ID: 02DEC2016C_016.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.25	Baseline	chandrase nas	12/04/16 18:28

Lab Sample ID: 320-23783-8 Client Sample ID: DPT-16-31-SO-14-15

Date Analyzed: 12/02/16 16:59 Lab File ID: 02DEC2016C_017.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.26	Baseline	chandrase nas	12/04/16 18:29

Lab Sample ID: 320-23783-9 Client Sample ID: DPT-16-30-SO-14-15

Date Analyzed: 12/02/16 17:07 Lab File ID: 02DEC2016C_018.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.26	Baseline	chandrase nas	12/04/16 18:29

Lab Sample ID: 320-23783-10 Client Sample ID: DPT-16-28-SO-14-15

Date Analyzed: 12/02/16 17:14 Lab File ID: 02DEC2016C_019.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.26	Baseline	chandrase nas	12/04/16 18:30

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140429

Lab Sample ID: 320-23783-11 Client Sample ID: DPT-16-15-SO-14-15

Date Analyzed: 12/02/16 17:22 Lab File ID: 02DEC2016C_020.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctane Sulfonate (PFOS)	3.26	Baseline	chandrase nas	12/04/16 18:31

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140564

Lab Sample ID: IC 320-140564/4 Client Sample ID: _____

Date Analyzed: 12/03/16 13:48 Lab File ID: 03DEC2016A_004.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorononanoic acid (PFNA)	3.28	Incomplete Integration	chandrase nas	12/05/16 09:42
Perfluoroundecanoic acid (PFUnA)	3.97	Incomplete Integration	chandrase nas	12/05/16 09:42

Lab Sample ID: IC 320-140564/7 Client Sample ID: _____

Date Analyzed: 12/03/16 14:11 Lab File ID: 03DEC2016A_007.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorohexanesulfonic acid (PFHxS)	2.53	Baseline	chandrase nas	12/05/16 09:41
Perfluorooctanoic acid (PFOA)	2.89	Baseline	chandrase nas	12/05/16 09:41

Lab Sample ID: IC 320-140564/9 Client Sample ID: _____

Date Analyzed: 12/03/16 14:26 Lab File ID: 03DEC2016A_009.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorotridecanoic Acid (PFTriA)	4.50	Incomplete Integration	chandrase nas	12/05/16 09:44

Lab Sample ID: 320-23783-3 Client Sample ID: DPT-16-35-GW-21-25

Date Analyzed: 12/03/16 16:26 Lab File ID: 03DEC2016B_004.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.85	Baseline	chandrase nas	12/05/16 10:31
Perfluorooctane Sulfonate (PFOS)	3.23	Baseline	chandrase nas	12/05/16 10:31

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 140564

Lab Sample ID: 320-23783-4 Client Sample ID: DPT-16-35-GW-31-35

Date Analyzed: 12/03/16 16:33 Lab File ID: 03DEC2016B_005.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.85	Baseline	chandrase nas	12/05/16 10:34

Lab Sample ID: 320-23783-12 Client Sample ID: DPT-16-15-GW-18-22

Date Analyzed: 12/03/16 16:41 Lab File ID: 03DEC2016B_006.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.82	Baseline	chandrase nas	12/05/16 10:35
Perfluorooctane Sulfonate (PFOS)	3.19	Baseline	chandrase nas	12/05/16 10:35

Lab Sample ID: 320-23783-13 Client Sample ID: DPT-16-15-GW-31-35

Date Analyzed: 12/03/16 16:48 Lab File ID: 03DEC2016B_007.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.86	Baseline	chandrase nas	12/05/16 10:36

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Analysis Batch Number: 141062

Lab Sample ID: 320-23783-5 DL Client Sample ID: DPT-16-19-SO-00-01 DL

Date Analyzed: 12/07/16 17:33 Lab File ID: 07DEC2016B_009.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorooctanoic acid (PFOA)	2.80	Baseline	westendor fc	12/08/16 08:55

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
LCMPFCSU_00046	03/01/17	11/03/16	Methanol, Lot Baker 144541	50000 uL	LCM2PFHxDA_00008	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00007	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00007	1000 uL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00008	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00011	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00008	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00011	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00008	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00012	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00008	1000 uL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00008	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00012	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00017	1000 uL	13C4 PFOS	0.956 ug/mL
LCMPFUDa_00009	1000 uL	13C2 PFUnA	1 ug/mL					
.LCM2PFHxDA_00008	01/07/21	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)	13C2-PFHxDA	50 ug/mL	
.LCM2PFTeDA_00007	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115			(Purchased Reagent)	13C2-PFTeDA	50 ug/mL	
.LCM4PFHPA_00007	05/27/21	Wellington Laboratories, Lot M4PFHpa0516			(Purchased Reagent)	13C4-PFHpa	50 ug/mL	
.LCM5PFPEA_00008	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)	13C5-PFPeA	50 ug/mL	
.LCM8FOSA_00011	12/22/17	Wellington Laboratories, Lot M8FOSA1215I			(Purchased Reagent)	13C8 FOSA	50 ug/mL	
.LCMPFBA_00008	05/24/21	Wellington Laboratories, Lot MPFBA0516			(Purchased Reagent)	13C4 PFBA	50 ug/mL	
.LCMPFDA_00011	08/19/20	Wellington Laboratories, Lot MPFDA0815			(Purchased Reagent)	13C2 PFDA	50 ug/mL	
.LCMPFDoA_00008	04/08/21	Wellington Laboratories, Lot MPFDoA0416			(Purchased Reagent)	13C2 PFDoA	50 ug/mL	
.LCMPFHxA_00012	04/08/21	Wellington Laboratories, Lot MPFHxA0416			(Purchased Reagent)	13C2 PFHxA	50 ug/mL	
.LCMPFHxS_00008	10/23/20	Wellington Laboratories, Lot MPFHxS1015			(Purchased Reagent)	1802 PFHxS	47.3 ug/mL	
.LCMPFNA_00008	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)	13C5 PFNA	50 ug/mL	
.LCMPFOA_00012	01/22/21	Wellington Laboratories, Lot MPFOA0116			(Purchased Reagent)	13C4 PFOA	50 ug/mL	
.LCMPFOS_00017	08/03/21	Wellington Laboratories, Lot MPFOS0816			(Purchased Reagent)	13C4 PFOS	47.8 ug/mL	
.LCMPFUDa_00009	02/12/21	Wellington Laboratories, Lot MPFUDa0216			(Purchased Reagent)	13C2 PFUnA	50 ug/mL	
LCPPFC-L0_00002	03/01/17	10/14/16	MeOH/H2O, Lot 144541	10 mL	LCMPFCSU_00045	500 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpa	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							1802 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
13C2 PFUnA	50 ng/mL							
.LCMPFCSU_00045	03/01/17	08/31/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00006	1000 uL	13C4-PFHpa	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCM5PFPEA 00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA 00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA 00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA 00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA 00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA 00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS 00007	1000 uL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA 00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA 00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS 00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa 00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA 00006	01/07/21	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA 00006	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA 00006	05/22/20	Wellington Laboratories, Lot M4PFHPA0515			(Purchased Reagent)		13C4-PFHpA	50 ug/mL
..LCM5PFPEA 00007	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA 00010	12/22/17	Wellington Laboratories, Lot M8FOSA1215I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA 00007	05/24/21	Wellington Laboratories, Lot MPFBA0516			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA 00010	08/19/20	Wellington Laboratories, Lot MPFDA0815			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA 00007	04/08/21	Wellington Laboratories, Lot MPFDoA0416			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA 00011	04/08/21	Wellington Laboratories, Lot MPFHxA0416			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS 00007	10/23/20	Wellington Laboratories, Lot MPFHxS1015			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA 00007	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA 00011	01/22/21	Wellington Laboratories, Lot MPFOA0116			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS 00015	01/22/21	Wellington Laboratories, Lot MPFOS0116			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa 00008	10/31/19	Wellington Laboratories, Lot MPFUDa1014			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
LCPFCL-L1_00021	12/28/16	08/03/16	MeOH/H2O, Lot 90285	5 mL	LCMPFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
					13C2 PFUnA	50 ng/mL		
					LCPFCLSP_00057	25 uL	Perfluorobutyric acid	0.5 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	0.442 ng/mL
							Perfluorodecanoic acid	0.5 ng/mL
							Perfluorododecanoic acid	0.5 ng/mL
							Perfluorodecane Sulfonic acid	0.482 ng/mL
							Perfluoroheptanoic acid	0.5 ng/mL
Perfluoroheptanesulfonic Acid	0.476 ng/mL							
Perfluorohexanoic acid	0.5 ng/mL							

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorohexadecanoic acid	0.5 ng/mL
							Perfluorohexanesulfonic acid	0.455 ng/mL
							Perfluorononanoic acid	0.5 ng/mL
							Perfluorooctanoic acid (PFOA)	0.5 ng/mL
							Perfluorooctadecanoic acid	0.5 ng/mL
							Perfluorooctane Sulfonate (PFOS)	0.464 ng/mL
							Perfluorooctane Sulfonamide	0.5 ng/mL
							Perfluoropentanoic acid	0.5 ng/mL
							Perfluorotetradecanoic acid	0.5 ng/mL
							Perfluorotridecanoic acid	0.5 ng/mL
							Perfluoroundecanoic acid	0.5 ng/mL
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHFA_00006	1000 uL	13C4-PFHFA	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00006	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHFA_00006	05/22/20	Wellington Laboratories, Lot M4PFHFA0515			(Purchased Reagent)		13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00007	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17	Wellington Laboratories, Lot M8FOSA1215I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21	Wellington Laboratories, Lot MPFBA0516			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20	Wellington Laboratories, Lot MPFDA0815			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21	Wellington Laboratories, Lot MPFDoA0416			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21	Wellington Laboratories, Lot MPFHxA0416			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20	Wellington Laboratories, Lot MPFHxS1015			(Purchased Reagent)		1802 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21	Wellington Laboratories, Lot MPFOA0116			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21	Wellington Laboratories, Lot MPFOS0116			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00008	10/31/19	Wellington Laboratories, Lot MPFUdA1014			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00057	02/01/17	08/03/16	Methanol, Lot 090285	10000 uL	LCPFCSP_00056	1000 uL	Perfluorobutyric acid	0.1 ug/mL
							Perfluorobutanesulfonic acid (PFBS)	0.0884 ug/mL
							Perfluorodecanoic acid	0.1 ug/mL
							Perfluorododecanoic acid	0.1 ug/mL
							Perfluorodecane Sulfonic acid	0.0964 ug/mL
							Perfluoroheptanoic acid	0.1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration	
					Reagent ID	Volume Added			
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL	
							Perfluorohexanoic acid	0.1 ug/mL	
							Perfluorohexadecanoic acid	0.1 ug/mL	
							Perfluorohexanesulfonic acid	0.091 ug/mL	
							Perfluorononanoic acid	0.1 ug/mL	
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL	
							Perfluorooctadecanoic acid	0.1 ug/mL	
							Perfluorooctane Sulfonate (PFOS)	0.0928 ug/mL	
							Perfluorooctane Sulfonamide	0.1 ug/mL	
							Perfluoropentanoic acid	0.1 ug/mL	
							Perfluorotetradecanoic acid	0.1 ug/mL	
							Perfluorotridecanoic acid	0.1 ug/mL	
							Perfluoroundecanoic acid	0.1 ug/mL	
..LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBFA_00004	200 uL	Perfluorobutyric acid	1 ug/mL	
					LCPFBFS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL	
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL	
					LCPFDaA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL	
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL	
					LCPFHpA_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL	
					LCPFHpS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL	
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL	
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL	
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL	
					LCPFNA_00005	200 uL	Perfluorononanoic acid	1 ug/mL	
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL	
					LCPFODA_00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL	
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL	
					LCPFOSA_00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL	
					LCPFPeA_00005	200 uL	Perfluoropentanoic acid	1 ug/mL	
					LCPFTeDA_00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL	
					LCPFTrDA_00004	200 uL	Perfluorotridecanoic acid	1 ug/mL	
					LCPFUdA_00004	200 uL	Perfluoroundecanoic acid	1 ug/mL	
...LCPFBFA_00004	01/30/20		Wellington Laboratories, Lot PFBA0115				(Purchased Reagent)	Perfluorobutyric acid	50 ug/mL
...LCPFBFS_00004	10/09/19		Wellington Laboratories, Lot LFPBFS1014				(Purchased Reagent)	Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDA_00005	07/02/20		Wellington Laboratories, Lot PFDA0615				(Purchased Reagent)	Perfluorodecanoic acid	50 ug/mL
...LCPFDaA_00005	01/30/20		Wellington Laboratories, Lot PFDoA0115				(Purchased Reagent)	Perfluorododecanoic acid	50 ug/mL
...LCPFDS_00005	07/02/20		Wellington Laboratories, Lot LFPDS0615				(Purchased Reagent)	Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00005	01/22/21		Wellington Laboratories, Lot PFHpA0116				(Purchased Reagent)	Perfluoroheptanoic acid	50 ug/mL
...LCPFHpS_00008	11/06/20		Wellington Laboratories, Lot LFPHpS1115				(Purchased Reagent)	Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA_00004	12/22/20		Wellington Laboratories, Lot PFHxA1215				(Purchased Reagent)	Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707				(Purchased Reagent)	Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxS-br_00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615				(Purchased Reagent)	Perfluorohexanesulfonic acid	45.5 ug/mL
...LCPFNA_00005	10/23/20		Wellington Laboratories, Lot PFNA1015				(Purchased Reagent)	Perfluorononanoic acid	50 ug/mL
...LCPFOA_00006	11/06/20		Wellington Laboratories, Lot PFOA1115				(Purchased Reagent)	Perfluorooctanoic acid (PFOA)	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
...LCPFODA 00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
...LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
...LCPFOSA 00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA 00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA 00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA 00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
...LCPFUDa 00004	08/19/20		Wellington Laboratories, Lot PFUDa0815		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L2_00022	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
							LCPFCSP_00057	50 uL
					Perfluorobutanesulfonic acid (PFBS)	0.884 ng/mL		
					Perfluorodecanoic acid	1 ng/mL		
					Perfluorododecanoic acid	1 ng/mL		
					Perfluorodecane Sulfonic acid	0.964 ng/mL		
					Perfluoroheptanoic acid	1 ng/mL		
					Perfluoroheptanesulfonic Acid	0.952 ng/mL		
					Perfluorohexanoic acid	1 ng/mL		
					Perfluorohexadecanoic acid	1 ng/mL		
					Perfluorohexanesulfonic acid	0.91 ng/mL		
					Perfluorononanoic acid	1 ng/mL		
Perfluorooctanoic acid (PFOA)	1 ng/mL							
Perfluorooctadecanoic acid	1 ng/mL							
Perfluorooctane Sulfonate (PFOS)	0.928 ng/mL							
Perfluorooctane Sulfonamide	1 ng/mL							
Perfluoropentanoic acid	1 ng/mL							
Perfluorotetradecanoic acid	1 ng/mL							
Perfluorotridecanoic acid	1 ng/mL							
Perfluoroundecanoic acid	1 ng/mL							
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00006	1000 uL	13C4-PFHpA	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTEda_00006	12/07/20		Wellington Laboratories, Lot M2PFTEda1115		(Purchased Reagent)		13C2-PFTEda	50 ug/mL
..LCM4PFHFA_00006	05/22/20		Wellington Laboratories, Lot M4PFHFA0515		(Purchased Reagent)		13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00007	05/22/20		Wellington Laboratories, Lot M5PFPEA0515		(Purchased Reagent)		13C5-PFPEA	50 ug/mL
..LCM8FOSA_00010	12/22/17		Wellington Laboratories, Lot M8FOSA1215I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21		Wellington Laboratories, Lot MPFBA0516		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20		Wellington Laboratories, Lot MPFDA0815		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21		Wellington Laboratories, Lot MPFDoA0416		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21		Wellington Laboratories, Lot MPFHxA0416		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20		Wellington Laboratories, Lot MPFHxS1015		(Purchased Reagent)		1802 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21		Wellington Laboratories, Lot MPFOA0116		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21		Wellington Laboratories, Lot MPFOS0116		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00008	10/31/19		Wellington Laboratories, Lot MPFUDa1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
..LCPFCSP_00057	02/01/17	08/03/16	Methanol, Lot 090285	10000 uL	LCPFCSP_00056	1000 uL	Perfluorobutyric acid	0.1 ug/mL
							Perfluorobutanesulfonic acid (PFBS)	0.0884 ug/mL
							Perfluorodecanoic acid	0.1 ug/mL
							Perfluorododecanoic acid	0.1 ug/mL
							Perfluorodecane Sulfonic acid	0.0964 ug/mL
							Perfluoroheptanoic acid	0.1 ug/mL
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL
							Perfluorohexanoic acid	0.1 ug/mL
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid	0.091 ug/mL
							Perfluorononanoic acid	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctane Sulfonate (PFOS)	0.0928 ug/mL
							Perfluorooctane Sulfonamide	0.1 ug/mL
							Perfluoropentanoic acid	0.1 ug/mL
							Perfluorotetradecanoic acid	0.1 ug/mL
							Perfluorotridecanoic acid	0.1 ug/mL
							Perfluoroundecanoic acid	0.1 ug/mL
..LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBA_00004	200 uL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA_00005	200 uL	Perfluorononanoic acid	1 ug/mL
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA_00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00004	200 uL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUDa_00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
...LCPFBA_00004	01/30/20		Wellington Laboratories, Lot PFBA0115		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFBS_00004	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDA_00005	07/02/20		Wellington Laboratories, Lot PFDA0615		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDoA_00005	01/30/20		Wellington Laboratories, Lot PFDoA0115		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFDS_00005	07/02/20		Wellington Laboratories, Lot LPFDS0615		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00005	01/22/21		Wellington Laboratories, Lot PFHpA0116		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
...LCPFHpS_00008	11/06/20		Wellington Laboratories, Lot LPFHpS1115		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA_00004	12/22/20		Wellington Laboratories, Lot PFHxA1215		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxS-br_00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615		(Purchased Reagent)		Perfluorohexanesulfonic acid	45.5 ug/mL
...LCPFNA_00005	10/23/20		Wellington Laboratories, Lot PFNA1015		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
...LCPFOA_00006	11/06/20		Wellington Laboratories, Lot PFOA1115		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFODA_00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
...LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
...LCPFOSA_00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA_00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA_00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA_00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
...LCPFUDa_00004	08/19/20		Wellington Laboratories, Lot PFUDa0815		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L3_00019	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFCSP_00057	250 uL	1802 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
							Perfluorobutyric acid	5 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	4.42 ng/mL
							Perfluorodecanoic acid	5 ng/mL
							Perfluorododecanoic acid	5 ng/mL
							Perfluorodecane Sulfonic acid	4.82 ng/mL
							Perfluoroheptanoic acid	5 ng/mL
							Perfluoroheptanesulfonic Acid	4.76 ng/mL
							Perfluorohexanoic acid	5 ng/mL
							Perfluorohexadecanoic acid	5 ng/mL
							Perfluorohexanesulfonic acid	4.55 ng/mL
							Perfluorononanoic acid	5 ng/mL
							Perfluorooctanoic acid (PFOA)	5 ng/mL
							Perfluorooctadecanoic acid	5 ng/mL
Perfluorooctane Sulfonate (PFOS)	4.64 ng/mL							
Perfluorooctane Sulfonamide	5 ng/mL							
Perfluoropentanoic acid	5 ng/mL							
Perfluorotetradecanoic acid	5 ng/mL							
Perfluorotridecanoic acid	5 ng/mL							
Perfluoroundecanoic acid	5 ng/mL							
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00006	1000 uL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00006	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00006	05/22/20	Wellington Laboratories, Lot M4PFHpA0515			(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA_00007	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17	Wellington Laboratories, Lot M8FOSA1215I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21	Wellington Laboratories, Lot MPFBA0516			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20	Wellington Laboratories, Lot MPFDA0815			(Purchased Reagent)		13C2 PFDA	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCMPFDoA_00007	04/08/21		Wellington Laboratories, Lot MPFDoA0416		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21		Wellington Laboratories, Lot MPFHxA0416		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20		Wellington Laboratories, Lot MPFHxS1015		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21		Wellington Laboratories, Lot MPFOA0116		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21		Wellington Laboratories, Lot MPFOS0116		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00008	10/31/19		Wellington Laboratories, Lot MPFUdA1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
..LCPFCSP_00057	02/01/17	08/03/16	Methanol, Lot 090285	10000 uL	LCPFCSP_00056	1000 uL	Perfluorobutyric acid	0.1 ug/mL
							Perfluorobutanesulfonic acid (PFBS)	0.0884 ug/mL
							Perfluorodecanoic acid	0.1 ug/mL
							Perfluorododecanoic acid	0.1 ug/mL
							Perfluorodecane Sulfonic acid	0.0964 ug/mL
							Perfluoroheptanoic acid	0.1 ug/mL
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL
							Perfluorohexanoic acid	0.1 ug/mL
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid	0.091 ug/mL
							Perfluorononanoic acid	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctane Sulfonate (PFOS)	0.0928 ug/mL
							Perfluorooctane Sulfonamide	0.1 ug/mL
							Perfluoropentanoic acid	0.1 ug/mL
							Perfluorotetradecanoic acid	0.1 ug/mL
							Perfluorotridecanoic acid	0.1 ug/mL
							Perfluoroundecanoic acid	0.1 ug/mL
..LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFDA_00004	200 uL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA_00005	200 uL	Perfluorononanoic acid	1 ug/mL
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA_00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTTrDA_00004	200 uL	Perfluorotridecanoic acid	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
...LCPFBA 00004	01/30/20		Wellington Laboratories, Lot PFBA0115		LCPFUdA 00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
...LCPFBS_00004	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFDA 00005	07/02/20		Wellington Laboratories, Lot PFDA0615		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDa 00005	01/30/20		Wellington Laboratories, Lot PFDoA0115		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDS 00005	07/02/20		Wellington Laboratories, Lot LPFDS0615		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFHpA 00005	01/22/21		Wellington Laboratories, Lot PFHpA0116		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHps 00008	11/06/20		Wellington Laboratories, Lot LPFHps1115		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
...LCPFHxA 00004	12/22/20		Wellington Laboratories, Lot PFHxA1215		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxS-br 00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFNA 00005	10/23/20		Wellington Laboratories, Lot PFNA1015		(Purchased Reagent)		Perfluorohexanesulfonic acid	45.5 ug/mL
...LCPFNA 00006	11/06/20		Wellington Laboratories, Lot PFNA1115		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
...LCPFODA 00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
...LCPFOSA 00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
...LCPFPeA 00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluorooctane Sulfonylamide	50 ug/mL
...LCPFTeDA 00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTrDA 00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFUdA 00004	08/19/20		Wellington Laboratories, Lot PFUdA0815		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
LCPFC-L4_00022	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCPFCFSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpa	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
							LCPFCSP_00056	100 uL
							Perfluorobutanesulfonic acid (PFBS)	17.68 ng/mL
							Perfluorodecanoic acid	20 ng/mL
							Perfluorododecanoic acid	20 ng/mL
							Perfluorodecane Sulfonic acid	19.28 ng/mL
							Perfluoroheptanoic acid	20 ng/mL
							Perfluoroheptanesulfonic Acid	19.04 ng/mL
							Perfluorohexanoic acid	20 ng/mL
							Perfluorohexadecanoic acid	20 ng/mL
		Perfluorohexanesulfonic acid	18.2 ng/mL					
		Perfluorononanoic acid	20 ng/mL					

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration	
					Reagent ID	Volume Added			
							Perfluorooctanoic acid (PFOA)	20 ng/mL	
							Perfluorooctadecanoic acid	20 ng/mL	
							Perfluorooctane Sulfonate (PFOS)	18.56 ng/mL	
							Perfluorooctane Sulfonamide	20 ng/mL	
							Perfluoropentanoic acid	20 ng/mL	
							Perfluorotetradecanoic acid	20 ng/mL	
							Perfluorotridecanoic acid	20 ng/mL	
							Perfluoroundecanoic acid	20 ng/mL	
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL	
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL	
					LCM4PFHFA_00006	1000 uL	13C4-PFHFA	1 ug/mL	
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL	
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL	
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL	
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL	
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL	
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL	
					LCMPFHxS_00007	1000 uL	18O2 PFHxS	0.946 ug/mL	
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL	
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL	
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL	
					LCMPFUDa_00008	1000 uL	13C2 PFUnA	1 ug/mL	
..LCM2PFHxDA_00006	01/07/21		Wellington Laboratories, Lot M2PFHxDA1112				(Purchased Reagent)	13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00006	12/07/20		Wellington Laboratories, Lot M2PFTeDA1115				(Purchased Reagent)	13C2-PFTeDA	50 ug/mL
..LCM4PFHFA_00006	05/22/20		Wellington Laboratories, Lot M4PFHFA0515				(Purchased Reagent)	13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00007	05/22/20		Wellington Laboratories, Lot M5PFPeA0515				(Purchased Reagent)	13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17		Wellington Laboratories, Lot M8FOSA1215I				(Purchased Reagent)	13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21		Wellington Laboratories, Lot MPFBA0516				(Purchased Reagent)	13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20		Wellington Laboratories, Lot MPFDA0815				(Purchased Reagent)	13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21		Wellington Laboratories, Lot MPFDoA0416				(Purchased Reagent)	13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21		Wellington Laboratories, Lot MPFHxA0416				(Purchased Reagent)	13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20		Wellington Laboratories, Lot MPFHxS1015				(Purchased Reagent)	18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19		Wellington Laboratories, Lot MPFNA0414				(Purchased Reagent)	13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21		Wellington Laboratories, Lot MPFOA0116				(Purchased Reagent)	13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21		Wellington Laboratories, Lot MPFOS0116				(Purchased Reagent)	13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00008	10/31/19		Wellington Laboratories, Lot MPFUDa1014				(Purchased Reagent)	13C2 PFUnA	50 ug/mL
.LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBA_00004	200 uL	Perfluorobutyric acid	1 ug/mL	
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL	
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL	
					LCPFDoA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL	
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL	
					LCPFHFA_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL	
					LCPFHFS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL	
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL	
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL	

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA 00005	200 uL	Perfluorononanoic acid	1 ug/mL
					LCPFOA 00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA 00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA 00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA 00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA 00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA 00004	200 uL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA 00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA 00004	01/30/20	Wellington Laboratories, Lot PFBA0115			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBS_00004	10/09/19	Wellington Laboratories, Lot LPFBS1014			(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA 00005	07/02/20	Wellington Laboratories, Lot PFDA0615			(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA 00005	01/30/20	Wellington Laboratories, Lot PFDoA0115			(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDS 00005	07/02/20	Wellington Laboratories, Lot LPFDS0615			(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA 00005	01/22/21	Wellington Laboratories, Lot PFHpA0116			(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS 00008	11/06/20	Wellington Laboratories, Lot LPFHps1115			(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA 00004	12/22/20	Wellington Laboratories, Lot PFHxA1215			(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA 00004	11/28/17	Wellington Laboratories, Lot PFHxDA0707			(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS-br_00001	07/03/20	Wellington Laboratories, Lot brPFHxSK0615			(Purchased Reagent)		Perfluorohexanesulfonic acid	45.5 ug/mL
..LCPFNA 00005	10/23/20	Wellington Laboratories, Lot PFNA1015			(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFOA 00006	11/06/20	Wellington Laboratories, Lot PFOA1115			(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA 00005	01/30/20	Wellington Laboratories, Lot PFODA0115			(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS-br_00001	10/14/20	Wellington Laboratories, Lot brPFOSK1015			(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
..LCPFOSA 00006	09/02/17	Wellington Laboratories, Lot FOSA0815I			(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA 00005	01/30/20	Wellington Laboratories, Lot PFPeA0115			(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA 00004	12/09/20	Wellington Laboratories, Lot PFTeDA1215			(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA 00004	12/10/18	Wellington Laboratories, Lot PFTrDA1213			(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA 00004	08/19/20	Wellington Laboratories, Lot PFUdA0815			(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L4_00023	02/01/17	09/30/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00045	250 uL	13C2 PFHxA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
					LCPFCSP_00056	100 uL	Perfluorobutanesulfonic acid (PFBS)	17.68 ng/mL
							Perfluorooctanoic acid (PFOA)	20 ng/mL
							Perfluorooctane Sulfonate (PFOS)	18.56 ng/mL
..LCMPFCSU_00045	03/01/17	08/31/16	Methanol, Lot Baker 115935	50000 uL	LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
..LCMPFHxA_00011	04/08/21	Wellington Laboratories, Lot MPFHxA0416			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFOA_00011	01/22/21	Wellington Laboratories, Lot MPFOA0116			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21	Wellington Laboratories, Lot MPFOS0116			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
..LCPFBS_00004	10/09/19	Wellington Laboratories, Lot LPFBS1014			(Purchased Reagent)	Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL	
..LCPFOA_00006	11/06/20	Wellington Laboratories, Lot PFOA1115			(Purchased Reagent)	Perfluorooctanoic acid (PFOA)	50 ug/mL	
..LCPFOS-br_00001	10/14/20	Wellington Laboratories, Lot brPFOSK1015			(Purchased Reagent)	Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL	
LCPFC-L5_00020	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDaA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
							LCPFCSP_00056	250 uL
					Perfluorobutanesulfonic acid (PFBS)	44.2 ng/mL		
					Perfluorodecanoic acid	50 ng/mL		
					Perfluorododecanoic acid	50 ng/mL		
					Perfluorodecane Sulfonic acid	48.2 ng/mL		
					Perfluoroheptanoic acid	50 ng/mL		
					Perfluoroheptanesulfonic Acid	47.6 ng/mL		
					Perfluorohexanoic acid	50 ng/mL		
					Perfluorohexadecanoic acid	50 ng/mL		
					Perfluorohexanesulfonic acid	45.5 ng/mL		
					Perfluorononanoic acid	50 ng/mL		
					Perfluorooctanoic acid (PFOA)	50 ng/mL		
					Perfluorooctadecanoic acid	50 ng/mL		
Perfluorooctane Sulfonate (PFOS)	46.4 ng/mL							
Perfluorooctane Sulfonamide	50 ng/mL							
Perfluoropentanoic acid	50 ng/mL							
Perfluorotetradecanoic acid	50 ng/mL							
Perfluorotridecanoic acid	50 ng/mL							
Perfluoroundecanoic acid	50 ng/mL							
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCM4PFHPA_00006	1000 uL	13C4-PFHpA	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTEdA_00006	12/07/20		Wellington Laboratories, Lot M2PFTEdA1115		(Purchased Reagent)		13C2-PFTEdA	50 ug/mL
..LCM4PFHPA_00006	05/22/20		Wellington Laboratories, Lot M4PFHPA0515		(Purchased Reagent)		13C4-PFHpA	50 ug/mL
..LCM5PFPEA_00007	05/22/20		Wellington Laboratories, Lot M5PFPeA0515		(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17		Wellington Laboratories, Lot M8FOSA1215I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21		Wellington Laboratories, Lot MPFBA0516		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20		Wellington Laboratories, Lot MPFDA0815		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21		Wellington Laboratories, Lot MPFDoA0416		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21		Wellington Laboratories, Lot MPFHxA0416		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20		Wellington Laboratories, Lot MPFHxS1015		(Purchased Reagent)		1802 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21		Wellington Laboratories, Lot MPFOA0116		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21		Wellington Laboratories, Lot MPFOS0116		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00008	10/31/19		Wellington Laboratories, Lot MPFUdA1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
..LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBA_00004	200 uL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA_00005	200 uL	Perfluorononanoic acid	1 ug/mL
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA_00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00004	200 uL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA_00004	01/30/20		Wellington Laboratories, Lot PFBA0115		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFBS_00004	10/09/19		Wellington Laboratories, Lot LPFBS1014			(Purchased Reagent)	Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA 00005	07/02/20		Wellington Laboratories, Lot PFDA0615			(Purchased Reagent)	Perfluorodecanoic acid	50 ug/mL
..LCPFDaA 00005	01/30/20		Wellington Laboratories, Lot PFDoA0115			(Purchased Reagent)	Perfluorododecanoic acid	50 ug/mL
..LCPFDS 00005	07/02/20		Wellington Laboratories, Lot LPFDS0615			(Purchased Reagent)	Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA 00005	01/22/21		Wellington Laboratories, Lot PFHpA0116			(Purchased Reagent)	Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS 00008	11/06/20		Wellington Laboratories, Lot LPFHpS1115			(Purchased Reagent)	Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA 00004	12/22/20		Wellington Laboratories, Lot PFHxA1215			(Purchased Reagent)	Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707			(Purchased Reagent)	Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS-br 00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615			(Purchased Reagent)	Perfluorohexanesulfonic acid	45.5 ug/mL
..LCPFNA 00005	10/23/20		Wellington Laboratories, Lot PFNA1015			(Purchased Reagent)	Perfluorononanoic acid	50 ug/mL
..LCPFOA 00006	11/06/20		Wellington Laboratories, Lot PFOA1115			(Purchased Reagent)	Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA 00005	01/30/20		Wellington Laboratories, Lot PFODA0115			(Purchased Reagent)	Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015			(Purchased Reagent)	Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
..LCPFOSA 00006	09/02/17		Wellington Laboratories, Lot FOSA0815I			(Purchased Reagent)	Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA 00005	01/30/20		Wellington Laboratories, Lot PFPeA0115			(Purchased Reagent)	Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA 00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215			(Purchased Reagent)	Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA 00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213			(Purchased Reagent)	Perfluorotridecanoic acid	50 ug/mL
..LCPFUDA 00004	08/19/20		Wellington Laboratories, Lot PFUDA0815			(Purchased Reagent)	Perfluoroundecanoic acid	50 ug/mL
LCPFC-L6_00019	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
					13C2 PFUnA	50 ng/mL		
					LCPFCSP_00056	1000 uL	Perfluorobutyric acid	200 ng/mL
					Perfluorobutanesulfonic acid (PFBS)		176.8 ng/mL	
					Perfluorodecanoic acid		200 ng/mL	
					Perfluorododecanoic acid		200 ng/mL	
					Perfluorodecane Sulfonic acid		192.8 ng/mL	
					Perfluoroheptanoic acid		200 ng/mL	
					Perfluoroheptanesulfonic Acid		190.4 ng/mL	
					Perfluorohexanoic acid		200 ng/mL	
					Perfluorohexadecanoic acid		200 ng/mL	
Perfluorohexanesulfonic acid	182 ng/mL							
Perfluorononanoic acid	200 ng/mL							
Perfluorooctanoic acid (PFOA)	200 ng/mL							
Perfluorooctadecanoic acid	200 ng/mL							

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorooctane Sulfonate (PFOS)	185.6 ng/mL
							Perfluorooctane Sulfonamide	200 ng/mL
							Perfluoropentanoic acid	200 ng/mL
							Perfluorotetradecanoic acid	200 ng/mL
							Perfluorotridecanoic acid	200 ng/mL
							Perfluoroundecanoic acid	200 ng/mL
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00006	1000 uL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00006	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00006	05/22/20	Wellington Laboratories, Lot M4PFHPA0515			(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA_00007	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17	Wellington Laboratories, Lot M8FOSA1215I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21	Wellington Laboratories, Lot MPFBA0516			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20	Wellington Laboratories, Lot MPFDA0815			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21	Wellington Laboratories, Lot MPFDoA0416			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21	Wellington Laboratories, Lot MPFHxA0416			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20	Wellington Laboratories, Lot MPFHxS1015			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21	Wellington Laboratories, Lot MPFOA0116			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21	Wellington Laboratories, Lot MPFOS0116			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00008	10/31/19	Wellington Laboratories, Lot MPFUDa1014			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBA_00004	200 uL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00005	200 uL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00005	200 uL	Perfluorododecanoic acid	1 ug/mL
					LCPFDS_00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpa_00005	200 uL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00004	200 uL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS-br_00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA_00005	200 uL	Perfluorononanoic acid	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFOA_00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA_00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPPeA_00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00004	200 uL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUDa_00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA_00004	01/30/20		Wellington Laboratories, Lot PFBA0115		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBS_00004	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA_00005	07/02/20		Wellington Laboratories, Lot PFDA0615		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA_00005	01/30/20		Wellington Laboratories, Lot PFDoA0115		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDS_00005	07/02/20		Wellington Laboratories, Lot LPFDS0615		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA_00005	01/22/21		Wellington Laboratories, Lot PFHpA0116		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHps_00008	11/06/20		Wellington Laboratories, Lot LPFHps1115		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA_00004	12/22/20		Wellington Laboratories, Lot PFHxA1215		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS-br_00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615		(Purchased Reagent)		Perfluorohexanesulfonic acid	45.5 ug/mL
..LCPFNA_00005	10/23/20		Wellington Laboratories, Lot PFNA1015		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFOA_00006	11/06/20		Wellington Laboratories, Lot PFOA1115		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
..LCPFOSA_00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA_00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA_00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUDa_00004	08/19/20		Wellington Laboratories, Lot PFUDa0815		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L7_00019	12/28/16	08/03/16	MeOH/H2O, Lot 090285	5 mL	LCPMFCSU_00044	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFTeDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL
							18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
					LCPFCSP_00056	2000 uL	Perfluorobutyric acid	400 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	353.6 ng/mL
							Perfluorodecanoic acid	400 ng/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorododecanoic acid	400 ng/mL
							Perfluorodecane Sulfonic acid	385.6 ng/mL
							Perfluoroheptanoic acid	400 ng/mL
							Perfluoroheptanesulfonic Acid	380.8 ng/mL
							Perfluorohexanoic acid	400 ng/mL
							Perfluorohexadecanoic acid	400 ng/mL
							Perfluorohexanesulfonic acid	364 ng/mL
							Perfluorononanoic acid	400 ng/mL
							Perfluorooctanoic acid (PFOA)	400 ng/mL
							Perfluorooctadecanoic acid	400 ng/mL
							Perfluorooctane Sulfonate (PFOS)	371.2 ng/mL
							Perfluorooctane Sulfonamide	400 ng/mL
							Perfluoropentanoic acid	400 ng/mL
							Perfluorotetradecanoic acid	400 ng/mL
							Perfluorotridecanoic acid	400 ng/mL
							Perfluoroundecanoic acid	400 ng/mL
.LCMPFCSU_00044	12/28/16	06/28/16	Methanol, Lot Baker 115935	50000 uL	LCM2PFHxDA_00006	1000 uL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00006	1000 uL	13C2-PFTeDA	1 ug/mL
					LCM4PFHFA_00006	1000 uL	13C4-PFHFA	1 ug/mL
					LCM5PFPEA_00007	1000 uL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00010	1000 uL	13C8 FOSA	1 ug/mL
					LCMPFBA_00007	1000 uL	13C4 PFBA	1 ug/mL
					LCMPFDA_00010	1000 uL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00007	1000 uL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00011	1000 uL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00007	1000 uL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00007	1000 uL	13C5 PFNA	1 ug/mL
					LCMPFOA_00011	1000 uL	13C4 PFOA	1 ug/mL
					LCMPFOS_00015	1000 uL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00008	1000 uL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00006	01/07/21		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00006	12/07/20		Wellington Laboratories, Lot M2PFTeDA1115		(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHFA_00006	05/22/20		Wellington Laboratories, Lot M4PFHFA0515		(Purchased Reagent)		13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00007	05/22/20		Wellington Laboratories, Lot M5PFPeA0515		(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00010	12/22/17		Wellington Laboratories, Lot M8FOSA1215I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00007	05/24/21		Wellington Laboratories, Lot MPFBA0516		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00010	08/19/20		Wellington Laboratories, Lot MPFDA0815		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00007	04/08/21		Wellington Laboratories, Lot MPFDoA0416		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00011	04/08/21		Wellington Laboratories, Lot MPFHxA0416		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00007	10/23/20		Wellington Laboratories, Lot MPFHxS1015		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00007	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00011	01/22/21		Wellington Laboratories, Lot MPFOA0116		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00015	01/22/21		Wellington Laboratories, Lot MPFOS0116		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00008	10/31/19		Wellington Laboratories, Lot MPFUDa1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00056	02/01/17	08/01/16	Methanol, Lot 090285	10000 uL	LCPFBA_00004	200 uL	Perfluorobutyric acid	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFBS_00004	200 uL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA 00005	200 uL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA 00005	200 uL	Perfluorododecanoic acid	1 ug/mL
					LCPFDS 00005	200 uL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA 00005	200 uL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS 00008	200 uL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA 00004	200 uL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA 00004	200 uL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS-br 00001	200 uL	Perfluorohexanesulfonic acid	0.91 ug/mL
					LCPFNA 00005	200 uL	Perfluorononanoic acid	1 ug/mL
					LCPFoA 00006	200 uL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA 00005	200 uL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS-br_00001	200 uL	Perfluorooctane Sulfonate (PFOS)	0.928 ug/mL
					LCPFOSA 00006	200 uL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA 00005	200 uL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA 00004	200 uL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA 00004	200 uL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUDA 00004	200 uL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA 00004	01/30/20		Wellington Laboratories, Lot PFBA0115		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBS_00004	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA 00005	07/02/20		Wellington Laboratories, Lot PFDA0615		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA 00005	01/30/20		Wellington Laboratories, Lot PFDoA0115		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDS 00005	07/02/20		Wellington Laboratories, Lot LPFDS0615		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA 00005	01/22/21		Wellington Laboratories, Lot PFHpA0116		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS 00008	11/06/20		Wellington Laboratories, Lot LPFHpS1115		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA 00004	12/22/20		Wellington Laboratories, Lot PFHxA1215		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS-br 00001	07/03/20		Wellington Laboratories, Lot brPFHxSK0615		(Purchased Reagent)		Perfluorohexanesulfonic acid	45.5 ug/mL
..LCPFNA 00005	10/23/20		Wellington Laboratories, Lot PFNA1015		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFoA 00006	11/06/20		Wellington Laboratories, Lot PFOA1115		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA 00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS-br_00001	10/14/20		Wellington Laboratories, Lot brPFOSK1015		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
..LCPFOSA 00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA 00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA 00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA 00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUDA 00004	08/19/20		Wellington Laboratories, Lot PFUDA0815		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC2-L1_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
							M2-8:2FTS	47.9 ng/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFC2SP_00014	25 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.474 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.479 ng/mL
							N-ethylperfluoro-1-octanesulfo namide	0.5 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	0.5 ng/mL
							MeFOSA	0.5 ng/mL
							N-methyl perfluorooctane sulfonamidoacetic acid	0.5 ng/mL
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NEtFOSA-M_00001	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M_00001	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA_00001	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NEtFOSAA_00001	200 uL	d5-NEtFOSAA	1 ug/mL
					LCM2-6:FtS_00001	200 uL	M2-6:2FtS	0.95 ug/mL
					LCM2-8:2FtS_00001	200 uL	M2-8:2FtS	0.958 ug/mL
						(Purchased Reagent)	d-N-EtFOSA-M	50 ug/mL
						(Purchased Reagent)	d-N-MeFOSA-M	50 ug/mL
						(Purchased Reagent)	d3-NMeFOSAA	50 ug/mL
						(Purchased Reagent)	d5-NEtFOSAA	50 ug/mL
						(Purchased Reagent)	M2-6:2FtS	47.5 ug/mL
						(Purchased Reagent)	M2-8:2FtS	47.9 ug/mL
.LCPFC2SP_00014	01/20/17	07/20/16	Methanol, Lot 104453	5000 uL	LCPFC2SP_00013	500 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.0948 ug/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.0958 ug/mL
							N-ethylperfluoro-1-octanesulfo namide	0.1 ug/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
							MeFOSA	0.1 ug/mL
							N-methyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
.LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL
					LC8:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfo namide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M_00001	200 uL	MeFOSA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
...LC6:2FTS_00001	10/03/17		WELLINGTON, Lot 62FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
...LC8:2FTS_00001	10/03/17		WELLINGTON, Lot 82FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
...LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NETFOSA0714M		(Purchased Reagent)		N-ethylperfluoro-1-octanesulfoamide	50 ug/mL
...LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NETFOSAA0113		(Purchased Reagent)		N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
...LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M		(Purchased Reagent)		MeFOSA	50 ug/mL
...LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214		(Purchased Reagent)		N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFC2-L2_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
					M2-8:2FTS	47.9 ng/mL		
					LCPFC2SP_00014	50 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ng/mL
							N-ethylperfluoro-1-octanesulfoamide	1 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	1 ng/mL
MeFOSA	1 ng/mL							
N-methyl perfluorooctane sulfonamidoacetic acid	1 ng/mL							
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NETFOSA-M_00001	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M_00001	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA_00001	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NETFOSAA_00001	200 uL	d5-NETFOSAA	1 ug/mL
					LCM2-6:FOSAA_00001	200 uL	M2-6:2FTS	0.95 ug/mL
					LCM2-8:2FOSAA_00001	200 uL	M2-8:2FTS	0.958 ug/mL
..LCd-NETFOSA-M_00001	03/10/19		WELLINGTON, Lot dNETFOSA0314M		(Purchased Reagent)		d-N-EtFOSA-M	50 ug/mL
..LCd-NMeFOSA-M_00001	01/28/19		WELLINGTON, Lot dNMeFOSA0114M		(Purchased Reagent)		d-N-MeFOSA-M	50 ug/mL
..LCd3-NMeFOSAA_00001	01/31/18		WELLINGTON, Lot d3NMeFOSAA0113		(Purchased Reagent)		d3-NMeFOSAA	50 ug/mL
..LCd5-NETFOSAA_00001	05/08/20		WELLINGTON, Lot d5NETFOSAA0515		(Purchased Reagent)		d5-NETFOSAA	50 ug/mL
..LCM2-6:FOSAA_00001	07/15/17		WELLINGTON, Lot M262FOSAA0714		(Purchased Reagent)		M2-6:2FTS	47.5 ug/mL
..LCM2-8:2FOSAA_00001	04/13/17		WELLINGTON, Lot M282FOSAA0414		(Purchased Reagent)		M2-8:2FTS	47.9 ug/mL
.LCPFC2SP_00014	01/20/17	07/20/16	Methanol, Lot 104453	5000 uL	LCPFC2SP_00013	500 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.0948 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.0958 ug/mL
							N-ethylperfluoro-1-octanesulfonamide	0.1 ug/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
							MeFOSA	0.1 ug/mL
							N-methyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
..LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FTS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL
					LC8:2FTS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfonamide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M_00001	200 uL	MeFOSA	1 ug/mL
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
...LC6:2FTS_00001	10/03/17		WELLINGTON, Lot 62FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
...LC8:2FTS_00001	10/03/17		WELLINGTON, Lot 82FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
...LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NETFOSA0714M		(Purchased Reagent)		N-ethylperfluoro-1-octanesulfonamide	50 ug/mL
...LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NETFOSAA0113		(Purchased Reagent)		N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
...LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M		(Purchased Reagent)		MeFOSA	50 ug/mL
...LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214		(Purchased Reagent)		N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFC2-L3_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
							M2-8:2FTS	47.9 ng/mL
					LCPFC2SP_00014	250 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	4.74 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	4.79 ng/mL
							N-ethylperfluoro-1-octanesulfonamide	5 ng/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							N-ethyl perfluorooctane sulfonamidoacetic acid	5 ng/mL
							MeFOSA	5 ng/mL
							N-methyl perfluorooctane sulfonamidoacetic acid	5 ng/mL
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NETfOSA-M 00001	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M 00001	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA 00001	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NETfOSAA 00001	200 uL	d5-NETfOSAA	1 ug/mL
					LCM2-6:FtS 00001	200 uL	M2-6:2FtS	0.95 ug/mL
					LCM2-8:2FtS 00001	200 uL	M2-8:2FtS	0.958 ug/mL
..LCd-NETfOSA-M 00001	03/10/19		WELLINGTON, Lot dNetFOSA0314M		(Purchased Reagent)		d-N-EtFOSA-M	50 ug/mL
..LCd-NMeFOSA-M 00001	01/28/19		WELLINGTON, Lot dNMeFOSA0114M		(Purchased Reagent)		d-N-MeFOSA-M	50 ug/mL
..LCd3-NMeFOSAA 00001	01/31/18		WELLINGTON, Lot d3NMeFOSAA0113		(Purchased Reagent)		d3-NMeFOSAA	50 ug/mL
..LCd5-NETfOSAA 00001	05/08/20		WELLINGTON, Lot d5NETfOSAA0515		(Purchased Reagent)		d5-NETfOSAA	50 ug/mL
..LCM2-6:FtS 00001	07/15/17		WELLINGTON, Lot M262FtS0714		(Purchased Reagent)		M2-6:2FtS	47.5 ug/mL
..LCM2-8:2FtS 00001	04/13/17		WELLINGTON, Lot M282FtS0414		(Purchased Reagent)		M2-8:2FtS	47.9 ug/mL
.LCPFC2SP_00014	01/20/17	07/20/16	Methanol, Lot 104453	5000 uL	LCPFC2SP_00013	500 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.0948 ug/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.0958 ug/mL
							N-ethylperfluoro-1-octanesulfoamide	0.1 ug/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
							MeFOSA	0.1 ug/mL
							N-methyl perfluorooctane sulfonamidoacetic acid	0.1 ug/mL
..LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL
					LC8:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfoamide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M 00001	200 uL	MeFOSA	1 ug/mL
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
...LC6:2FtS_00001	10/03/17		WELLINGTON, Lot 62FtS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
...LC8:2FtS_00001	10/03/17		WELLINGTON, Lot 82FtS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
...LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NETFOSA0714M		(Purchased Reagent)		N-ethylperfluoro-1-octanesulfo namide	50 ug/mL
...LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NETFOSAA0113		(Purchased Reagent)		N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
...LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M		(Purchased Reagent)		MeFOSA	50 ug/mL
...LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214		(Purchased Reagent)		N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFC2-L4_00003	02/26/17	09/22/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00008	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
					LCPFC2SP_00017	200 uL	M2-8:2FTS	47.9 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	18.96 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	19.16 ng/mL
							N-ethylperfluoro-1-octanesulfo namide	20 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	20 ng/mL
.LCMPFC2SU_00008	02/26/17	08/26/16	Methanol, Lot 104453	10000 uL	LCd-NEtFOSA-M_00002	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M_00002	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA_00002	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NETFOSAA_00002	200 uL	d5-NETFOSAA	1 ug/mL
					LCM2-6:Fts_00002	200 uL	M2-6:2FTS	0.95 ug/mL
					LCM2-8:2Fts_00002	200 uL	M2-8:2FTS	0.958 ug/mL
					..LCd-NEtFOSA-M_00002	03/10/19		WELLINGTON, Lot dNEtFOSA0314M
..LCd-NMeFOSA-M_00002	06/10/21		WELLINGTON, Lot dNMeFOSA0616M		(Purchased Reagent)	d-N-MeFOSA-M	50 ug/mL	
..LCd3-NMeFOSAA_00002	01/20/21		WELLINGTON, Lot d3NMeFOSAA0116		(Purchased Reagent)	d3-NMeFOSAA	50 ug/mL	
..LCd5-NETFOSAA_00002	12/07/20		WELLINGTON, Lot d5NETFOSAA1115		(Purchased Reagent)	d5-NETFOSAA	50 ug/mL	
..LCM2-6:Fts_00002	01/08/21		WELLINGTON, Lot M262Fts0116		(Purchased Reagent)	M2-6:2Fts	47.5 ug/mL	
..LCM2-8:2Fts_00002	01/08/21		WELLINGTON, Lot M282Fts0116		(Purchased Reagent)	M2-8:2Fts	47.9 ug/mL	
.LCPFC2SP_00017	03/02/17	09/02/16	Methanol, Lot 104453	10000 uL	LC6:2Fts_00002	100 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.474 ug/mL
					LC8:2Fts_00002	100 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.479 ug/mL
					LCN-EtFOSA-M_00003	100 uL	N-ethylperfluoro-1-octanesulfo namide	0.5 ug/mL
					LCN-EtFOSAA_00002	100 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	0.5 ug/mL
					LCN-MeFOSA-M_00002	100 uL	MeFOSA	0.5 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCN-MeFOSAA_00003	100 uL	N-methyl perfluorooctane sulfonamidoacetic acid	0.5 ug/mL
..LC6:2FTS_00002	06/25/21		WELLINGTON, Lot 62FTS0616		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
..LC8:2FTS_00002	10/23/20		WELLINGTON, Lot 82FTS1015		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
..LCN-EtFOSA-M_00003	05/24/21		WELLINGTON, Lot NETFOSA0516M		(Purchased Reagent)		N-ethylperfluoro-1-octanesulfo namide	50 ug/mL
..LCN-EtFOSAA_00002	01/20/21		WELLINGTON, Lot NETFOSAA0116		(Purchased Reagent)		N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
..LCN-MeFOSA-M_00002	05/24/21		WELLINGTON, Lot NMeFOSA0714M		(Purchased Reagent)		MeFOSA	50 ug/mL
..LCN-MeFOSAA_00003	01/20/21		WELLINGTON, Lot NMeFOSAA0116		(Purchased Reagent)		N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFC2-L5_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
					M2-8:2FTS	47.9 ng/mL		
					LCPFC2SP_00013	250 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ng/mL
							N-ethylperfluoro-1-octanesulfo namide	50 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	50 ng/mL
MeFOSA	50 ng/mL							
N-methyl perfluorooctane sulfonamidoacetic acid	50 ng/mL							
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NETFOSA-M_00001	200 uL	d-N-EtFOSA-M	1 ug/mL
							LCd-NMeFOSA-M_00001	1 ug/mL
							LCd3-NMeFOSAA_00001	1 ug/mL
							LCd5-NETFOSAA_00001	1 ug/mL
							LCM2-6:FOS_00001	0.95 ug/mL
							LCM2-8:2FOS_00001	0.958 ug/mL
..LCd-NETFOSA-M_00001	03/10/19		WELLINGTON, Lot dNETFOSA0314M		(Purchased Reagent)		d-N-EtFOSA-M	50 ug/mL
..LCd-NMeFOSA-M_00001	01/28/19		WELLINGTON, Lot dNMeFOSA0114M		(Purchased Reagent)		d-N-MeFOSA-M	50 ug/mL
..LCd3-NMeFOSAA_00001	01/31/18		WELLINGTON, Lot d3NMeFOSAA0113		(Purchased Reagent)		d3-NMeFOSAA	50 ug/mL
..LCd5-NETFOSAA_00001	05/08/20		WELLINGTON, Lot d5NETFOSAA0515		(Purchased Reagent)		d5-NETFOSAA	50 ug/mL
..LCM2-6:FOS_00001	07/15/17		WELLINGTON, Lot M262FOS0714		(Purchased Reagent)		M2-6:2FOS	47.5 ug/mL
..LCM2-8:2FOS_00001	04/13/17		WELLINGTON, Lot M282FOS0414		(Purchased Reagent)		M2-8:2FOS	47.9 ug/mL
.LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FOS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LC8:2FTS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfo namide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M_00001	200 uL	MeFOSA	1 ug/mL
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
..LC6:2FTS_00001	10/03/17		WELLINGTON, Lot 62FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
..LC8:2FTS_00001	10/03/17		WELLINGTON, Lot 82FTS1014		(Purchased Reagent)		Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
..LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NETFOSA0714M		(Purchased Reagent)		N-ethylperfluoro-1-octanesulfo namide	50 ug/mL
..LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NETFOSAA0113		(Purchased Reagent)		N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
..LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M		(Purchased Reagent)		MeFOSA	50 ug/mL
..LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214		(Purchased Reagent)		N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPF2-L6_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NETFOSAA	50 ng/mL
							M2-6:2FTS	47.5 ng/mL
							M2-8:2FTS	47.9 ng/mL
					LCPFC2SP_00013	1000 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	189.6 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	191.6 ng/mL
							N-ethylperfluoro-1-octanesulfo namide	200 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	200 ng/mL
						MeFOSA	200 ng/mL	
						N-methyl perfluorooctane sulfonamidoacetic acid	200 ng/mL	
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NETFOSA-M_00001	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M_00001	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA_00001	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NETFOSAA_00001	200 uL	d5-NETFOSAA	1 ug/mL
					LCM2-6:FTS_00001	200 uL	M2-6:2FTS	0.95 ug/mL
					LCM2-8:2FTS_00001	200 uL	M2-8:2FTS	0.958 ug/mL
..LCd-NETFOSA-M_00001	03/10/19		WELLINGTON, Lot dNETFOSA0314M		(Purchased Reagent)		d-N-EtFOSA-M	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCd-NMeFOSA-M_00001	01/28/19		WELLINGTON, Lot dNMeFOSA0114M			(Purchased Reagent)	d-N-MeFOSA-M	50 ug/mL
..LCd3-NMeFOSAA_00001	01/31/18		WELLINGTON, Lot d3NMeFOSAA0113			(Purchased Reagent)	d3-NMeFOSAA	50 ug/mL
..LCd5-NEtFOSAA_00001	05/08/20		WELLINGTON, Lot d5NEtFOSAA0515			(Purchased Reagent)	d5-NEtFOSAA	50 ug/mL
..LCM2-6:FtS_00001	07/15/17		WELLINGTON, Lot M262FtS0714			(Purchased Reagent)	M2-6:2FtS	47.5 ug/mL
..LCM2-8:2FtS_00001	04/13/17		WELLINGTON, Lot M282FtS0414			(Purchased Reagent)	M2-8:2FtS	47.9 ug/mL
.LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL
					LC8:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfo namide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M_00001	200 uL	MeFOSA	1 ug/mL
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
..LC6:2FtS_00001	10/03/17		WELLINGTON, Lot 62FtS1014			(Purchased Reagent)	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
..LC8:2FtS_00001	10/03/17		WELLINGTON, Lot 82FtS1014			(Purchased Reagent)	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
..LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NEtFOSA0714M			(Purchased Reagent)	N-ethylperfluoro-1-octanesulfo namide	50 ug/mL
..LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NEtFOSAA0113			(Purchased Reagent)	N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
..LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M			(Purchased Reagent)	MeFOSA	50 ug/mL
..LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214			(Purchased Reagent)	N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFC2-L7_00002	01/08/17	07/20/16	MeOH/H2O, Lot 104453	5 mL	LCMPFC2SU_00005	250 uL	d-N-EtFOSA-M	50 ng/mL
							d-N-MeFOSA-M	50 ng/mL
							d3-NMeFOSAA	50 ng/mL
							d5-NEtFOSAA	50 ng/mL
							M2-6:2FtS	47.5 ng/mL
							M2-8:2FtS	47.9 ng/mL
					LCPFC2SP_00013	2000 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	379.2 ng/mL
							Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	383.2 ng/mL
							N-ethylperfluoro-1-octanesulfo namide	400 ng/mL
							N-ethyl perfluorooctane sulfonamidoacetic acid	400 ng/mL
MeFOSA	400 ng/mL							

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							N-methyl perfluorooctane sulfonamidoacetic acid	400 ng/mL
.LCMPFC2SU_00005	01/08/17	07/08/16	Methanol, Lot 104453	10000 uL	LCd-NEtFOSA-M 00001	200 uL	d-N-EtFOSA-M	1 ug/mL
					LCd-NMeFOSA-M 00001	200 uL	d-N-MeFOSA-M	1 ug/mL
					LCd3-NMeFOSAA 00001	200 uL	d3-NMeFOSAA	1 ug/mL
					LCd5-NEtFOSAA 00001	200 uL	d5-NEtFOSAA	1 ug/mL
					LCM2-6:FtS 00001	200 uL	M2-6:2FtS	0.95 ug/mL
					LCM2-8:2FtS 00001	200 uL	M2-8:2FtS	0.958 ug/mL
..LCd-NEtFOSA-M 00001	03/10/19		WELLINGTON, Lot dNEtFOSA0314M			(Purchased Reagent)	d-N-EtFOSA-M	50 ug/mL
..LCd-NMeFOSA-M 00001	01/28/19		WELLINGTON, Lot dNMeFOSA0114M			(Purchased Reagent)	d-N-MeFOSA-M	50 ug/mL
..LCd3-NMeFOSAA 00001	01/31/18		WELLINGTON, Lot d3NMeFOSAA0113			(Purchased Reagent)	d3-NMeFOSAA	50 ug/mL
..LCd5-NEtFOSAA 00001	05/08/20		WELLINGTON, Lot d5NEtFOSAA0515			(Purchased Reagent)	d5-NEtFOSAA	50 ug/mL
..LCM2-6:FtS 00001	07/15/17		WELLINGTON, Lot M262FtS0714			(Purchased Reagent)	M2-6:2FtS	47.5 ug/mL
..LCM2-8:2FtS 00001	04/13/17		WELLINGTON, Lot M282FtS0414			(Purchased Reagent)	M2-8:2FtS	47.9 ug/mL
.LCPFC2SP_00013	01/20/17	07/20/16	Methanol, Lot 104453	10000 uL	LC6:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	0.948 ug/mL
					LC8:2FtS_00001	200 uL	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	0.958 ug/mL
					LCN-EtFOSA-M_00002	200 uL	N-ethylperfluoro-1-octanesulfonamide	1 ug/mL
					LCN-EtFOSAA_00001	200 uL	N-ethyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
					LCN-MeFOSA-M_00001	200 uL	MeFOSA	1 ug/mL
					LCN-MeFOSAA_00001	200 uL	N-methyl perfluorooctane sulfonamidoacetic acid	1 ug/mL
..LC6:2FtS_00001	10/03/17		WELLINGTON, Lot 62FtS1014			(Purchased Reagent)	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2)	47.4 ug/mL
..LC8:2FtS_00001	10/03/17		WELLINGTON, Lot 82FtS1014			(Purchased Reagent)	Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (8:2)	47.9 ug/mL
..LCN-EtFOSA-M_00002	07/14/19		WELLINGTON, Lot NEtFOSA0714M			(Purchased Reagent)	N-ethylperfluoro-1-octanesulfonamide	50 ug/mL
..LCN-EtFOSAA_00001	01/29/18		WELLINGTON, Lot NEtFOSAA0113			(Purchased Reagent)	N-ethyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
..LCN-MeFOSA-M_00001	07/15/19		WELLINGTON, Lot NMeFOSA0714M			(Purchased Reagent)	MeFOSA	50 ug/mL
..LCN-MeFOSAA_00001	12/09/19		WELLINGTON, Lot NMeFOSAA1214			(Purchased Reagent)	N-methyl perfluorooctane sulfonamidoacetic acid	50 ug/mL
LCPFCIC_00020	03/01/17	12/01/16	MeOH/H2O, Lot 09285	5 mL	LCMPFCSU_00046	250 uL	13C2-PFHxDA	50 ng/mL
							13C2-PFtEDA	50 ng/mL
							13C4-PFHpA	50 ng/mL
							13C5-PFPeA	50 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	50 ng/mL
							13C2 PFDA	50 ng/mL
							13C2 PFDoA	50 ng/mL
							13C2 PFHxA	50 ng/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFACMXB_00007	125 uL	18O2 PFHxS	47.3 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	50 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	44.25 ng/mL
							Perfluorooctane Sulfonate (PFOS)	47.75 ng/mL
.LCMPFCSU_00046	03/01/17	11/03/16	Methanol, Lot Baker 144541	50000 uL	LCM2PFHxDA_00008	1000 uL	13C2-PFHxDA	1 ug/mL
							LCM2PFTeDA_00007	1 ug/mL
							LCM4PFHPA_00007	1 ug/mL
							LCM5PFPEA_00008	1 ug/mL
							LCM8FOSA_00011	1 ug/mL
							LCMPFBA_00008	1 ug/mL
							LCMPFDA_00011	1 ug/mL
							LCMPFDoA_00008	1 ug/mL
							LCMPFHxA_00012	1 ug/mL
							LCMPFHxS_00008	0.946 ug/mL
							LCMPFNA_00008	1 ug/mL
							LCMPFOA_00012	1 ug/mL
							LCMPFOS_00017	0.956 ug/mL
							LCMPFUDa_00009	1 ug/mL
							..LCM2PFHxDA_00008	01/07/21
..LCM2PFTeDA_00007	12/07/20	Wellington Laboratories, Lot M2PFTeDA1115	(Purchased Reagent)	13C2-PFTeDA	50 ug/mL			
..LCM4PFHPA_00007	05/27/21	Wellington Laboratories, Lot M4PFHPa0516	(Purchased Reagent)	13C4-PFHpa	50 ug/mL			
..LCM5PFPEA_00008	05/22/20	Wellington Laboratories, Lot M5PFPeA0515	(Purchased Reagent)	13C5-PFPeA	50 ug/mL			
..LCM8FOSA_00011	12/22/17	Wellington Laboratories, Lot M8FOSA1215I	(Purchased Reagent)	13C8 FOSA	50 ug/mL			
..LCMPFBA_00008	05/24/21	Wellington Laboratories, Lot MPFBA0516	(Purchased Reagent)	13C4 PFBA	50 ug/mL			
..LCMPFDA_00011	08/19/20	Wellington Laboratories, Lot MPFDA0815	(Purchased Reagent)	13C2 PFDA	50 ug/mL			
..LCMPFDoA_00008	04/08/21	Wellington Laboratories, Lot MPFDoA0416	(Purchased Reagent)	13C2 PFDoA	50 ug/mL			
..LCMPFHxA_00012	04/08/21	Wellington Laboratories, Lot MPFHxA0416	(Purchased Reagent)	13C2 PFHxA	50 ug/mL			
..LCMPFHxS_00008	10/23/20	Wellington Laboratories, Lot MPFHxS1015	(Purchased Reagent)	18O2 PFHxS	47.3 ug/mL			
..LCMPFNA_00008	04/13/19	Wellington Laboratories, Lot MPFNA0414	(Purchased Reagent)	13C5 PFNA	50 ug/mL			
..LCMPFOA_00012	01/22/21	Wellington Laboratories, Lot MPFOA0116	(Purchased Reagent)	13C4 PFOA	50 ug/mL			
..LCMPFOS_00017	08/03/21	Wellington Laboratories, Lot MPFOS0816	(Purchased Reagent)	13C4 PFOS	47.8 ug/mL			
..LCMPFUDa_00009	02/12/21	Wellington Laboratories, Lot MPFUDa0216	(Purchased Reagent)	13C2 PFUnA	50 ug/mL			
.LCPFACMXB_00007	11/06/20	Wellington Laboratories, Lot PFACMXB1115	(Purchased Reagent)	Perfluorobutanesulfonic acid (PFBS)	1.77 ug/mL			
					Perfluorooctane Sulfonate (PFOS)	1.91 ug/mL		
					Perfluorooctanoic acid (PFOA)	2 ug/mL		
LCPFCSP_00070	05/15/17	11/15/16	Methanol, Lot 090285	10000 uL	LCPFBA_00005	100 uL	Perfluorobutyric acid	0.5 ug/mL
					LCPFBS_00005	100 uL	Perfluorobutane Sulfonate	0.442 ug/mL
							Perfluorobutanesulfonic acid (PFBS)	0.442 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-23783-1

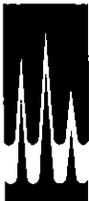
SDG No.: _____

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFDA_00005	100 uL	Perfluorodecanoic acid	0.5 ug/mL
					LCPFDoA_00005	100 uL	Perfluorododecanoic acid	0.5 ug/mL
					LCPFDS_00006	100 uL	Perfluorodecane Sulfonate	0.482 ug/mL
							Perfluorodecane Sulfonic acid	0.482 ug/mL
					LCPFHpA_00005	100 uL	Perfluoroheptanoic acid	0.5 ug/mL
					LCPFHpS_00009	100 uL	Perfluoroheptane Sulfonate	0.476 ug/mL
							Perfluoroheptanesulfonic Acid	0.476 ug/mL
					LCPFHxA_00004	100 uL	Perfluorohexanoic acid	0.5 ug/mL
					LCPFHxDA_00006	100 uL	Perfluorohexadecanoic acid	0.5 ug/mL
					LCPFHxS-br_00002	100 uL	Perfluorohexane Sulfonate	0.455 ug/mL
							Perfluorohexanesulfonic acid	0.455 ug/mL
					LCPFNA_00005	100 uL	Perfluorononanoic acid	0.5 ug/mL
					LCPFOA_00006	100 uL	Perfluorooctanoic acid (PFOA)	0.5 ug/mL
					LCPFODA_00005	100 uL	Perfluorooctadecanoic acid	0.5 ug/mL
					LCPFOS-br_00002	100 uL	Perfluorooctane Sulfonate (PFOS)	0.464 ug/mL
					LCPFOSA_00006	100 uL	Perfluorooctane Sulfonamide	0.5 ug/mL
					LCPFPeA_00005	100 uL	Perfluoropentanoic acid	0.5 ug/mL
					LCPFTeDA_00004	100 uL	Perfluorotetradecanoic acid	0.5 ug/mL
					LCPFTrDA_00004	100 uL	Perfluorotridecanoic acid	0.5 ug/mL
					LCPFUDa_00005	100 uL	Perfluoroundecanoic acid	0.5 ug/mL
.LCPFBA_00005	05/27/21		Wellington Laboratories, Lot PFBA0516		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
.LCPFBS_00005	03/15/21		Wellington Laboratories, Lot LPFBS0316		(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
							Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
.LCPFDA_00005	07/02/20		Wellington Laboratories, Lot PFDA0615		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
.LCPFDoA_00005	01/30/20		Wellington Laboratories, Lot PFDoA0115		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
.LCPFDS_00006	05/24/21		Wellington Laboratories, Lot LPFDS0516		(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
							Perfluorodecane Sulfonic acid	48.2 ug/mL
.LCPFHpA_00005	01/22/21		Wellington Laboratories, Lot PFHpA0116		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
.LCPFHpS_00009	11/06/20		Wellington Laboratories, Lot LPFHpS1115		(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
							Perfluoroheptanesulfonic Acid	47.6 ug/mL
.LCPFHxA_00004	12/22/20		Wellington Laboratories, Lot PFHxA1215		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
.LCPFHxDA_00006	05/25/21		Wellington Laboratories, Lot PFHxDA0516		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
.LCPFHxS-br_00002	07/03/20		Wellington Laboratories, Lot brPFHxSK0615		(Purchased Reagent)		Perfluorohexane Sulfonate	45.5 ug/mL
							Perfluorohexanesulfonic acid	45.5 ug/mL
.LCPFNA_00005	10/23/20		Wellington Laboratories, Lot PFNA1015		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
.LCPFOA_00006	11/06/20		Wellington Laboratories, Lot PFOA1115		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
.LCPFODA_00005	01/30/20		Wellington Laboratories, Lot PFODA0115		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
.LCPFOS-br_00002	10/14/20		Wellington Laboratories, Lot brPFOS1015		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	46.4 ug/mL
.LCPFOSA_00006	09/02/17		Wellington Laboratories, Lot FOSA0815I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
.LCPFPeA_00005	01/30/20		Wellington Laboratories, Lot PFPeA0115		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
.LCPFTeDA_00004	12/09/20		Wellington Laboratories, Lot PFTeDA1215		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
.LCPFTrDA_00004	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
.LCPFUDa_00005	08/19/20		Wellington Laboratories, Lot PFUDa0815		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL

Reagent

LC6:2FTS_00001

r: 7hclis ev
S: 7hclis sw

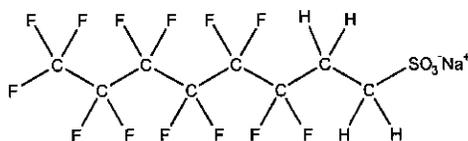


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: 6:2FTS **LOT NUMBER:** 62FTS1014
COMPOUND: Sodium 1H,1H,2H,2H-perfluorooctane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: $C_8H_4F_{13}SO_3Na$ **MOLECULAR WEIGHT:** 450.15
CONCENTRATION: $50.0 \pm 2.5 \mu\text{g/ml}$ (Na salt) **SOLVENT(S):** Methanol
 $47.4 \pm 2.4 \mu\text{g/ml}$ (6:2FTS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 10/03/2014
EXPIRY DATE: (mm/dd/yyyy) 10/03/2017
RECOMMENDED STORAGE: Refrigerate ampoule

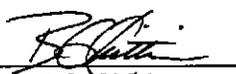
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 03/27/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

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EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

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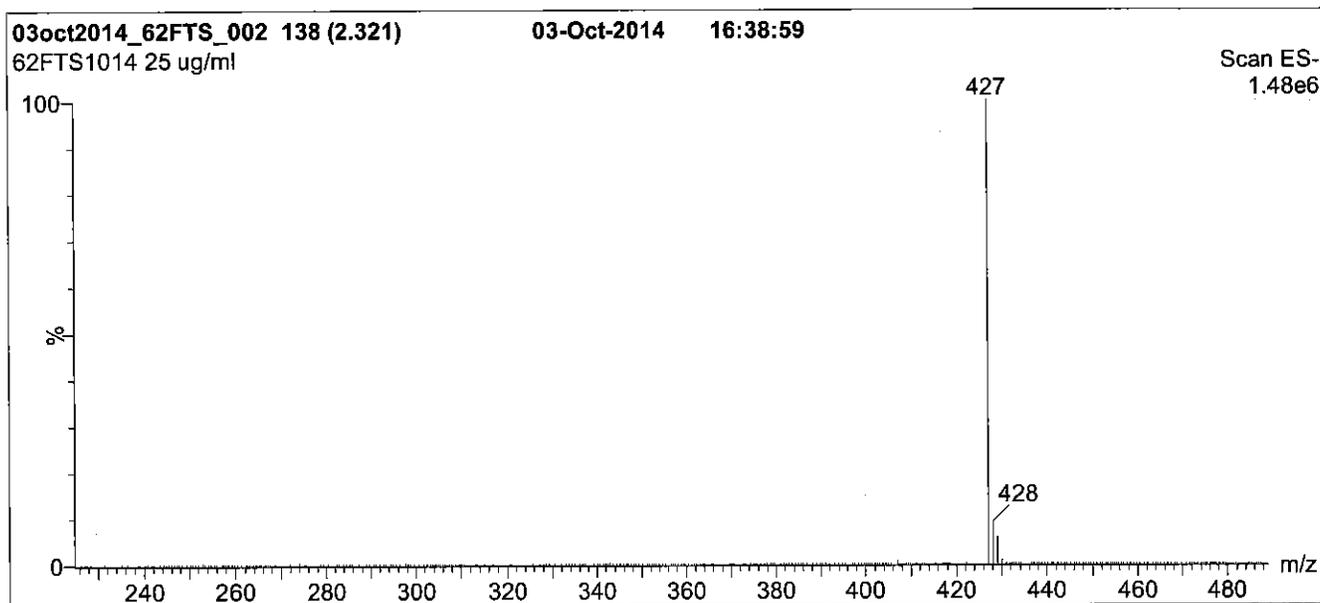
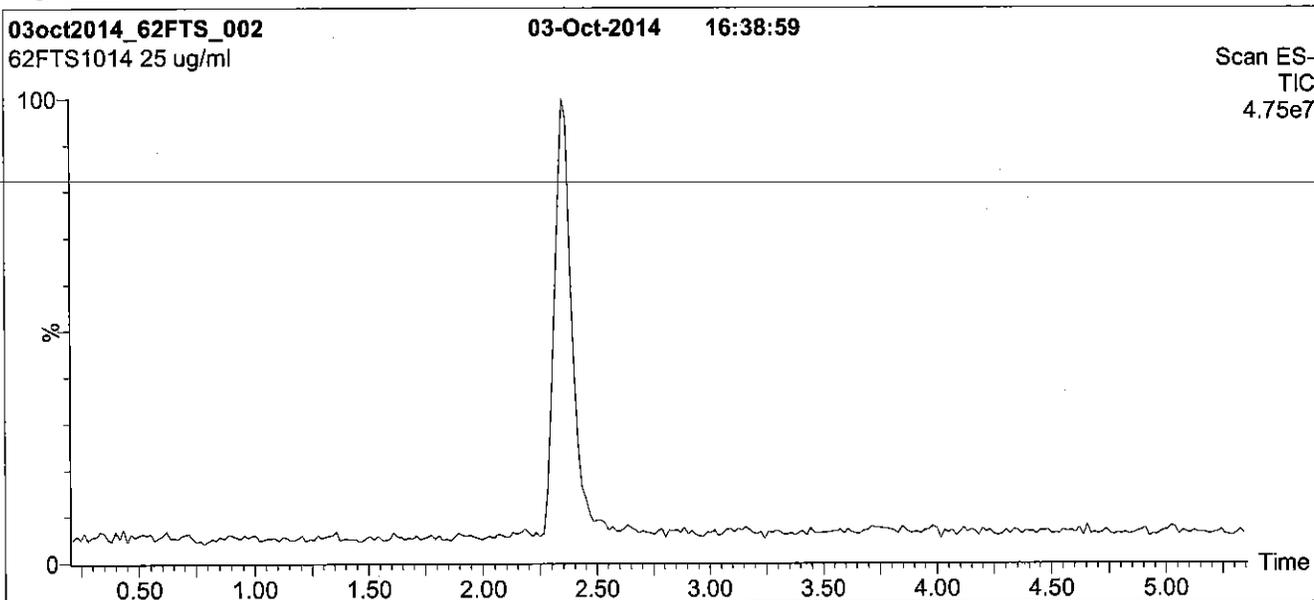
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: 6:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 55% (80:20 MeOH:ACN) / 45% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min
before returning to initial conditions in 0.5 min.
Time: 10 min

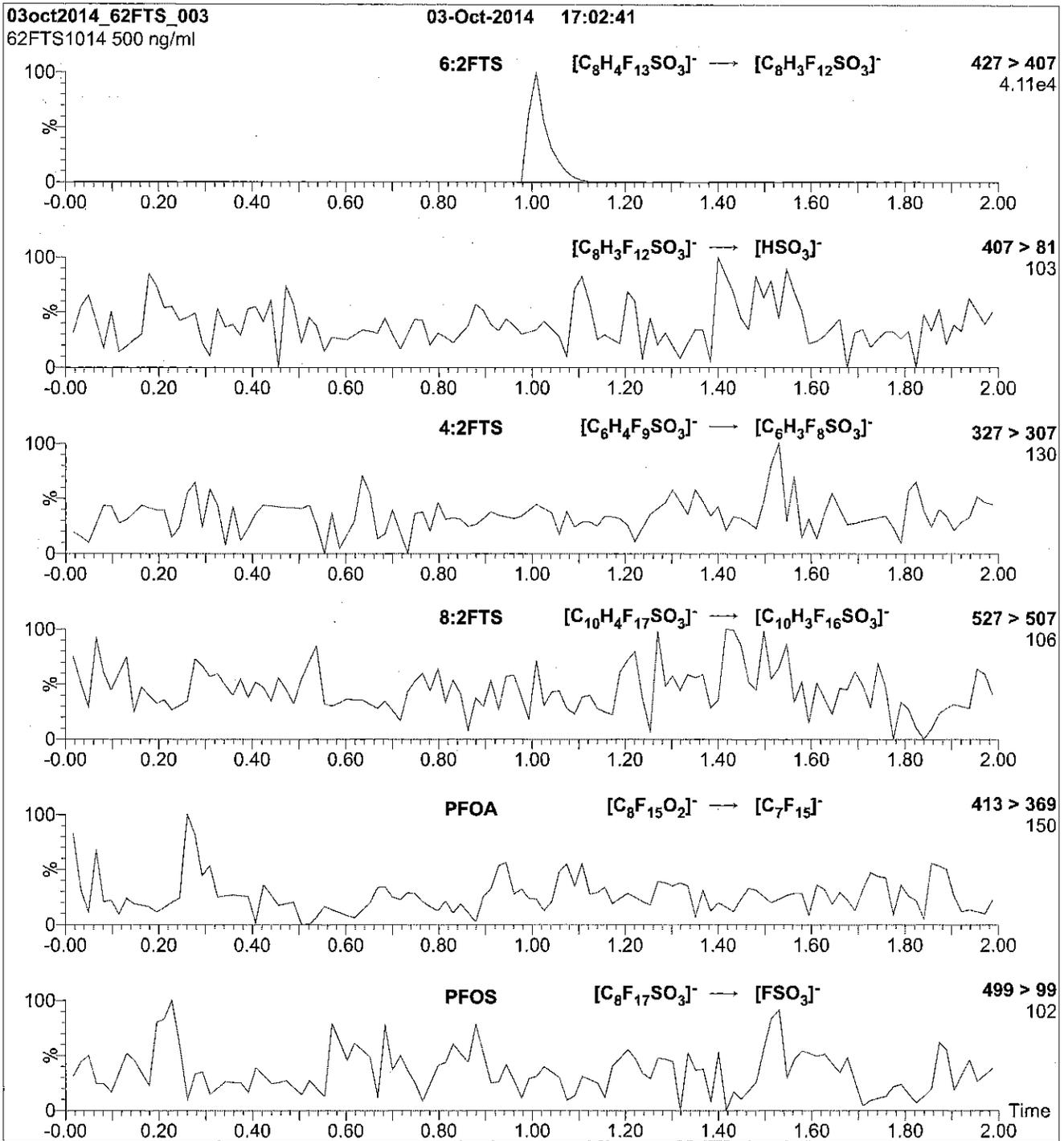
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: 6:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml 6:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.50e-3
Collision Energy (eV) = 25

Reagent

LC6:2FTS_00002

R: 8/23/16 SBC



715544

ID: LC6:2FTS_00002

Exp: 06/25/21 Prod: SBC
6:2FTS

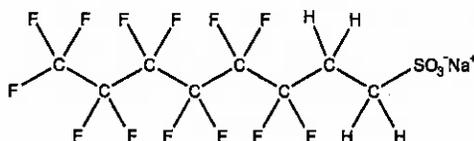


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: 6:2FTS **LOT NUMBER:** 62FTS0616
COMPOUND: Sodium 1H,1H,2H,2H-perfluorooctane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: $C_8H_4F_{13}SO_3Na$ **MOLECULAR WEIGHT:** 450.15
CONCENTRATION: $50.0 \pm 2.5 \mu\text{g/ml}$ (Na salt) **SOLVENT(S):** Methanol
 $47.4 \pm 2.4 \mu\text{g/ml}$ (6:2FTS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 06/25/2016
EXPIRY DATE: (mm/dd/yyyy) 06/25/2021
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

• See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 06/29/2016

(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • Info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

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SYNTHESIS / CHARACTERIZATION:

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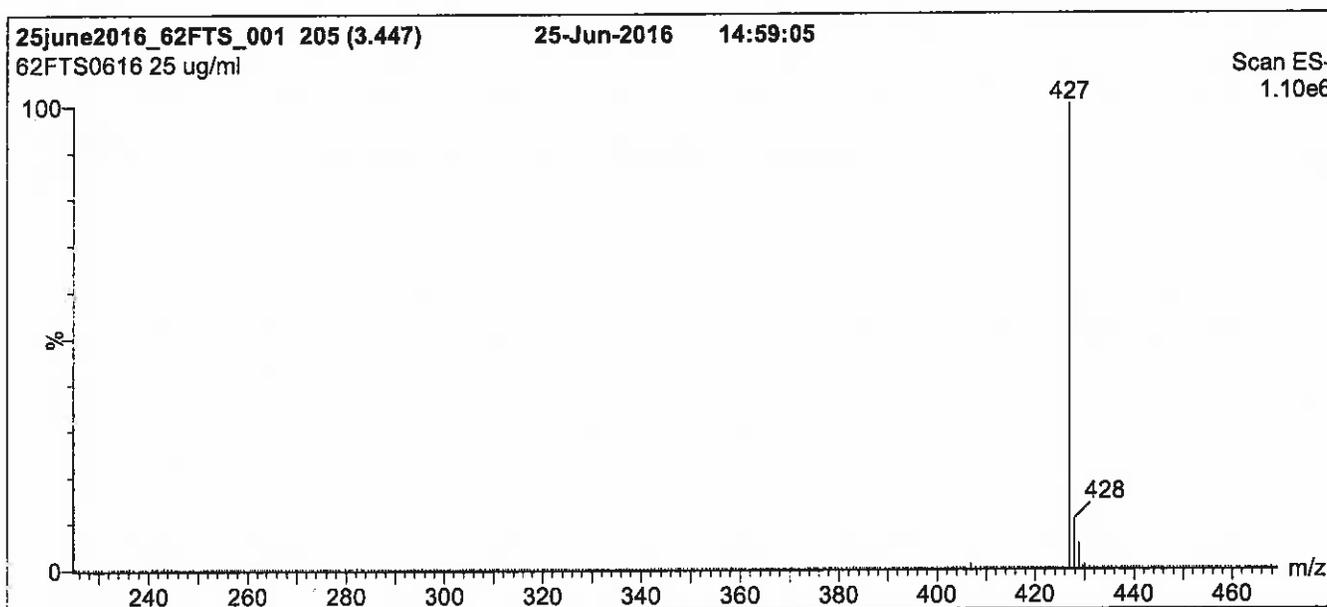
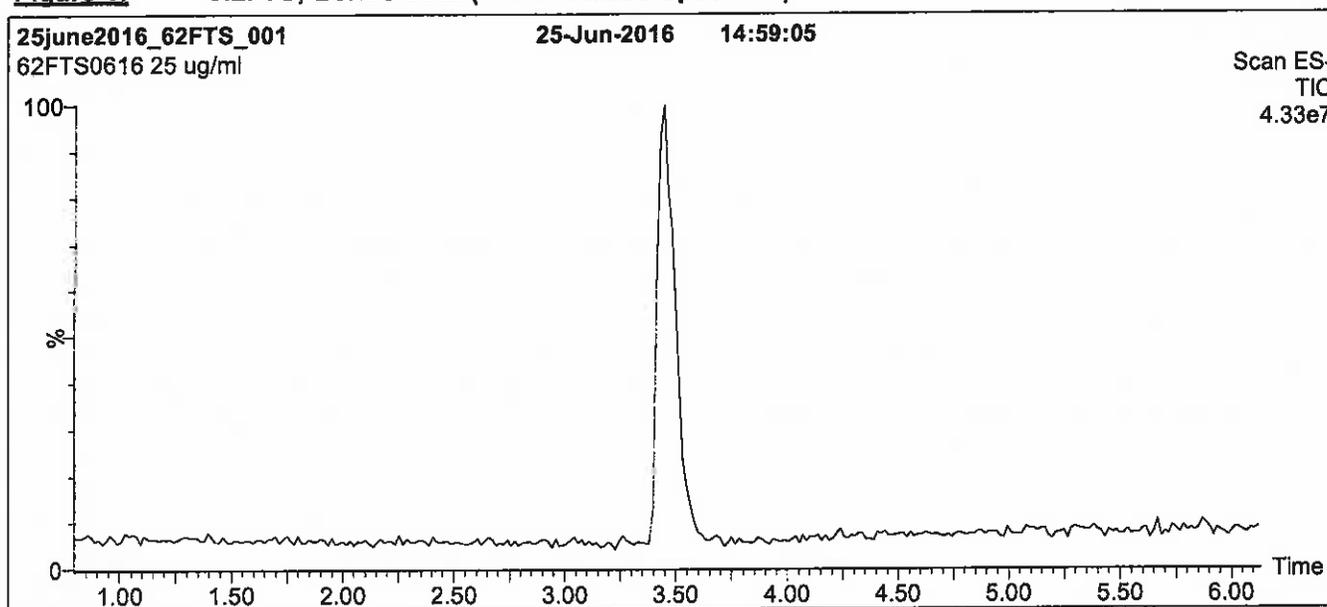
QUALITY MANAGEMENT:

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Figure 1: 6:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

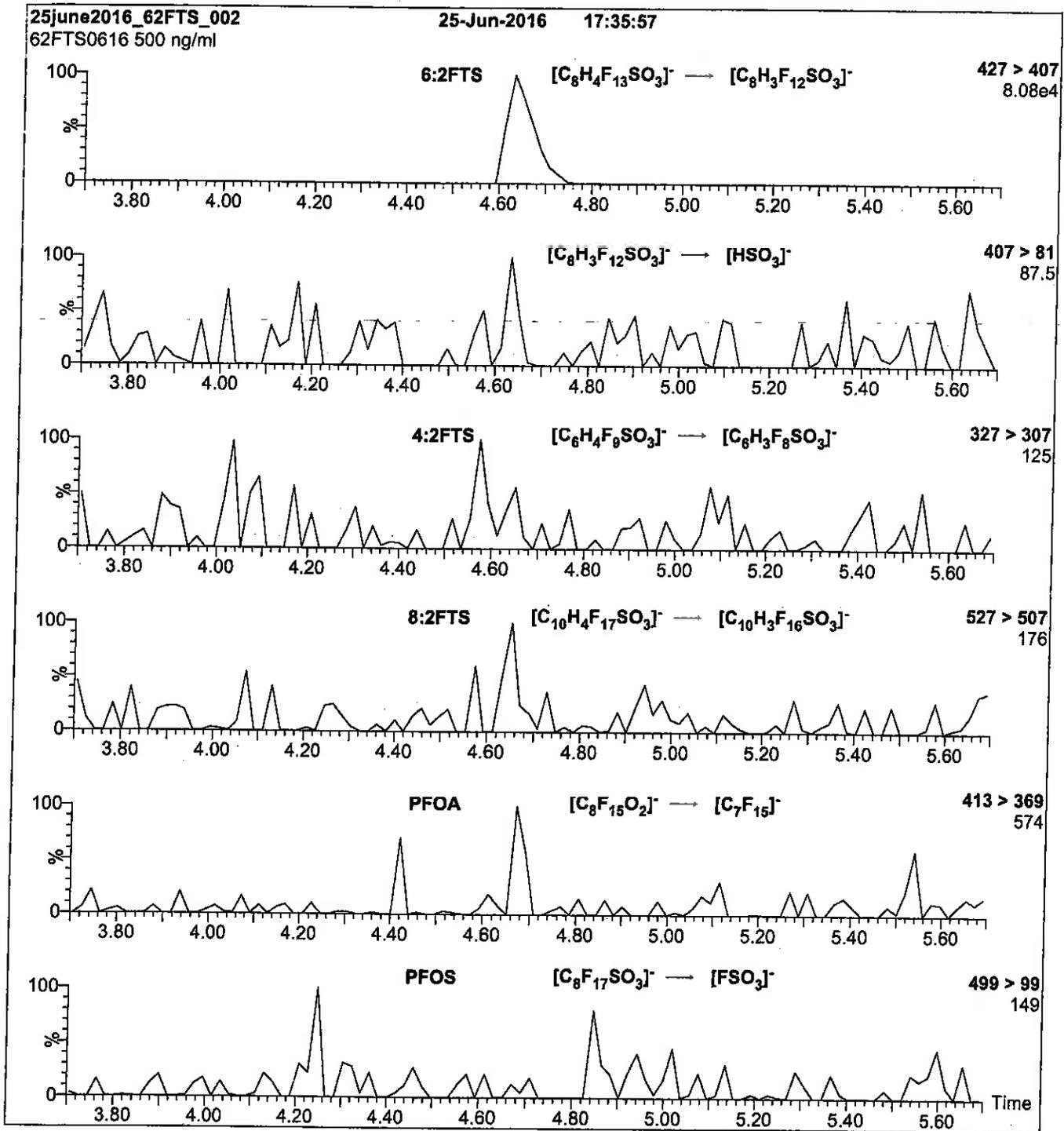
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: 6:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml 6:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.46e-3
Collision Energy (eV) = 25

Reagent

LC8 : 2FTS _ 00001

r: 71615 8V
S: 71215 8V

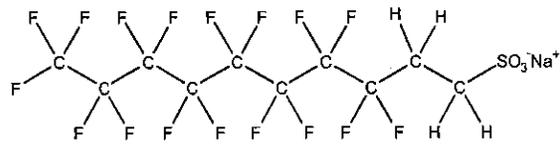


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: 8:2FTS **LOT NUMBER:** 82FTS1014
COMPOUND: Sodium 1H,1H,2H,2H-perfluorodecane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₀H₄F₁₇SO₃Na **MOLECULAR WEIGHT:** 550.16
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.9 ± 2.4 µg/ml (8:2FTS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 10/03/2014
EXPIRY DATE: (mm/dd/yyyy) 10/03/2017
RECOMMENDED STORAGE: Refrigerate ampoule

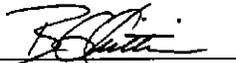
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 03/27/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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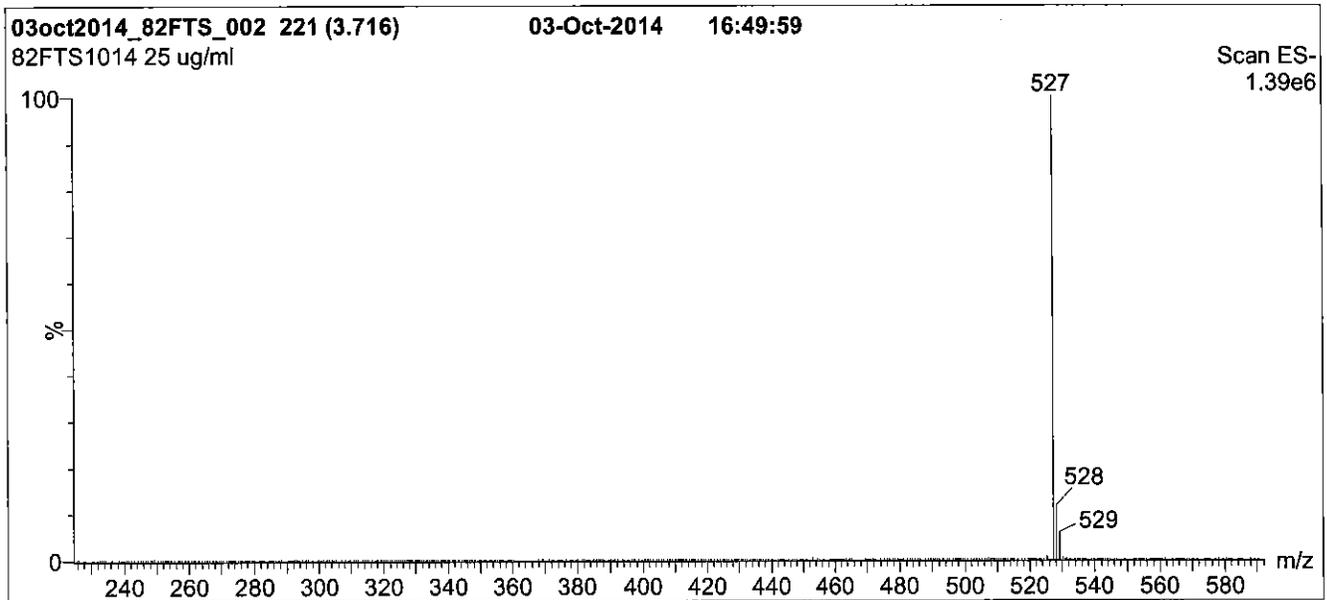
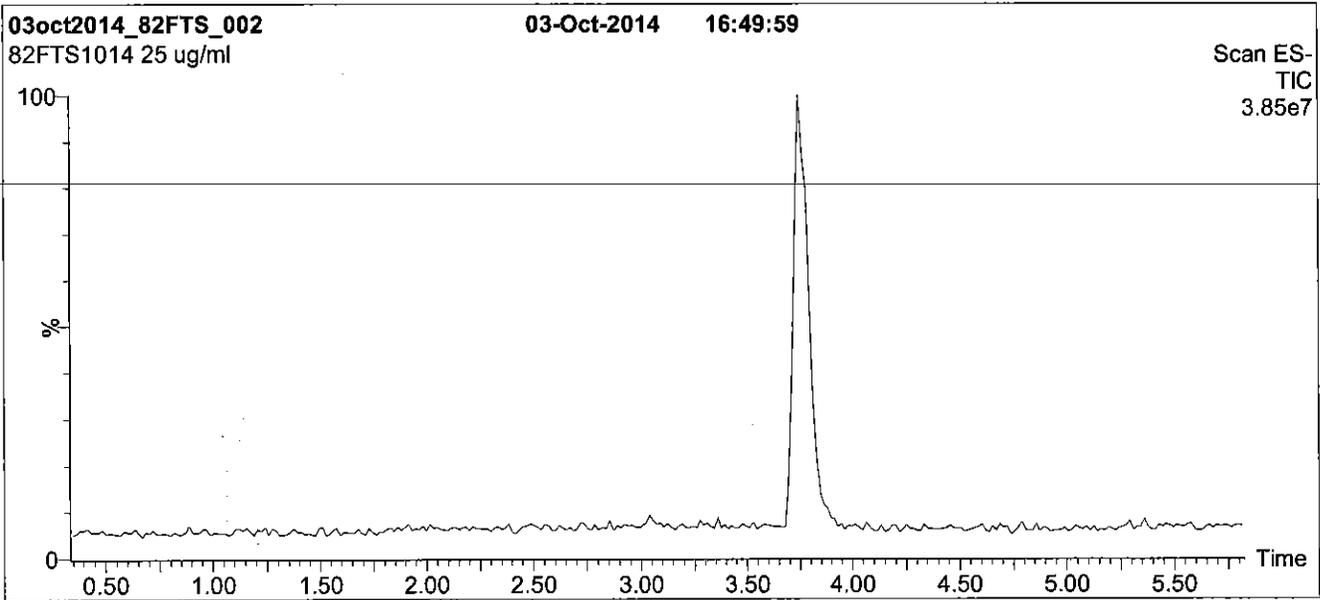
QUALITY MANAGEMENT:

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Figure 1: 8:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro micro API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μm, 2.1 x 100 mm

Mobile phase: Gradient
Start: 55% (80:20 MeOH:ACN) / 45% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min.
Return to initial conditions in 0.5 min.
Time: 10 min

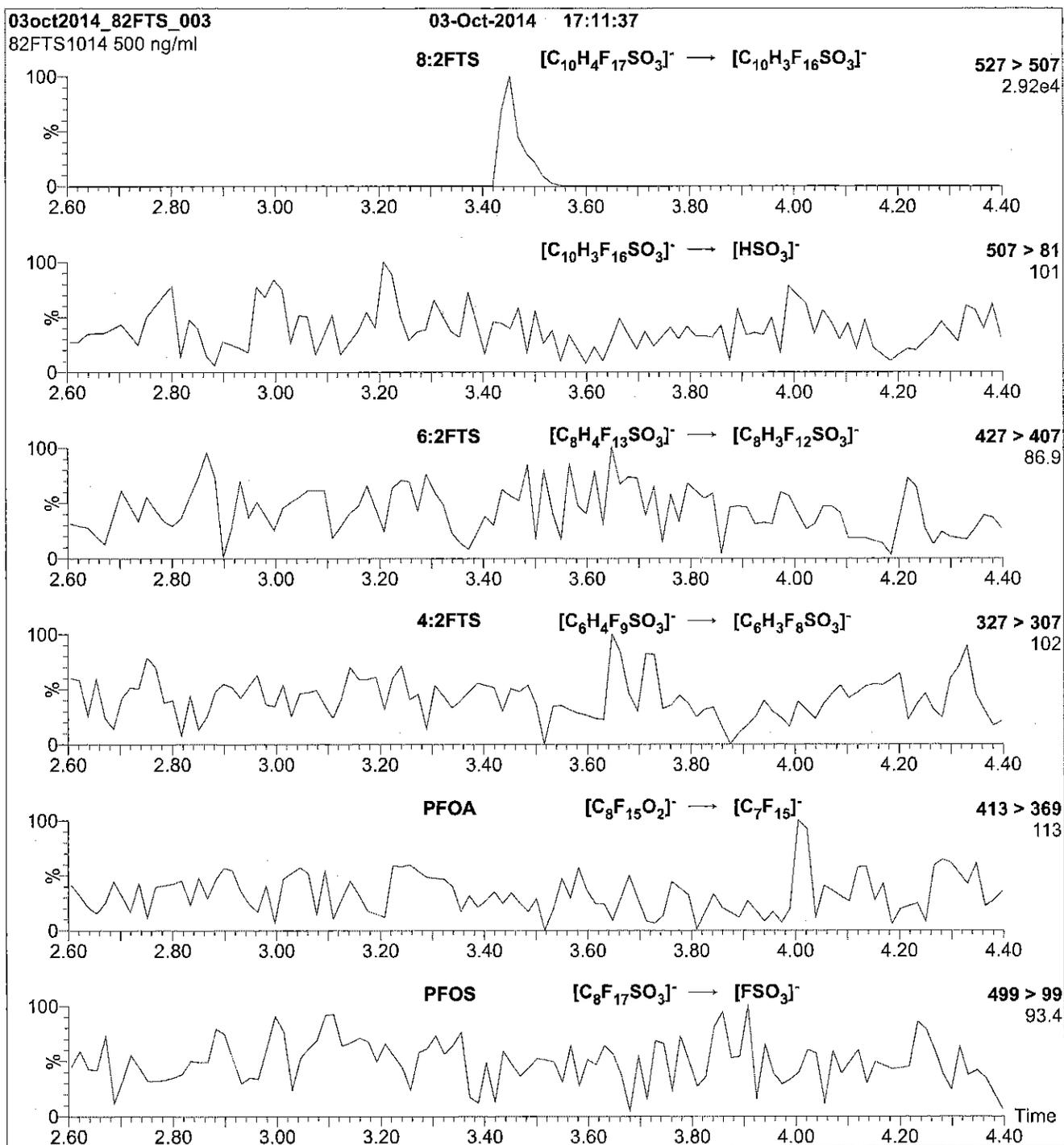
Flow: 300 μl/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: 8:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml 8:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.50e-3
Collision Energy (eV) = 30

Reagent

LC8 : 2FTS _ 00002

R: 8/23/16 SBC

715545
ID: LC8:2FTS_00002
Exp: 10/23/20 Prod: SBC
8:2FTS

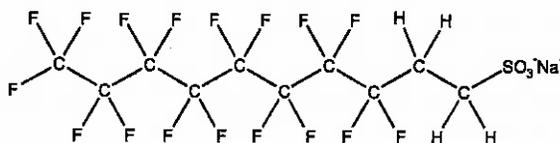


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: 8:2FTS **LOT NUMBER:** 82FTS1015
COMPOUND: Sodium 1H,1H,2H,2H-perfluorodecane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₀H₄F₁₇SO₃Na **MOLECULAR WEIGHT:** 550.16
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.9 ± 2.4 µg/ml (8:2FTS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 10/23/2015
EXPIRY DATE: (mm/dd/yyyy) 10/23/2020
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 10/27/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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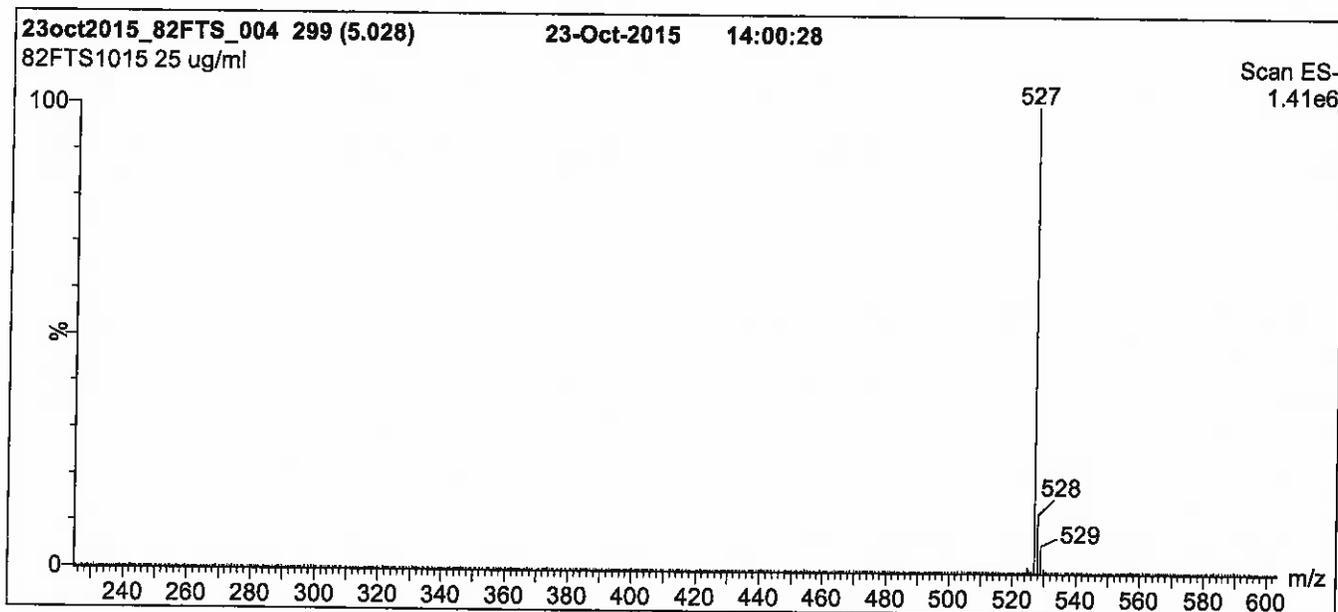
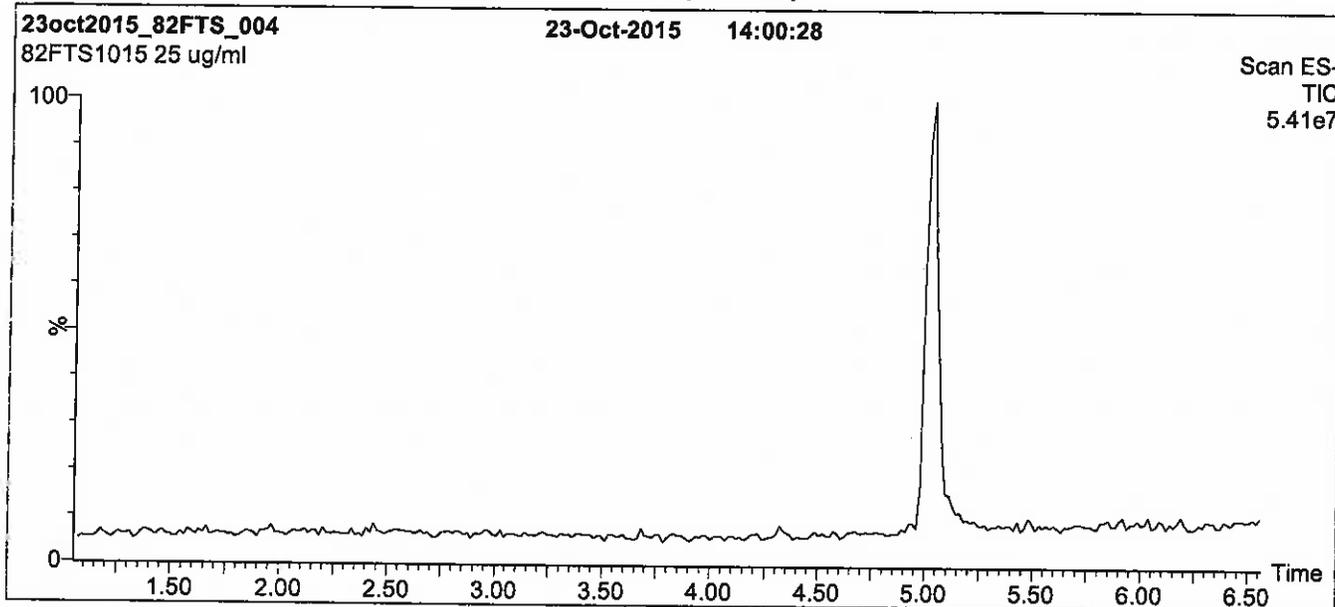
QUALITY MANAGEMENT:

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Figure 1: 8:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min.
Return to Initial conditions in 0.5 min.
Time: 10 min

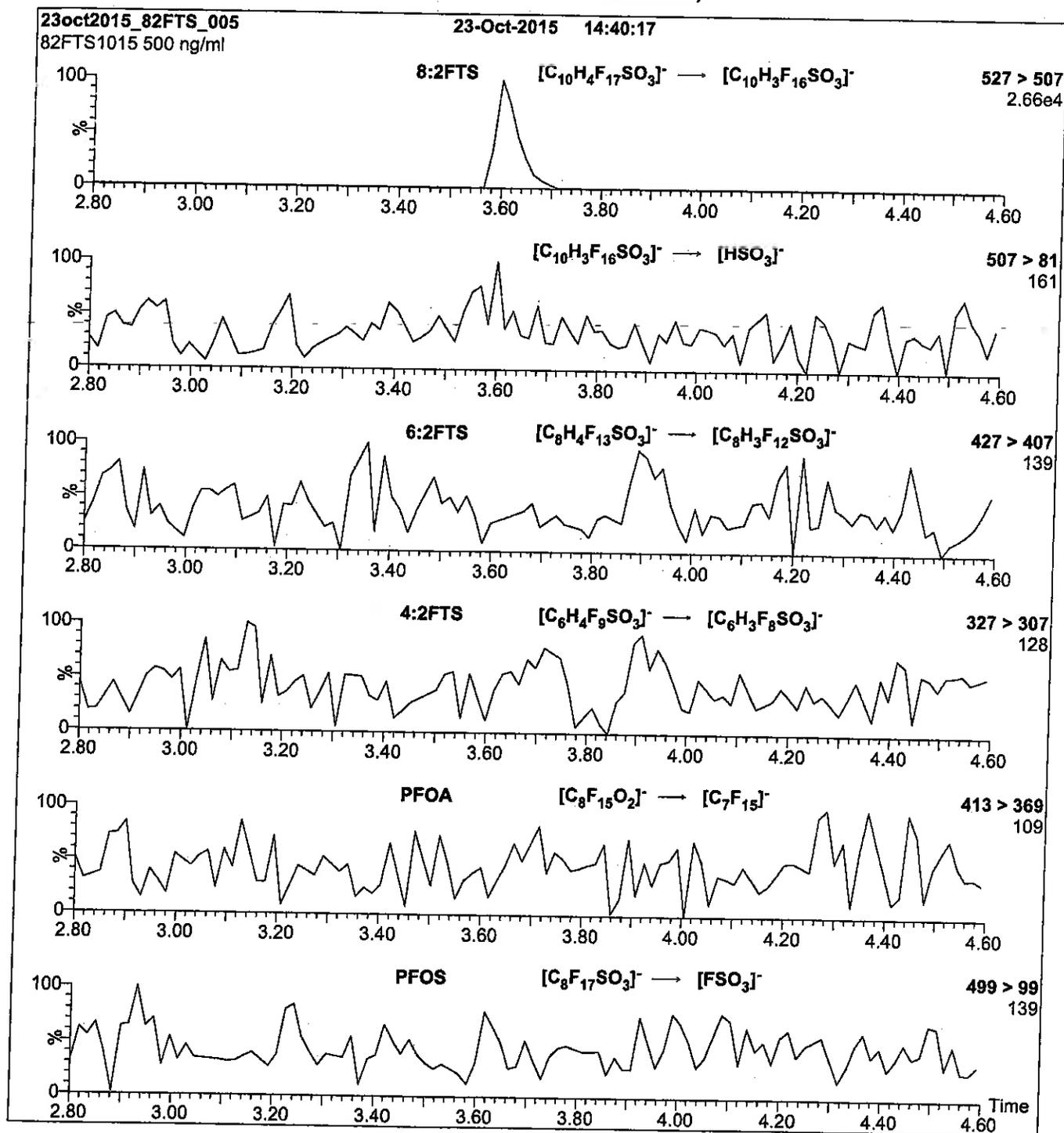
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: 8:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml 8:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.28e-3
Collision Energy (eV) = 30

Reagent

LCd-NEtFOSA-M_00001

C: 7/16/15 8/



WELLINGTON LABORATORIES

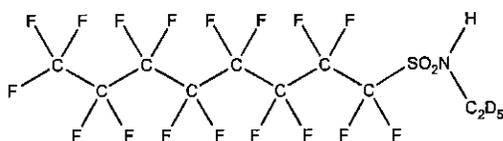
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d-N-EtFOSA-M
COMPOUND: N-ethyl-d₅-perfluoro-1-octanesulfonamide

LOT NUMBER: dNEtFOSA0314M

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: C₁₀D₅HF₁₇NO₂S
CONCENTRATION: 50 ± 2.5 µg/ml
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 03/10/2014
EXPIRY DATE: (mm/dd/yyyy) 03/10/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 532.23
SOLVENT(S): Methanol
ISOTOPIC PURITY: ≥98% ²H₅

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:
B.G. Chittim

Date: 04/01/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

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UNCERTAINTY:

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$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

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EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

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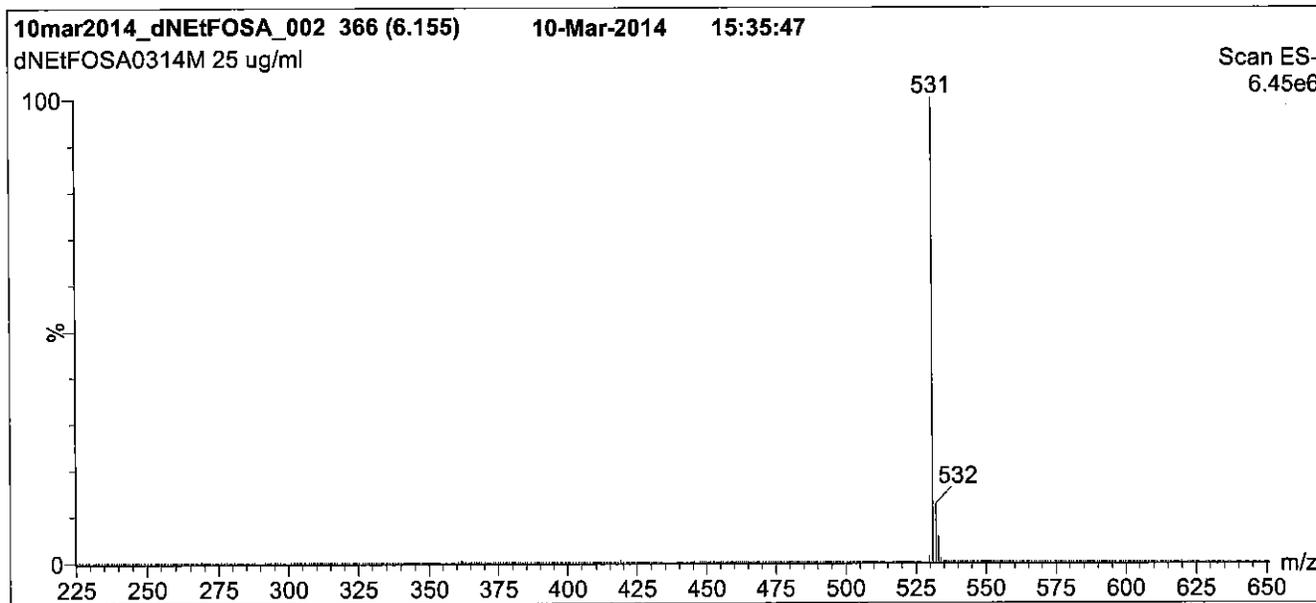
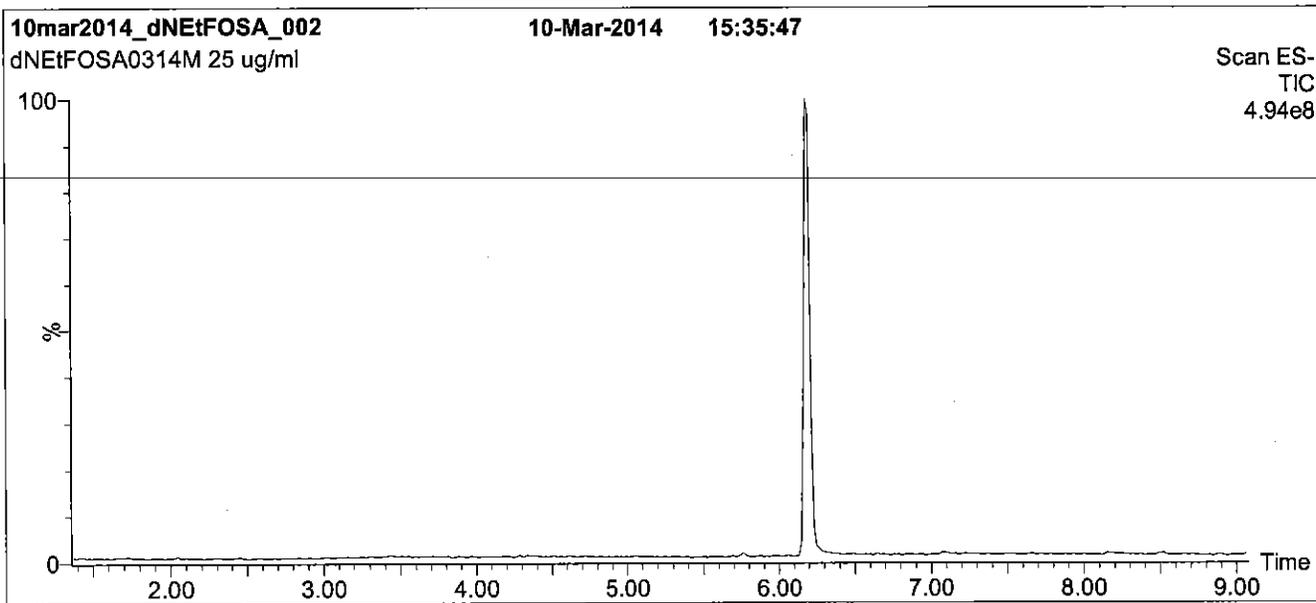
QUALITY MANAGEMENT:

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Figure 1: d-N-EtFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 40% H₂O / 60% (80:20 MeOH:ACN)
 (both with 10mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

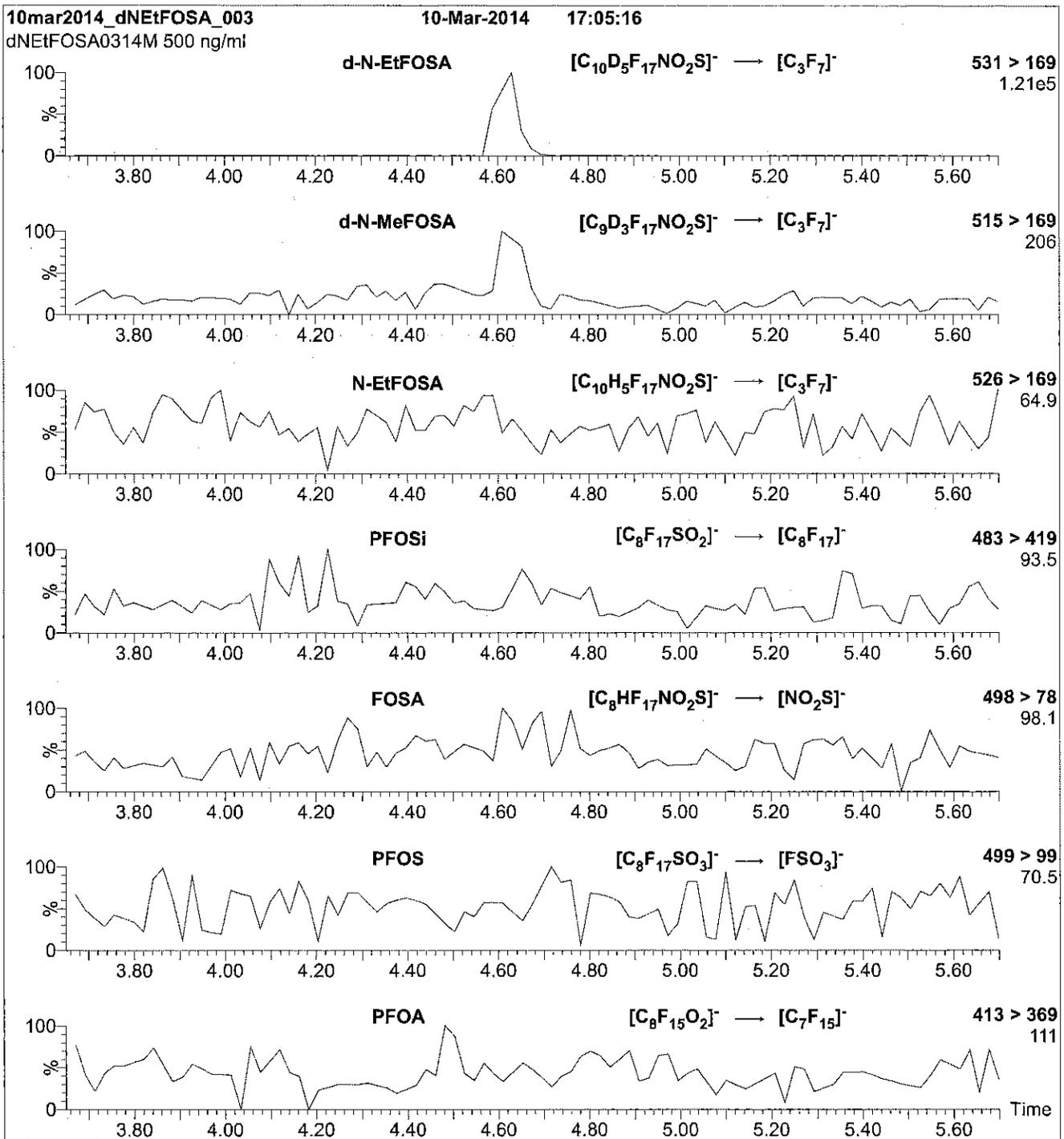
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 950 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d-N-EtFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d-N-EtFOSA-M)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
Collision Energy (eV) = 25

Reagent

LCd-NEtFOSA-M_00002

R-7/6/16 CAW



671571
ID: LCd-NEtFOSA-M_00002
Exp: 03/10/19 Pipd: CBW
d-N-EtFOSA-M



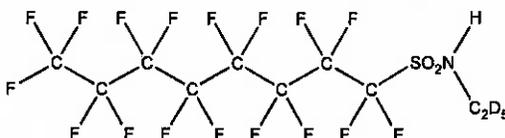
WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d-N-EtFOSA-M
COMPOUND: N-ethyl-d₅-perfluoro-1-octanesulfonamide

LOT NUMBER: dNEtFOSA0314M

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₀D₅HF₁₇NO₂S
CONCENTRATION: 50 ± 2.5 µg/ml
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 03/10/2014
EXPIRY DATE: (mm/dd/yyyy) 03/10/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 532.23
SOLVENT(S): Methanol
ISOTOPIC PURITY: ≥98% ²H₅

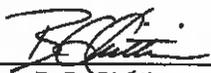
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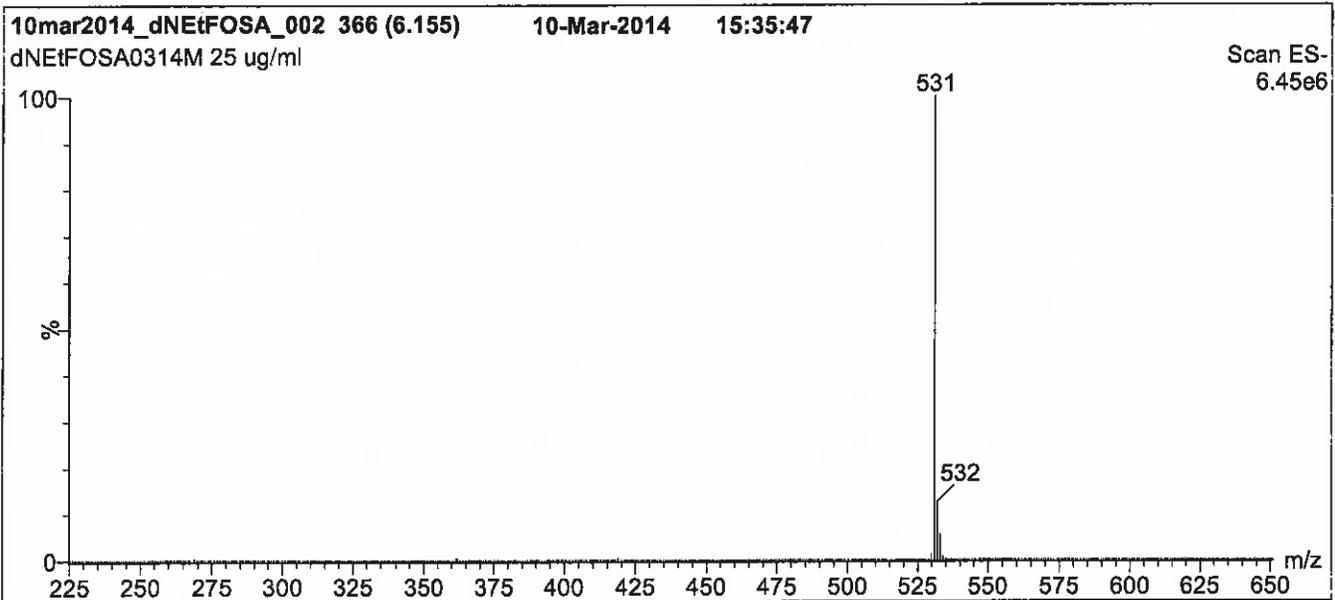
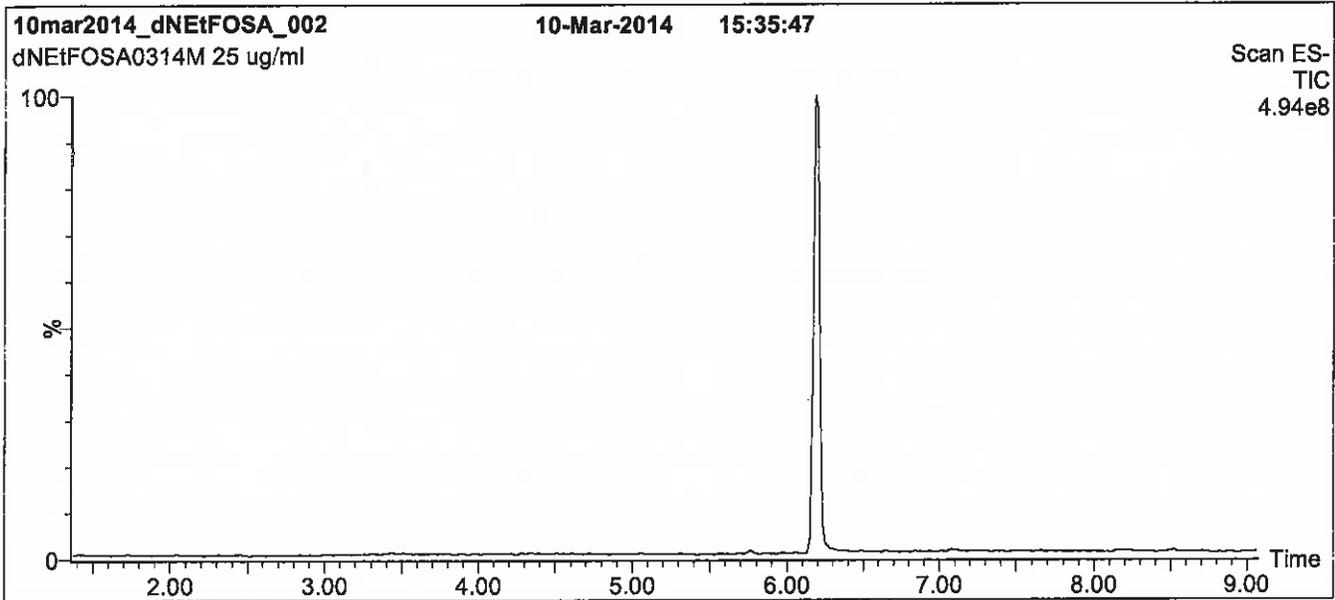
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Figure 1: d-N-EtFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gracient
 Start: 40% H₂O / 60% (80:20 MeOH:ACN)
 (both with 10mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

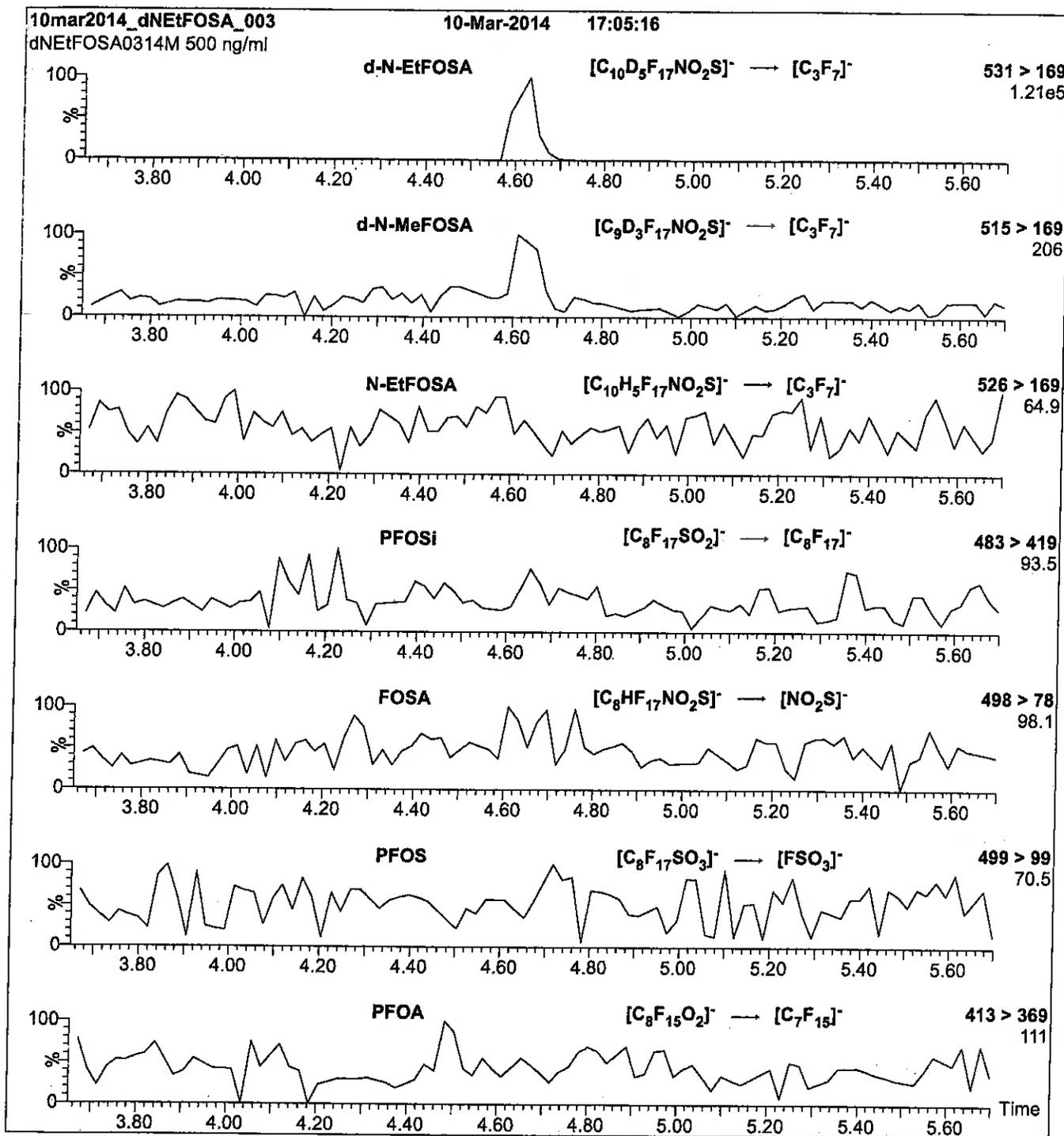
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 950 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d-N-EtFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d-N-EtFOSA-M)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
Collision Energy (eV) = 25

Reagent

LCd-NMeFOSA-M_00001

r: 7/16/15 SKW



WELLINGTON LABORATORIES

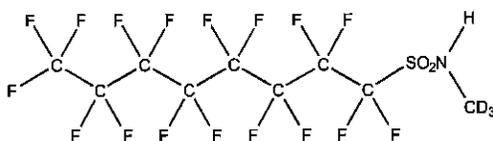
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d-N-MeFOSA-M
COMPOUND: N-methyl-d₃-perfluoro-1-octanesulfonamide

LOT NUMBER: dNMeFOSA0114M

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: C₉D₃HF₁₇NO₂S
CONCENTRATION: 50 ± 2.5 µg/ml
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/28/2014
EXPIRY DATE: (mm/dd/yyyy) 01/28/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 516.19
SOLVENT(S): Methanol
ISOTOPIC PURITY: ≥98% ²H₃

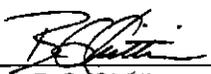
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Certified By: 
B.G. Chittim
Date: 04/01/2015
(mm/dd/yyyy)

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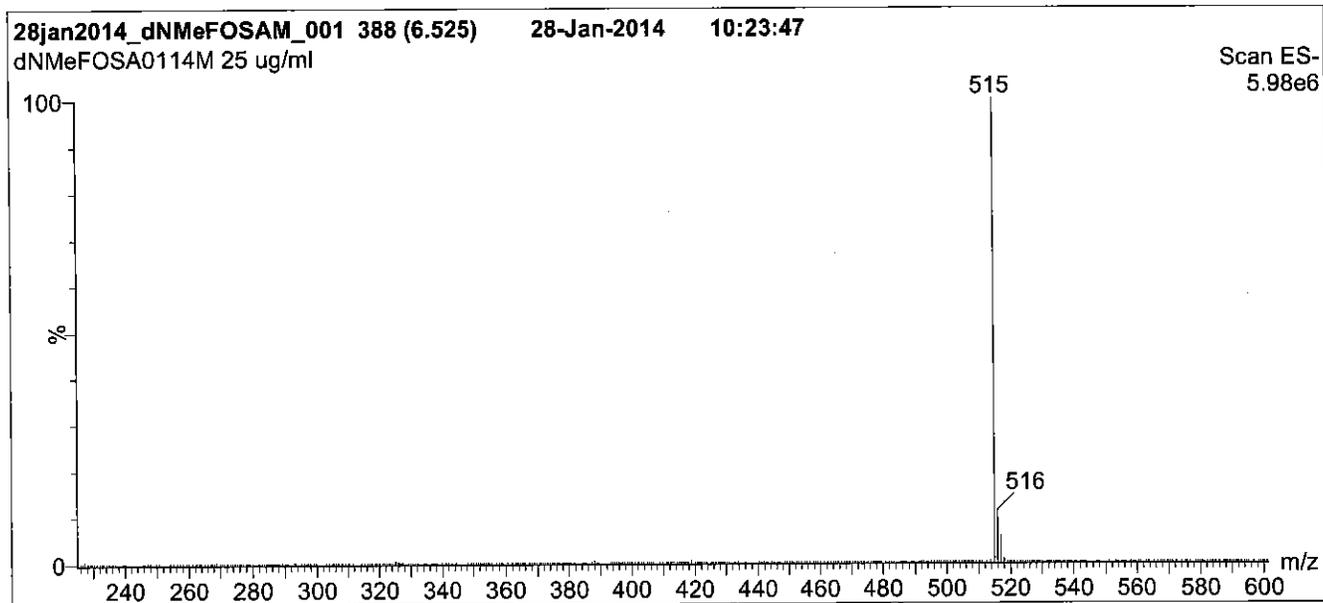
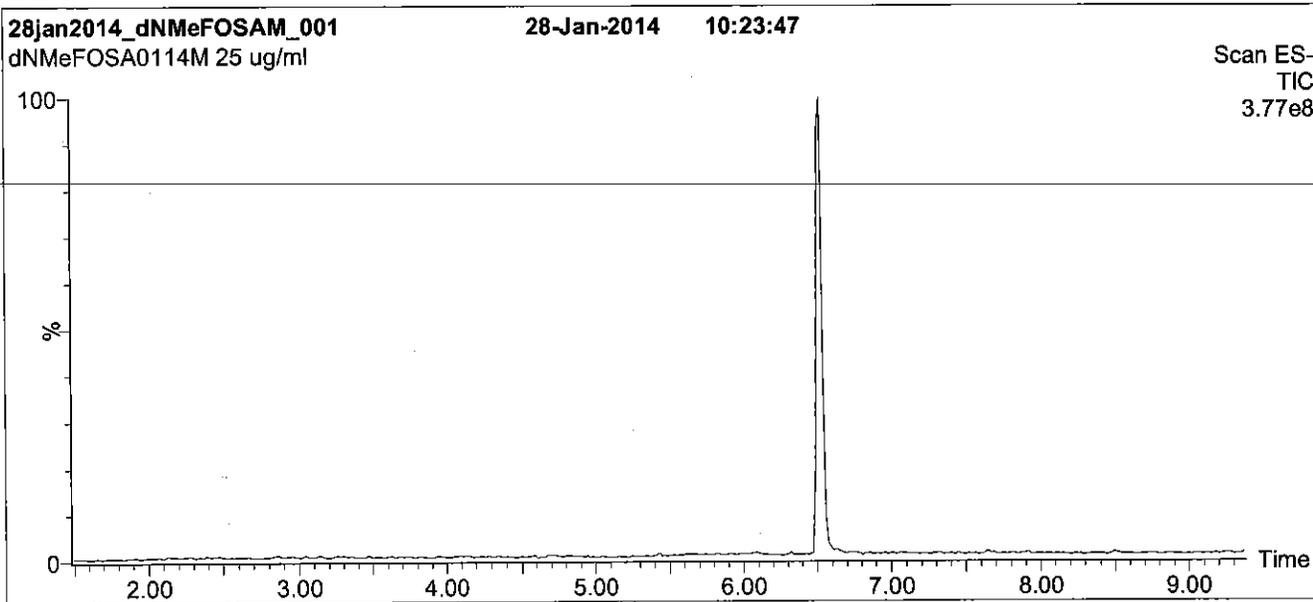
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Figure 1: d-N-MeFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% H₂O / 50% (80:20 MeOH:ACN)
 (both with 10mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 1.5 min. Return to initial conditions over 0.5 min.
 Time: 10 min

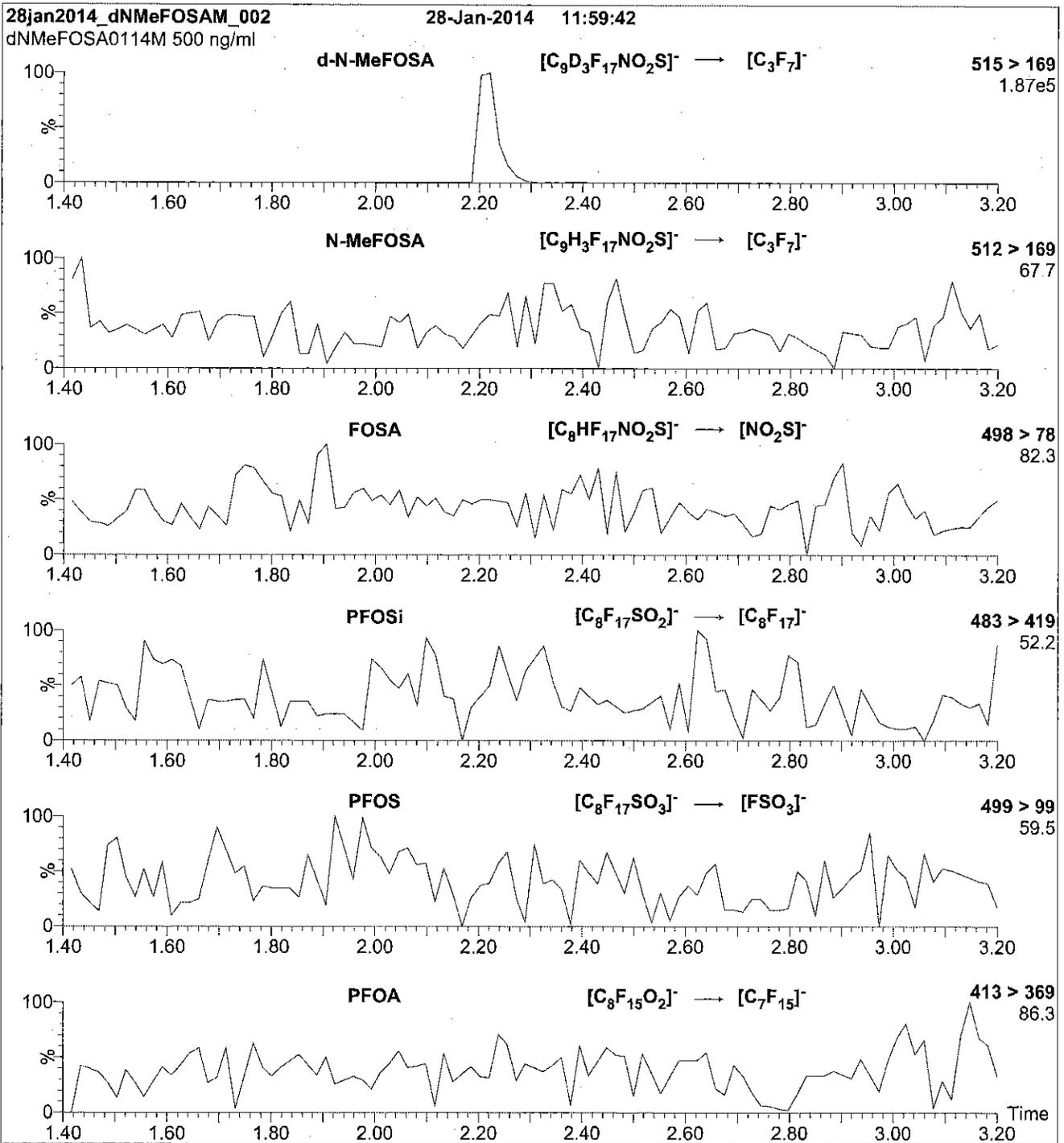
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.50
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d-N-MeFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d-N-MeFOSA-M)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.50e-3
Collision Energy (eV) = 30

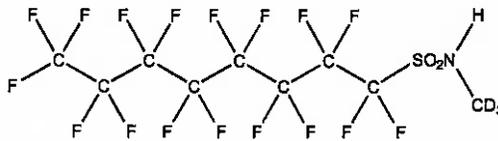
Reagent

LCd-NMeFOSA-M_00002

**WELLINGTON
LABORATORIES****CERTIFICATE OF ANALYSIS
DOCUMENTATION**

PRODUCT CODE: d-N-MeFOSA-M **LOT NUMBER:** dNMeFOSA0616M
COMPOUND: N-methyl-d₃-perfluoro-1-octanesulfonamide

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₈D₃HF₁₇NO₂S **MOLECULAR WEIGHT:** 516.19
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥98% ²H₃
LAST TESTED: (mm/dd/yyyy) 06/10/2016
EXPIRY DATE: (mm/dd/yyyy) 06/10/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

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Certified By: _____

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Date: 06/16/2016

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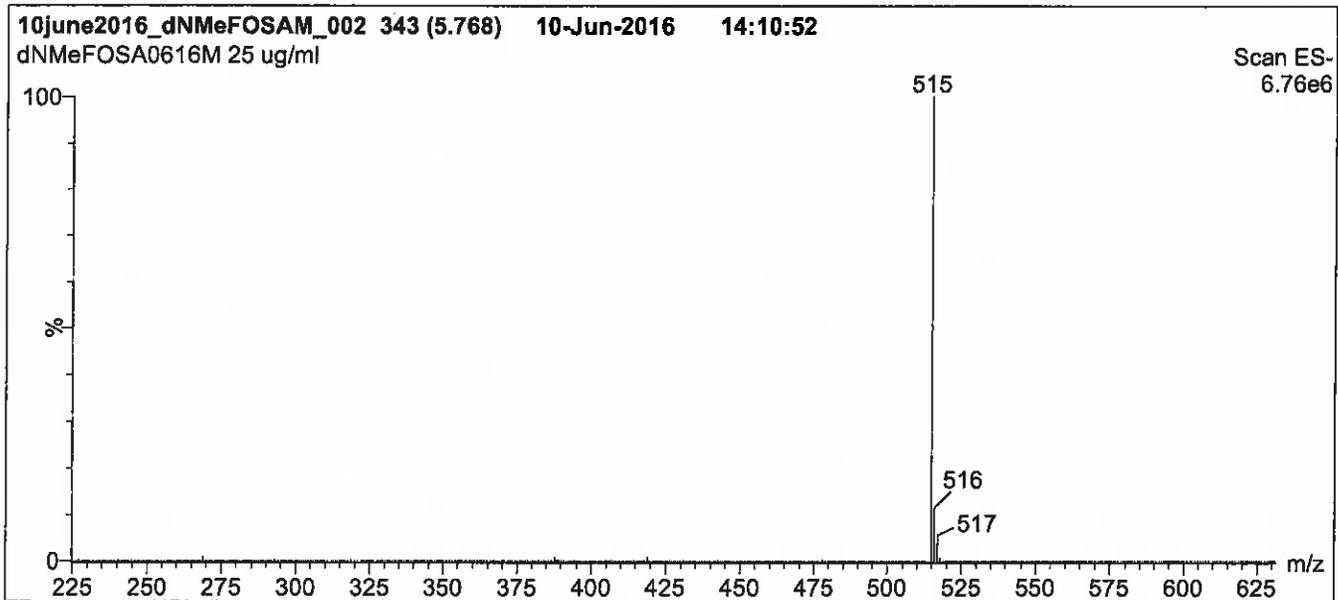
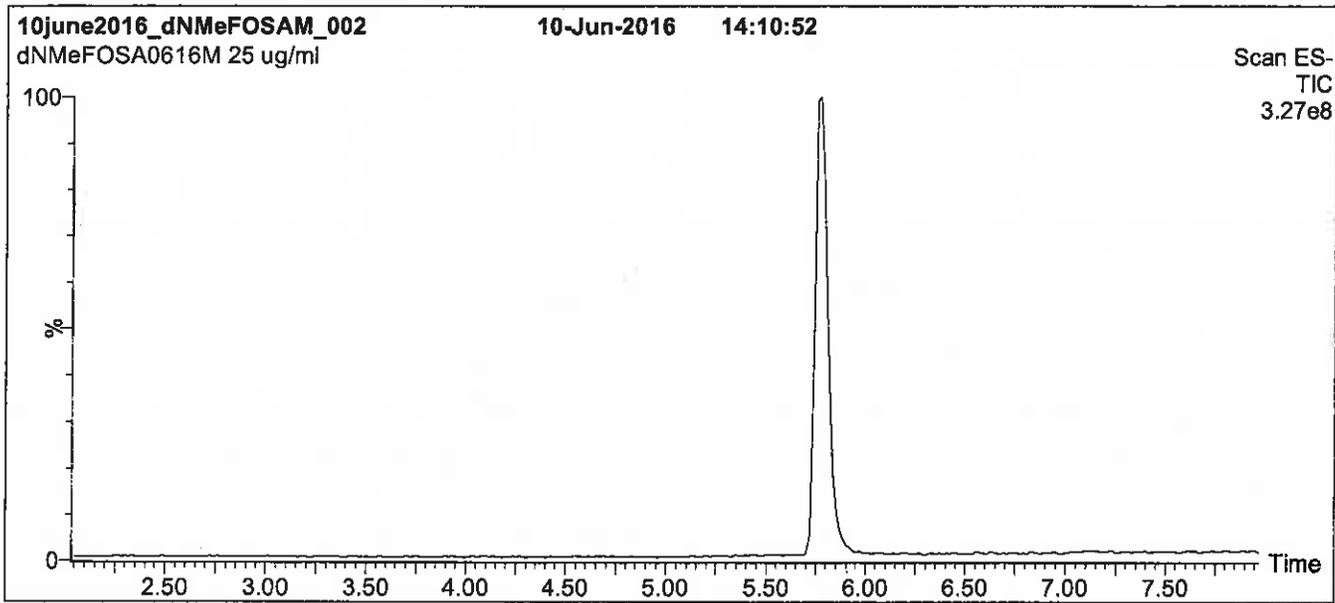
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: d-N-MeFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 40% H₂O / 60% (80:20 MeOH:ACN)
(both with 10mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

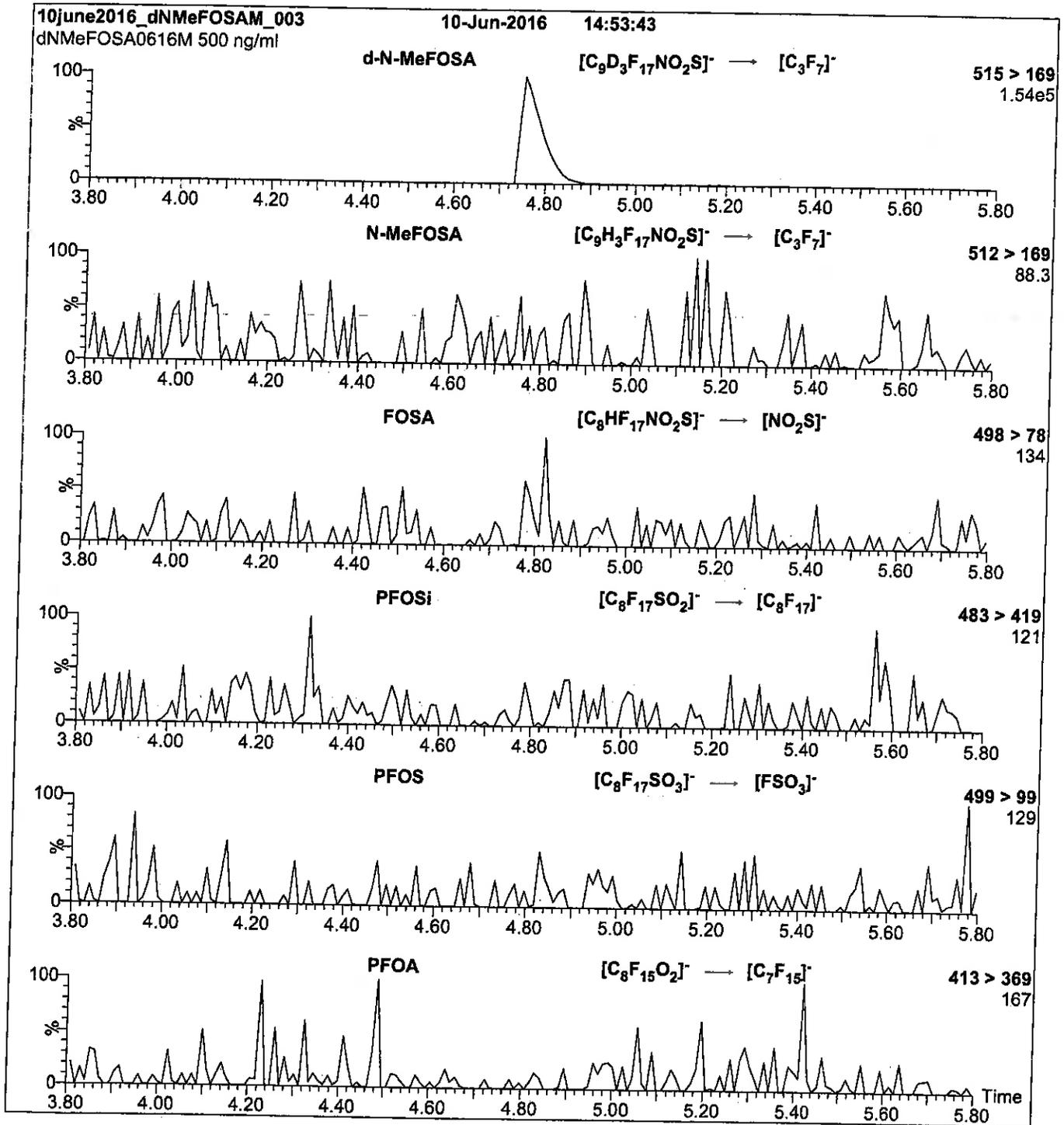
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.50
Cone Voltage (V) = 40.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: d-N-MeFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d-N-MeFOSA-M)

Mobile phase: isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.39e-3
Collision Energy (eV) = 25

Reagent

LCd3-NMeFOSAA_00001

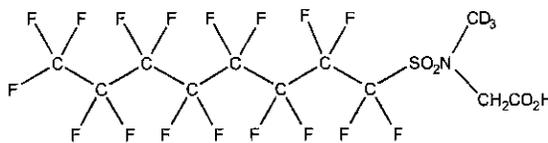


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d3-N-MeFOSAA **LOT NUMBER:** d3NMeFOSAA0113
COMPOUND: N-methyl-d3-perfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: $C_{11}D_3H_3F_{17}NO_4S$ **MOLECULAR WEIGHT:** 574.23
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Methanol
 Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** $\geq 98\% \text{ } ^2\text{H}_3$
LAST TESTED: (mm/dd/yyyy) 01/31/2013
EXPIRY DATE: (mm/dd/yyyy) 01/31/2018
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: _____


 B.G. Chittim

Date: 04/06/2015
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

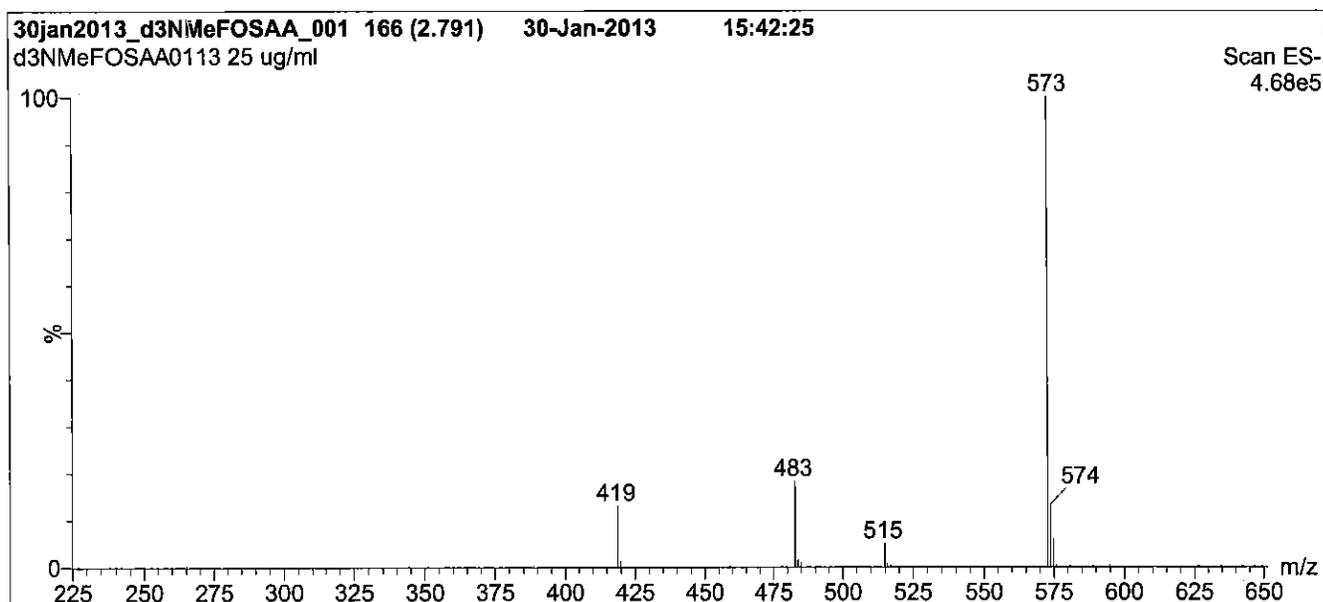
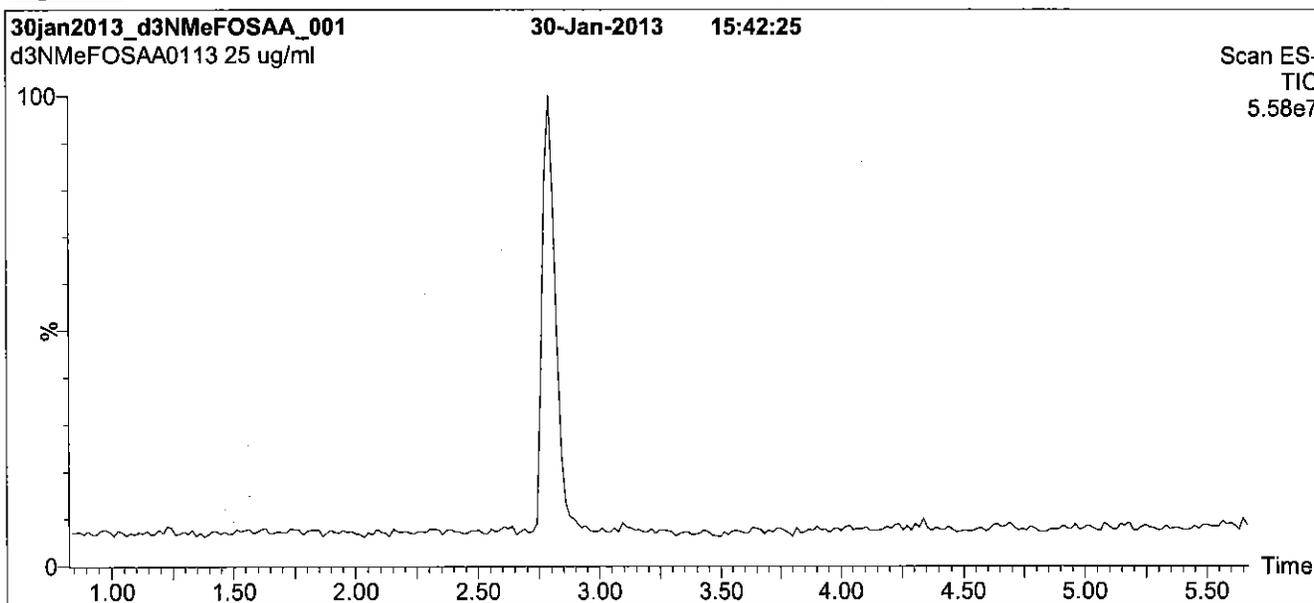
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: d3-N-MeFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 65% (80:20 MeOH:ACN) / 35% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

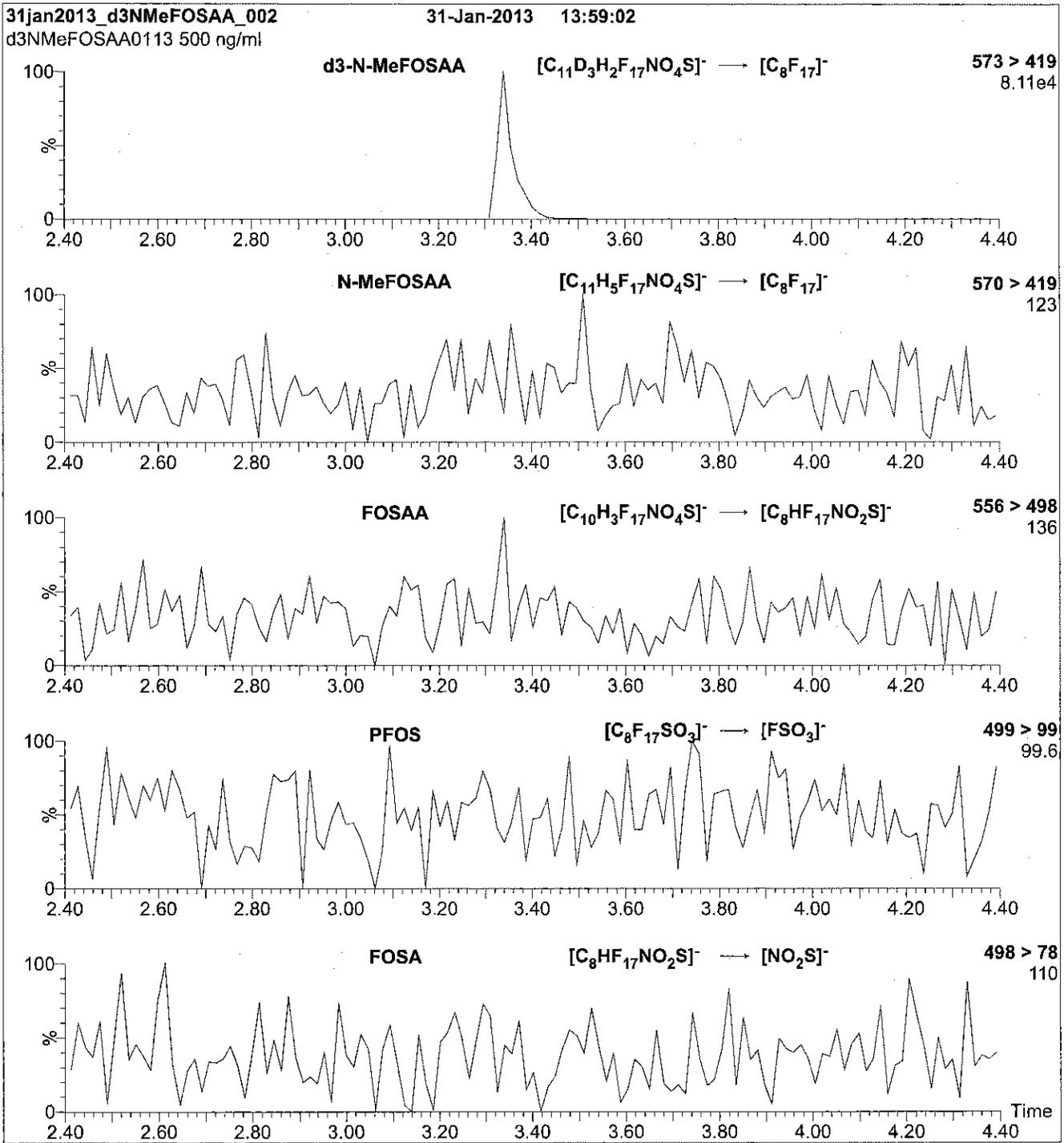
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 35.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: d3-N-MeFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d3-N-MeFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
Collision Energy (eV) = 25

Reagent

LCd3-NMeFOSAA_00002

R: 7/6/16 CBW

671572
ID: LCd3-NMeFOSAA_00002
Exp: 01/2021 Prpd: CBW
d3-N-MeFOSAA

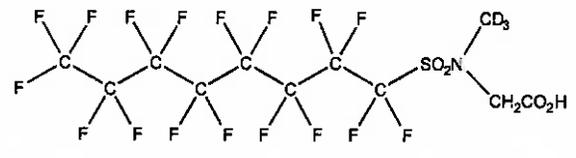


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d3-N-MeFOSAA **LOT NUMBER:** d3NMeFOSAA0116
COMPOUND: N-methyl-d3-perfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₁D₃H₃F₁₇NO₄S **MOLECULAR WEIGHT:** 574.23
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥98% ²H₃
LAST TESTED: (mm/dd/yyyy) 01/20/2016
EXPIRY DATE: (mm/dd/yyyy) 01/20/2021
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 01/25/2016
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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EXPIRY DATE / PERIOD OF VALIDITY:

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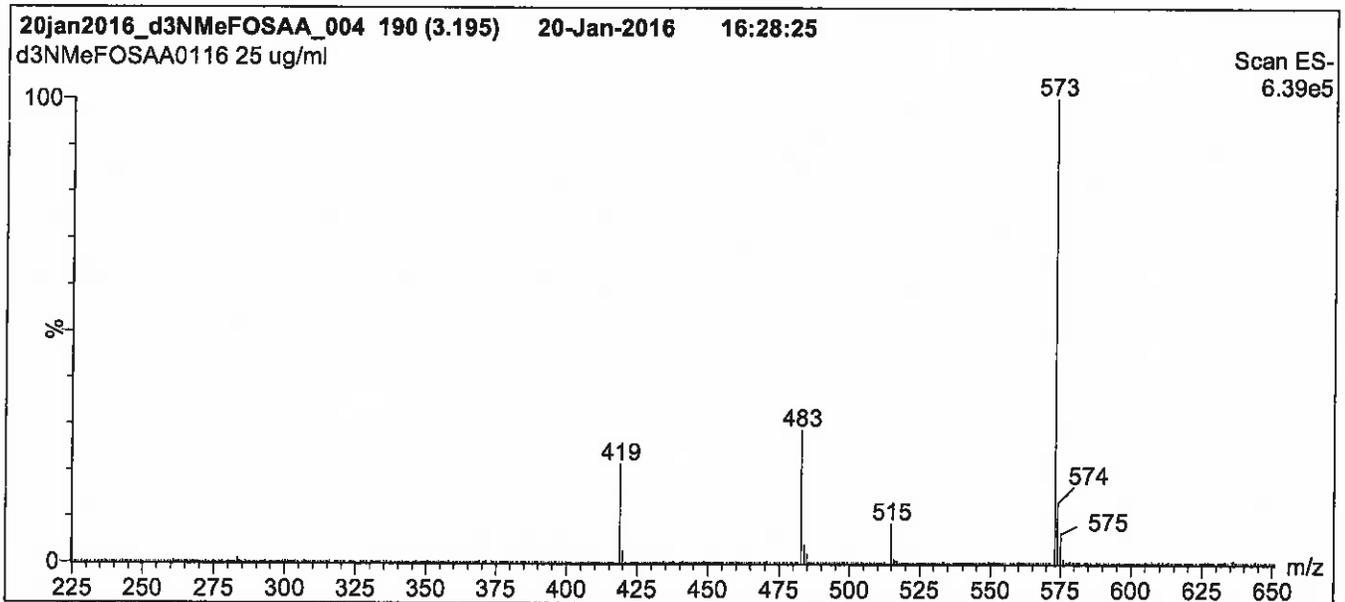
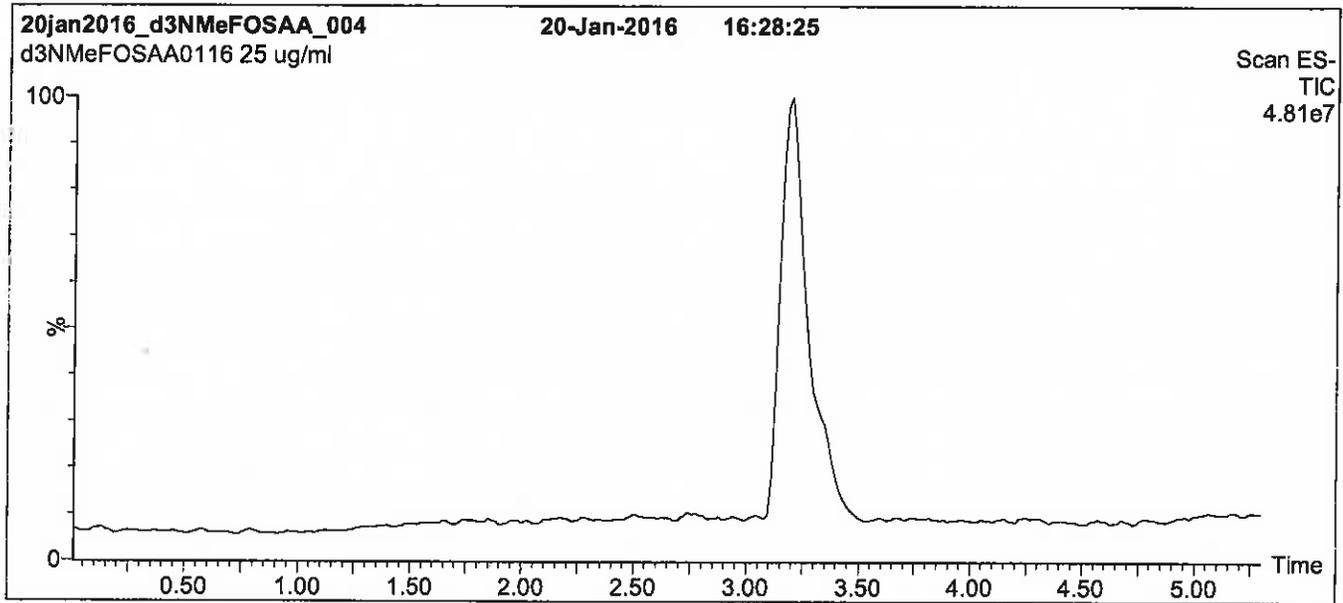
QUALITY MANAGEMENT:

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Figure 1: d3-N-MeFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

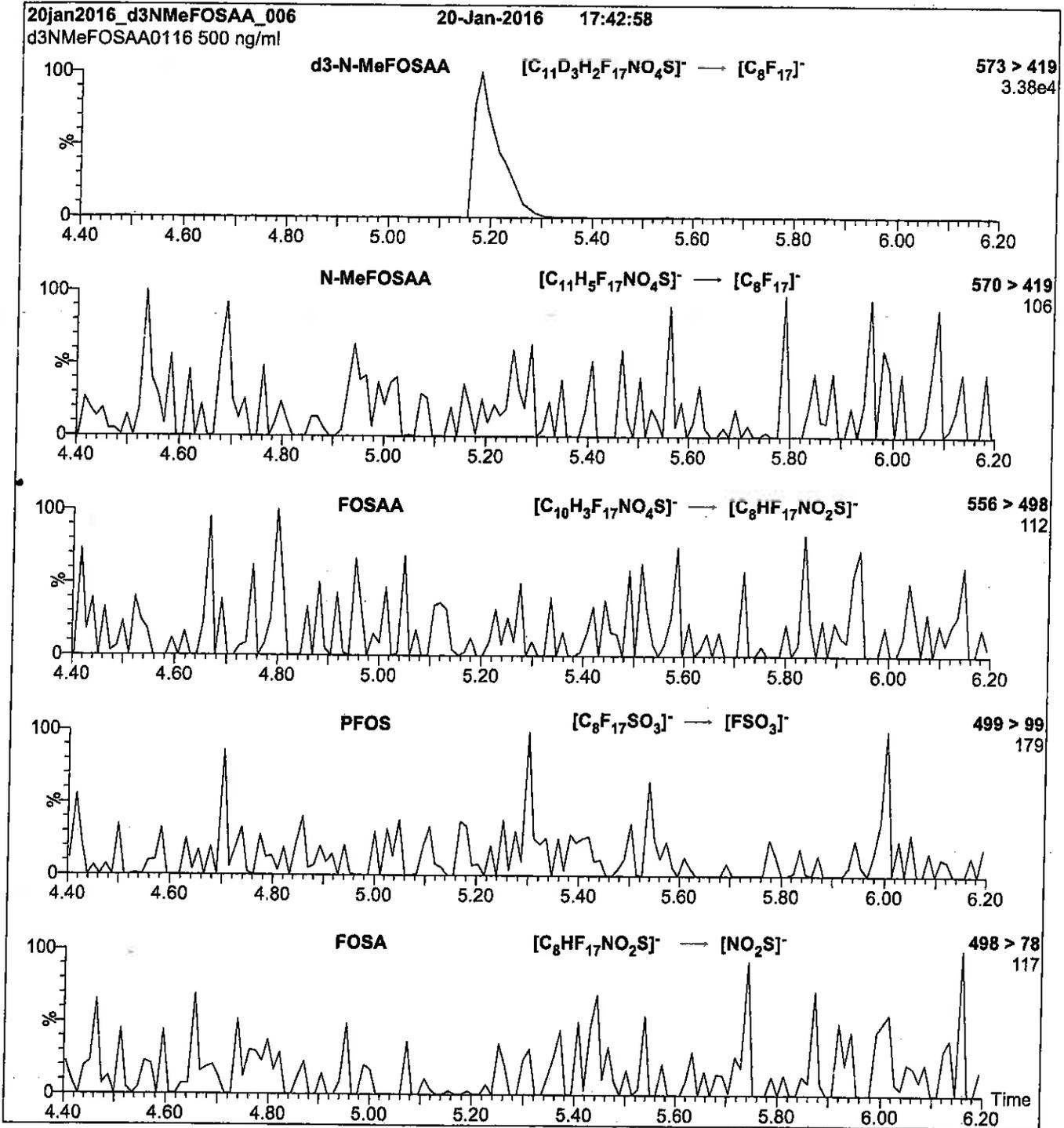
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d3-N-MeFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d3-N-MeFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.66e-3
Collision Energy (eV) = 25

Reagent

LCd5-NEtFOSAA_00001

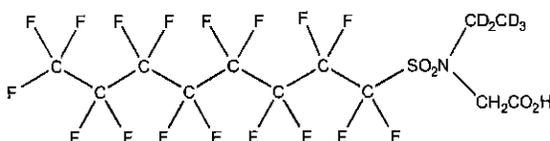


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d5-N-EtFOSAA **LOT NUMBER:** d5NEtFOSAA0515
COMPOUND: N-ethyl-d5-perfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₂D₅H₃F₁₇NO₄S
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 590.27
SOLVENT(S): Methanol
 Water (<1%)

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 05/08/2015
EXPIRY DATE: (mm/dd/yyyy) 05/08/2020
RECOMMENDED STORAGE: Refrigerate ampoule

ISOTOPIC PURITY: ≥98% ²H₅

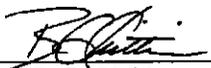
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 05/11/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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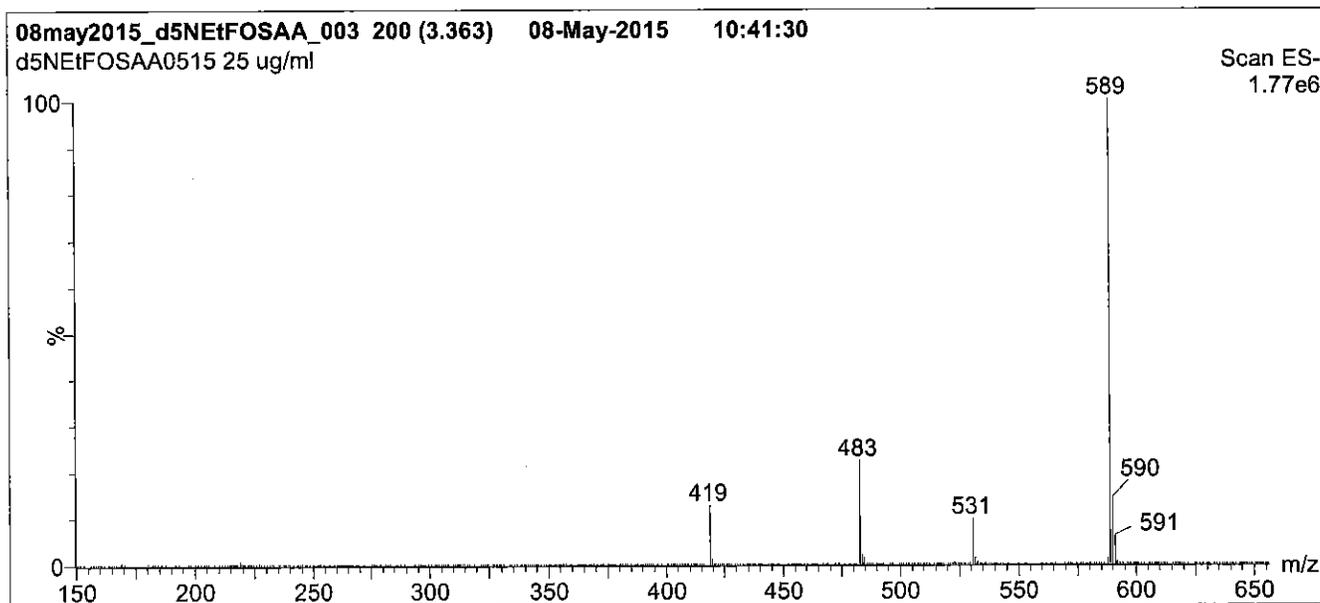
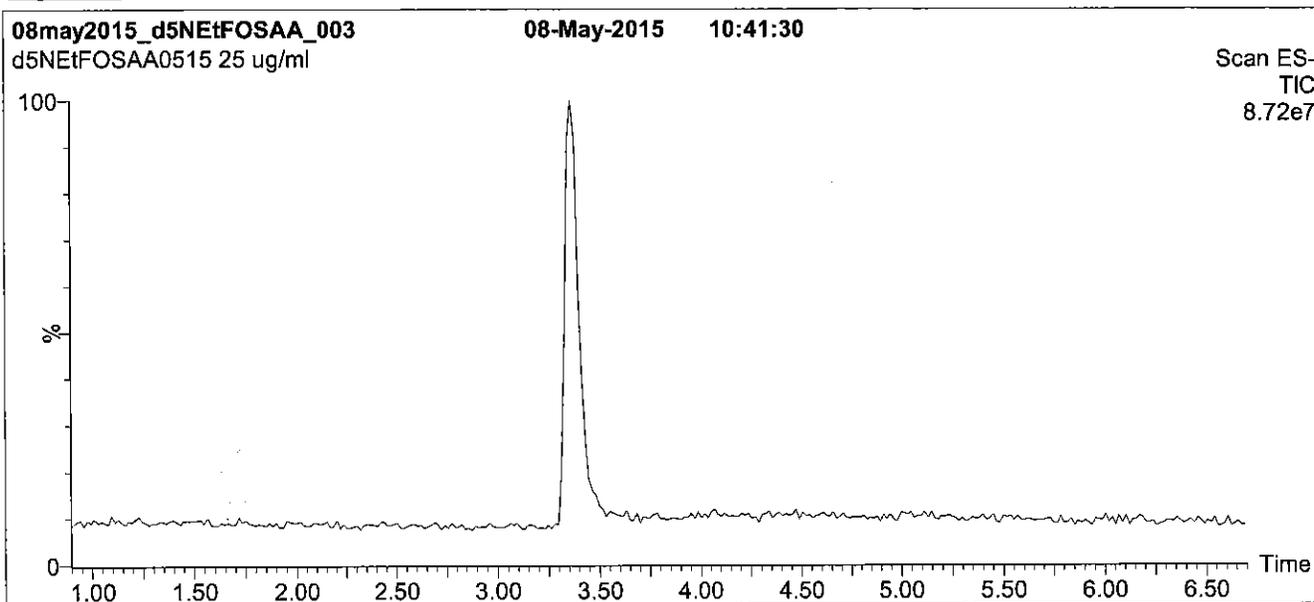
QUALITY MANAGEMENT:

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Figure 1: d5-N-EtFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 65% (80:20 MeOH:ACN) / 35% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

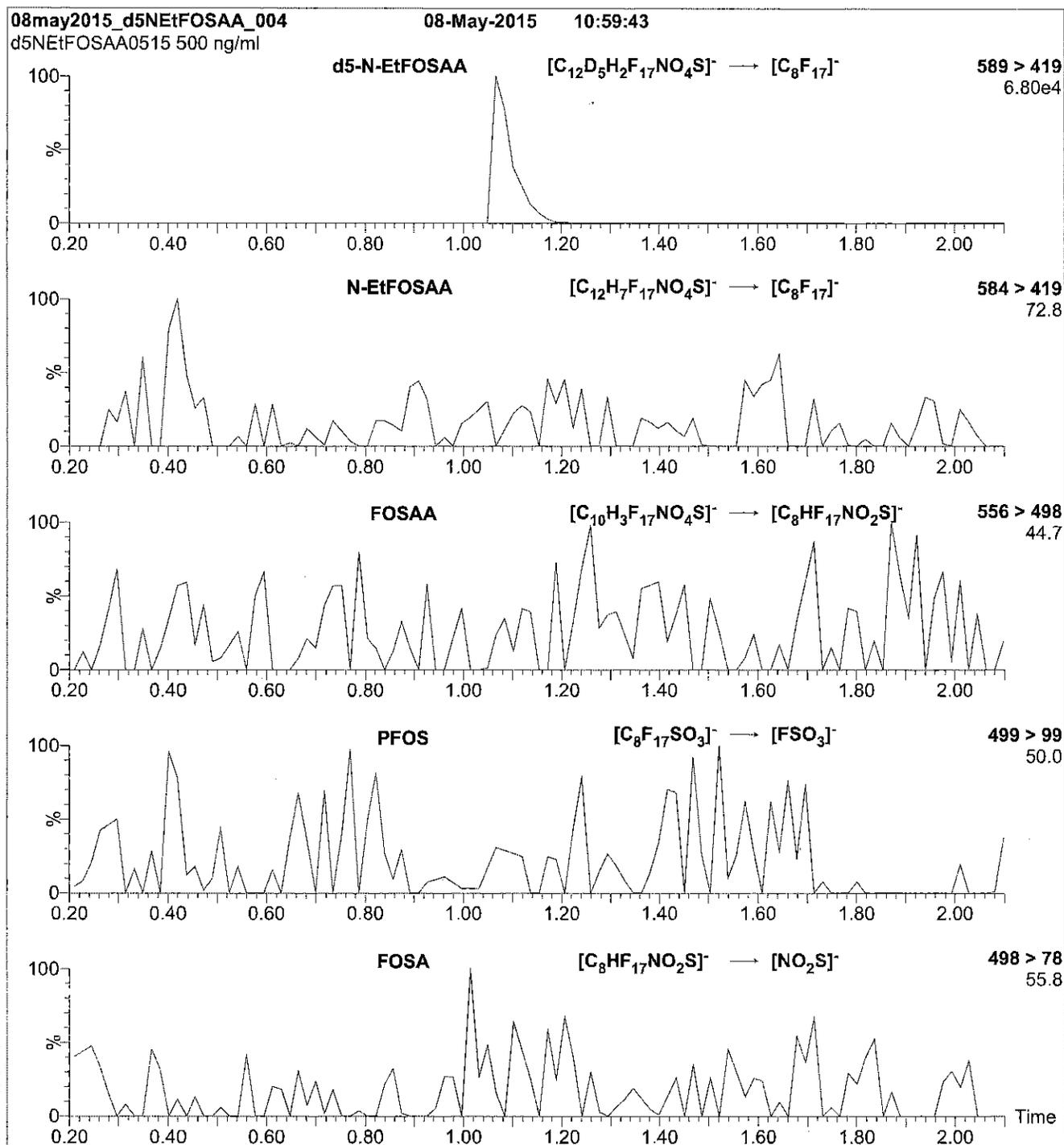
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d5-N-EtFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d5-N-EtFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.24e-3
Collision Energy (eV) = 25

Reagent

LCd5-NEtFOSAA_00002

R: 7/6/16 CBW



671603
ID: LCd5-NEtFOSAA_00002
Exp: 12/07/20 Prep: CBW
d5-N-EtFOSAA

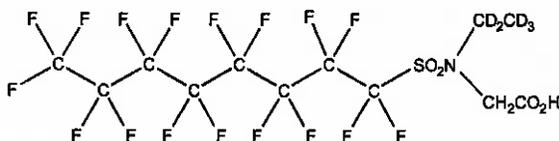


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: d5-N-EtFOSAA **LOT NUMBER:** d5NEtFOSAA1115
COMPOUND: N-ethyl-d5-perfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₁₂D₅H₃F₁₇NO₄S
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 590.27
SOLVENT(S): Methanol
Water (<1%)
ISOTOPIC PURITY: ≥98% ²H₆

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 12/07/2015
EXPIRY DATE: (mm/dd/yyyy) 12/07/2020
RECOMMENDED STORAGE: Refrigerate ampoule

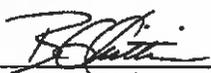
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 12/07/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

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HOMOGENEITY:

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UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

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EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

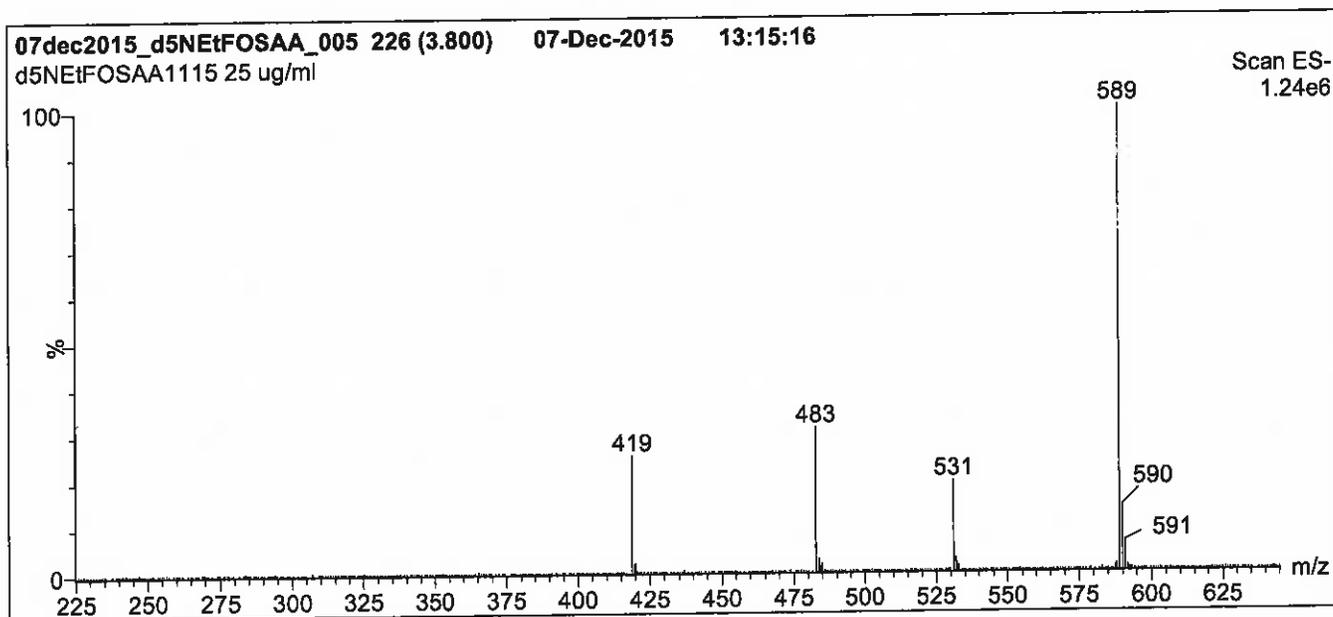
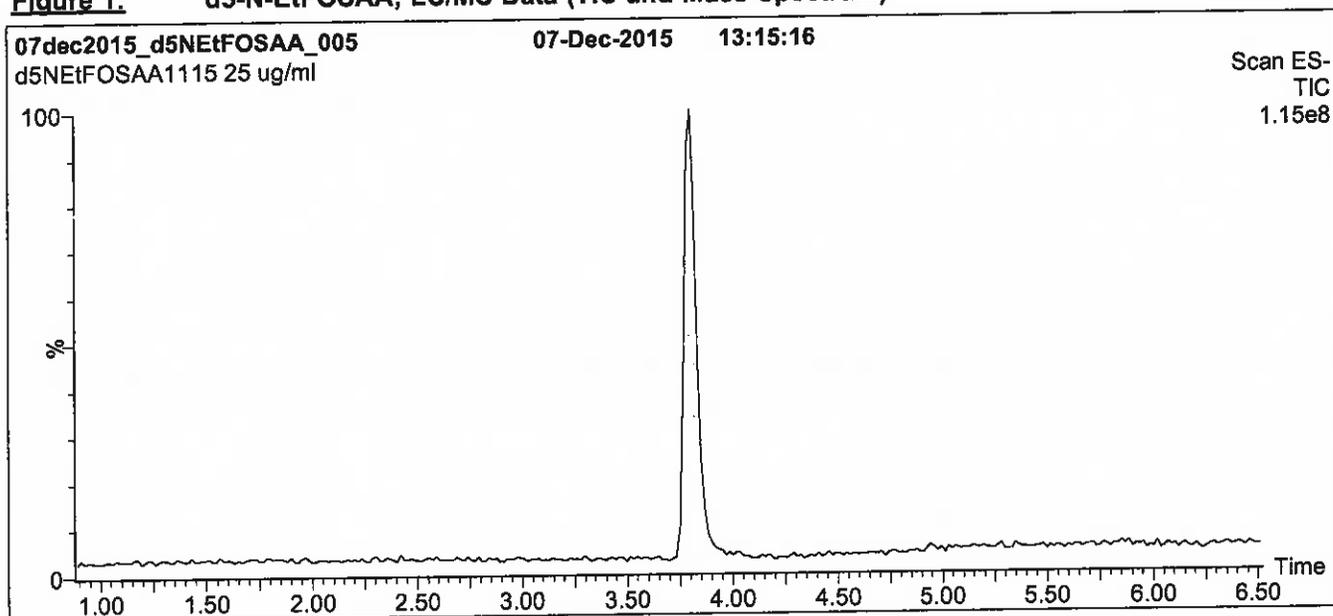
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: d5-N-EtFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 65% (80:20 MeOH:ACN) / 35% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min.
 before returning to initial conditions in 0.5 min.
 Time: 10 min

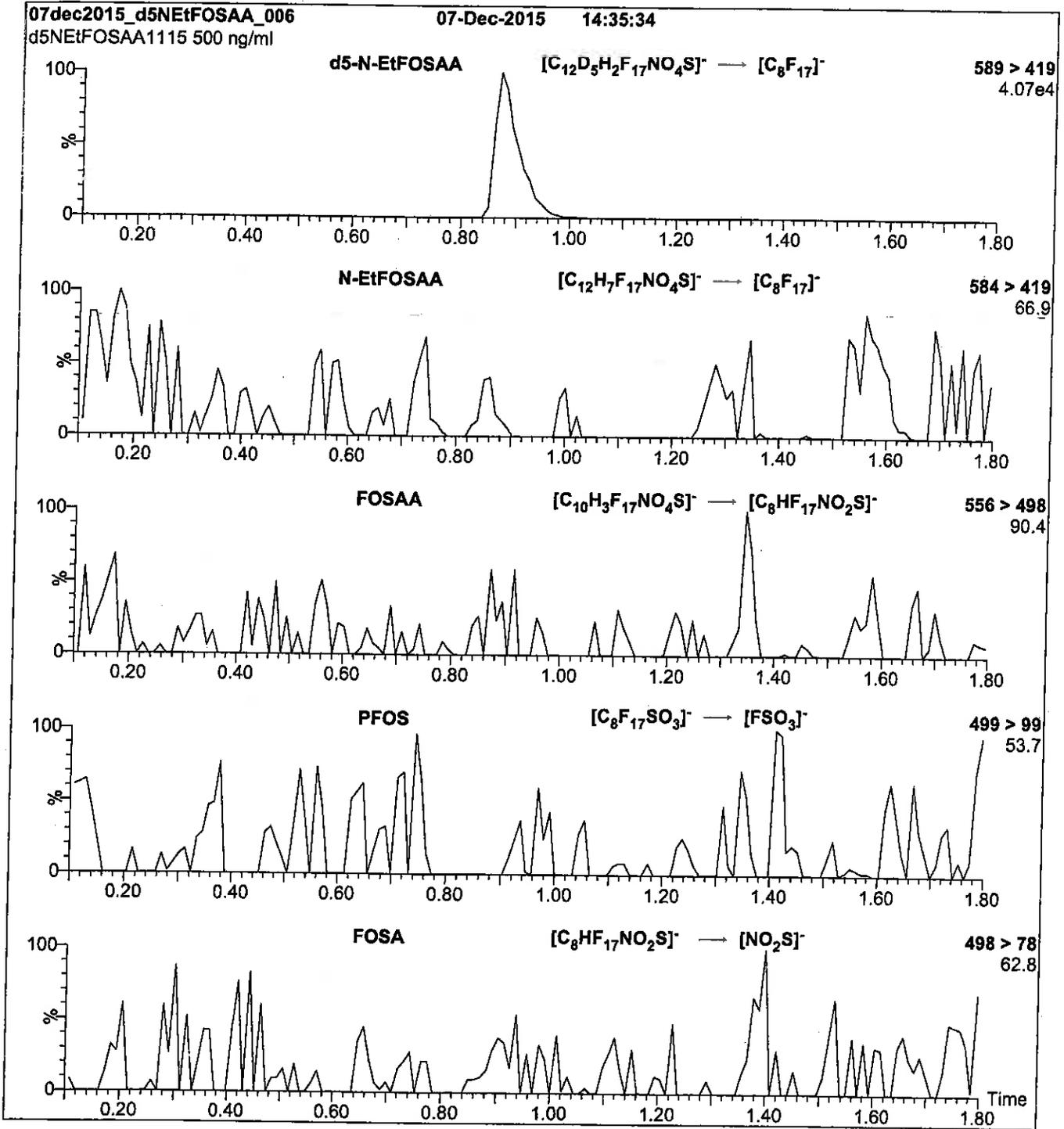
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: d5-N-EtFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml d5-N-EtFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
Collision Energy (eV) = 25

Reagent

LCM2-6:FTS_00001

R: 7/16/15 SW
S: 7/20/15 SW

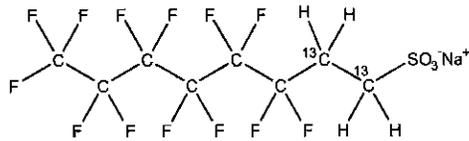


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2-6:2FTS **LOT NUMBER:** M262FTS0714
COMPOUND: Sodium 1H,1H,2H,2H-perfluoro-[1,2-¹³C₂]octane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₆H₄F₁₃SO₃Na **MOLECULAR WEIGHT:** 452.13
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.5 ± 2.4 µg/ml (M2-6:2FTS anion)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 07/15/2014 (1,2-¹³C₂)
EXPIRY DATE: (mm/dd/yyyy) 07/15/2017
RECOMMENDED STORAGE: Refrigerate ampoule

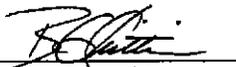
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The native 6:2FTS contains 4.22% of ³⁴S (due to natural isotopic abundance) therefore both native 6:2FTS and M2-6:2FTS will produce signals in the m/z 429 to m/z 409 channel during SRM analysis. We recommend using the m/z 429 to m/z 81 transition to monitor for M2-6:2FTS during quantitative analysis as it will be free of any native contribution (see Figure 2).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 03/27/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

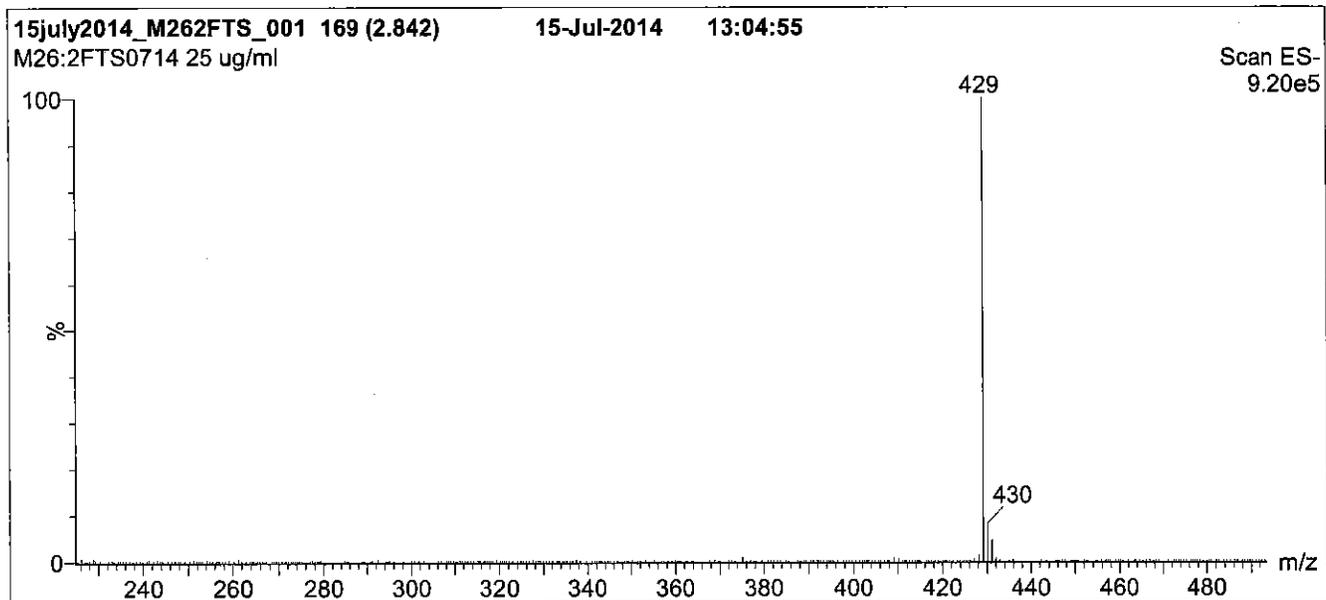
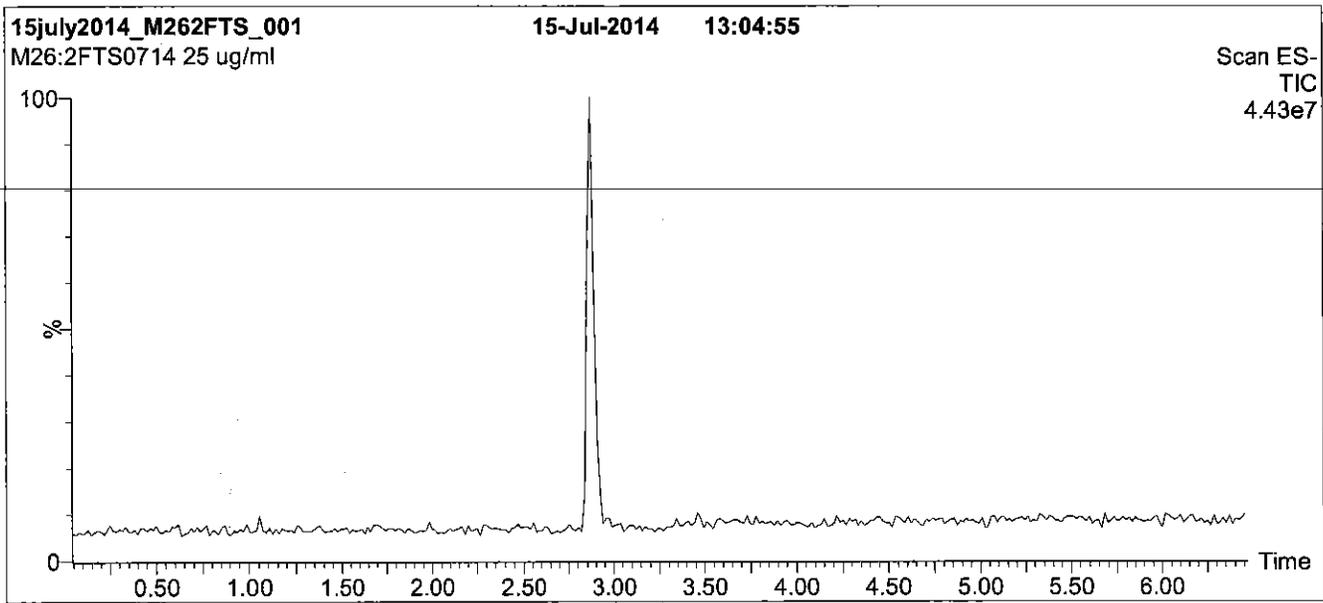
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: M2-6:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 55% (80:20 MeOH:ACN) / 45% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min
and hold for 2 min before returning
to initial conditions in 0.5 min.
Time: 10 min

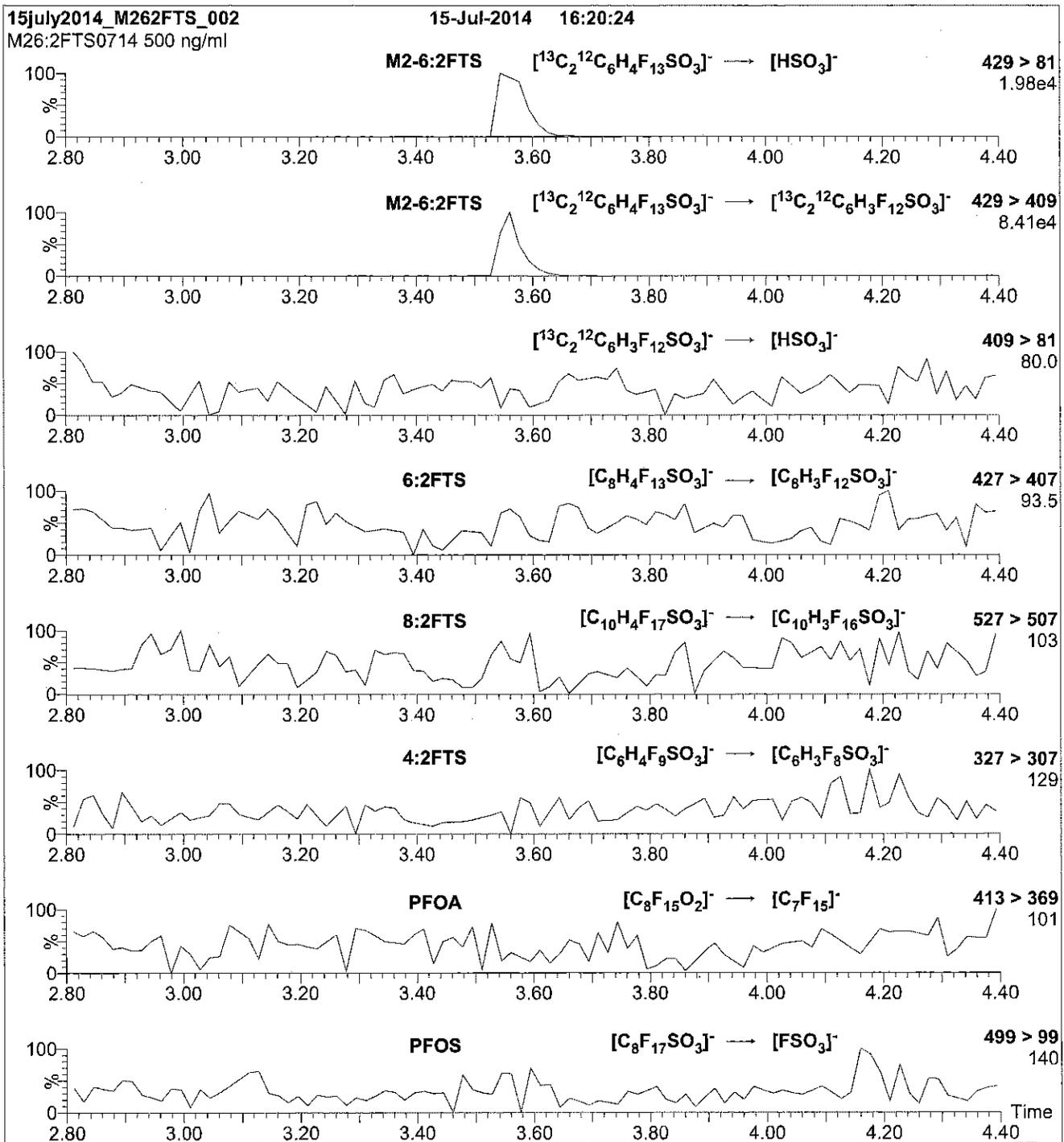
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 950 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2-6:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2-6:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.43e-3
Collision Energy (eV) = 25

Reagent

LCM2-6:FTS_00002

R: 7/6/16 CSW

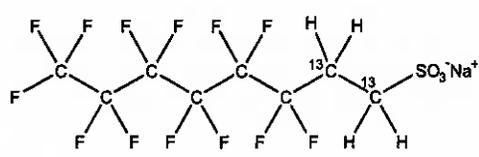
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Exp: 01/08/21 Prod: CSW
M2-6:2F2S

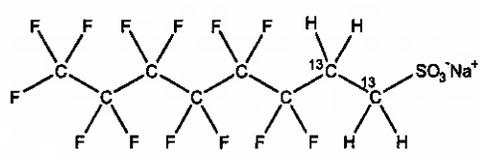


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2-6:2F2S **LOT NUMBER:** M262F2S0116
COMPOUND: Sodium 1H,1H,2H,2H-perfluoro-[1,2-¹³C₂]octane sulfonate

STRUCTURE:  **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₆H₄F₁₃SO₃Na **MOLECULAR WEIGHT:** 452.13
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
 47.5 ± 2.4 µg/ml (M2-6:2F2S anion)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 01/08/2016 (1,2-¹³C₂)
EXPIRY DATE: (mm/dd/yyyy) 01/08/2021
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The native 6:2F2S contains 4.22% of ³⁴S (due to natural isotopic abundance) therefore both native 6:2F2S and M2-6:2F2S will produce signals in the m/z 429 to m/z 409 channel during SRM analysis. We recommend using the m/z 429 to m/z 81 transition to monitor for M2-6:2F2S during quantitative analysis as it will be free of any native contribution (see Figure 2).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 01/11/2016
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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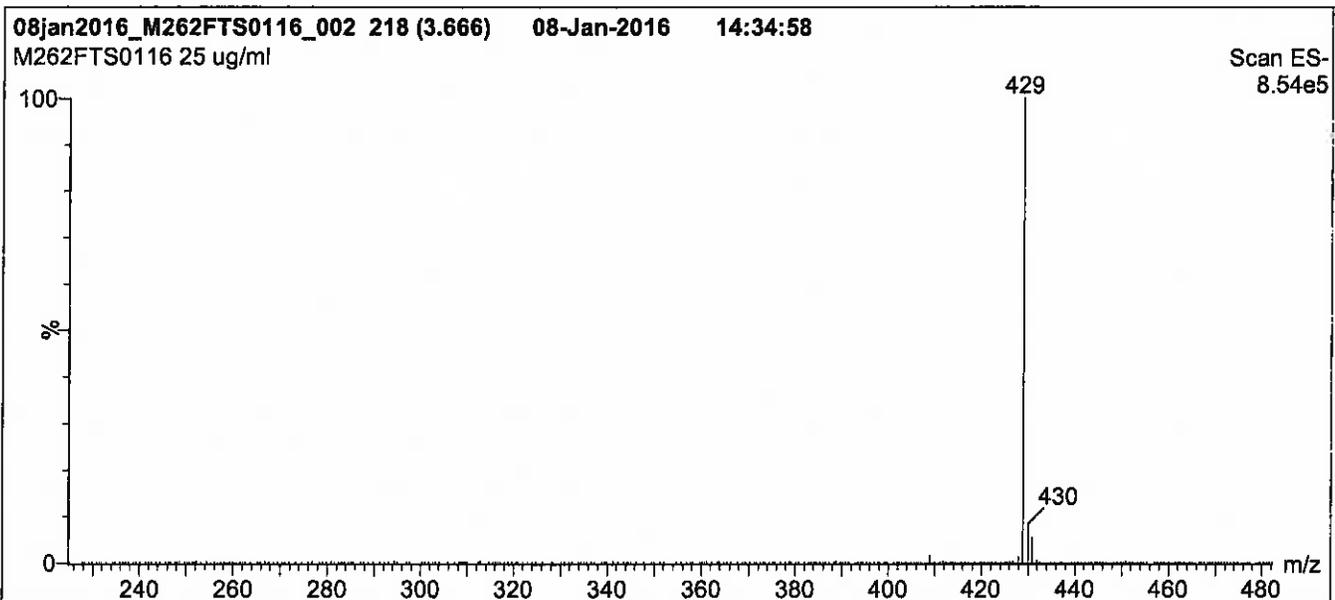
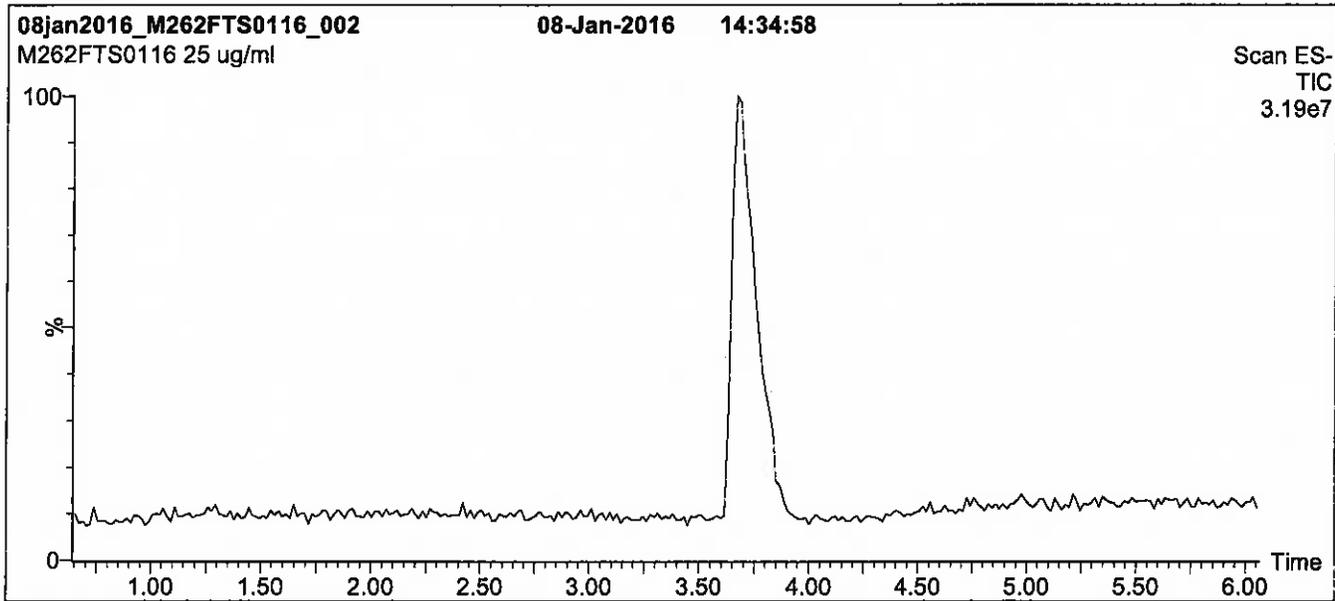
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Figure 1: M2-6:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min
and hold for 2 min before returning
to initial conditions in 0.5 min.
Time: 10 min

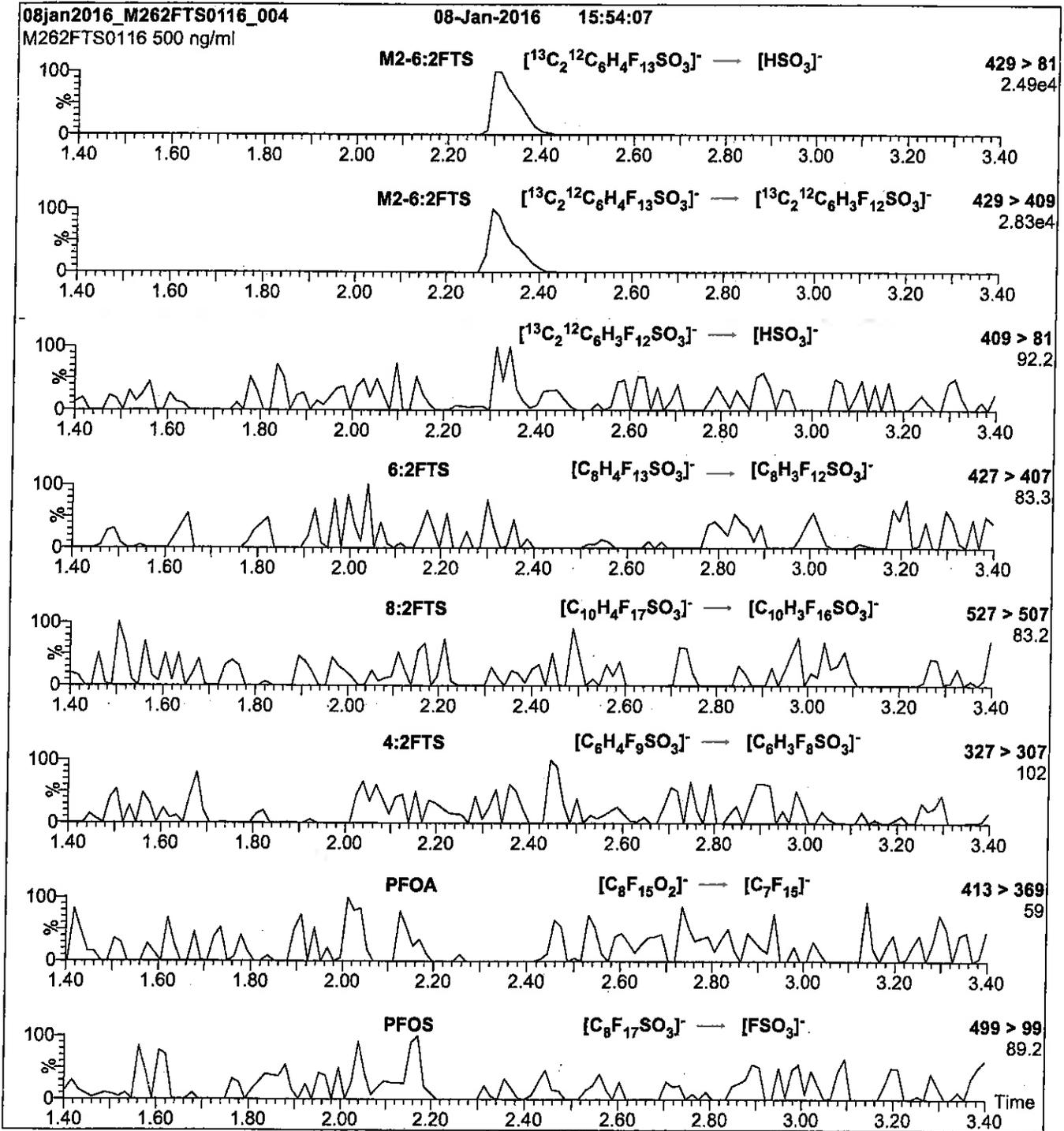
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2-6:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2-6:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.28e-3
Collision Energy (eV) = 25

Reagent

LCM2-8:2FTS_00001

r: 7/16/15 ✓
s: 7/22/15 STV

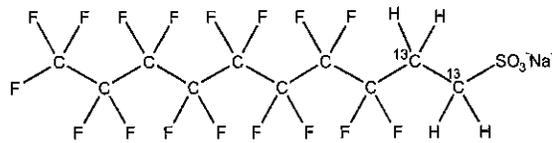


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2-8:2FTS **LOT NUMBER:** M282FTS0414
COMPOUND: Sodium 1H,1H,2H,2H-perfluoro-[1,2-¹³C₂]decane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₈H₄F₁₇SO₃Na **MOLECULAR WEIGHT:** 552.15
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.9 ± 2.4 µg/ml (M2-8:2FTS anion)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 04/13/2014 (1,2-¹³C₂)
EXPIRY DATE: (mm/dd/yyyy) 04/13/2017
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The native 8:2FTS contains 4.22% of ³⁴S (due to natural isotopic abundance) therefore both native 8:2FTS and M2-8:2FTS will produce signals in the m/z 529 to m/z 509 channel during SRM analysis. We recommend using the m/z 529 to m/z 81 transition to monitor for M2-8:2FTS during quantitative analysis as it will be free of any native contribution (see Figure 2).

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The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

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Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

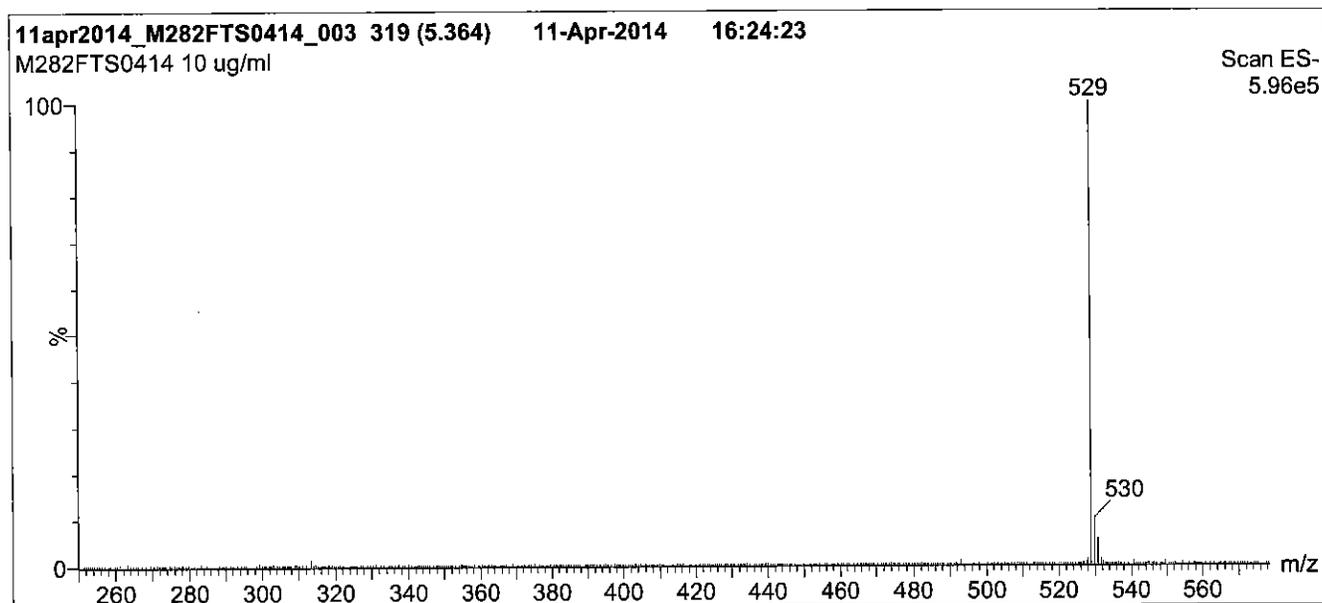
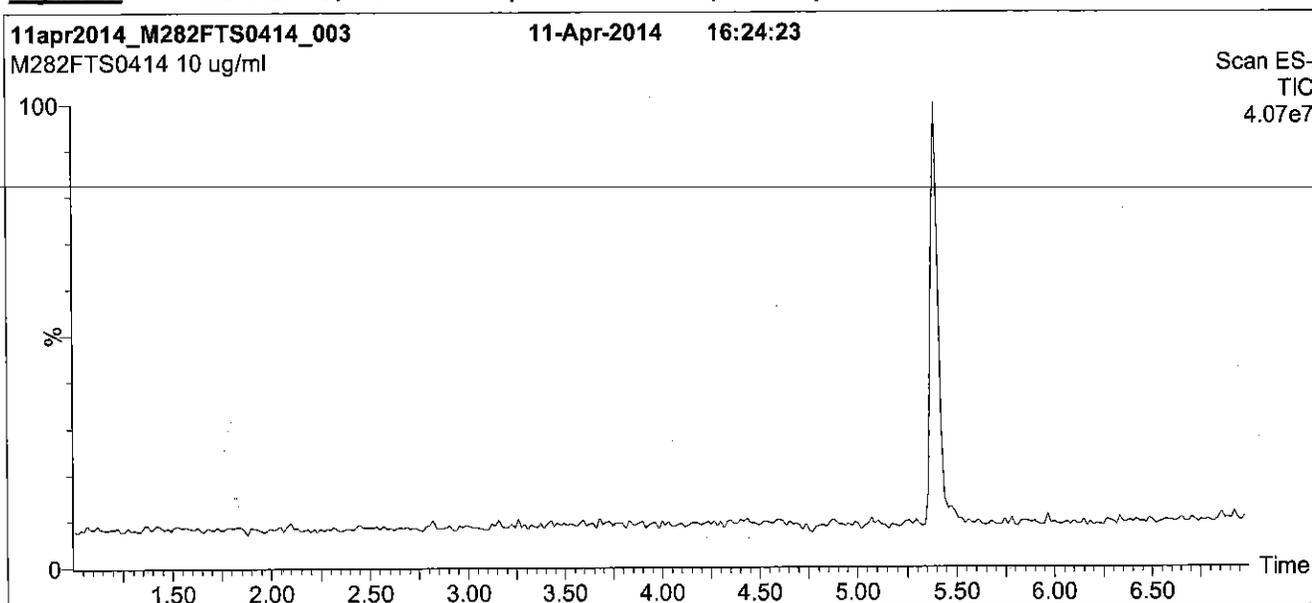
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: M2-8:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min
and hold for 2 min before returning
to initial conditions in 0.5 min.
Time: 10 min

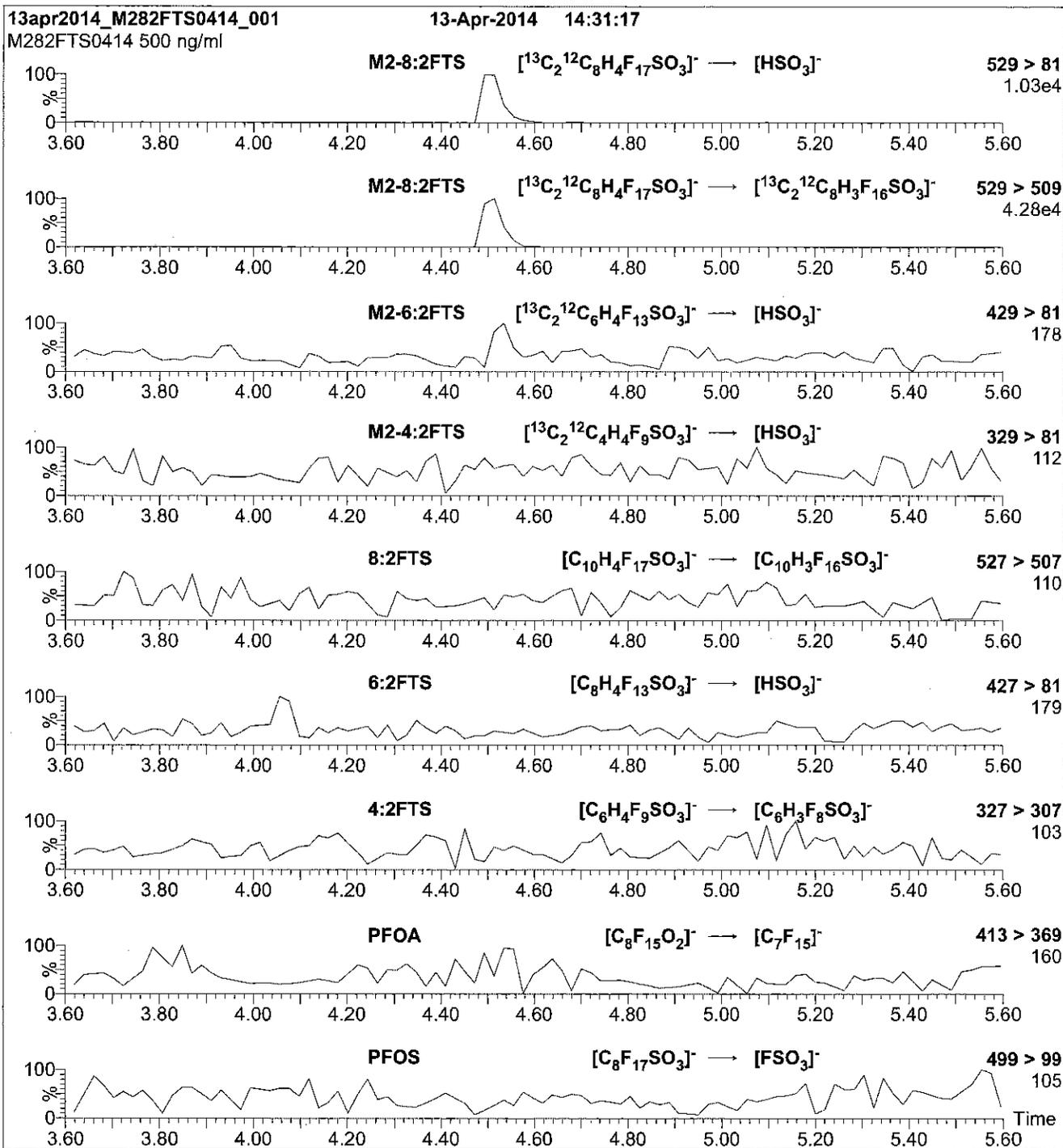
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2-8:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μl (500 ng/ml M2-8:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
 (both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

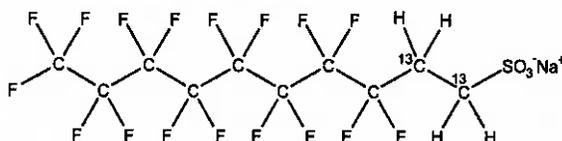
MS Parameters

Collision Gas (mbar) = 3.24e-3
 Collision Energy (eV) = 25

Reagent

LCM2-8:2FTS_00002

R: 7/6/16 CBW

671602
ID: LCM2-8:2FTS_00002
Exp: 01/08/21 Prod: CBW
M2-8:2FTS**WELLINGTON**
LABORATORIES**CERTIFICATE OF ANALYSIS**
DOCUMENTATION**PRODUCT CODE:** M2-8:2FTS **LOT NUMBER:** M282FTS0116
COMPOUND: Sodium 1H,1H,2H,2H-perfluoro-[1,2-¹³C₂]decane sulfonate**STRUCTURE:** **CAS #:** Not available

MOLECULAR FORMULA:	¹³ C ₂ ¹² C ₈ H ₄ F ₁₇ SO ₃ Na	MOLECULAR WEIGHT:	552.15
CONCENTRATION:	50.0 ± 2.5 µg/ml (Na salt)	SOLVENT(S):	Methanol
	47.9 ± 2.4 µg/ml (M2-8:2FTS anion)	ISOTOPIC PURITY:	≥99% ¹³ C
CHEMICAL PURITY:	>98%		(1,2- ¹³ C ₂)
LAST TESTED: (mm/dd/yyyy)	01/08/2016		
EXPIRY DATE: (mm/dd/yyyy)	01/08/2021		
RECOMMENDED STORAGE:	Refrigerate ampoule		

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The native 8:2FTS contains 4.22% of ³⁴S (due to natural isotopic abundance) therefore both native 8:2FTS and M2-8:2FTS will produce signals in the m/z 529 to m/z 509 channel during SRM analysis. We recommend using the m/z 529 to m/z 81 transition to monitor for M2-8:2FTS during quantitative analysis as it will be free of any native contribution (see Figure 2).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim
Date: 01/18/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

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HOMOGENEITY:

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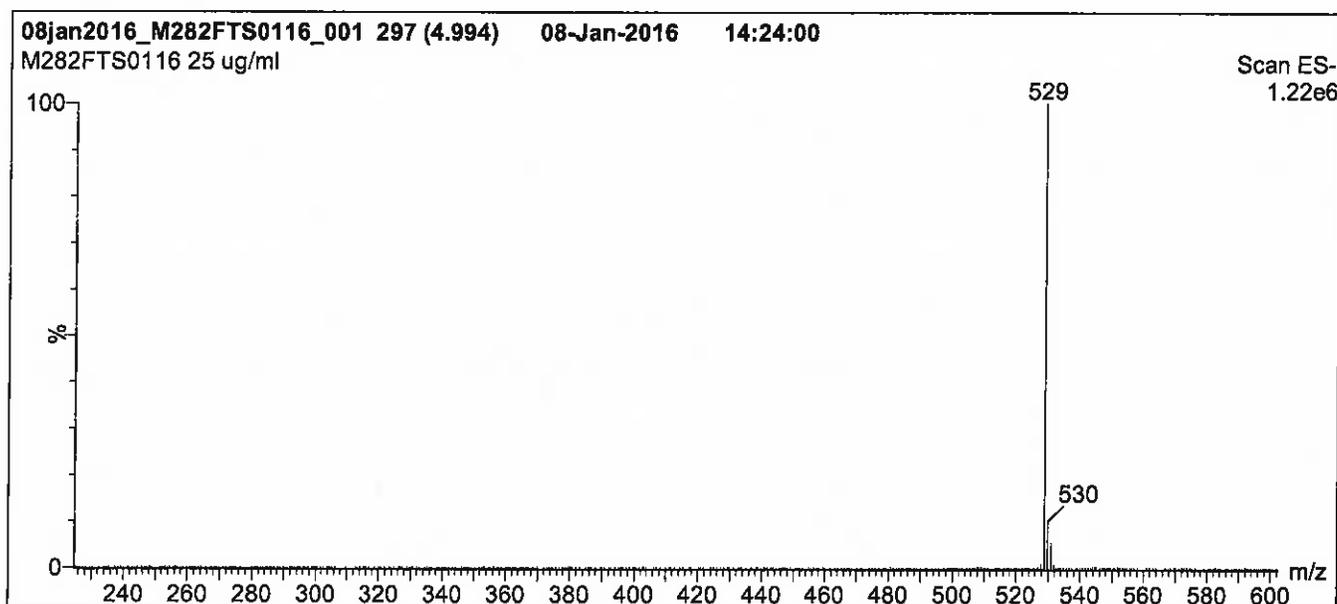
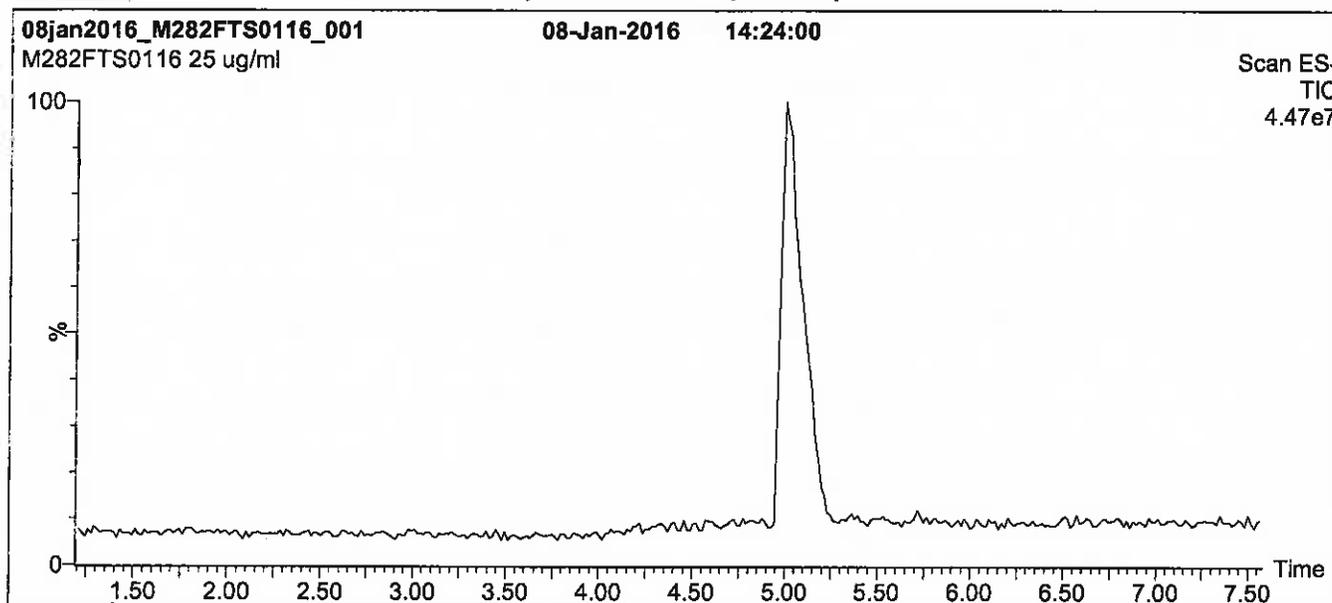
QUALITY MANAGEMENT:

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Figure 1: M2-8:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min
and hold for 2 min before returning
to initial conditions in 0.5 min.
Time: 10 min

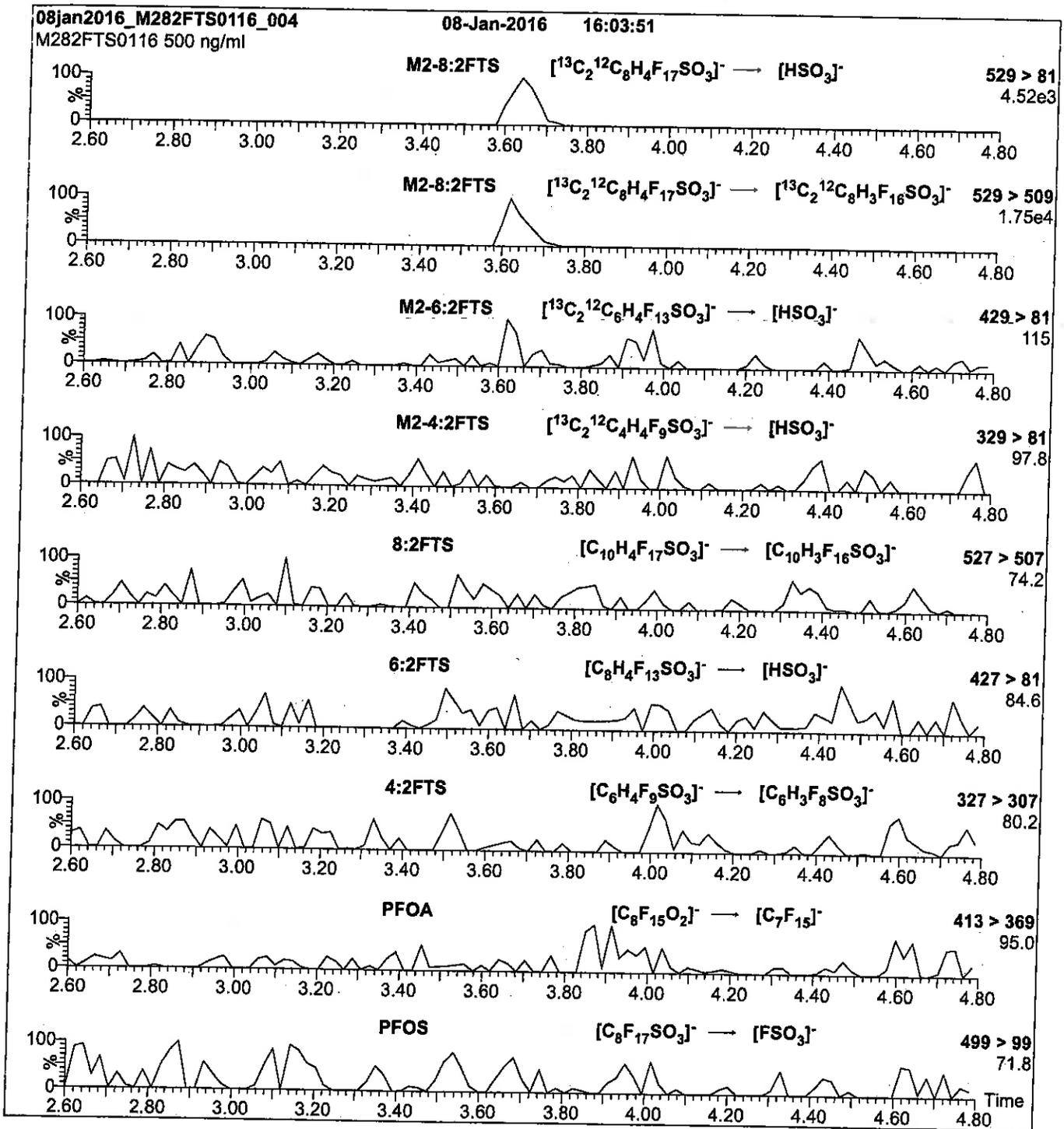
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2-8:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2-8:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = $3.20\text{e-}3$
Collision Energy (eV) = 30

Reagent

LCM2PFHxDA_00008

R: SBC 9/22/16

739512
ID: LCM2PFHxDA_00008
Exp: 01/07/21 Prod: SBC
13C2-PFHxDA at 50ug/mL

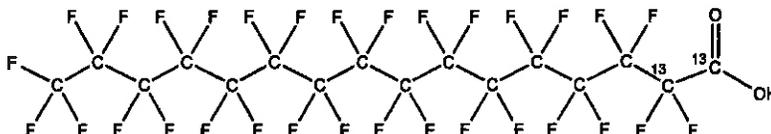


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2PFHxDA **LOT NUMBER:** M2PFHxDA1112
COMPOUND: Perfluoro-n-[1,2-¹³C₂]hexadecanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA:	¹³ C ₂ ¹² C ₁₄ HF ₃₁ O ₂	MOLECULAR WEIGHT:	816.11
CONCENTRATION:	50 ± 2.5 µg/ml	SOLVENT(S):	Methanol Water (<1%)
CHEMICAL PURITY:	>98%	ISOTOPIC PURITY:	≥99% ¹³ C (1,2- ¹³ C ₂)
LAST TESTED: (mm/dd/yyyy)	01/07/2016		
EXPIRY DATE: (mm/dd/yyyy)	01/07/2021		
RECOMMENDED STORAGE:	Store ampoule in a cool, dark place		

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.3% of native perfluoro-n-hexadecanoic acid.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 01/11/2016
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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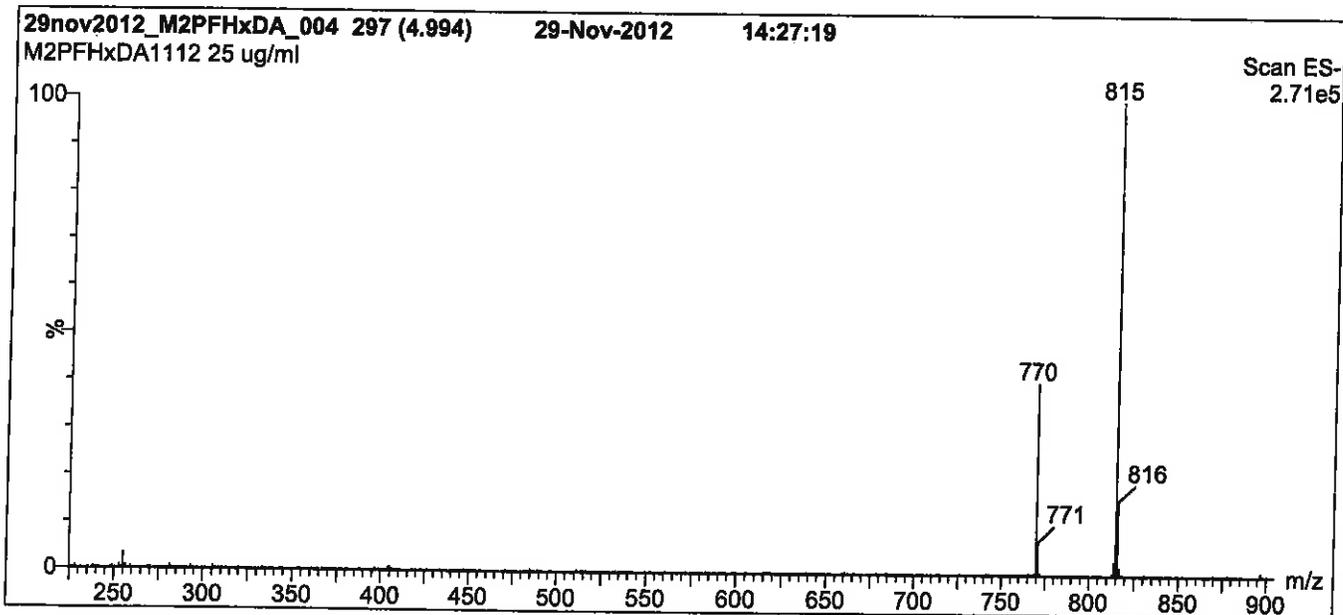
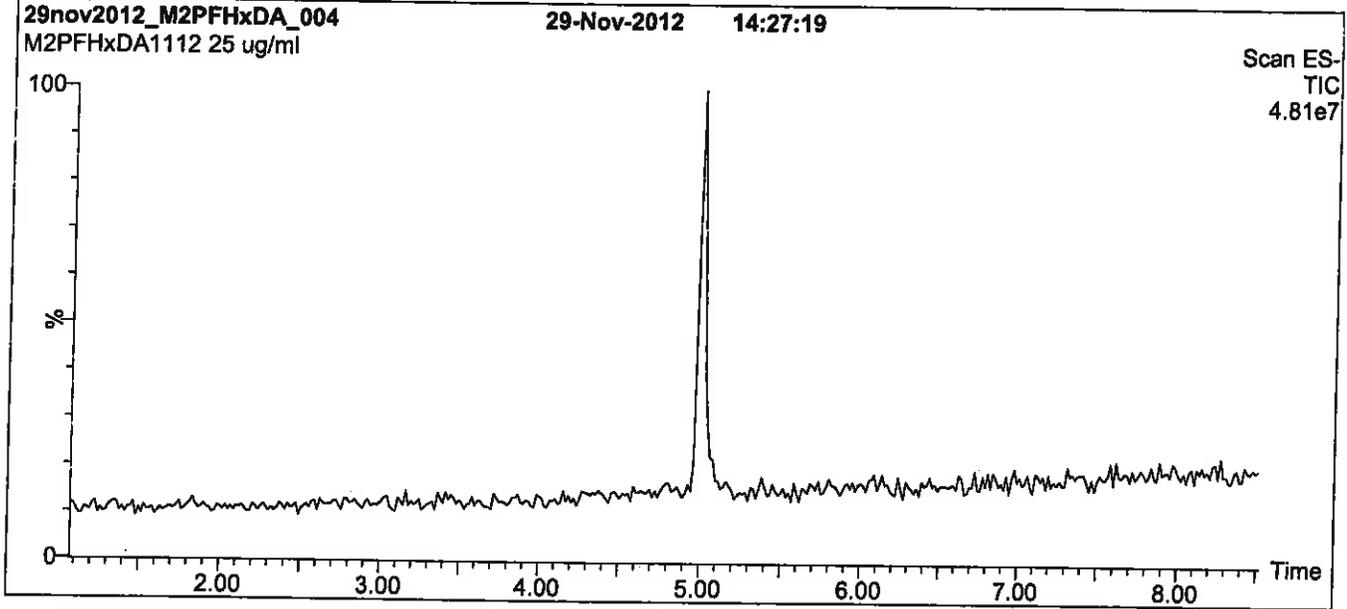
QUALITY MANAGEMENT:

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Figure 1: M2PFHxDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 60% (80:20 MeOH:ACN) / 40% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 100% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

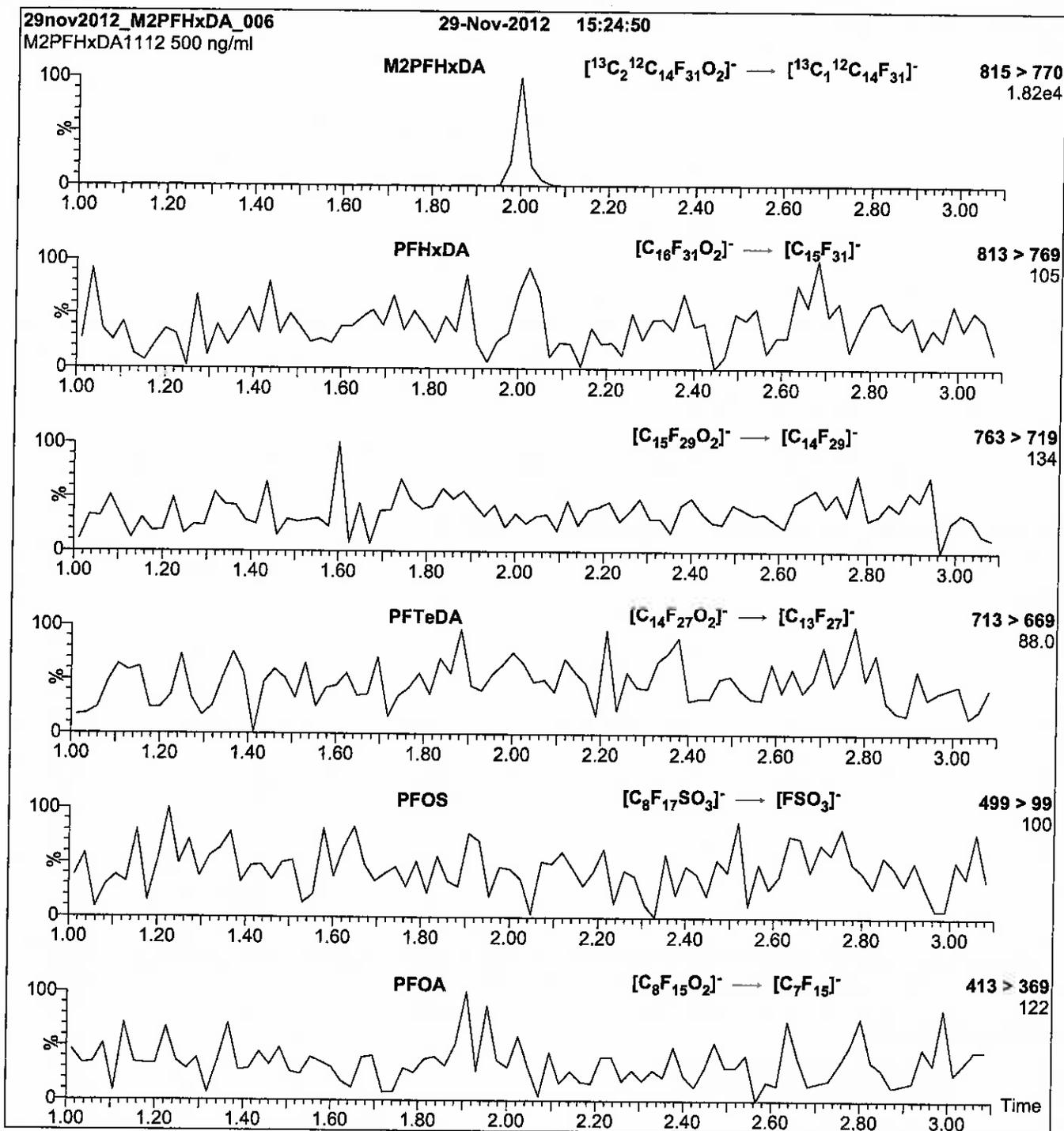
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 1200 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 25.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2PFHxDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2PFHxDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.39e-3
Collision Energy (eV) = 15

Reagent

LCM2PFTeDA_00007

Scanned 10/14/16 R: Soc 9/22/16

739563
ID: LCM2PFTeDA_00007
Exp: 12/07/20 Pppl: SBC
13C2-PFTeDA at 50ug/mL

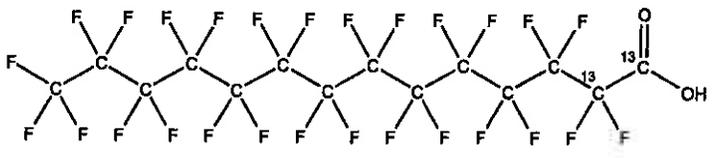


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2PFTeDA **LOT NUMBER:** M2PFTeDA1115
COMPOUND: Perfluoro-n-[1,2-¹³C₂]tetradecanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₁₂HF₂₇O₂ **MOLECULAR WEIGHT:** 716.10
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 12/07/2015 **Water (<1%)**
EXPIRY DATE: (mm/dd/yyyy) 12/07/2020 **(1,2-¹³C₂)**
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: B.G. Chittim Date: 12/08/2015
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • Info@well-labs.com

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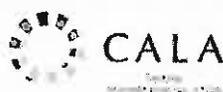
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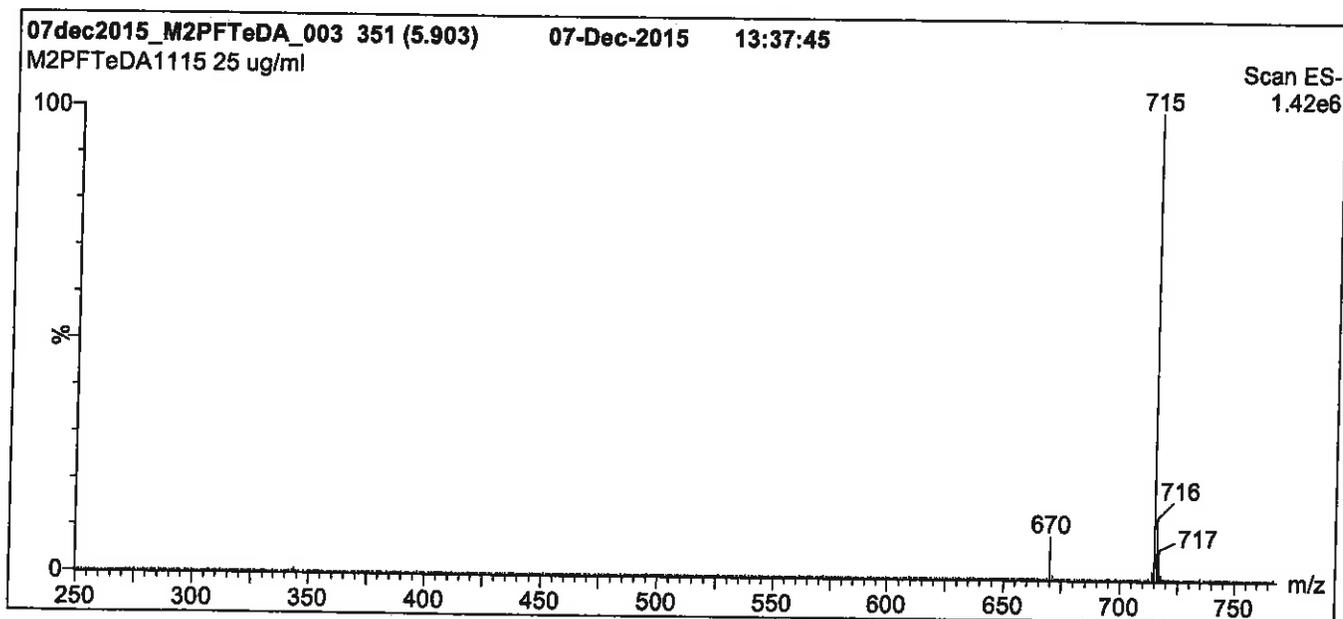
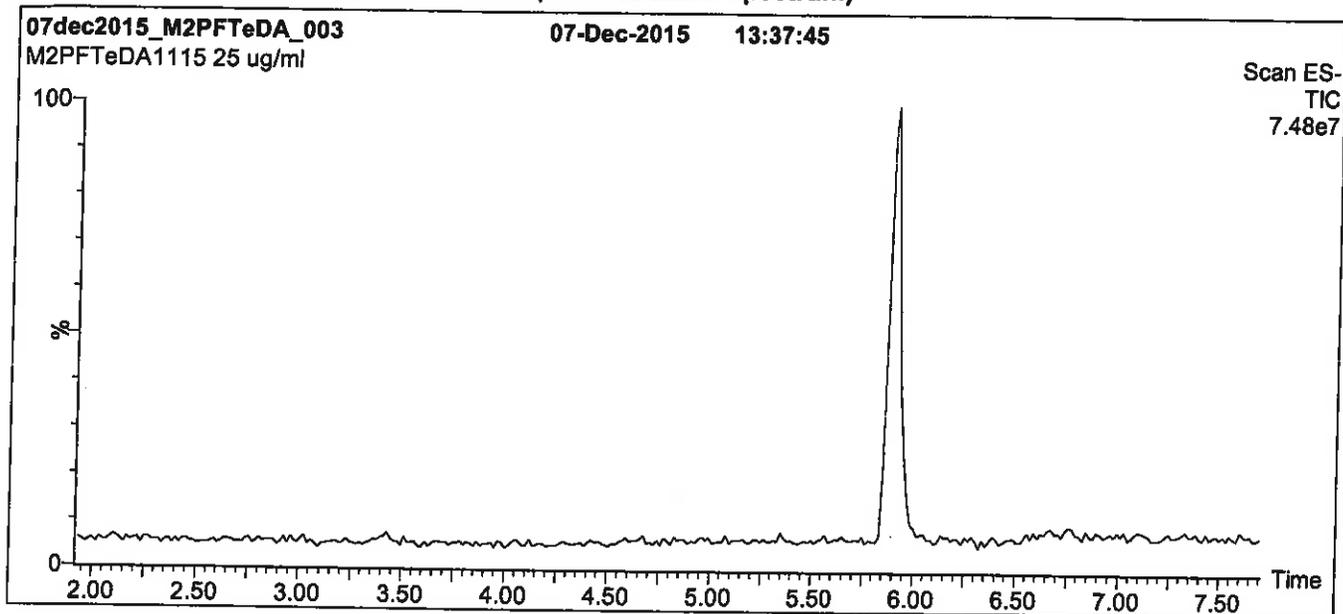
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Figure 1: M2PFTeDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 65% (80:20 MeOH:ACN) / 35% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

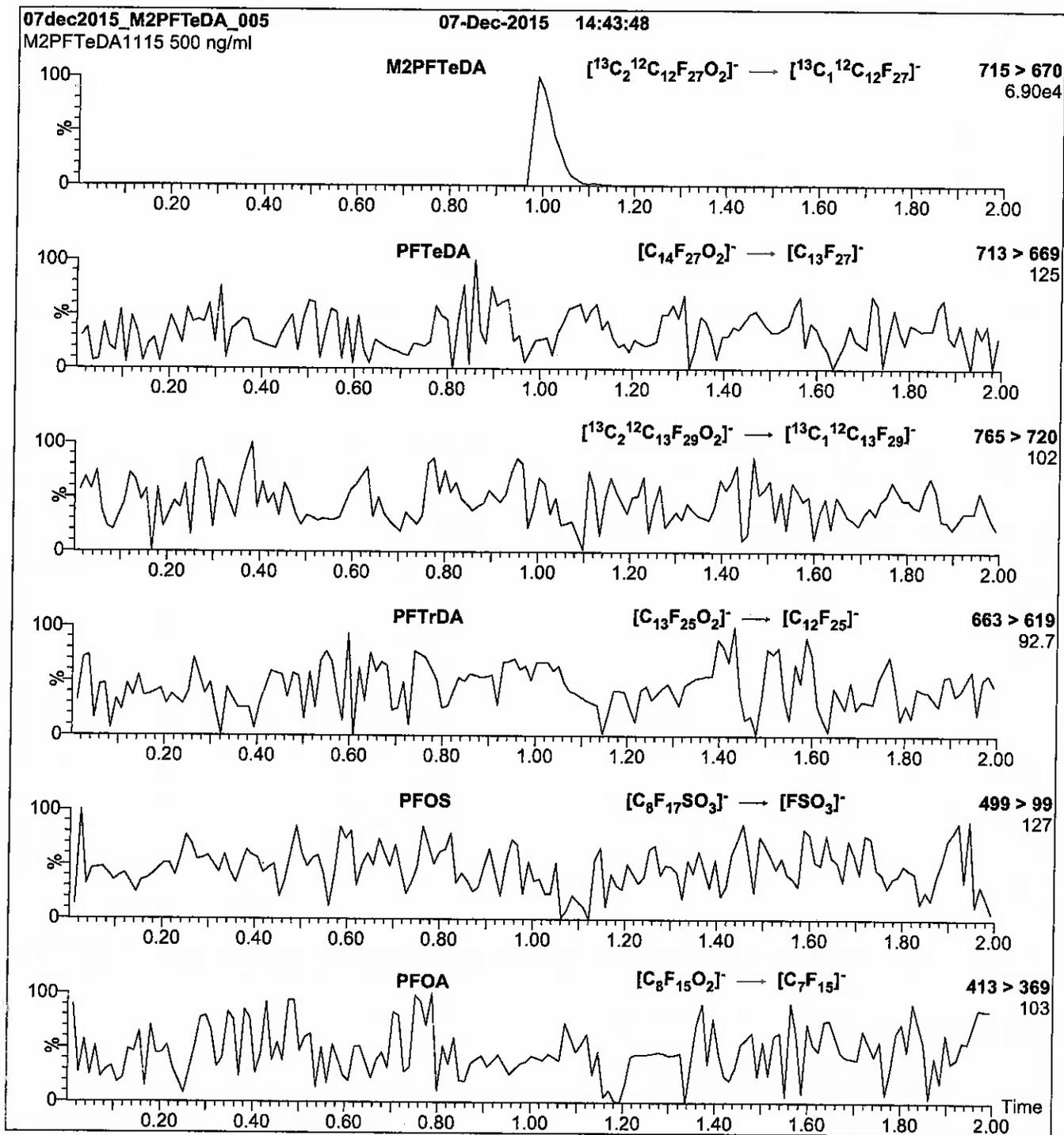
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 1250 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2PFTeDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2PFTeDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.28e-3
Collision Energy (eV) = 14

Reagent

LCM4PFHPA_00007

f: SBC a/22/16

739567
ID: LCM4PFHPA_00007
Exp: 05/27/21 Prpd: SBC
13C4-Perfluoroheptanoic a



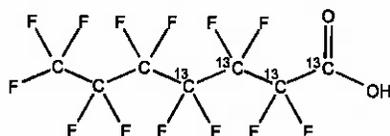
WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

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PRODUCT CODE: M4PFHpA **LOT NUMBER:** M4PFHpA0516
COMPOUND: Perfluoro-n-[1,2,3,4-¹³C₄]heptanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₄¹²C₃HF₁₃O₂ **MOLECULAR WEIGHT:** 368.03
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99%¹³C
(1,2,3,4-¹³C₄)
LAST TESTED: (mm/dd/yyyy) 05/27/2016
EXPIRY DATE: (mm/dd/yyyy) 05/27/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

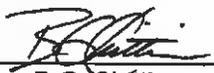
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 07/05/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

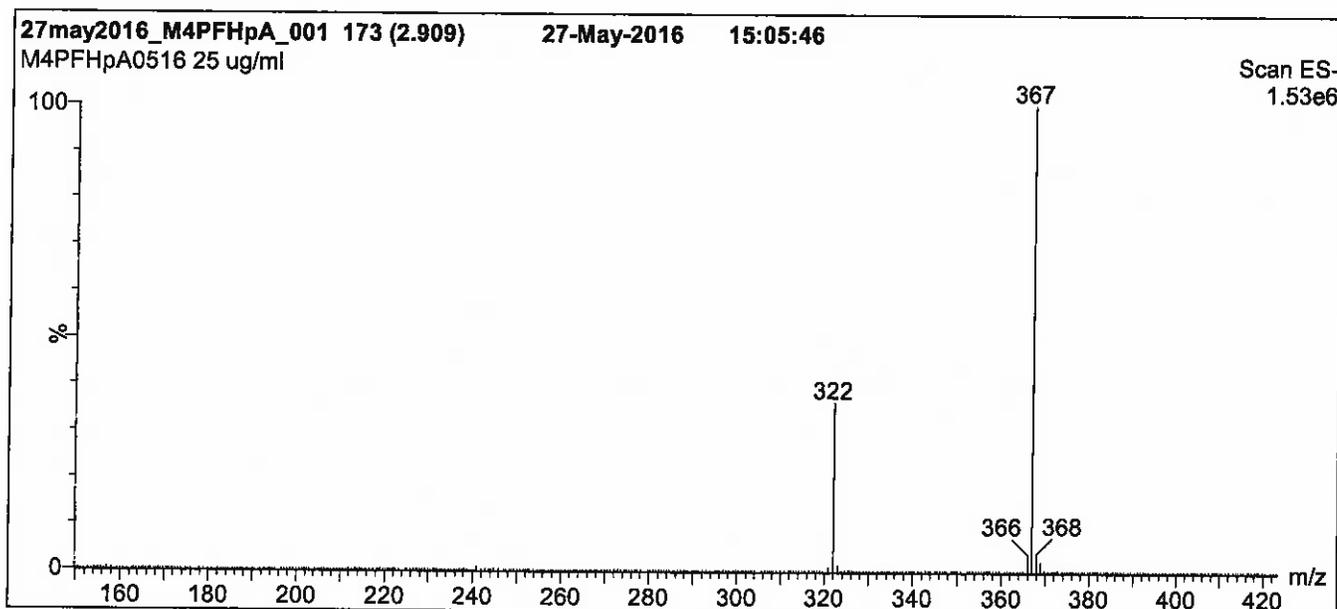
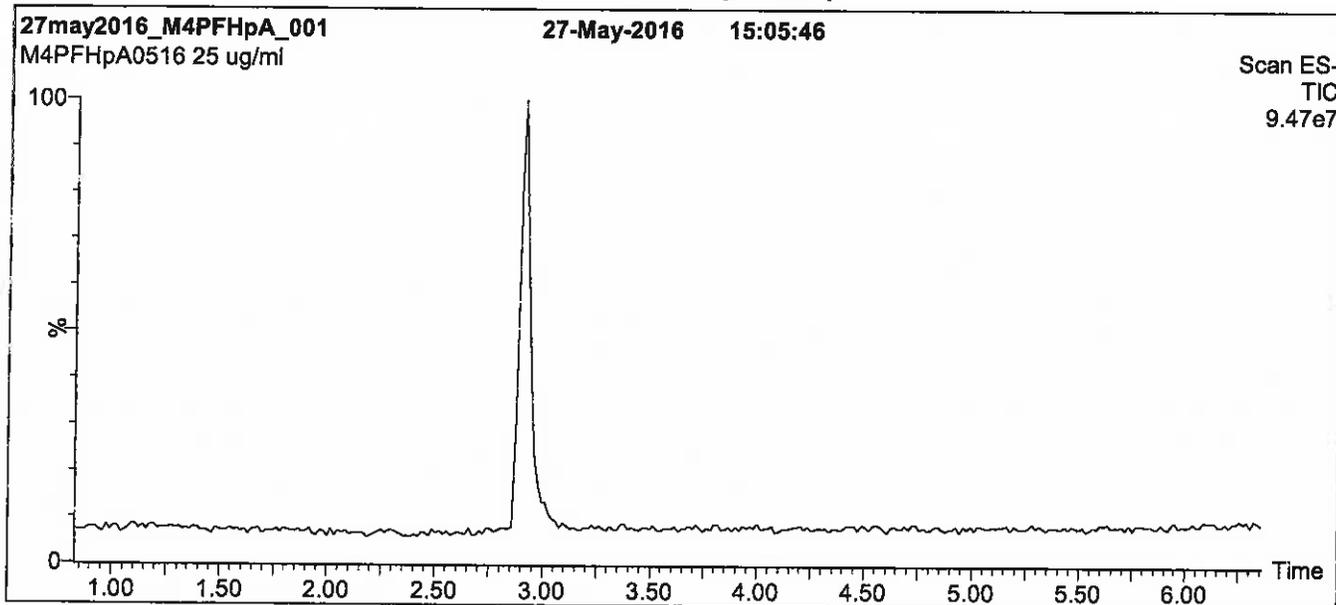
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: M4PFHpA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7.5 min and hold for 1.5 min before returning to initial conditions in 0.5 min.
Time: 10 min

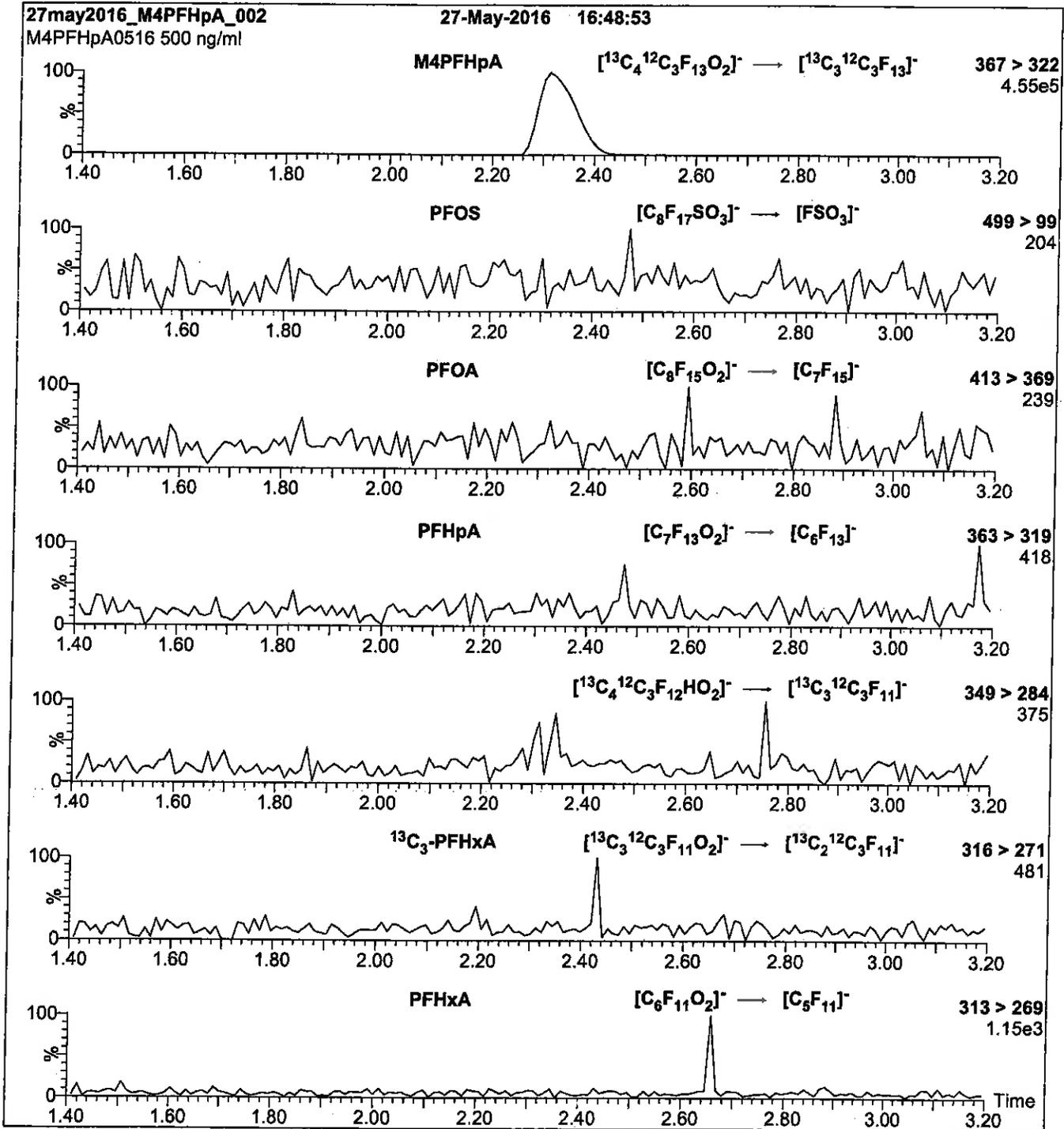
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: M4PFHpA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M4PFHpA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.35e-3
Collision Energy (eV) = 11

Reagent

LCM5PFPEA_00008

R: 8BC 9/22/16



739590
ID: LCM5PFPEA_00008
Exp: 05/22/20 Prpt: SAC
13C5-Perfluoropentanoic a

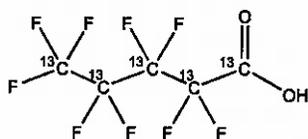


WELLINGTON
LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

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PRODUCT CODE: M5PFPeA **LOT NUMBER:** M5PFPeA0515
COMPOUND: Perfluoro-n-[¹³C₅]pentanoic acid
STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₅HF₉O₂ **MOLECULAR WEIGHT:** 269.01
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 05/22/2015 (¹³C₅)
EXPIRY DATE: (mm/dd/yyyy) 05/22/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains < 0.1% of perfluoro-n-pentanoic acid.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: *B.G. Chittim* **Date:** 05/25/2015
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

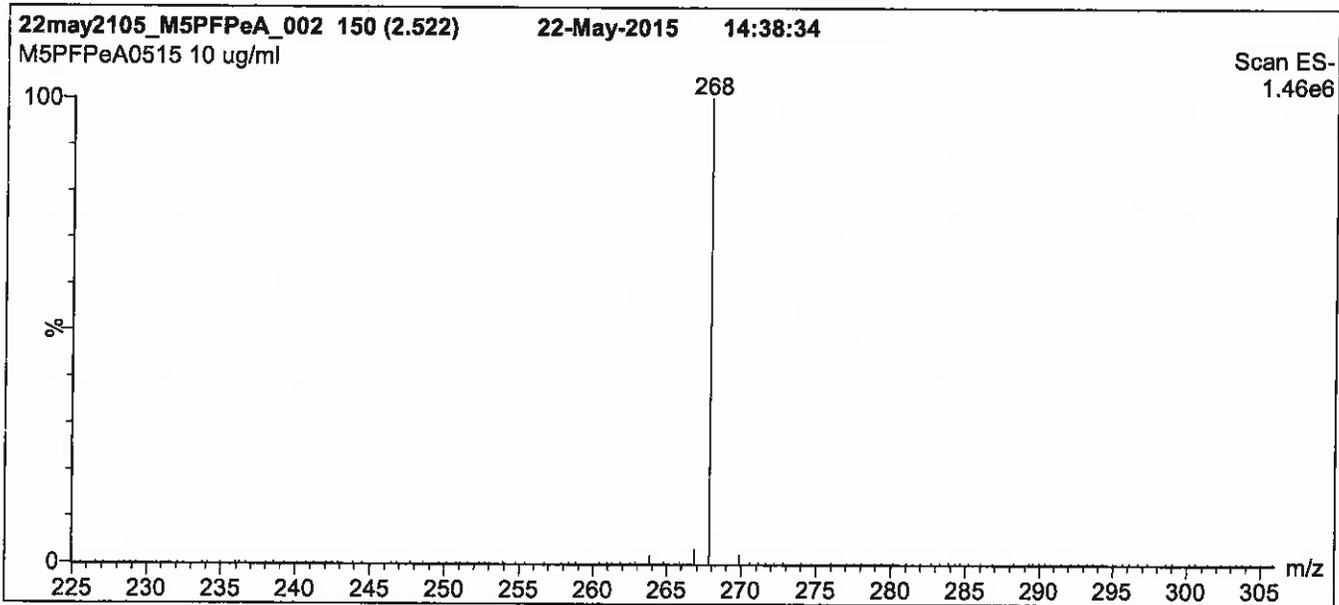
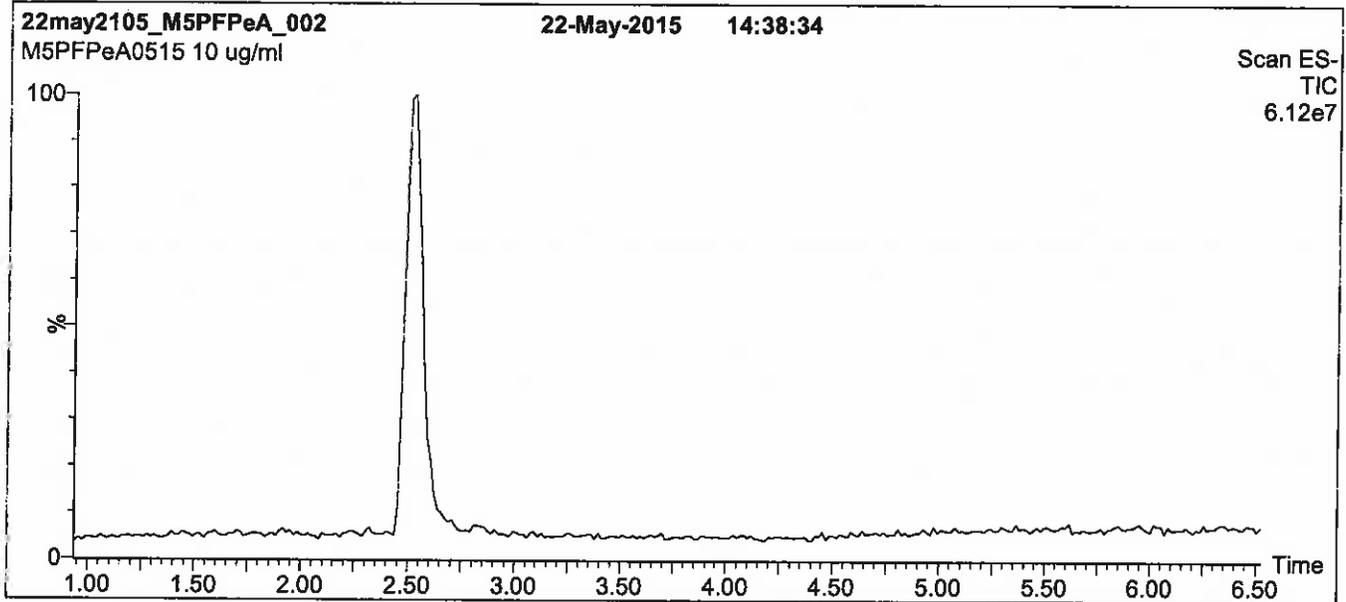
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: M5PFPeA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 40% (80:20 MeOH:ACN) / 60% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for
1.5 min before returning to initial conditions in 0.5 min.
Time: 10 min

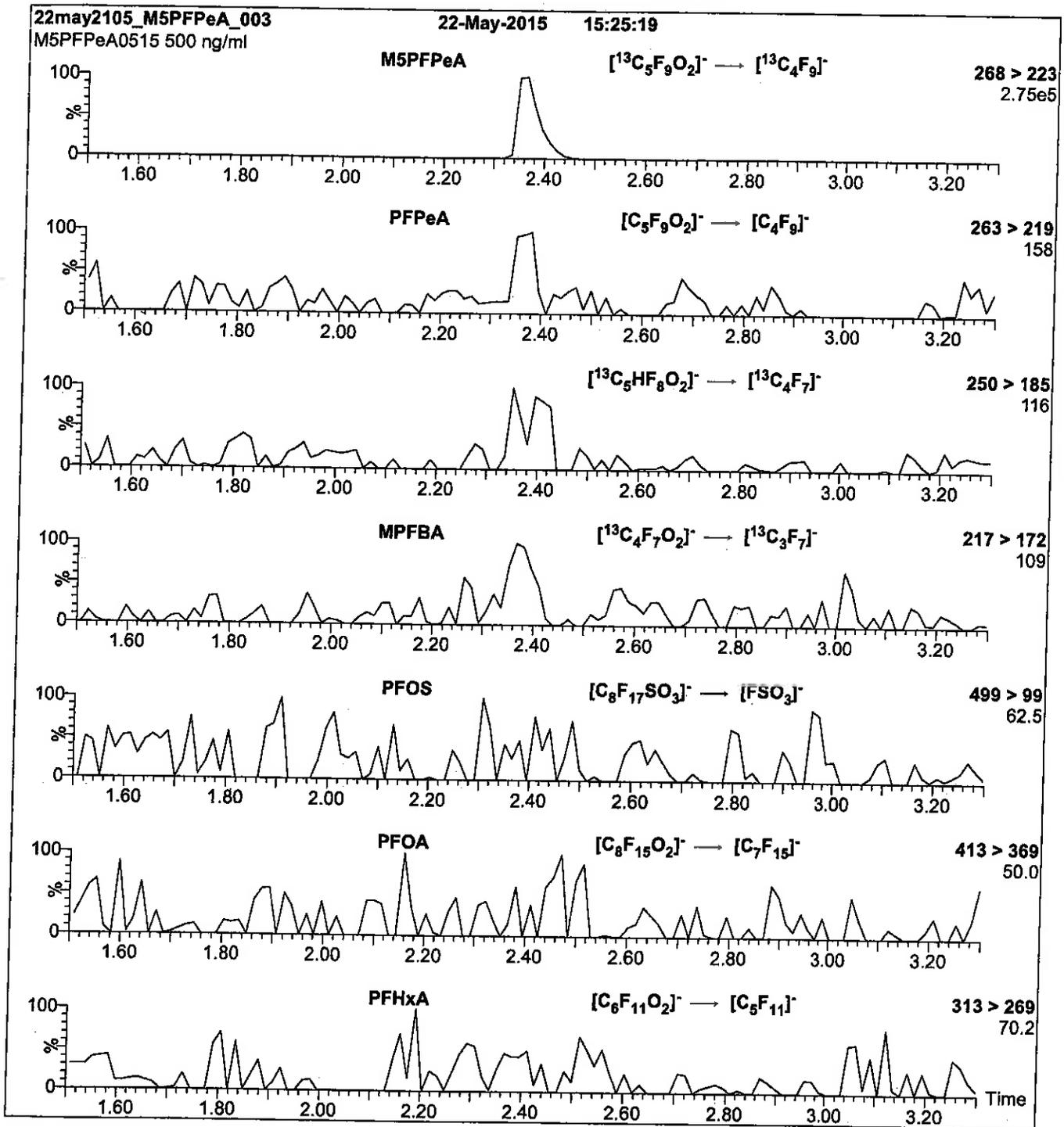
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: M5PFPeA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M5PFPeA)

MS Parameters

Collision Gas (mbar) = 3.35e-3
Collision Energy (eV) = 9

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

Reagent

LCM8FOSA_00011

Scanned 10/14/16 R: SBC 9/22/16



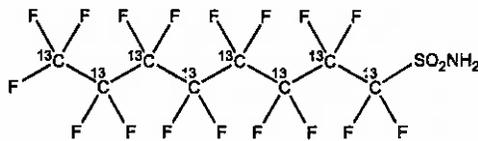
739615
ID: LCM8FOSA_00011
Exp: 12/22/17 Prod: SBC
13C8-Perfluorooctanesulfo



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M8FOSA-I **LOT NUMBER:** M8FOSA1215I
COMPOUND: Perfluoro-1-[¹³C]₈octanesulfonamide
STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₈H₂F₁₇NO₂S **MOLECULAR WEIGHT:** 507.09
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Isopropanol
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 12/22/2015 **EXPIRY DATE:** (mm/dd/yyyy) 12/22/2017
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 01/14/2016
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

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HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

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TRACEABILITY:

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EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

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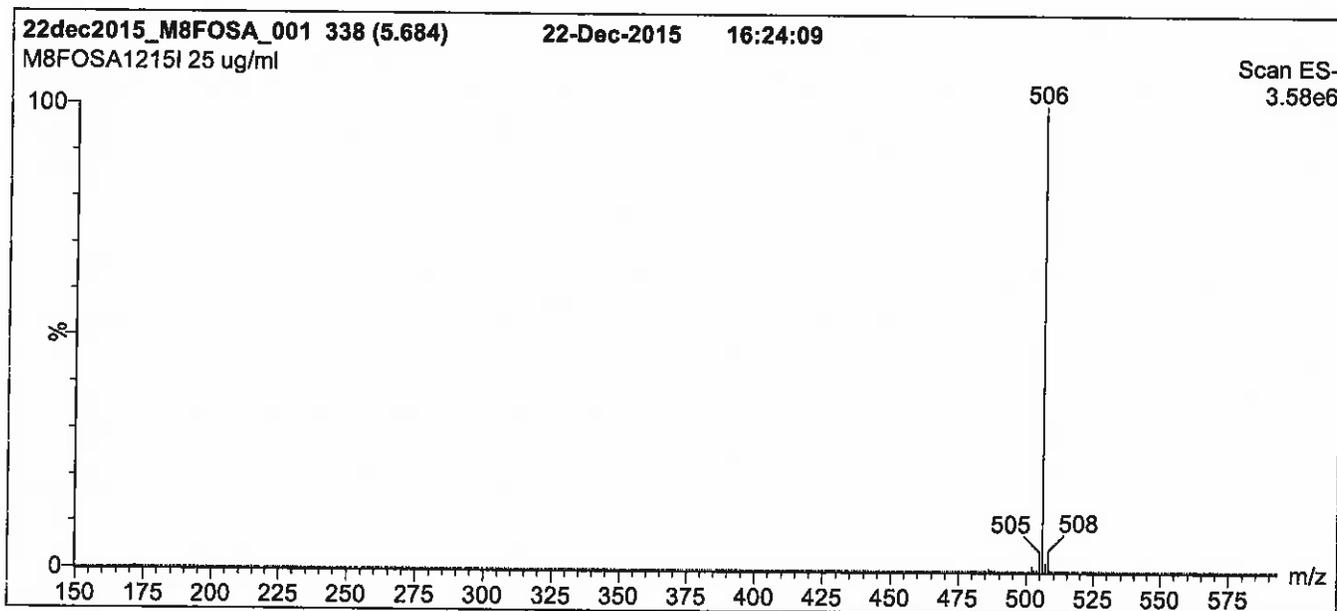
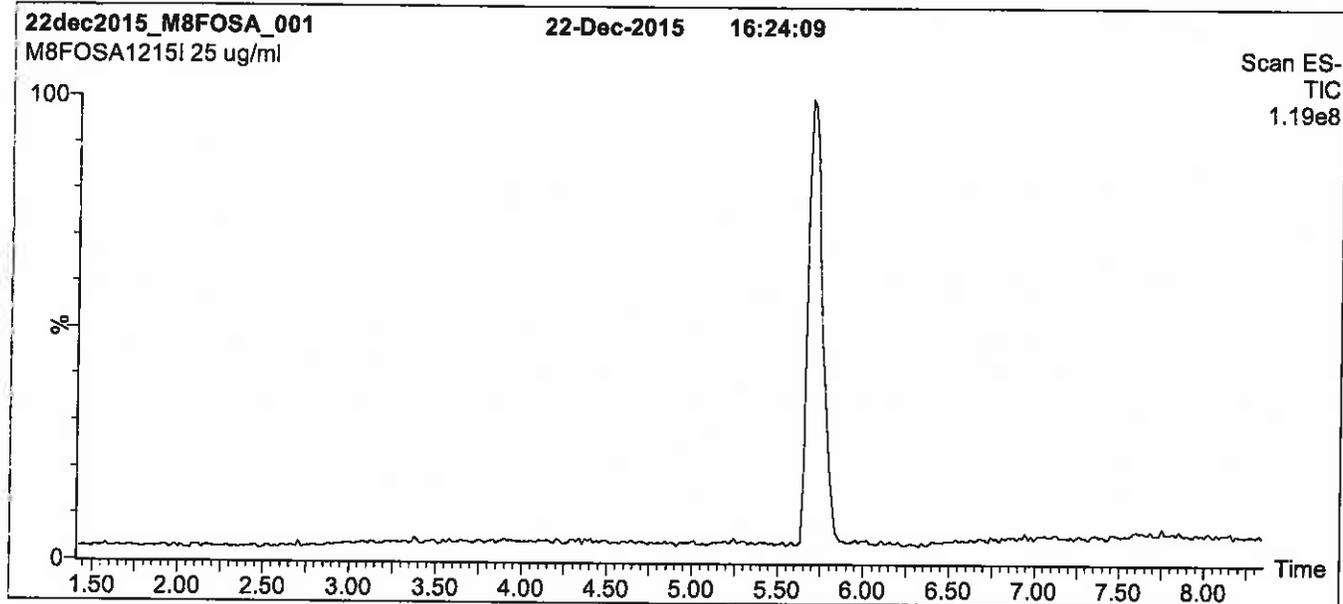
QUALITY MANAGEMENT:

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For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: M8FOSA-I; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min
before returning to initial conditions in 0.5 min.
Time: 10 min

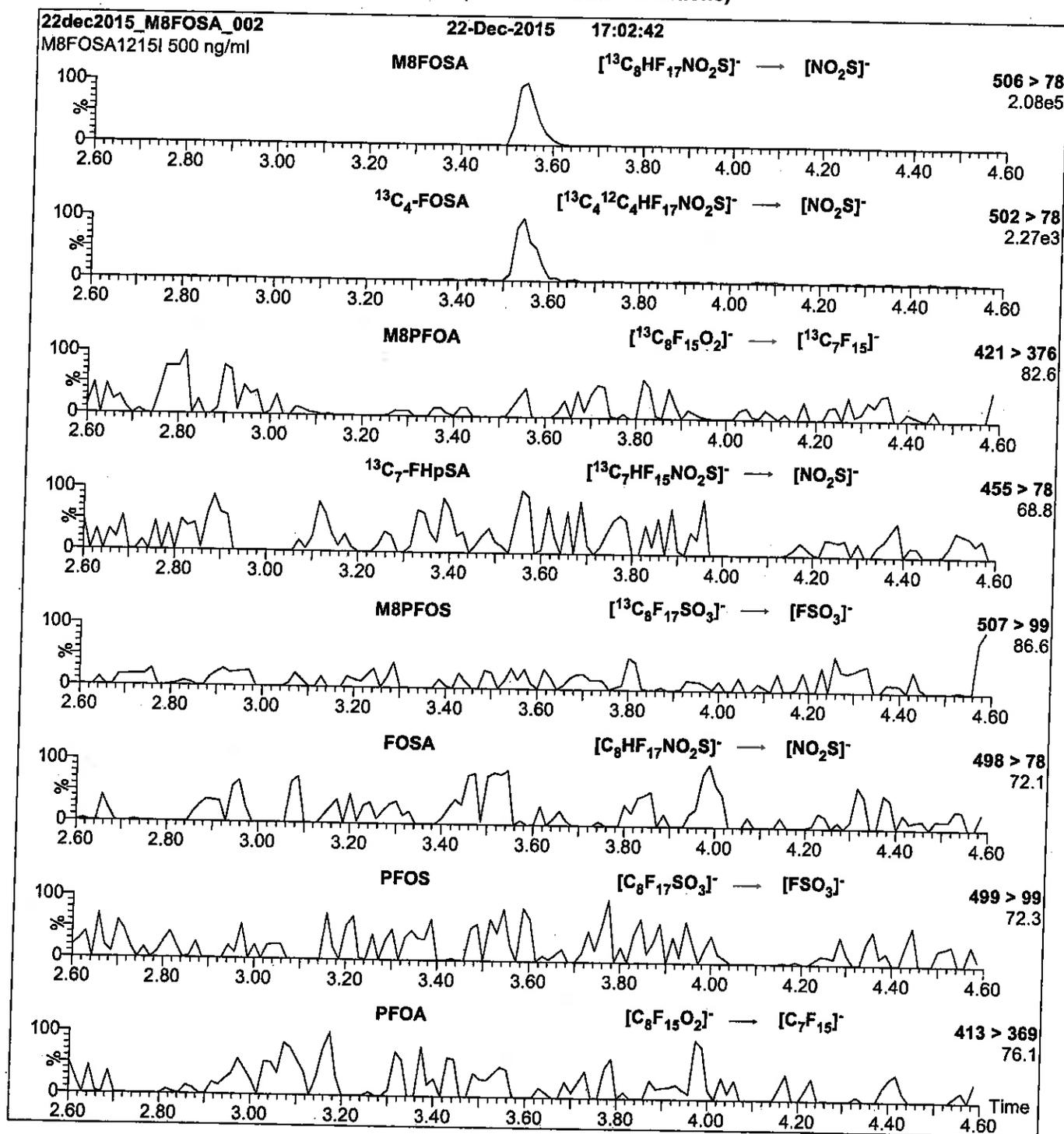
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.50
Cone Voltage (V) = 40.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: M8FOSA-I; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M8FOSA-I)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.39e-3
Collision Energy (eV) = 30

Reagent

LCMPFBA_00008

R: 8BC 9/22/16



739593
ID: LCMFBA_00008
Exp: 05/24/21 Prp4: SEC
13C4-Perfluorobutanoic ac

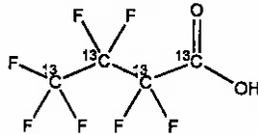


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

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PRODUCT CODE: MPFBA **LOT NUMBER:** MPFBA0516
COMPOUND: Perfluoro-n-[1,2,3,4-¹³C₄]butanoic acid
STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₄HF₇O₂ **MOLECULAR WEIGHT:** 218.01
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99%¹³C
(1,2,3,4-¹³C₄)
LAST TESTED: (mm/dd/yyyy) 05/24/2016
EXPIRY DATE: (mm/dd/yyyy) 05/24/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

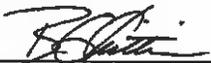
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 05/30/2016
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

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HOMOGENEITY:

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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

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EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

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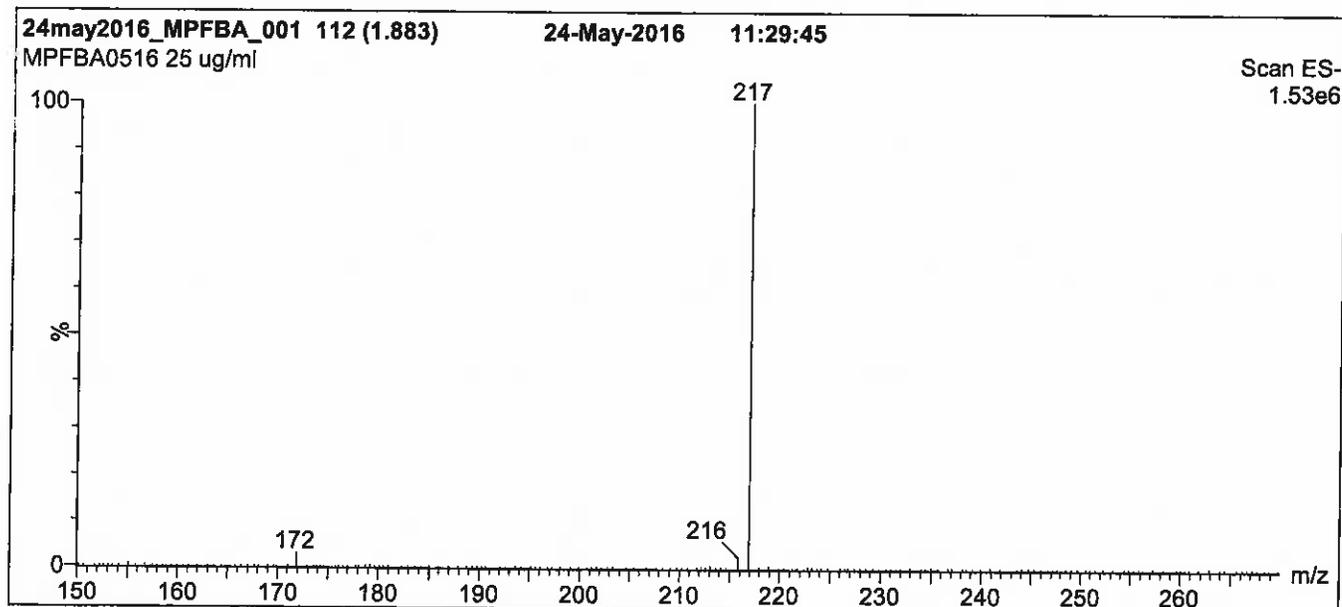
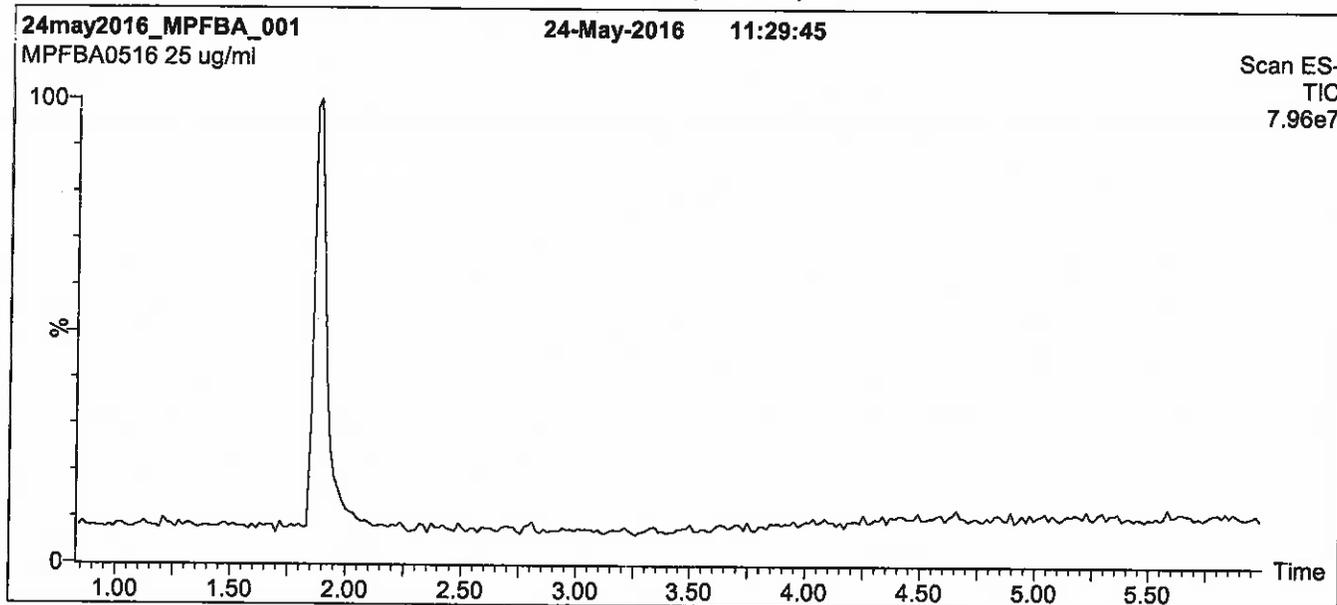
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: MPFBA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 30% (80:20 MeOH:ACN) / 70% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

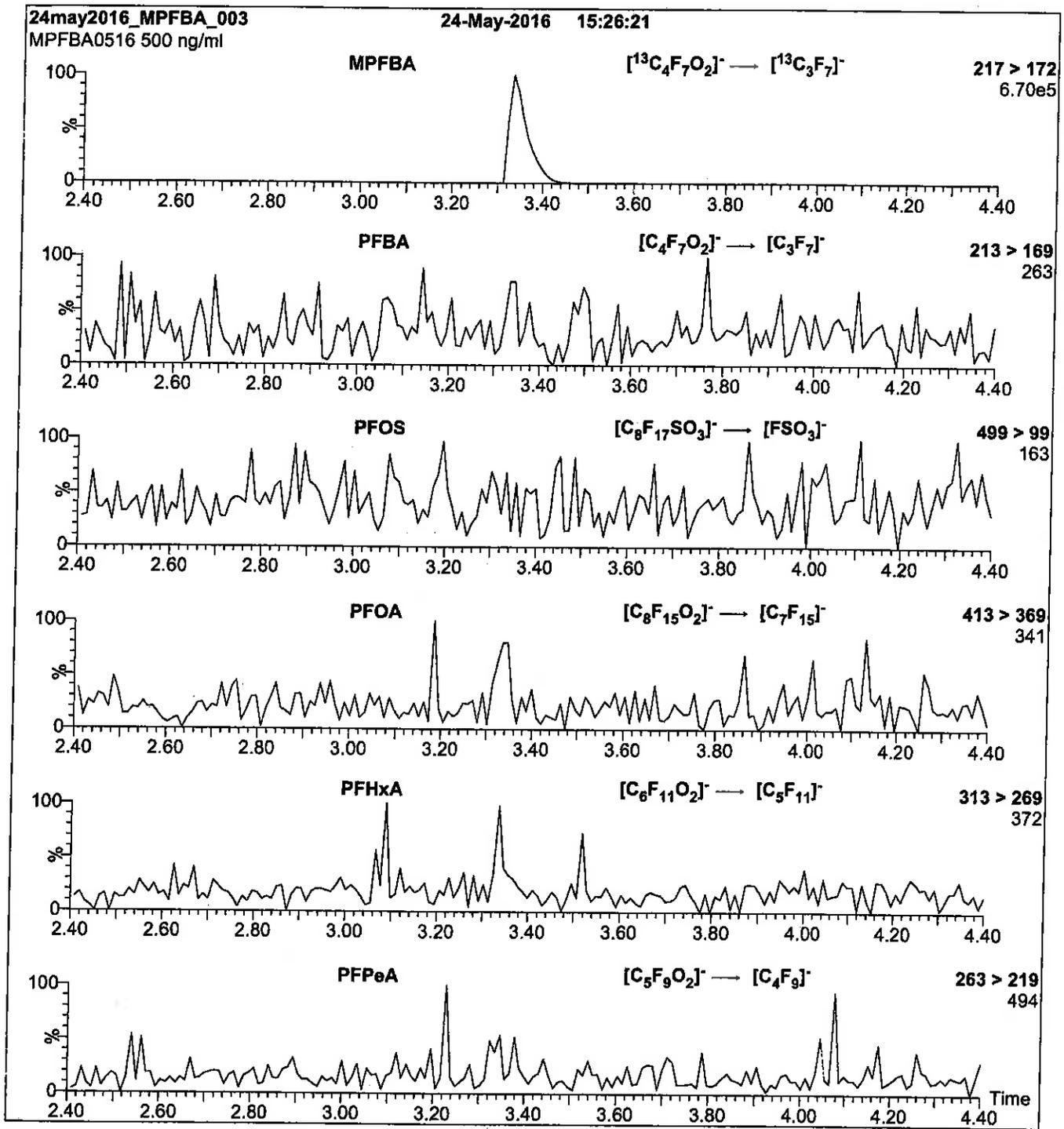
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 10.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFBA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop Injection
10 μl (500 ng/ml MPFBA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.50e-3
Collision Energy (eV) = 10

Reagent

LCMPFDA_00011

Scanned 10/14/16 R: SBC 9/22/16

739609
ID: LCMFDA_00011
Exp: 08/19/20 Prep: SBC
13C2-Perfluorodecanoic a

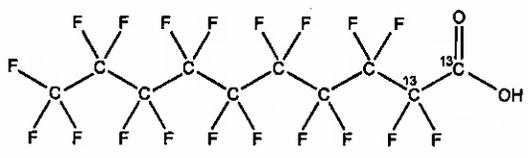


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFDA **LOT NUMBER:** MPFDA0815
COMPOUND: Perfluoro-n-[1,2-¹³C₂]decanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₈HF₁₈O₂ **MOLECULAR WEIGHT:** 516.07
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
(1,2-¹³C₂)
LAST TESTED: (mm/dd/yyyy) 08/19/2015
EXPIRY DATE: (mm/dd/yyyy) 08/19/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

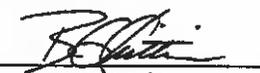
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains < 0.1% of ¹³C₁-PFNA.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 08/21/2015
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

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LIMITED WARRANTY:

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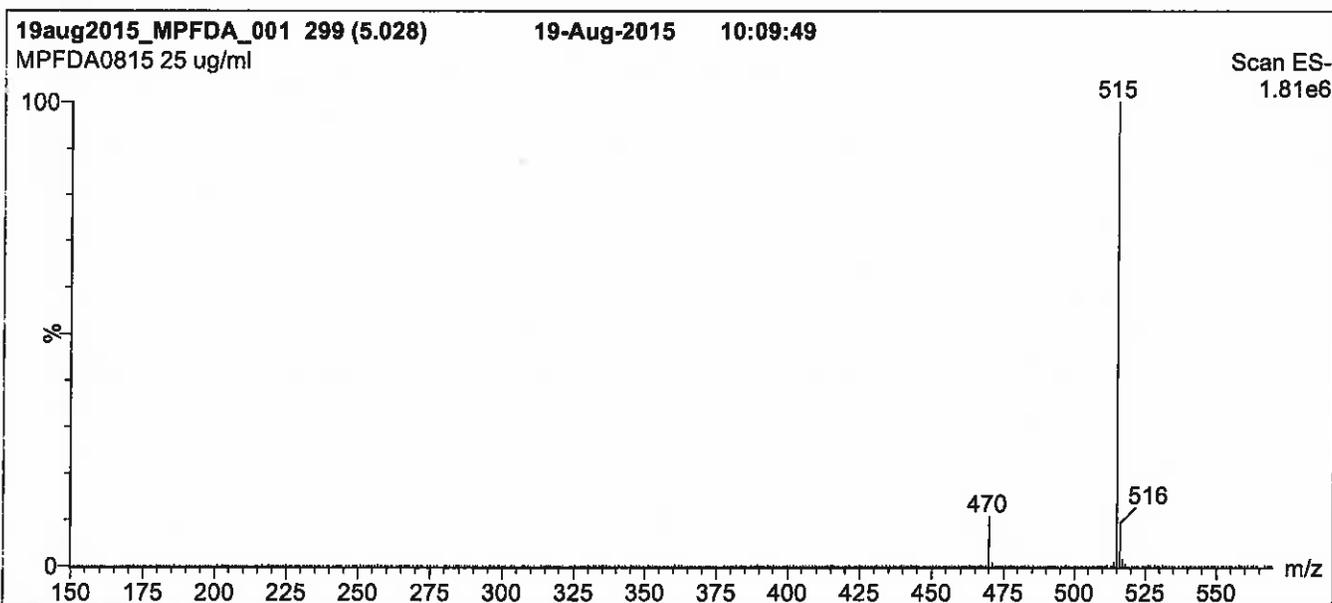
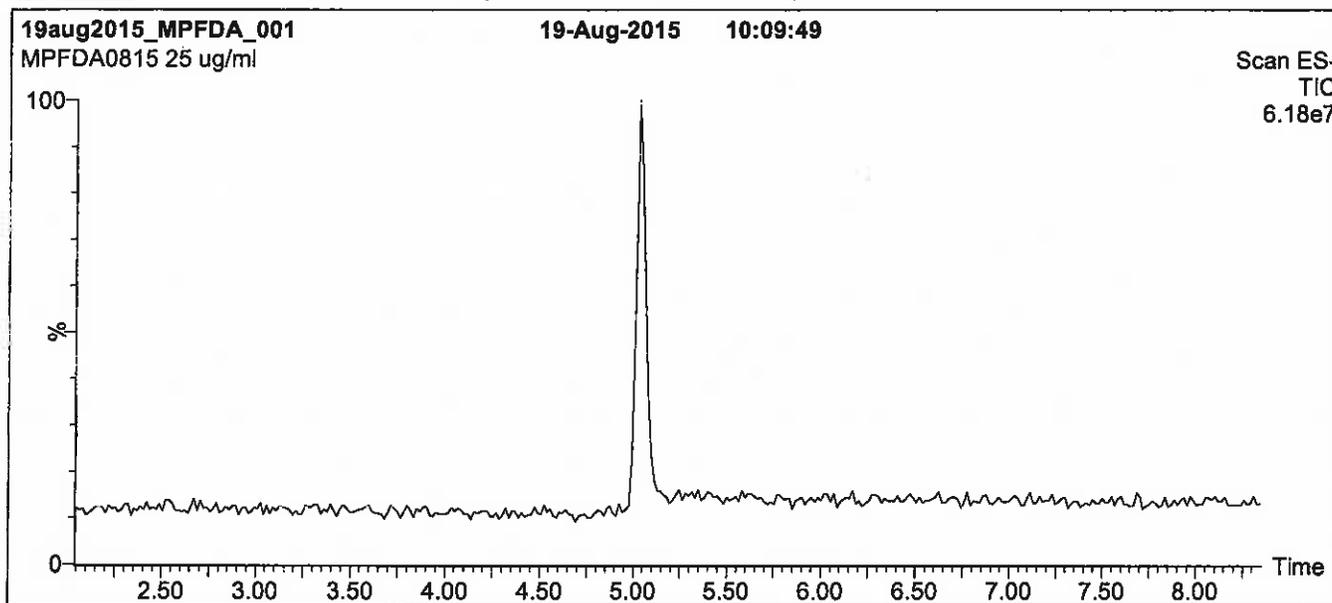
QUALITY MANAGEMENT:

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Figure 1: MPFDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

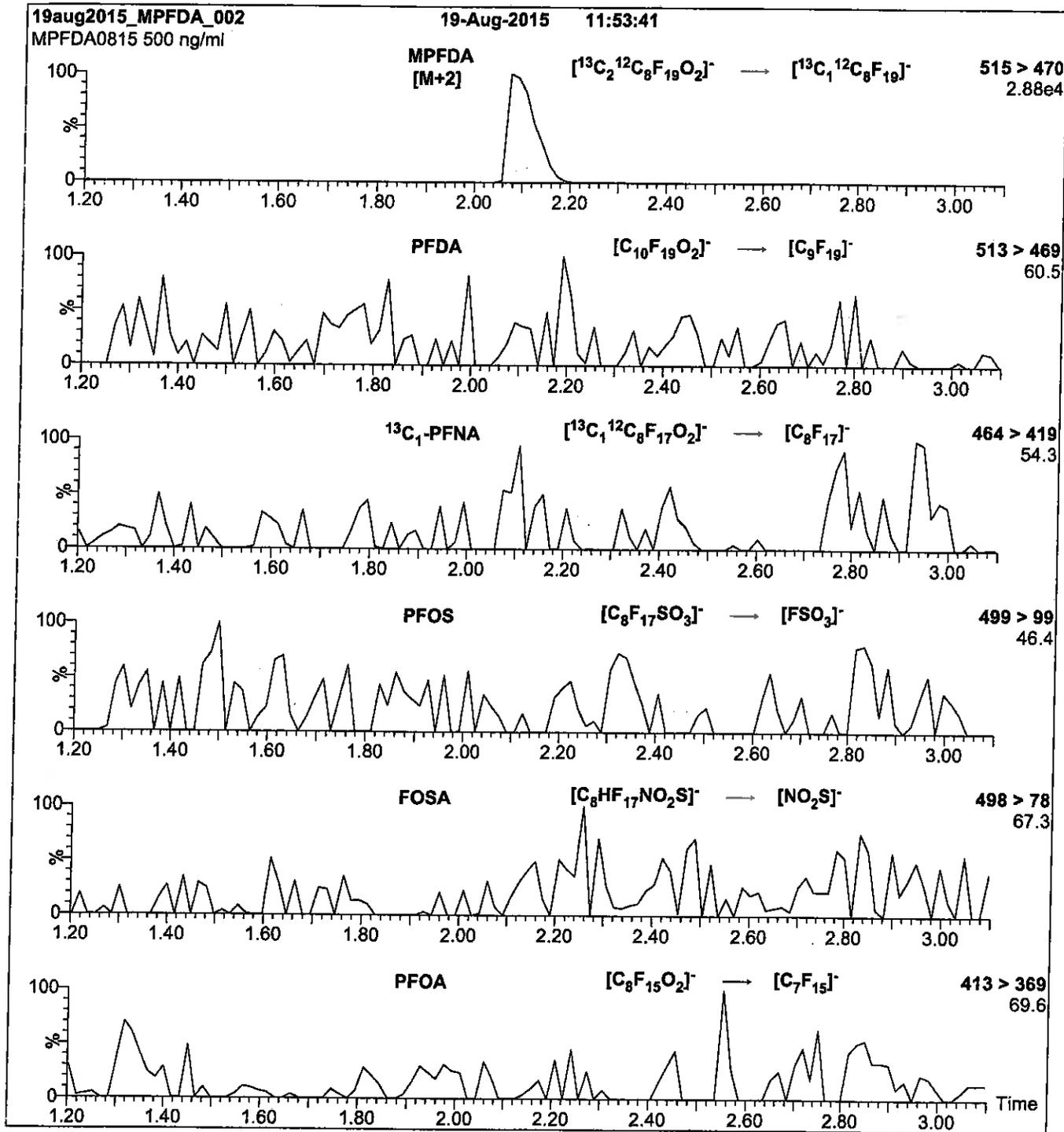
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml MPFDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.35e-3
Collision Energy (eV) = 13

Reagent

LCMPFD_oA_00008

R: 882 9/22/16



739598
ID: LCMFDoA_00008
Exp: 04/08/21 Prod: SBC
13C2-Perfluorododecanoic



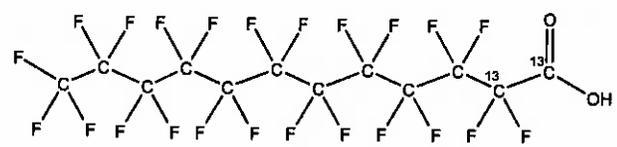
WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

Scanned 10/14/16 SR

PRODUCT CODE: MPFDoA **LOT NUMBER:** MPFDoA0416
COMPOUND: Perfluoro-n-[1,2-¹³C₂]dodecanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₁₀HF₂₃O₂ **MOLECULAR WEIGHT:** 616.08
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 04/08/2016 (1,2-¹³C₂)
EXPIRY DATE: (mm/dd/yyyy) 04/08/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 04/15/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • Info@well-labs.com

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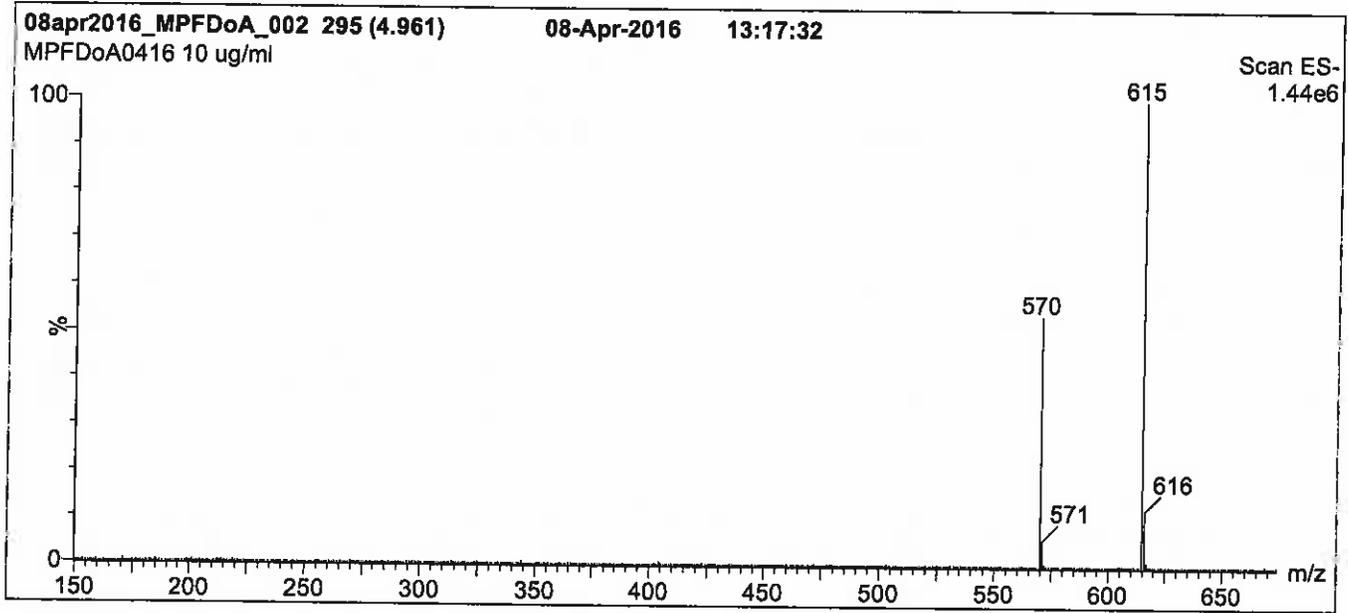
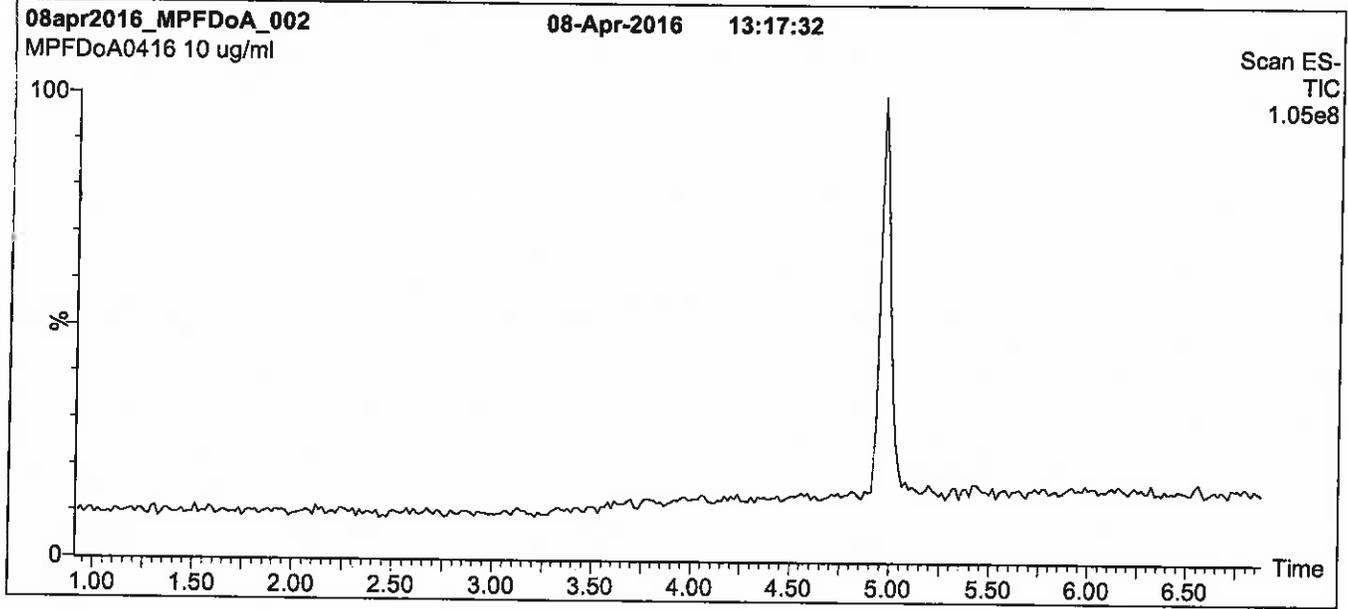
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Figure 1: MPFDoA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

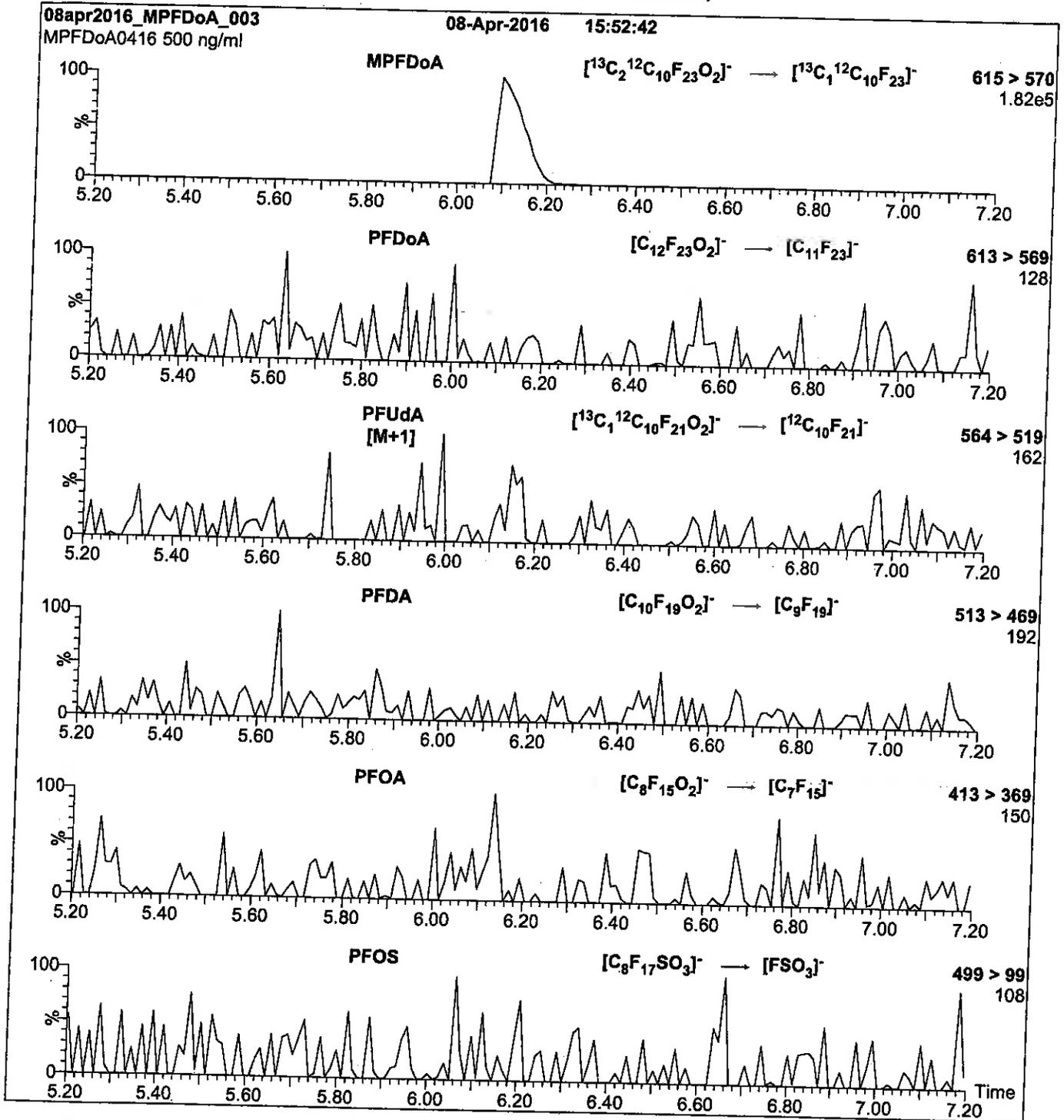
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 20.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFDoA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFDoA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.24e-3
Collision Energy (eV) = 13

Reagent

LCMPFHxA_00012

Scanned 10/11/16 R: SBC 9/22/16

739612
ID: LCMPFHxA_00012
Exp: 04/08/21 Prpd: SBC
13C2-Perfluorohexanoic ac



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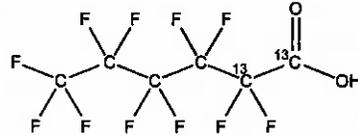
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFHxA
COMPOUND: Perfluoro-n-[1,2-¹³C₂]hexanoic acid

LOT NUMBER: MPFHxA0416

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: ¹³C₂¹²C₄HF₁₁O₂
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 316.04
SOLVENT(S): Methanol
Water (<1%)

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 04/08/2016
EXPIRY DATE: (mm/dd/yyyy) 04/08/2021

ISOTOPIC PURITY: ≥99%¹³C
(1,2-¹³C₂)

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

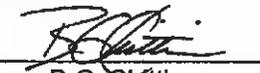
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains < 0.1% of perfluoro-n-hexanoic acid and ~ 0.3% of perfluoro-n-octanoic acid.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim
Date: 04/29/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
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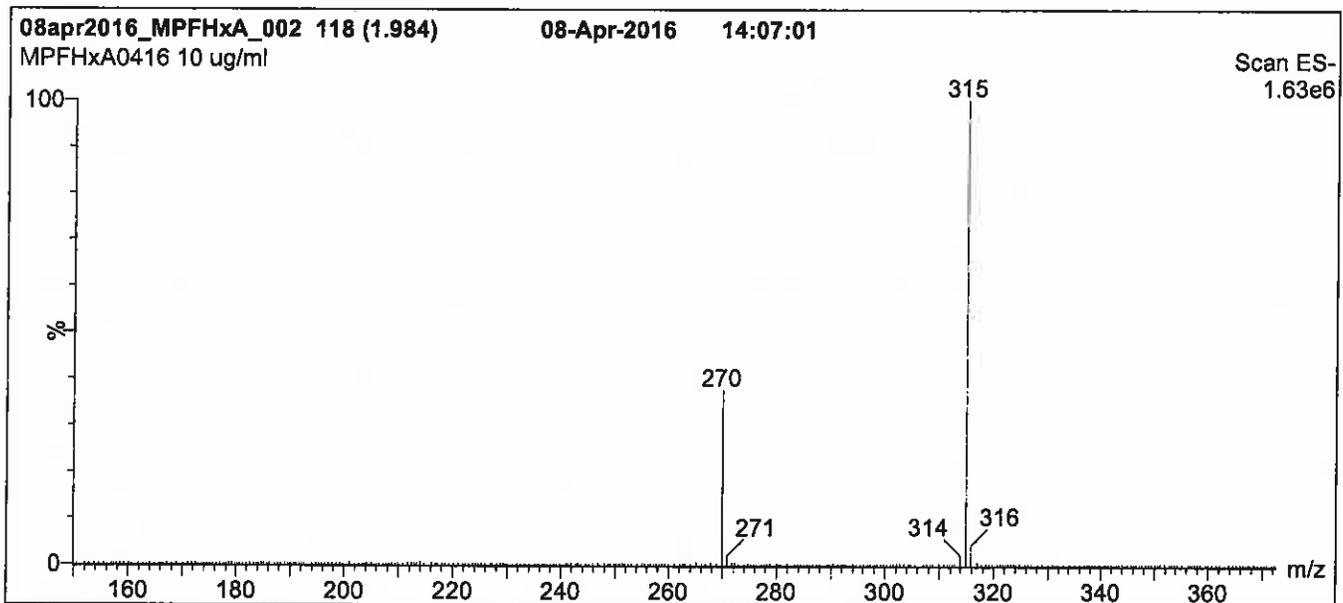
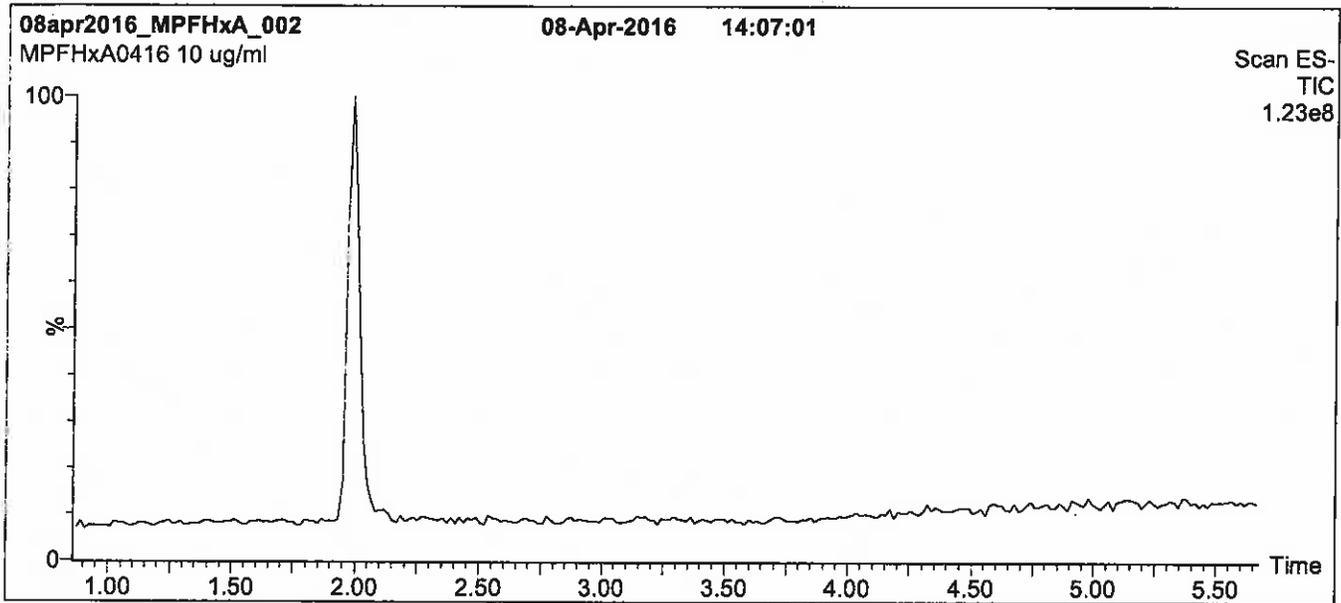
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Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

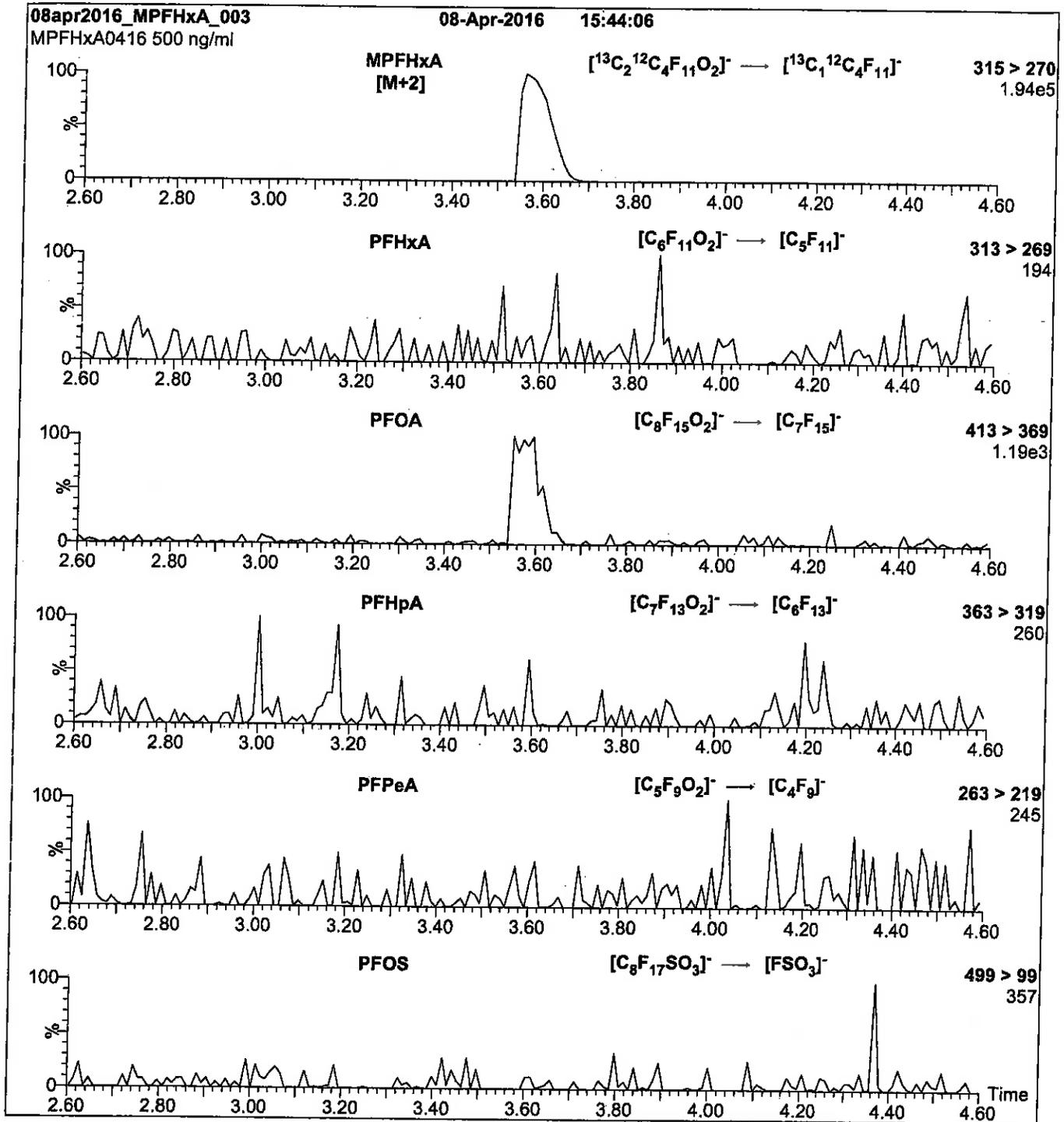
Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7.5 min and hold for 1.5 min
 before returning to initial conditions over 0.5 min.
 Time: 10 min

Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)
 Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFHxA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml MPFHxA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.39e-3
Collision Energy (eV) = 10

Reagent

LCMPFHXS_00008

R: 800 9/22/16



739601

ID: LCMPFHxS_00008

Exp: 10/23/20 Prod: SBC

18O2-Perfluorohexanesulfo



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

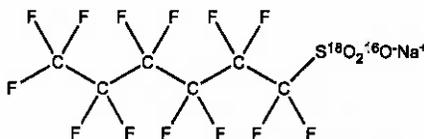
Scanned 10/14/16 SK

PRODUCT CODE: MPFHxS
COMPOUND: Sodium perfluoro-1-hexane[¹⁸O₂]sulfonate

LOT NUMBER: MPFHxS1015

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: C₆F₁₃S¹⁸O₂¹⁶ONa
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt)
47.3 ± 2.4 µg/ml (MPFHxS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 10/23/2015
EXPIRY DATE: (mm/dd/yyyy) 10/23/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 426.10
SOLVENT(S): Methanol
ISOTOPIC PURITY: >94% (¹⁸O₂)

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The response factor for MPFHxS (C₆F₁₃S¹⁸O₂¹⁶O) has been observed to be up to 10% lower than for PFHxS (C₆F₁₃S¹⁶O₃) when both compounds are injected together. This difference may vary between instruments.
- Due to the isotopic purity of the starting material (¹⁸O₂ >94%), MPFHxS contains ~ 0.3% of PFHxS. This value agrees with the theoretical percent relative abundance that is expected based on the stated isotopic purity.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 10/28/2015

(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

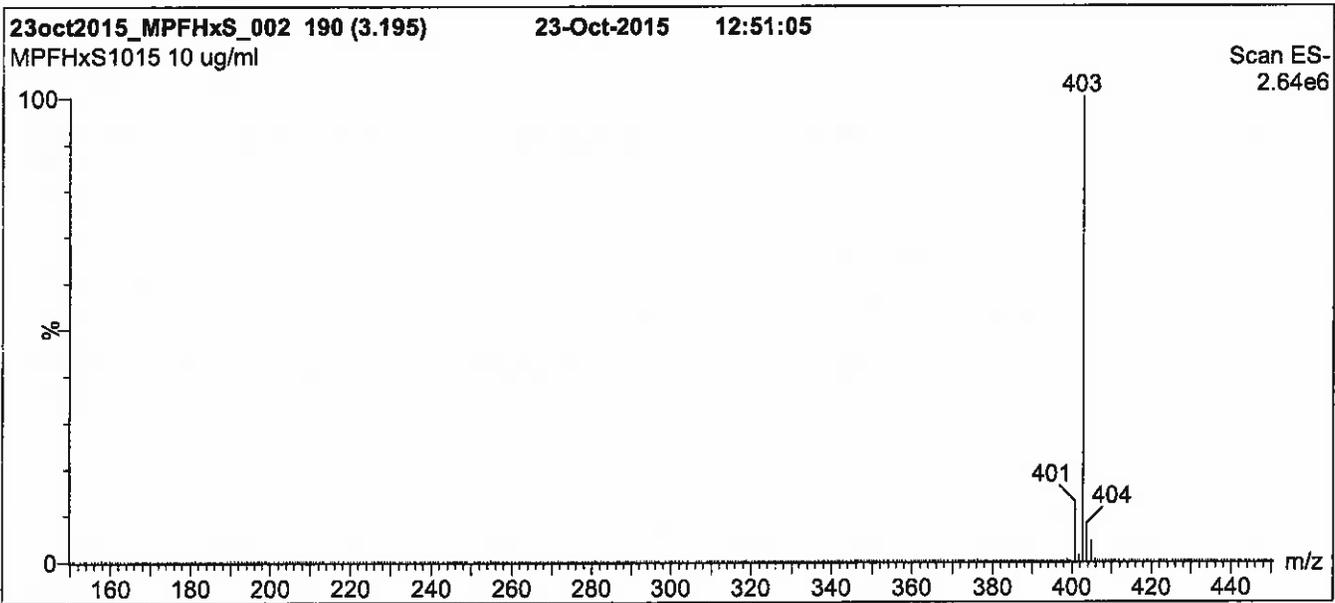
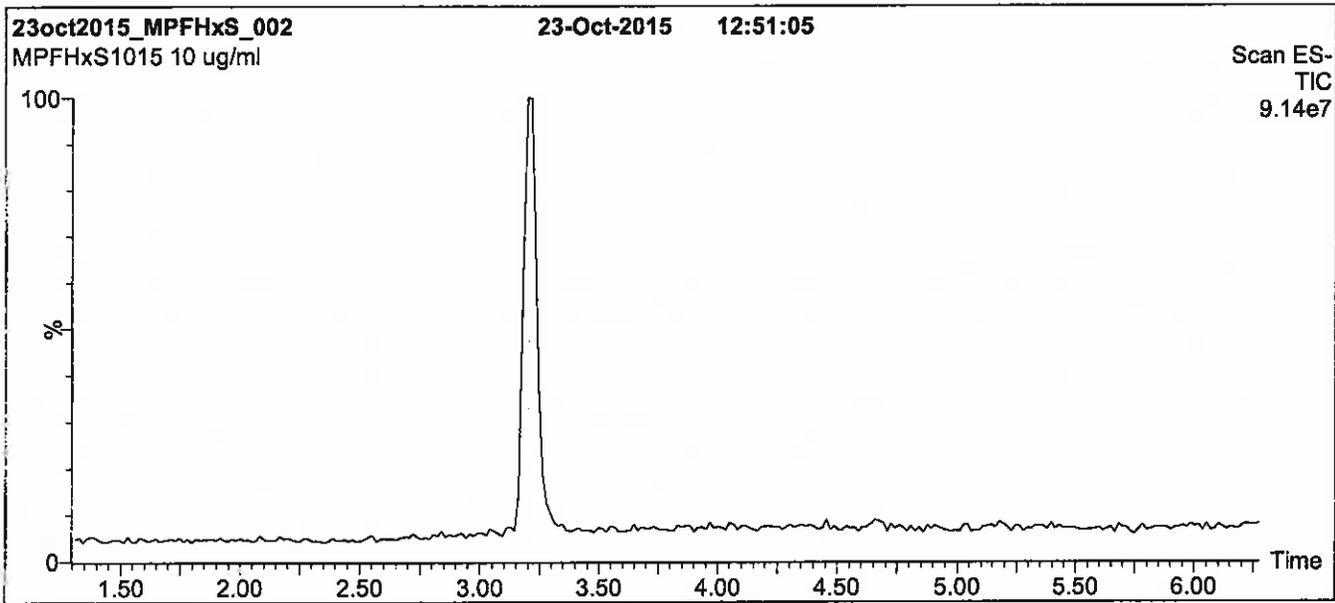
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: MPFHxS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

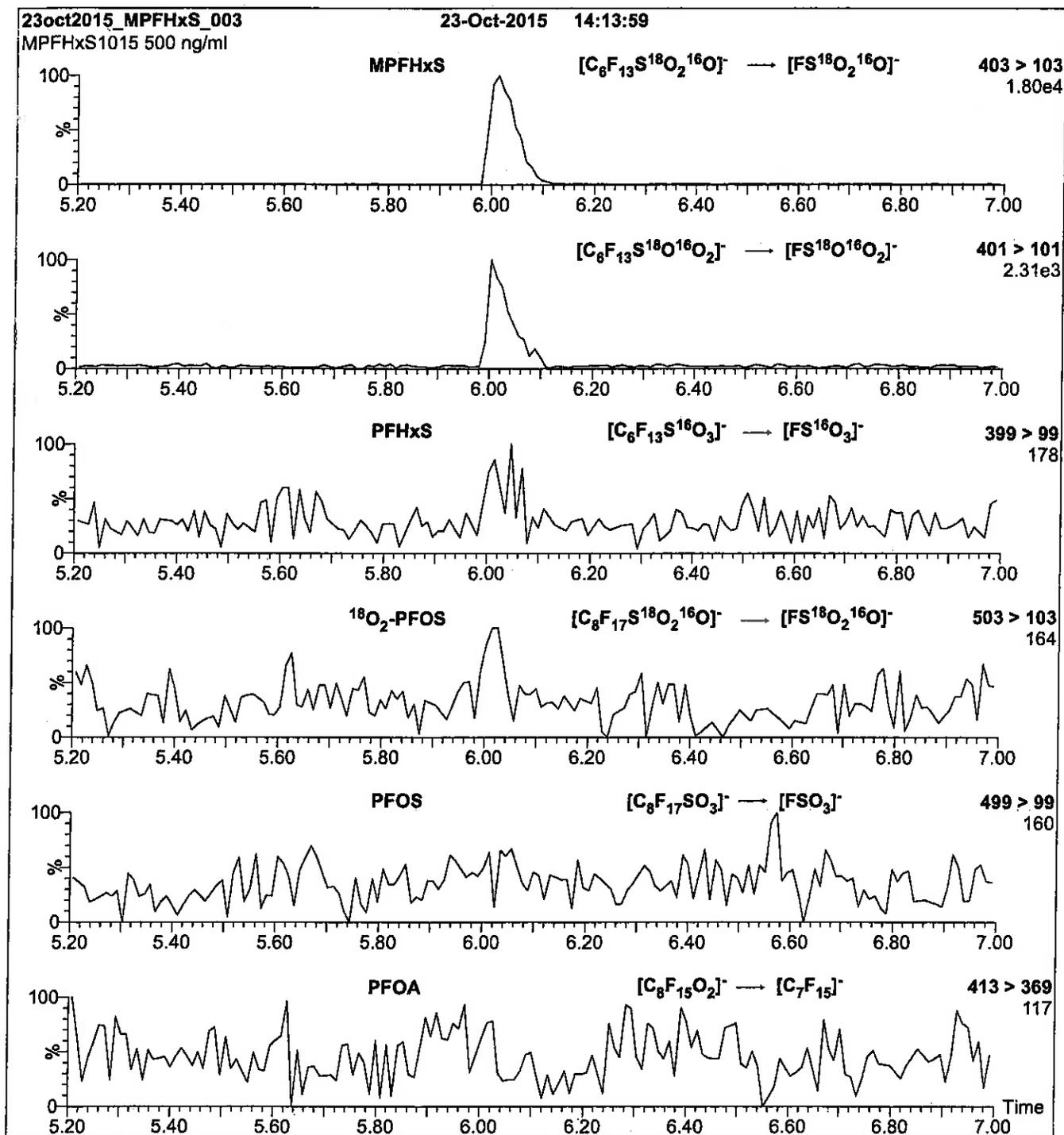
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 50.00
 Cone Gas Flow (l/hr) = 60
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFHxS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml MPFHxS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.35e-3
 Collision Energy (eV) = 30

Reagent

LCMPFNA_00008

Scanned 10/14/16 R: SBC 9/22/16



739637
ID: LCM:PFNA_0008
Exp: 04/13/19 Pppl: SBC
13C5-Perfluoronoic aci

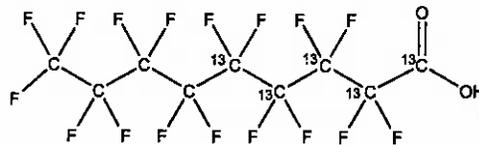


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFNA **LOT NUMBER:** MPFNA0414
COMPOUND: Perfluoro-n-[1,2,3,4,5-¹³C₅]nonanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₅¹²C₄HF₁₇O₂ **MOLECULAR WEIGHT:** 469.04
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99%¹³C
(1,2,3,4,5-¹³C₅)
LAST TESTED: (mm/dd/yyyy) 04/13/2014
EXPIRY DATE: (mm/dd/yyyy) 04/13/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 04/01/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

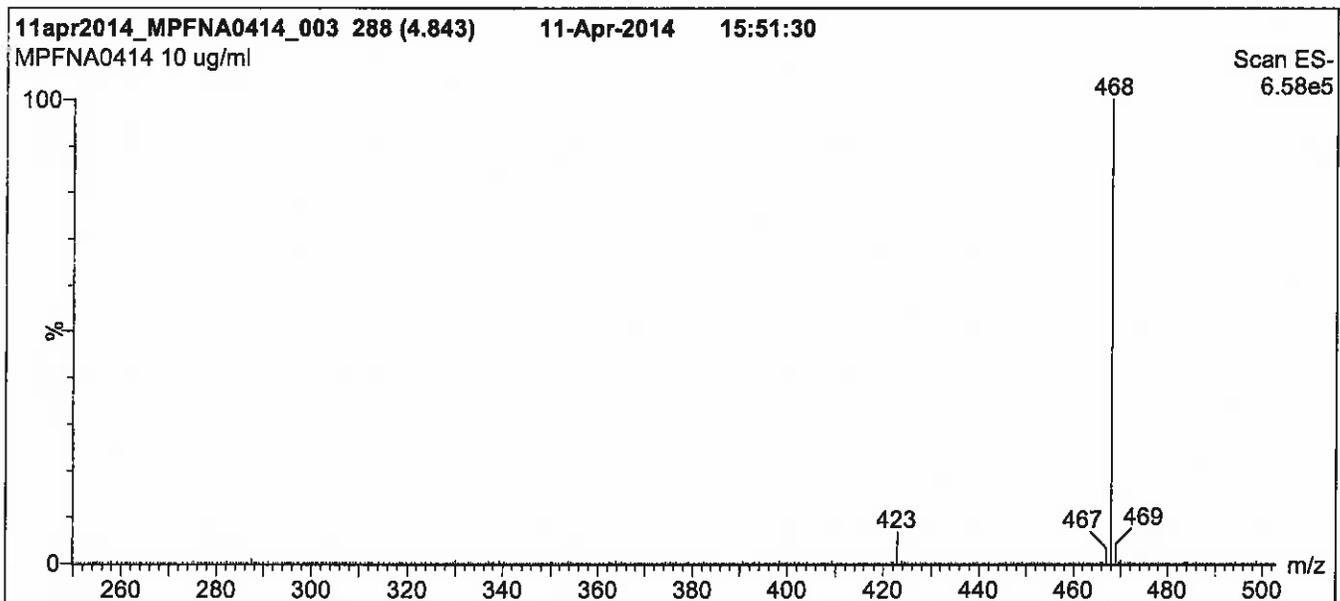
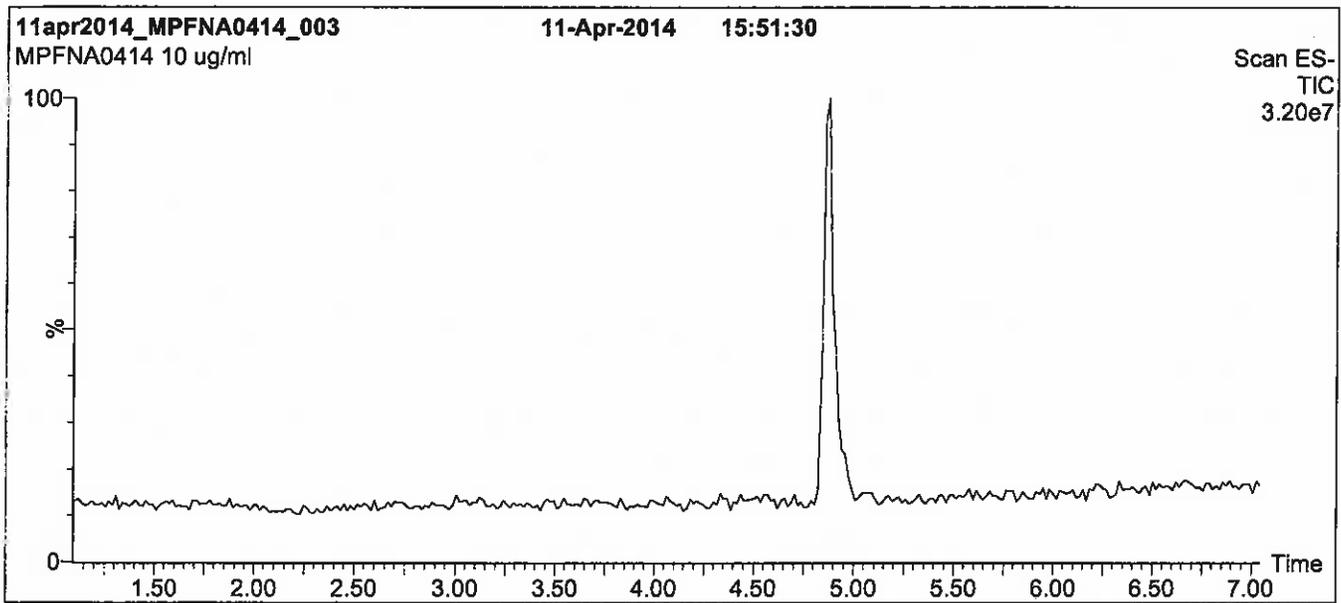
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: MPFNA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

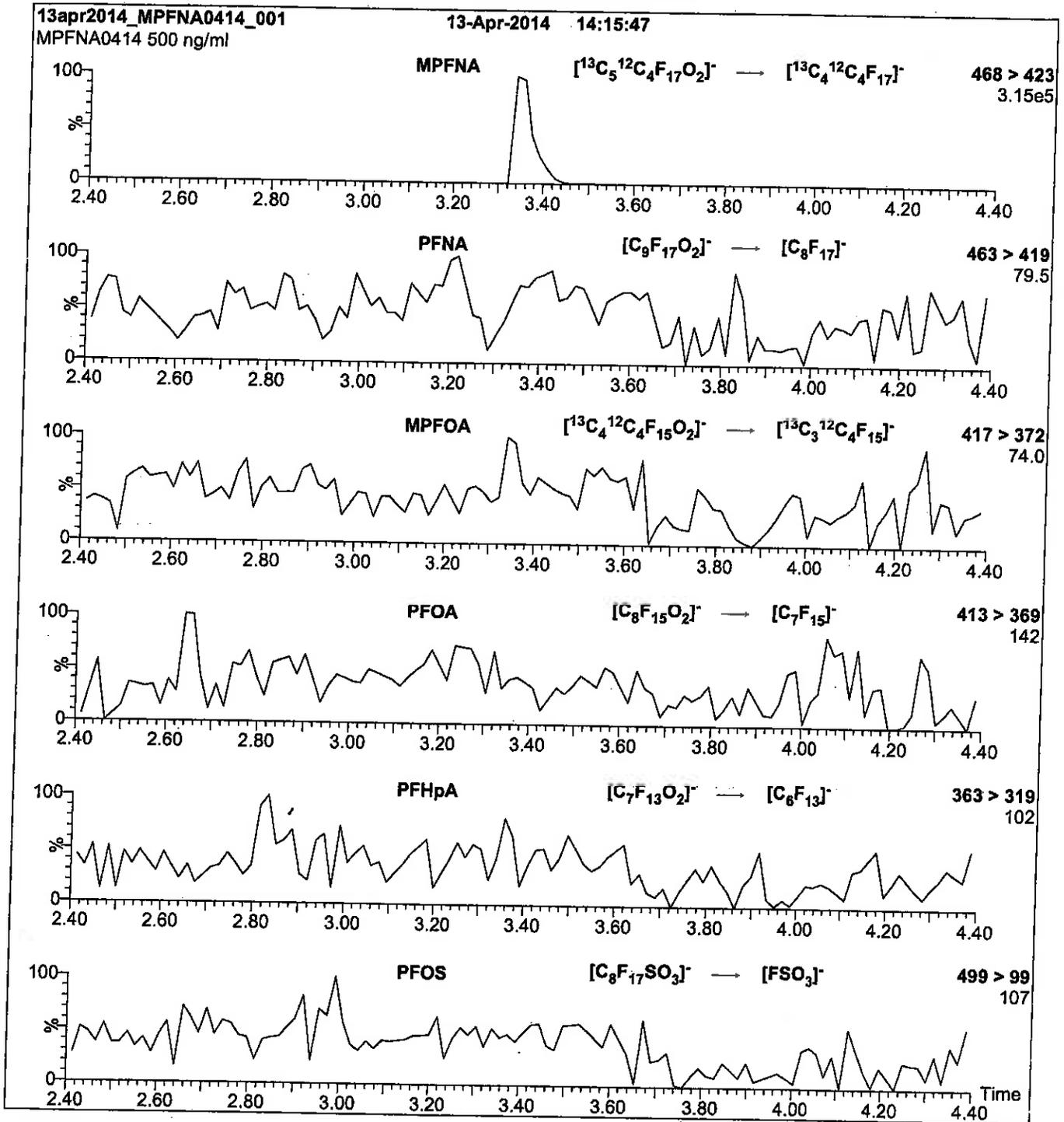
Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm
 Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 850 amu)
 Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFNA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFNA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.28e-3
Collision Energy (eV) = 11

Reagent

LCMPFOA_00012

R: SBC 9/22/16



738683
ID: LCMFOA_00012
Exp: 01/22/21 Prep: SBC
13C4-Perfluorooctanoic ac



WELLINGTON LABORATORIES

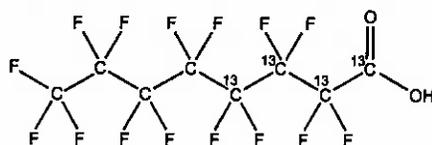
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFOA
COMPOUND: Perfluoro-n-[1,2,3,4-¹³C₄]octanoic acid

LOT NUMBER: MPFOA0116

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: ¹³C₄¹²C₄HF₁₅O₂
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 418.04
SOLVENT(S): Methanol
Water (<1%)

CHEMICAL PURITY: >98%

ISOTOPIC PURITY: ≥99% ¹³C
(1,2,3,4-¹³C₄)

LAST TESTED: (mm/dd/yyyy) 01/22/2016

EXPIRY DATE: (mm/dd/yyyy) 01/22/2021

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.1% of native perfluoro-n-octanoic acid (PFOA).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:
B.G. Chittim

Date: 02/01/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

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LIMITED WARRANTY:

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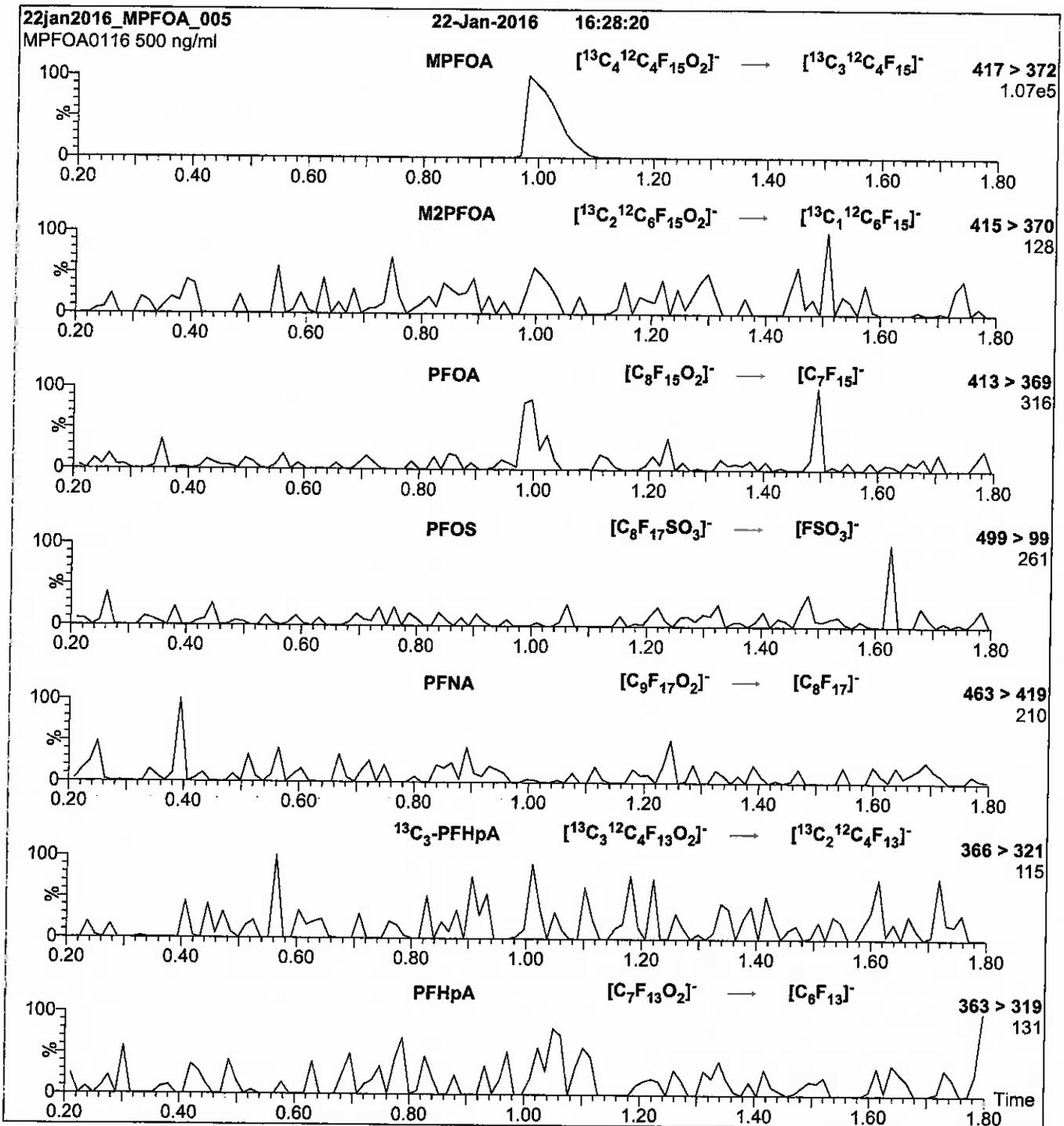
QUALITY MANAGEMENT:

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For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 2: MPFOA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFOA)

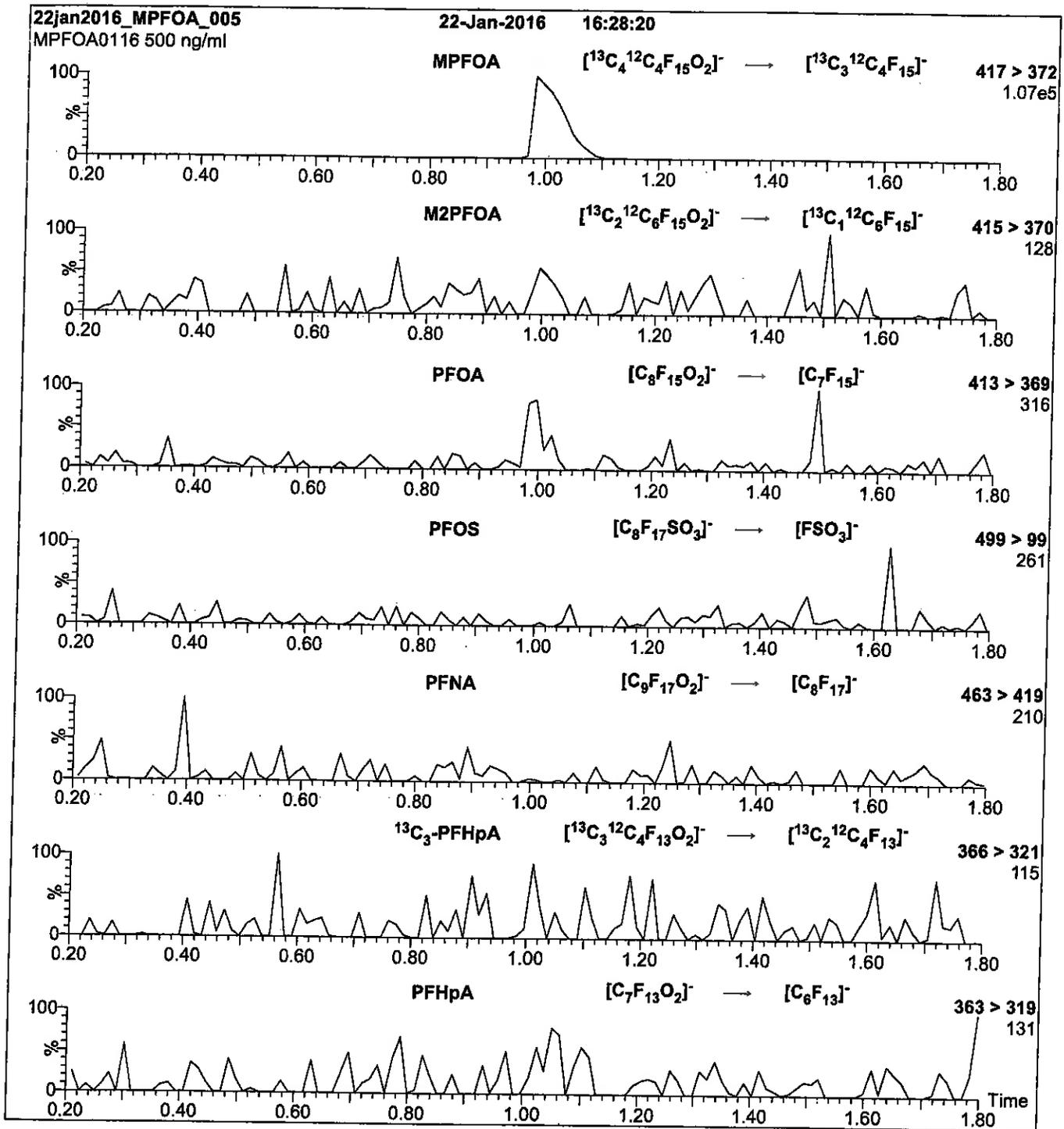
Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.58e-3
Collision Energy (eV) = 10

Figure 2: MPFOA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFOA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.58e-3
Collision Energy (eV) = 10

Reagent

LCMPFOS_00017

R: 9/9/16 802

728309
ID: LCMPPFOS_00017
Exp: 08/03/21 Prpd: SBC
13C4-Perfluorooctanesulfo

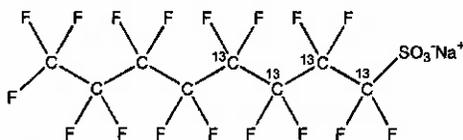


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFOS **LOT NUMBER:** MPFOS0816
COMPOUND: Sodium perfluoro-1-[1,2,3,4-¹³C₄]octanesulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₄¹²C₄F₁₇SO₃Na **MOLECULAR WEIGHT:** 526.08
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.8 ± 2.4 µg/ml (MPFOS anion)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 08/03/2016 (1,2,3,4-¹³C₄)
EXPIRY DATE: (mm/dd/yyyy) 08/03/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

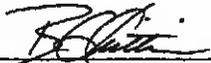
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.8% Sodium perfluoro-1-[1,2,3-¹³C₃]heptanesulfonate.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 08/05/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

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SYNTHESIS / CHARACTERIZATION:

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HOMOGENEITY:

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where x is expressed as a relative standard uncertainty of the individual parameter.

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LIMITED WARRANTY:

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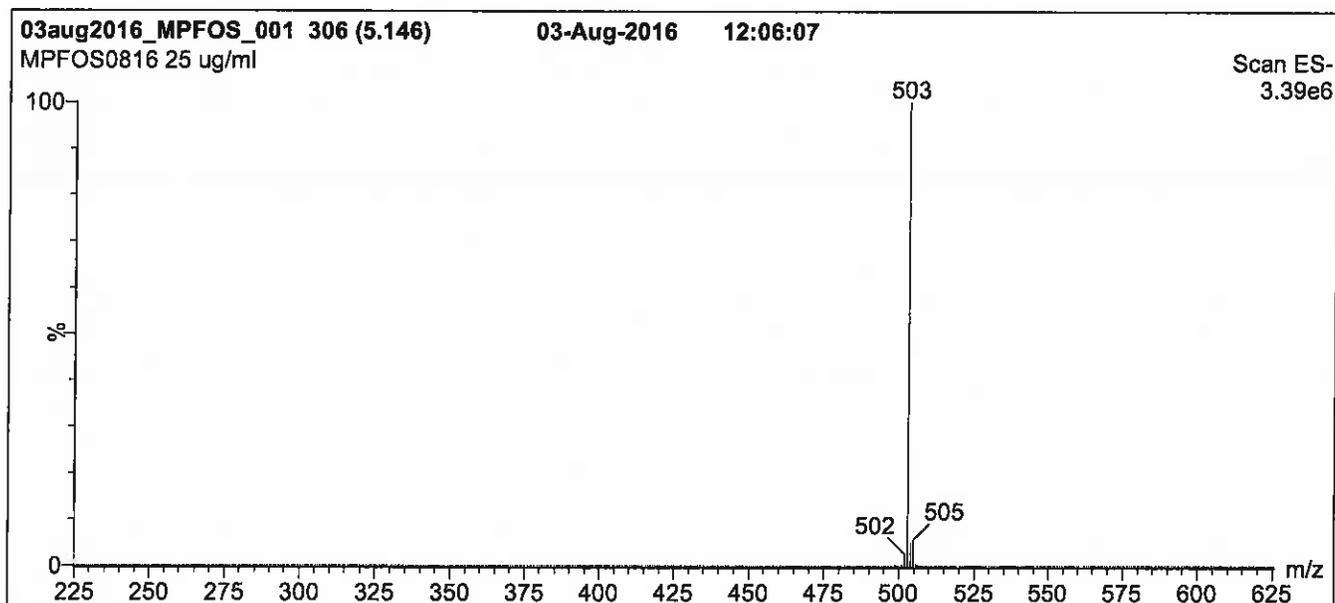
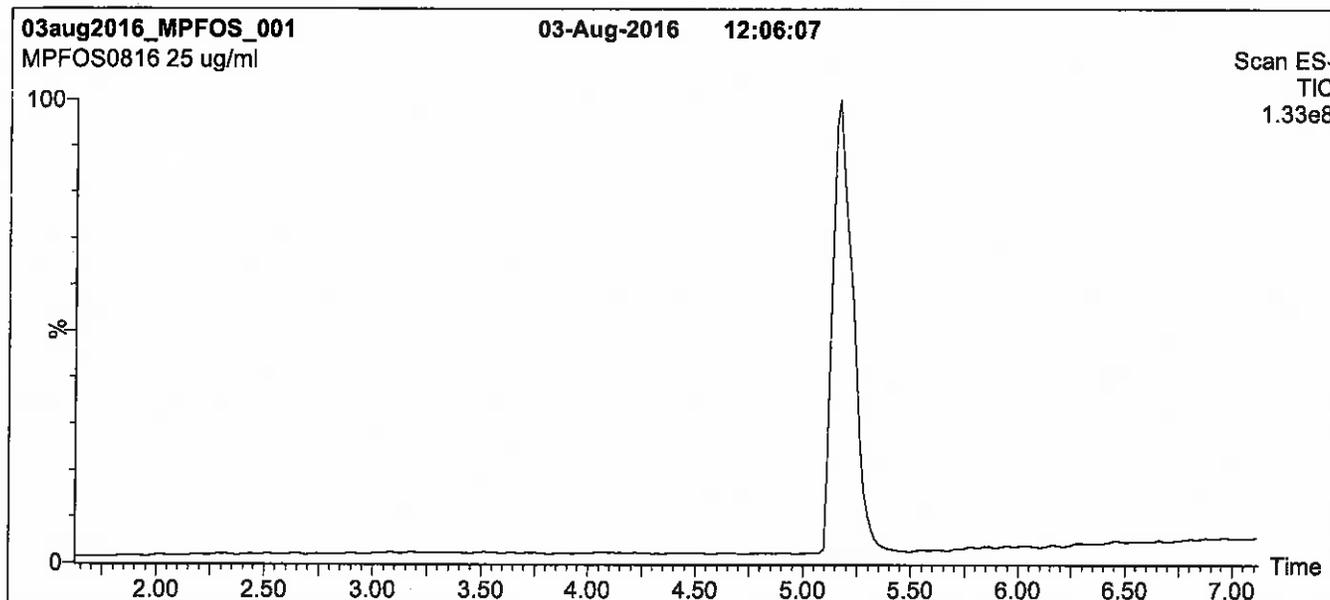
QUALITY MANAGEMENT:

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Figure 1: MPFOS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 45% (80:20 MeOH:ACN) / 55% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

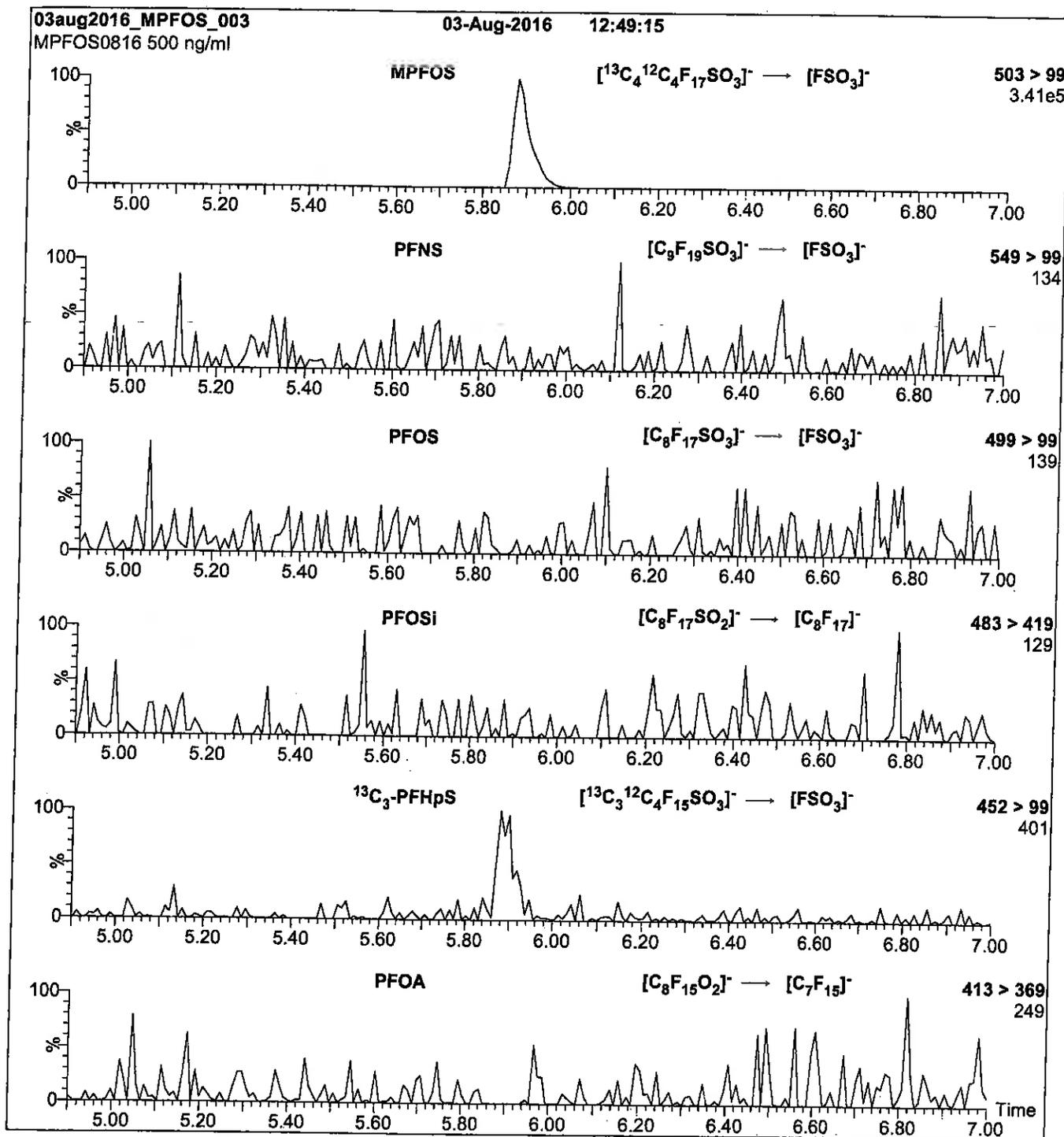
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 60.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFOS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFOS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.46e-3
Collision Energy (eV) = 40

Reagent

LCMPFUdA_00009

R: SBC 9/22/16

739604
ID: LCMPFUdA_00009
Exp: 02/12/21 Prod: SBC
13C2-Perfluoroundecanoic



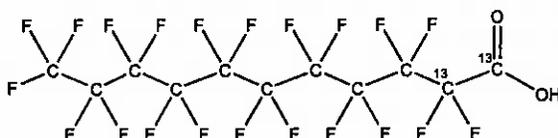
WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

Scanned 10/14/16 SK

PRODUCT CODE: MPFUdA **LOT NUMBER:** MPFUdA0216
COMPOUND: Perfluoro-n-[1,2-¹³C₂]undecanoic acid

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₉HF₂₁O₂ **MOLECULAR WEIGHT:** 566.08
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
(1,2-¹³C₂)
LAST TESTED: (mm/dd/yyyy) 02/12/2016
EXPIRY DATE: (mm/dd/yyyy) 02/12/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

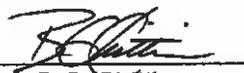
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Presence of 1-¹³C₁-PFUdA (~1%; see Figure 2), 2-¹³C₁-PFUdA (~1%), and PFUdA (~0.2%; see Figure 2) are due to the isotopic purity of the ¹³C-precursor.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 02/24/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

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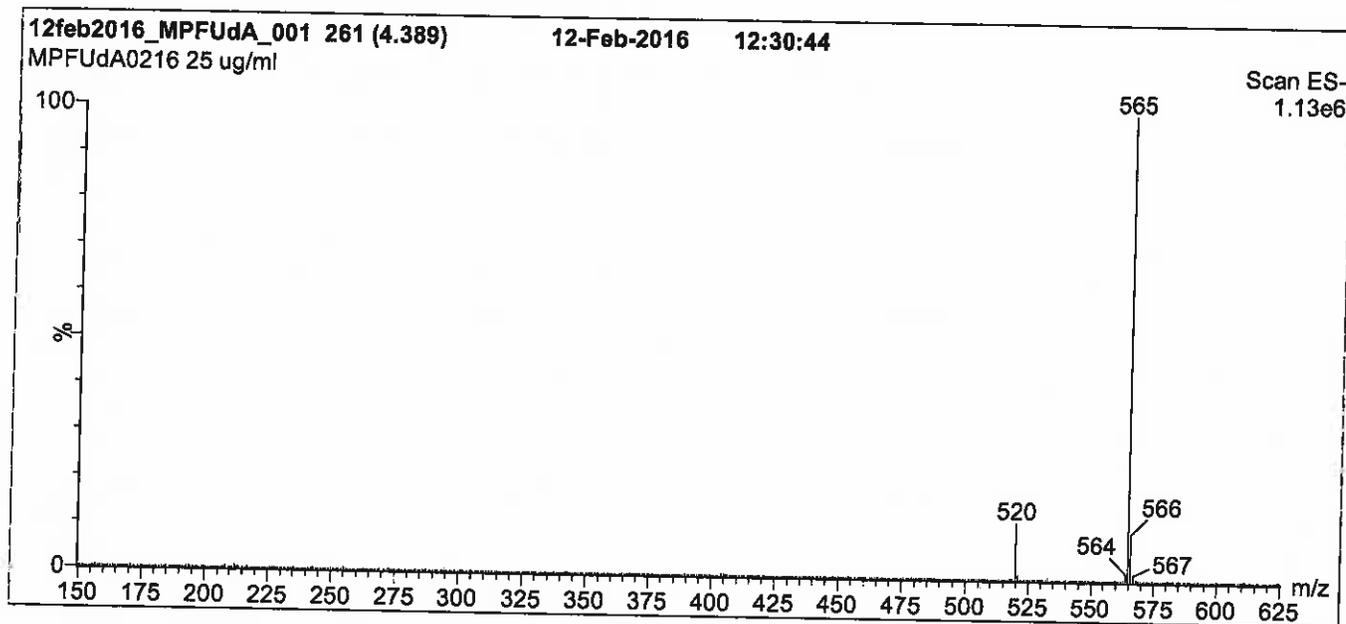
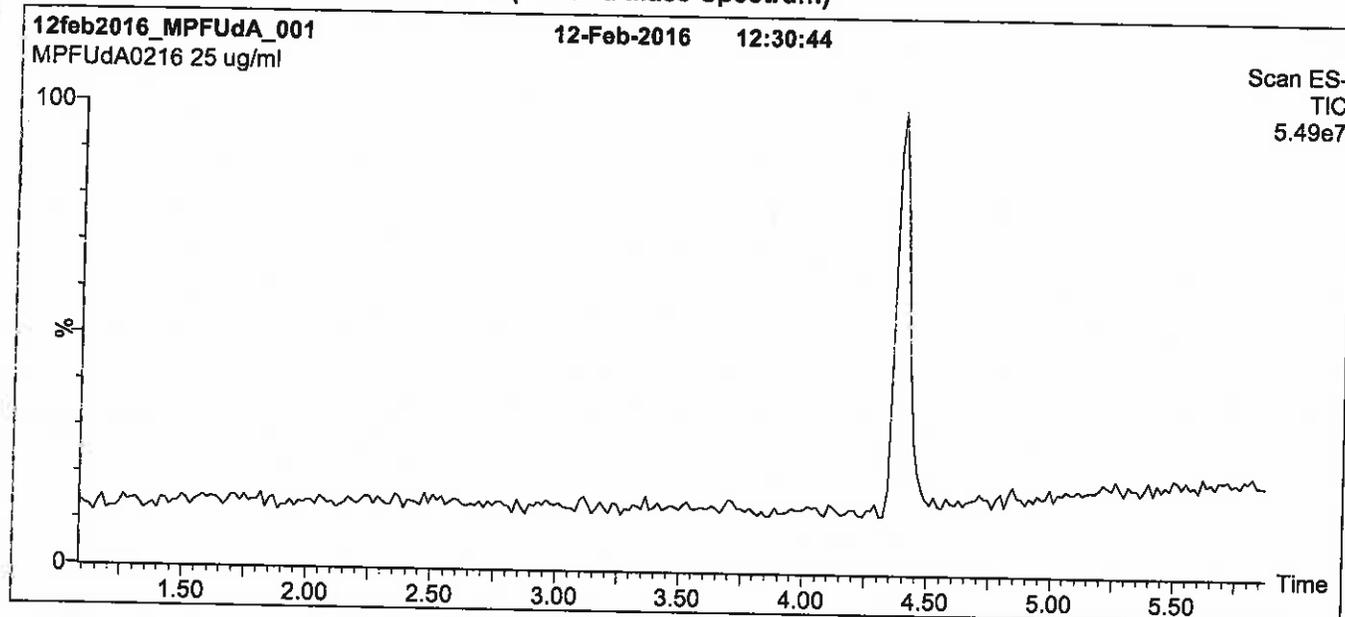
QUALITY MANAGEMENT:

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Figure 1: MPFUdA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 1.5 min before returning to initial conditions in 0.5 min.
 Time: 10 min

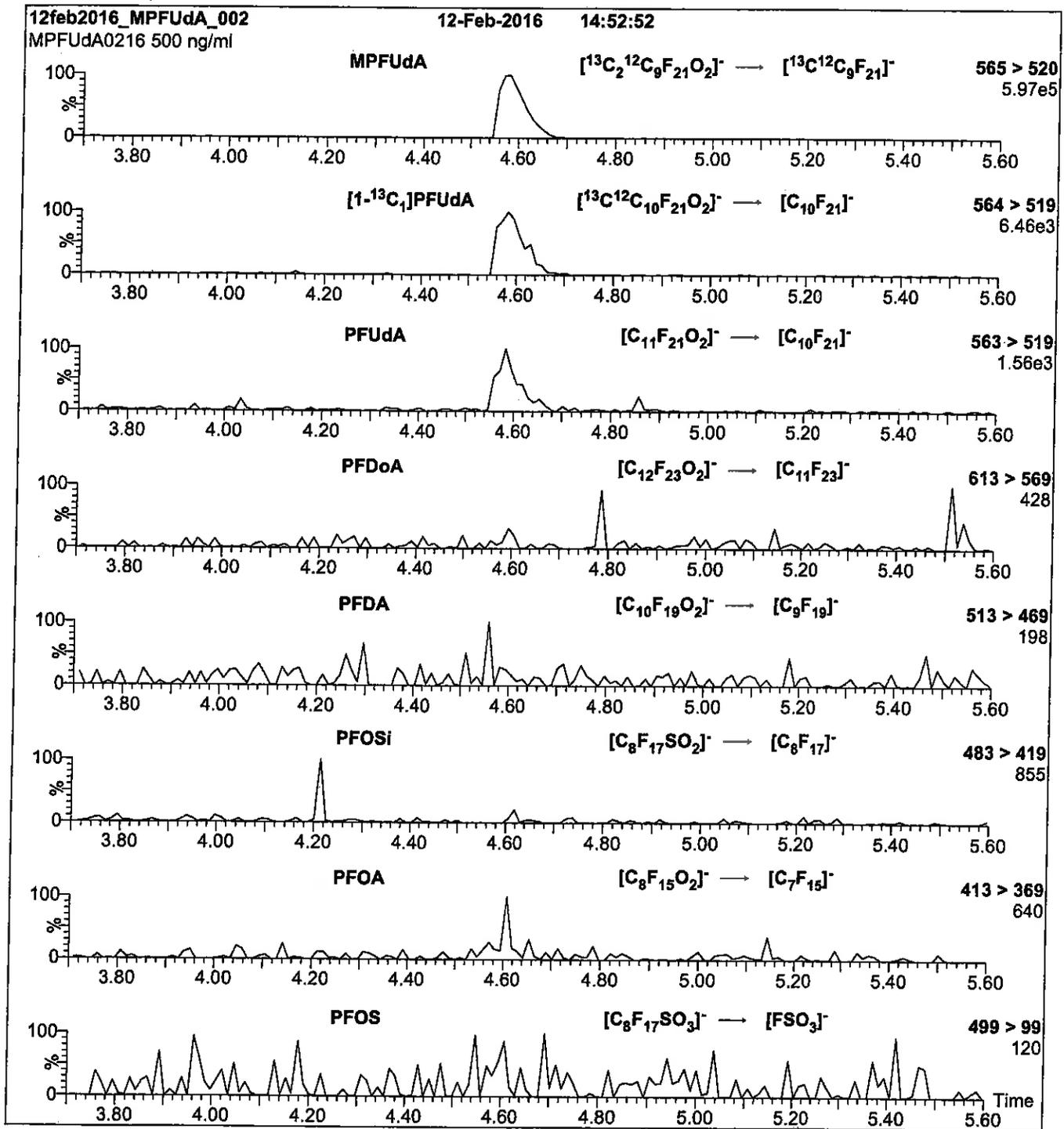
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 65
 Desolvation Gas Flow (l/hr) = 750

Figure 2: MPFUdA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml MPFUdA)

Mobile phase: Isocratic 80% MeOH / 20% H_2O

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.35e-3
Collision Energy (eV) = 11

Reagent

LCN-EtFOSA-M_00002

P: 7/16/15 SW



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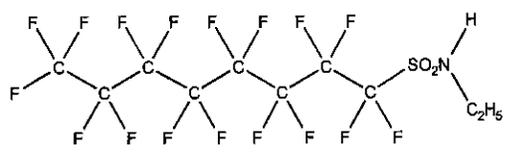
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-EtFOSA-M
COMPOUND: N-ethylperfluoro-1-octanesulfonamide

LOT NUMBER: NEtFOSA0714M

STRUCTURE:

CAS #: 4151-50-2



MOLECULAR FORMULA: C₁₀H₆F₁₇NO₂S
CONCENTRATION: 50 ± 2.5 µg/ml
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 07/14/2014
EXPIRY DATE: (mm/dd/yyyy) 07/14/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 527.20
SOLVENT(S): Methanol

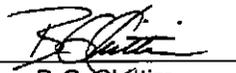
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ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim

Date: 04/01/2015
(mm/dd/yyyy)

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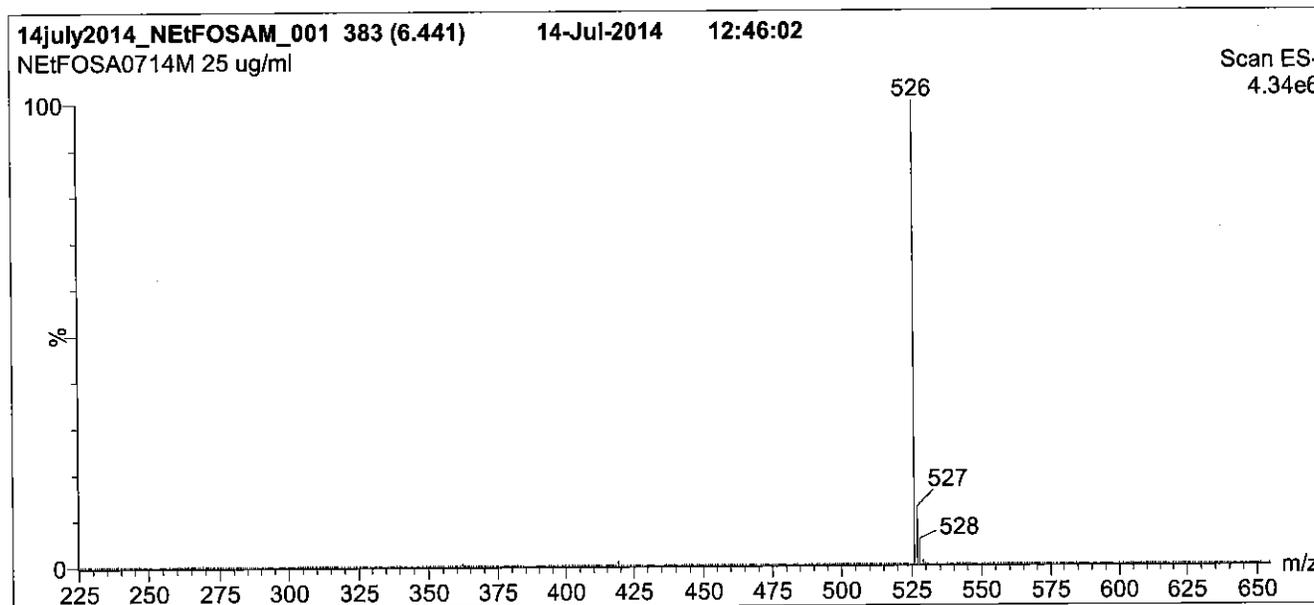
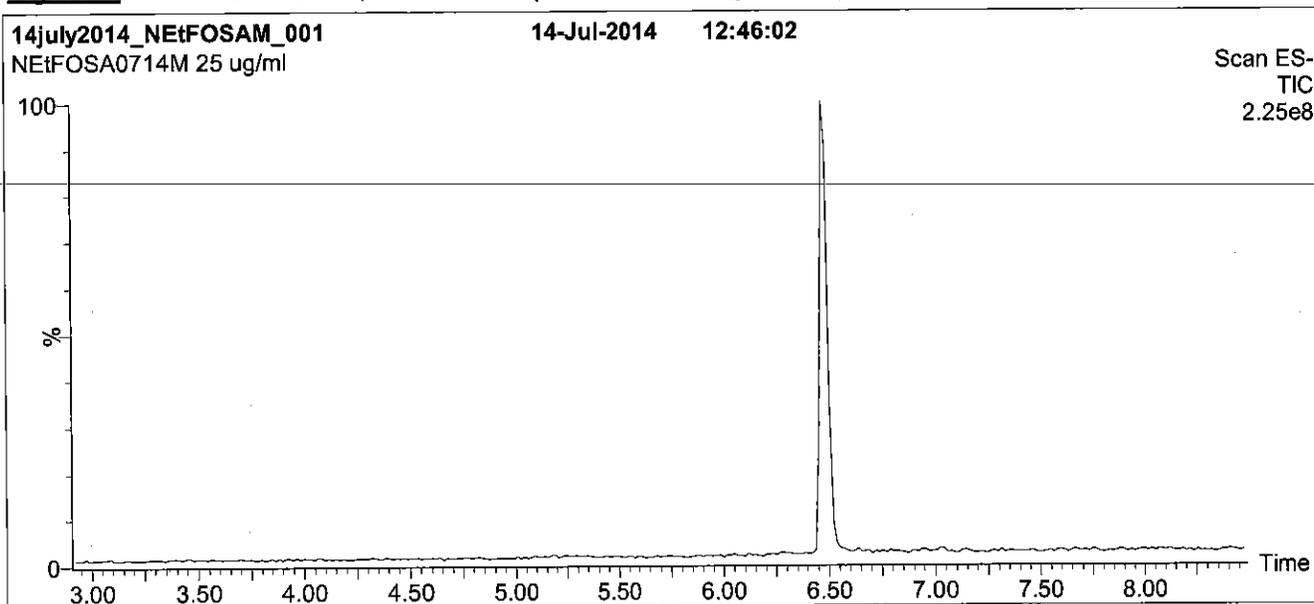
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Figure 1: N-EtFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 45% H₂O / 55% (80:20 MeOH:ACN)
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

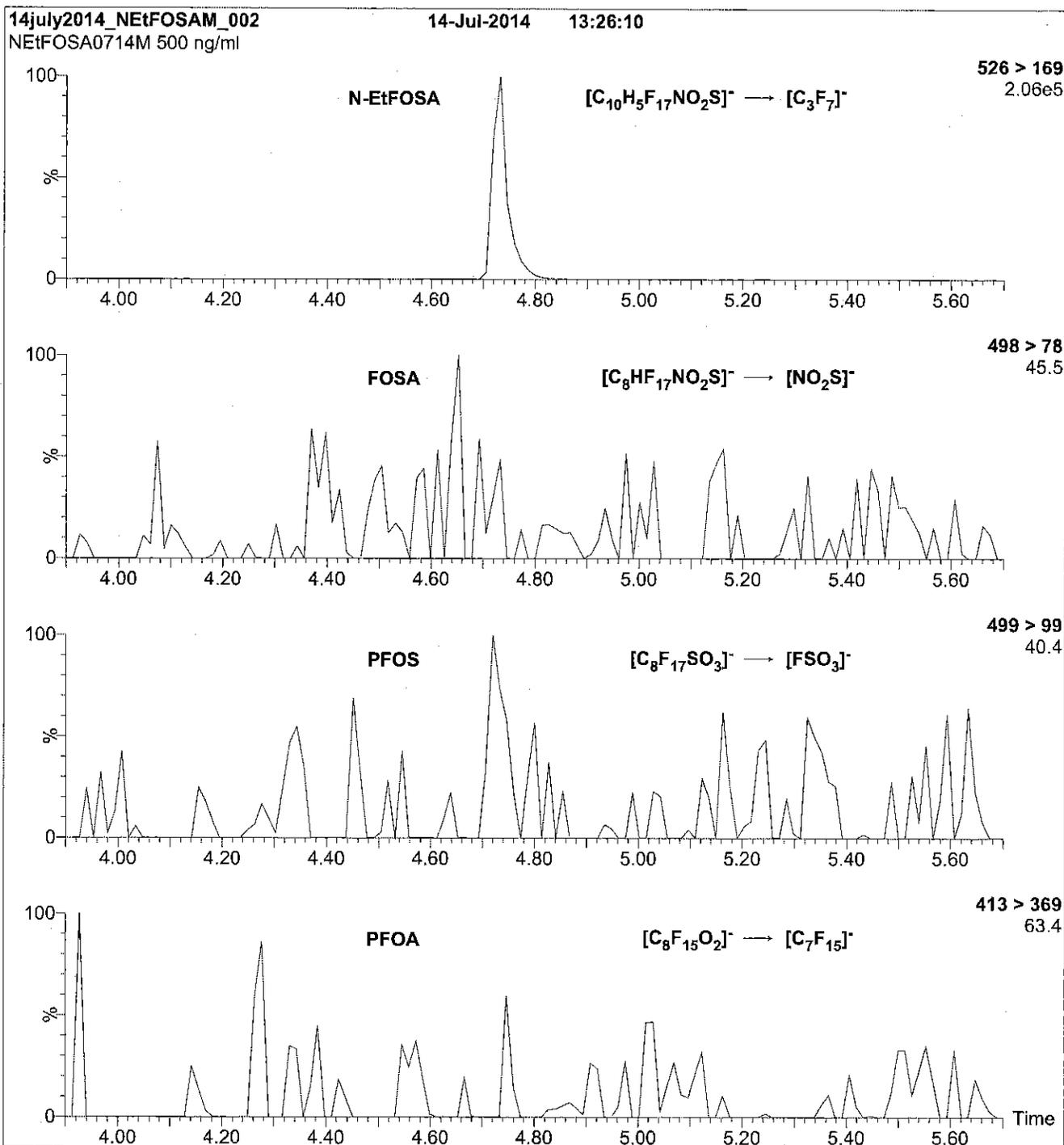
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 950 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.50
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-EtFOA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml N-EtFOA-M)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.50e-3
 Collision Energy (eV) = 30

Reagent

LCN-EtFOSA-M_00003

R: 8/23/16 SBC



715563
ID: LCN-EtFOSA-M_00003
Exp: 05/24/21 Prpt: SBC
N-EtFOSA-M

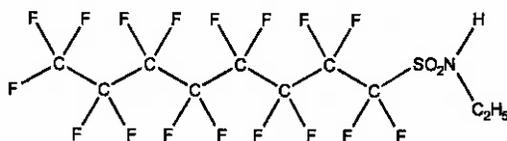


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CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-EtFOSA-M **LOT NUMBER:** NEtFOSA0516M
COMPOUND: N-ethylperfluoro-1-octanesulfonamide

STRUCTURE: **CAS #:** 4151-50-2



MOLECULAR FORMULA: C₁₀H₈F₁₇NO₂S **MOLECULAR WEIGHT:** 527.20
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 05/24/2016
EXPIRY DATE: (mm/dd/yyyy) 05/24/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 05/27/2016

(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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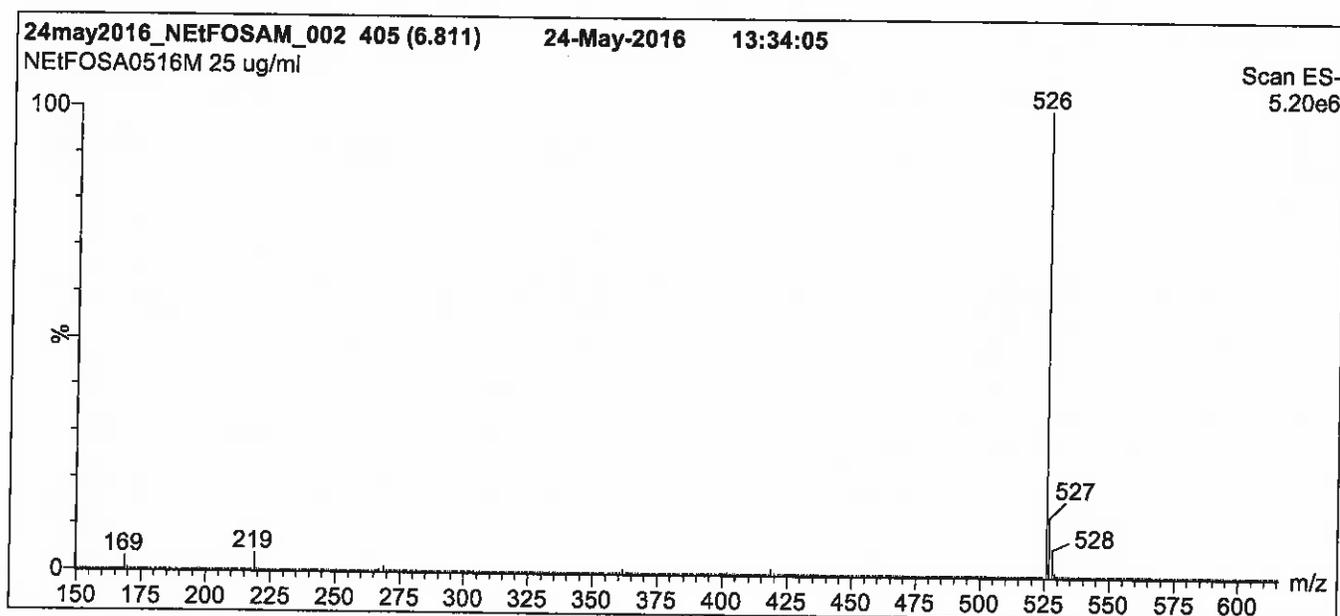
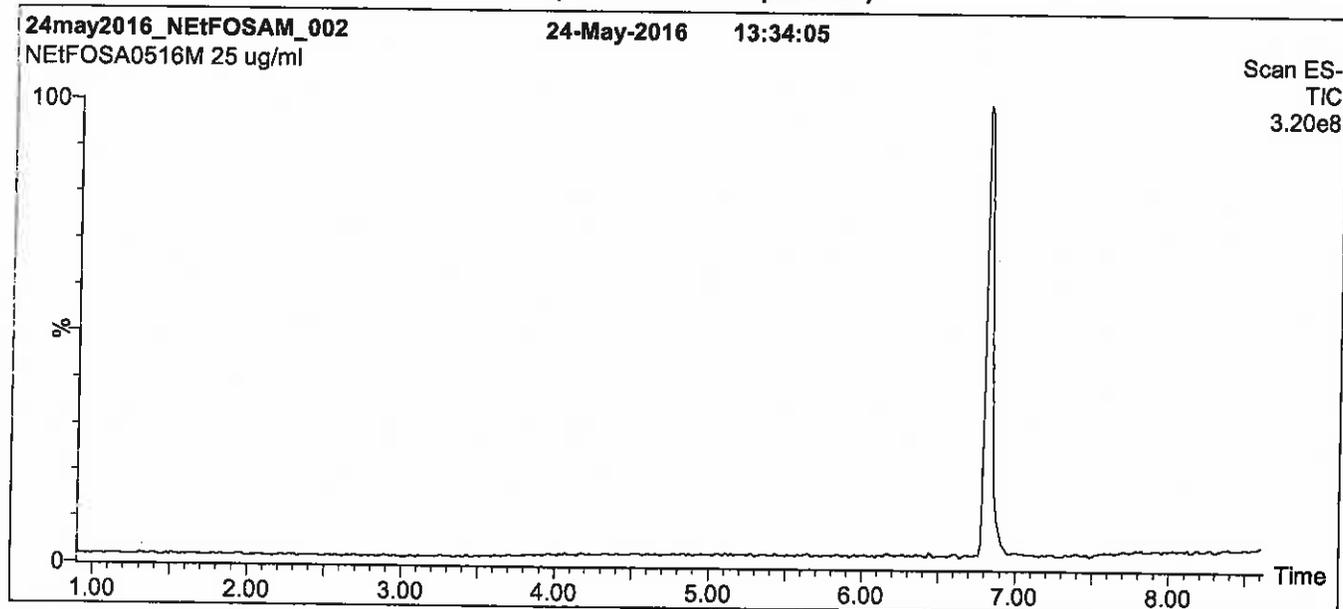
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MS: Micromass Quattro *micro* API MS

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 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 45% H₂O / 55% (80:20 MeOH:ACN)
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 Ramp to 90% organic over 7.5 min and hold for 1.5 min before returning to initial conditions in 0.5 min.
 Time: 10 min

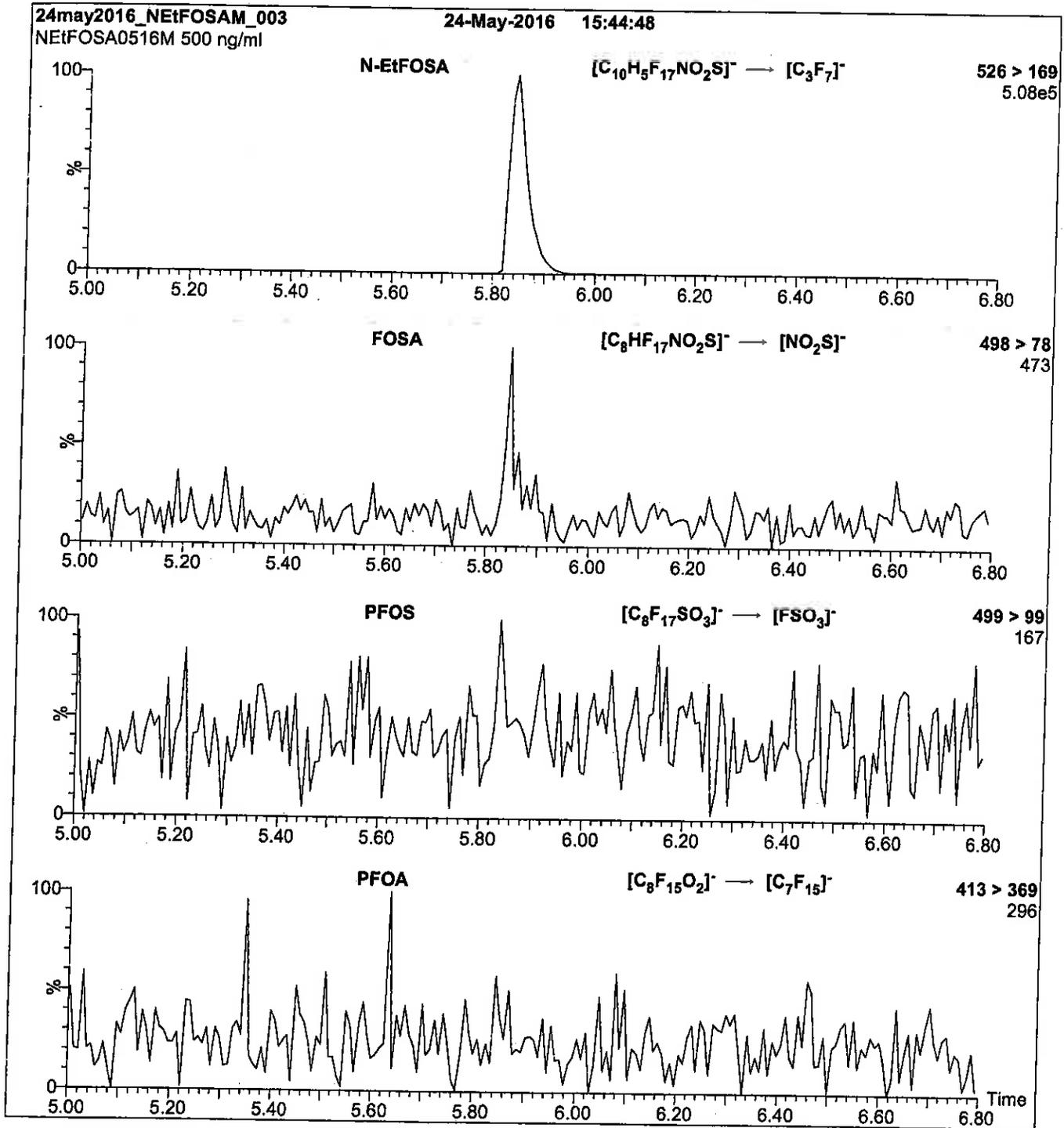
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.50
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-EtFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml N-EtFOSA-M)

MS Parameters

Collision Gas (mbar) = 3.54e-3
 Collision Energy (eV) = 30

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

Reagent

LCN-EtFOSAA_00001

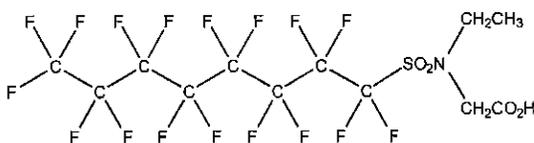


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-EtFOSAA **LOT NUMBER:** NEtFOSAA0113
COMPOUND: N-ethylperfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** 2991-50-6



MOLECULAR FORMULA: C₁₂H₈F₁₇NO₄S
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 585.23
SOLVENT(S): Methanol
 Water (<1%)

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/29/2013
EXPIRY DATE: (mm/dd/yyyy) 01/29/2018
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 04/06/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

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UNCERTAINTY:

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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

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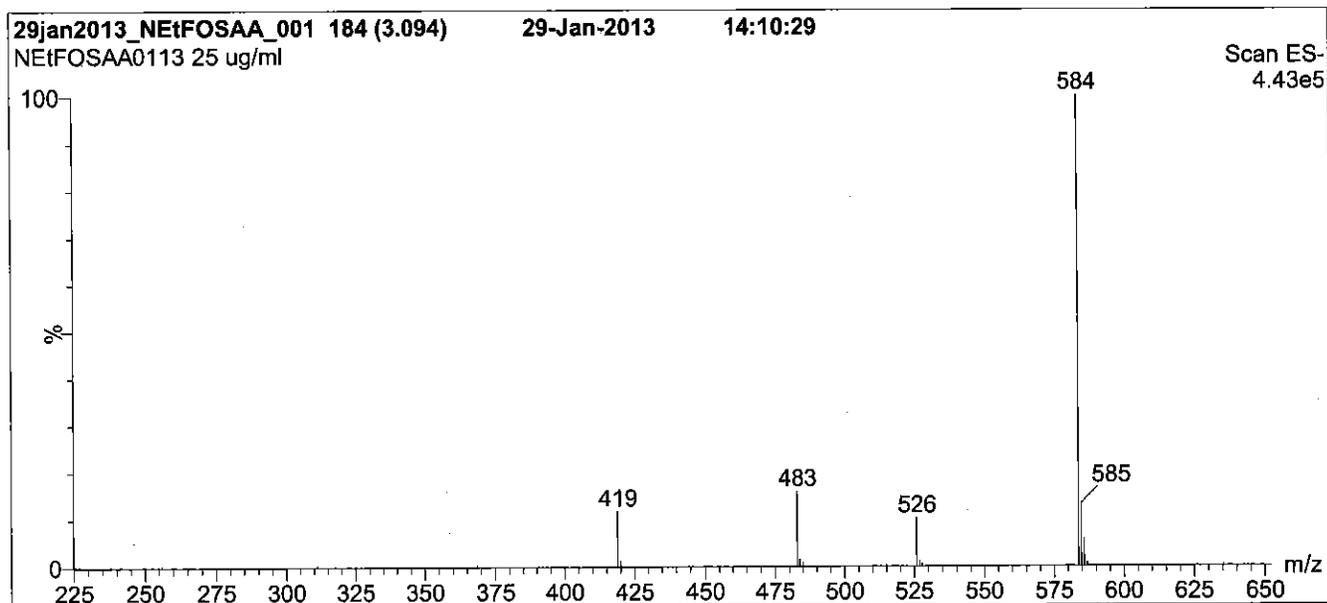
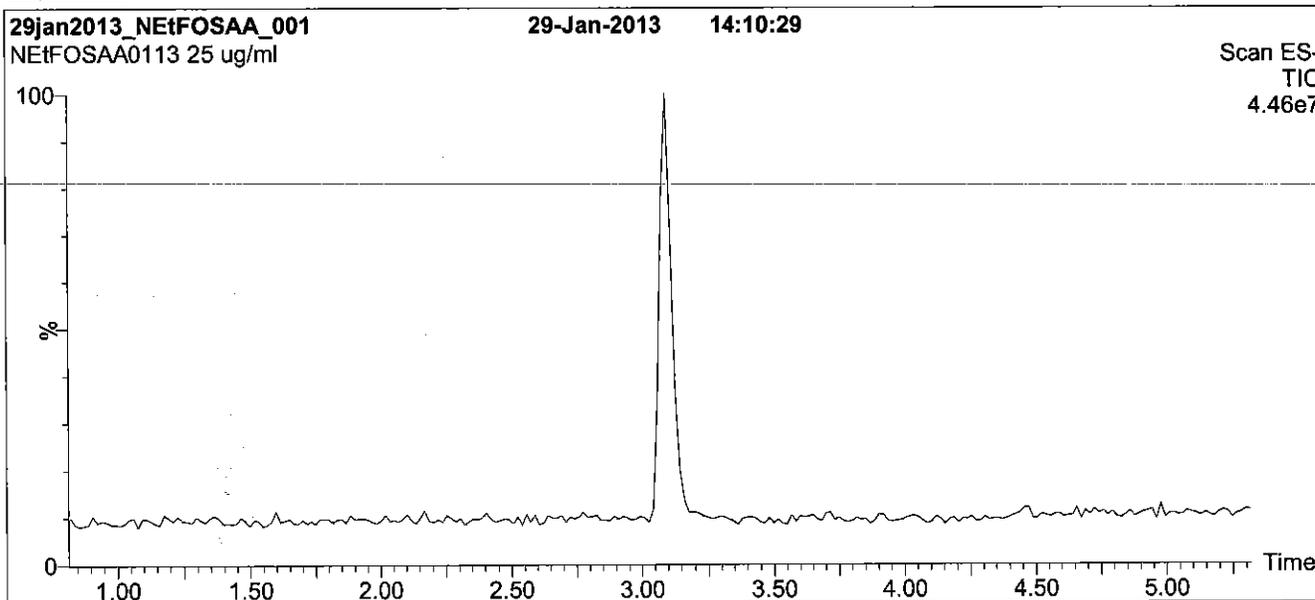
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: N-EtFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 65% (80:20 MeOH:ACN) / 35% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

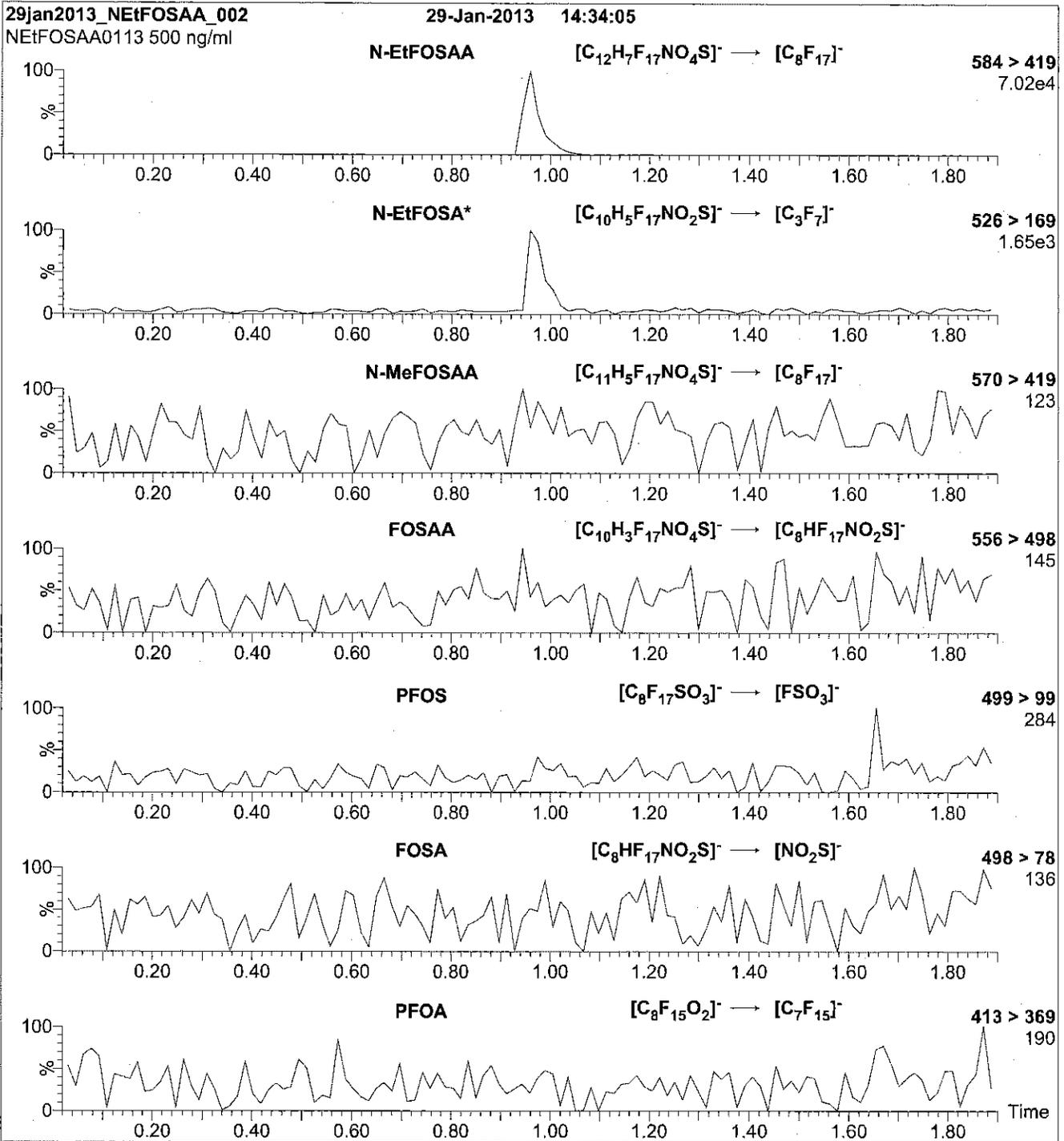
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-EtFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Note: N-EtFOSA is formed by fragmentation of N-EtFOSAA.

Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml N-EtFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
 Collision Energy (eV) = 25

Reagent

LCN-ETFOSAA_00002

R: 8/23/16 SBC



715561
ID: LCN-EiFOSAA_00002
Exp: 01/2021 Pp# 98C
N-EiFOSAA

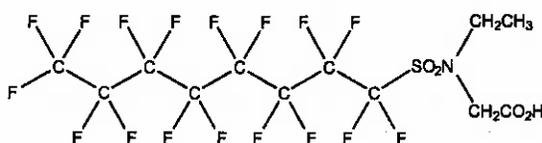


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-EtFOSAA **LOT NUMBER:** NEiFOSAA0116
COMPOUND: N-ethylperfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** 2991-50-6



MOLECULAR FORMULA:	C ₁₂ H ₈ F ₁₇ NO ₄ S	MOLECULAR WEIGHT:	585.23
CONCENTRATION:	50 ± 2.5 µg/ml	SOLVENT(S):	Methanol Water (<1%)
CHEMICAL PURITY:	>98%		
LAST TESTED: (mm/dd/yyyy)	01/20/2016		
EXPIRY DATE: (mm/dd/yyyy)	01/20/2021		
RECOMMENDED STORAGE:	Refrigerate ampoule		

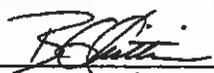
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim

Date: 01/21/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • Info@well-labs.com

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EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

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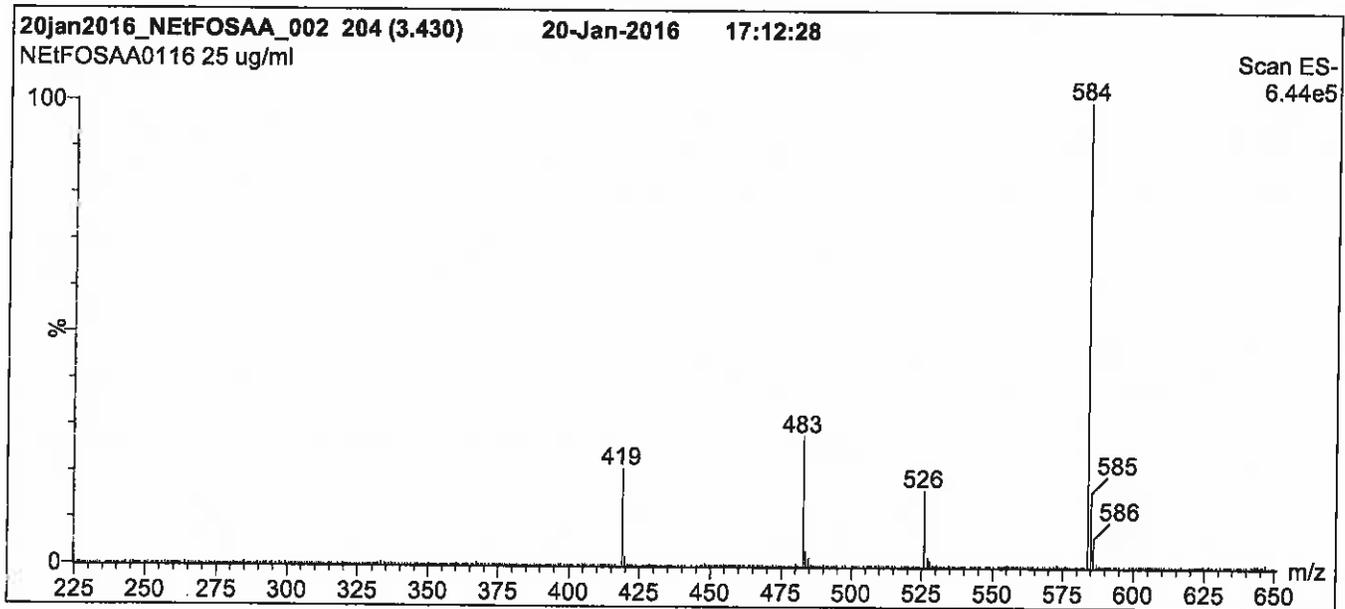
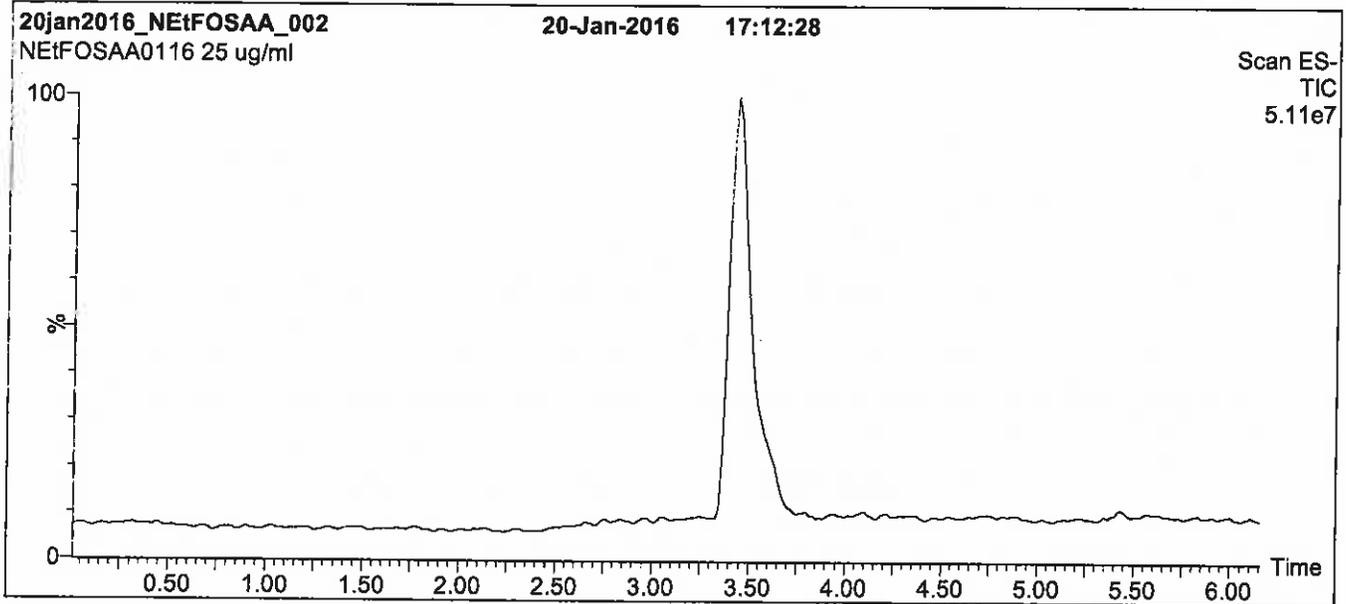
QUALITY MANAGEMENT:

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Figure 1: N-EtFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

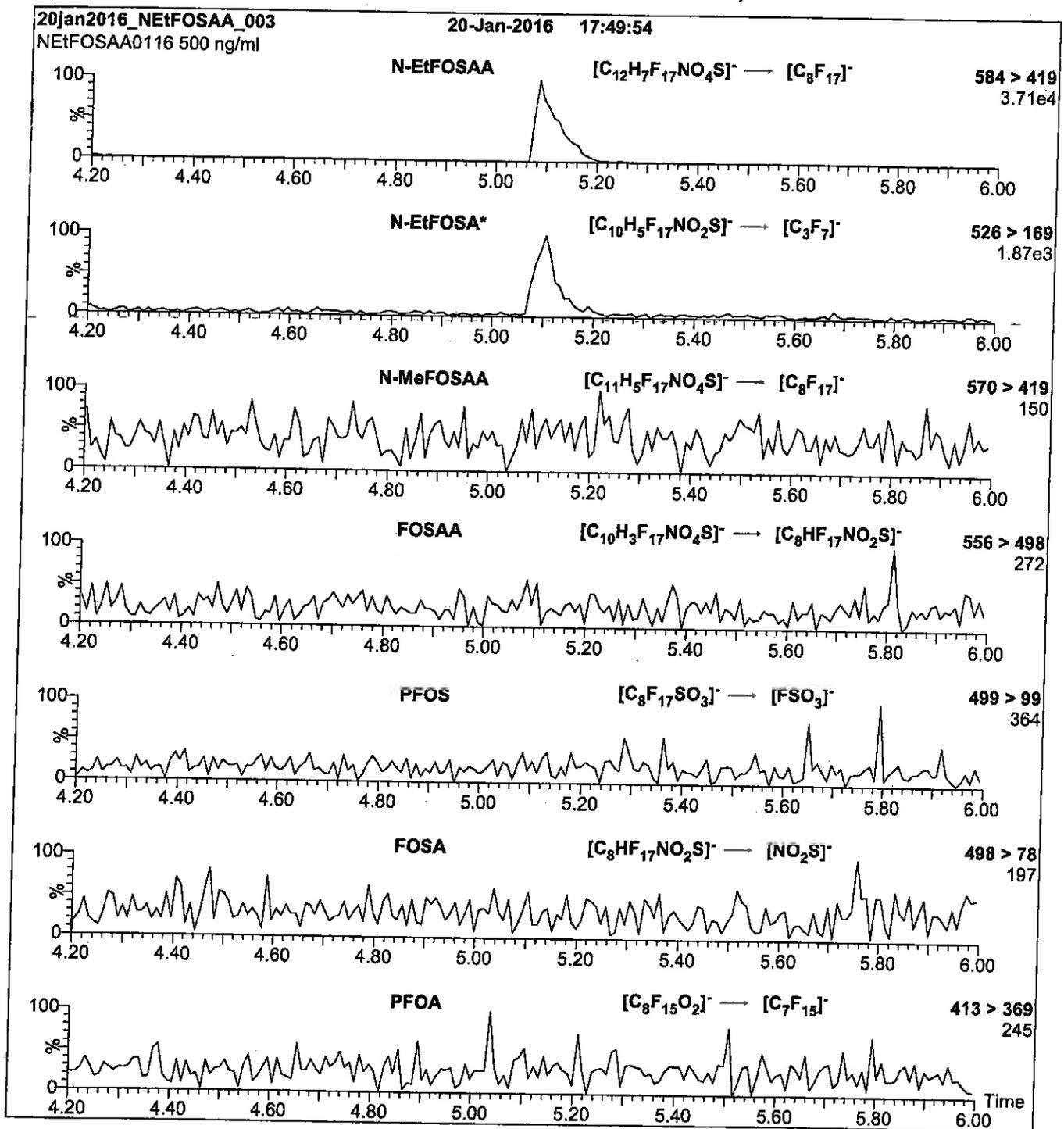
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-EtFOSAA; LC/MS/MS Data (Selected MRM Transitions)



Note: N-EtFOSA is formed by fragmentation of N-EtFOSAA.

Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml N-EtFOSAA)

MS Parameters

Collision Gas (mbar) = 3.66e-3
Collision Energy (eV) = 25

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

Reagent

LCN-MeFOSA-M_00001

V: 7/16/15 SPW



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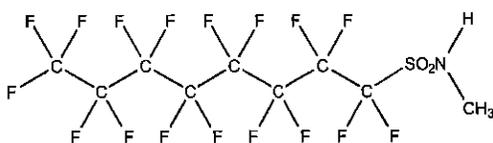
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-MeFOSA-M
COMPOUND: N-methylperfluoro-1-octanesulfonamide

LOT NUMBER: NMeFOSA0714M

STRUCTURE:

CAS #: 31506-32-8



MOLECULAR FORMULA: C₉H₄F₁₇NO₂S
CONCENTRATION: 50 ± 2.5 µg/ml
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 07/15/2014
EXPIRY DATE: (mm/dd/yyyy) 07/15/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 513.17
SOLVENT(S): Methanol

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 04/01/2015

(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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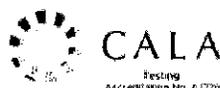
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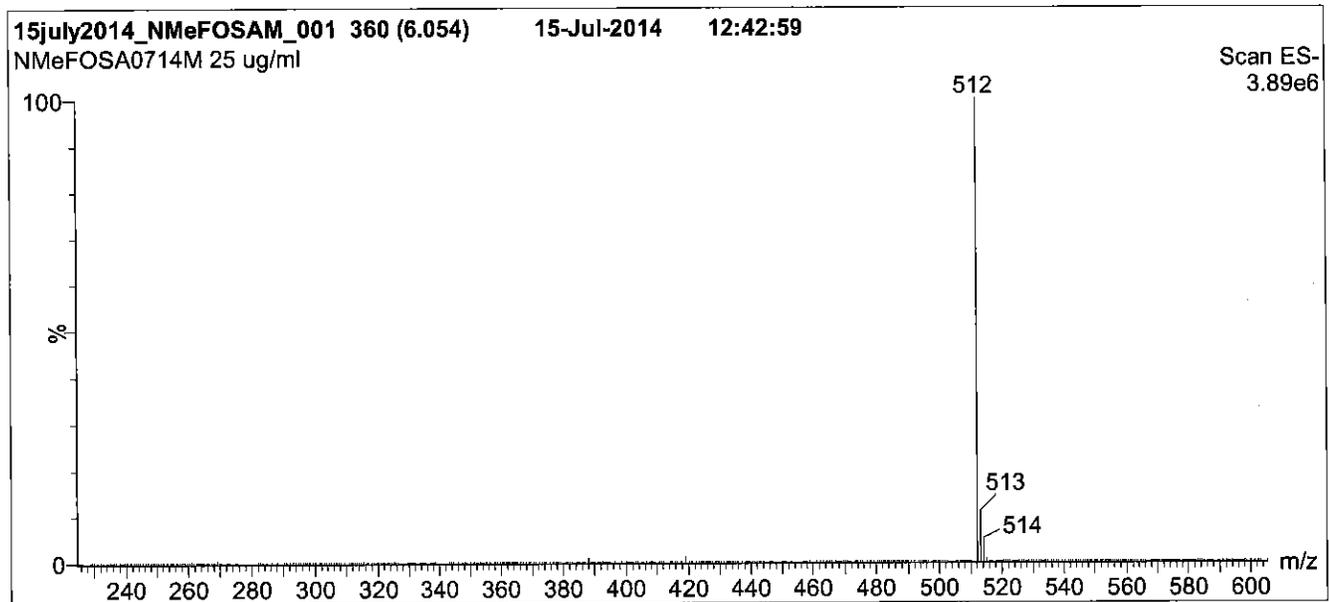
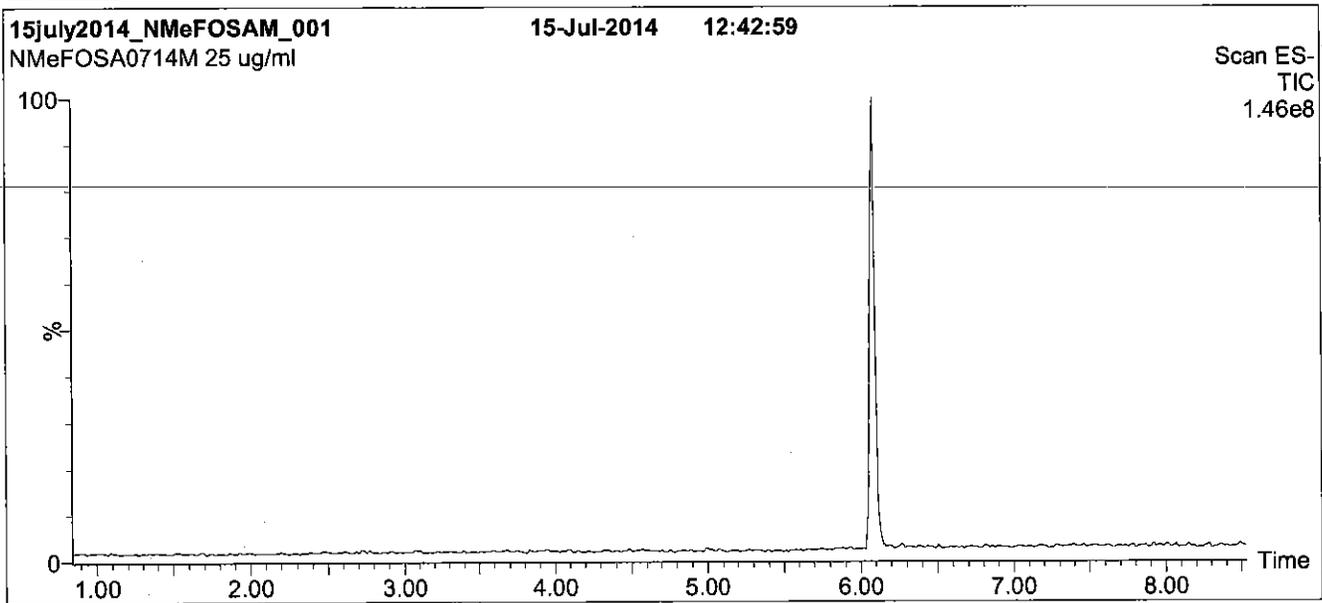
QUALITY MANAGEMENT:

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Figure 1: N-MeFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 45% H₂O / 55% (80:20 MeOH:ACN)
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 2 min before returning to initial conditions in 0.5 min.
 Time: 10 min

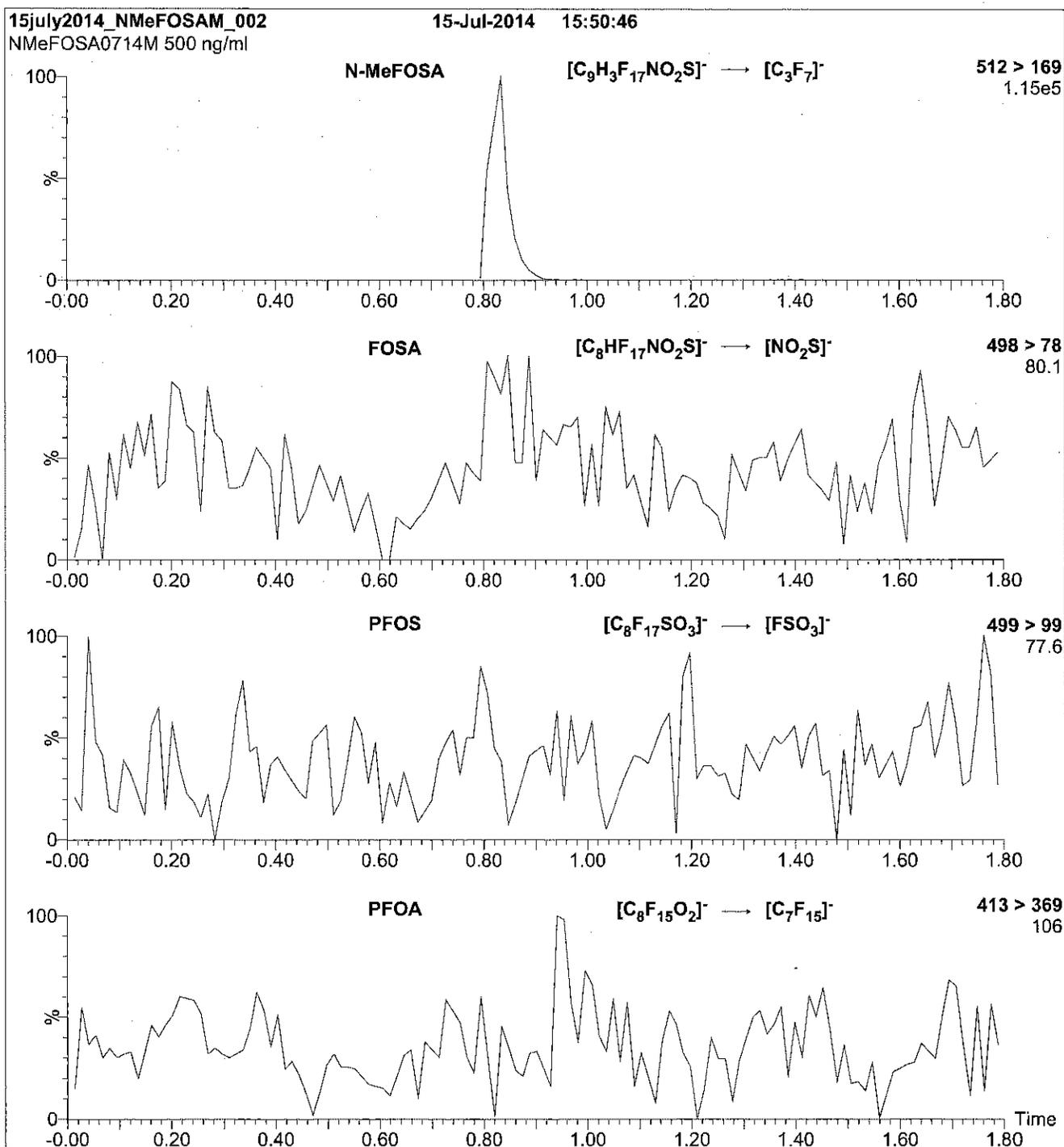
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 950 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.50
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-MeFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml N-MeFOSA-M)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 30

Reagent

LCN-MeFOSA-M_00002

R: 8/23/16 SBC



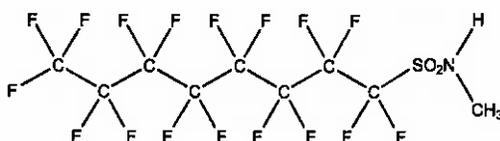
715564
ID: LCN-MeFOSA-M_00002
Exp: 05/24/21 Pppl: SBC
N-MeFOSA-M



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-MeFOSA-M **LOT NUMBER:** NMeFOSA0516M
COMPOUND: N-methylperfluoro-1-octanesulfonamide
STRUCTURE: **CAS #:** 31506-32-8



MOLECULAR FORMULA: C₉H₄F₁₇NO₂S **MOLECULAR WEIGHT:** 513.17
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 05/24/2016
EXPIRY DATE: (mm/dd/yyyy) 05/24/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 05/26/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

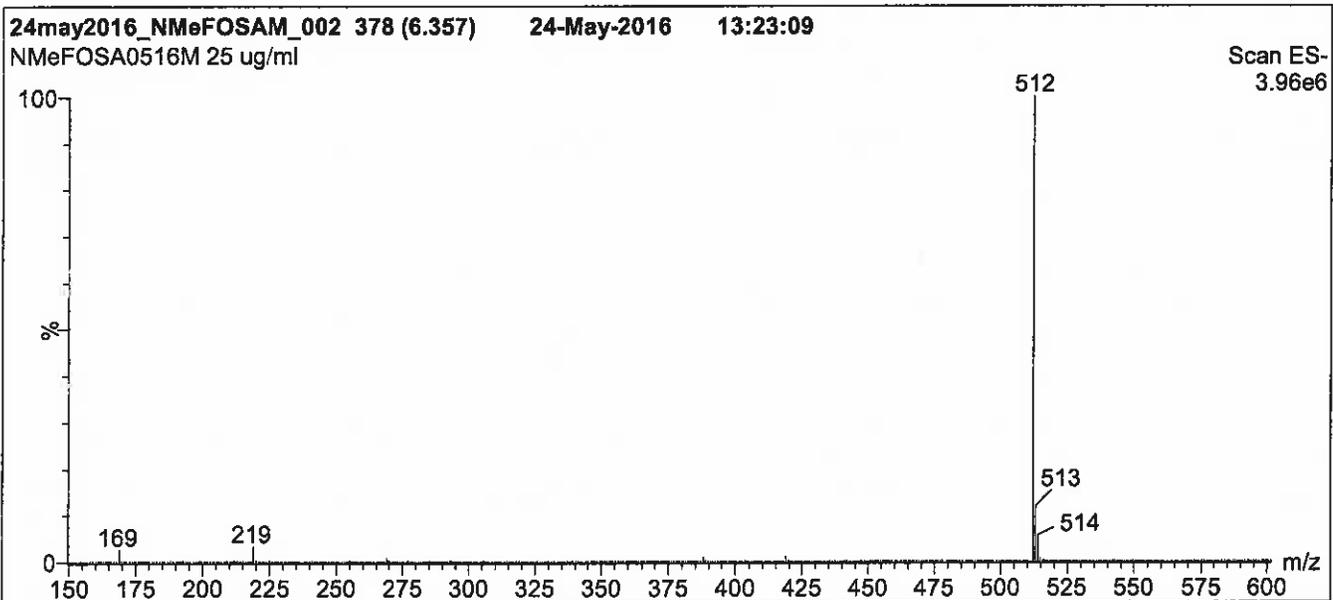
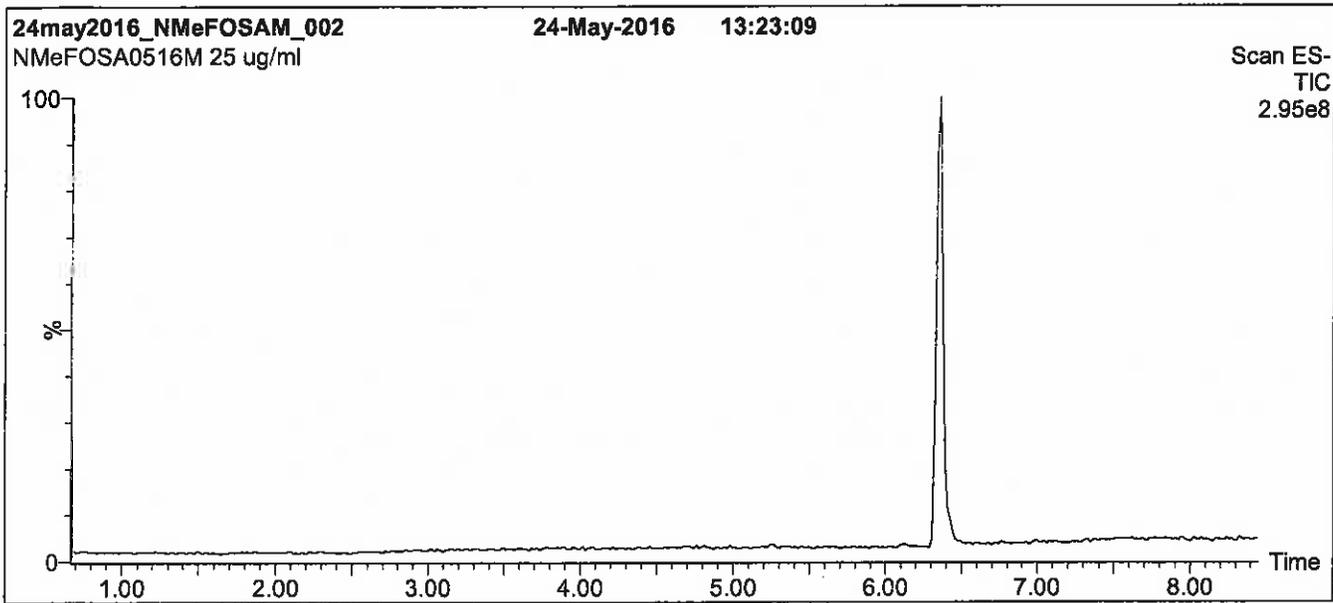
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: N-MeFOSA-M; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 45% H₂O / 55% (80:20 MeOH:ACN)
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7.5 min and hold for
 1.5 min before returning to initial conditions in 0.5 min.
 Time: 10 min

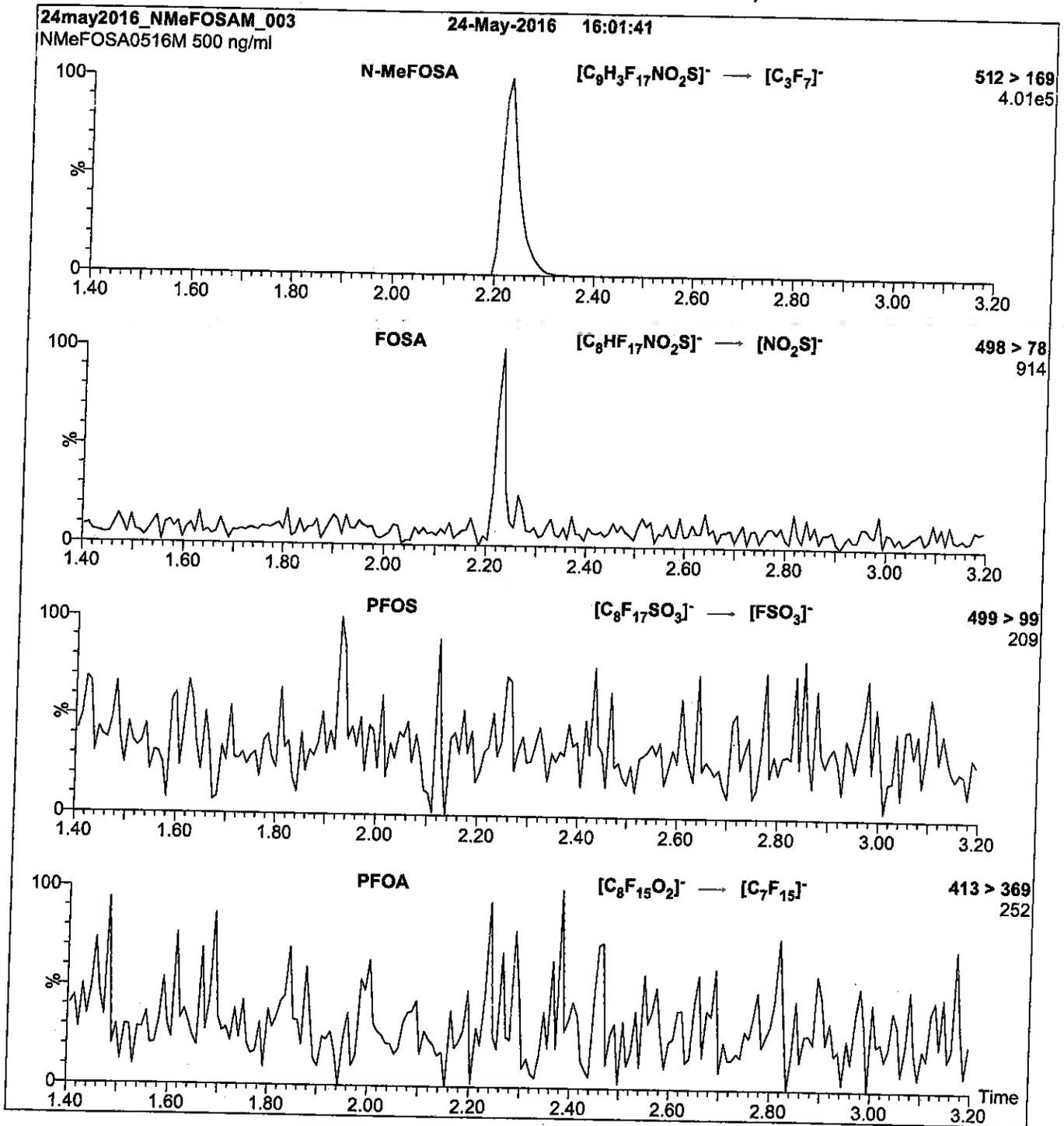
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.50
 Cone Voltage (V) = 40.00
 Core Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-MeFOSA-M; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml N-MeFOSA-M)

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 30

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

Reagent

LCN-MeFOSAA_00001

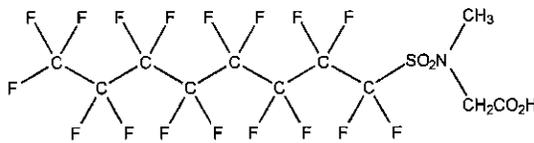


WELLINGTON
LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

PRODUCT CODE: N-MeFOSAA **LOT NUMBER:** NMeFOSAA1214
COMPOUND: N-methylperfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** 2355-31-9



MOLECULAR FORMULA: C₁₁H₆F₁₇NO₄S **MOLECULAR WEIGHT:** 571.21
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
 Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 12/09/2014
EXPIRY DATE: (mm/dd/yyyy) 12/09/2019
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 04/06/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

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$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

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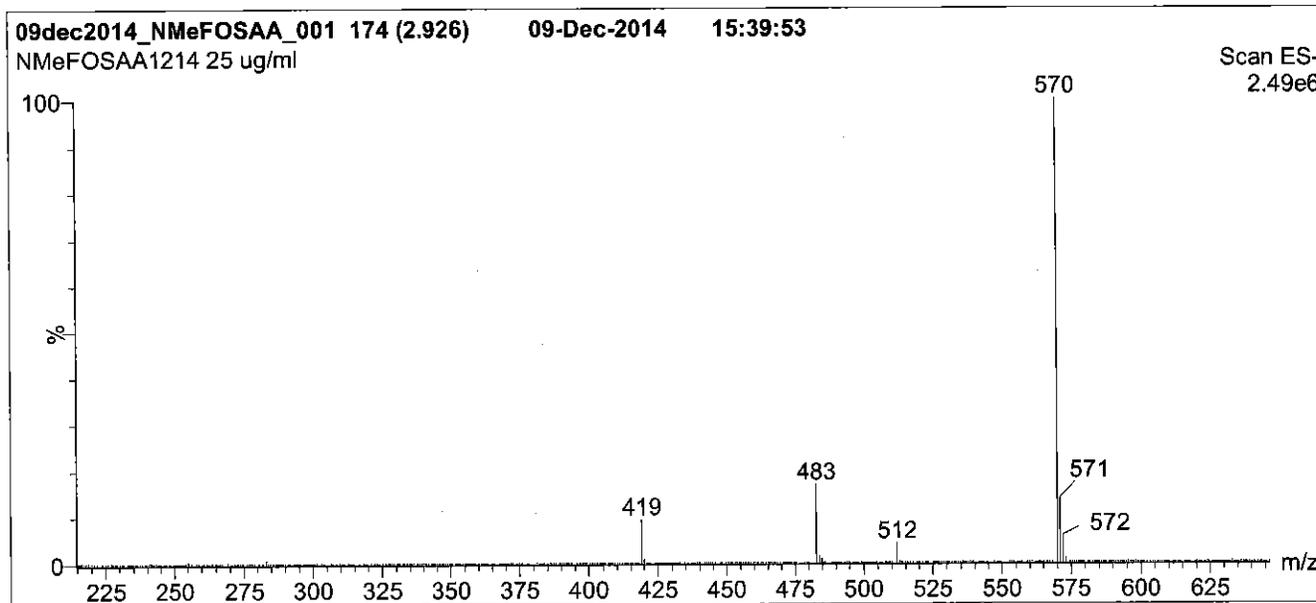
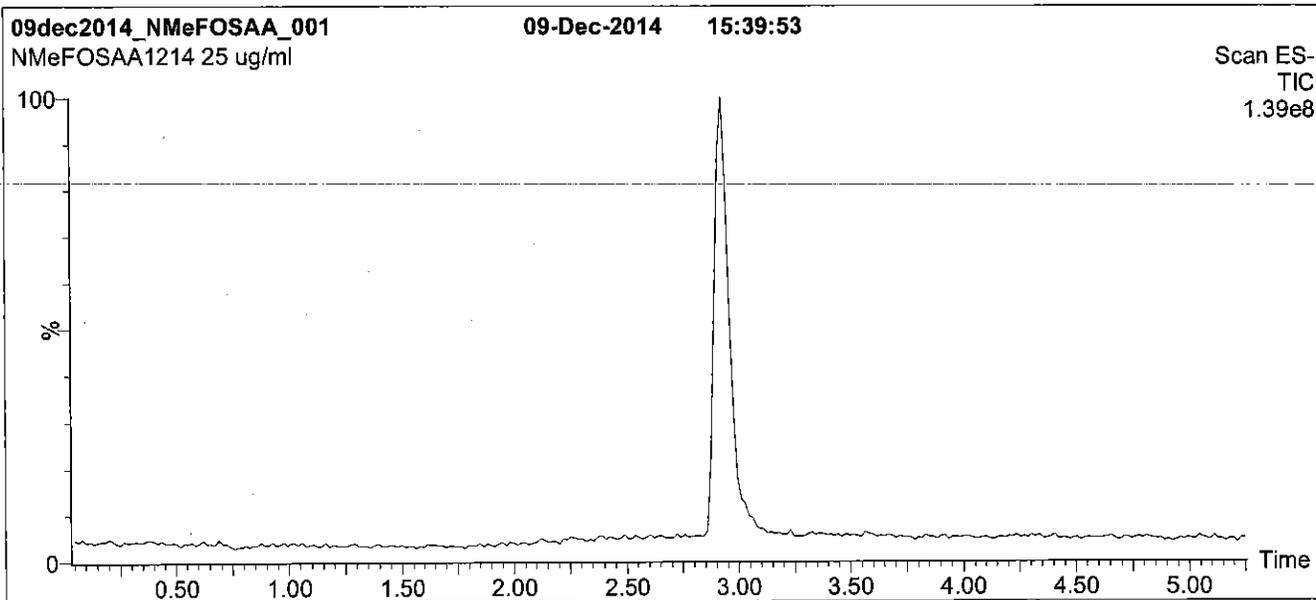
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: N-MeFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 65% (80:20 MeOH:ACN) / 35% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

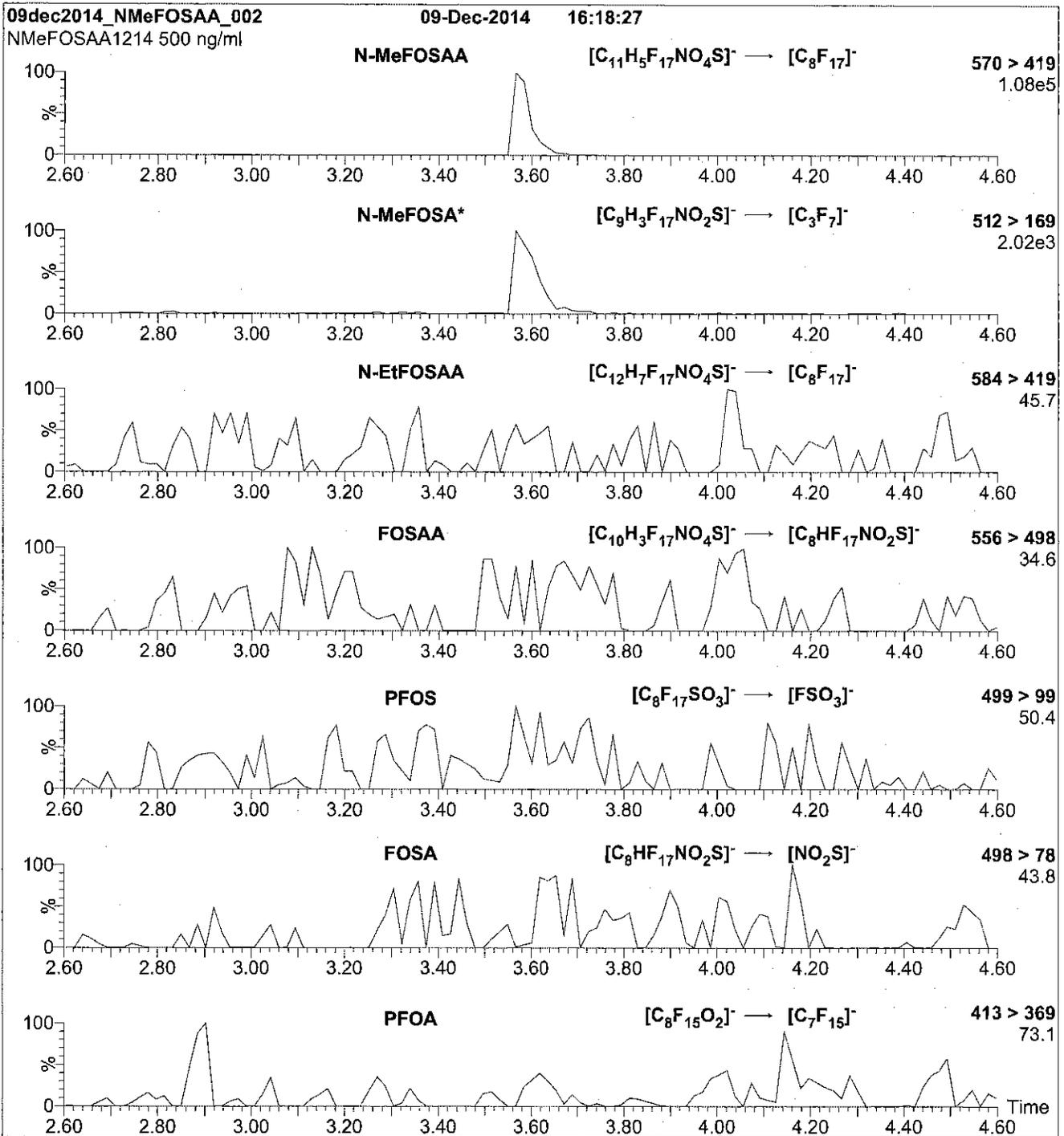
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (215 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-MeFOSAA; LC/MS/MS Data (Selected MRM Transitions)



*Note: N-MeFOSA is formed by fragmentation of N-MeFOSAA.

Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml N-MeFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
Collision Energy (eV) = 25

Reagent

LCN-MeFOSAA_00003

R: 8/23/16 *SAE*

715562
ID: LCN-MeFOSAA_00003
Exp: 01/20/21 Prpd: SEC
N-MeFOSAA

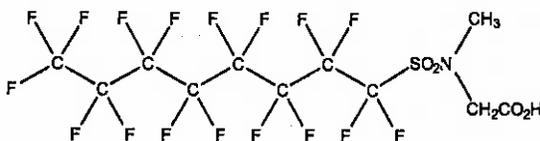


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: N-MeFOSAA **LOT NUMBER:** NMeFOSAA0116
COMPOUND: N-methylperfluoro-1-octanesulfonamidoacetic acid

STRUCTURE: **CAS #:** 2355-31-9



MOLECULAR FORMULA: C₁₁H₈F₁₇NO₄S **MOLECULAR WEIGHT:** 571.21
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/20/2016
EXPIRY DATE: (mm/dd/yyyy) 01/20/2021
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent the conversion of the acetic acid moiety to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 01/21/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

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where x is expressed as a relative standard uncertainty of the individual parameter.

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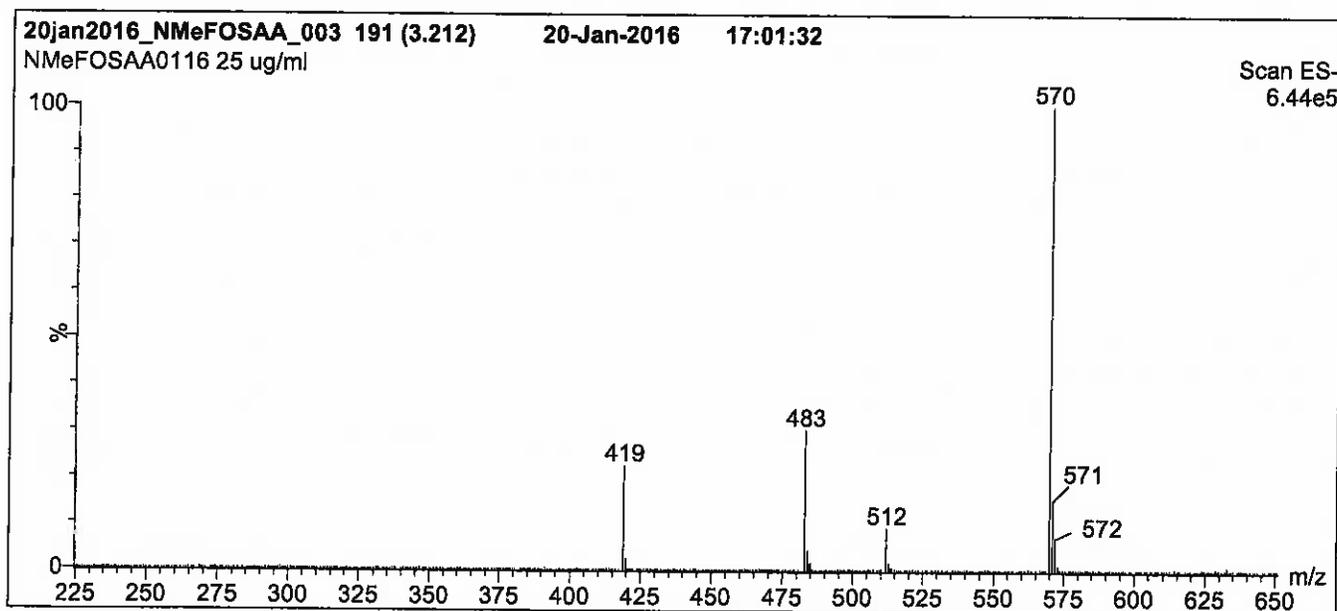
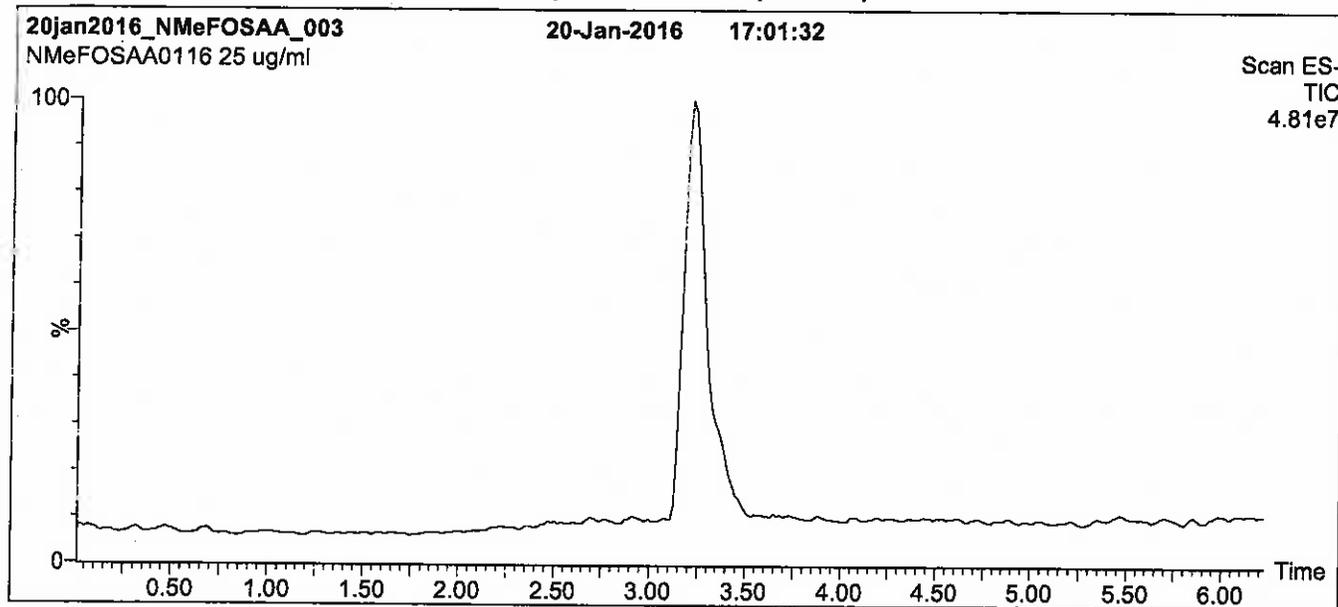
QUALITY MANAGEMENT:

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Figure 1: N-MeFOSAA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

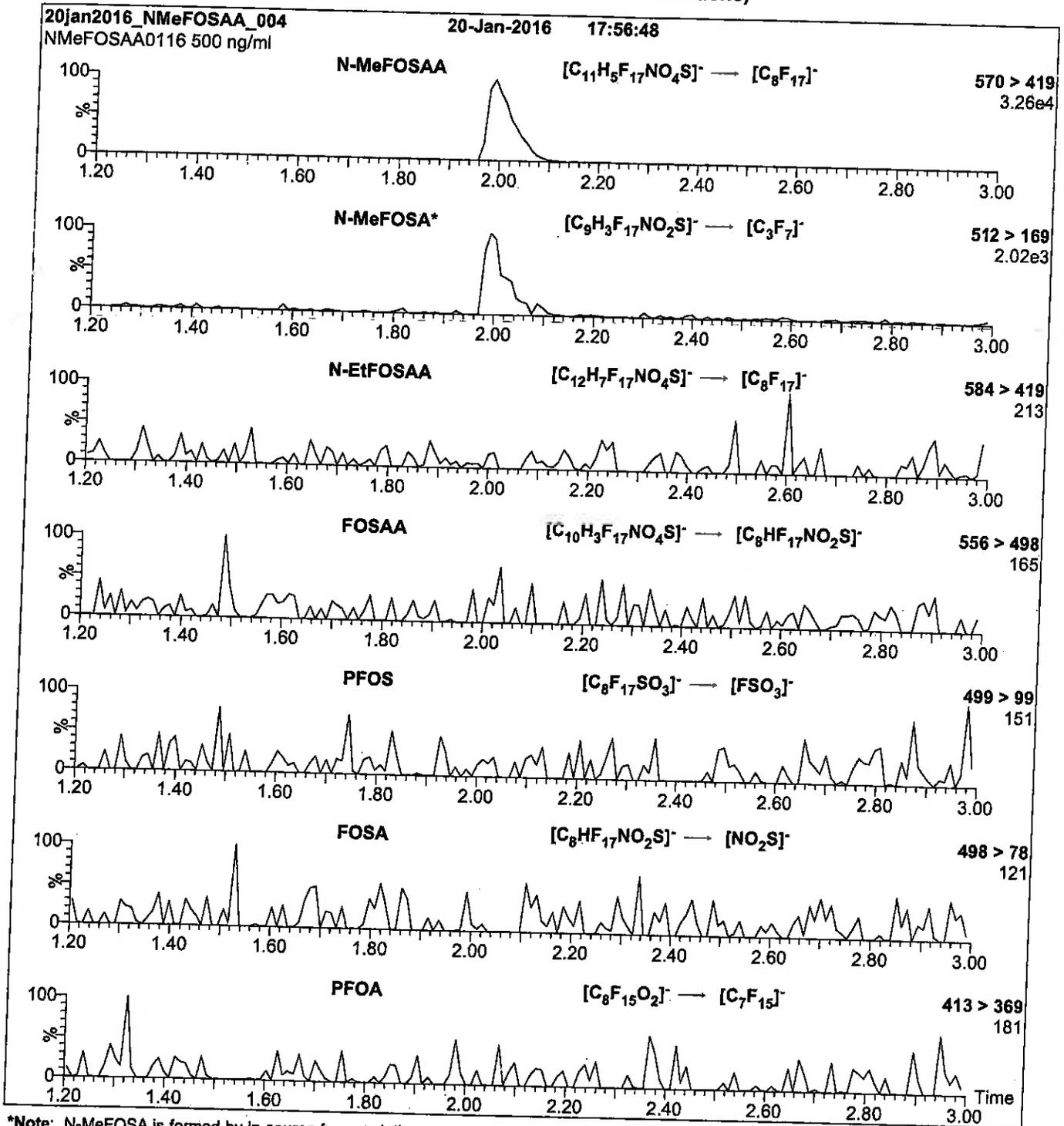
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 35.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: N-MeFOSAA; LC/MS/MS Data (Selected MRM Transitions)



*Note: N-MeFOSA is formed by in-source fragmentation.

Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml N-MeFOSAA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.66e-3
Collision Energy (eV) = 25

Reagent

LCPFACMXB_00007



WELLINGTON
LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

PFAC-MXB

**Solution/Mixture of Native
Perfluoroalkylcarboxylic Acids and
Native Perfluoroalkylsulfonates**

PRODUCT CODE: PFAC-MXB
LOT NUMBER: PFACMXB1115
SOLVENT(S): Methanol / Water (<1%)
DATE PREPARED: (mm/dd/yyyy) 11/04/2015
LAST TESTED: (mm/dd/yyyy) 11/06/2015
EXPIRY DATE: (mm/dd/yyyy) 11/06/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DESCRIPTION:

PFAC-MXB is a solution/mixture of thirteen native perfluoroalkylcarboxylic acids (C₄-C₁₄, C₁₆, and C₁₈) and four native perfluoroalkylsulfonates (C₄, C₆, C₈ and C₁₀). The full name, abbreviation and concentration for each of the components are given in Table A.

The individual perfluoroalkylcarboxylic acids and perfluoroalkylsulfonates all have chemical purities of >98%.

DOCUMENTATION/ DATA ATTACHED:

Table A: Components and Concentrations of the Solution/Mixture
 Figure 1: LC/MS Data (SiR)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)
 Figure 3: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acids to their respective methyl esters.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HAZARDS:

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SYNTHESIS / CHARACTERIZATION:

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HOMOGENEITY:

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TRACEABILITY:

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LIMITED WARRANTY:

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QUALITY MANAGEMENT:

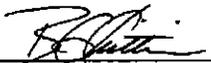
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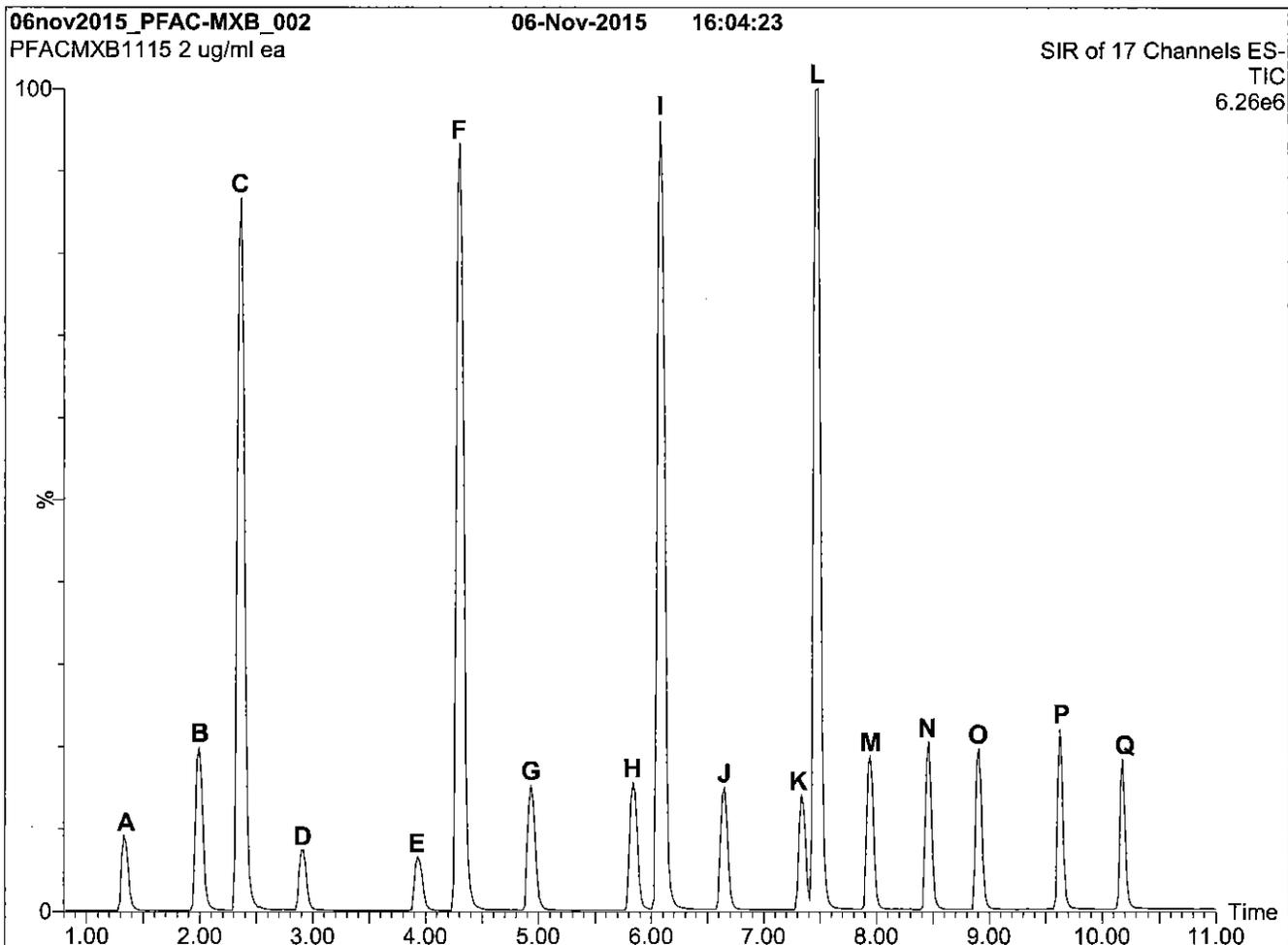
Table A: PFAC-MXB; Components and Concentrations (ng/ml, ± 5% in Methanol / Water (<1%))

Name	Abbreviation	Concentration (ng/ml)		Peak Assignment in Figure 1
		as the salt	as the anion	
Perfluoro-n-butanoic acid	PFBA	2000		A
Perfluoro-n-pentanoic acid	PFPeA	2000		B
Perfluoro-n-hexanoic acid	PFHxA	2000		D
Perfluoro-n-heptanoic acid	PFHpA	2000		E
Perfluoro-n-octanoic acid	PFOA	2000		G
Perfluoro-n-nonanoic acid	PFNA	2000		H
Perfluoro-n-decanoic acid	PFDA	2000		J
Perfluoro-n-undecanoic acid	PFUdA	2000		K
Perfluoro-n-dodecanoic acid	PFDoA	2000		M
Perfluoro-n-tridecanoic acid	PFTrDA	2000		N
Perfluoro-n-tetradecanoic acid	PFTeDA	2000		O
Perfluoro-n-hexadecanoic acid	PFHxDA	2000		P
Perfluoro-n-octadecanoic acid	PFODA	2000		Q
Name	Abbreviation	Concentration (ng/ml)		Peak Assignment in Figure 1
		as the salt	as the anion	
Potassium perfluoro-1-butanesulfonate	L-PFBS	2000	1770	C
Sodium perfluoro-1-hexanesulfonate	L-PFHxS	2000	1890	F
Sodium perfluoro-1-octanesulfonate	L-PFOS	2000	1910	I
Sodium perfluoro-1-decanesulfonate	L-PFDS	2000	1930	L

Certified By: 
B.G. Chittim

Date: 11/11/2015
(mm/dd/yyyy)

Figure 1: PFAC-MXB; LC/MS Data (Total Ion Current Chromatogram; SIR)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 55% H₂O / 45% (80:20 MeOH:ACN)
(both with 10 mM NH₄OAc buffer)
Ramp to 95% organic over 10 min and hold for 1 min
before returning to initial conditions in 0.5 min.

Time: 12 min

Flow: 300 μ l/min

MS Parameters

Experiment: SIR of 17 Channels

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = variable (10-70)
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFAC-MXB; LC/MS/MS Data (Selected MRM Transitions)

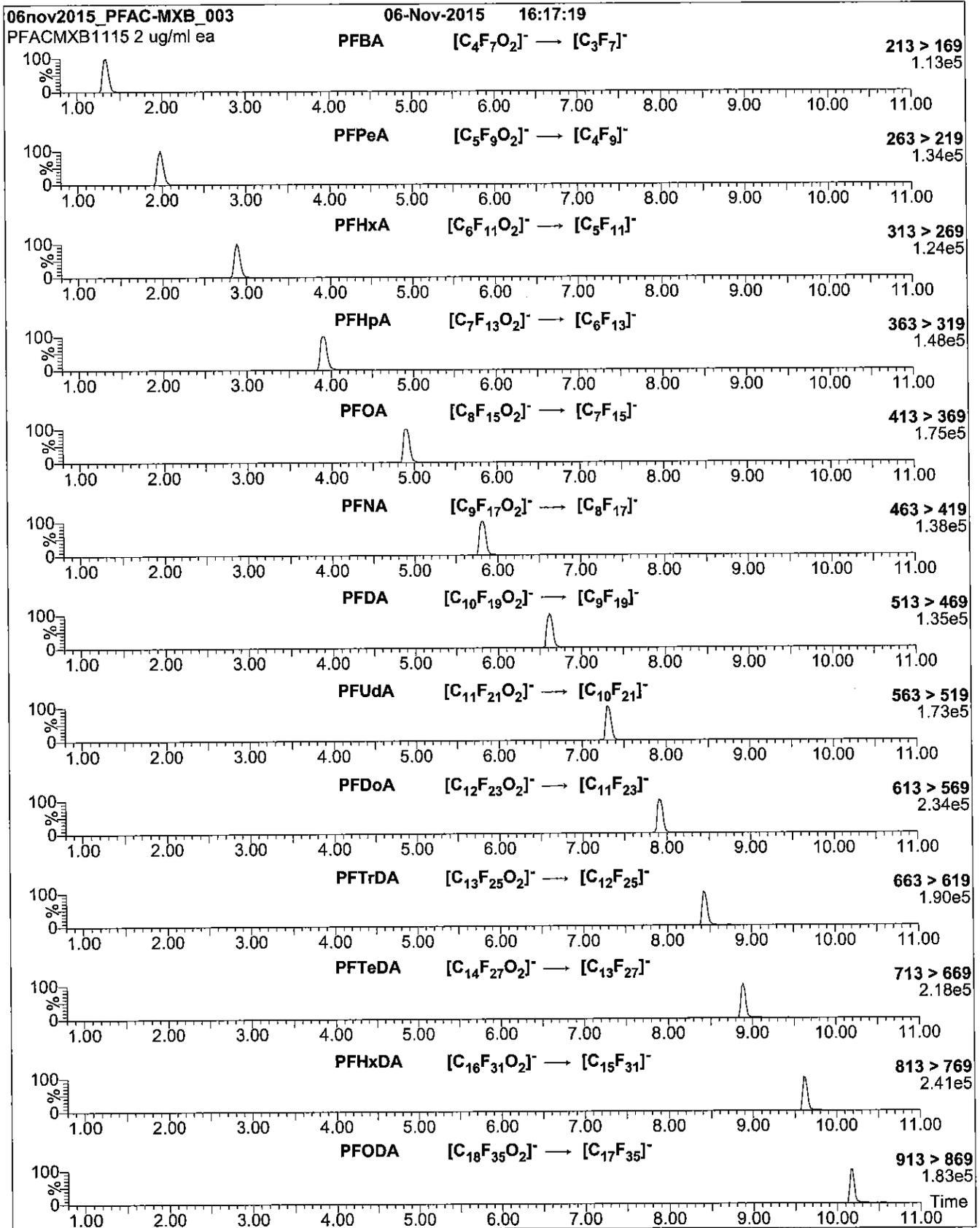
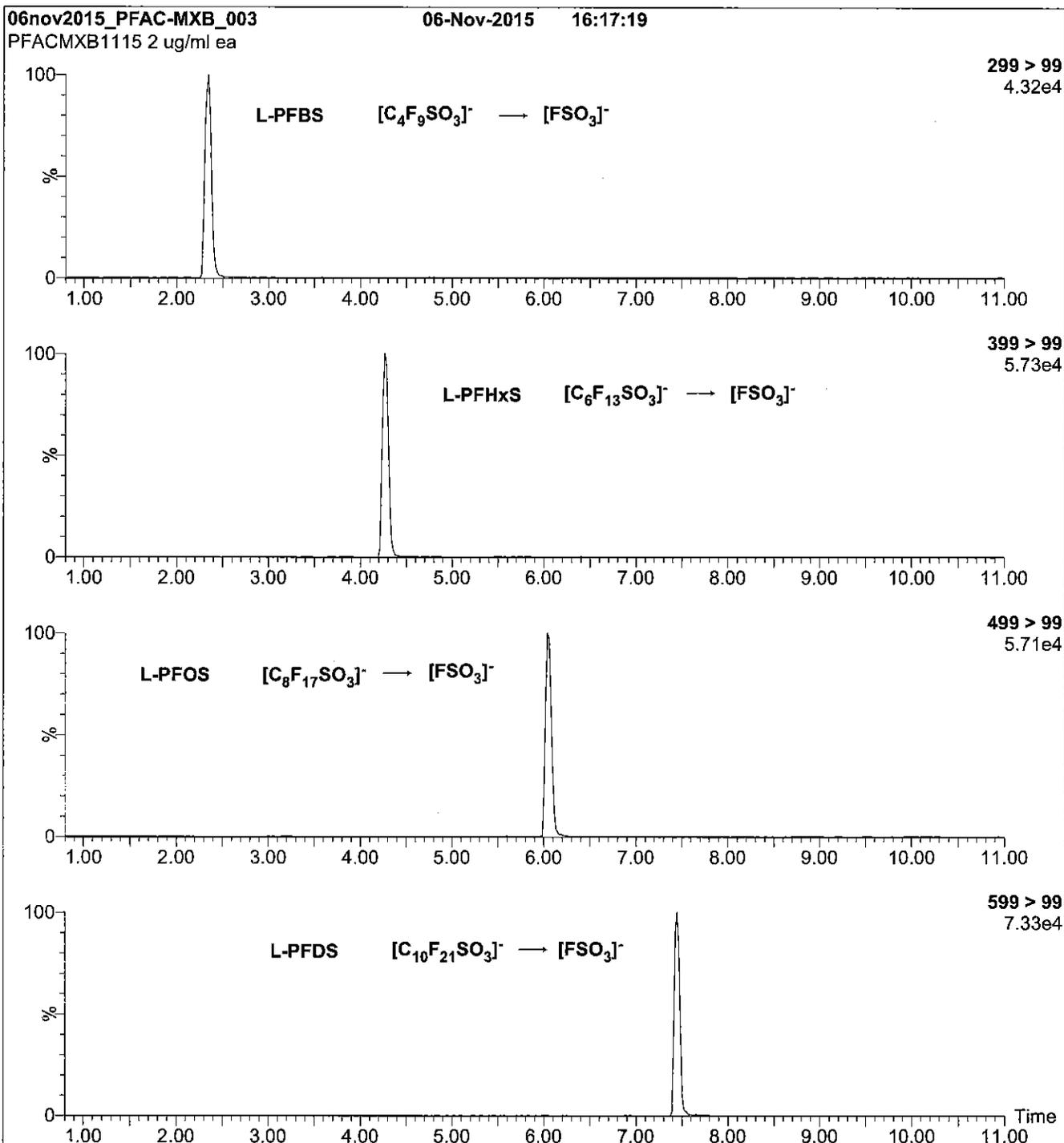


Figure 3: PFAC-MXB; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figures 2 and 3:

Injection: on-column (PFAC-MXB)
 Mobile phase: Same as Figure 1
 Flow: 300 μ /min

MS Parameters
 Collision Gas (mbar) = 3.24e-3
 Collision Energy (eV) = 8-50 (variable)

Reagent

LCPFBA_00004



R: 2125/16 CBW

587895

ID: LCPFBA_00004

Exp: 01/30/20 Prep: CBW

PF-n-butanoic acid

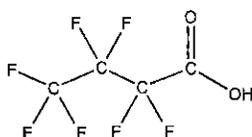


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFBA **LOT NUMBER:** PFBA0115
COMPOUND: Perfluoro-n-butanoic acid

STRUCTURE: **CAS #:** 375-22-4



MOLECULAR FORMULA: C₄HF₇O₂ **MOLECULAR WEIGHT:** 214.04
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
 Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/30/2015
EXPIRY DATE: (mm/dd/yyyy) 01/30/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 

B.G. Chittim

Date: 03/25/2015
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

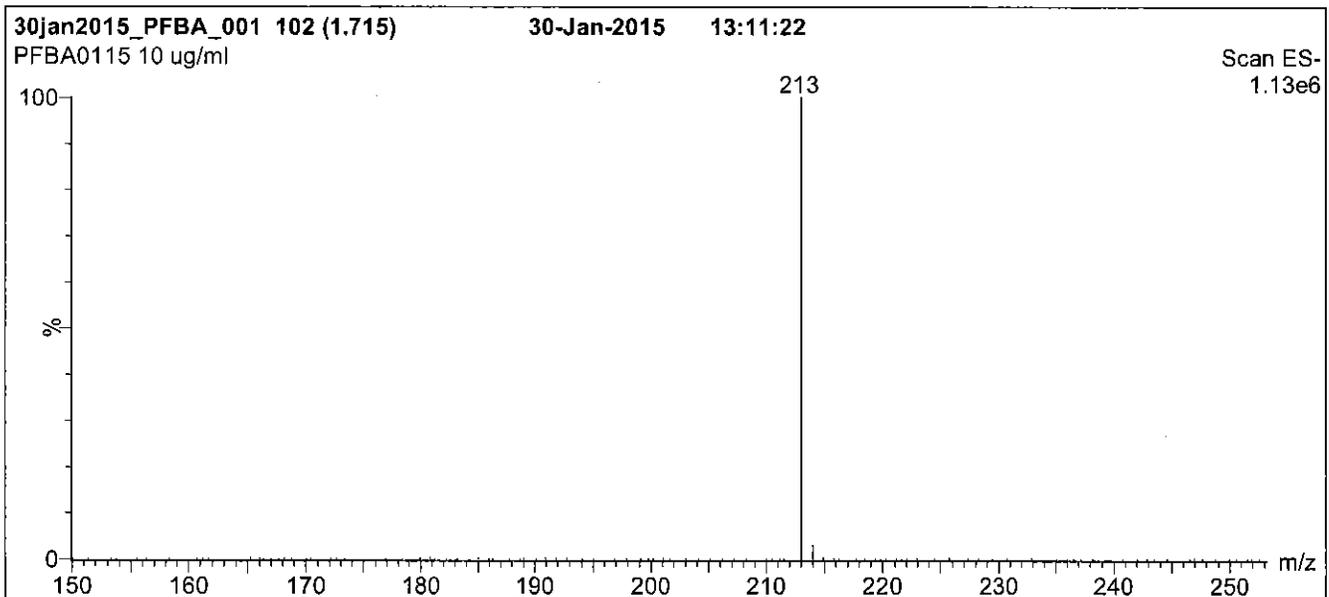
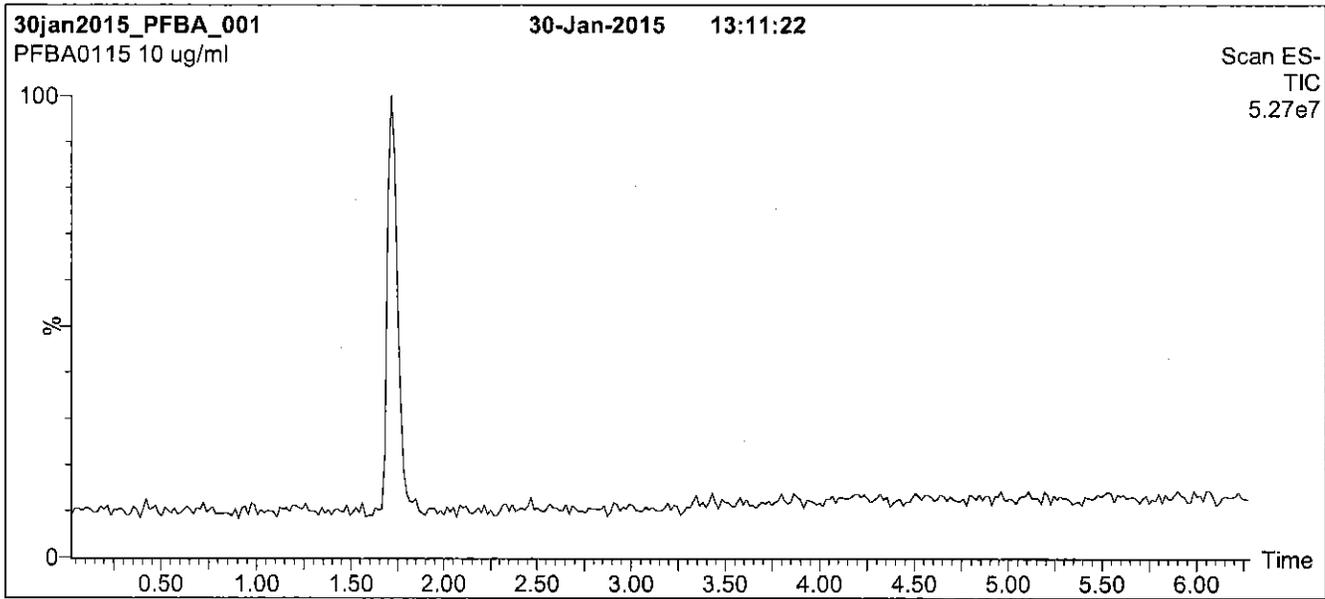
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: PFBA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 30% (80:20 MeOH:ACN) / 70% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7.5 min and hold for 1 min
before returning to initial conditions in 0.5 min.
Time: 10 min

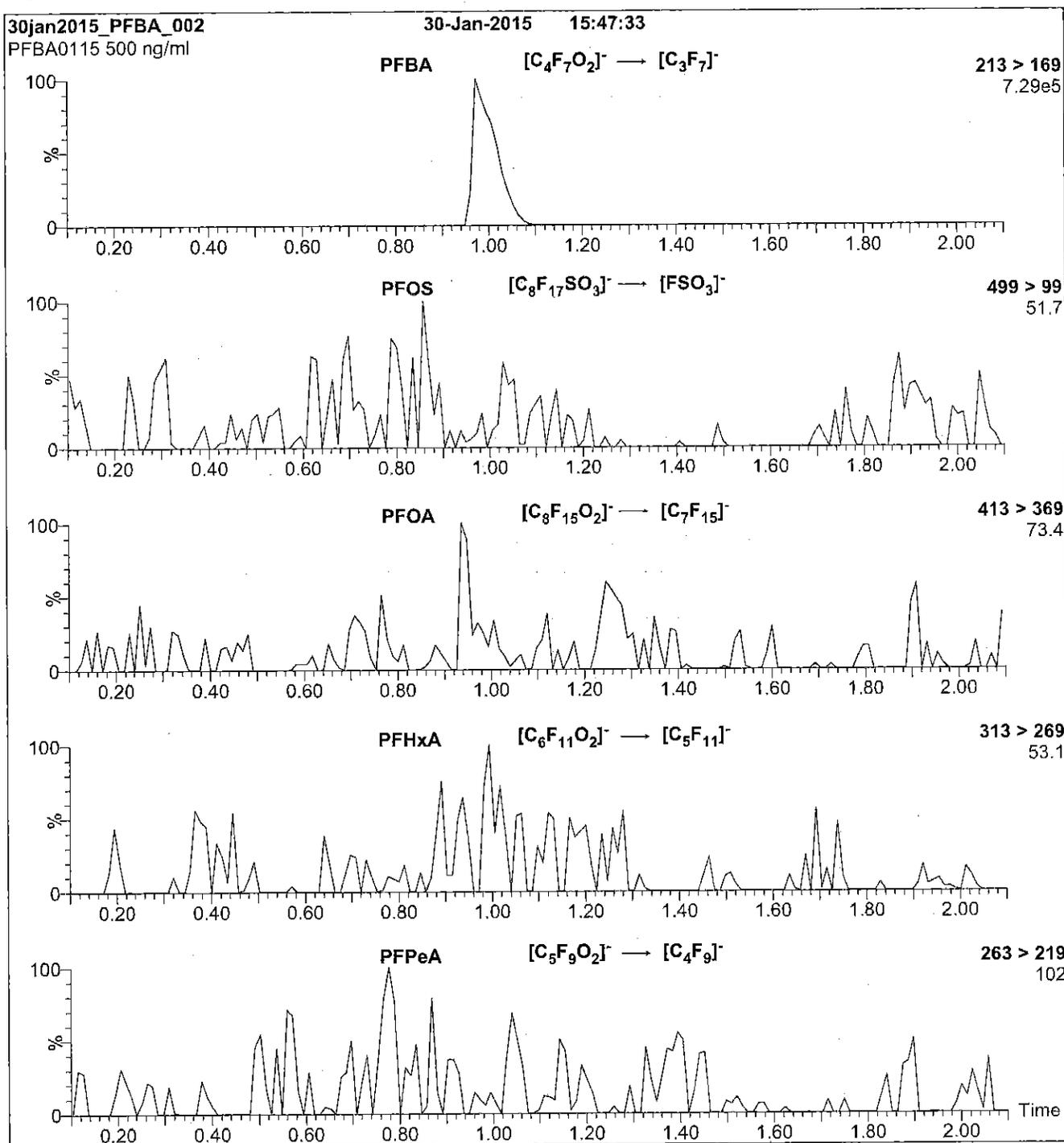
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 8.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFBA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFBA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.35e-3
 Collision Energy (eV) = 10

Reagent

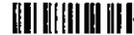
LCPFBA_00005

Scanned
10/16/14

R: SBC 9/13/16



730531
ID: LCPFBA_00005
Exp: 05/27/21 Prpd: SBC
PF-n-butanolic acid



730532
ID: LCPFBA_00006
Exp: 05/27/21 Prpd: SBC
PF-n-butanolic acid



WELLINGTON LABORATORIES

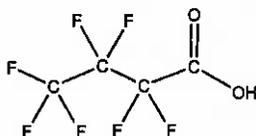
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFBA
COMPOUND: Perfluoro-n-butanolic acid

LOT NUMBER: PFBA0516

STRUCTURE:

CAS #: 375-22-4



MOLECULAR FORMULA: C₄HF₇O₂
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 214.04
SOLVENT(S): Methanol
Water (<1%)

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 05/27/2016
EXPIRY DATE: (mm/dd/yyyy) 05/27/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole.eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 05/31/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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where x is expressed as a relative standard uncertainty of the individual parameter.

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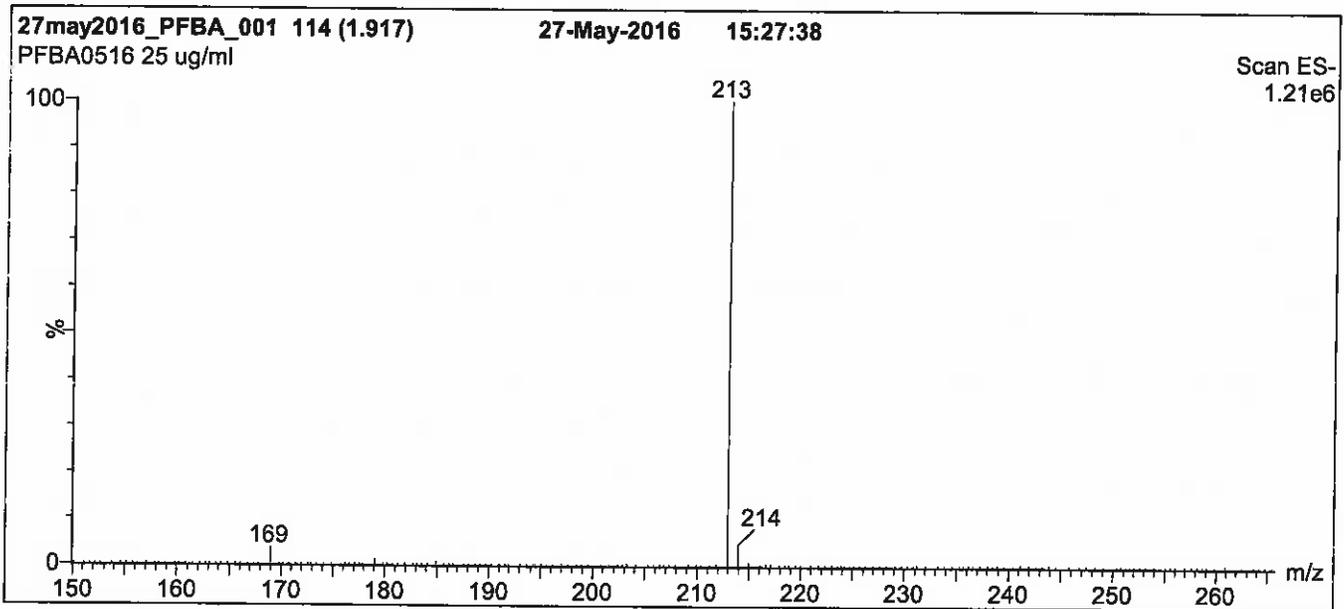
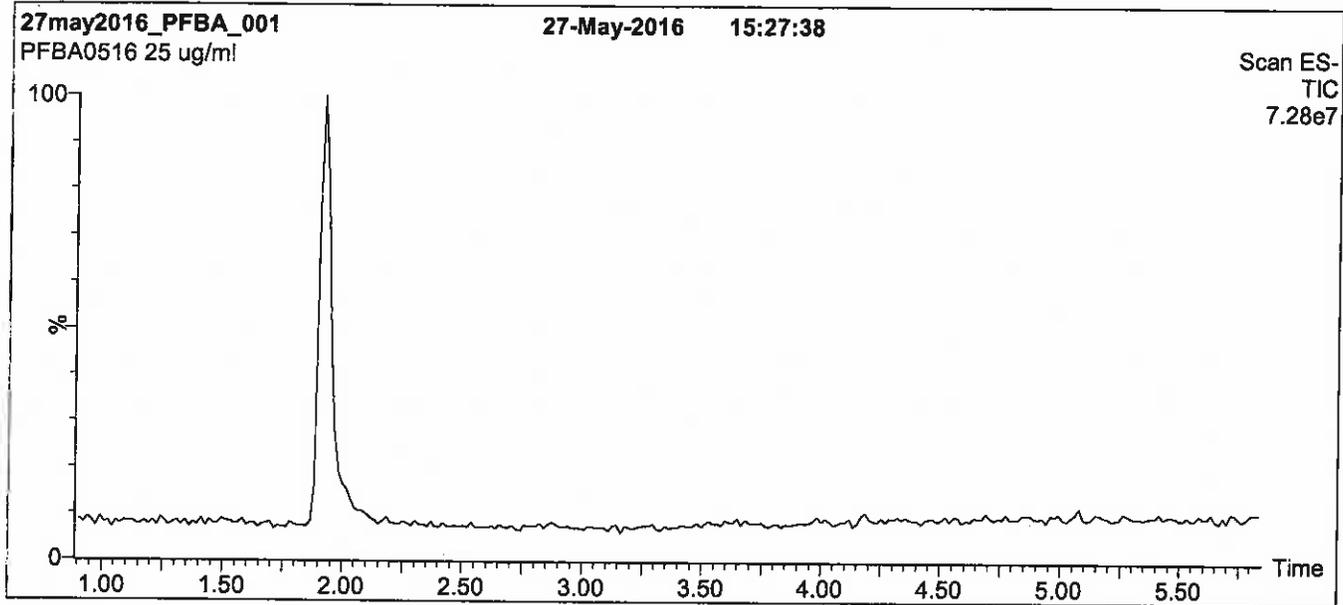
QUALITY MANAGEMENT:

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Figure 1: PFBA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 30% (80:20 MeOH:ACN) / 70% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min before returning to initial conditions in 0.5 min.
Time: 10 min

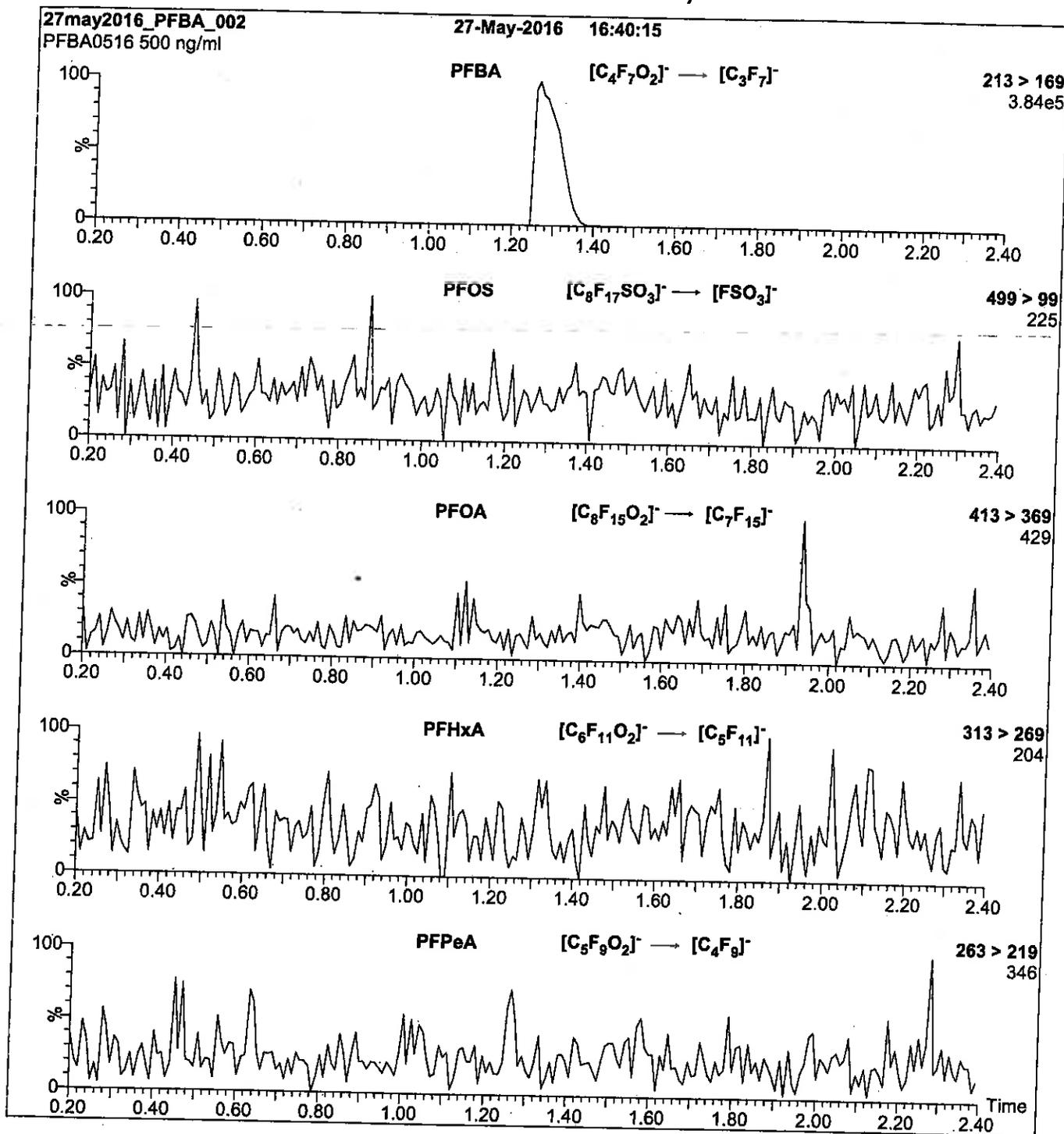
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 10.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFBA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFBA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.62e-3
Collision Energy (eV) = 10

Reagent

LCPFBS_00004



Rec. 3/29/16 JRB ✓

605236

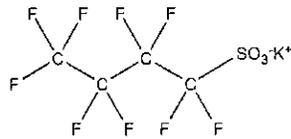
ID: LCPFBS_00004

Exp: 10/09/19 Prpd: CBW

PF-1-butanesulfonate K sa

**WELLINGTON
LABORATORIES****CERTIFICATE OF ANALYSIS
DOCUMENTATION**

PRODUCT CODE: L-PFBS **LOT NUMBER:** LPFBS1014
COMPOUND: Potassium perfluoro-1-butanesulfonate
STRUCTURE: **CAS #:** 29420-49-3



MOLECULAR FORMULA: C₄F₉SO₃K **MOLECULAR WEIGHT:** 338.19
CONCENTRATION: 50.0 ± 2.5 µg/ml (K salt) **SOLVENT(S):** Methanol
44.2 ± 2.2 µg/ml (PFBS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 10/09/2014
EXPIRY DATE: (mm/dd/yyyy) 10/09/2019
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 04/02/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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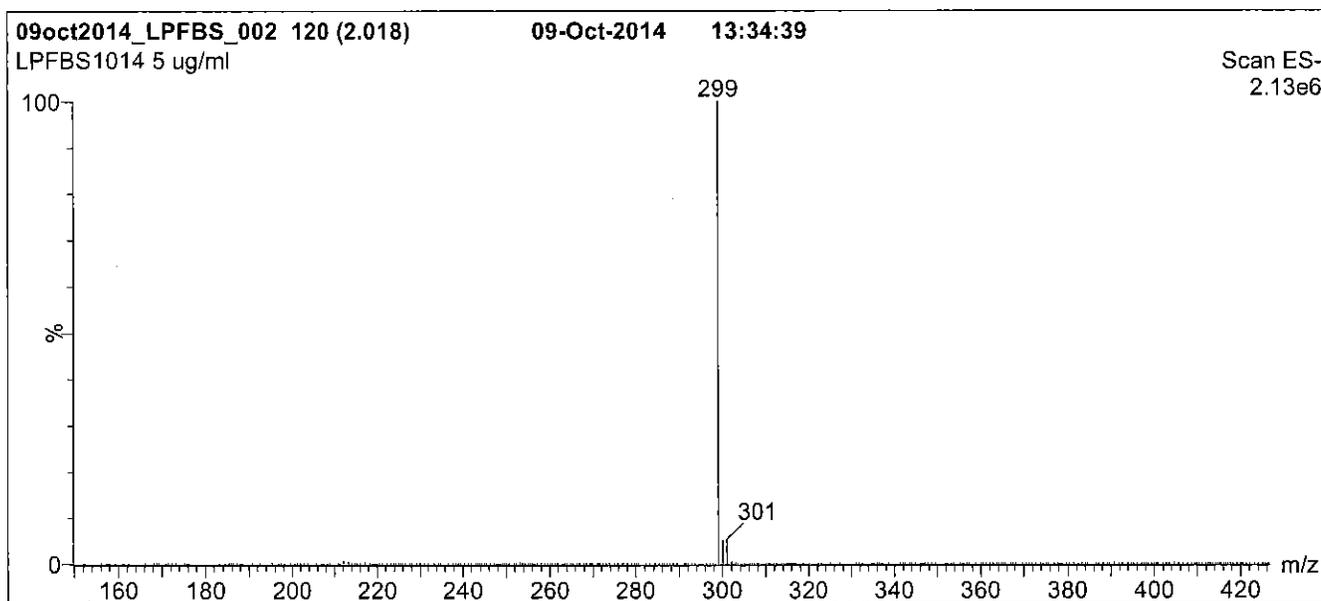
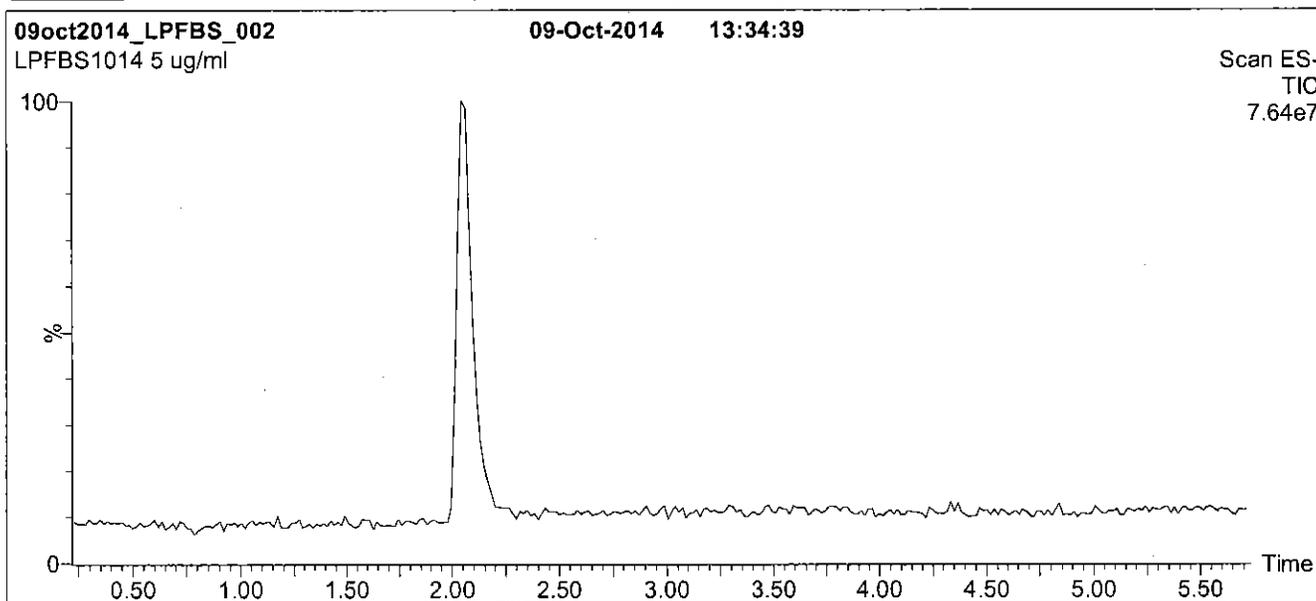
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Figure 1: L-PFBS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 40% (80:20 MeOH:ACN) / 60% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

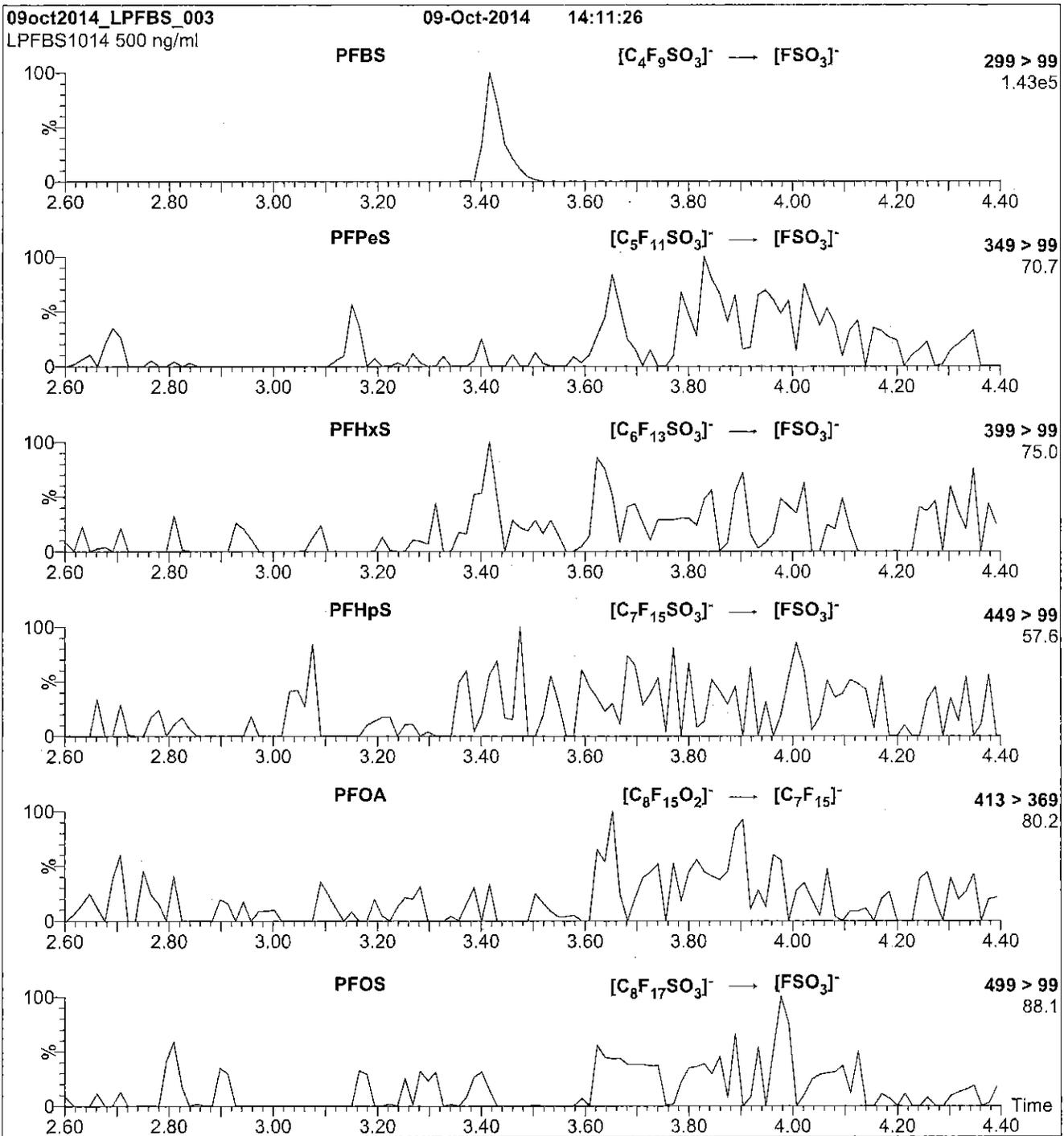
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 40.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: L-PFBS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml L-PFBS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
 Collision Energy (eV) = 25

Reagent

LCPFBS_00005

R: 9/9/16 gbe



728306
ID: LCM2-8:2FTS_00003
Exp: 01/08/21 Prpd: SBC
M2-8:2FTS

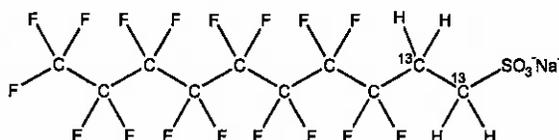


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: M2-8:2FTS **LOT NUMBER:** M282FTS0116
COMPOUND: Sodium 1H,1H,2H,2H-perfluoro-[1,2-¹³C₂]decane sulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: ¹³C₂¹²C₈H₄F₁₇SO₃Na **MOLECULAR WEIGHT:** 552.15
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
47.9 ± 2.4 µg/ml (M2-8:2FTS anion)
CHEMICAL PURITY: >98% **ISOTOPIC PURITY:** ≥99% ¹³C
LAST TESTED: (mm/dd/yyyy) 01/08/2016 (1,2-¹³C₂)
EXPIRY DATE: (mm/dd/yyyy) 01/08/2021
RECOMMENDED STORAGE: Refrigerate ampoule

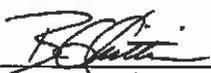
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- The native 8:2FTS contains 4.22% of ³⁴S (due to natural isotopic abundance) therefore both native 8:2FTS and M2-8:2FTS will produce signals in the m/z 529 to m/z 509 channel during SRM analysis. We recommend using the m/z 529 to m/z 81 transition to monitor for M2-8:2FTS during quantitative analysis as it will be free of any native contribution (see Figure 2).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 01/18/2016
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

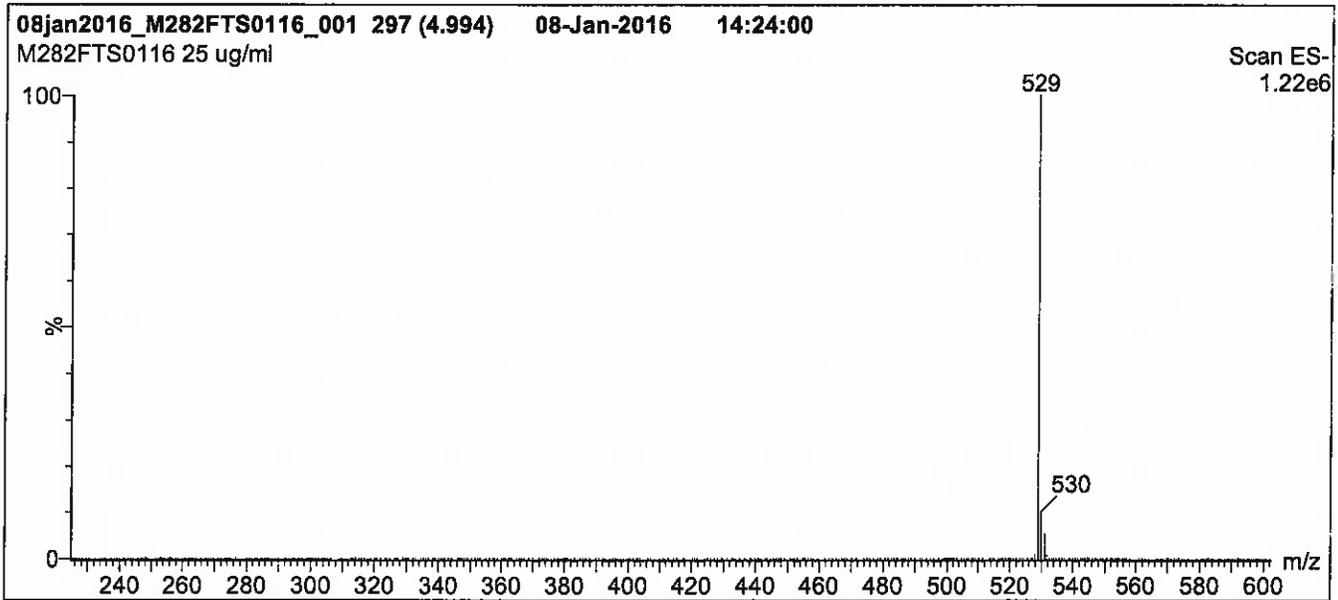
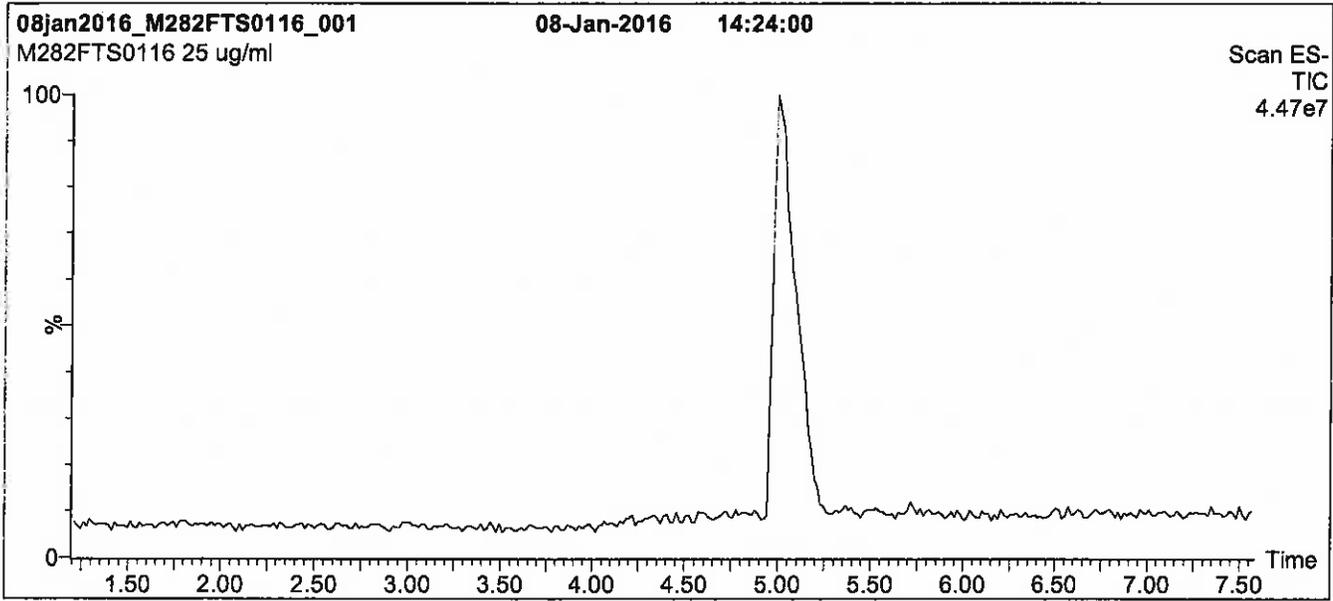
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: M2-8:2FTS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min
and hold for 2 min before returning
to initial conditions in 0.5 min.
Time: 10 min

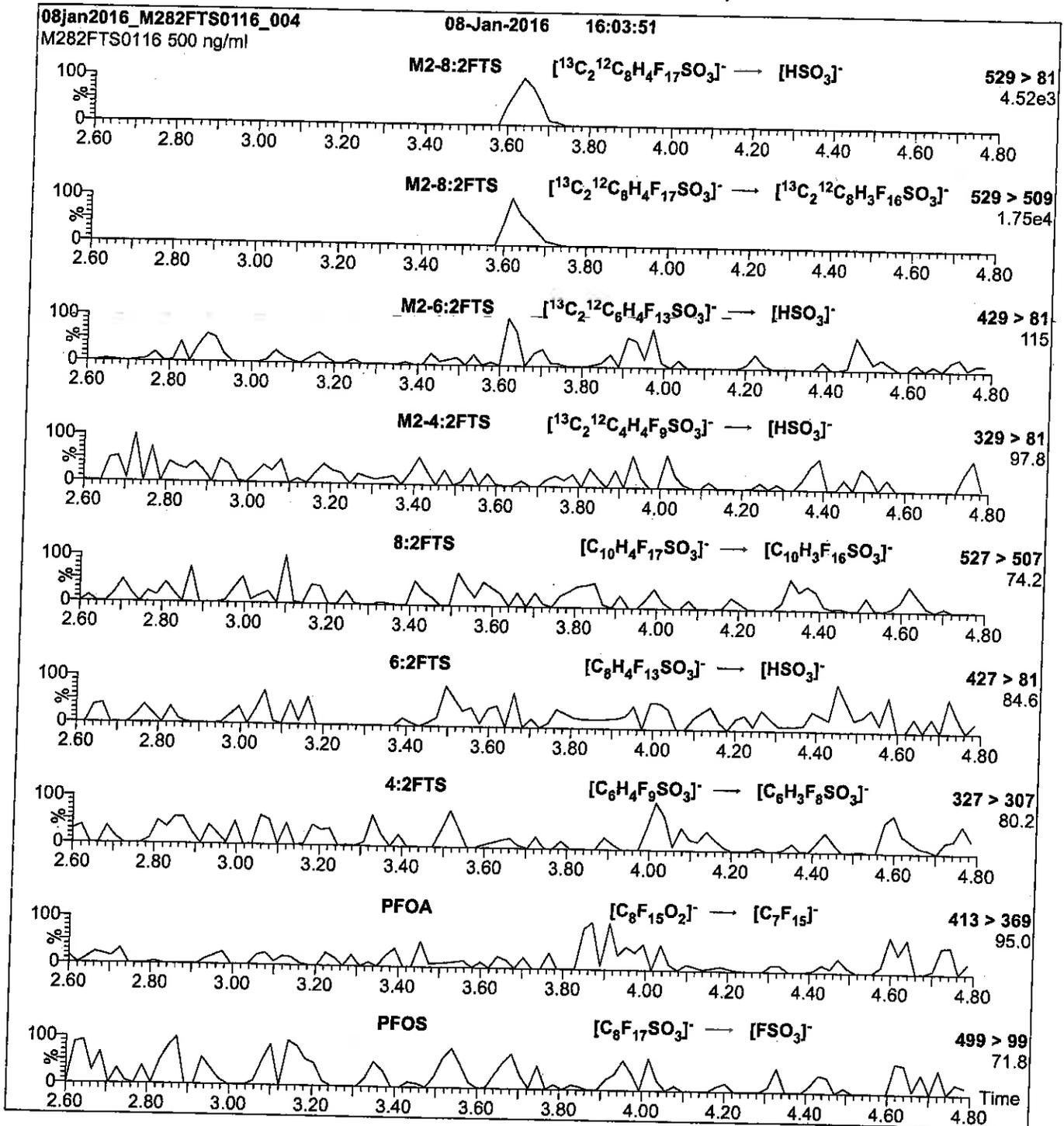
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 30.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: M2-8:2FTS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μl (500 ng/ml M2-8:2FTS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H_2O
(both with 10 mM NH_4OAc buffer)

Flow: 300 $\mu\text{l}/\text{min}$

MS Parameters

Collision Gas (mbar) = 3.20e-3
Collision Energy (eV) = 30

R: SBC 9/13/16



730511
ID: LCPFBS_00005
Exp: 03/15/21 Pripd: SBC
PF-1-butanesulfonate K sa



730512
ID: LCPFBS_00006
Exp: 03/15/21 Pripd: SBC
PF-1-butanesulfonate K sa

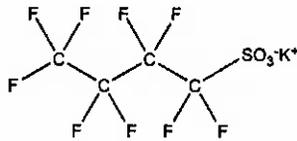


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CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: L-PFBS **LOT NUMBER:** LPFBS0316
COMPOUND: Potassium perfluoro-1-butanesulfonate

STRUCTURE: **CAS #:** 29420-49-3



MOLECULAR FORMULA: C₄F₉SO₃K **MOLECULAR WEIGHT:** 338.19
CONCENTRATION: 50.0 ± 2.5 µg/ml (K salt) **SOLVENT(S):** Methanol
44.2 ± 2.2 µg/ml (PFBS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 03/15/2016
EXPIRY DATE: (mm/dd/yyyy) 03/15/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

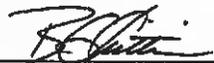
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 03/21/2016
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

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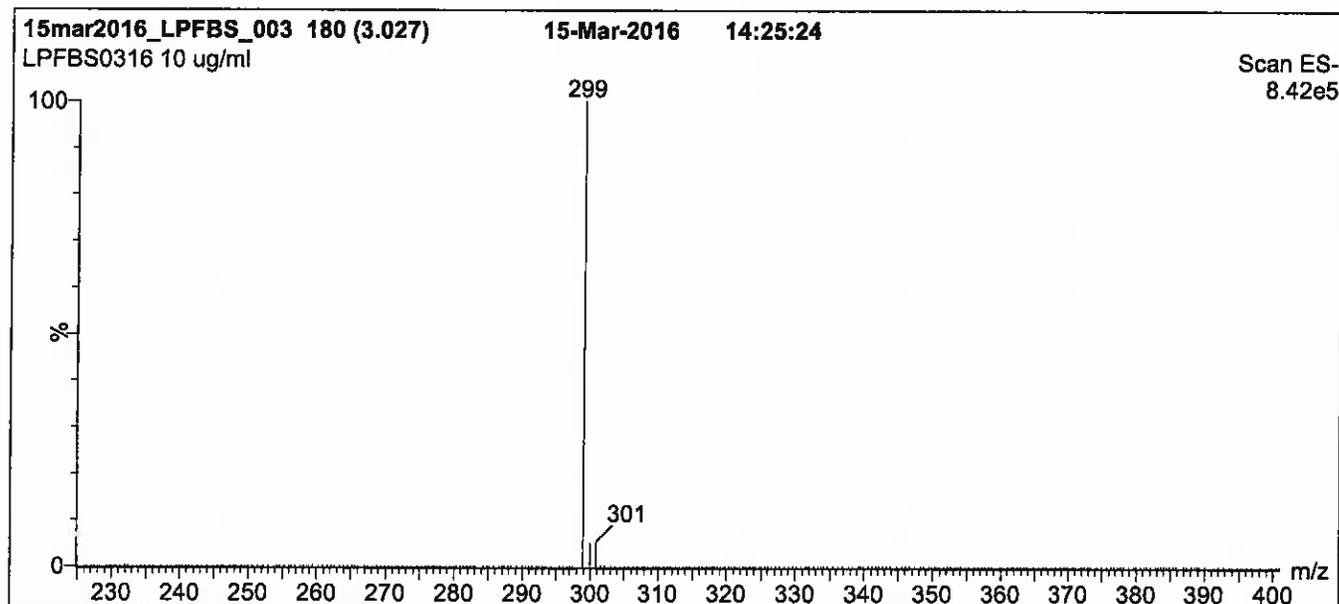
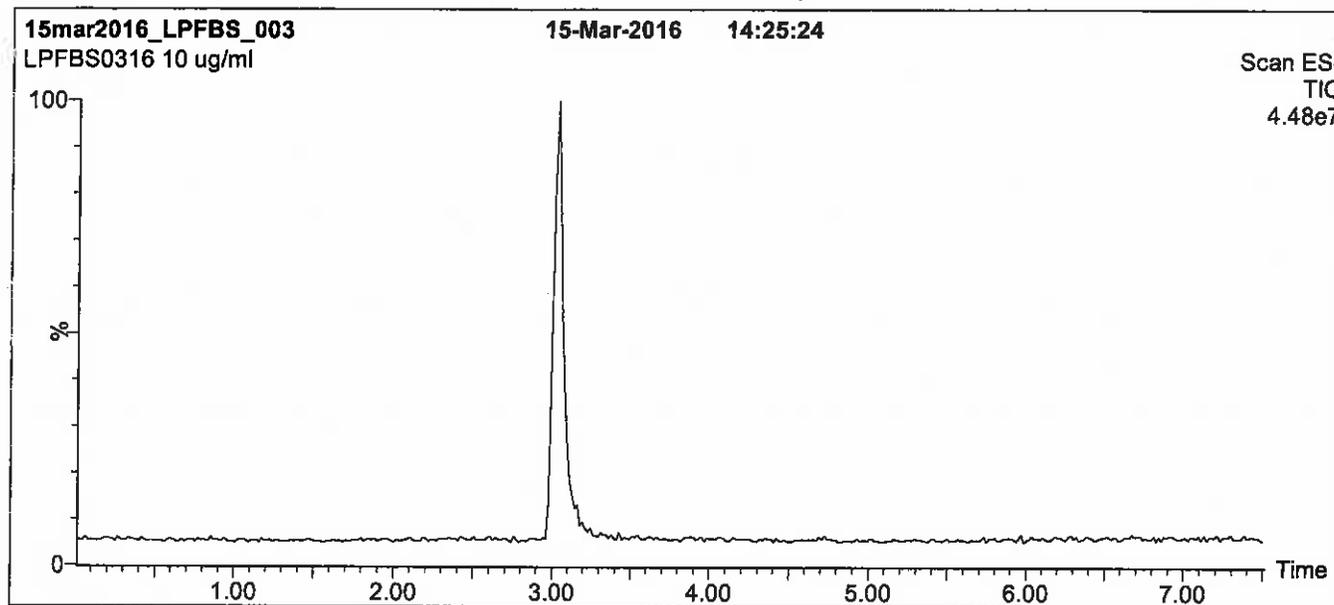
QUALITY MANAGEMENT:

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Figure 1: L-PFBS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 40% (80:20 MeOH:ACN) / 60% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

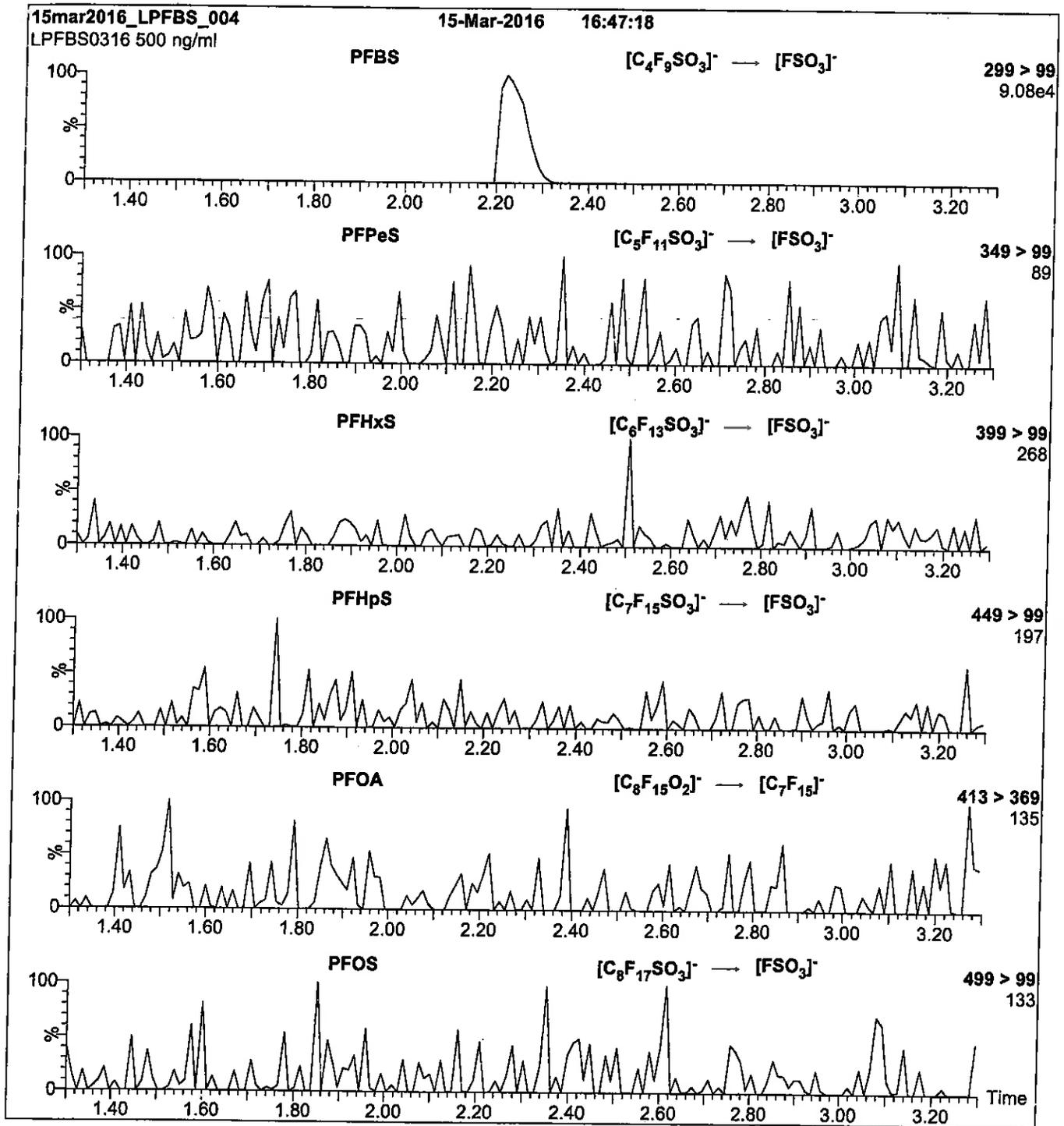
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 40.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: L-PFBS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml L-PFBS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.20e-3
 Collision Energy (eV) = 25

Reagent

LCPFDA_00005

R: 7/16/16 CBW



671576
ID: LCPFDA_00305
Exp: 07/02/20 Pipd: CBW
PF-n-decanoic acid

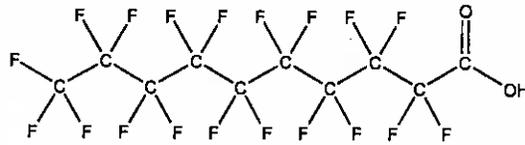


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CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFDA **LOT NUMBER:** PFDA0615
COMPOUND: Perfluoro-n-decanoic acid

STRUCTURE: **CAS #:** 335-76-2



MOLECULAR FORMULA: C₁₀H₁₉O₂ **MOLECULAR WEIGHT:** 514.08
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 07/02/2015
EXPIRY DATE: (mm/dd/yyyy) 07/02/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.6% PFNA and ~ 0.3% PFOA.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 07/24/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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LIMITED WARRANTY:

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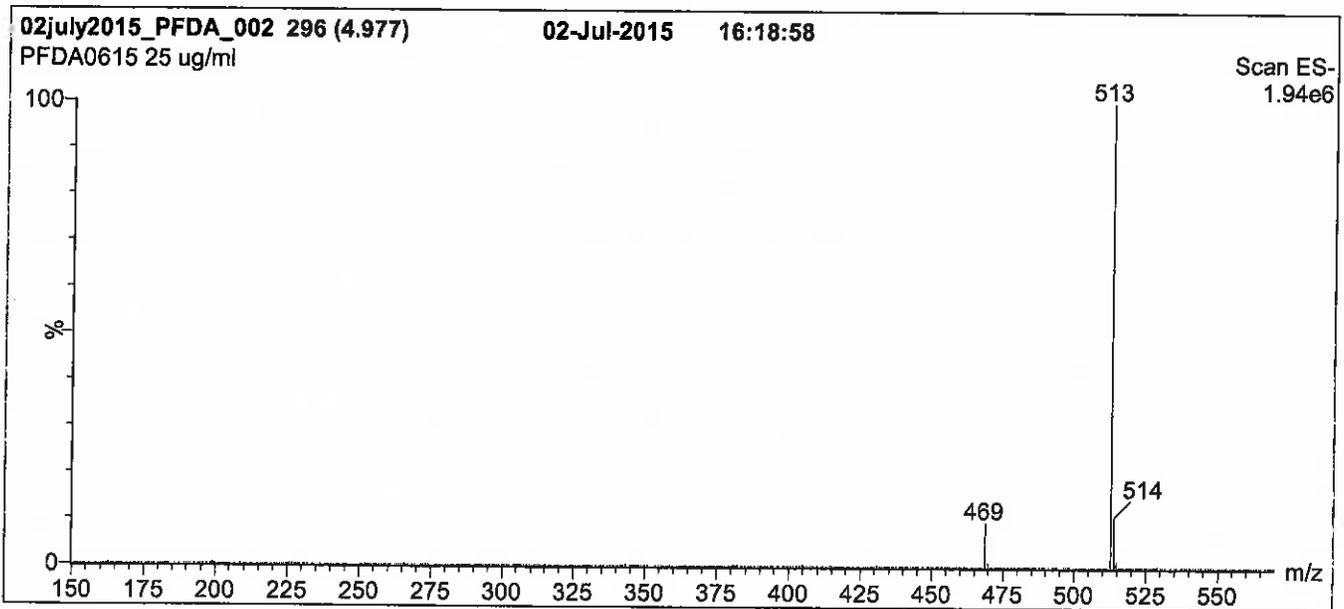
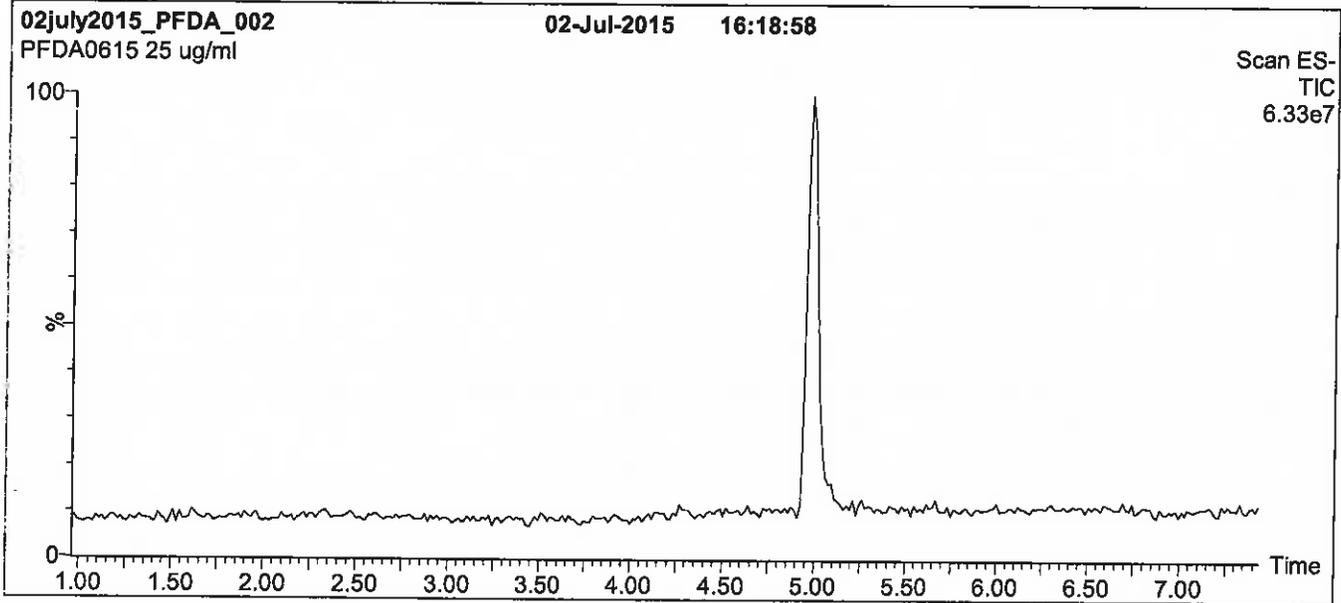
QUALITY MANAGEMENT:

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Figure 1: PFDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 2 min before returning to initial conditions in 0.5 min.
 Time: 10 min

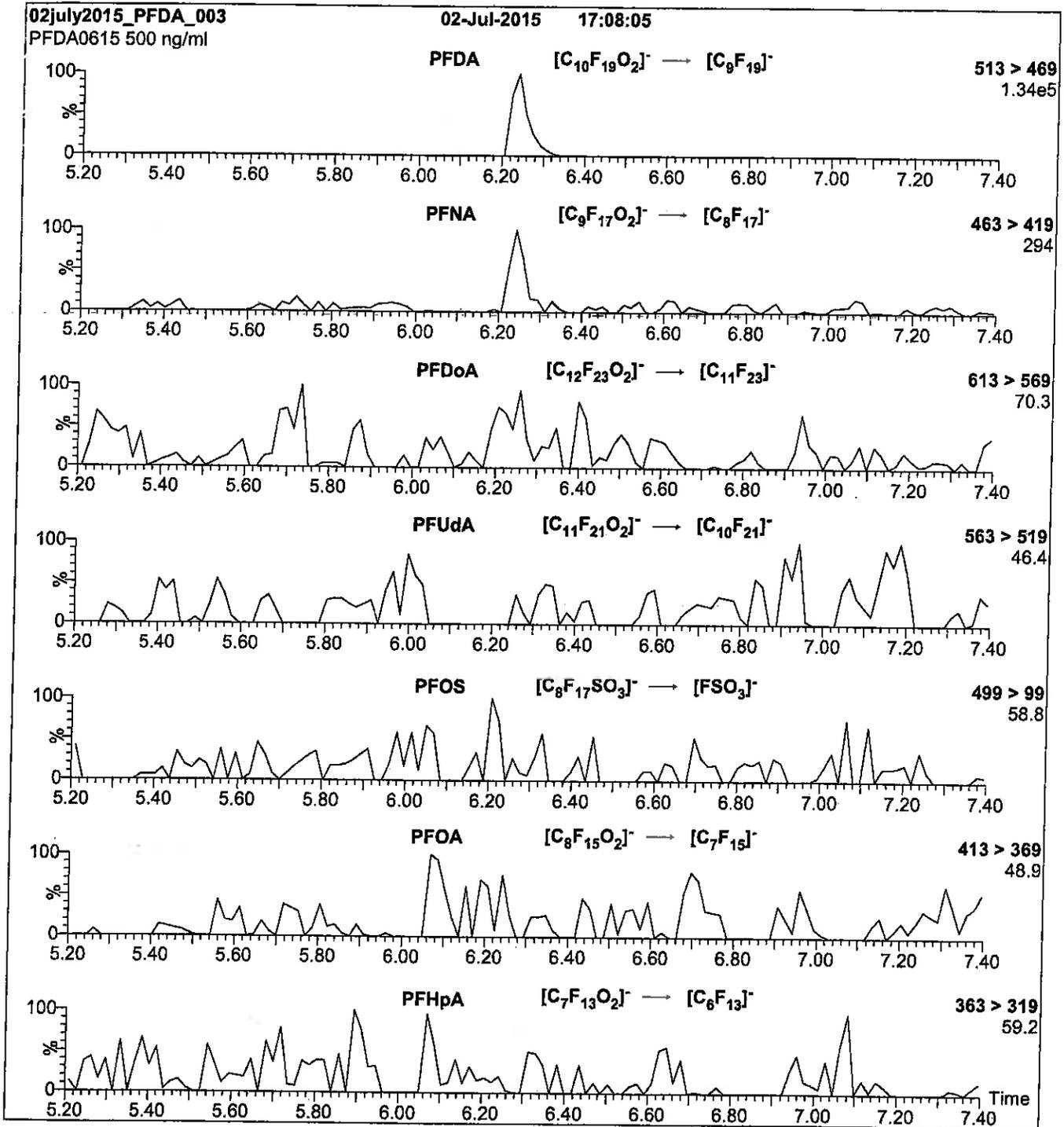
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: PFDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.62e-3
Collision Energy (eV) = 13

Reagent

LCPFDoA_00005

R: 7/6/16 can

671601
ID: LCPFD0A_00005
Exp: 01/30/20 Prod: CBW
PF-n-dodecanoic acid

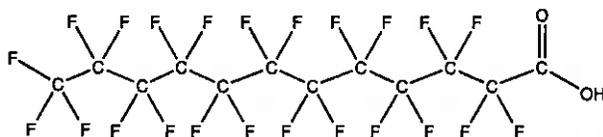


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LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

PRODUCT CODE: PFD0A **LOT NUMBER:** PFD0A0115
COMPOUND: Perfluoro-n-dodecanoic acid

STRUCTURE: **CAS #:** 307-55-1



MOLECULAR FORMULA: $C_{12}HF_{23}O_2$ **MOLECULAR WEIGHT:** 614.10
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/30/2015
EXPIRY DATE: (mm/dd/yyyy) 01/30/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 03/25/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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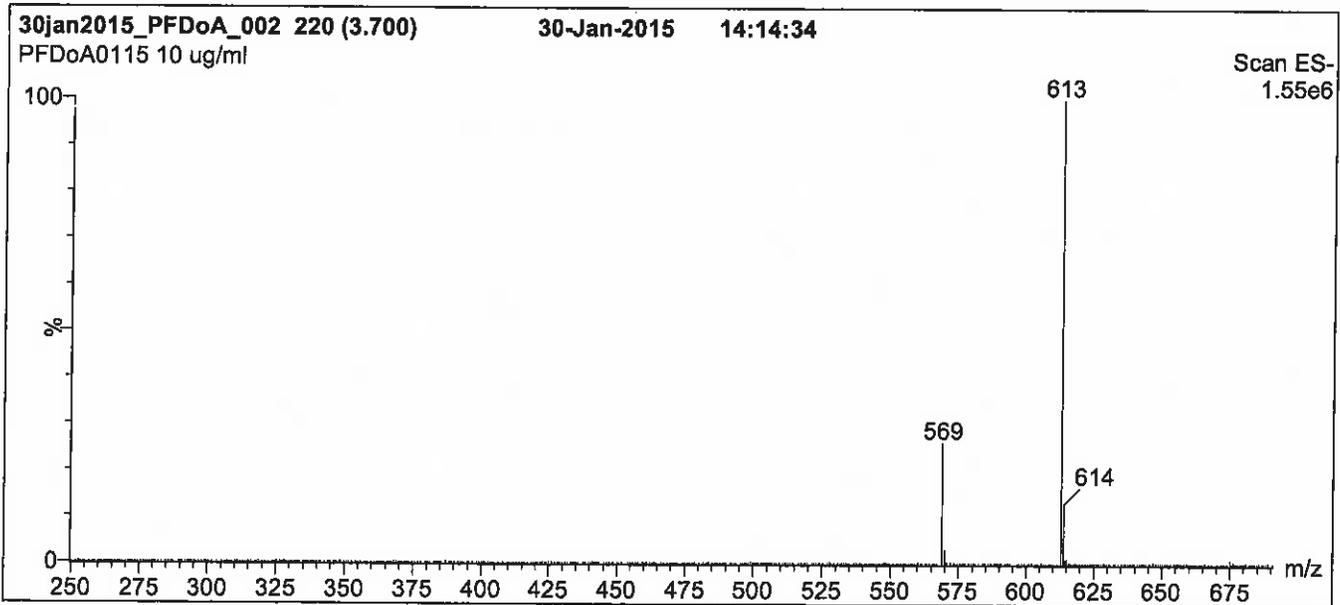
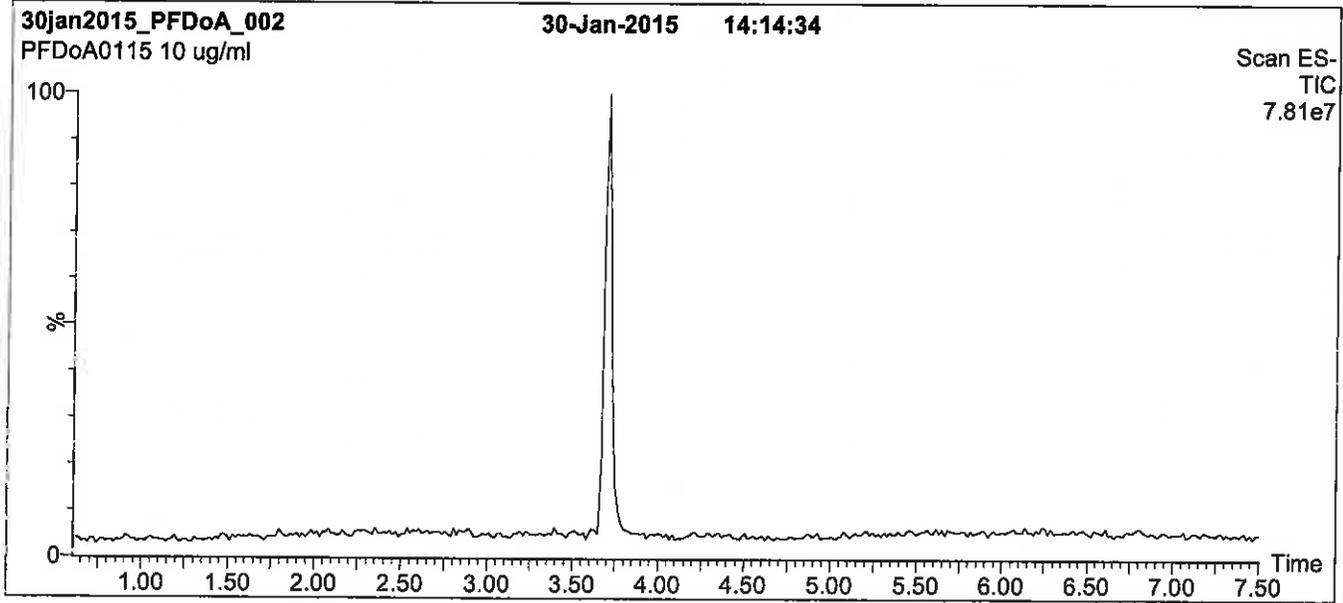
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Figure 1: PFD_oA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 60% (80:20 MeOH:ACN) / 40% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

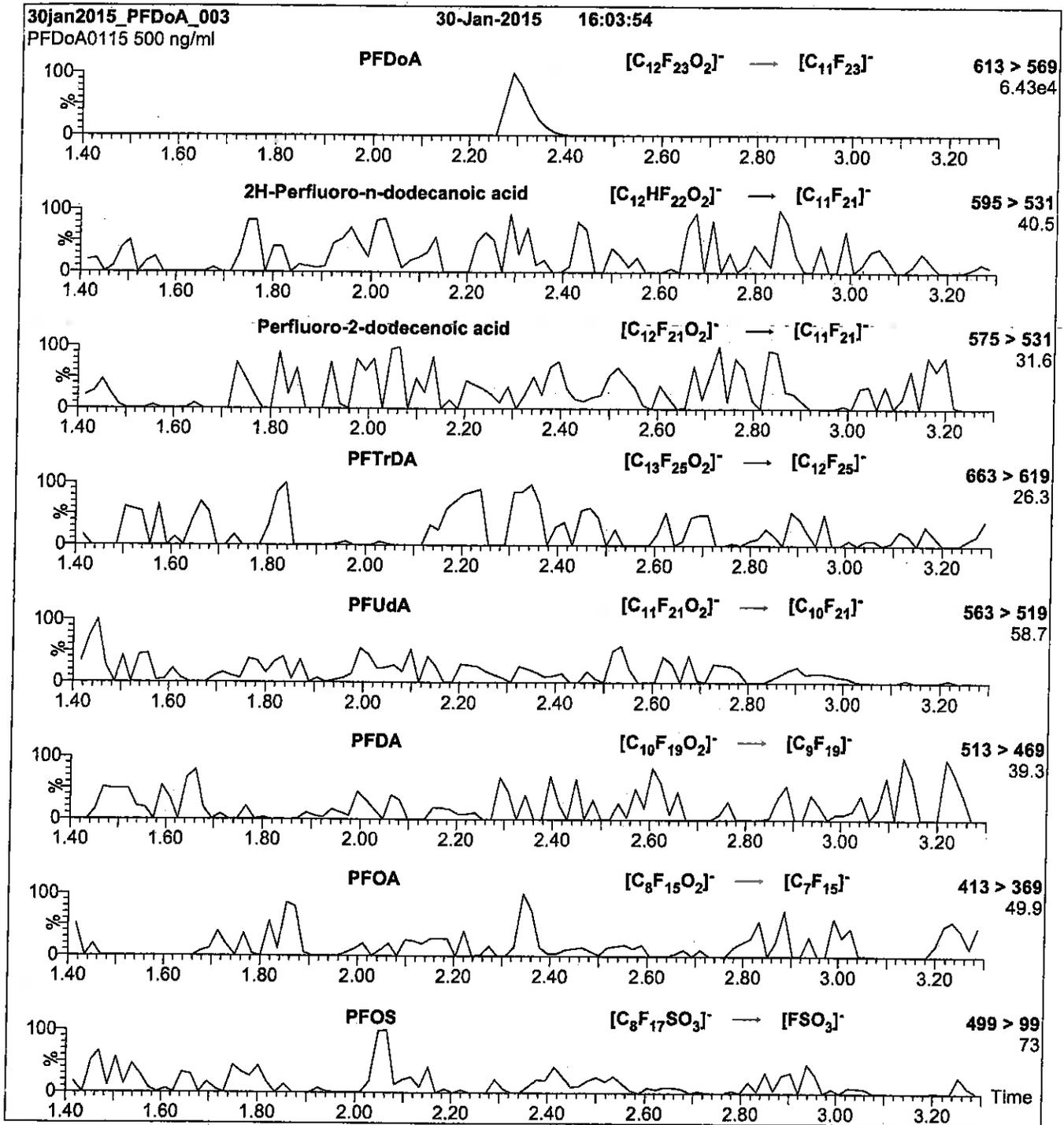
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 1000 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 20.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFDoA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFDoA)

MS Parameters

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Collision Gas (mbar) = 3.28e-3
 Collision Energy (eV) = 13

Flow: 300 μ l/min

Reagent

LCPFDS_00005



605240
 ID: LCPFDS_00005
 Exp: 07/02/20 Prep: CBW
 PF-1-decanesulfonate sodi

Rec. 3/29/16 JRB

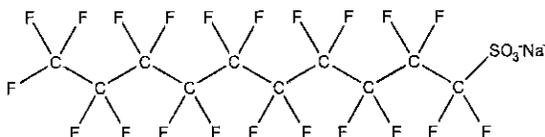


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CERTIFICATE OF ANALYSIS
 DOCUMENTATION

PRODUCT CODE: L-PFDS **LOT NUMBER:** LPFDS0615
COMPOUND: Sodium perfluoro-1-decanesulfonate

STRUCTURE: **CAS #:** 2806-15-7



MOLECULAR FORMULA: C₁₀F₂₁SO₃Na **MOLECULAR WEIGHT:** 622.13
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
 48.2 ± 2.4 µg/ml (PFDS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 07/02/2015
EXPIRY DATE: (mm/dd/yyyy) 07/02/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

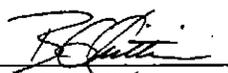
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.9% of sodium perfluoro-1-dodecanesulfonate (L-PFDoS).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 12/07/2015
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

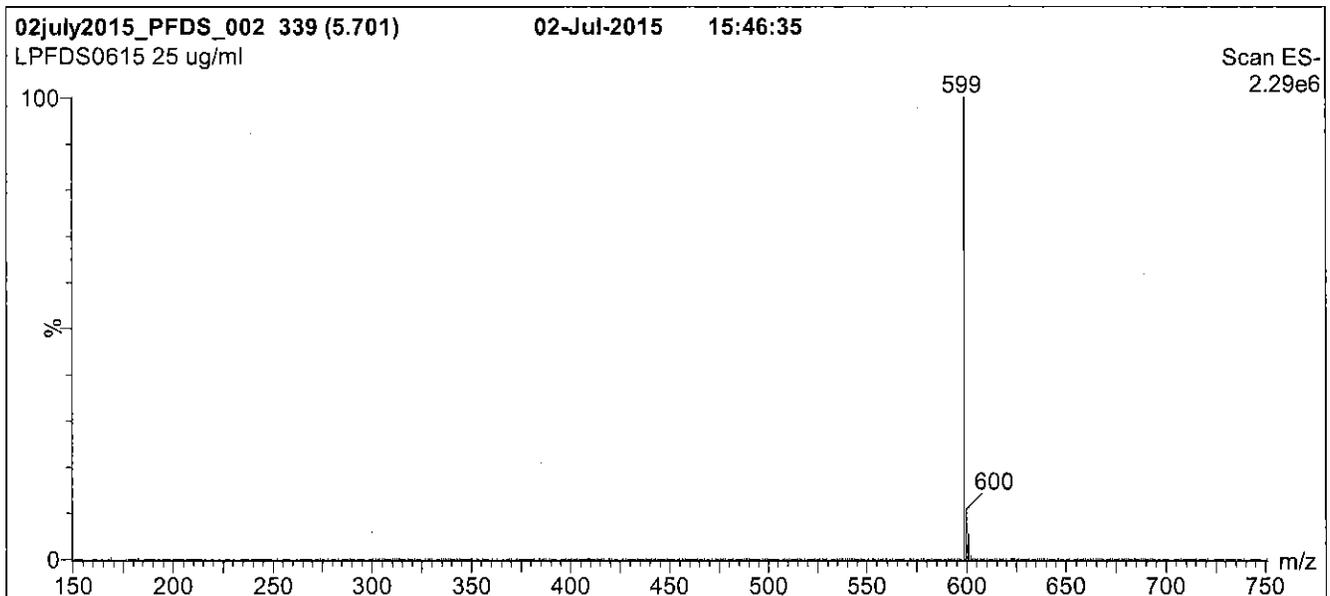
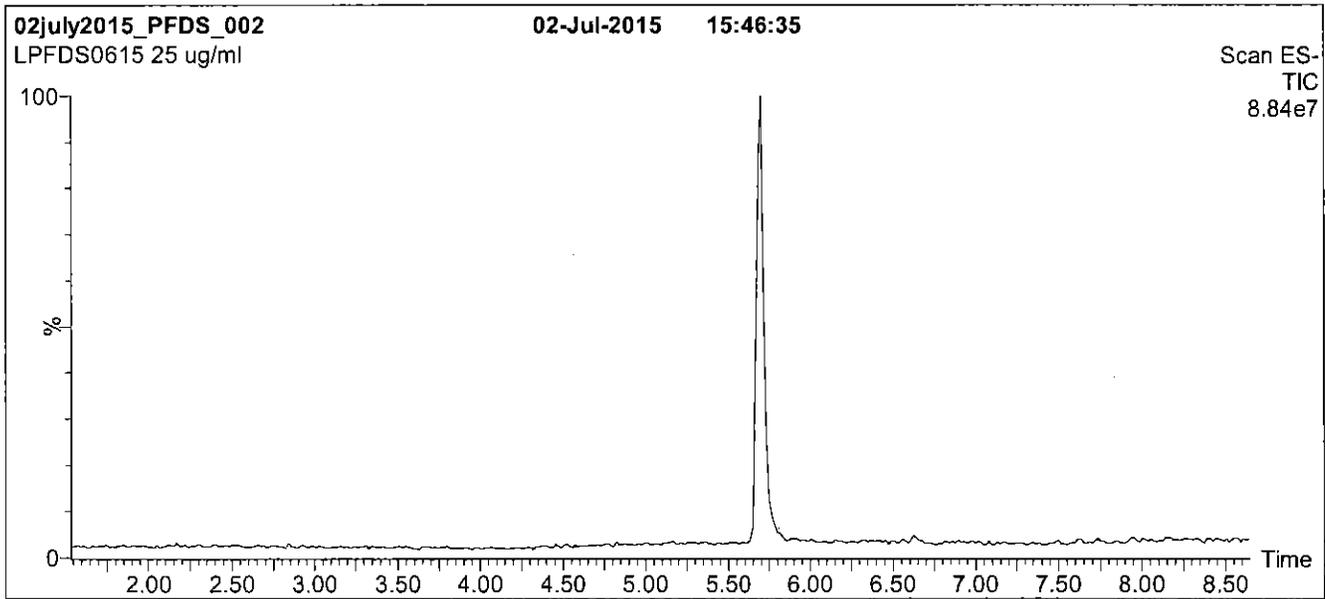
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: L-PFDS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for
2 min before returning to initial conditions in 0.5 min.
Time: 10 min

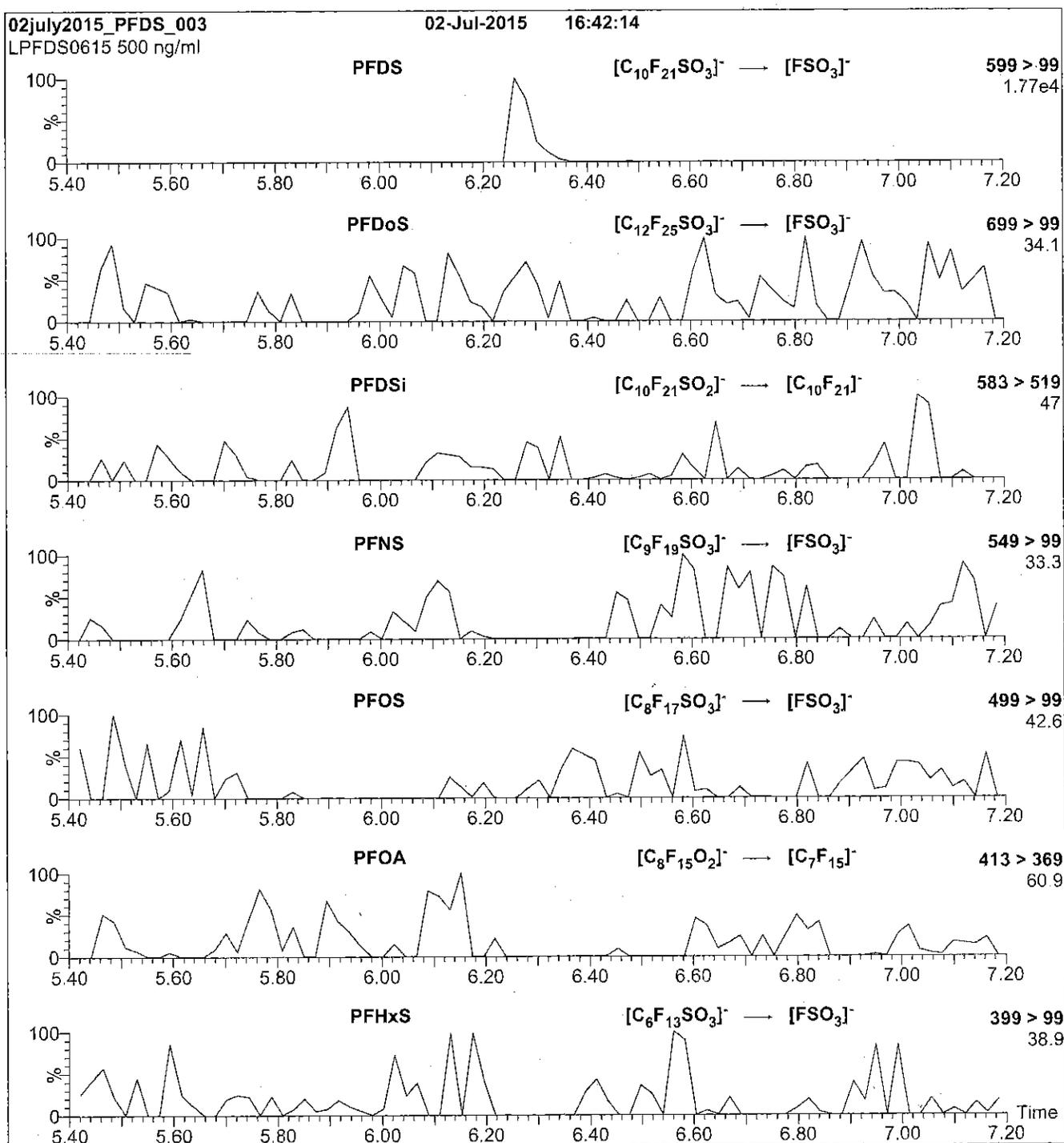
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 70.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: L-PFDS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml L-PFDS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 50

Reagent

LCPFHpA_00005



609639

ID: LCPFHpA_00005

Exp: 01/22/21 Prpd: CBW

PF-n-heptanoic acid

R: 4/7/16 CBW



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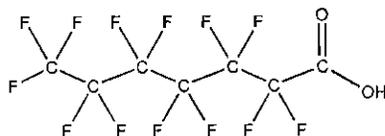
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFHpA
COMPOUND: Perfluoro-n-heptanoic acid

LOT NUMBER: PFHpA0116

STRUCTURE:

CAS #: 375-85-9



MOLECULAR FORMULA: C₇H₁₃O₂
CONCENTRATION: 50 ± 2.5 µg/ml

MOLECULAR WEIGHT: 364.06
SOLVENT(S): Methanol
Water (<1%)

CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/22/2016
EXPIRY DATE: (mm/dd/yyyy) 01/22/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 02/02/2016

(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON 'N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

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The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

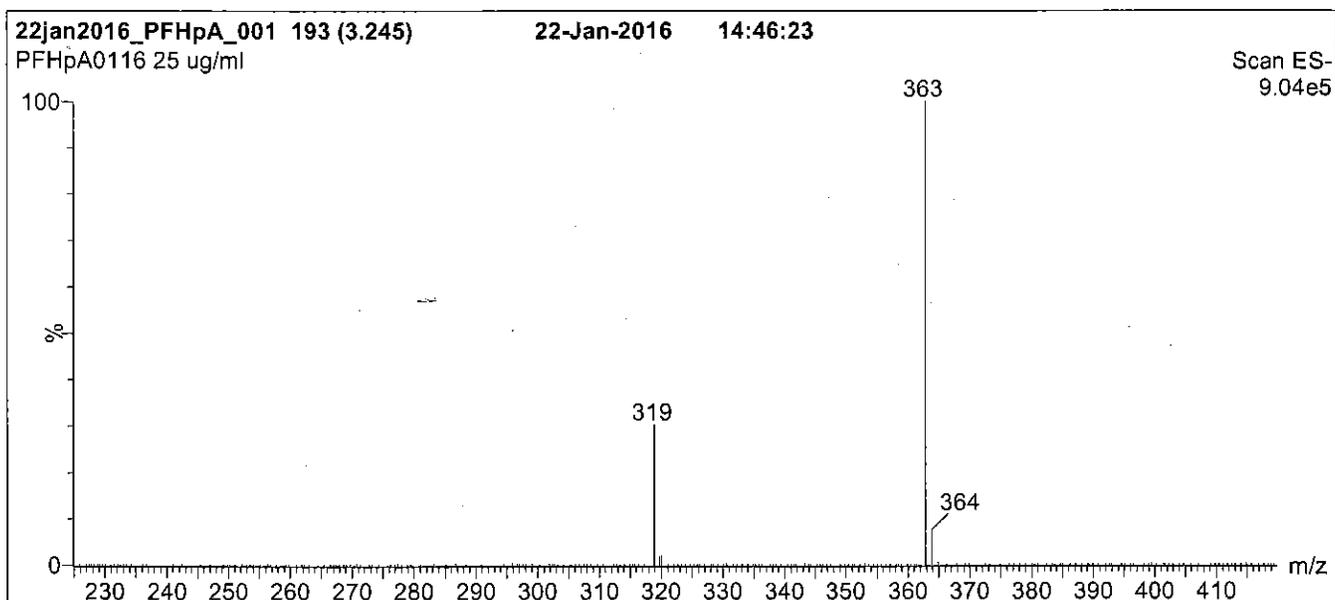
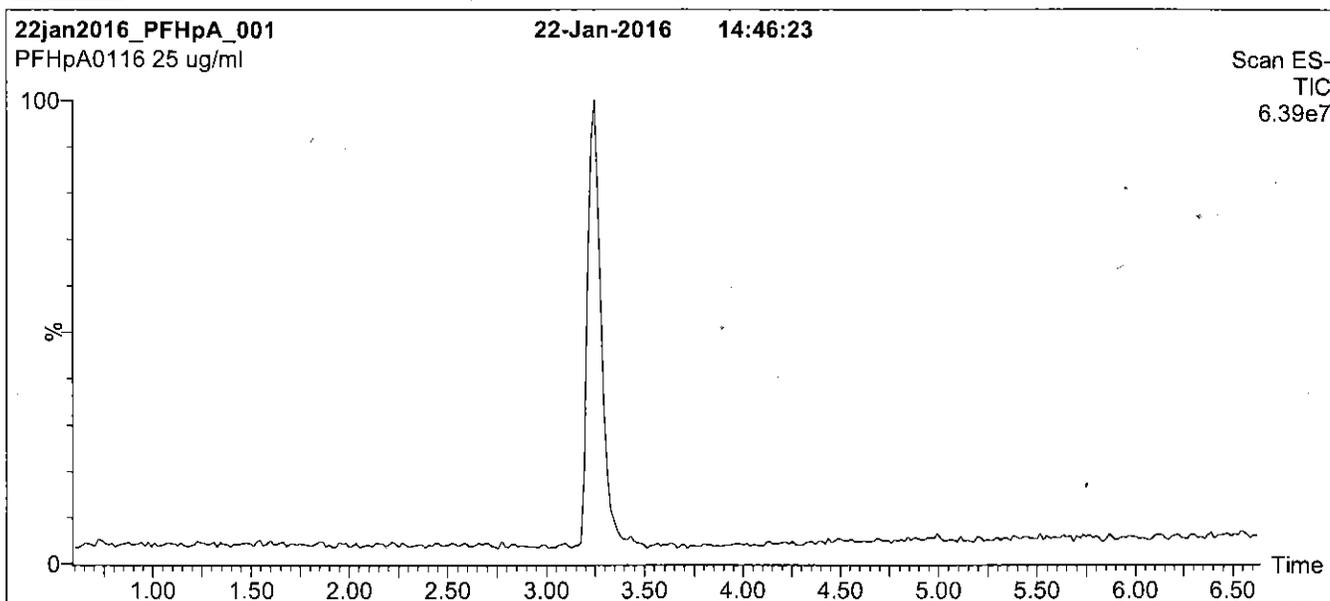
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: PFHpA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 55% (80:20 MeOH:ACN) / 45% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for
2 min before returning to initial conditions in 0.5 min.
Time: 10 min

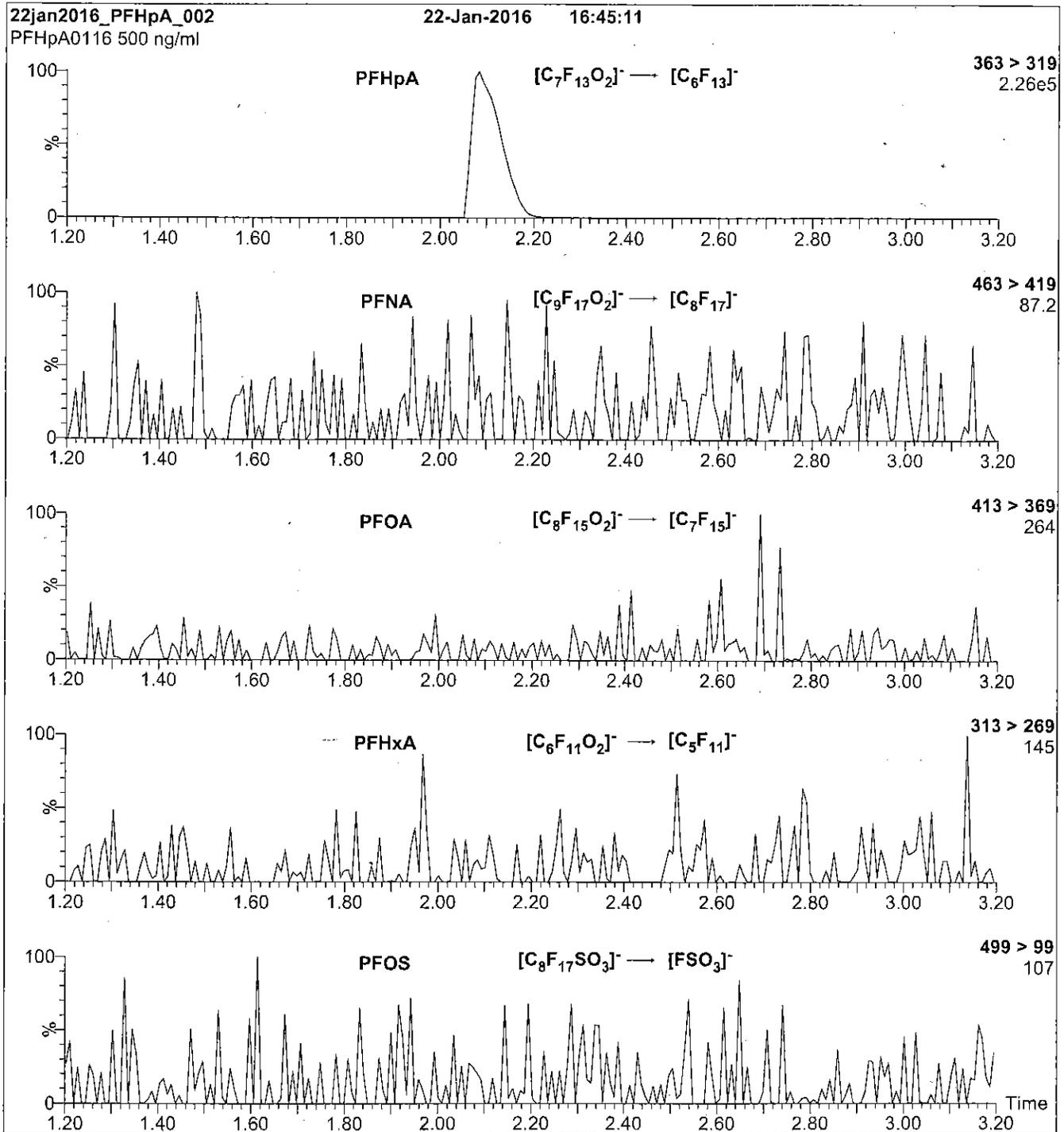
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFHpA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFHpA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.50e-3
Collision Energy (eV) = 11

Reagent

LCPFHpS_00008



627751
 ID: LCPFHPS_00008
 Exp: 11/06/20 Ppt: CBW
 PFHpS at 47.6ug/ml

R: 5/10/16 CBW

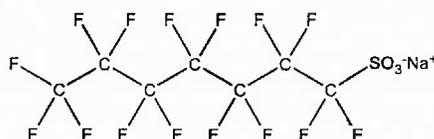


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CERTIFICATE OF ANALYSIS
 DOCUMENTATION

PRODUCT CODE: L-PFHpS **LOT NUMBER:** LPFHpS1115
COMPOUND: Sodium perfluoro-1-heptanesulfonate

STRUCTURE: **CAS #:** Not available



MOLECULAR FORMULA: C₇F₁₅SO₃Na **MOLECULAR WEIGHT:** 472.10
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol
 47.6 ± 2.4 µg/ml (PFHpS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 11/06/2015
EXPIRY DATE: (mm/dd/yyyy) 11/06/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

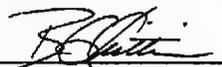
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.1% of L-PFHxS (C₆F₁₃SO₃Na) and ~ 0.2% of L-PFOS (C₈F₁₇SO₃Na).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
 B.G. Chittim **Date:** 11/09/2015
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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HOMOGENEITY:

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UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

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where x is expressed as a relative standard uncertainty of the individual parameter.

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EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

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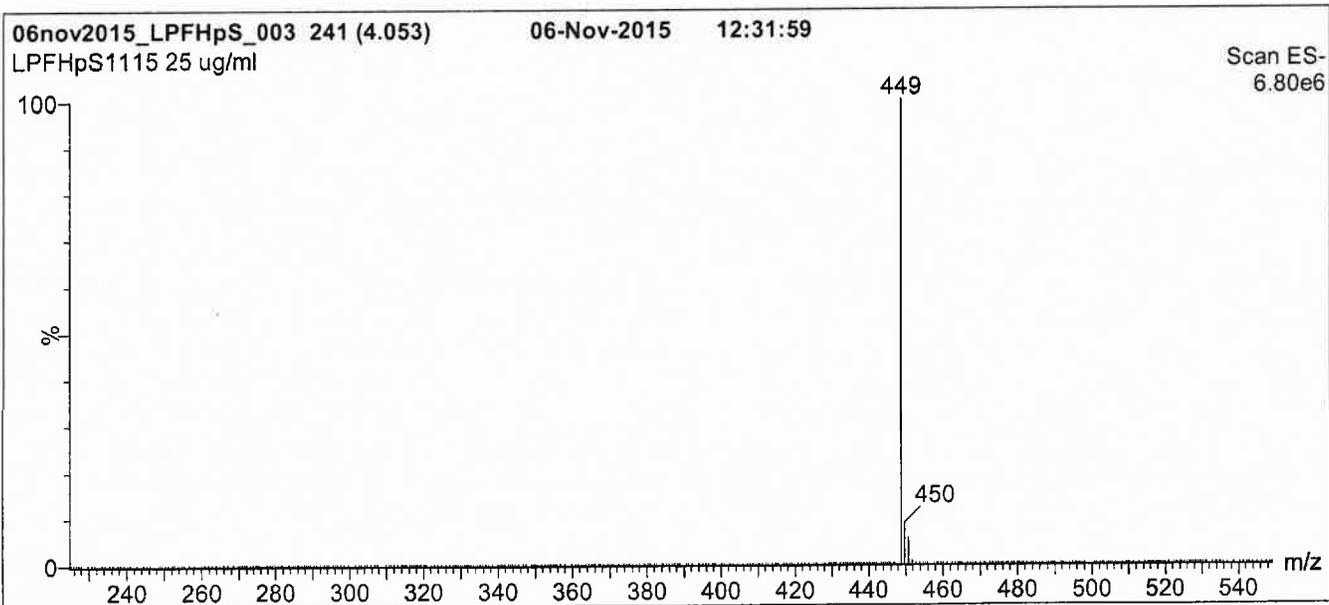
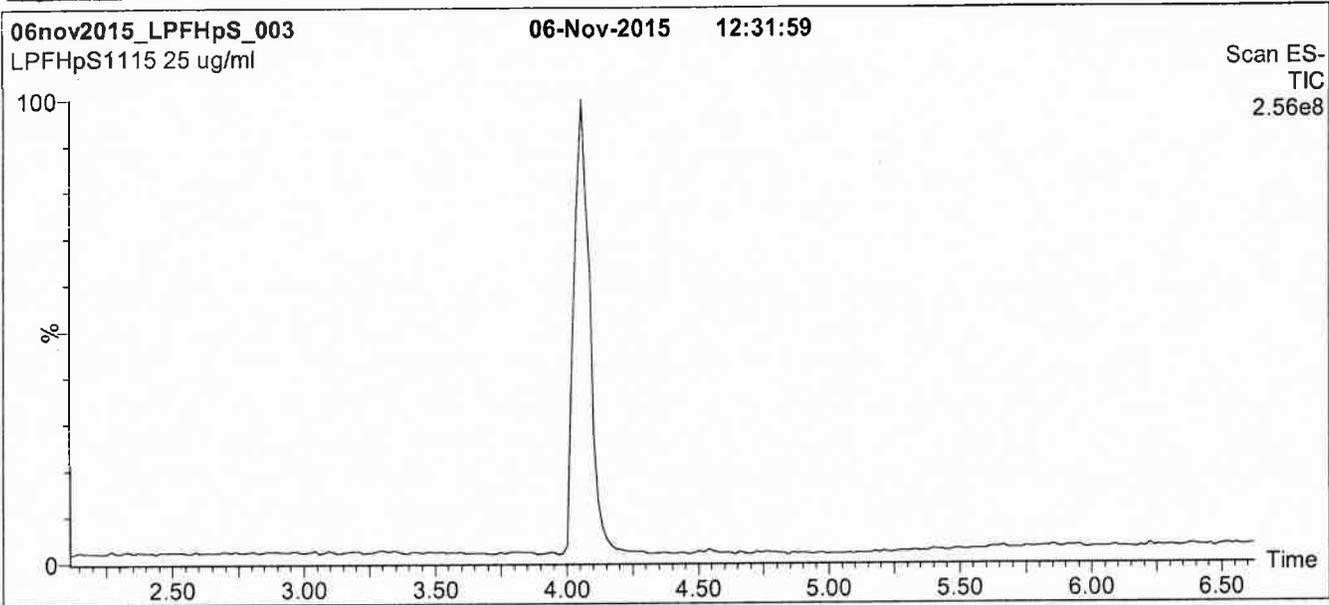
QUALITY MANAGEMENT:

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Figure 1: L-PFHpS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μm, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold
 for 2 min before returning to initial conditions in 0.5 min.
 Time: 10 min

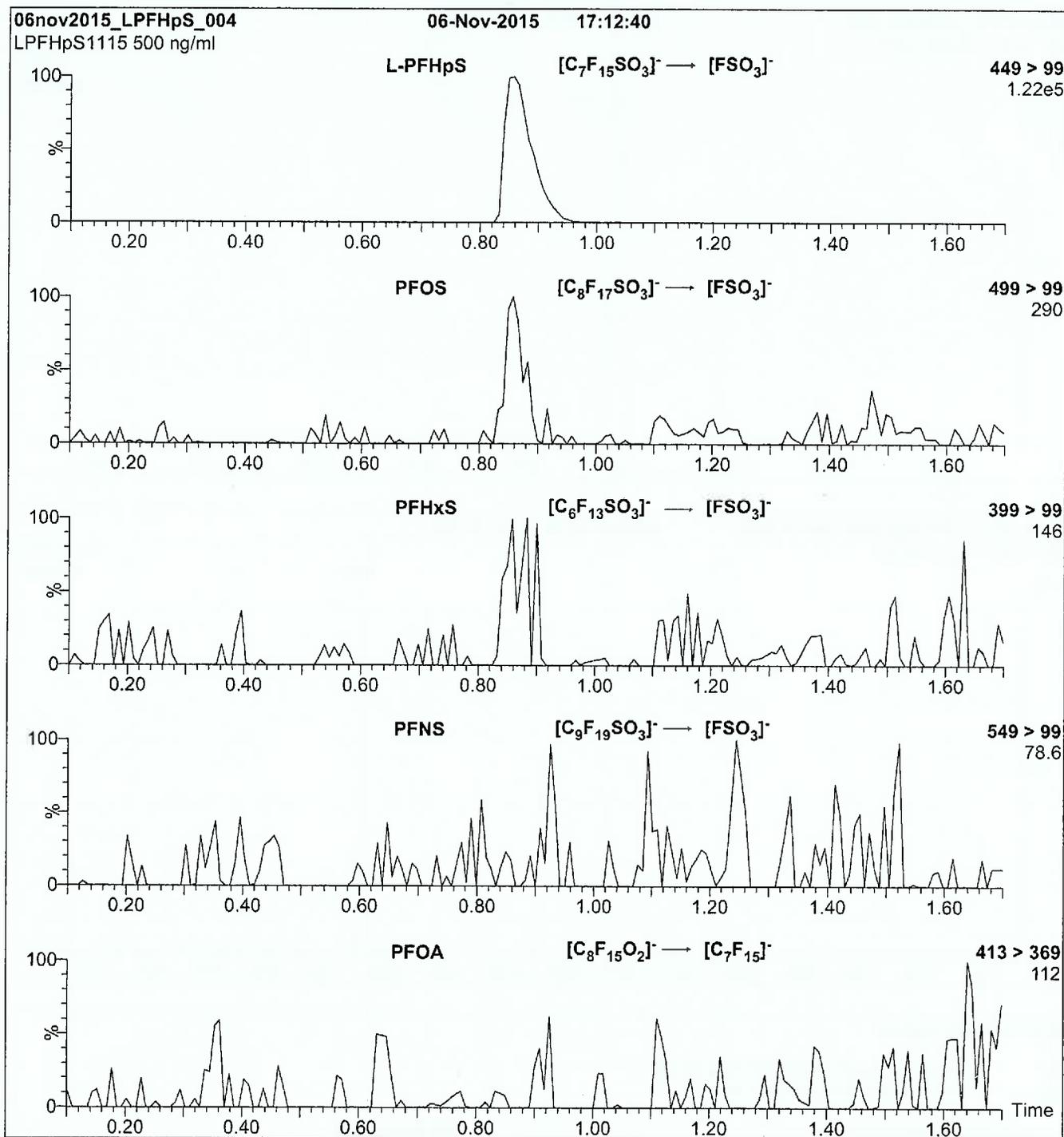
Flow: 300 μl/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 60.00
 Cone Gas Flow (l/hr) = 60
 Desolvation Gas Flow (l/hr) = 750

Figure 2: L-PFHpS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml L-PFHpS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
 Collision Energy (eV) = 35

Reagent

LCPFHpS_00009

Scanned
10/14/16 SP
R: 8BC 9/13/16



730635
ID: LCPFHpS_00009
Exp: 11/06/20 Prpd: SBC
PFHpS at 47.6ug/mL



730639
ID: LCPFHpS_00010
Exp: 11/06/20 Prpd: SBC
PFHpS at 47.6ug/mL



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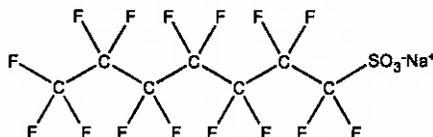
CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: L-PFHpS
COMPOUND: Sodium perfluoro-1-heptanesulfonate

LOT NUMBER: LPFHpS1115

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA: C₇F₁₅SO₃Na
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt)
47.6 ± 2.4 µg/ml (PFHpS anion)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 11/06/2015
EXPIRY DATE: (mm/dd/yyyy) 11/06/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

MOLECULAR WEIGHT: 472.10
SOLVENT(S): Methanol

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.1% of L-PFHxS (C₆F₁₃SO₃Na) and ~ 0.2% of L-PFOS (C₈F₁₇SO₃Na).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 11/09/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

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EXPIRY DATE / PERIOD OF VALIDITY:

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LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

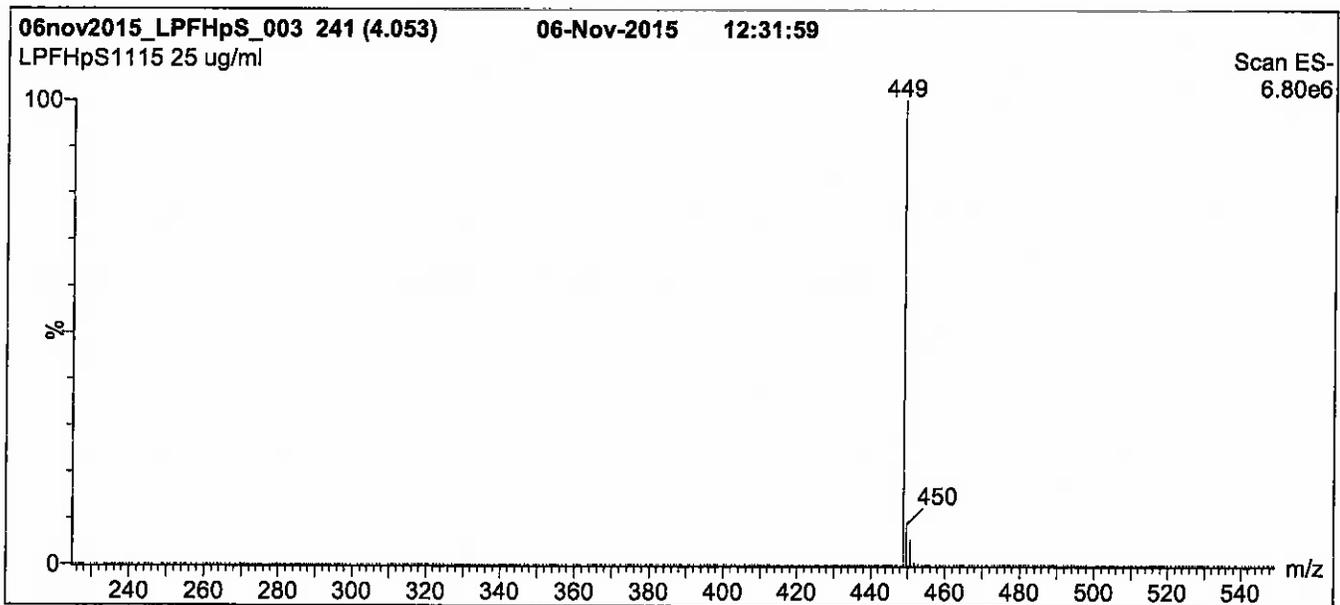
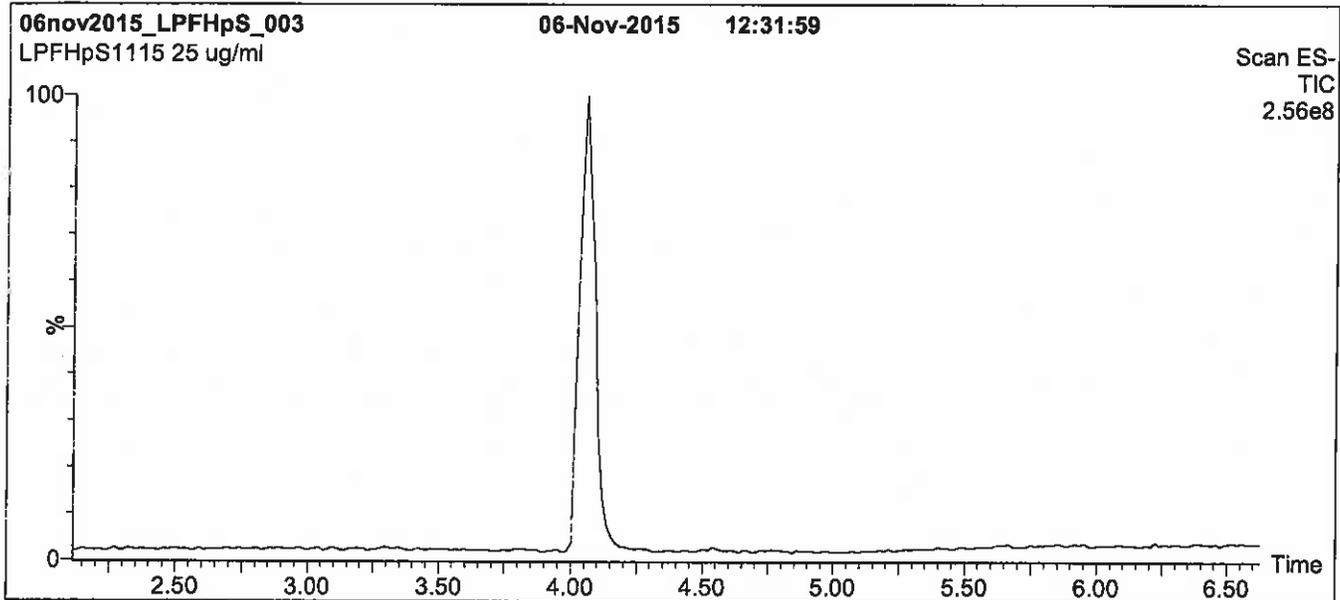
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: L-PFHpS; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold
for 2 min before returning to initial conditions in 0.5 min.
Time: 10 min

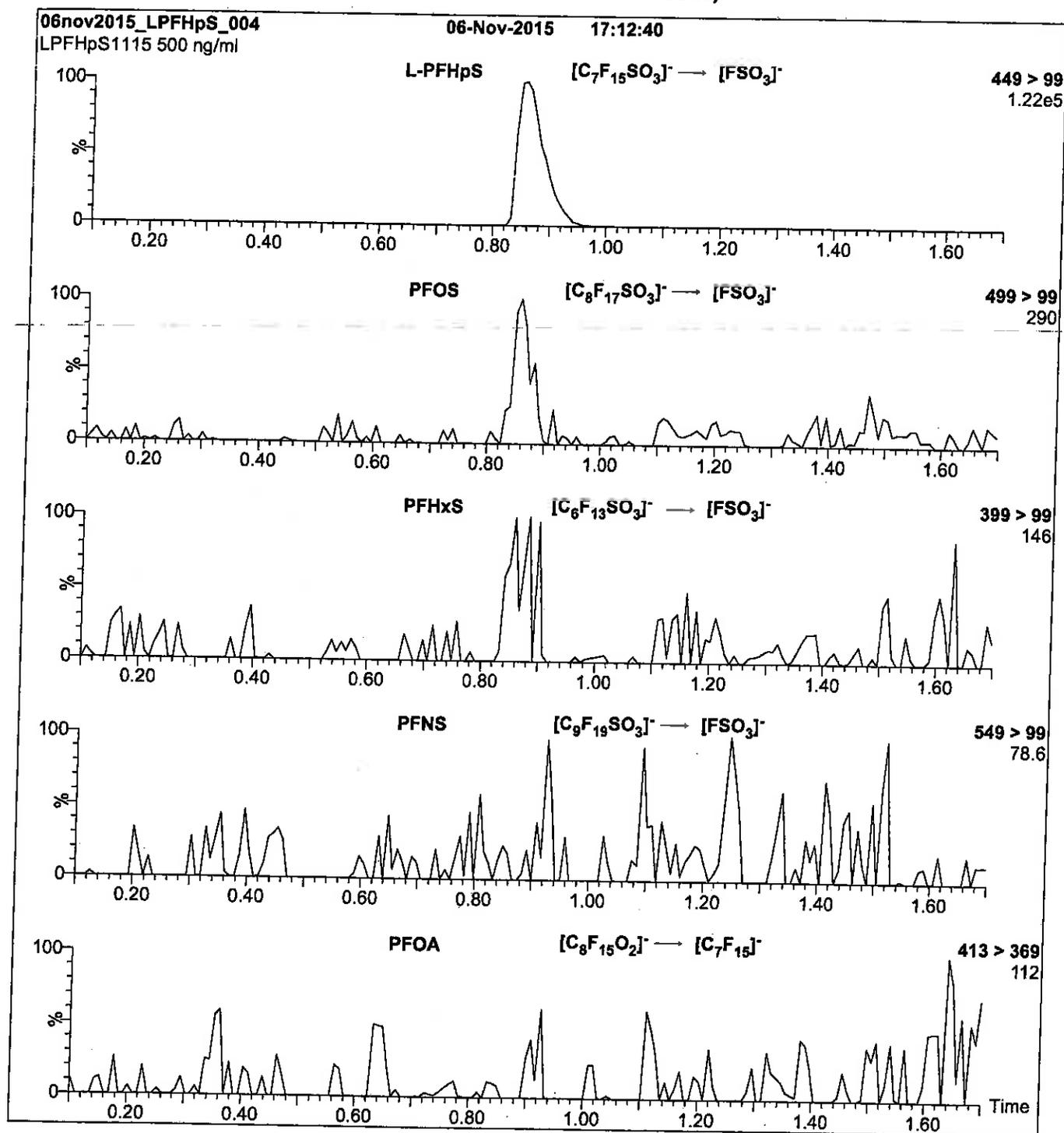
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 60.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: L-PFHpS; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml L-PFHpS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
 Collision Energy (eV) = 35

Reagent

LCPFHxA_00004



609702
 ID: LCPFHxA_00004
 Exp: 12/22/20 Prpd: CBW
 PF-n-hexanoic acid

R: 4/7/16 CBW

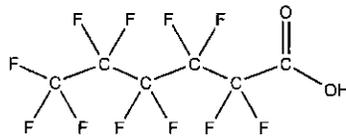


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CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFHxA **LOT NUMBER:** PFHxA1215
COMPOUND: Perfluoro-n-hexanoic acid

STRUCTURE: **CAS #:** 307-24-4



MOLECULAR FORMULA: C₆HF₁₁O₂ **MOLECULAR WEIGHT:** 314.05
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
 Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 12/22/2015
EXPIRY DATE: (mm/dd/yyyy) 12/22/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.2% of Perfluoro-n-pentanoic acid (PFPeA).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: B.G. Crittitt Date: 12/23/2015
 B.G. Crittitt (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

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LIMITED WARRANTY:

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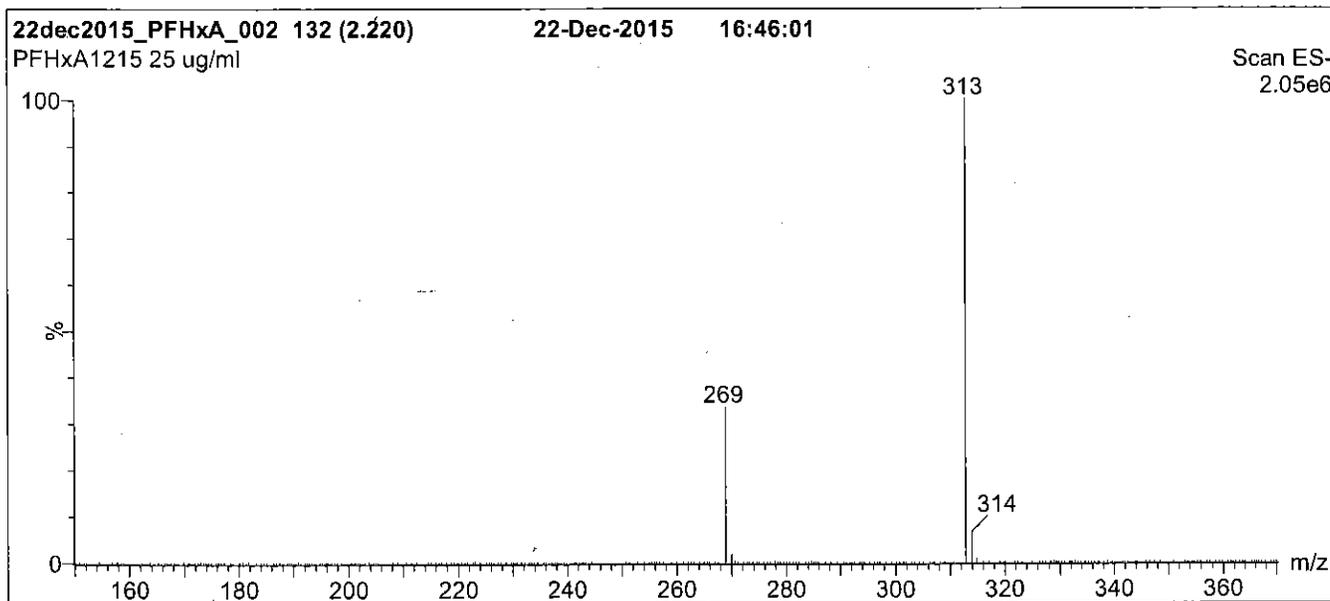
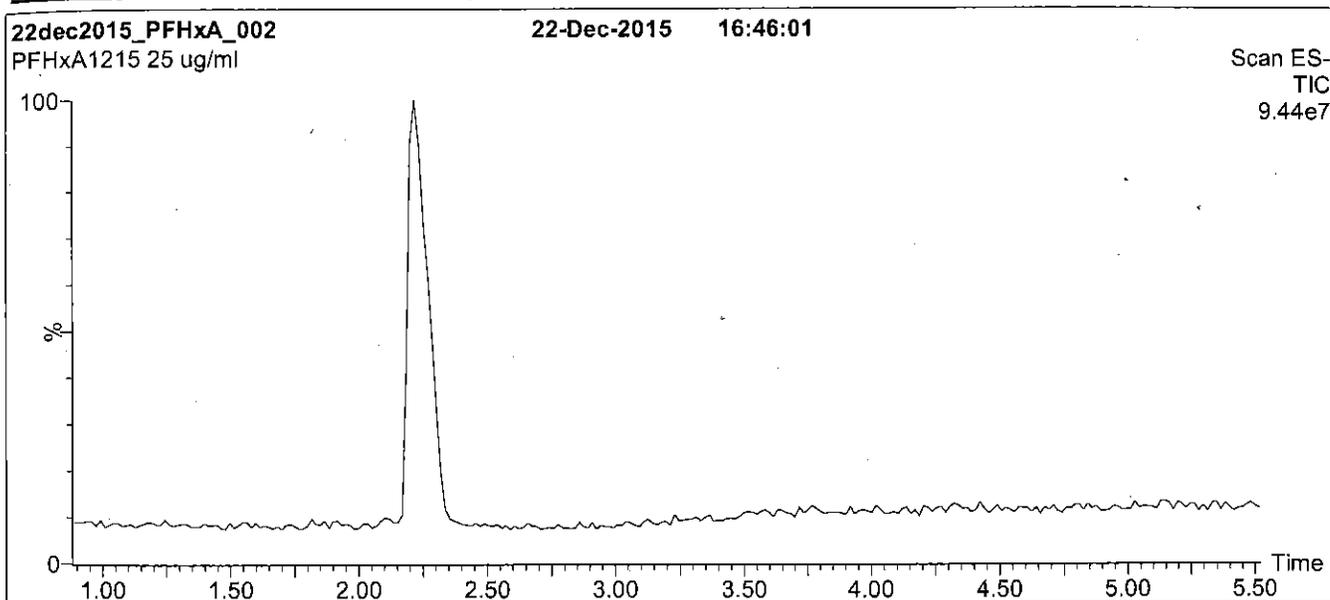
QUALITY MANAGEMENT:

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Figure 1: PFHxA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

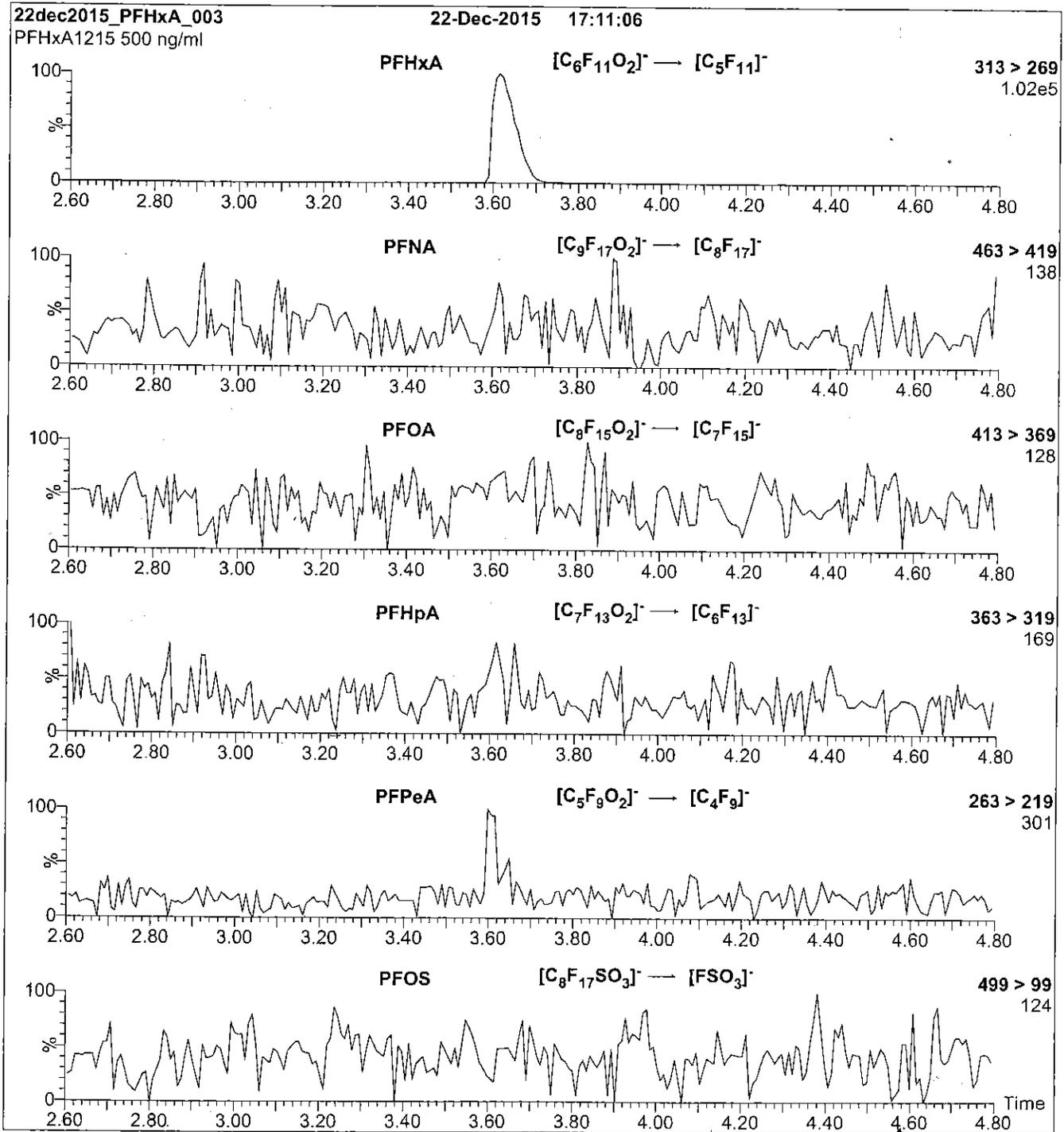
Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm
Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min
before returning to initial conditions in 0.5 min.
Time: 10 min

Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)
Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 100
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFHxA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFHxA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
 Collision Energy (eV) = 10

Reagent

LCPFHxDA_00006

R: SBC 9/13/16

Scanned 10/14/16



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730630
ID: LCPFHxDA_00006
Exp: 05/25/21 Prod: SBC
PFHxDA stock 50ug/mL

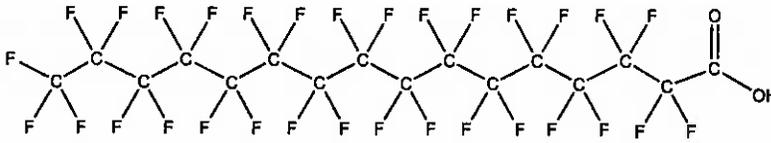


730631
ID: LCPFHxDA_00007
Exp: 05/25/21 Prod: SBC
PFHxDA stock 50ug/mL

**CERTIFICATE OF ANALYSIS
DOCUMENTATION**

PRODUCT CODE: PFHxDA **LOT NUMBER:** PFHxDA0516
COMPOUND: Perfluoro-n-hexadecanoic acid

STRUCTURE: **CAS #:** 67905-19-5



MOLECULAR FORMULA: C₁₆H₃₁O₂ **MOLECULAR WEIGHT:** 814.13
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 05/25/2016
EXPIRY DATE: (mm/dd/yyyy) 05/25/2021
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.4% of PFODA.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chrifim

Date: 05/27/2016
(mm/dd/yyyy)

**Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com**

INTENDED USE:

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HAZARDS:

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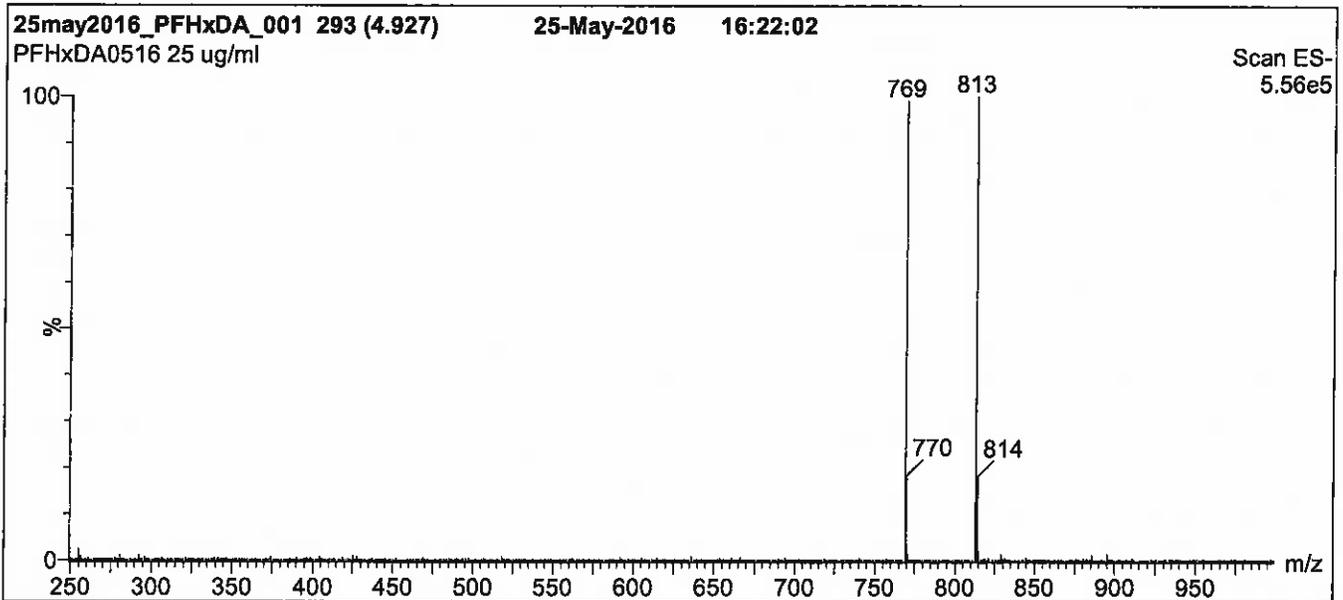
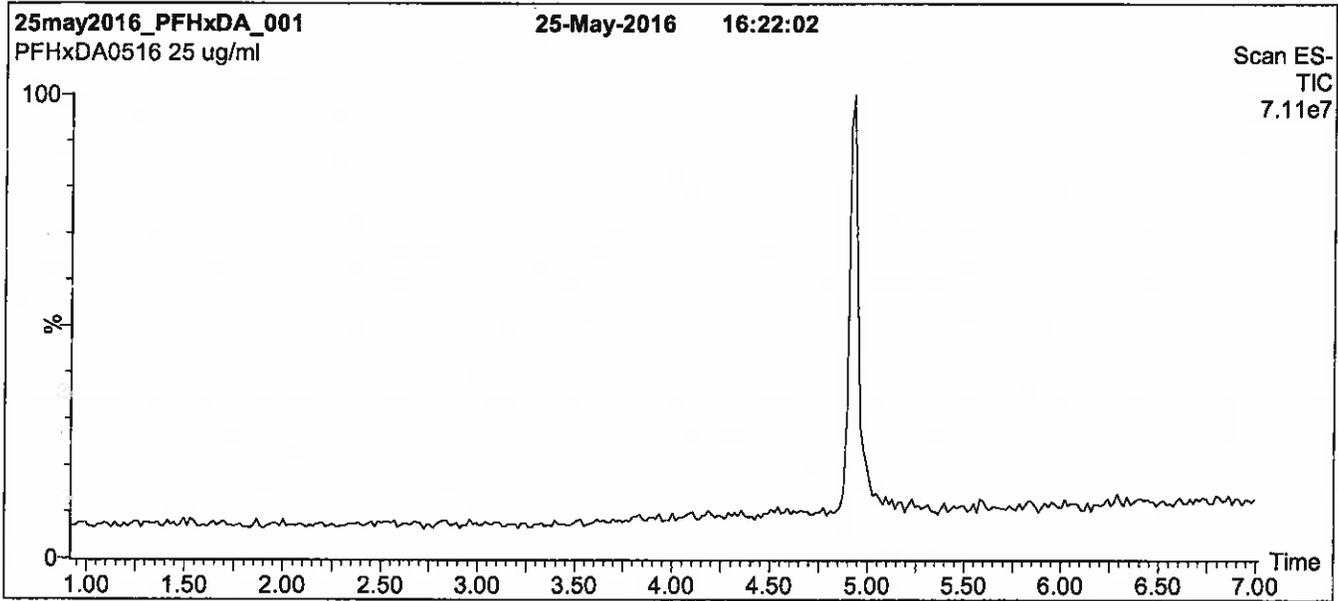
QUALITY MANAGEMENT:

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Figure 1: PFHxDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 70% (80:20 MeOH:ACN) / 30% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 95% organic over 6 min and hold for 2.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

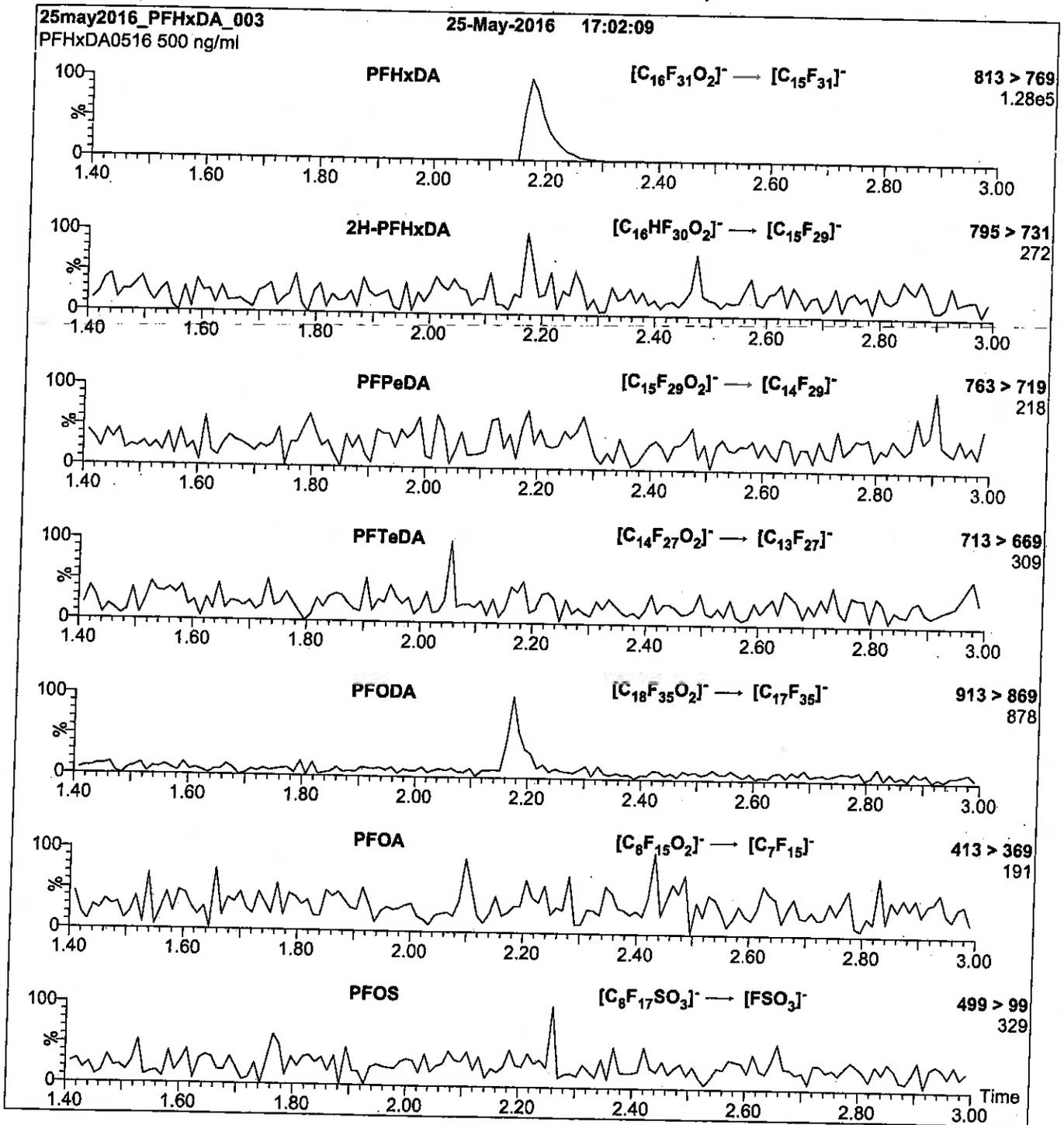
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 1250 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 25.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFHxDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFHxDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.66e-3
 Collision Energy (eV) = 15

Reagent

LCPFHxS-br_00001



PS 12/9/15 SW

566007
ID: LCPFHxS-br_00001
Exp: 07/03/20 Pppl: CBW
Potassium Perfluorohexane



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CERTIFICATE OF ANALYSIS DOCUMENTATION

br-PFHxSK

Potassium Perfluorohexanesulfonate Solution/Mixture of Linear and Branched Isomers

PRODUCT CODE: br-PFHxSK
LOT NUMBER: brPFHxSK0615
CONCENTRATION: 50.0 ± 2.5 µg/ml (total potassium salt)
 45.5 ± 2.3 µg/ml (total PFHxS anion)
SOLVENT(S): Methanol
DATE PREPARED: (mm/dd/yyyy) 06/29/2015
LAST TESTED: (mm/dd/yyyy) 07/03/2015
EXPIRY DATE: (mm/dd/yyyy) 07/03/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DESCRIPTION:

The chemical purity has been determined to be ≥98% perfluorohexanesulfonate linear and branched isomers. The full name, structure and percent composition for each of the identified isomeric components are given in Table A.

DOCUMENTATION/ DATA ATTACHED:

Table A: Isomeric Components and Percent Composition by ¹⁹F-NMR
Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS Data
Figure 3: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.5% of perfluoro-1-pentanesulfonate and ~ 0.2% of perfluoro-1-octanesulfonate.
- CAS#: 3871-99-6 (for linear isomer; potassium salt).

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INTENDED USE:

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Table A: br-PFHxSK; Isomeric Components and Percent Composition (by ¹⁹F-NMR)*

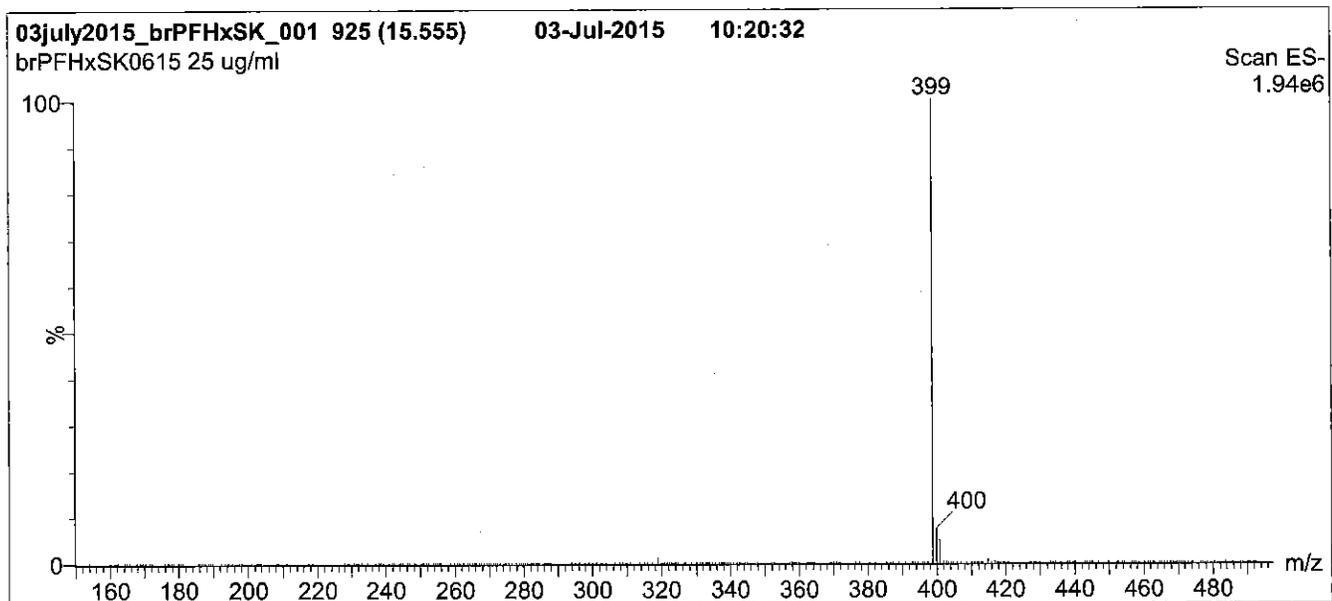
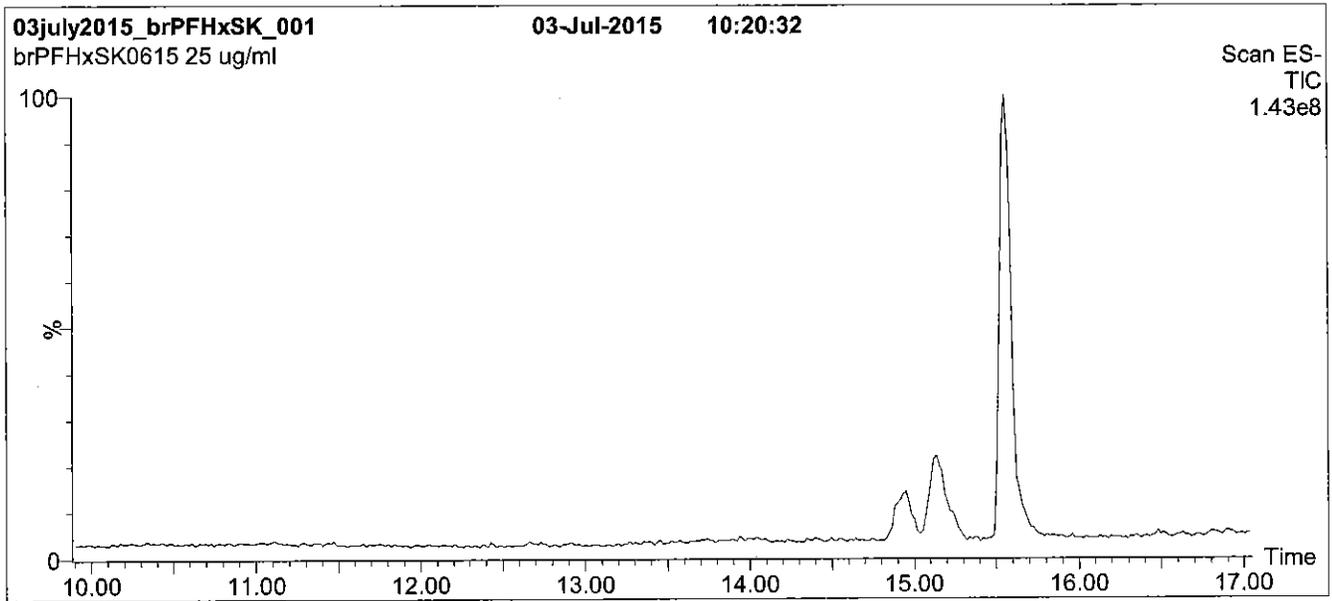
Isomer	Name	Structure	Percent Composition by ¹⁹ F-NMR
1	Potassium perfluoro-1-hexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ ⁻ K ⁺	81.1
2	Potassium 1-trifluoromethylperfluoropentanesulfonate**	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	2.9
3	Potassium 2-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	1.4
4	Potassium 3-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	5.0
5	Potassium 4-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	8.9
6	Potassium 3,3-di(trifluoromethyl)perfluorobutanesulfonate	$\begin{array}{c} \text{CF}_3 \\ \\ \text{CF}_3\text{CCF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	0.2
7	Other Unidentified Isomers		0.5

* Percent of total perfluorohexanesulfonate isomers only.
 ** Systematic Name: Potassium perfluorohexane-2-sulfonate.

Certified By: 
 B.G. Chittim

Date: 07/15/2015
(mm/dd/yyyy)

Figure 1: br-PFHxSK; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 20% (80:20 MeOH:ACN) / 80% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 50% organic over 14 min. Ramp to
90% organic over 3 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 20 min

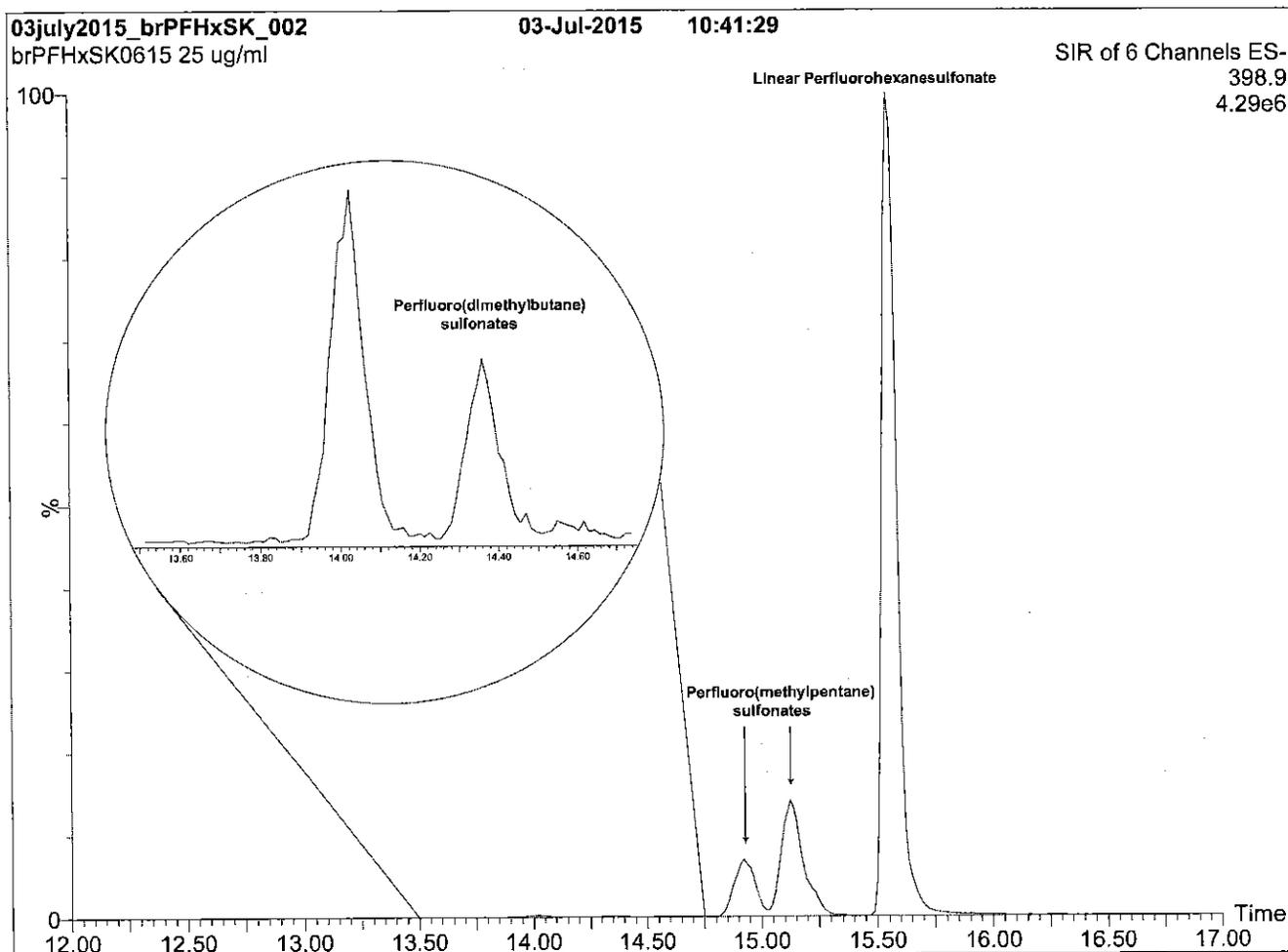
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 50.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: br-PFHxSK; LC/MS Data



Conditions for Figure 2:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μm, 2.1 x 100 mm

Mobile phase: Gradient
Start: 20% (80:20 MeOH:ACN) / 80% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 50% organic over 14 min. Ramp to
90% organic over 3 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 20 min

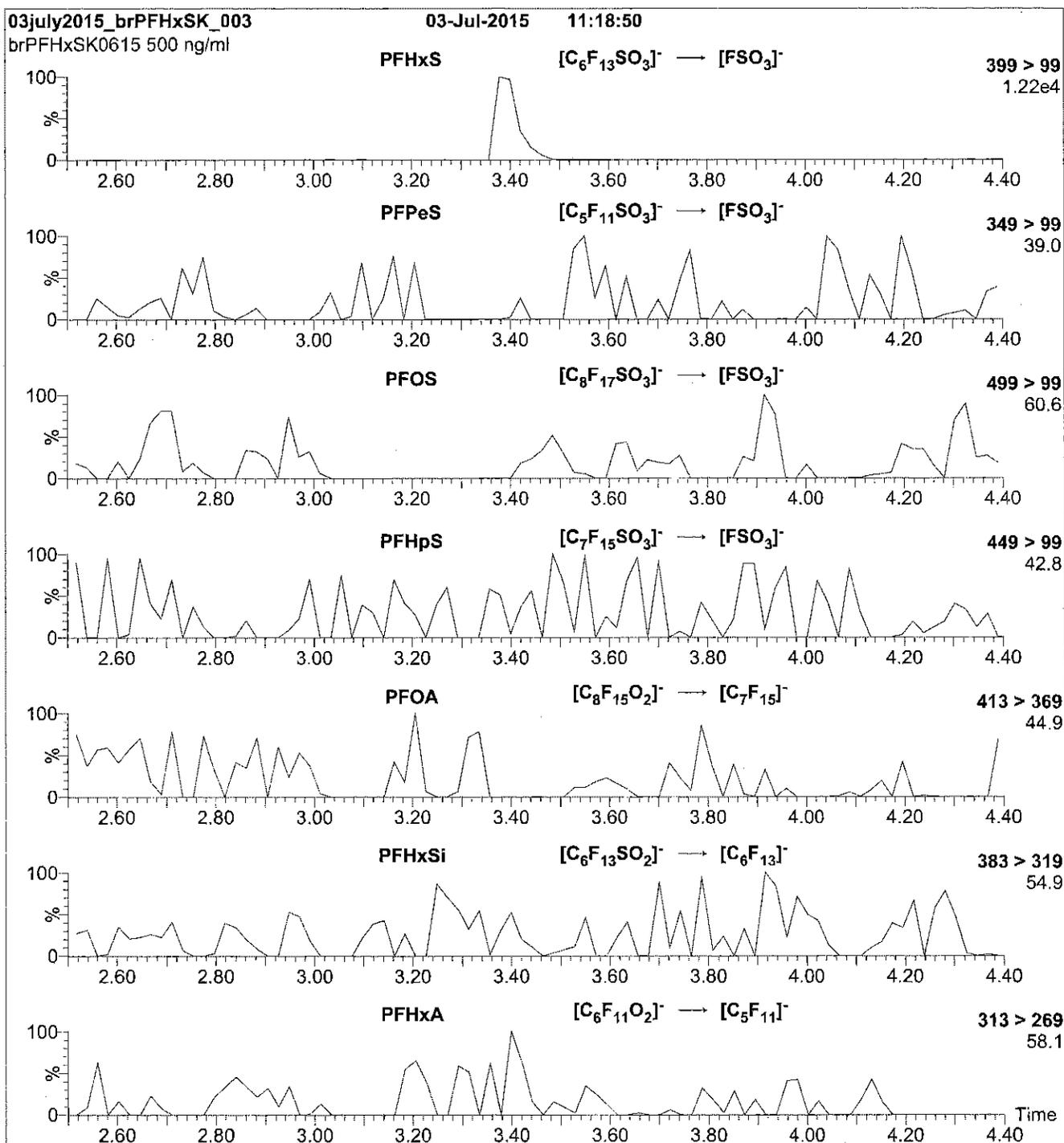
Flow: 300 μl/min

MS Parameters

Experiment: SIR (6 channels)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 50.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 3: br-PFHxSK; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 3:

Injection: Direct loop injection
10 μ l (500 ng/ml br-PFHxSK)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 30

Reagent

LCPFHxS-br_00002

SBC
R: 9/13/16



730513
ID: LCPFHxS-br_00002
Exp: 07/03/20 Ppfd: SBC
Potassium Perfluorohexane



730514
ID: LCPFHxS-br_00003
Exp: 07/03/20 Ppfd: SBC
Potassium Perfluorohexane



WELLINGTON
LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

br-PFHxSK

**Potassium Perfluorohexanesulfonate
Solution/Mixture of Linear and
Branched Isomers**

PRODUCT CODE: br-PFHxSK
LOT NUMBER: brPFHxSK0615
CONCENTRATION: 50.0 ± 2.5 µg/ml (total potassium salt)
45.5 ± 2.3 µg/ml (total PFHxS anion)
SOLVENT(S): Methanol
DATE PREPARED: (mm/dd/yyyy) 06/29/2015
LAST TESTED: (mm/dd/yyyy) 07/03/2015
EXPIRY DATE: (mm/dd/yyyy) 07/03/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DESCRIPTION:

The chemical purity has been determined to be ≥98% perfluorohexanesulfonate linear and branched isomers. The full name, structure and percent composition for each of the identified isomeric components are given in Table A.

DOCUMENTATION/ DATA ATTACHED:

Table A: Isomeric Components and Percent Composition by ¹⁹F-NMR
Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS Data
Figure 3: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains ~ 0.5% of perfluoro-1-pentanesulfonate and ~ 0.2% of perfluoro-1-octanesulfonate.
- CAS#: 3871-99-6 (for linear isomer; potassium salt).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compounds it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Table A: br-PFHxSK; Isomeric Components and Percent Composition (by ¹⁹F-NMR)*

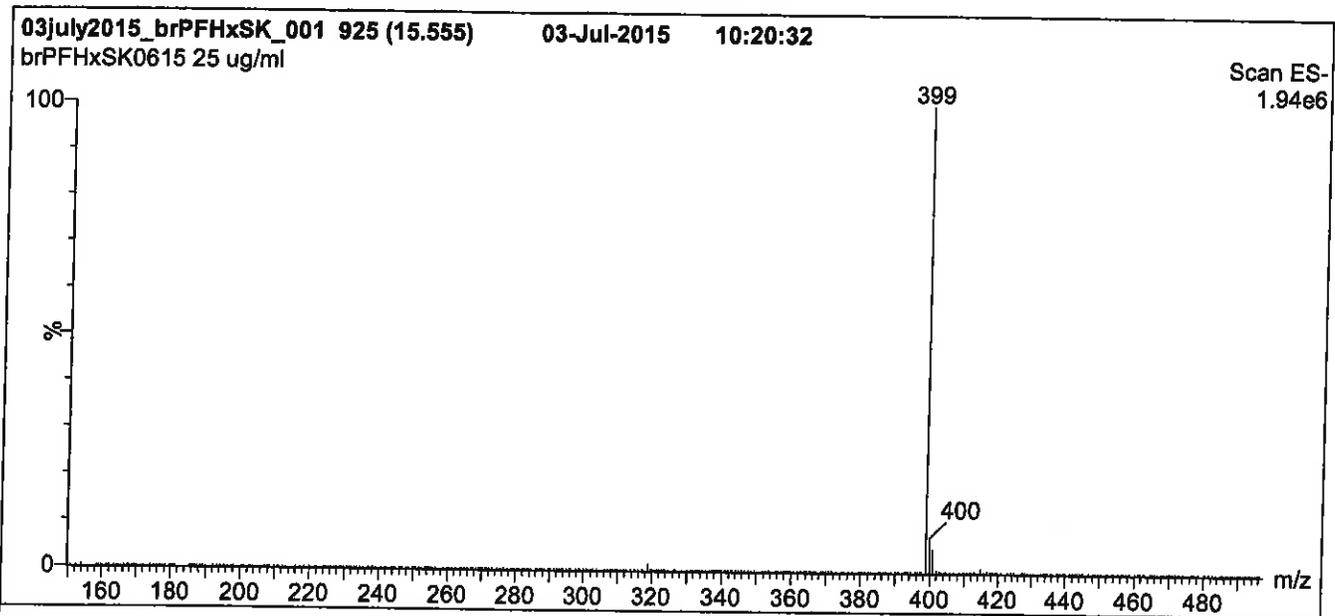
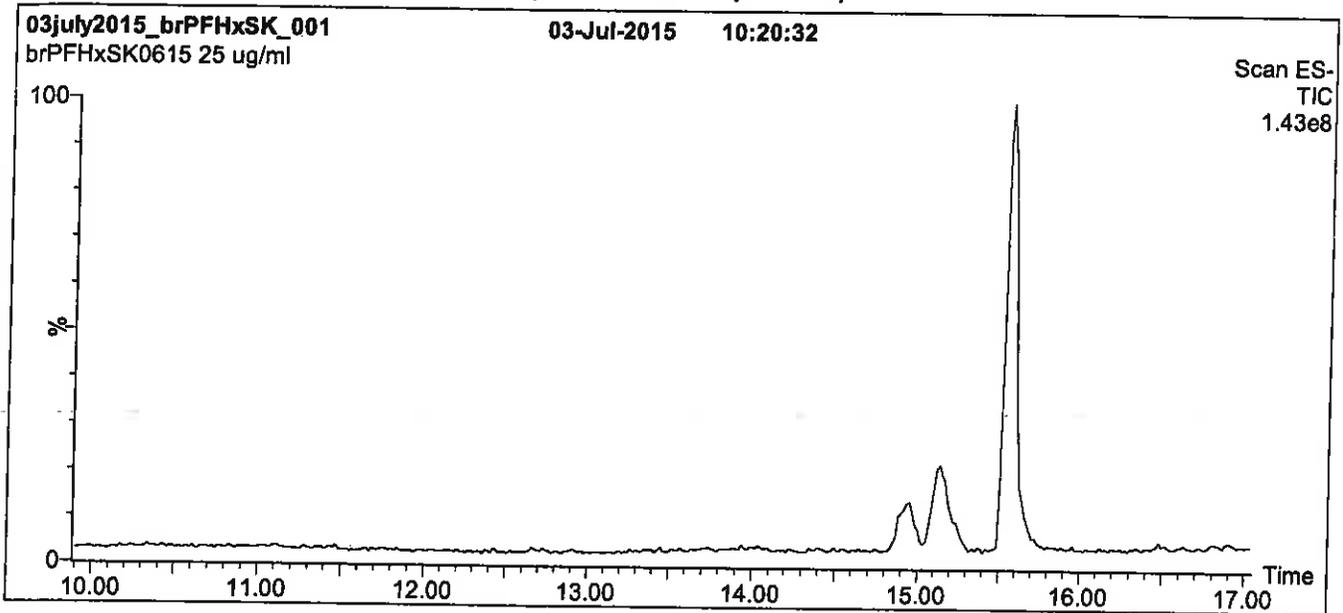
Isomer	Name	Structure	Percent Composition by ¹⁹ F-NMR
1	Potassium perfluoro-1-hexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ ⁻ K ⁺	81.1
2	Potassium 1-trifluoromethylperfluoropentanesulfonate**	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	2.9
3	Potassium 2-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	1.4
4	Potassium 3-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	5.0
5	Potassium 4-trifluoromethylperfluoropentanesulfonate	$\begin{array}{c} \text{CF}_3\text{CF}_2\text{CF}_2\text{CF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	8.9
6	Potassium 3,3-di(trifluoromethyl)perfluorobutanesulfonate	$\begin{array}{c} \text{CF}_3 \\ \\ \text{CF}_3\text{CCF}_2\text{CF}_2\text{SO}_3^-\text{K}^+ \\ \\ \text{CF}_3 \end{array}$	0.2
7	Other Unidentified Isomers		0.5

* Percent of total perfluorohexanesulfonate isomers only.
 ** Systematic Name: Potassium perfluorohexane-2-sulfonate.

Certified By: 
 B.G. Chittim

Date: 07/15/2015
(mm/dd/yyyy)

Figure 1: br-PFHxSK; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 20% (80:20 MeOH:ACN) / 80% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 50% organic over 14 min. Ramp to
90% organic over 3 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 20 min

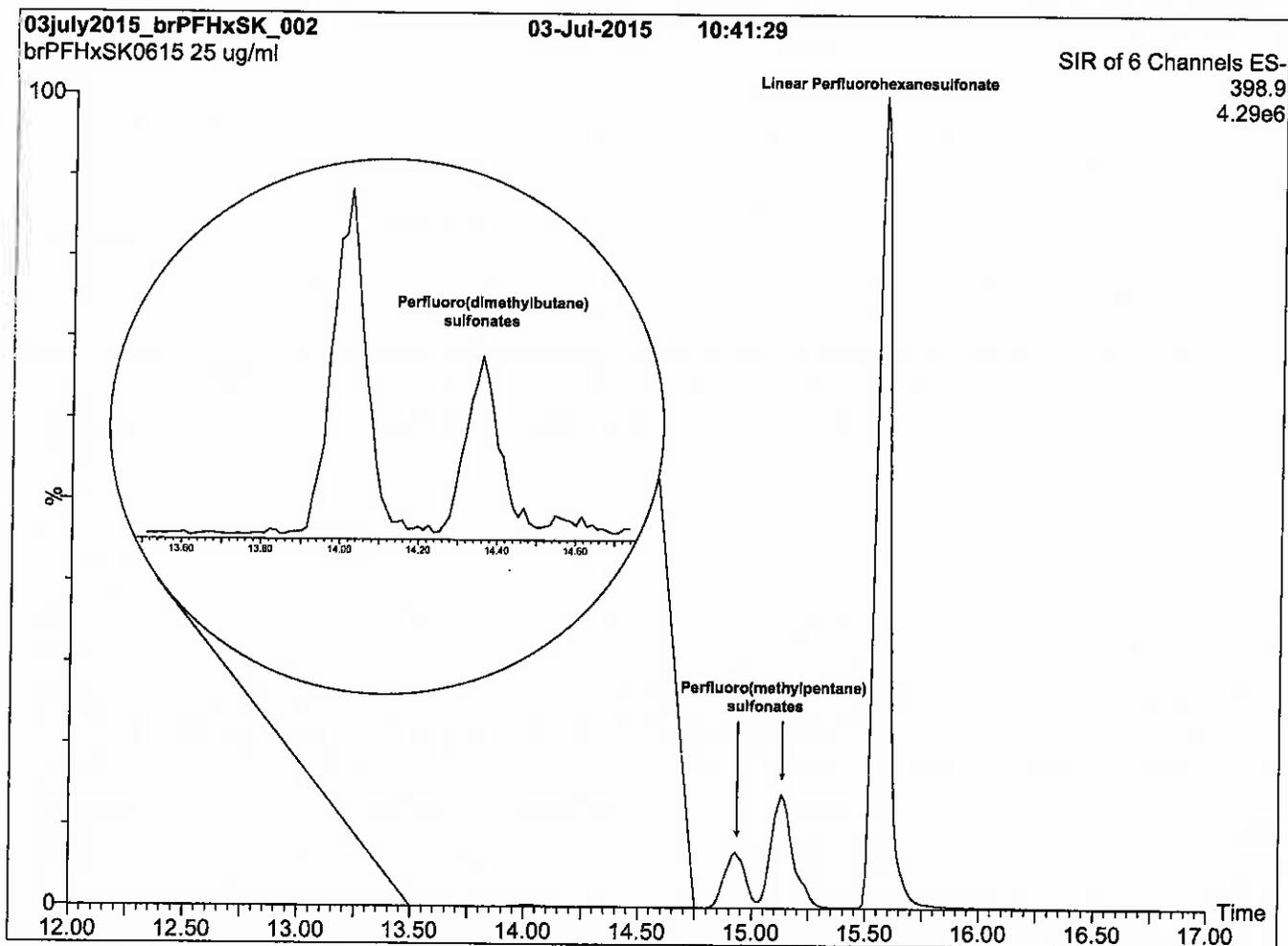
MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 50.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Flow: 300 μ l/min

Figure 2: br-PFHxSK; LC/MS Data



Conditions for Figure 2:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

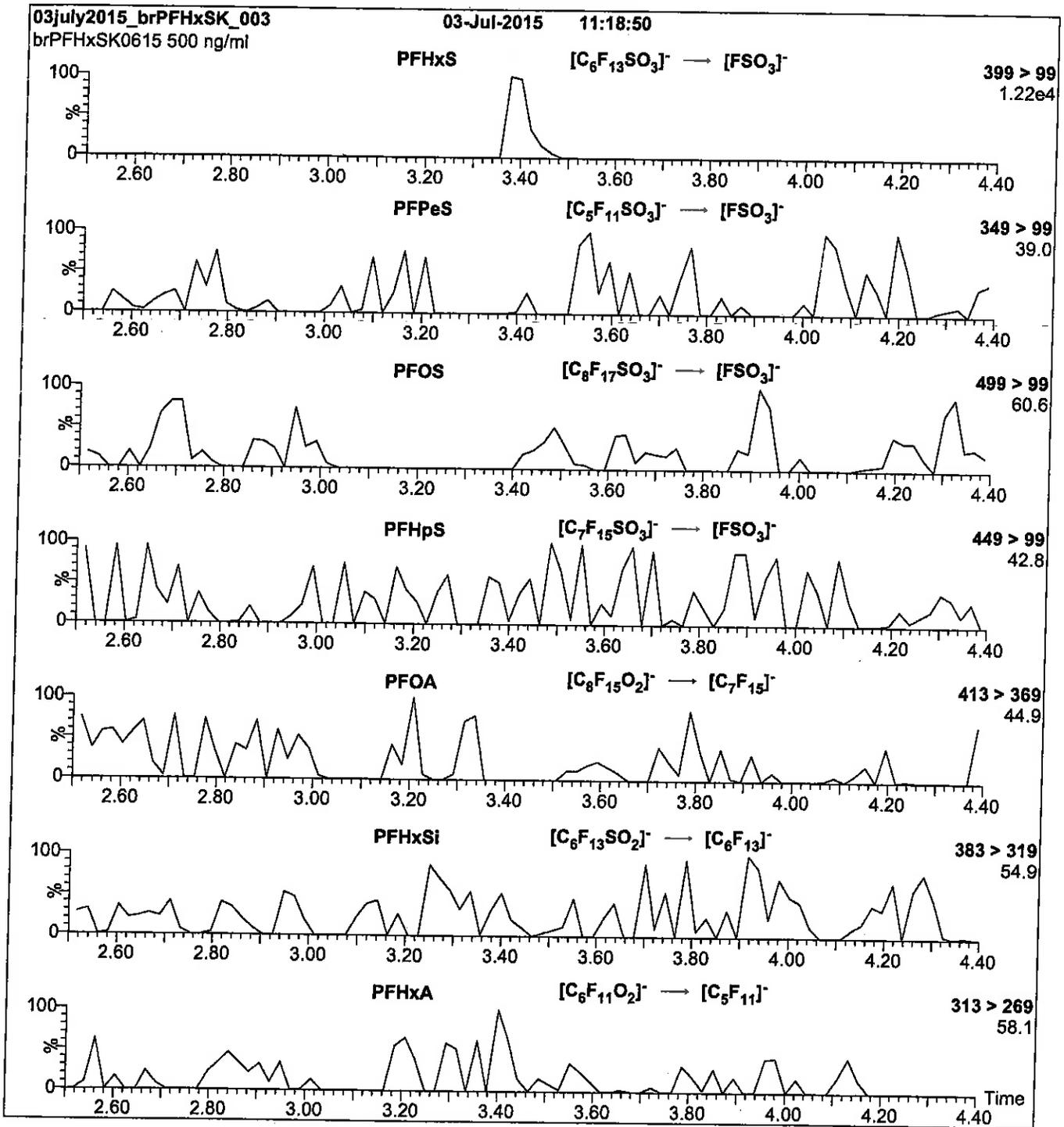
Mobile phase: Gradient
 Start: 20% (80:20 MeOH:ACN) / 80% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 50% organic over 14 min. Ramp to
 90% organic over 3 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 20 min

Flow: 300 μ l/min

MS Parameters

Experiment: SIR (6 channels)
 Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 50.00
 Cone Gas Flow (l/hr) = 60
 Desolvation Gas Flow (l/hr) = 750

Figure 3: br-PFHxSK; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 3:

Injection: Direct loop injection
10 μ l (500 ng/ml br-PFHxSK)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 30

Reagent

LCPFNA_00005



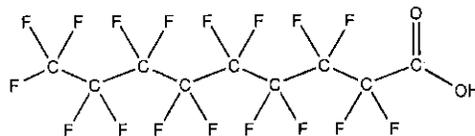
R: 4/7/16 CBW

609703

ID: LCPFNA_00005

Exp: 10/23/20 Prod: CBW

PF-n-nonanoic acid

**WELLINGTON**
LABORATORIES**CERTIFICATE OF ANALYSIS**
DOCUMENTATION**PRODUCT CODE:** PFNA
COMPOUND: Perfluoro-n-nonanoic acid**LOT NUMBER:** PFNA1015**STRUCTURE:****CAS #:** 375-95-1**MOLECULAR FORMULA:** C₉H_{F₁₇}O₂
CONCENTRATION: 50 ± 2.5 µg/ml**MOLECULAR WEIGHT:** 464.08
SOLVENT(S): Methanol
Water (<1%)**CHEMICAL PURITY:** >98%
LAST TESTED: (mm/dd/yyyy) 10/23/2015
EXPIRY DATE: (mm/dd/yyyy) 10/23/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place**DOCUMENTATION/ DATA ATTACHED:**Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)**ADDITIONAL INFORMATION:**

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.1% of perfluoro-n-octanoic acid (PFOA) and < 0.1% of perfluoro-n-heptanoic acid (PFHpA).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**Certified By:**

B.G. Chittim
Date: 10/30/2015
(mm/dd/yyyy)Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON 'N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

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$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

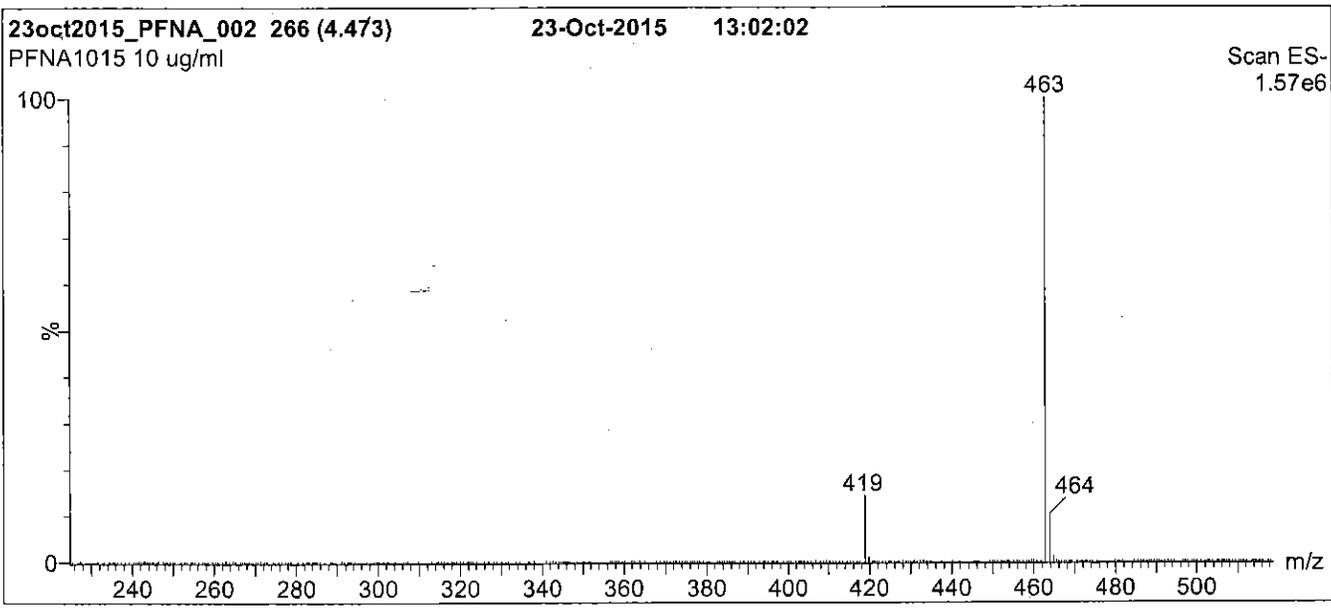
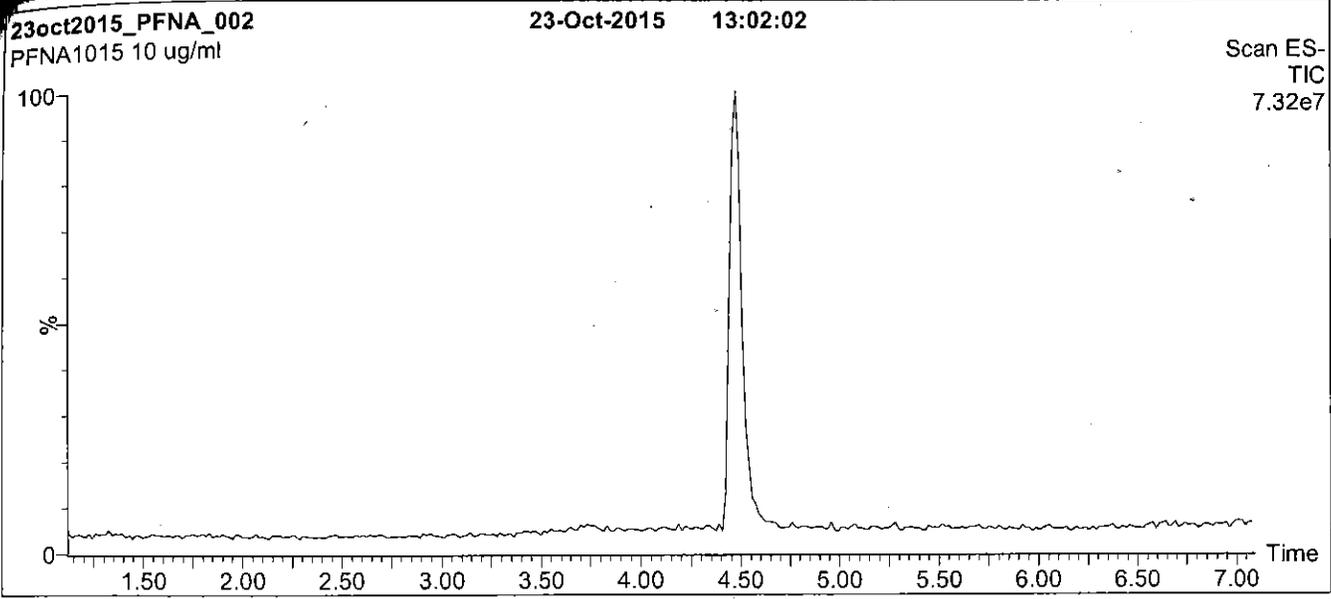
QUALITY MANAGEMENT:

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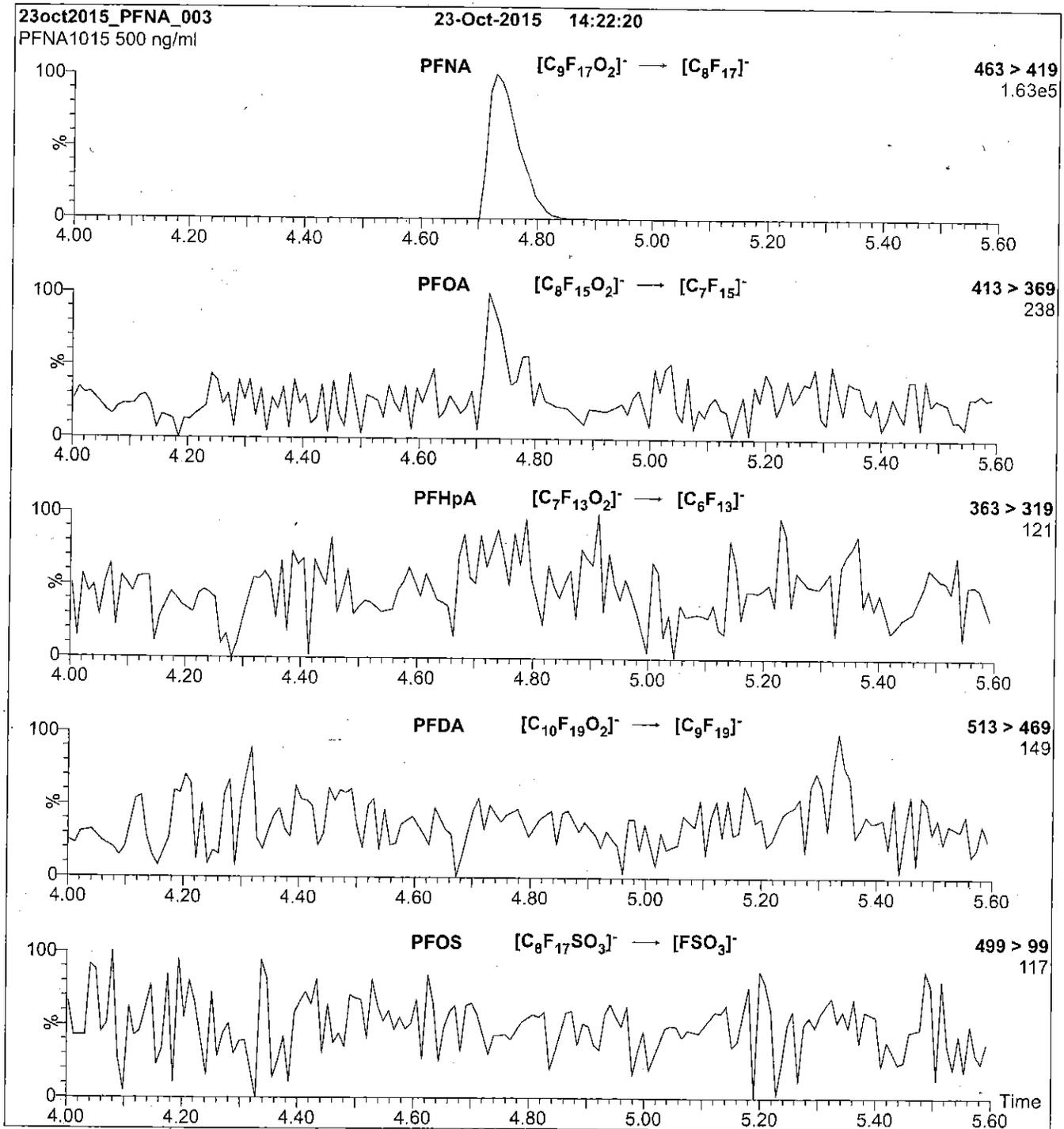
Figure 1: PFNA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC:	Waters Acquity Ultra Performance LC
MS:	Micromass Quattro <i>micro</i> API MS
Chromatographic Conditions	
Column:	Acquity UPLC BEH Shield RP ₁₈ 1.7 μ m, 2.1 x 100 mm
Mobile phase:	Gradient Start: 50% (80:20 MeOH:ACN) / 50% H ₂ O (both with 10 mM NH ₄ OAc buffer) Ramp to 90% organic over 7 min and hold for 2 min before returning to initial conditions in 0.5 min. Time: 10 min
Flow:	300 μ l/min
MS Parameters	
Experiment:	Full Scan (225 - 850 amu)
Source:	Electrospray (negative)
Capillary Voltage (kV):	2.00
Cone Voltage (V):	15.00
Cone Gas Flow (l/hr):	50
Desolvation Gas Flow (l/hr):	750

Figure 2: PFNA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFNA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.28e-3
 Collision Energy (eV) = 11

Reagent

LCPFOA_00006

R-716/16 CBW



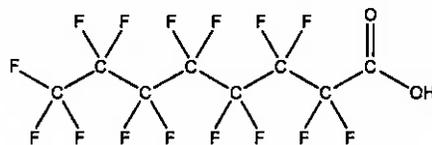
671577
ID: LCPFOA_00006
Exp: 11/06/20 Prod: CBW
PF-n-octanoic acid



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFOA **LOT NUMBER:** PFOA1115
COMPOUND: Perfluoro-n-octanoic acid
STRUCTURE: **CAS #:** 335-67-1



MOLECULAR FORMULA: $C_8HF_{15}O_2$ **MOLECULAR WEIGHT:** 414.07
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 11/06/2015
EXPIRY DATE: (mm/dd/yyyy) 11/06/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 11/11/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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where x is expressed as a relative standard uncertainty of the individual parameter.

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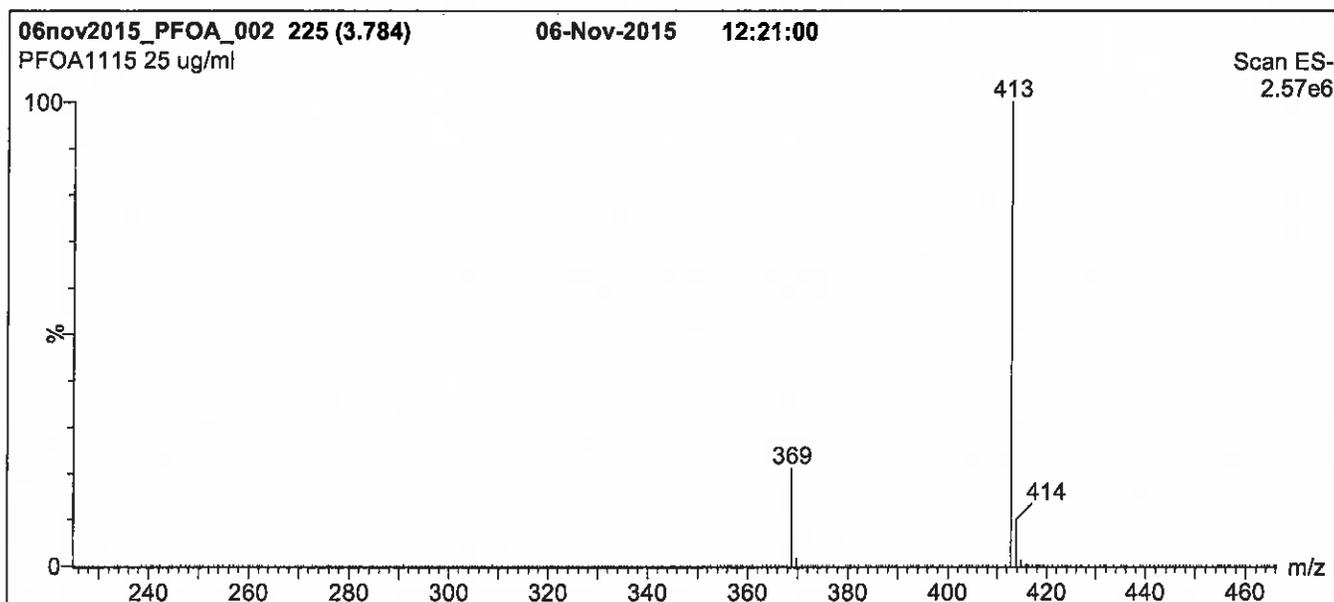
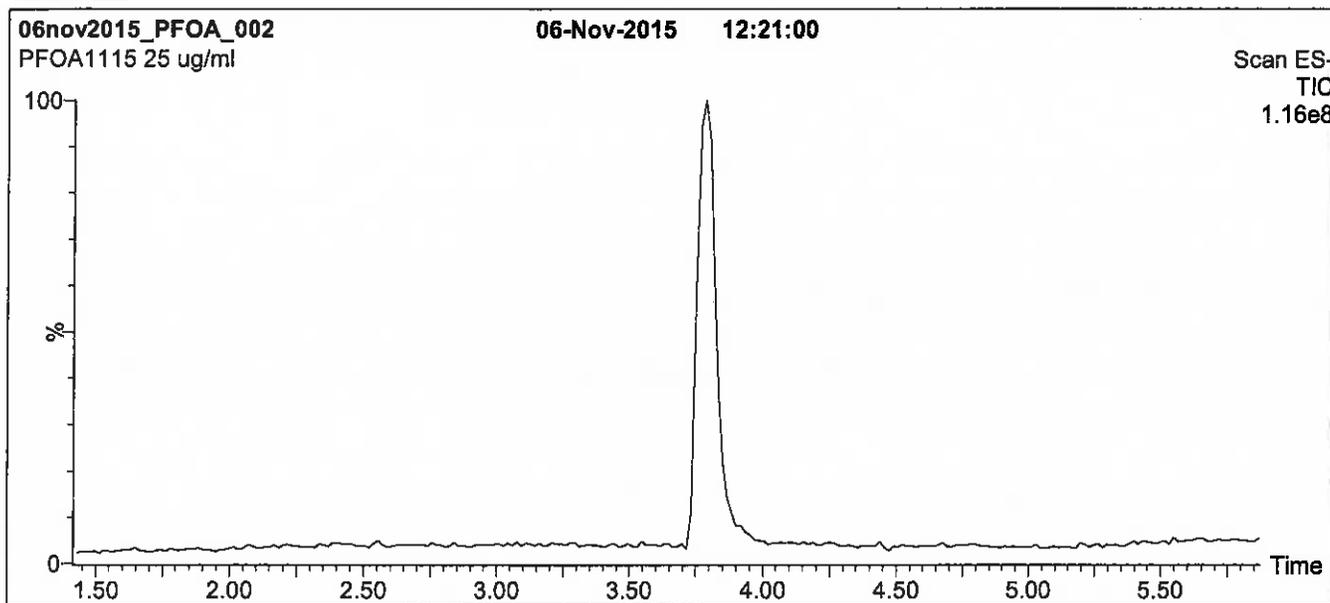
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: PFOA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro micro API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 2 min before returning to initial conditions in 0.5 min.
 Time: 10 min

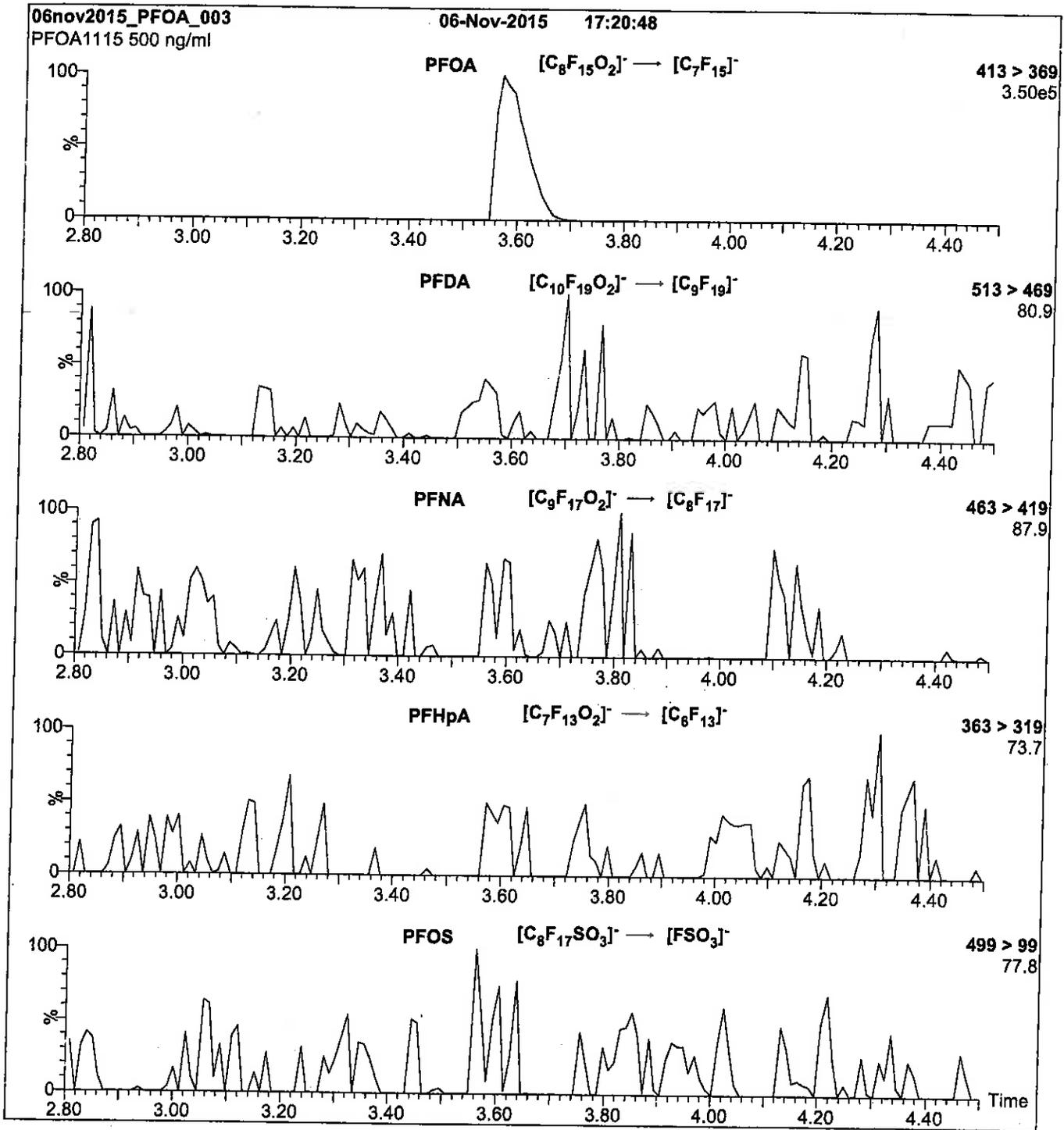
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 100
 Desolvation Gas Flow (l/hr) = 750

Figure 2: PFOA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFOA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.17e-3
Collision Energy (eV) = 10

Reagent

LCPFODA_00005

605234
ID: LCPFODA_00005
Exp: 01/30/20 Prod: CBW
PFODA stock 50ug/ml

Rec. 3/20/16 JRB

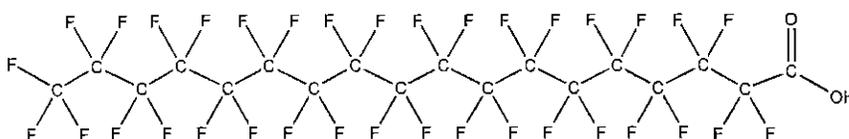


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CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFODA **LOT NUMBER:** PFODA0115
COMPOUND: Perfluoro-n-octadecanoic acid

STRUCTURE: **CAS #:** 16517-11-6



MOLECULAR FORMULA: $C_{18}H_{35}O_2$ **MOLECULAR WEIGHT:** 914.14
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/30/2015
EXPIRY DATE: (mm/dd/yyyy) 01/30/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

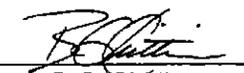
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 03/25/2015
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

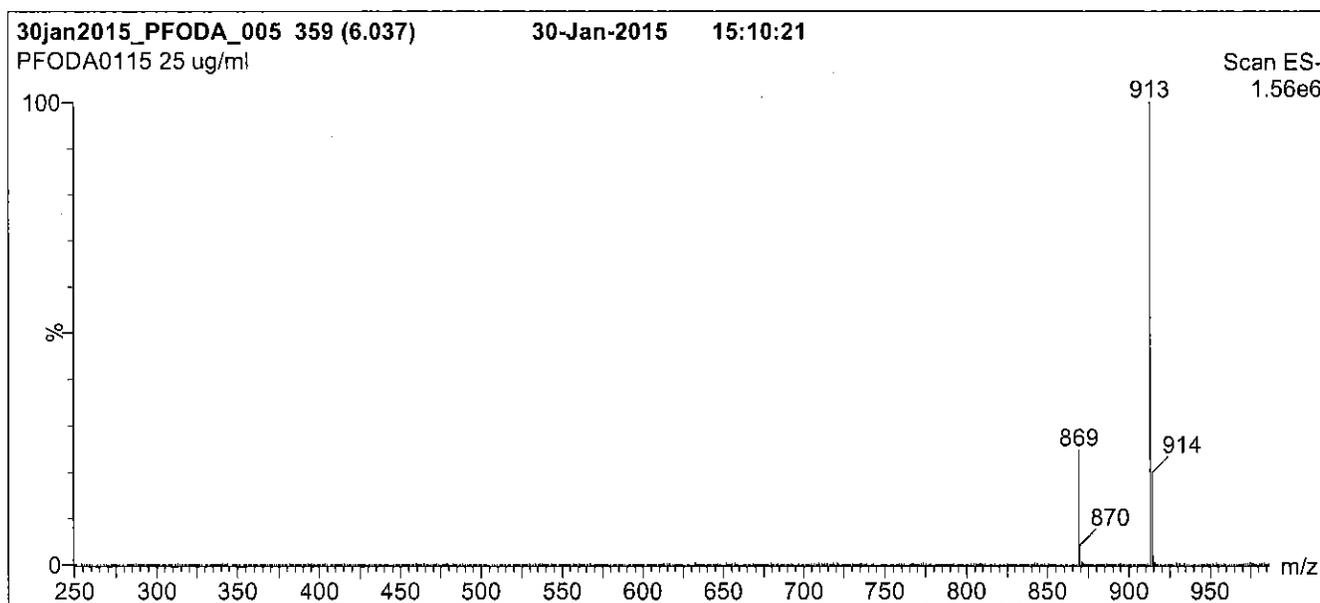
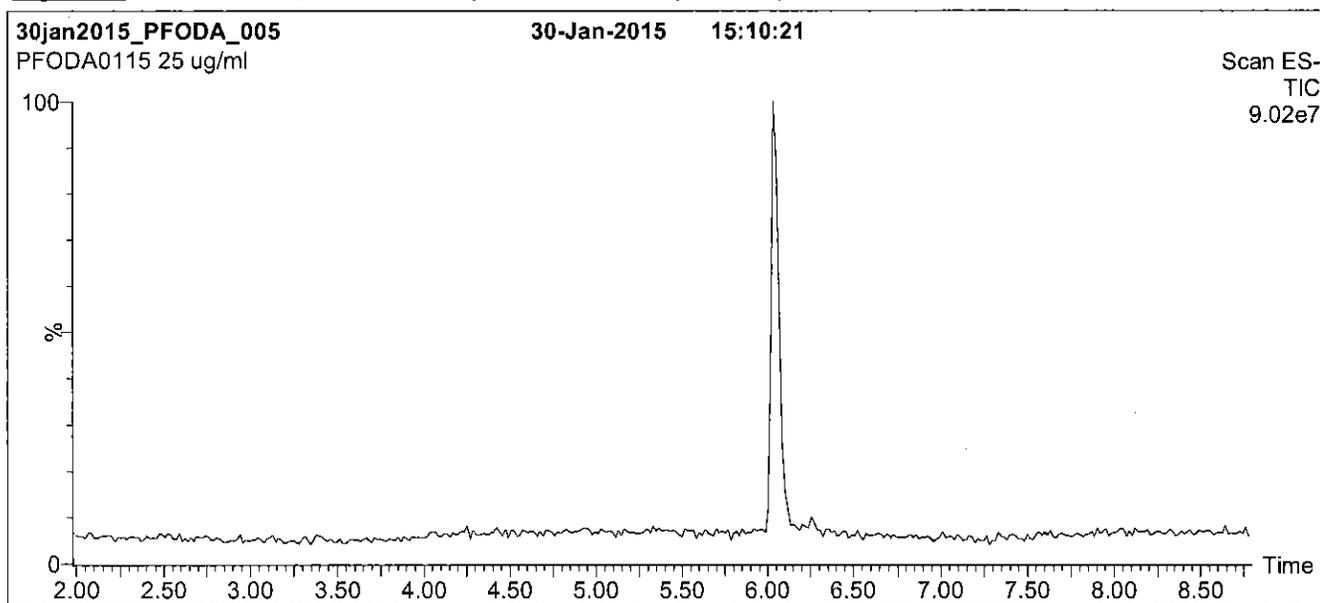
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Figure 1: PFODA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for
 1.5 min before returning to initial conditions in 0.5 min.
 Time: 10 min

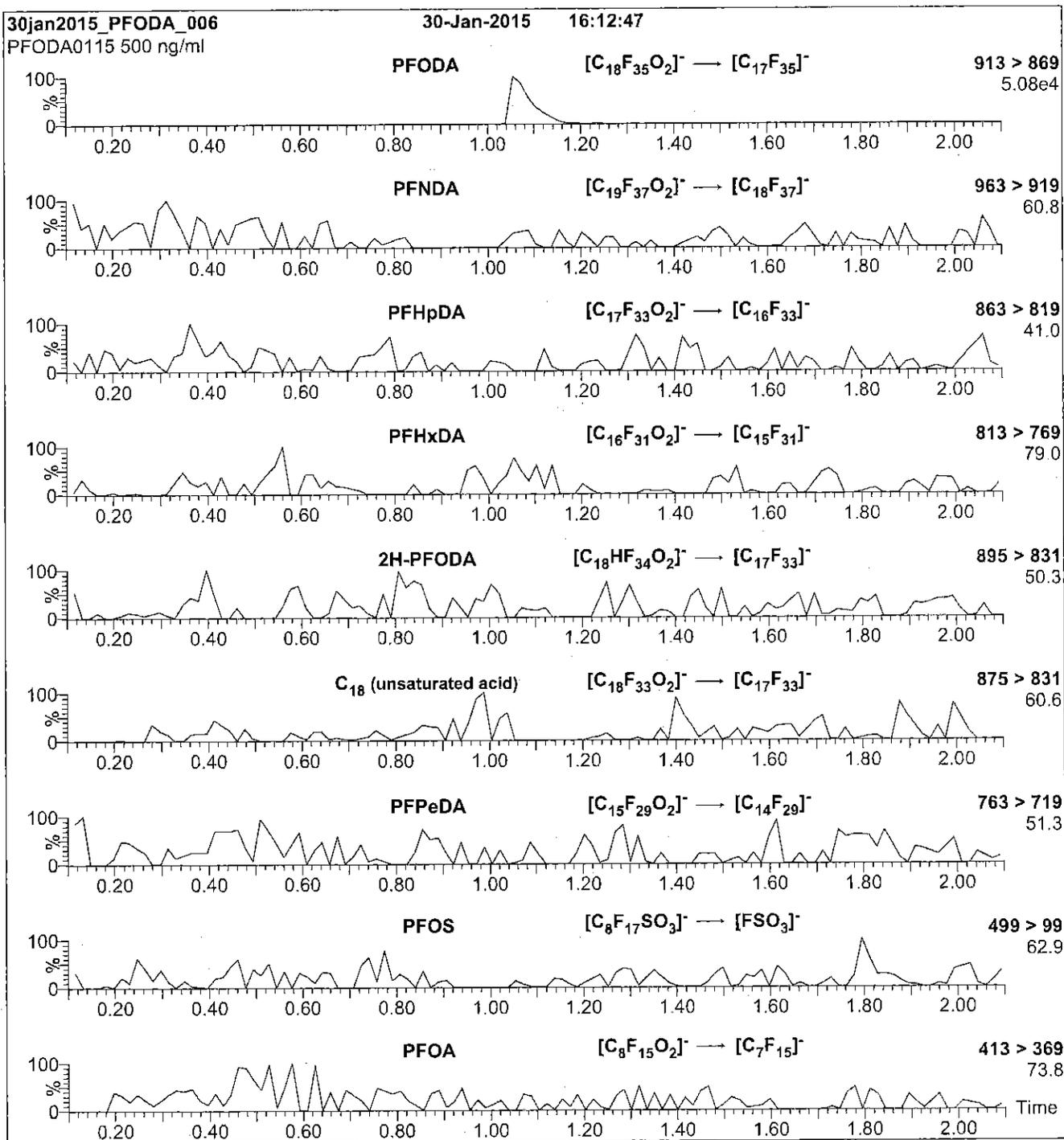
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (250 - 1000 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 25.00
 Cone Gas Flow (l/hr) = 50
 Desolvation Gas Flow (l/hr) = 750

Figure 2: PFODA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 µl (500 ng/ml PFODA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 µl/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
 Collision Energy (eV) = 15

Reagent

LCPFOS-br_00001



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

br-PFOSK

Potassium Perfluorooctanesulfonate Solution/Mixture of Linear and Branched Isomers

PRODUCT CODE: br-PFOSK
LOT NUMBER: brPFOSK1015
CONCENTRATION: 50 ± 2.5 µg/ml (total potassium salt)
46.4 ± 2.3 µg/ml (total PFOS anion)
SOLVENT(S): Methanol
DATE PREPARED: (mm/dd/yyyy) 10/13/2015
LAST TESTED: (mm/dd/yyyy) 10/14/2015
EXPIRY DATE: (mm/dd/yyyy) 10/14/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DESCRIPTION:

The chemical purity has been determined to be ≥98% perfluorooctanesulfonate linear and branched isomers. The full name, structure and percent composition for each of the isomeric components are given in Table A.

DOCUMENTATION/ DATA ATTACHED:

Table A: Isomeric Components and Percent Composition by ¹⁹F-NMR
Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS Data (SIR)
Figure 3: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- A 5-point calibration curve was generated using linear PFOS (potassium salt) and mass-labelled PFOS as an internal standard to enable quantitation of br-PFOSK using isotopic dilution.
- CAS#: 2795-39-3 (for linear isomer; potassium salt).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compounds it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



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Table A: br-PFOSK; Isomeric Components and Percent Composition (by ¹⁹F-NMR)*

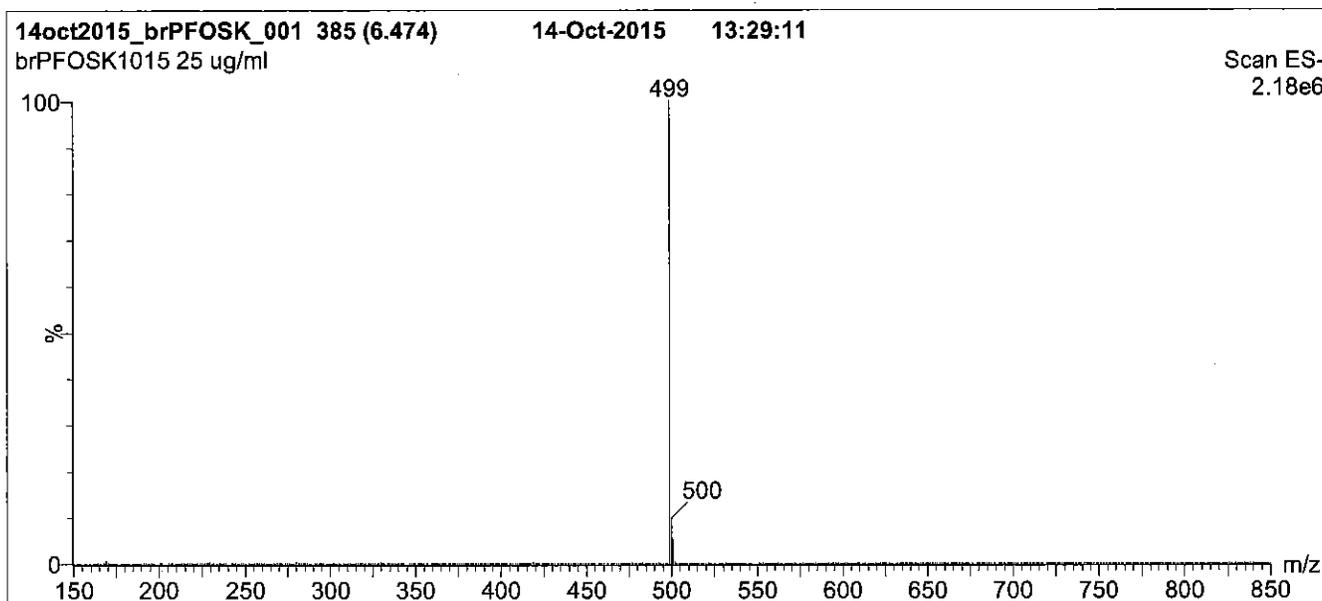
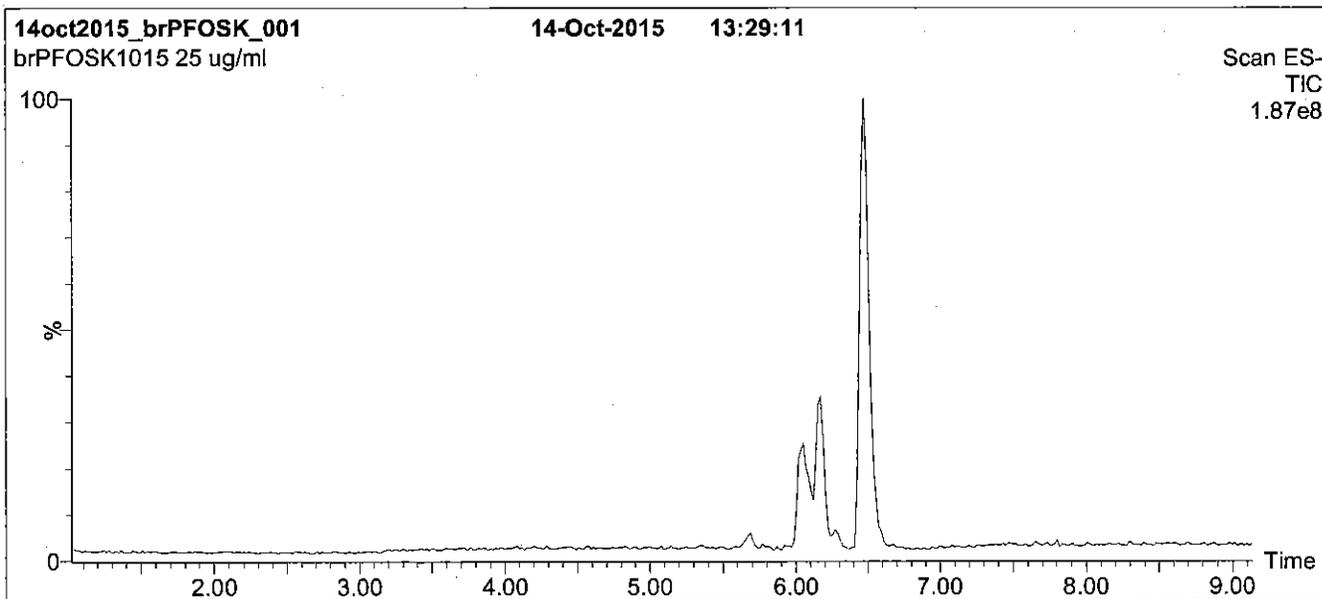
Isomer	Name	Structure	Percent Composition by ¹⁹ F-NMR
1	Potassium perfluoro-1-octanesulfonate	CF ₃ CF ₂ SO ₃ K ⁺	78.8
2	Potassium 1-trifluoromethylperfluoroheptanesulfonate**	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ C(SO ₃)K ⁺ CF ₃	1.2
3	Potassium 2-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF(SO ₃)K ⁺ CF ₃	0.6
4	Potassium 3-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF(SO ₃)K ⁺ CF ₃	1.9
5	Potassium 4-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF(SO ₃)K ⁺ CF ₃	2.2
6	Potassium 5-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF(SO ₃)K ⁺ CF ₃	4.5
7	Potassium 6-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF(SO ₃)K ⁺ CF ₃	10.0
8	Potassium 5,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ -C(CF ₃)CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	0.2
9	Potassium 4,4-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ CF ₂ -C(CF ₃)CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	0.03
10	Potassium 4,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ -CF(CF ₃)-CF(CF ₃)-CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.4
11	Potassium 3,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ -CF(CF ₃)-CF ₂ -CF(CF ₃)-CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.07

* Percent of total perfluorooctanesulfonate isomers only. Isomers are labelled in Figure 2.

** Systematic Name: Potassium perfluorooctane-2-sulfonate.

Certified By:  **Date:** 10/15/2015
B.G. Chittim (mm/dd/yyyy)

Figure 1: br-PFOSK; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 45% (80:20 MeOH:ACN) / 55% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 12 min and hold for 2 min.
Return to initial conditions over 0.5 min.
Time: 16 min

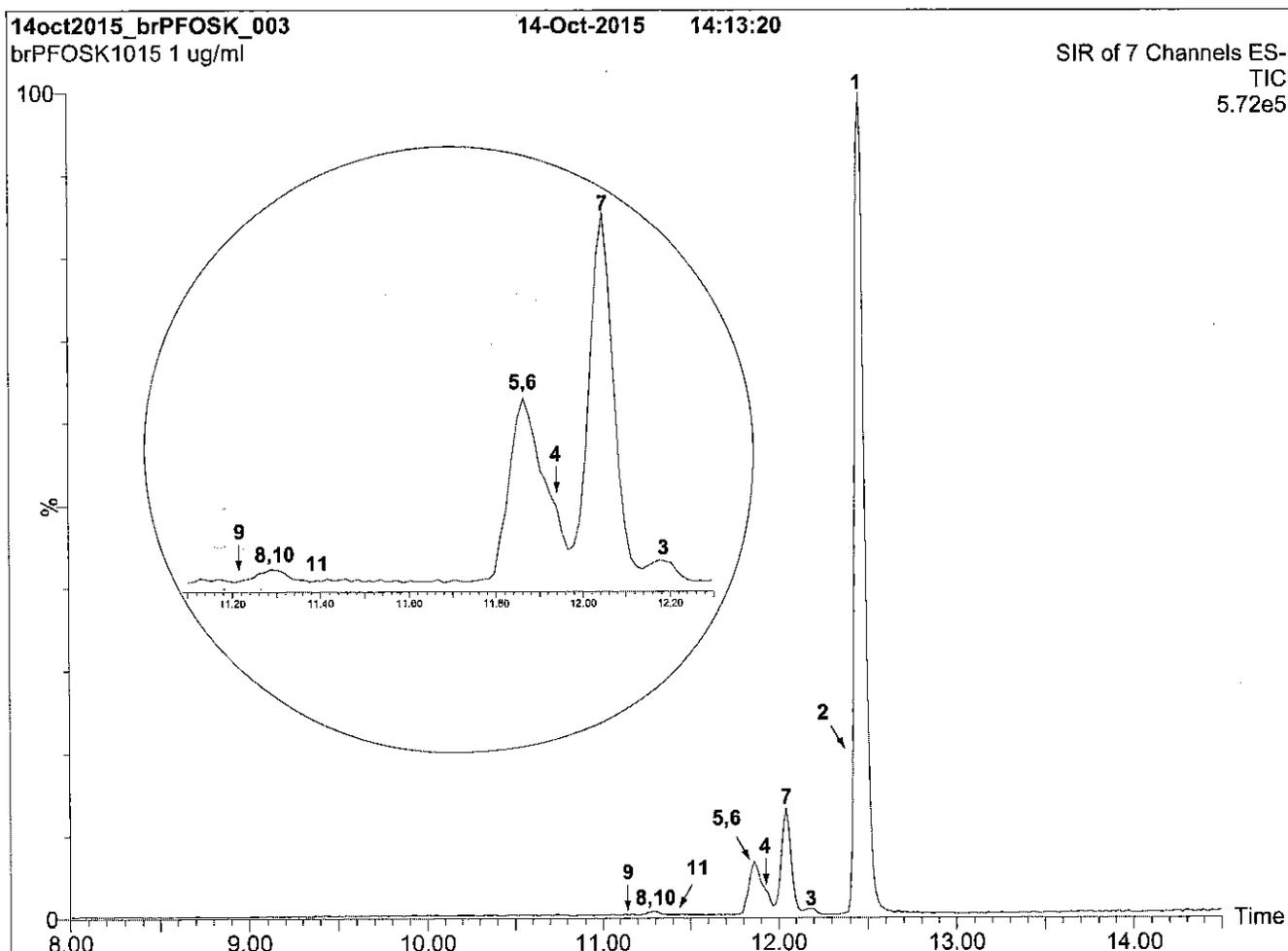
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 60.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: br-PFOSK; LC/MS Data (SIR)



Conditions for Figure 2:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

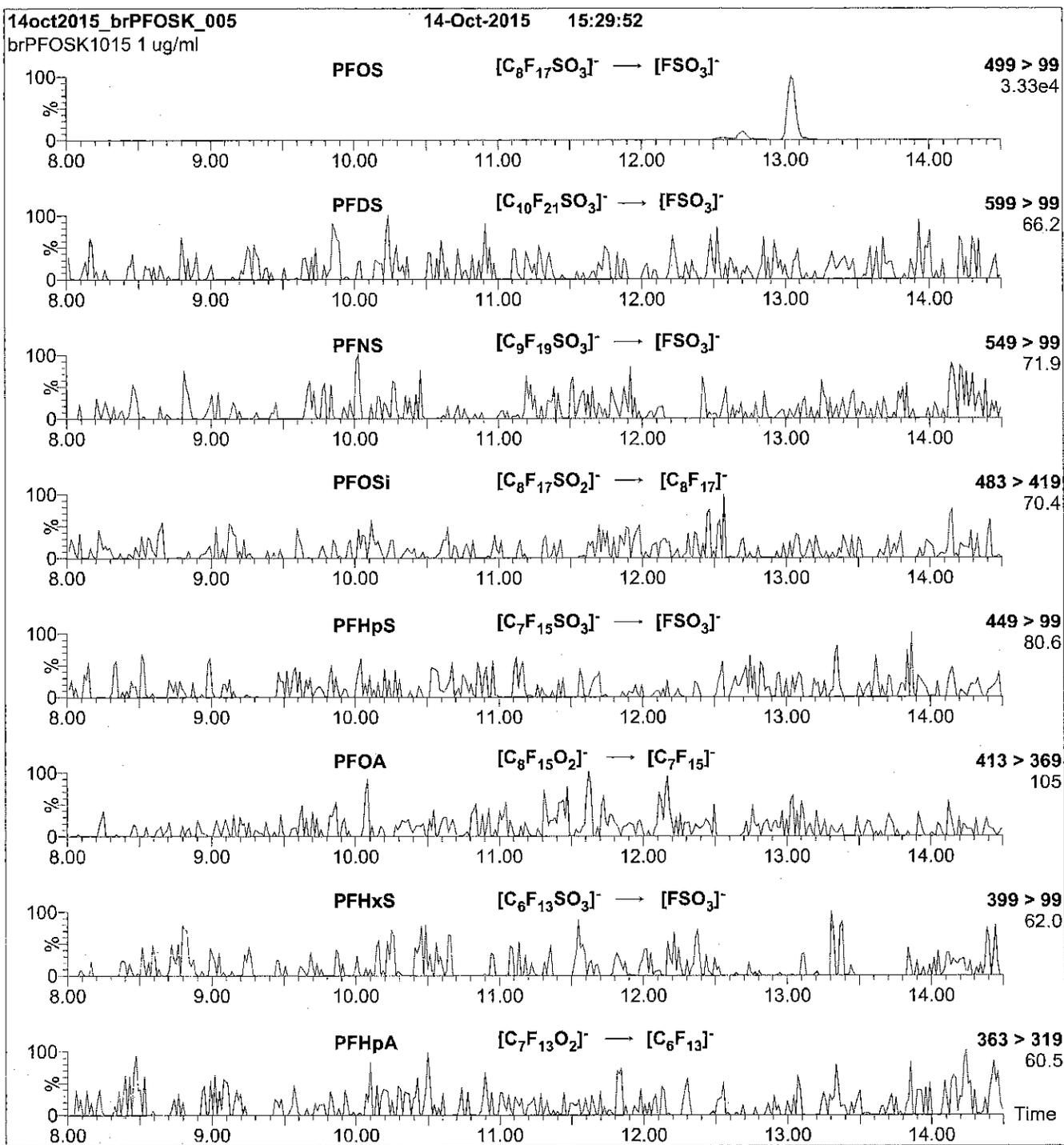
Chromatographic Conditions:

Column: Acquity UPLC BEH Shield RP₁₈ (1.7 μ m, 2.1 x 100 mm)
Injection: 1.0 μ g/ml of br-PFOSK
Mobile Phase: Gradient
45% (80:20 MeOH:ACN) / 55% H₂O (both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 15 min and hold for 3 min.
Return to initial conditions over 1 min.
Time: 20 min
Flow: 300 μ l/min

MS Conditions:

SIR (ES⁻)
Source = 110 °C
Desolvation = 325 °C
Cone Voltage = 60V

Figure 3: br-PFOSK; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 3:

Injection: On-column
 Mobile phase: Same as Figure 2
 Flow: 300 µl/min

MS Parameters

Collision Gas (mbar) = 3.06e-3
 Collision Energy (eV) = 11-50 (variable)

Reagent

LCPFOS-br_00002

Scanned
10/14/16 SR

R: SBC 9/13/16



730515
ID: LCPFOS-br_00002
Exp: 10/14/20 Prpt: SBC
Potassium Perfluorooctane



730516
ID: LCPFOS-br_00003
Exp: 10/14/20 Prpt: SBC
Potassium Perfluorooctane



WELLINGTON
LABORATORIES

CERTIFICATE OF ANALYSIS
DOCUMENTATION

br-PFOSK

Potassium Perfluorooctanesulfonate
Solution/Mixture of Linear and
Branched Isomers

PRODUCT CODE: br-PFOSK
LOT NUMBER: brPFOSK1015
CONCENTRATION: 50 ± 2.5 µg/ml (total potassium salt)
46.4 ± 2.3 µg/ml (total PFOS anion)
SOLVENT(S): Methanol
DATE PREPARED: (mm/dd/yyyy) 10/13/2015
LAST TESTED: (mm/dd/yyyy) 10/14/2015
EXPIRY DATE: (mm/dd/yyyy) 10/14/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DESCRIPTION:

The chemical purity has been determined to be ≥98% perfluorooctanesulfonate linear and branched isomers. The full name, structure and percent composition for each of the isomeric components are given in Table A.

DOCUMENTATION/ DATA ATTACHED:

Table A: Isomeric Components and Percent Composition by ¹⁹F-NMR
Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS Data (SIR)
Figure 3: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- A 5-point calibration curve was generated using linear PFOS (potassium salt) and mass-labelled PFOS as an internal standard to enable quantitation of br-PFOSK using isotopic dilution.
- CAS#: 2795-39-3 (for linear isomer; potassium salt).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

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HAZARDS:

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HOMOGENEITY:

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UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

QUALITY MANAGEMENT:

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Table A: br-PFOSK; Isomeric Components and Percent Composition (by ¹⁹F-NMR)*

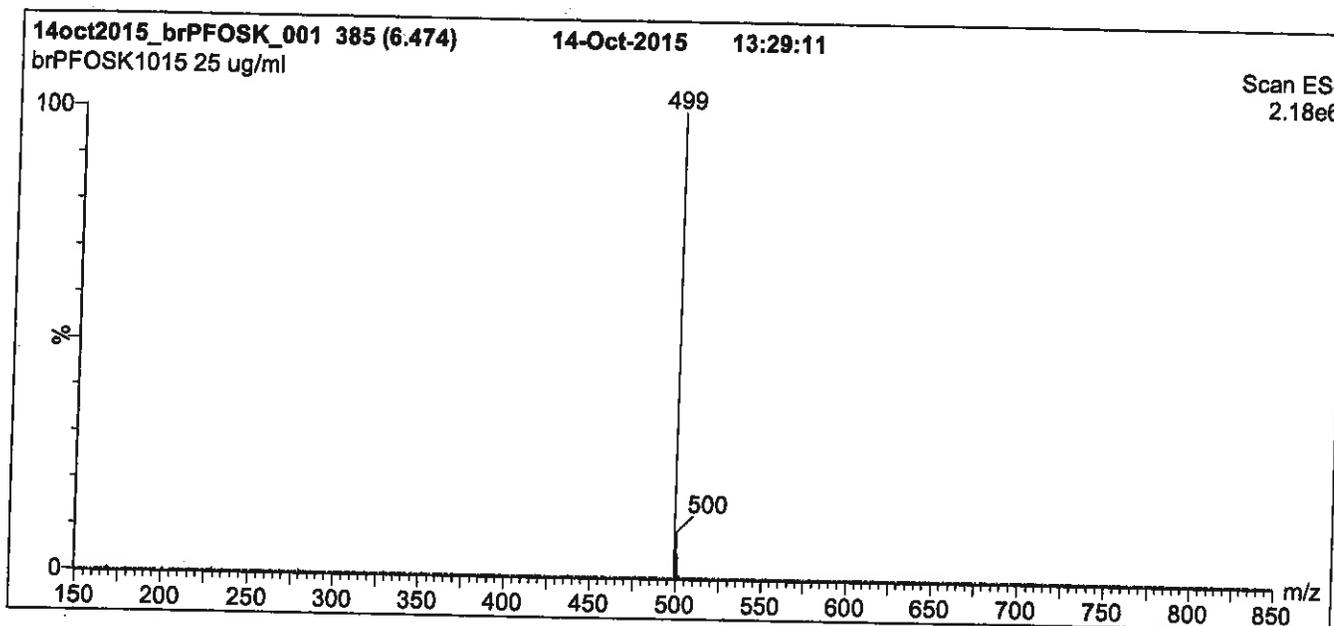
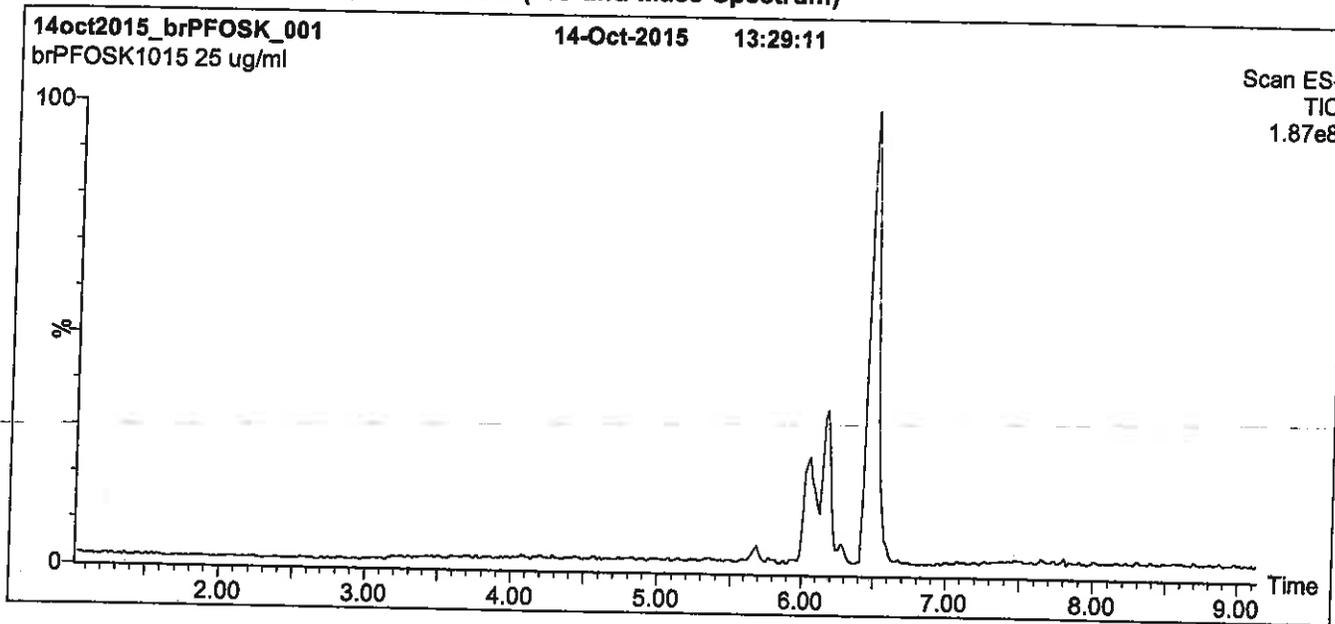
Isomer	Name	Structure	Percent Composition by ¹⁹ F-NMR
1	Potassium perfluoro-1-octanesulfonate	CF ₃ CF ₂ SO ₃ K ⁺	78.8
2	Potassium 1-trifluoromethylperfluoroheptanesulfonate**	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	1.2
3	Potassium 2-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	0.6
4	Potassium 3-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	1.9
5	Potassium 4-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	2.2
6	Potassium 5-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	4.5
7	Potassium 6-trifluoromethylperfluoroheptanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃	10.0
8	Potassium 5,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.2
9	Potassium 4,4-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.03
10	Potassium 4,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.4
11	Potassium 3,5-di(trifluoromethyl)perfluorohexanesulfonate	CF ₃ CF ₂ CF ₂ CF ₂ CF ₂ CF ₂ SO ₃ K ⁺ CF ₃ CF ₃	0.07

* Percent of total perfluorooctanesulfonate isomers only. Isomers are labeled in Figure 2.
 ** Systematic Name: Potassium perfluorooctane-2-sulfonate.

Certified By: 
 B.G. Chittim

Date: 10/15/2015
(mm/dd/yyyy)

Figure 1: br-PFOSK; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈,
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 45% (80:20 MeOH:ACN) / 55% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 12 min and hold for 2 min.
Return to initial conditions over 0.5 min.
Time: 16 min

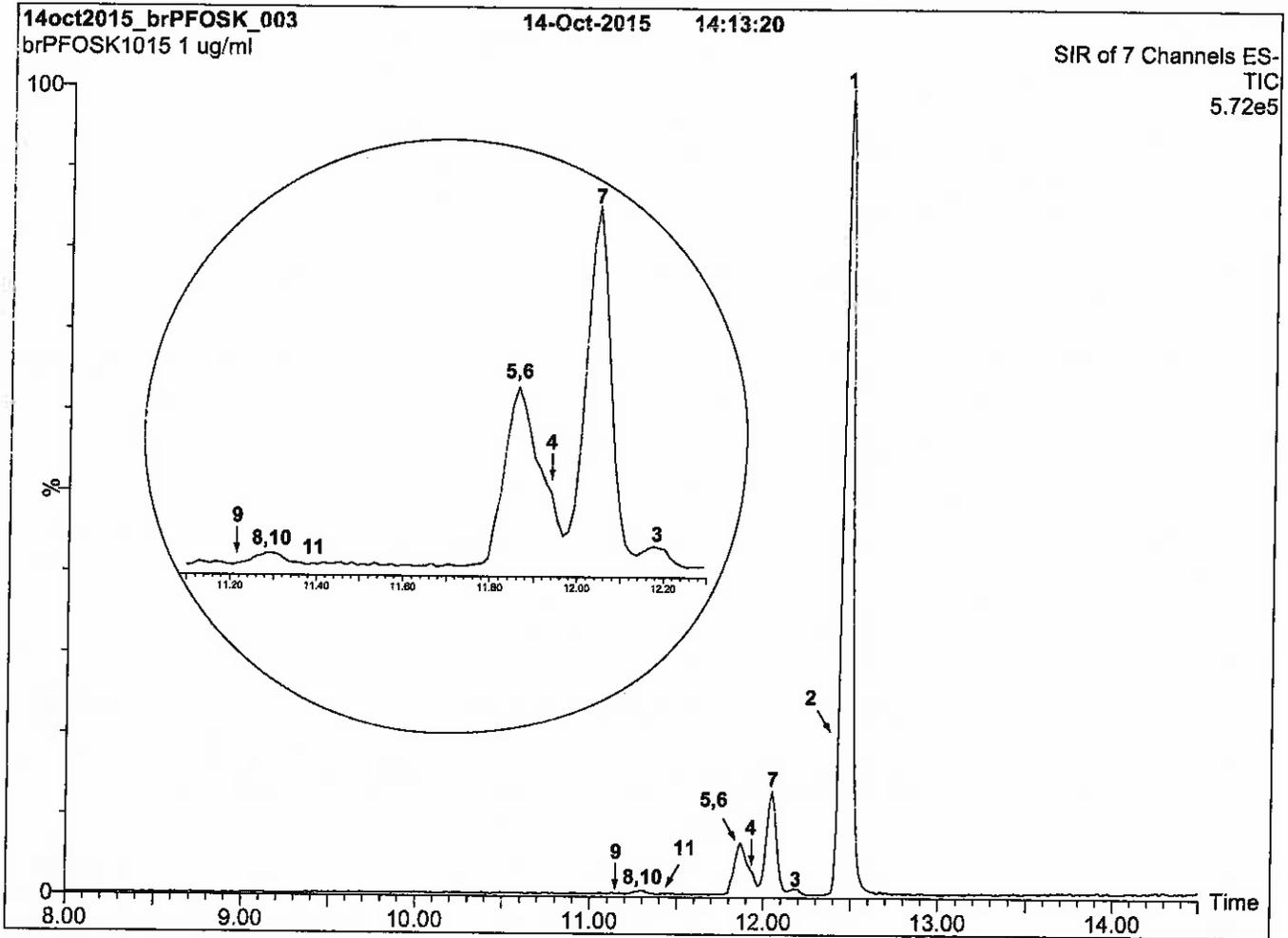
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.00
Cone Voltage (V) = 60.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: br-PFOSK; LC/MS Data (SIR)



Conditions for Figure 2:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

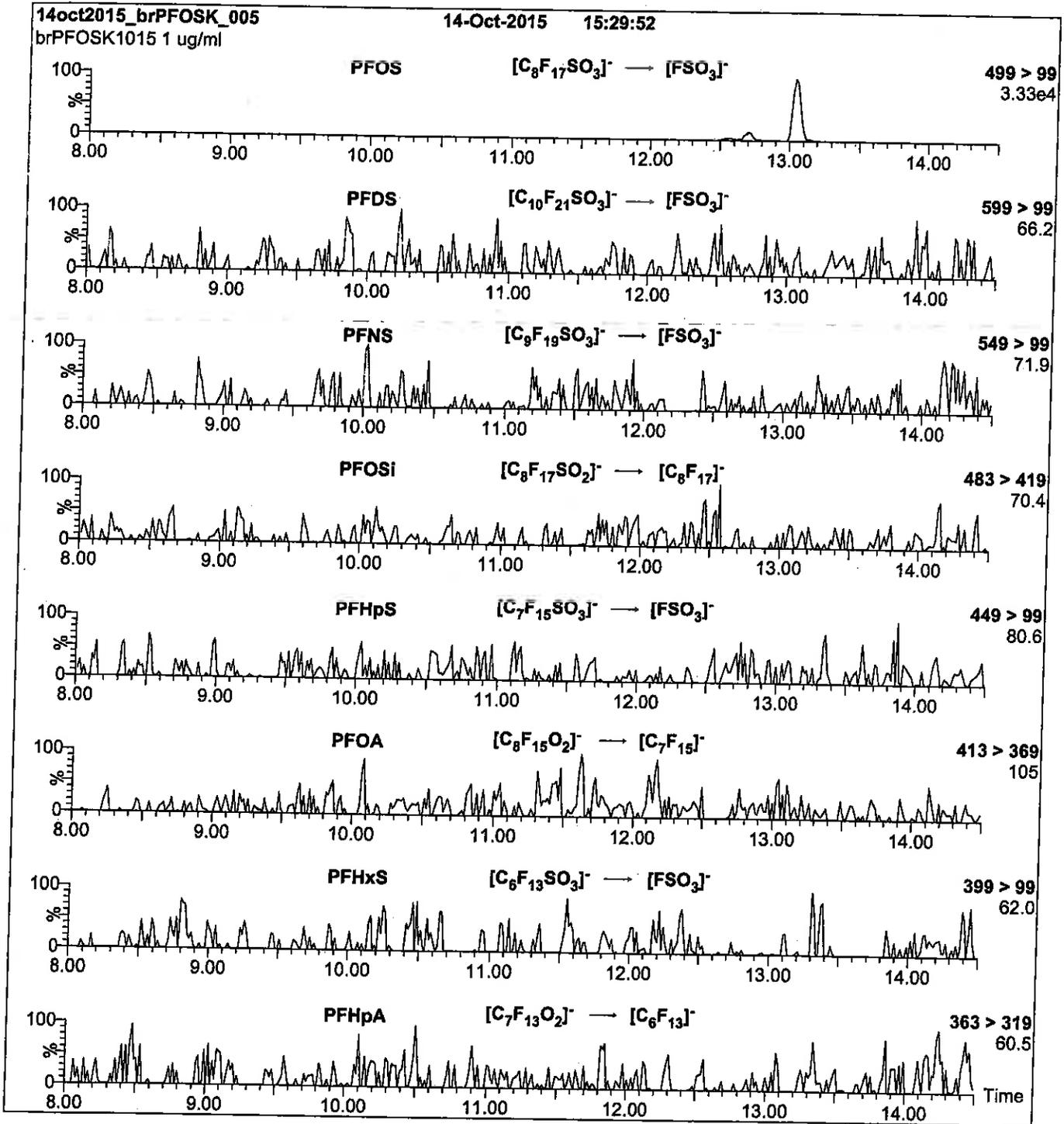
Chromatographic Conditions:

Column: Acquity UPLC BEH Shield RP₁₈ (1.7 μ m, 2.1 x 100 mm)
Injection: 1.0 μ g/ml of br-PFOSK
Mobile Phase: Gradient
45% (80:20 MeOH:ACN) / 55% H₂O (both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 15 min and hold for 3 min.
Return to initial conditions over 1 min.
Time: 20 min
Flow: 300 μ l/min

MS Conditions:

SIR (ES)
Source = 110 $^{\circ}$ C
Desolvation = 325 $^{\circ}$ C
Cone Voltage = 60V

Figure 3: br-PFOSK; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 3:

Injection: On-column

Mobile phase: Same as Figure 2

Flow: 300 μ /min

MS Parameters

Collision Gas (mbar) = 3.06e-3

Collision Energy (eV) = 11-50 (variable)

Reagent

LCPFOSA_00006

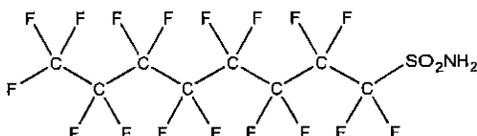


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: FOSA-I **LOT NUMBER:** FOSA0815I
COMPOUND: Perfluoro-1-octanesulfonamide

STRUCTURE: **CAS #:** 754-91-6



MOLECULAR FORMULA: $C_8H_2F_{17}NO_2S$ **MOLECULAR WEIGHT:** 499.14
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Isopropanol
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 09/02/2015
EXPIRY DATE: (mm/dd/yyyy) 09/02/2017
RECOMMENDED STORAGE: Refrigerate ampoule

DOCUMENTATION/ DATA ATTACHED:

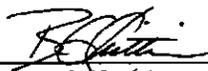
Figure 1: LC/MS Data (TIC and Mass Spectrum)
 Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: _____


 B.G. Chittim

Date: 09/11/2015
 (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
 519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

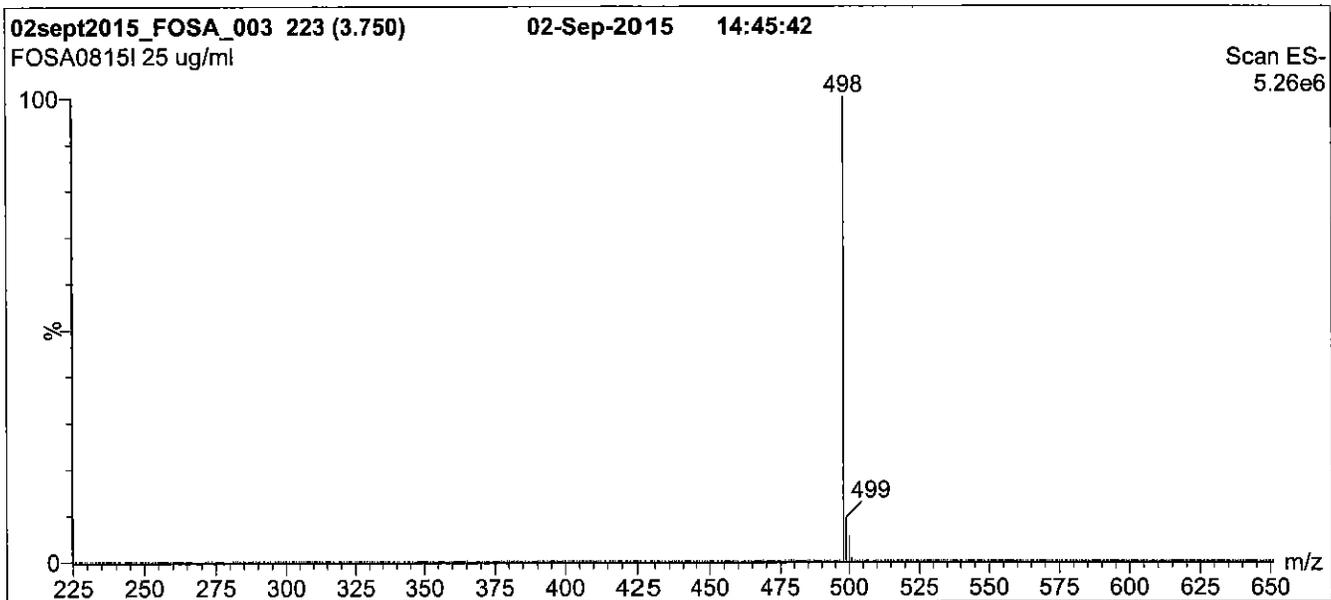
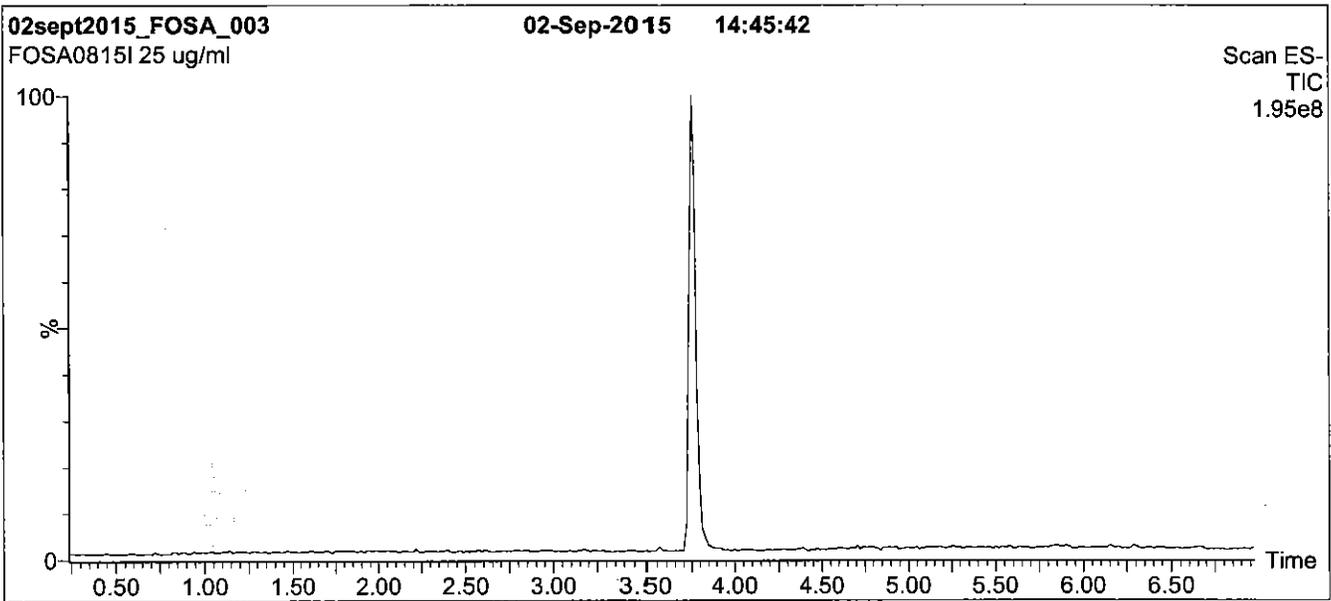
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: FOSA-I; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP_{1a}
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 60% (80:20 MeOH:ACN) / 40% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

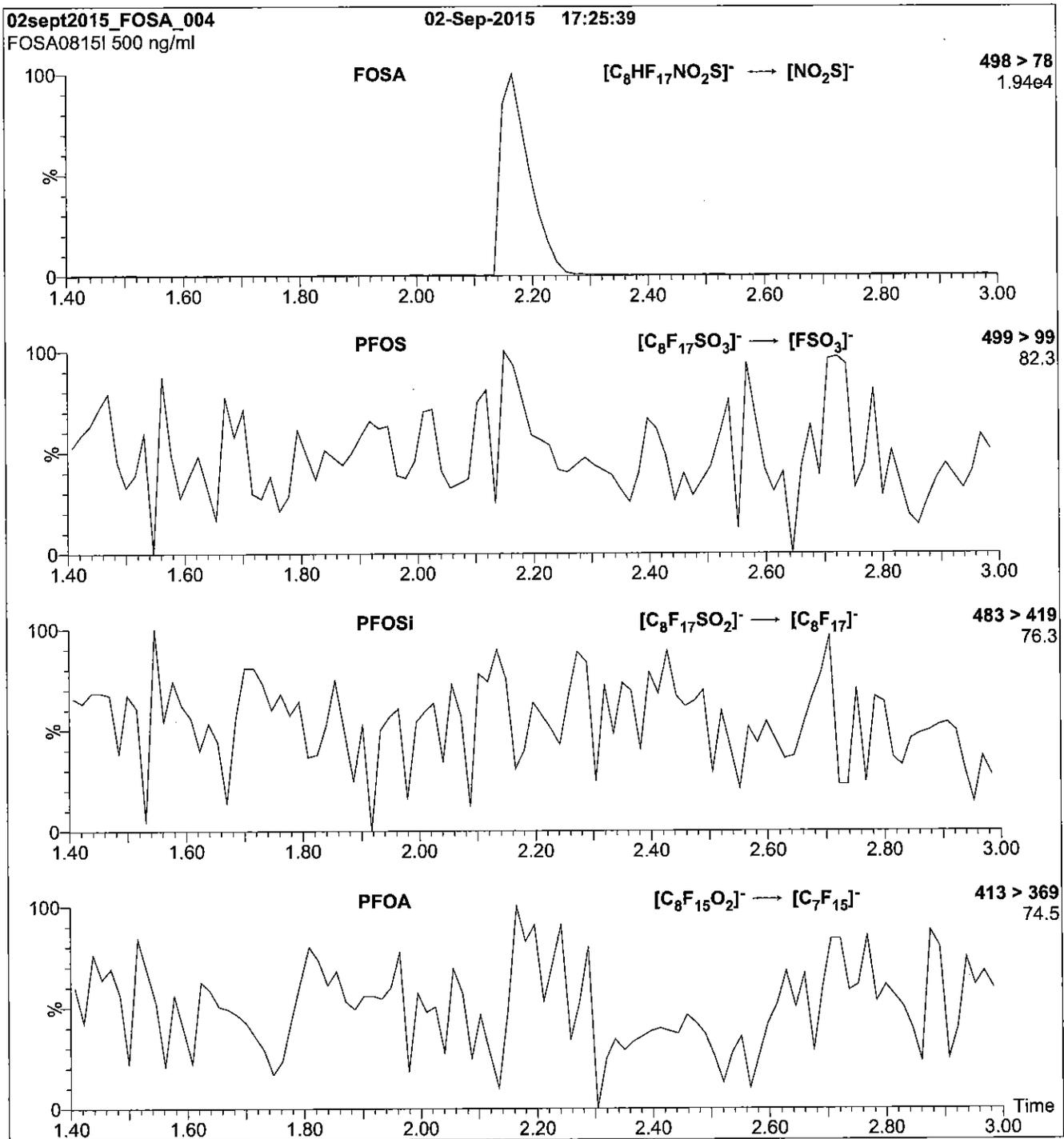
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 2.50
Cone Voltage (V) = 40.00
Cone Gas Flow (l/hr) = 50
Desolvation Gas Flow (l/hr) = 750

Figure 2: FOSA-I; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml FOSA-I)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.54e-3
Collision Energy (eV) = 30

Reagent

LCFPeA_00005

R: 7/6/16 CBW



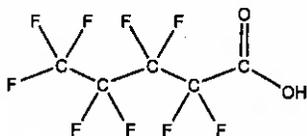
671579
ID: LCPFPeA_00005
Exp: 01/30/20 Prod: CBW
PF-n-pentanoic acid



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFPeA **LOT NUMBER:** PFPeA0115
COMPOUND: Perfluoro-n-pentanoic acid
STRUCTURE: **CAS #:** 2706-90-3



MOLECULAR FORMULA: $C_5HF_9O_2$ **MOLECULAR WEIGHT:** 264.05
CONCENTRATION: $50 \pm 2.5 \mu\text{g/ml}$ **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 01/30/2015
EXPIRY DATE: (mm/dd/yyyy) 01/30/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

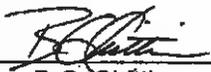
DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.3% of Perfluoro-n-heptanoic acid (PFHpA) and ~ 0.2% of $C_5H_2F_8O_2$ (hydrido - derivative) as measured by ^{19}F NMR.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By: 
B.G. Chittim **Date:** 03/26/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

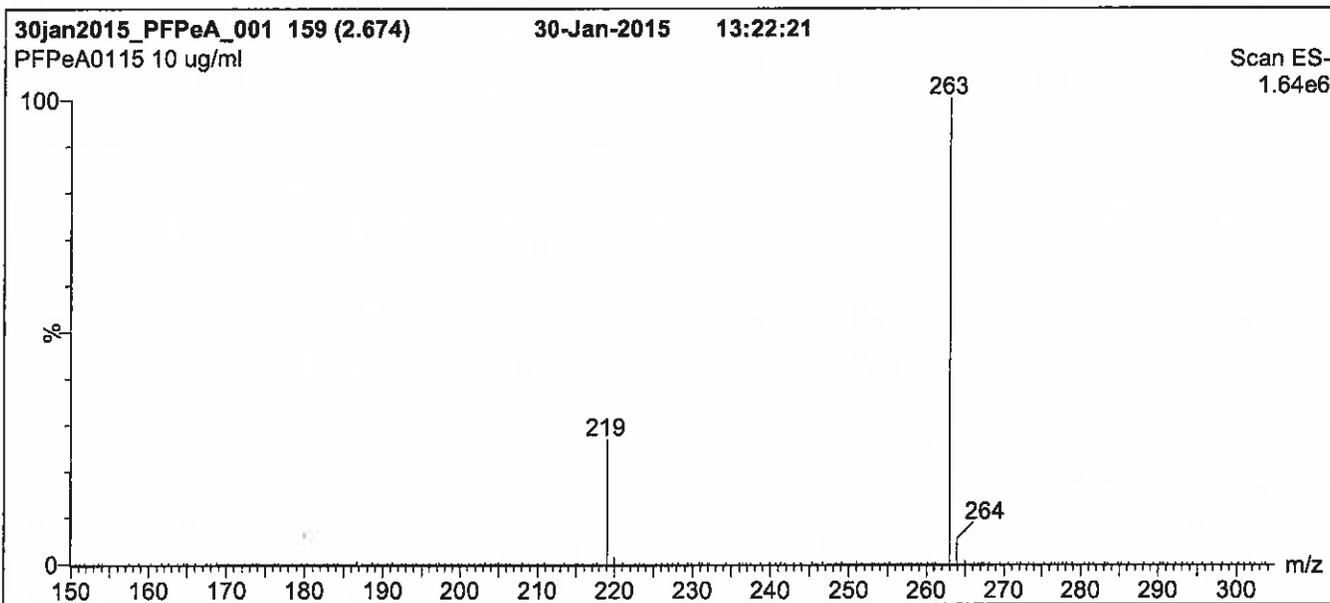
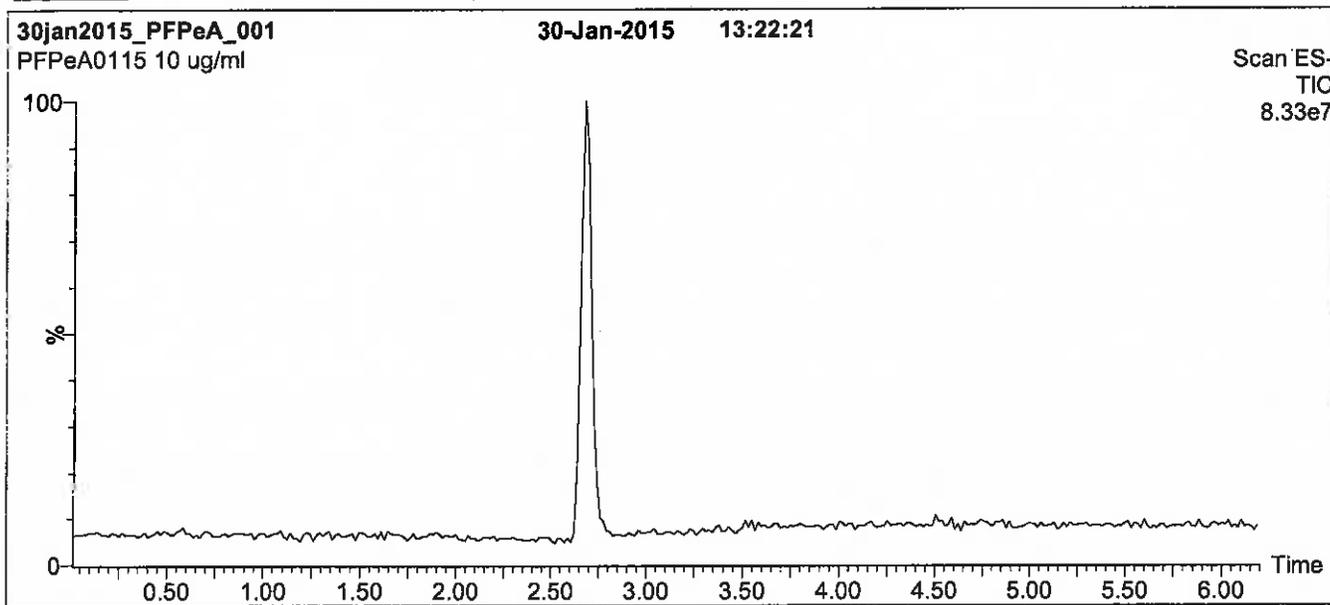
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: PFPeA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 30% (80:20 MeOH:ACN) / 70% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7.5 min and hold for 1 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

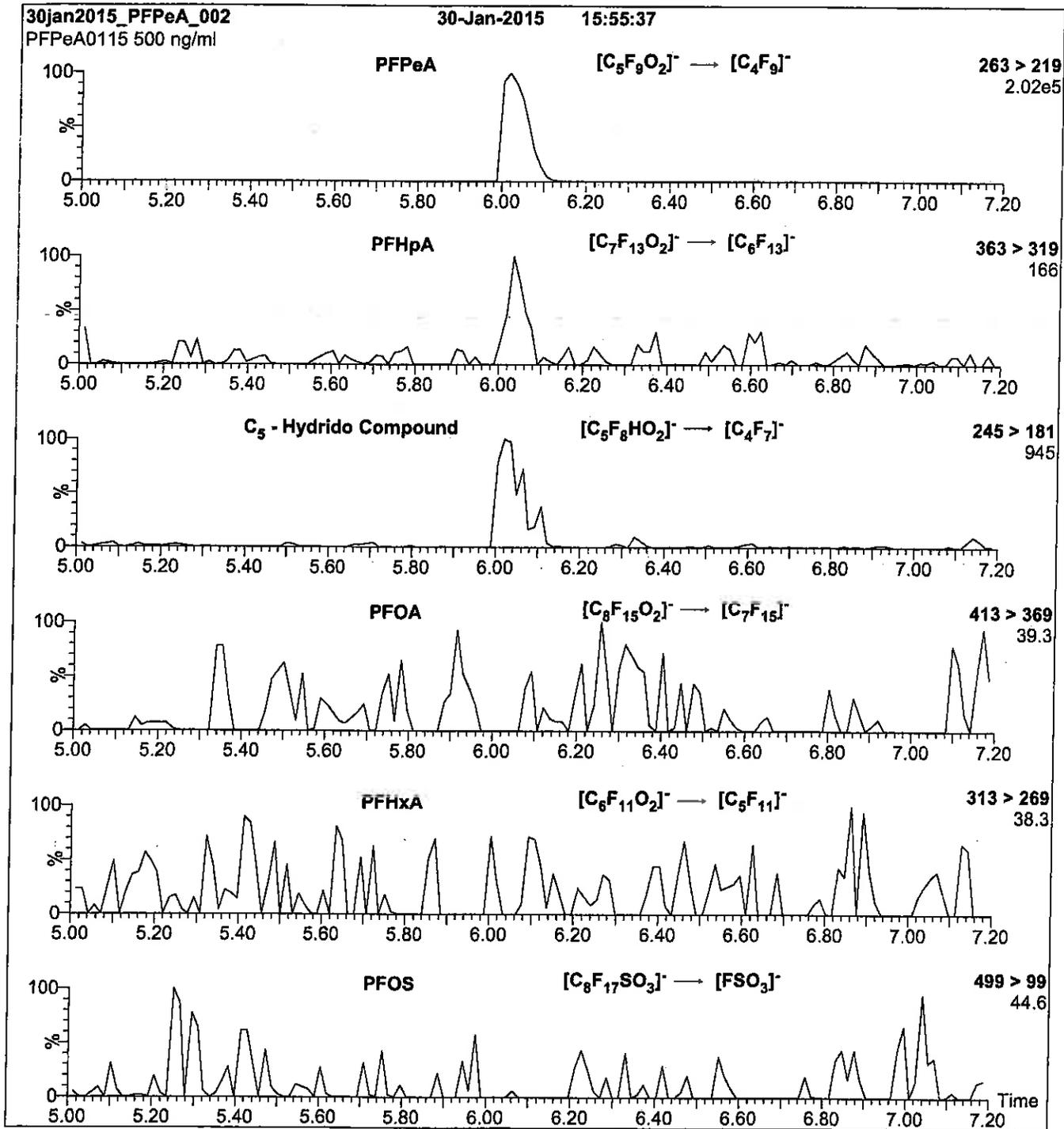
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 60
 Desolvation Gas Flow (l/hr) = 750

Figure 2: PFPeA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFPeA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.35e-3
 Collision Energy (eV) = 9

Reagent

LCPFTeDA_00004



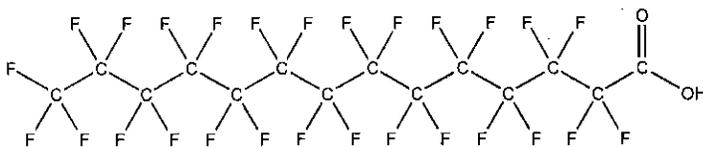
R: 4/7/16 CBW

609636

ID: LCPFTeDA_00004

Exp: 12/09/20 Pripd: CBW

PF-n-tetradecanoic acid

**WELLINGTON**
LABORATORIES**CERTIFICATE OF ANALYSIS**
DOCUMENTATION**PRODUCT CODE:** PFTeDA **LOT NUMBER:** PFTeDA1215
COMPOUND: Perfluoro-n-tetradecanoic acid**STRUCTURE:** **CAS #:** 376-06-7

MOLECULAR FORMULA:	$C_{14}H_{27}O_2$	MOLECULAR WEIGHT:	714.11
CONCENTRATION:	$50 \pm 2.5 \mu\text{g/ml}$	SOLVENT(S):	Methanol Water (<1%)
CHEMICAL PURITY:	>98%		
LAST TESTED: (mm/dd/yyyy)	12/09/2015		
EXPIRY DATE: (mm/dd/yyyy)	12/09/2020		
RECOMMENDED STORAGE:	Store ampoule in a cool, dark place		

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.2% of PFDa ($C_{12}H_{23}O_2$) and ~ 0.2% of PFPeDA ($C_{15}H_{29}O_2$).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim
Date: 12/09/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

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where x is expressed as a relative standard uncertainty of the individual parameter.

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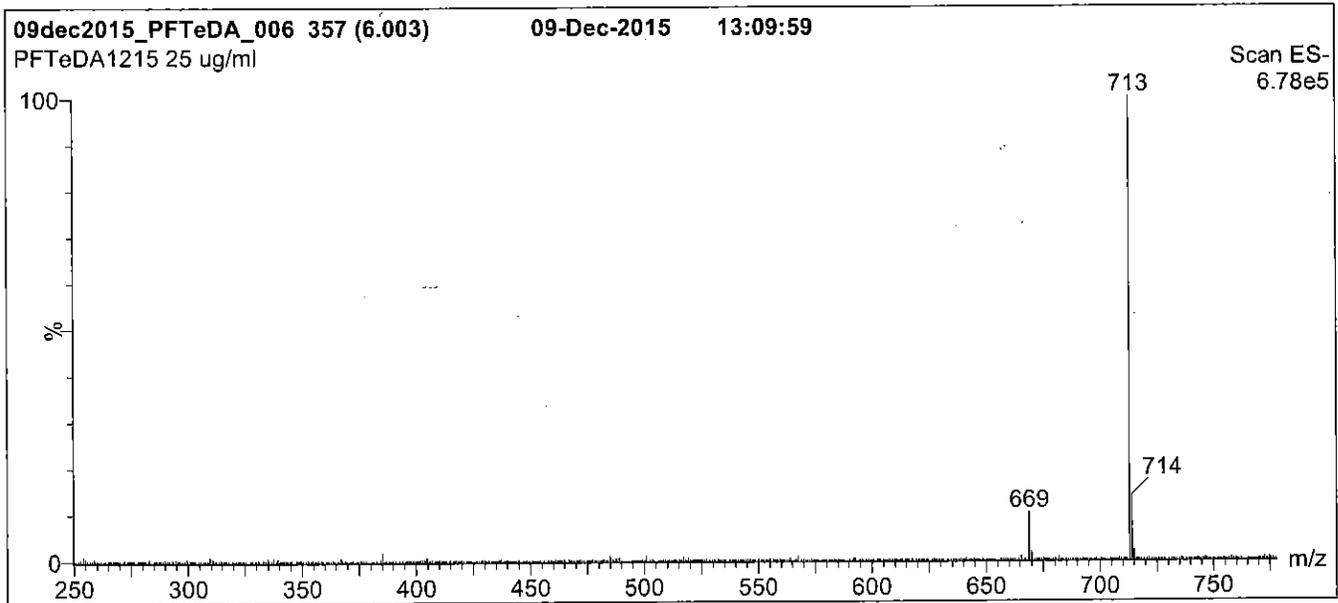
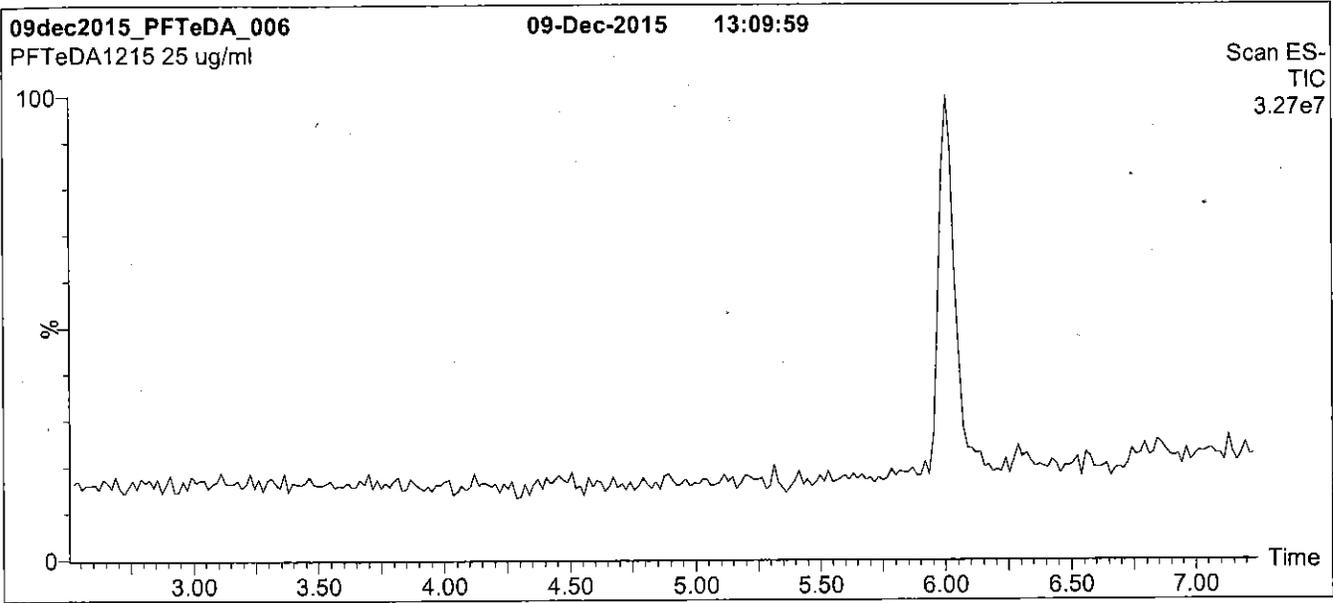
QUALITY MANAGEMENT:

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Figure 1: PFTeDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro micro API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 µm, 2.1 x 100 mm

Mobile phase: Gradient
Start: 65% (80:20 MeOH:ACN) / 35% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7.5 min and hold for 1.5 min
before returning to initial conditions in 0.5 min.
Time: 10 min

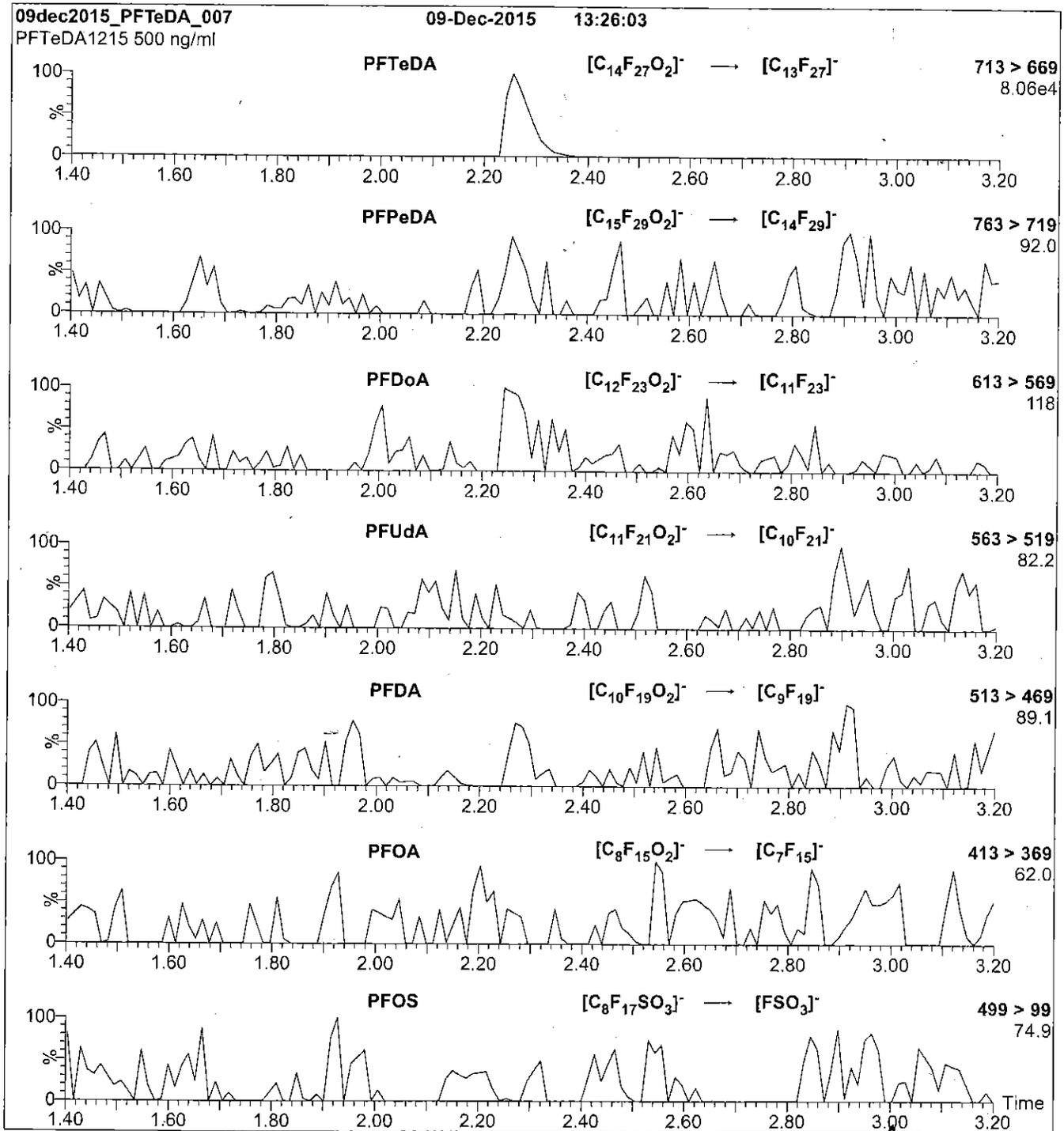
Flow: 300 µl/min

MS Parameters

Experiment: Full Scan (250 - 1250 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 60
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFTeDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFTeDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.43e-3
Collision Energy (eV) = 14

Reagent

LCPFT_rDA_00004



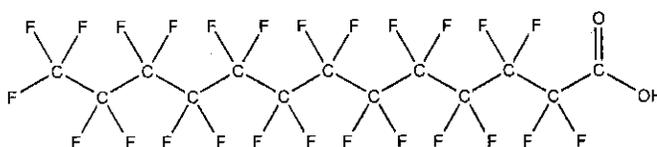
R: 4/7/16 CBW

609697

ID: LCPFTrDA_00004

Exp: 12/10/18 Prep: CBW

PF-n-tridecanoic acid

**WELLINGTON**
LABORATORIES**CERTIFICATE OF ANALYSIS**
DOCUMENTATION**PRODUCT CODE:** PFTTrDA **LOT NUMBER:** PFTTrDA1213
COMPOUND: Perfluoro-n-tridecanoic acid**STRUCTURE:** **CAS #:** 72629-94-8

MOLECULAR FORMULA:	$C_{13}H_1F_{25}O_2$	MOLECULAR WEIGHT:	664.11
CONCENTRATION:	$50 \pm 2.5 \mu\text{g/ml}$	SOLVENT(S):	Methanol Water (<1%)
CHEMICAL PURITY:	>98%		
LAST TESTED: (mm/dd/yyyy)	12/10/2013		
EXPIRY DATE: (mm/dd/yyyy)	12/10/2018		
RECOMMENDED STORAGE:	Store ampoule in a cool, dark place		

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains ~ 0.1% of PFUdA ($C_{11}H_9F_{21}O_2$); ~ 0.4% of PFDaA ($C_{12}H_7F_{23}O_2$), and ~ 0.1% of PFTeDA ($C_{14}H_5F_{27}O_2$).

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim
Date: 03/25/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON 'N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the stated technical and purity specifications.

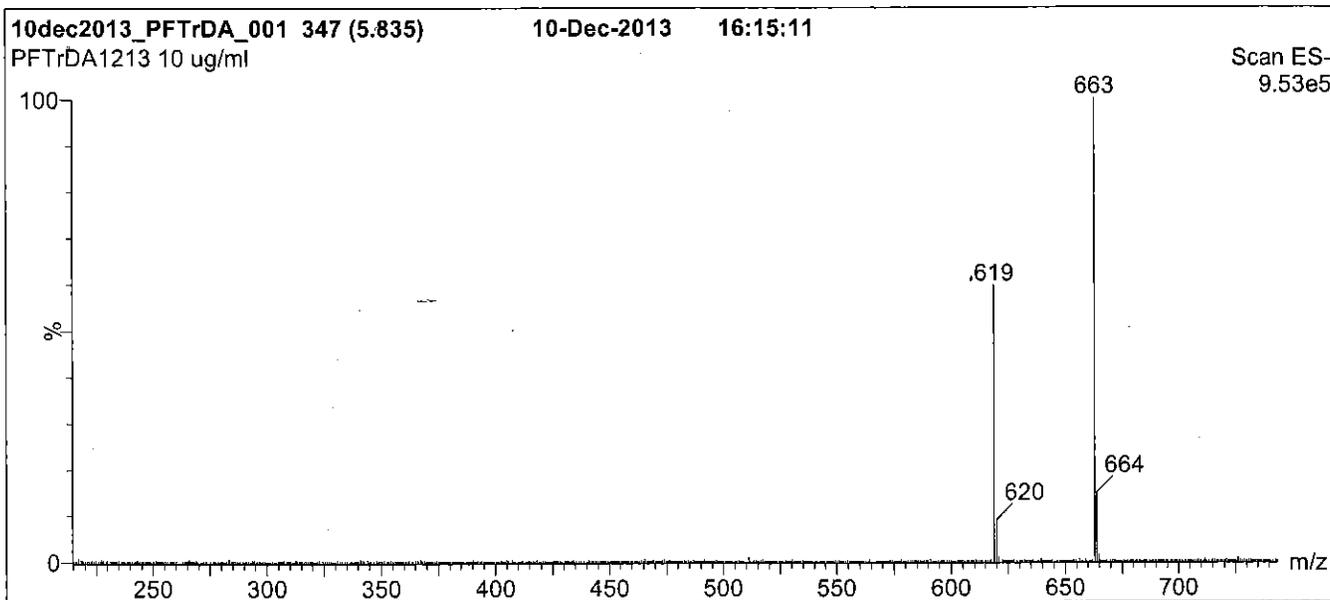
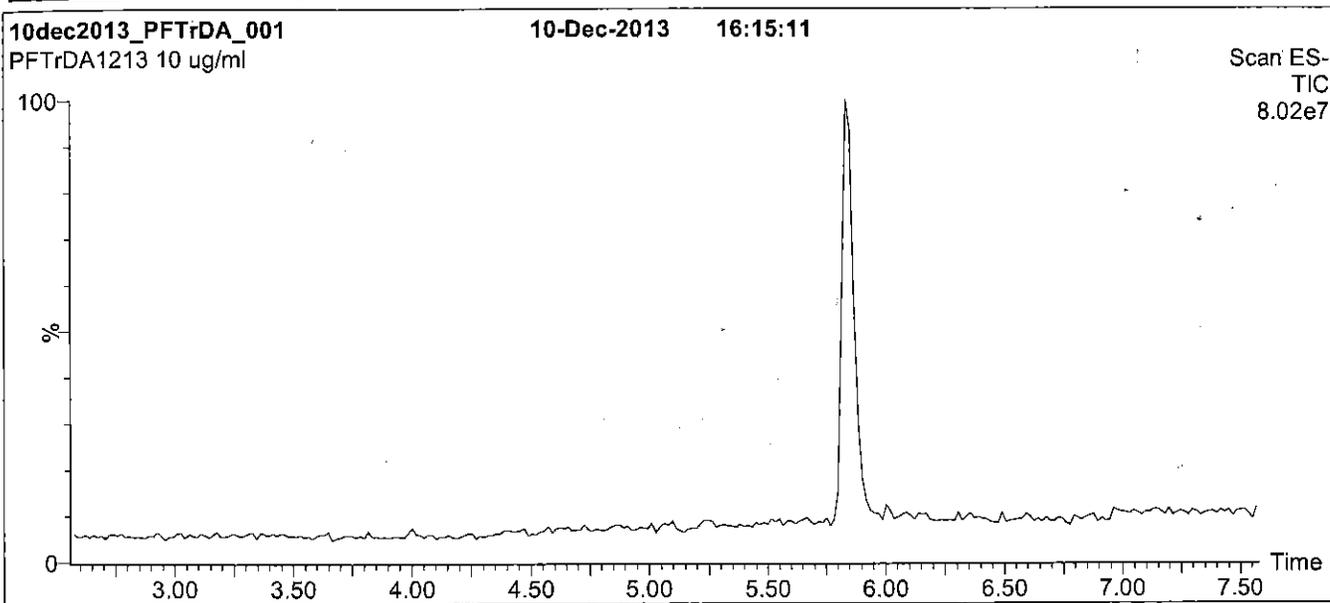
QUALITY MANAGEMENT:

This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at www.well-labs.com or contact us directly at info@well-labs.com

Figure 1: PFTrDA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 60% (80:20 MeOH:ACN) / 40% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 1.5 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

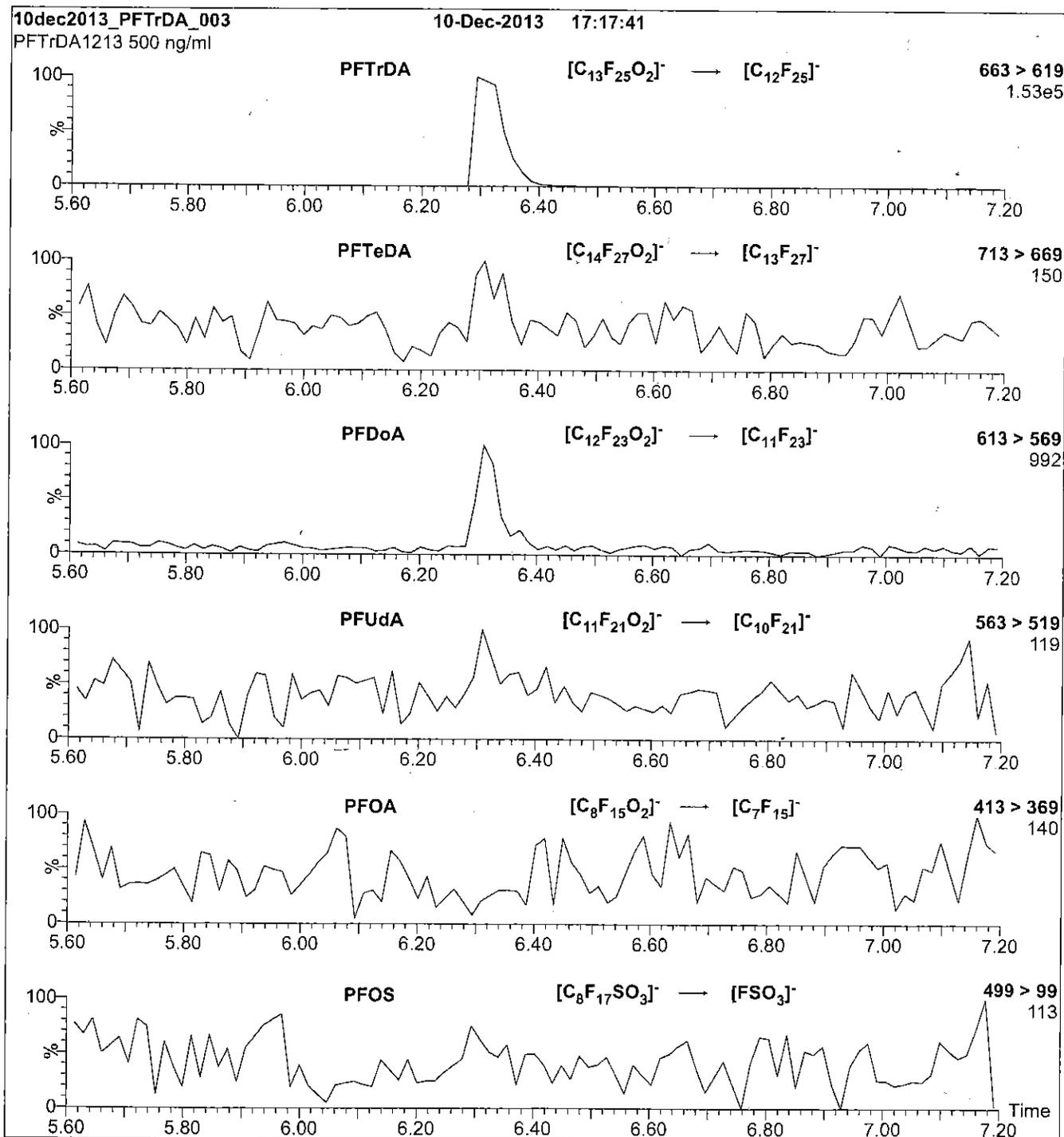
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (215 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 2.00
 Cone Voltage (V) = 22.00
 Cone Gas Flow (l/hr) = 60
 Desolvation Gas Flow (l/hr) = 650

Figure 2: PFTrDA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
10 μ l (500 ng/ml PFTrDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
(both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.28e-3
Collision Energy (eV) = 15

Reagent

LCPFUdA_00004

605242
ID: LCPFUdA_00004
Exp: 08/19/20 Pppe: CBW
PF-n-undecanoic acid

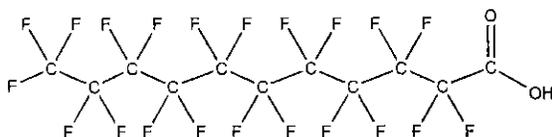
Rec. 3/29/16 JRB ✓



WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFUdA **LOT NUMBER:** PFUdA0815
COMPOUND: Perfluoro-n-undecanoic acid
STRUCTURE: **CAS #:** 2058-94-8



MOLECULAR FORMULA: C₁₁H₂₁O₂ **MOLECULAR WEIGHT:** 564.09
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 08/19/2015
EXPIRY DATE: (mm/dd/yyyy) 08/19/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

Figure 1: LC/MS Data (TIC and Mass Spectrum)
Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:

B.G. Chittim

Date: 08/21/2015
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

INTENDED USE:

The products prepared by Wellington Laboratories Inc. are for laboratory use only. This certified reference material (CRM) was designed to be used as a standard for the identification and/or quantification of the specific chemical compound it contains.

HAZARDS:

This product should only be used by qualified personnel familiar with its potential hazards and trained in the handling of hazardous chemicals. Due care should be exercised to prevent unnecessary human contact or ingestion. All procedures should be carried out in a well-functioning fume hood and suitable gloves, eye protection, and clothing should be worn at all times. Waste should be disposed of according to national and regional regulations. Safety Data Sheets (SDSs) are available upon request.

SYNTHESIS / CHARACTERIZATION:

Where possible, all of our products are synthesized using single-product unambiguous routes. They are then characterized, and their structures and purities confirmed, using a combination of the most relevant techniques, such as NMR, GC/MS, LC/MS/MS, SFC/UV/MS/MS, x-ray crystallography, and melting point. Isotopic purities of mass-labelled compounds are also confirmed using HRGC/HRMS and/or LC/MS/MS.

HOMOGENEITY:

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS, LC/MS/MS and/or SFC/UV/MS/MS. The relative response factors of the analyte of interest in each solution are required to be <5% RSD. New solution lots of existing products are compared to older lots in the same manner, which further confirms the homogeneity of the crystalline material as well as the stability and homogeneity of the solutions in the storage containers.

UNCERTAINTY:

The maximum combined relative standard uncertainty of our reference standard solutions is calculated using the following equation:

The combined relative standard uncertainty, $u_c(y)$, of a value y and the uncertainty of the independent parameters x_1, x_2, \dots, x_n on which it depends is:

$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

where x is expressed as a relative standard uncertainty of the individual parameter.

The individual uncertainties taken into account include those associated with weights (calibration of the balance) and volumes (calibration of the volumetric glassware). An expanded maximum combined percent relative uncertainty of $\pm 5\%$ (calculated with a coverage factor of 2 and a level of confidence of 95%) is stated on the Certificate of Analysis for all of our products.

TRACEABILITY:

All reference standard solutions are traceable to specific crystalline lots. The microbalances used for solution preparation are regularly tested by an external ISO/IEC 17025 accredited calibration company. In addition, their calibration is verified prior to each weighing using NIST and/or NRC traceable external weights. All volumetric glassware used is of Class A tolerance and has been tested according to the appropriate ASTM procedures, which are ultimately traceable to NIST. For certain products, traceability to international interlaboratory studies has also been established.

EXPIRY DATE / PERIOD OF VALIDITY:

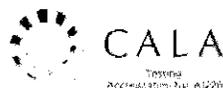
Ongoing stability studies of this product have demonstrated stability in its composition and concentration, until the specified expiry date, in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

LIMITED WARRANTY:

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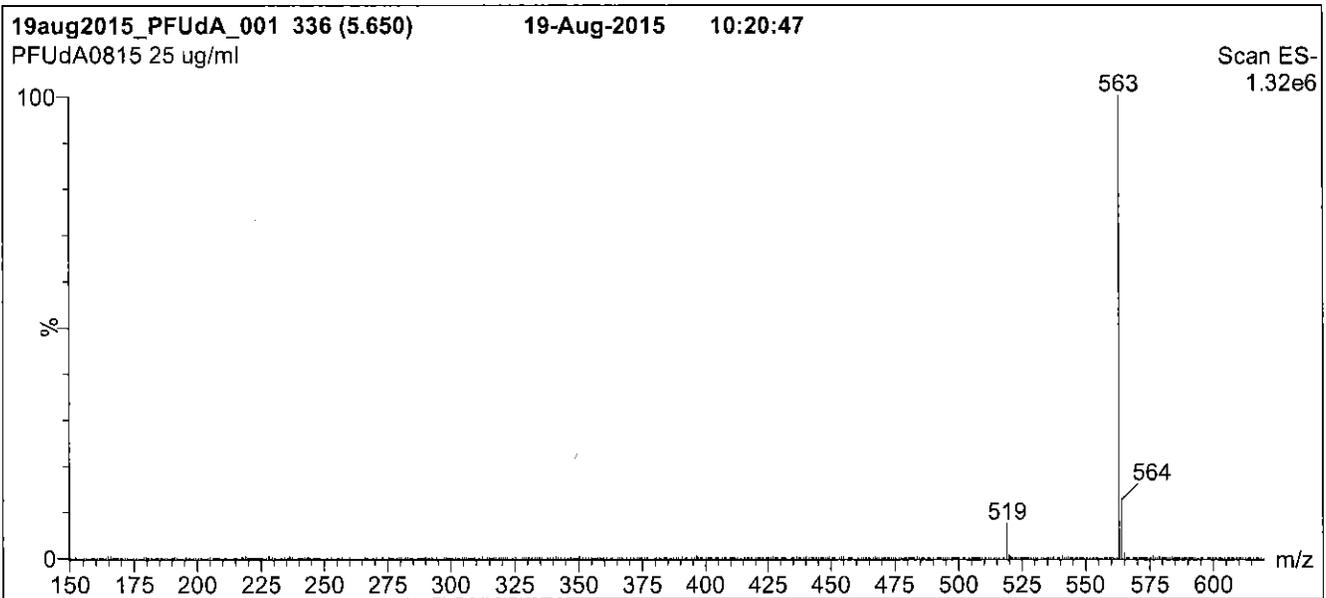
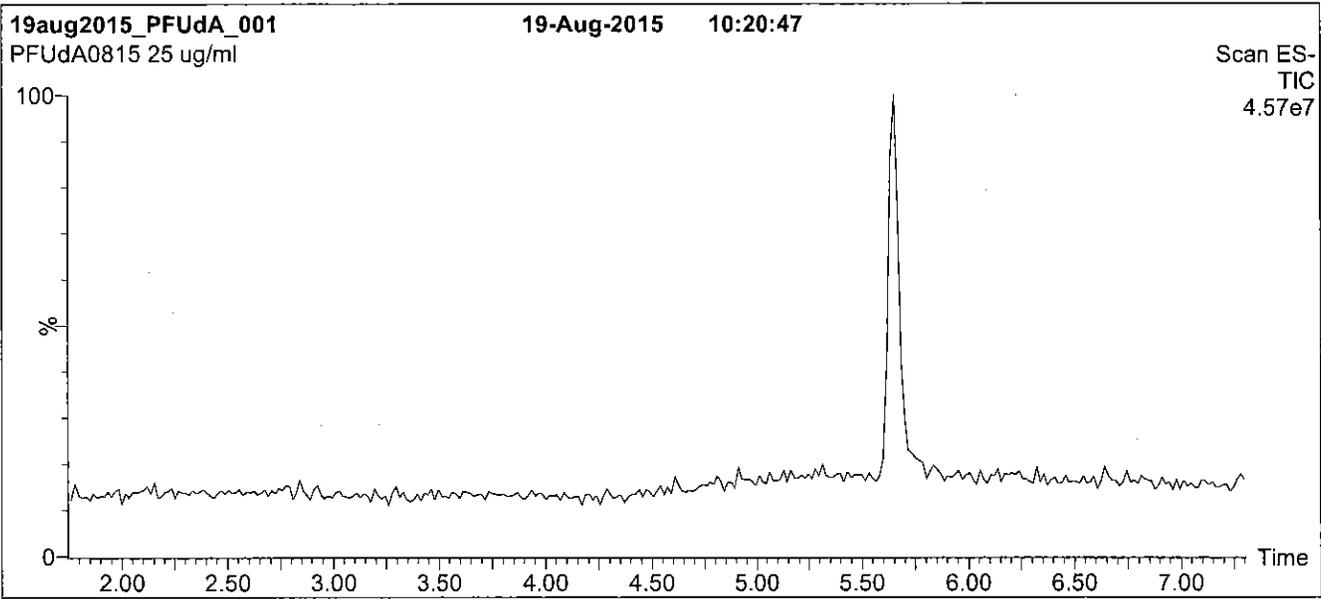
QUALITY MANAGEMENT:

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Figure 1: PFUdA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro *micro* API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
 1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
 Start: 50% (80:20 MeOH:ACN) / 50% H₂O
 (both with 10 mM NH₄OAc buffer)
 Ramp to 90% organic over 7 min and hold for 2 min
 before returning to initial conditions in 0.5 min.
 Time: 10 min

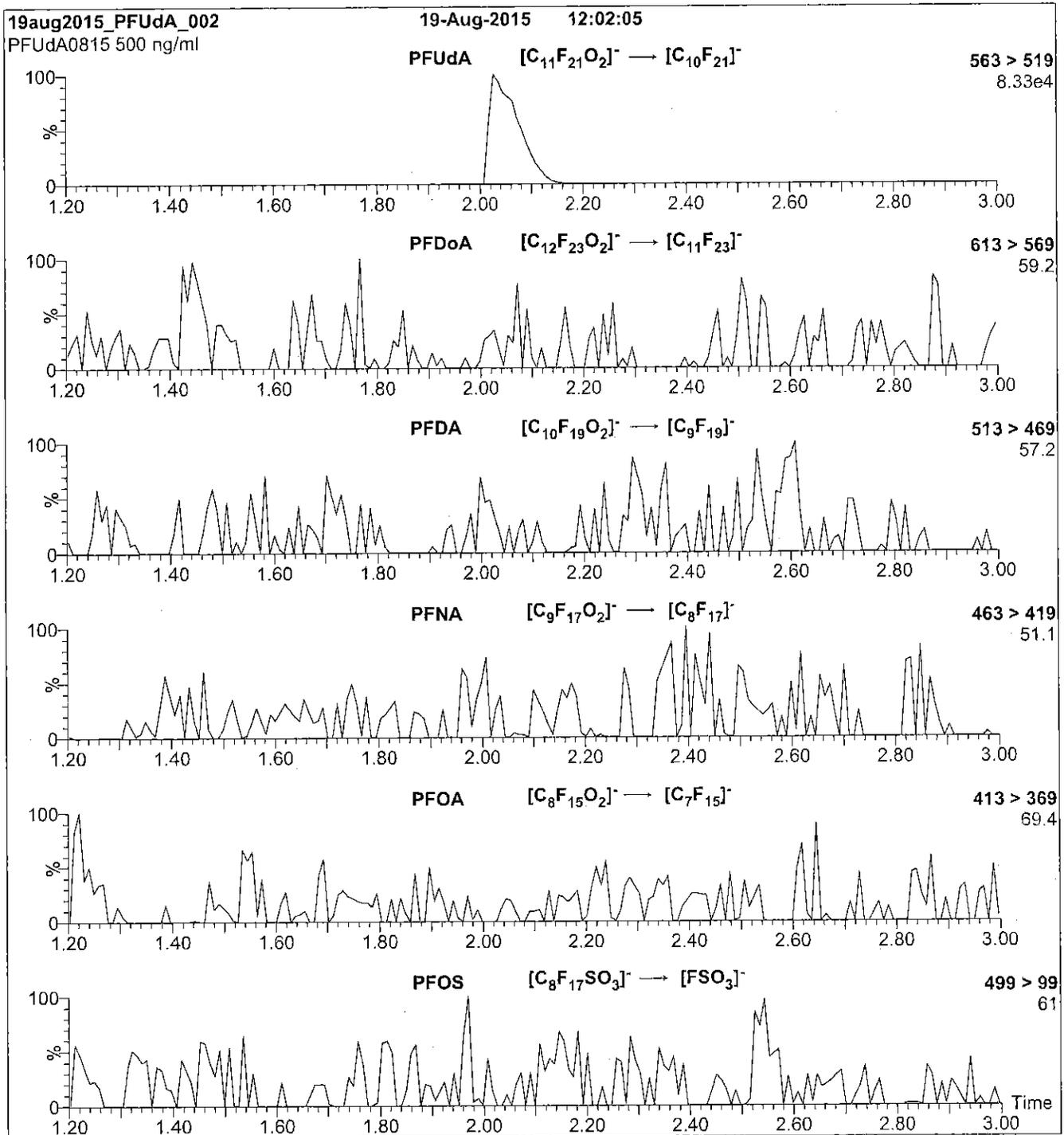
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
 Capillary Voltage (kV) = 3.00
 Cone Voltage (V) = 15.00
 Cone Gas Flow (l/hr) = 65
 Desolvation Gas Flow (l/hr) = 750

Figure 2: PFUdA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFUdA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
 Collision Energy (eV) = 11

Reagent

LCPFUdA_00005

Scanned
10/14/16 R: SBC 9/13/16

730535
ID: LCPFUdA_00005
Exp: 08/19/20 Prj: SBC
PF-n-undecanoic acid

730536
ID: LCPFUdA_00006
Exp: 08/19/20 Prj: SBC
PF-n-undecanoic acid

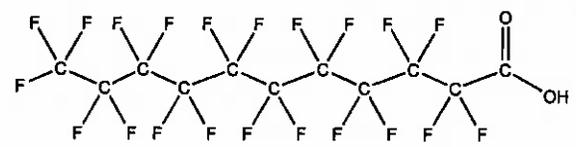


WELLINGTON LABORATORIES

CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFUdA **LOT NUMBER:** PFUdA0815
COMPOUND: Perfluoro-n-undecanoic acid

STRUCTURE: **CAS #:** 2058-94-8



MOLECULAR FORMULA: C₁₁H₂₁O₂ **MOLECULAR WEIGHT:** 564.09
CONCENTRATION: 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol
Water (<1%)
CHEMICAL PURITY: >98%
LAST TESTED: (mm/dd/yyyy) 08/19/2015
EXPIRY DATE: (mm/dd/yyyy) 08/19/2020
RECOMMENDED STORAGE: Store ampoule in a cool, dark place

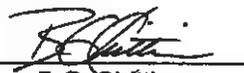
DOCUMENTATION/ DATA ATTACHED:

- Figure 1: LC/MS Data (TIC and Mass Spectrum)
- Figure 2: LC/MS/MS Data (Selected MRM Transitions)

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE

Certified By:  **Date:** 08/21/2015
B.G. Chittim (mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA
519-822-2436 • Fax: 519-822-2849 • info@well-labs.com

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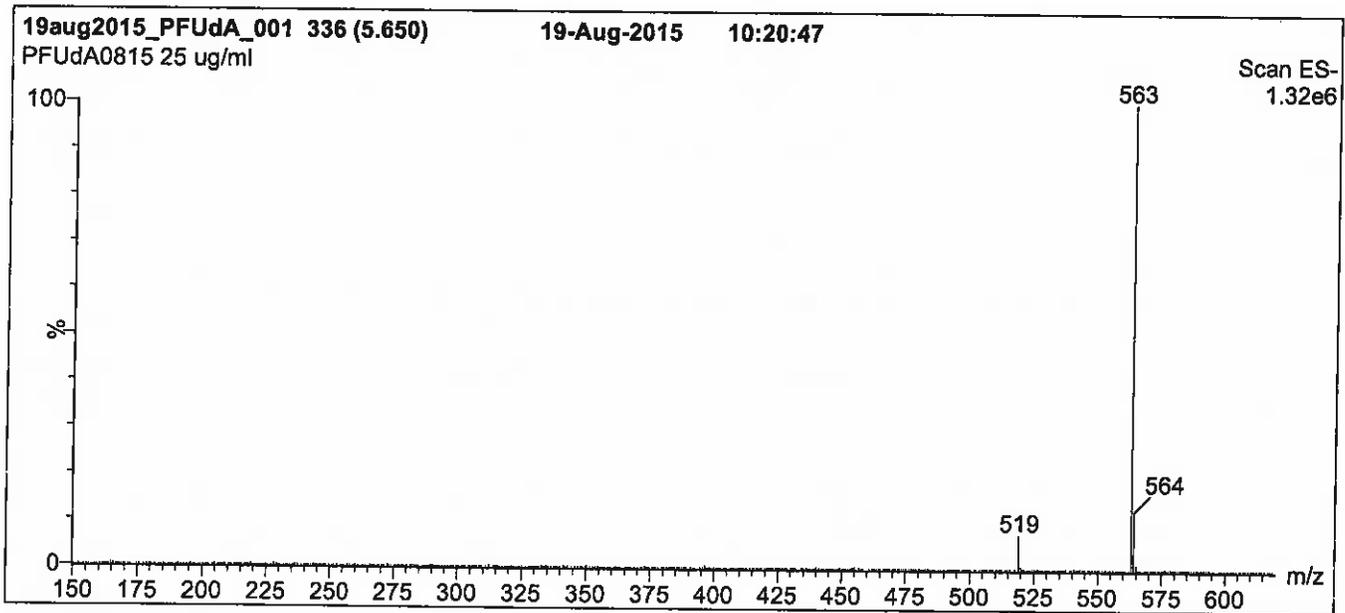
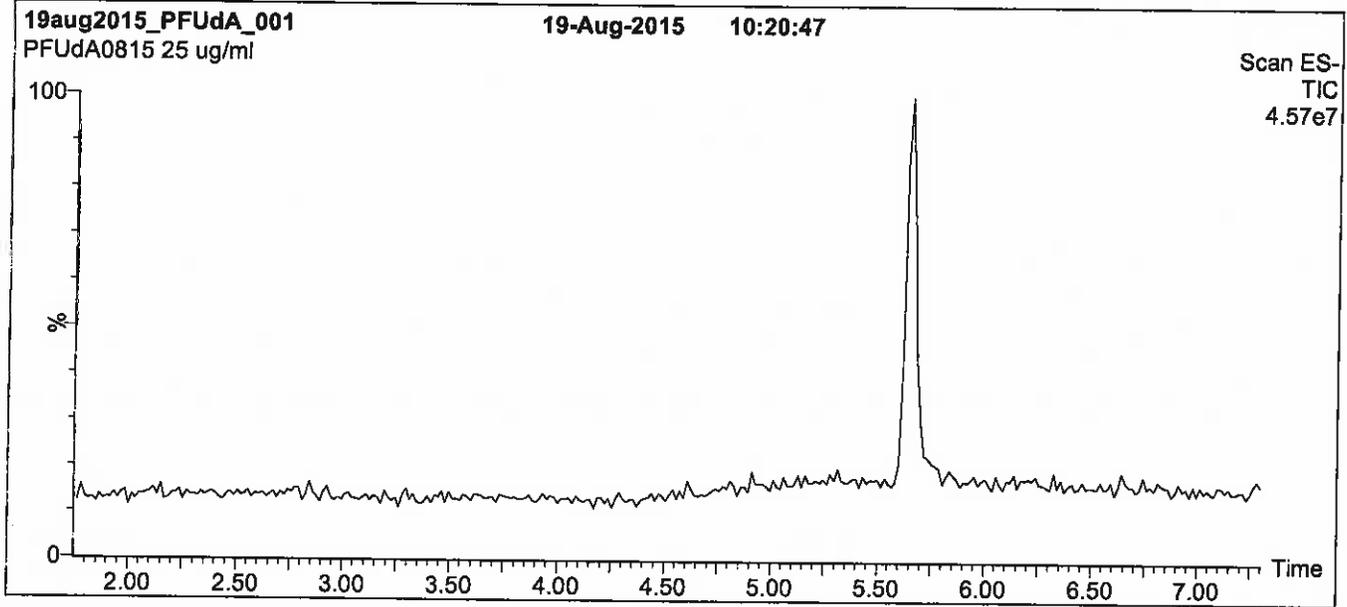
QUALITY MANAGEMENT:

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Figure 1: PFUdA; LC/MS Data (TIC and Mass Spectrum)



Conditions for Figure 1:

LC: Waters Acquity Ultra Performance LC
MS: Micromass Quattro micro API MS

Chromatographic Conditions

Column: Acquity UPLC BEH Shield RP₁₈
1.7 μ m, 2.1 x 100 mm

Mobile phase: Gradient
Start: 50% (80:20 MeOH:ACN) / 50% H₂O
(both with 10 mM NH₄OAc buffer)
Ramp to 90% organic over 7 min and hold for 2 min
before returning to initial conditions in 0.5 min.
Time: 10 min

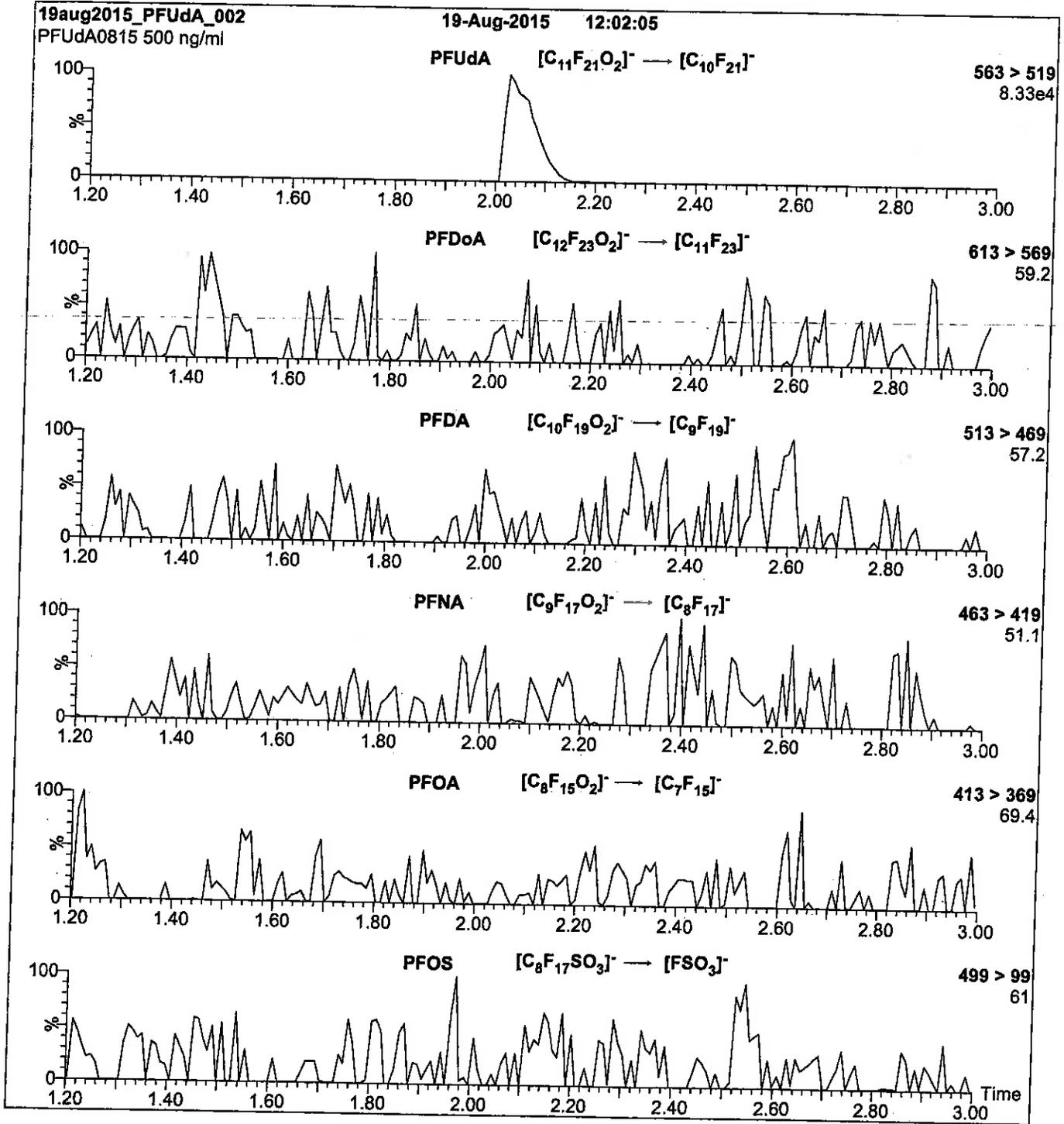
Flow: 300 μ l/min

MS Parameters

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)
Capillary Voltage (kV) = 3.00
Cone Voltage (V) = 15.00
Cone Gas Flow (l/hr) = 65
Desolvation Gas Flow (l/hr) = 750

Figure 2: PFUdA; LC/MS/MS Data (Selected MRM Transitions)



Conditions for Figure 2:

Injection: Direct loop injection
 10 μ l (500 ng/ml PFUdA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H₂O
 (both with 10 mM NH₄OAc buffer)

Flow: 300 μ l/min

MS Parameters

Collision Gas (mbar) = 3.31e-3
 Collision Energy (eV) = 11

Method PFC DOD

Perfluronated Hydrocarbons (LC/MS)
by Method PFC_DOD

FORM II
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Matrix: Solid Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFHxS #	PFOA #	PFOS #
DPT-16-35-SO-00-01	320-23783-1	86	93	73
DPT-16-35-SO-19-20	320-23783-2	108	112	101
DPT-16-19-SO-00-01	320-23783-5	85	96	54
DPT-16-19-SO-00-01 DL	320-23783-5 DL	106	136	90
DPT-16-19-SO-17-18	320-23783-6	73	69	64
DPT-16-32-SO-14-15	320-23783-7	70	96	62
DPT-16-31-SO-14-15	320-23783-8	93	108	82
DPT-16-30-SO-14-15	320-23783-9	90	101	63
DPT-16-28-SO-14-15	320-23783-10	89	105	61
DPT-16-15-SO-14-15	320-23783-11	89	101	66
	MB 320-139627/1-A	104	108	94
	LCS 320-139627/2-A	120	119	112

PFHxS = 1802 PFHxS	<u>QC LIMITS</u>
PFOA = 13C4 PFOA	25-150
PFOS = 13C4 PFOS	25-150
	25-150

Column to be used to flag recovery values

FORM II 537 (Modified)

FORM II
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Matrix: Water Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFHxS #	PFOA #	PFOS #
DPT-16-35-GW-21-25	320-23783-3	43	70	58
DPT-16-35-GW-21-25 DL	320-23783-3 DL	83	91	87
DPT-16-35-GW-31-35	320-23783-4	30	52	76
DPT-16-35-GW-31-35 DL	320-23783-4 DL	73	78	91
DPT-16-15-GW-18-22	320-23783-12	21 Q	23 Q	29
DPT-16-15-GW-18-22 DL	320-23783-12 DL	104	83	94
DPT-16-15-GW-31-35	320-23783-13	66	22 Q	82
DPT-16-15-GW-31-35 DL	320-23783-13 DL	101	33	95
	MB 320-139749/1-A	86	93	85
	LCS 320-139749/2-A	105	108	104
	LCSD 320-139749/3-A	104	116	104

PFHxS = 1802 PFHxS
PFOA = 13C4 PFOA
PFOS = 13C4 PFOS

QC LIMITS
25-150
25-150
25-150

Column to be used to flag recovery values

FORM II 537 (Modified)

FORM III
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Matrix: Solid Level: Low Lab File ID: 02DEC2016C_002.d
 Lab ID: LCS 320-139627/2-A Client ID: _____

COMPOUND	SPIKE ADDED (ug/Kg)	LCS CONCENTRATION (ug/Kg)	LCS % REC	QC LIMITS REC	#
Perfluorooctanoic acid (PFOA)	4.00	4.25	106	60-140	
Perfluorooctane Sulfonate (PFOS)	3.71	3.66	99	60-140	
13C4 PFOA	10.0	11.9	119	25-150	
13C4 PFOS	9.56	10.7	112	25-150	
Perfluorobutanesulfonic acid (PFBS)	3.54	4.03	114	50-150	
18O2 PFHxS	9.46	11.3	120	25-150	

Column to be used to flag recovery and RPD values
 FORM III 537 (Modified)

FORM III
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Matrix: Water Level: Low Lab File ID: 03DEC2016B_002.d
 Lab ID: LCS 320-139749/2-A Client ID: _____

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Perfluorooctanoic acid (PFOA)	0.0400	0.0296	74	60-140	
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0261	70	60-140	
13C4 PFOA	0.100	0.108	108	25-150	
13C4 PFOS	0.0956	0.0995	104	25-150	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0289	82	50-150	
18O2 PFHxS	0.0946	0.0995	105	25-150	

Column to be used to flag recovery and RPD values
 FORM III 537 (Modified)

FORM III
LCMS LAB CONTROL SAMPLE DUPLICATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Matrix: Water Level: Low Lab File ID: 03DEC2016B_003.d
 Lab ID: LCSD 320-139749/3-A Client ID: _____

COMPOUND	SPIKE ADDED (ug/L)	LCSD CONCENTRATION (ug/L)	LCSD % REC	% RPD	QC LIMITS		#
					RPD	REC	
Perfluorooctanoic acid (PFOA)	0.0400	0.0314	79	6	30	60-140	
Perfluorooctane Sulfonate (PFOS)	0.0371	0.0289	78	10	30	60-140	
13C4 PFOA	0.100	0.116	116			25-150	
13C4 PFOS	0.0956	0.0990	104			25-150	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0325	92	12	30	50-150	
18O2 PFHxS	0.0946	0.0981	104			25-150	

Column to be used to flag recovery and RPD values
 FORM III 537 (Modified)

FORM IV
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab File ID: 02DEC2016C_001.d Lab Sample ID: MB 320-139627/1-A
 Matrix: Solid Date Extracted: 11/28/2016 11:03
 Instrument ID: A8_N Date Analyzed: 12/02/2016 14:59
 Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-139627/2-A	02DEC2016C_002.d	12/02/2016 15:07
DPT-16-35-SO-00-01	320-23783-1	02DEC2016C_008.d	12/02/2016 15:52
DPT-16-35-SO-19-20	320-23783-2	02DEC2016C_009.d	12/02/2016 15:59
DPT-16-19-SO-00-01	320-23783-5	02DEC2016C_010.d	12/02/2016 16:07
DPT-16-19-SO-17-18	320-23783-6	02DEC2016C_015.d	12/02/2016 16:44
DPT-16-32-SO-14-15	320-23783-7	02DEC2016C_016.d	12/02/2016 16:52
DPT-16-31-SO-14-15	320-23783-8	02DEC2016C_017.d	12/02/2016 16:59
DPT-16-30-SO-14-15	320-23783-9	02DEC2016C_018.d	12/02/2016 17:07
DPT-16-28-SO-14-15	320-23783-10	02DEC2016C_019.d	12/02/2016 17:14
DPT-16-15-SO-14-15	320-23783-11	02DEC2016C_020.d	12/02/2016 17:22
DPT-16-19-SO-00-01 DL	320-23783-5 DL	07DEC2016B_009.d	12/07/2016 17:33

FORM IV
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab File ID: 03DEC2016B_001.d Lab Sample ID: MB 320-139749/1-A
 Matrix: Water Date Extracted: 11/28/2016 16:55
 Instrument ID: A8_N Date Analyzed: 12/03/2016 16:03
 Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-139749/2-A	03DEC2016B_002.d	12/03/2016 16:11
	LCSD 320-139749/3-A	03DEC2016B_003.d	12/03/2016 16:18
DPT-16-35-GW-21-25	320-23783-3	03DEC2016B_004.d	12/03/2016 16:26
DPT-16-35-GW-31-35	320-23783-4	03DEC2016B_005.d	12/03/2016 16:33
DPT-16-15-GW-18-22	320-23783-12	03DEC2016B_006.d	12/03/2016 16:41
DPT-16-15-GW-31-35	320-23783-13	03DEC2016B_007.d	12/03/2016 16:48
DPT-16-15-GW-18-22 DL	320-23783-12 DL	07DEC2016B_005.d	12/07/2016 17:03
DPT-16-35-GW-21-25 DL	320-23783-3 DL	07DEC2016B_006.d	12/07/2016 17:10
DPT-16-35-GW-31-35 DL	320-23783-4 DL	07DEC2016B_007.d	12/07/2016 17:18
DPT-16-15-GW-31-35 DL	320-23783-13 DL	07DEC2016B_008.d	12/07/2016 17:25

FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-SO-00-01 Lab Sample ID: 320-23783-1
 Matrix: Solid Lab File ID: 02DEC2016C_008.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 09:24
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.01(g) Date Analyzed: 12/02/2016 15:52
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 13.5 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	12	M	0.58	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	28		0.58	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.21	J	0.46	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	93		25-150
STL00991	13C4 PFOS	73		25-150
STL00994	18O2 PFHxS	86		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_008.d
 Lims ID: 320-23783-A-1-A
 Client ID: DPT-16-35-SO-00-01
 Sample Type: Client
 Inject. Date: 02-Dec-2016 15:52:08 ALS Bottle#: 8 Worklist Smp#: 12
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-1-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:25:25 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:22:28

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.900	1.907	-0.007	1.000	376131	0.9011				
298.90 > 99.00	1.900	1.907	-0.007	1.000	161964		2.32(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.527	2.534	-0.007		12236591	40.5		85.5	1572000	
D 14 13C4 PFOA										
417.00 > 372.00	2.886	2.886	0.0		10158301	46.3		92.6	967819	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.886	2.894	-0.008	1.000	10474209	50.5			59538	M
413.00 > 169.00	2.886	2.894	-0.008	1.000	6995834		1.50(0.90-1.10)		279169	M
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.268	3.267	0.001	1.000	23227692	119.5			623949	
499.00 > 99.00	3.268	3.267	0.001	1.000	4818749		4.82(0.90-1.10)		175438	
D 17 13C4 PFOS										
503.00 > 80.00	3.268	3.267	0.001		8546516	34.9		73.1	153218	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_008.d

Injection Date: 02-Dec-2016 15:52:08

Instrument ID: A8_N

Lims ID: 320-23783-A-1-A

Lab Sample ID: 320-23783-1

Client ID: DPT-16-35-SO-00-01

Operator ID: A8-PC\A8

ALS Bottle#: 8

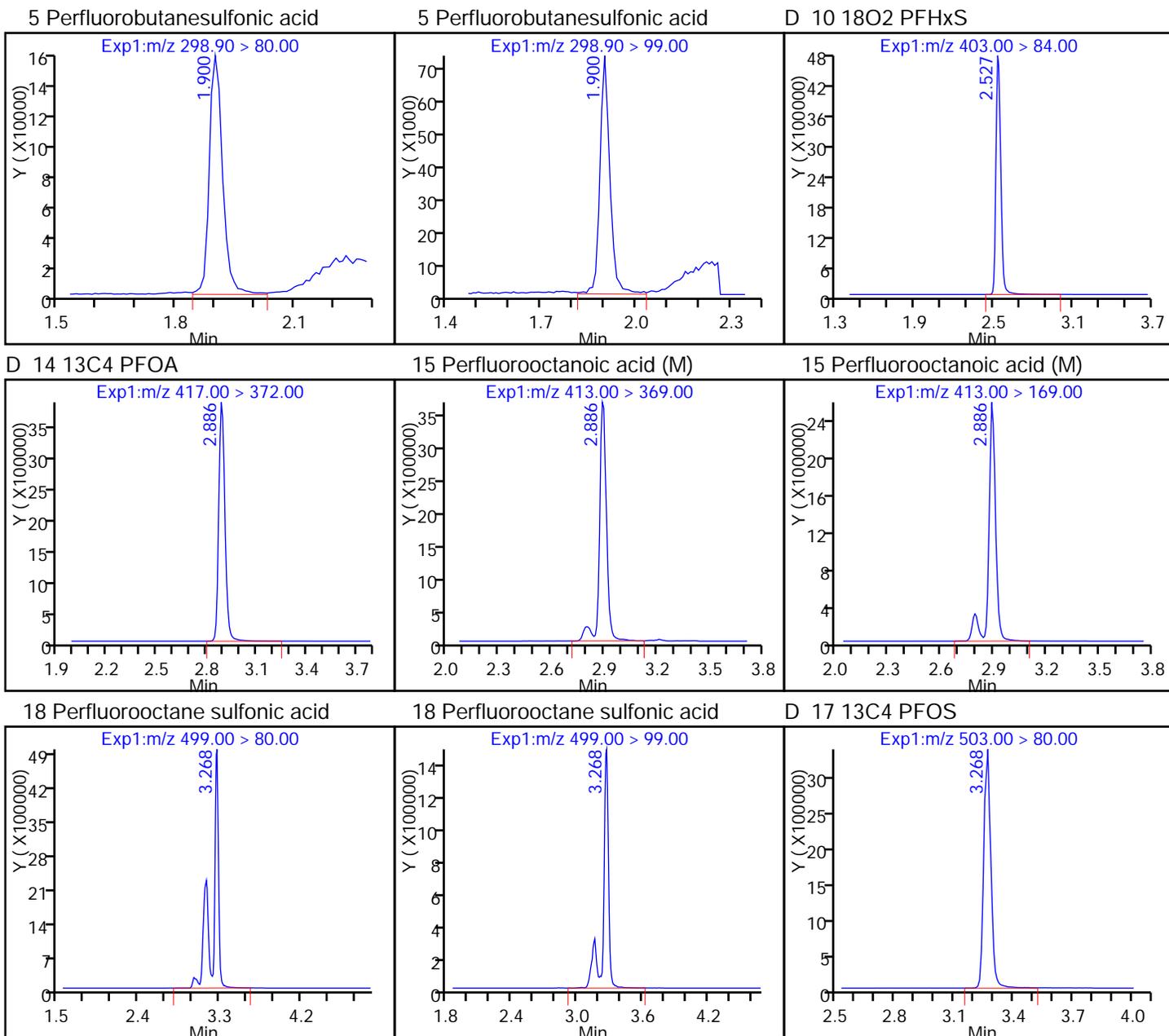
Worklist Smp#: 12

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

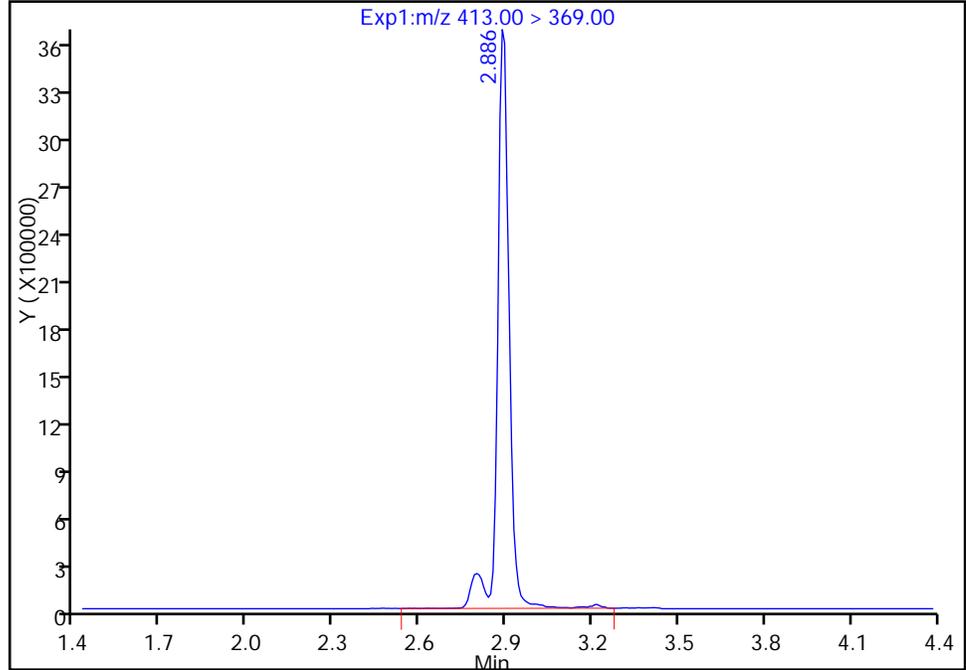
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Injection Date: 02-Dec-2016 15:52:08 Instrument ID: A8_N
Lims ID: 320-23783-A-1-A Lab Sample ID: 320-23783-1
Client ID: DPT-16-35-SO-00-01
Operator ID: A8-PC\A8 ALS Bottle#: 8 Worklist Smp#: 12
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

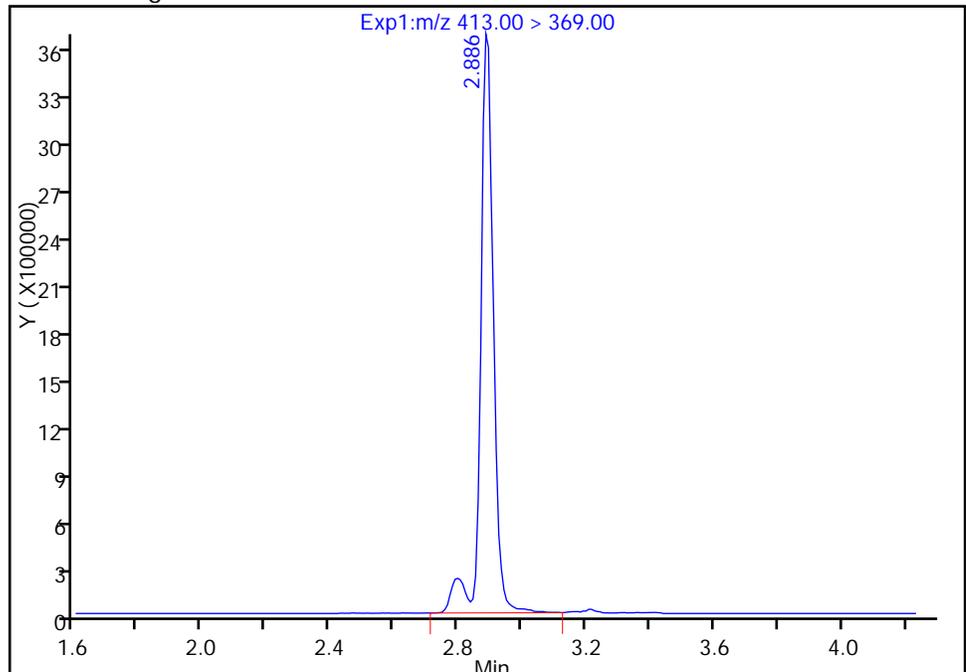
RT: 2.89
Area: 10606743
Amount: 51.168180
Amount Units: ng/ml

Processing Integration Results



RT: 2.89
Area: 10474209
Amount: 50.528821
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:22:28
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

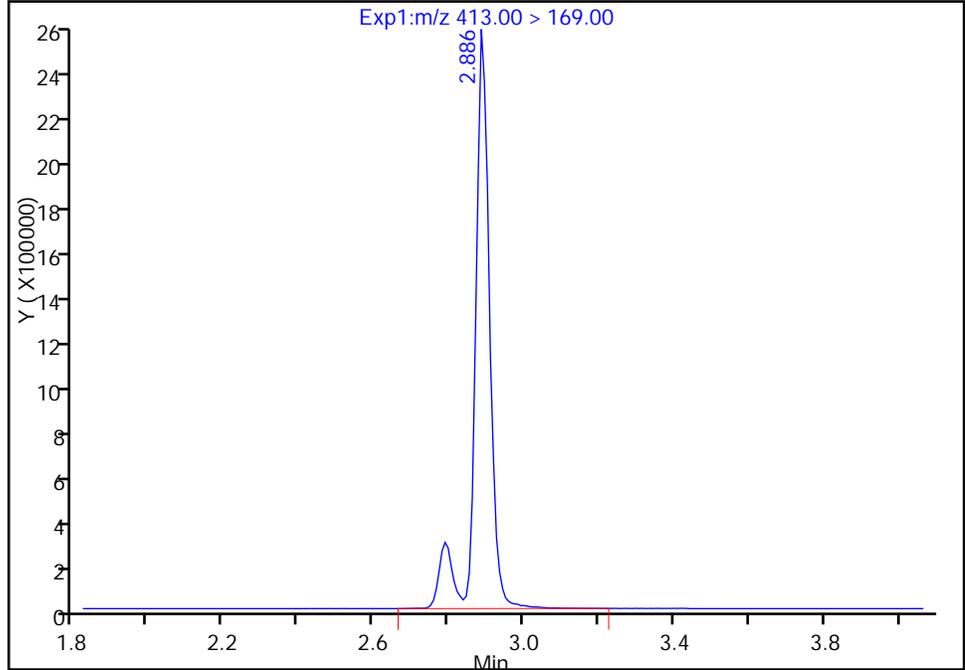
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Injection Date: 02-Dec-2016 15:52:08 Instrument ID: A8_N
Lims ID: 320-23783-A-1-A Lab Sample ID: 320-23783-1
Client ID: DPT-16-35-SO-00-01
Operator ID: A8-PC\A8 ALS Bottle#: 8 Worklist Smp#: 12
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 2

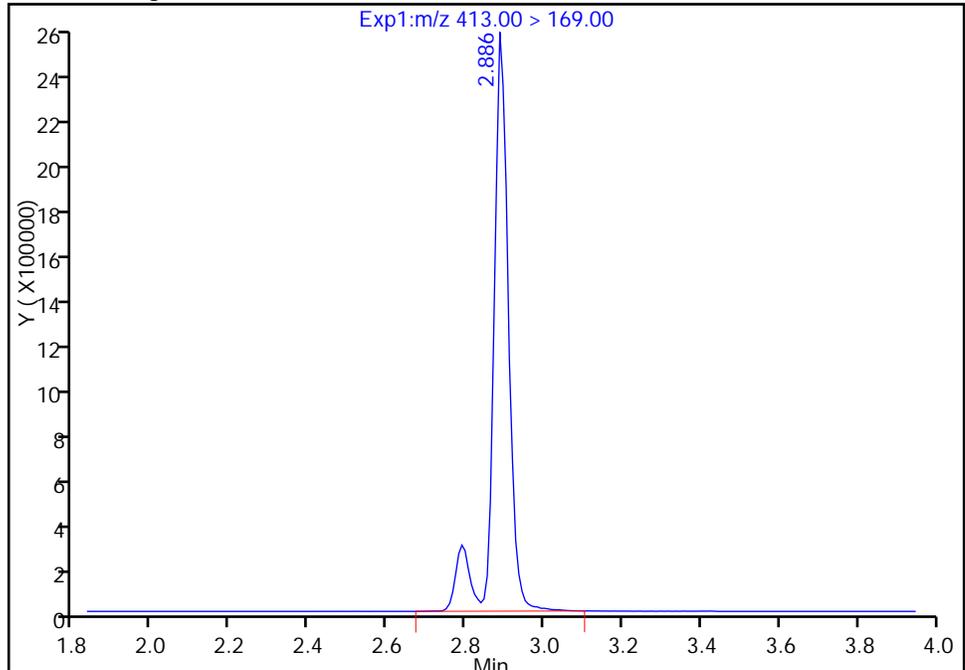
RT: 2.89
Area: 7020249
Amount: 51.168180
Amount Units: ng/ml

Processing Integration Results



RT: 2.89
Area: 6995834
Amount: 50.528821
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:22:28

Audit Action: Manually Integrated

Audit Reason: Baseline

FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-SO-19-20 Lab Sample ID: 320-23783-2
 Matrix: Solid Lab File ID: 02DEC2016C_009.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 09:36
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.07(g) Date Analyzed: 12/02/2016 15:59
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 17.2 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.56	J	0.60	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	4.0	M	0.60	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	112		25-150
STL00991	13C4 PFOS	101		25-150
STL00994	18O2 PFHxS	108		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_009.d
 Lims ID: 320-23783-A-2-A
 Client ID: DPT-16-35-SO-19-20
 Sample Type: Client
 Inject. Date: 02-Dec-2016 15:59:39 ALS Bottle#: 9 Worklist Smp#: 13
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-2-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:25:25 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:23:39

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
--------	----	--------	--------	--------	----------	--------------	---------------	------	-----	-------

5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.897	1.907	-0.010	1.000	139624	0.2652				
298.90 > 99.00	1.897	1.907	-0.010	1.000	63391		2.20(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.538	2.534	0.004		15434355	51.0		108	736230	
D 14 13C4 PFOA										
417.00 > 372.00	2.888	2.886	0.002		12258453	55.9		112	1106294	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.888	2.894	-0.006	1.000	591556	2.36			7002	
413.00 > 169.00	2.888	2.894	-0.006	1.000	387219		1.53(0.90-1.10)		22352	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.270	3.267	0.003	1.000	4485030	16.7			140297	M
499.00 > 99.00	3.270	3.267	0.003	1.000	1003003		4.47(0.90-1.10)		39467	M
D 17 13C4 PFOS										
503.00 > 80.00	3.262	3.267	-0.005		11777207	48.1		101	763283	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_009.d

Injection Date: 02-Dec-2016 15:59:39

Instrument ID: A8_N

Lims ID: 320-23783-A-2-A

Lab Sample ID: 320-23783-2

Client ID: DPT-16-35-SO-19-20

Operator ID: A8-PC\A8

ALS Bottle#: 9

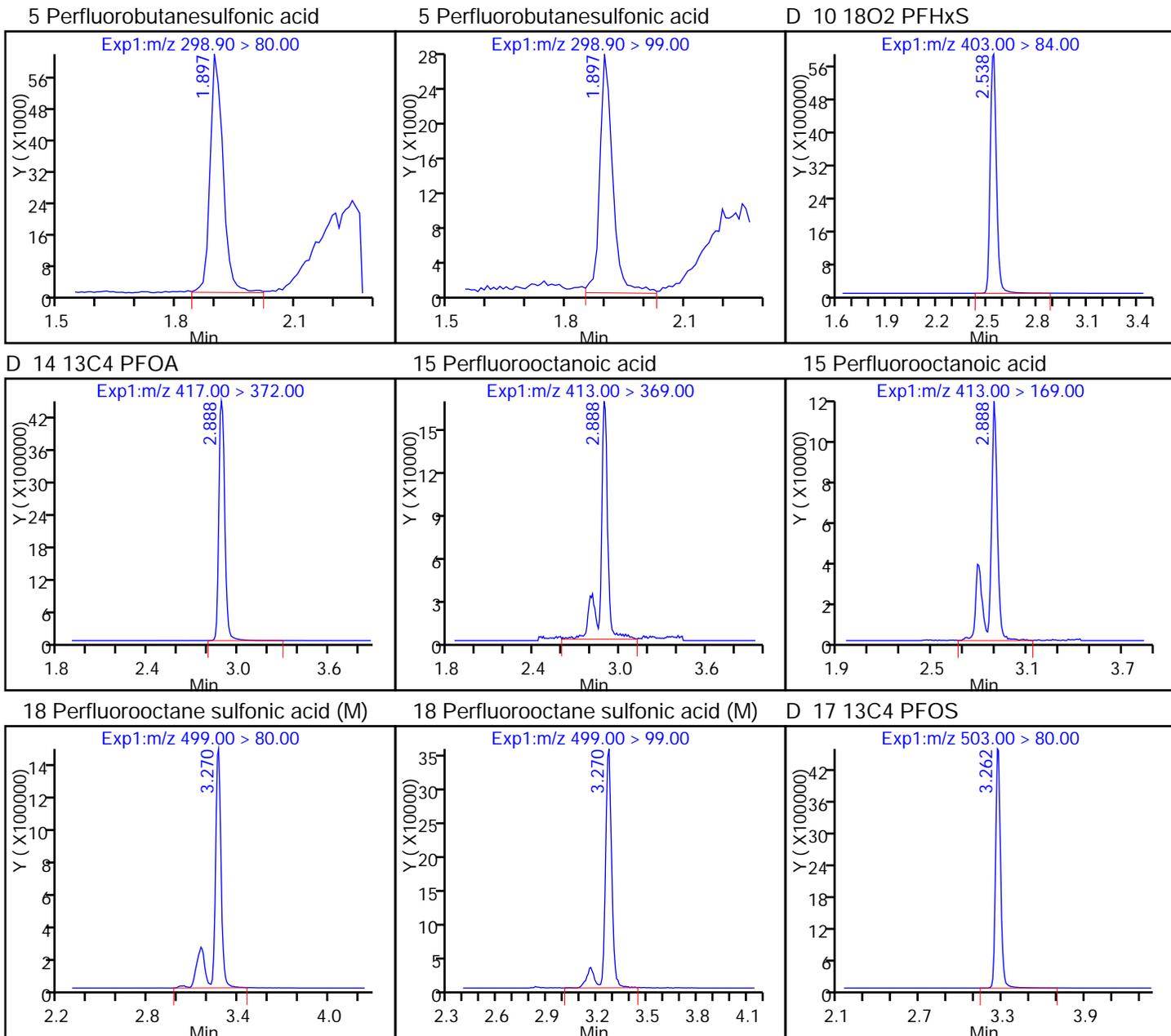
Worklist Smp#: 13

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

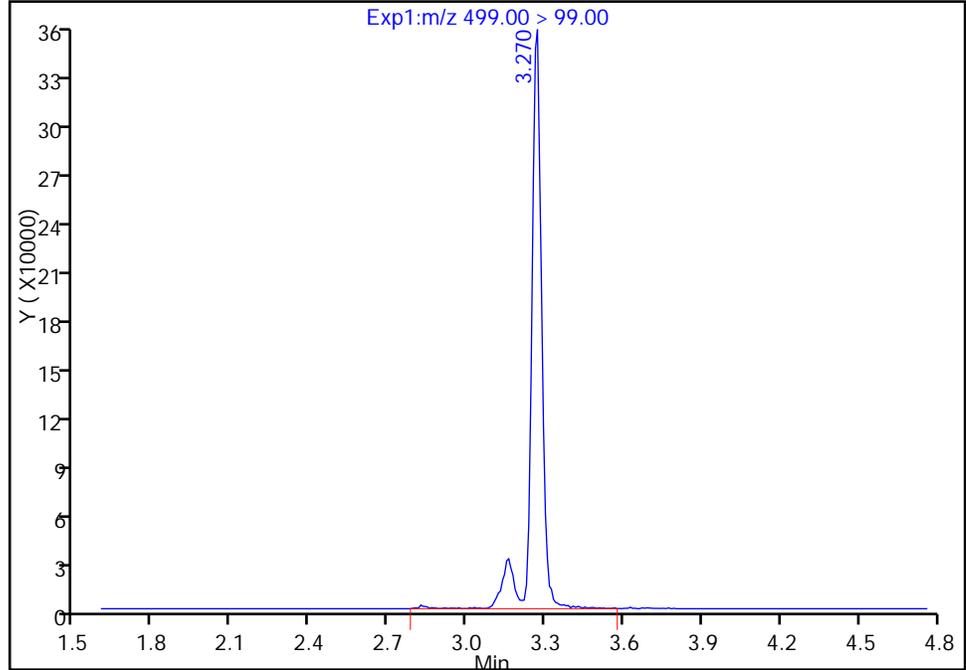
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Injection Date: 02-Dec-2016 15:59:39 Instrument ID: A8_N
Lims ID: 320-23783-A-2-A Lab Sample ID: 320-23783-2
Client ID: DPT-16-35-SO-19-20
Operator ID: A8-PC\A8 ALS Bottle#: 9 Worklist Smp#: 13
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

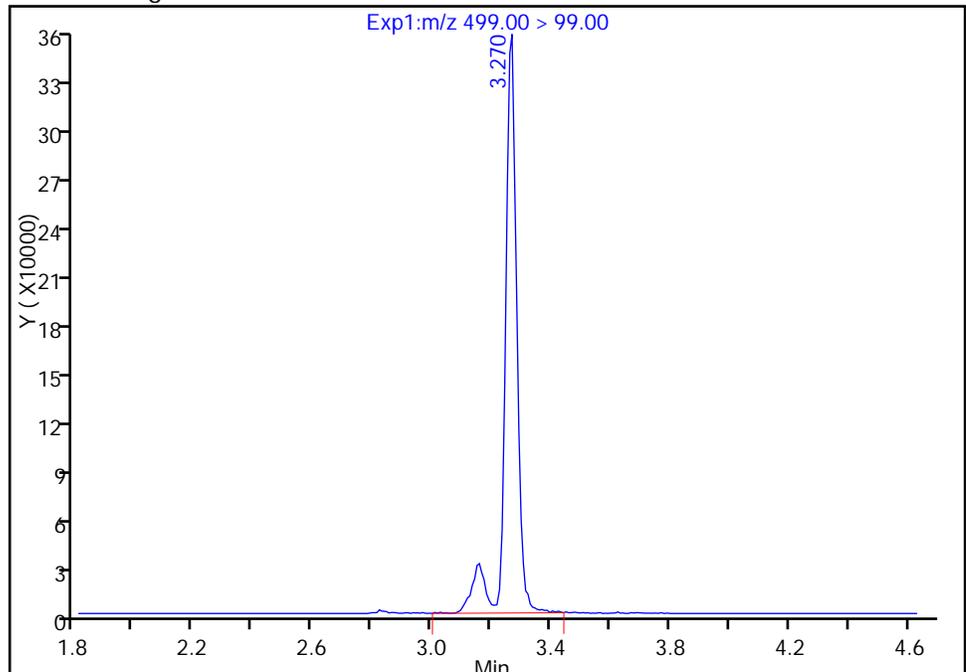
RT: 3.27
Area: 1020091
Amount: 16.875521
Amount Units: ng/ml

Processing Integration Results



RT: 3.27
Area: 1003003
Amount: 16.741126
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:23:39

Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

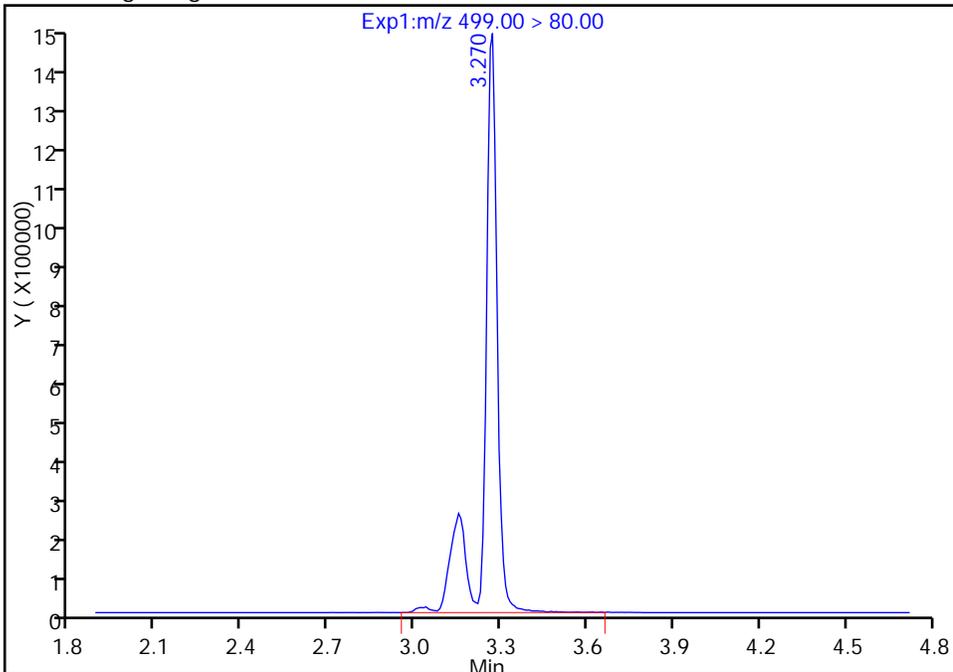
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_009.d
Injection Date: 02-Dec-2016 15:59:39 Instrument ID: A8_N
Lims ID: 320-23783-A-2-A Lab Sample ID: 320-23783-2
Client ID: DPT-16-35-SO-19-20
Operator ID: A8-PC\A8 ALS Bottle#: 9 Worklist Smp#: 13
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

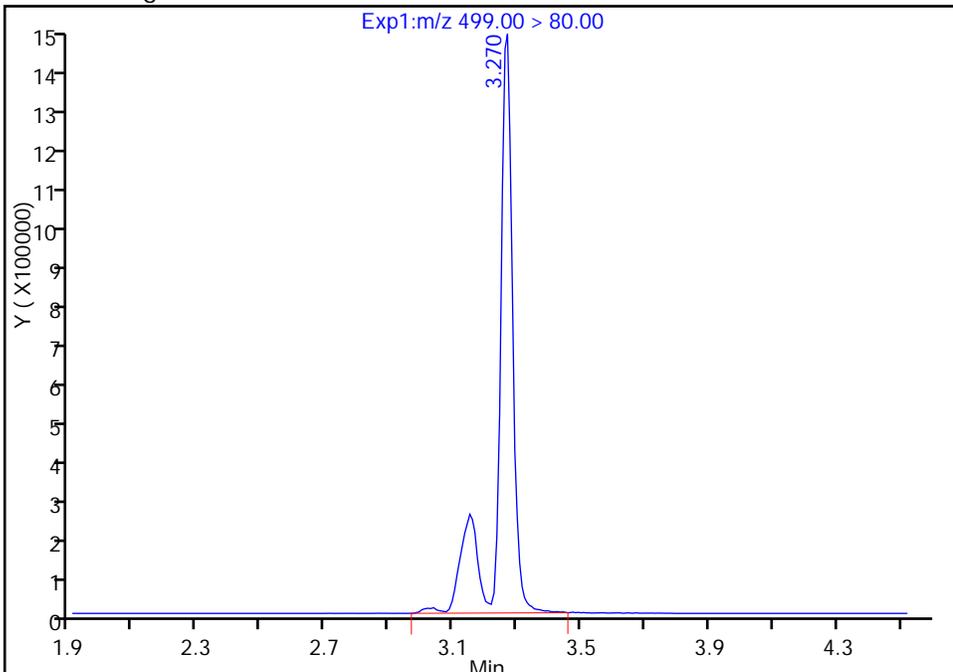
RT: 3.27
Area: 4521035
Amount: 16.875521
Amount Units: ng/ml

Processing Integration Results



RT: 3.27
Area: 4485030
Amount: 16.741126
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-GW-21-25 Lab Sample ID: 320-23783-3
 Matrix: Water Lab File ID: 03DEC2016B_004.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 10:10
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 250.2 (mL) Date Analyzed: 12/03/2016 16:26
 Con. Extract Vol.: 0.5 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.60	E M	0.0025	0.00075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.9	E M	0.0040	0.0013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.28		0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	70		25-150
STL00991	13C4 PFOS	58		25-150
STL00994	18O2 PFHxS	43		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d
 Lims ID: 320-23783-A-3-A
 Client ID: DPT-16-35-GW-21-25
 Sample Type: Client
 Inject. Date: 03-Dec-2016 16:26:12 ALS Bottle#: 35 Worklist Smp#: 25
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-3-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:45:18 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:31:29

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.878	1.900	-0.022	1.000	29884773	138.9				
298.90 > 99.00	1.878	1.900	-0.022	1.000	13466947		2.22(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.496	2.531	-0.035		6375608	20.4		43.2	910398	
D 14 13C4 PFOA										
417.00 > 372.00	2.848	2.880	-0.032		7643048	34.9		69.7	414375	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.848	2.887	-0.039	1.000	49480944	302.0			255484	EM
413.00 > 169.00	2.848	2.887	-0.039	1.000	37577807		1.32(0.90-1.10)		0.0	M
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.225	3.258	-0.033	1.000	142928519	926.0			0.0	EM
499.00 > 99.00	3.225	3.258	-0.033	1.000	38827184		3.68(0.90-1.10)		0.0	M
D 17 13C4 PFOS										
503.00 > 80.00	3.225	3.259	-0.034		6798776	27.6		57.8	26873	

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d

Injection Date: 03-Dec-2016 16:26:12

Instrument ID: A8_N

Lims ID: 320-23783-A-3-A

Lab Sample ID: 320-23783-3

Client ID: DPT-16-35-GW-21-25

Operator ID: A8-PC\A8

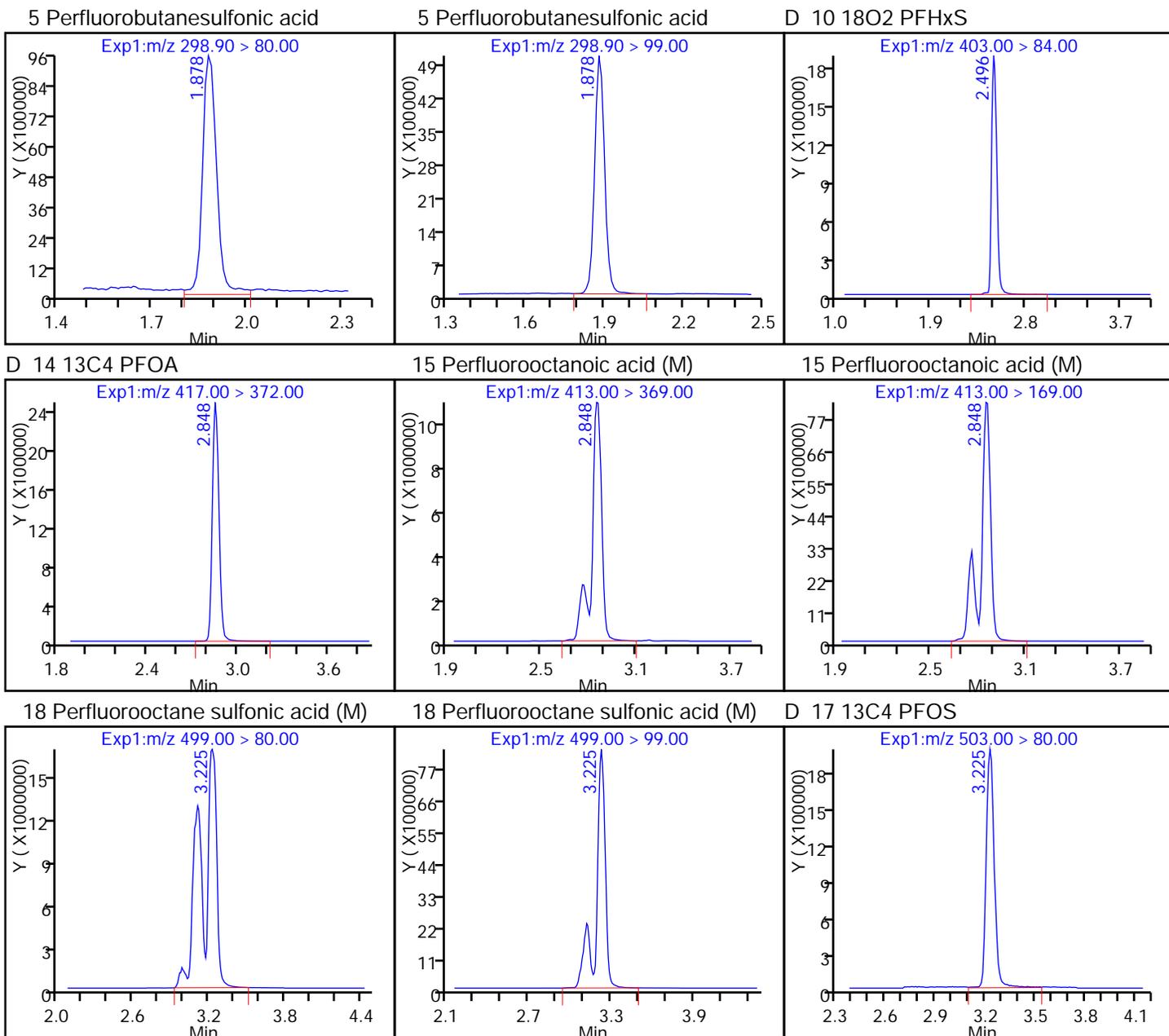
ALS Bottle#: 35 Worklist Smp#: 25

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

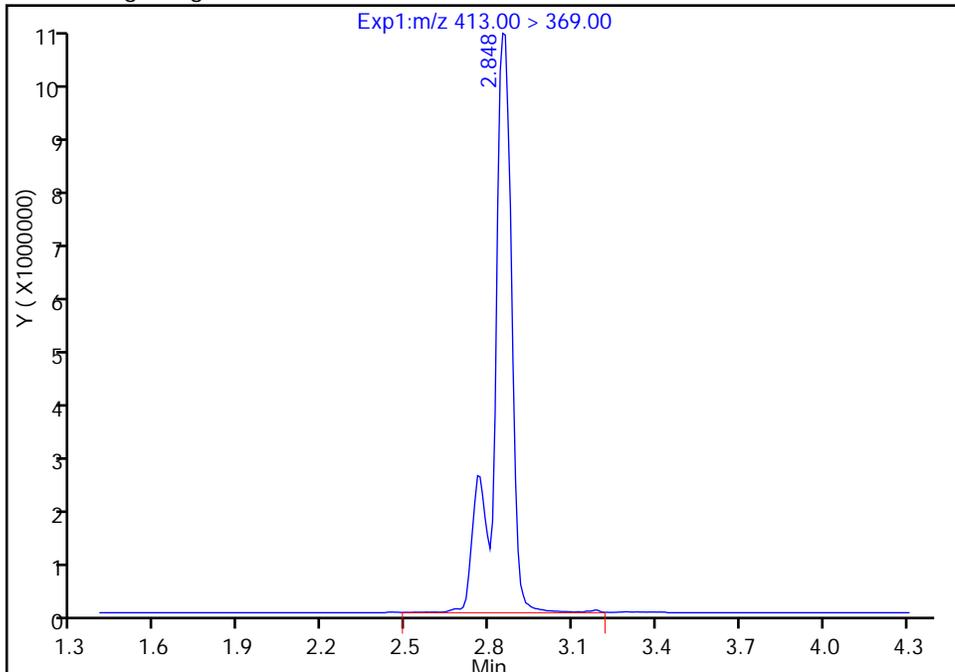
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d
Injection Date: 03-Dec-2016 16:26:12 Instrument ID: A8_N
Lims ID: 320-23783-A-3-A Lab Sample ID: 320-23783-3
Client ID: DPT-16-35-GW-21-25
Operator ID: A8-PC\A8 ALS Bottle#: 35 Worklist Smp#: 25
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

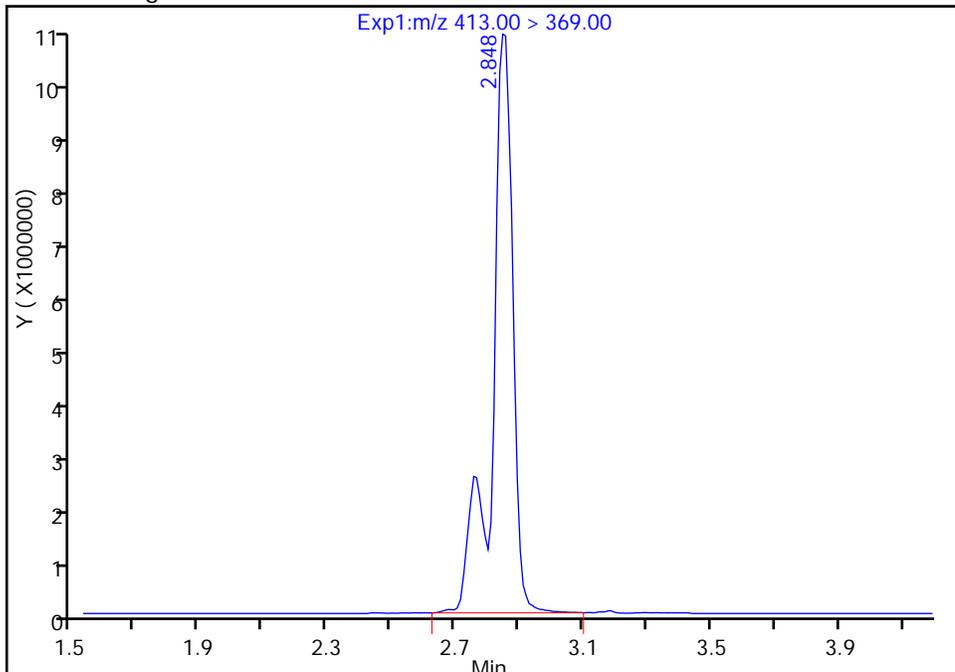
RT: 2.85
Area: 50072994
Amount: 305.5952
Amount Units: ng/ml

Processing Integration Results



RT: 2.85
Area: 49480944
Amount: 301.9820
Amount Units: ng/ml

Manual Integration Results



TestAmerica Sacramento

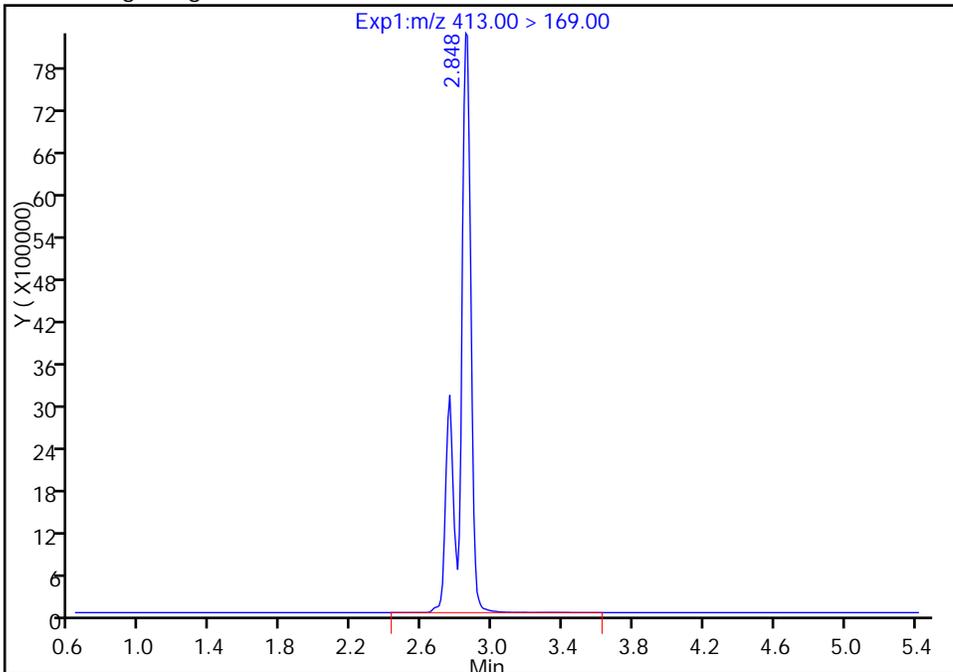
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d
Injection Date: 03-Dec-2016 16:26:12 Instrument ID: A8_N
Lims ID: 320-23783-A-3-A Lab Sample ID: 320-23783-3
Client ID: DPT-16-35-GW-21-25
Operator ID: A8-PC\A8 ALS Bottle#: 35 Worklist Smp#: 25
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 2

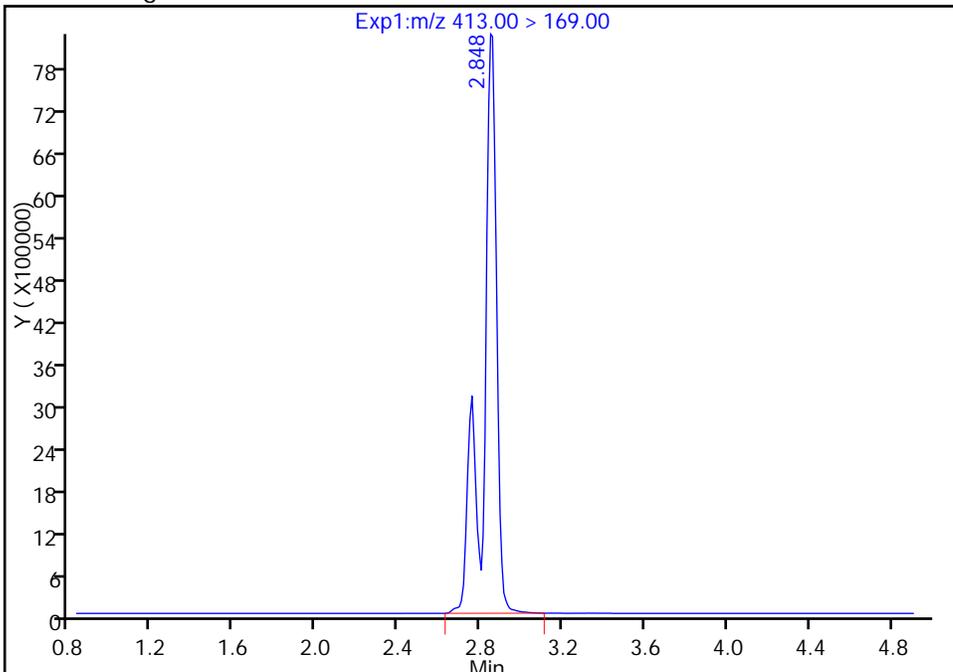
RT: 2.85
Area: 37697521
Amount: 305.5952
Amount Units: ng/ml

Processing Integration Results



RT: 2.85
Area: 37577807
Amount: 301.9820
Amount Units: ng/ml

Manual Integration Results



TestAmerica Sacramento

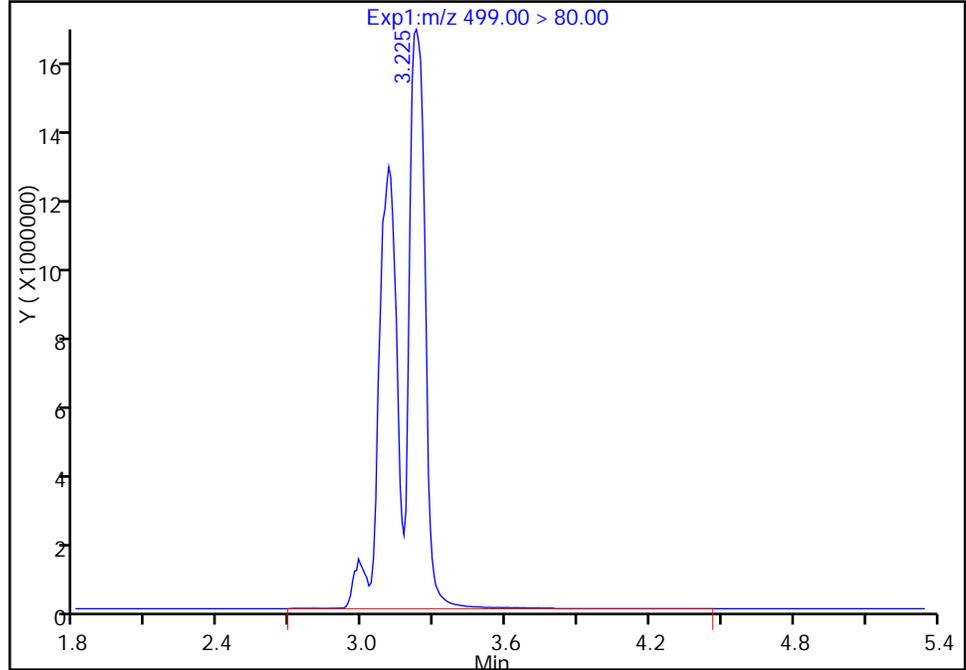
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d
Injection Date: 03-Dec-2016 16:26:12 Instrument ID: A8_N
Lims ID: 320-23783-A-3-A Lab Sample ID: 320-23783-3
Client ID: DPT-16-35-GW-21-25
Operator ID: A8-PC\A8 ALS Bottle#: 35 Worklist Smp#: 25
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

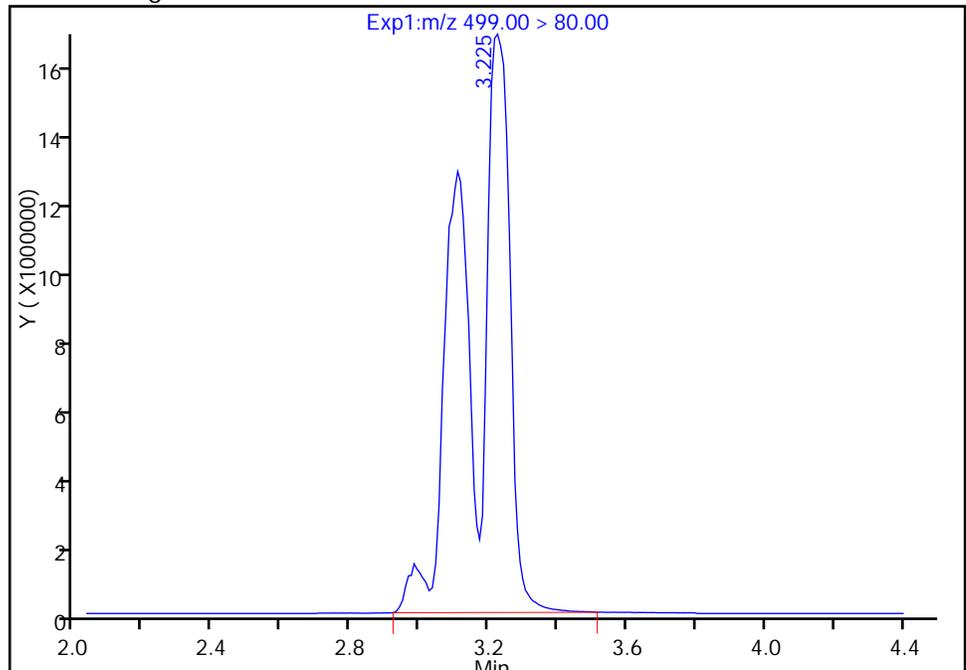
RT: 3.22
Area: 144567877
Amount: 936.5901
Amount Units: ng/ml

Processing Integration Results



RT: 3.22
Area: 142928519
Amount: 925.9694
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 10:31:29

Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

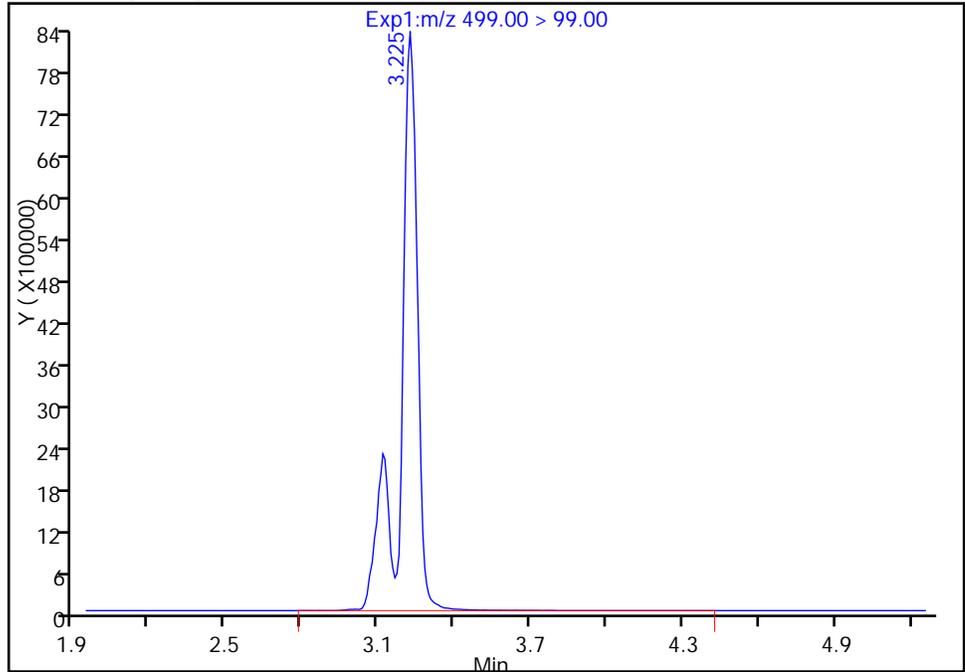
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_004.d
Injection Date: 03-Dec-2016 16:26:12 Instrument ID: A8_N
Lims ID: 320-23783-A-3-A Lab Sample ID: 320-23783-3
Client ID: DPT-16-35-GW-21-25
Operator ID: A8-PC\A8 ALS Bottle#: 35 Worklist Smp#: 25
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

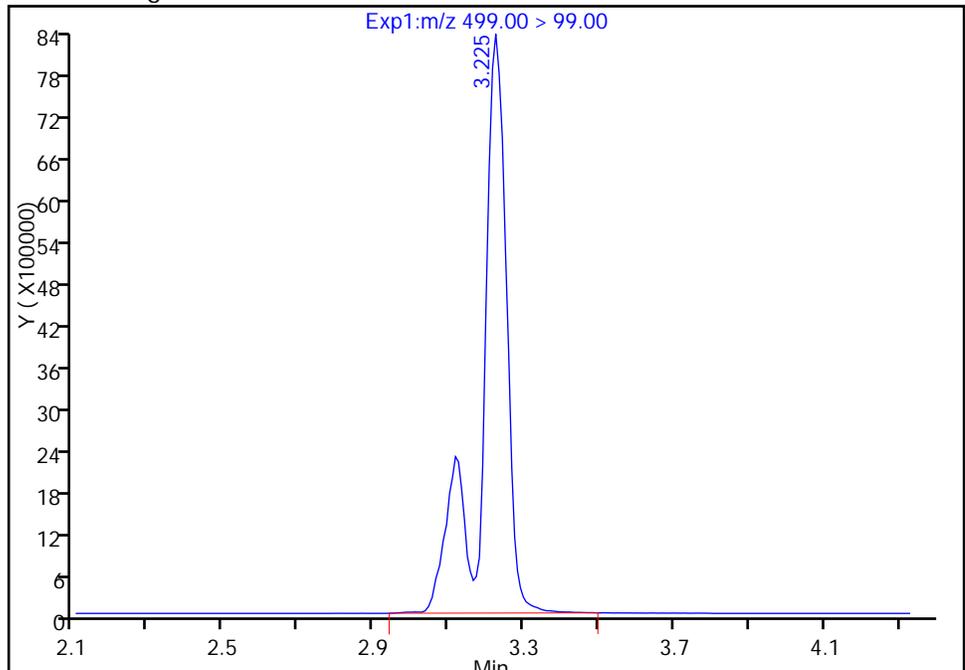
RT: 3.22
Area: 39077186
Amount: 936.5901
Amount Units: ng/ml

Processing Integration Results



RT: 3.22
Area: 38827184
Amount: 925.9694
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-GW-21-25 DL Lab Sample ID: 320-23783-3 DL
 Matrix: Water Lab File ID: 07DEC2016B_006.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 10:10
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 250.2 (mL) Date Analyzed: 12/07/2016 17:10
 Con. Extract Vol.: 0.5 (mL) Dilution Factor: 10
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 141062 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.76	D	0.025	0.0075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	2.4	D	0.040	0.013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.19	D	0.025	0.0092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	91		25-150
STL00991	13C4 PFOS	87		25-150
STL00994	18O2 PFHxS	83		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_006.d
 Lims ID: 320-23783-A-3-A
 Client ID: DPT-16-35-GW-21-25
 Sample Type: Client
 Inject. Date: 07-Dec-2016 17:10:55 ALS Bottle#: 19 Worklist Smp#: 6
 Injection Vol: 2.0 ul Dil. Factor: 10.0000
 Sample Info: 320-23783-a-3-a 10X
 Misc. Info.: Plate: 1 Rack: 3
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:54 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:54:57

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.848	1.848	0.0	1.000	3907991	9.41				
298.90 > 99.00	1.848	1.848	0.0	1.000	1604262		2.44(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.457	2.460	-0.003		1230623	3.94		8.3	107443	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.803	2.802	0.001	1.000	8183022	38.2			208567	
413.00 > 169.00	2.803	2.802	0.001	1.000	5434289		1.51(0.90-1.10)		201585	
D 14 13C4 PFOA										
417.00 > 372.00	2.803	2.802	0.001		998595	4.55		9.1	120423	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.170	3.061	0.109	1.000	27465853	118.2			0.0	
499.00 > 99.00	3.062	3.061	0.001	0.966	5756275		4.77(0.90-1.10)		40061	
D 17 13C4 PFOS										
503.00 > 80.00	3.170	3.169	0.001		1023919	4.16		8.7	16170	

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_006.d

Injection Date: 07-Dec-2016 17:10:55

Instrument ID: A8_N

Lims ID: 320-23783-A-3-A

Lab Sample ID: 320-23783-3

Client ID: DPT-16-35-GW-21-25

Operator ID: A8-PC\A8

ALS Bottle#: 19

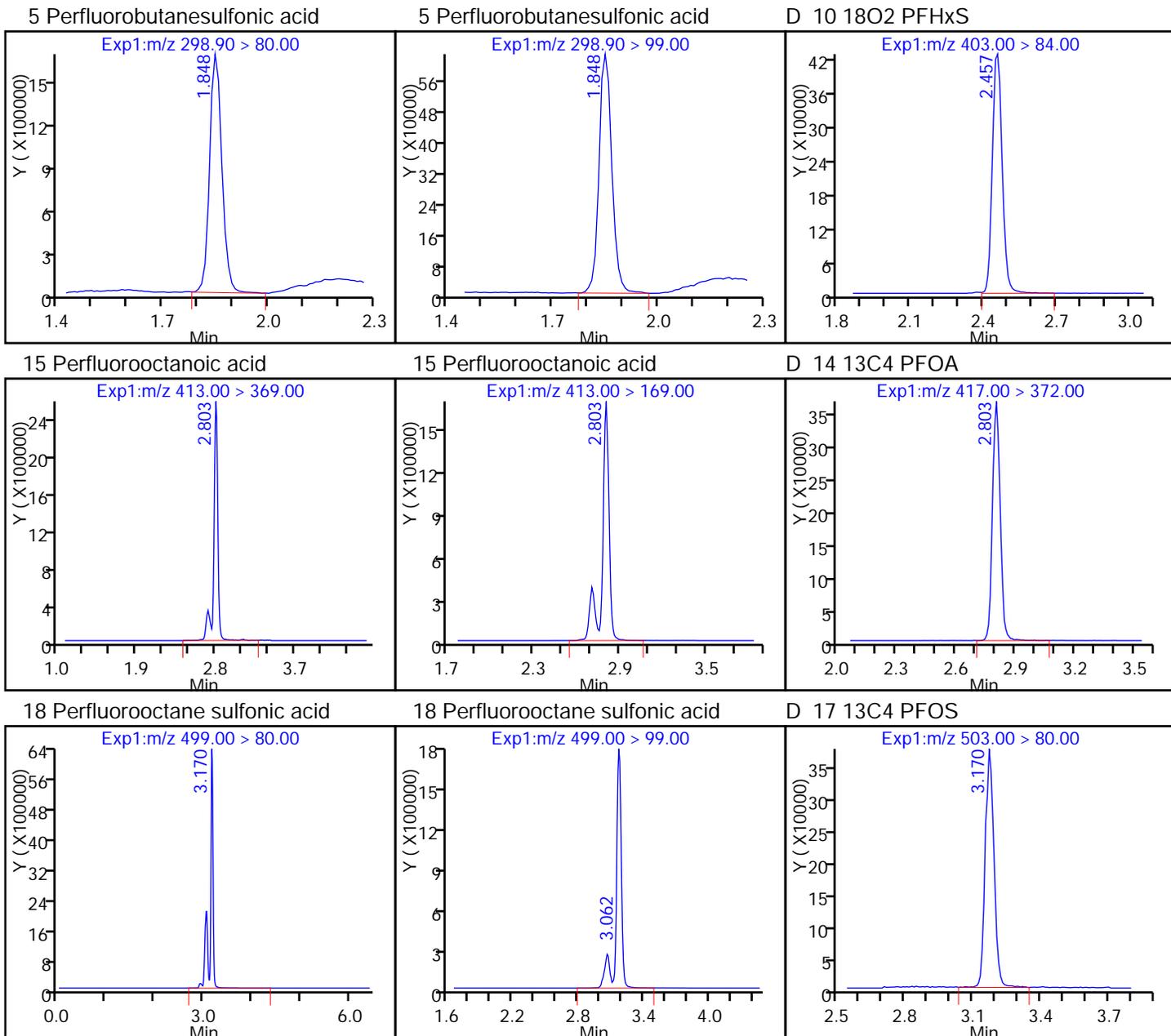
Worklist Smp#: 6

Injection Vol: 2.0 ul

Dil. Factor: 10.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-GW-31-35 Lab Sample ID: 320-23783-4
 Matrix: Water Lab File ID: 03DEC2016B_005.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 10:30
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 249.6(mL) Date Analyzed: 12/03/2016 16:33
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	<i>Perfluorooctanoic acid (PFOA)</i>	1.1	E M	0.0025	0.00075
1763-23-1	<i>Perfluorooctane Sulfonate (PFOS)</i>	0.72	E	0.0040	0.0013
375-73-5	<i>Perfluorobutanesulfonic acid (PFBS)</i>	0.58	E	0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	52		25-150
STL00991	13C4 PFOS	76		25-150
STL00994	18O2 PFHxS	30		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_005.d
 Lims ID: 320-23783-A-4-A
 Client ID: DPT-16-35-GW-31-35
 Sample Type: Client
 Inject. Date: 03-Dec-2016 16:33:42 ALS Bottle#: 36 Worklist Smp#: 26
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-4-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:34:03 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:34:03

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.878	1.900	-0.022	1.000	42955471	291.6				E
298.90 > 99.00	1.878	1.900	-0.022	1.000	21977968		1.95(0.00-0.00)			E
D 10 18O2 PFHxS										
403.00 > 84.00	2.496	2.531	-0.035		4364938	14.0		29.5	61505	
D 14 13C4 PFOA										
417.00 > 372.00	2.846	2.880	-0.034		5746936	26.2		52.4	436796	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.846	2.887	-0.041	1.000	65382761	530.7			276814	EM
413.00 > 169.00	2.846	2.887	-0.041	1.000	52263596		1.25(0.90-1.10)		154214	M
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.220	3.258	-0.038	1.000	72785937	358.1			2044570	E
499.00 > 99.00	3.220	3.258	-0.038	1.000	16679335		4.36(0.90-1.10)		2131667	
D 17 13C4 PFOS										
503.00 > 80.00	3.220	3.259	-0.039		8953037	36.4		76.1	114840	

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_005.d

Injection Date: 03-Dec-2016 16:33:42

Instrument ID: A8_N

Lims ID: 320-23783-A-4-A

Lab Sample ID: 320-23783-4

Client ID: DPT-16-35-GW-31-35

Operator ID: A8-PC\A8

ALS Bottle#: 36

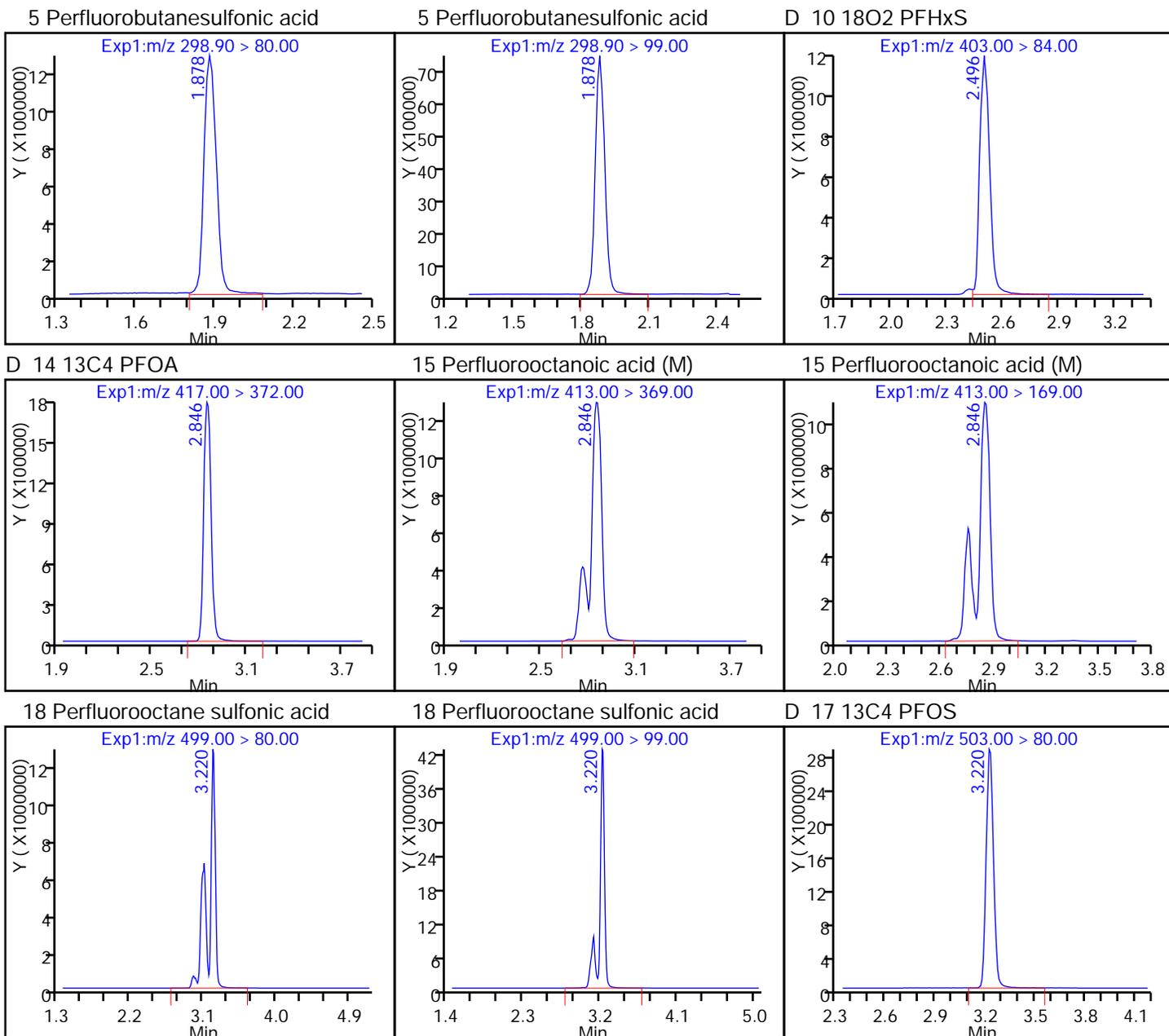
Worklist Smp#: 26

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

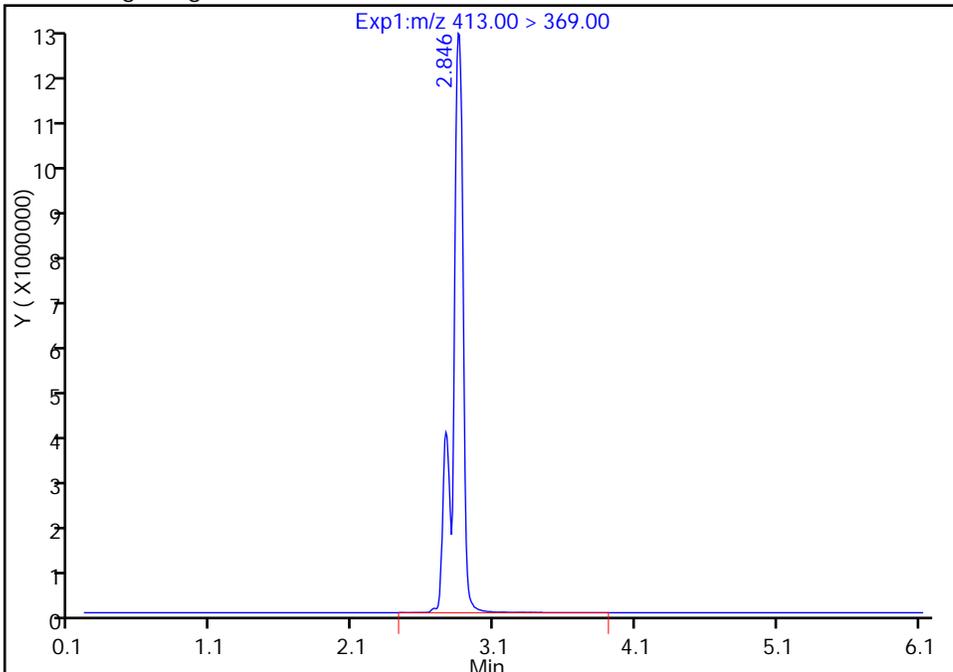
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_005.d
Injection Date: 03-Dec-2016 16:33:42 Instrument ID: A8_N
Lims ID: 320-23783-A-4-A Lab Sample ID: 320-23783-4
Client ID: DPT-16-35-GW-31-35
Operator ID: A8-PC\A8 ALS Bottle#: 36 Worklist Smp#: 26
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

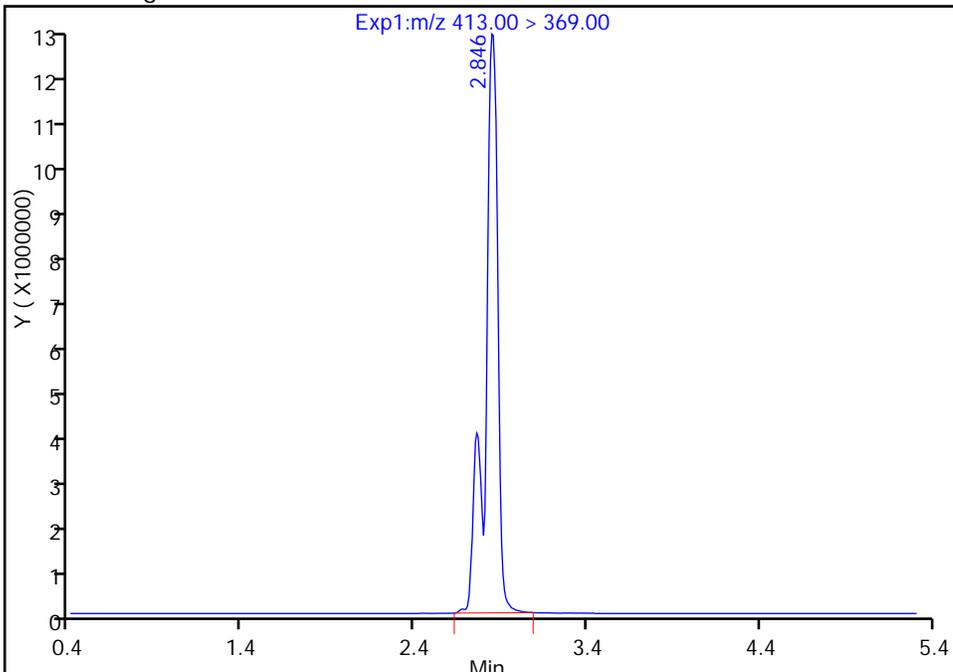
RT: 2.85
Area: 66001828
Amount: 535.7093
Amount Units: ng/ml

Processing Integration Results



RT: 2.85
Area: 65382761
Amount: 530.6846
Amount Units: ng/ml

Manual Integration Results



TestAmerica Sacramento

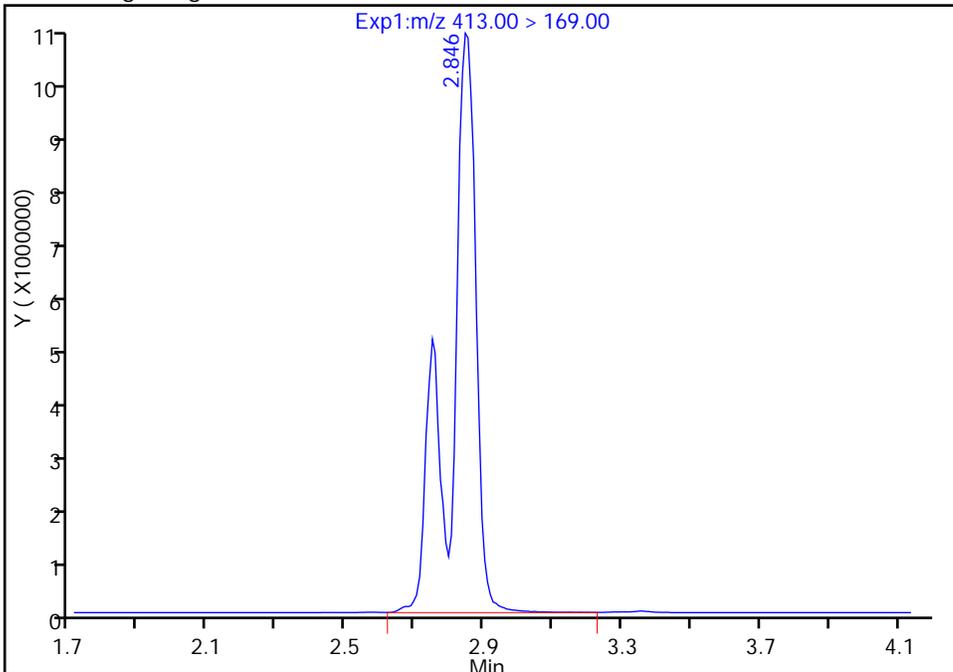
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_005.d
Injection Date: 03-Dec-2016 16:33:42 Instrument ID: A8_N
Lims ID: 320-23783-A-4-A Lab Sample ID: 320-23783-4
Client ID: DPT-16-35-GW-31-35
Operator ID: A8-PC\A8 ALS Bottle#: 36 Worklist Smp#: 26
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 2

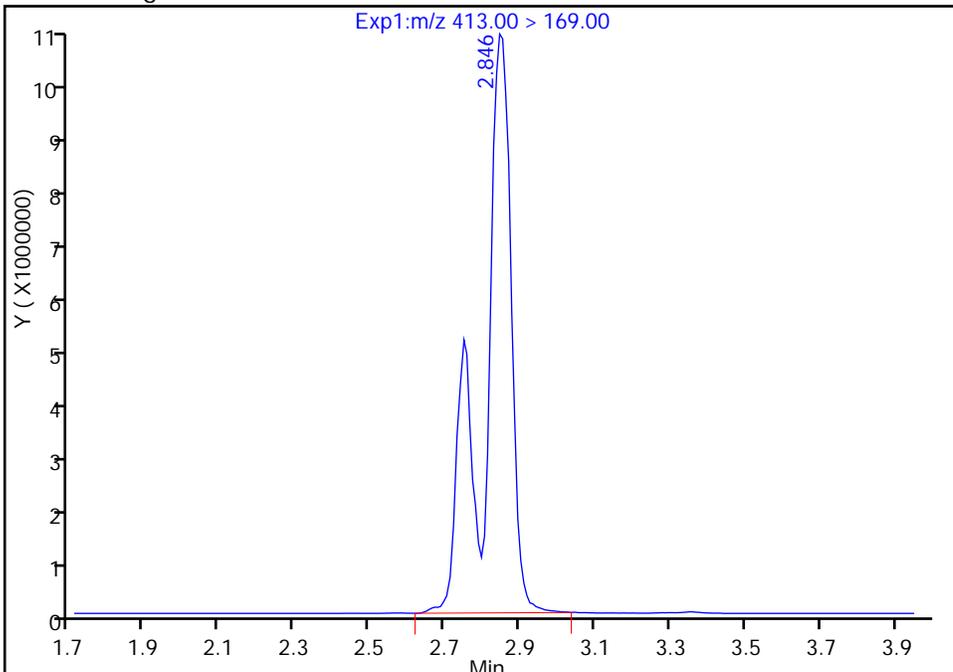
RT: 2.85
Area: 52579475
Amount: 535.7093
Amount Units: ng/ml

Processing Integration Results



RT: 2.85
Area: 52263596
Amount: 530.6846
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-35-GW-31-35 DL Lab Sample ID: 320-23783-4 DL
 Matrix: Water Lab File ID: 07DEC2016B_007.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 10:30
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 249.6(mL) Date Analyzed: 12/07/2016 17:18
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 10
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 141062 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	1.4	D	0.025	0.0075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.80	D	0.040	0.013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.38	D	0.025	0.0092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	78		25-150
STL00991	13C4 PFOS	91		25-150
STL00994	18O2 PFHxS	73		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_007.d
 Lims ID: 320-23783-A-4-A
 Client ID: DPT-16-35-GW-31-35
 Sample Type: Client
 Inject. Date: 07-Dec-2016 17:18:25 ALS Bottle#: 20 Worklist Smp#: 7
 Injection Vol: 2.0 ul Dil. Factor: 10.0000
 Sample Info: 320-23783-a-4-a 10X
 Misc. Info.: Plate: 1 Rack: 3
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:54 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:55:07

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.848	1.848	0.0	1.000	6973833	19.1				
298.90 > 99.00	1.848	1.848	0.0	1.000	3038662		2.30(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.455	2.460	-0.005		1083716	3.47		7.3	88167	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.804	2.802	0.002	1.000	13051856	71.3			290557	
413.00 > 169.00	2.804	2.802	0.002	1.000	8917503		1.46(0.90-1.10)		215086	
D 14 13C4 PFOA										
417.00 > 372.00	2.804	2.802	0.002		853859	3.89		7.8	100091	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.179	3.061	0.118	1.000	9766998	40.1			458205	
499.00 > 99.00	3.171	3.061	0.110	0.997	1917058		5.09(0.90-1.10)		144974	
D 17 13C4 PFOS										
503.00 > 80.00	3.171	3.169	0.002		1073837	4.37		9.1	40552	

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_007.d

Injection Date: 07-Dec-2016 17:18:25

Instrument ID: A8_N

Lims ID: 320-23783-A-4-A

Lab Sample ID: 320-23783-4

Client ID: DPT-16-35-GW-31-35

Operator ID: A8-PC\A8

ALS Bottle#: 20

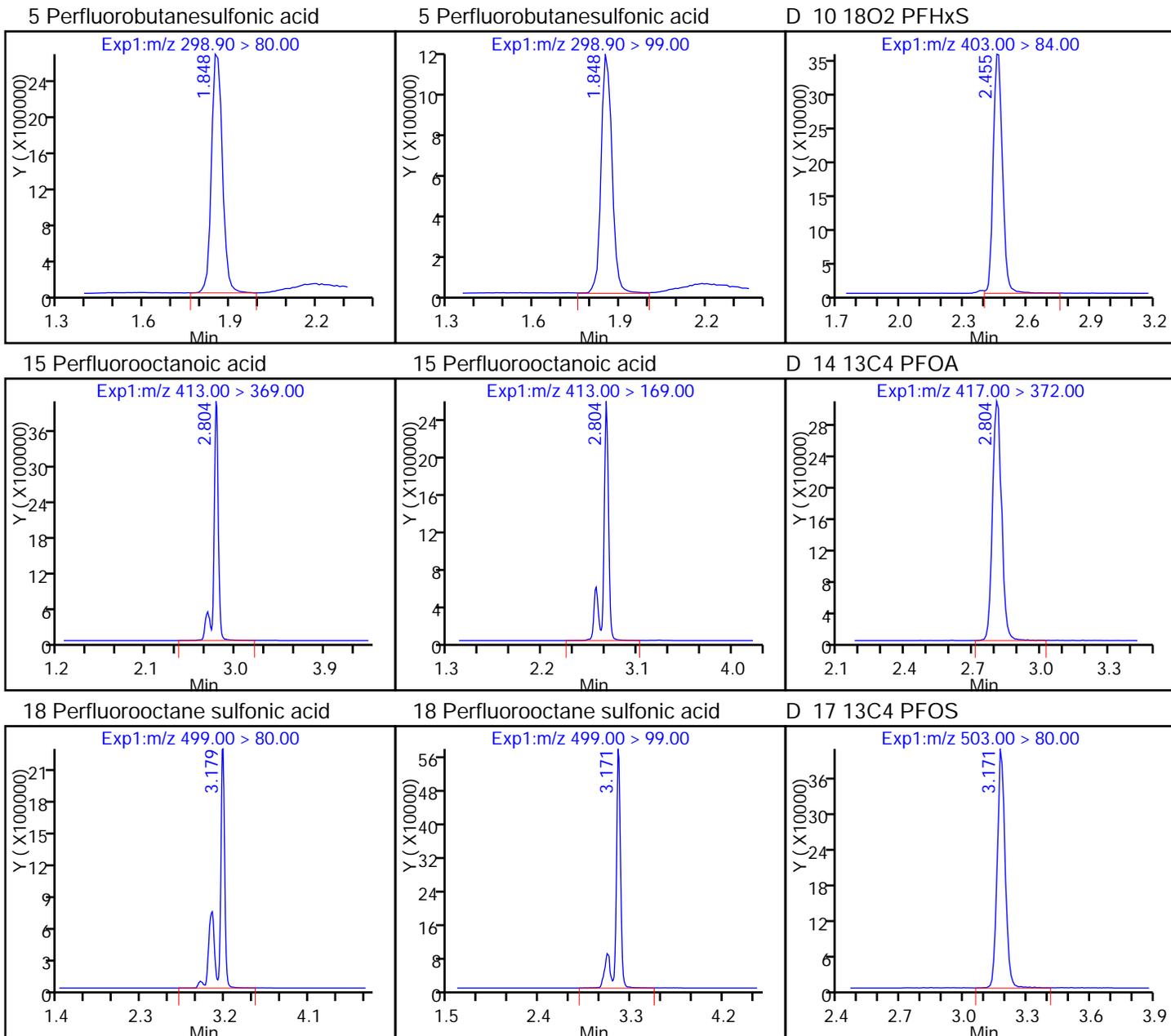
Worklist Smp#: 7

Injection Vol: 2.0 ul

Dil. Factor: 10.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-19-SO-00-01 Lab Sample ID: 320-23783-5
 Matrix: Solid Lab File ID: 02DEC2016C_010.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 11:20
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.02(g) Date Analyzed: 12/02/2016 16:07
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: 13.3 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	3.8		0.57	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	130	M E	0.57	0.14
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.18	J	0.46	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	96		25-150
STL00991	13C4 PFOS	54		25-150
STL00994	18O2 PFHxS	85		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_010.d
 Lims ID: 320-23783-A-5-A
 Client ID: DPT-16-19-SO-00-01
 Sample Type: Client
 Inject. Date: 02-Dec-2016 16:07:10 ALS Bottle#: 10 Worklist Smp#: 14
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-5-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:25:25 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:25:16

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.900	1.907	-0.007	1.000	322239	0.7750				
298.90 > 99.00	1.900	1.907	-0.007	1.000	154730		2.08(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.528	2.534	-0.006		12188867	40.3		85.2	600398	
D 14 13C4 PFOA										
417.00 > 372.00	2.887	2.886	0.0		10568981	48.2		96.3	1059712	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.887	2.894	-0.008	1.000	3598653	16.7			37425	
413.00 > 169.00	2.887	2.894	-0.008	1.000	2481983		1.45(0.90-1.10)		115185	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.268	3.267	0.001	1.000	82856204	581.8			366942	EM
499.00 > 99.00	3.268	3.267	0.001	1.000	22573580		3.67(0.90-1.10)		245829	M
D 17 13C4 PFOS										
503.00 > 80.00	3.261	3.267	-0.006		6260717	25.6		53.5	208105	

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_010.d

Injection Date: 02-Dec-2016 16:07:10

Instrument ID: A8_N

Lims ID: 320-23783-A-5-A

Lab Sample ID: 320-23783-5

Client ID: DPT-16-19-SO-00-01

Operator ID: A8-PC\A8

ALS Bottle#: 10

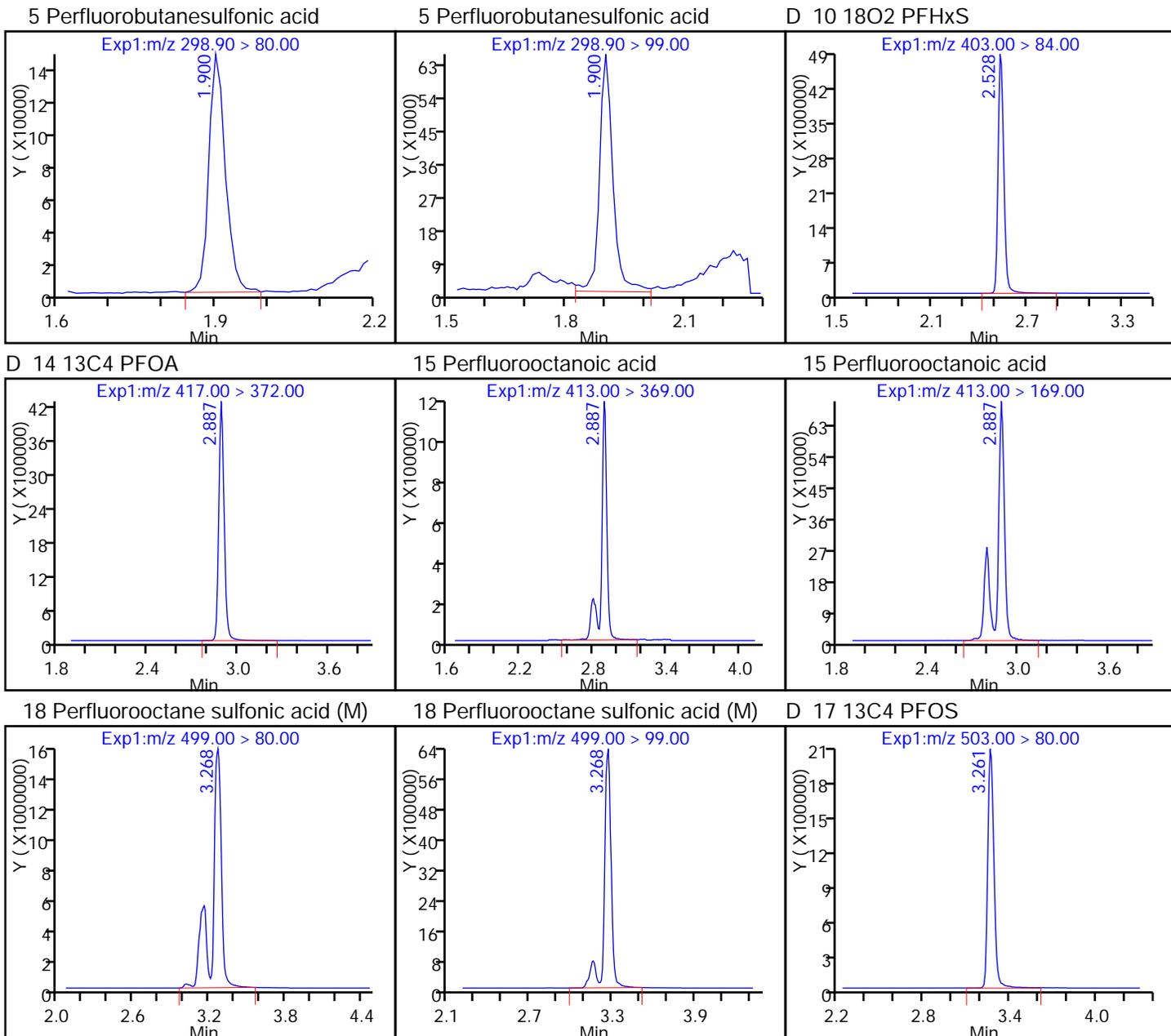
Worklist Smp#: 14

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

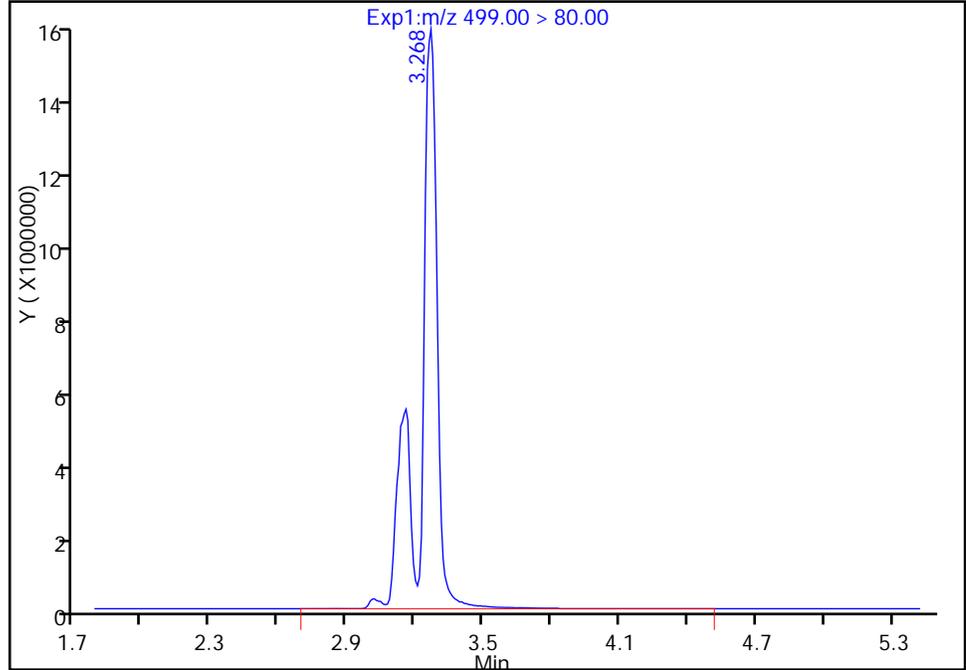
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_010.d
Injection Date: 02-Dec-2016 16:07:10 Instrument ID: A8_N
Lims ID: 320-23783-A-5-A Lab Sample ID: 320-23783-5
Client ID: DPT-16-19-SO-00-01
Operator ID: A8-PC\A8 ALS Bottle#: 10 Worklist Smp#: 14
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

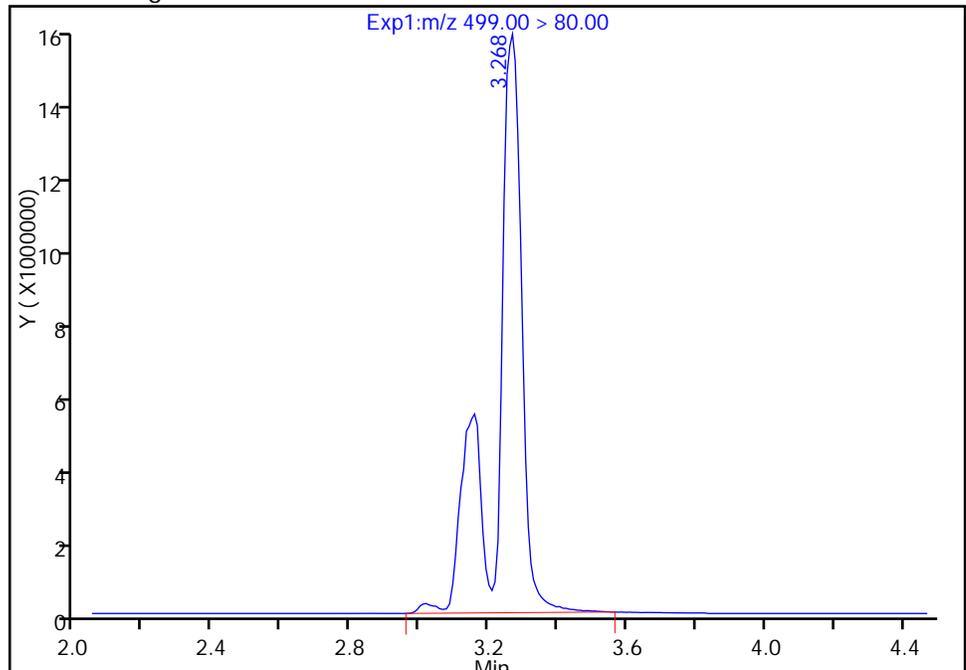
RT: 3.27
Area: 84017279
Amount: 589.9377
Amount Units: ng/ml

Processing Integration Results



RT: 3.27
Area: 82856204
Amount: 581.7851
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:25:16
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

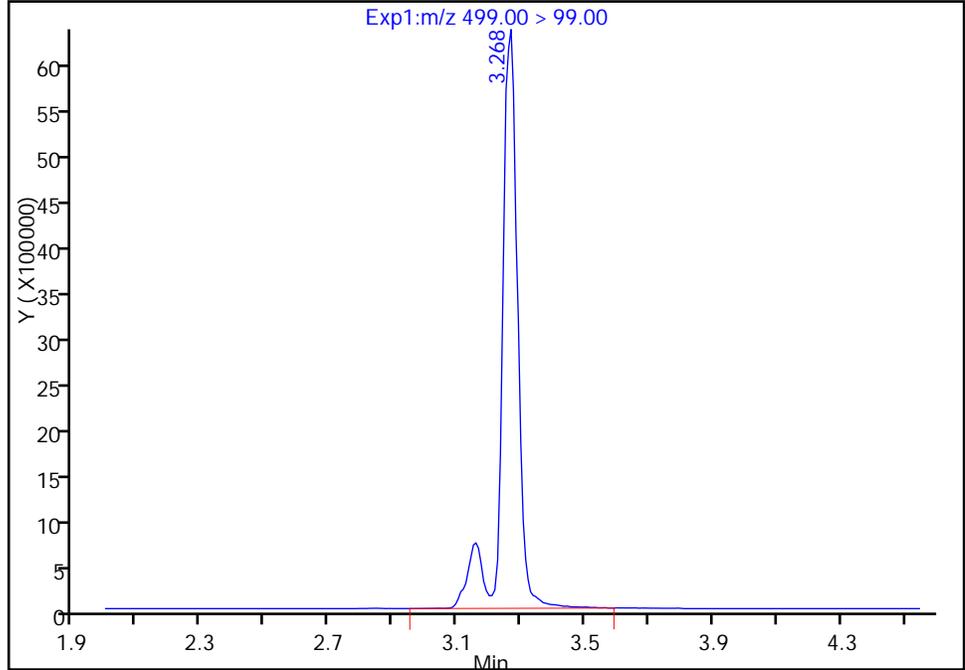
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_010.d
Injection Date: 02-Dec-2016 16:07:10 Instrument ID: A8_N
Lims ID: 320-23783-A-5-A Lab Sample ID: 320-23783-5
Client ID: DPT-16-19-SO-00-01
Operator ID: A8-PC\A8 ALS Bottle#: 10 Worklist Smp#: 14
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

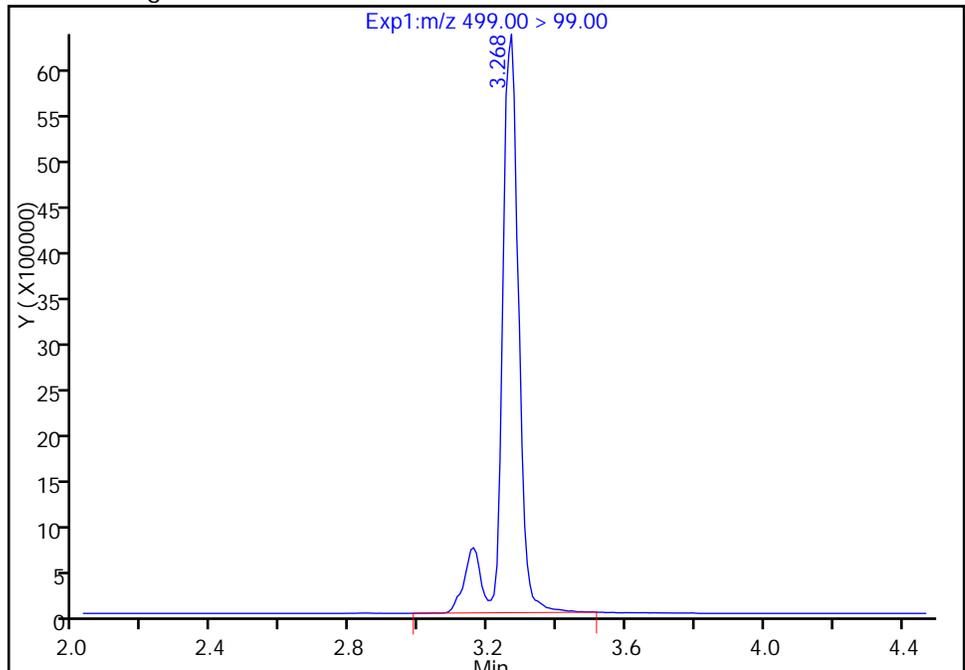
RT: 3.27
Area: 22733991
Amount: 589.9377
Amount Units: ng/ml

Processing Integration Results



RT: 3.27
Area: 22573580
Amount: 581.7851
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-19-SO-00-01 DL Lab Sample ID: 320-23783-5 DL
 Matrix: Solid Lab File ID: 07DEC2016B_009.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 11:20
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.02(g) Date Analyzed: 12/07/2016 17:33
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 10
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 13.3 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 141062 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	<i>Perfluorooctanoic acid (PFOA)</i>	3.5	J D M	5.7	1.2
1763-23-1	Perfluorooctane Sulfonate (PFOS)	170	D	5.7	1.4
375-73-5	<i>Perfluorobutanesulfonic acid (PFBS)</i>	3.4	U	4.6	1.2

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	136		25-150
STL00991	13C4 PFOS	90		25-150
STL00994	18O2 PFHxS	106		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_009.d
 Lims ID: 320-23783-A-5-A
 Client ID: DPT-16-19-SO-00-01
 Sample Type: Client
 Inject. Date: 07-Dec-2016 17:33:26 ALS Bottle#: 22 Worklist Smp#: 9
 Injection Vol: 2.0 ul Dil. Factor: 10.0000
 Sample Info: 320-23783-a-5-a 10X
 Misc. Info.: Plate: 1 Rack: 3
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:54 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:55:45

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.849	1.848	0.001	1.000	36343	0.0690				
298.90 > 99.00	1.839	1.848	-0.009	0.995	14796		2.46(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.448	2.460	-0.012		1561968	5.00		10.6	195821	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.800	2.802	-0.002	1.000	487830	1.53			4567	M
413.00 > 169.00	2.776	2.802	-0.026	0.991	333150		1.46(0.90-1.10)		5791	M
D 14 13C4 PFOA										
417.00 > 372.00	2.800	2.802	-0.002		1490656	6.80		13.6	147470	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.169	3.061	0.108	1.000	17658015	73.6			19687	
499.00 > 99.00	3.169	3.061	0.108	1.000	3944310		4.48(0.90-1.10)		337659	
D 17 13C4 PFOS										
503.00 > 80.00	3.169	3.169	0.0		1056967	4.30		9.0	64846	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_009.d

Injection Date: 07-Dec-2016 17:33:26

Instrument ID: A8_N

Lims ID: 320-23783-A-5-A

Lab Sample ID: 320-23783-5

Client ID: DPT-16-19-SO-00-01

Operator ID: A8-PC\A8

ALS Bottle#: 22

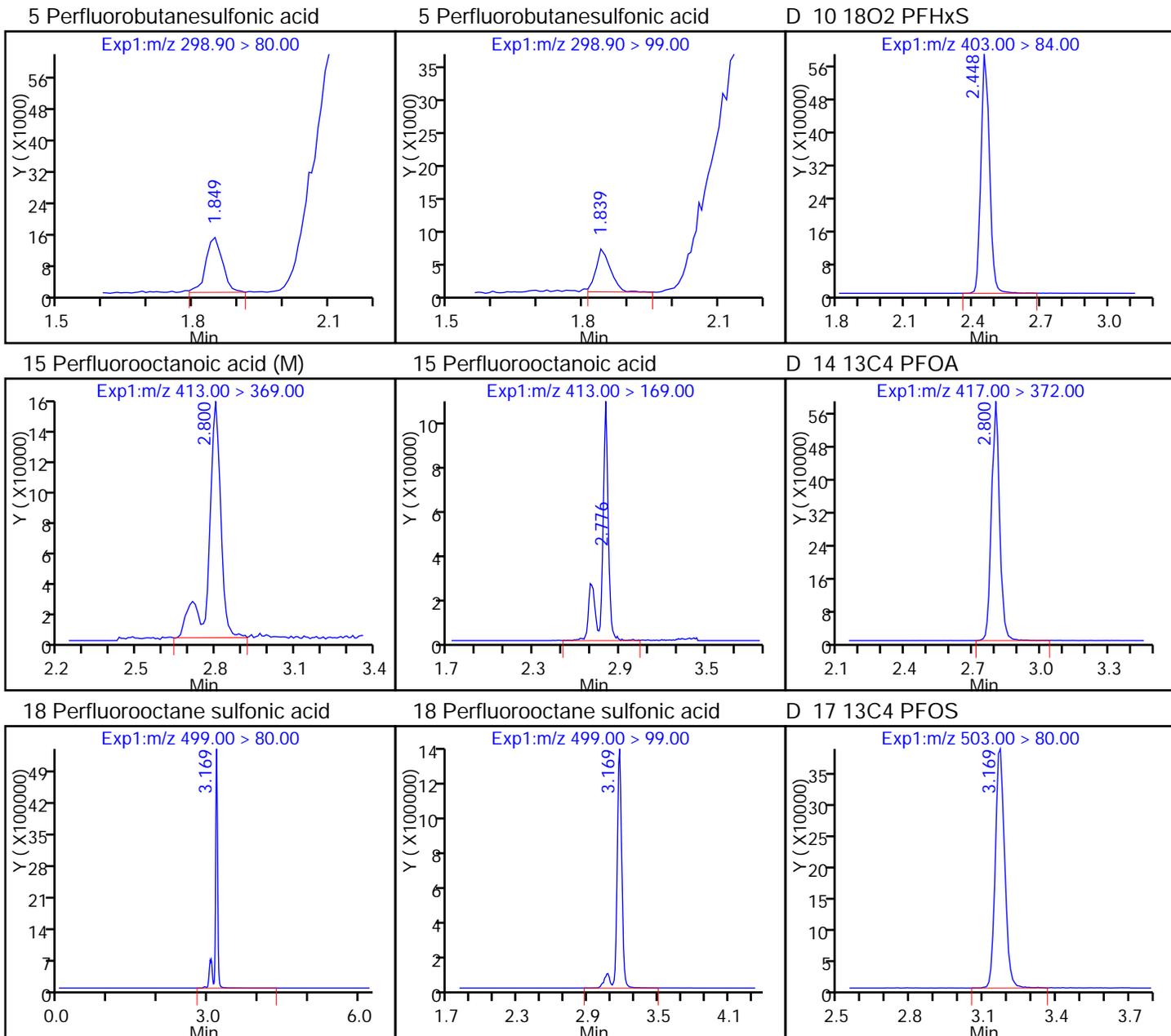
Worklist Smp#: 9

Injection Vol: 2.0 ul

Dil. Factor: 10.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

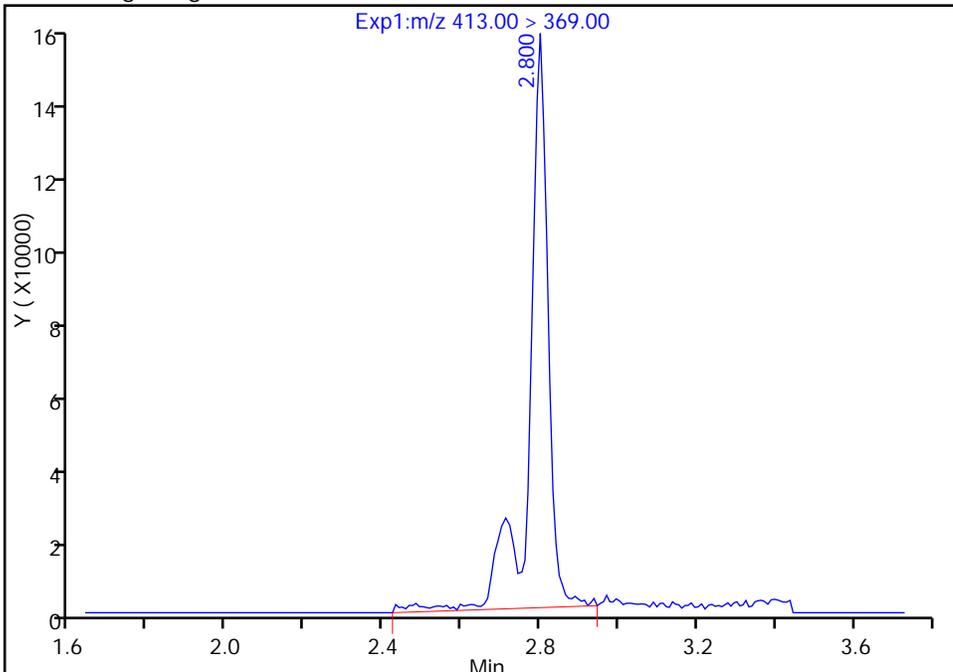
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_009.d
Injection Date: 07-Dec-2016 17:33:26 Instrument ID: A8_N
Lims ID: 320-23783-A-5-A Lab Sample ID: 320-23783-5
Client ID: DPT-16-19-SO-00-01
Operator ID: A8-PC\A8 ALS Bottle#: 22 Worklist Smp#: 9
Injection Vol: 2.0 ul Dil. Factor: 10.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

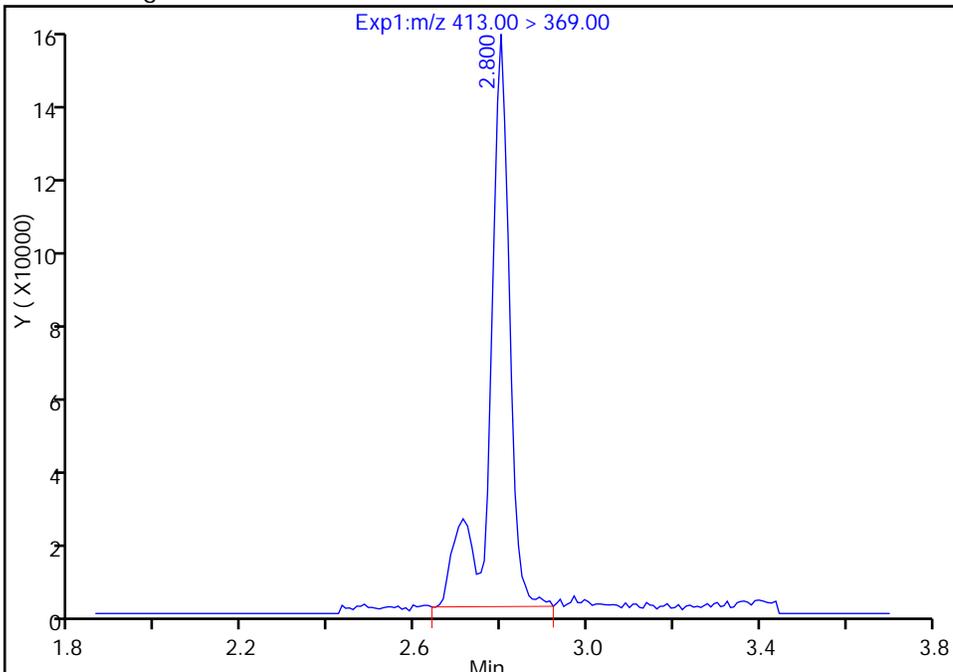
RT: 2.80
Area: 514676
Amount: 1.610520
Amount Units: ng/ml

Processing Integration Results



RT: 2.80
Area: 487830
Amount: 1.526514
Amount Units: ng/ml

Manual Integration Results



Reviewer: westendorfc, 08-Dec-2016 08:55:45

Audit Action: Manually Integrated

Audit Reason: Baseline

FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-19-SO-17-18 Lab Sample ID: 320-23783-6
 Matrix: Solid Lab File ID: 02DEC2016C_015.d
 Analysis Method: 537 (Modified) Date Collected: 11/22/2016 11:30
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.06(g) Date Analyzed: 12/02/2016 16:44
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 20.3 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.37	J	0.62	0.13
1763-23-1	Perfluorooctane Sulfonate (PFOS)	6.5	M	0.62	0.16
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.50	0.13

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	69		25-150
STL00991	13C4 PFOS	64		25-150
STL00994	18O2 PFHxS	73		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_015.d
 Lims ID: 320-23783-A-6-A
 Client ID: DPT-16-19-SO-17-18
 Sample Type: Client
 Inject. Date: 02-Dec-2016 16:44:43 ALS Bottle#: 11 Worklist Smp#: 19
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-6-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:32:28 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:27:06

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.900	1.907	-0.007	1.000	37817	0.1060				
298.90 > 99.00	1.890	1.907	-0.017	0.995	17034		2.22(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.529	2.534	-0.005		10457609	34.6		73.1	1389657	
D 14 13C4 PFOA										
417.00 > 372.00	2.878	2.886	-0.008		7593353	34.6		69.2	606548	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.886	2.894	-0.008	1.000	233612	1.51			3299	
413.00 > 169.00	2.886	2.894	-0.008	1.000	151520		1.54(0.90-1.10)		6224	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.259	3.267	-0.008	1.000	4405095	26.1			1503701M	
499.00 > 99.00	3.259	3.267	-0.008	1.000	989788		4.45(0.90-1.10)		13953	M
D 17 13C4 PFOS										
503.00 > 80.00	3.259	3.267	-0.008		7431648	30.4		63.6	309964	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_015.d

Injection Date: 02-Dec-2016 16:44:43

Instrument ID: A8_N

Lims ID: 320-23783-A-6-A

Lab Sample ID: 320-23783-6

Client ID: DPT-16-19-SO-17-18

Operator ID: A8-PC\A8

ALS Bottle#: 11

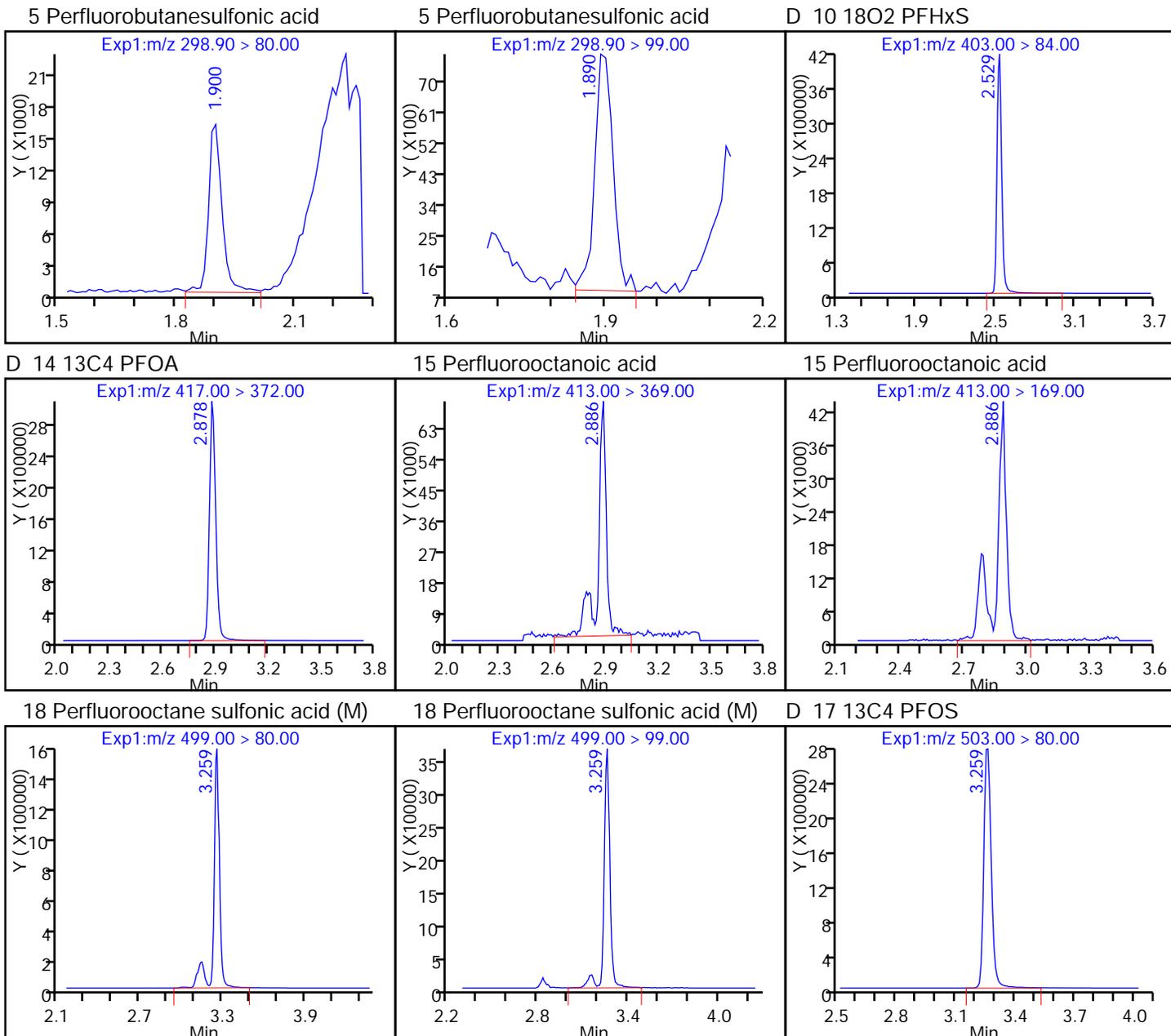
Worklist Smp#: 19

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

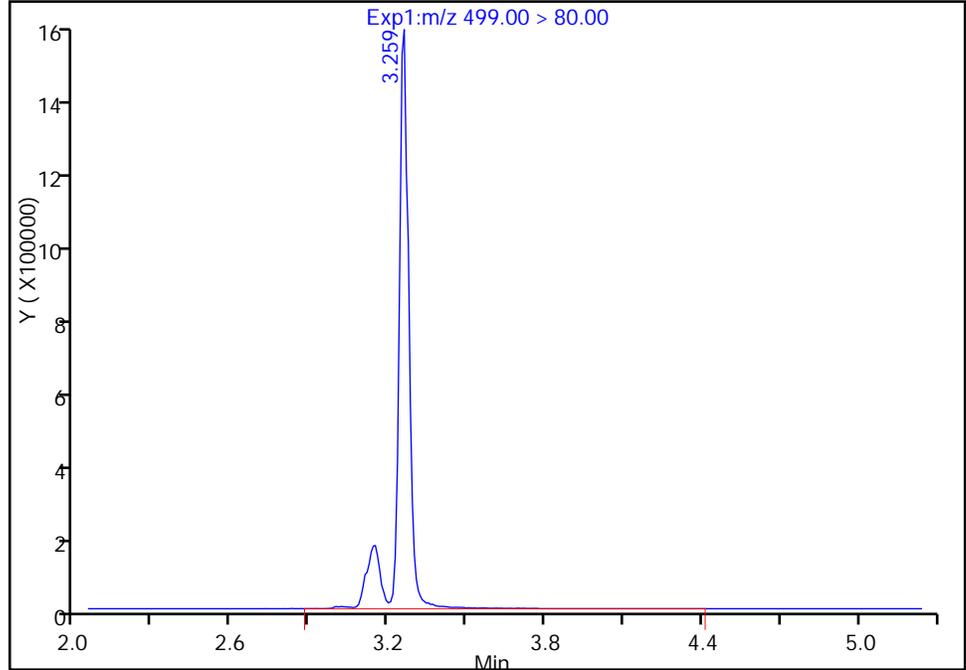
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_015.d
Injection Date: 02-Dec-2016 16:44:43 Instrument ID: A8_N
Lims ID: 320-23783-A-6-A Lab Sample ID: 320-23783-6
Client ID: DPT-16-19-SO-17-18
Operator ID: A8-PC\A8 ALS Bottle#: 11 Worklist Smp#: 19
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

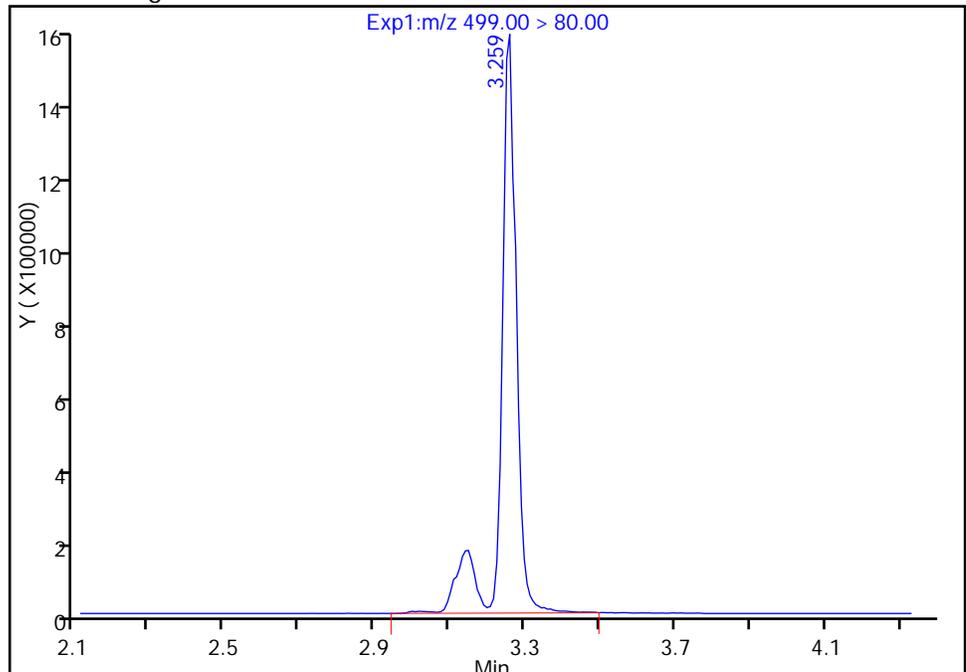
RT: 3.26
Area: 4460025
Amount: 26.382369
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 4405095
Amount: 26.057442
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:27:06
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

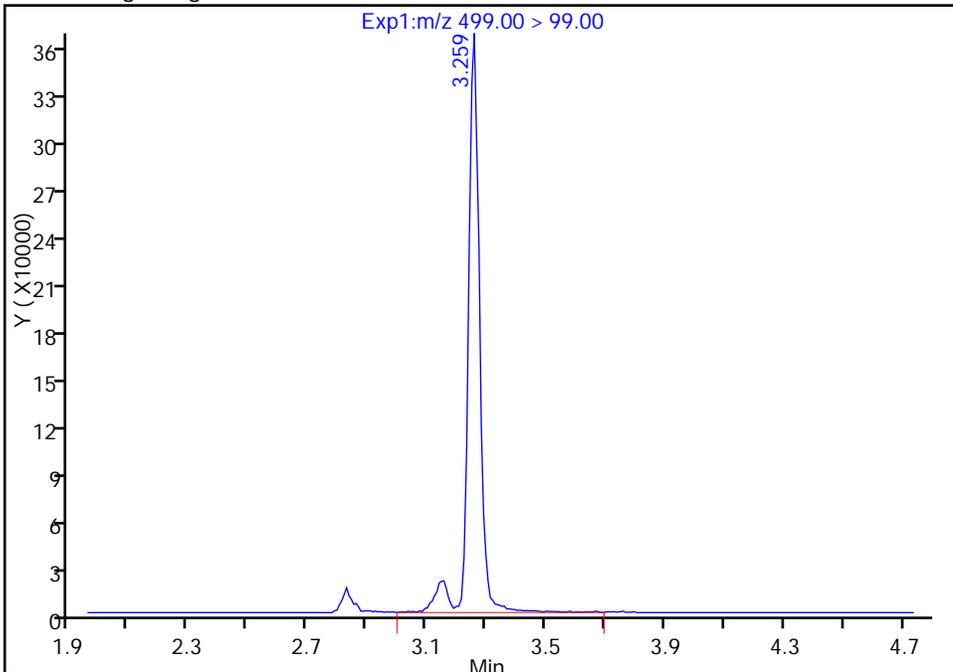
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_015.d
Injection Date: 02-Dec-2016 16:44:43 Instrument ID: A8_N
Lims ID: 320-23783-A-6-A Lab Sample ID: 320-23783-6
Client ID: DPT-16-19-SO-17-18
Operator ID: A8-PC\A8 ALS Bottle#: 11 Worklist Smp#: 19
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

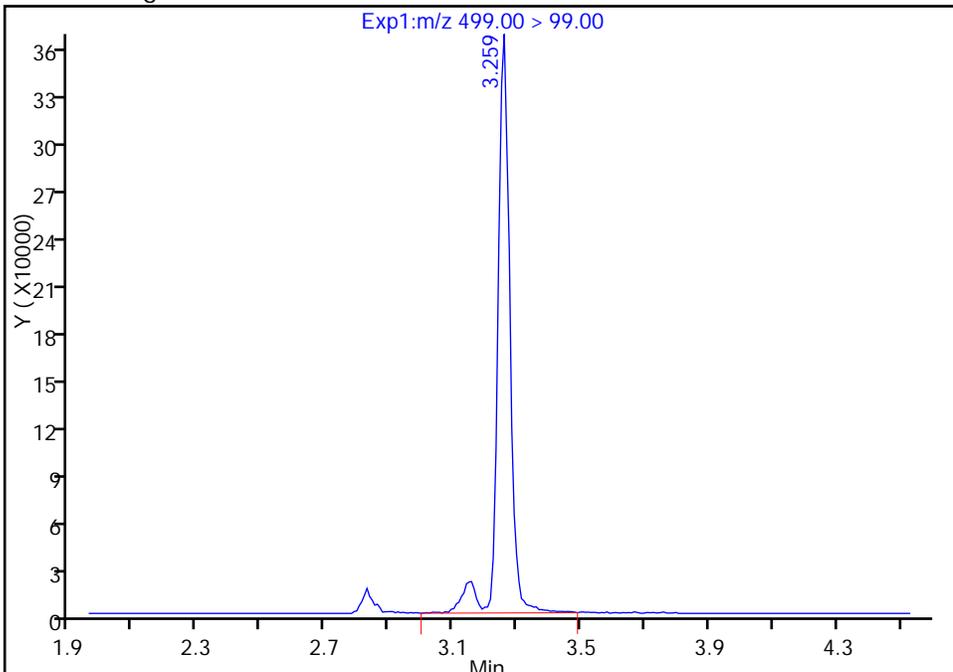
RT: 3.26
Area: 1001013
Amount: 26.382369
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 989788
Amount: 26.057442
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-32-SO-14-15 Lab Sample ID: 320-23783-7
 Matrix: Solid Lab File ID: 02DEC2016C_016.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 09:36
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.03(g) Date Analyzed: 12/02/2016 16:52
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 17.7 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	1.1		0.60	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	46	M	0.60	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	96		25-150
STL00991	13C4 PFOS	62		25-150
STL00994	18O2 PFHxS	70		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_016.d
 Lims ID: 320-23783-A-7-A
 Client ID: DPT-16-32-SO-14-15
 Sample Type: Client
 Inject. Date: 02-Dec-2016 16:52:14 ALS Bottle#: 12 Worklist Smp#: 20
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-7-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:28:23 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:28:23

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.897	1.907	-0.010	1.000	112764	0.3307				
298.90 > 99.00	1.897	1.907	-0.010	1.000	49175		2.29(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.523	2.534	-0.012		9995081	33.0		69.9	312906	
D 14 13C4 PFOA										
417.00 > 372.00	2.879	2.886	-0.007		10547358	48.1		96.1	1344938	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.879	2.894	-0.015	1.000	968867	4.50			18515	
413.00 > 169.00	2.879	2.894	-0.015	1.000	619103		1.56(0.90-1.10)		36207	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.254	3.267	-0.013	1.000	31467926	190.2			148739	M
499.00 > 99.00	3.262	3.267	-0.005	1.003	8099323		3.89(0.90-1.10)		115582	M
D 17 13C4 PFOS										
503.00 > 80.00	3.262	3.267	-0.005		7274549	29.7		62.2	237038	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_016.d

Injection Date: 02-Dec-2016 16:52:14

Instrument ID: A8_N

Lims ID: 320-23783-A-7-A

Lab Sample ID: 320-23783-7

Client ID: DPT-16-32-SO-14-15

Operator ID: A8-PC\A8

ALS Bottle#: 12

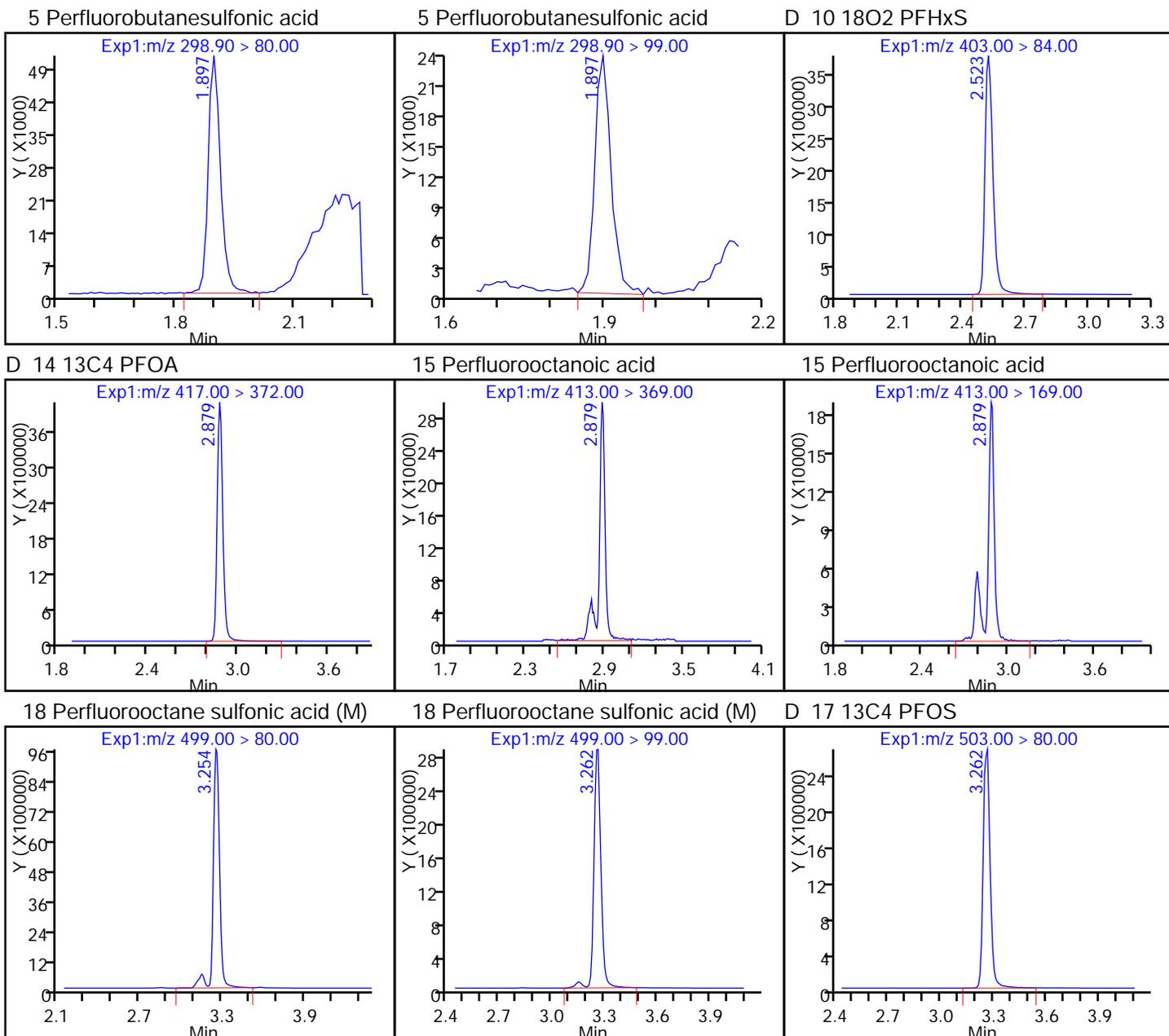
Worklist Smp#: 20

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

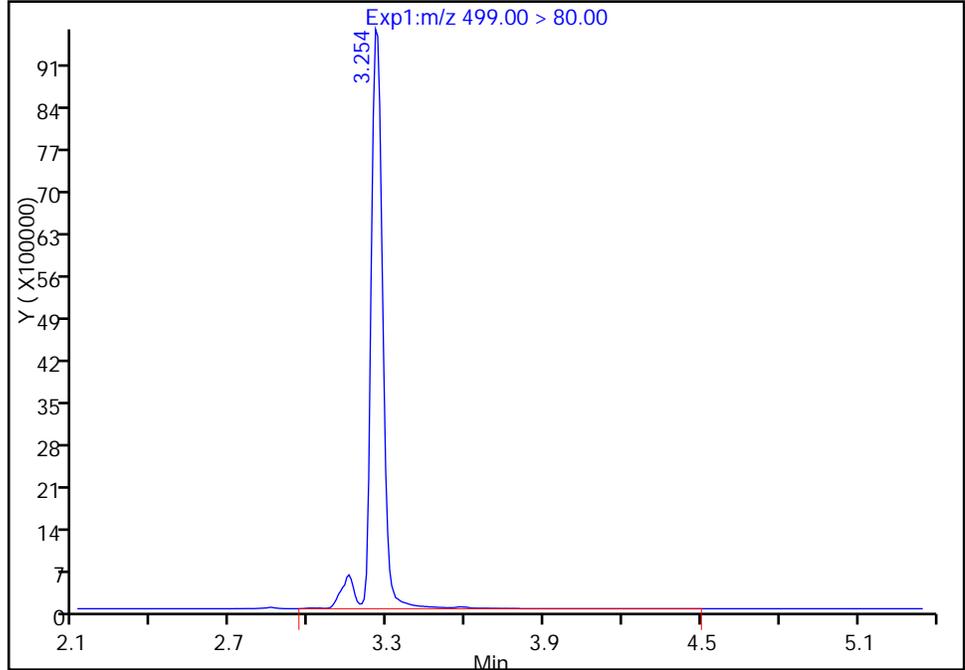
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_016.d
Injection Date: 02-Dec-2016 16:52:14 Instrument ID: A8_N
Lims ID: 320-23783-A-7-A Lab Sample ID: 320-23783-7
Client ID: DPT-16-32-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 12 Worklist Smp#: 20
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

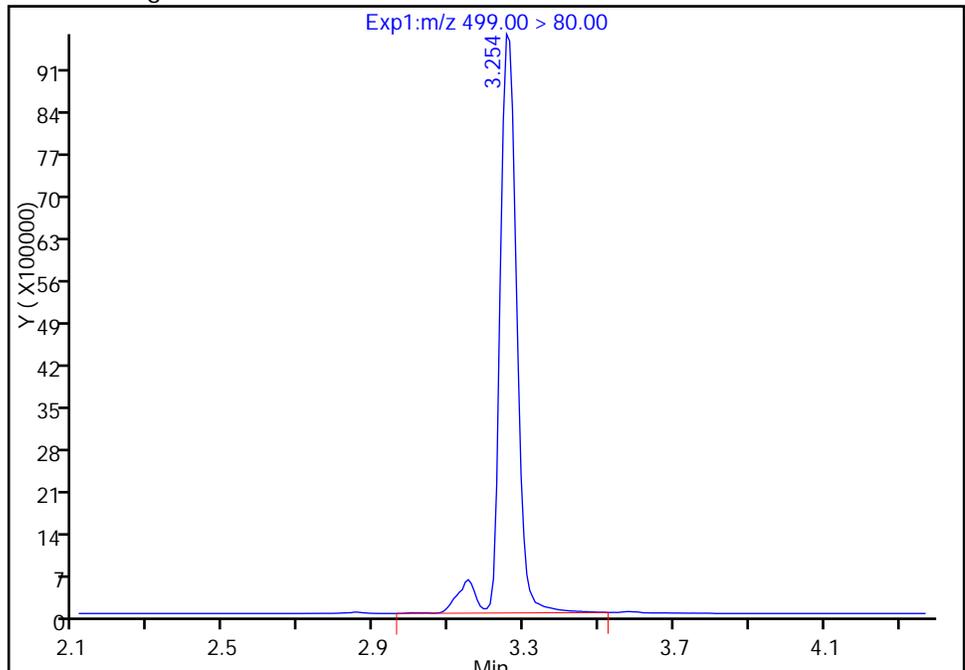
RT: 3.25
Area: 31952647
Amount: 193.0912
Amount Units: ng/ml

Processing Integration Results



RT: 3.25
Area: 31467926
Amount: 190.1620
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:28:23
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

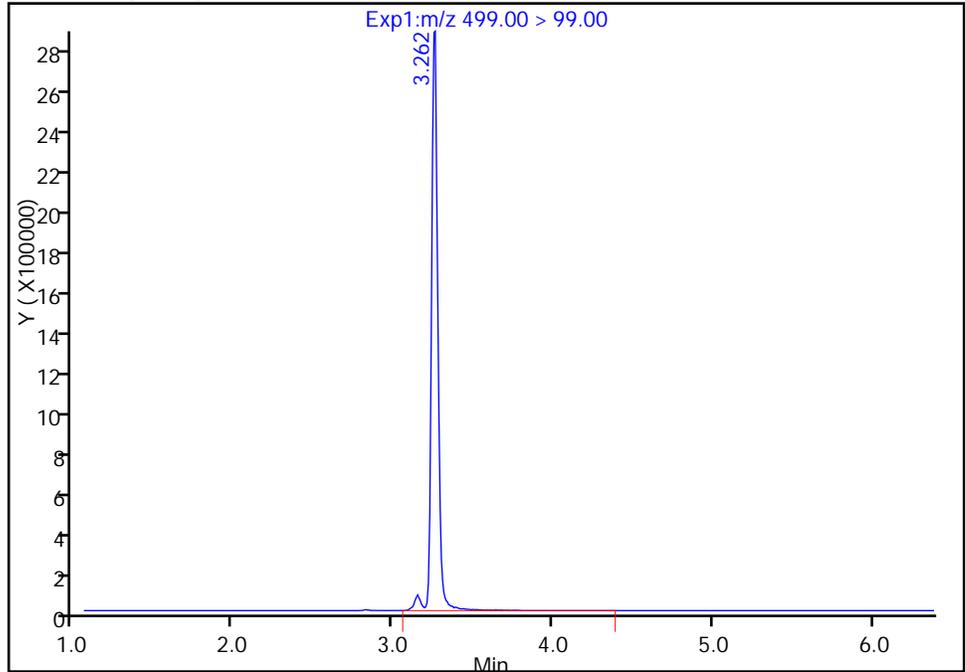
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_016.d
Injection Date: 02-Dec-2016 16:52:14 Instrument ID: A8_N
Lims ID: 320-23783-A-7-A Lab Sample ID: 320-23783-7
Client ID: DPT-16-32-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 12 Worklist Smp#: 20
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

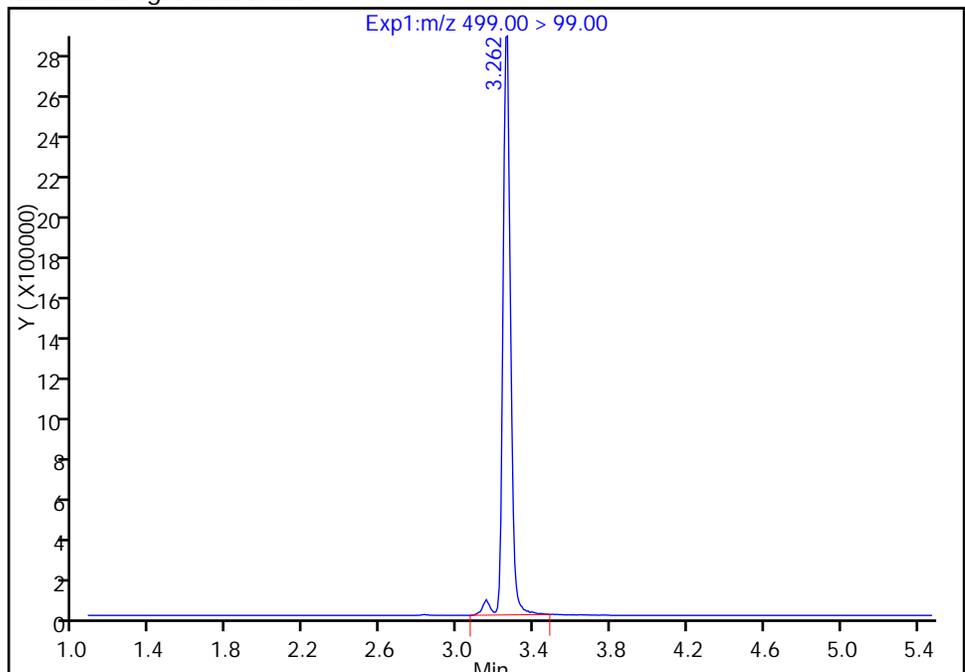
RT: 3.26
Area: 8210883
Amount: 193.0912
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 8099323
Amount: 190.1620
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:28:23

Audit Action: Manually Integrated

Audit Reason: Baseline

FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-31-SO-14-15 Lab Sample ID: 320-23783-8
 Matrix: Solid Lab File ID: 02DEC2016C_017.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 10:12
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.00(g) Date Analyzed: 12/02/2016 16:59
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: 16.2 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.27	J	0.60	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	11	M	0.60	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	108		25-150
STL00991	13C4 PFOS	82		25-150
STL00994	18O2 PFHxS	93		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_017.d
 Lims ID: 320-23783-A-8-A
 Client ID: DPT-16-31-SO-14-15
 Sample Type: Client
 Inject. Date: 02-Dec-2016 16:59:44 ALS Bottle#: 13 Worklist Smp#: 21
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-8-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:29:01 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:29:01

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.906	1.907	-0.001	1.000	12175	0.0270				
298.90 > 99.00	1.906	1.907	-0.001	1.000	6063		2.01(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.529	2.534	-0.005		13238095	43.8		92.5	640197	
D 14 13C4 PFOA										
417.00 > 372.00	2.886	2.886	0.0		11860760	54.1		108	1429539	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.886	2.894	-0.008	1.000	273450	1.13			3347	
413.00 > 169.00	2.886	2.894	-0.008	1.000	180855		1.51(0.90-1.10)		9027	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.259	3.267	-0.008	1.000	10232948	46.9			149415	M
499.00 > 99.00	3.259	3.267	-0.008	1.000	2262395		4.52(0.90-1.10)		70060	M
D 17 13C4 PFOS										
503.00 > 80.00	3.259	3.267	-0.008		9582384	39.2		82.0	477814	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_017.d

Injection Date: 02-Dec-2016 16:59:44

Instrument ID: A8_N

Lims ID: 320-23783-A-8-A

Lab Sample ID: 320-23783-8

Client ID: DPT-16-31-SO-14-15

Operator ID: A8-PC\A8

ALS Bottle#: 13

Worklist Smp#: 21

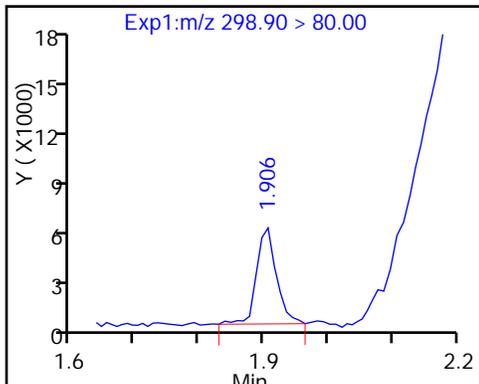
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

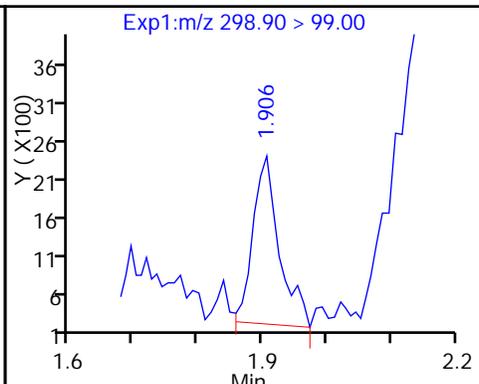
Method: A8_N

Limit Group: LC PFC_DOD ICAL

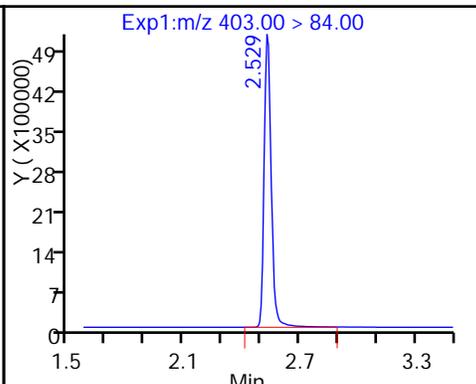
5 Perfluorobutanesulfonic acid



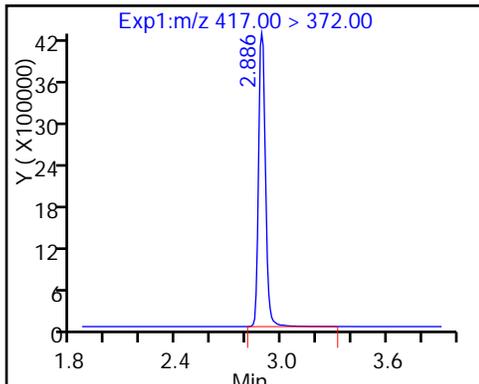
5 Perfluorobutanesulfonic acid



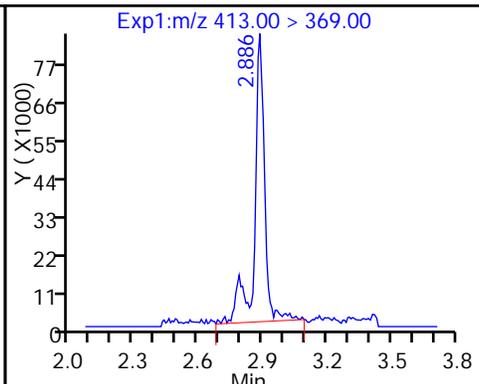
D 10 18O2 PFHxS



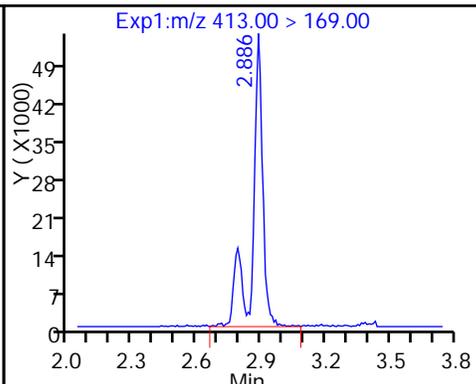
D 14 13C4 PFOA



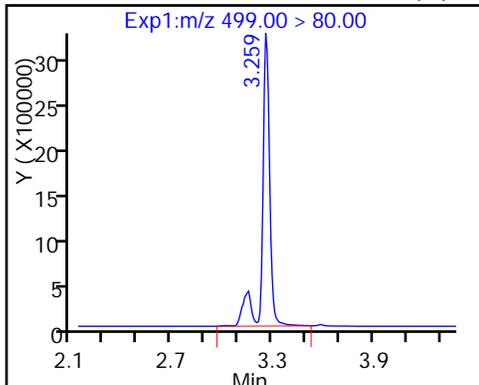
15 Perfluorooctanoic acid



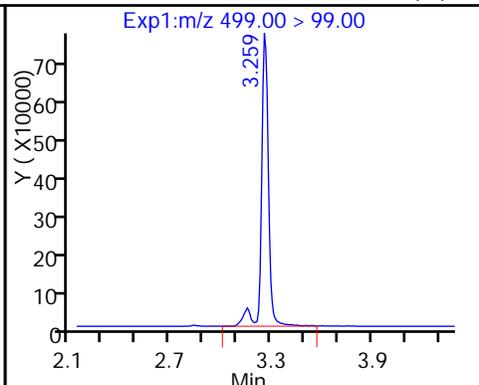
15 Perfluorooctanoic acid



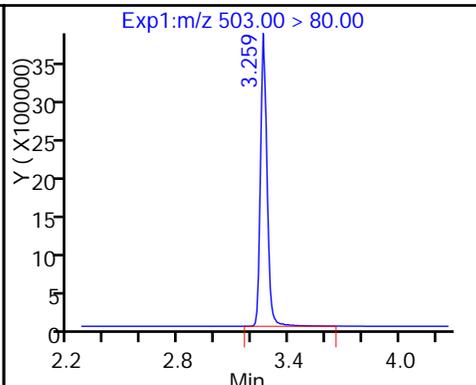
18 Perfluorooctane sulfonic acid (M)



18 Perfluorooctane sulfonic acid (M)



D 17 13C4 PFOS



TestAmerica Sacramento

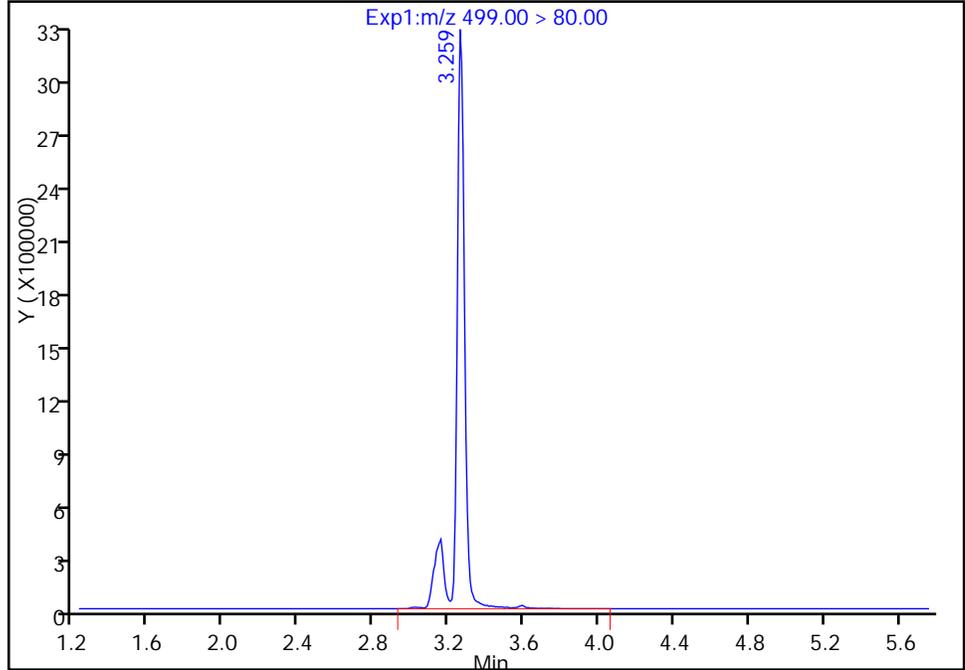
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_017.d
Injection Date: 02-Dec-2016 16:59:44 Instrument ID: A8_N
Lims ID: 320-23783-A-8-A Lab Sample ID: 320-23783-8
Client ID: DPT-16-31-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 13 Worklist Smp#: 21
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

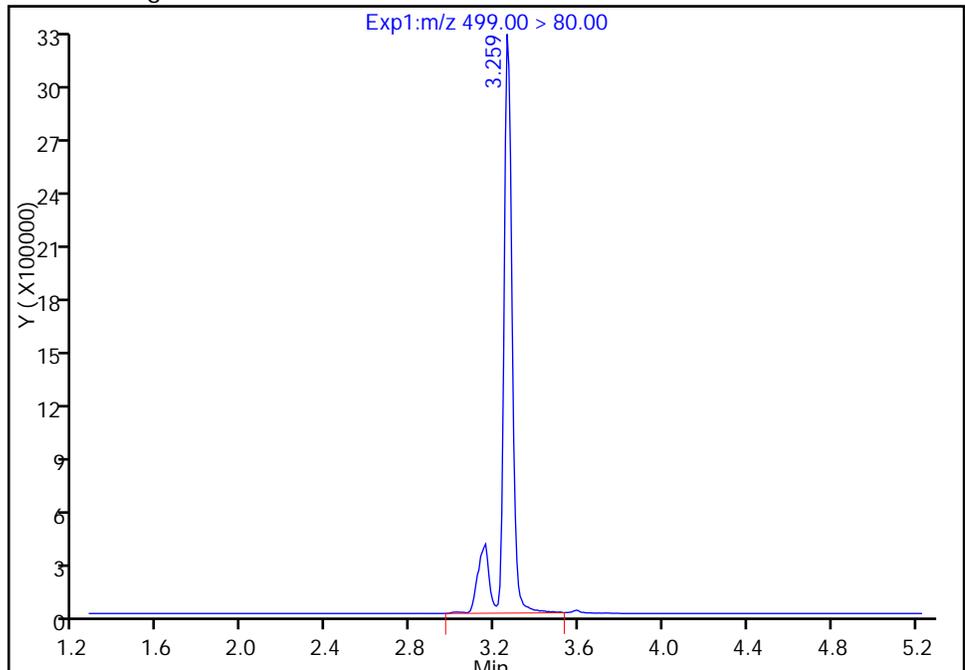
RT: 3.26
Area: 10384466
Amount: 47.640052
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 10232948
Amount: 46.944944
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:29:01
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

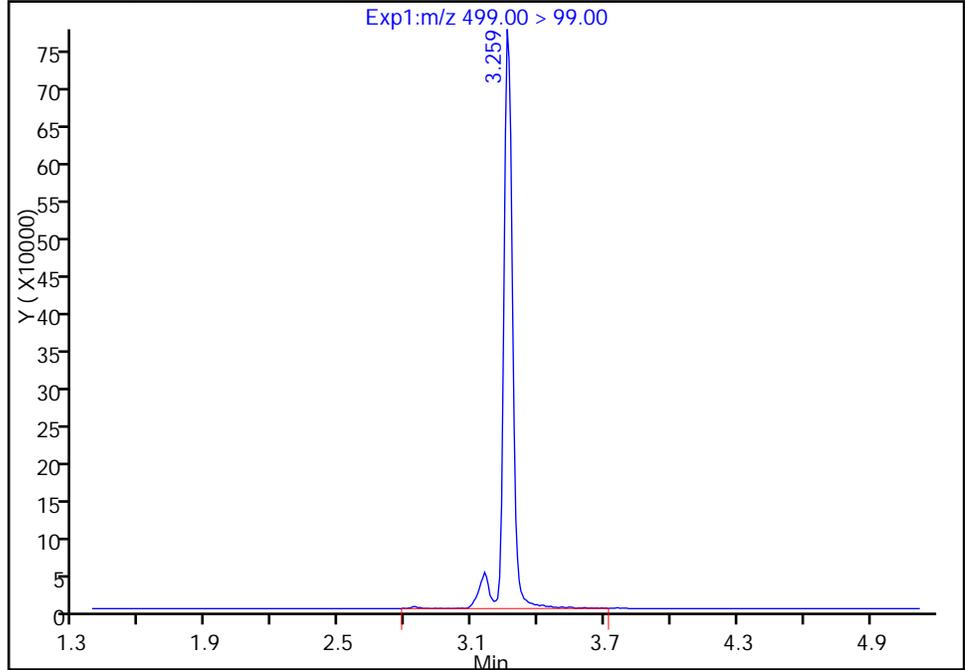
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_017.d
Injection Date: 02-Dec-2016 16:59:44 Instrument ID: A8_N
Lims ID: 320-23783-A-8-A Lab Sample ID: 320-23783-8
Client ID: DPT-16-31-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 13 Worklist Smp#: 21
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

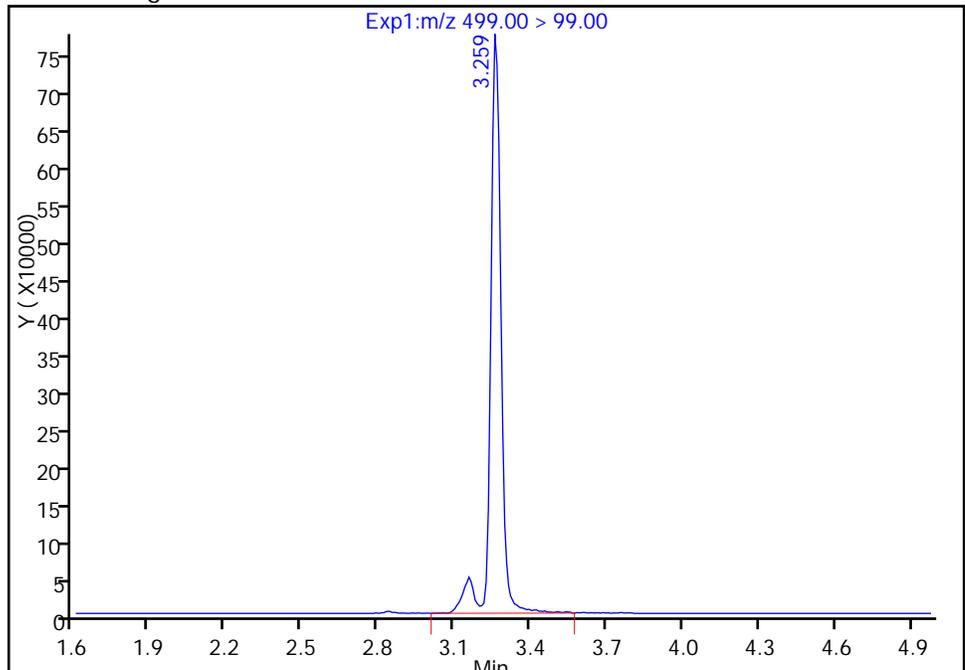
RT: 3.26
Area: 2281253
Amount: 47.640052
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 2262395
Amount: 46.944944
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-30-SO-14-15 Lab Sample ID: 320-23783-9
 Matrix: Solid Lab File ID: 02DEC2016C_018.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 10:46
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.02(g) Date Analyzed: 12/02/2016 17:07
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: 16.2 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	5.5		0.59	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	57	M	0.59	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.36	U	0.48	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	101		25-150
STL00991	13C4 PFOS	63		25-150
STL00994	18O2 PFHxS	90		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_018.d
 Lims ID: 320-23783-A-9-A
 Client ID: DPT-16-30-SO-14-15
 Sample Type: Client
 Inject. Date: 02-Dec-2016 17:07:17 ALS Bottle#: 14 Worklist Smp#: 22
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-9-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:29:56 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:29:56

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.906	1.907	-0.001	1.000	83824	0.1911				
298.90 > 99.00	1.906	1.907	-0.001	1.000	39255		2.14(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.531	2.534	-0.003		12857523	42.5		89.9	435506	
D 14 13C4 PFOA										
417.00 > 372.00	2.888	2.886	0.002		11077577	50.5		101	1065418	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.888	2.894	-0.006	1.000	5268737	23.3			80849	
413.00 > 169.00	2.888	2.894	-0.006	1.000	3495011		1.51(0.90-1.10)		139772	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.262	3.267	-0.005	1.000	40173597	240.8			358921	M
499.00 > 99.00	3.262	3.267	-0.005	1.000	8876431		4.53(0.90-1.10)		281528	M
D 17 13C4 PFOS										
503.00 > 80.00	3.262	3.267	-0.005		7333704	30.0		62.7	567209	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_018.d

Injection Date: 02-Dec-2016 17:07:17

Instrument ID: A8_N

Lims ID: 320-23783-A-9-A

Lab Sample ID: 320-23783-9

Client ID: DPT-16-30-SO-14-15

Operator ID: A8-PC\A8

ALS Bottle#: 14

Worklist Smp#: 22

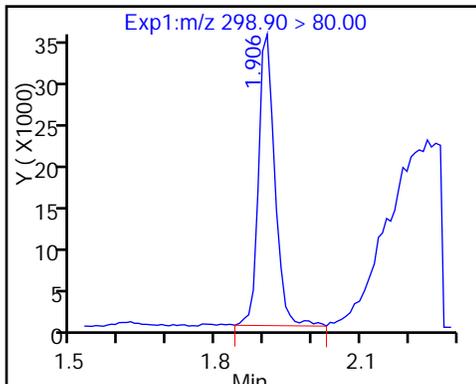
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

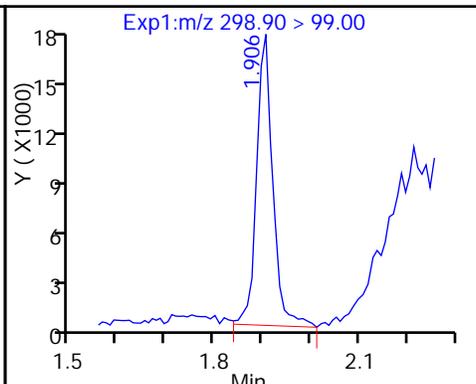
Method: A8_N

Limit Group: LC PFC_DOD ICAL

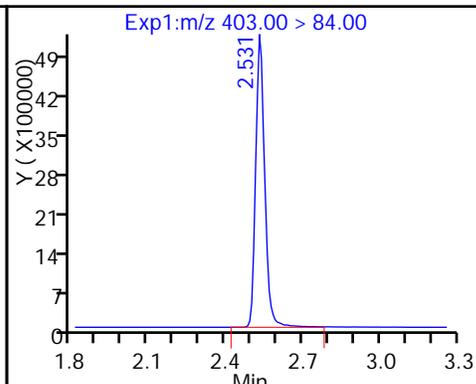
5 Perfluorobutanesulfonic acid



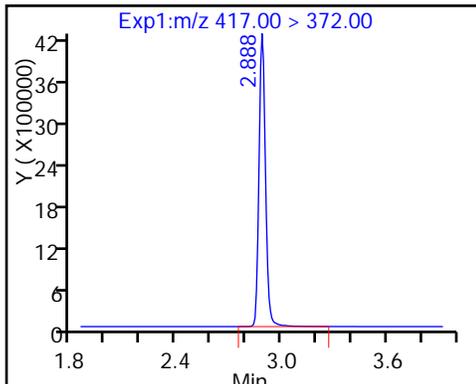
5 Perfluorobutanesulfonic acid



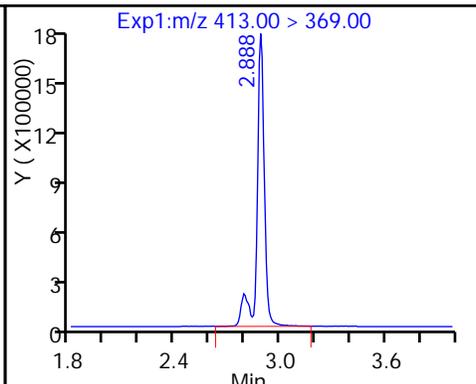
D 10 18O2 PFHxS



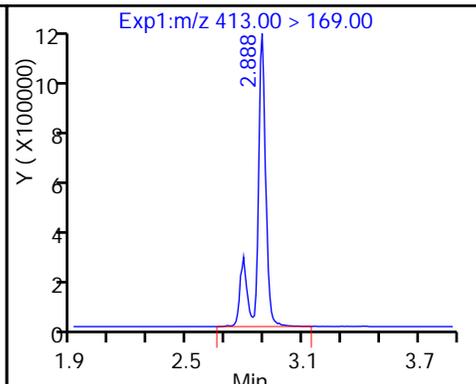
D 14 13C4 PFOA



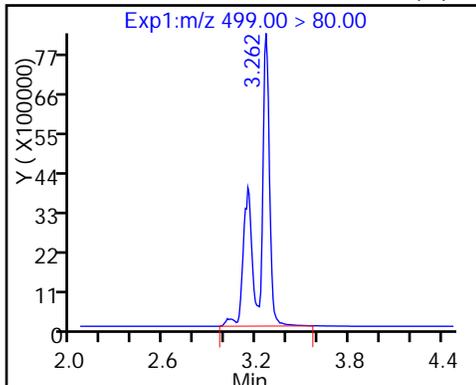
15 Perfluorooctanoic acid



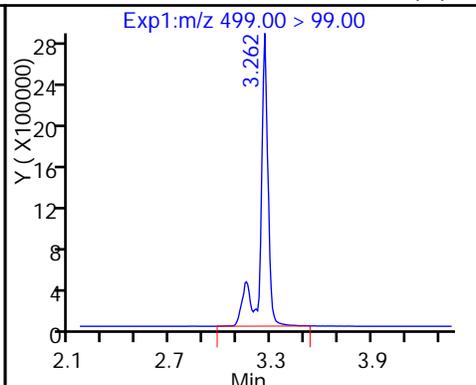
15 Perfluorooctanoic acid



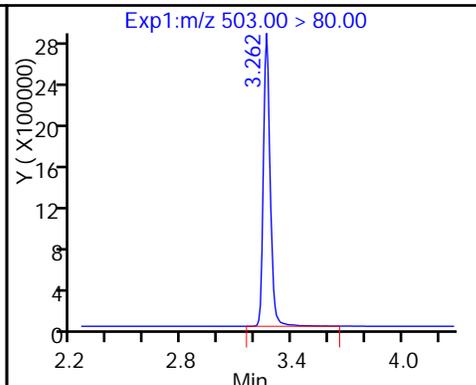
18 Perfluorooctane sulfonic acid (M)



18 Perfluorooctane sulfonic acid (M)



D 17 13C4 PFOS



TestAmerica Sacramento

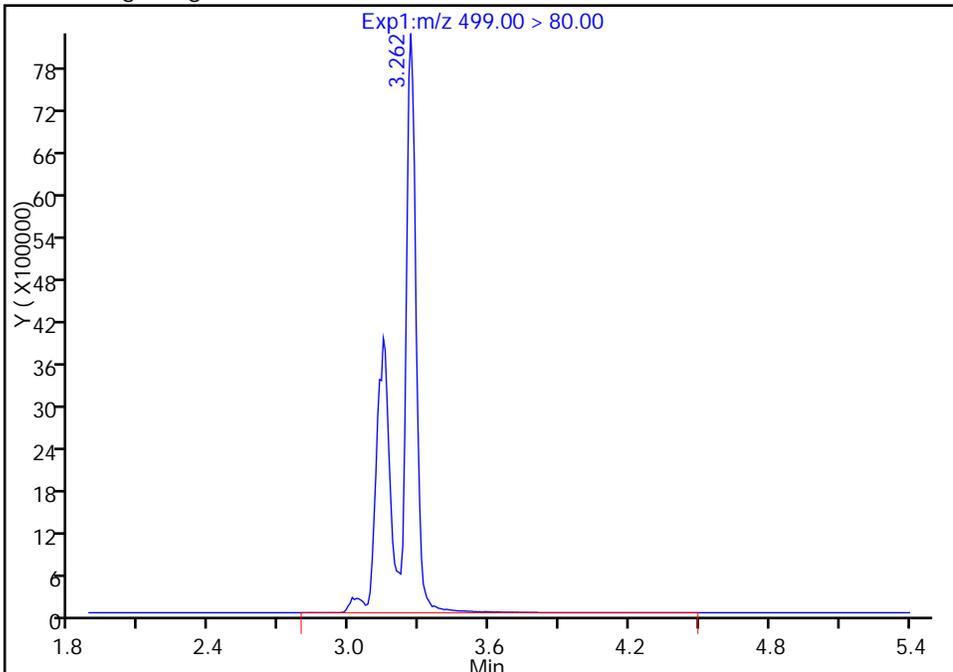
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_018.d
Injection Date: 02-Dec-2016 17:07:17 Instrument ID: A8_N
Lims ID: 320-23783-A-9-A Lab Sample ID: 320-23783-9
Client ID: DPT-16-30-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 14 Worklist Smp#: 22
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

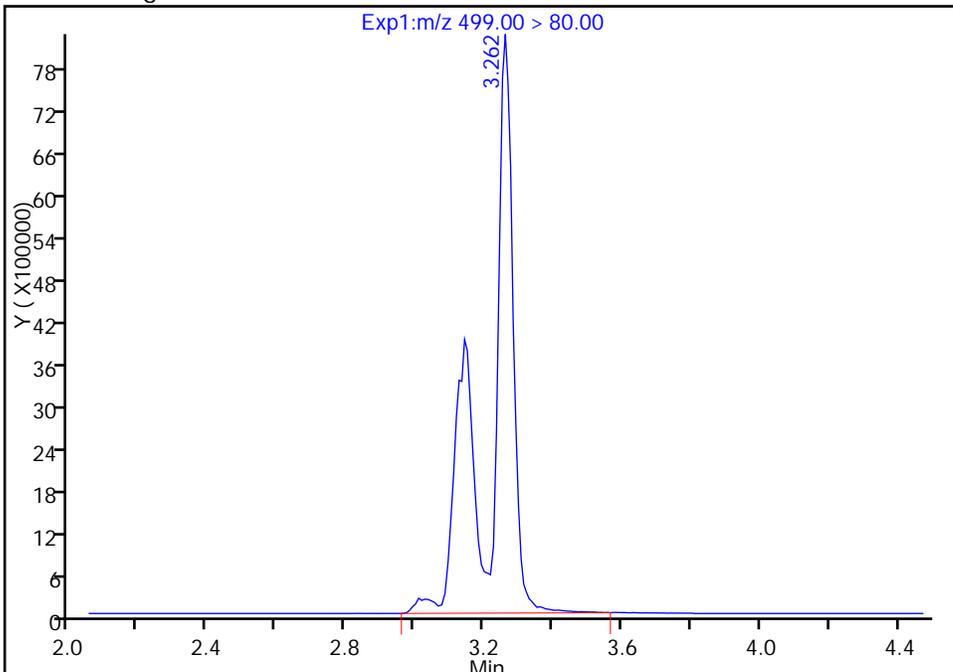
RT: 3.26
Area: 40507569
Amount: 242.8144
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 40173597
Amount: 240.8125
Amount Units: ng/ml

Manual Integration Results



TestAmerica Sacramento

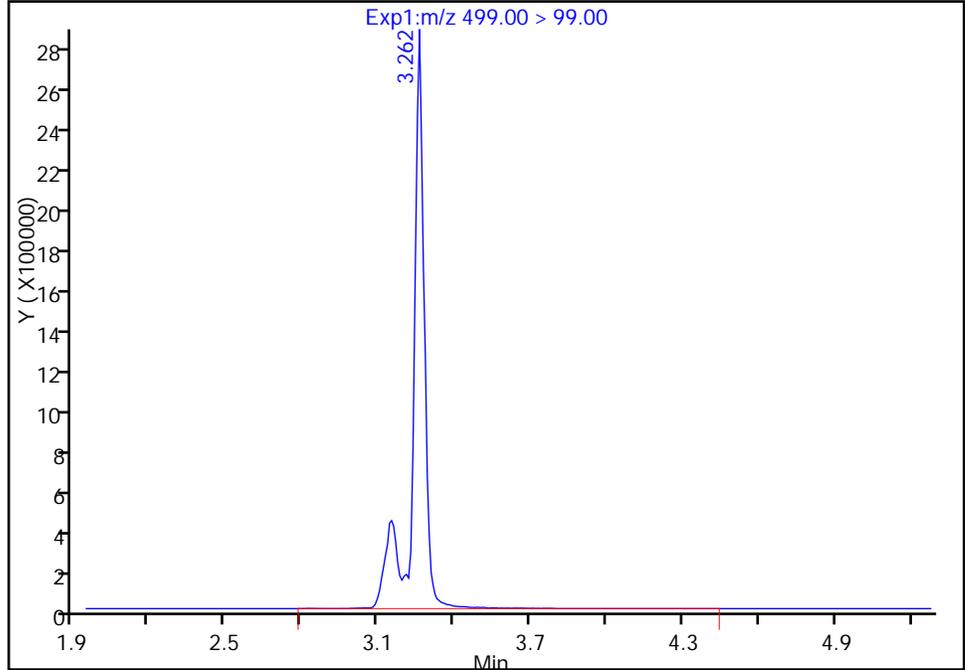
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_018.d
Injection Date: 02-Dec-2016 17:07:17 Instrument ID: A8_N
Lims ID: 320-23783-A-9-A Lab Sample ID: 320-23783-9
Client ID: DPT-16-30-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 14 Worklist Smp#: 22
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

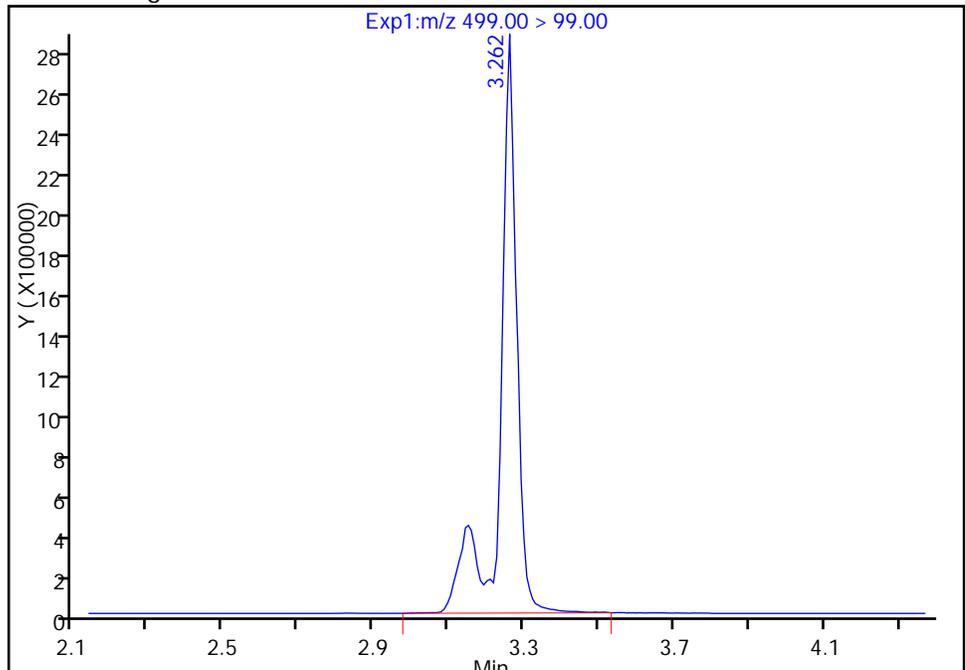
RT: 3.26
Area: 8968373
Amount: 242.8144
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 8876431
Amount: 240.8125
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-28-SO-14-15 Lab Sample ID: 320-23783-10
 Matrix: Solid Lab File ID: 02DEC2016C_019.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 11:54
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.03(g) Date Analyzed: 12/02/2016 17:14
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: 14.8 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.45	J	0.58	0.12
1763-23-1	Perfluorooctane Sulfonate (PFOS)	10	M	0.58	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.35	U	0.47	0.12

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	105		25-150
STL00991	13C4 PFOS	61		25-150
STL00994	18O2 PFHxS	89		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_019.d
 Lims ID: 320-23783-A-10-A
 Client ID: DPT-16-28-SO-14-15
 Sample Type: Client
 Inject. Date: 02-Dec-2016 17:14:47 ALS Bottle#: 15 Worklist Smp#: 23
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-10-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:30:24 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:30:24

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.897	1.907	-0.010	1.000	35638	0.0818				
298.90 > 99.00	1.897	1.907	-0.010	1.000	18199		1.96(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.522	2.534	-0.012		12769908	42.2		89.3	822686	
D 14 13C4 PFOA										
417.00 > 372.00	2.879	2.886	-0.007		11510268	52.5		105	884090	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.879	2.894	-0.015	1.000	453417	1.93			7493	
413.00 > 169.00	2.888	2.894	-0.006	1.003	277329		1.63(0.90-1.10)		14351	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.262	3.267	-0.005	1.000	7177023	44.4			141636	M
499.00 > 99.00	3.262	3.267	-0.005	1.000	1486211		4.83(0.90-1.10)		54507	M
D 17 13C4 PFOS										
503.00 > 80.00	3.262	3.267	-0.005		7102884	29.0		60.7	332839	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_019.d

Injection Date: 02-Dec-2016 17:14:47

Instrument ID: A8_N

Lims ID: 320-23783-A-10-A

Lab Sample ID: 320-23783-10

Client ID: DPT-16-28-SO-14-15

Operator ID: A8-PC\A8

ALS Bottle#: 15

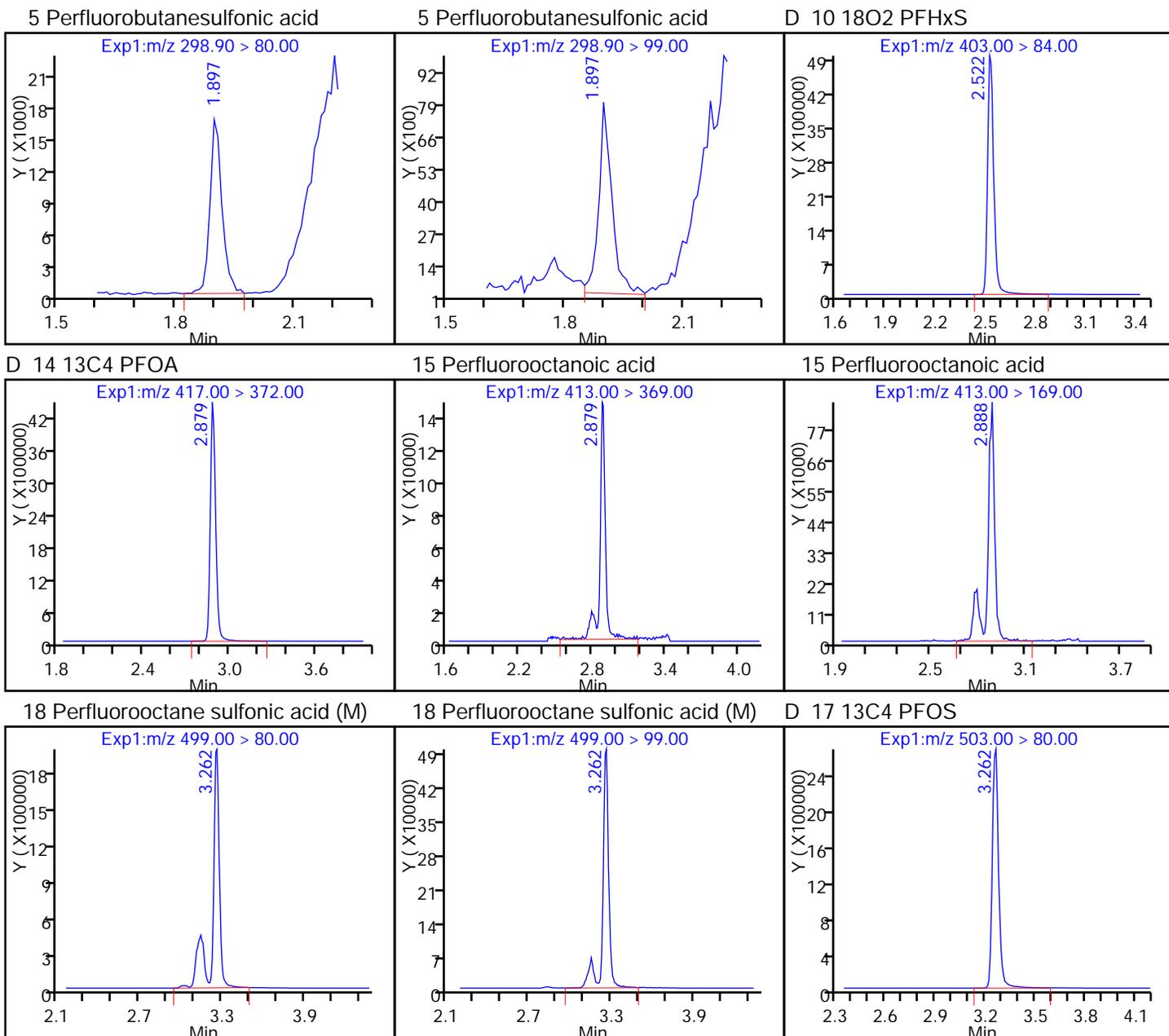
Worklist Smp#: 23

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

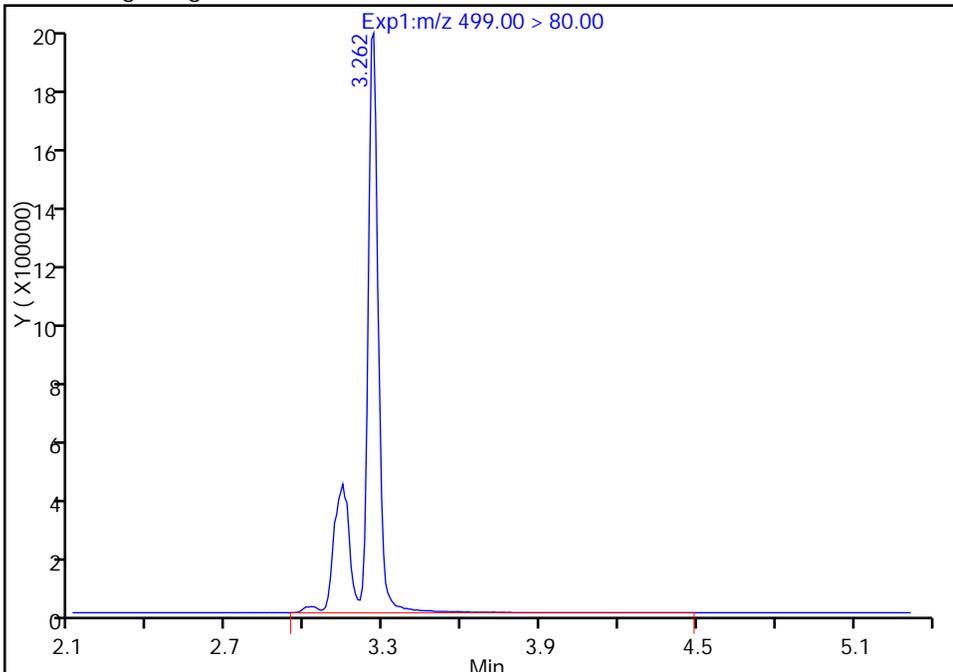
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_019.d
Injection Date: 02-Dec-2016 17:14:47 Instrument ID: A8_N
Lims ID: 320-23783-A-10-A Lab Sample ID: 320-23783-10
Client ID: DPT-16-28-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 15 Worklist Smp#: 23
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

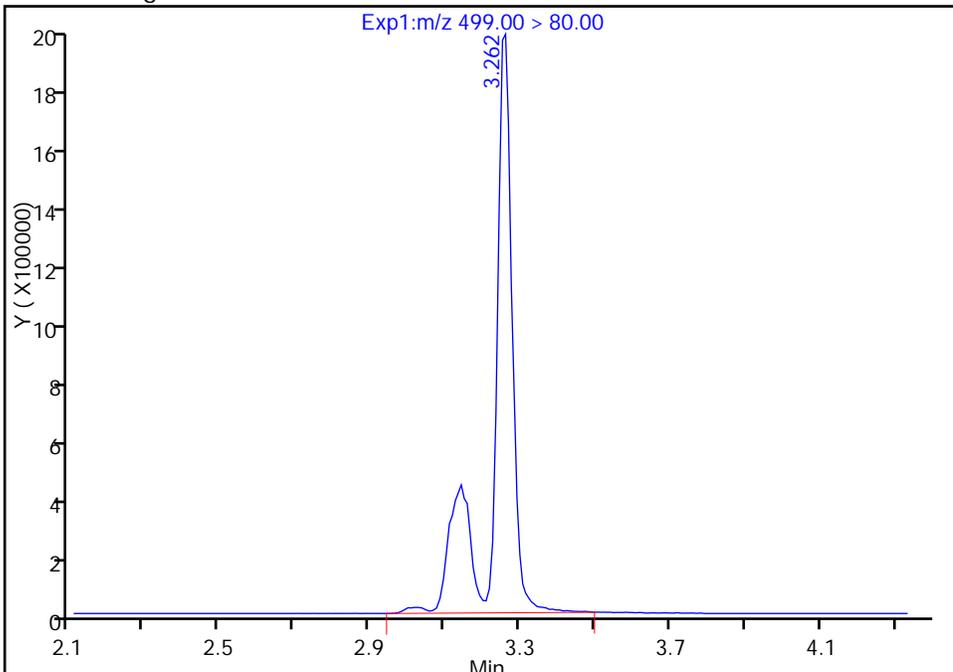
RT: 3.26
Area: 7288147
Amount: 45.107009
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 7177023
Amount: 44.419252
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:30:24
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

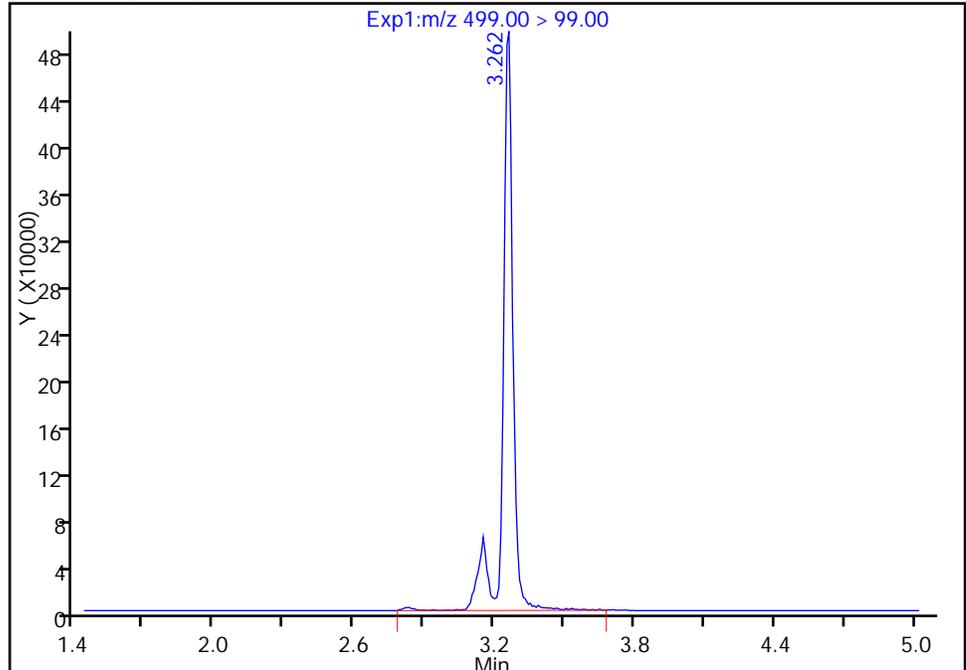
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_019.d
Injection Date: 02-Dec-2016 17:14:47 Instrument ID: A8_N
Lims ID: 320-23783-A-10-A Lab Sample ID: 320-23783-10
Client ID: DPT-16-28-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 15 Worklist Smp#: 23
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

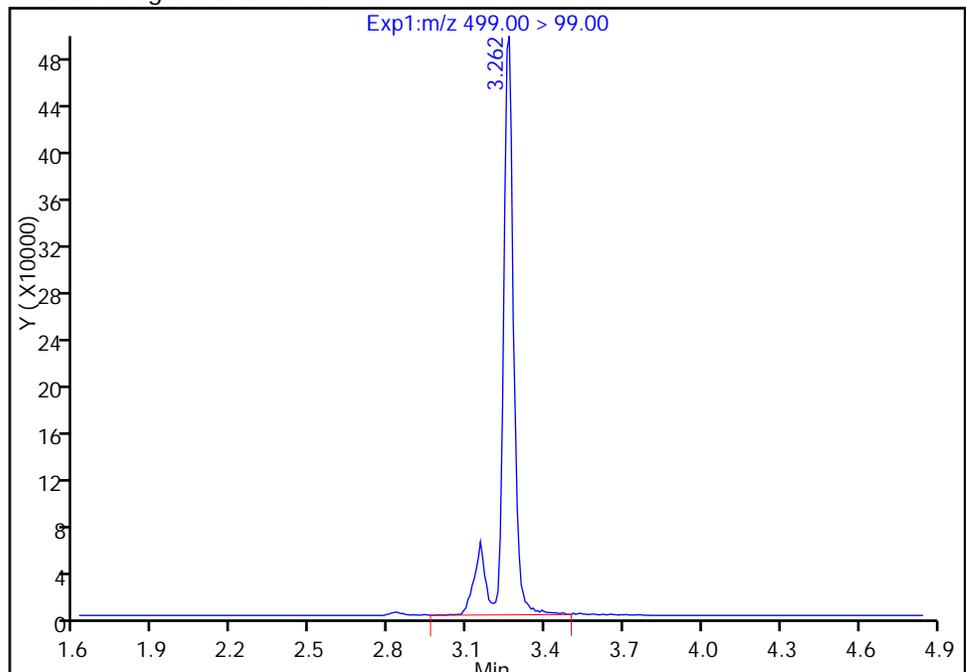
RT: 3.26
Area: 1513264
Amount: 45.107009
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 1486211
Amount: 44.419252
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-15-SO-14-15 Lab Sample ID: 320-23783-11
 Matrix: Solid Lab File ID: 02DEC2016C_020.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 12:50
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.05(g) Date Analyzed: 12/02/2016 17:22
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: 19.4 GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.51	J	0.61	0.13
1763-23-1	Perfluorooctane Sulfonate (PFOS)	11	M	0.61	0.15
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.37	U	0.49	0.13

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	101		25-150
STL00991	13C4 PFOS	66		25-150
STL00994	18O2 PFHxS	89		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_020.d
 Lims ID: 320-23783-A-11-A
 Client ID: DPT-16-15-SO-14-15
 Sample Type: Client
 Inject. Date: 02-Dec-2016 17:22:19 ALS Bottle#: 16 Worklist Smp#: 24
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-11-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:31:06 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:31:06

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.906	1.907	-0.001	1.000	66604	0.1533				
298.90 > 99.00	1.897	1.907	-0.010	0.995	29112		2.29(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.531	2.534	-0.003		12739222	42.1		89.0	636115	
D 14 13C4 PFOA										
417.00 > 372.00	2.888	2.886	0.002		11061350	50.4		101	1057568	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.880	2.894	-0.014	1.000	465088	2.06			7440	
413.00 > 169.00	2.880	2.894	-0.014	1.000	288775		1.61(0.90-1.10)		14333	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.262	3.267	-0.005	1.000	7622896	43.1			269368	M
499.00 > 99.00	3.262	3.267	-0.005	1.000	1717205		4.44(0.90-1.10)		67501	M
D 17 13C4 PFOS										
503.00 > 80.00	3.262	3.267	-0.005		7772512	31.8		66.5	373209	

QC Flag Legend

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_020.d

Injection Date: 02-Dec-2016 17:22:19

Instrument ID: A8_N

Lims ID: 320-23783-A-11-A

Lab Sample ID: 320-23783-11

Client ID: DPT-16-15-SO-14-15

Operator ID: A8-PC\A8

ALS Bottle#: 16

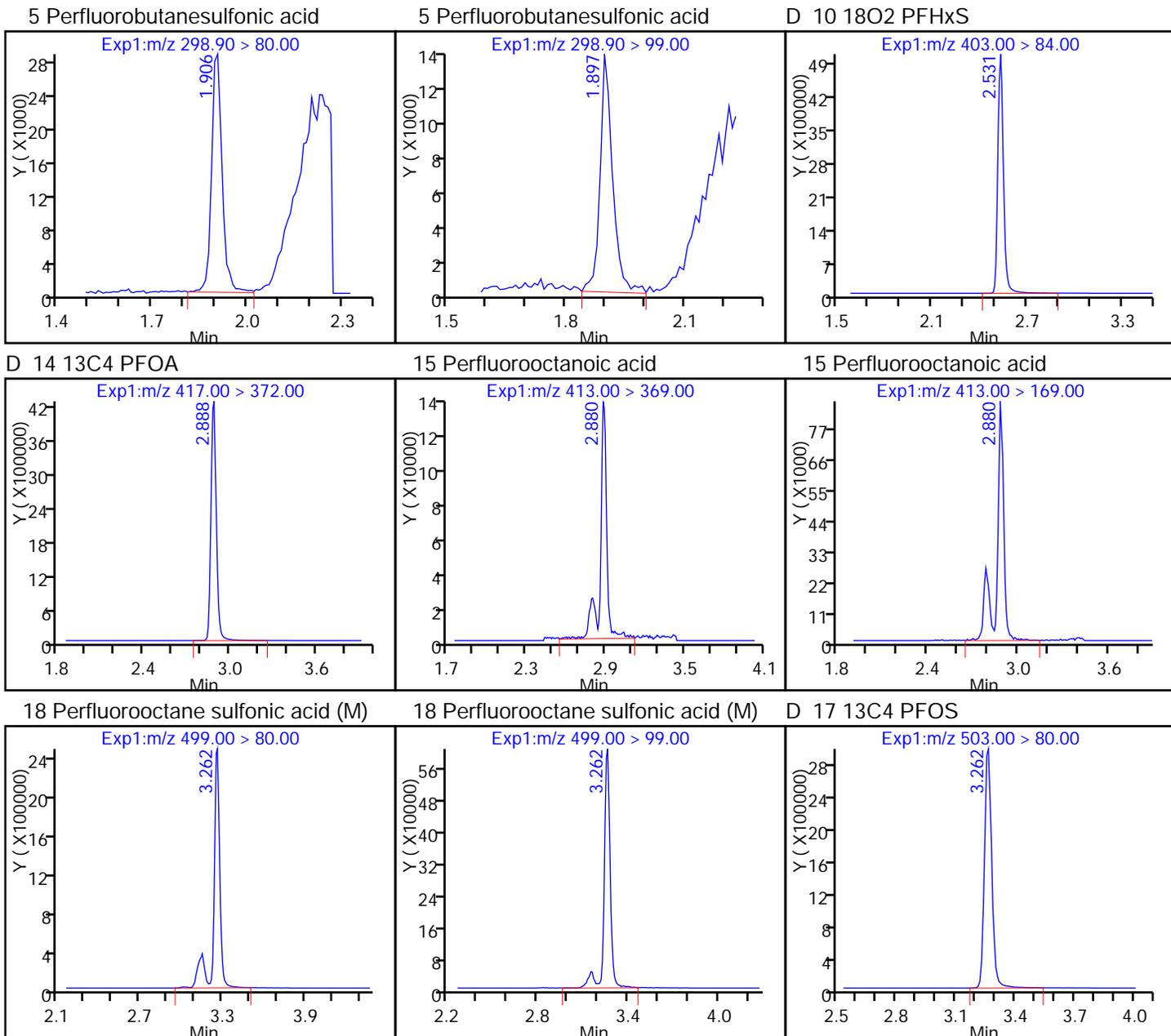
Worklist Smp#: 24

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

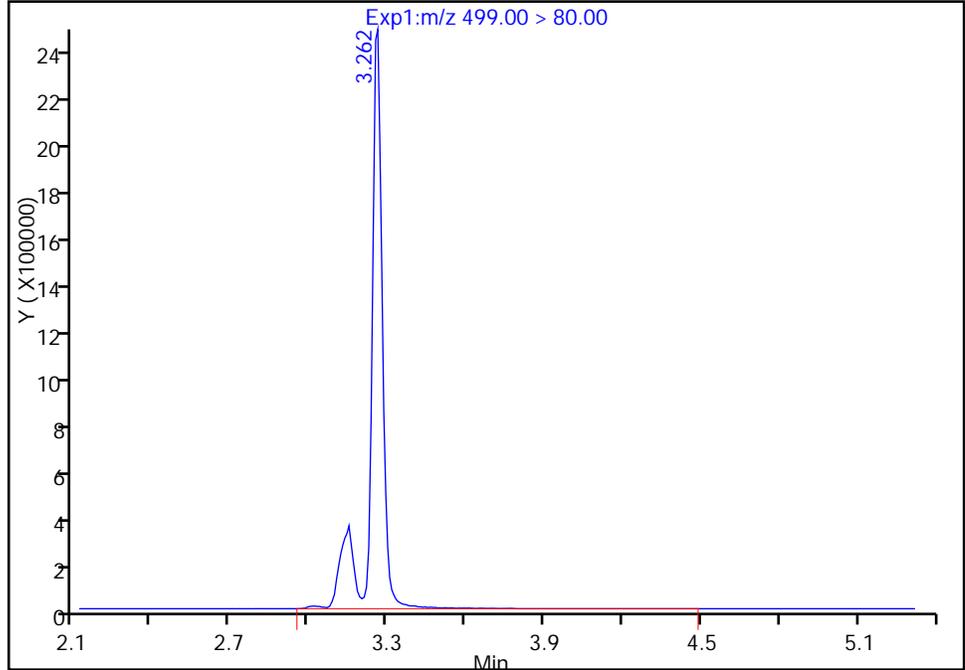
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_020.d
Injection Date: 02-Dec-2016 17:22:19 Instrument ID: A8_N
Lims ID: 320-23783-A-11-A Lab Sample ID: 320-23783-11
Client ID: DPT-16-15-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 16 Worklist Smp#: 24
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

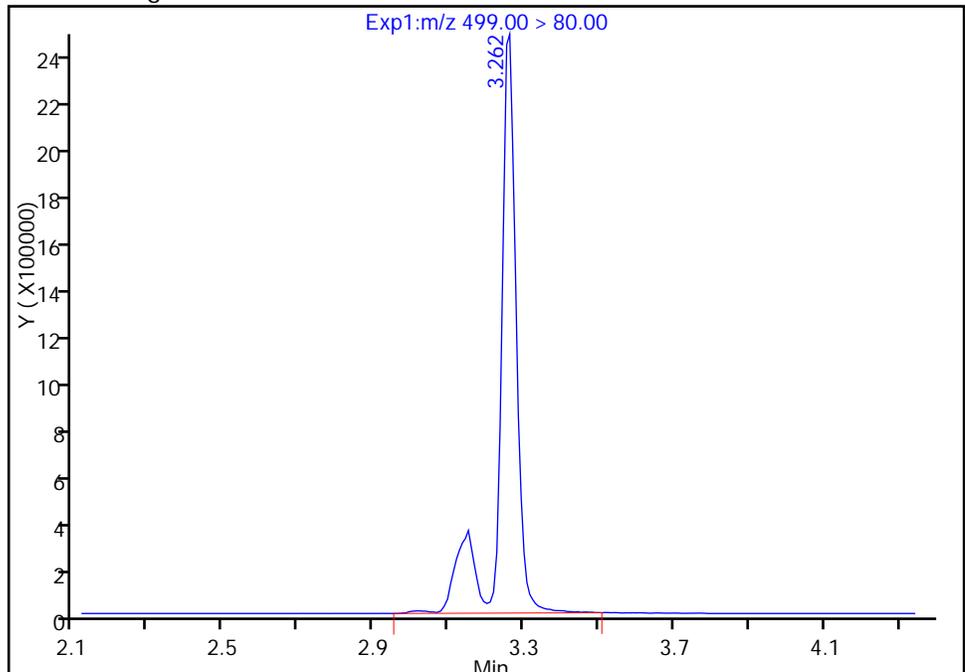
RT: 3.26
Area: 7709603
Amount: 43.604593
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 7622896
Amount: 43.114189
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 04-Dec-2016 18:31:06
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

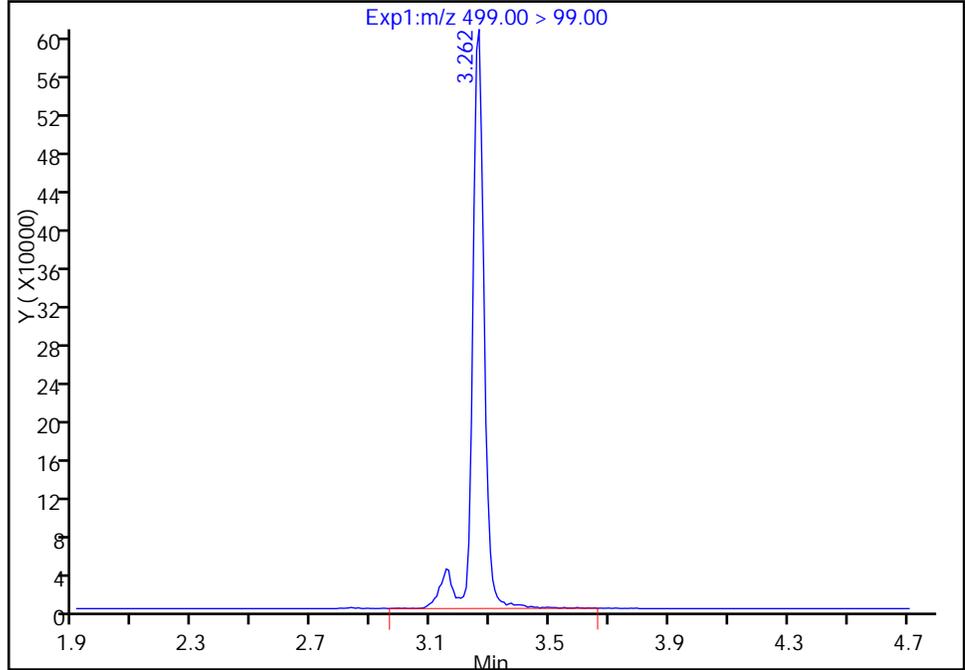
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_020.d
Injection Date: 02-Dec-2016 17:22:19 Instrument ID: A8_N
Lims ID: 320-23783-A-11-A Lab Sample ID: 320-23783-11
Client ID: DPT-16-15-SO-14-15
Operator ID: A8-PC\A8 ALS Bottle#: 16 Worklist Smp#: 24
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

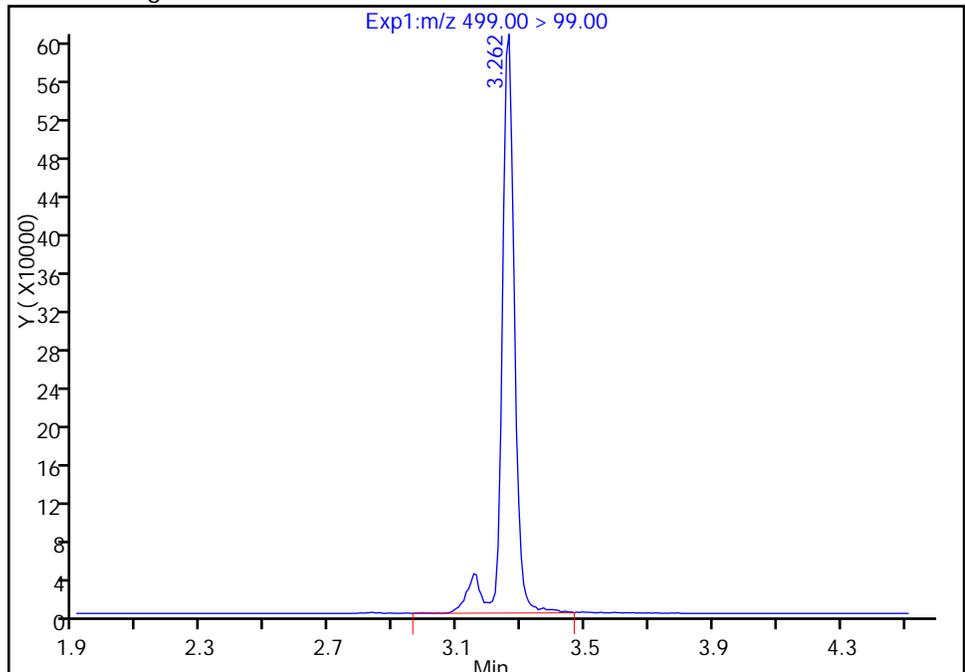
RT: 3.26
Area: 1729991
Amount: 43.604593
Amount Units: ng/ml

Processing Integration Results



RT: 3.26
Area: 1717205
Amount: 43.114189
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-15-GW-18-22 Lab Sample ID: 320-23783-12
 Matrix: Water Lab File ID: 03DEC2016B_006.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 13:30
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 249.7(mL) Date Analyzed: 12/03/2016 16:41
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	<i>Perfluorooctanoic acid (PFOA)</i>	5.8	E M	0.0025	0.00075
1763-23-1	<i>Perfluorooctane Sulfonate (PFOS)</i>	6.5	E M	0.0040	0.0013
375-73-5	<i>Perfluorobutanesulfonic acid (PFBS)</i>	1.2	E	0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	23	Q	25-150
STL00991	13C4 PFOS	29		25-150
STL00994	18O2 PFHxS	21	Q	25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d
 Lims ID: 320-23783-A-12-A
 Client ID: DPT-16-15-GW-18-22
 Sample Type: Client
 Inject. Date: 03-Dec-2016 16:41:11 ALS Bottle#: 37 Worklist Smp#: 27
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-12-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:35:31 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:35:31

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.868	1.900	-0.032	1.000	63043653	598.1				E
298.90 > 99.00	1.878	1.900	-0.022	1.005	38170744		1.65(0.00-0.00)			E
D 10 18O2 PFHxS										
403.00 > 84.00	2.489	2.531	-0.042		3123671	10.0		21.1	34160	
D 14 13C4 PFOA										
417.00 > 372.00	2.841	2.880	-0.039		2480157	11.3		22.6	166976	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.817	2.887	-0.070	1.000	152985407	2877.3			85801	EM
413.00 > 169.00	2.825	2.887	-0.062	1.003	150895964		1.01(0.90-1.10)		144575	M
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.191	3.258	-0.067	1.000	255054939	3244.2			125297	EM
499.00 > 99.00	3.217	3.258	-0.041	1.008	89493448		2.85(0.90-1.10)		438551	M
D 17 13C4 PFOS										
503.00 > 80.00	3.217	3.259	-0.042		3462821	14.1		29.4	43644	

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d

Injection Date: 03-Dec-2016 16:41:11

Instrument ID: A8_N

Lims ID: 320-23783-A-12-A

Lab Sample ID: 320-23783-12

Client ID: DPT-16-15-GW-18-22

Operator ID: A8-PC\A8

ALS Bottle#: 37

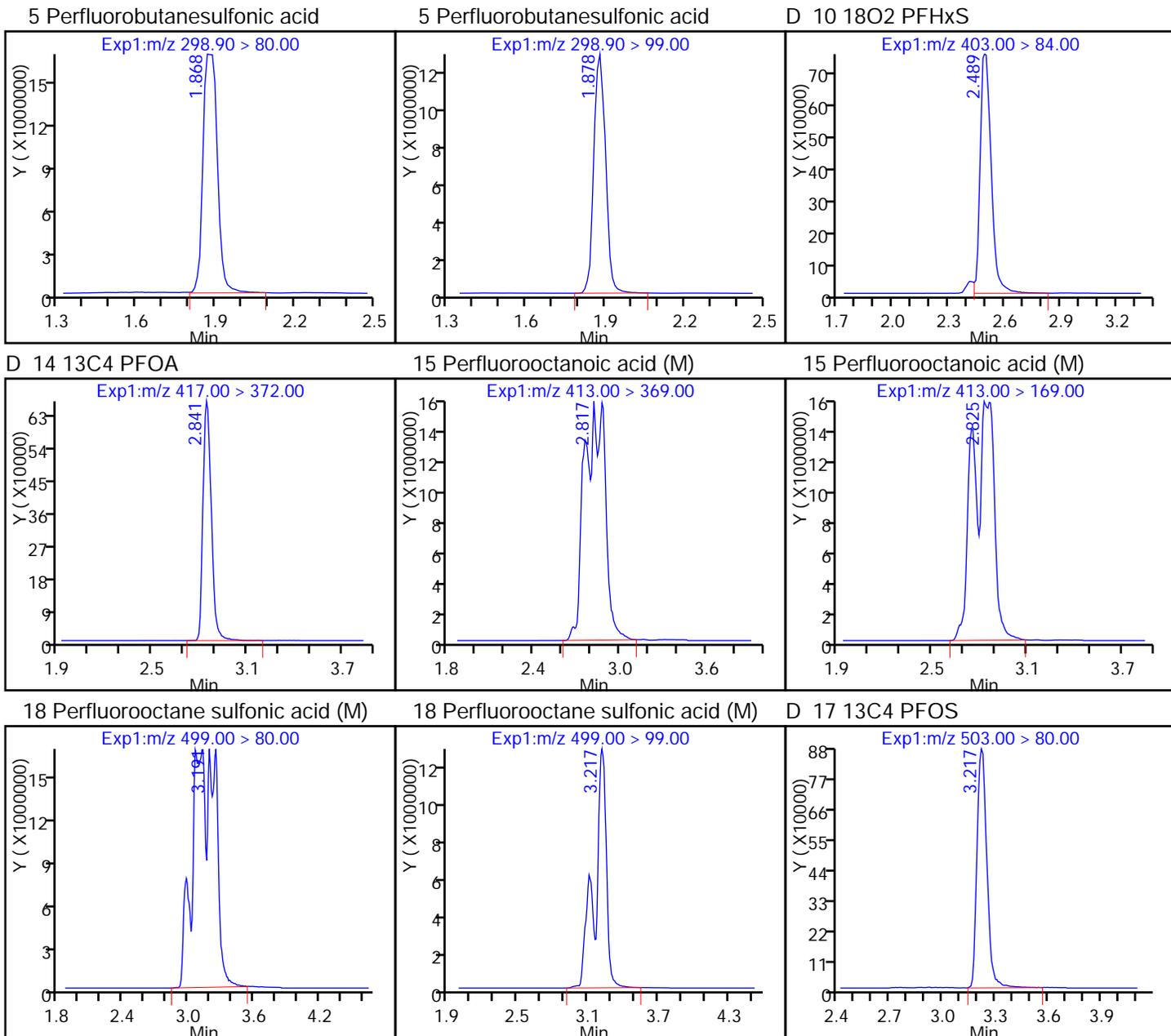
Worklist Smp#: 27

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

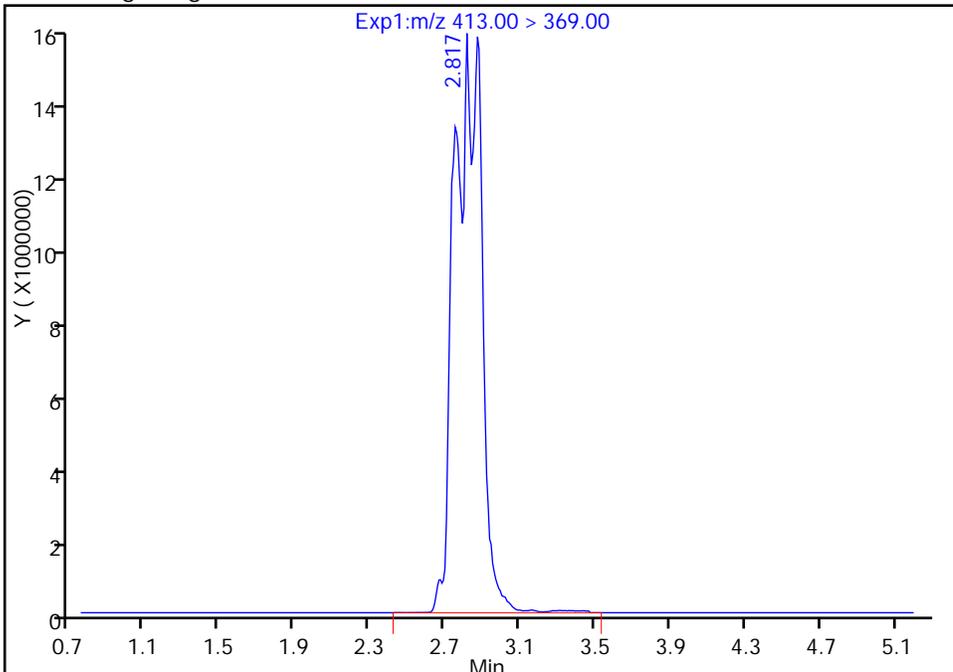
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d
Injection Date: 03-Dec-2016 16:41:11 Instrument ID: A8_N
Lims ID: 320-23783-A-12-A Lab Sample ID: 320-23783-12
Client ID: DPT-16-15-GW-18-22
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 27
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

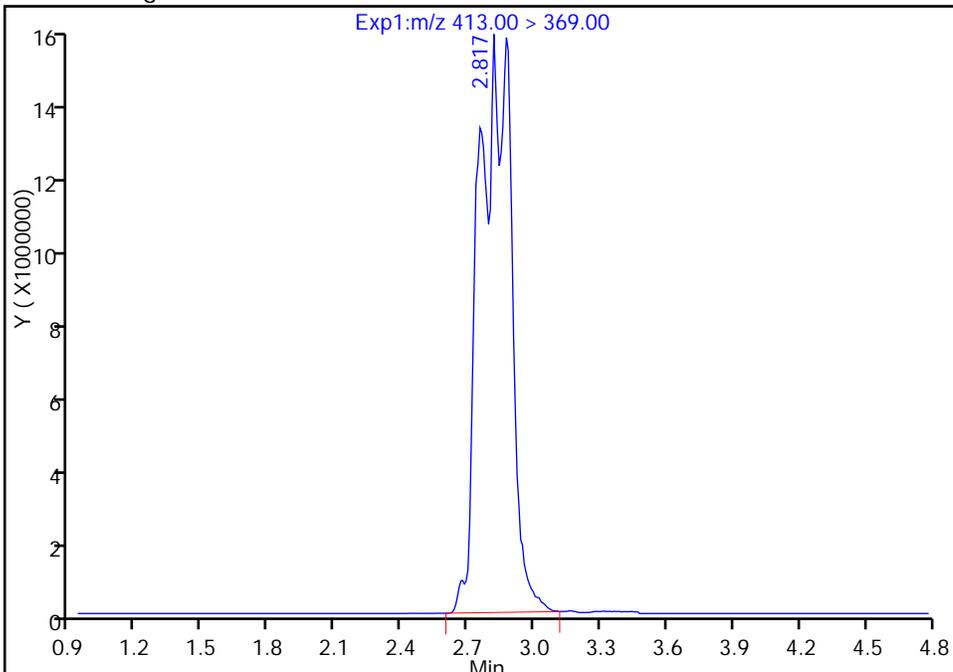
RT: 2.82
Area: 154922470
Amount: 2913.7001
Amount Units: ng/ml

Processing Integration Results



RT: 2.82
Area: 152985407
Amount: 2877.2689
Amount Units: ng/ml

Manual Integration Results



TestAmerica Sacramento

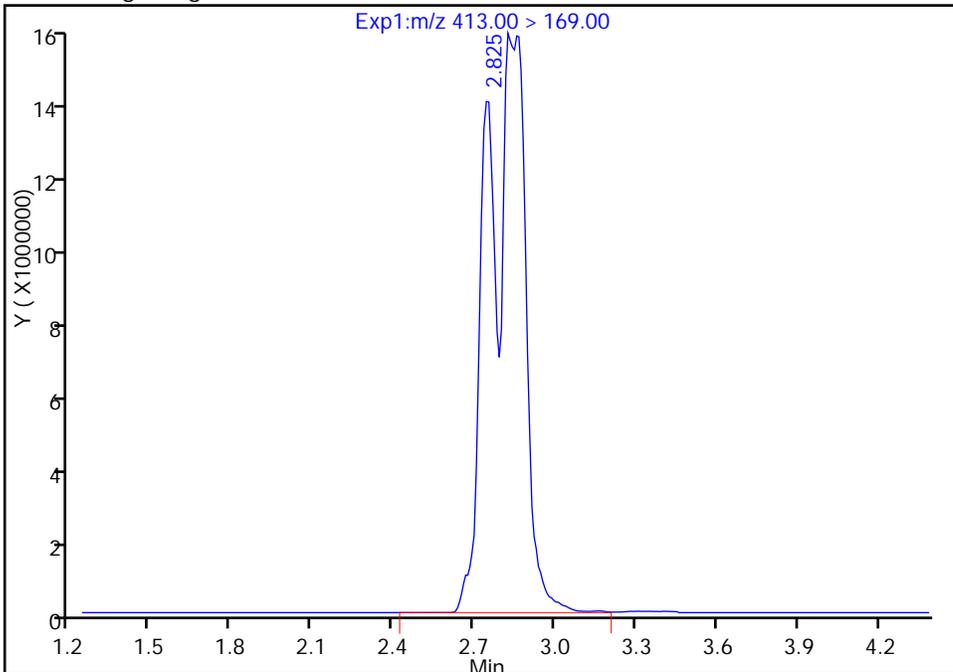
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d
Injection Date: 03-Dec-2016 16:41:11 Instrument ID: A8_N
Lims ID: 320-23783-A-12-A Lab Sample ID: 320-23783-12
Client ID: DPT-16-15-GW-18-22
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 27
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 2

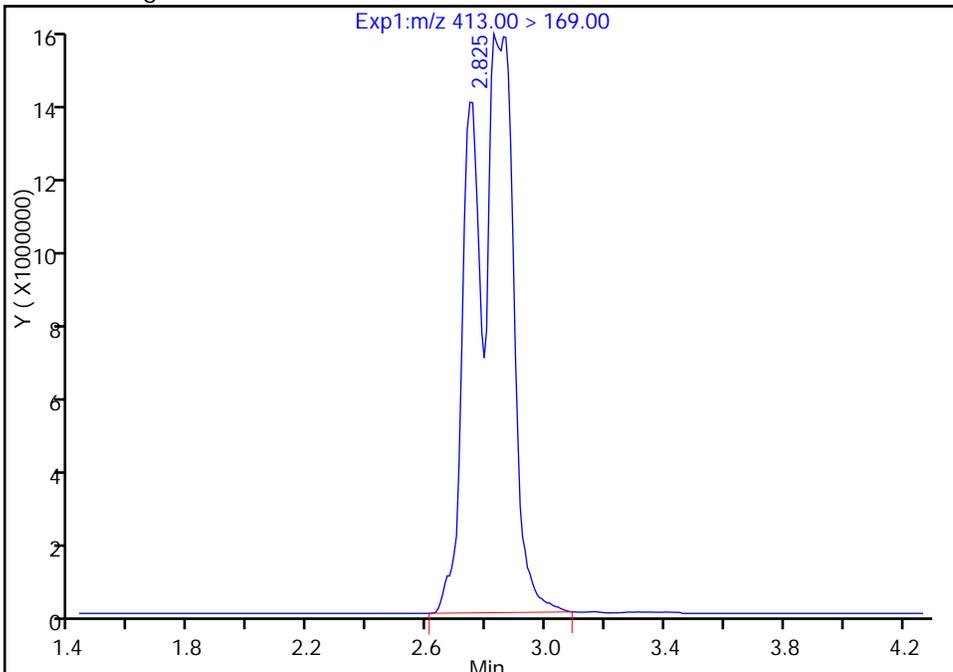
RT: 2.83
Area: 151742676
Amount: 2913.7001
Amount Units: ng/ml

Processing Integration Results



RT: 2.83
Area: 150895964
Amount: 2877.2689
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 10:35:31

Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

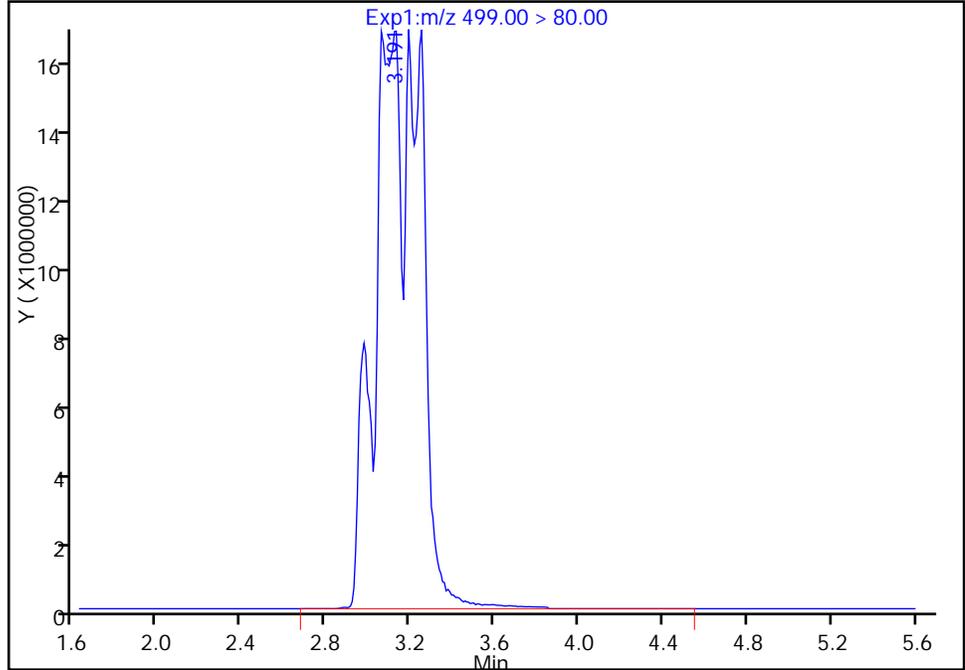
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d
Injection Date: 03-Dec-2016 16:41:11 Instrument ID: A8_N
Lims ID: 320-23783-A-12-A Lab Sample ID: 320-23783-12
Client ID: DPT-16-15-GW-18-22
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 27
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 1

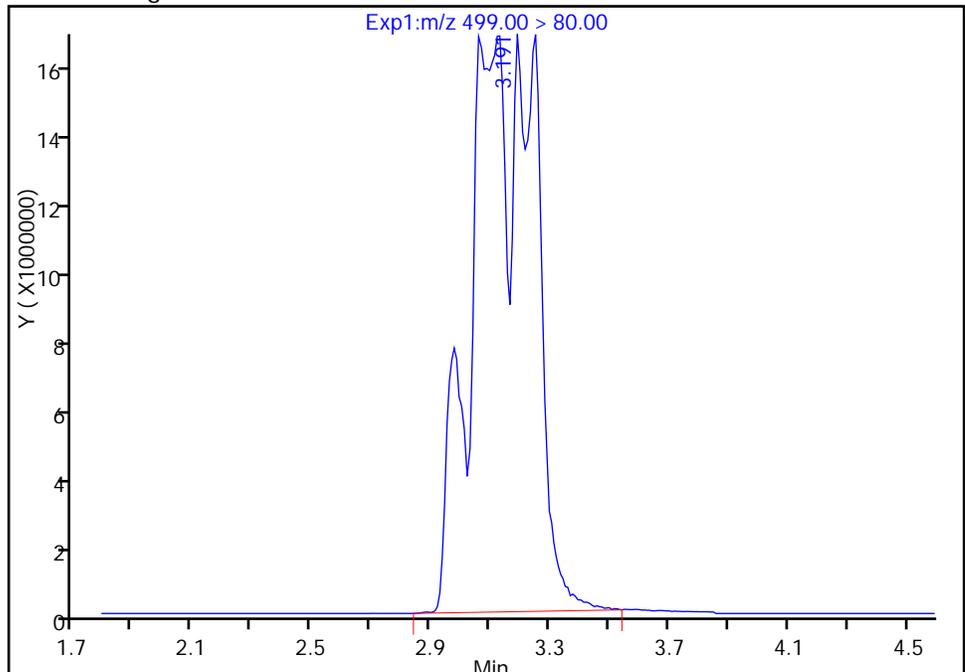
RT: 3.19
Area: 258759726
Amount: 3291.3581
Amount Units: ng/ml

Processing Integration Results



RT: 3.19
Area: 255054939
Amount: 3244.2341
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 10:35:31
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

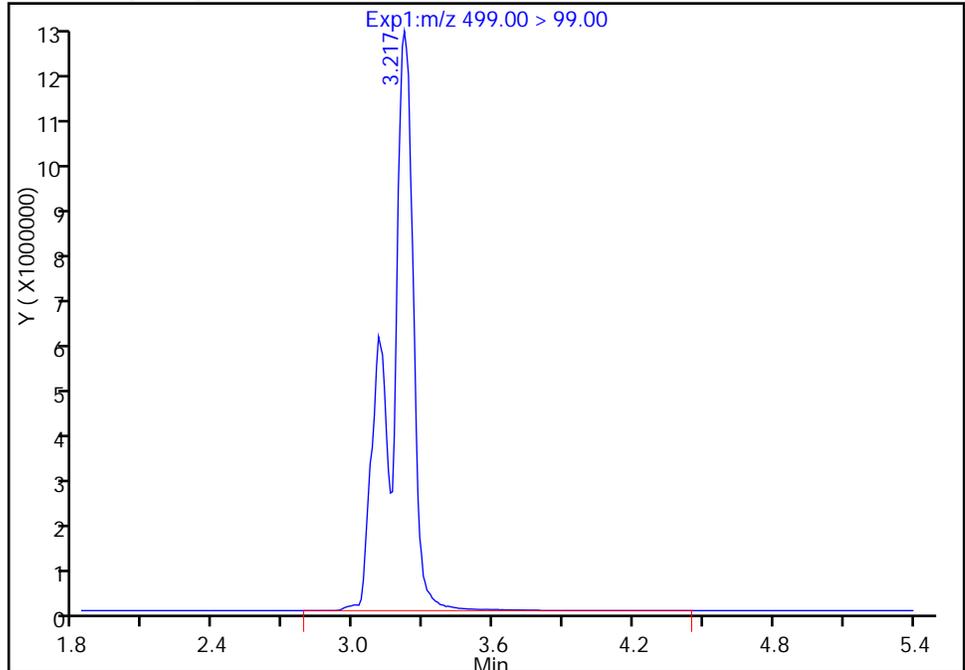
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_006.d
Injection Date: 03-Dec-2016 16:41:11 Instrument ID: A8_N
Lims ID: 320-23783-A-12-A Lab Sample ID: 320-23783-12
Client ID: DPT-16-15-GW-18-22
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 27
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

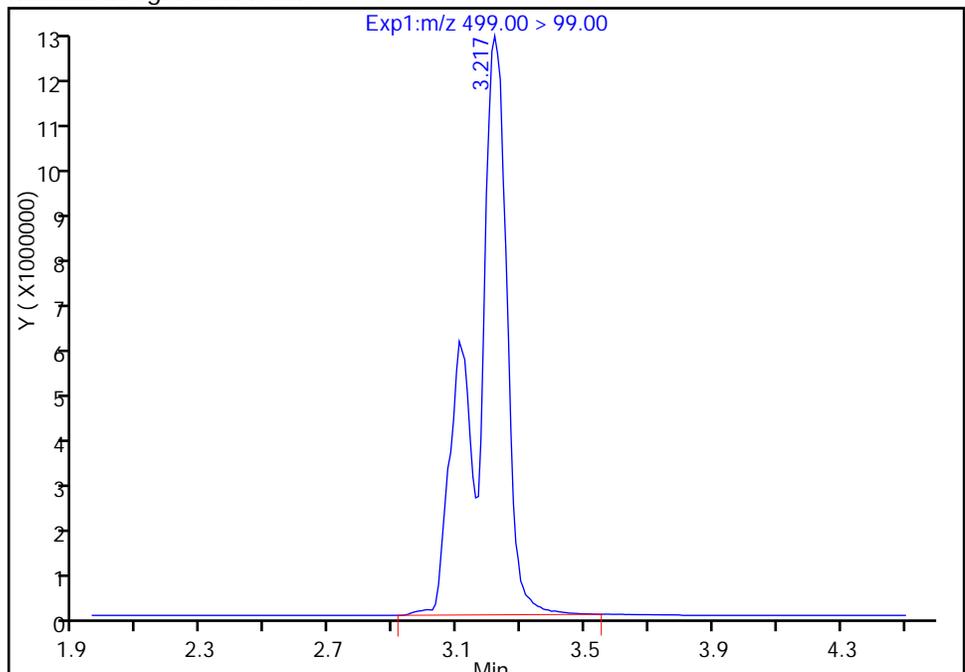
RT: 3.22
Area: 90233657
Amount: 3291.3581
Amount Units: ng/ml

Processing Integration Results



RT: 3.22
Area: 89493448
Amount: 3244.2341
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-15-GW-18-22 DL Lab Sample ID: 320-23783-12 DL
 Matrix: Water Lab File ID: 07DEC2016B_005.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 13:30
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 249.7(mL) Date Analyzed: 12/07/2016 17:03
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 100
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 141062 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	15	D	0.25	0.075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	12	D	0.40	0.13
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.54	D	0.25	0.092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	83		25-150
STL00991	13C4 PFOS	94		25-150
STL00994	18O2 PFHxS	104		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_005.d
 Lims ID: 320-23783-A-12-A
 Client ID: DPT-16-15-GW-18-22
 Sample Type: Client
 Inject. Date: 07-Dec-2016 17:03:25 ALS Bottle#: 18 Worklist Smp#: 5
 Injection Vol: 2.0 ul Dil. Factor: 100.0000
 Sample Info: 320-23783-a-12-a 100X
 Misc. Info.: Plate: 1 Rack: 3
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:54 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:54:48

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.852	1.848	0.004	1.000	1388458	2.68				
298.90 > 99.00	1.852	1.848	0.004	1.000	589024		2.36(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.452	2.460	-0.008		153817	0.4925		1.0	28373	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.802	2.802	0.0	1.000	14553741	74.9			293900	
413.00 > 169.00	2.802	2.802	0.0	1.000	9910045		1.47(0.90-1.10)		410094	
D 14 13C4 PFOA										
417.00 > 372.00	2.810	2.802	0.008		90597	0.4132		0.8	10591	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.162	3.061	0.101	1.000	14701138	58.6			349461	
499.00 > 99.00	3.170	3.061	0.109	1.003	2913043		5.05(0.90-1.10)		226375	
D 17 13C4 PFOS										
503.00 > 80.00	3.170	3.169	0.001		110529	0.4493		0.9	10146	

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_005.d

Injection Date: 07-Dec-2016 17:03:25

Instrument ID: A8_N

Lims ID: 320-23783-A-12-A

Lab Sample ID: 320-23783-12

Client ID: DPT-16-15-GW-18-22

Operator ID: A8-PC\A8

ALS Bottle#: 18

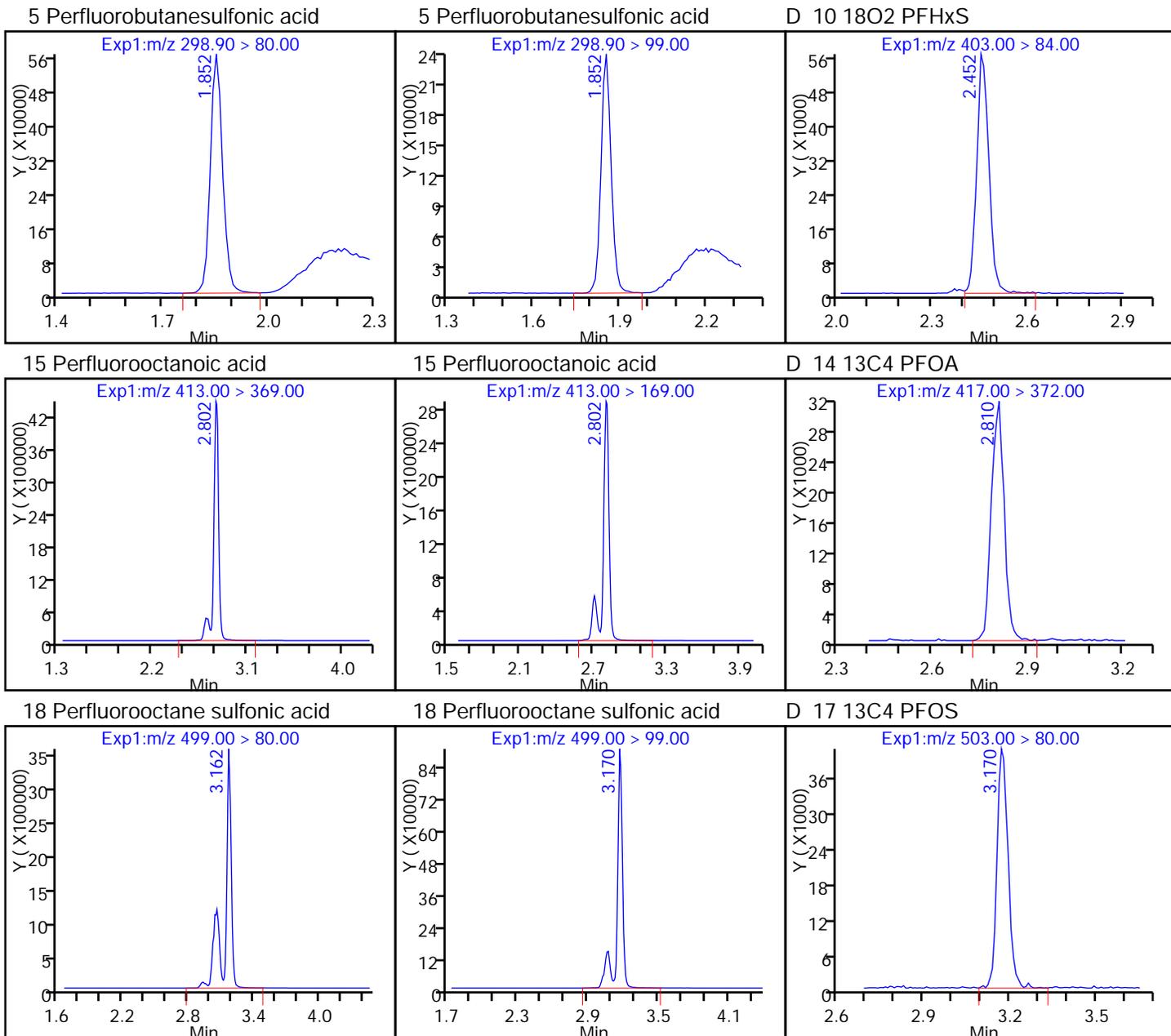
Worklist Smp#: 5

Injection Vol: 2.0 ul

Dil. Factor: 100.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-15-GW-31-35 Lab Sample ID: 320-23783-13
 Matrix: Water Lab File ID: 03DEC2016B_007.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 14:00
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 245.9(mL) Date Analyzed: 12/03/2016 16:48
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.27	M	0.0025	0.00076
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.46	E	0.0041	0.0013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.048		0.0025	0.00093

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	22	Q	25-150
STL00991	13C4 PFOS	82		25-150
STL00994	18O2 PFHxS	66		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_007.d
 Lims ID: 320-23783-A-13-A
 Client ID: DPT-16-15-GW-31-35
 Sample Type: Client
 Inject. Date: 03-Dec-2016 16:48:42 ALS Bottle#: 38 Worklist Smp#: 28
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: 320-23783-a-13-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:36:41 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:36:41

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.877	1.900	-0.023	1.000	7767443	23.5				
298.90 > 99.00	1.877	1.900	-0.023	1.000	3342231		2.32(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.510	2.531	-0.021		9782375	31.3		66.2	382293	
D 14 13C4 PFOA										
417.00 > 372.00	2.846	2.880	-0.034		2458776	11.2		22.4	137087	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.862	2.887	-0.025	1.000	6930971	131.5			41657	M
413.00 > 169.00	2.862	2.887	-0.025	1.000	4477342		1.55(0.90-1.10)		24092	M
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.225	3.258	-0.033	1.000	49255849	225.4			955116	E
499.00 > 99.00	3.225	3.258	-0.033	1.000	11461351		4.30(0.90-1.10)		205350	E
D 17 13C4 PFOS										
503.00 > 80.00	3.225	3.259	-0.034		9623480	39.1		81.8	247467	

QC Flag Legend

Processing Flags

E - Exceeded Maximum Amount

Review Flags

M - Manually Integrated

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_007.d

Injection Date: 03-Dec-2016 16:48:42

Instrument ID: A8_N

Lims ID: 320-23783-A-13-A

Lab Sample ID: 320-23783-13

Client ID: DPT-16-15-GW-31-35

Operator ID: A8-PC\A8

ALS Bottle#: 38

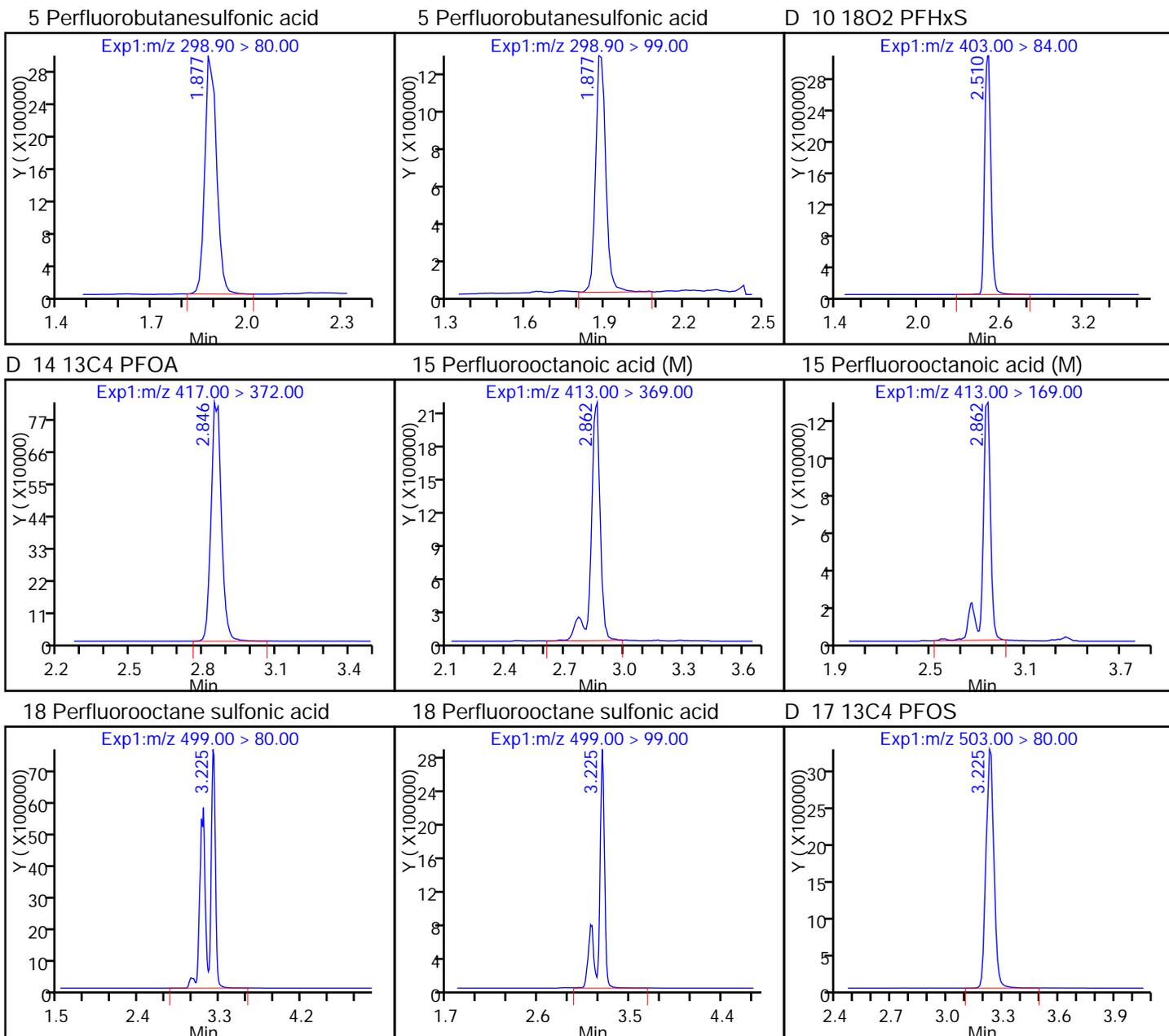
Worklist Smp#: 28

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



TestAmerica Sacramento

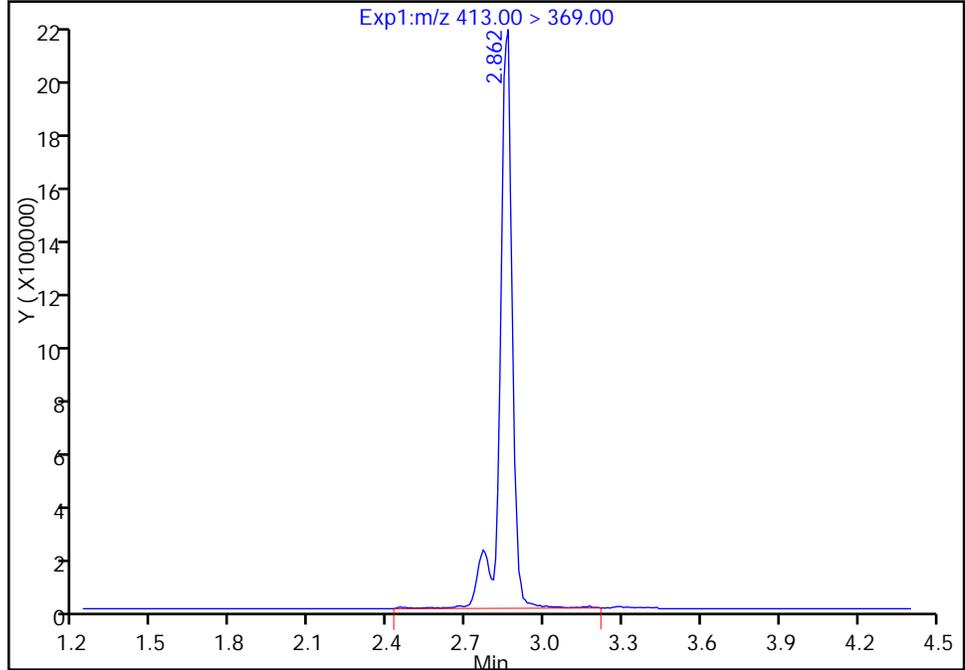
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_007.d
Injection Date: 03-Dec-2016 16:48:42 Instrument ID: A8_N
Lims ID: 320-23783-A-13-A Lab Sample ID: 320-23783-13
Client ID: DPT-16-15-GW-31-35
Operator ID: A8-PC\A8 ALS Bottle#: 38 Worklist Smp#: 28
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

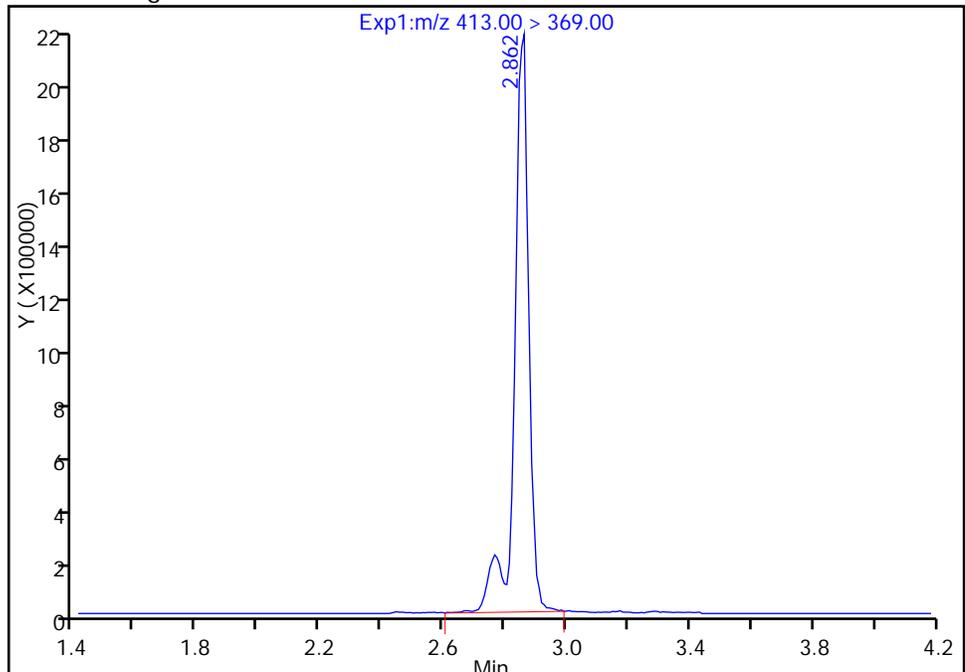
RT: 2.86
Area: 7098853
Amount: 134.6725
Amount Units: ng/ml

Processing Integration Results



RT: 2.86
Area: 6930971
Amount: 131.4876
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 10:36:41
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

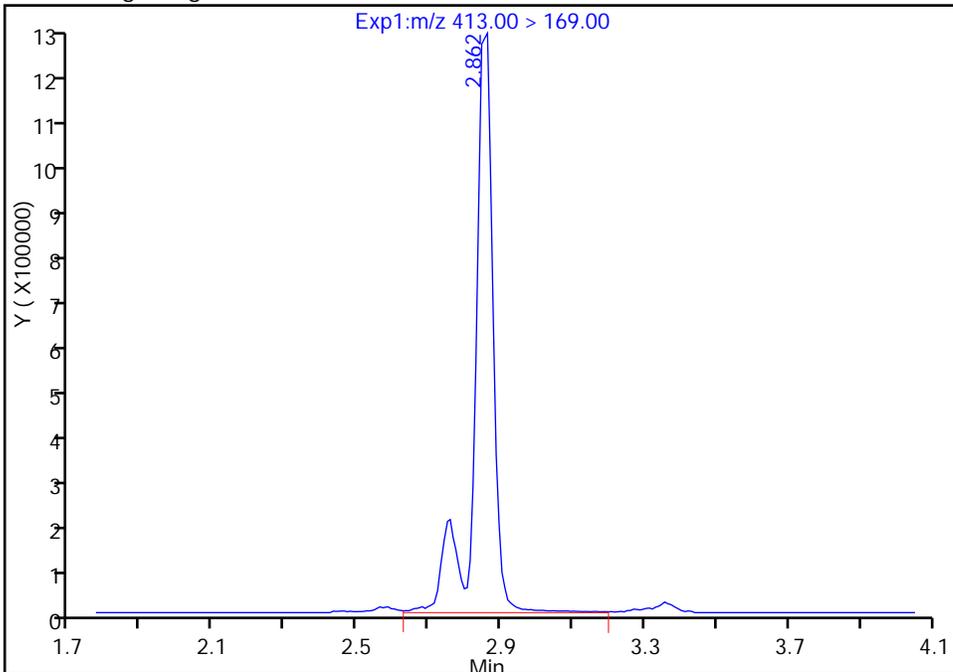
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_007.d
Injection Date: 03-Dec-2016 16:48:42 Instrument ID: A8_N
Lims ID: 320-23783-A-13-A Lab Sample ID: 320-23783-13
Client ID: DPT-16-15-GW-31-35
Operator ID: A8-PC\A8 ALS Bottle#: 38 Worklist Smp#: 28
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 2

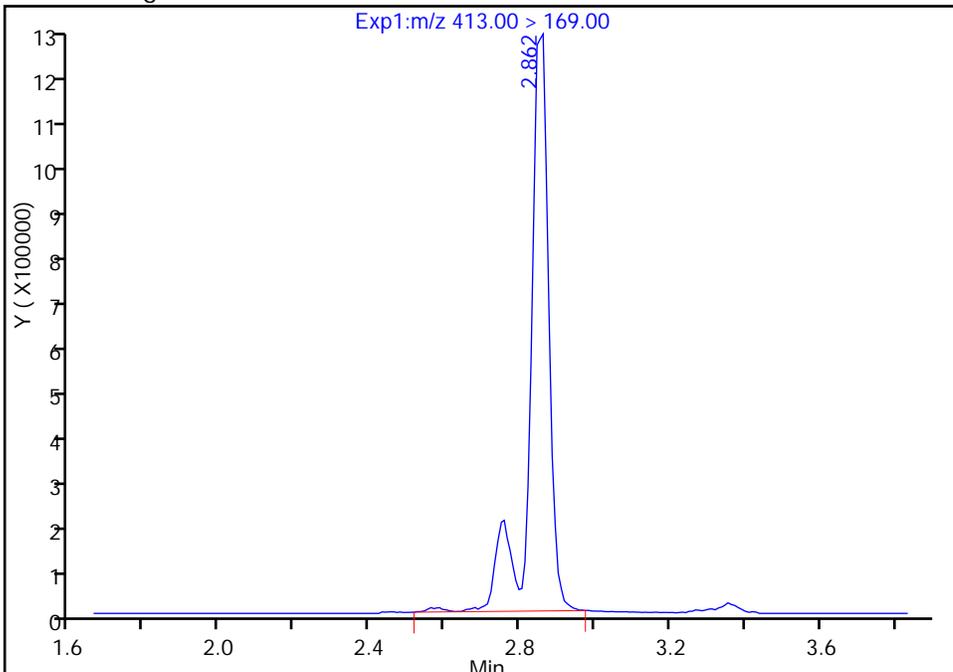
RT: 2.86
Area: 4593402
Amount: 134.6725
Amount Units: ng/ml

Processing Integration Results



RT: 2.86
Area: 4477342
Amount: 131.4876
Amount Units: ng/ml

Manual Integration Results



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: DPT-16-15-GW-31-35 DL Lab Sample ID: 320-23783-13 DL
 Matrix: Water Lab File ID: 07DEC2016B_008.d
 Analysis Method: 537 (Modified) Date Collected: 11/21/2016 14:00
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 245.9(mL) Date Analyzed: 12/07/2016 17:25
 Con. Extract Vol.: 0.5(mL) Dilution Factor: 10
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 141062 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	<i>Perfluorooctanoic acid (PFOA)</i>	0.27	D	0.025	0.0076
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.45	D	0.041	0.013
375-73-5	<i>Perfluorobutanesulfonic acid (PFBS)</i>	0.035	D	0.025	0.0093

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	33		25-150
STL00991	13C4 PFOS	95		25-150
STL00994	18O2 PFHxS	101		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_008.d
 Lims ID: 320-23783-A-13-A
 Client ID: DPT-16-15-GW-31-35
 Sample Type: Client
 Inject. Date: 07-Dec-2016 17:25:55 ALS Bottle#: 21 Worklist Smp#: 8
 Injection Vol: 2.0 ul Dil. Factor: 10.0000
 Sample Info: 320-23783-a-13-a 10X
 Misc. Info.: Plate: 1 Rack: 3
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:54 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:55:20

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.850	1.848	0.002	1.000	858416	1.71				
298.90 > 99.00	1.850	1.848	0.002	1.000	395855		2.17(0.00-0.00)			
D 10 18O2 PFHxS										
403.00 > 84.00	2.449	2.460	-0.011		1484791	4.75		10.1	181205	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.797	2.802	-0.005	1.000	1028052	13.1			14401	
413.00 > 169.00	2.797	2.802	-0.005	1.000	654698		1.57(0.90-1.10)		22473	
D 14 13C4 PFOA										
417.00 > 372.00	2.797	2.802	-0.005		365581	1.67		3.3	70562	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.166	3.061	0.105	1.000	5564937	21.9			511379	
499.00 > 99.00	3.047	3.061	-0.014	0.962	1233383		4.51(0.90-1.10)		24472	
D 17 13C4 PFOS										
503.00 > 80.00	3.166	3.169	-0.003		1117563	4.54		9.5	78635	

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_008.d

Injection Date: 07-Dec-2016 17:25:55

Instrument ID: A8_N

Lims ID: 320-23783-A-13-A

Lab Sample ID: 320-23783-13

Client ID: DPT-16-15-GW-31-35

Operator ID: A8-PC\A8

ALS Bottle#: 21

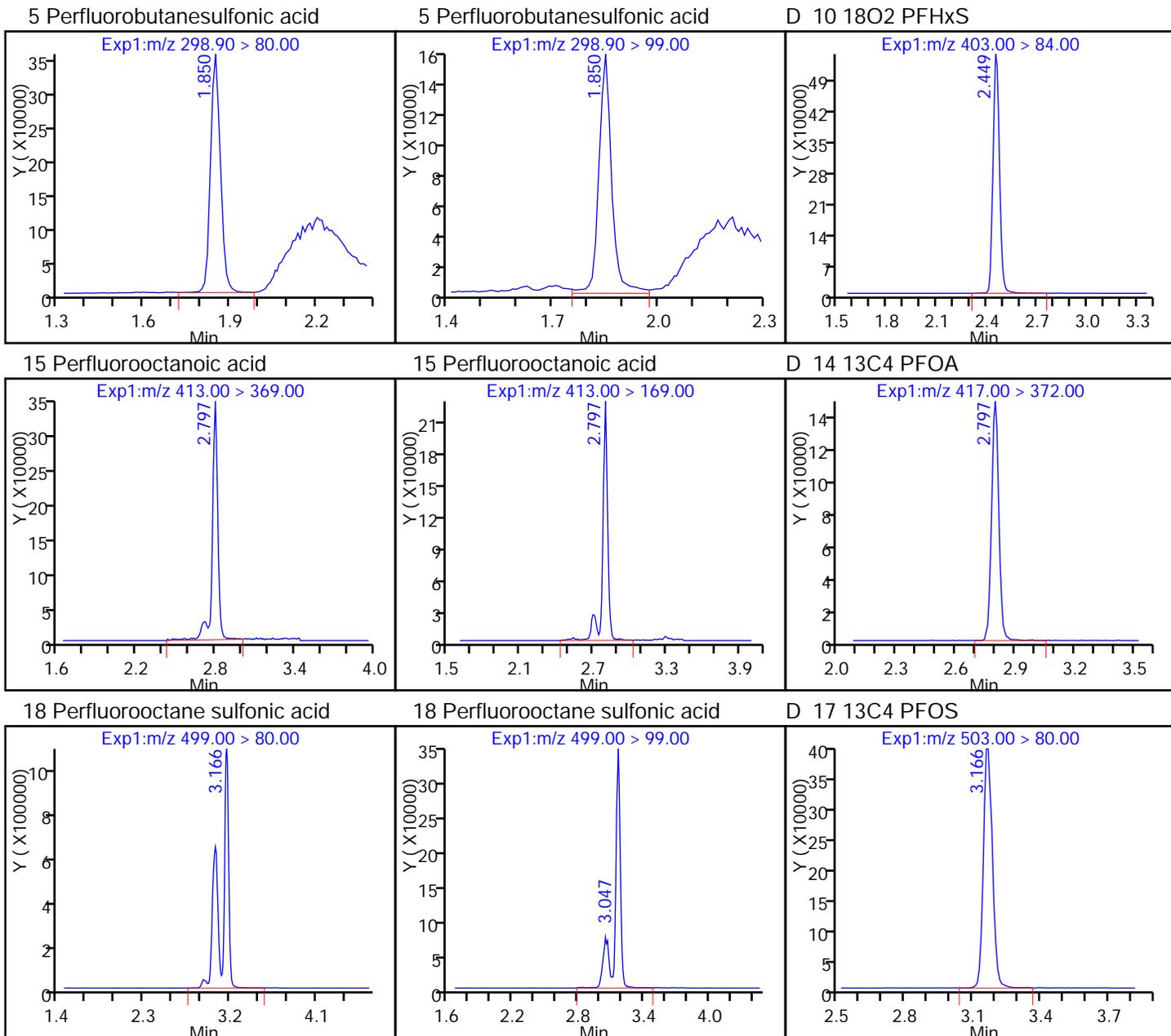
Worklist Smp#: 8

Injection Vol: 2.0 ul

Dil. Factor: 10.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL



FORM VI
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
 RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140382/4	02DEC2016A_004.d
Level 2	IC 320-140382/14	02DEC2016A_014.d
Level 3	IC 320-140382/5	02DEC2016A_005.d
Level 4	IC 320-140382/15	02DEC2016A_015.d
Level 5	IC 320-140382/6	02DEC2016A_006.d
Level 6	IC 320-140382/16	02DEC2016A_016.d
Level 7	IC 320-140382/7	02DEC2016A_007.d
Level 8	IC 320-140382/17	02DEC2016A_017.d
Level 9	IC 320-140382/8	02DEC2016A_008.d
Level 10	IC 320-140382/18	02DEC2016A_018.d
Level 11	IC 320-140382/9	02DEC2016A_009.d
Level 12	IC 320-140382/19	02DEC2016A_019.d
Level 13	IC 320-140382/10	02DEC2016A_010.d
Level 14	IC 320-140382/20	02DEC2016A_020.d

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12	LVL 13	LVL 14								
Perfluorobutanoic acid (PFBA)	1.630 1.610		1.622 ++++		1.622		1.614		1.614		1.367 - 1.867	1.619
Perfluoropentanoic acid (PFPeA)	1.944 1.910		1.916 ++++		1.926		1.916		1.916		1.670 - 2.170	1.921
Perfluorobutanesulfonic acid (PFBS)	1.983 1.949		1.964 ++++		1.964		1.954		1.945		1.781 - 2.141	1.960
Perfluorohexanoic acid (PFHxA)	2.281 2.225		2.239 ++++		2.247		2.239		2.226		1.991 - 2.491	2.243
Perfluoroheptanoic acid (PFHpA)	2.648 2.578		2.604 ++++		2.600		2.593		2.593		2.350 - 2.850	2.603
Perfluoroheptanesulfonic acid (PFHxS)	++++ 2.601		2.619 2.598		2.616		2.608		2.608		2.365 - 2.865	2.608
6:2FTS		++++ 2.904		2.920 ++++		2.910		2.914		2.919	2.663 - 3.163	2.913
Perfluorooctanoic acid (PFOA)	++++ 2.958		2.993 2.967		2.981		2.982		2.972		2.734 - 3.234	2.976
Perfluoroheptanesulfonic Acid (PFHpS)	3.032 2.966		2.993 ++++		2.989		2.982		2.972		2.736 - 3.236	2.989
Perfluorooctane Sulfonate (PFOS)	++++ 3.340		3.380 3.345		3.366		3.362		3.352		3.116 - 3.616	3.358
Perfluorononanoic acid (PFNA)	3.416 3.348		3.380 ++++		3.383		3.371		3.361		3.123 - 3.623	3.377
Perfluorooctane Sulfonamide (FOSA)	3.686 3.643		3.657 ++++		3.652		3.648		3.648		3.406 - 3.906	3.656
8:2FTS		3.668 3.667		3.684 ++++		3.673		3.678		3.674	3.424 - 3.924	3.674

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12	LVL 13	LVL 14								
Perfluorodecanoic acid (PFDA)	3.786 3.721		3.746 ++++		3.740		3.738		3.727		3.488 - 3.988	3.743
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)		3.851 3.832		3.857 3.849		3.847		3.843		3.847	3.597 - 4.097	3.847
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)		4.020 4.009		4.026 4.018		4.015		4.011		4.016	3.766 - 4.266	4.016
Perfluorodecanesulfonic acid (PFDS)	4.096 4.028		4.061 4.028		4.046		4.048		4.039		3.799 - 4.299	4.049
Perfluoroundecanoic acid (PFUnA)	4.120 4.054		4.086 ++++		4.071		4.065		4.056		3.821 - 4.321	4.075
MeFOSA		4.115 4.102		4.114 4.118		4.104		4.112		4.108	3.860 - 4.360	4.110
N-EtFOSA-M		4.293 4.296		4.301 4.309		4.297		4.302		4.295	4.049 - 4.549	4.299
Perfluorododecanoic acid (PFDoA)	4.420 4.343		4.381 ++++		4.376		4.365		4.354		4.120 - 4.620	4.373
Perfluorotridecanoic Acid (PFTriA)	4.681 4.613		4.653 ++++		4.637		4.631		4.629		4.386 - 4.886	4.641
Perfluorotetradecanoic acid (PFTeA)	4.935 4.864		4.893 ++++		4.888		4.883		4.867		4.634 - 5.134	4.888
Perfluoro-n-hexadecanoic acid (PFHxDA)	5.370 5.290		5.334 ++++		5.326		5.315		5.306		5.070 - 5.570	5.324
Perfluoro-n-octadecanoic acid (PFODA)	5.784 5.688		5.738 5.682		5.726		5.717		5.702		5.470 - 5.970	5.720
13C4 PFBA	1.630 1.602		1.614 1.609		1.622		1.614		1.614		1.365 - 1.865	1.615
13C5-PFPeA	1.944 1.910		1.916 ++++		1.916		1.916		1.916		1.668 - 2.168	1.920
13C2 PFHxA	2.281 2.225		2.239 ++++		2.239		2.231		2.226		1.989 - 2.489	2.240
13C4-PFHpA	2.641 2.578		2.604 ++++		2.600		2.593		2.593		2.349 - 2.849	2.602
18O2 PFHxS	2.656 2.594		2.619 2.598		2.616		2.608		2.608		2.364 - 2.864	2.614
M2-6:2FTS		2.925 2.912		2.912 2.917		2.918		2.906		2.911	2.665 - 3.165	2.914
13C4 PFOA	3.032 2.966		2.993 2.958		2.989		2.974		2.972		2.734 - 3.234	2.983
13C4 PFOS	3.416 3.340		3.371 3.345		3.366		3.362		3.352		3.115 - 3.615	3.365
13C5 PFNA	3.425 3.355		3.380 ++++		3.375		3.371		3.361		3.124 - 3.624	3.378
13C8 FOSA	3.686 3.634		3.648 ++++		3.652		3.648		3.648		3.401 - 3.901	3.653

FORM VI
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
 RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382
 SDG No.: _____
 Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N
 Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12	LVL 13	LVL 14								
M2-8:2FTS		3.678 3.667		3.684 3.676		3.673		3.678		3.674	3.426 - 3.926	3.676
13C2 PFDA	3.786 3.710		3.746 ++++		3.740		3.738		3.727		3.488 - 3.988	3.741
d3-NMeFOSAA		3.843 3.824		3.848 3.841		3.838		3.843		3.839	3.589 - 4.089	3.839
d5-NEtFOSAA		4.011 4.000		4.017 4.009		4.006		4.011		4.007	3.759 - 4.259	4.009
13C2 PFUnA	4.120 4.045		4.078 ++++		4.079		4.065		4.056		3.820 - 4.320	4.074
d-N-MeFOSA-M		4.108 4.102		4.106 4.110		4.104		4.104		4.100	3.855 - 4.355	4.105
d-N-EtFOSA-M		4.293 4.286		4.292 4.299		4.287		4.293		4.285	4.040 - 4.540	4.291
13C2 PFDoA	4.420 4.343		4.381 ++++		4.376		4.365		4.354		4.120 - 4.620	4.373
13C2-PFTeDA	4.935 4.855		4.893 4.860		4.880		4.883		4.867		4.632 - 5.132	4.882
13C2-PFHxDA	5.370 5.290		5.334 5.300		5.326		5.315		5.306		5.070 - 5.570	5.320

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140382/4	02DEC2016A_004.d
Level 2	IC 320-140382/14	02DEC2016A_014.d
Level 3	IC 320-140382/5	02DEC2016A_005.d
Level 4	IC 320-140382/15	02DEC2016A_015.d
Level 5	IC 320-140382/6	02DEC2016A_006.d
Level 6	IC 320-140382/16	02DEC2016A_016.d
Level 7	IC 320-140382/7	02DEC2016A_007.d
Level 8	IC 320-140382/17	02DEC2016A_017.d
Level 9	IC 320-140382/8	02DEC2016A_008.d
Level 10	IC 320-140382/18	02DEC2016A_018.d
Level 11	IC 320-140382/9	02DEC2016A_009.d
Level 12	IC 320-140382/19	02DEC2016A_019.d
Level 13	IC 320-140382/10	02DEC2016A_010.d
Level 14	IC 320-140382/20	02DEC2016A_020.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 5 LVL 9 LVL 13	LVL 2 LVL 6 LVL 10 LVL 14	LVL 3 LVL 7 LVL 11	LVL 4 LVL 8 LVL 12		B	M1	M2								
13C4 PFBA	315621 417432 329797 296657		338632 378720 305244		Ave		340300.229			12.7			50.0			
13C5-PFPeA	252783 331097 252000 ++++		264694 302516 218073		Ave		270193.930			14.9			50.0			
13C2 PFHxA	230991 294886 230866 ++++		246623 269968 205545		Ave		246479.820			12.9			50.0			
13C4-PFHpA	202776 256831 196027 ++++		215578 228495 163352		Ave		210510.050			15.0			50.0			
18O2 PFHxS	293951 366456 304899 239385		310687 347463 254328		Ave		302452.643			15.1			50.0			

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 5 LVL 9 LVL 13	LVL 2 LVL 6 LVL 10 LVL 14	LVL 3 LVL 7 LVL 11	LVL 4 LVL 8 LVL 12		B	M1	M2								
M2-6:2FTS		124739 140185 117602 164705		120558 166448 140694	Ave		139275.699			14.4		50.0				
13C4 PFOA	223843 294113 211325 144514		246973 252700 162409		Ave		219410.943			23.8		50.0				
13C4 PFOS	227748 298992 246737 190884		263387 280702 203865		Ave		244616.441			16.2		50.0				
13C5 PFNA	167208 216529 165500 ++++		188629 194611 136179		Ave		178109.173			15.7		50.0				
13C8 FOSA	366949 471533 404926 ++++		431176 451020 359915		Ave		414252.957			10.9		50.0				
M2-8:2FTS		117628 124671 122166 166617		108840 154273 140315	Ave		133501.446			15.7		50.0				
13C2 PFDA	151221 196057 159286 ++++		168367 182254 139218		Ave		166066.990			12.5		50.0				
d3-NMeFOSAA		77601 82804 68846 79183		72124 99067 80331	Ave		79993.7857			12.1		50.0				
d5-NETfOSAA		89291 89787 77393 89956		82682 104841 89500	Ave		89064.1286			9.5		50.0				

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 5 LVL 9 LVL 13	LVL 2 LVL 6 LVL 10 LVL 14	LVL 3 LVL 7 LVL 11	LVL 4 LVL 8 LVL 12		B	M1	M2								
13C2 PFUnA	122340 154918 114588 ++++		133394 138267 91167		Ave		125779.243			17.4			50.0			
d-N-MeFOSA-M		93620 107246 99342 124175		103638 117800 98551	Ave		106338.894			10.4			50.0			
d-N-EtFOSA-M		86581 104104 97144 118377		100979 112224 93938	Ave		101906.806			10.6			50.0			
13C2 PFDoA	108677 140992 109985 ++++		119003 126396 96377		Ave		116904.873			13.3			50.0			
13C2-PFTeDA	221880 299085 231293 207930		259139 271923 202045		Ave		241899.200			14.8			50.0			
13C2-PFHxDA	113226 160713 124217 120034		135127 148580 110713		Ave		130372.880			14.3			50.0			

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_NGC Column: Acquity ID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/02/2016 10:29Calibration End Date: 12/02/2016 12:29Calibration ID: 26849

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10												
Perfluorobutanoic acid (PFBA)	287036		298633		387562	AveID		0.8853			10.0		35.0				
	219376	371759	++++	294183													
Perfluoropentanoic acid (PFPeA)	321460		290628		336046	AveID		1.0408			15.3		35.0				
	171488	327954	++++	249486													
Perfluorobutanesulfonic acid (PFBS)	481176		500216		668327	AveID		1.6135			15.6		50.0				
	291824	642971	++++	491581													
Perfluorohexanoic acid (PFHxA)	235306		236370		291709	AveID		0.9559			8.0		35.0				
	166899	274051	++++	217418													
Perfluoroheptanoic acid (PFHpA)	233630		216223		259601	AveID		1.0342			8.4		35.0				
	146715	250282	++++	204959													
Perfluorohexanesulfonic acid (PFHxS)	++++		376003		406797	AveID		1.0596			10.2		35.0				
	251770	380131		323081													
6:2FTS		++++		93245		AveID		0.8391			14.3		35.0				
	103949	105654	152954	++++	118849												
Perfluorooctanoic acid (PFOA)	++++		290666		312182	AveID		1.0203			12.7		35.0				
	148950	280154	118726		218968												
Perfluoroheptanesulfonic Acid (PFHpS)	277206		299283		372775	AveID		1.1690			9.7		50.0				
	199679	364908	++++	279830													
Perfluorooctane Sulfonate (PFOS)	++++		275533		316788	AveID		1.0873			3.0		35.0				
	225419	309544	216190		265826												
Perfluorononanoic acid (PFNA)	170148		184877		217102	AveID		0.9912			3.7		35.0				
	126022	200322	++++	164187													
Perfluorooctane Sulfonamide (FOSA)	360690		403791		474054	AveID		0.9353			12.5		35.0				
	255544	468609	++++	379917													

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_NGC Column: Acquity ID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/02/2016 10:29Calibration End Date: 12/02/2016 12:29Calibration ID: 26849

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10												
8:2FTS	90169	91935	146110	98719	120115	AveID	0.8444				13.7		35.0				
		101586		++++													
Perfluorodecanoic acid (PFDA)	155778	183568	159778	151623	187771	AveID	0.9693				4.2		35.0				
	128074		++++														
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)	57976	61790	91493	59295	73940	AveID	0.8686				13.5		35.0				
		68867	71812														
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)	56792	60032	94072	56004	70934	AveID	0.7671				14.6		35.0				
		70327	70891														
Perfluorodecanesulfonic acid (PFDS)	138425	156537	162158	201706		AveID	0.6271				6.9		50.0				
	126426	189172	107269														
Perfluoroundecanoic acid (PFUnA)	152116	148792	111549	152861		AveID	1.0596				9.9		35.0				
	89994	145410	++++														
MeFOSA	72052	74554	106455	73327	96734	AveID	0.7963				13.8		35.0				
		79073	89254														
N-EtFOSA-M	73147	71070	109355	78846	100899	AveID	0.8497				14.4		35.0				
		82751	88734														
Perfluorododecanoic acid (PFDoA)	109058	109023	101710	133861		AveID	0.9465				3.6		35.0				
	88491	122257	++++														
Perfluorotridecanoic Acid (PFTriA)	109674	126175	114082	147998		AveID	1.0402				5.1		50.0				
	92748	141902	++++														
Perfluorotetradecanoic acid (PFTeA)	228612	237250	206027	270118		AveID	1.9323				8.3		50.0				
	159620	259265	++++														
Perfluoro-n-hexadecanoic acid (PFHxDA)	267566	218894	113089	160135		LlID	0.8109	0.9850						0.9990		0.9900	
	93727	142551	++++														
Perfluoro-n-octadecanoic acid (PFODA)	80380	118037	107852	105867		AveID	0.7863				19.0		50.0				
	68360	72886	77747														

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
 RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140382/4	02DEC2016A_004.d
Level 2	IC 320-140382/14	02DEC2016A_014.d
Level 3	IC 320-140382/5	02DEC2016A_005.d
Level 4	IC 320-140382/15	02DEC2016A_015.d
Level 5	IC 320-140382/6	02DEC2016A_006.d
Level 6	IC 320-140382/16	02DEC2016A_016.d
Level 7	IC 320-140382/7	02DEC2016A_007.d
Level 8	IC 320-140382/17	02DEC2016A_017.d
Level 9	IC 320-140382/8	02DEC2016A_008.d
Level 10	IC 320-140382/18	02DEC2016A_018.d
Level 11	IC 320-140382/9	02DEC2016A_009.d
Level 12	IC 320-140382/19	02DEC2016A_019.d
Level 13	IC 320-140382/10	02DEC2016A_010.d
Level 14	IC 320-140382/20	02DEC2016A_020.d

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
		LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
13C4 PFBA	Ave	15781027	18935985	16931601	16489860	20871580	50.0	50.0	50.0	50.0	50.0
		15262180		14832847		50.0	50.0				
		12639170		13234715		16554836	50.0		50.0		50.0
13C5-PFPeA	Ave	10903650	15125796	+++++	12600012	16554836	50.0	50.0	50.0	50.0	50.0
		11549542		12331127		14744313	50.0		50.0		50.0
		10277273		11543313		12841573	50.0		50.0		50.0
13C2 PFHxA	Ave	10138817	11424765	10778897	9801352	12841573	50.0	50.0	50.0	50.0	50.0
		8167611		+++++		50.0	+++++		50.0		
		13903865		14695505		17333346	47.3		47.3		47.3
12029706	16435009	14421709	47.3	47.3	47.3						
5925112	7906261	5726516	47.5	47.5	47.5	47.5					
6658764	6682949	5586093	47.5		47.5	47.5					
11192131	12348635	14705649	50.0		50.0	50.0	50.0				
8120442	7225722	10566267	50.0	50.0		50.0					
12634984	12348635	14705649	50.0	50.0		50.0					

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
		LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
		LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	
13C4 PFOS	Ave	10886366	13417569	12589922	11794026	14291797	47.8	47.8	47.8	47.8	47.8
		9744748		9124233			47.8		47.8		
13C5 PFNA	Ave	8360391	9730546	9431435	8274989	10826461	50.0	50.0	50.0	50.0	50.0
		6808930		+++++			50.0		+++++		
13C8 FOSA	Ave	18347473	22550983	21558794	20246280	23576629	50.0	50.0	50.0	50.0	50.0
		17995728		+++++			50.0		+++++		
M2-8:2FTS	Ave	5971752	5634375	7389674	5213436	5851766	47.9	47.9	47.9	47.9	47.9
			6721067		7980965		47.9		47.9		
13C2 PFDA	Ave	7561029	9112679	8418346	7964307	9802839	50.0	50.0	50.0	50.0	50.0
		6960897		+++++			50.0		+++++		
d3-NMeFOSAA	Ave	4140205	3880040	4953346	3606216	3442292	50.0	50.0	50.0	50.0	50.0
			4016571		3959155		50.0		50.0		
d5-NEtFOSAA	Ave	4489327	4464530	5242030	4497788	3869672	50.0	50.0	50.0	50.0	50.0
			4474975				50.0		50.0		
13C2 PFUnA	Ave	6117015	6913372	6669716	5729423	7745917	50.0	50.0	50.0	50.0	50.0
		4558330		+++++			50.0		+++++		
d-N-MeFOSA-M	Ave	5362314	4681022	5889991	5181881	4967092	50.0	50.0	50.0	50.0	50.0
			4927555		6208758		50.0		50.0		
d-N-EtFOSA-M	Ave	5205212	4329074	5611197	5048951	4857179	50.0	50.0	50.0	50.0	50.0
			4696896		5918873		50.0		50.0		
13C2 PFDoA	Ave	5433841	6319787	5950161	5499233	7049582	50.0	50.0	50.0	50.0	50.0
		4818858		+++++			50.0		+++++		
13C2-PFtEDA	Ave	11093975	13596161	12956938	11564641	14954236	50.0	50.0	50.0	50.0	50.0
		10102255		10396514			50.0		50.0		
13C2-PFHxDA	Ave	5661315	7428976	6756331	6210852	8035665	50.0	50.0	50.0	50.0	50.0
		5535655		6001714			50.0		50.0		

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140382
SDG No.: _____
Instrument ID: A8_N GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N
Calibration Start Date: 12/02/2016 10:29 Calibration End Date: 12/02/2016 12:29 Calibration ID: 26849

Curve Type Legend:

Ave = Average

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_NGC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/02/2016 10:29Calibration End Date: 12/02/2016 12:29Calibration ID: 26849

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140382/4	02DEC2016A_004.d
Level 2	IC 320-140382/14	02DEC2016A_014.d
Level 3	IC 320-140382/5	02DEC2016A_005.d
Level 4	IC 320-140382/15	02DEC2016A_015.d
Level 5	IC 320-140382/6	02DEC2016A_006.d
Level 6	IC 320-140382/16	02DEC2016A_016.d
Level 7	IC 320-140382/7	02DEC2016A_007.d
Level 8	IC 320-140382/17	02DEC2016A_017.d
Level 9	IC 320-140382/8	02DEC2016A_008.d
Level 10	IC 320-140382/18	02DEC2016A_018.d
Level 11	IC 320-140382/9	02DEC2016A_009.d
Level 12	IC 320-140382/19	02DEC2016A_019.d
Level 13	IC 320-140382/10	02DEC2016A_010.d
Level 14	IC 320-140382/20	02DEC2016A_020.d

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
Perfluorobutanoic acid (PFBA)		AveID	143518		298633		1937809	0.500		1.00		5.00
			43875101	7435187	++++	14709131	200	20.0	++++	50.0		
Perfluoropentanoic acid (PFPeA)		AveID	160730		290628		1680231	0.500		1.00		5.00
			34297516	6559082	++++	12474289	200	20.0	++++	50.0		
Perfluorobutanesulfonic acid (PFBS)		AveID	212680		442191		2954006	0.442		0.884		4.42
			51594571	11367724	++++	21727893	177	17.7	++++	44.2		
Perfluorohexanoic acid (PFHxA)		AveID	117653		236370		1458544	0.500		1.00		5.00
			33379792	5481010	++++	10870881	200	20.0	++++	50.0		
Perfluoroheptanoic acid (PFHpA)		AveID	116815		216223		1298003	0.500		1.00		5.00
			29343029	5005633	++++	10247953	200	20.0	++++	50.0		
Perfluorohexanesulfonic acid (PFHxS)		AveID	++++		342163		1850926	++++		0.910		4.55
			45822208	6918390	77870104	14700167	182	18.2	364	45.5		
6:2FTS		AveID	492719	++++	2900008		88396	4.74	++++	19.0		0.948
				20031946			5633441	190	190	++++	47.4	

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_NGC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/02/2016 10:29Calibration End Date: 12/02/2016 12:29Calibration ID: 26849

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
Perfluorooctanoic acid (PFOA)		AveID	+++++	5603089	290666	10948383	1560911	+++++	20.0	1.00	50.0	5.00
			29789966		47490545			200		400		
Perfluoroheptanesulfonic Acid (PFHpS)		AveID	131950	6947857	284917	13319919	1774408	0.476	19.0	0.952	47.6	4.76
			38018954		+++++			190		+++++		
Perfluorooctane Sulfonate (PFOS)		AveID	+++++	5745134	255695	12334342	1469896	+++++	18.6	0.928	46.4	4.64
			41837729		80249842			186		371		
Perfluorononanoic acid (PFNA)		AveID	85074	4006430	184877	8209339	1085508	0.500	20.0	1.00	50.0	5.00
			25204338		+++++			200		+++++		
Perfluorooctane Sulfonamide (FOSA)		AveID	180345	9372172	403791	18995833	2370271	0.500	20.0	1.00	50.0	5.00
			51108754		+++++			200		+++++		
8:2FTS		AveID	431909	44037	2799477	94573	5753486	4.79	0.479	19.2	0.958	47.9
				19463965		+++++			192		+++++	
Perfluorodecanoic acid (PFDA)		AveID	77889	3671358	159778	7581125	938853	0.500	20.0	1.00	50.0	5.00
			25614867		+++++			200		+++++		
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)		AveID	289882	30895	1829869	59295	3696988	5.00	0.500	20.0	1.00	50.0
				13773335		28724886		200		400		
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)		AveID	283960	30016	1881436	56004	3546696	5.00	0.500	20.0	1.00	50.0
				14065497		28356327		200		400		
Perfluorodecanesulfonic acid (PFDS)		AveID	66721	3647236	150902	7816007	972225	0.482	19.3	0.964	48.2	4.82
			24374906		41362935			193		386		
Perfluoroundecanoic acid (PFUnA)		AveID	76058	2908196	148792	5577456	764307	0.500	20.0	1.00	50.0	5.00
			17998762		+++++			200		+++++		
MeFOSA		AveID	360259	37277	2129108	73327	4836683	5.00	0.500	20.0	1.00	50.0
				15814534		35701676		200		400		
N-EtFOSA-M		AveID	365734	35535	2187091	78846	5044937	5.00	0.500	20.0	1.00	50.0
				16550291		35493503			200		400	

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140382

SDG No.: _____

Instrument ID: A8_NGC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/02/2016 10:29Calibration End Date: 12/02/2016 12:29Calibration ID: 26849

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
Perfluorododecanoic acid (PFDoA)		AveID	54529		109023		669306	0.500		1.00		5.00
			17698276	2445148	++++	5085511		200	20.0	++++	50.0	
			LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	
Perfluorotridecanoic Acid (PFTriA)		AveID	54837		126175		739992	0.500		1.00		5.00
			18549592	2838032	++++	5704096		200	20.0	++++	50.0	
			LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	
Perfluorotetradecanoic acid (PFTeA)		AveID	114306		237250		1350591	0.500		1.00		5.00
			31923969	5185293	++++	10301361		200	20.0	++++	50.0	
			LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	
Perfluoro-n-hexadecanoic acid (PFHxDA)		L1ID	133783		218894		800676	0.500		1.00		5.00
			18745470	2851018	++++	5654463		200	20.0	++++	50.0	
			LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	
Perfluoro-n-octadecanoic acid (PFODA)		AveID	40190		118037		529337	0.500		1.00		5.00
			13671969	1457725		5392609		200	20.0		50.0	
			LVL 11	LVL 12	LVL 13	LVL 14	LVL 11	LVL 12	LVL 13	LVL 14	LVL 10	

Curve Type Legend:

AveID = Average isotope dilution
L1ID = Linear 1/conc IsoDil

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_004.d
 Lims ID: IC L1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 02-Dec-2016 10:29:38 ALS Bottle#: 37 Worklist Smp#: 4
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L1_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:29 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:37:48

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.630	1.615	0.015	15781027	46.4		92.7	1736709	
1 Perfluorobutyric acid	212.90 > 169.00	1.630	1.617	0.013	143518	0.5136		103	1140	
D 4 13C5-PFPeA	267.90 > 223.00	1.944	1.918	0.026	12639170	46.8		93.6	942300	
3 Perfluoropentanoic acid	262.90 > 219.00	1.944	1.920	0.024	160730	0.6109		122	1551	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.983	1.961	0.022	212680	0.4484		101		
	298.90 > 99.00	1.983	1.961	0.022	94809		2.24(0.00-0.00)	101		
D 6 13C2 PFHxA	315.00 > 270.00	2.281	2.239	0.042	11549542	46.9		93.7	865664	
7 Perfluorohexanoic acid	313.00 > 269.00	2.281	2.241	0.040	117653	0.5329		107	3266	
D 11 13C4-PFHpA	367.00 > 322.00	2.641	2.599	0.042	10138817	48.2		96.3	926585	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.648	2.600	0.048	116815	0.5570		111	2490	
D 10 18O2 PFHxS	403.00 > 84.00	2.656	2.614	0.042	13903865	46.0		97.2	1297944	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.656	2.615	0.041	196748	0.6317		139		
15 Perfluorooctanoic acid	413.00 > 369.00	3.032	2.984	0.048	158739	0.6950		139	1573	
	413.00 > 169.00	3.032	2.984	0.048	89616		1.77(0.90-1.10)	139	8158	

D 14 13C4 PFOA

417.00 > 372.00 3.032 2.984 0.048 11192131 51.0 102 1037140
 Page 505 of 834 12/09/2016

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	3.032	2.986	0.046	1.000	131950	0.4956	104		
D 17 13C4 PFOS	503.00	> 80.00	3.416	3.365	0.051		10886366	44.5	93.1	481501	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.416	3.366	0.050	1.000	120361	0.4860	105	16520	
	499.00	> 99.00	3.425	3.366	0.059	1.003	28366	4.24(0.90-1.10)	105	2029	
20 Perfluorononanoic acid	463.00	> 419.00	3.416	3.373	0.043	1.000	85074	0.5133	103	1441	
D 19 13C5 PFNA	468.00	> 423.00	3.425	3.374	0.051		8360391	46.9	93.9	601596	
D 21 13C8 FOSA	506.00	> 78.00	3.686	3.651	0.035		18347473	44.3	88.6	659525	
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.686	3.656	0.030	1.000	180345	0.5254	105	13418	
24 Perfluorodecanoic acid	513.00	> 469.00	3.786	3.738	0.048	1.000	77889	0.5314	106	2050	
D 23 13C2 PFDA	515.00	> 470.00	3.786	3.738	0.048		7561029	45.5	91.1	167715	
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	4.096	4.049	0.047	1.000	66721	0.4671	96.9		
D 27 13C2 PFUnA	565.00	> 520.00	4.120	4.070	0.050		6117015	48.6	97.3	398481	
28 Perfluoroundecanoic acid	563.00	> 519.00	4.120	4.071	0.049	1.000	76058	0.5867	117	2233	
D 30 13C2 PFDoA	615.00	> 570.00	4.420	4.370	0.050		5433841	46.5	93.0	245704	
29 Perfluorododecanoic acid	613.00	> 569.00	4.420	4.370	0.050	1.000	54529	0.5301	106	74.3	
31 Perfluorotridecanoic acid	663.00	> 619.00	4.681	4.636	0.045	1.000	54837	0.4851	97.0	81.1	
D 32 13C2-PFTeDA	715.00	> 670.00	4.935	4.882	0.053		11093975	45.9	91.7	599425	
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.935	4.884	0.051	1.000	114306	0.5443	109	74.0	
	713.00	> 169.00	4.918	4.884	0.034	0.997	21224	5.39(0.00-0.00)	109	2088	M
D 34 13C2-PFHxDA	815.00	> 770.00	5.370	5.320	0.050		5661315	43.4	86.8	129845	
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.370	5.320	0.050	1.000	133783	0.4265	85.3	239	
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.784	5.720	0.064	1.000	40190	0.4703	94.1	105	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L1_00021

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_004.d

Injection Date: 02-Dec-2016 10:29:38

Instrument ID: A8_N

Lims ID: IC L1

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 37

Worklist Smp#: 4

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

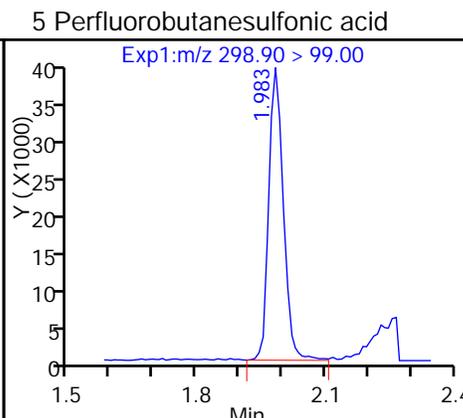
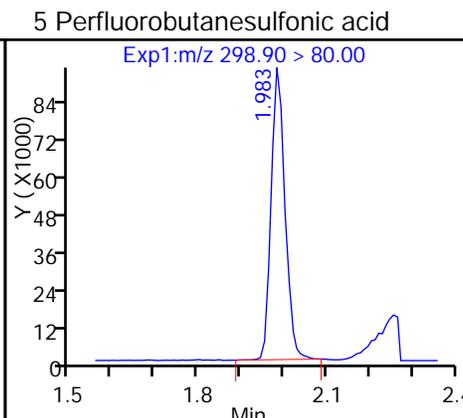
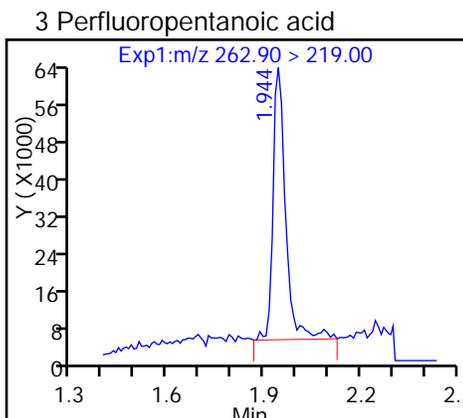
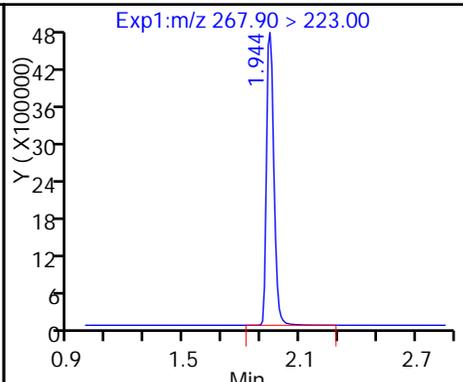
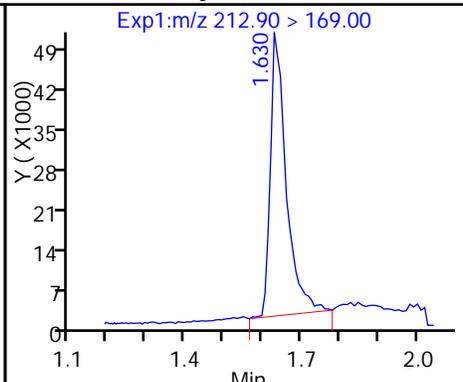
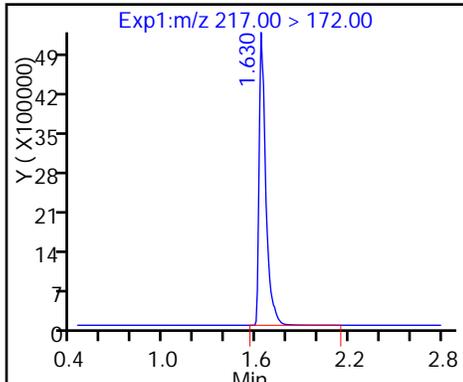
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

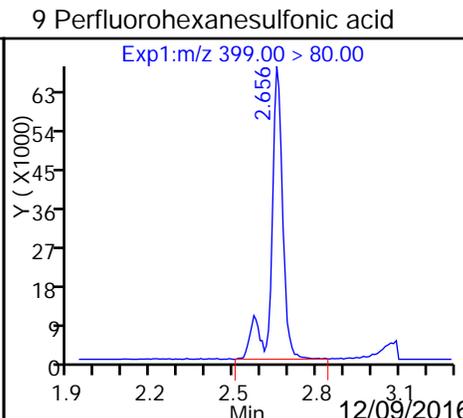
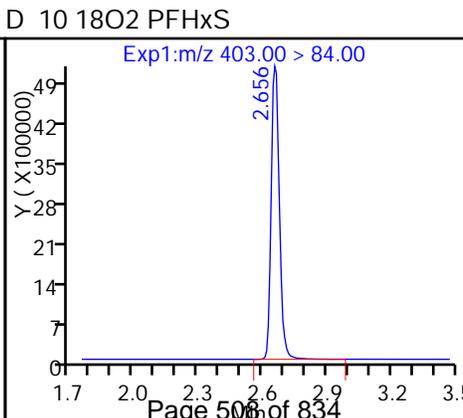
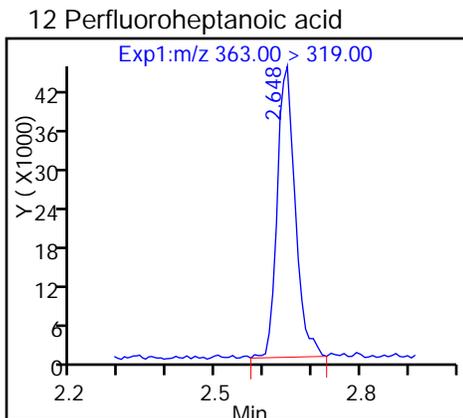
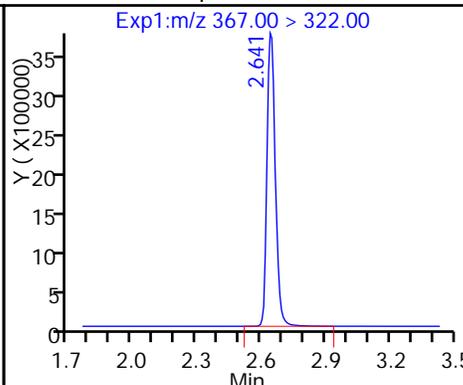
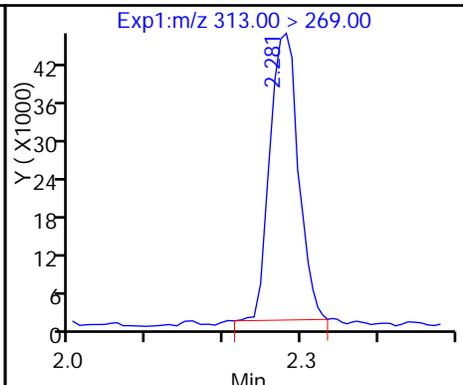
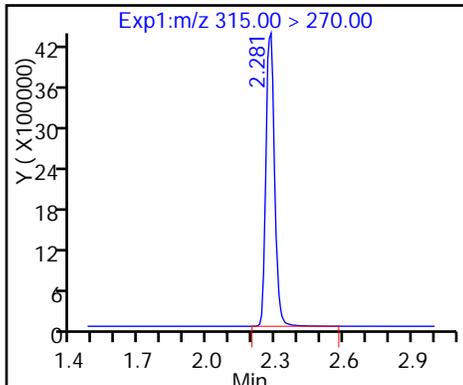
D 4 13C5-PFPeA

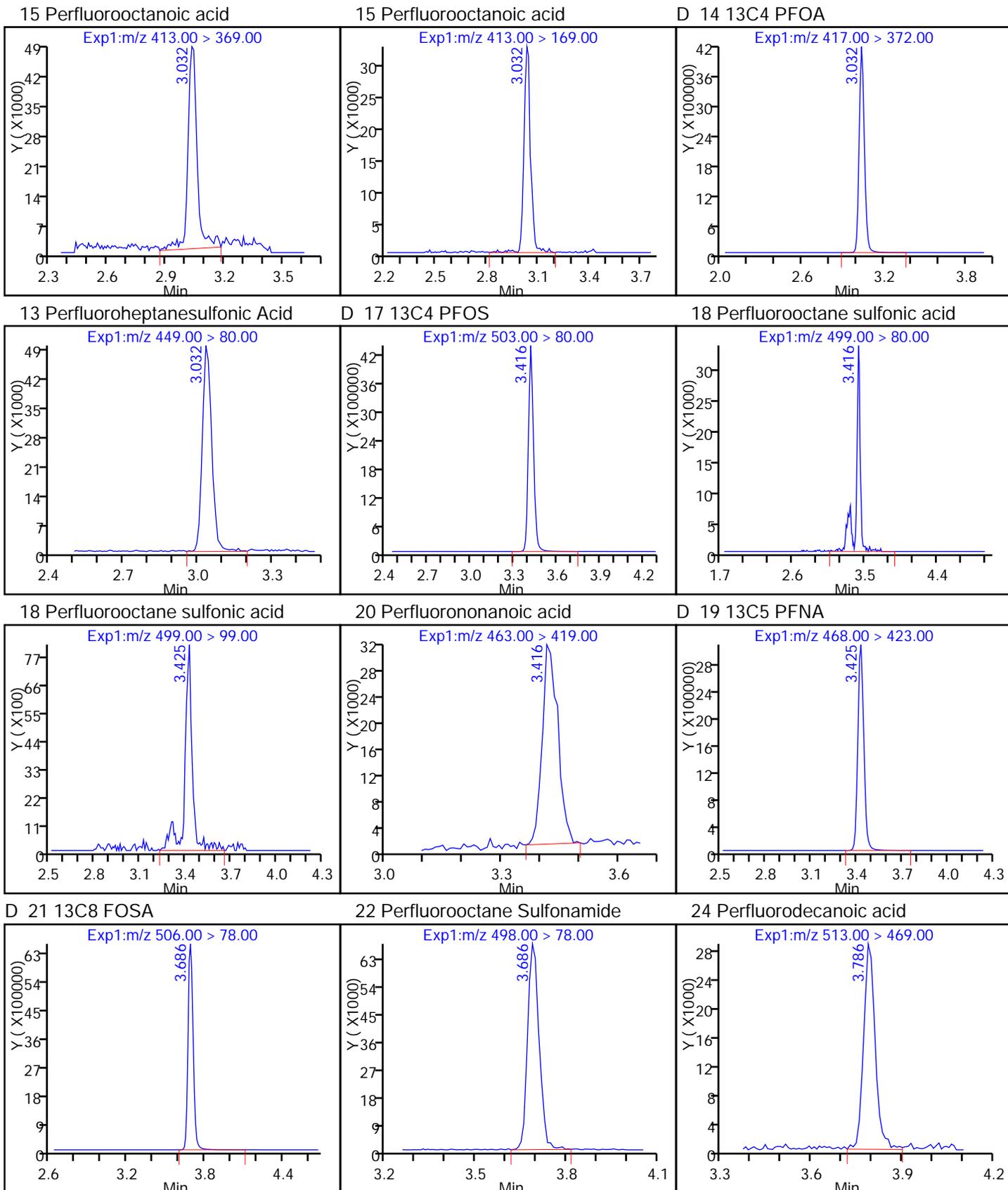


D 6 13C2 PFHxA

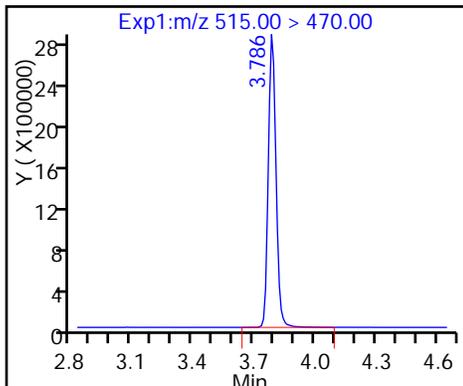
7 Perfluorohexanoic acid

D 11 13C4-PFHpA

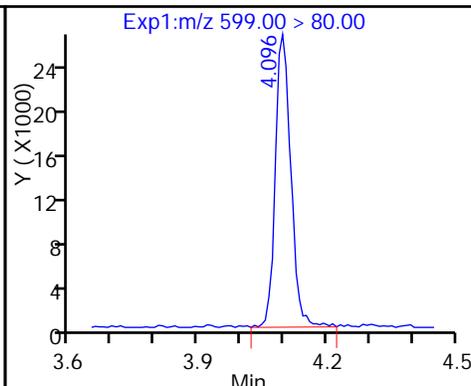




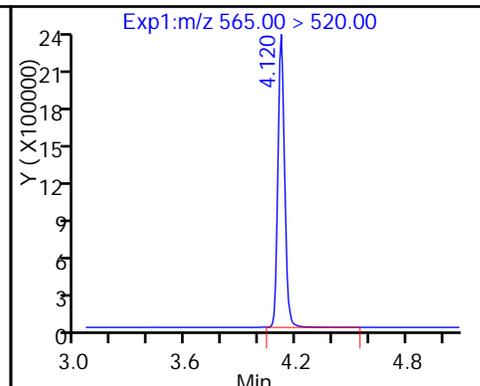
D 23 13C2 PFDA



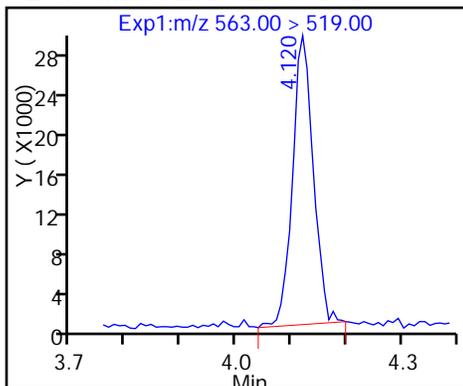
26 Perfluorodecane Sulfonic acid



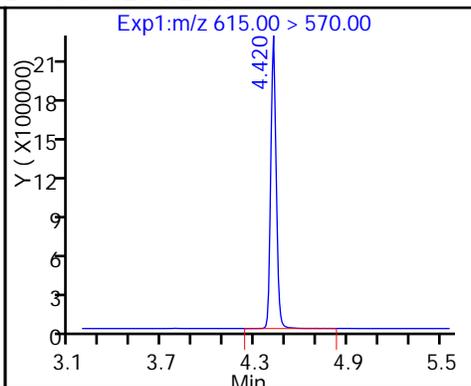
D 27 13C2 PFUnA



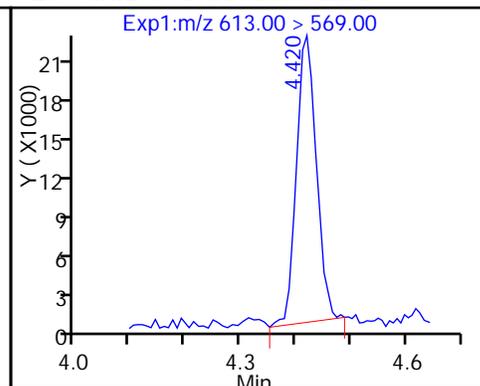
28 Perfluoroundecanoic acid



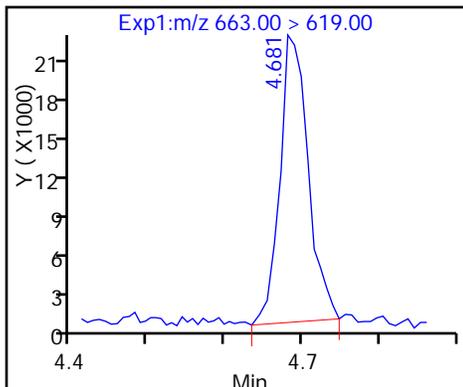
D 30 13C2 PFDaA



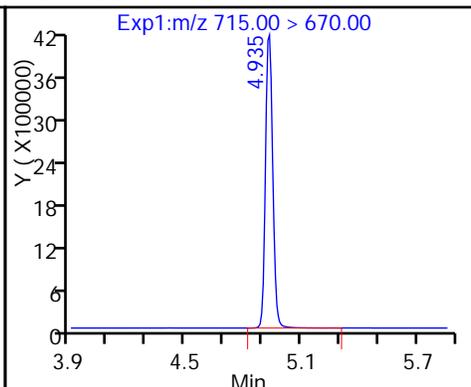
29 Perfluorododecanoic acid



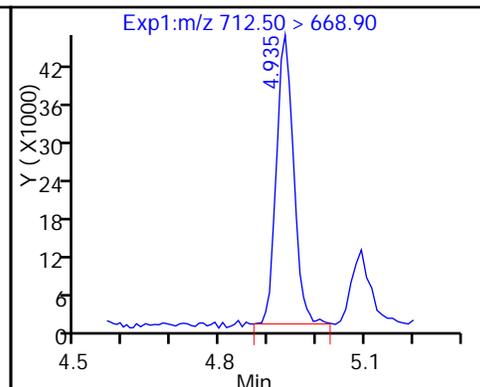
31 Perfluorotridecanoic acid



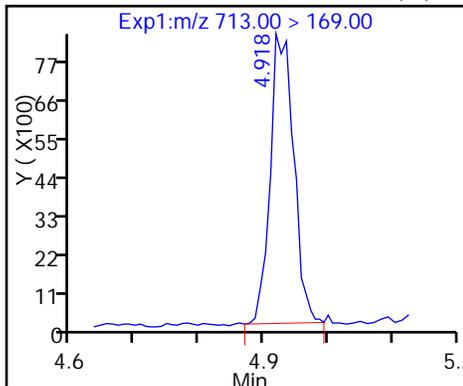
D 32 13C2-PFTeDA



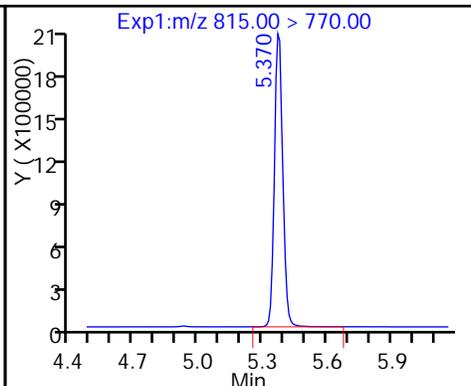
33 Perfluorotetradecanoic acid



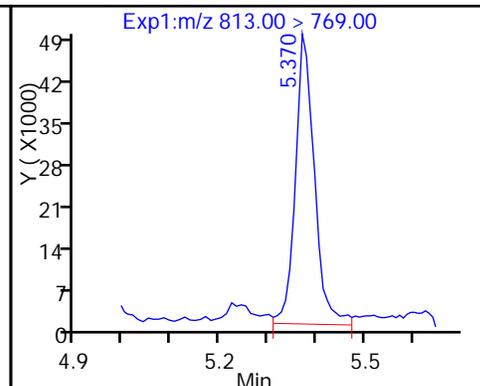
33 Perfluorotetradecanoic acid (M)



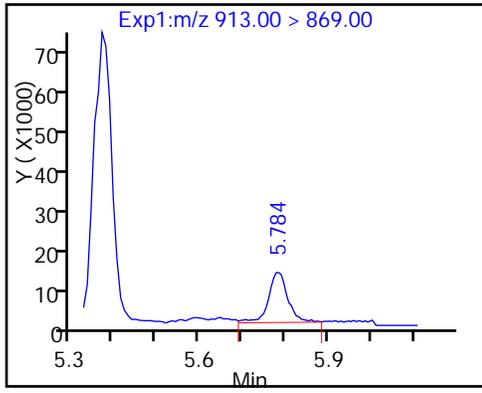
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento

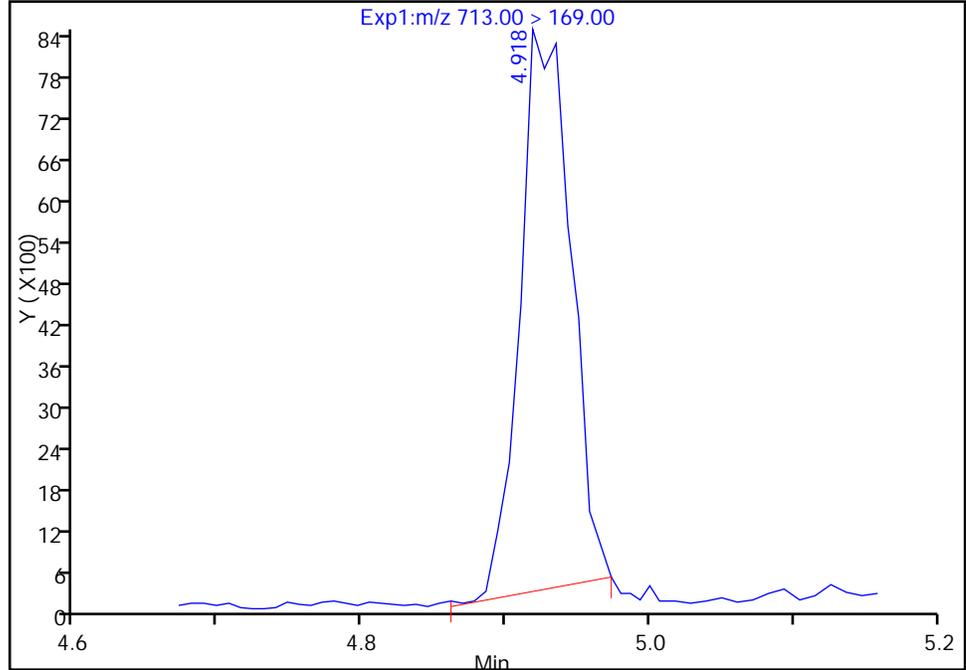
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_004.d
Injection Date: 02-Dec-2016 10:29:38 Instrument ID: A8_N
Lims ID: IC L1
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 4
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

33 Perfluorotetradecanoic acid, CAS: 376-06-7

Signal: 2

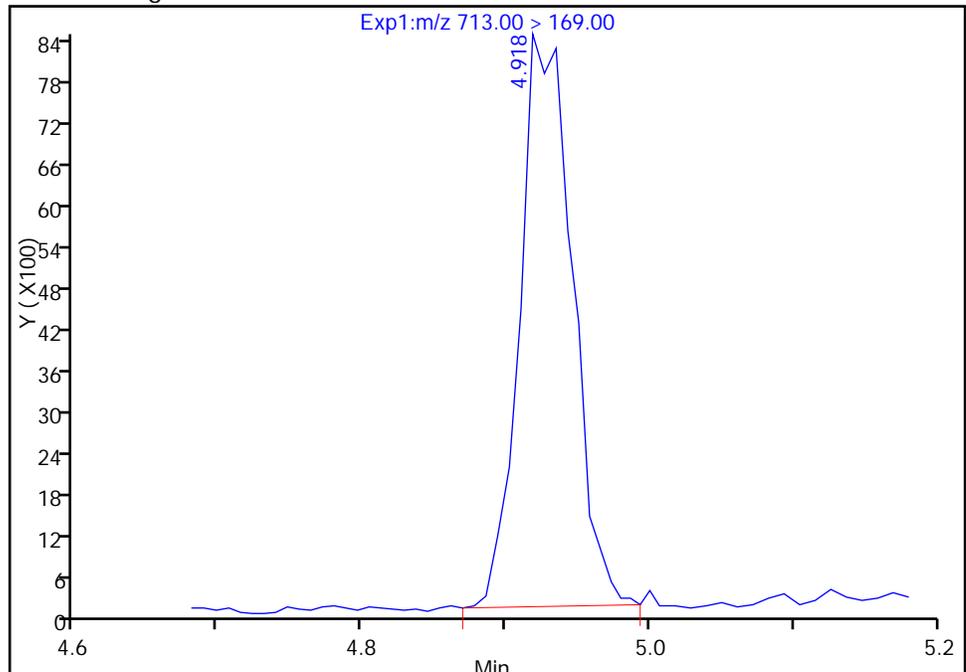
RT: 4.92
Area: 20090
Amount: 0.574454
Amount Units: ng/ml

Processing Integration Results



RT: 4.92
Area: 21224
Amount: 0.544327
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 02-Dec-2016 13:37:48
Audit Action: Manually Integrated

Audit Reason: Incomplete Integration

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_005.d
 Lims ID: IC L2
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 02-Dec-2016 10:37:07 ALS Bottle#: 38 Worklist Smp#: 5
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L2_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:31 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:38:35

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.614	1.615	-0.001	16931601	49.8		99.5	861260	
1 Perfluorobutyric acid	212.90 > 169.00	1.622	1.617	0.005	298633	1.00		99.6	2205	
D 4 13C5-PFPeA	267.90 > 223.00	1.916	1.918	-0.002	13234715	49.0		98.0	1668802	
3 Perfluoropentanoic acid	262.90 > 219.00	1.916	1.920	-0.004	290628	1.05		105	2905	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.964	1.961	0.003	442191	0.8821		99.8		
	298.90 > 99.00	1.954	1.961	-0.007	183888		2.40(0.00-0.00)	99.8		
D 6 13C2 PFHxA	315.00 > 270.00	2.239	2.239	0.0	12331127	50.0		100	1160245	
7 Perfluorohexanoic acid	313.00 > 269.00	2.239	2.241	-0.002	236370	1.00		100	7007	
D 11 13C4-PFHpA	367.00 > 322.00	2.604	2.599	0.005	10778897	51.2		102	1033732	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.604	2.600	0.004	216223	0.9699		97.0	3288	
D 10 18O2 PFHxS	403.00 > 84.00	2.619	2.614	0.005	14695505	48.6		103	816827	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.619	2.615	0.004	342163	1.04		114		
15 Perfluorooctanoic acid	413.00 > 369.00	2.993	2.984	0.009	290666	1.15		115	3791	
	413.00 > 169.00	2.993	2.984	0.009	166261		1.75(0.90-1.10)	115	12537	

D 14 13C4 PFOA

417.00 > 372.00 2.993 2.984 0.009 12348635 56.3 113 1464087
 Page 513 of 834 12/09/2016

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.993	2.986	0.007	1.000	284917	0.9254	97.2		
D 17 13C4 PFOS	503.00	> 80.00	3.371	3.365	0.006		12589922	51.5	108	793078	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.380	3.366	0.014	1.000	255695	0.8928	96.2	33794	
	499.00	> 99.00	3.371	3.366	0.005	0.997	58595	4.36(0.90-1.10)	96.2	6167	M
20 Perfluorononanoic acid	463.00	> 419.00	3.380	3.373	0.007	1.000	184877	0.9888	98.9	3375	
D 19 13C5 PFNA	468.00	> 423.00	3.380	3.374	0.006		9431435	53.0	106	1129875	
D 21 13C8 FOSA	506.00	> 78.00	3.648	3.651	-0.003		21558794	52.0	104	974406	
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.657	3.656	0.001	1.000	403791	1.00	100	30165	
24 Perfluorodecanoic acid	513.00	> 469.00	3.746	3.738	0.008	1.000	159778	0.9790	97.9	5715	
D 23 13C2 PFDA	515.00	> 470.00	3.746	3.738	0.008		8418346	50.7	101	248297	
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	4.061	4.049	0.012	1.000	150902	0.9136	94.8		
D 27 13C2 PFUnA	565.00	> 520.00	4.078	4.070	0.008		6669716	53.0	106	287118	
28 Perfluoroundecanoic acid	563.00	> 519.00	4.086	4.071	0.015	1.000	148792	1.05	105	3690	
D 30 13C2 PFDoA	615.00	> 570.00	4.381	4.370	0.011		5950161	50.9	102	144810	
29 Perfluorododecanoic acid	613.00	> 569.00	4.381	4.370	0.011	1.000	109023	0.9679	96.8	130	
31 Perfluorotridecanoic acid	663.00	> 619.00	4.653	4.636	0.017	1.000	126175	1.02	102	142	
D 32 13C2-PFTeDA	715.00	> 670.00	4.893	4.882	0.011		12956938	53.6	107	632063	
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.893	4.884	0.009	1.000	237250	1.03	103	104	
	713.00	> 169.00	4.885	4.884	0.001	0.998	41035	5.78(0.00-0.00)	103	4184	
D 34 13C2-PFHxDA	815.00	> 770.00	5.334	5.320	0.014		6756331	51.8	104	147007	
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.334	5.320	0.014	1.000	218894	1.04	104	287	
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.738	5.720	0.018	1.000	118037	1.26	126	233	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L2_00022

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_005.d

Injection Date: 02-Dec-2016 10:37:07

Instrument ID: A8_N

Lims ID: IC L2

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 38

Worklist Smp#: 5

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

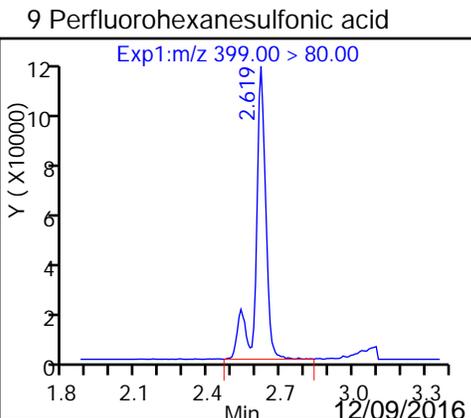
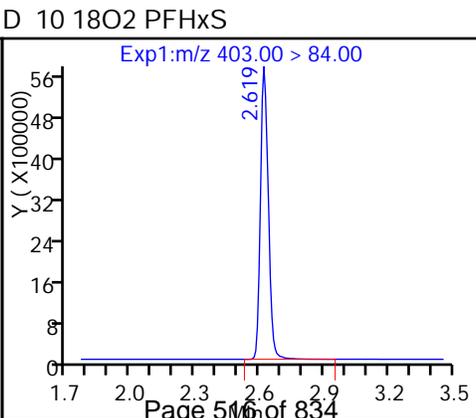
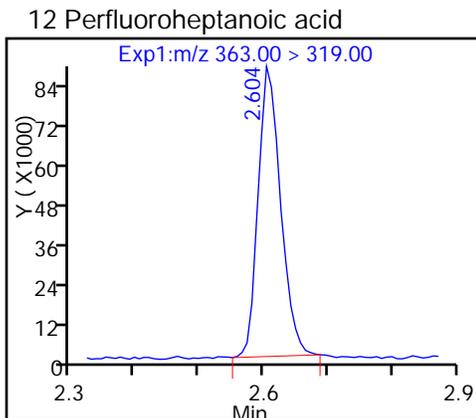
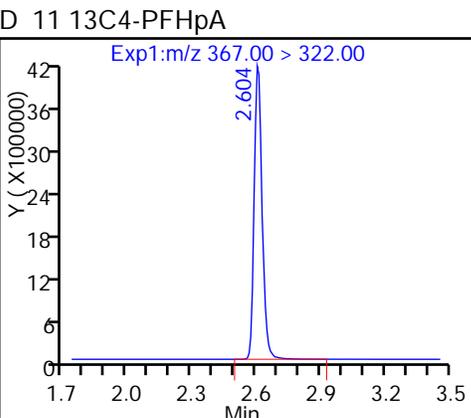
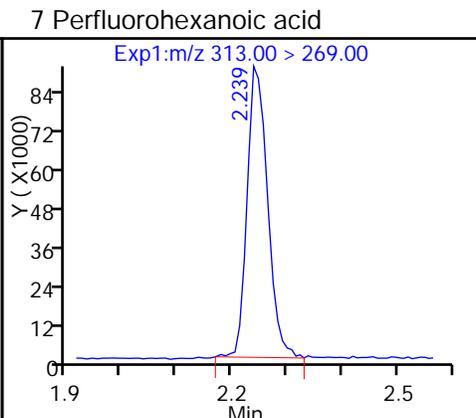
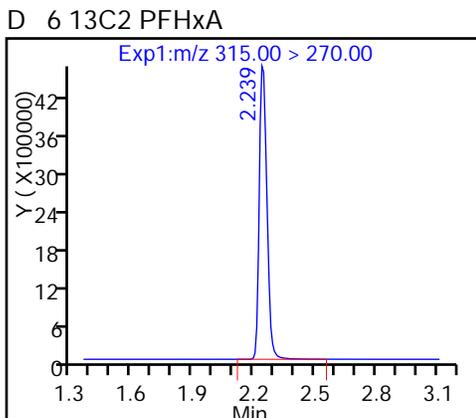
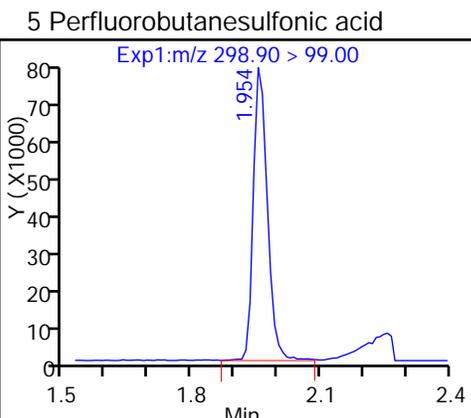
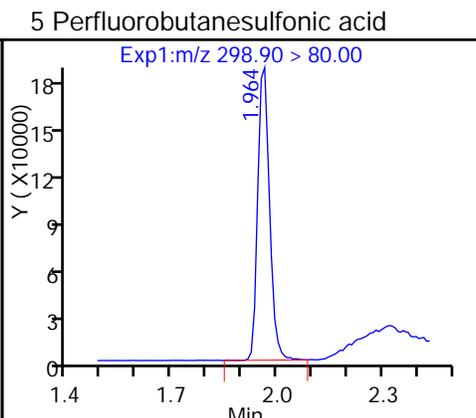
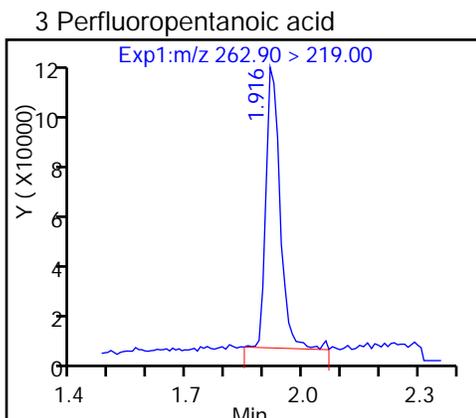
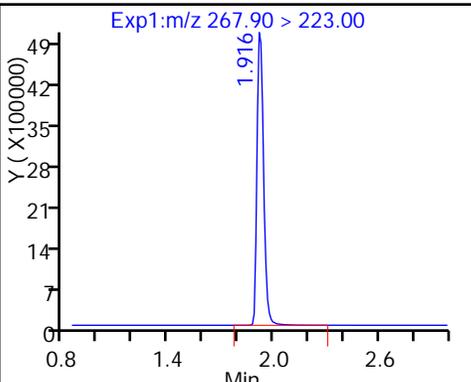
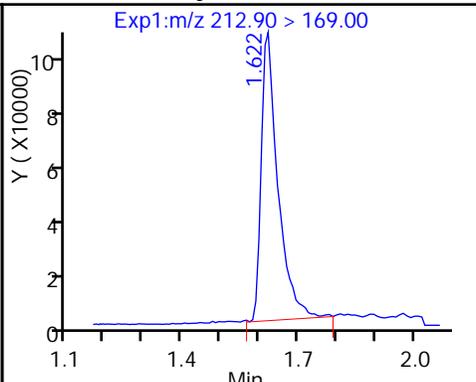
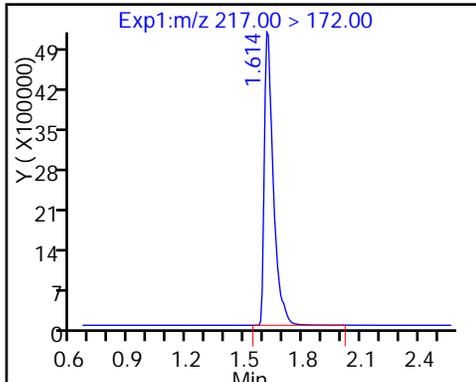
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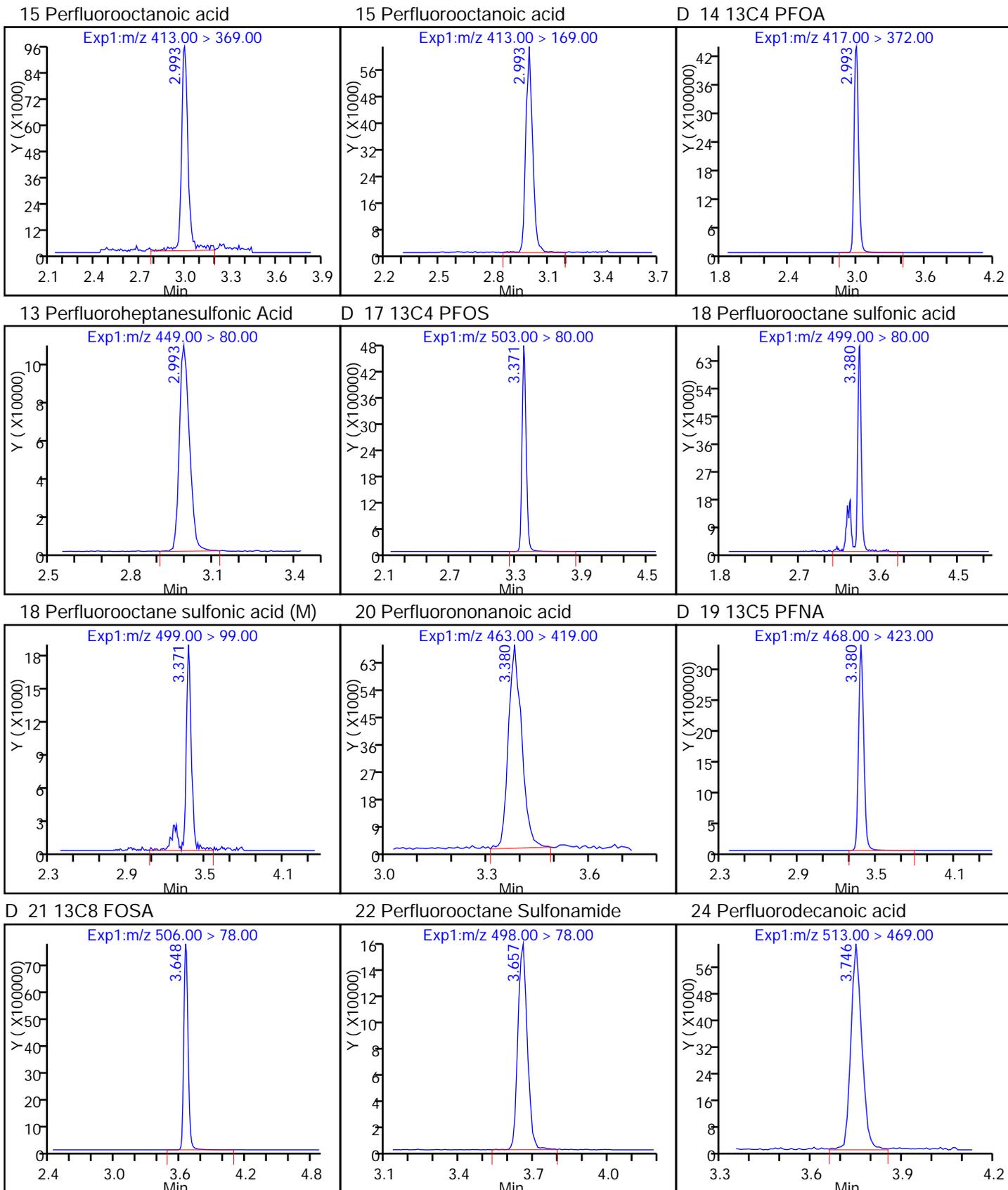
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

D 4 13C5-PFPeA

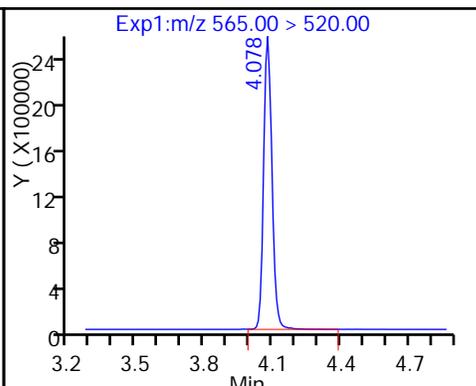
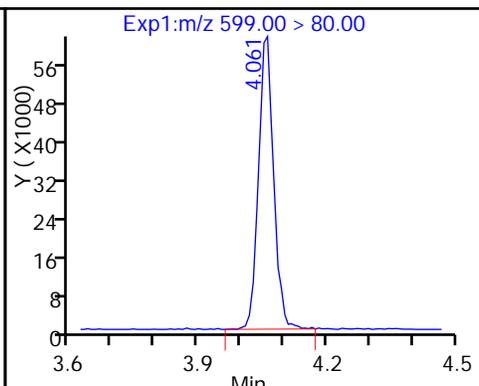
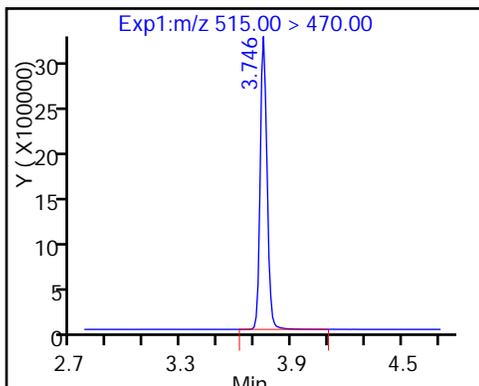




D 23 13C2 PFDA

26 Perfluorodecane Sulfonic acid

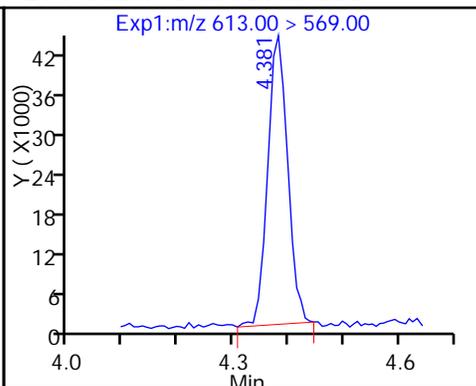
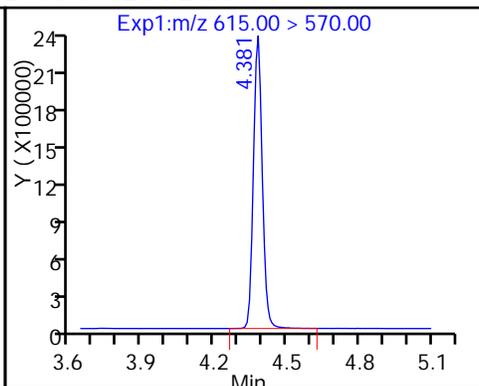
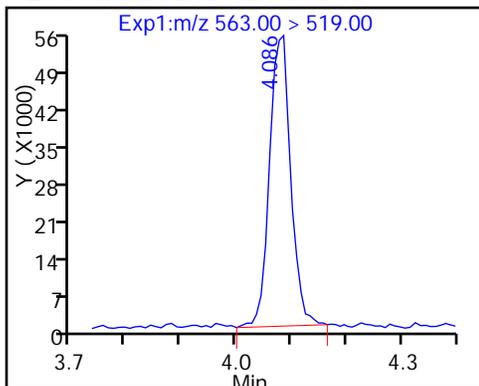
D 27 13C2 PFUnA



28 Perfluoroundecanoic acid

D 30 13C2 PFDaA

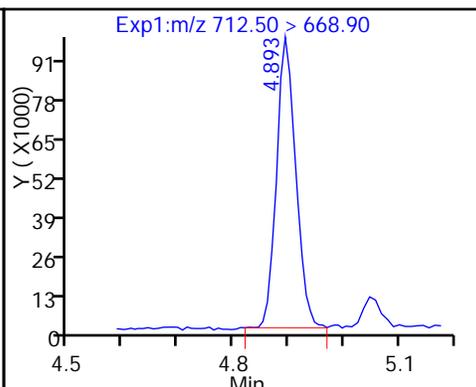
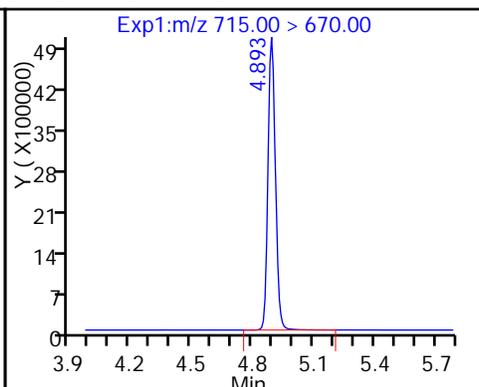
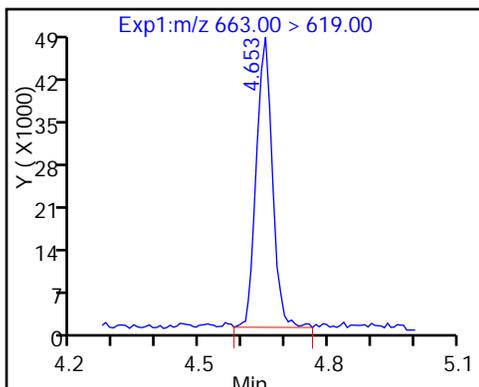
29 Perfluorododecanoic acid



31 Perfluorotridecanoic acid

D 32 13C2-PFTeDA

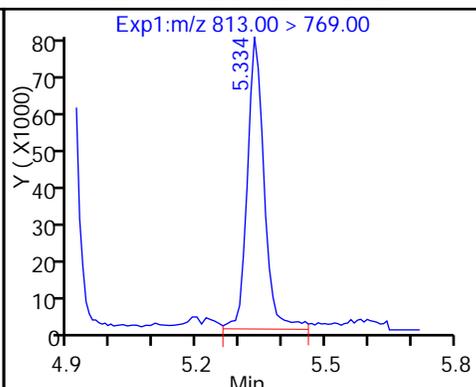
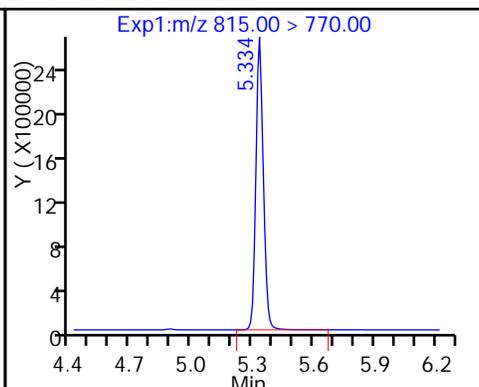
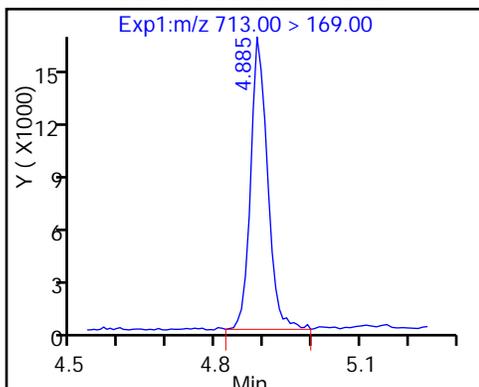
33 Perfluorotetradecanoic acid



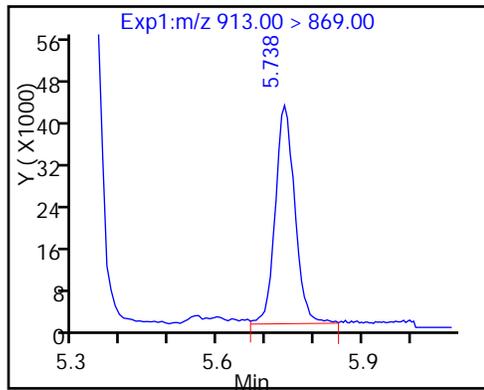
33 Perfluorotetradecanoic acid

D 34 13C2-PFHxD A

35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento

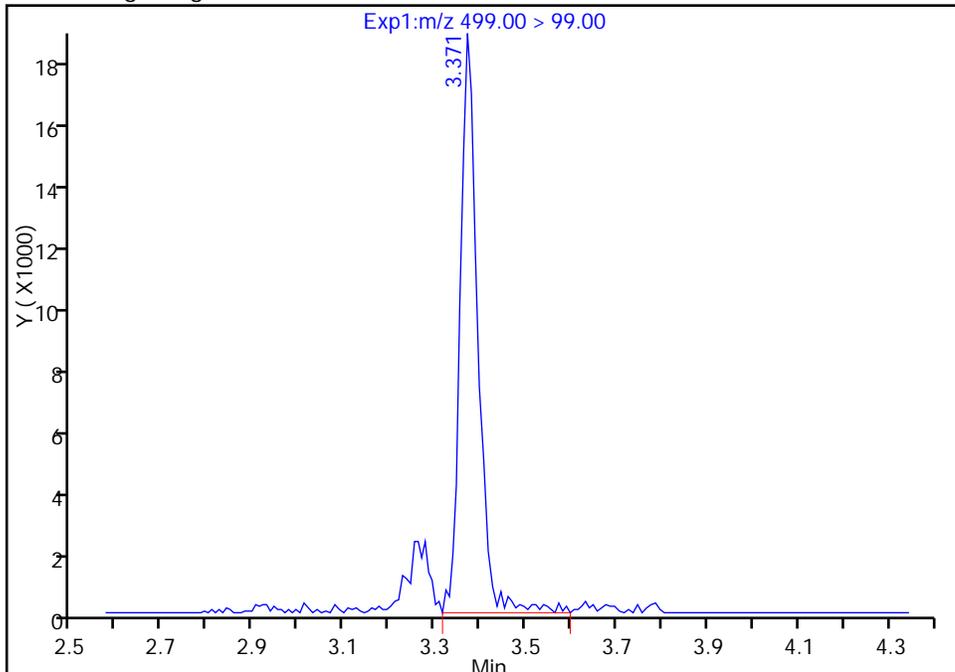
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Injection Date: 02-Dec-2016 10:37:07 Instrument ID: A8_N
Lims ID: IC L2
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 38 Worklist Smp#: 5
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

18 Perfluorooctane sulfonic acid, CAS: 1763-23-1

Signal: 2

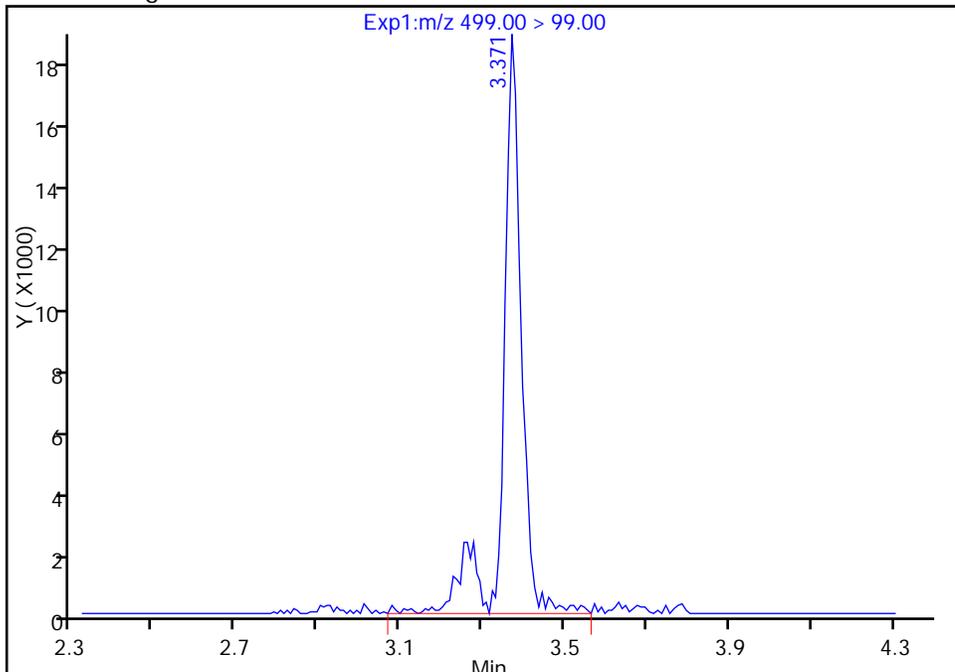
RT: 3.37
Area: 50423
Amount: 0.892814
Amount Units: ng/ml

Processing Integration Results



RT: 3.37
Area: 58595
Amount: 0.892814
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 02-Dec-2016 13:38:35
Audit Action: Manually Integrated

Audit Reason: Isomers

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_006.d
 Lims ID: IC L3
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 02-Dec-2016 10:44:37 ALS Bottle#: 39 Worklist Smp#: 6
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L3_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:32 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:38:56

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.622	1.615	0.007	20871580	61.3		123	1270902	
1 Perfluorobutyric acid	212.90 > 169.00	1.622	1.617	0.005	1937809	5.24		105	16704	
D 4 13C5-PFPeA	267.90 > 223.00	1.916	1.918	-0.002	16554836	61.3		123	1251857	
3 Perfluoropentanoic acid	262.90 > 219.00	1.926	1.920	0.006	1680231	4.88		97.5	18184	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.964	1.961	0.003	2954006	5.00		113		
	298.90 > 99.00	1.955	1.961	-0.007	1239873		2.38(0.00-0.00)	113		
D 6 13C2 PFHxA	315.00 > 270.00	2.239	2.239	-0.001	14744313	59.8		120	1821555	
7 Perfluorohexanoic acid	313.00 > 269.00	2.247	2.241	0.006	1458544	5.17		103	41938	
D 11 13C4-PFHpA	367.00 > 322.00	2.600	2.599	0.001	12841573	61.0		122	826915	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.600	2.600	0.0	1298003	4.89		97.7	25441	
D 10 18O2 PFHxS	403.00 > 84.00	2.616	2.614	0.002	17333346	57.3		121	1088243	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.616	2.615	0.001	1850926	4.77		105		
15 Perfluorooctanoic acid	413.00 > 369.00	2.981	2.984	-0.003	1560911	5.20		104	22497	
	413.00 > 169.00	2.989	2.984	0.005	958311		1.63(0.90-1.10)	104	88434	
D 14 13C4 PFOA	417.00 > 372.00	2.989	2.984	0.005	14705649	67.0		134	862824	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.989	2.986	0.003	1.000	1774408	5.08		107		
D 17 13C4 PFOS										
503.00 > 80.00	3.366	3.365	0.001		14291797	58.4		122	925059	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.366	3.366	0.0	1.000	1469896	4.52		97.4	134789	
499.00 > 99.00	3.366	3.366	0.0	1.000	312360		4.71(0.90-1.10)	97.4	26246	
20 Perfluorononanoic acid										
463.00 > 419.00	3.383	3.373	0.010	1.000	1085508	5.06		101	19854	
D 19 13C5 PFNA										
468.00 > 423.00	3.375	3.374	0.001		10826461	60.8		122	367418	
D 21 13C8 FOSA										
506.00 > 78.00	3.652	3.651	0.001		23576629	56.9		114	722968	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.652	3.656	-0.004	1.000	2370271	5.37		107	305013	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.740	3.738	0.002	1.000	938853	4.94		98.8	39325	
D 23 13C2 PFDA										
515.00 > 470.00	3.740	3.738	0.002		9802839	59.0		118	248176	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	4.046	4.049	-0.003	1.000	972225	5.18		108		
D 27 13C2 PFUnA										
565.00 > 520.00	4.079	4.070	0.009		7745917	61.6		123	312996	
28 Perfluoroundecanoic acid										
563.00 > 519.00	4.071	4.071	0.0	1.000	764307	4.66		93.1	21162	
D 30 13C2 PFDoA										
615.00 > 570.00	4.376	4.370	0.006		7049582	60.3		121	194268	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.376	4.370	0.006	1.000	669306	5.02		100	638	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.637	4.636	0.001	1.000	739992	5.05		101	761	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.880	4.882	-0.002		14954236	61.8		124	563118	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.888	4.884	0.004	1.000	1350591	4.96		99.1	554	
713.00 > 169.00	4.880	4.884	-0.004	0.998	220454		6.13(0.00-0.00)	99.1	17882	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.326	5.320	0.006		8035665	61.6		123	159038	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.326	5.320	0.006	1.000	800676	4.94		98.8	999	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.726	5.720	0.006	1.000	529337	4.77		95.5	857	

Reagents:

LCPFC-L3_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_006.d

Injection Date: 02-Dec-2016 10:44:37

Instrument ID: A8_N

Lims ID: IC L3

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 39

Worklist Smp#: 6

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

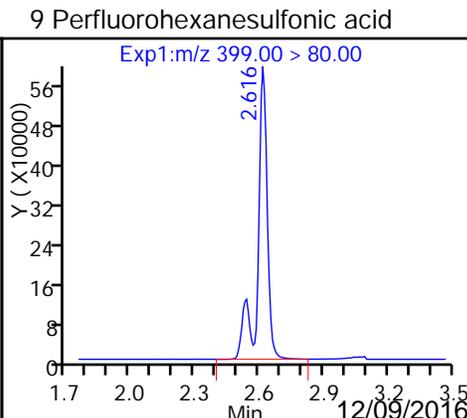
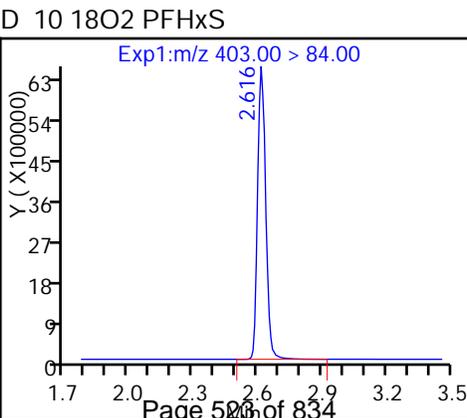
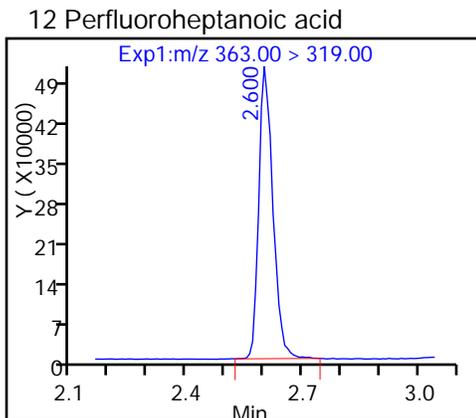
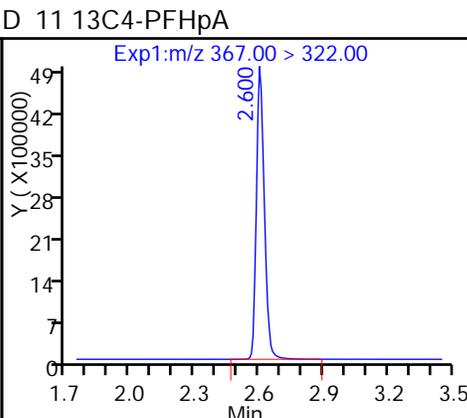
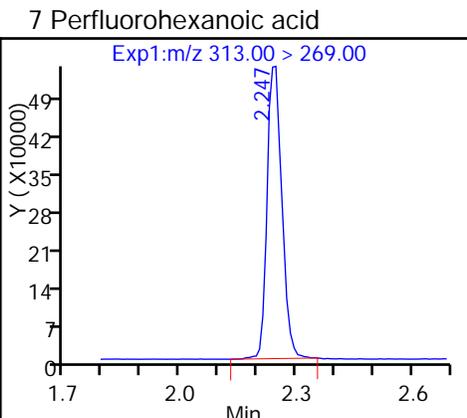
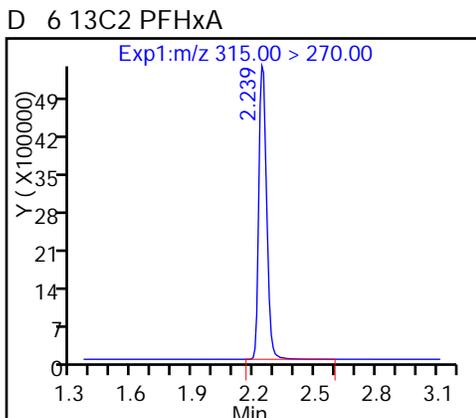
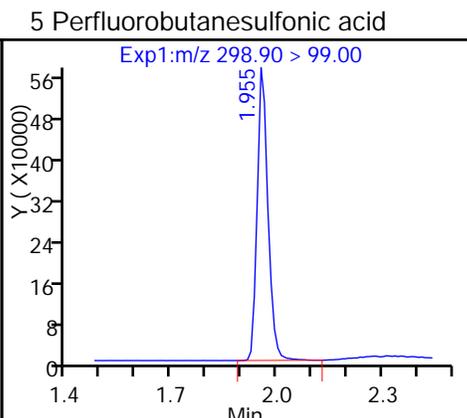
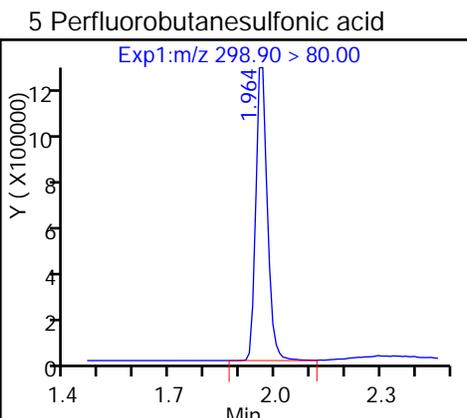
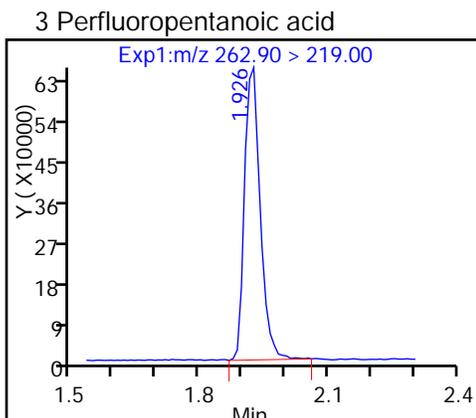
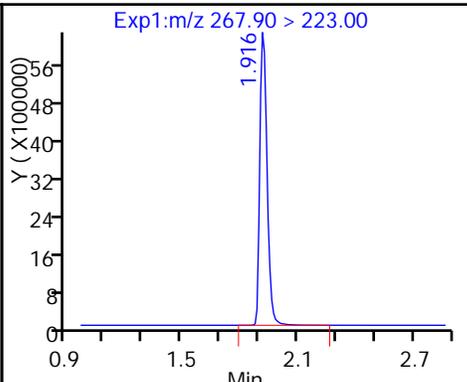
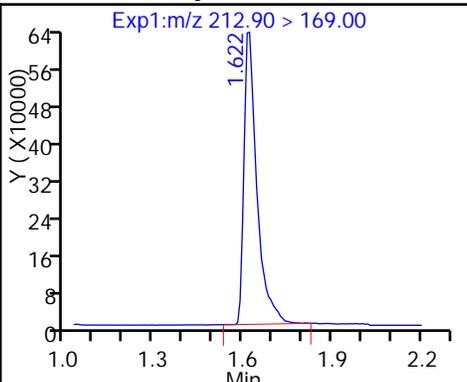
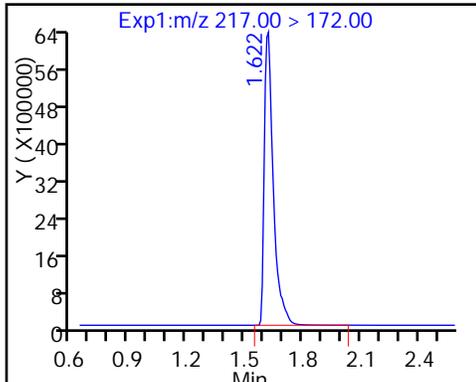
Method: A8_N

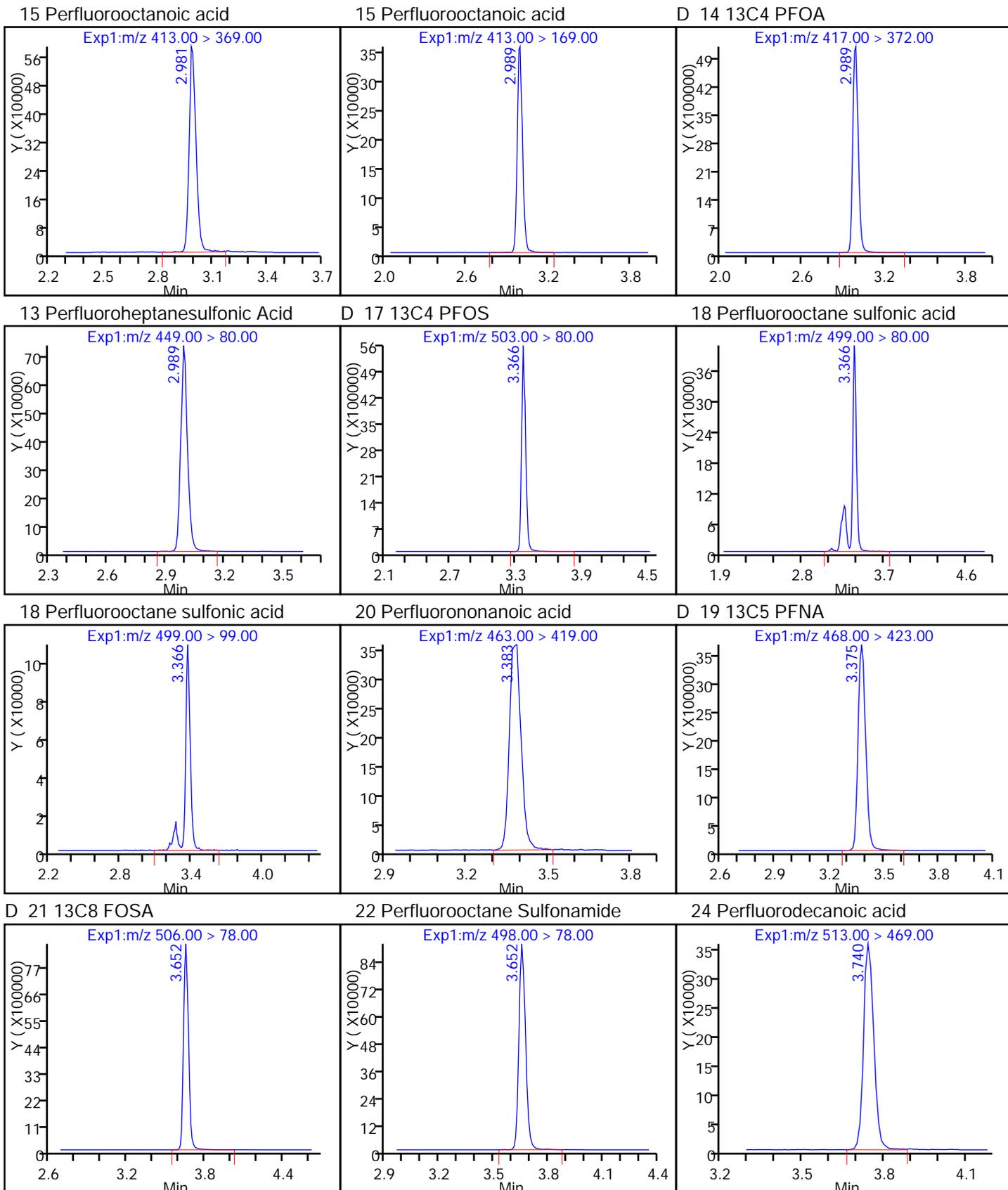
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

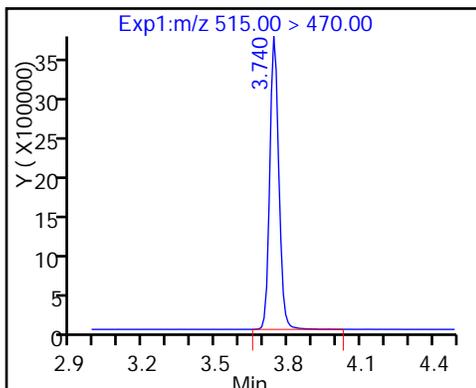
1 Perfluorobutyric acid

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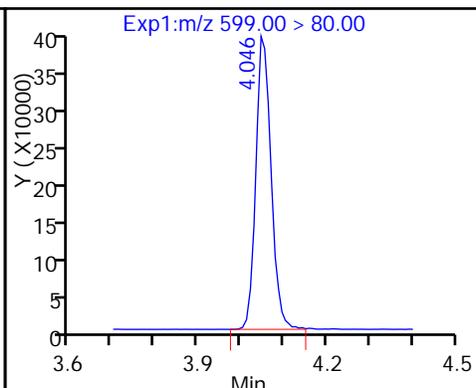




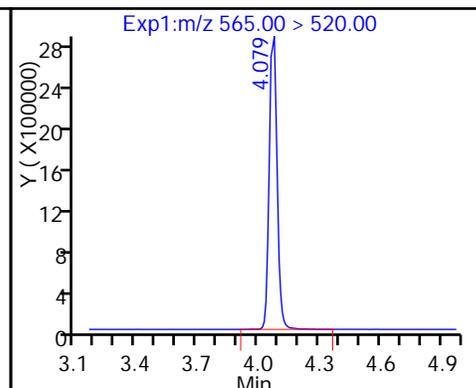
D 23 13C2 PFDA



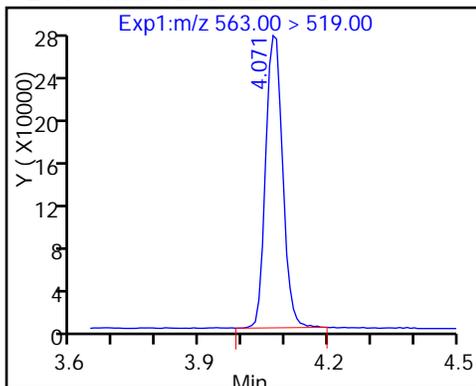
26 Perfluorodecane Sulfonic acid



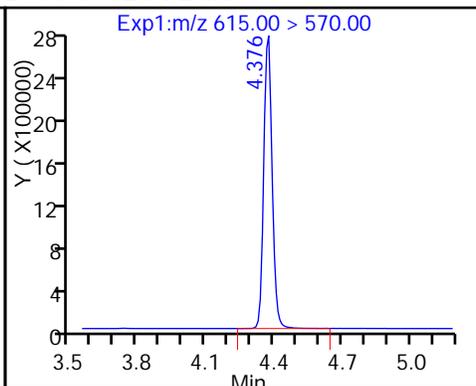
D 27 13C2 PFUnA



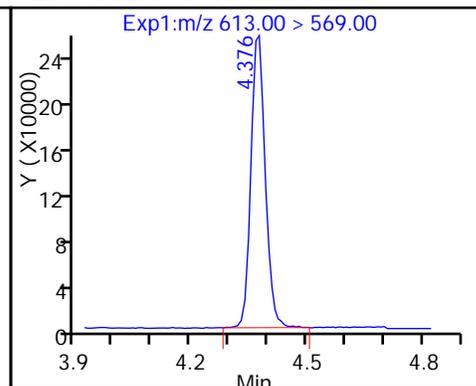
28 Perfluoroundecanoic acid



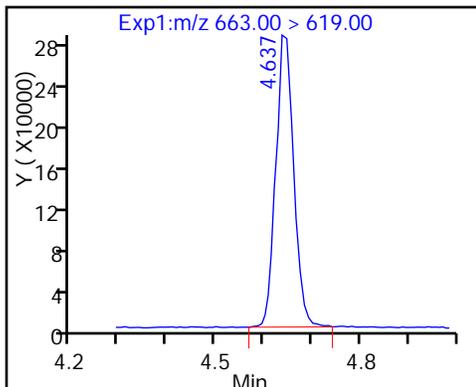
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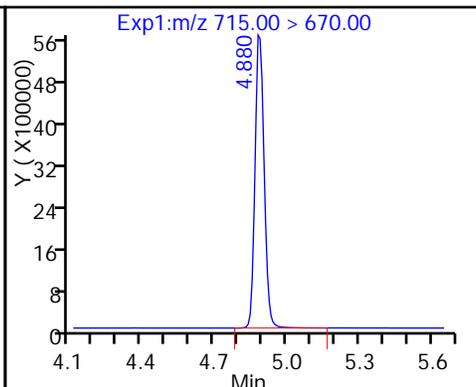
29 Perfluorododecanoic acid



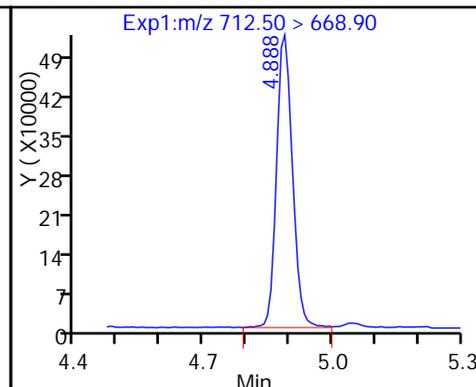
31 Perfluorotridecanoic acid



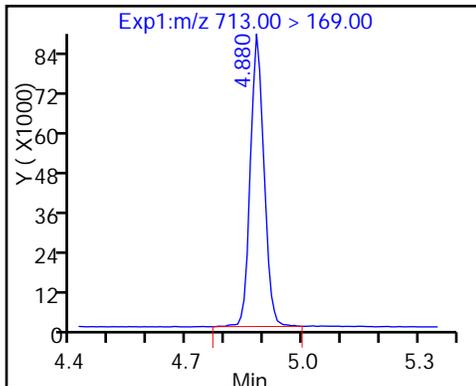
D 32 13C2-PFTeDA



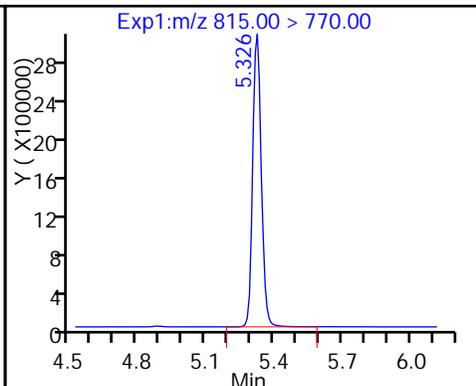
33 Perfluorotetradecanoic acid



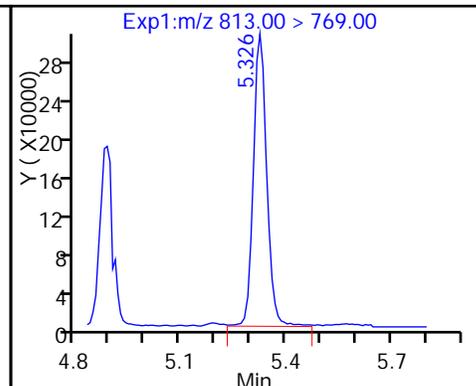
33 Perfluorotetradecanoic acid



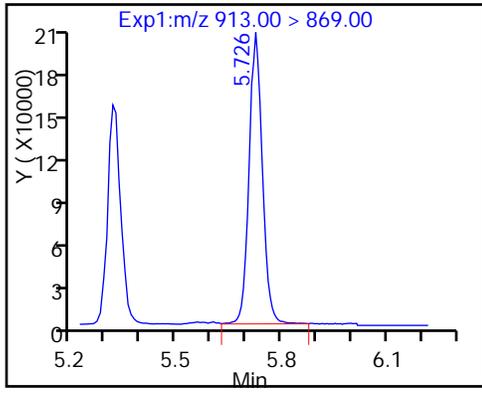
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_007.d
 Lims ID: IC L4
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 02-Dec-2016 10:52:06 ALS Bottle#: 40 Worklist Smp#: 7
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L4_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:34 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:36:08

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.614	1.615	-0.001	18935985	55.6		111	1406021	
1 Perfluorobutyric acid	212.90 > 169.00	1.614	1.617	-0.003	7435187	22.2		111	60389	
D 4 13C5-PFPeA	267.90 > 223.00	1.916	1.918	-0.002	15125796	56.0		112	1182515	
3 Perfluoropentanoic acid	262.90 > 219.00	1.916	1.920	-0.004	6559082	20.8		104	50558	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.954	1.961	-0.007	11367724	20.3		115		
	298.90 > 99.00	1.954	1.961	-0.007	4784706		2.38(0.00-0.00)	115		
D 6 13C2 PFHxA	315.00 > 270.00	2.231	2.239	-0.008	13498378	54.8		110	1276286	
7 Perfluorohexanoic acid	313.00 > 269.00	2.239	2.241	-0.002	5481010	21.2		106	245929	
D 11 13C4-PFHpA	367.00 > 322.00	2.593	2.599	-0.006	11424765	54.3		109	4354289	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.593	2.600	-0.007	5005633	21.2		106	82737	
D 10 18O2 PFHxS	403.00 > 84.00	2.608	2.614	-0.006	16435009	54.3		115	2029890	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.608	2.615	-0.007	6918390	18.8		103		
15 Perfluorooctanoic acid	413.00 > 369.00	2.982	2.984	-0.002	5603089	21.7		109	76131	
	413.00 > 169.00	2.974	2.984	-0.010	3413139		1.64(0.90-1.10)	109	244664	
D 14 13C4 PFOA	417.00 > 372.00	2.974	2.984	-0.010	12634984	57.6		115	1129431	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.982	2.986	-0.004	1.000	6947857	21.2		111	
D 17 13C4 PFOS	503.00	> 80.00	3.362	3.365	-0.003		13417569	54.9		115	370037
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.362	3.366	-0.004	1.000	5745134	18.8		101	258473
	499.00	> 99.00	3.362	3.366	-0.004	1.000	1279643		4.49(0.90-1.10)	101	205142
20 Perfluorononanoic acid	463.00	> 419.00	3.371	3.373	-0.002	1.000	4006430	20.8		104	50680
D 19 13C5 PFNA	468.00	> 423.00	3.371	3.374	-0.003		9730546	54.6		109	1069619
D 21 13C8 FOSA	506.00	> 78.00	3.648	3.651	-0.003		22550983	54.4		109	565745
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.648	3.656	-0.008	1.000	9372172	22.2		111	355431
24 Perfluorodecanoic acid	513.00	> 469.00	3.738	3.738	0.0	1.000	3671358	20.8		104	97996
D 23 13C2 PFDA	515.00	> 470.00	3.738	3.738	0.0		9112679	54.9		110	235961
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	4.048	4.049	-0.001	1.000	3647236	20.7		107	
D 27 13C2 PFUnA	565.00	> 520.00	4.065	4.070	-0.005		6913372	55.0		110	380701
28 Perfluoroundecanoic acid	563.00	> 519.00	4.065	4.071	-0.006	1.000	2908196	19.8		99.2	75091
D 30 13C2 PFDoA	615.00	> 570.00	4.365	4.370	-0.005		6319787	54.1		108	160074
29 Perfluorododecanoic acid	613.00	> 569.00	4.365	4.370	-0.005	1.000	2445148	20.4		102	2507
31 Perfluorotridecanoic acid	663.00	> 619.00	4.631	4.636	-0.005	1.000	2838032	21.6		108	3011
D 32 13C2-PFTeDA	715.00	> 670.00	4.883	4.882	0.001		13596161	56.2		112	754431
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.883	4.884	-0.001	1.000	5185293	21.2		106	2202
	713.00	> 169.00	4.875	4.884	-0.009	0.998	789160		6.57(0.00-0.00)	106	63701
D 34 13C2-PFHxDA	815.00	> 770.00	5.315	5.320	-0.005		7428976	57.0		114	148527
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.315	5.320	-0.005	1.000	2851018	22.1		110	3329
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.717	5.720	-0.003	1.000	1457725	14.7		73.3	2373

Reagents:

LCPFC-L4_00022

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_007.d

Injection Date: 02-Dec-2016 10:52:06

Instrument ID: A8_N

Lims ID: IC L4

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 40

Worklist Smp#: 7

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

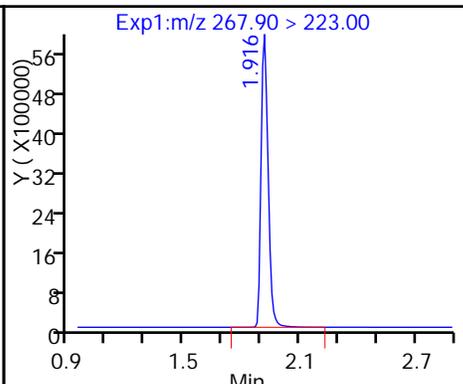
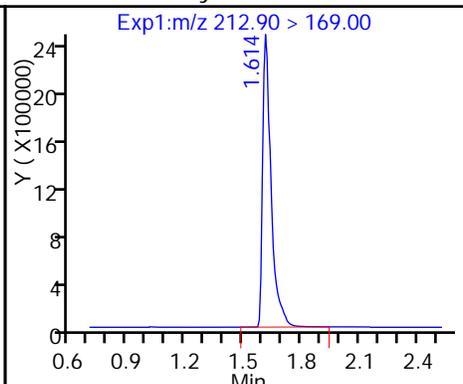
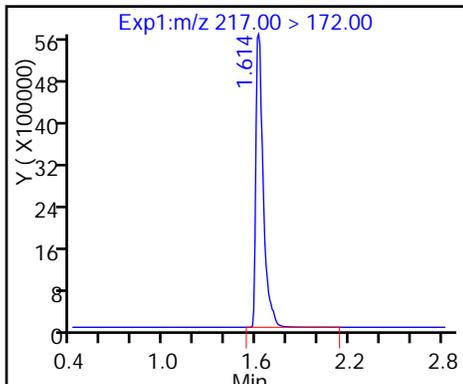
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

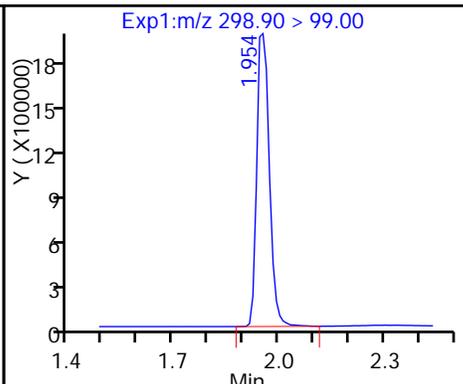
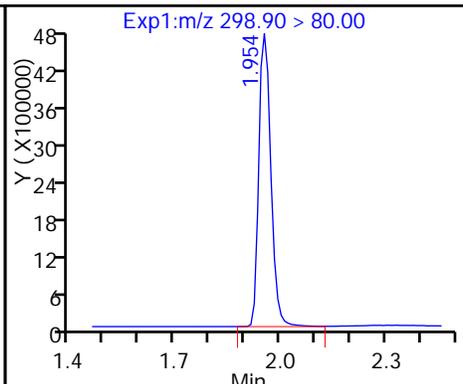
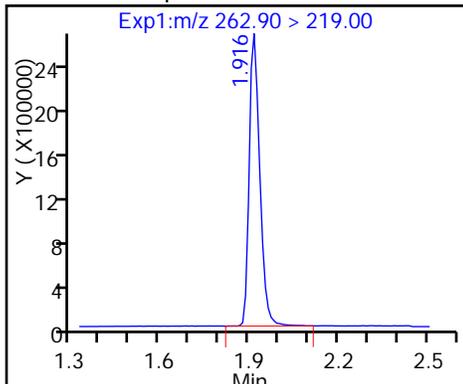
D 4 13C5-PFPeA



3 Perfluoropentanoic acid

5 Perfluorobutanesulfonic acid

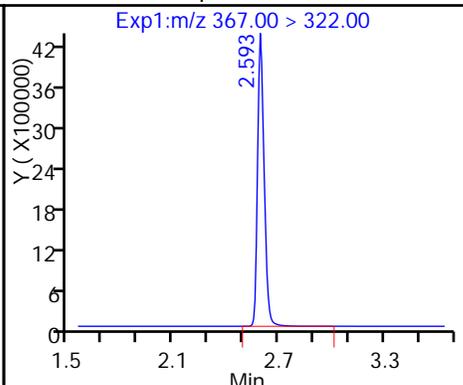
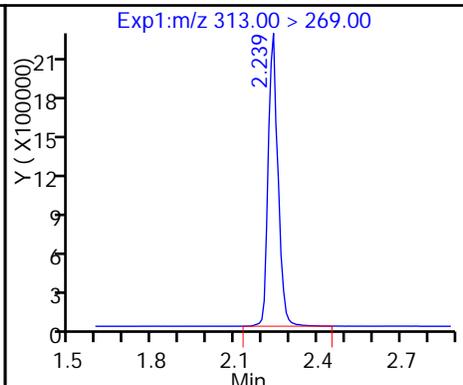
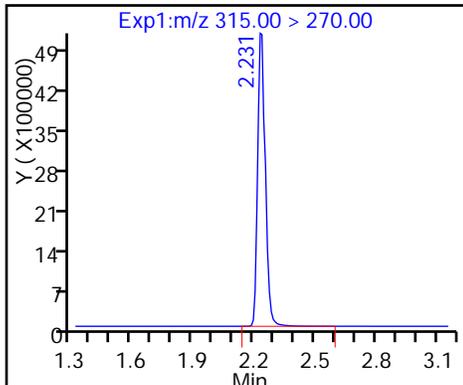
5 Perfluorobutanesulfonic acid



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

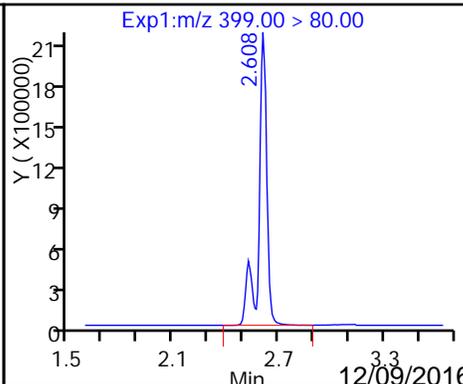
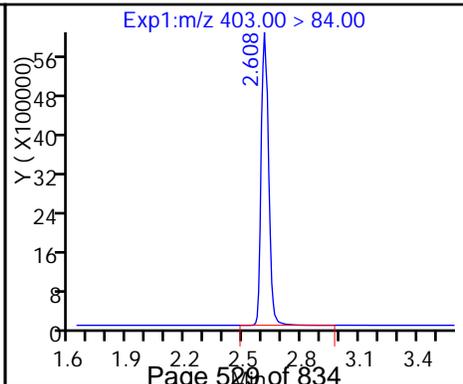
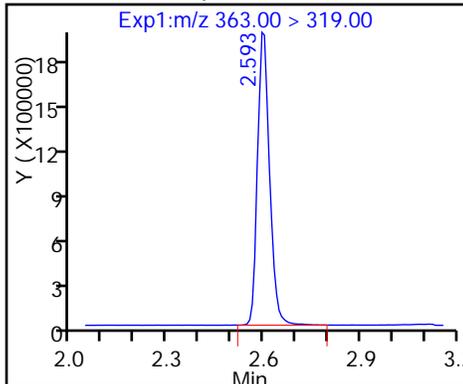
D 11 13C4-PFHpA

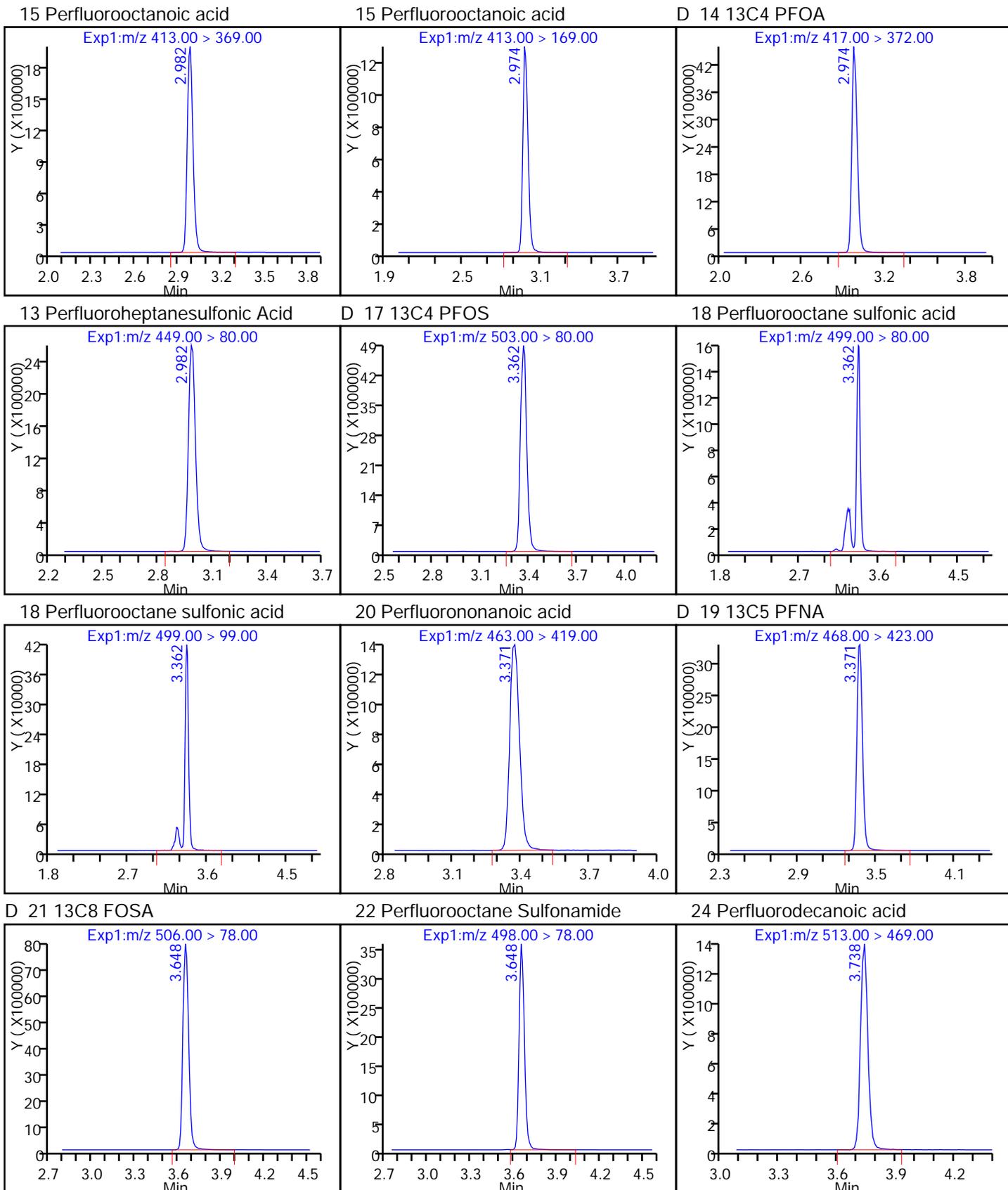


12 Perfluoroheptanoic acid

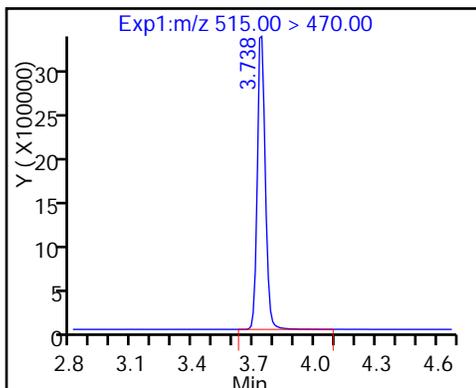
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9 Perfluorohexanesulfonic acid

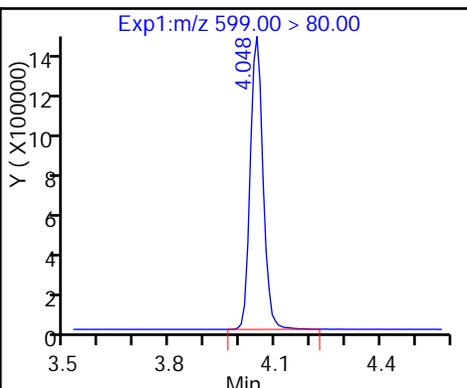




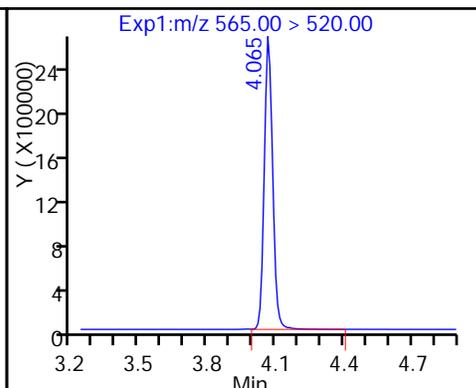
D 23 13C2 PFDA



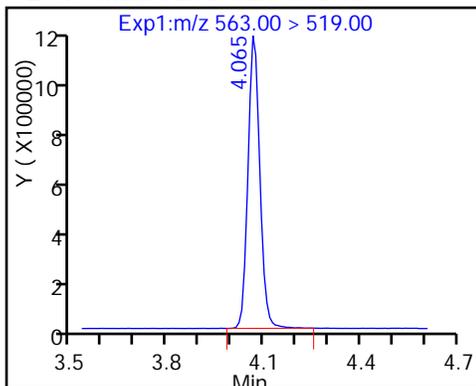
26 Perfluorodecane Sulfonic acid



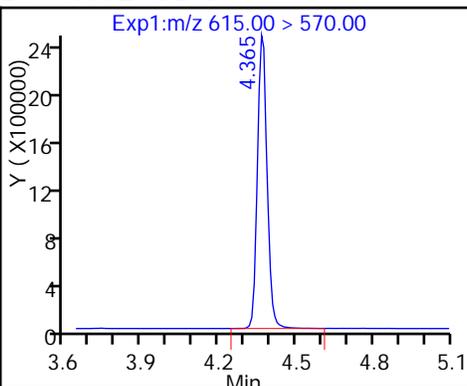
D 27 13C2 PFUnA



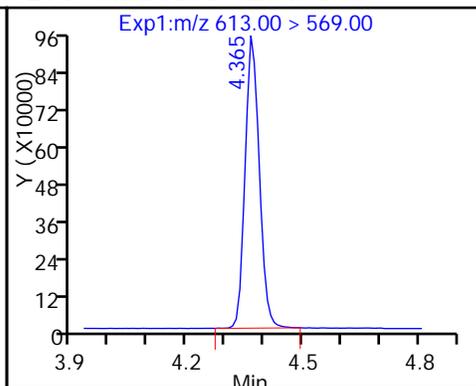
28 Perfluoroundecanoic acid



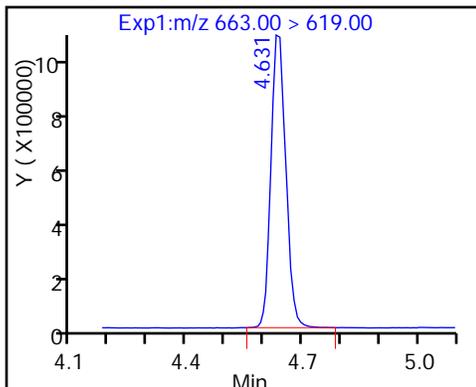
D 30 13C2 PFDaA



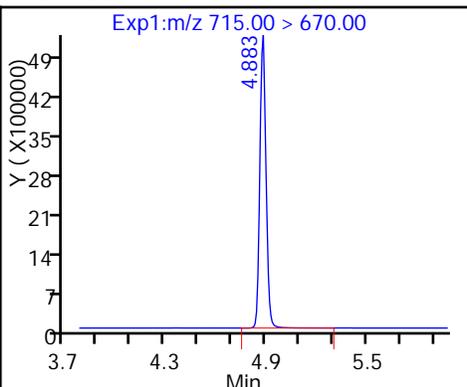
29 Perfluorododecanoic acid



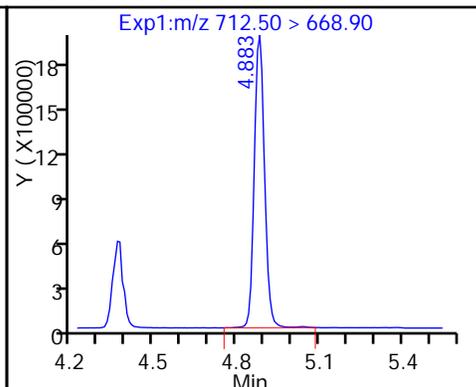
31 Perfluorotridecanoic acid



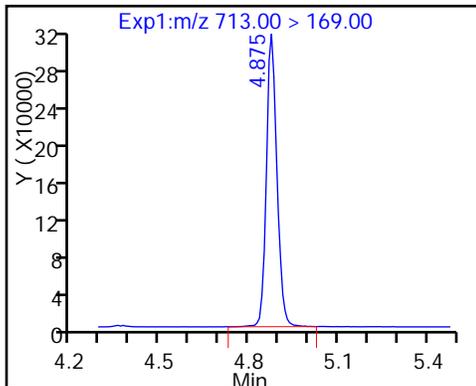
D 32 13C2-PFTeDA



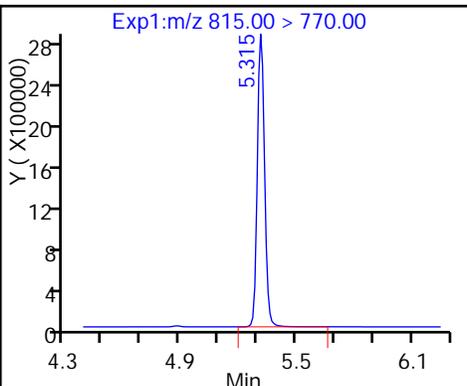
33 Perfluorotetradecanoic acid



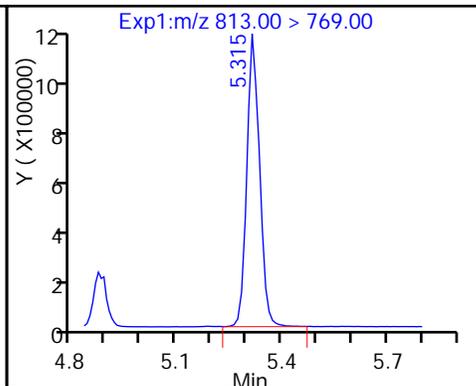
33 Perfluorotetradecanoic acid



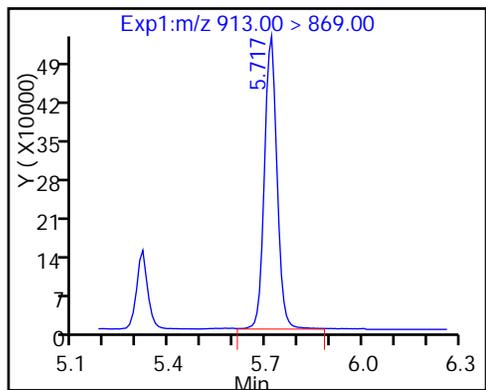
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_008.d
 Lims ID: IC L5
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 02-Dec-2016 10:59:32 ALS Bottle#: 41 Worklist Smp#: 8
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L5_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:36 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:39:38

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.614	1.615	-0.001	16489860	48.5		96.9	1638772	
1 Perfluorobutyric acid	212.90 > 169.00	1.614	1.617	-0.003	14709131	50.4		101	160438	
D 4 13C5-PFPeA	267.90 > 223.00	1.916	1.918	-0.002	12600012	46.6		93.3	943576	
3 Perfluoropentanoic acid	262.90 > 219.00	1.916	1.920	-0.004	12474289	47.6		95.1	93266	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.945	1.961	-0.016	21727893	44.2		99.9		
	298.90 > 99.00	1.945	1.961	-0.016	10461832		2.08(0.00-0.00)	99.9		
D 6 13C2 PFHxA	315.00 > 270.00	2.226	2.239	-0.013	11543313	46.8		93.7	1429849	
7 Perfluorohexanoic acid	313.00 > 269.00	2.226	2.241	-0.015	10870881	49.3		98.5	324820	
D 11 13C4-PFHpA	367.00 > 322.00	2.593	2.599	-0.006	9801352	46.6		93.1	1194438	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.593	2.600	-0.007	10247953	50.6		101	131961	
D 10 18O2 PFHxS	403.00 > 84.00	2.608	2.614	-0.006	14421709	47.7		101	804399	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.608	2.615	-0.007	14700167	45.5		100		
15 Perfluorooctanoic acid	413.00 > 369.00	2.972	2.984	-0.012	10948383	50.8		102	135691	
	413.00 > 169.00	2.972	2.984	-0.012	6763939		1.62(0.90-1.10)	102	350881	
D 14 13C4 PFOA	417.00 > 372.00	2.972	2.984	-0.012	10566267	48.2		96.3	1210292	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.972	2.986	-0.014	1.000	13319919	46.2		97.0		
D 17 13C4 PFOS										
503.00 > 80.00	3.352	3.365	-0.013		11794026	48.2		101	334101	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.352	3.366	-0.014	1.000	12334342	46.0		99.1	3198173	
499.00 > 99.00	3.352	3.366	-0.014	1.000	2699806		4.57(0.90-1.10)	99.1	276987	
20 Perfluorononanoic acid										
463.00 > 419.00	3.361	3.373	-0.012	1.000	8209339	50.0		100	106298	
D 19 13C5 PFNA										
468.00 > 423.00	3.361	3.374	-0.013		8274989	46.5		92.9	464531	
D 21 13C8 FOSA										
506.00 > 78.00	3.648	3.651	-0.003		20246280	48.9		97.7	412589	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.648	3.656	-0.008	1.000	18995833	50.2		100	525604	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.727	3.738	-0.011	1.000	7581125	49.1		98.2	195732	
D 23 13C2 PFDA										
515.00 > 470.00	3.727	3.738	-0.011		7964307	48.0		95.9	183252	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	4.039	4.049	-0.010	1.000	7816007	50.5		105		
D 27 13C2 PFUnA										
565.00 > 520.00	4.056	4.070	-0.014		5729423	45.6		91.1	366101	
28 Perfluoroundecanoic acid										
563.00 > 519.00	4.056	4.071	-0.015	1.000	5577456	45.9		91.9	89442	
D 30 13C2 PFDoA										
615.00 > 570.00	4.354	4.370	-0.016		5499233	47.0		94.1	178583	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.354	4.370	-0.016	1.000	5085511	48.8		97.7	5770	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.629	4.636	-0.007	1.000	5704096	49.9		99.7	5914	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.867	4.882	-0.015		11564641	47.8		95.6	896262	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.867	4.884	-0.017	1.000	10301361	48.5		96.9	4268	
713.00 > 169.00	4.867	4.884	-0.017	1.000	1659281		6.21(0.00-0.00)	96.9	129903	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.306	5.320	-0.014		6210852	47.6		95.3	147641	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.306	5.320	-0.014	1.000	5654463	51.4		103	6481	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.702	5.720	-0.018	1.000	5392609	62.4		125	9388	

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_008.d

Injection Date: 02-Dec-2016 10:59:32

Instrument ID: A8_N

Lims ID: IC L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 8

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

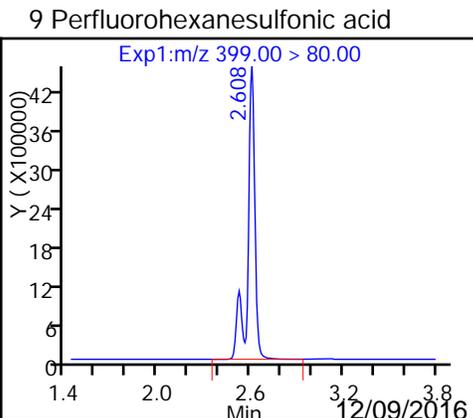
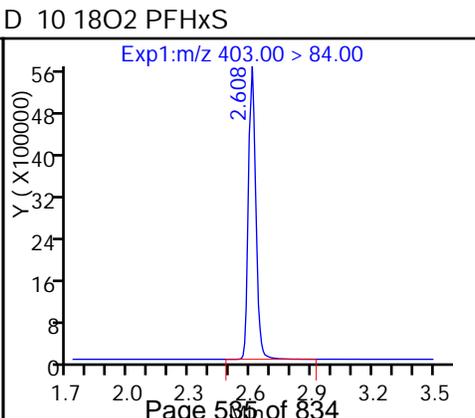
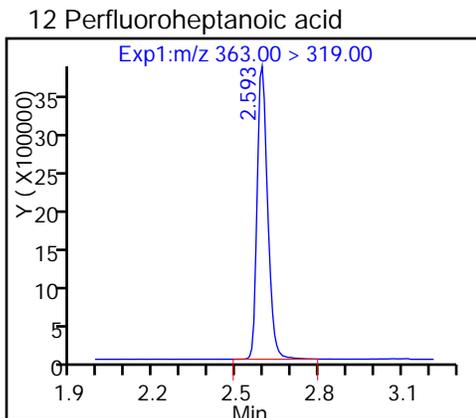
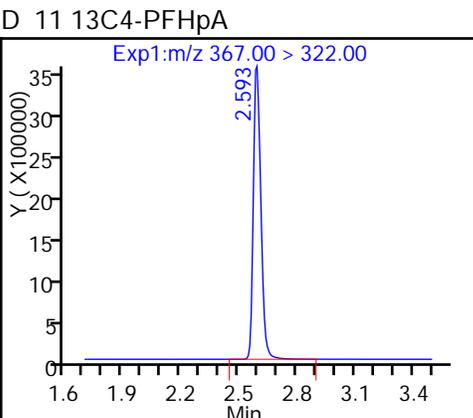
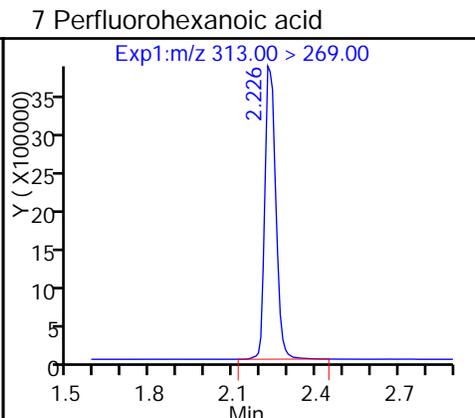
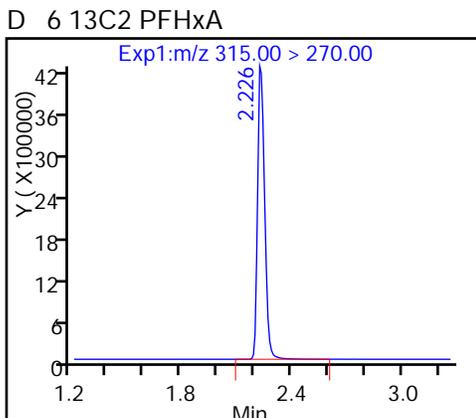
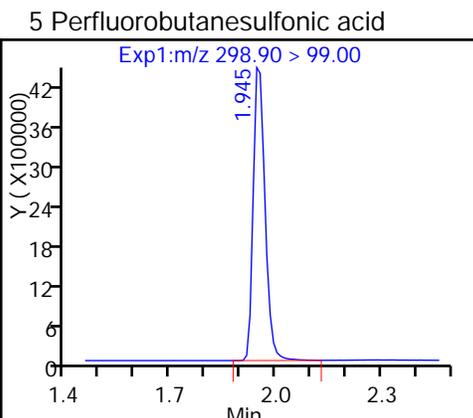
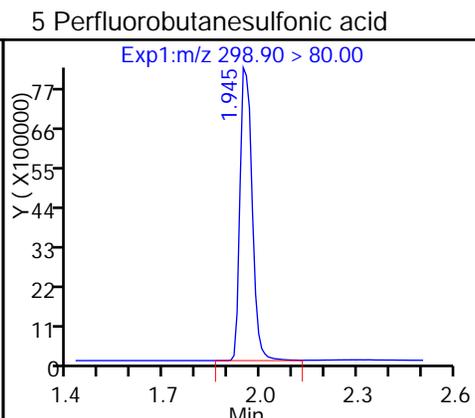
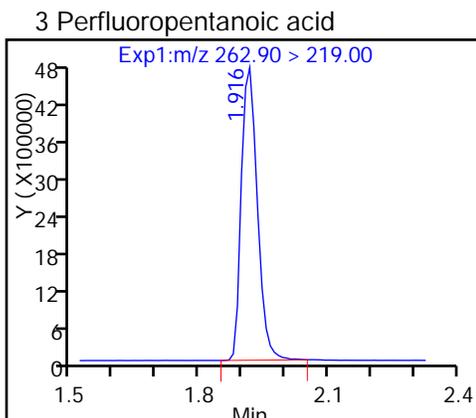
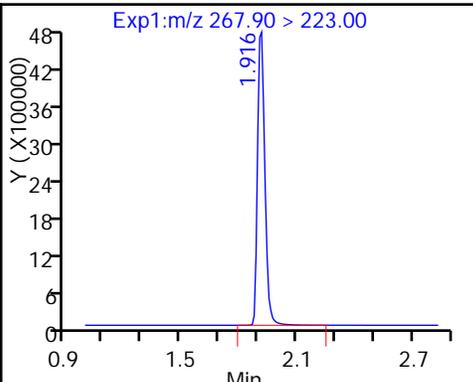
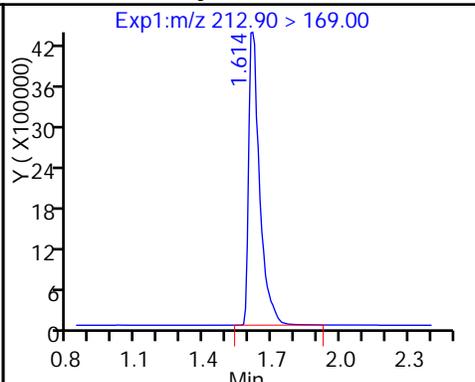
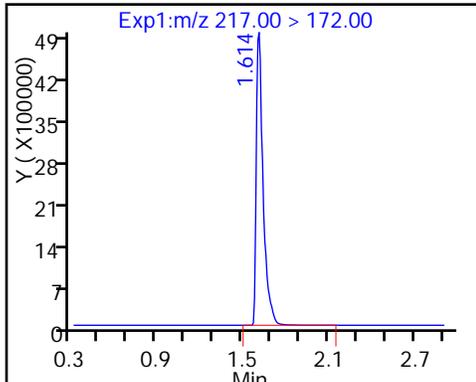
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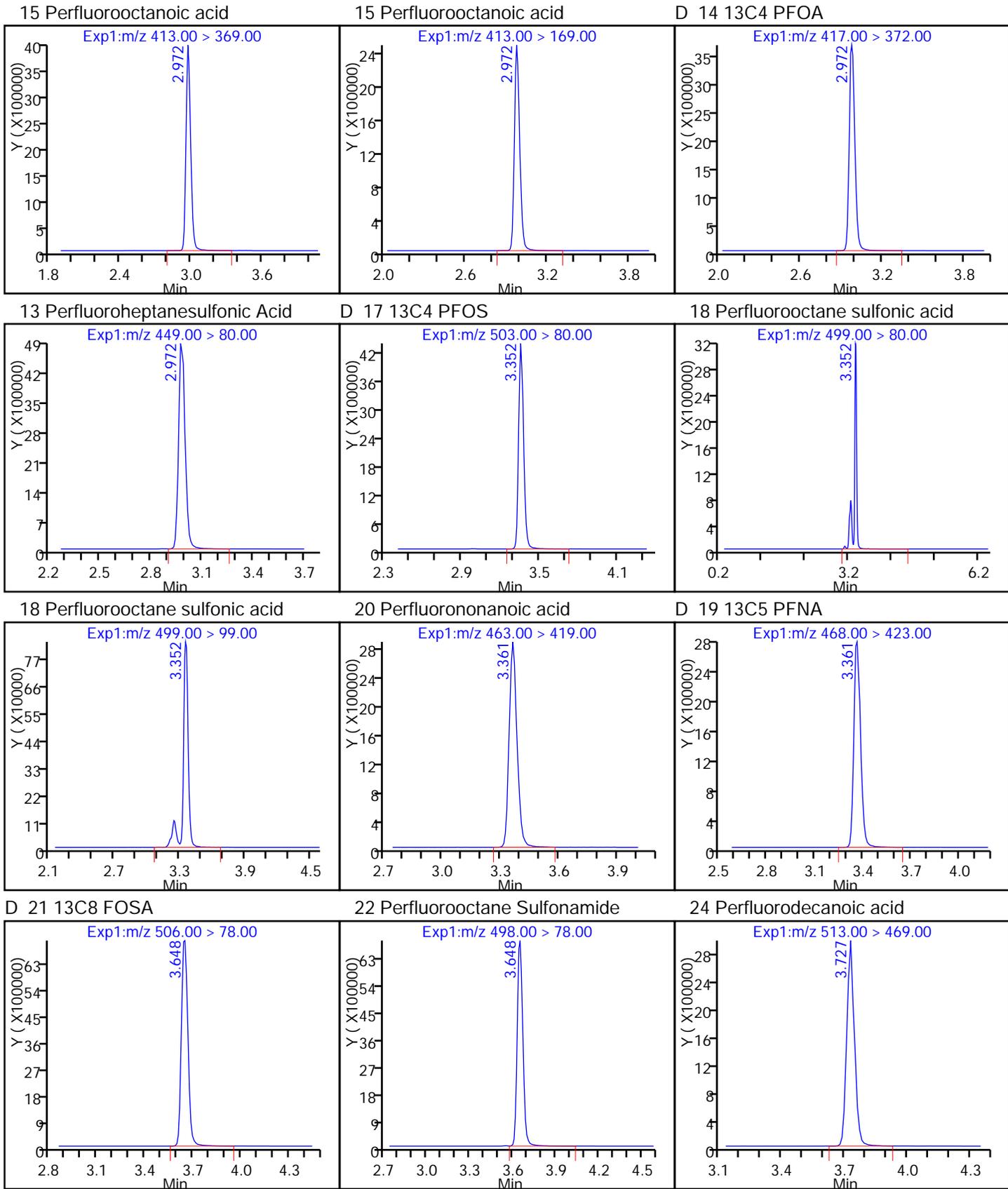
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

D 4 13C5-PFPeA

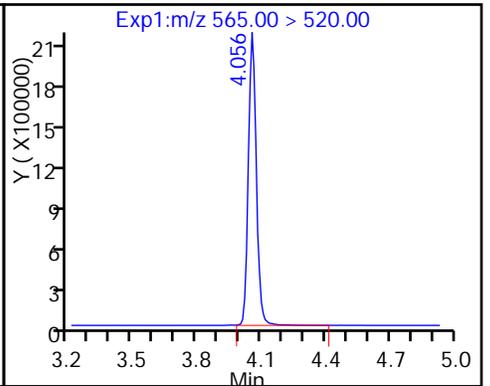
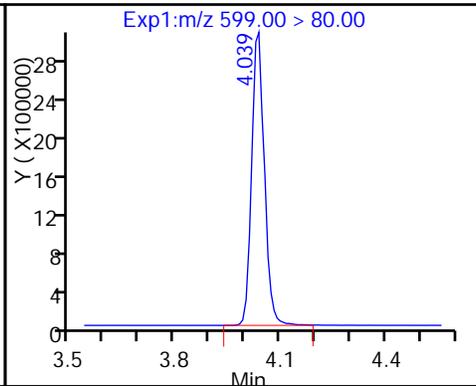
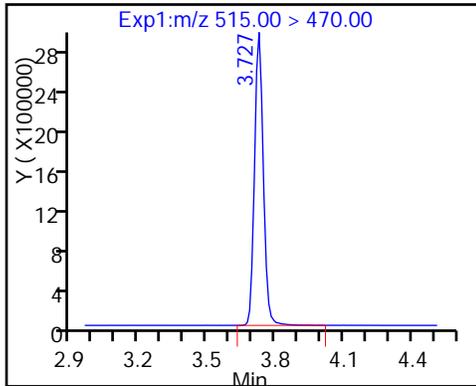




D 23 13C2 PFDA

26 Perfluorodecane Sulfonic acid

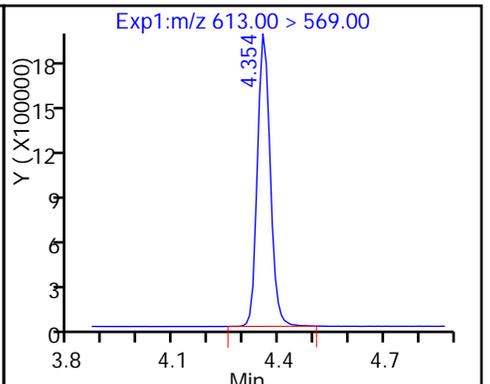
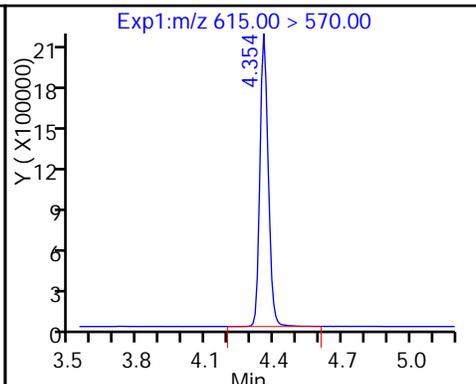
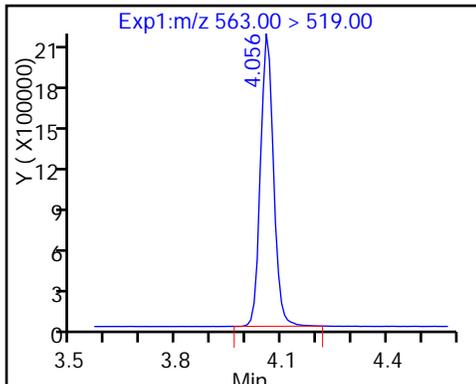
D 27 13C2 PFUnA



28 Perfluoroundecanoic acid

D 30 13C2 PFDaA

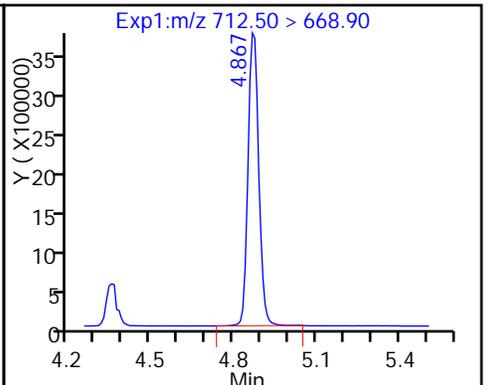
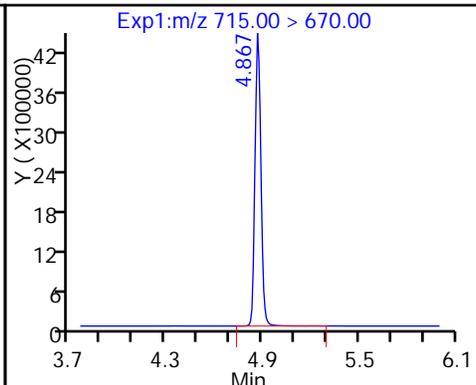
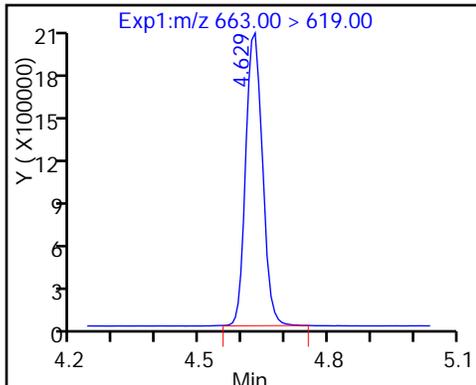
29 Perfluorododecanoic acid



31 Perfluorotridecanoic acid

D 32 13C2-PFTeDA

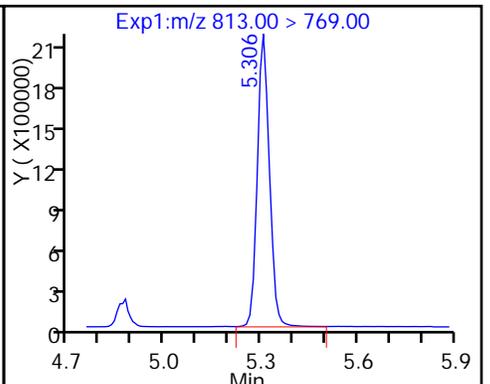
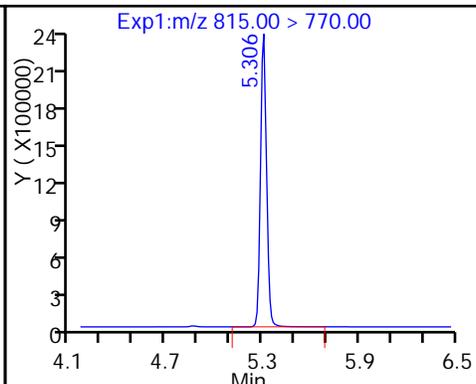
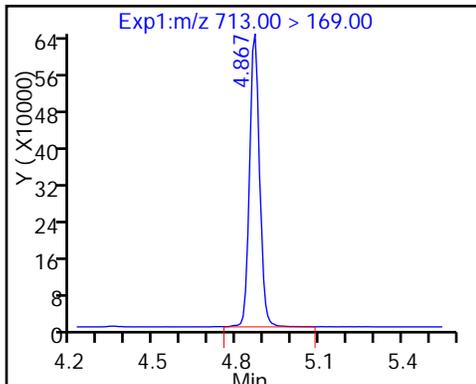
33 Perfluorotetradecanoic acid



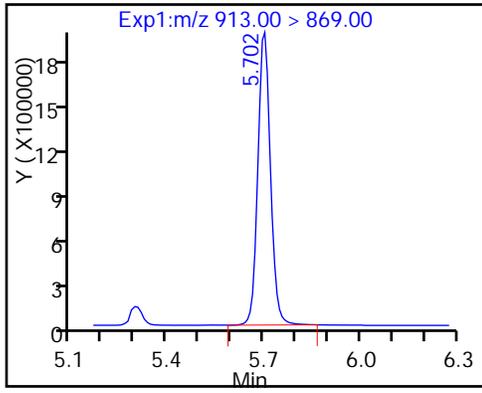
33 Perfluorotetradecanoic acid

D 34 13C2-PFHxDA

35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_009.d
 Lims ID: IC L6
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 02-Dec-2016 11:07:02 ALS Bottle#: 42 Worklist Smp#: 9
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L6_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:37 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:41:33

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.602	1.615	-0.013	15262180	44.8		89.7	765913	
1 Perfluorobutyric acid	212.90 > 169.00	1.610	1.617	-0.007	1.000	43875101	162.4	81.2	412834	
D 4 13C5-PFPeA	267.90 > 223.00	1.910	1.918	-0.008	10903650	40.4		80.7	1365253	
3 Perfluoropentanoic acid	262.90 > 219.00	1.910	1.920	-0.010	1.000	34297516	151.1	75.6	270259	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.949	1.961	-0.012	1.000	51594571	125.7	71.1		
	298.90 > 99.00	1.939	1.961	-0.022	0.995	28882906	1.79(0.00-0.00)	71.1		
D 6 13C2 PFHxA	315.00 > 270.00	2.225	2.239	-0.014	10277273	41.7		83.4	928195	
7 Perfluorohexanoic acid	313.00 > 269.00	2.225	2.241	-0.016	1.000	33379792	169.9	84.9	455453	
D 11 13C4-PFHpA	367.00 > 322.00	2.578	2.599	-0.021	8167611	38.8		77.6	1010737	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.578	2.600	-0.022	1.000	29343029	173.7	86.8	246277	
D 10 18O2 PFHxS	403.00 > 84.00	2.594	2.614	-0.020	12029706	39.8		84.1	822262	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.601	2.615	-0.014	1.000	45822208	170.0	93.4		
15 Perfluorooctanoic acid	413.00 > 369.00	2.958	2.984	-0.026	1.000	29789966	179.8	89.9	348106	
	413.00 > 169.00	2.966	2.984	-0.018	1.003	20146226	1.48(0.90-1.10)	89.9	805973	
D 14 13C4 PFOA	417.00 > 372.00	2.966	2.984	-0.018	8120442	37.0		74.0	869612	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.966	2.986	-0.020	1.000	38018954	159.5		83.8	
D 17 13C4 PFOS	503.00	> 80.00	3.340	3.365	-0.025		9744748	39.8		83.3	81434
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.340	3.366	-0.026	1.000	41837729	188.7		102	8979223
	499.00	> 99.00	3.340	3.366	-0.026	1.000	9665788		4.33(0.90-1.10)	102	925115
20 Perfluorononanoic acid	463.00	> 419.00	3.348	3.373	-0.025	1.000	25204338	186.7		93.4	245903
D 19 13C5 PFNA	468.00	> 423.00	3.355	3.374	-0.019		6808930	38.2		76.5	557624
D 21 13C8 FOSA	506.00	> 78.00	3.634	3.651	-0.017		17995728	43.4		86.9	430609
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.643	3.656	-0.013	1.000	51108754	151.8		75.9	319115
24 Perfluorodecanoic acid	513.00	> 469.00	3.721	3.738	-0.017	1.000	25614867	189.8		94.9	333805
D 23 13C2 PFDA	515.00	> 470.00	3.710	3.738	-0.028		6960897	41.9		83.8	168207
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	4.028	4.049	-0.021	1.000	24374906	190.6		98.9	
D 27 13C2 PFUnA	565.00	> 520.00	4.045	4.070	-0.025		4558330	36.2		72.5	316100
28 Perfluoroundecanoic acid	563.00	> 519.00	4.054	4.071	-0.017	1.000	17998762	186.3		93.2	327205
D 30 13C2 PFDoA	615.00	> 570.00	4.343	4.370	-0.027		4818858	41.2		82.4	222562
29 Perfluorododecanoic acid	613.00	> 569.00	4.343	4.370	-0.027	1.000	17698276	194.0		97.0	23046
31 Perfluorotridecanoic acid	663.00	> 619.00	4.613	4.636	-0.023	1.000	18549592	185.0		92.5	33008
D 32 13C2-PFTeDA	715.00	> 670.00	4.855	4.882	-0.027		10102255	41.8		83.5	643461
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.864	4.884	-0.020	1.000	31923969	171.4		85.7	17512
	713.00	> 169.00	4.855	4.884	-0.029	0.998	6059732		5.27(0.00-0.00)	85.7	373056
D 34 13C2-PFHxDA	815.00	> 770.00	5.290	5.320	-0.030		5535655	42.5		84.9	147627
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.290	5.320	-0.030	1.000	18745470	196.6		98.3	24259
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.688	5.720	-0.032	1.000	13671969	180.4		90.2	26246

Reagents:

LCPFC-L6_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_009.d

Injection Date: 02-Dec-2016 11:07:02

Instrument ID: A8_N

Lims ID: IC L6

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 42

Worklist Smp#: 9

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

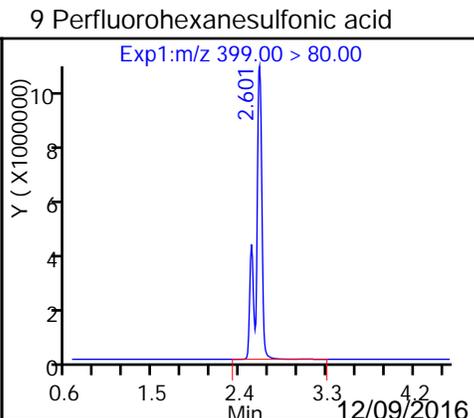
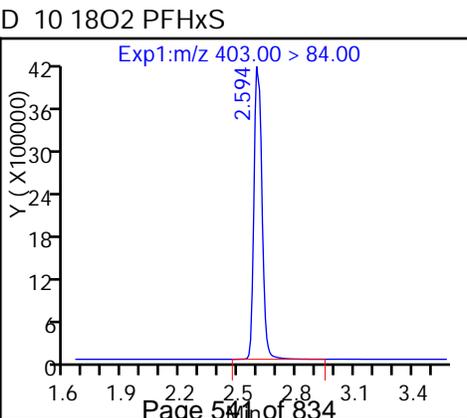
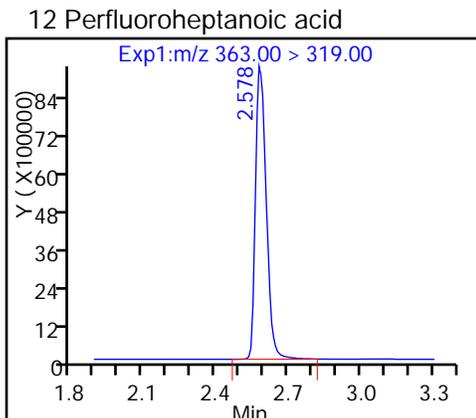
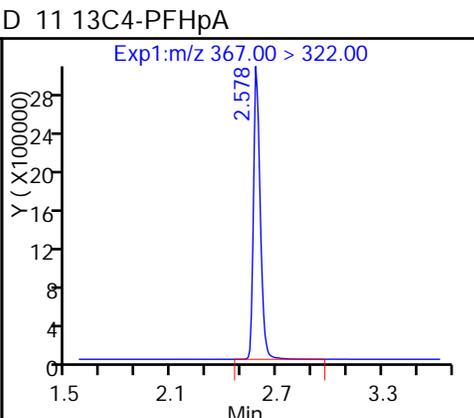
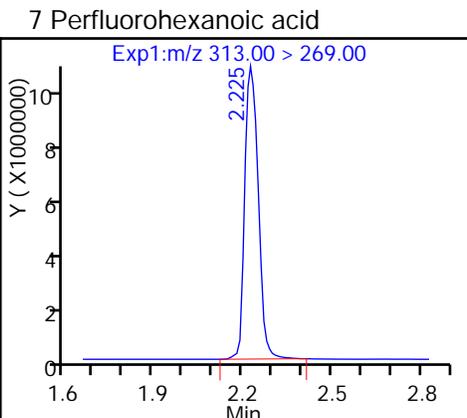
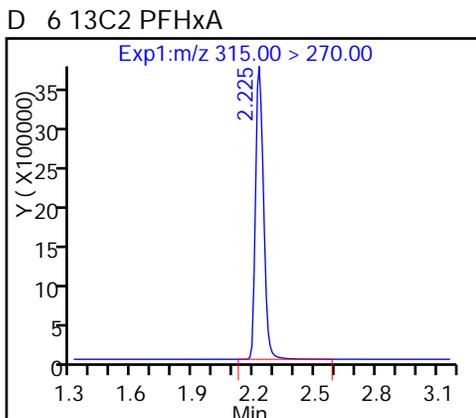
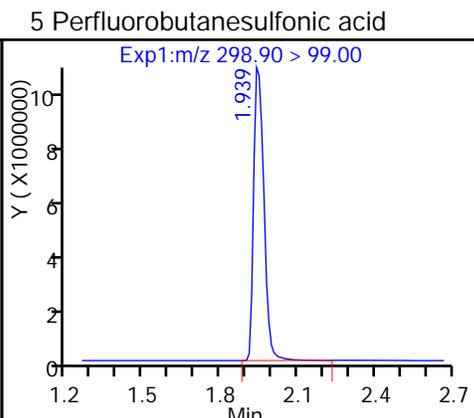
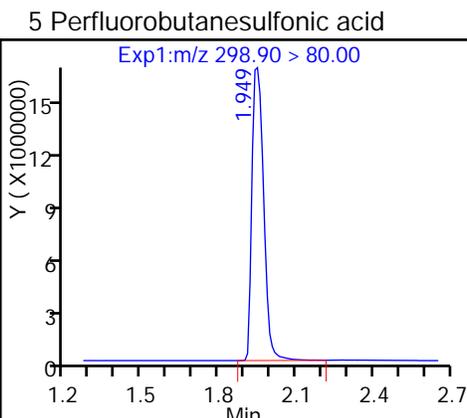
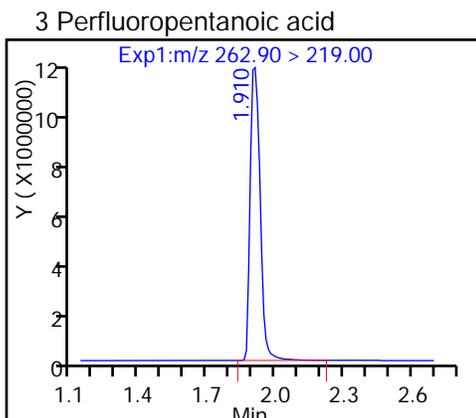
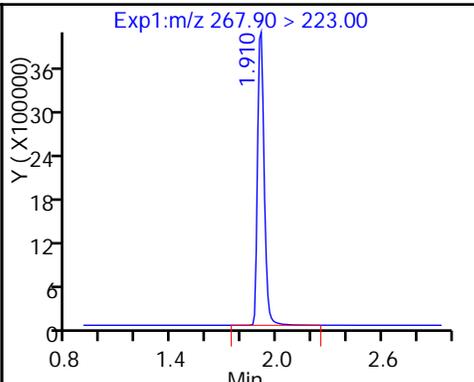
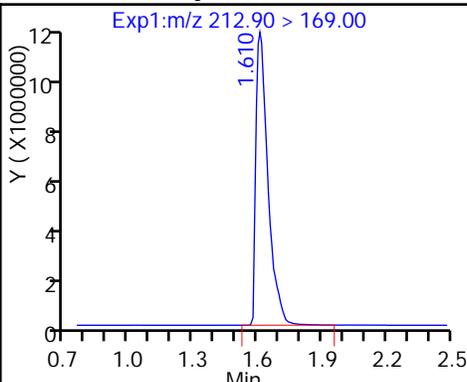
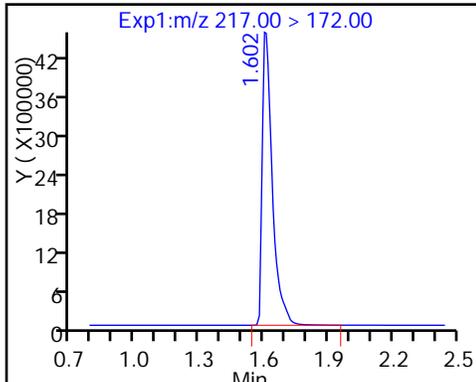
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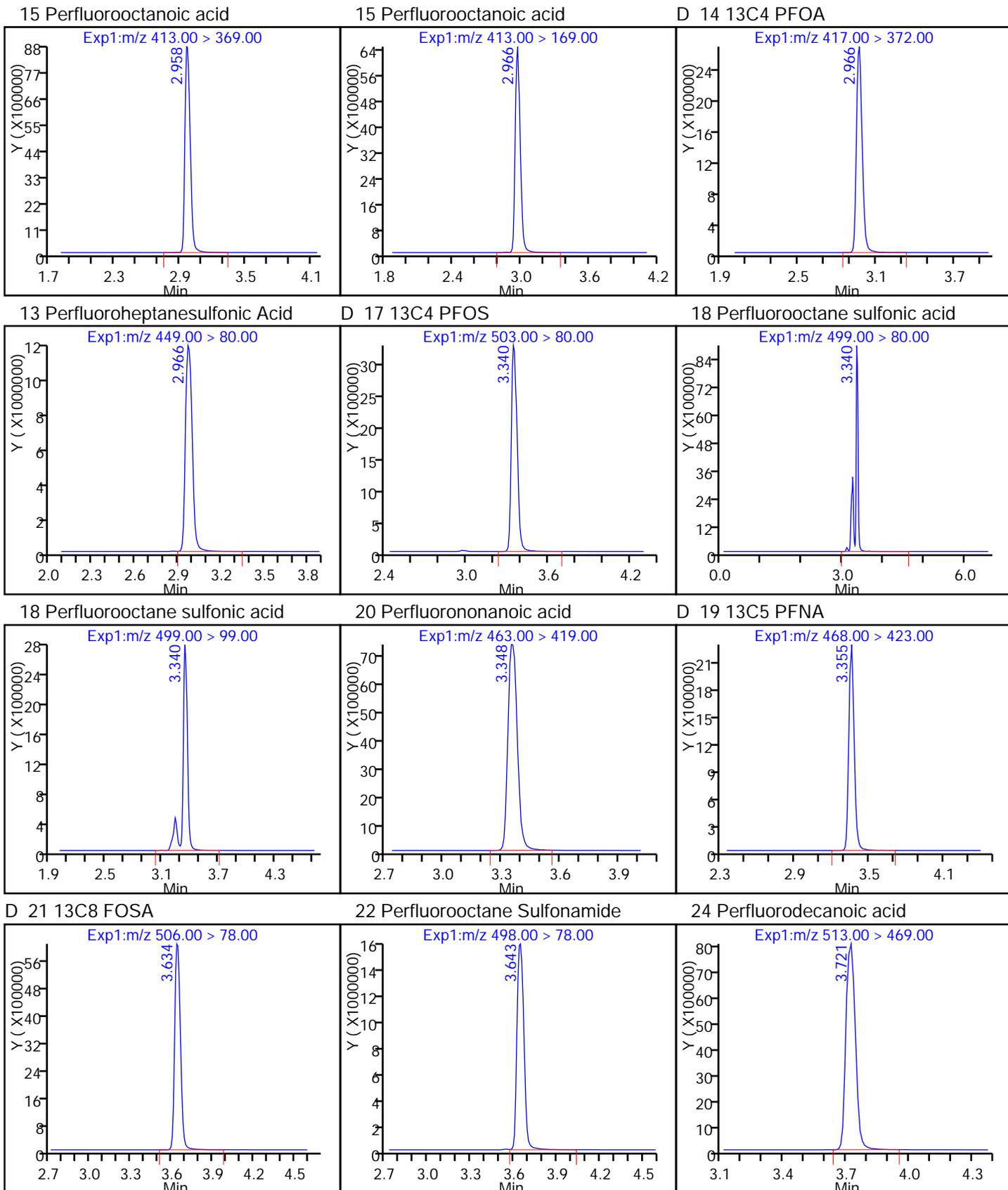
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

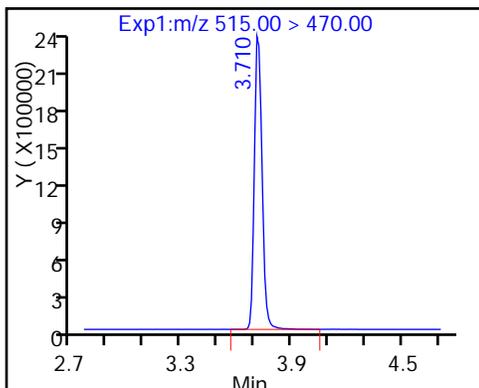
1 Perfluorobutyric acid

D 4 13C5-PFPeA

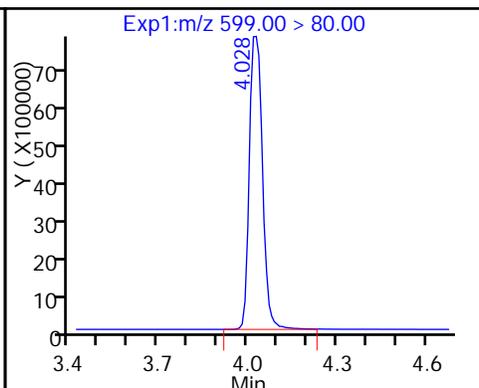




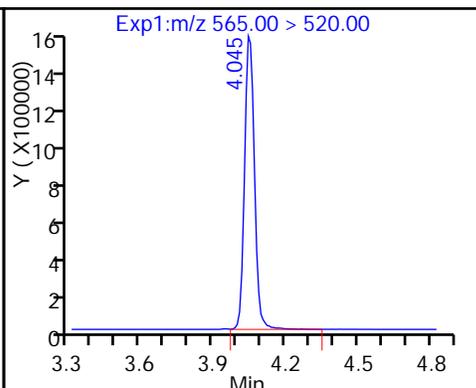
D 23 13C2 PFDA



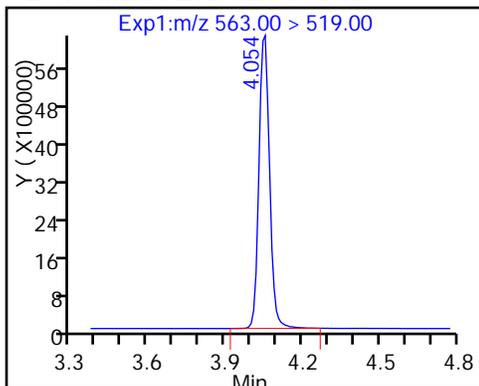
26 Perfluorodecane Sulfonic acid



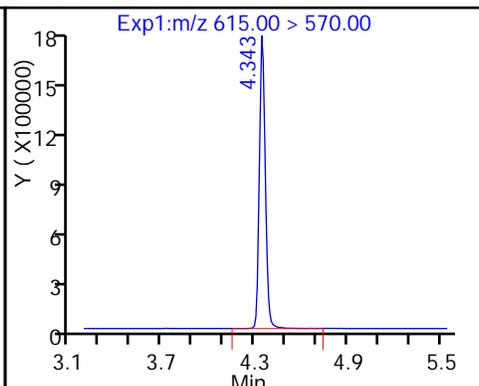
D 27 13C2 PFUnA



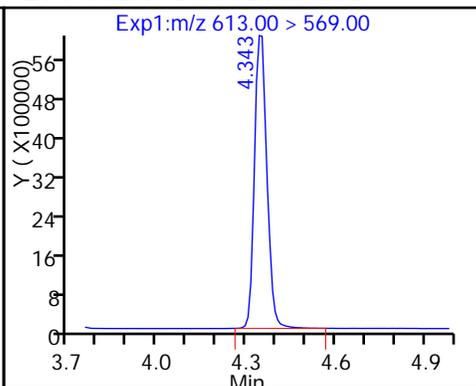
28 Perfluoroundecanoic acid



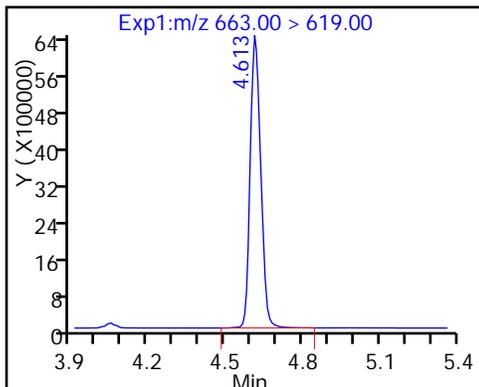
D 30 13C2 PFDaA



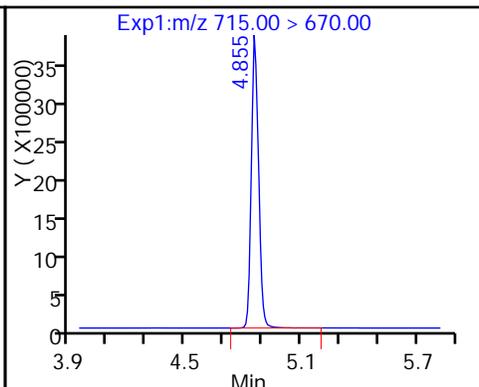
29 Perfluorododecanoic acid



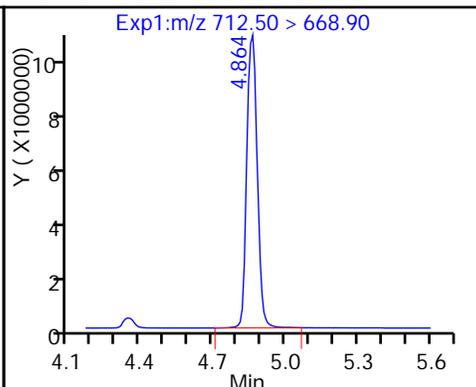
31 Perfluorotridecanoic acid



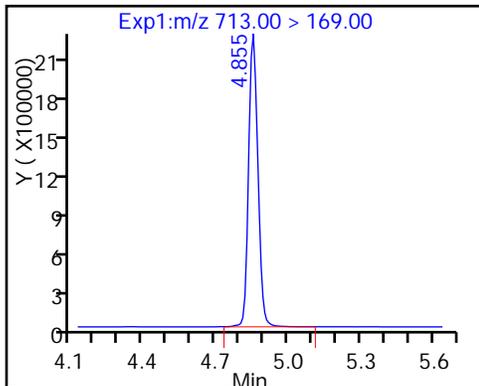
D 32 13C2-PFTeDA



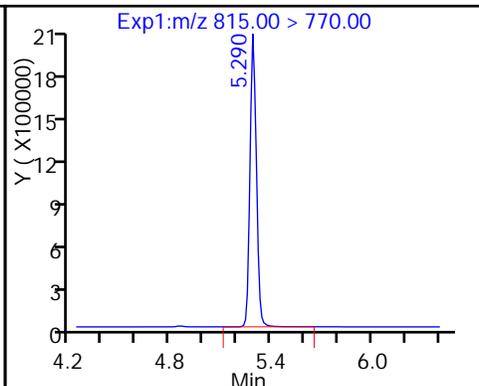
33 Perfluorotetradecanoic acid



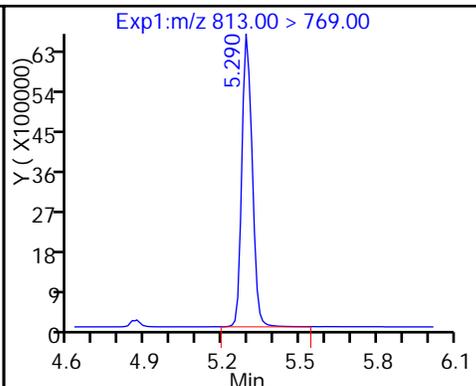
33 Perfluorotetradecanoic acid



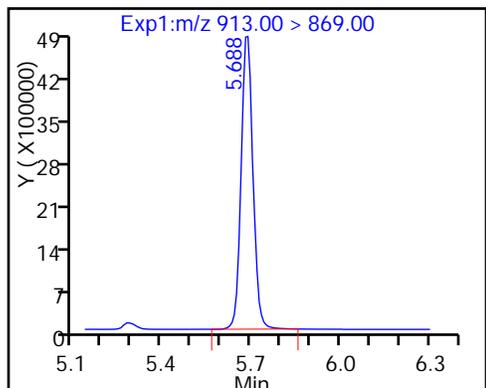
D 34 13C2-PFHxD A



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_010.d
 Lims ID: IC L7
 Client ID:
 Sample Type: IC Calib Level: 7
 Inject. Date: 02-Dec-2016 11:14:32 ALS Bottle#: 43 Worklist Smp#: 10
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L7_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:51:39 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:42:13

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.609	1.615	-0.006	14832847	43.6		87.2	714054	
1 Perfluorobutyric acid	212.90 > 169.00	1.609	1.617	-0.008	69781585	265.7		66.4	349641	
D 4 13C5-PFPeA	267.90 > 223.00	1.910	1.918	-0.008	10386976	38.4		76.9	960901	
3 Perfluoropentanoic acid	262.90 > 219.00	1.910	1.920	-0.010	50360659	232.9		58.2	296264	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.968	1.961	0.007	41429650	107.3		30.3		
	298.90 > 99.00	1.948	1.961	-0.013	44560969		0.93(0.00-0.00)	30.3		
D 6 13C2 PFHxA	315.00 > 270.00	2.230	2.239	-0.009	9976646	40.5		81.0	859246	
7 Perfluorohexanoic acid	313.00 > 269.00	2.230	2.241	-0.011	51025463	267.5		66.9	759942	
D 11 13C4-PFHpA	367.00 > 322.00	2.581	2.599	-0.019	7479568	35.5		71.1	874763	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.581	2.600	-0.020	45359019	293.2		73.3	320883	
D 10 18O2 PFHxS	403.00 > 84.00	2.598	2.614	-0.016	11322930	37.4		79.1	1230464	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.598	2.615	-0.017	77870104	307.0		84.3		
15 Perfluorooctanoic acid	413.00 > 369.00	2.967	2.984	-0.017	47490545	322.1		80.5	2406009	
	413.00 > 169.00	2.967	2.984	-0.017	33779619		1.41(0.90-1.10)	80.5	3107357	
D 14 13C4 PFOA	417.00 > 372.00	2.958	2.984	-0.026	7225722	32.9		65.9	438118	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.967	2.986	-0.019	1.000	59371135	266.1		69.9	
D 17 13C4 PFOS	503.00	> 80.00	3.345	3.365	-0.020		9124233	37.3		78.0	67315
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.345	3.366	-0.021	1.000	80249842	386.6		104	13058140
	499.00	> 99.00	3.345	3.366	-0.021	1.000	19443677		4.13(0.90-1.10)	104	690162
20 Perfluorononanoic acid	463.00	> 419.00	3.353	3.373	-0.020	1.000	41884345	335.5		83.9	392400
D 19 13C5 PFNA	468.00	> 423.00	3.353	3.374	-0.021		6297115	35.4		70.7	371230
D 21 13C8 FOSA	506.00	> 78.00	3.642	3.651	-0.009		18384094	44.4		88.8	592069
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.660	3.656	0.004	1.000	73164187	212.7		53.2	145203
24 Perfluorodecanoic acid	513.00	> 469.00	3.709	3.738	-0.029	1.000	42979124	313.4		78.4	353734
D 23 13C2 PFDA	515.00	> 470.00	3.720	3.738	-0.018		7073000	42.6		85.2	160460
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	4.028	4.049	-0.021	1.000	41362935	345.5		89.6	
D 27 13C2 PFUnA	565.00	> 520.00	4.045	4.070	-0.025		4373854	34.8		69.5	211760
28 Perfluoroundecanoic acid	563.00	> 519.00	4.045	4.071	-0.026	1.000	31200994	336.6		84.2	378305
D 30 13C2 PFDoA	615.00	> 570.00	4.353	4.370	-0.017		5148784	44.0		88.1	149077
29 Perfluorododecanoic acid	613.00	> 569.00	4.353	4.370	-0.017	1.000	32927149	337.8		84.5	44886
31 Perfluorotridecanoic acid	663.00	> 619.00	4.610	4.636	-0.026	1.000	34277350	320.0		80.0	59163 M
D 32 13C2-PFTeDA	715.00	> 670.00	4.860	4.882	-0.022		10396514	43.0		86.0	969333
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.860	4.884	-0.024	1.000	50372831	253.2		63.3	26628
	713.00	> 169.00	4.851	4.884	-0.033	0.998	12426058		4.05(0.00-0.00)	63.3	473842
D 34 13C2-PFHxDA	815.00	> 770.00	5.300	5.320	-0.020		6001714	46.0		92.1	121116
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.300	5.320	-0.020	1.000	35113710	345.4		86.3	34228
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.682	5.720	-0.038	1.000	31098706	384.1		96.0	54097

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L7_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_010.d

Injection Date: 02-Dec-2016 11:14:32

Instrument ID: A8_N

Lims ID: IC L7

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 43

Worklist Smp#: 10

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

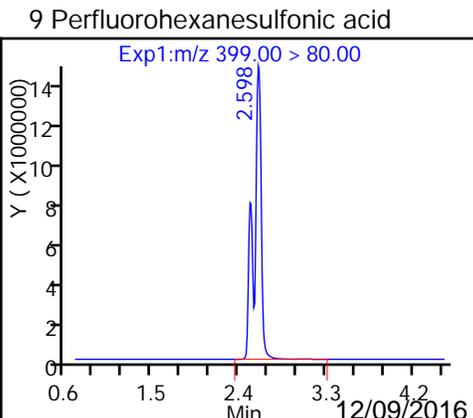
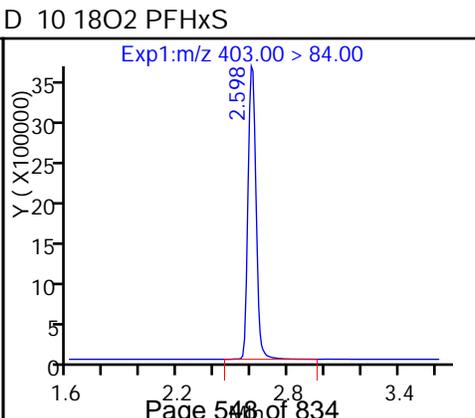
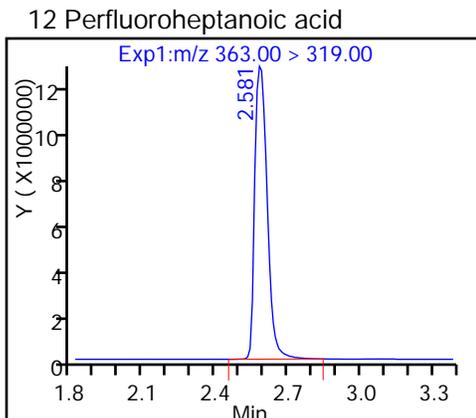
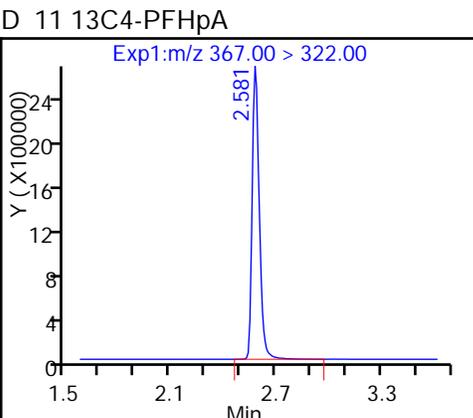
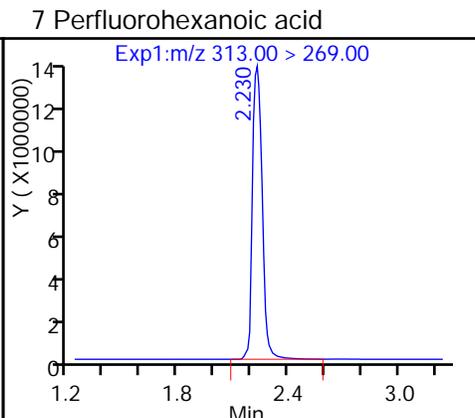
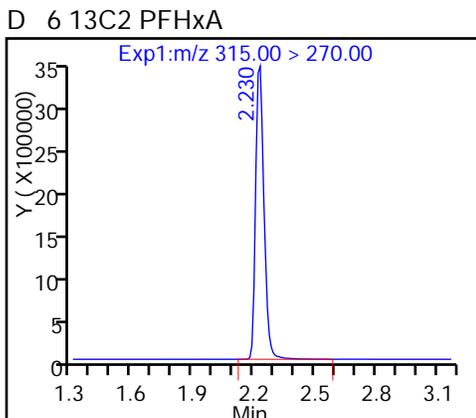
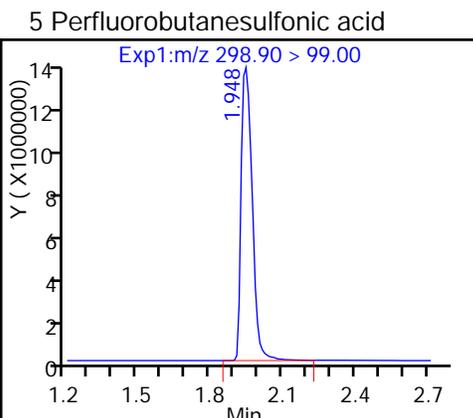
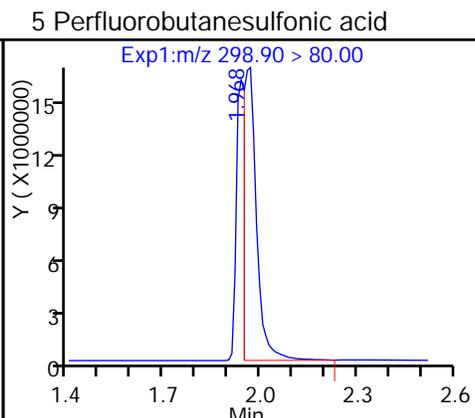
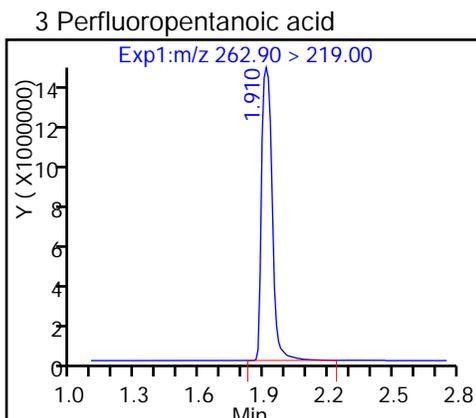
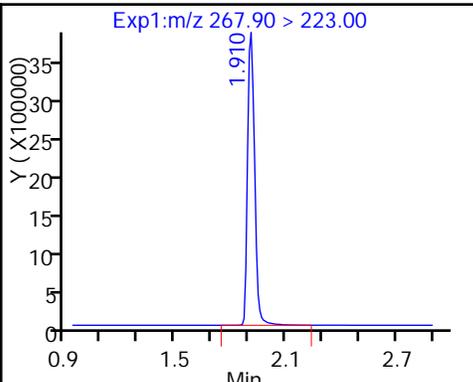
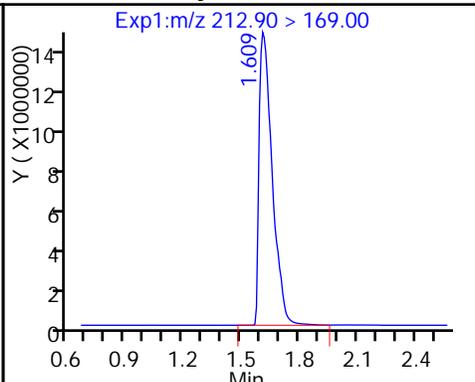
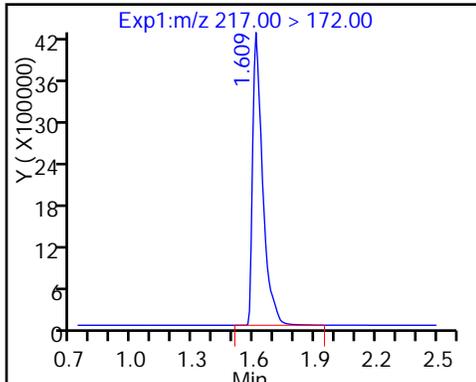
Method: A8_N

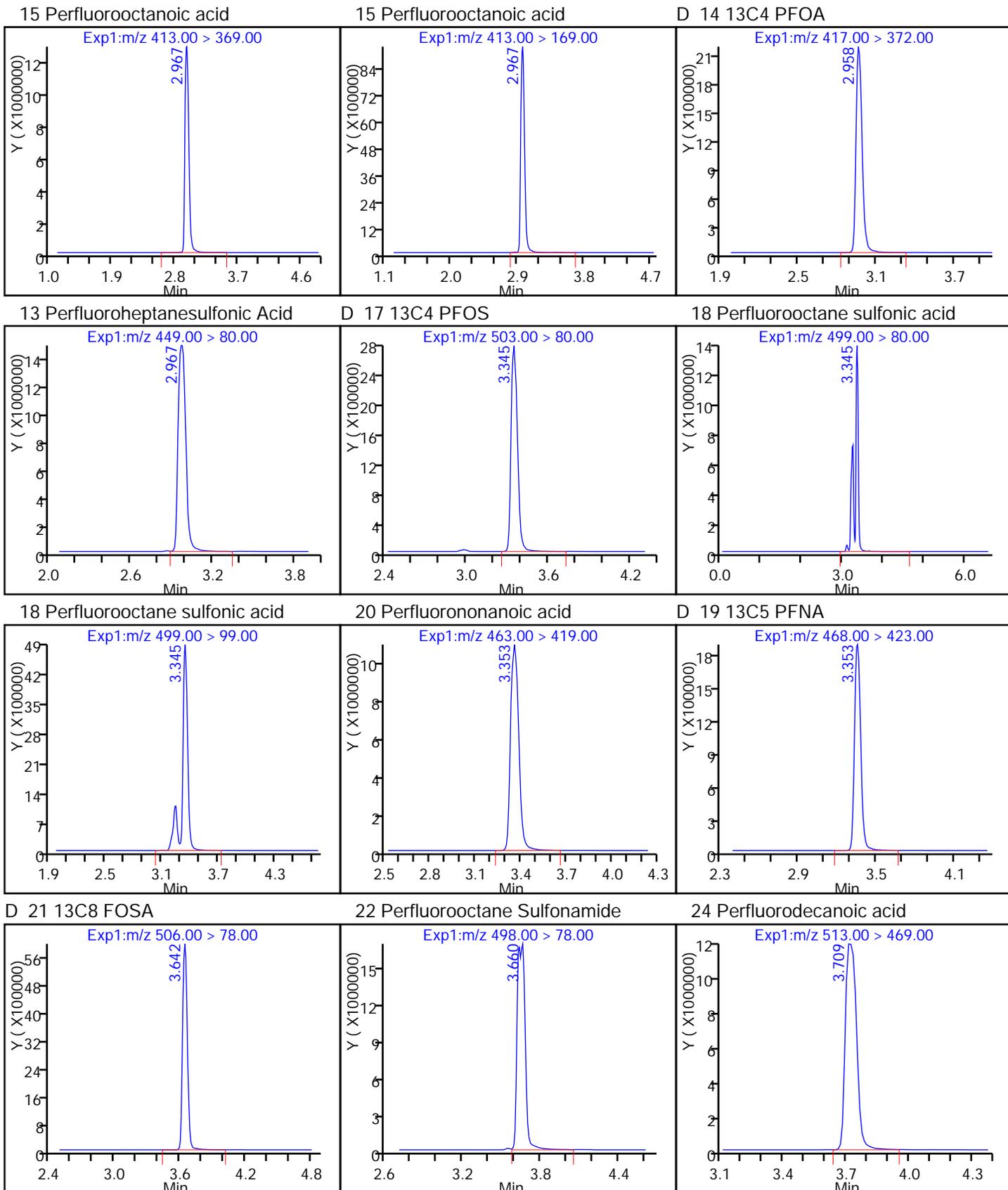
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

D 4 13C5-PFPeA

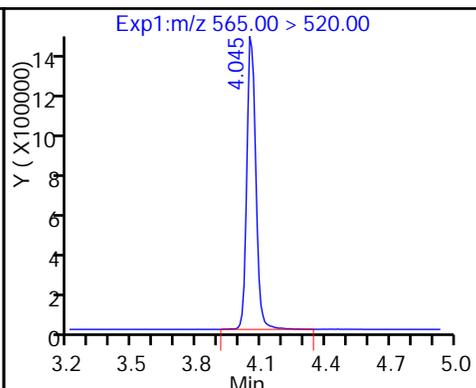
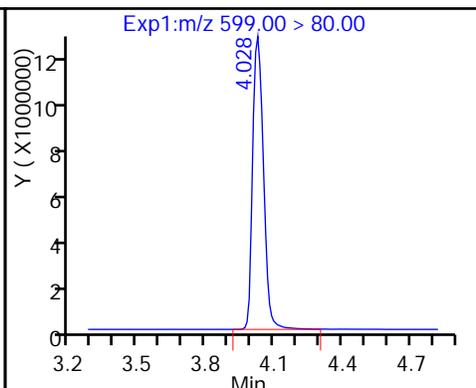
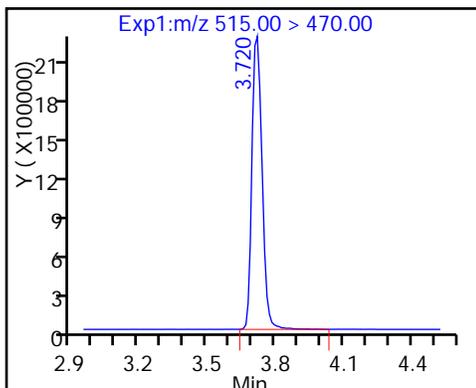




D 23 13C2 PFDA

26 Perfluorodecane Sulfonic acid

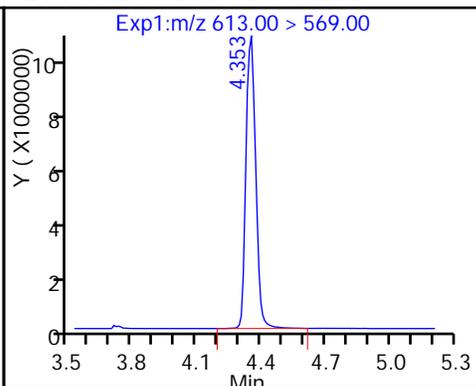
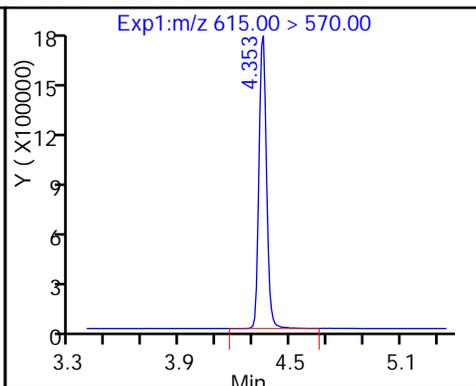
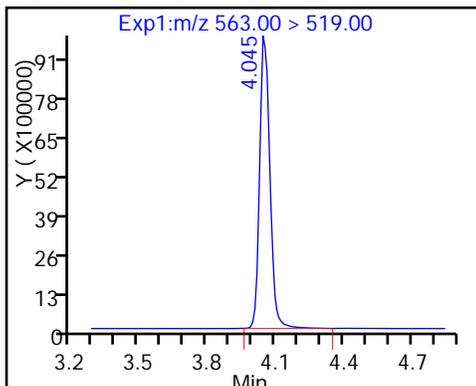
D 27 13C2 PFUa



28 Perfluoroundecanoic acid

D 30 13C2 PFDa

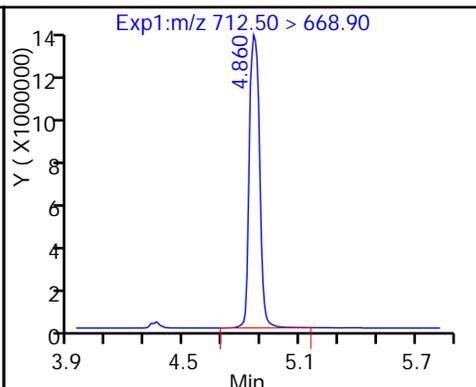
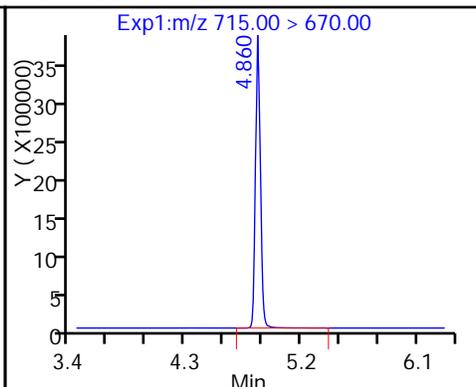
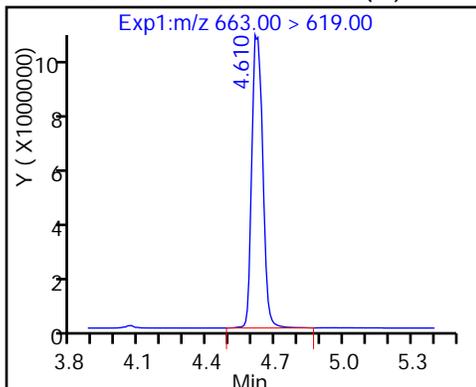
29 Perfluorododecanoic acid



31 Perfluorotridecanoic acid (M)

D 32 13C2-PFTeDA

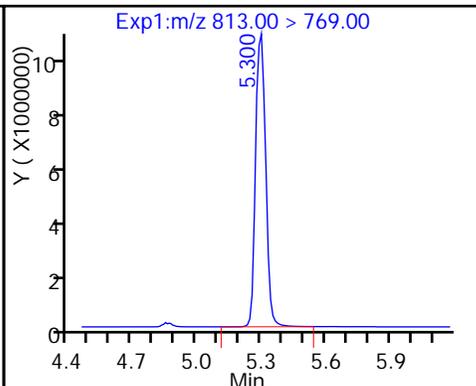
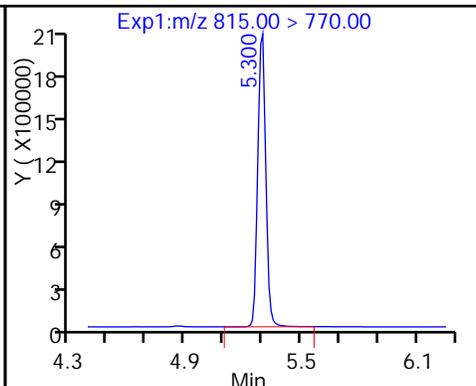
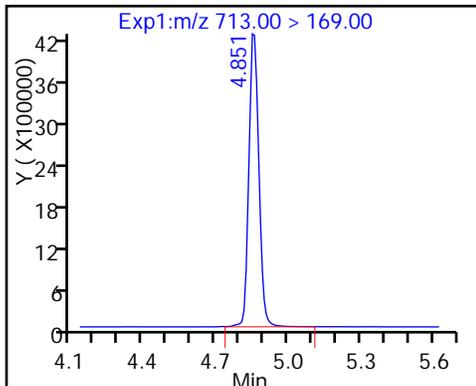
33 Perfluorotetradecanoic acid



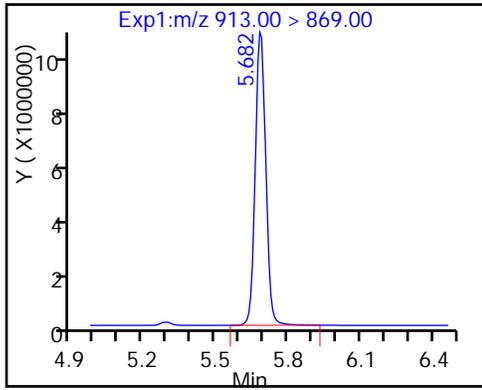
33 Perfluorotetradecanoic acid

D 34 13C2-PFHxDA

35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento

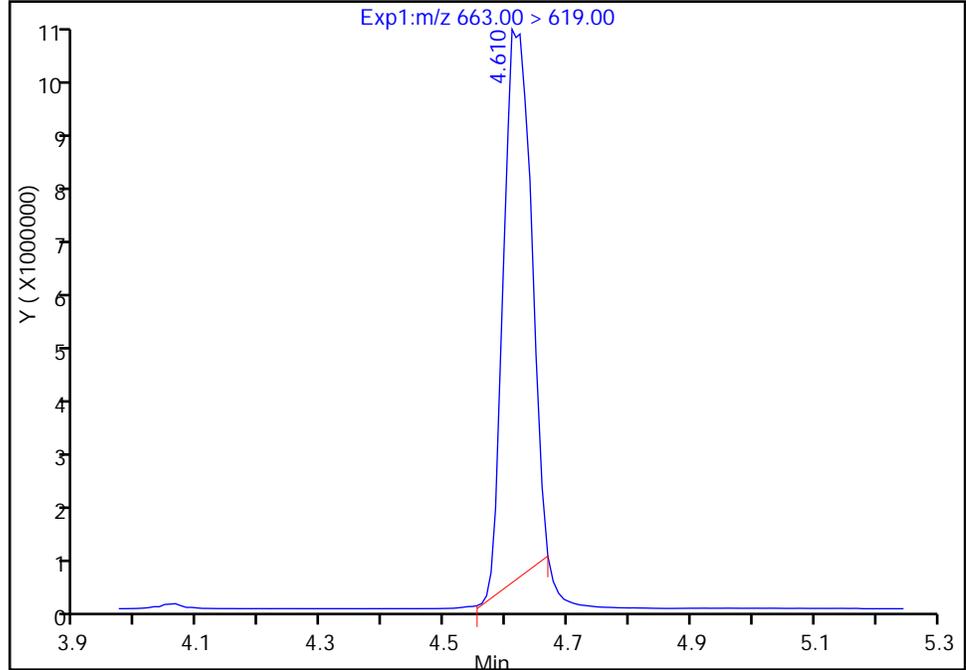
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Injection Date: 02-Dec-2016 11:14:32 Instrument ID: A8_N
Lims ID: IC L7
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 43 Worklist Smp#: 10
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

31 Perfluorotridecanoic acid, CAS: 72629-94-8

Signal: 1

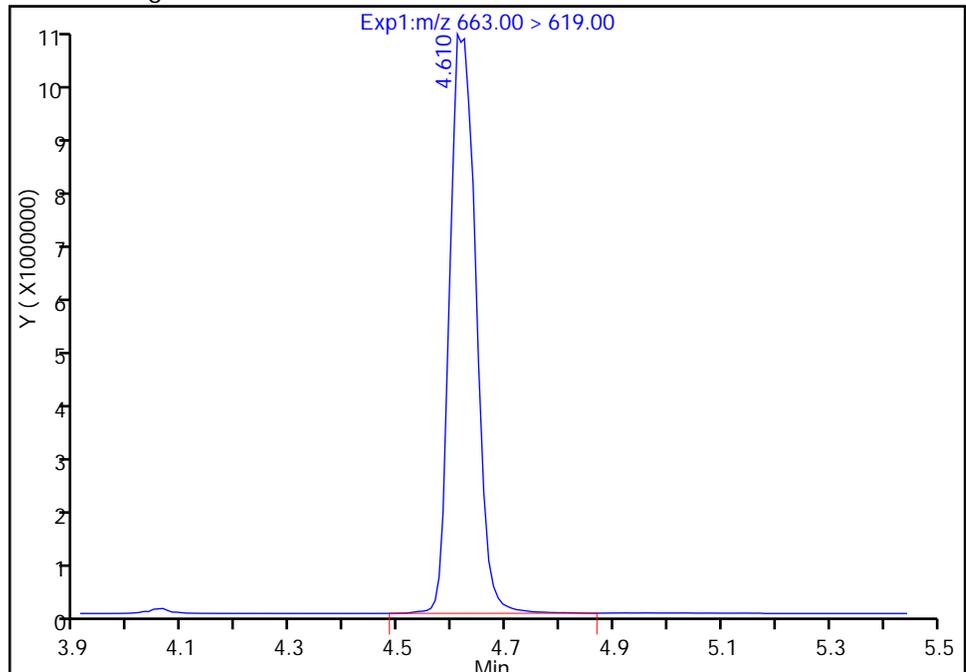
RT: 4.61
Area: 30068675
Amount: 293.1955
Amount Units: ng/ml

Processing Integration Results



RT: 4.61
Area: 34277350
Amount: 319.9934
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 02-Dec-2016 13:42:13

Audit Action: Manually Integrated

Audit Reason: Incomplete Integration

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_014.d
 Lims ID: IC L1 Add-on
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 02-Dec-2016 11:44:30 ALS Bottle#: 46 Worklist Smp#: 14
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L1 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:41 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:56:57

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.909	2.913	-0.004	1.000	51838	0.4953	104		
D 47 M2-6:2FTS	429.00 > 409.00	2.925	2.915	0.010		5925112	42.5	89.6		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.668	3.674	-0.006	0.997	44037	0.4434	92.6		
D 42 M2-8:2FTS	529.00 > 509.00	3.678	3.676	0.002		5634375	42.2	88.1		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.843	3.839	0.004		3880040	48.5	97.0		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.851	3.847	0.004	1.002	30895	0.4583	91.7		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.011	4.009	0.002		4464530	50.1	100		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.020	4.016	0.004	1.002	30016	0.4382	87.6		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.108	4.105	0.003		4681022	44.0	88.0		
54 MeFOSA	512.00 > 169.00	4.115	4.110	0.005	1.000	37277	0.5000	100		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.293	4.290	0.003		4329074	42.5	85.0		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.293	4.299	-0.006	1.000	35535	0.4830	96.6		

Reagents:

LCPFC2-L1_00002

Amount Added: 1.00

Units: mL

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_014.d

Injection Date: 02-Dec-2016 11:44:30

Instrument ID: A8_N

Lims ID: IC L1 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 46

Worklist Smp#: 14

Injection Vol: 2.0 ul

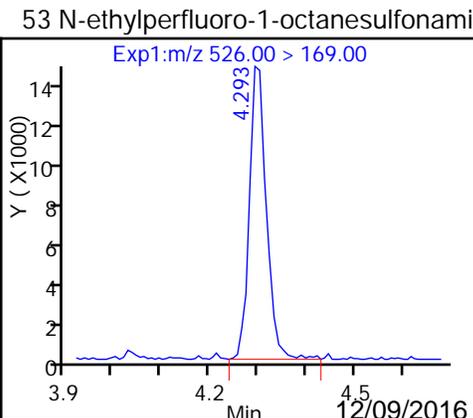
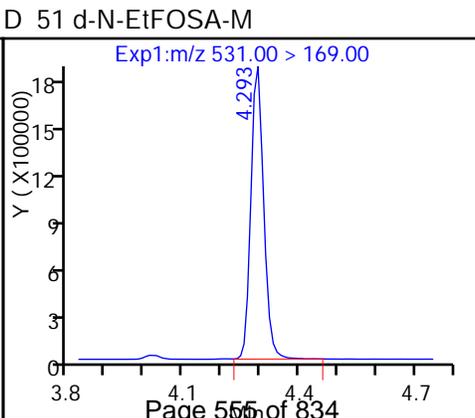
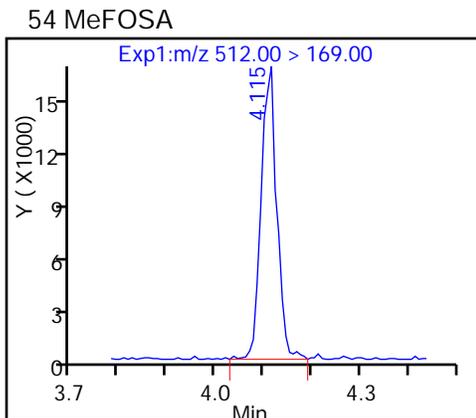
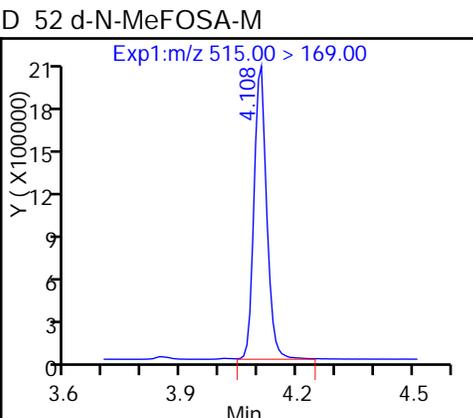
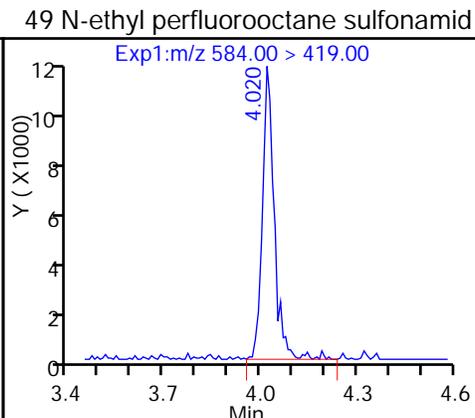
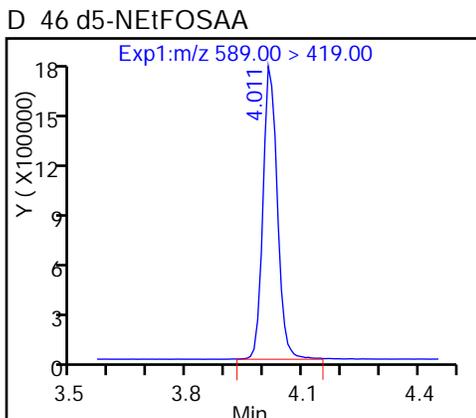
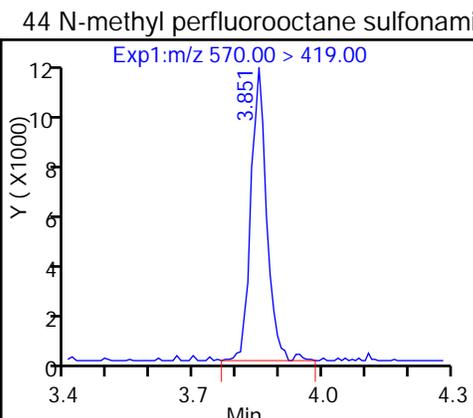
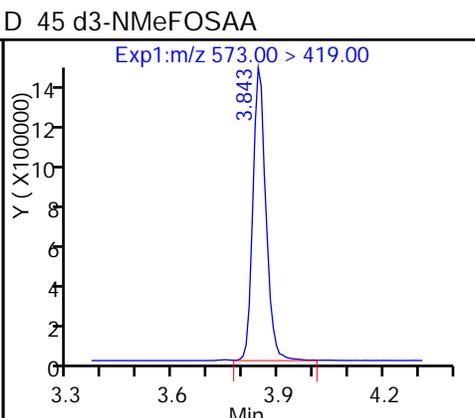
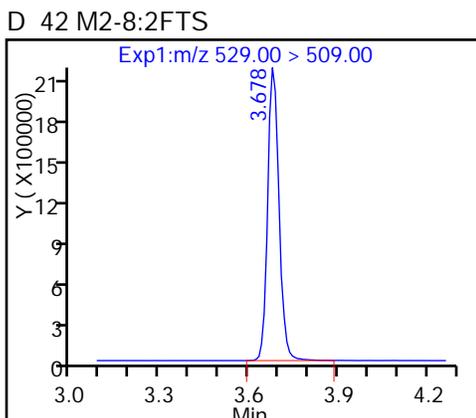
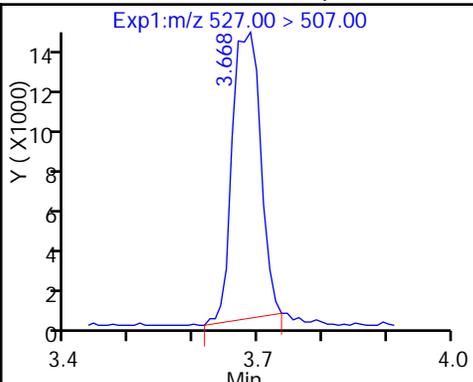
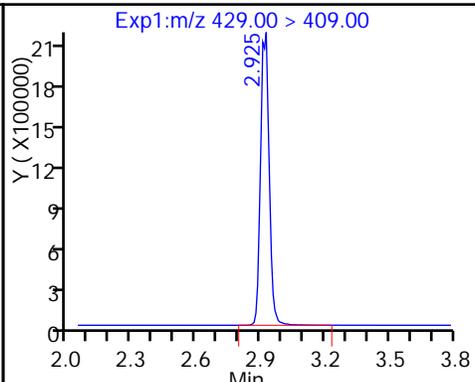
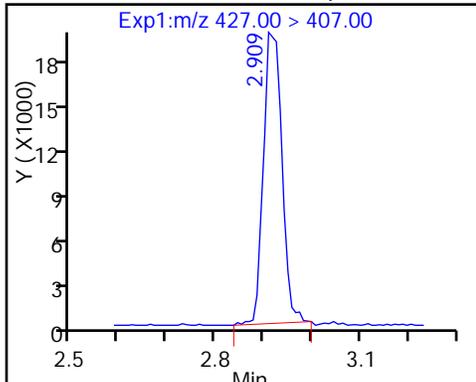
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_015.d
 Lims ID: IC L2 Add-on
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 02-Dec-2016 11:52:00 ALS Bottle#: 47 Worklist Smp#: 15
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L2 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:42 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:57:07

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.920	2.913	0.007	1.000	88396	0.8738	92.2		
D 47 M2-6:2FTS	429.00 > 409.00	2.912	2.915	-0.003		5726516	41.1	86.6		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.684	3.674	0.010	1.000	94573	1.03	107		
D 42 M2-8:2FTS	529.00 > 509.00	3.684	3.676	0.008		5213436	39.1	81.5		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.848	3.839	0.009		3606216	45.1	90.2		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.857	3.847	0.010	1.002	59295	0.9465	94.6		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.017	4.009	0.008		4134123	46.4	92.8		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.026	4.016	0.010	1.002	56004	0.8830	88.3		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.106	4.105	0.001		5181881	48.7	97.5		
54 MeFOSA	512.00 > 169.00	4.114	4.110	0.004	1.000	73327	0.8885	88.8		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.292	4.290	0.002		5048951	49.5	99.1		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.301	4.299	0.002	1.000	78846	0.9189	91.9		

Reagents:

LCPFC2-L2_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_015.d

Injection Date: 02-Dec-2016 11:52:00

Instrument ID: A8_N

Lims ID: IC L2 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 47

Worklist Smp#: 15

Injection Vol: 2.0 ul

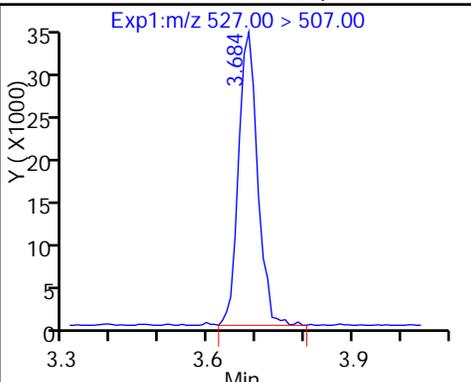
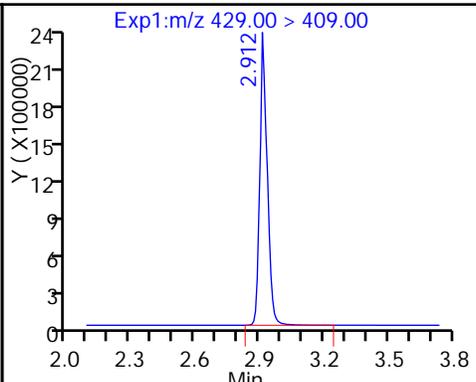
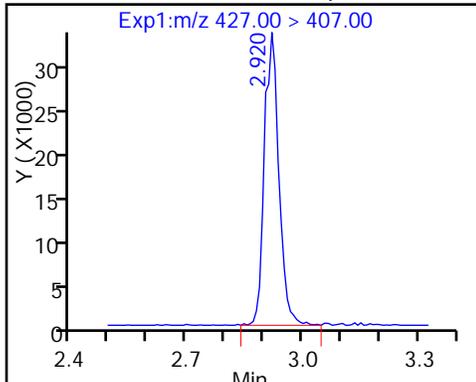
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

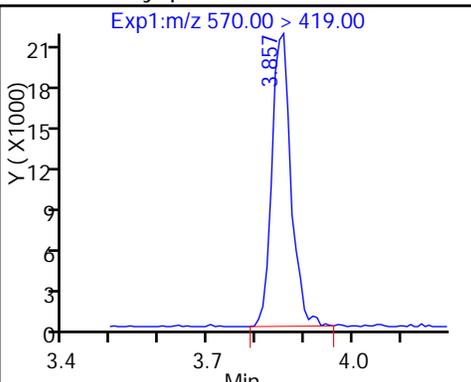
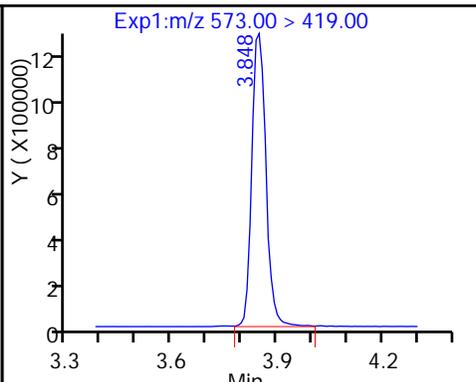
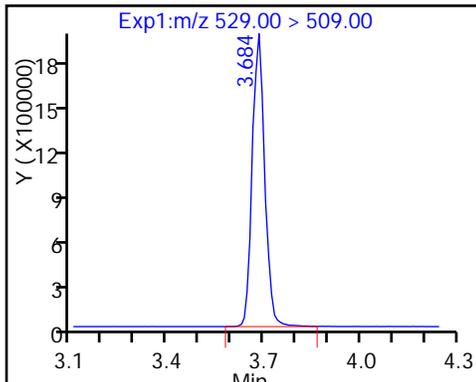
43 Sodium 1H,1H,2H,2H-perfluorooctane



D 42 M2-8:2FTS

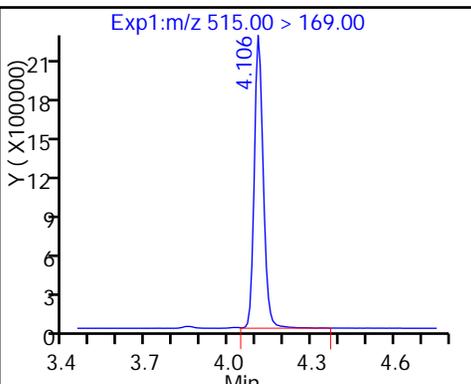
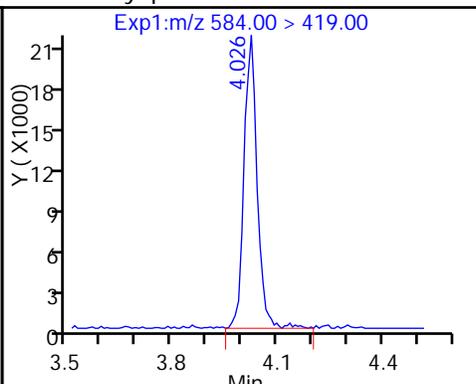
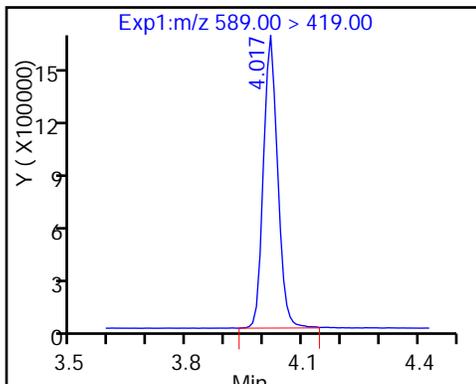
D 45 d3-NMeFOSAA

44 N-methyl perfluorooctane sulfonami



D 46 d5-NEtFOSAA

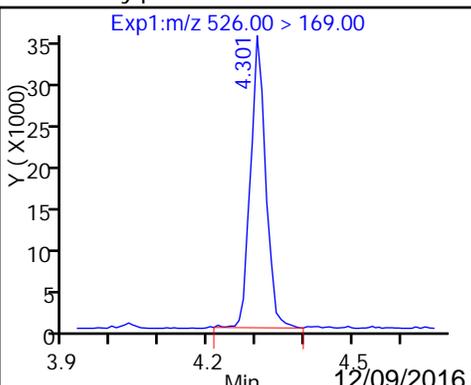
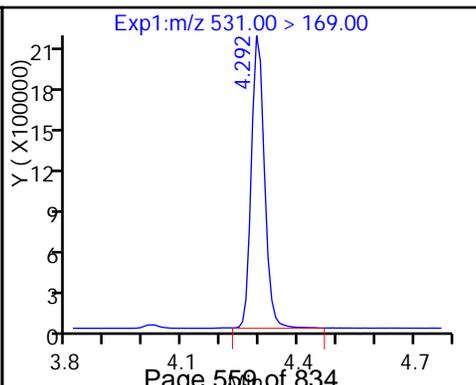
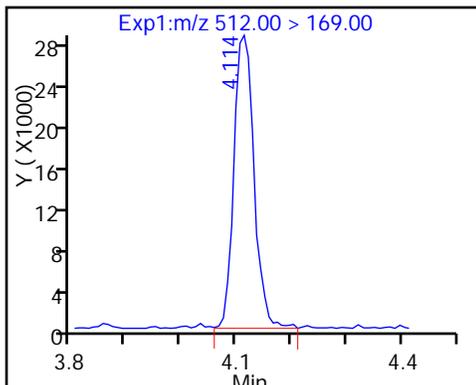
49 N-ethyl perfluorooctane sulfonamid D 52 d-N-MeFOSA-M



54 MeFOSA

D 51 d-N-EtFOSA-M

53 N-ethylperfluoro-1-octanesulfonami



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_016.d
 Lims ID: IC L3 Add-on
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 02-Dec-2016 11:59:31 ALS Bottle#: 48 Worklist Smp#: 16
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L3 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:44 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:57:18

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.910	2.913	-0.003	1.000	492719	4.19	88.4		
D 47 M2-6:2FTS	429.00 > 409.00	2.918	2.915	0.003		6658764	47.8	101		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.673	3.674	-0.001	1.000	431909	4.10	85.7		
D 42 M2-8:2FTS	529.00 > 509.00	3.673	3.676	-0.003		5971752	44.7	93.4		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.838	3.839	-0.001		4140205	51.8	104		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.847	3.847	0.0	1.002	289882	4.03	80.6		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.006	4.009	-0.003		4489327	50.4	101		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.015	4.016	-0.001	1.002	283960	4.12	82.5		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.104	4.105	-0.001		5362314	50.4	101		
54 MeFOSA	512.00 > 169.00	4.104	4.110	-0.006	1.000	360259	4.22	84.4		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.287	4.290	-0.003		5205212	51.1	102		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.297	4.299	-0.002	1.000	365734	4.13	82.7		

Reagents:

LCPFC2-L3_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_016.d

Injection Date: 02-Dec-2016 11:59:31

Instrument ID: A8_N

Lims ID: IC L3 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 48

Worklist Smp#: 16

Injection Vol: 2.0 ul

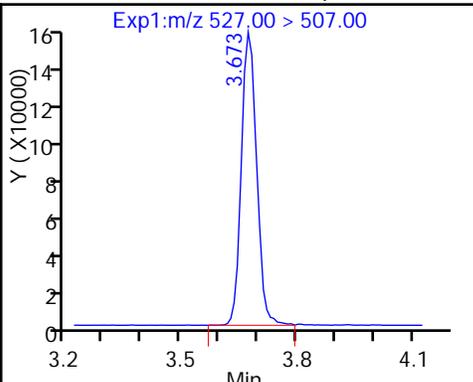
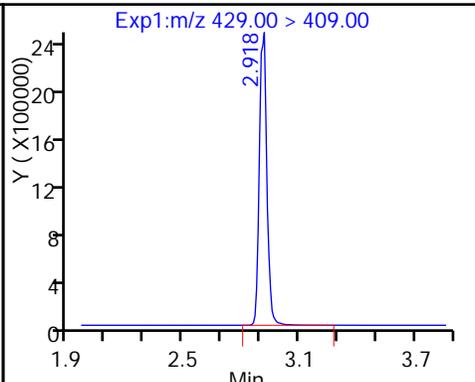
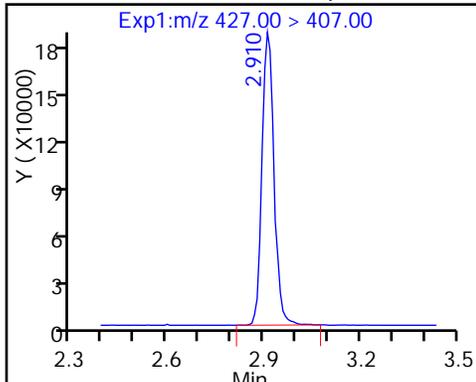
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

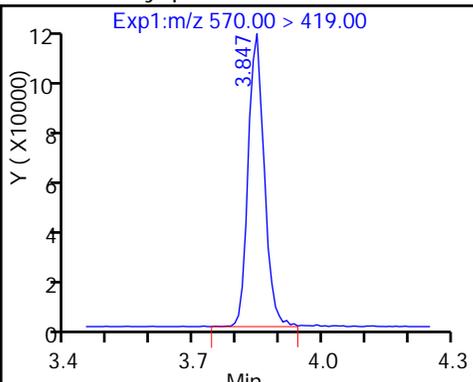
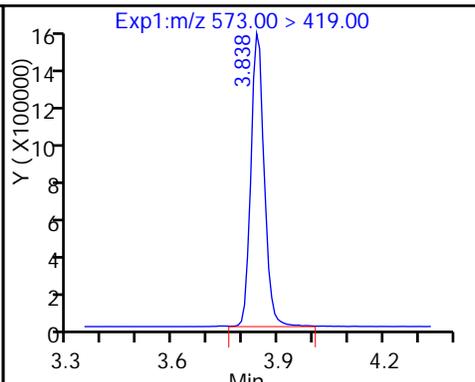
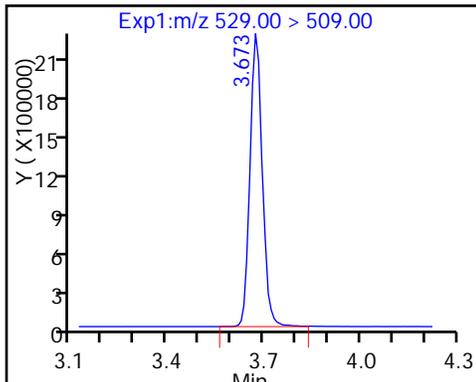
43 Sodium 1H,1H,2H,2H-perfluorooctane



D 42 M2-8:2FTS

D 45 d3-NMeFOSAA

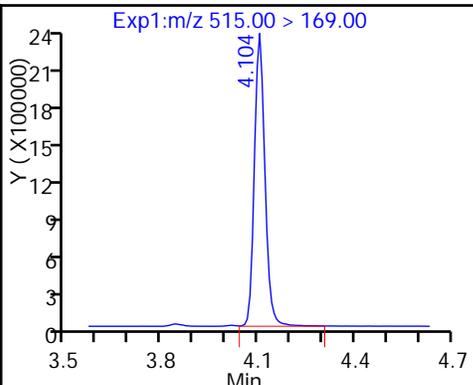
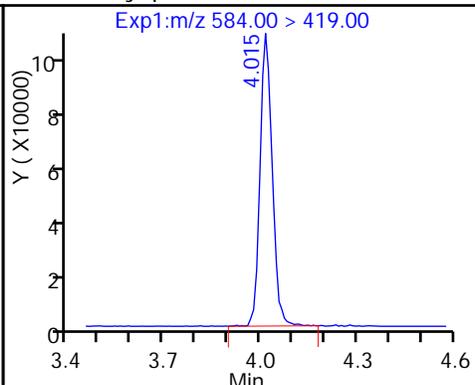
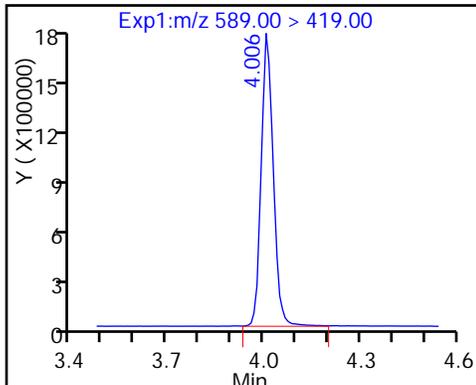
44 N-methyl perfluorooctane sulfonami



D 46 d5-NEtFOSAA

49 N-ethyl perfluorooctane sulfonamid

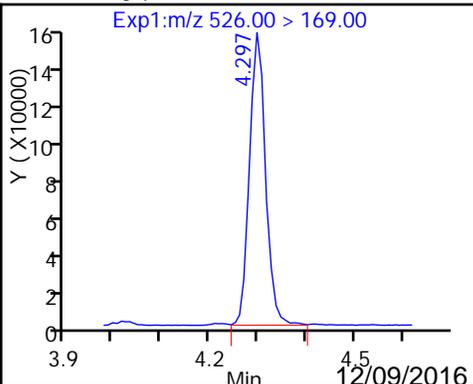
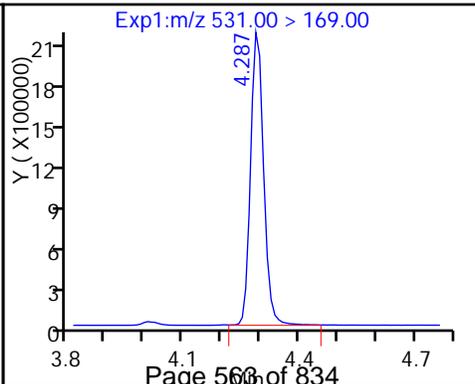
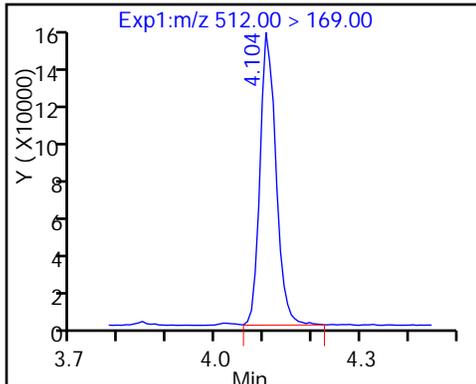
D 52 d-N-MeFOSA-M



54 MeFOSA

D 51 d-N-EtFOSA-M

53 N-ethylperfluoro-1-octanesulfonami



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_017.d
 Lims ID: IC L4 Add-on
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 02-Dec-2016 12:07:00 ALS Bottle#: 49 Worklist Smp#: 17
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L4 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:46 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:56:04

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.914	2.913	0.001	1.000	2900008	20.8	110		
D 47 M2-6:2FTS	429.00 > 409.00	2.906	2.915	-0.009		7906261	56.8	120		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.678	3.674	0.004	1.000	2799477	21.5	112		
D 42 M2-8:2FTS	529.00 > 509.00	3.678	3.676	0.002		7389674	55.4	116		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.843	3.839	0.004		4953346	61.9	124		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.843	3.847	-0.004	1.000	1829869	21.3	106		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.011	4.009	0.002		5242030	58.9	118		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.011	4.016	-0.005	1.000	1881436	23.4	117		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.104	4.105	-0.001		5889991	55.4	111		
54 MeFOSA	512.00 > 169.00	4.112	4.110	0.002	1.000	2129108	22.7	113		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.293	4.290	0.003		5611197	55.1	110		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.302	4.299	0.003	1.000	2187091	22.9	115		

Reagents:

LCPFC2-L4_00003

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_017.d

Injection Date: 02-Dec-2016 12:07:00

Instrument ID: A8_N

Lims ID: IC L4 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 49

Worklist Smp#: 17

Injection Vol: 2.0 ul

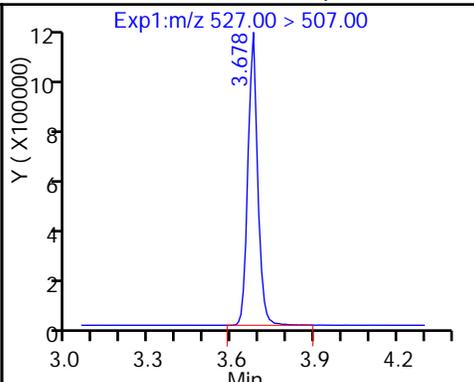
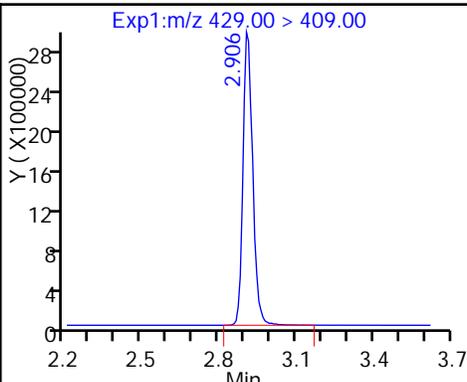
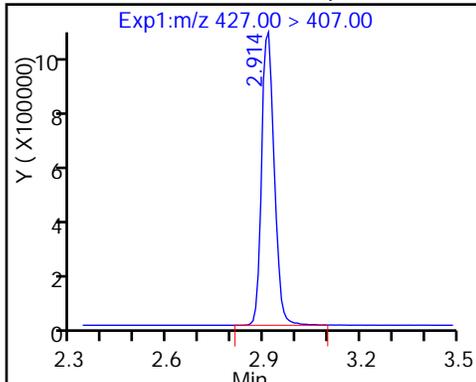
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

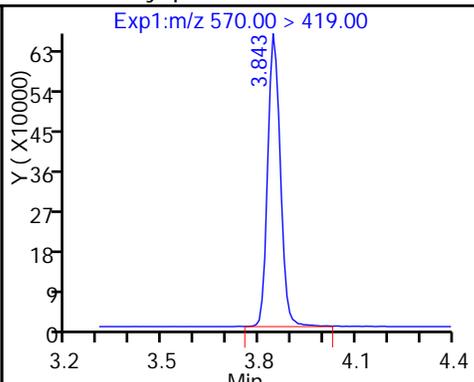
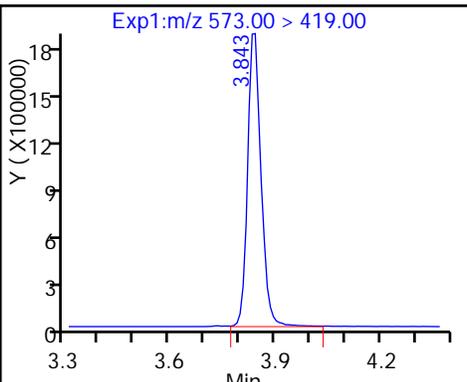
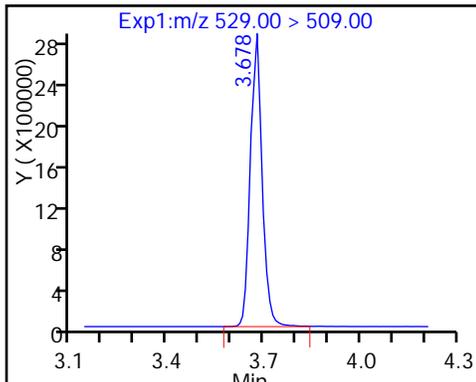
43 Sodium 1H,1H,2H,2H-perfluorooctane



D 42 M2-8:2FTS

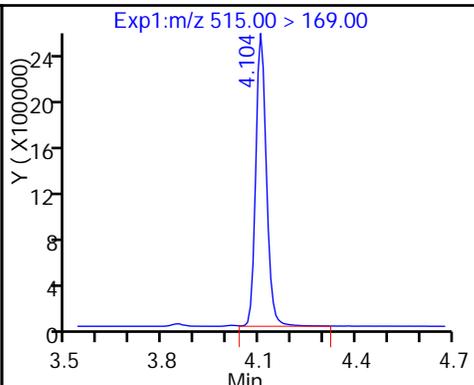
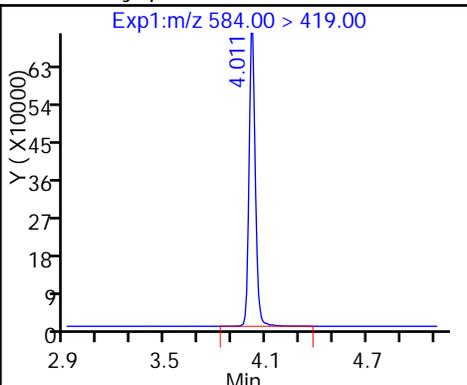
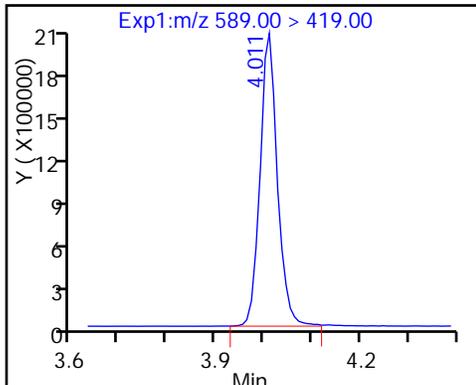
D 45 d3-NMeFOSAA

44 N-methyl perfluorooctane sulfonami



D 46 d5-NEtFOSAA

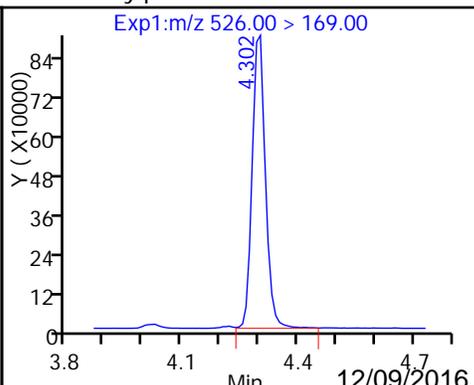
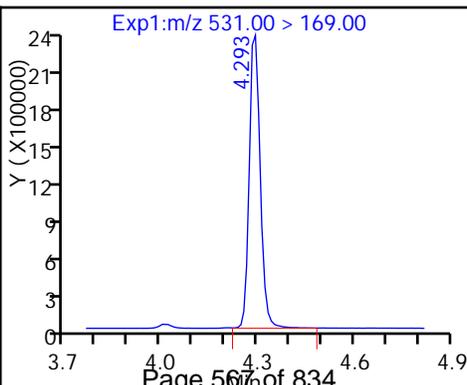
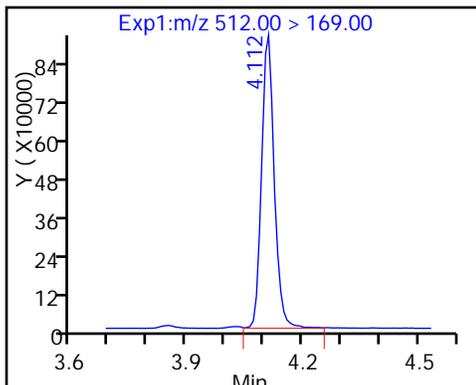
49 N-ethyl perfluorooctane sulfonamid D 52 d-N-MeFOSA-M



54 MeFOSA

D 51 d-N-EtFOSA-M

53 N-ethylperfluoro-1-octanesulfonami



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_018.d
 Lims ID: IC L5 Add-on
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 02-Dec-2016 12:14:30 ALS Bottle#: 50 Worklist Smp#: 18
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L5 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:47 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:57:38

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.919	2.913	0.006	1.000	5633441	57.1	120		
D 47 M2-6:2FTS	429.00 > 409.00	2.911	2.915	-0.004		5586093	40.1	84.4		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.674	3.674	0.0	1.000	5753486	55.8	116		
D 42 M2-8:2FTS	529.00 > 509.00	3.674	3.676	-0.002		5851766	43.8	91.5		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.839	3.839	0.0		3442292	43.0	86.1		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.847	3.847	0.0	1.002	3696988	61.8	124		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.007	4.009	-0.002		3869672	43.4	86.9		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.016	4.016	0.0	1.002	3546696	59.7	119		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.100	4.105	-0.005		4967092	46.7	93.4		
54 MeFOSA	512.00 > 169.00	4.108	4.110	-0.002	1.000	4836683	61.1	122		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.285	4.290	-0.005		4857179	47.7	95.3		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.295	4.299	-0.004	1.000	5044937	61.1	122		

Reagents:

LCPFC2-L5_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_018.d

Injection Date: 02-Dec-2016 12:14:30

Instrument ID: A8_N

Lims ID: IC L5 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 50

Worklist Smp#: 18

Injection Vol: 2.0 ul

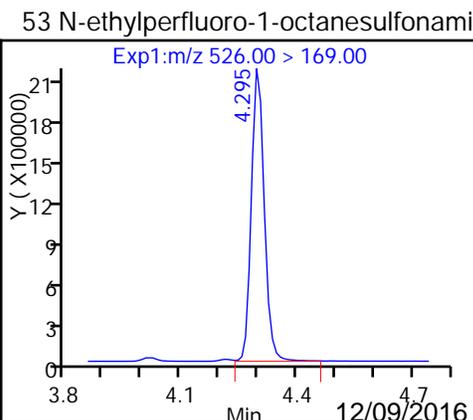
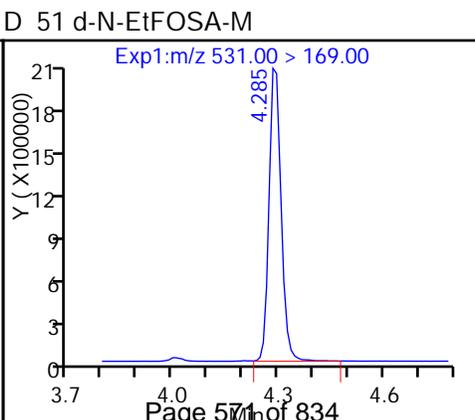
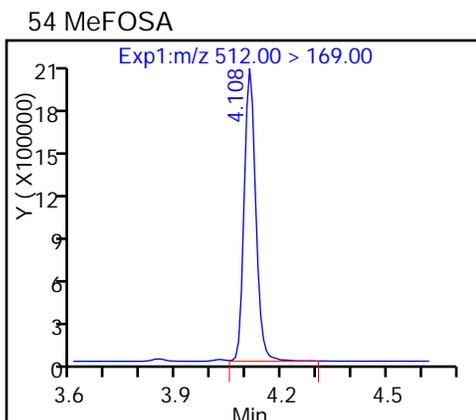
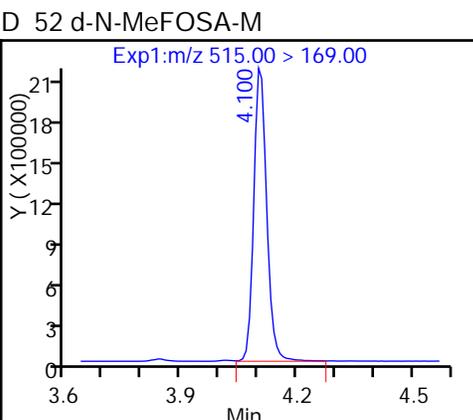
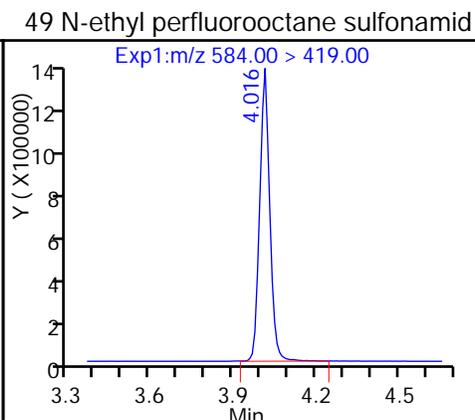
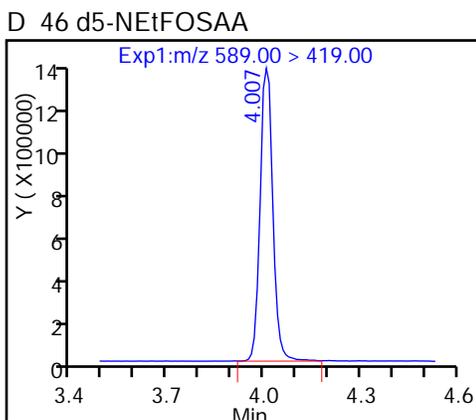
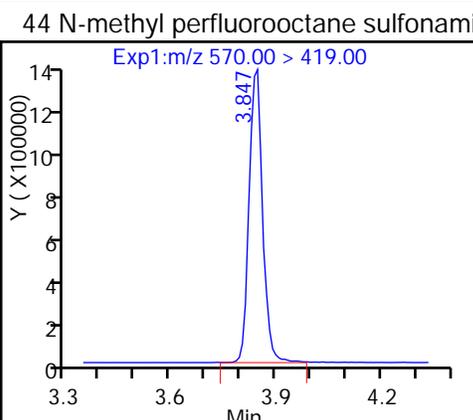
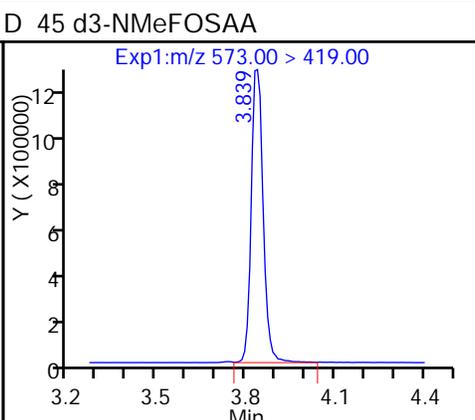
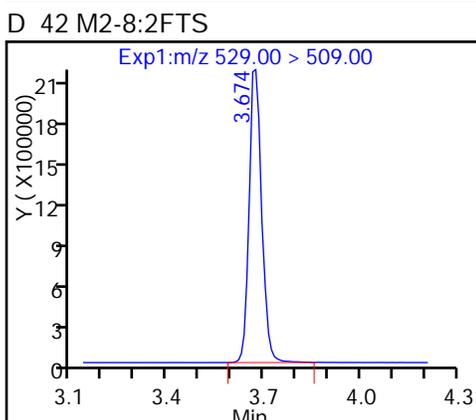
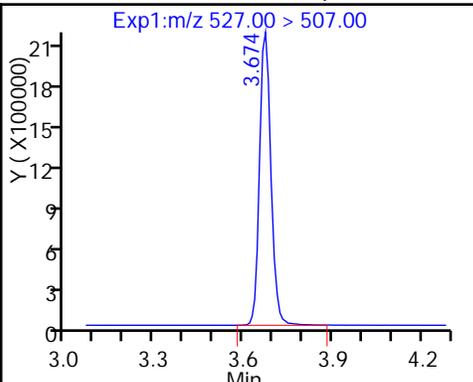
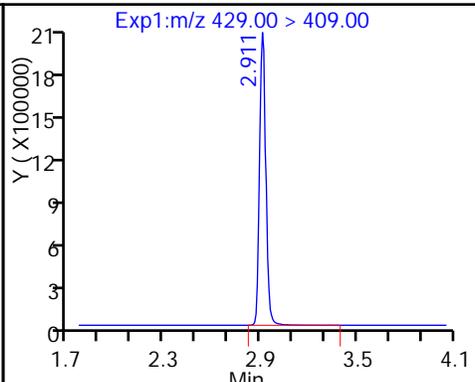
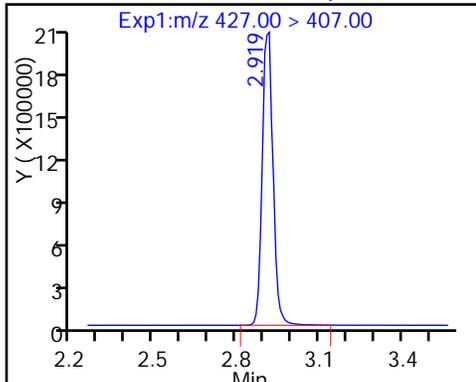
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_019.d
 Lims ID: IC L6 Add-on
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 02-Dec-2016 12:22:00 ALS Bottle#: 51 Worklist Smp#: 19
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L6 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:49 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:57:48

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.904	2.913	-0.009	1.000	20031946	169.7	89.5		
D 47 M2-6:2FTS	429.00 > 409.00	2.912	2.915	-0.003		6682949	48.0	101		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.667	3.674	-0.007	1.000	19463965	164.3	85.7		
D 42 M2-8:2FTS	529.00 > 509.00	3.667	3.676	-0.009		6721067	50.3	105		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.824	3.839	-0.015		4016571	50.2	100		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.832	3.847	-0.015	1.002	13773335	197.4	98.7		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.000	4.009	-0.009		4474975	50.2	100		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.009	4.016	-0.007	1.002	14065497	204.9	102		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.102	4.105	-0.003		4927555	46.3	92.7		
54 MeFOSA	512.00 > 169.00	4.102	4.110	-0.008	1.000	15814534	201.5	101		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.286	4.290	-0.004		4696896	46.1	92.2		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.296	4.299	-0.003	1.000	16550291	207.3	104		

Reagents:

LCPFC2-L6_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_019.d

Injection Date: 02-Dec-2016 12:22:00

Instrument ID: A8_N

Lims ID: IC L6 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 51

Worklist Smp#: 19

Injection Vol: 2.0 ul

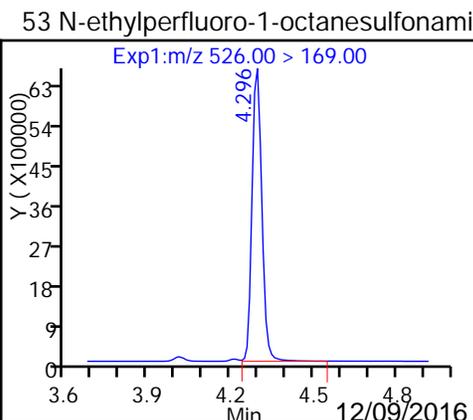
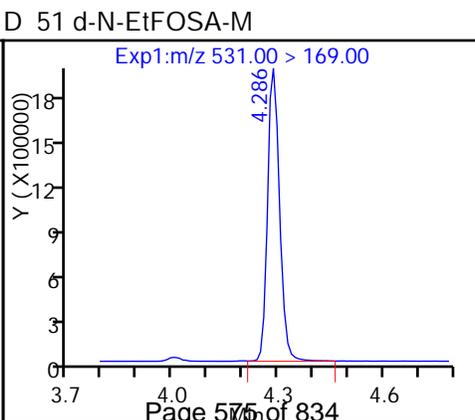
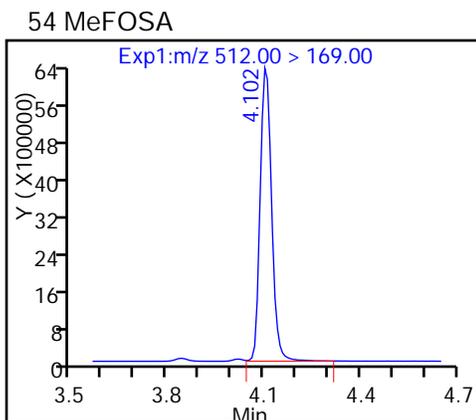
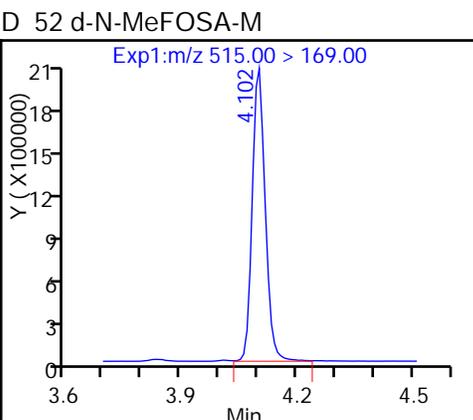
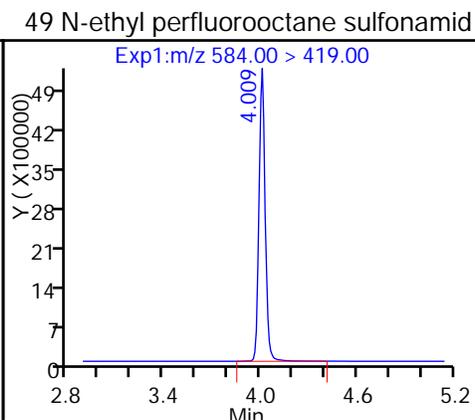
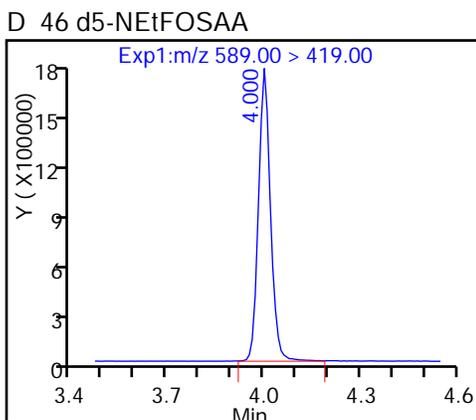
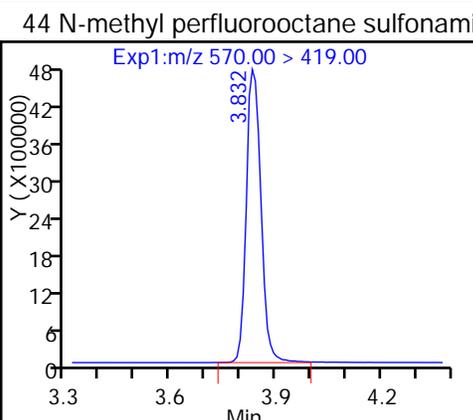
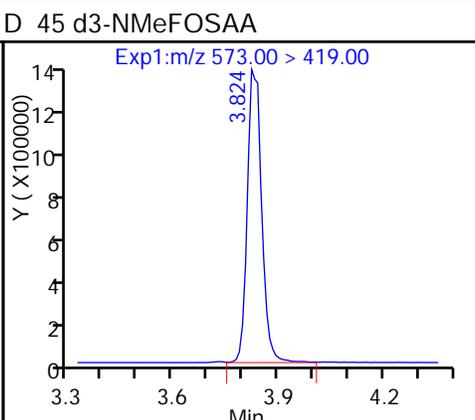
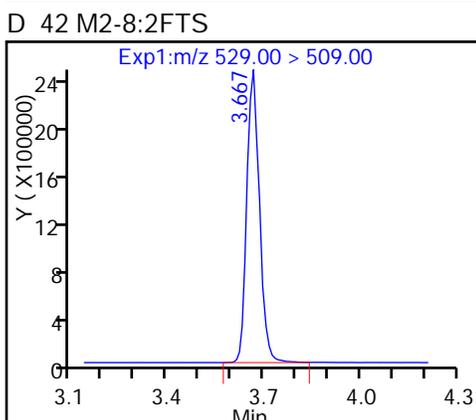
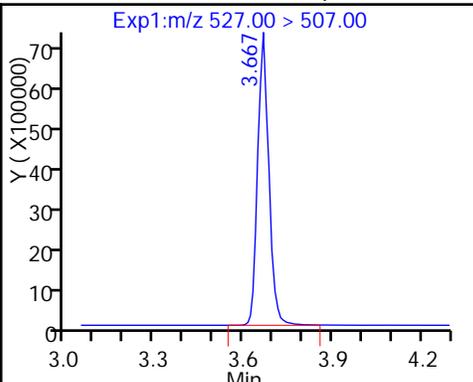
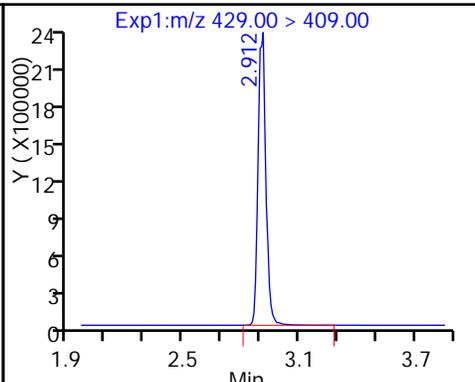
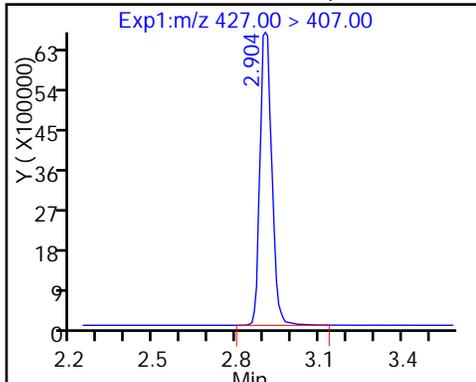
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Lims ID: IC L7 Add-on
 Client ID:
 Sample Type: IC Calib Level: 7
 Inject. Date: 02-Dec-2016 12:29:29 ALS Bottle#: 52 Worklist Smp#: 20
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L7 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 14:52:50 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK011

First Level Reviewer: chandrasenas Date: 02-Dec-2016 13:58:03

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.917	2.913	0.004	1.000	37224925	269.4	71.0		
D 47 M2-6:2FTS	429.00 > 409.00	2.917	2.915	0.002		7823475	56.2	118		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.676	3.674	0.002	1.000	37119838	263.9	68.9		
D 42 M2-8:2FTS	529.00 > 509.00	3.676	3.676	0.0		7980965	59.8	125		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.841	3.839	0.002		3959155	49.5	99.0		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.849	3.847	0.002	1.002	28724886	417.6	104		
D 46 d5-NEtFOSAA	589.00 > 419.00	4.009	4.009	0.0		4497788	50.5	101		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	4.018	4.016	0.002	1.002	28356327	410.9	103		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.110	4.105	0.005		6208758	58.4	117		
54 MeFOSA	512.00 > 169.00	4.118	4.110	0.008	1.000	35701676	361.0	90.3		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.299	4.290	0.009		5918873	58.1	116		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.309	4.299	0.010	1.000	35493503	352.9	88.2		

Reagents:

LCPFC2-L7_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Injection Date: 02-Dec-2016 12:29:29

Instrument ID: A8_N

Lims ID: IC L7 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 52

Worklist Smp#: 20

Injection Vol: 2.0 ul

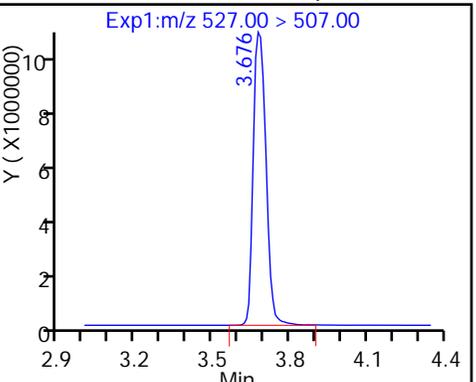
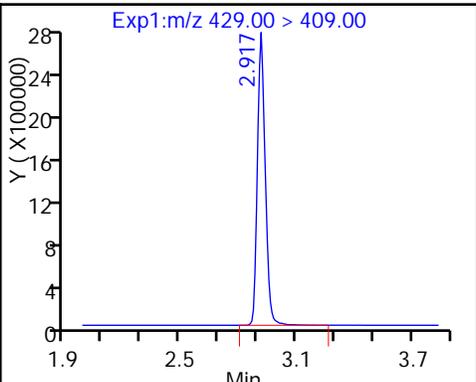
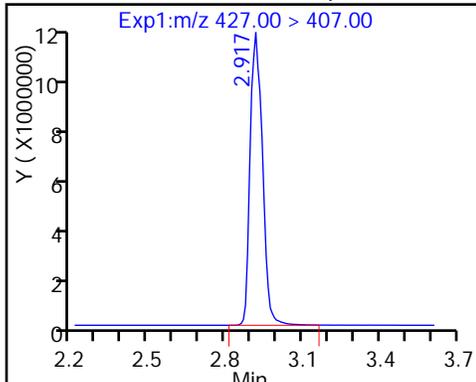
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

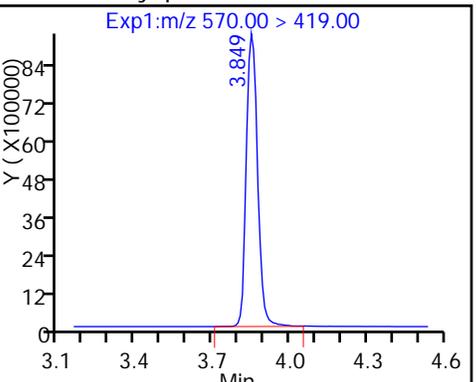
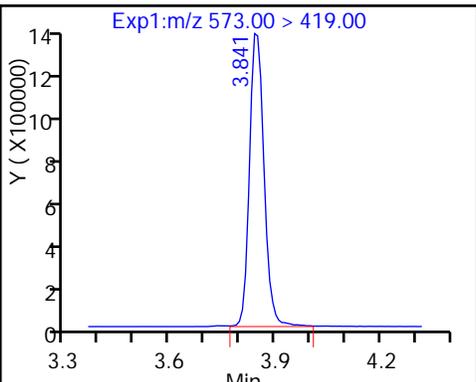
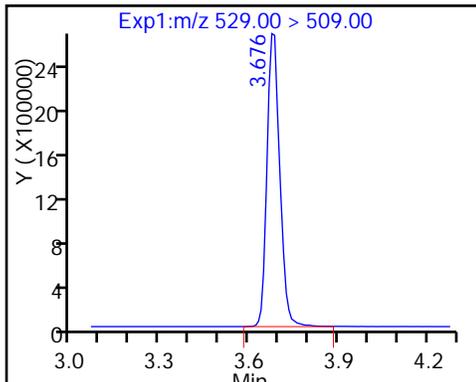
43 Sodium 1H,1H,2H,2H-perfluorooctane



D 42 M2-8:2FTS

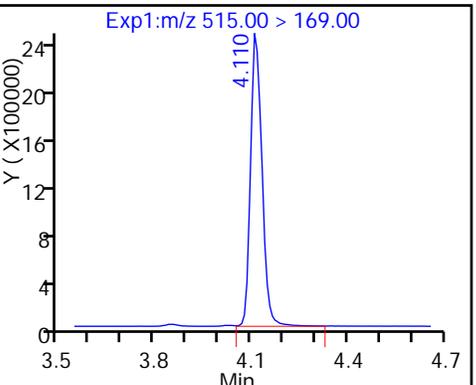
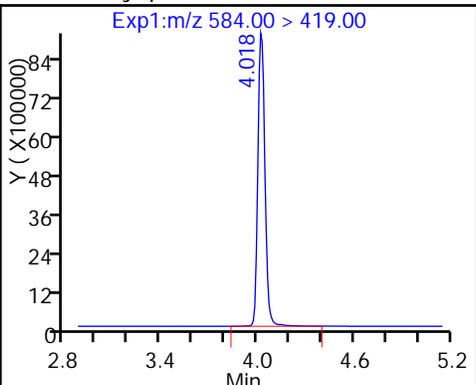
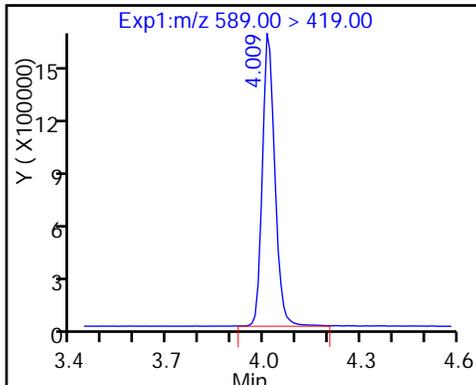
D 45 d3-NMeFOSAA

44 N-methyl perfluorooctane sulfonami



D 46 d5-NEtFOSAA

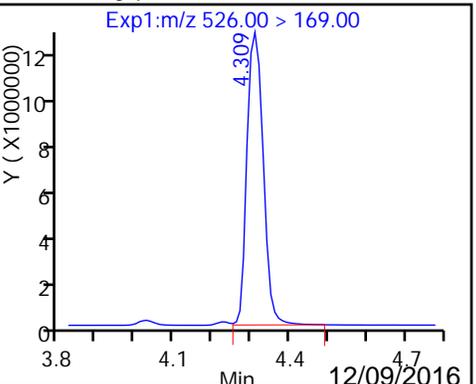
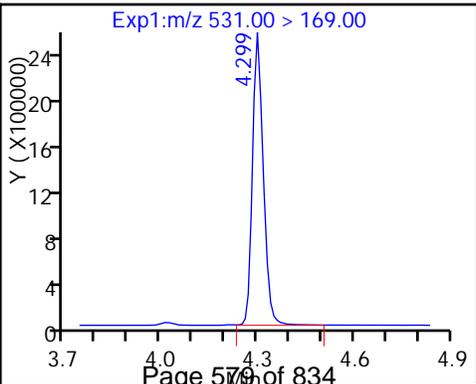
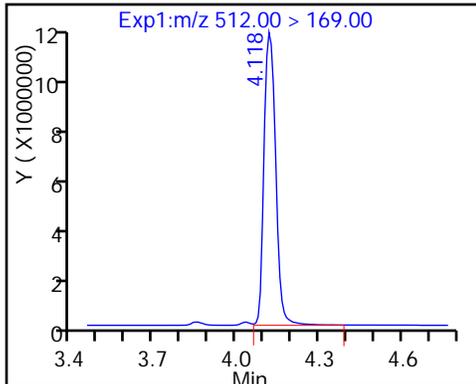
49 N-ethyl perfluorooctane sulfonamid D 52 d-N-MeFOSA-M



54 MeFOSA

D 51 d-N-EtFOSA-M

53 N-ethylperfluoro-1-octanesulfonami



FORM VI
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
 RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140564/4	03DEC2016A_004.d
Level 2	IC 320-140564/13	03DEC2016A_013.d
Level 3	IC 320-140564/5	03DEC2016A_005.d
Level 4	IC 320-140564/14	03DEC2016A_014.d
Level 5	IC 320-140564/6	03DEC2016A_006.d
Level 6	IC 320-140564/15	03DEC2016A_015.d
Level 7	IC 320-140564/7	03DEC2016A_007.d
Level 8	IC 320-140564/16	03DEC2016A_016.d
Level 9	IC 320-140564/8	03DEC2016A_008.d
Level 10	IC 320-140564/17	03DEC2016A_017.d
Level 11	IC 320-140564/9	03DEC2016A_009.d
Level 12	IC 320-140564/18	03DEC2016A_018.d

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12										
Perfluorobutanoic acid (PFBA)	1.582 1.574		1.574		1.574		1.582		1.574		1.327 - 1.827	1.577
Perfluoropentanoic acid (PFPeA)	1.868 1.858		1.858		1.858		1.868		1.858		1.611 - 2.111	1.861
Perfluorobutanesulfonic acid (PFBS)	1.906 1.896		1.897		1.897		1.907		1.897		1.720 - 2.080	1.900
Perfluorohexanoic acid (PFHxA)	2.170 2.157		2.164		2.163		2.171		2.160		1.914 - 2.414	2.164
Perfluoroheptanoic acid (PFHpA)	2.521 2.501		2.512		2.516		2.512		2.510		2.262 - 2.762	2.512
Perfluorohexanesulfonic acid (PFHxS)	++++ 2.526		2.535		2.531		2.527		2.525		2.281 - 2.781	2.529
6:2FTS		++++ 2.838		2.830		2.838		2.840		2.830	2.586 - 3.086	2.835
Perfluorooctanoic acid (PFOA)	++++ 2.868		2.894		2.880		2.894		2.884		2.637 - 3.137	2.884
Perfluoroheptanesulfonic Acid (PFHpS)	2.902 2.884		2.886		2.888		2.886		2.884		2.638 - 3.138	2.888
Perfluorooctane Sulfonate (PFOS)	++++ 3.249		3.259		3.255		3.259		3.257		3.008 - 3.508	3.256
Perfluorononanoic acid (PFNA)	3.275 3.249		3.268		3.263		3.267		3.257		3.013 - 3.513	3.263
Perfluorooctane Sulfonamide (FOSA)	3.581 3.571		3.573		3.576		3.573		3.570		3.324 - 3.824	3.574
8:2FTS		3.601 3.589		3.579		3.588		3.592		3.579	3.338 - 3.838	3.588
Perfluorodecanoic acid (PFDA)	3.632 3.613		3.632		3.618		3.623		3.620		3.373 - 3.873	3.623

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12										
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)		3.774 3.752		3.760		3.760		3.756		3.750	3.509 - 4.009	3.759
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)		3.931 3.927		3.935		3.926		3.931		3.925	3.679 - 4.179	3.929
Perfluorodecanesulfonic acid (PFDS)	3.945 3.921		3.944		3.938		3.939		3.928		3.686 - 4.186	3.936
Perfluoroundecanoic acid (PFUnA)	3.971 3.947		3.953		3.956		3.956		3.946		3.705 - 4.205	3.955
MeFOSA		4.063 4.061		4.051		4.051		4.063		4.051	3.807 - 4.307	4.057
N-EtFOSA-M		4.253 4.249		4.239		4.240		4.244		4.245	3.995 - 4.495	4.245
Perfluorododecanoic acid (PFDoA)	4.261 4.236		4.254		4.253		4.253		4.242		4.000 - 4.500	4.250
Perfluorotridecanoic Acid (PFTriA)	4.535 4.501		4.515		4.521		4.524		4.510		4.268 - 4.768	4.518
Perfluorotetradecanoic acid (PFTeA)	4.778 4.746		4.770		4.761		4.764		4.750		4.511 - 5.011	4.762
Perfluoro-n-hexadecanoic acid (PFHxDA)	5.208 5.160		5.196		5.191		5.187		5.177		4.936 - 5.436	5.187
Perfluoro-n-octadecanoic acid (PFODA)	5.584 5.536		5.565		5.561		5.563		5.547		5.309 - 5.809	5.559
13C4 PFBA	1.582 1.565		1.574		1.574		1.574		1.574		1.324 - 1.824	1.574
13C5-PFPeA	1.868 1.858		1.858		1.858		1.868		1.858		1.611 - 2.111	1.861
13C2 PFHxA	2.179 2.157		2.164		2.163		2.162		2.160		1.914 - 2.414	2.164
13C4-PFHpA	2.521 2.501		2.512		2.508		2.512		2.510		2.261 - 2.761	2.511
18O2 PFHxS	2.544 2.518		2.535		2.531		2.535		2.525		2.281 - 2.781	2.531
M2-6:2FTS		2.840 2.838		2.830		2.838		2.840		2.830	2.586 - 3.086	2.836
13C4 PFOA	2.886 2.868		2.886		2.880		2.886		2.876		2.630 - 3.130	2.880
13C4 PFOS	3.267 3.249		3.259		3.263		3.259		3.257		3.009 - 3.509	3.259
13C5 PFNA	3.275 3.249		3.268		3.263		3.267		3.257		3.013 - 3.513	3.263
13C8 FOSA	3.581 3.562		3.565		3.576		3.573		3.570		3.321 - 3.821	3.571
M2-8:2FTS		3.601 3.589		3.579		3.596		3.592		3.587	3.341 - 3.841	3.591

FORM VI
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
 RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	RT WINDOW	AVG RT
	LVL 11	LVL 12										
13C2 PFDA	3.640 3.621		3.624		3.627		3.623		3.620		3.376 - 3.876	3.626
d3-NMeFOSAA		3.765 3.752		3.760		3.751		3.756		3.750	3.506 - 4.006	3.756
d5-NEtFOSAA		3.931 3.919		3.918		3.918		3.931		3.917	3.672 - 4.172	3.922
13C2 PFUnA	3.971 3.947		3.961		3.956		3.965		3.946		3.708 - 4.208	3.958
d-N-MeFOSA-M		4.054 4.052		4.051		4.051		4.055		4.042	3.801 - 4.301	4.051
d-N-EtFOSA-M		4.244 4.239		4.231		4.240		4.244		4.236	3.989 - 4.489	4.239
13C2 PFDoA	4.261 4.236		4.262		4.253		4.253		4.242		4.001 - 4.501	4.251
13C2-PFTeDA	4.778 4.738		4.762		4.761		4.764		4.750		4.509 - 5.009	4.759
13C2-PFHxDA	5.208 5.171		5.196		5.180		5.187		5.177		4.936 - 5.436	5.187

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140564/4	03DEC2016A_004.d
Level 2	IC 320-140564/13	03DEC2016A_013.d
Level 3	IC 320-140564/5	03DEC2016A_005.d
Level 4	IC 320-140564/14	03DEC2016A_014.d
Level 5	IC 320-140564/6	03DEC2016A_006.d
Level 6	IC 320-140564/15	03DEC2016A_015.d
Level 7	IC 320-140564/7	03DEC2016A_007.d
Level 8	IC 320-140564/16	03DEC2016A_016.d
Level 9	IC 320-140564/8	03DEC2016A_008.d
Level 10	IC 320-140564/17	03DEC2016A_017.d
Level 11	IC 320-140564/9	03DEC2016A_009.d
Level 12	IC 320-140564/18	03DEC2016A_018.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4		B	M1	M2								
	LVL 5	LVL 6	LVL 7	LVL 8												
	LVL 9	LVL 10	LVL 11	LVL 12												
13C4 PFBA	310185 395986 320580		331347 365089 291784		Ave		335828.500			11.4		50.0				
13C5-PFPeA	254875 312687 249553		258388 288538 223232		Ave		264545.467			11.9		50.0				
13C2 PFHxA	230194 279196 219710		235263 254328 206225		Ave		237485.877			10.9		50.0				
13C4-PFHpA	200488 249965 193077		209795 229496 161658		Ave		207413.233			14.7		50.0				
18O2 PFHxS	293993 367326 302399		309858 340372 260105		Ave		312342.301			12.0		50.0				
M2-6:2FTS		91389 101069 97657		86632 132540 123901	Ave		105531.519			17.5		50.0				
13C4 PFOA	219399 265152 198189		227804 241383 163620		Ave		219257.923			16.1		50.0				

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4		B	M1	M2								
	LVL 5	LVL 6	LVL 7	LVL 8												
LVL 9	LVL 10	LVL 11	LVL 12													
13C4 PFOS	228339		232908		Ave		246009.066			12.7		50.0				
	292548		275934													
	236006		210320													
13C5 PFNA	162836		167249		Ave		166415.157			14.2		50.0				
	202161		178407													
	157087		130751													
13C8 FOSA	382785		396984		Ave		402279.103			11.3		50.0				
	462525		446533													
	385863		338985													
M2-8:2FTS		89009		81776	Ave		97142.4774			15.9		50.0				
		92918		112285												
		86527		120341												
13C2 PFDA	153345		157229		Ave		157817.020			12.7		50.0				
	184278		175727													
	148117		128207													
d3-NMeFOSAA		66435		67366	Ave		72529.4733			12.4		50.0				
		73365		86475												
		62459		79076												
d5-NEtFOSAA		76443		76375	Ave		79657.5967			10.3		50.0				
		81901		93004												
		68389		81834												
13C2 PFUnA	116625		121719		Ave		118762.217			14.6		50.0				
	141738		129709													
	112505		90276													
d-N-MeFOSA-M		93925		97701	Ave		105101.257			10.8		50.0				
		110531		123254												
		95888		109309												
d-N-EtFOSA-M		87118		94000	Ave		99165.2500			9.9		50.0				
		101086		114956												
		93356		104475												
13C2 PFDoA	107060		108978		Ave		112083.837			11.2		50.0				
	131990		121360													
	106770		96346													
13C2-PFTEdA	220709		228021		Ave		231173.483			12.4		50.0				
	269763		257560													
	220622		190365													
13C2-PFHxDA	122044		129673		Ave		129725.210			12.9		50.0				
	151839		147280													
	117867		109649													

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_NGC Column: Acquity ID: 2.1 (mm)Heated Purge: (Y/N) NCalibration Start Date: 12/03/2016 13:48Calibration End Date: 12/03/2016 15:33Calibration ID: 26875

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R ² OR COD	#	MIN R ² OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6 LVL 11	LVL 7 LVL 12	LVL 8	LVL 9	LVL 10												
Perfluorobutanoic acid (PFBA)	285568 214032	350423	286663	280973	351800	AveID	0.8740				8.8		35.0				
Perfluoropentanoic acid (PFPeA)	300210 177640	305783	267561	251941	316000	AveID	1.0148				12.2		35.0				
Perfluorobutanesulfonic acid (PFBS)	504086 300541	615867	490253	486734	626278	AveID	1.5960				14.4		50.0				
Perfluorohexanoic acid (PFHxA)	235776 165590	261205	221801	208409	271706	AveID	0.9531				8.6		35.0				
Perfluoroheptanoic acid (PFHpA)	221302 150068	245911	215700	198314	250864	AveID	1.0271				5.9		35.0				
Perfluorohexanesulfonic acid (PFHxS)	++++ 259457	382897	373893	318646	405919	AveID	1.0976				7.2		35.0				
6:2FTS	74414	++++ 95077	114411	73191	96558	AveID	0.8401				11.7		35.0				
Perfluorooctanoic acid (PFOA)	++++ 154270	264619	266474	208682	291075	AveID	1.0719				7.8		35.0				
Perfluoroheptanesulfonic Acid (PFHpS)	288964 206923	350751	264304	290052	344790	AveID	1.1771				9.2		50.0				
Perfluorooctane Sulfonate (PFOS)	++++ 232098	303437	248775	257626	311027	AveID	1.0852				1.7		35.0				
Perfluorononanoic acid (PFNA)	173474 124497	190216	158219	153501	196312	AveID	0.9963				5.5		35.0				
Perfluorooctane Sulfonamide (FOSA)	370100 253291	455589	375193	365380	452435	AveID	0.9341				10.2		35.0				

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_NGC Column: Acquity ID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/03/2016 13:48Calibration End Date: 12/03/2016 15:33Calibration ID: 26875

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10												
8:2FTS	68103	78395 87944	105918	66735	89333	AveID		0.8560			14.0		35.0				
Perfluorodecanoic acid (PFDA)	156640 117857	171719	147102	143186	173721	AveID		0.9605			3.8		35.0				
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)	52006	47238 67834	82772	53845	69682	AveID		0.8583			18.3		35.0				
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)	53051	50866 65238	81472	49948	65231	AveID		0.7657			17.0		35.0				
Perfluorodecanesulfonic acid (PFDS)	142463 129604	184058	155987	152444	180118	AveID		0.6398			3.9		50.0				
Perfluoroundecanoic acid (PFUnA)	154250 89091	133679	130643	111061	140855	AveID		1.0657			12.2		35.0				
MeFOSA	71380	74884 88180	109492	72155	93127	AveID		0.8080			14.0		35.0				
N-EtFOSA-M	73063	72158 91173	110071	72107	94707	AveID		0.8605			12.9		35.0				
Perfluorododecanoic acid (PFDoA)	106404 88867	120759	100943	100016	121390	AveID		0.9490			3.8		35.0				
Perfluorotridecanoic Acid (PFTriA)	106898 87470	118653	100348	100682	125535	AveID		0.9498			3.6		50.0				
Perfluorotetradecanoic acid (PFTeA)	219098 152968	234056	215694	192030	235101	AveID		1.8536			8.9		50.0				
Perfluoro-n-hexadecanoic acid (PFHxDA)	254818 94562	143517	184956	107642	153296	L1ID	0.7490	0.9930						0.9980		0.9900	
Perfluoro-n-octadecanoic acid (PFODA)	100080 89220	110258	119152	114581	134794	AveID		0.9929			8.1		50.0				

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140564/4	03DEC2016A_004.d
Level 2	IC 320-140564/13	03DEC2016A_013.d
Level 3	IC 320-140564/5	03DEC2016A_005.d
Level 4	IC 320-140564/14	03DEC2016A_014.d
Level 5	IC 320-140564/6	03DEC2016A_006.d
Level 6	IC 320-140564/15	03DEC2016A_015.d
Level 7	IC 320-140564/7	03DEC2016A_007.d
Level 8	IC 320-140564/16	03DEC2016A_016.d
Level 9	IC 320-140564/8	03DEC2016A_008.d
Level 10	IC 320-140564/17	03DEC2016A_017.d
Level 11	IC 320-140564/9	03DEC2016A_009.d
Level 12	IC 320-140564/18	03DEC2016A_018.d

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
		LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
13C4 PFBA	Ave	15509259	18254473	16567343	16029004	19799279	50.0	50.0	50.0	50.0	50.0
		14589192					50.0				
13C5-PFPeA	Ave	12743752	14426922	12919402	12477636	15634346	50.0	50.0	50.0	50.0	50.0
		11161582					50.0				
13C2 PFHxA	Ave	11509685	12716423	11763139	10985481	13959775	50.0	50.0	50.0	50.0	50.0
		10311260					50.0				
13C4-PFHpA	Ave	10024418	11474824	10489748	9653835	12498264	50.0	50.0	50.0	50.0	50.0
		8082881					50.0				
1802 PFHxS	Ave	13905891	16099619	14656306	14303487	17374498	47.3	47.3	47.3	47.3	47.3
		12302944					47.3				
M2-6:2FTS	Ave	4800798	4340981	6295656	4115043	4638709	47.5	47.5	47.5	47.5	47.5
		5885296					47.5	47.5			
13C4 PFOA	Ave	10969960	12069172	11390215	9909434	13257596	50.0	50.0	50.0	50.0	50.0
		8181000					50.0				
13C4 PFOS	Ave	10914613	13189624	11132983	11281105	13983781	47.8	47.8	47.8	47.8	47.8
		10053294					47.8				

FORM VI
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1 LVL 6 LVL 11	LVL 2 LVL 7 LVL 12	LVL 3 LVL 8	LVL 4 LVL 9	LVL 5 LVL 10	LVL 1 LVL 6 LVL 11	LVL 2 LVL 7 LVL 12	LVL 3 LVL 8	LVL 4 LVL 9	LVL 5 LVL 10
13C5 PFNA	Ave	8141793 6537549	8920352	8362456	7854327	10108070	50.0 50.0	50.0	50.0	50.0	50.0
13C8 FOSA	Ave	19139237 16949269	22326638	19849207	19293148	23126232	50.0 50.0	50.0	50.0	50.0	50.0
M2-8:2FTS	Ave	4450751	4263521 5764339	5378444	3917070	4144623	47.9	47.9 47.9	47.9	47.9	47.9
13C2 PFDA	Ave	7667226 6410340	8786342	7861453	7405870	9213875	50.0 50.0	50.0	50.0	50.0	50.0
d3-NMeFOSAA	Ave	3668263	3321758 3953776	4323749	3368323	3122973	50.0	50.0 50.0	50.0	50.0	50.0
d5-NEtFOSAA	Ave	4095032	3822164 4091711	4650188	3818745	3419439	50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFUnA	Ave	5831263 4513820	6485474	6085968	5625237	7086903	50.0 50.0	50.0	50.0	50.0	50.0
d-N-MeFOSA-M	Ave	5526561	4696243 5465444	6162693	4885054	4794382	50.0	50.0 50.0	50.0	50.0	50.0
d-N-EtFOSA-M	Ave	5054297	4355916 5223764	5747792	4699997	4667809	50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFDoA	Ave	5353024 4817286	6067987	5448882	5338479	6599493	50.0 50.0	50.0	50.0	50.0	50.0
13C2-PFTeDA	Ave	11035471 9518260	12877999	11401035	11031119	13488161	50.0 50.0	50.0	50.0	50.0	50.0
13C2-PFHxDA	Ave	6102203 5482429	7363997	6483658	5893329	7591947	50.0 50.0	50.0	50.0	50.0	50.0

Curve Type Legend:

Ave = Average

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_NGC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/03/2016 13:48Calibration End Date: 12/03/2016 15:33Calibration ID: 26875

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	IC 320-140564/4	03DEC2016A_004.d
Level 2	IC 320-140564/13	03DEC2016A_013.d
Level 3	IC 320-140564/5	03DEC2016A_005.d
Level 4	IC 320-140564/14	03DEC2016A_014.d
Level 5	IC 320-140564/6	03DEC2016A_006.d
Level 6	IC 320-140564/15	03DEC2016A_015.d
Level 7	IC 320-140564/7	03DEC2016A_007.d
Level 8	IC 320-140564/16	03DEC2016A_016.d
Level 9	IC 320-140564/8	03DEC2016A_008.d
Level 10	IC 320-140564/17	03DEC2016A_017.d
Level 11	IC 320-140564/9	03DEC2016A_009.d
Level 12	IC 320-140564/18	03DEC2016A_018.d

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
Perfluorobutanoic acid (PFBA)		AveID	142784	7008467	286663	14048645	1758998	0.500	20.0	1.00	50.0	5.00
			42806432					200				
Perfluoropentanoic acid (PFPeA)		AveID	150105	6115663	267561	12597041	1580002	0.500	20.0	1.00	50.0	5.00
			35527936					200				
Perfluorobutanesulfonic acid (PFBS)		AveID	222806	10888534	433384	21513660	2768149	0.442	17.7	0.884	44.2	4.42
			53135654					177				
Perfluorohexanoic acid (PFHxA)		AveID	117888	5224102	221801	10420431	1358530	0.500	20.0	1.00	50.0	5.00
			33117925					200				
Perfluoroheptanoic acid (PFHpA)		AveID	110651	4918226	215700	9915679	1254320	0.500	20.0	1.00	50.0	5.00
			30013620					200				
Perfluorohexanesulfonic acid (PFHxS)		AveID	+++++	6968734	340243	14498414	1846932	+++++	18.2	0.910	45.5	4.55
			47221104					182				
6:2FTS		AveID	352721	+++++	2169230	69385	4576863	4.74	+++++	19.0	0.948	47.4
			18026595	190								
Perfluorooctanoic acid (PFOA)		AveID	+++++	5292388	266474	10434085	1455373	+++++	20.0	1.00	50.0	5.00
			30853962					200				

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica SacramentoJob No.: 320-23783-1Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_NGC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 12/03/2016 13:48Calibration End Date: 12/03/2016 15:33Calibration ID: 26875

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
Perfluoroheptanesulfonic Acid (PFHpS)		AveID	137547 39398056	6678298	251617	13806464	1641199	0.476 190	19.0	0.952	47.6	4.76
Perfluorooctane Sulfonate (PFOS)		AveID	++++ 43077470	5631788	230863	11953842	1443163	++++ 186	18.6	0.928	46.4	4.64
Perfluorononanoic acid (PFNA)		AveID	86737 24899337	3804327	158219	7675034	981559	0.500 200	20.0	1.00	50.0	5.00
Perfluorooctane Sulfonamide (FOSA)		AveID	185050 50658186	9111782	375193	18269014	2262174	0.500 200	20.0	1.00	50.0	5.00
8:2FTS		AveID	326213	37551 16850029	2029390	63932	4279048	4.79	0.479 192	19.2	0.958	47.9
Perfluorodecanoic acid (PFDA)		AveID	78320 23571490	3434382	147102	7159278	868604	0.500 200	20.0	1.00	50.0	5.00
N-methyl perfluorooctane sulfonamidoacetic acid (NMeFOSAA)		AveID	260030	23619 13566790	1655435	53845	3484088	5.00	0.500 200	20.0	1.00	50.0
N-ethyl perfluorooctane sulfonamidoacetic acid (NEtFOSAA)		AveID	265257	25433 13047697	1629443	49948	3261537	5.00	0.500 200	20.0	1.00	50.0
Perfluorodecanesulfonic acid (PFDS)		AveID	68667 24987612	3548629	150371	7347809	868168	0.482 193	19.3	0.964	48.2	4.82
Perfluoroundecanoic acid (PFUnA)		AveID	77125 17818228	2673576	130643	5553051	704273	0.500 200	20.0	1.00	50.0	5.00
MeFOSA		AveID	356898	37442 17635948	2189848	72155	4656344	5.00	0.500 200	20.0	1.00	50.0
N-EtFOSA-M		AveID	365313	36079 18234509	2201423	72107	4735360	5.00	0.500 200	20.0	1.00	50.0
Perfluorododecanoic acid (PFDoA)		AveID	53202 17773370	2415187	100943	5000795	606951	0.500 200	20.0	1.00	50.0	5.00

RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1 Analy Batch No.: 140564

SDG No.: _____

Instrument ID: A8_N GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) NCalibration Start Date: 12/03/2016 13:48 Calibration End Date: 12/03/2016 15:33 Calibration ID: 26875

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5
			LVL 6	LVL 7	LVL 8	LVL 9	LVL 10	LVL 6	LVL 7	LVL 8	LVL 9	LVL 10
			LVL 11	LVL 12				LVL 11	LVL 12			
Perfluorotridecanoic Acid (PFTriA)		AveID	53449	2373060	100348	5034109	627673	0.500	20.0	1.00	50.0	5.00
			17494025					200				
Perfluorotetradecanoic acid (PFTeA)		AveID	109549	4681126	215694	9601518	1175506	0.500	20.0	1.00	50.0	5.00
			30593527					200				
Perfluoro-n-hexadecanoic acid (PFHxDA)		L1ID	127409	2870345	184956	5382109	766480	0.500	20.0	1.00	50.0	5.00
			18912385					200				
Perfluoro-n-octadecanoic acid (PFODA)		AveID	50040	2205153	119152	5729058	673971	0.500	20.0	1.00	50.0	5.00
			17844066					200				

Curve Type Legend:

AveID = Average isotope dilution
L1ID = Linear 1/conc IsoDil

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_004.d
 Lims ID: IC L1
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 03-Dec-2016 13:48:41 ALS Bottle#: 37 Worklist Smp#: 4
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L1_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:05 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:42:37

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.582	1.574	0.008	15509259	46.2		92.4	1670611	
1 Perfluorobutyric acid	212.90 > 169.00	1.582	1.577	0.005	142784	0.5267		105	902	
3 Perfluoropentanoic acid	262.90 > 219.00	1.868	1.861	0.007	150105	0.5803		116	1311	
D 4 13C5-PFPeA	267.90 > 223.00	1.868	1.861	0.007	12743752	48.2		96.3	785537	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.906	1.900	0.006	222806	0.4748		107		
	298.90 > 99.00	1.906	1.900	0.006	90815		2.45(0.00-0.00)	107		
7 Perfluorohexanoic acid	313.00 > 269.00	2.170	2.164	0.006	117888	0.5373		107	2866	
D 6 13C2 PFHxA	315.00 > 270.00	2.179	2.164	0.015	11509685	48.5		96.9	1035144	
D 11 13C4-PFHpA	367.00 > 322.00	2.521	2.511	0.010	10024418	48.3		96.7	941002	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.521	2.512	0.009	110651	0.5374		107	1512	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.544	2.531	0.013	195583	0.6061		133		
D 10 18O2 PFHxS	403.00 > 84.00	2.544	2.531	0.013	13905891	44.5		94.1	1599849	
D 14 13C4 PFOA	417.00 > 372.00	2.886	2.880	0.006	10969960	50.0		100	702523	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.902	2.887	0.015	1.000	176658	0.7512		150	2006	
413.00 > 169.00	2.902	2.887	0.015	1.000	83124		2.13(0.90-1.10)	150	4894	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.902	2.888	0.014	1.000	137547	0.5117		108		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.267	3.258	0.009	1.000	118203	0.4770		103	15866	
499.00 > 99.00	3.250	3.258	-0.008	0.995	0		0.00(0.90-1.10)	103		
D 17 13C4 PFOS										
503.00 > 80.00	3.267	3.259	0.008		10914613	44.4		92.8	963774	
D 19 13C5 PFNA										
468.00 > 423.00	3.275	3.263	0.012		8141793	48.9		97.8	521938	
20 Perfluorononanoic acid										
463.00 > 419.00	3.275	3.263	0.012	1.000	86737	0.5346		107	1741	M
D 21 13C8 FOSA										
506.00 > 78.00	3.581	3.571	0.010		19139237	47.6		95.2	555083	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.581	3.574	0.007	1.000	185050	0.5175		104	16767	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.632	3.623	0.009	1.000	78320	0.5318		106	2431	
D 23 13C2 PFDA										
515.00 > 470.00	3.640	3.626	0.014		7667226	48.6		97.2	219644	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.945	3.936	0.009	1.000	68667	0.4701		97.5		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.971	3.955	0.016	1.000	77125	0.6205		124	2113	M
D 27 13C2 PFUnA										
565.00 > 520.00	3.971	3.958	0.013		5831263	49.1		98.2	341809	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.261	4.250	0.011	1.000	53202	0.5236		105	752	
D 30 13C2 PFDaA										
615.00 > 570.00	4.261	4.251	0.010		5353024	47.8		95.5	191441	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.535	4.518	0.017	1.000	53449	0.5256		105	91.0	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.778	4.759	0.019		11035471	47.7		95.5	488569	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.778	4.761	0.017	1.000	109549	0.5520		110	60.5	
713.00 > 169.00	4.770	4.761	0.009	0.998	20451		5.36(0.00-0.00)	110	3502	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.208	5.186	0.022		6102203	47.0		94.1	113568	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.208	5.186	0.022	1.000	127409	0.4441		88.8	179	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.584	5.559	0.025	1.000	50040	0.4708		94.2	82.2	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L1_00021

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_004.d

Injection Date: 03-Dec-2016 13:48:41

Instrument ID: A8_N

Lims ID: IC L1

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 37

Worklist Smp#: 4

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

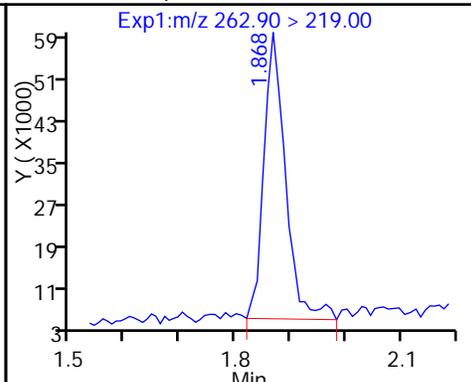
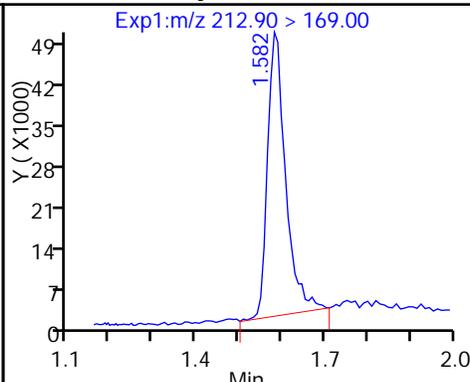
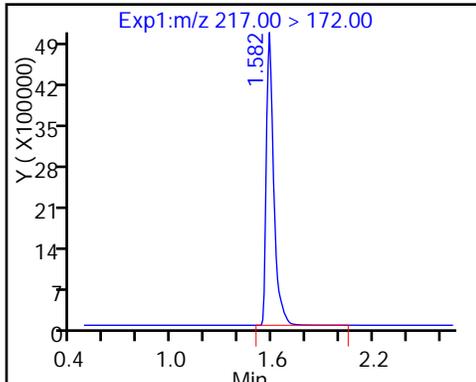
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

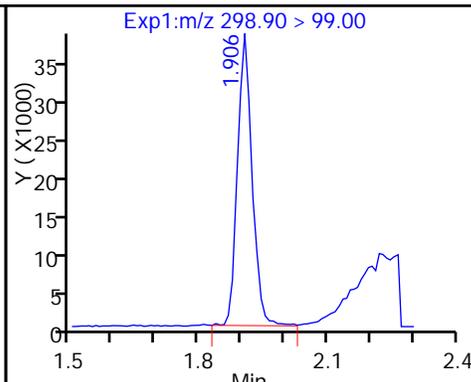
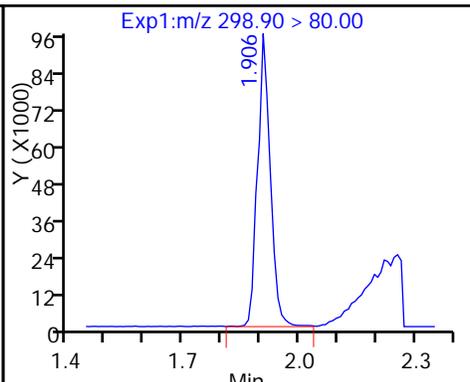
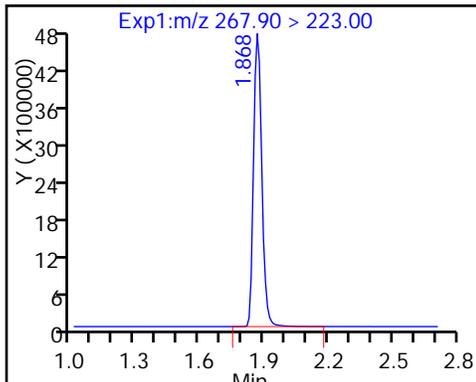
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

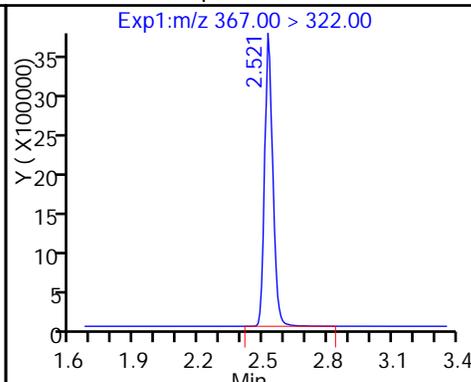
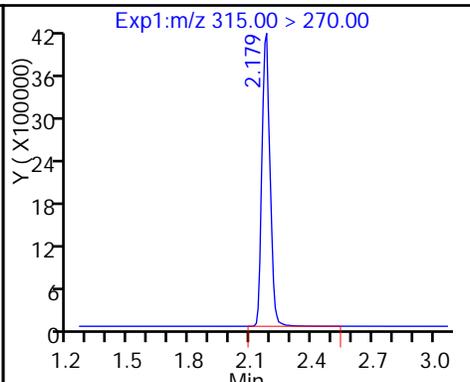
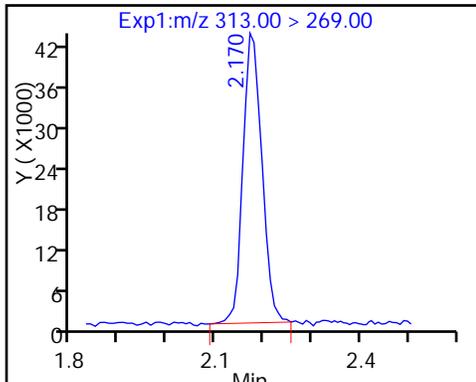
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

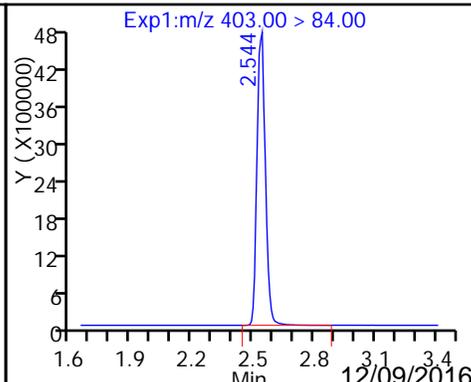
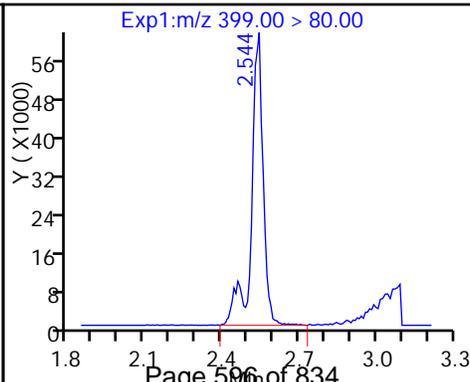
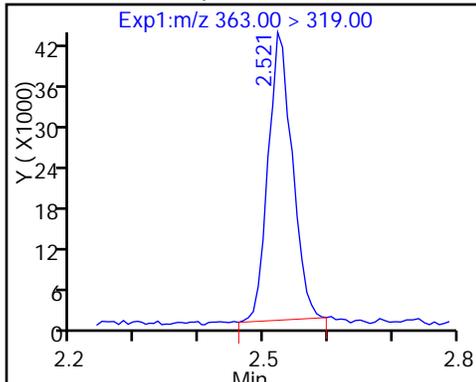
D 11 13C4-PFHpA



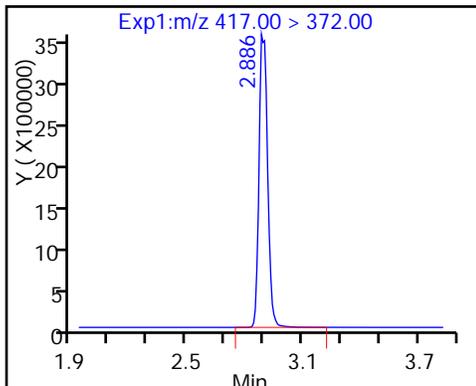
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

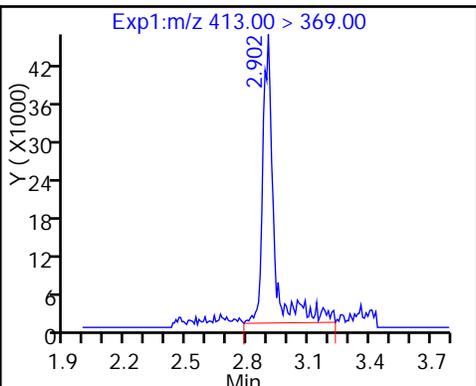
D 10 18O2 PFHxS



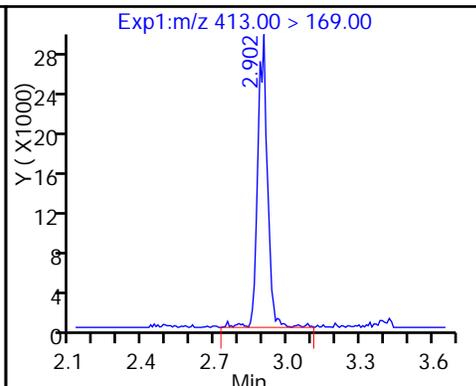
D 14 13C4 PFOA



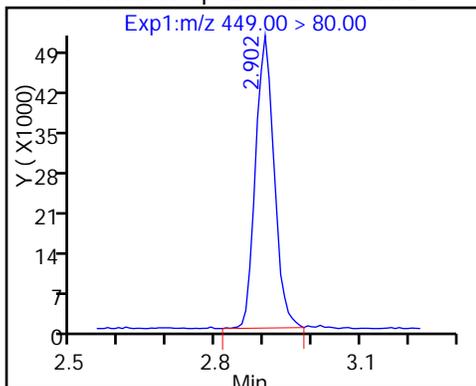
15 Perfluorooctanoic acid



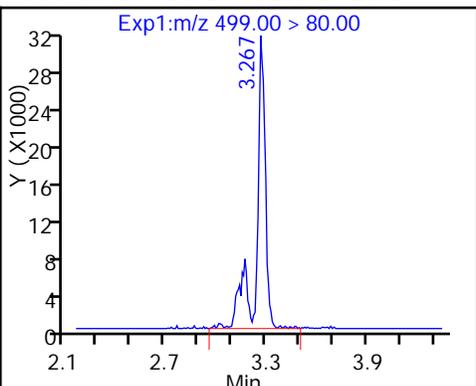
15 Perfluorooctanoic acid



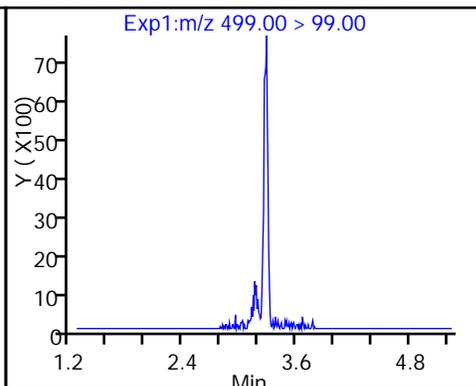
13 Perfluoroheptanesulfonic Acid



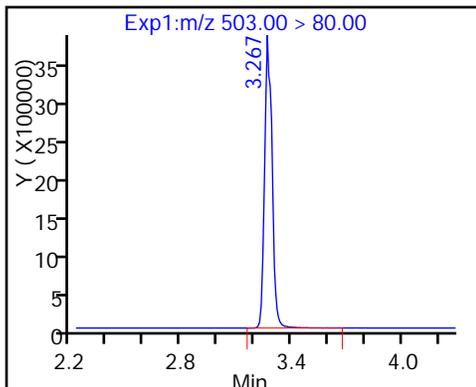
18 Perfluorooctane sulfonic acid



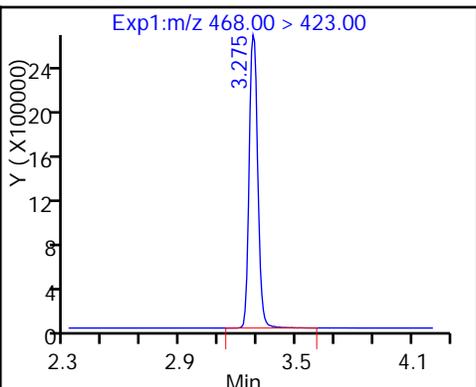
18 Perfluorooctane sulfonic acid



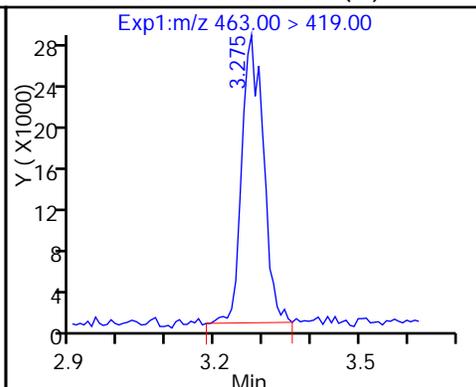
D 17 13C4 PFOS



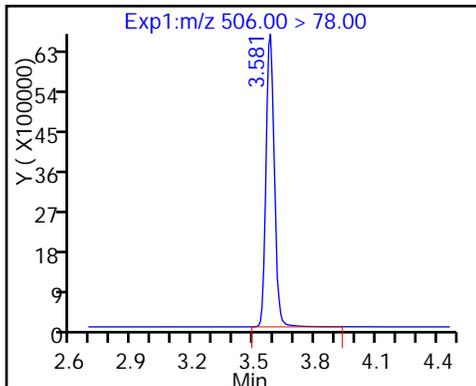
D 19 13C5 PFNA



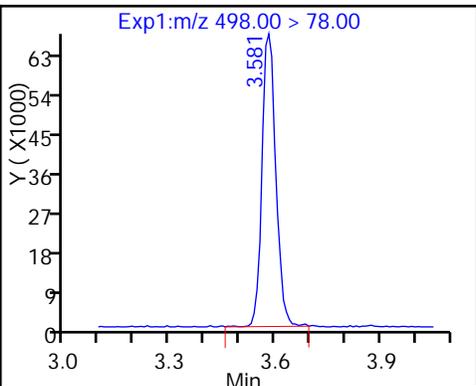
20 Perfluorononanoic acid (M)



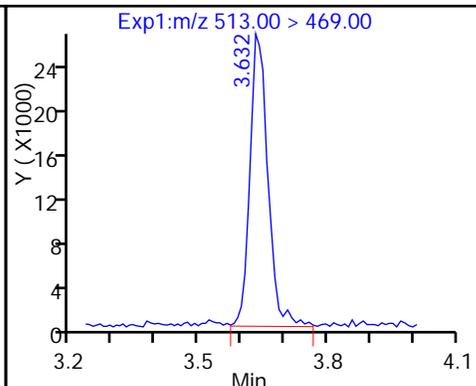
D 21 13C8 FOSA



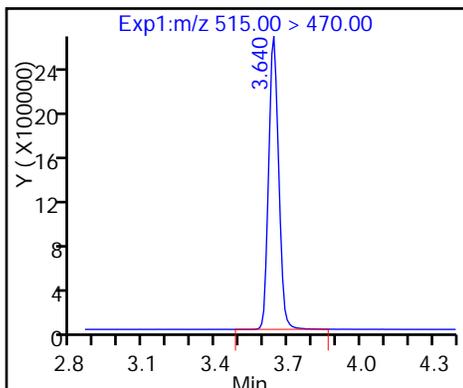
22 Perfluorooctane Sulfonamide



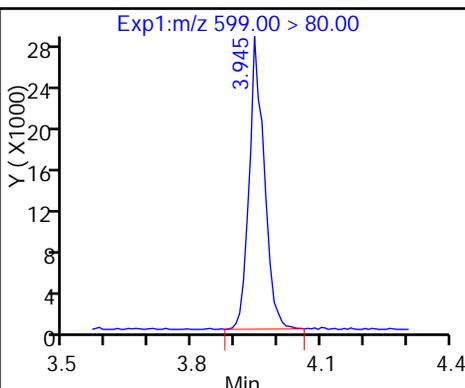
24 Perfluorodecanoic acid



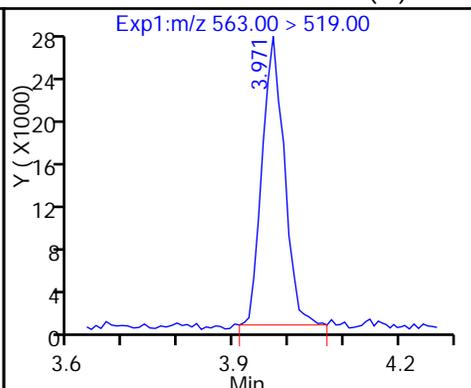
D 23 13C2 PFDA



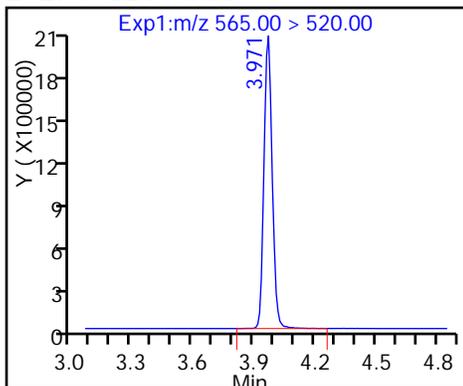
26 Perfluorodecane Sulfonic acid



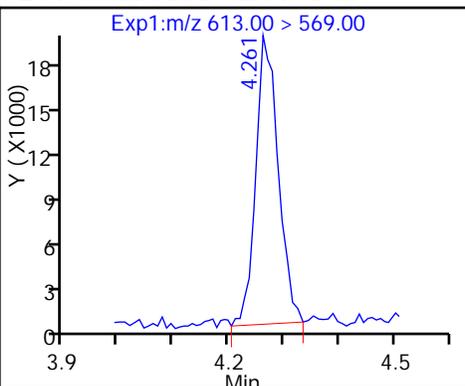
28 Perfluoroundecanoic acid (M)



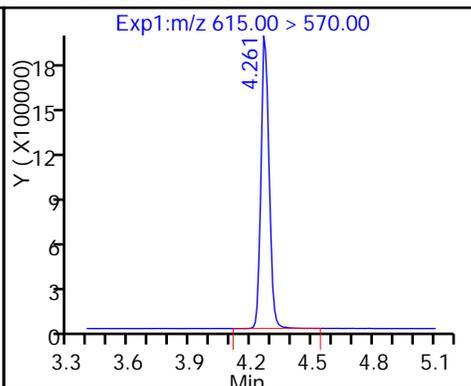
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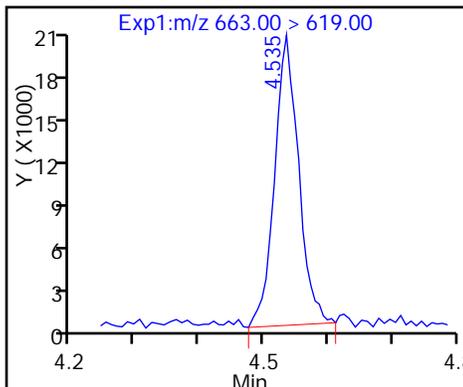
29 Perfluorododecanoic acid



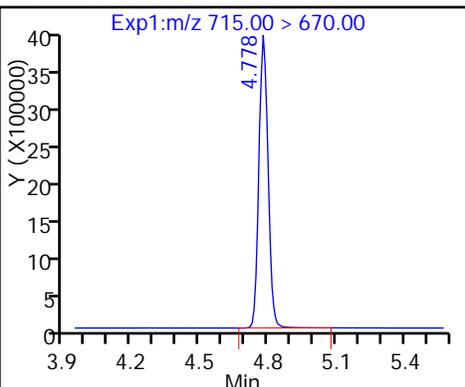
D 30 13C2 PFDa



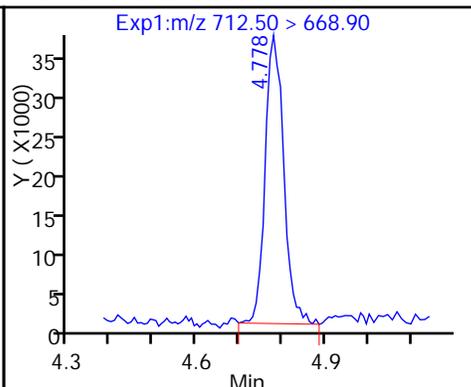
31 Perfluorotridecanoic acid



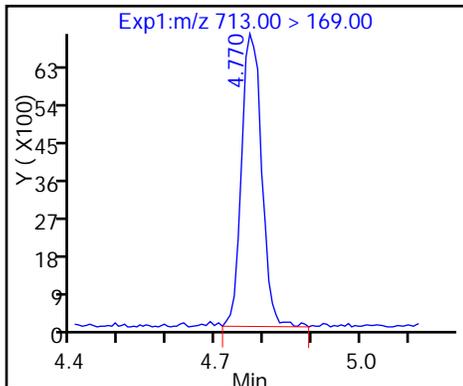
D 32 13C2-PFTeDa



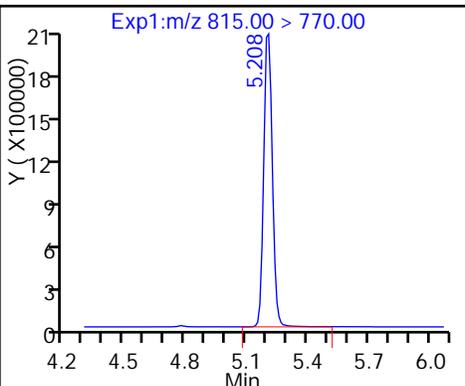
33 Perfluorotetradecanoic acid



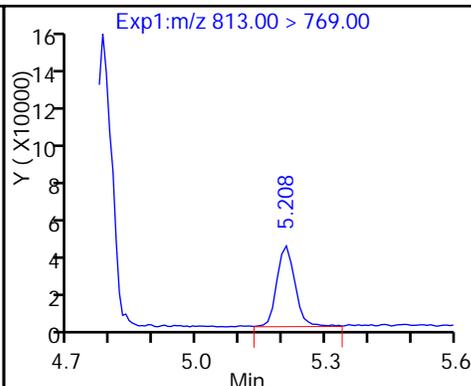
33 Perfluorotetradecanoic acid



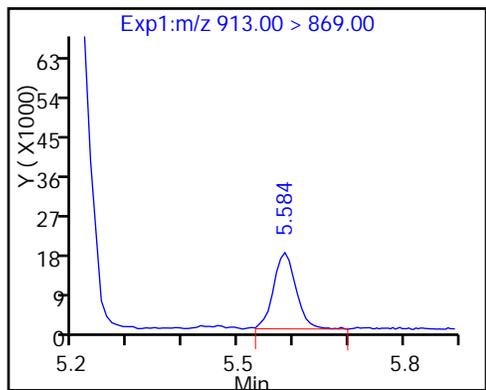
D 34 13C2-PFHxDa



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



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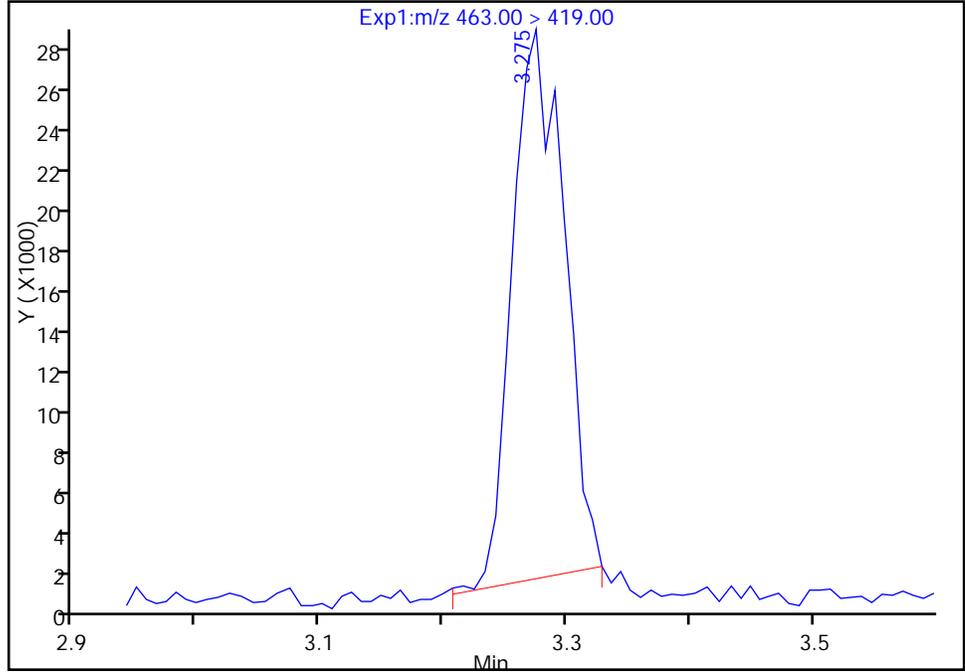
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_004.d
Injection Date: 03-Dec-2016 13:48:41 Instrument ID: A8_N
Lims ID: IC L1
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 4
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

20 Perfluorononanoic acid, CAS: 375-95-1

Signal: 1

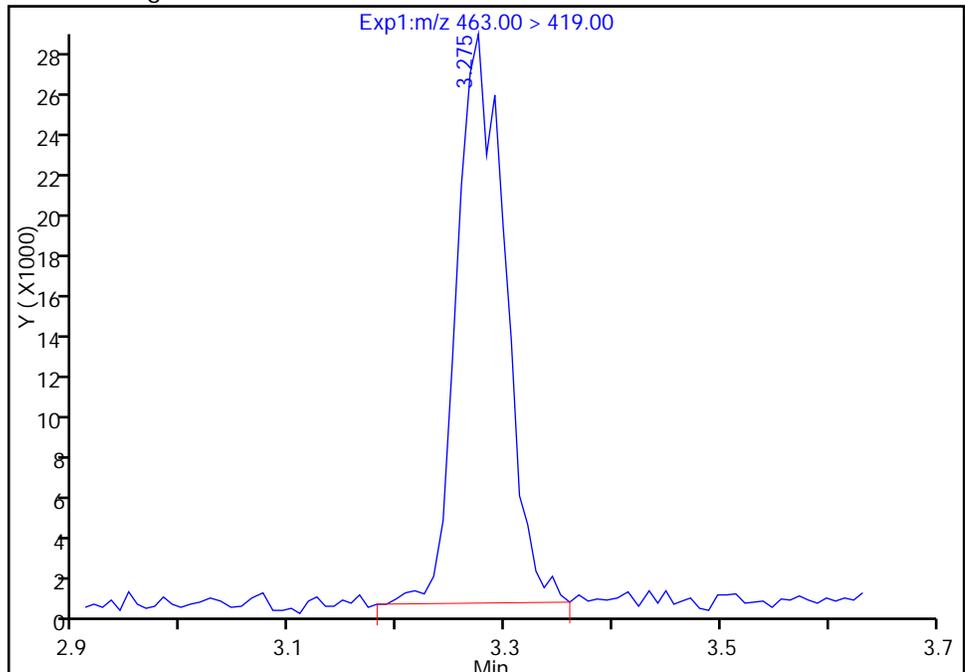
RT: 3.28
Area: 78587
Amount: 0.492646
Amount Units: ng/ml

Processing Integration Results



RT: 3.28
Area: 86737
Amount: 0.534631
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 09:42:37
Audit Action: Manually Integrated

Audit Reason: Incomplete Integration

TestAmerica Sacramento

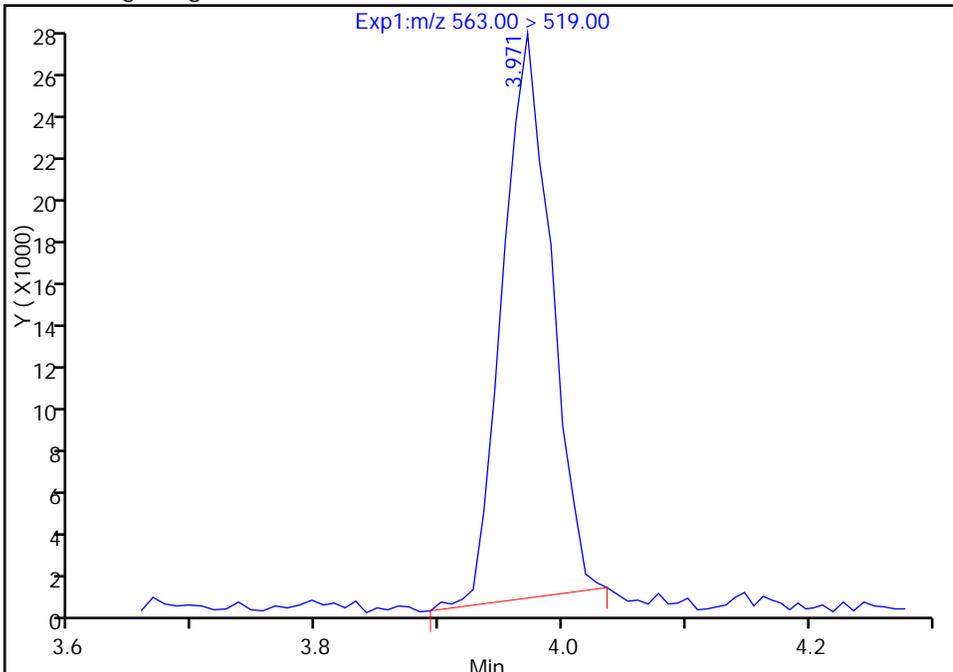
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Injection Date: 03-Dec-2016 13:48:41 Instrument ID: A8_N
Lims ID: IC L1
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 37 Worklist Smp#: 4
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

28 Perfluoroundecanoic acid, CAS: 2058-94-8

Signal: 1

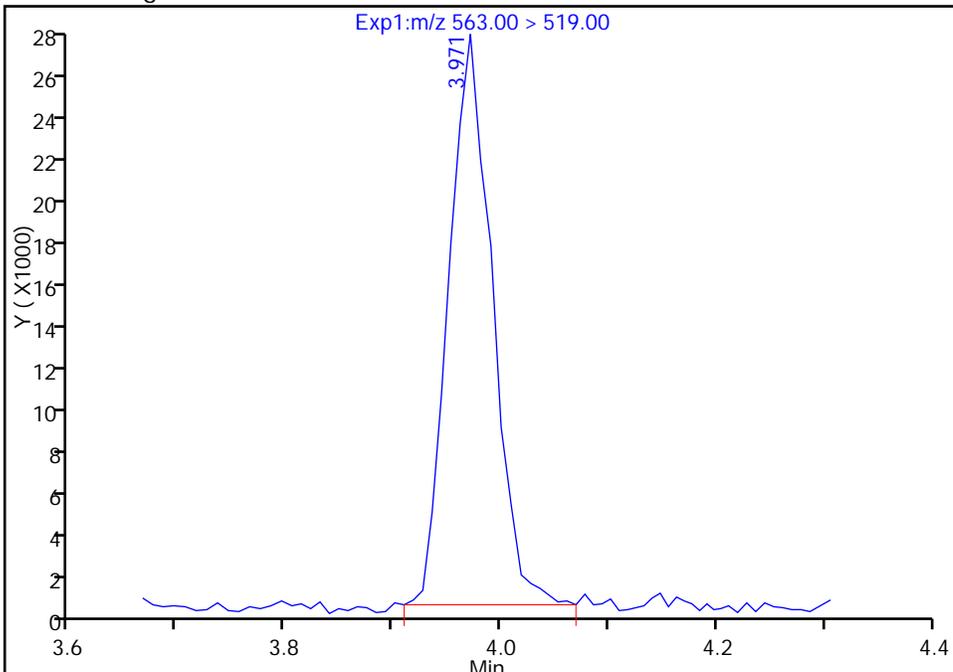
RT: 3.97
Area: 74499
Amount: 0.603647
Amount Units: ng/ml

Processing Integration Results



RT: 3.97
Area: 77125
Amount: 0.620524
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 09:42:37

Audit Action: Manually Integrated

Audit Reason: Incomplete Integration

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_005.d
 Lims ID: IC L2
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 03-Dec-2016 13:56:13 ALS Bottle#: 38 Worklist Smp#: 5
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L2_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:08 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:43:09

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	16567343	49.3		98.7	1253121	
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.577	-0.003	1.000	286663	0.9899	99.0	1783	
3 Perfluoropentanoic acid	262.90 > 219.00	1.858	1.861	-0.003	1.000	267561	1.02	102	2233	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.861	-0.003		12919402	48.8	97.7	738841	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.897	1.900	-0.003	1.000	433384	0.8763	99.1		
	298.90 > 99.00	1.897	1.900	-0.003	1.000	173533	2.50(0.00-0.00)	99.1		
7 Perfluorohexanoic acid	313.00 > 269.00	2.164	2.164	0.0	1.000	221801	0.9891	98.9	4857	
D 6 13C2 PFHxA	315.00 > 270.00	2.164	2.164	0.0		11763139	49.5	99.1	702628	
D 11 13C4-PFHpA	367.00 > 322.00	2.512	2.511	0.001		10489748	50.6	101	757325	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.512	2.512	0.0	1.000	215700	1.00	100	2505	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.535	2.531	0.004	1.000	340243	1.00	110		
D 10 18O2 PFHxS	403.00 > 84.00	2.535	2.531	0.004		14656306	46.9	99.2	569002	
D 14 13C4 PFOA	417.00 > 372.00	2.886	2.880	0.006		11390215	51.9	104	1255788	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.894	2.887	0.007	1.000	266474	1.09		109	2568	
413.00 > 169.00	2.886	2.887	-0.001	0.997	153584		1.74(0.90-1.10)	109	8517	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.886	2.888	-0.002	1.000	251617	0.9178		96.4		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.259	3.258	0.001	1.000	230863	0.9134		98.4	60496	
499.00 > 99.00	3.268	3.258	0.010	1.003	49554		4.66(0.90-1.10)	98.4	7390	
D 17 13C4 PFOS										
503.00 > 80.00	3.259	3.259	0.0		11132983	45.3		94.7	553508	
D 19 13C5 PFNA										
468.00 > 423.00	3.268	3.263	0.005		8362456	50.3		101	686636	
20 Perfluorononanoic acid										
463.00 > 419.00	3.268	3.263	0.005	1.000	158219	0.9495		94.9	2139	
D 21 13C8 FOSA										
506.00 > 78.00	3.565	3.571	-0.006		19849207	49.3		98.7	435087	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.573	3.574	-0.001	1.000	375193	1.01		101	37696	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.632	3.623	0.009	1.000	147102	0.9741		97.4	6264	
D 23 13C2 PFDA										
515.00 > 470.00	3.624	3.626	-0.002		7861453	49.8		99.6	210924	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.944	3.936	0.008	1.000	150371	1.01		105		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.953	3.955	-0.003	1.000	130643	1.01		101	3153	
D 27 13C2 PFUnA										
565.00 > 520.00	3.961	3.958	0.003		6085968	51.2		102	355525	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.254	4.250	0.004	1.000	100943	0.9761		97.6	1508	
D 30 13C2 PFDaA										
615.00 > 570.00	4.262	4.251	0.011		5448882	48.6		97.2	155924	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.515	4.518	-0.003	1.000	100348	0.9695		96.9	172	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.762	4.759	0.003		11401035	49.3		98.6	776058	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.770	4.761	0.009	1.000	215694	1.07		107	105	
713.00 > 169.00	4.762	4.761	0.001	0.998	35526		6.07(0.00-0.00)	107	12035	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.196	5.186	0.010		6483658	50.0		100.0	170445	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.196	5.186	0.010	1.000	184956	0.9549		95.5	215	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.565	5.559	0.006	1.000	119152	1.10		110	244	

Reagents:

LCPFC-L2_00022

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_005.d

Injection Date: 03-Dec-2016 13:56:13

Instrument ID: A8_N

Lims ID: IC L2

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 38

Worklist Smp#: 5

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

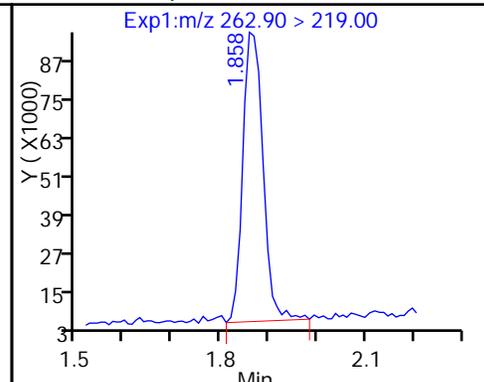
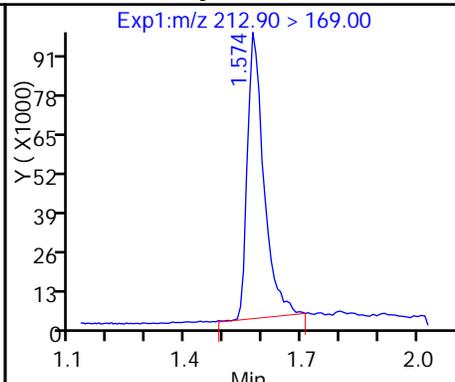
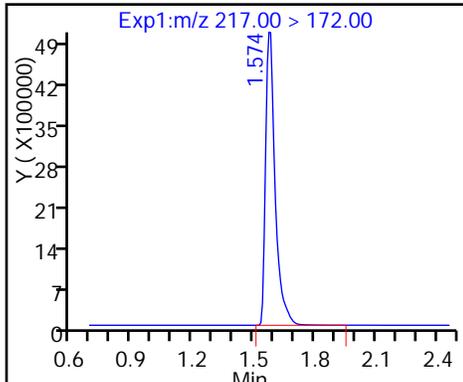
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

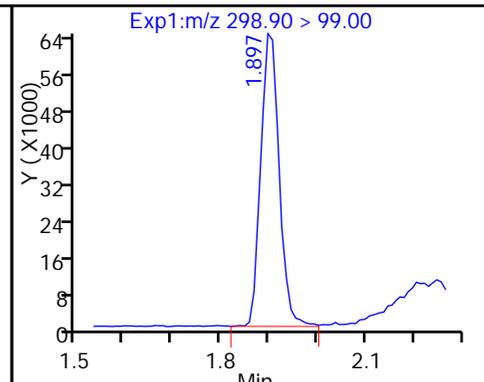
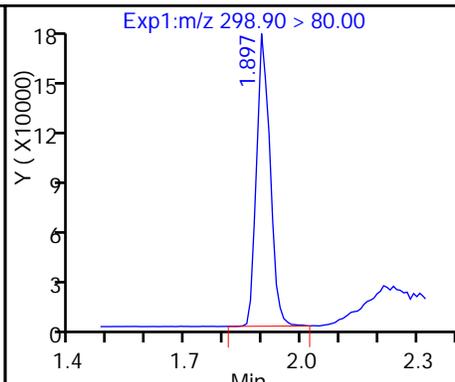
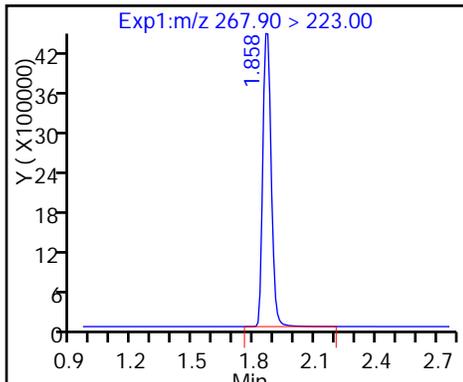
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

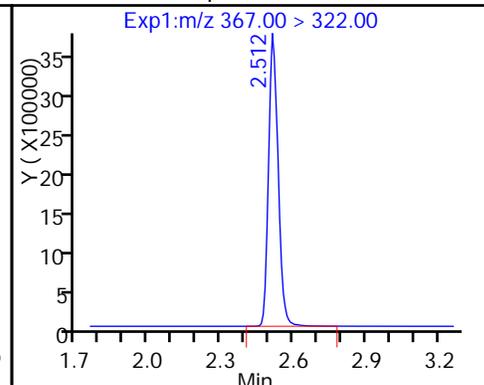
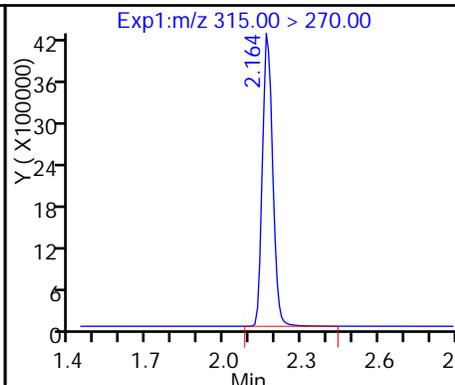
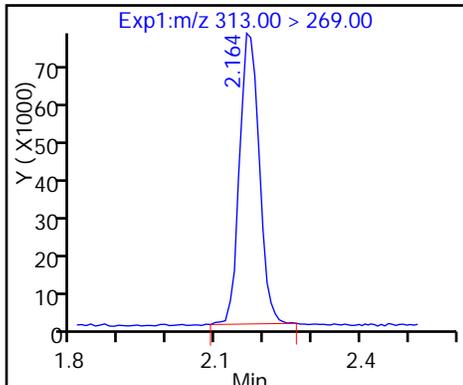
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

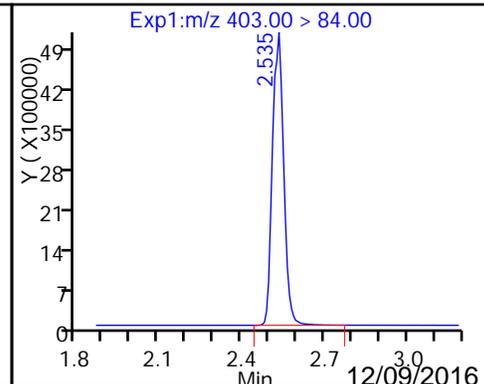
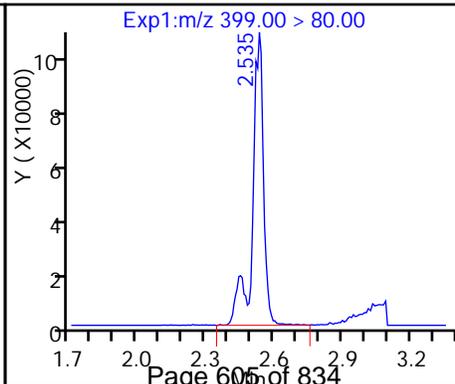
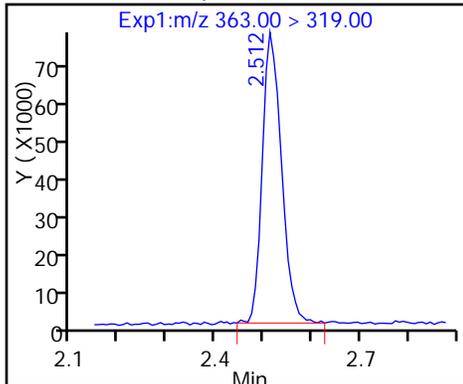
D 11 13C4-PFHpA



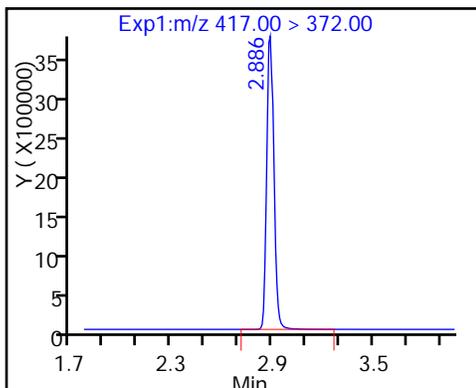
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

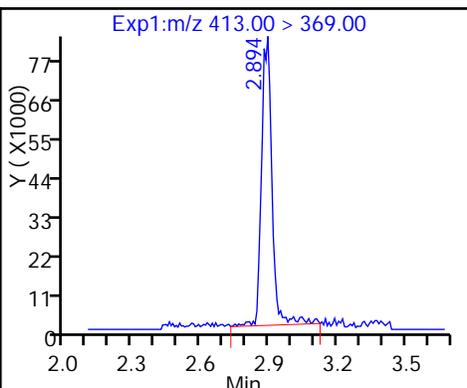
D 10 18O2 PFHxS



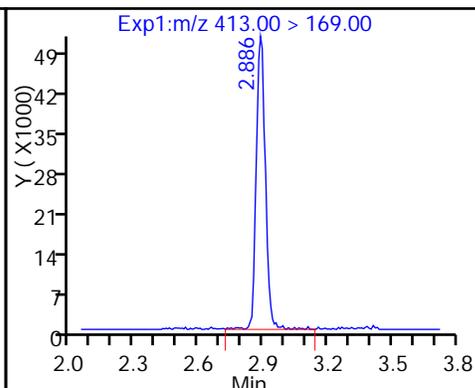
D 14 13C4 PFOA



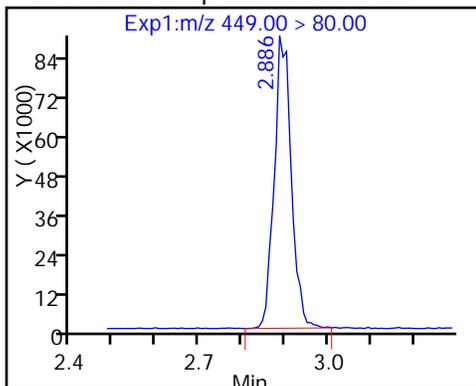
15 Perfluorooctanoic acid



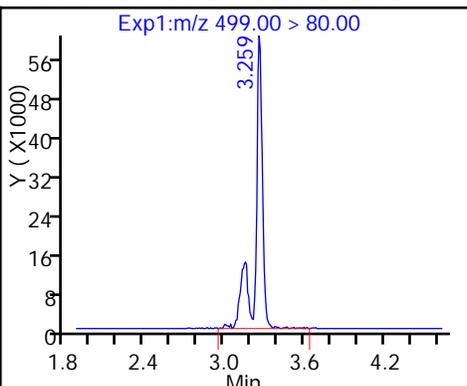
15 Perfluorooctanoic acid



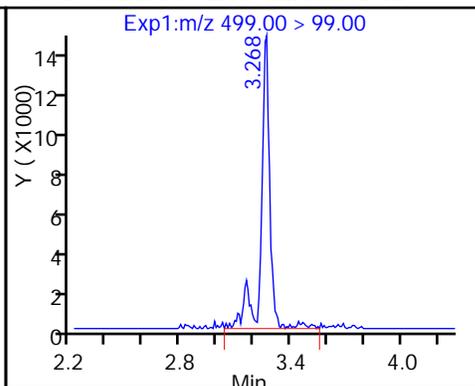
13 Perfluoroheptanesulfonic Acid



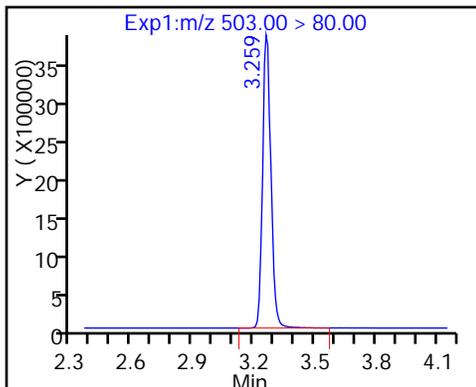
18 Perfluorooctane sulfonic acid



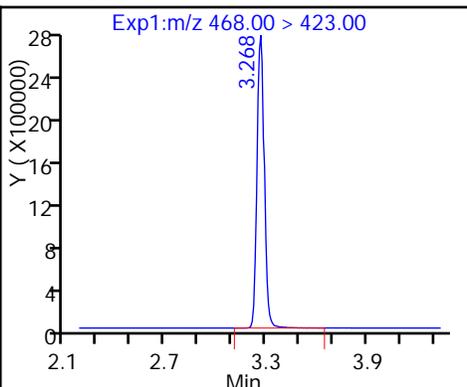
18 Perfluorooctane sulfonic acid



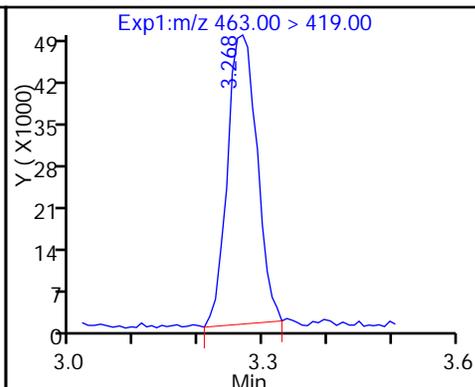
D 17 13C4 PFOS



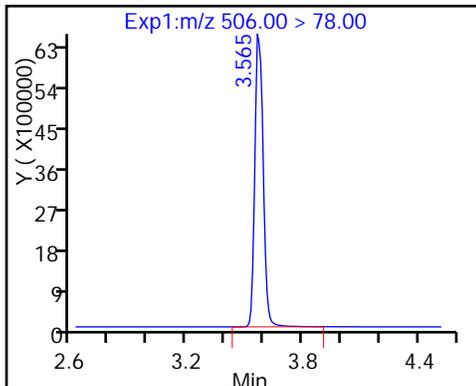
D 19 13C5 PFNA



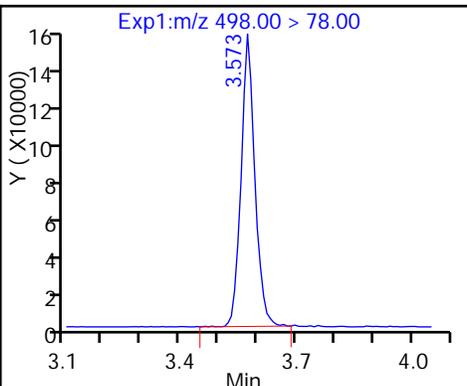
20 Perfluorononanoic acid



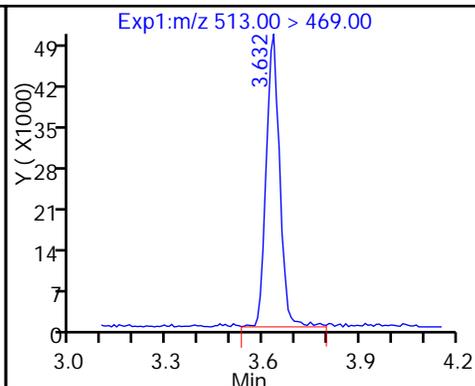
D 21 13C8 FOSA



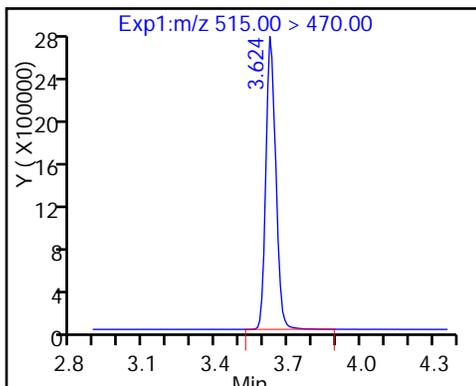
22 Perfluorooctane Sulfonamide



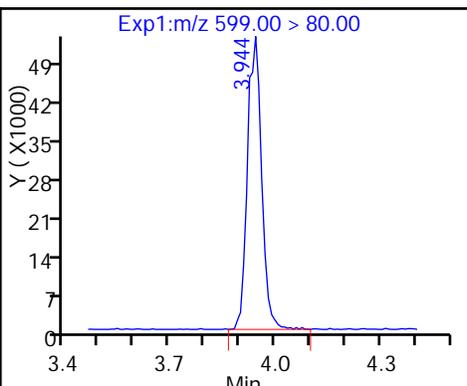
24 Perfluorodecanoic acid



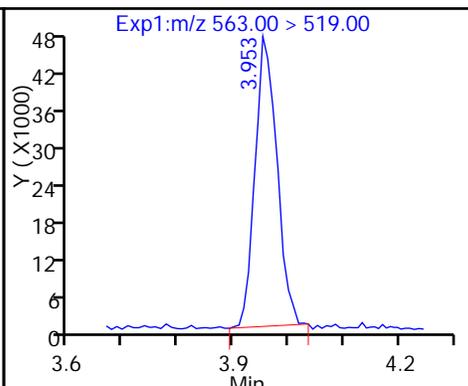
D 23 13C2 PFDA



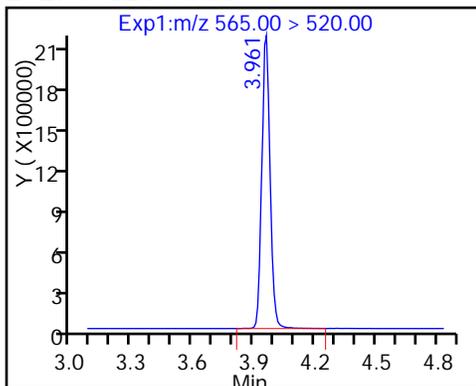
26 Perfluorodecane Sulfonic acid



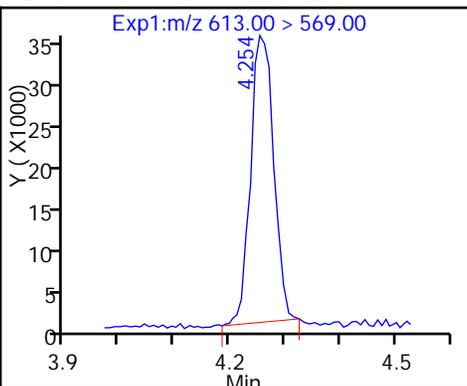
28 Perfluoroundecanoic acid



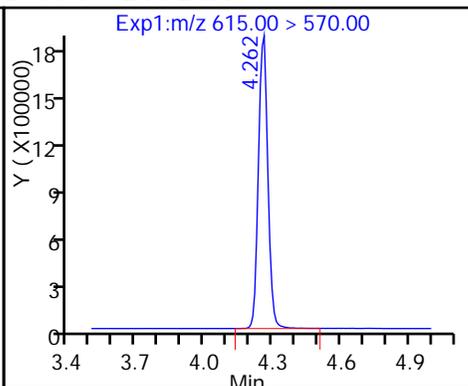
D 27 13C2 PFUa



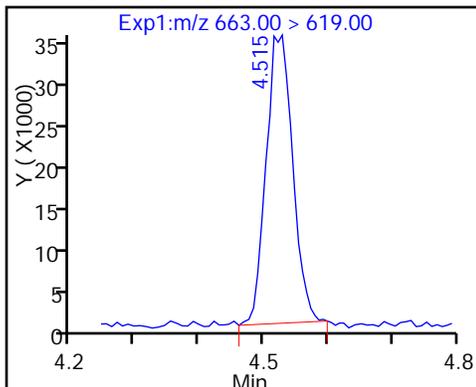
29 Perfluorododecanoic acid



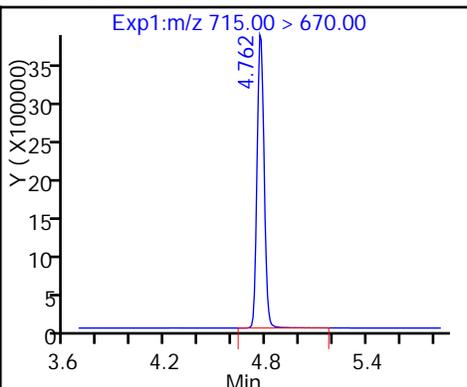
D 30 13C2 PFDa



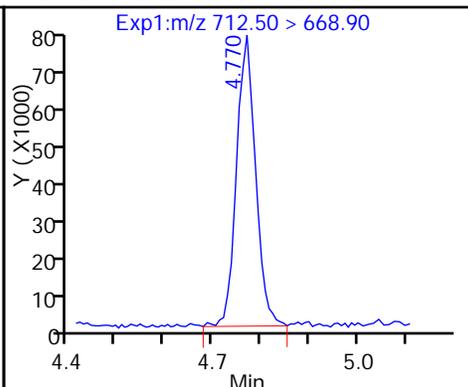
31 Perfluorotridecanoic acid



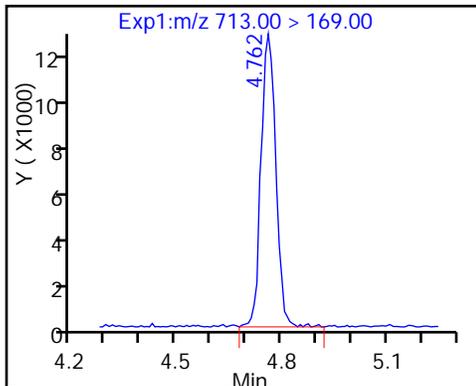
D 32 13C2-PFTeDa



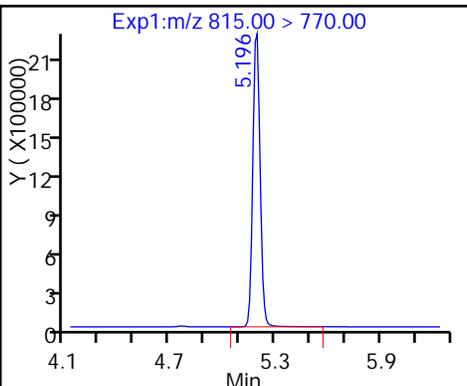
33 Perfluorotetradecanoic acid



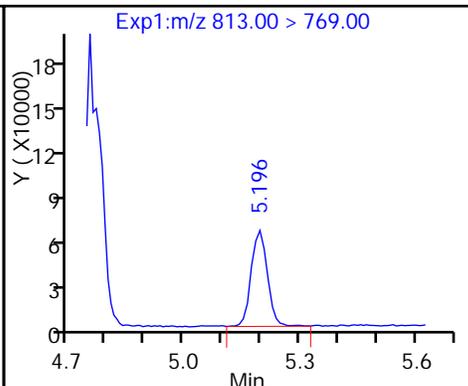
33 Perfluorotetradecanoic acid



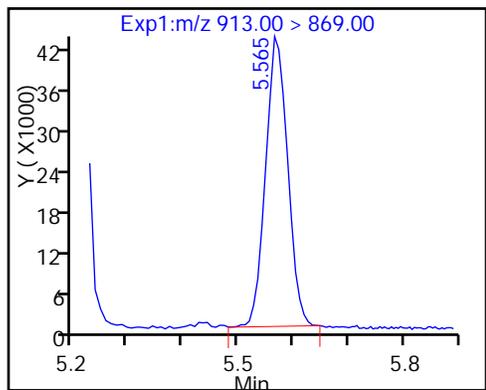
D 34 13C2-PFHxDa



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_006.d
 Lims ID: IC L3
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 03-Dec-2016 14:03:44 ALS Bottle#: 39 Worklist Smp#: 6
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L3_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:10 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:43:24

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	19799279	59.0		118	1507463	
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.577	-0.003	1758998	5.08		102	10276	
3 Perfluoropentanoic acid	262.90 > 219.00	1.858	1.861	-0.003	1580002	4.98		99.6	13754	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.861	-0.003	15634346	59.1		118	910049	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.897	1.900	-0.003	2768149	4.72		107		
	298.90 > 99.00	1.897	1.900	-0.003	1152600		2.40(0.00-0.00)	107		
7 Perfluorohexanoic acid	313.00 > 269.00	2.163	2.164	-0.001	1358530	5.11		102	38617	
D 6 13C2 PFHxA	315.00 > 270.00	2.163	2.164	-0.001	13959775	58.8		118	866927	
D 11 13C4-PFHpA	367.00 > 322.00	2.508	2.511	-0.003	12498264	60.3		121	1070818	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.516	2.512	0.004	1254320	4.89		97.7	16816	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.531	2.531	0.0	1846932	4.58		101		
D 10 18O2 PFHxS	403.00 > 84.00	2.531	2.531	0.0	17374498	55.6		118	872275	
D 14 13C4 PFOA	417.00 > 372.00	2.880	2.880	0.0	13257596	60.5		121	935997	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.880	2.887	-0.007	1.000	1455373	5.12		102	23006	
413.00 > 169.00	2.888	2.887	0.001	1.003	828495		1.76(0.90-1.10)	102	49224	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.888	2.888	0.0	1.000	1641199	4.77		100		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.255	3.258	-0.003	1.000	1443163	4.55		98.0	374092	
499.00 > 99.00	3.255	3.258	-0.003	1.000	319030		4.52(0.90-1.10)	98.0	32937	
D 17 13C4 PFOS										
503.00 > 80.00	3.263	3.259	0.004		13983781	56.8		119	604878	
D 19 13C5 PFNA										
468.00 > 423.00	3.263	3.263	0.0		10108070	60.7		121	545062	
20 Perfluorononanoic acid										
463.00 > 419.00	3.263	3.263	0.0	1.000	981559	4.87		97.5	16647	
D 21 13C8 FOSA										
506.00 > 78.00	3.576	3.571	0.005		23126232	57.5		115	745838	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.576	3.574	0.002	1.000	2262174	5.24		105	133644	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.618	3.623	-0.005	1.000	868604	4.91		98.1	29437	
D 23 13C2 PFDA										
515.00 > 470.00	3.627	3.626	0.001		9213875	58.4		117	291561	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.938	3.936	0.002	1.000	868168	4.64		96.2		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.956	3.955	0.001	1.000	704273	4.66		93.2	19079	
D 27 13C2 PFUnA										
565.00 > 520.00	3.956	3.958	-0.002		7086903	59.7		119	336150	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.253	4.250	0.003	1.000	606951	4.85		96.9	11284	
D 30 13C2 PFDoA										
615.00 > 570.00	4.253	4.251	0.002		6599493	58.9		118	172819	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.521	4.518	0.003	1.000	627673	5.01		100	1096	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.761	4.759	0.002		13488161	58.3		117	1155339	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.761	4.761	0.0	1.000	1175506	4.80		96.1	484	
713.00 > 169.00	4.761	4.761	0.0	1.000	194335		6.05(0.00-0.00)	96.1	68892	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.180	5.186	-0.006		7591947	58.5		117	144235	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.191	5.186	0.005	1.000	766480	5.09		102	693	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.561	5.559	0.002	1.000	673971	5.14		103	887	

Reagents:

LCPFC-L3_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_006.d

Injection Date: 03-Dec-2016 14:03:44

Instrument ID: A8_N

Lims ID: IC L3

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 39

Worklist Smp#: 6

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

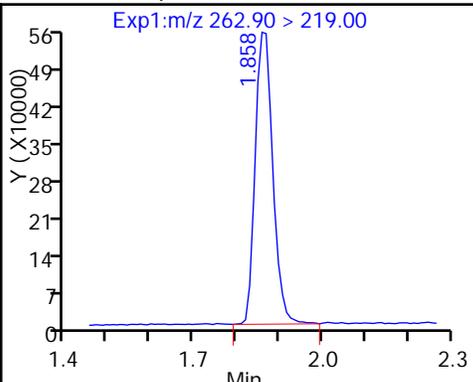
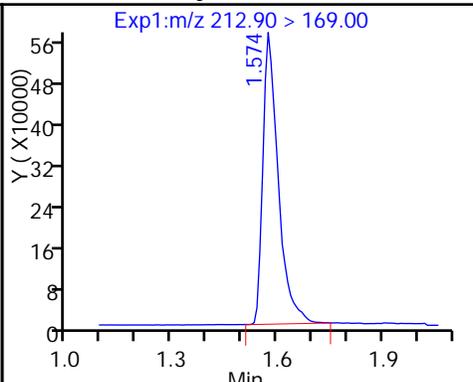
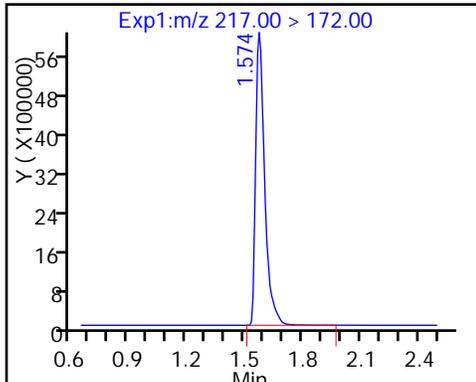
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

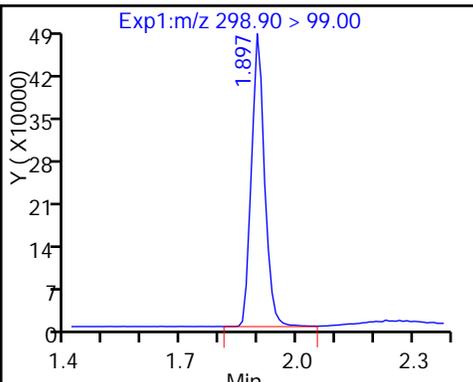
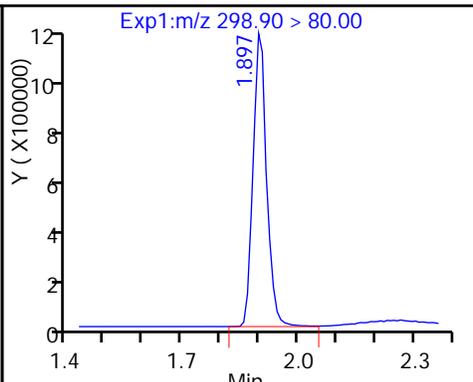
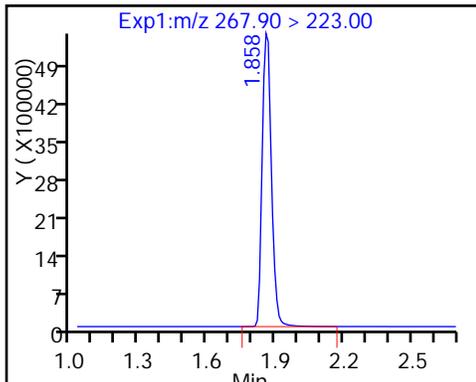
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

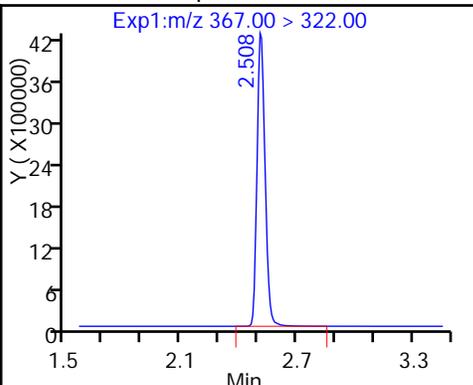
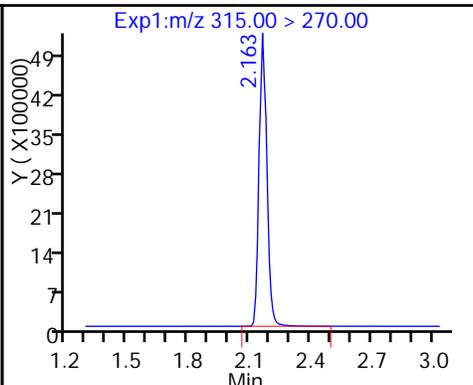
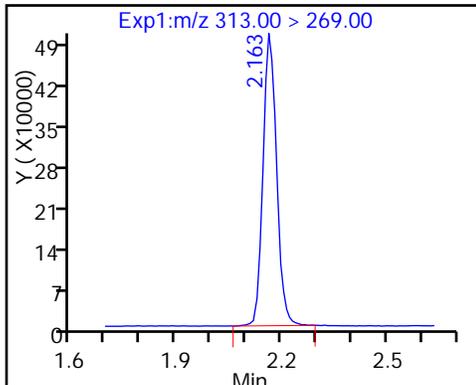
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

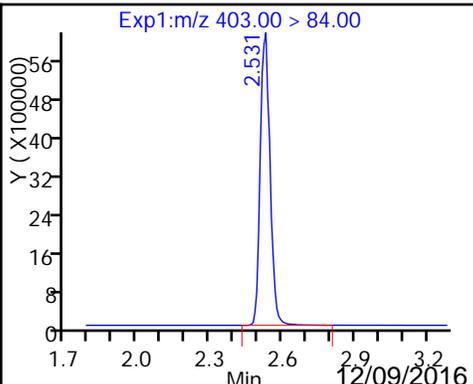
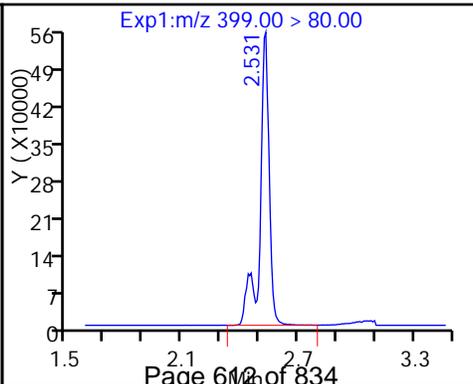
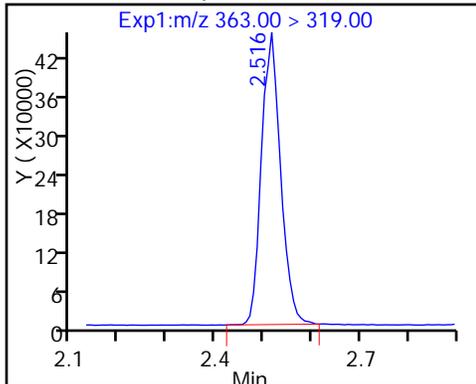
D 11 13C4-PFHpA



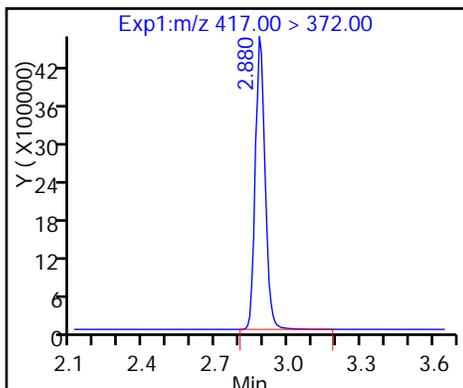
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

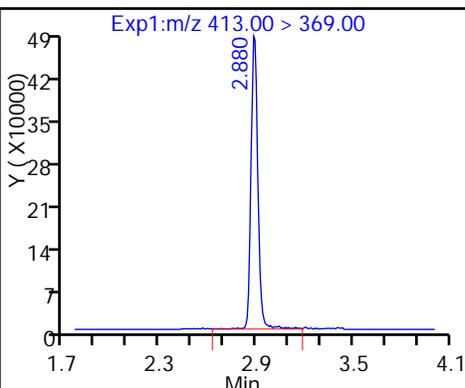
D 10 18O2 PFHxS



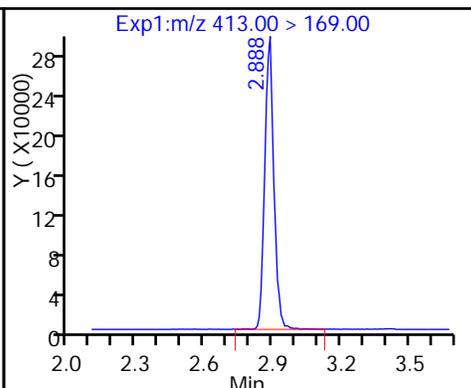
D 14 13C4 PFOA



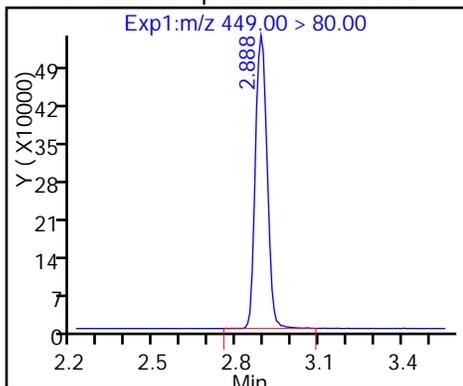
15 Perfluorooctanoic acid



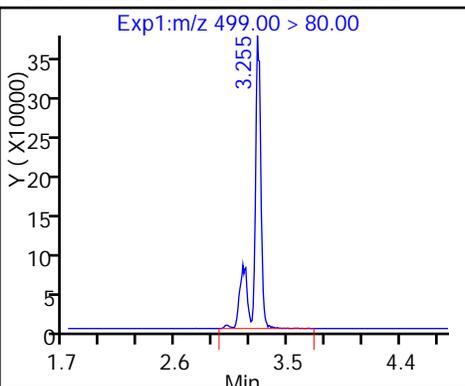
15 Perfluorooctanoic acid



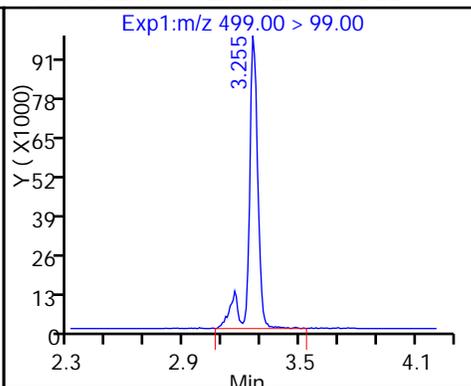
13 Perfluoroheptanesulfonic Acid



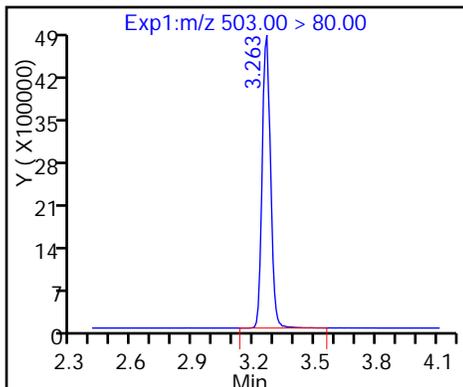
18 Perfluorooctane sulfonic acid



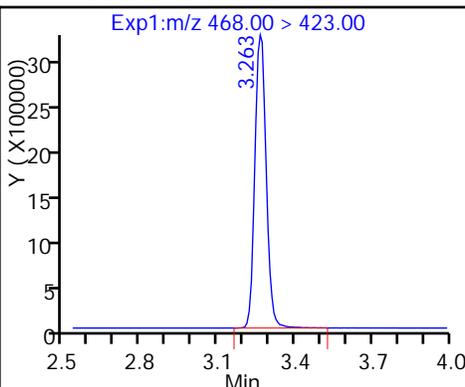
18 Perfluorooctane sulfonic acid



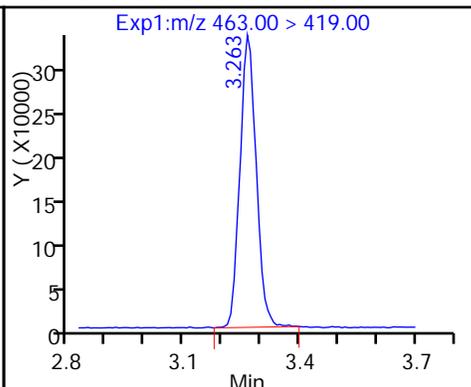
D 17 13C4 PFOS



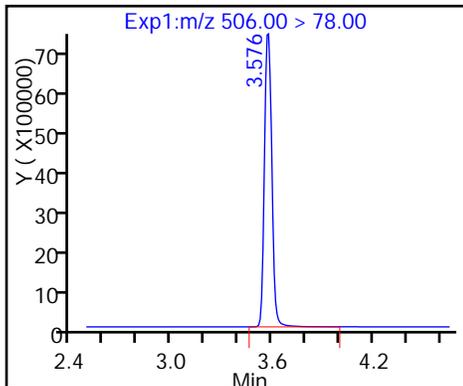
D 19 13C5 PFNA



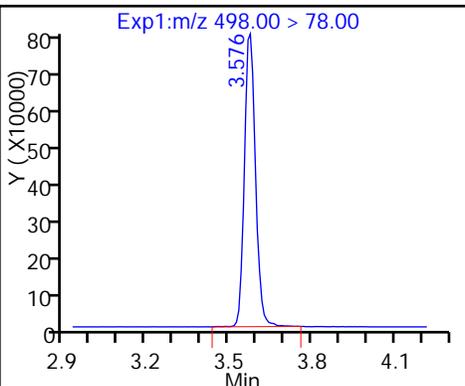
20 Perfluorononanoic acid



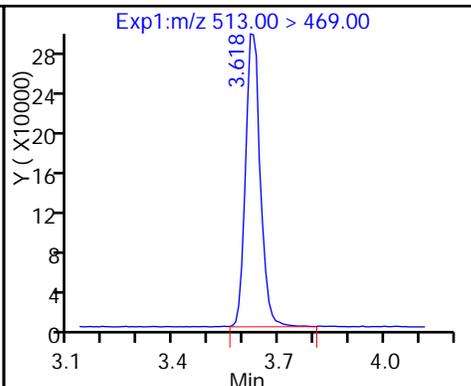
D 21 13C8 FOSA



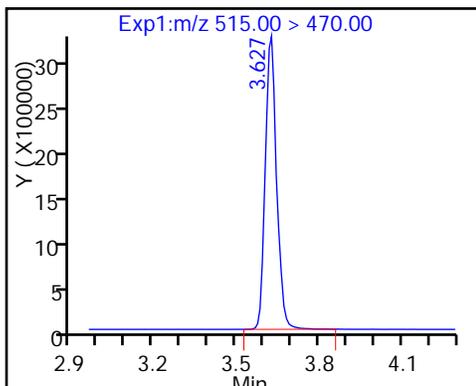
22 Perfluorooctane Sulfonamide



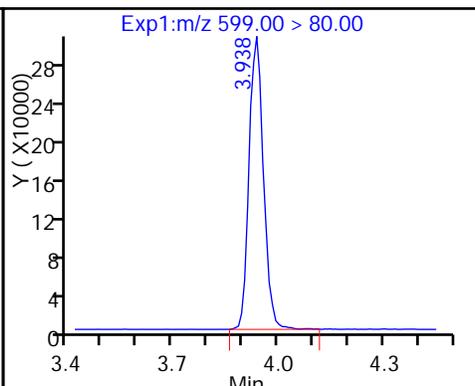
24 Perfluorodecanoic acid



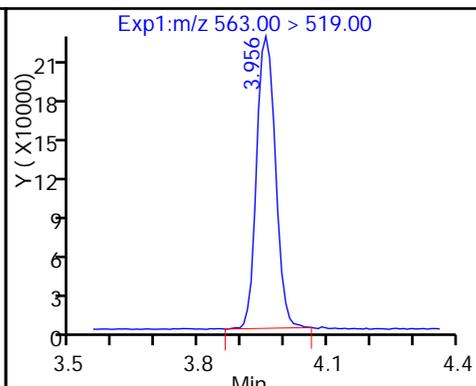
D 23 13C2 PFDA



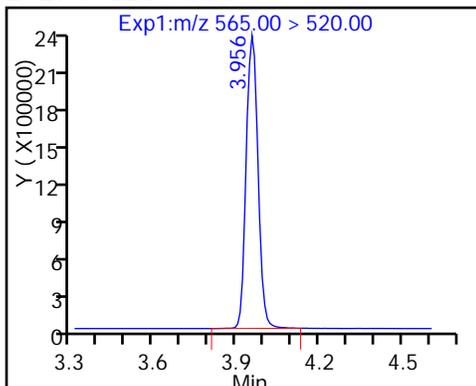
26 Perfluorodecane Sulfonic acid



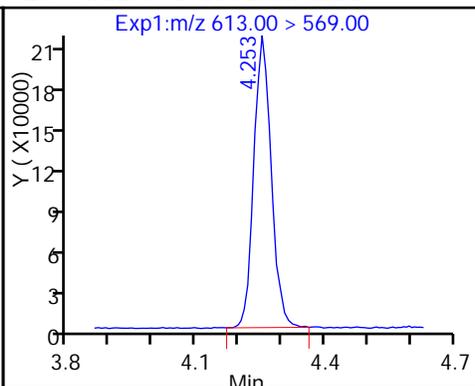
28 Perfluoroundecanoic acid



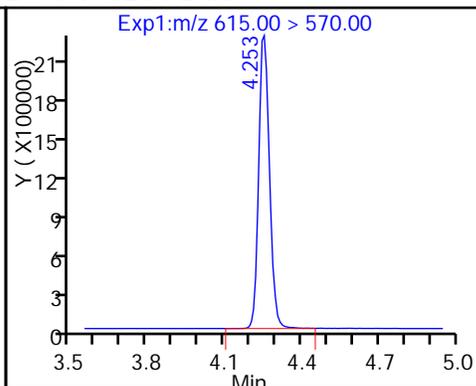
D 27 13C2 PFUa



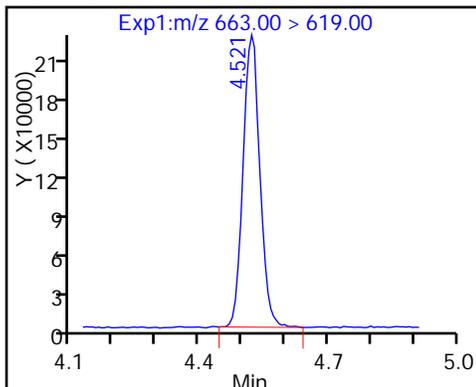
29 Perfluorododecanoic acid



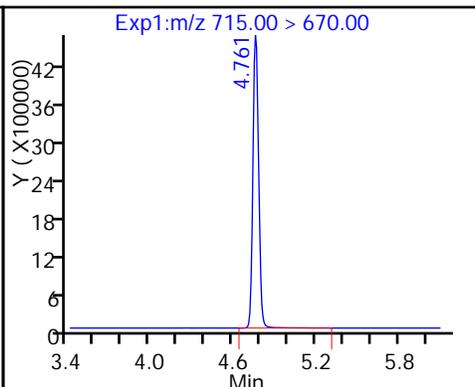
D 30 13C2 PFDa



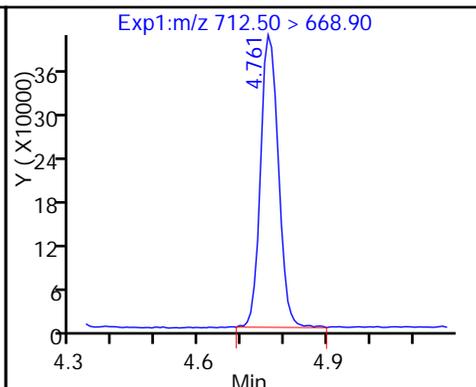
31 Perfluorotridecanoic acid



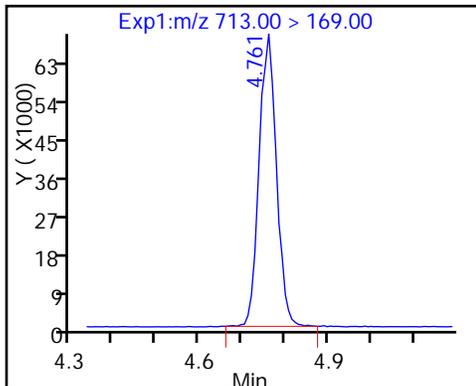
D 32 13C2-PFTeDA



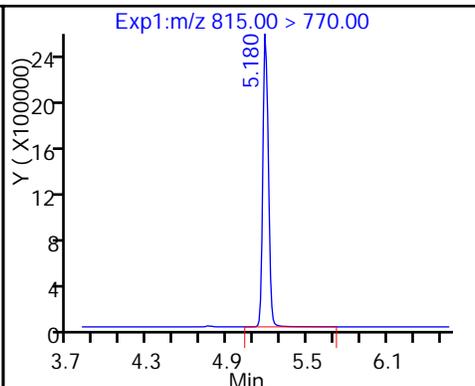
33 Perfluorotetradecanoic acid



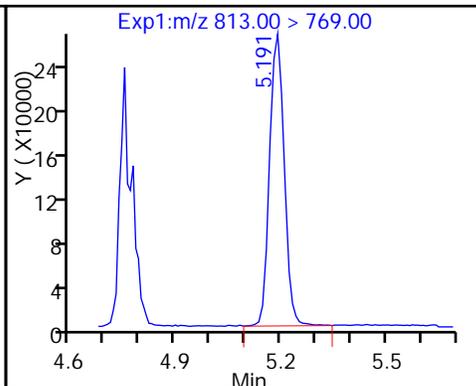
33 Perfluorotetradecanoic acid



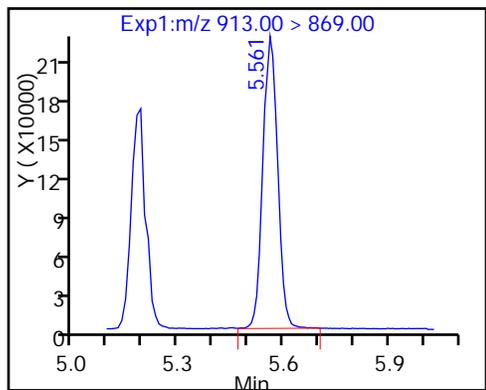
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_007.d
 Lims ID: IC L4
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 03-Dec-2016 14:11:14 ALS Bottle#: 40 Worklist Smp#: 7
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L4_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:13 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:41:57

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	18254473	54.4		109	916244	
1 Perfluorobutyric acid	212.90 > 169.00	1.582	1.577	0.005	7008467	22.0		110	42848	
3 Perfluoropentanoic acid	262.90 > 219.00	1.868	1.861	0.007	6115663	20.9		104	53754	
D 4 13C5-PFPeA	267.90 > 223.00	1.868	1.861	0.007	14426922	54.5		109	1030803	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.907	1.900	0.007	10888534	20.0		113		
	298.90 > 99.00	1.897	1.900	-0.003	4709173		2.31(0.00-0.00)	113		
7 Perfluorohexanoic acid	313.00 > 269.00	2.171	2.164	0.007	5224102	21.6		108	141802	
D 6 13C2 PFHxA	315.00 > 270.00	2.162	2.164	-0.002	12716423	53.5		107	709396	
D 11 13C4-PFHpA	367.00 > 322.00	2.512	2.511	0.001	11474824	55.3		111	830208	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.512	2.512	0.0	4918226	20.9		104	66176	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.527	2.531	-0.004	6968734	18.7		102		M M
D 10 18O2 PFHxS	403.00 > 84.00	2.535	2.531	0.004	16099619	51.5		109	1076833	
D 14 13C4 PFOA	417.00 > 372.00	2.886	2.880	0.006	12069172	55.0		110	656359	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										M
413.00 > 369.00	2.894	2.887	0.007	1.000	5292388	20.5		102	50246	M
413.00 > 169.00	2.886	2.887	-0.001	0.997	3252983		1.63(0.90-1.10)	102	188817	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.886	2.888	-0.002	1.000	6678298	20.6		108		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.259	3.258	0.001	1.000	5631788	18.8		101	741165	
499.00 > 99.00	3.259	3.258	0.001	1.000	1257953		4.48(0.90-1.10)	101	196382	
D 17 13C4 PFOS										
503.00 > 80.00	3.259	3.259	0.0		13189624	53.6		112	527854	
D 19 13C5 PFNA										
468.00 > 423.00	3.267	3.263	0.004		8920352	53.6		107	500218	
20 Perfluorononanoic acid										
463.00 > 419.00	3.267	3.263	0.004	1.000	3804327	21.4		107	70301	
D 21 13C8 FOSA										
506.00 > 78.00	3.573	3.571	0.002		22326638	55.5		111	379770	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.573	3.574	-0.001	1.000	9111782	21.8		109	354493	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.623	3.623	0.0	1.000	3434382	20.3		102	104771	
D 23 13C2 PFDA										
515.00 > 470.00	3.623	3.626	-0.003		8786342	55.7		111	333180	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.939	3.936	0.003	1.000	3548629	20.1		104		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.956	3.955	0.001	1.000	2673576	19.3		96.7	68734	
D 27 13C2 PFUnA										
565.00 > 520.00	3.965	3.958	0.007		6485474	54.6		109	683577	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.253	4.250	0.003	1.000	2415187	21.0		105	47125	
D 30 13C2 PFDoA										
615.00 > 570.00	4.253	4.251	0.002		6067987	54.1		108	181188	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.524	4.518	0.006	1.000	2373060	20.6		103	5080	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.764	4.759	0.005		12877999	55.7		111	442175	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.764	4.761	0.003	1.000	4681126	20.8		104	1697	
713.00 > 169.00	4.756	4.761	-0.005	0.998	771098		6.07(0.00-0.00)	104	90317	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.187	5.186	0.001		7363997	56.8		114	132680	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.187	5.186	0.001	1.000	2870345	23.1		115	2476	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.563	5.559	0.004	1.000	2205153	18.3		91.5	3091	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L4_00022

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_007.d

Injection Date: 03-Dec-2016 14:11:14

Instrument ID: A8_N

Lims ID: IC L4

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 40

Worklist Smp#: 7

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

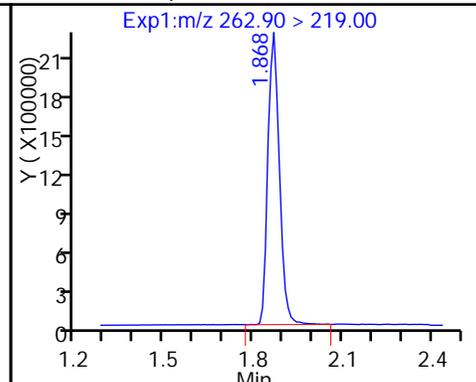
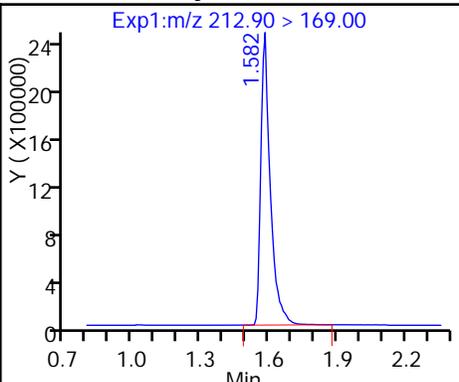
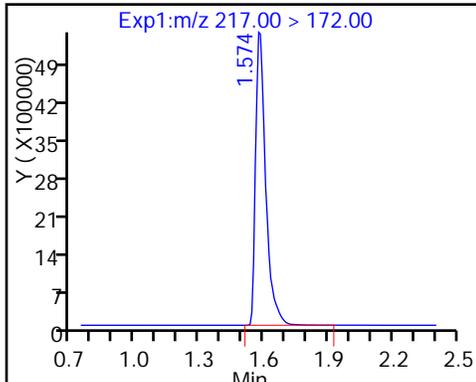
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

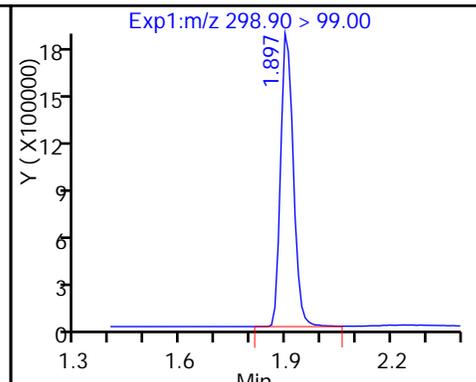
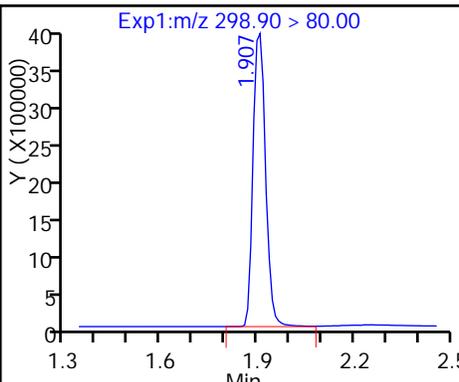
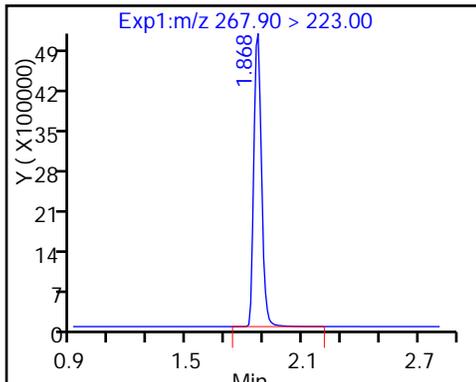
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

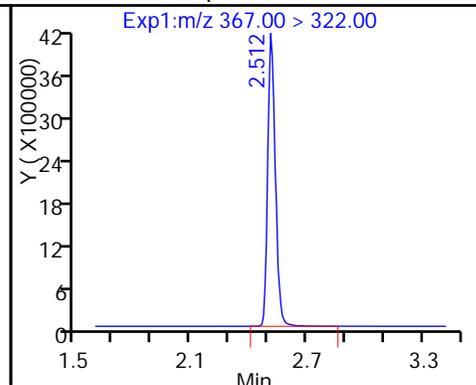
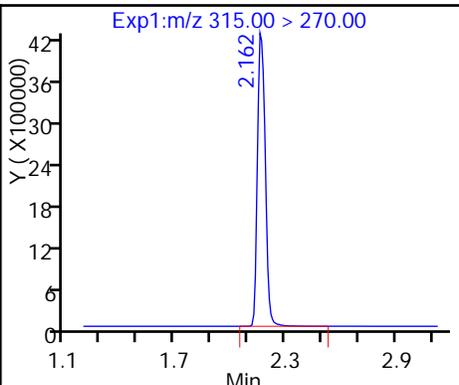
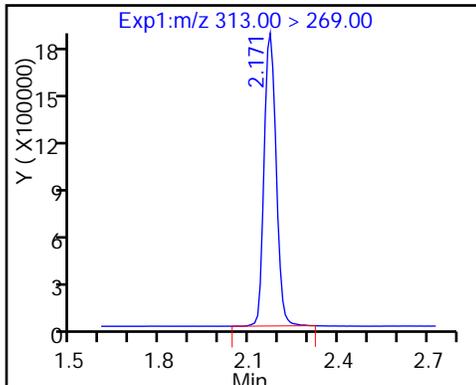
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

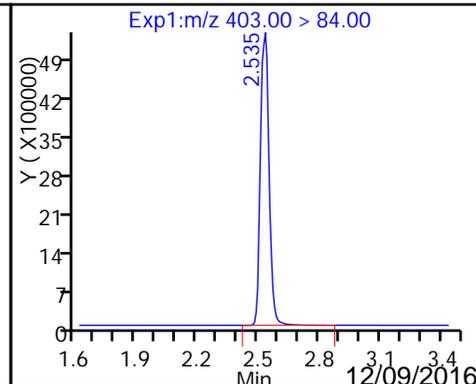
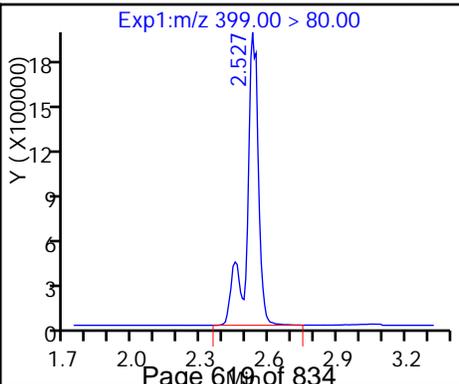
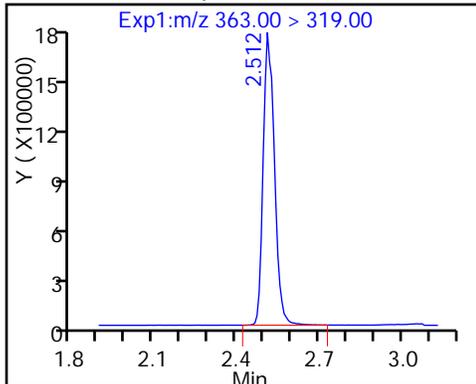
D 11 13C4-PFHpA



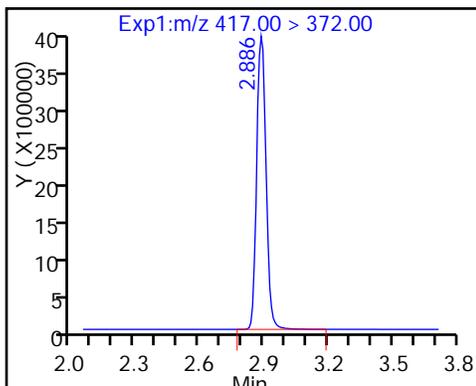
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid (M)

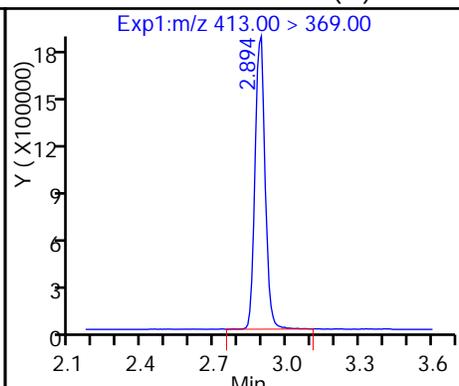
D 10 18O2 PFHxS



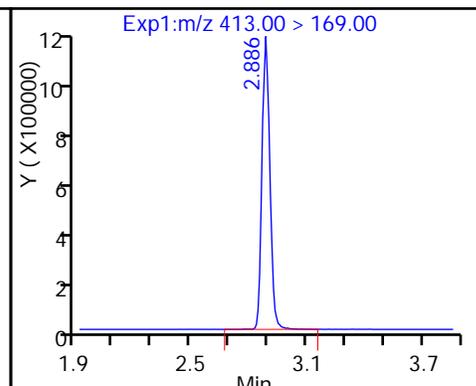
D 14 13C4 PFOA



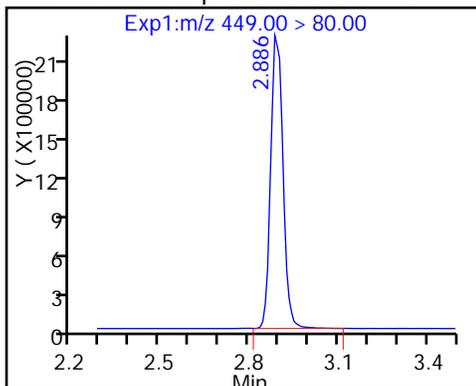
15 Perfluorooctanoic acid (M)



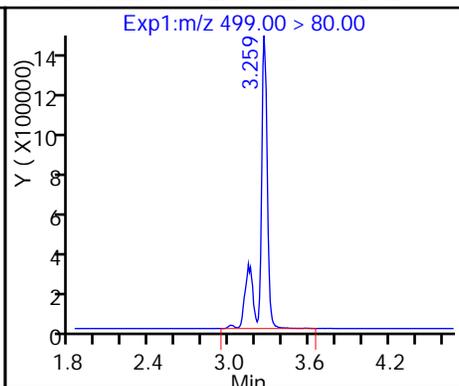
15 Perfluorooctanoic acid



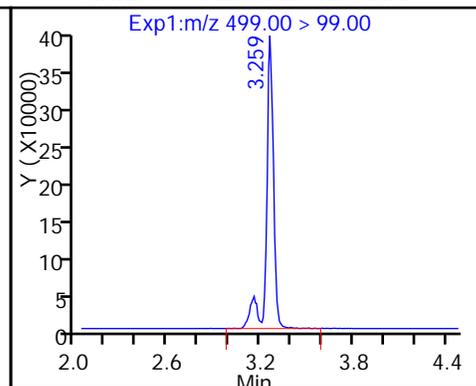
13 Perfluoroheptanesulfonic Acid



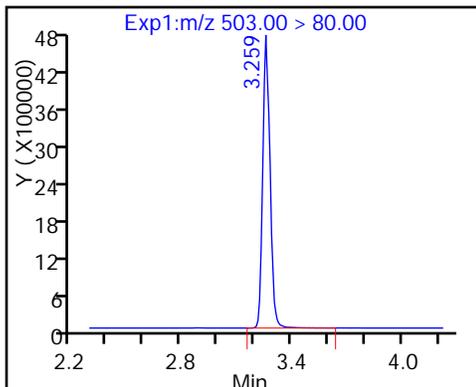
18 Perfluorooctane sulfonic acid



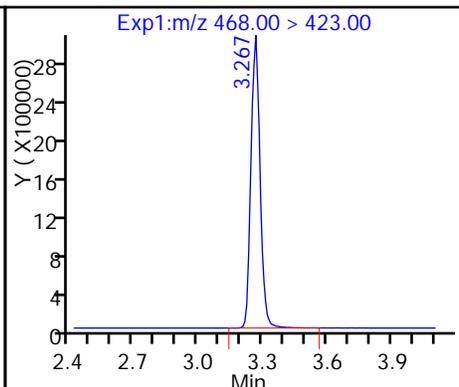
18 Perfluorooctane sulfonic acid



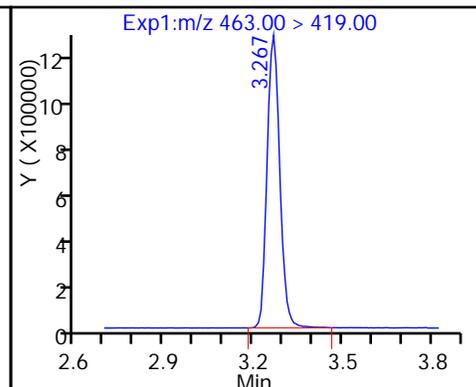
D 17 13C4 PFOS



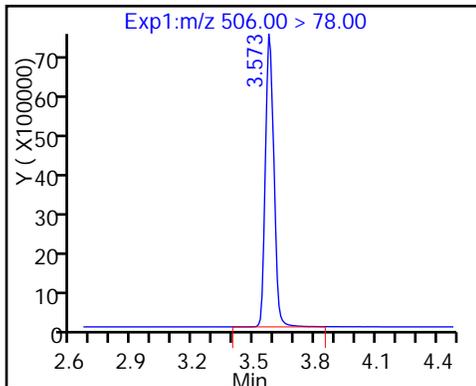
D 19 13C5 PFNA



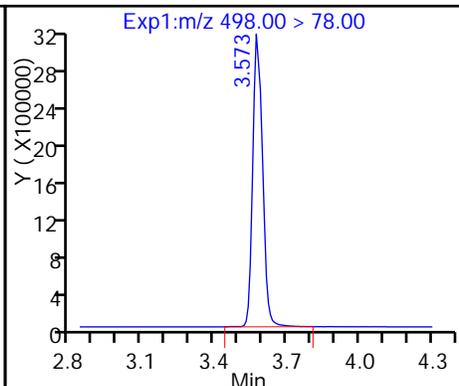
20 Perfluorononanoic acid



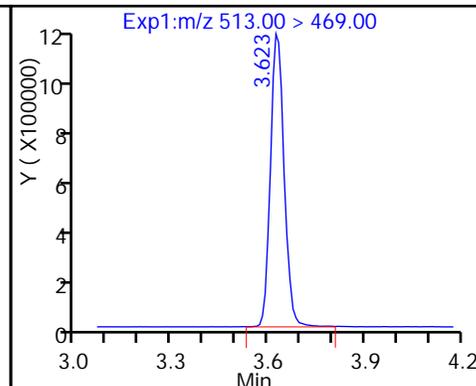
D 21 13C8 FOSA



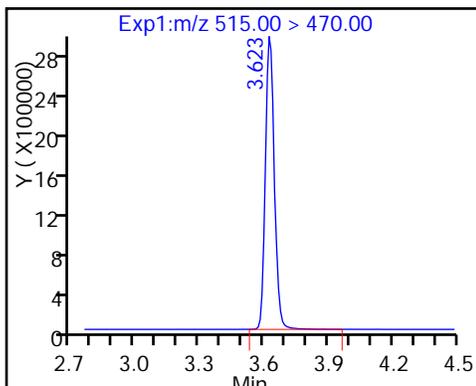
22 Perfluorooctane Sulfonamide



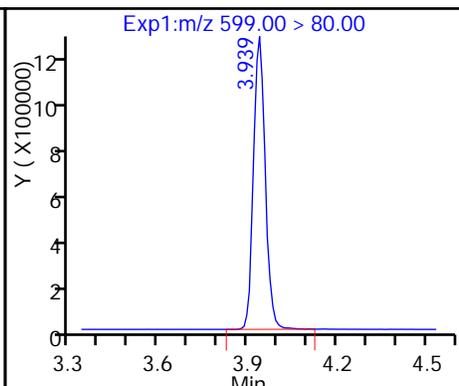
24 Perfluorodecanoic acid



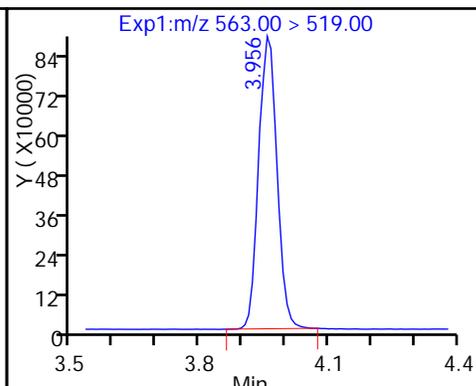
D 23 13C2 PFDA



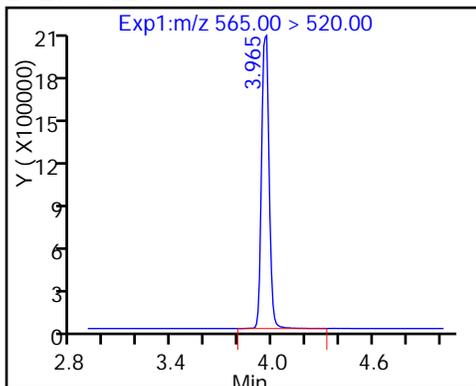
26 Perfluorodecane Sulfonic acid



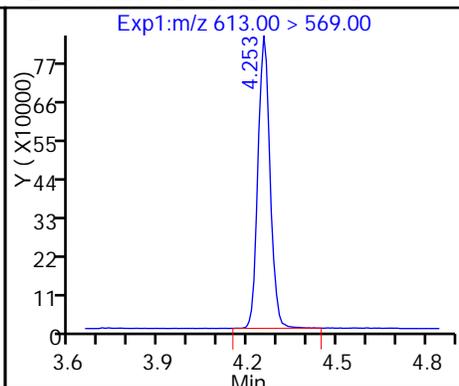
28 Perfluoroundecanoic acid



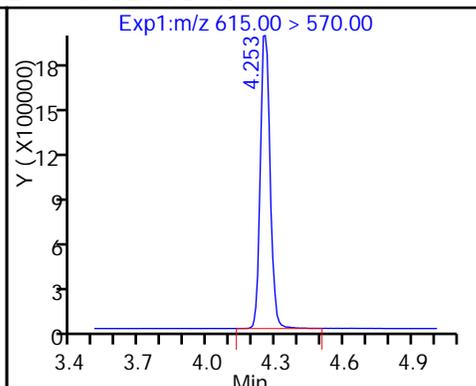
D 27 13C2 PFUa



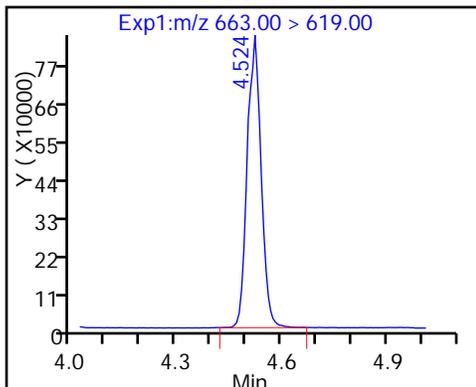
29 Perfluorododecanoic acid



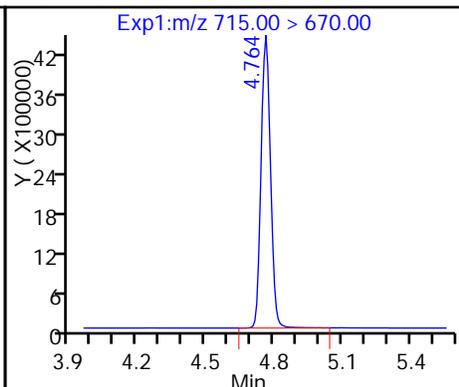
D 30 13C2 PFDa



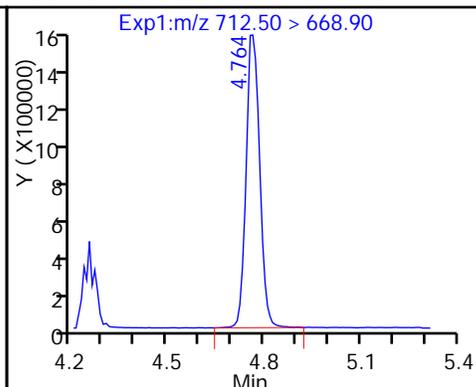
31 Perfluorotridecanoic acid



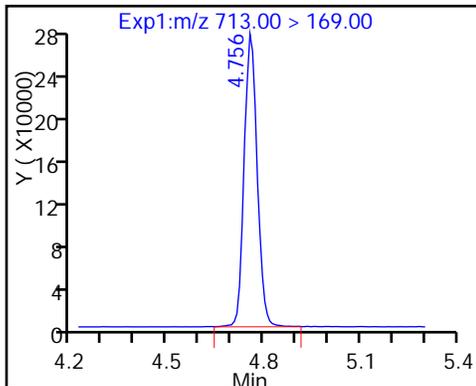
D 32 13C2-PFTeDA



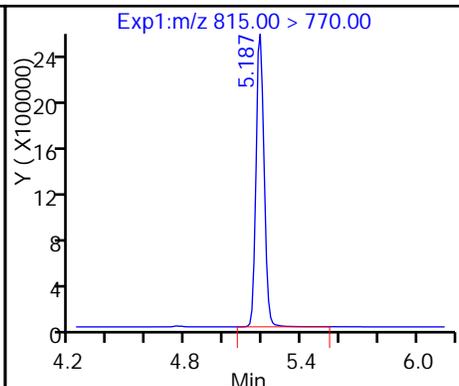
33 Perfluorotetradecanoic acid



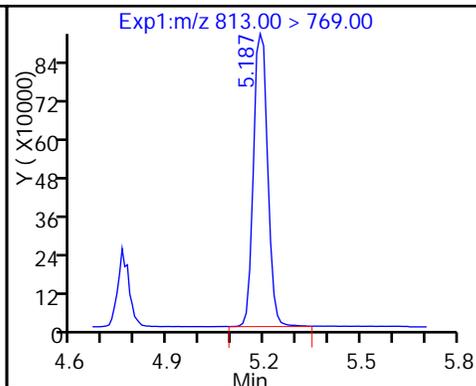
33 Perfluorotetradecanoic acid



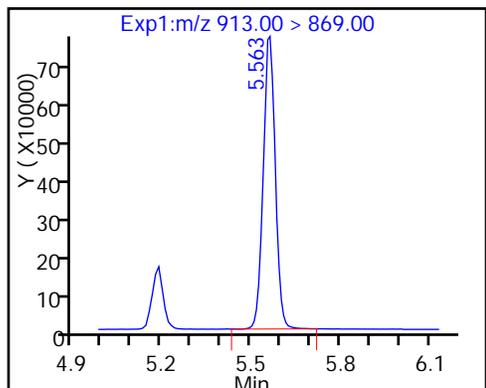
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



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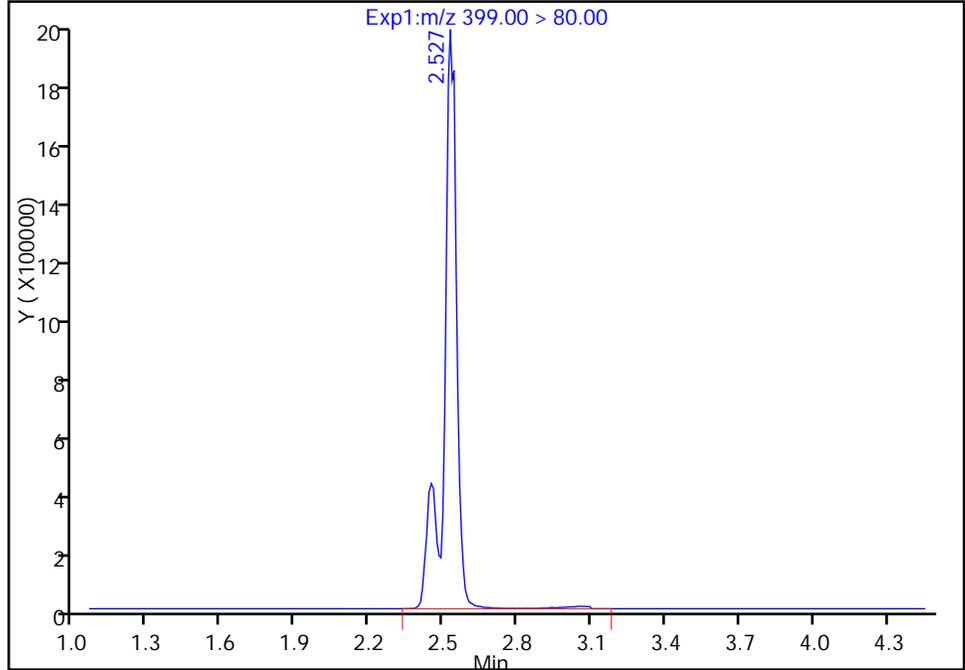
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_007.d
Injection Date: 03-Dec-2016 14:11:14 Instrument ID: A8_N
Lims ID: IC L4
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 40 Worklist Smp#: 7
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

9 Perfluorohexanesulfonic acid, CAS: 355-46-4

Signal: 1

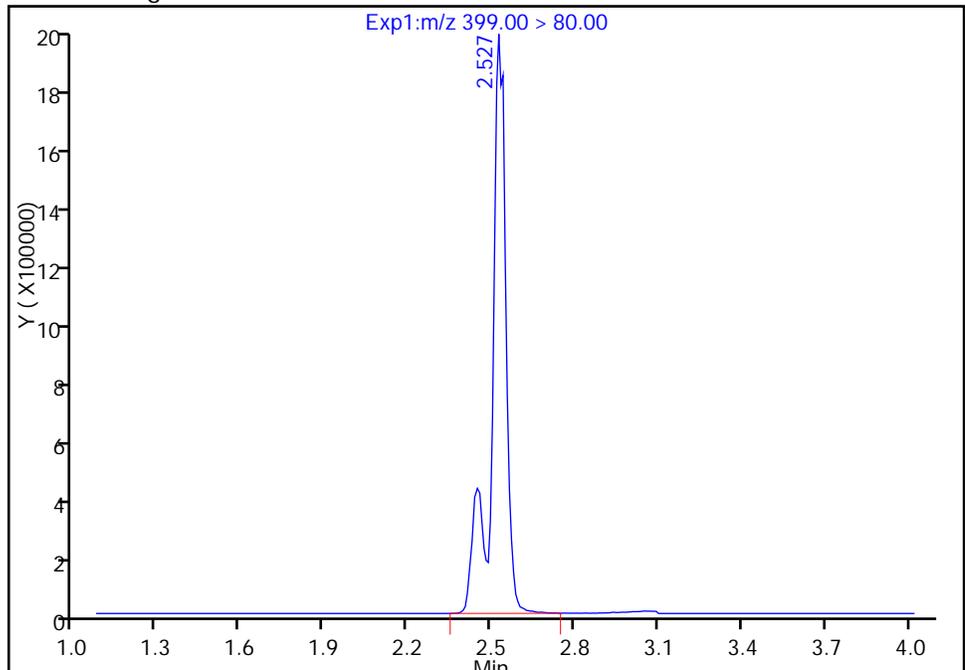
RT: 2.53
Area: 7050184
Amount: 18.826548
Amount Units: ng/ml

Processing Integration Results



RT: 2.53
Area: 6968734
Amount: 18.653631
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 09:41:57
Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento

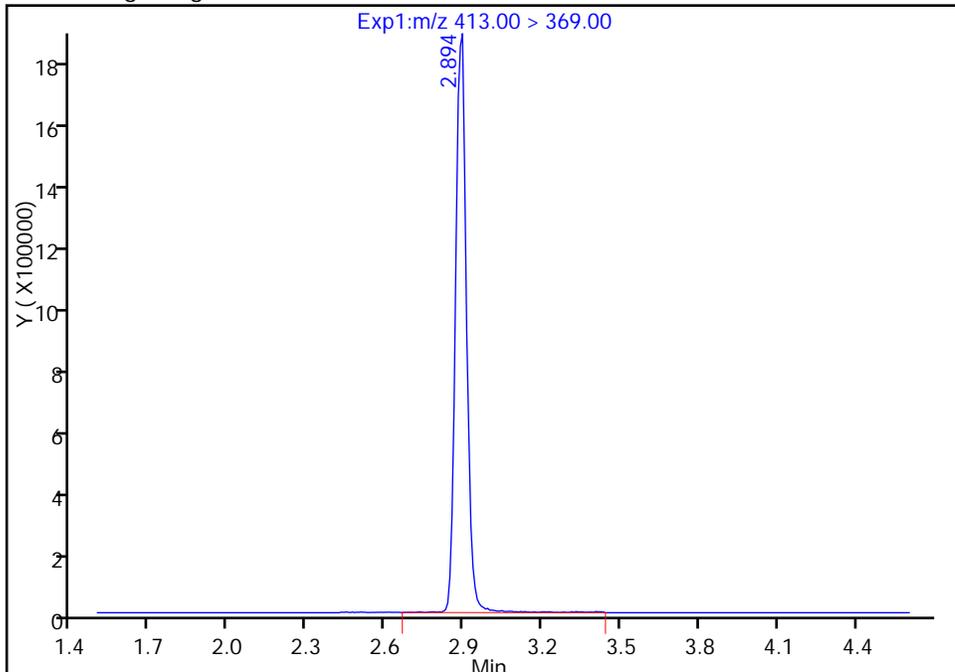
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_007.d
Injection Date: 03-Dec-2016 14:11:14 Instrument ID: A8_N
Lims ID: IC L4
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 40 Worklist Smp#: 7
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

15 Perfluorooctanoic acid, CAS: 335-67-1

Signal: 1

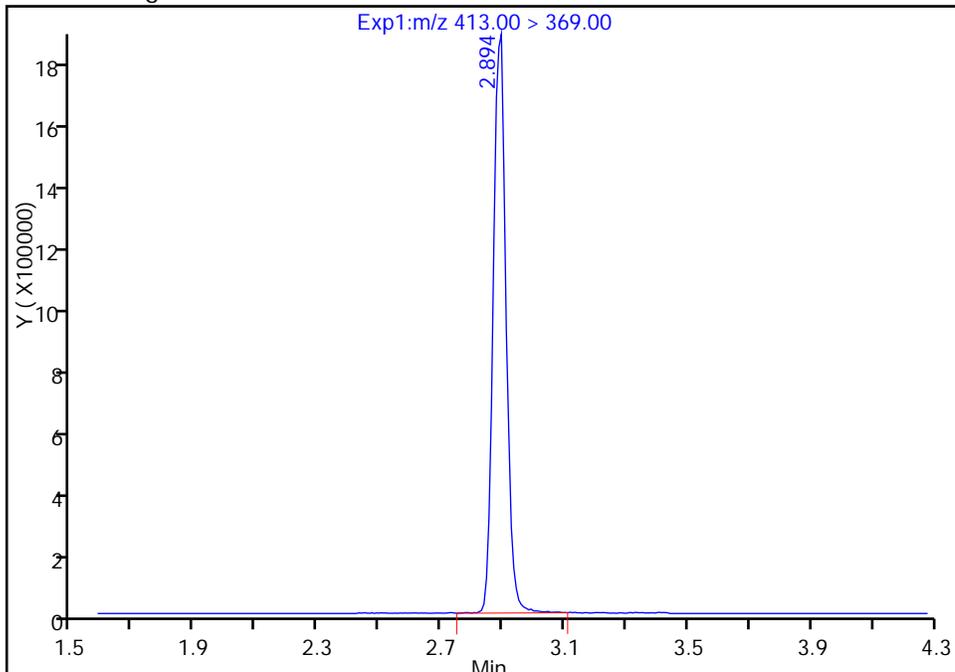
RT: 2.89
Area: 5360428
Amount: 20.662890
Amount Units: ng/ml

Processing Integration Results



RT: 2.89
Area: 5292388
Amount: 20.454262
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 09:41:57

Audit Action: Manually Integrated

Audit Reason: Baseline

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_008.d
 Lims ID: IC L5
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 03-Dec-2016 14:18:44 ALS Bottle#: 41 Worklist Smp#: 8
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L5_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:15 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:43:57

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	16029004	47.7		95.5	1234710	
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.577	-0.003	14048645	50.1		100	74155	
3 Perfluoropentanoic acid	262.90 > 219.00	1.858	1.861	-0.003	12597041	49.7		99.5	116129	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.861	-0.003	12477636	47.2		94.3	1170062	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.897	1.900	-0.003	21513660	44.6		101		
	298.90 > 99.00	1.897	1.900	-0.003	9875959		2.18(0.00-0.00)	101		
7 Perfluorohexanoic acid	313.00 > 269.00	2.160	2.164	-0.004	10420431	49.8		99.5	246256	
D 6 13C2 PFHxA	315.00 > 270.00	2.160	2.164	-0.004	10985481	46.3		92.5	787845	
D 11 13C4-PFHpA	367.00 > 322.00	2.510	2.511	-0.001	9653835	46.5		93.1	803660	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.510	2.512	-0.002	9915679	50.0		100	147911	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.525	2.531	-0.006	14498414	43.7		96.0		
D 10 18O2 PFHxS	403.00 > 84.00	2.525	2.531	-0.006	14303487	45.8		96.8	5055434	
D 14 13C4 PFOA	417.00 > 372.00	2.876	2.880	-0.004	9909434	45.2		90.4	408580	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.884	2.887	-0.003	1.000	10434085	49.1		98.2	197146	
413.00 > 169.00	2.876	2.887	-0.011	0.997	6358650		1.64(0.90-1.10)	98.2	316749	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.884	2.888	-0.004	1.000	13806464	49.7		104		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.257	3.258	-0.001	1.000	11953842	46.7		101	39071	
499.00 > 99.00	3.248	3.258	-0.010	0.997	2570510		4.65(0.90-1.10)	101	263704	
D 17 13C4 PFOS										
503.00 > 80.00	3.257	3.259	-0.002		11281105	45.9		95.9	268992	
D 19 13C5 PFNA										
468.00 > 423.00	3.257	3.263	-0.006		7854327	47.2		94.4	841834	
20 Perfluorononanoic acid										
463.00 > 419.00	3.257	3.263	-0.006	1.000	7675034	49.0		98.1	142675	
D 21 13C8 FOSA										
506.00 > 78.00	3.570	3.571	-0.001		19293148	48.0		95.9	389510	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.570	3.574	-0.004	1.000	18269014	50.7		101	626245	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.620	3.623	-0.003	1.000	7159278	50.3		101	175242	
D 23 13C2 PFDA										
515.00 > 470.00	3.620	3.626	-0.006		7405870	46.9		93.9	311468	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.928	3.936	-0.008	1.000	7347809	48.7		101		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.946	3.955	-0.009	1.000	5553051	46.3		92.6	133337	
D 27 13C2 PFUnA										
565.00 > 520.00	3.946	3.958	-0.012		5625237	47.4		94.7	281129	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.242	4.250	-0.008	1.000	5000795	49.4		98.7	86155	
D 30 13C2 PFDoA										
615.00 > 570.00	4.242	4.251	-0.009		5338479	47.6		95.3	147109	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.510	4.518	-0.008	1.000	5034109	49.6		99.3	21741	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.750	4.759	-0.009		11031119	47.7		95.4	553742	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.750	4.761	-0.011	1.000	9601518	48.5		97.0	4445	
713.00 > 169.00	4.750	4.761	-0.011	1.000	1523287		6.30(0.00-0.00)	97.0	181298	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.177	5.186	-0.009		5893329	45.4		90.9	104400	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.177	5.186	-0.009	1.000	5382109	50.0		100	6406	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.547	5.559	-0.012	1.000	5729058	54.0		108	9772	

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_008.d

Injection Date: 03-Dec-2016 14:18:44

Instrument ID: A8_N

Lims ID: IC L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 8

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

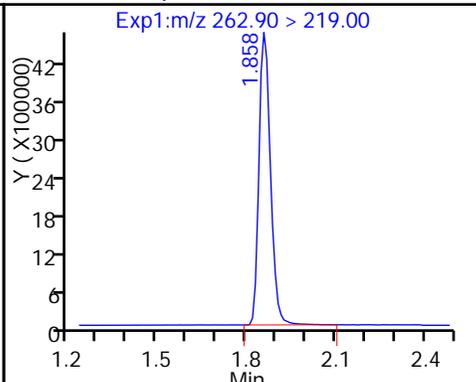
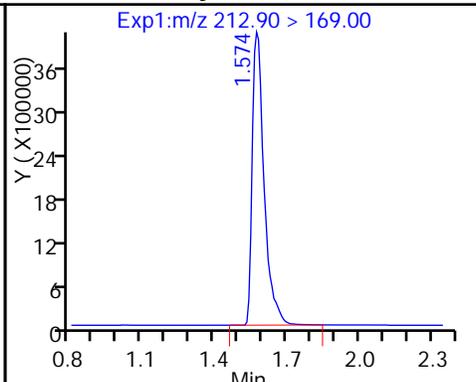
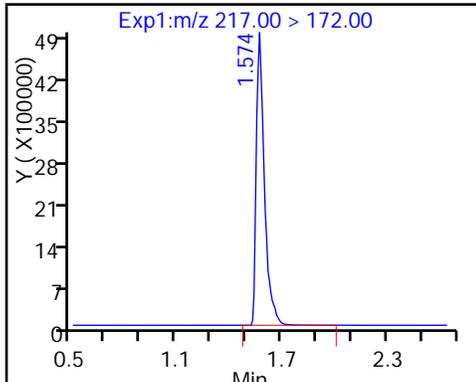
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

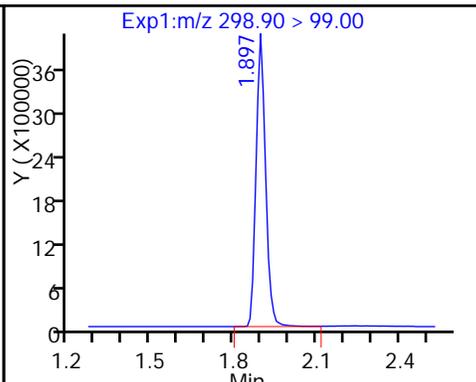
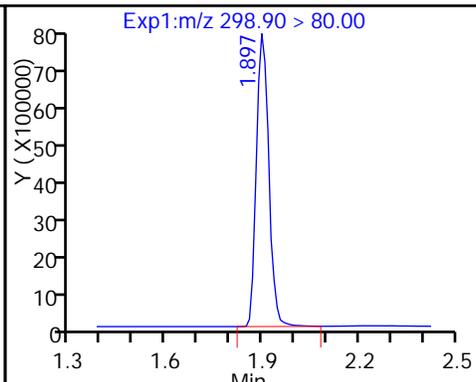
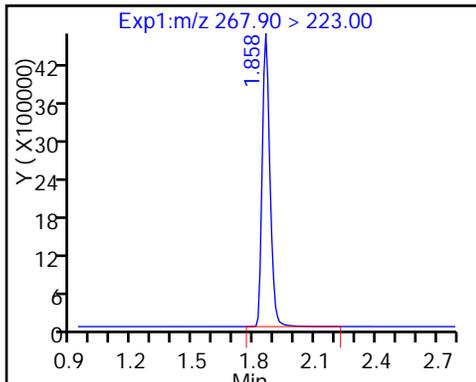
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

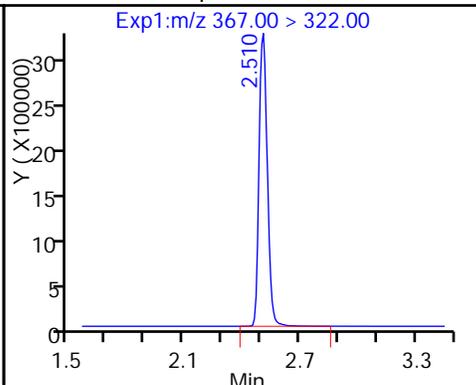
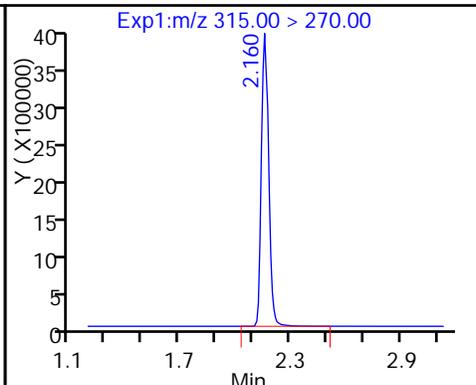
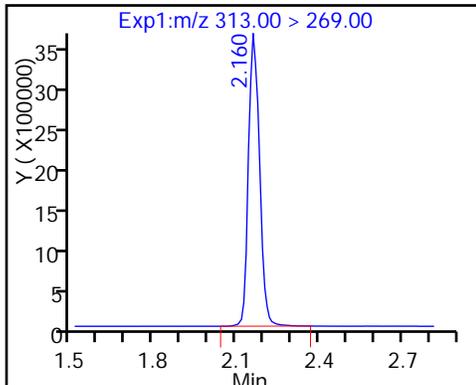
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

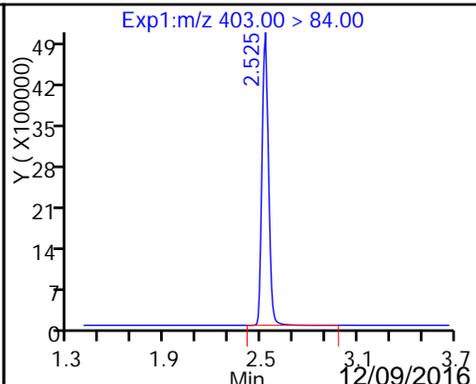
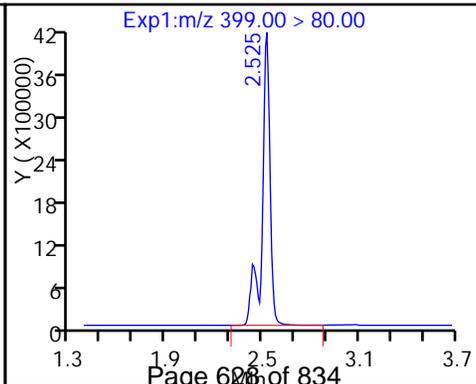
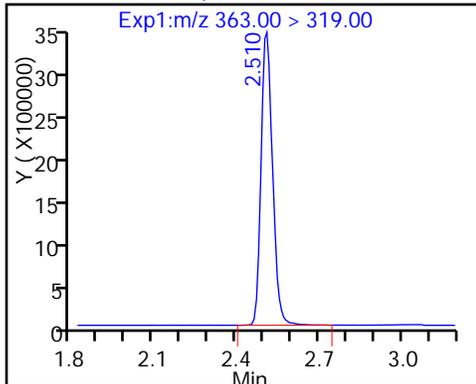
D 11 13C4-PFHpA



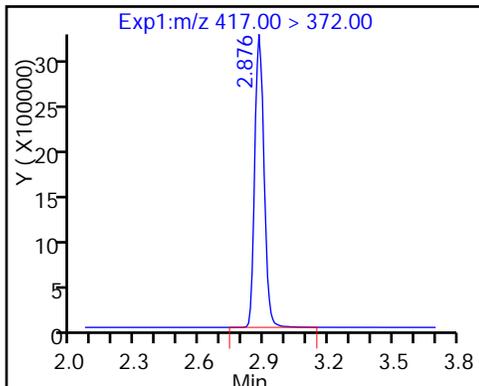
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

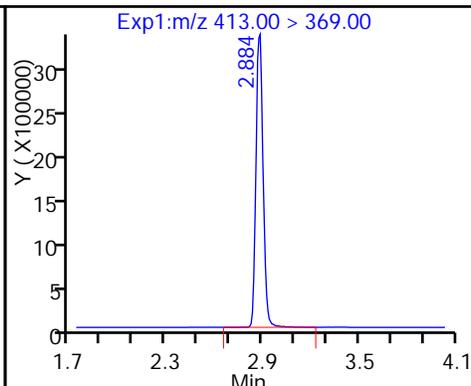
D 10 18O2 PFHxS



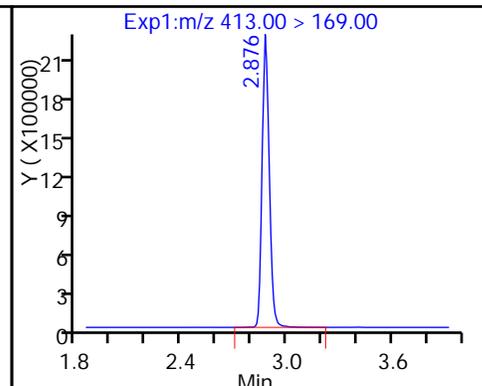
D 14 13C4 PFOA



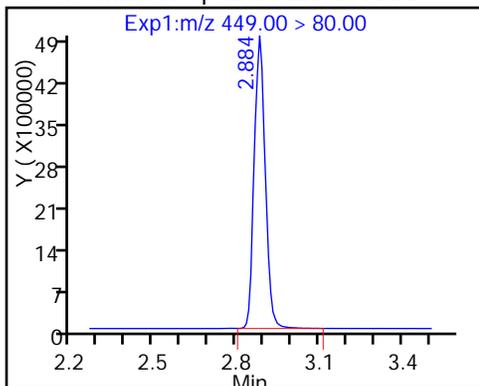
15 Perfluorooctanoic acid



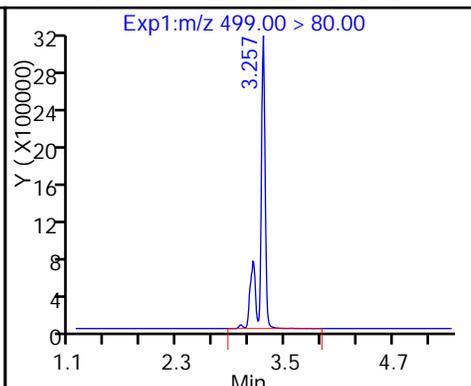
15 Perfluorooctanoic acid



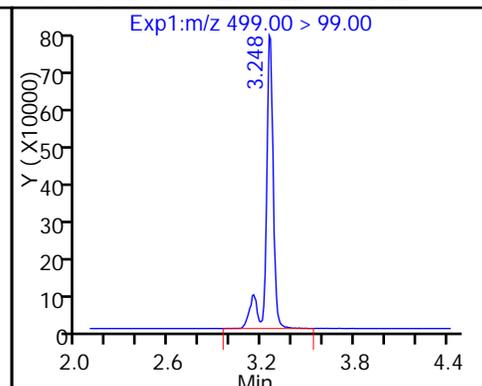
13 Perfluoroheptanesulfonic Acid



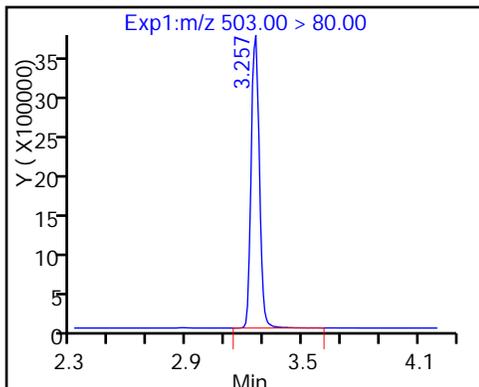
18 Perfluorooctane sulfonic acid



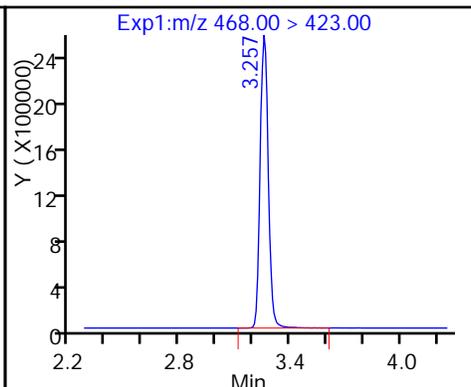
18 Perfluorooctane sulfonic acid



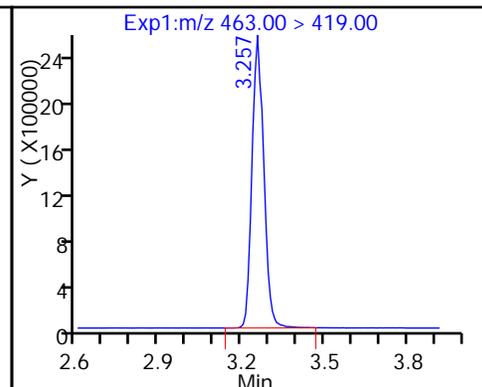
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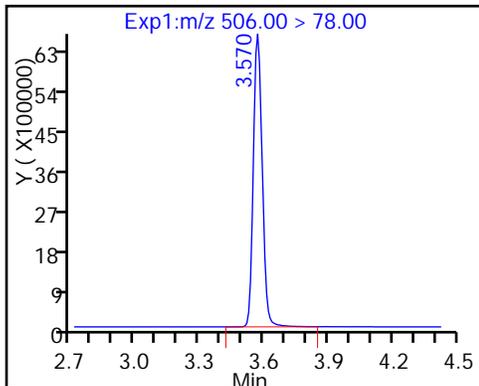
D 19 13C5 PFNA



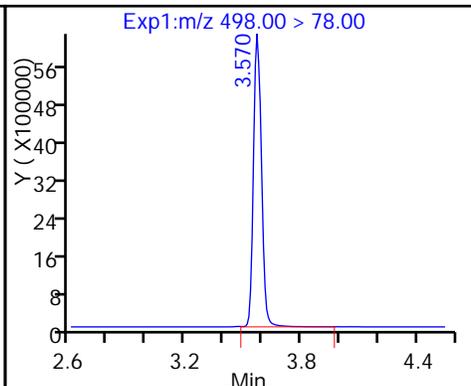
20 Perfluorononanoic acid



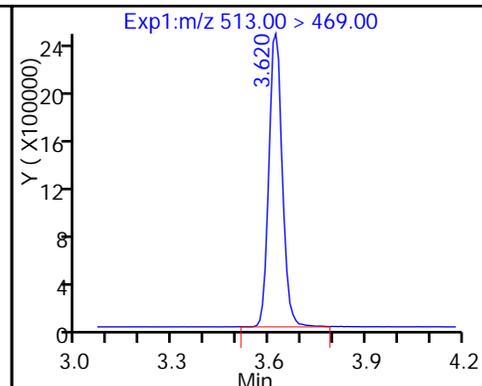
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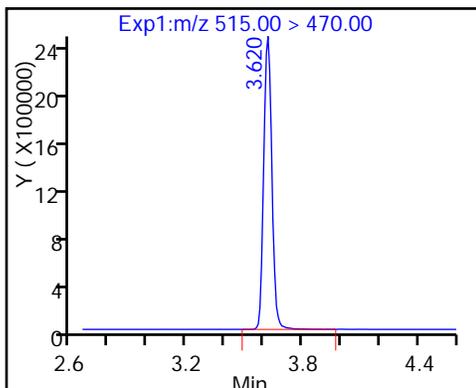
22 Perfluorooctane Sulfonamide



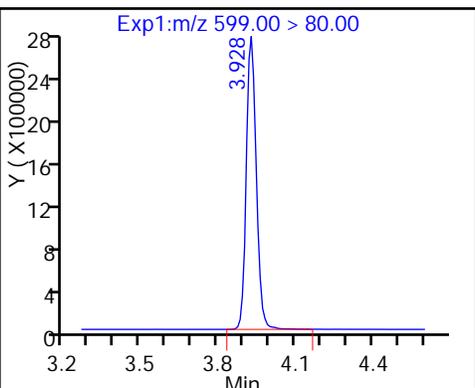
24 Perfluorodecanoic acid



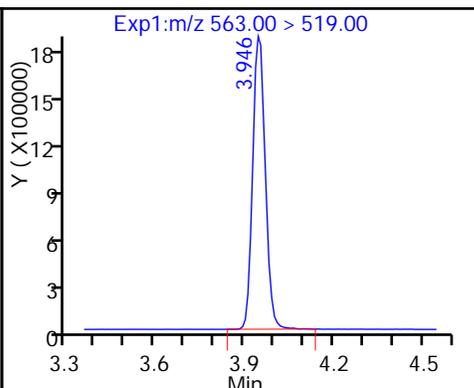
D 23 13C2 PFDA



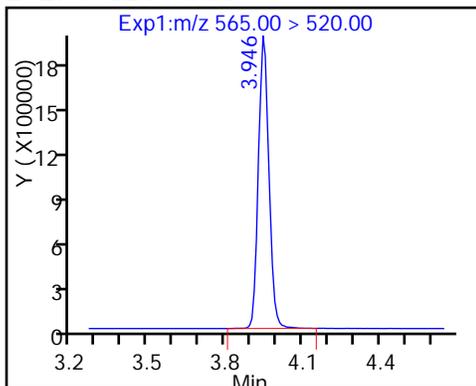
26 Perfluorodecane Sulfonic acid



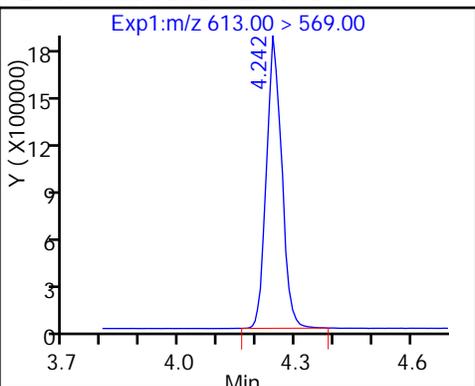
28 Perfluoroundecanoic acid



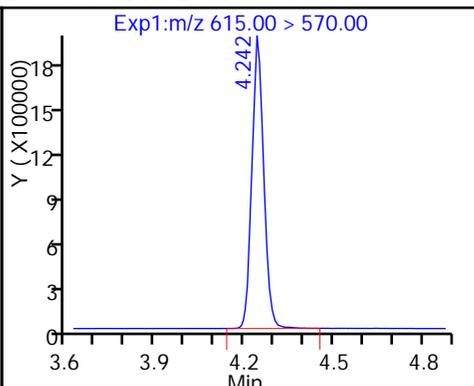
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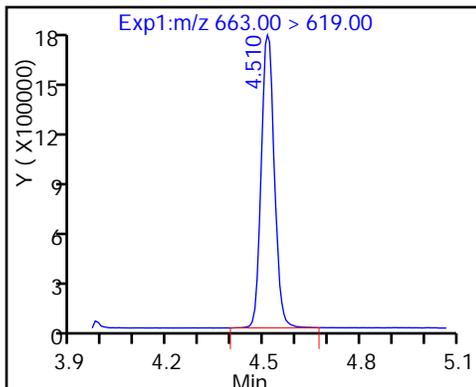
29 Perfluorododecanoic acid



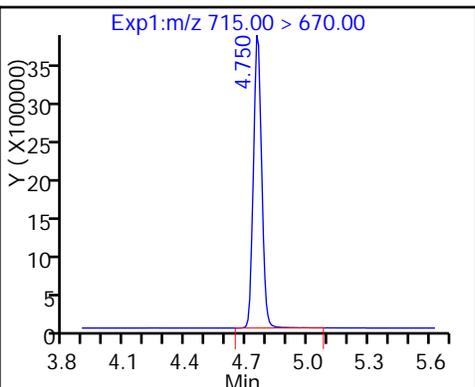
D 30 13C2 PFDaA



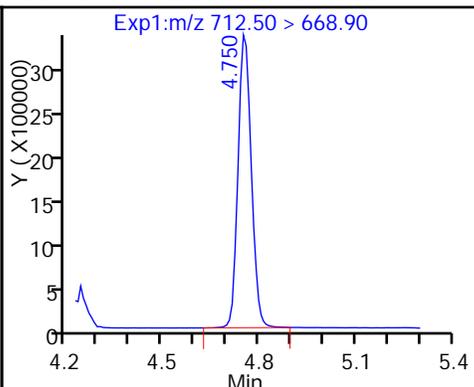
31 Perfluorotridecanoic acid



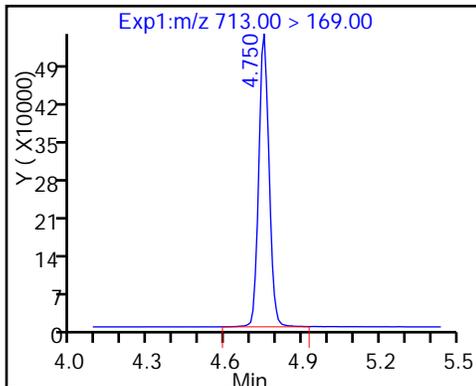
D 32 13C2-PFTeDA



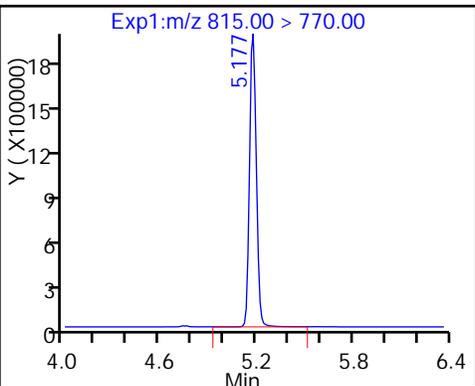
33 Perfluorotetradecanoic acid



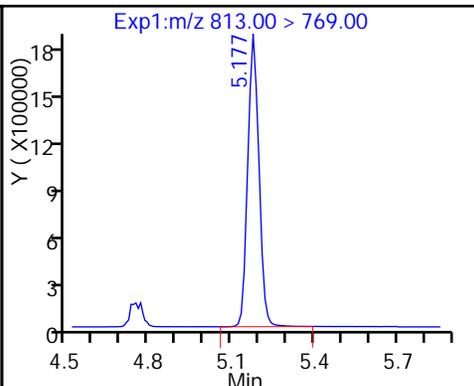
33 Perfluorotetradecanoic acid



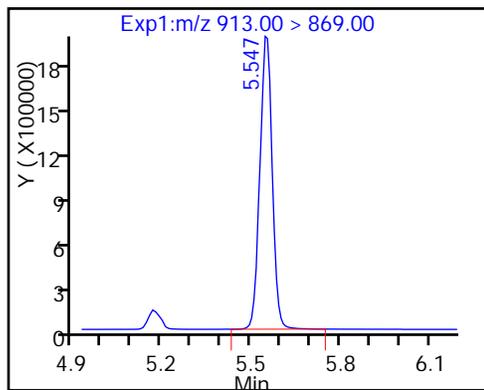
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_009.d
 Lims ID: IC L6
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 03-Dec-2016 14:26:13 ALS Bottle#: 42 Worklist Smp#: 9
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L6_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:24:17 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:44:26

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.565	1.574	-0.009	14589192	43.4		86.9	865034	
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.577	-0.003	1.000	42806432	167.9	83.9	195743	
3 Perfluoropentanoic acid	262.90 > 219.00	1.858	1.861	-0.003	1.000	35527936	156.8	78.4	261830	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.861	-0.003		11161582	42.2	84.4	966643	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.896	1.900	-0.004	1.000	53135654	128.0	72.4		
	298.90 > 99.00	1.896	1.900	-0.004	1.000	29815085	1.78(0.00-0.00)	72.4		
7 Perfluorohexanoic acid	313.00 > 269.00	2.157	2.164	-0.007	1.000	33117925	168.5	84.2	581515	
D 6 13C2 PFHxA	315.00 > 270.00	2.157	2.164	-0.007		10311260	43.4	86.8	509947	
D 11 13C4-PFHpA	367.00 > 322.00	2.501	2.511	-0.010		8082881	39.0	77.9	826093	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.501	2.512	-0.011	1.000	30013620	180.8	90.4	294610	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.526	2.531	-0.005	1.000	47221104	165.4	90.9		
D 10 18O2 PFHxS	403.00 > 84.00	2.518	2.531	-0.013		12302944	39.4	83.3	1330084	
D 14 13C4 PFOA	417.00 > 372.00	2.868	2.880	-0.012		8181000	37.3	74.6	486964	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.868	2.887	-0.019	1.000	30853962	175.9		88.0	291973	
413.00 > 169.00	2.876	2.887	-0.011	1.003	20504099		1.50(0.90-1.10)	88.0	540708	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.884	2.888	-0.004	1.000	39398056	159.1		83.6		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.249	3.258	-0.009	1.000	43077470	188.7		102	75165	
499.00 > 99.00	3.249	3.258	-0.009	1.000	9992099		4.31(0.90-1.10)	102	2648708	
D 17 13C4 PFOS										
503.00 > 80.00	3.249	3.259	-0.010		10053294	40.9		85.5	90966	
D 19 13C5 PFNA										
468.00 > 423.00	3.249	3.263	-0.014		6537549	39.3		78.6	482383	
20 Perfluorononanoic acid										
463.00 > 419.00	3.249	3.263	-0.014	1.000	24899337	191.1		95.6	358972	
D 21 13C8 FOSA										
506.00 > 78.00	3.562	3.571	-0.009		16949269	42.1		84.3	560152	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.571	3.574	-0.003	1.000	50658186	160.0		80.0	335690	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.613	3.623	-0.010	1.000	23571490	191.4		95.7	488396	
D 23 13C2 PFDA										
515.00 > 470.00	3.621	3.626	-0.005		6410340	40.6		81.2	207594	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.921	3.936	-0.015	1.000	24987612	185.7		96.3		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.947	3.955	-0.008	1.000	17818228	185.2		92.6	329189	
D 27 13C2 PFUnA										
565.00 > 520.00	3.947	3.958	-0.011		4513820	38.0		76.0	155727	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.236	4.250	-0.014	1.000	17773370	194.4		97.2	287464	
D 30 13C2 PFDoA										
615.00 > 570.00	4.236	4.251	-0.015		4817286	43.0		86.0	195663	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.501	4.518	-0.017	1.000	17494025	191.2		95.6	67382	M
D 32 13C2-PFTeDA										
715.00 > 670.00	4.738	4.759	-0.021		9518260	41.2		82.3	367157	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.746	4.761	-0.015	1.000	30593527	171.3		85.7	17849	
713.00 > 169.00	4.746	4.761	-0.015	1.000	5666745		5.40(0.00-0.00)	85.7	258419	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.171	5.186	-0.015		5482429	42.3		84.5	117464	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.160	5.186	-0.026	1.000	18912385	196.9		98.5	16230	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.536	5.559	-0.023	1.000	17844066	186.5		93.3	33443	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L6_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_009.d

Injection Date: 03-Dec-2016 14:26:13

Instrument ID: A8_N

Lims ID: IC L6

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 42

Worklist Smp#: 9

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

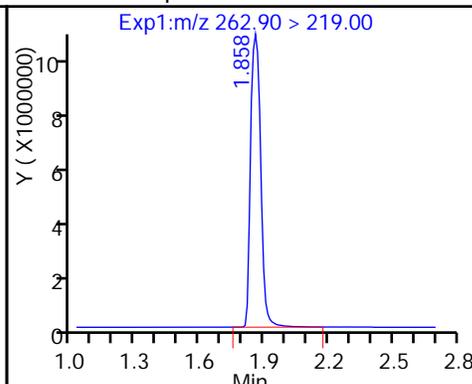
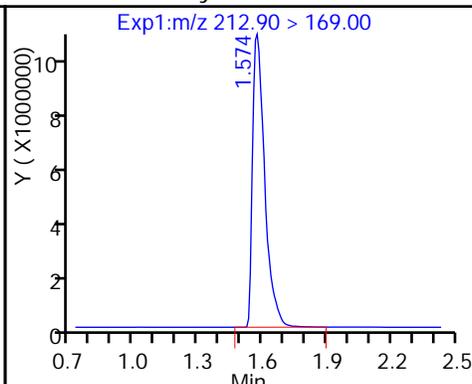
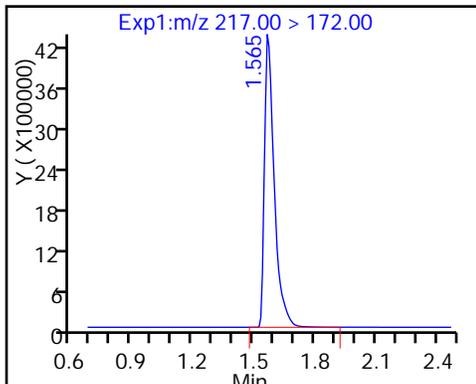
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

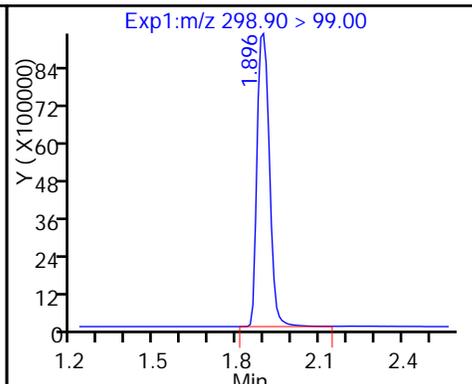
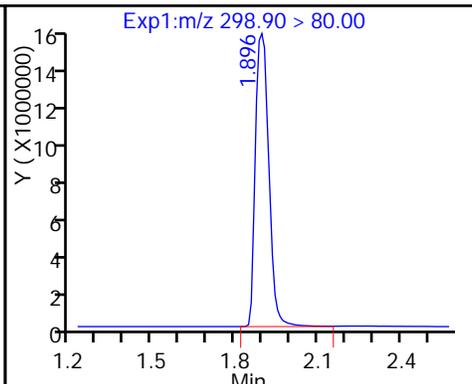
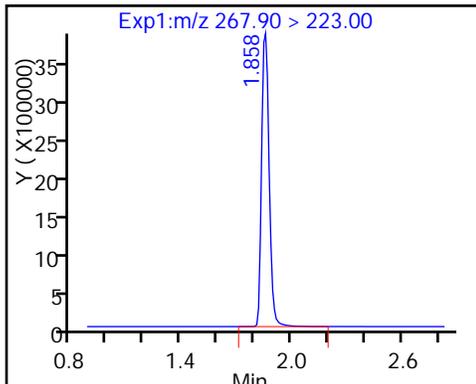
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

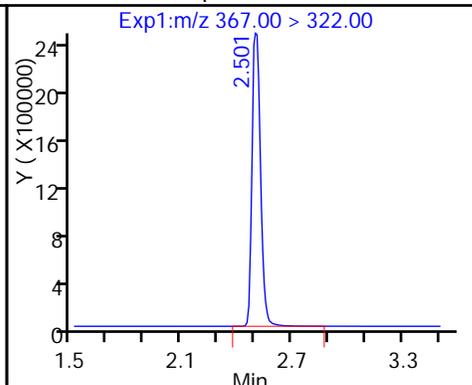
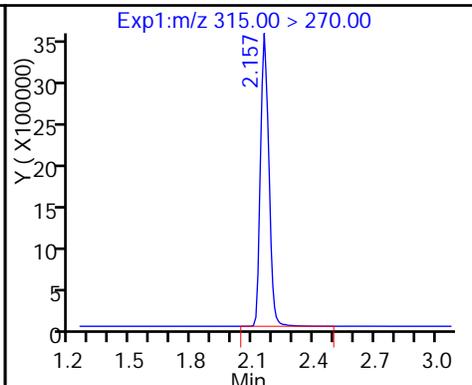
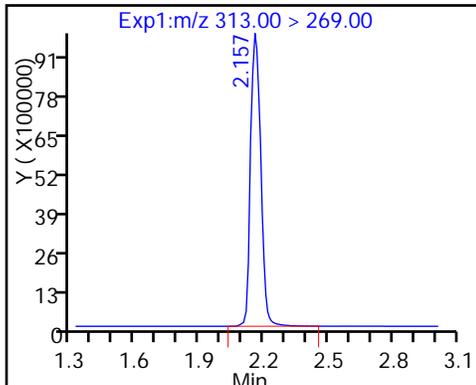
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

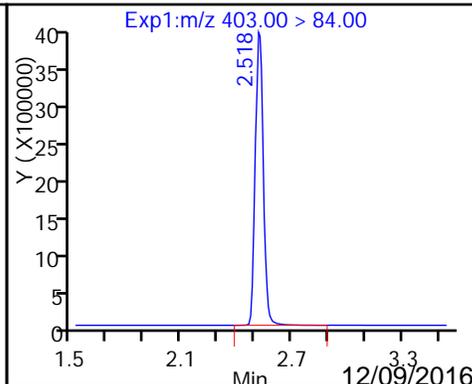
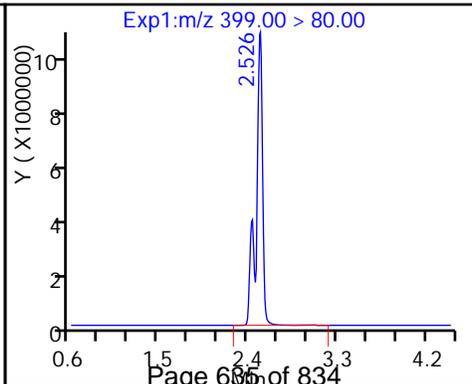
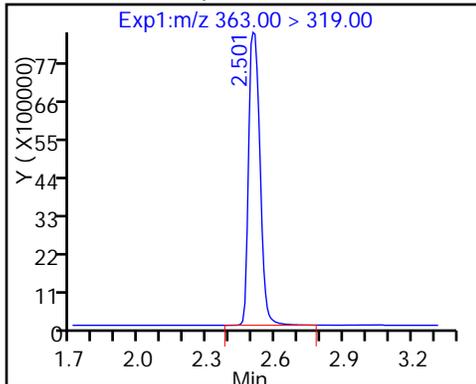
D 11 13C4-PFHpA



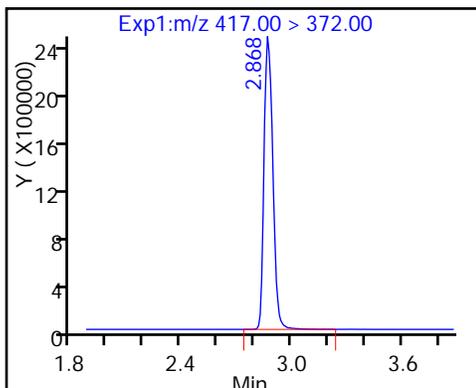
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

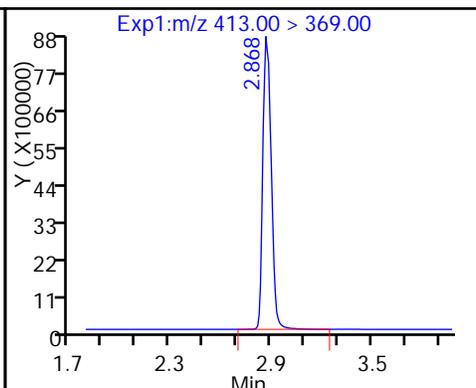
D 10 18O2 PFHxS



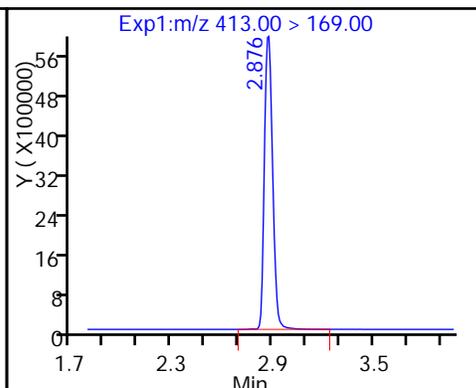
D 14 13C4 PFOA



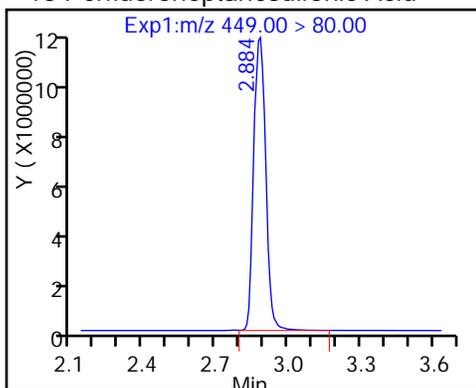
15 Perfluorooctanoic acid



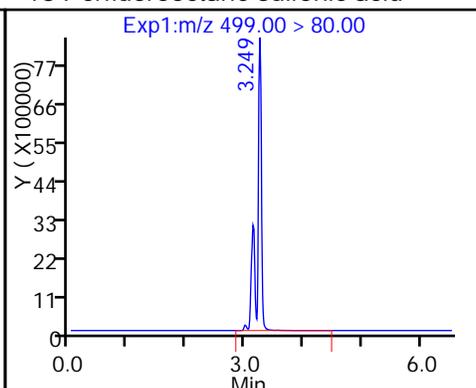
15 Perfluorooctanoic acid



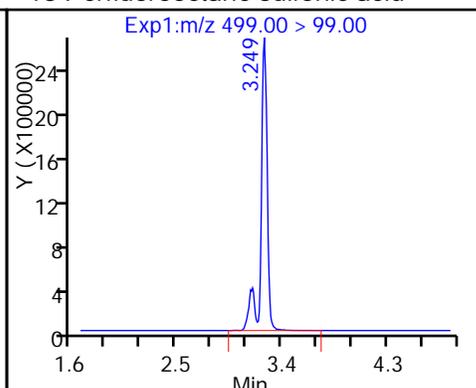
13 Perfluoroheptanesulfonic Acid



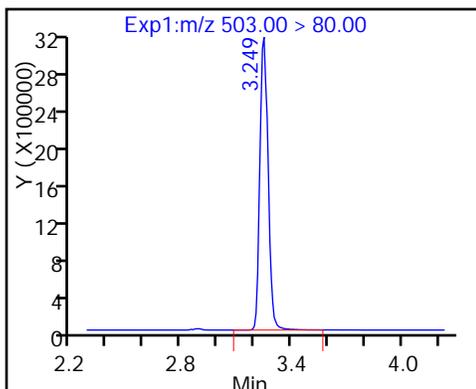
18 Perfluorooctane sulfonic acid



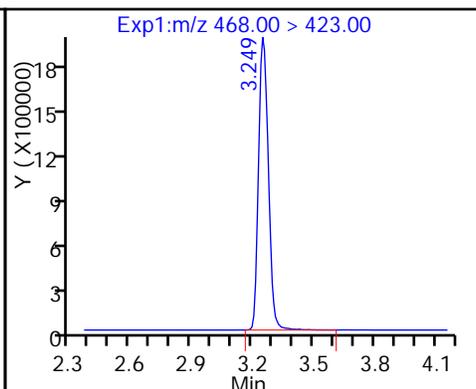
18 Perfluorooctane sulfonic acid



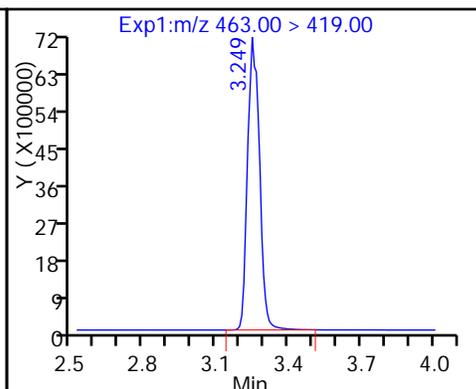
D 17 13C4 PFOS



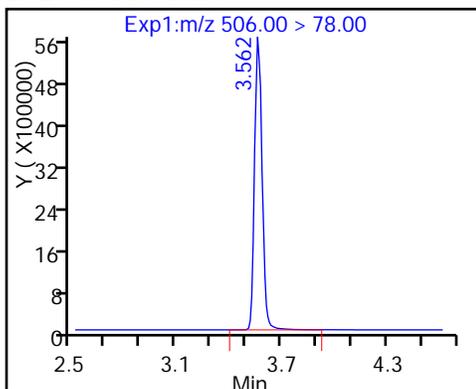
D 19 13C5 PFNA



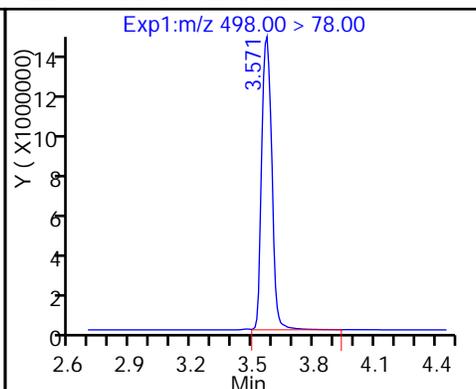
20 Perfluorononanoic acid



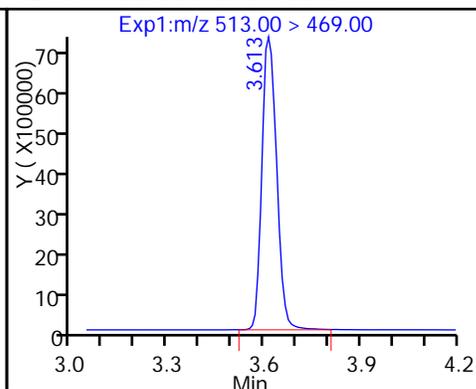
D 21 13C8 FOSA



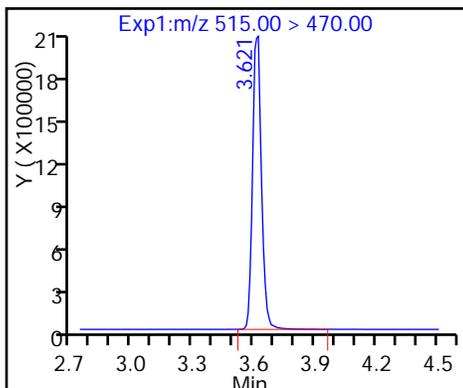
22 Perfluorooctane Sulfonamide



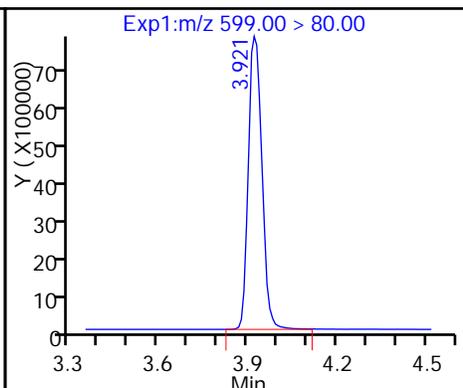
24 Perfluorodecanoic acid



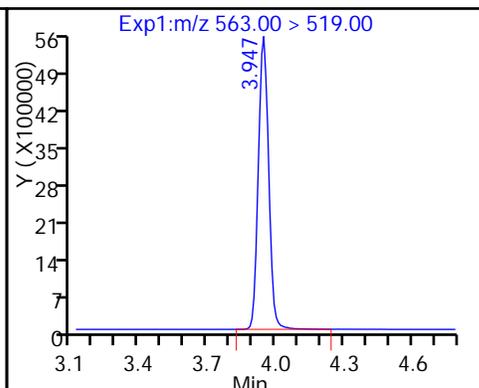
D 23 13C2 PFDA



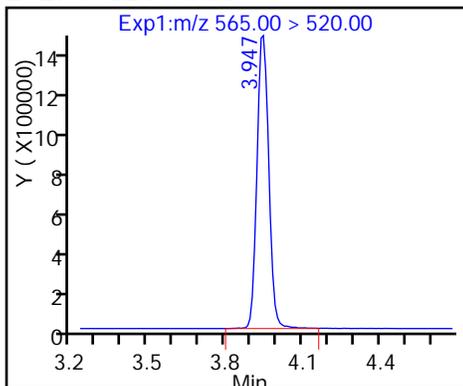
26 Perfluorodecane Sulfonic acid



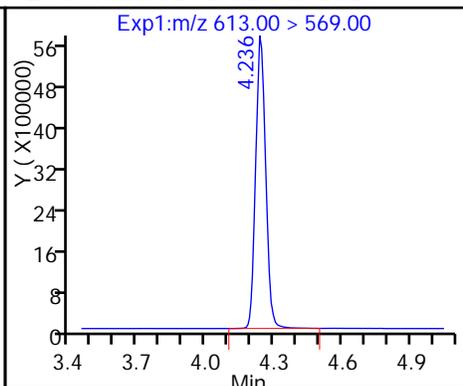
28 Perfluoroundecanoic acid



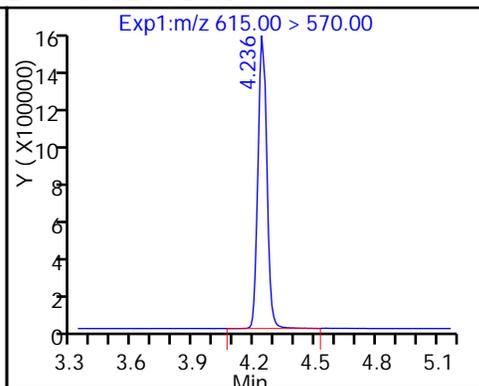
D 27 13C2 PFUa



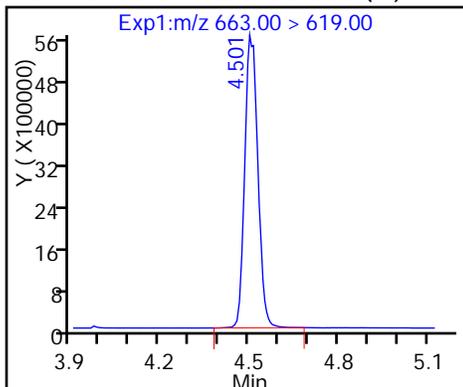
29 Perfluorododecanoic acid



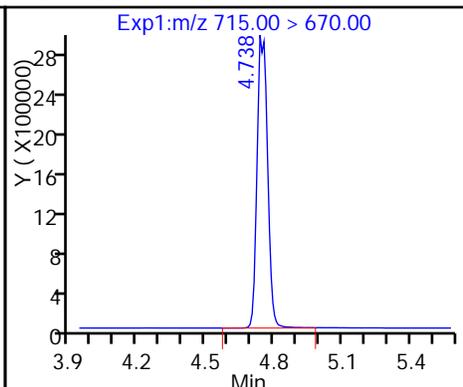
D 30 13C2 PFDa



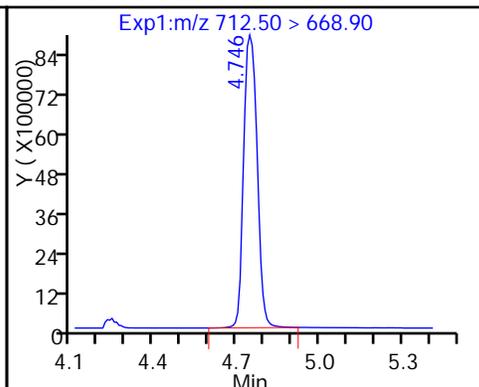
31 Perfluorotridecanoic acid (M)



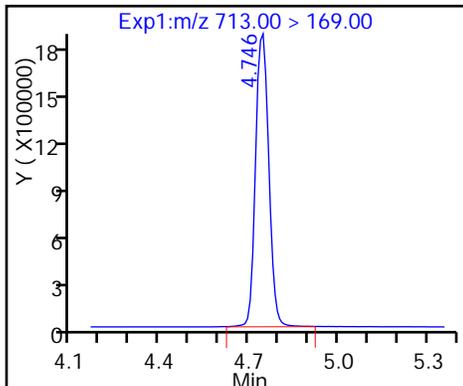
D 32 13C2-PFTeDA



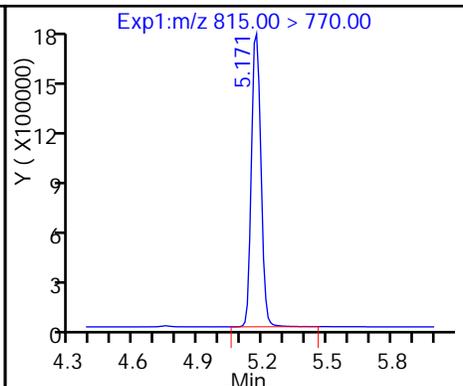
33 Perfluorotetradecanoic acid



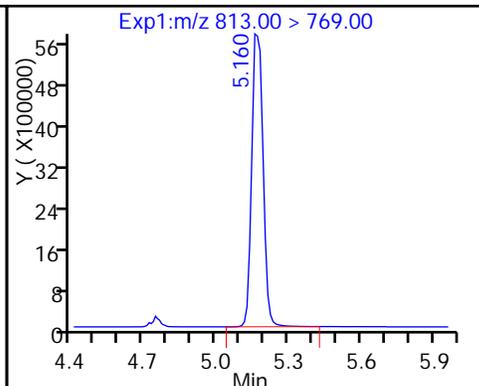
33 Perfluorotetradecanoic acid



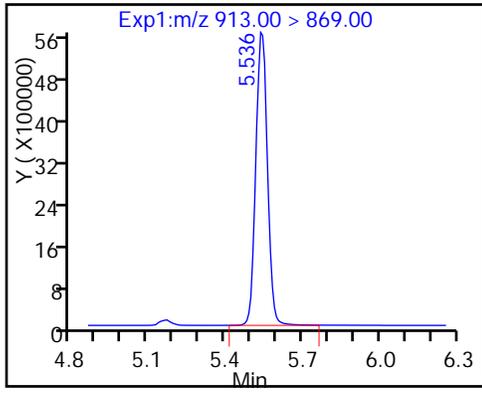
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



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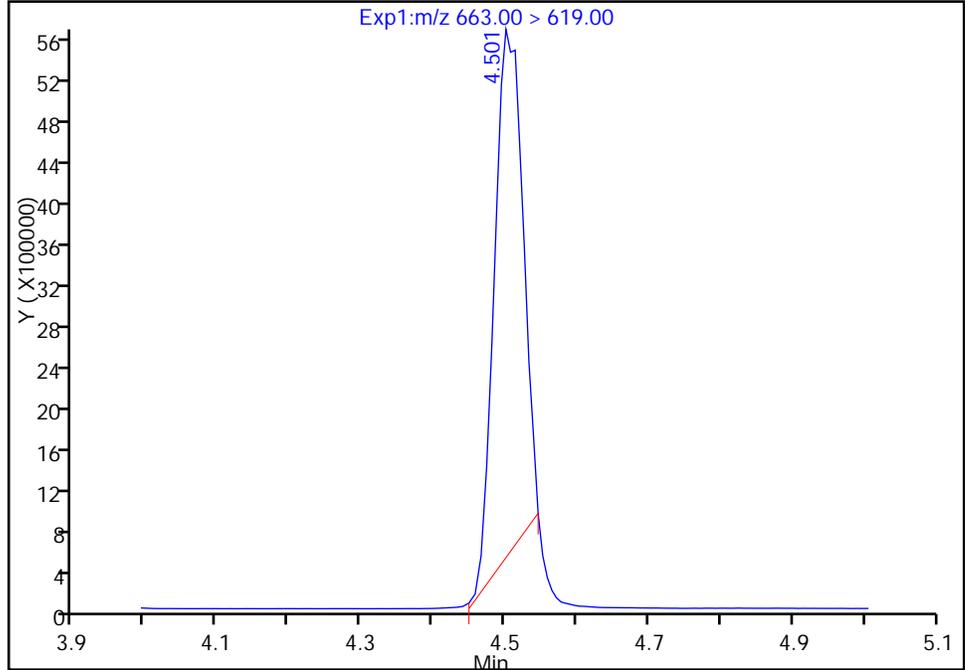
Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_009.d
Injection Date: 03-Dec-2016 14:26:13 Instrument ID: A8_N
Lims ID: IC L6
Client ID:
Operator ID: A8-PC\A8 ALS Bottle#: 42 Worklist Smp#: 9
Injection Vol: 2.0 ul Dil. Factor: 1.0000
Method: A8_N Limit Group: LC PFC_DOD ICAL
Column: Detector EXP1

31 Perfluorotridecanoic acid, CAS: 72629-94-8

Signal: 1

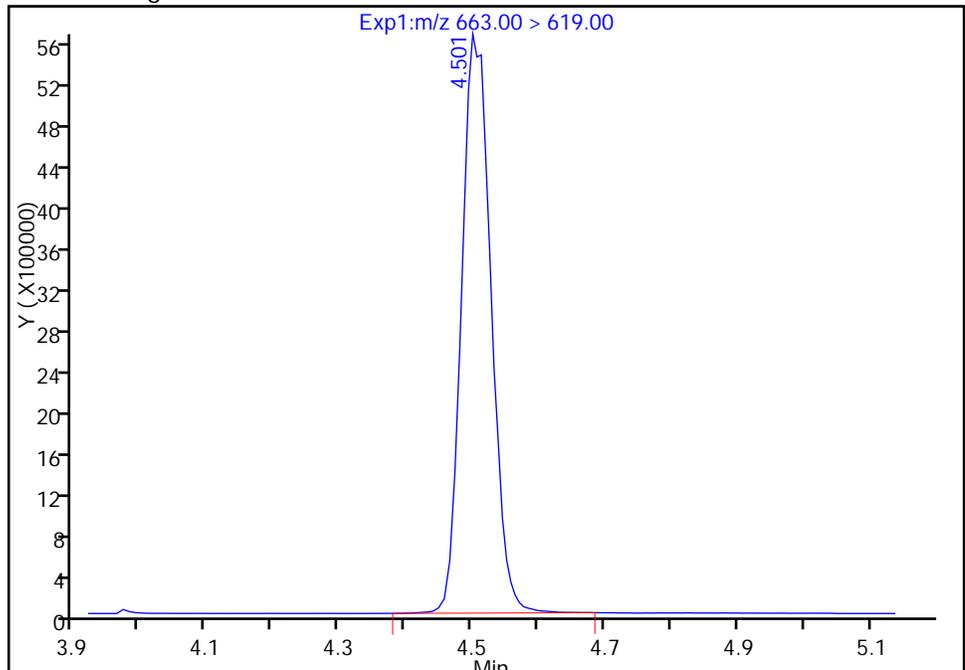
RT: 4.50
Area: 14077685
Amount: 158.7746
Amount Units: ng/ml

Processing Integration Results



RT: 4.50
Area: 17494025
Amount: 191.1674
Amount Units: ng/ml

Manual Integration Results



Reviewer: chandrasenas, 05-Dec-2016 09:44:26
Audit Action: Manually Integrated

Audit Reason: Incomplete Integration

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_013.d
 Lims ID: IC L1 Add-on
 Client ID:
 Sample Type: IC Calib Level: 1
 Inject. Date: 03-Dec-2016 14:56:13 ALS Bottle#: 46 Worklist Smp#: 13
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L1 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:00 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:47:53

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.840	2.836	0.004	1.000	40088	0.5222	110		
D 47 M2-6:2FTS	429.00 > 409.00	2.840	2.836	0.004		4340981	41.1	86.6		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.601	3.588	0.013	1.000	37551	0.4928	103		
D 42 M2-8:2FTS	529.00 > 509.00	3.601	3.591	0.010		4263521	43.9	91.6		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.765	3.756	0.009		3321758	45.8	91.6		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.774	3.759	0.015	1.002	23619	0.4142	82.8		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.931	3.922	0.009		3822164	48.0	96.0		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.931	3.929	0.002	1.000	25433	0.4345	86.9		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.054	4.051	0.003		4696243	44.7	89.4		
54 MeFOSA	512.00 > 169.00	4.063	4.057	0.006	1.000	37442	0.4934	98.7		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.244	4.239	0.005		4355916	43.9	87.9		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.253	4.245	0.008	1.000	36079	0.4813	96.3		

Reagents:

LCPFC2-L1_00002

Amount Added: 1.00

Units: mL

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_013.d

Injection Date: 03-Dec-2016 14:56:13

Instrument ID: A8_N

Lims ID: IC L1 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 46

Worklist Smp#: 13

Injection Vol: 2.0 ul

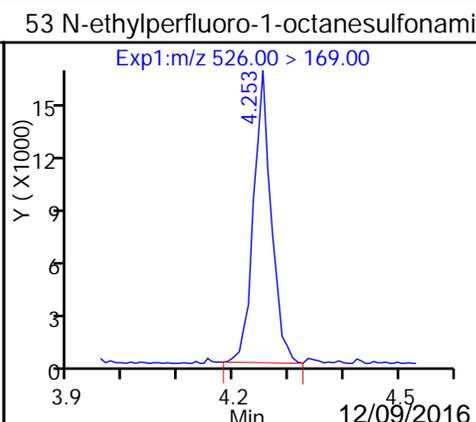
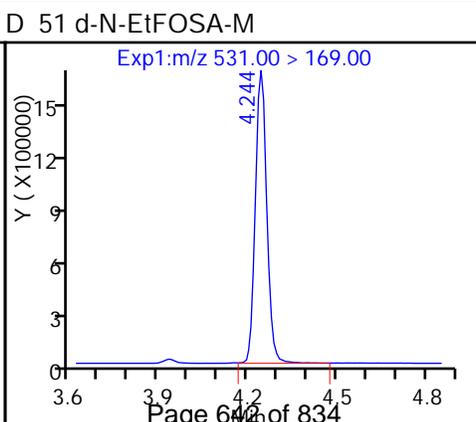
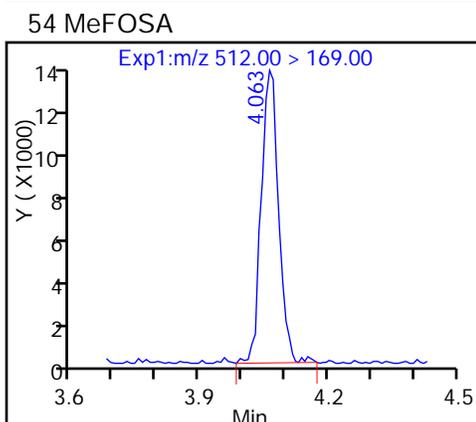
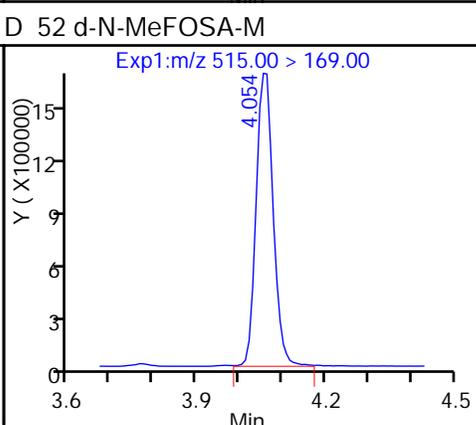
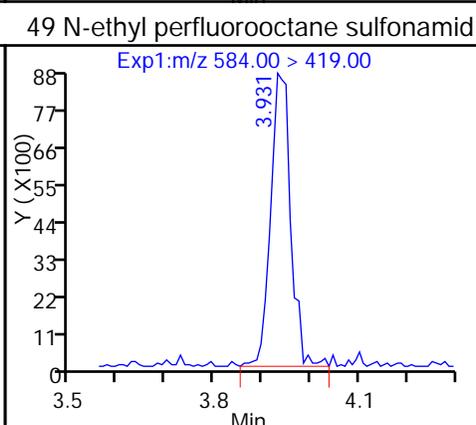
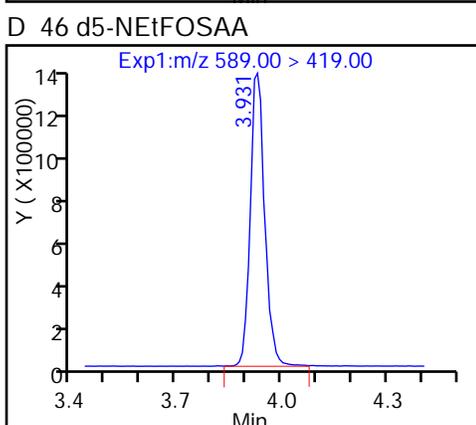
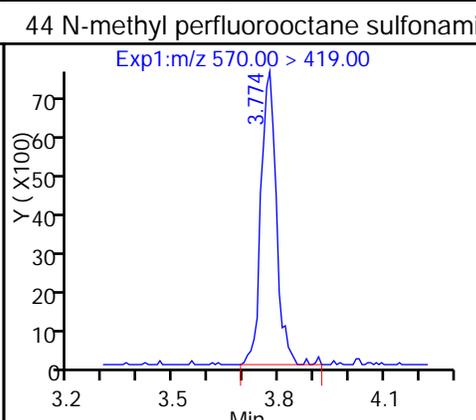
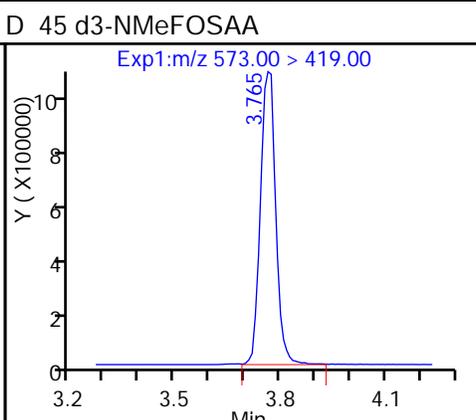
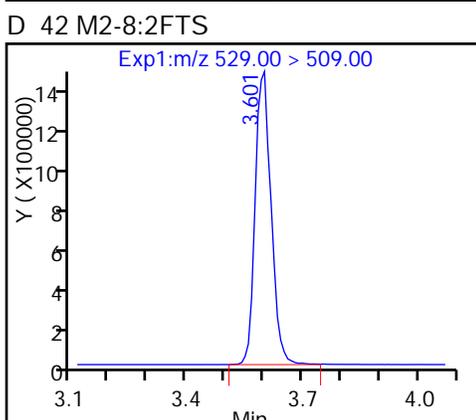
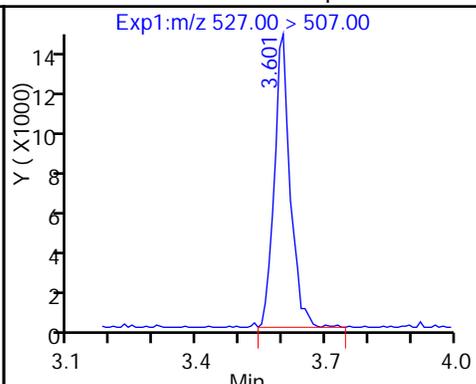
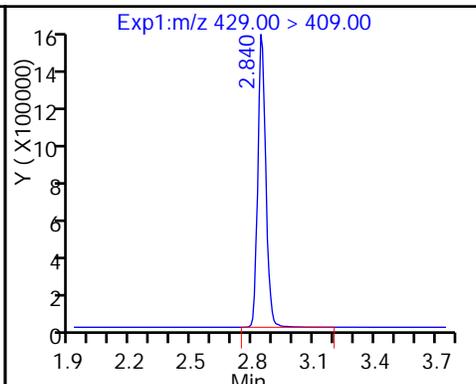
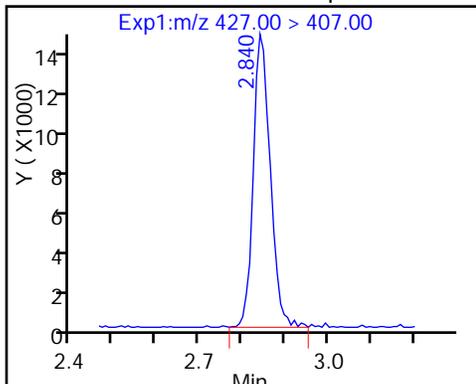
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_014.d
 Lims ID: IC L2 Add-on
 Client ID:
 Sample Type: IC Calib Level: 2
 Inject. Date: 03-Dec-2016 15:03:43 ALS Bottle#: 47 Worklist Smp#: 14
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L2 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:01 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:48:07

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.830	2.836	-0.006	1.000	69385	0.9534	101		
D 47 M2-6:2FTS	429.00 > 409.00	2.830	2.836	-0.006		4115043	39.0	82.1		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.579	3.588	-0.009	1.000	63932	0.9133	95.3		
D 42 M2-8:2FTS	529.00 > 509.00	3.579	3.591	-0.012		3917070	40.3	84.2		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.760	3.756	0.004		3368323	46.4	92.9		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.760	3.759	0.001	1.000	53845	0.9312	93.1		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.918	3.922	-0.004		3818745	47.9	95.9		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.935	3.929	0.006	1.004	49948	0.8541	85.4		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.051	4.051	0.0		4885054	46.5	93.0		
54 MeFOSA	512.00 > 169.00	4.051	4.057	-0.006	1.000	72155	0.9140	91.4		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.231	4.239	-0.008		4699997	47.4	94.8		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.239	4.245	-0.006	1.000	72107	0.8915	89.1		

Reagents:

LCPFC2-L2_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_014.d

Injection Date: 03-Dec-2016 15:03:43

Instrument ID: A8_N

Lims ID: IC L2 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 47

Worklist Smp#: 14

Injection Vol: 2.0 ul

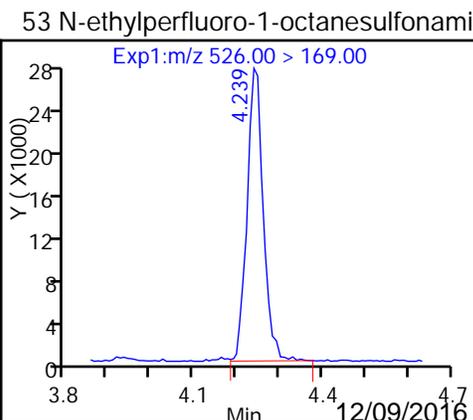
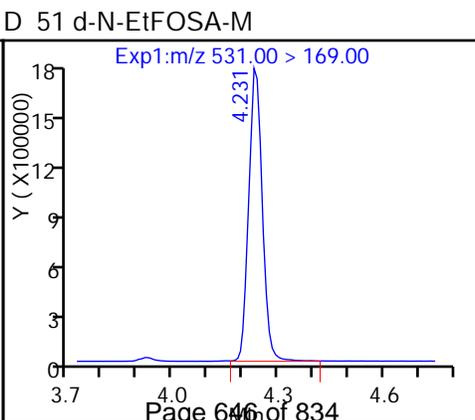
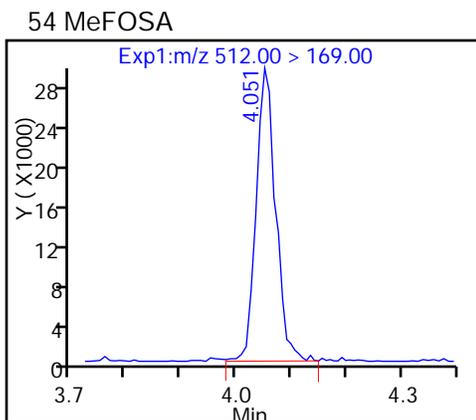
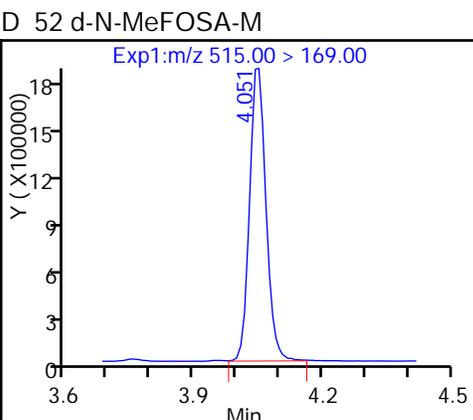
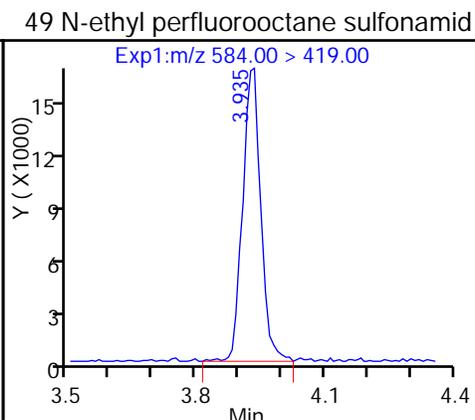
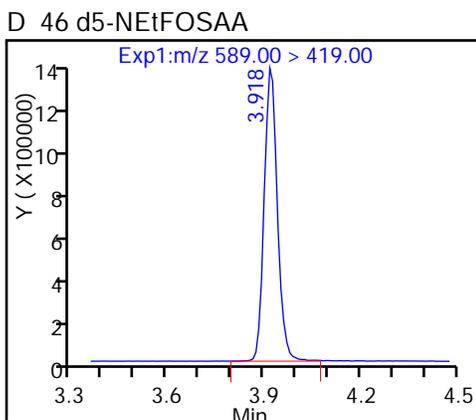
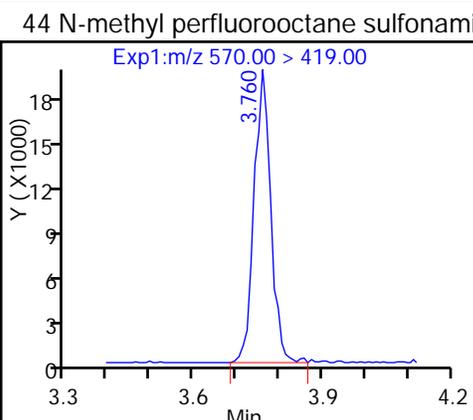
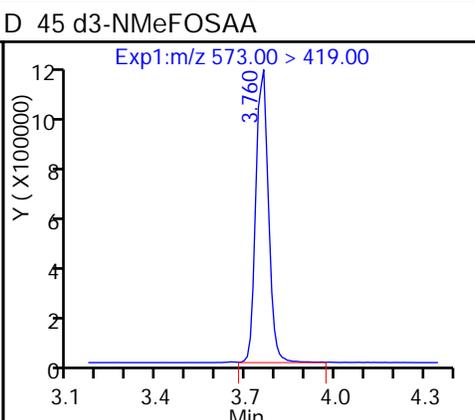
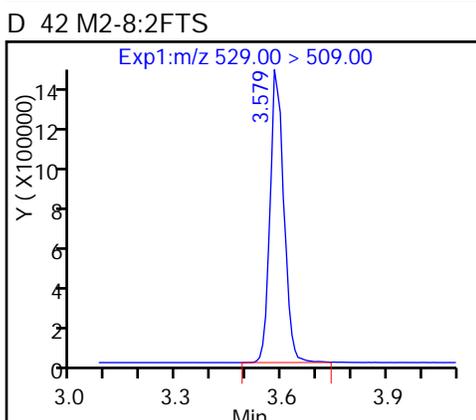
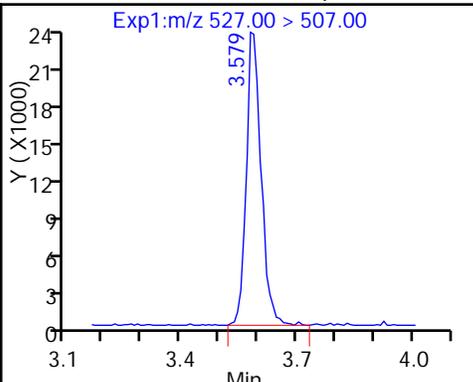
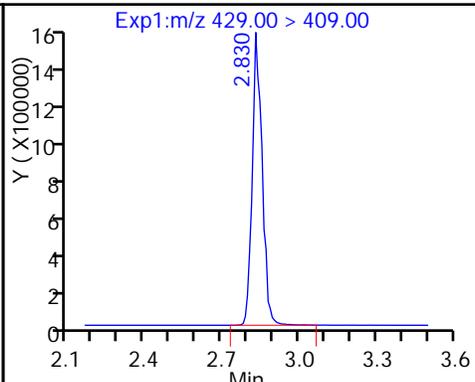
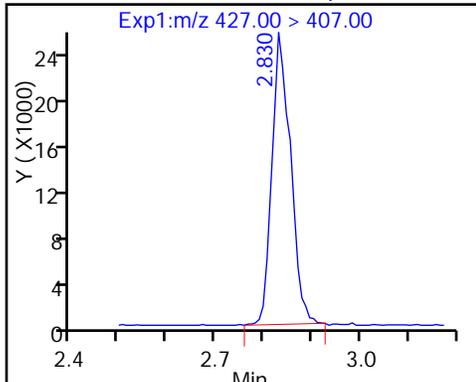
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_015.d
 Lims ID: IC L3 Add-on
 Client ID:
 Sample Type: IC Calib Level: 3
 Inject. Date: 03-Dec-2016 15:11:13 ALS Bottle#: 48 Worklist Smp#: 15
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L3 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:04 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:48:18

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.838	2.836	0.002	1.000	352721	4.15	87.6		
D 47 M2-6:2FTS	429.00 > 409.00	2.838	2.836	0.002		4800798	45.5	95.8		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.588	3.588	0.0	0.998	326213	4.10	85.6		
D 42 M2-8:2FTS	529.00 > 509.00	3.596	3.591	0.005		4450751	45.8	95.7		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.751	3.756	-0.005		3668263	50.6	101		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.760	3.759	0.001	1.002	260030	4.13	82.6		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.918	3.922	-0.004		4095032	51.4	103		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.926	3.929	-0.003	1.002	265257	4.23	84.6		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.051	4.051	0.0		5526561	52.6	105		
54 MeFOSA	512.00 > 169.00	4.051	4.057	-0.006	1.000	356898	4.00	79.9		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.240	4.239	0.001		5054297	51.0	102		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.240	4.245	-0.005	1.000	365313	4.20	84.0		

Reagents:

LCPFC2-L3_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_015.d

Injection Date: 03-Dec-2016 15:11:13

Instrument ID: A8_N

Lims ID: IC L3 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 48

Worklist Smp#: 15

Injection Vol: 2.0 ul

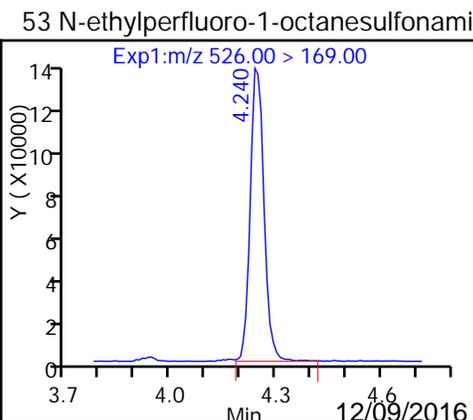
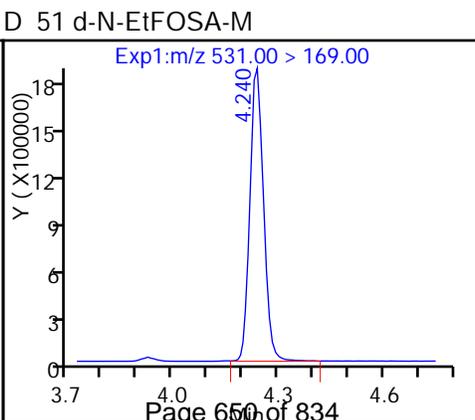
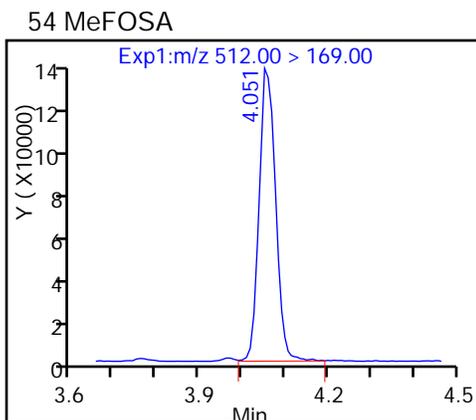
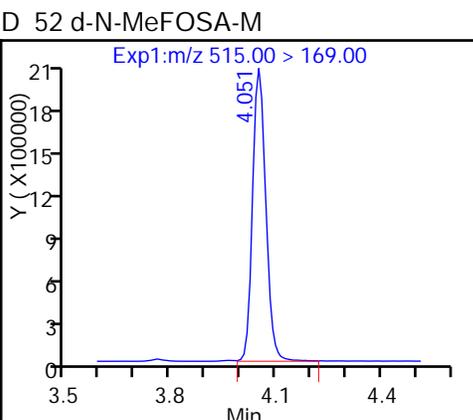
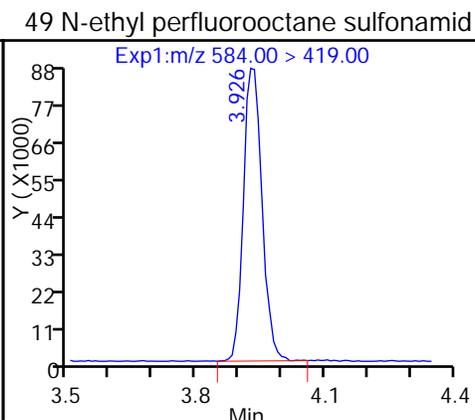
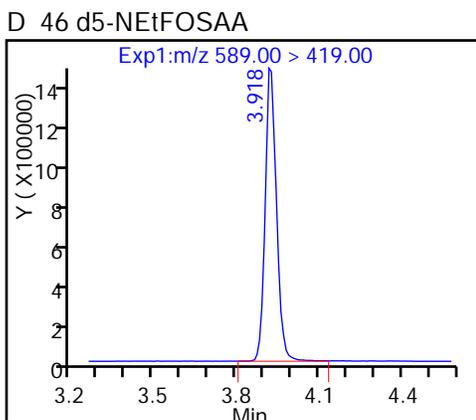
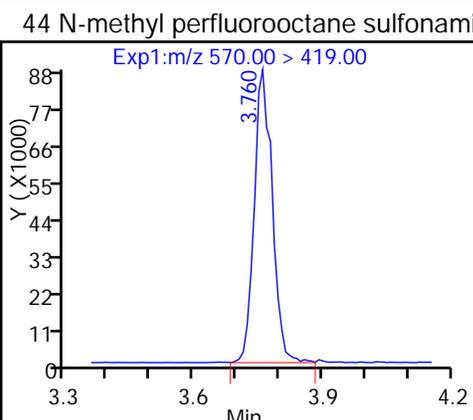
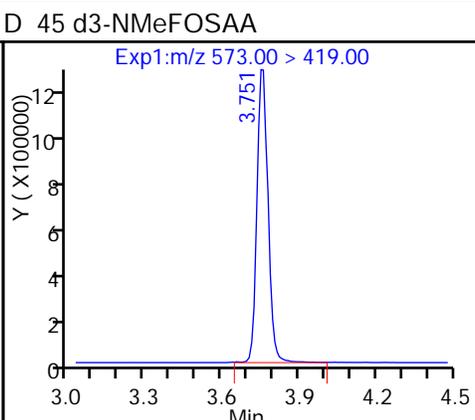
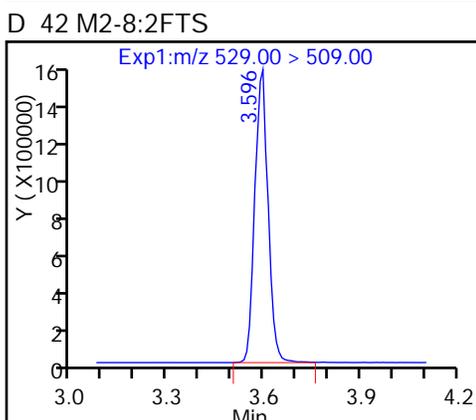
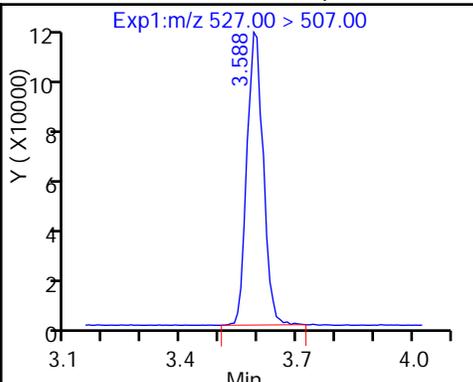
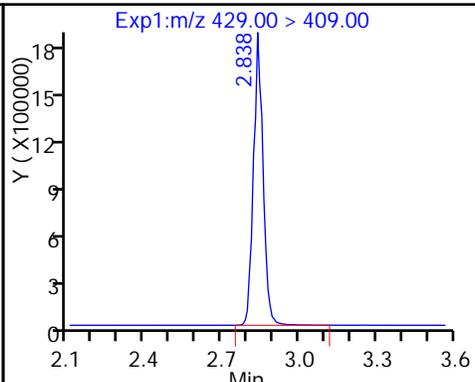
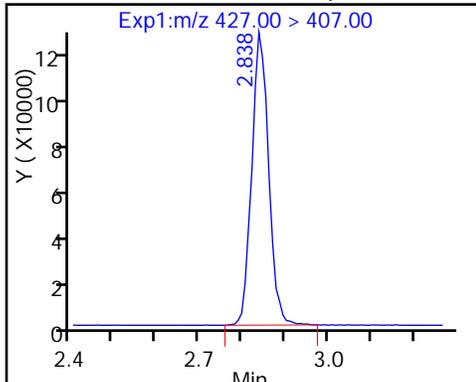
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_016.d
 Lims ID: IC L4 Add-on
 Client ID:
 Sample Type: IC Calib Level: 4
 Inject. Date: 03-Dec-2016 15:18:43 ALS Bottle#: 49 Worklist Smp#: 16
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L4 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:06 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:47:40

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.840	2.836	0.004	1.000	2169230	19.5	103		
D 47 M2-6:2FTS	429.00 > 409.00	2.840	2.836	0.004		6295656	59.7	126		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.592	3.588	0.004	1.000	2029390	21.1	110		
D 42 M2-8:2FTS	529.00 > 509.00	3.592	3.591	0.001		5378444	55.4	116		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.756	3.756	0.0		4323749	59.6	119		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.756	3.759	-0.003	1.000	1655435	22.3	112		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.931	3.922	0.009		4650188	58.4	117		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.931	3.929	0.002	1.000	1629443	22.9	114		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.055	4.051	0.004		6162693	58.6	117		
54 MeFOSA	512.00 > 169.00	4.063	4.057	0.006	1.000	2189848	22.0	110		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.244	4.239	0.005		5747792	58.0	116		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.244	4.245	-0.001	1.000	2201423	22.3	111		

Reagents:

LCPFC2-L4_00003

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_016.d

Injection Date: 03-Dec-2016 15:18:43

Instrument ID: A8_N

Lims ID: IC L4 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 49

Worklist Smp#: 16

Injection Vol: 2.0 ul

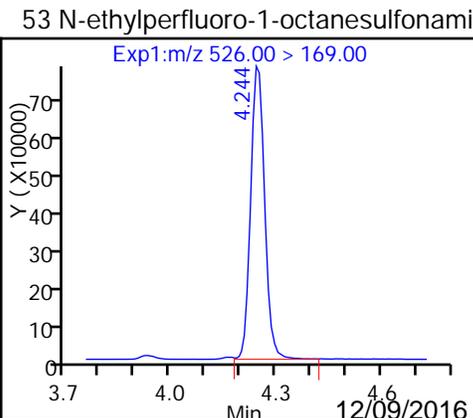
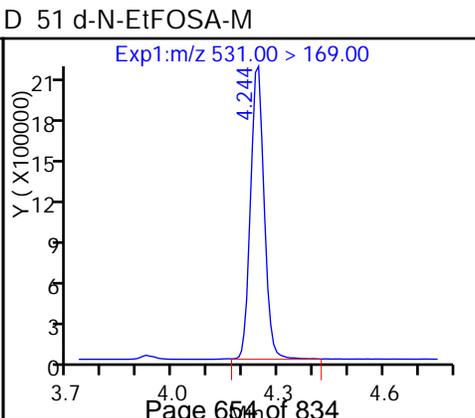
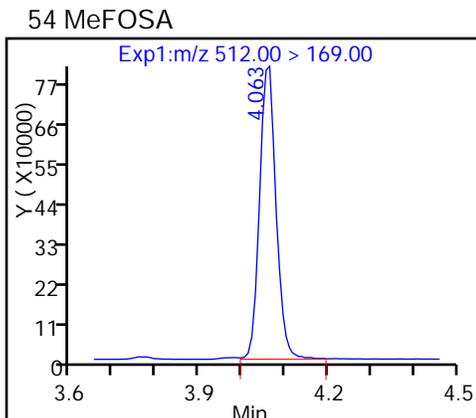
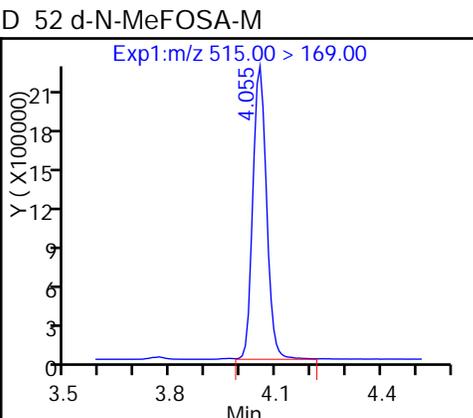
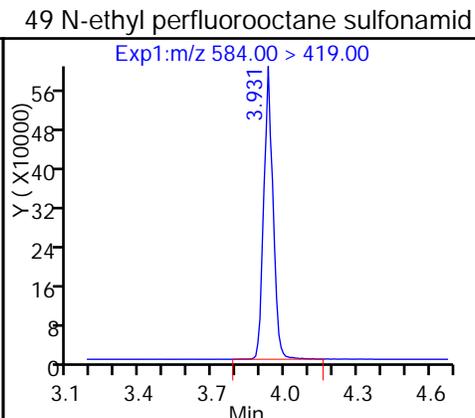
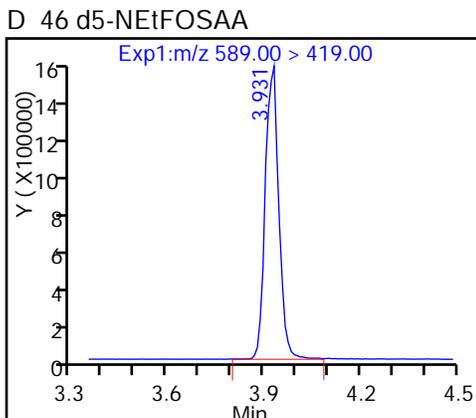
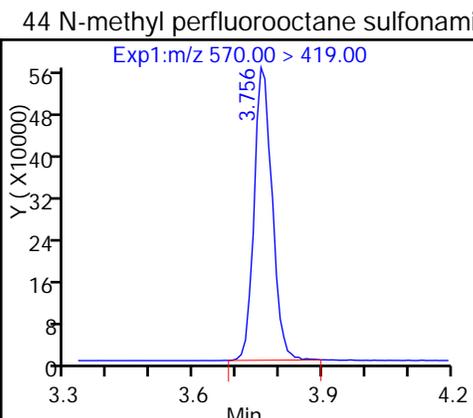
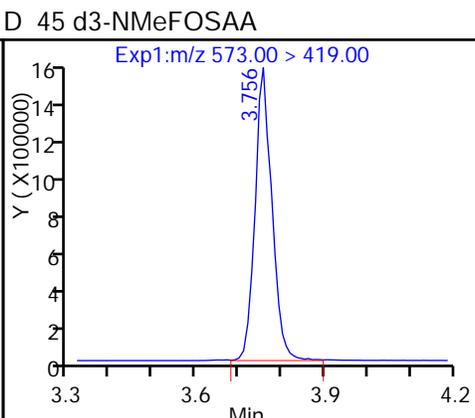
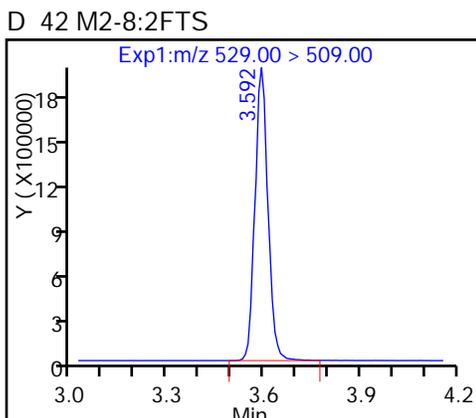
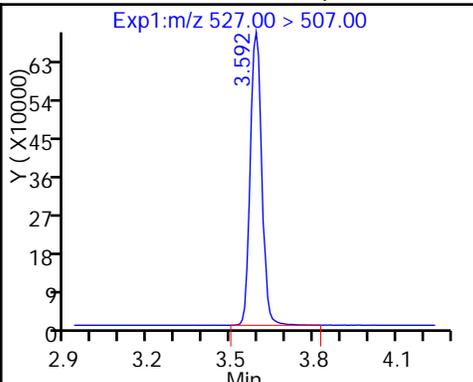
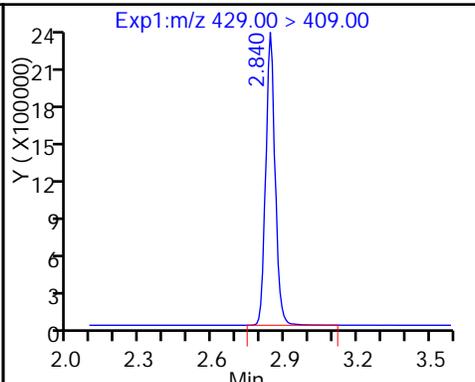
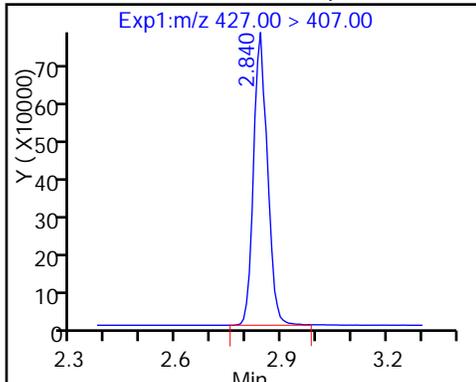
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_017.d
 Lims ID: IC L5 Add-on
 Client ID:
 Sample Type: IC Calib Level: 5
 Inject. Date: 03-Dec-2016 15:26:12 ALS Bottle#: 50 Worklist Smp#: 17
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L5 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:07 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:48:43

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.830	2.836	-0.006	1.000	4576863	55.8	118		
D 47 M2-6:2FTS	429.00 > 409.00	2.830	2.836	-0.006		4638709	44.0	92.5		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.579	3.588	-0.009	0.998	4279048	57.8	121		
D 42 M2-8:2FTS	529.00 > 509.00	3.587	3.591	-0.004		4144623	42.7	89.1		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.750	3.756	-0.006		3122973	43.1	86.1		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.750	3.759	-0.009	1.000	3484088	65.0	130		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.917	3.922	-0.005		3419439	42.9	85.9		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.925	3.929	-0.004	1.002	3261537	62.3	125		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.042	4.051	-0.009		4794382	45.6	91.2		
54 MeFOSA	512.00 > 169.00	4.051	4.057	-0.006	1.000	4656344	60.1	120		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.236	4.239	-0.003		4667809	47.1	94.1		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.245	4.245	0.0	1.000	4735360	58.9	118		

Reagents:

LCPFC2-L5_00002

Amount Added: 1.00

Units: mL

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_017.d

Injection Date: 03-Dec-2016 15:26:12

Instrument ID: A8_N

Lims ID: IC L5 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 50

Worklist Smp#: 17

Injection Vol: 2.0 ul

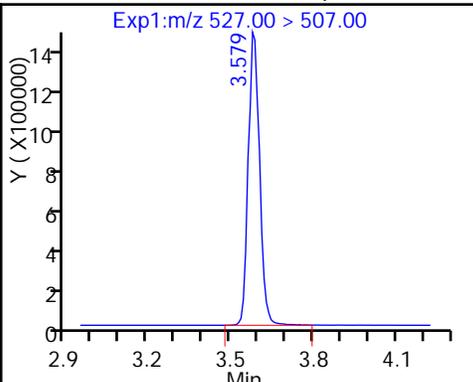
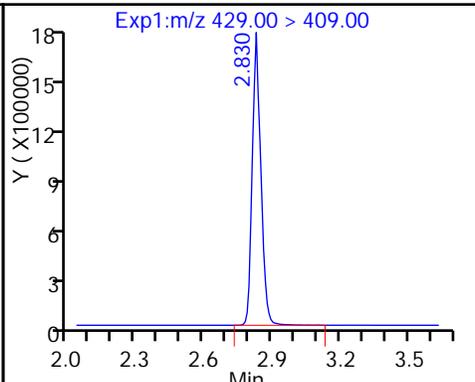
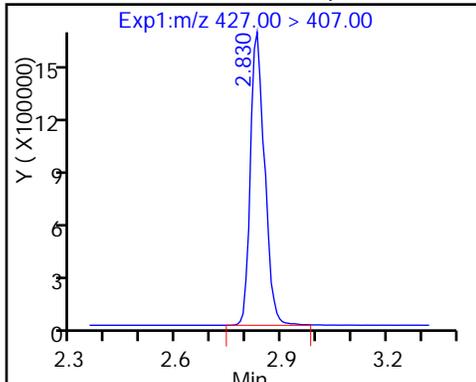
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane D 47 M2-6:2FTS

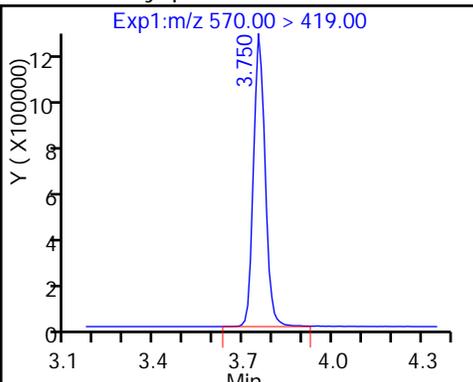
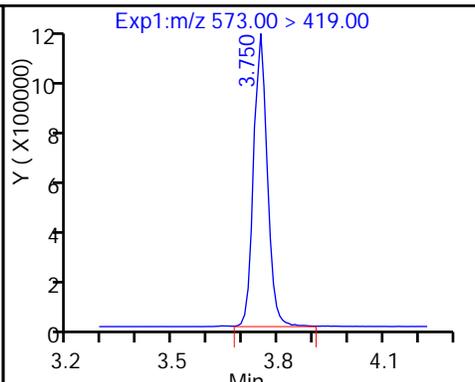
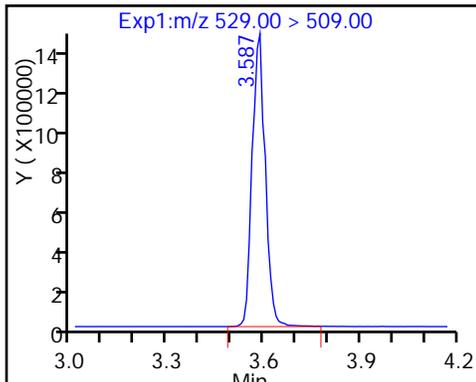
43 Sodium 1H,1H,2H,2H-perfluorooctane



D 42 M2-8:2FTS

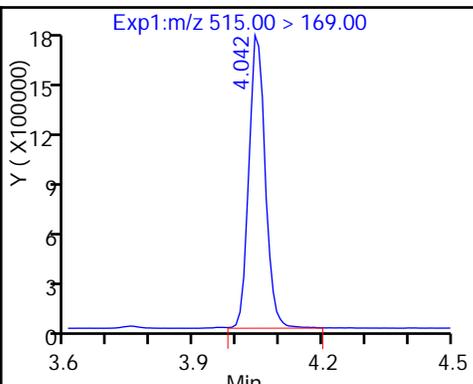
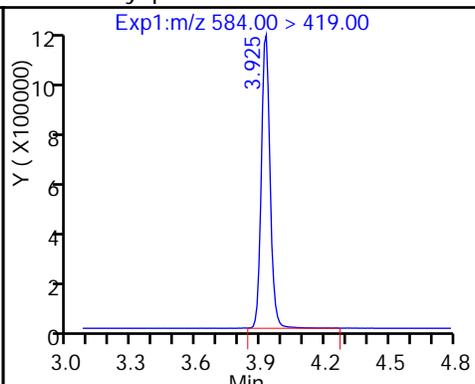
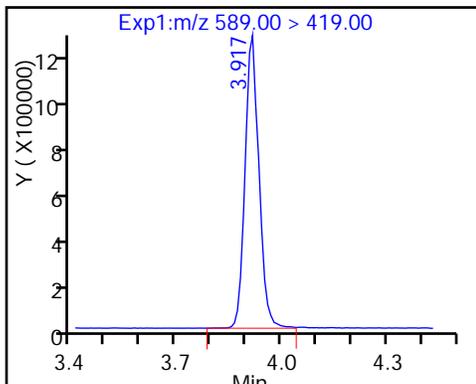
D 45 d3-NMeFOSAA

44 N-methyl perfluorooctane sulfonami



D 46 d5-NEtFOSAA

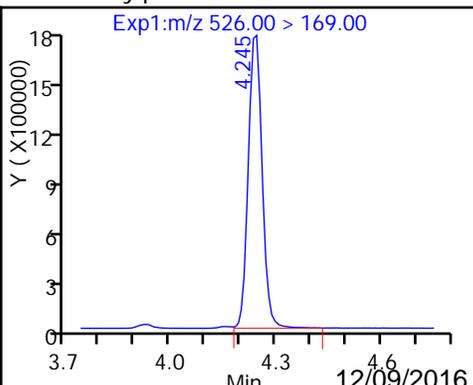
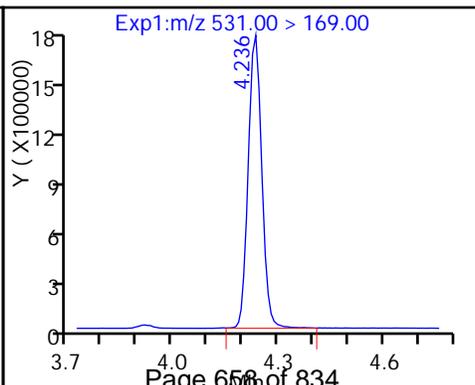
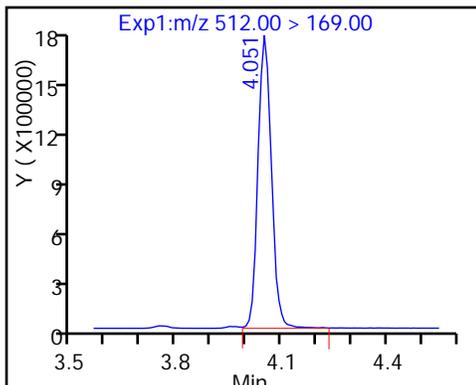
49 N-ethyl perfluorooctane sulfonamid D 52 d-N-MeFOSA-M



54 MeFOSA

D 51 d-N-EtFOSA-M

53 N-ethylperfluoro-1-octanesulfonami



TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Lims ID: IC L6 Add-on
 Client ID:
 Sample Type: IC Calib Level: 6
 Inject. Date: 03-Dec-2016 15:33:40 ALS Bottle#: 51 Worklist Smp#: 18
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: L6 ADD ON
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub3
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:28:09 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:49:51

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.838	2.836	0.002	1.000	18026595	173.2	91.3		
D 47 M2-6:2FTS	429.00 > 409.00	2.838	2.836	0.002		5885296	55.8	117		
43 Sodium 1H,1H,2H,2H-perfluorooctane	527.00 > 507.00	3.589	3.588	0.001	1.000	16850029	163.6	85.4		
D 42 M2-8:2FTS	529.00 > 509.00	3.589	3.591	-0.002		5764339	59.3	124		
D 45 d3-NMeFOSAA	573.00 > 419.00	3.752	3.756	-0.004		3953776	54.5	109		
44 N-methyl perfluorooctane sulfonami	570.00 > 419.00	3.752	3.759	-0.007	1.000	13566790	199.9	99.9		
D 46 d5-NEtFOSAA	589.00 > 419.00	3.919	3.922	-0.003		4091711	51.4	103		
49 N-ethyl perfluorooctane sulfonamid	584.00 > 419.00	3.927	3.929	-0.002	1.002	13047697	208.2	104		
D 52 d-N-MeFOSA-M	515.00 > 169.00	4.052	4.051	0.001		5465444	52.0	104		
54 MeFOSA	512.00 > 169.00	4.061	4.057	0.004	1.000	17635948	199.7	99.8		
D 51 d-N-EtFOSA-M	531.00 > 169.00	4.239	4.239	0.0		5223764	52.7	105		
53 N-ethylperfluoro-1-octanesulfonami	526.00 > 169.00	4.249	4.245	0.003	1.000	18234509	202.8	101		

Reagents:

LCPFC2-L6_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Injection Date: 03-Dec-2016 15:33:40

Instrument ID: A8_N

Lims ID: IC L6 Add-on

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 51

Worklist Smp#: 18

Injection Vol: 2.0 ul

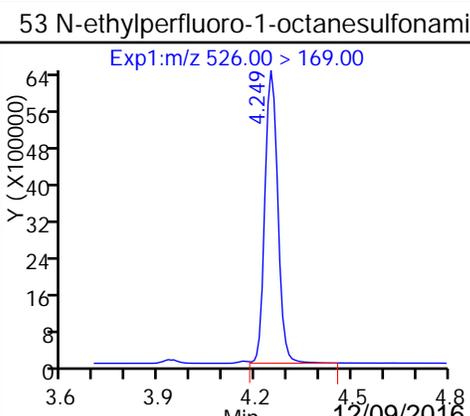
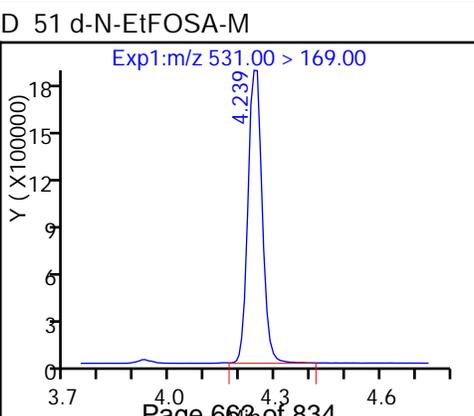
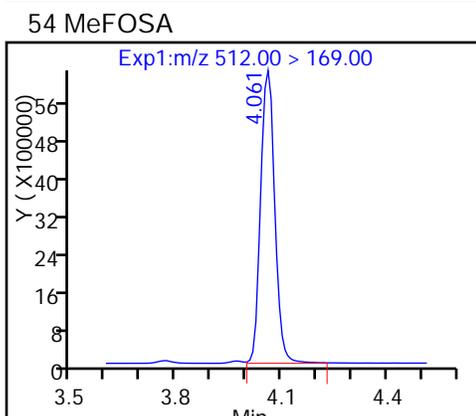
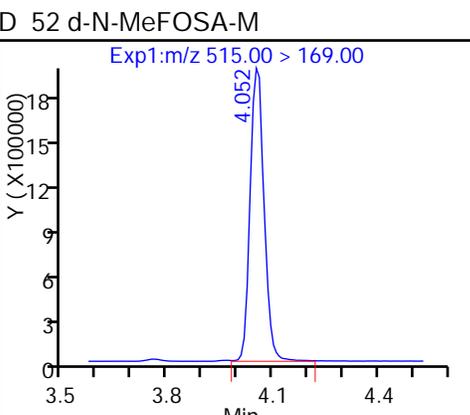
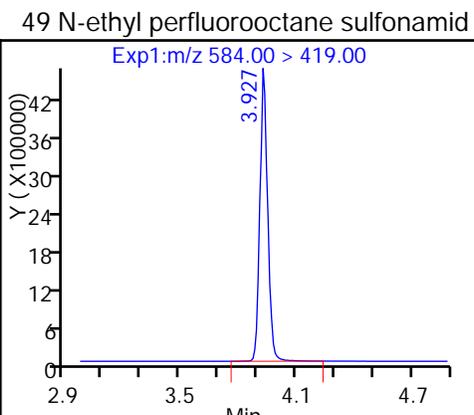
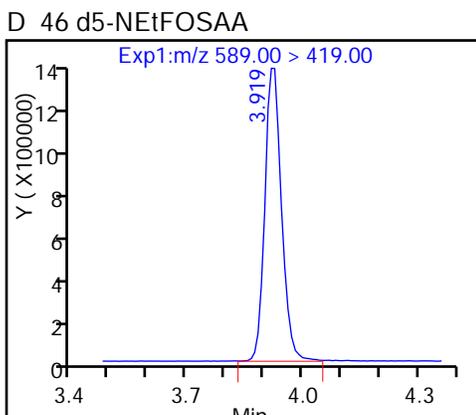
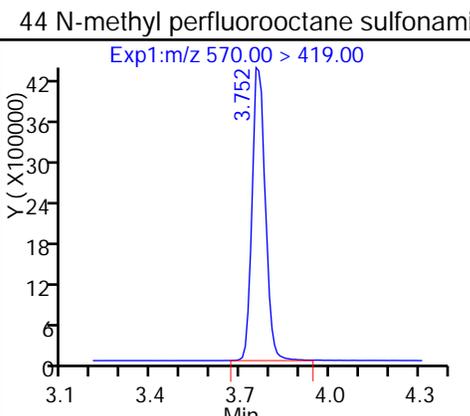
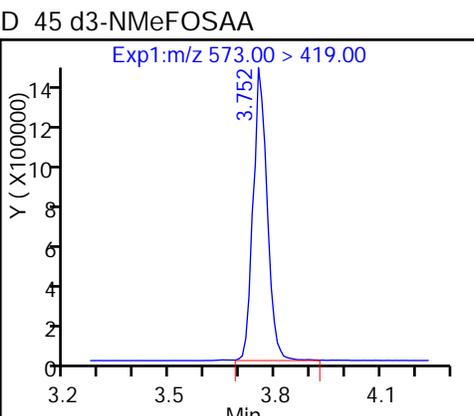
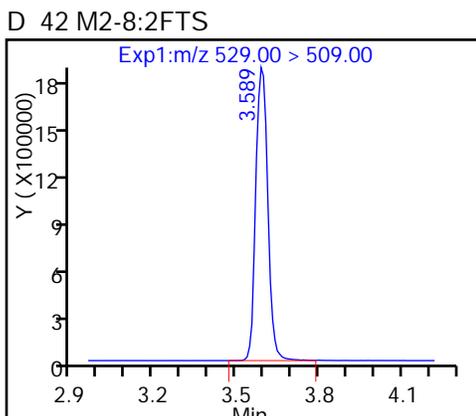
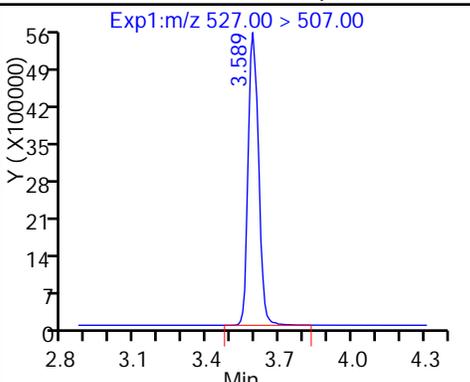
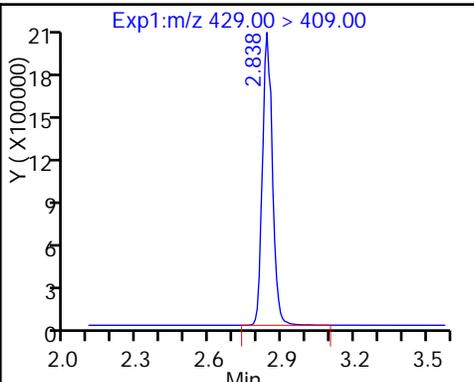
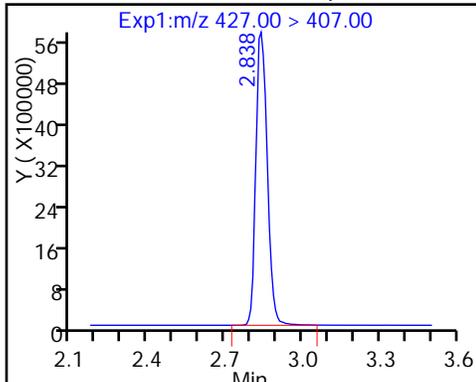
Dil. Factor: 1.0000

Method: A8_N

Limit Group: LC PFC_DOD ICAL

48 Sodium 1H,1H,2H,2H-perfluorooctane

43 Sodium 1H,1H,2H,2H-perfluorooctane



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: ICV 320-140382/12 Calibration Date: 12/02/2016 11:29
 Instrument ID: A8_N Calib Start Date: 12/02/2016 10:29
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/02/2016 12:29
 Lab File ID: 02DEC2016A_012.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.613	1.563		42.9	44.3	-3.1	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.020	0.9852		48.3	50.0	-3.4	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.087	0.9062		39.8	47.8	-16.7	25.0
13C4 PFBA	Ave	340300	315108		46.3	50.0	-7.4	50.0
13C5-PFPeA	Ave	270194	250588		46.4	50.0	-7.3	50.0
13C2 PFHxA	Ave	246480	226803		46.0	50.0	-8.0	50.0
13C4-PFHpA	Ave	210510	197868		47.0	50.0	-6.0	50.0
18O2 PFHxS	Ave	302453	298386		46.7	47.3	-1.3	50.0
13C4 PFOA	Ave	219411	208024		47.4	50.0	-5.2	50.0
13C4 PFOS	Ave	244616	241692		47.2	47.8	-1.2	50.0
13C5 PFNA	Ave	178109	162087		45.5	50.0	-9.0	50.0
13C8 FOSA	Ave	414253	389267		47.0	50.0	-6.0	50.0
13C2 PFDA	Ave	166067	160722		48.4	50.0	-3.2	50.0
13C2 PFUnA	Ave	125779	115580		45.9	50.0	-8.1	50.0
13C2 PFDoA	Ave	116905	110755		47.4	50.0	-5.3	50.0
13C2-PFTeDA	Ave	241899	239157		49.4	50.0	-1.1	50.0
13C2-PFHxDA	Ave	130373	130244		50.0	50.0	-0.1	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_012.d
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 02-Dec-2016 11:29:31 ALS Bottle#: 44 Worklist Smp#: 12
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: ICV_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist:
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 02-Dec-2016 16:10:33 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: barnettj Date: 02-Dec-2016 16:10:33

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.598	1.615	-0.017	15755416	46.3		92.6	1226463	
1 Perfluorobutyric acid	212.90 > 169.00	1.606	1.617	-0.011	13527367	48.5			149480	
D 4 13C5-PFPeA	267.90 > 223.00	1.897	1.918	-0.021	12529404	46.4		92.7	998019	
3 Perfluoropentanoic acid	262.90 > 219.00	1.897	1.920	-0.023	12023152	46.1			109623	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.936	1.961	-0.025	20638621	42.9				
	298.90 > 99.00	1.936	1.961	-0.025	9604105		2.15(0.00-0.00)			
D 6 13C2 PFHxA	315.00 > 270.00	2.213	2.239	-0.026	11340128	46.0		92.0	872752	
7 Perfluorohexanoic acid	313.00 > 269.00	2.213	2.241	-0.028	10367089	47.8			388218	
D 11 13C4-PFHpA	367.00 > 322.00	2.566	2.599	-0.033	9893393	47.0		94.0	1263133	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.573	2.600	-0.027	9494636	46.4			159398	
D 10 18O2 PFHxS	403.00 > 84.00	2.589	2.614	-0.025	14113680	46.7		98.7	888579	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.589	2.615	-0.026	13727834	43.4				
15 Perfluorooctanoic acid	413.00 > 369.00	2.951	2.984	-0.033	10247185	48.3			136835	
	413.00 > 169.00	2.943	2.984	-0.041	6439506		1.59(0.90-1.10)		319423	
D 14 13C4 PFOA	417.00 > 372.00	2.951	2.984	-0.033	10401182	47.4		94.8	602068	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.960	2.986	-0.026	1.000	13266199	47.0				
D 17 13C4 PFOS										
503.00 > 80.00	3.329	3.365	-0.036		11552855	47.2		98.8	248324	
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.329	3.366	-0.037	1.000	10458074	39.8			1887102	
499.00 > 99.00	3.329	3.366	-0.037	1.000	2551302		4.10(0.90-1.10)		466085	
20 Perfluorononanoic acid										
463.00 > 419.00	3.336	3.373	-0.037	1.000	7761968	48.3			97406	
D 19 13C5 PFNA										
468.00 > 423.00	3.336	3.374	-0.038		8104346	45.5		91.0	462400	
D 21 13C8 FOSA										
506.00 > 78.00	3.633	3.651	-0.018		19463355	47.0		94.0	341722	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.633	3.656	-0.023	1.000	18143701	49.8			283379	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.698	3.738	-0.040	1.000	7312192	46.9			156765	
D 23 13C2 PFDA										
515.00 > 470.00	3.698	3.738	-0.040		8036093	48.4		96.8	321643	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	4.010	4.049	-0.039	1.000	6928568	45.7				
D 27 13C2 PFUnA										
565.00 > 520.00	4.029	4.070	-0.041		5778978	45.9		91.9	312840	
28 Perfluoroundecanoic acid										
563.00 > 519.00	4.029	4.071	-0.042	1.000	5413472	44.2			130245	
D 30 13C2 PFDaA										
615.00 > 570.00	4.330	4.370	-0.040		5537768	47.4		94.7	163474	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.330	4.370	-0.040	1.000	4970310	47.4			6635	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.600	4.636	-0.036	1.000	5107276	44.3			7225	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.837	4.882	-0.045		11957860	49.4		98.9	638807	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.837	4.884	-0.047	1.000	10218574	47.7			5085	
713.00 > 169.00	4.837	4.884	-0.047	1.000	1568084		6.52(0.00-0.00)		100324	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.274	5.320	-0.046		6512177	50.0		99.9	123170	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.274	5.320	-0.046	1.000	5617133	50.7			6502	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.664	5.720	-0.056	1.000	5682438	65.3			7600	

Reagents:

LCPFCIC_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_012.d

Injection Date: 02-Dec-2016 11:29:31

Instrument ID: A8_N

Lims ID: ICV

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 44

Worklist Smp#: 12

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

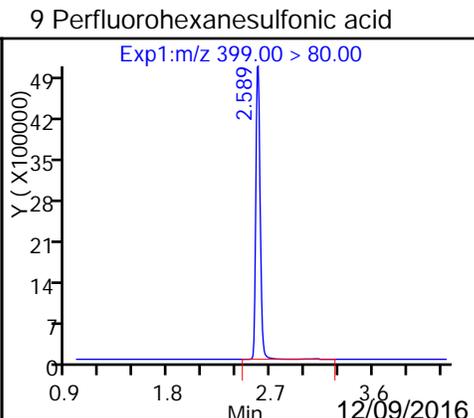
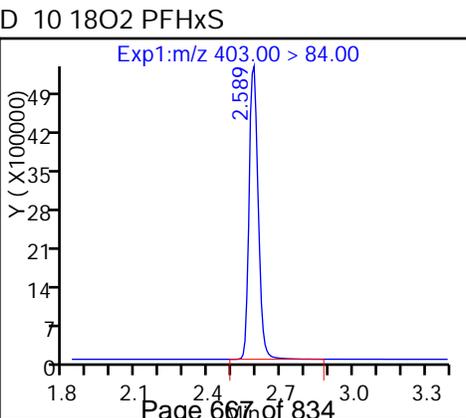
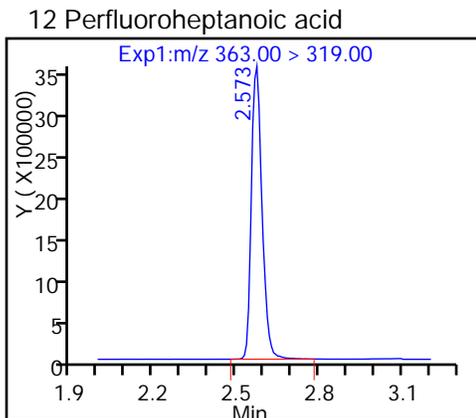
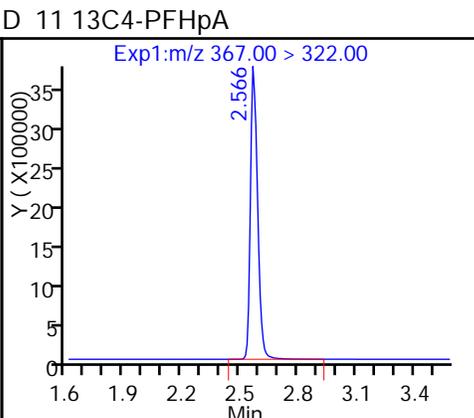
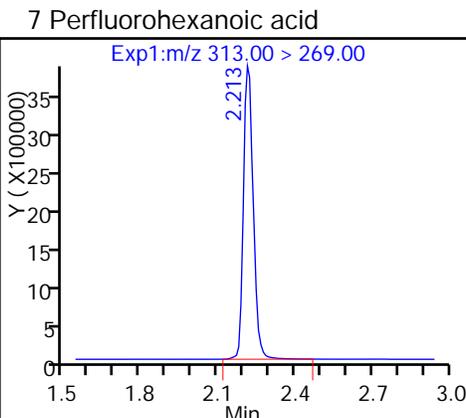
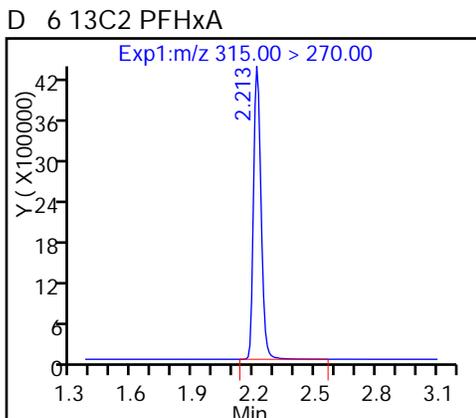
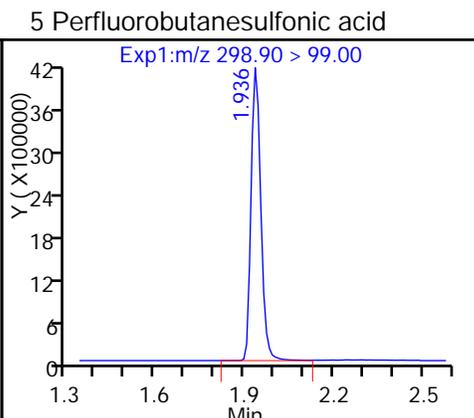
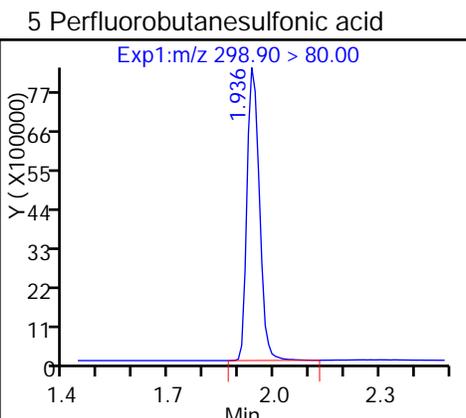
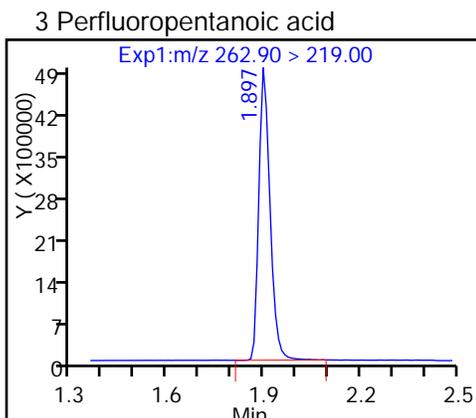
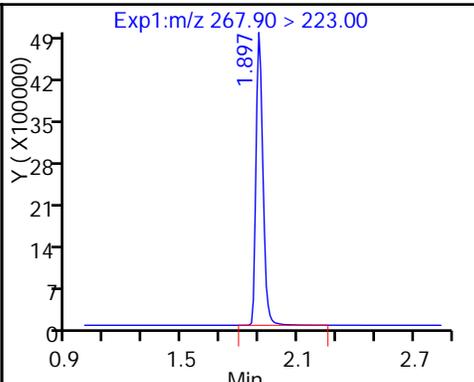
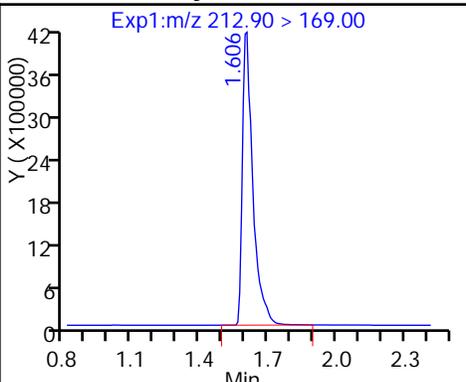
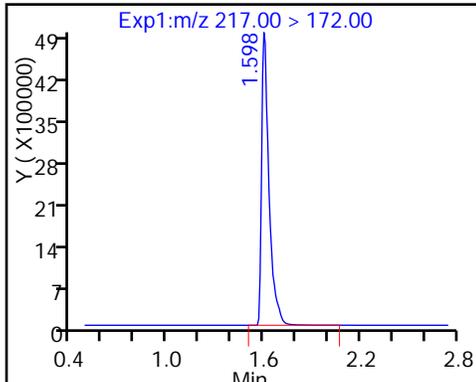
Method: A8_N

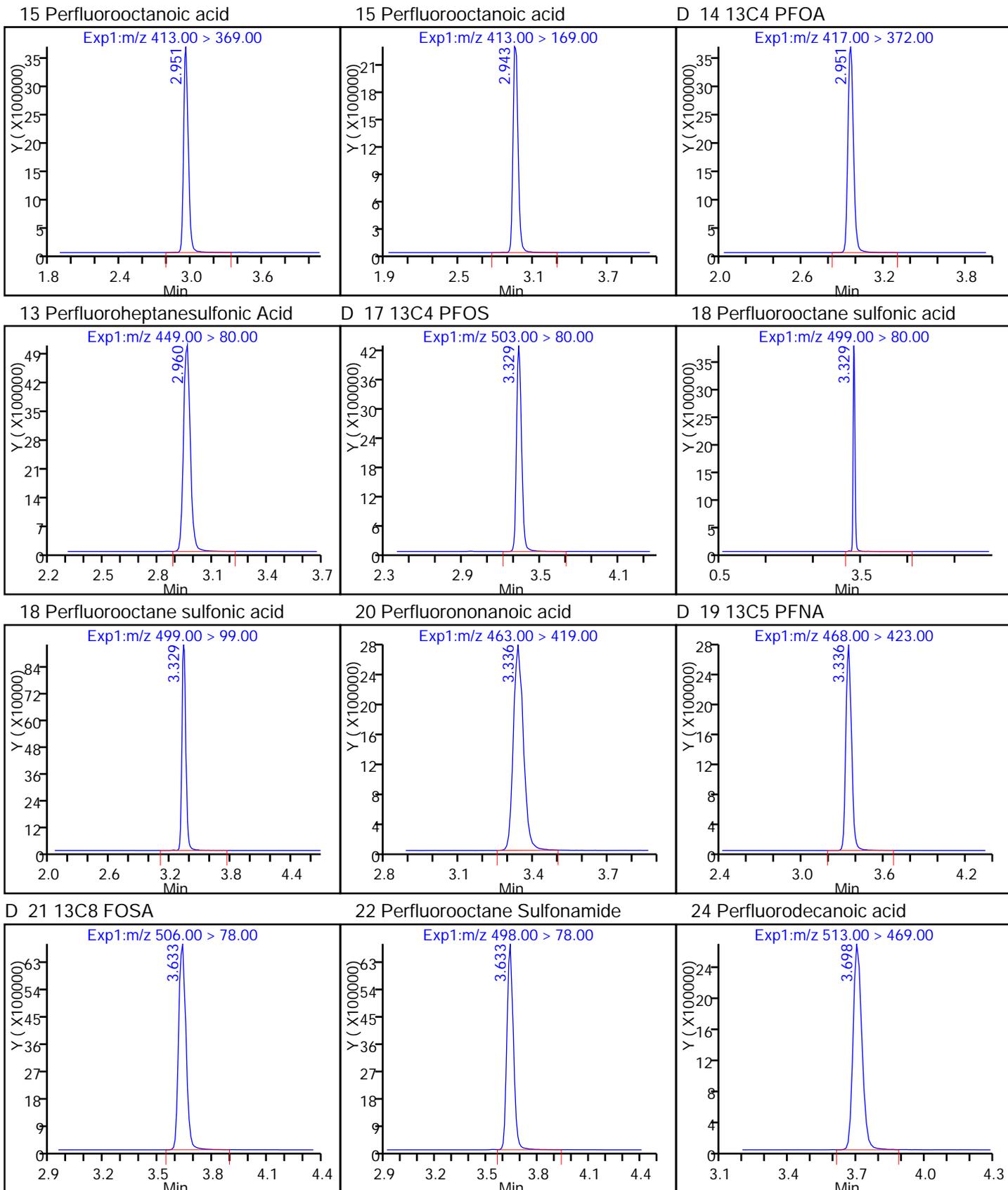
Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

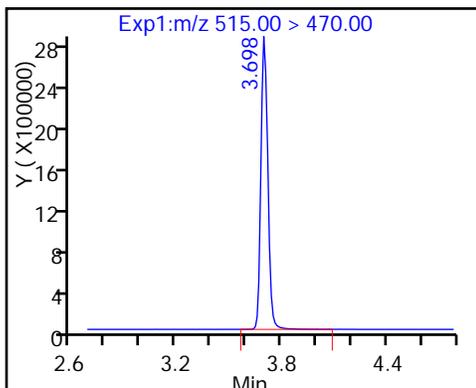
1 Perfluorobutyric acid

D 4 13C5-PFPeA

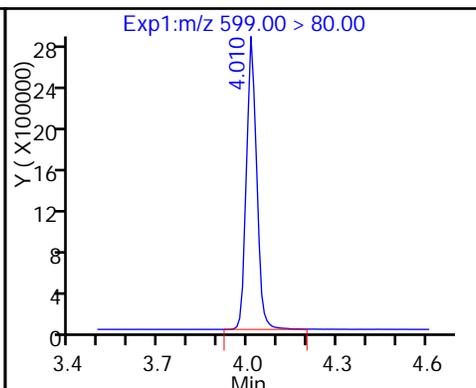




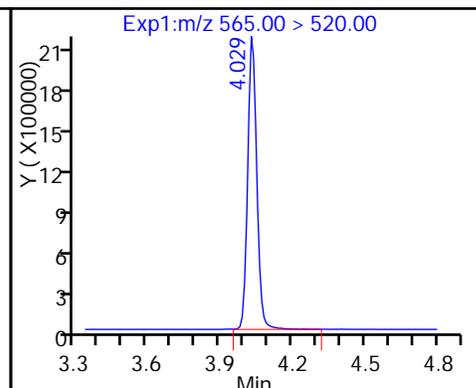
D 23 13C2 PFDA



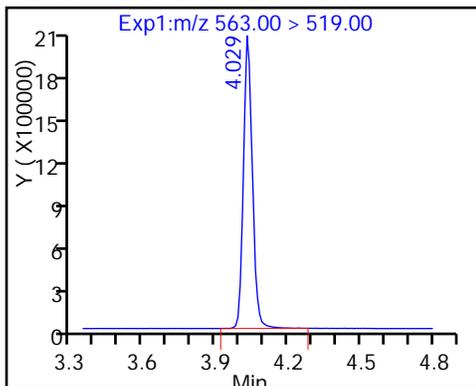
26 Perfluorodecane Sulfonic acid



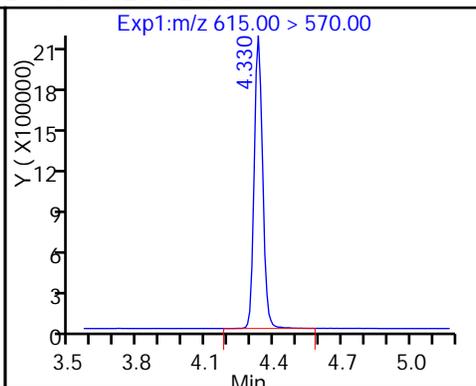
D 27 13C2 PFUnA



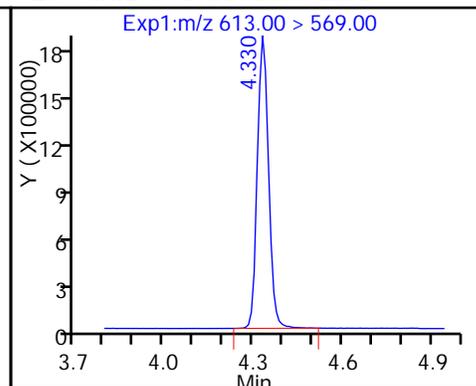
28 Perfluoroundecanoic acid



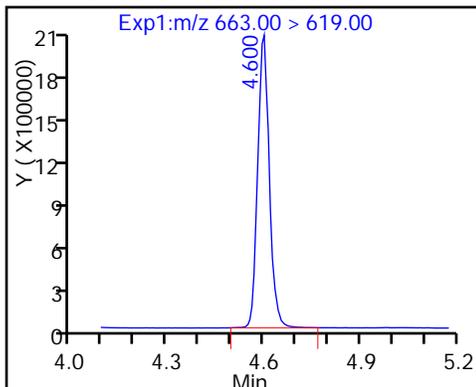
D 30 13C2 PFDaA



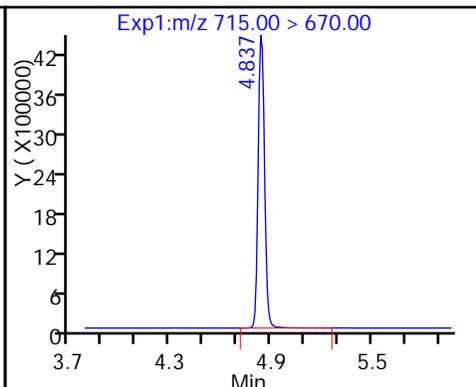
29 Perfluorododecanoic acid



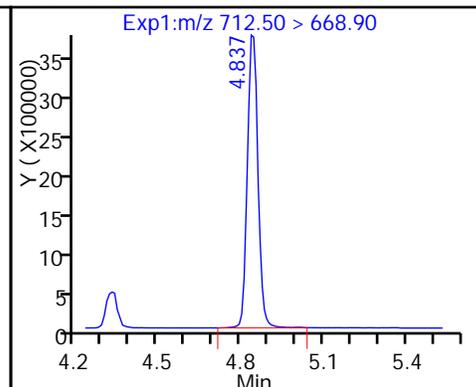
31 Perfluorotridecanoic acid



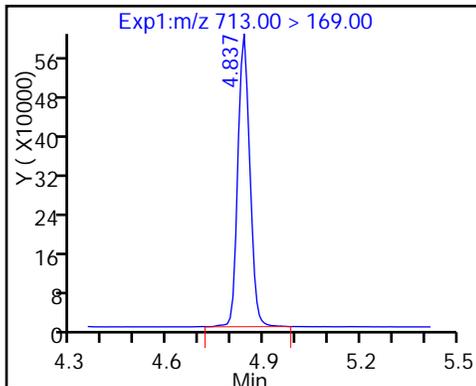
D 32 13C2-PFTeDA



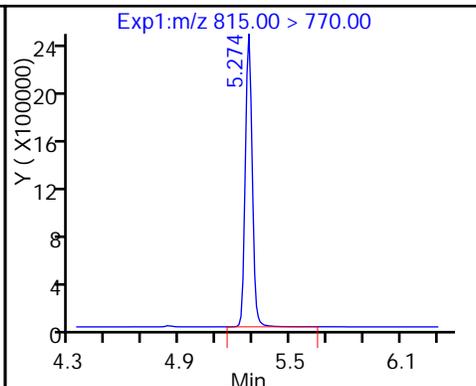
33 Perfluorotetradecanoic acid



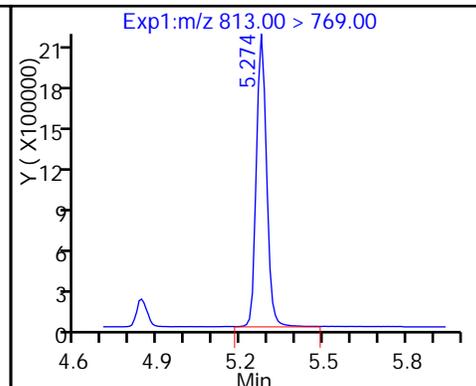
33 Perfluorotetradecanoic acid



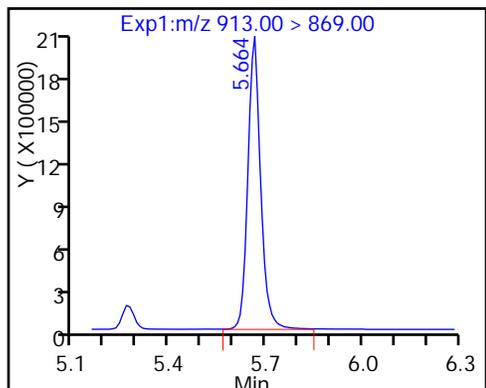
D 34 13C2-PFHxD A



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-140429/2 Calibration Date: 12/02/2016 14:37
 Instrument ID: A8_N Calib Start Date: 12/02/2016 10:29
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/02/2016 12:29
 Lab File ID: 02DEC2016B_014.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.613	1.603		43.9	44.2	-0.6	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.020	1.035		50.7	50.0	1.5	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.087	1.084		46.3	46.4	-0.3	25.0
13C4 PFBA	Ave	340300	339545		49.9	50.0	-0.2	50.0
13C5-PFPeA	Ave	270194	261302		48.4	50.0	-3.3	50.0
13C2 PFHxA	Ave	246480	229526		46.6	50.0	-6.9	50.0
13C4-PFHpA	Ave	210510	202011		48.0	50.0	-4.0	50.0
18O2 PFHxS	Ave	302453	317474		49.6	47.3	5.0	50.0
13C4 PFOA	Ave	219411	216146		49.3	50.0	-1.5	50.0
13C4 PFOS	Ave	244616	251364		49.1	47.8	2.8	50.0
13C5 PFNA	Ave	178109	169399		47.6	50.0	-4.9	50.0
13C8 FOSA	Ave	414253	402881		48.6	50.0	-2.7	50.0
13C2 PFDA	Ave	166067	161124		48.5	50.0	-3.0	50.0
13C2 PFUnA	Ave	125779	118770		47.2	50.0	-5.6	50.0
13C2 PFDoA	Ave	116905	109471		46.8	50.0	-6.4	50.0
13C2-PFTeDA	Ave	241899	223925		46.3	50.0	-7.4	50.0
13C2-PFHxDA	Ave	130373	118805		45.6	50.0	-8.9	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016B_014.d
 Lims ID: CCV L5
 Client ID:
 Sample Type: CCV
 Inject. Date: 02-Dec-2016 14:37:03 ALS Bottle#: 41 Worklist Smp#: 2
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L5
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:06:11 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d

Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:06:11

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.574	0.0	1.000	15171458	50.5	101	155330	
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0		16977232	49.9	99.8	2771435	
3 Perfluoropentanoic acid	262.90 > 219.00	1.868	1.868	0.0	1.000	12970874	47.7	95.4	120574	
D 4 13C5-PFPeA	267.90 > 223.00	1.868	1.868	0.0		13065075	48.4	96.7	2568563	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.907	1.907	0.0	1.000	22493915	43.9	99.4		
	298.90 > 99.00	1.907	1.907	0.0	1.000	10590632	2.12(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.173	2.173	0.0	1.000	10840520	49.4	98.8	348131	
D 6 13C2 PFHxA	315.00 > 270.00	2.173	2.173	0.0		11476282	46.6	93.1	919645	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.519	2.519	0.0	1.000	10487509	50.2	100	129458	
D 11 13C4-PFHpA	367.00 > 322.00	2.519	2.519	0.0		10100555	48.0	96.0	962415	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.534	2.534	0.0	1.000	15406543	45.8	101		
D 10 18O2 PFHxS	403.00 > 84.00	2.534	2.534	0.0		15016500	49.6	105	1172460	
D 14 13C4 PFOA	417.00 > 372.00	2.886	2.886	0.0		10807294	49.3	98.5	772353	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.894	2.894	0.0	1.000	11190759	50.7		101	205894	
413.00 > 169.00	2.894	2.894	0.0	1.000	6841143		1.64(0.90-1.10)		311496	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.902	2.902	0.0	1.000	14021365	47.7		100		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.267	3.267	0.0	1.000	12644938	46.3		99.7	53281	
499.00 > 99.00	3.267	3.267	0.0	1.000	2940349		4.30(0.90-1.10)		339746	
D 17 13C4 PFOS										
503.00 > 80.00	3.267	3.267	0.0		12015188	49.1		103	275554	
20 Perfluorononanoic acid										
463.00 > 419.00	3.275	3.275	0.0	1.000	8262291	49.2		98.4	153747	
D 19 13C5 PFNA										
468.00 > 423.00	3.275	3.275	0.0		8469930	47.6		95.1	605451	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.589	3.589	0.0	1.000	18261930	48.5		96.9	733539	
D 21 13C8 FOSA										
506.00 > 78.00	3.589	3.589	0.0		20144056	48.6		97.3	630987	
D 23 13C2 PFDA										
515.00 > 470.00	3.631	3.631	0.0		8056201	48.5		97.0	192938	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.631	3.631	0.0	1.000	7623608	48.8		97.6	194009	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.941	3.941	0.0	1.000	8048082	51.1		106		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.959	3.959	0.0	1.000	5737327	45.6		91.2	103207	
D 27 13C2 PFUnA										
565.00 > 520.00	3.967	3.967	0.0		5938511	47.2		94.4	310740	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.259	4.259	0.0	1.000	5083644	49.1		98.1	71841	
D 30 13C2 PFDaA										
615.00 > 570.00	4.259	4.259	0.0		5473570	46.8		93.6	186491	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.520	4.520	0.0	1.000	5281655	46.4		92.8	6905	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.772	4.772	0.0	1.000	10203271	48.2		96.5	4031	
713.00 > 169.00	4.763	4.772	-0.009	0.998	1624426		6.28(0.00-0.00)		120314	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.772	4.772	0.0		11196256	46.3		92.6	372028	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.184	5.184	0.0		5940244	45.6		91.1	99394	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.194	5.194	0.0	1.000	5316696	48.5		97.0	3968	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.562	5.562	0.0	1.000	5707122	66.3		133	7784	

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016B_014.d

Injection Date: 02-Dec-2016 14:37:03

Instrument ID: A8_N

Lims ID: CCV L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 2

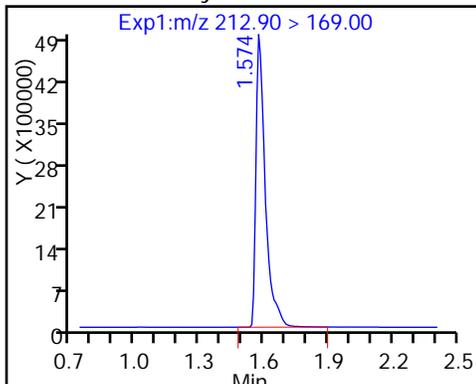
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

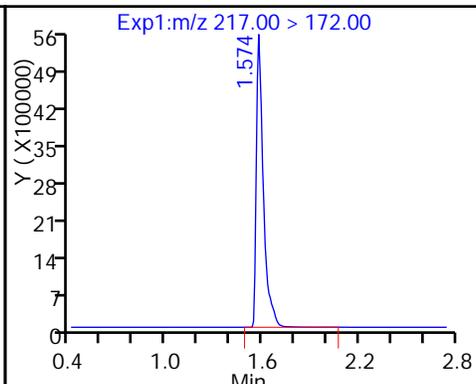
Method: A8_N

Limit Group: LC PFC_DOD ICAL

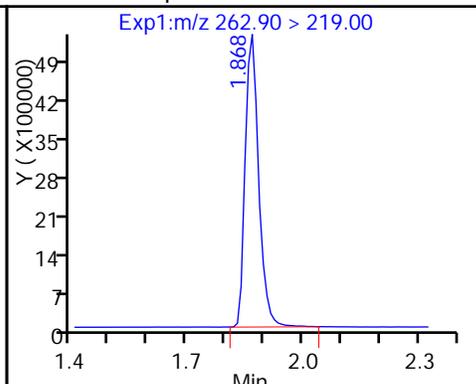
1 Perfluorobutyric acid



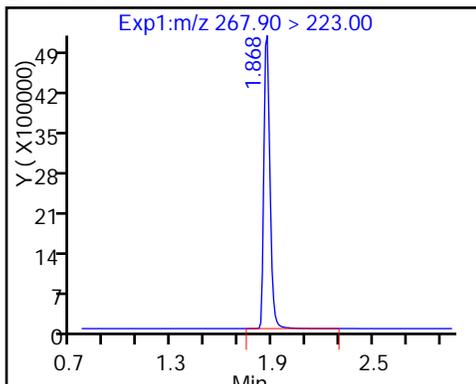
D 2 13C4 PFBA



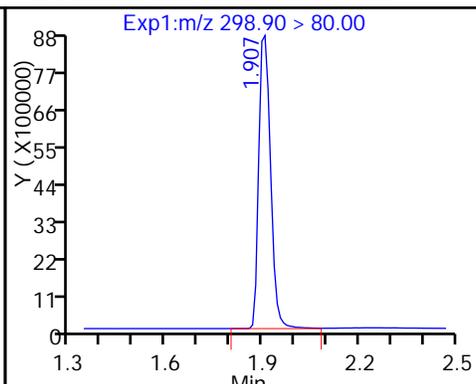
3 Perfluoropentanoic acid



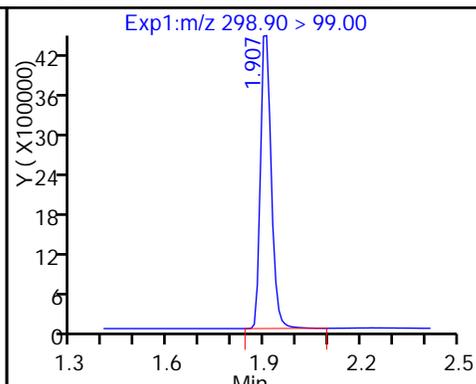
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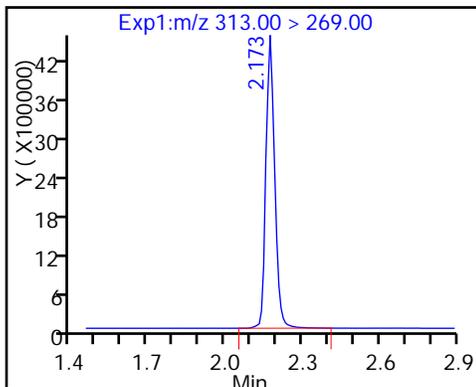
5 Perfluorobutanesulfonic acid



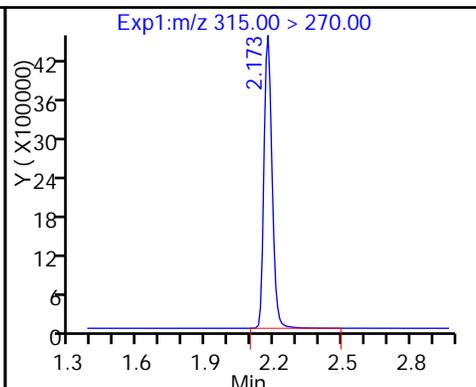
5 Perfluorobutanesulfonic acid



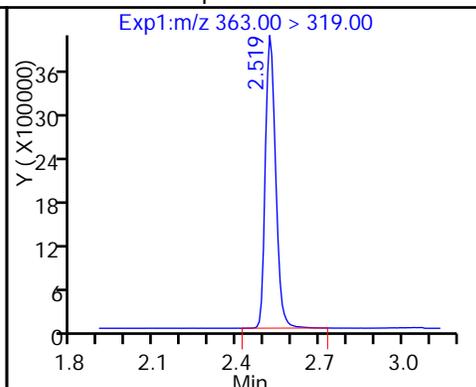
7 Perfluorohexanoic acid



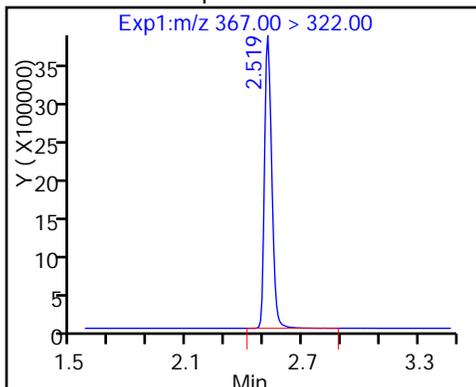
D 6 13C2 PFHxA



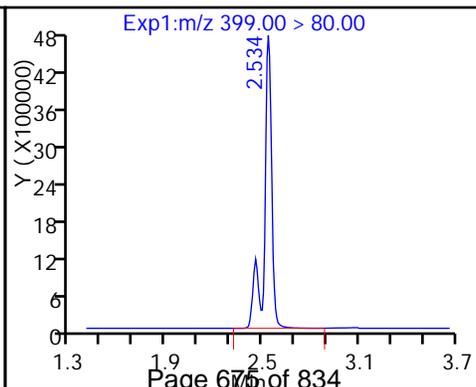
12 Perfluoroheptanoic acid



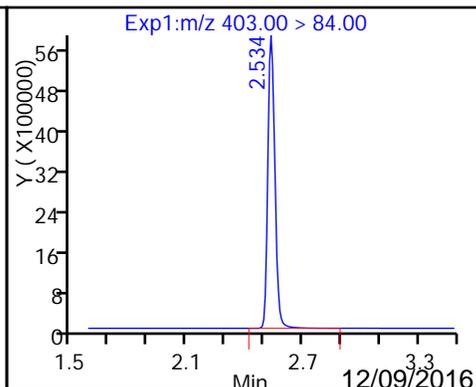
D 11 13C4-PFHpA



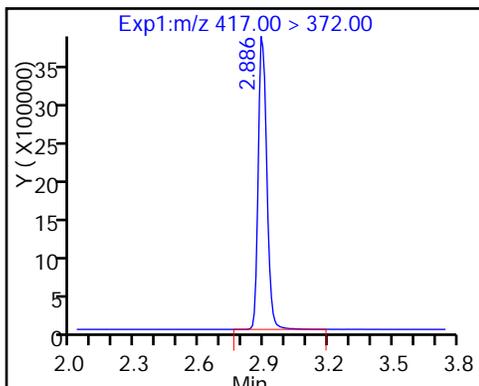
9 Perfluorohexanesulfonic acid



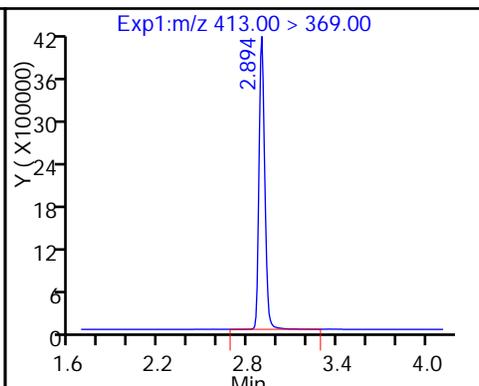
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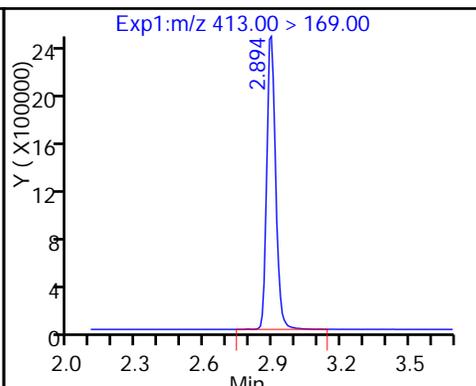
D 14 13C4 PFOA



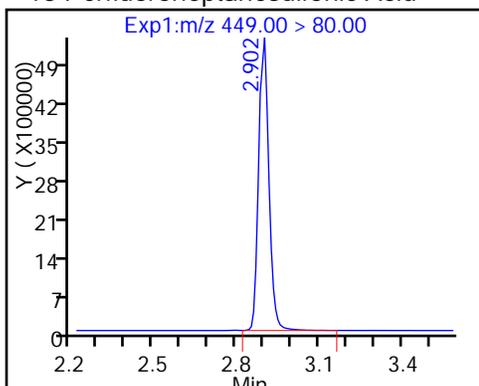
15 Perfluorooctanoic acid



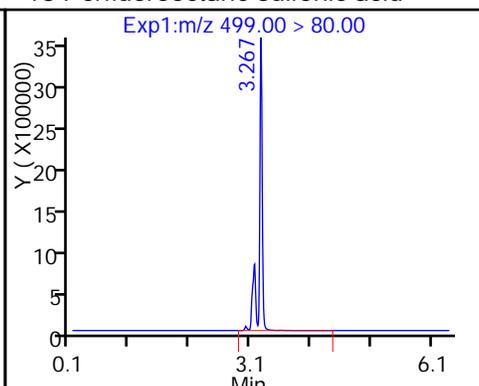
15 Perfluorooctanoic acid



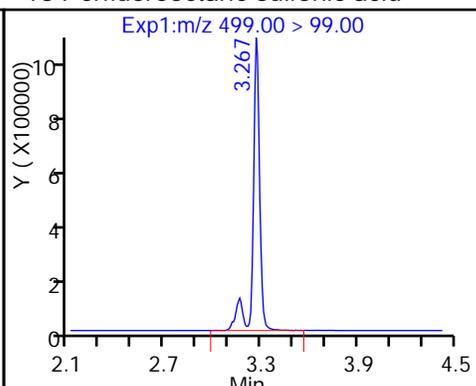
13 Perfluoroheptanesulfonic Acid



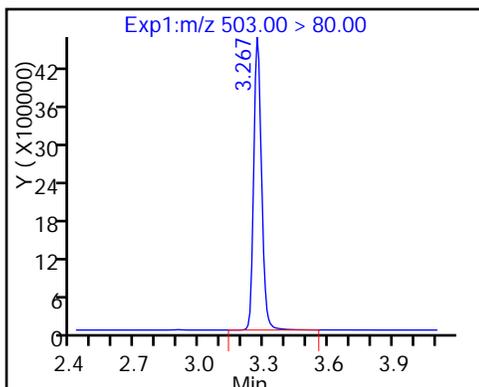
18 Perfluorooctane sulfonic acid



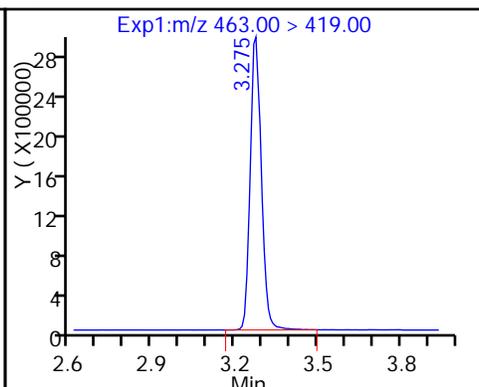
18 Perfluorooctane sulfonic acid



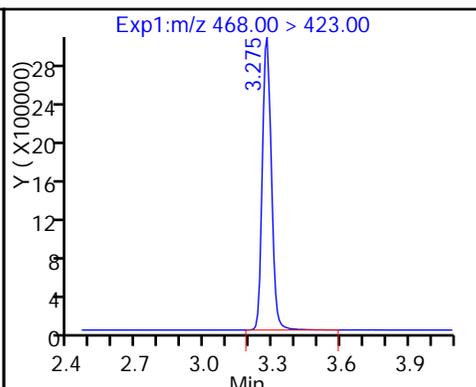
D 17 13C4 PFOS



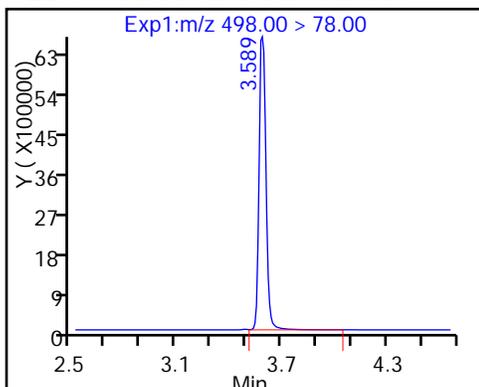
20 Perfluorononanoic acid



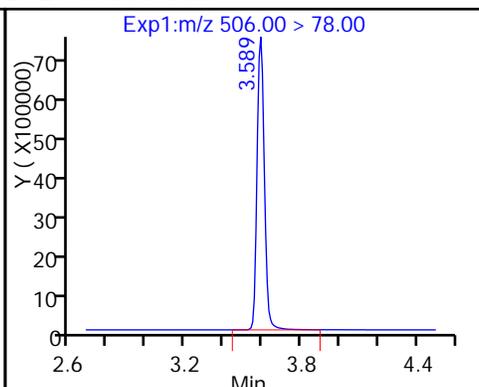
D 19 13C5 PFNA



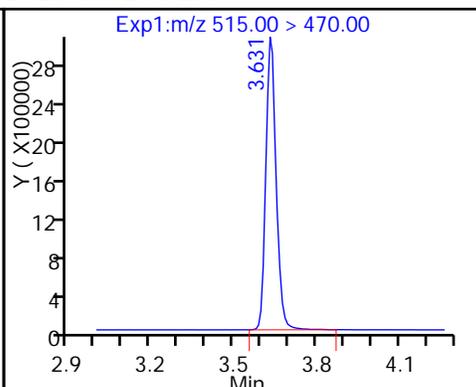
22 Perfluorooctane Sulfonamide

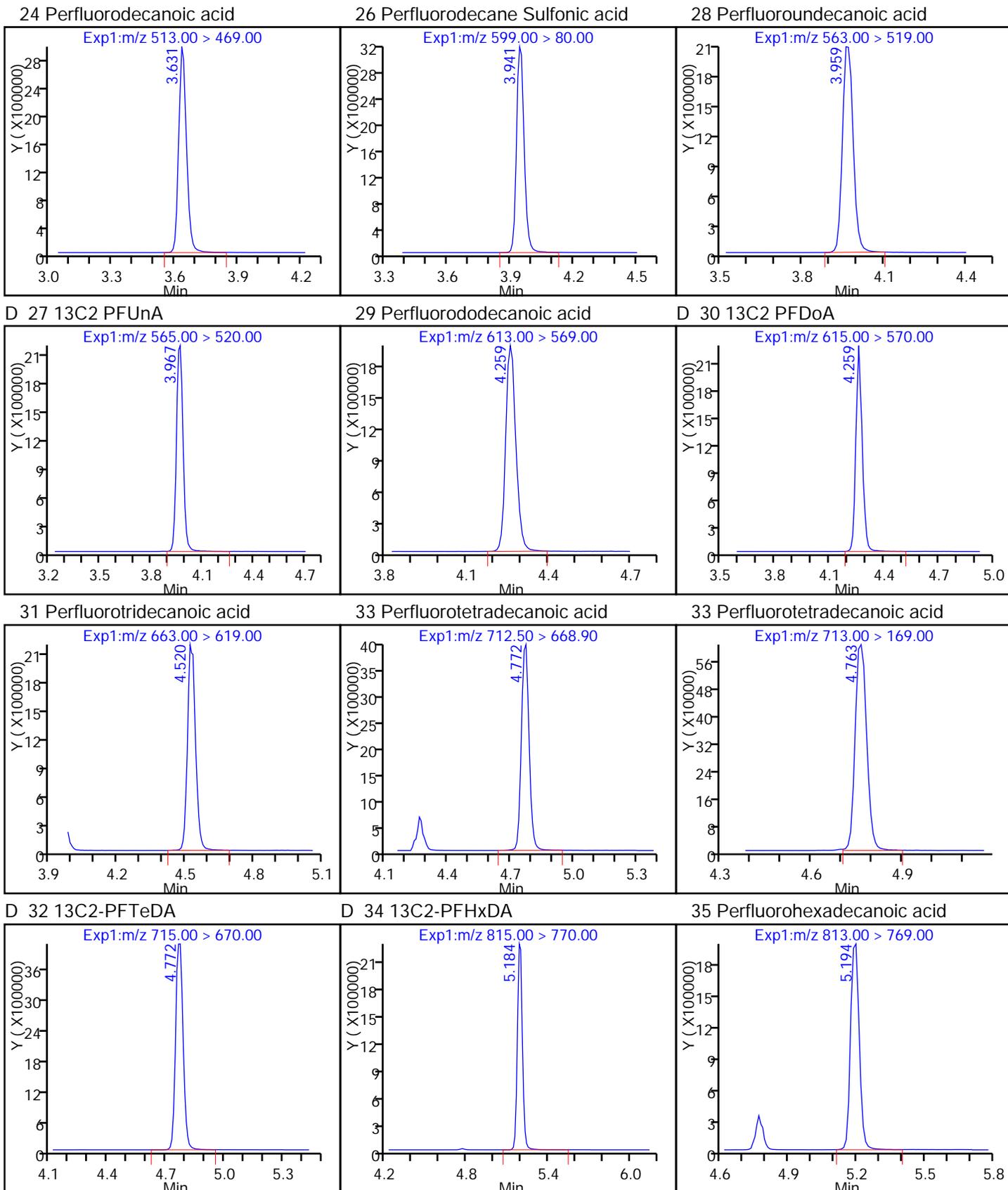


D 21 13C8 FOSA

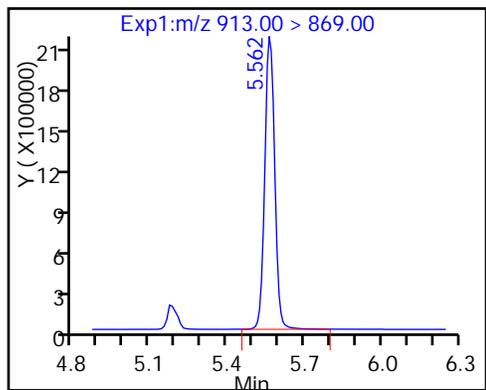


D 23 13C2 PFDA





36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-140429/16 Calibration Date: 12/02/2016 16:22
 Instrument ID: A8_N Calib Start Date: 12/02/2016 10:29
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/02/2016 12:29
 Lab File ID: 02DEC2016C_012.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.613	1.597		43.7	44.2	-1.0	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.020	1.019		49.9	50.0	-0.1	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.087	1.065		45.5	46.4	-2.0	25.0
13C4 PFBA	Ave	340300	332191		48.8	50.0	-2.4	50.0
13C5-PFPeA	Ave	270194	252615		46.7	50.0	-6.5	50.0
13C2 PFHxA	Ave	246480	228018		46.3	50.0	-7.5	50.0
13C4-PFHpA	Ave	210510	196748		46.7	50.0	-6.5	50.0
18O2 PFHxS	Ave	302453	307137		48.0	47.3	1.5	50.0
13C4 PFOA	Ave	219411	211412		48.2	50.0	-3.6	50.0
13C4 PFOS	Ave	244616	252256		49.3	47.8	3.1	50.0
13C5 PFNA	Ave	178109	159978		44.9	50.0	-10.2	50.0
13C8 FOSA	Ave	414253	392650		47.4	50.0	-5.2	50.0
13C2 PFDA	Ave	166067	151711		45.7	50.0	-8.6	50.0
13C2 PFUnA	Ave	125779	110399		43.9	50.0	-12.2	50.0
13C2 PFDoA	Ave	116905	109030		46.6	50.0	-6.7	50.0
13C2-PFTeDA	Ave	241899	217472		45.0	50.0	-10.1	50.0
13C2-PFHxDA	Ave	130373	114036		43.7	50.0	-12.5	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_012.d
 Lims ID: CCV L5
 Client ID:
 Sample Type: CCV
 Inject. Date: 02-Dec-2016 16:22:13 ALS Bottle#: 41 Worklist Smp#: 16
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L5
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:06:22 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:06:22

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.574	0.0	1.000	14622963	49.7	99.4	136701	
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0		16609540	48.8	97.6	1760564	
3 Perfluoropentanoic acid	262.90 > 219.00	1.859	1.868	-0.009	1.000	12389685	47.1	94.2	143388	
D 4 13C5-PFPeA	267.90 > 223.00	1.859	1.868	-0.009		12630763	46.7	93.5	1231239	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.907	1.907	0.0	1.000	21678790	43.7	99.0		
	298.90 > 99.00	1.897	1.907	-0.010	0.995	10208437		2.12(0.00-0.00)		
7 Perfluorohexanoic acid	313.00 > 269.00	2.171	2.173	-0.003	1.000	10711219	49.1	98.3	288186	
D 6 13C2 PFHxA	315.00 > 270.00	2.171	2.173	-0.003		11400916	46.3	92.5	2220908	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.516	2.519	-0.003	1.000	10025294	49.3	98.5	148580	
D 11 13C4-PFHpA	367.00 > 322.00	2.516	2.519	-0.003		9837421	46.7	93.5	979084	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.532	2.534	-0.002	1.000	14850728	45.6	100		
D 10 18O2 PFHxS	403.00 > 84.00	2.532	2.534	-0.002		14527568	48.0	102	761217	
D 14 13C4 PFOA	417.00 > 372.00	2.890	2.886	0.004		10570593	48.2	96.4	786274	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.890	2.894	-0.004	1.000	10770410	49.9		99.9	198921	
413.00 > 169.00	2.890	2.894	-0.004	1.000	6718323		1.60(0.90-1.10)		592810	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.890	2.902	-0.012	1.000	13560651	46.0		96.6		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.265	3.267	-0.002	1.000	12467043	45.5		98.0	3525927	
499.00 > 99.00	3.265	3.267	-0.002	1.000	2716230		4.59(0.90-1.10)		228232	
D 17 13C4 PFOS										
503.00 > 80.00	3.265	3.267	-0.002		12057843	49.3		103	232224	
20 Perfluorononanoic acid										
463.00 > 419.00	3.272	3.275	-0.003	1.000	7963240	50.2		100	144737	
D 19 13C5 PFNA										
468.00 > 423.00	3.265	3.275	-0.010		7998877	44.9		89.8	353617	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.587	3.589	-0.002	1.000	18194106	49.5		99.1	536694	
D 21 13C8 FOSA										
506.00 > 78.00	3.587	3.589	-0.002		19632521	47.4		94.8	564736	
D 23 13C2 PFDA										
515.00 > 470.00	3.629	3.631	-0.002		7585560	45.7		91.4	208800	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.629	3.631	-0.002	1.000	7277017	49.5		99.0	199518	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.938	3.941	-0.003	1.000	7729327	48.9		101		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.956	3.959	-0.003	1.000	5431519	46.4		92.9	112110	
D 27 13C2 PFUnA										
565.00 > 520.00	3.956	3.967	-0.011		5519927	43.9		87.8	528013	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.255	4.259	-0.004	1.000	5067396	49.1		98.2	77314	
D 30 13C2 PFDaA										
615.00 > 570.00	4.246	4.259	-0.013		5451503	46.6		93.3	157285	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.521	4.520	0.001	1.000	5122725	45.2		90.3	11584	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.757	4.772	-0.015	1.000	9633033	45.7		91.4	4134	
713.00 > 169.00	4.748	4.772	-0.024	0.998	1613548		5.97(0.00-0.00)		215079	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.757	4.772	-0.015		10873598	45.0		89.9	570141	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.173	5.184	-0.011		5701819	43.7		87.5	107923	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.184	5.194	-0.010	1.000	5253333	48.1		96.2	4932	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.545	5.562	-0.017	1.000	5747404	67.0		134	6736	

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_012.d

Injection Date: 02-Dec-2016 16:22:13

Instrument ID: A8_N

Lims ID: CCV L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 16

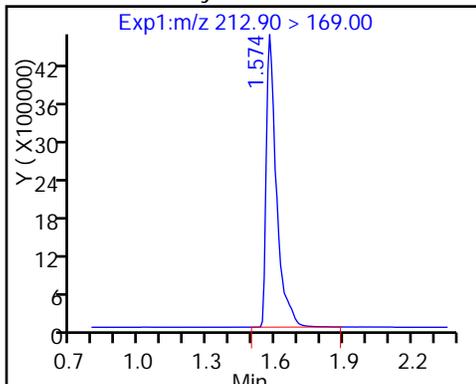
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

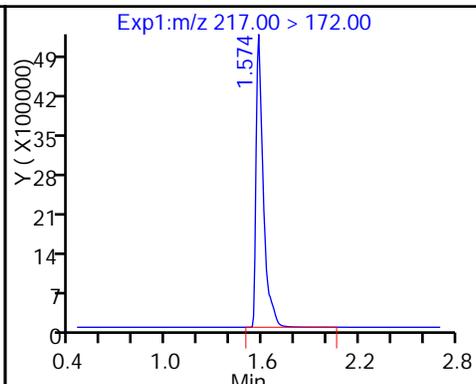
Method: A8_N

Limit Group: LC PFC_DOD ICAL

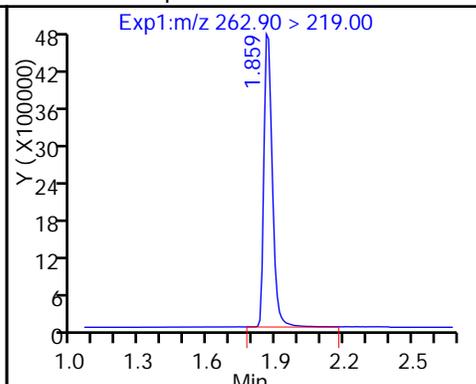
1 Perfluorobutyric acid



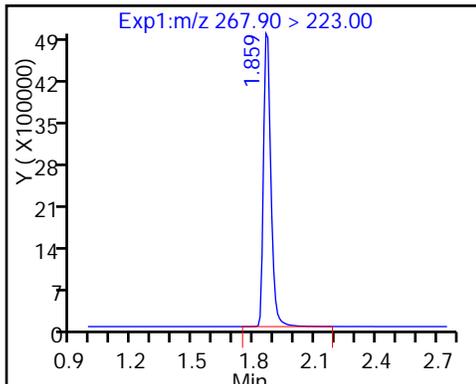
D 2 13C4 PFBA



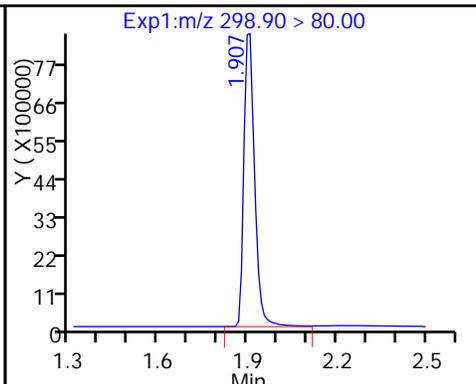
3 Perfluoropentanoic acid



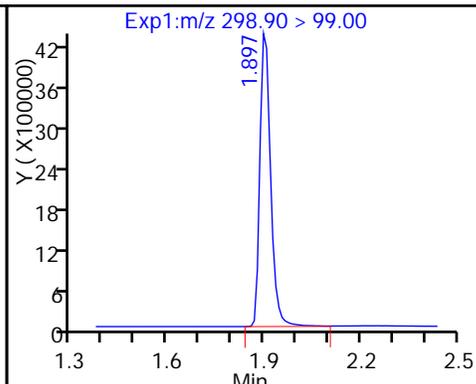
D 4 13C5-PFPeA



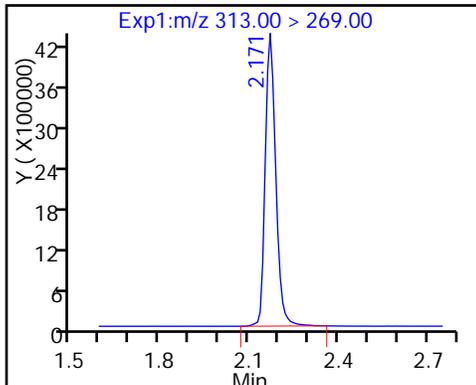
5 Perfluorobutanesulfonic acid



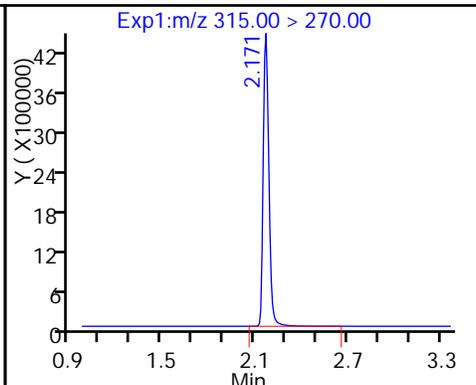
5 Perfluorobutanesulfonic acid



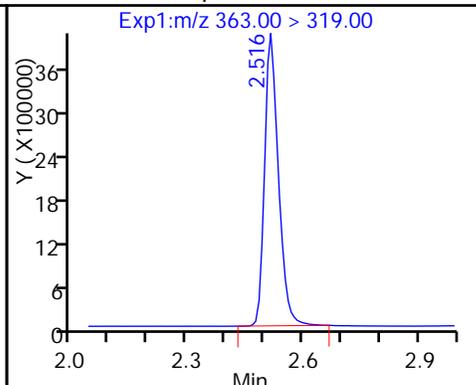
7 Perfluorohexanoic acid



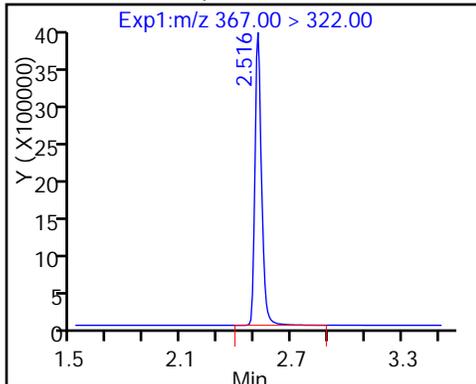
D 6 13C2 PFHxA



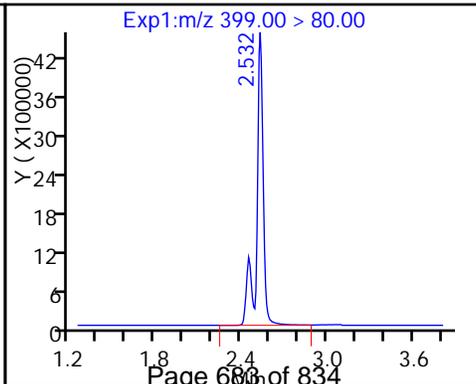
12 Perfluoroheptanoic acid



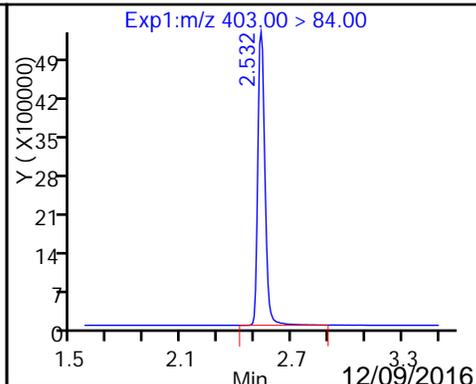
D 11 13C4-PFHpA



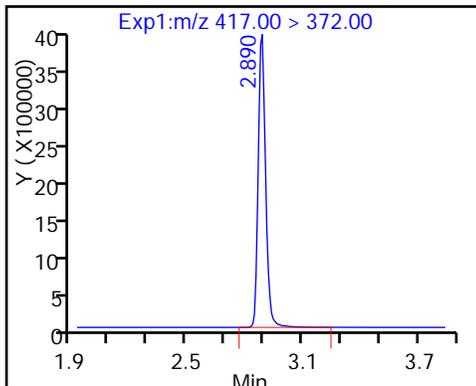
9 Perfluorohexanesulfonic acid



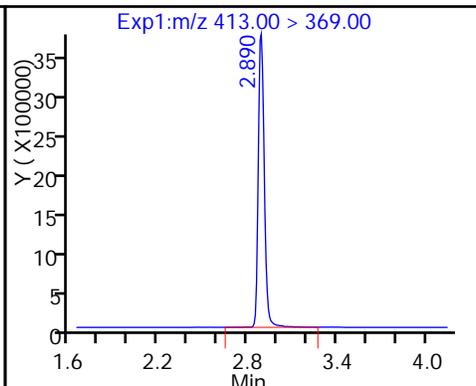
D 10 18O2 PFHxS



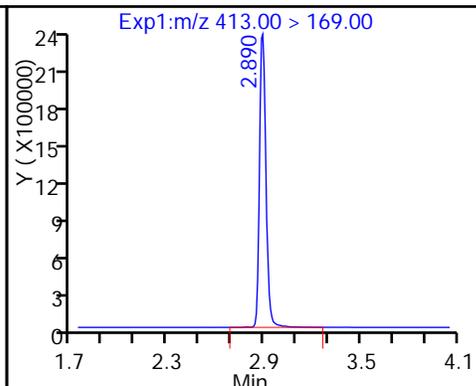
D 14 13C4 PFOA



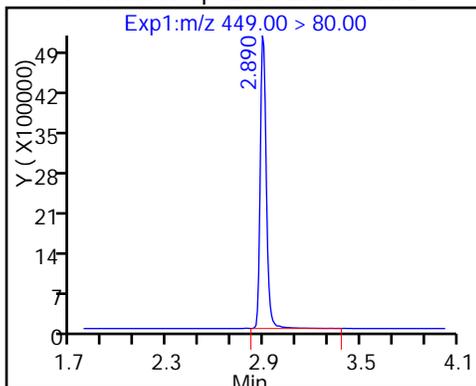
15 Perfluorooctanoic acid



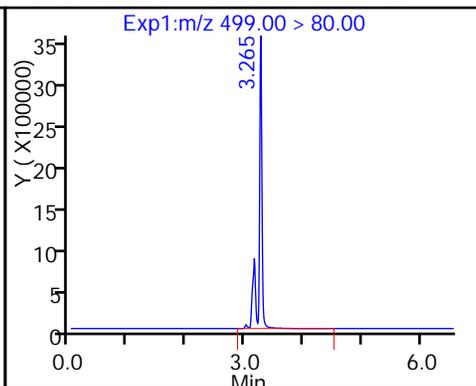
15 Perfluorooctanoic acid



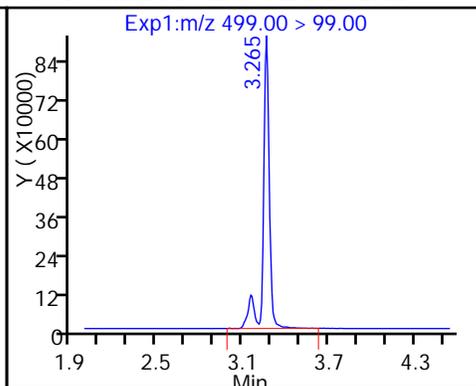
13 Perfluoroheptanesulfonic Acid



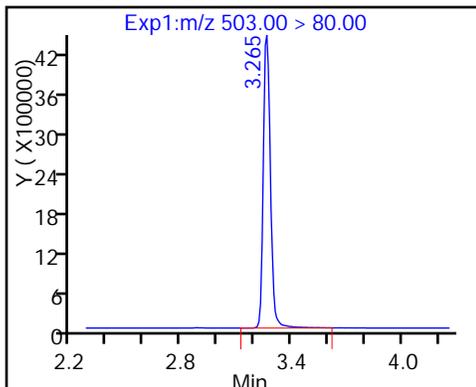
18 Perfluorooctane sulfonic acid



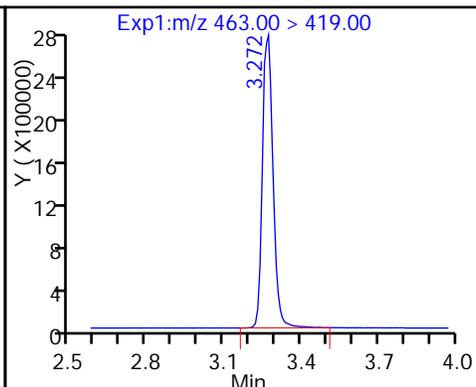
18 Perfluorooctane sulfonic acid



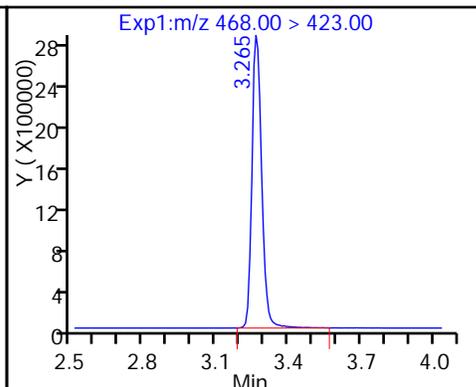
D 17 13C4 PFOS



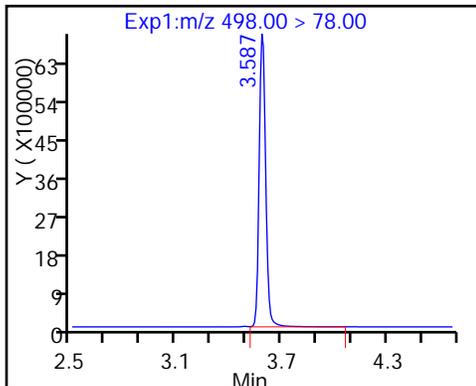
20 Perfluorononanoic acid



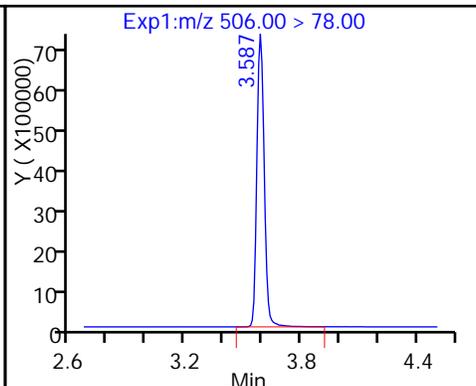
D 19 13C5 PFNA



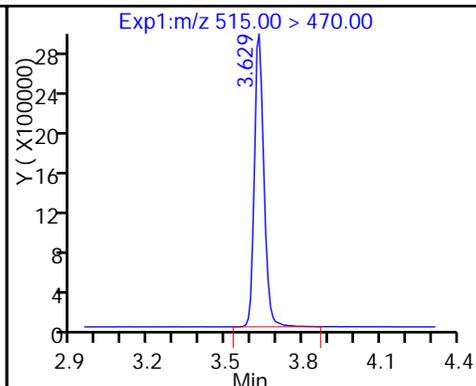
22 Perfluorooctane Sulfonamide

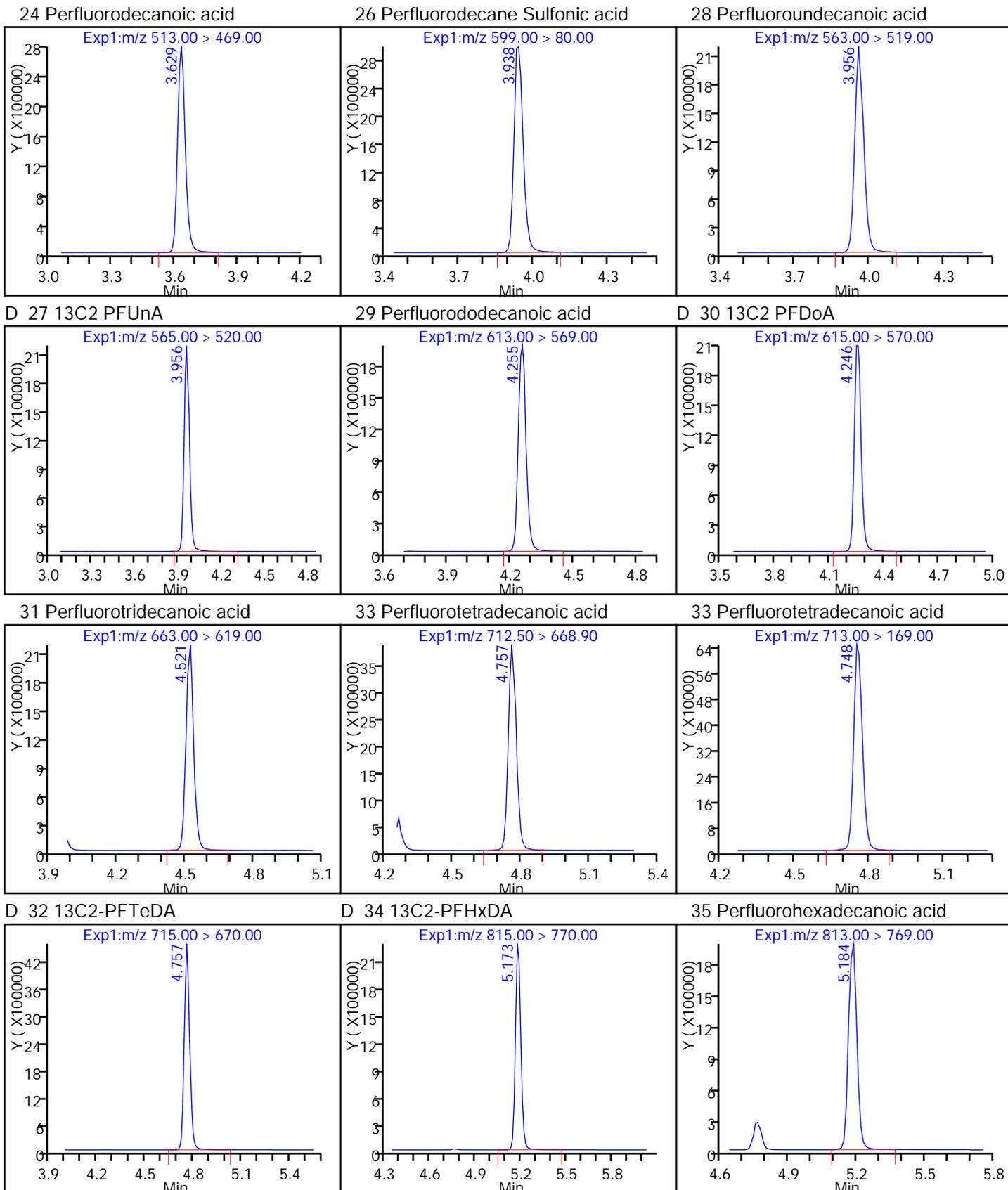


D 21 13C8 FOSA

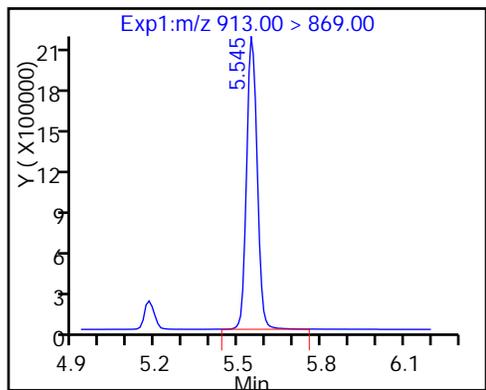


D 23 13C2 PFDA





36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-140429/26 Calibration Date: 12/02/2016 17:37
 Instrument ID: A8_N Calib Start Date: 12/02/2016 10:29
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/02/2016 12:29
 Lab File ID: 02DEC2016C_022.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.613	1.598		43.8	44.2	-1.0	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.020	1.046		51.3	50.0	2.5	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.087	1.089		46.5	46.4	0.2	25.0
13C4 PFBA	Ave	340300	325436		47.8	50.0	-4.4	50.0
13C5-PFPeA	Ave	270194	253322		46.9	50.0	-6.2	50.0
13C2 PFHxA	Ave	246480	222118		45.1	50.0	-9.9	50.0
13C4-PFHpA	Ave	210510	196792		46.7	50.0	-6.5	50.0
18O2 PFHxS	Ave	302453	292688		45.8	47.3	-3.2	50.0
13C4 PFOA	Ave	219411	207834		47.4	50.0	-5.3	50.0
13C4 PFOS	Ave	244616	231762		45.3	47.8	-5.3	50.0
13C5 PFNA	Ave	178109	162211		45.5	50.0	-8.9	50.0
13C8 FOSA	Ave	414253	370800		44.8	50.0	-10.5	50.0
13C2 PFDA	Ave	166067	151478		45.6	50.0	-8.8	50.0
13C2 PFUnA	Ave	125779	112787		44.8	50.0	-10.3	50.0
13C2 PFDoA	Ave	116905	108715		46.5	50.0	-7.0	50.0
13C2-PFTeDA	Ave	241899	220421		45.6	50.0	-8.9	50.0
13C2-PFHxDA	Ave	130373	118772		45.6	50.0	-8.9	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_022.d
 Lims ID: CCV L5
 Client ID:
 Sample Type: CCV
 Inject. Date: 02-Dec-2016 17:37:22 ALS Bottle#: 41 Worklist Smp#: 26
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L5
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:31:59 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:31:59

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
1 Perfluorobutyric acid	212.90 > 169.00	1.575	1.574	0.001	1.000	14279480	49.6	99.1	139417	
D 2 13C4 PFBA	217.00 > 172.00	1.575	1.574	0.001		16271795	47.8	95.6	1290471	
3 Perfluoropentanoic acid	262.90 > 219.00	1.859	1.868	-0.009	1.000	12234509	46.4	92.8	122207	
D 4 13C5-PFPeA	267.90 > 223.00	1.859	1.868	-0.009		12666104	46.9	93.8	1256391	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.898	1.907	-0.009	1.000	20670435	43.8	99.0		
	298.90 > 99.00	1.898	1.907	-0.009	1.000	9724349		2.13(0.00-0.00)		
7 Perfluorohexanoic acid	313.00 > 269.00	2.164	2.173	-0.009	1.000	10659778	50.2	100	311210	
D 6 13C2 PFHxA	315.00 > 270.00	2.164	2.173	-0.009		11105875	45.1	90.1	1390903	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.511	2.519	-0.008	1.000	9965121	49.0	97.9	129553	
D 11 13C4-PFHpA	367.00 > 322.00	2.511	2.519	-0.008		9839617	46.7	93.5	606816	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.526	2.534	-0.008	1.000	13863696	44.7	98.2		M
D 10 18O2 PFHxS	403.00 > 84.00	2.526	2.534	-0.008		13844166	45.8	96.8	284370	
D 14 13C4 PFOA	417.00 > 372.00	2.877	2.886	-0.009		10391697	47.4	94.7	923492	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.877	2.894	-0.017	1.000	10869296	51.3		103	351138	
413.00 > 169.00	2.877	2.894	-0.017	1.000	6481914		1.68(0.90-1.10)		594276	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.885	2.902	-0.017	1.000	13016744	48.0		101		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.258	3.267	-0.009	1.000	11713855	46.5		100	833502	
499.00 > 99.00	3.258	3.267	-0.009	1.000	2574140		4.55(0.90-1.10)		854286	
D 17 13C4 PFOS										
503.00 > 80.00	3.258	3.267	-0.009		11078221	45.3		94.7	225908	
20 Perfluorononanoic acid										
463.00 > 419.00	3.258	3.275	-0.017	1.000	7994712	49.7		99.4	111942	
D 19 13C5 PFNA										
468.00 > 423.00	3.258	3.275	-0.017		8110550	45.5		91.1	448870	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.579	3.589	-0.010	1.000	17622504	50.8		102	244916	
D 21 13C8 FOSA										
506.00 > 78.00	3.579	3.589	-0.010		18539988	44.8		89.5	782482	
D 23 13C2 PFDA										
515.00 > 470.00	3.621	3.631	-0.010		7573901	45.6		91.2	477033	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.621	3.631	-0.010	1.000	7456893	50.8		102	159874	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.929	3.941	-0.012	1.000	7073516	48.7		101		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.947	3.959	-0.012	1.000	5654902	47.3		94.6	116847	
D 27 13C2 PFUnA										
565.00 > 520.00	3.947	3.967	-0.020		5639344	44.8		89.7	266119	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.243	4.259	-0.016	1.000	4947736	48.1		96.2	64070	
D 30 13C2 PFDaA										
615.00 > 570.00	4.243	4.259	-0.016		5435760	46.5		93.0	206048	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.511	4.520	-0.009	1.000	5155981	45.6		91.2	17473	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.750	4.772	-0.022	1.000	9673132	46.0		92.1	3806	
713.00 > 169.00	4.742	4.772	-0.030	0.998	1561384		6.20(0.00-0.00)		311529	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.750	4.772	-0.022		11021074	45.6		91.1	497299	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.167	5.184	-0.017		5938609	45.6		91.1	101514	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.177	5.194	-0.017	1.000	5386746	49.5		99.0	4133	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.540	5.562	-0.022	1.000	5609402	65.6		131	7160	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_022.d

Injection Date: 02-Dec-2016 17:37:22

Instrument ID: A8_N

Lims ID: CCV L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 26

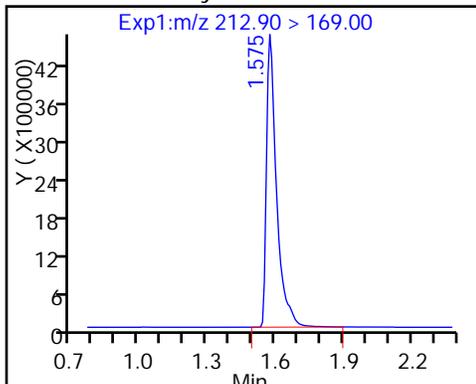
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

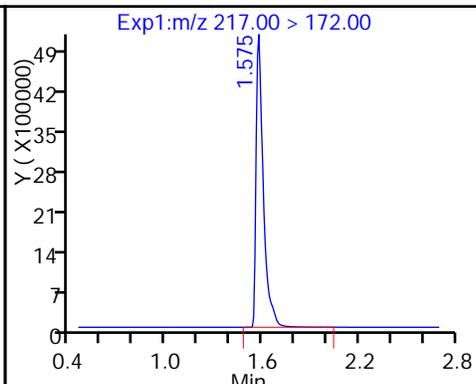
Method: A8_N

Limit Group: LC PFC_DOD ICAL

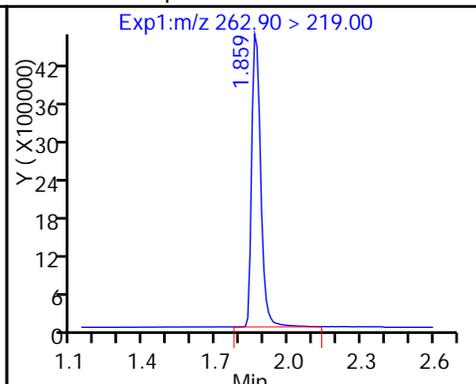
1 Perfluorobutyric acid



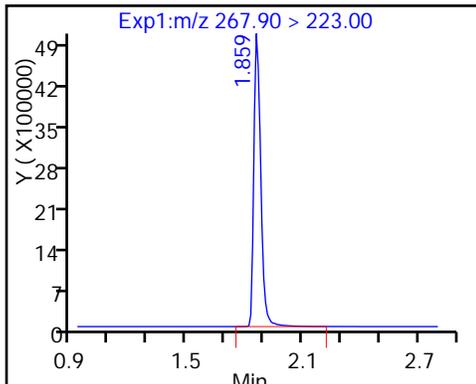
D 2 13C4 PFBA



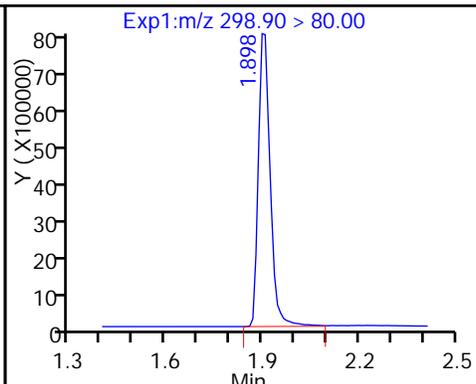
3 Perfluoropentanoic acid



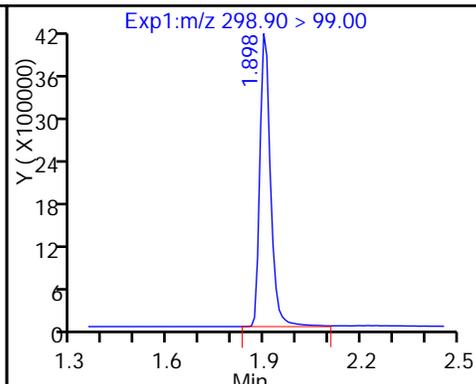
D 4 13C5-PFPeA



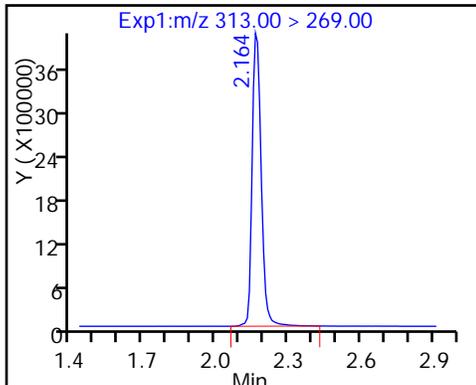
5 Perfluorobutanesulfonic acid



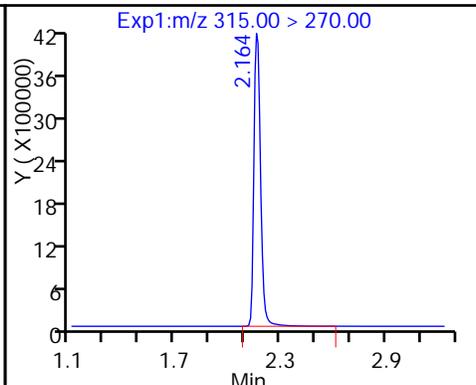
5 Perfluorobutanesulfonic acid



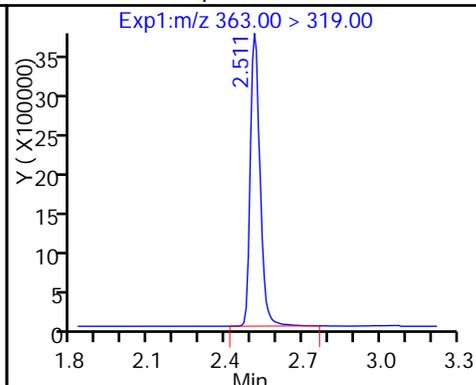
7 Perfluorohexanoic acid



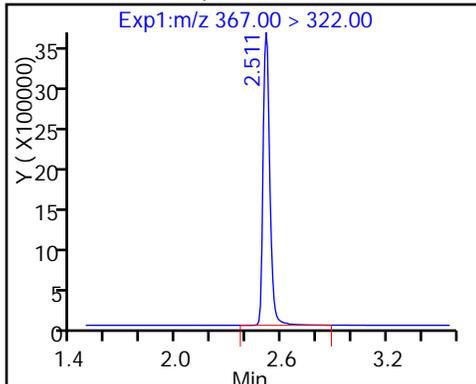
D 6 13C2 PFHxA



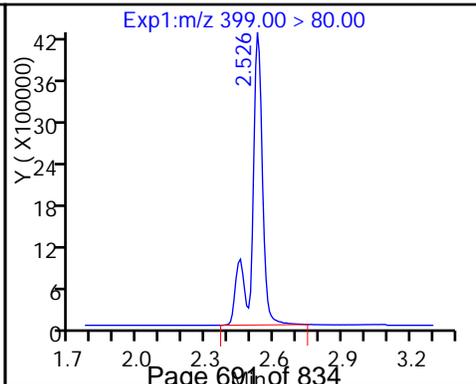
12 Perfluoroheptanoic acid



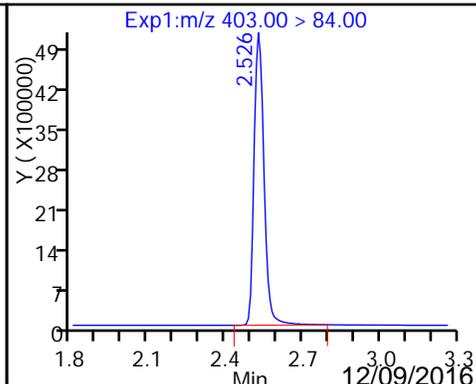
D 11 13C4-PFHpA



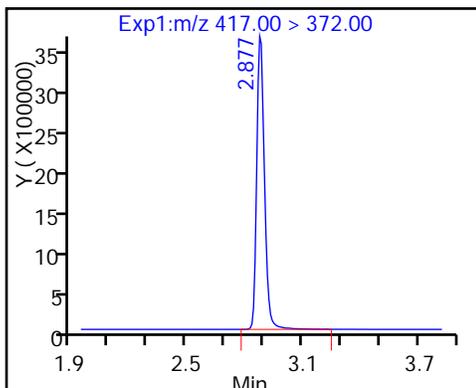
9 Perfluorohexanesulfonic acid (M)



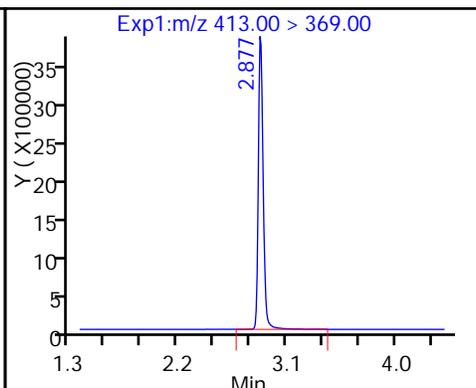
D 10 18O2 PFHxS



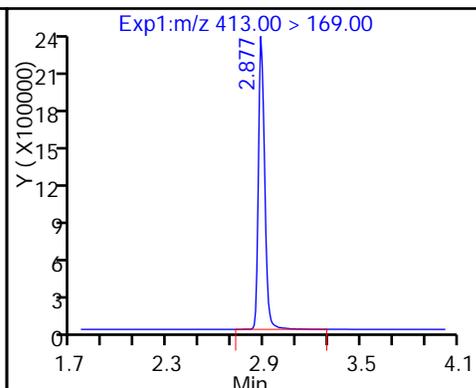
D 14 13C4 PFOA



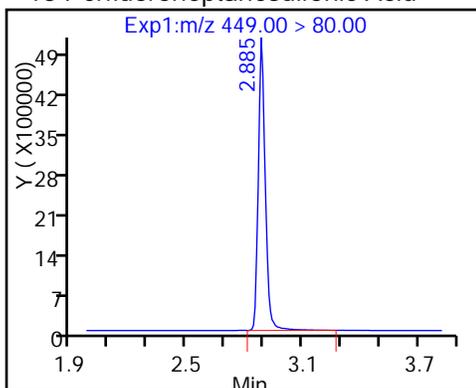
15 Perfluorooctanoic acid



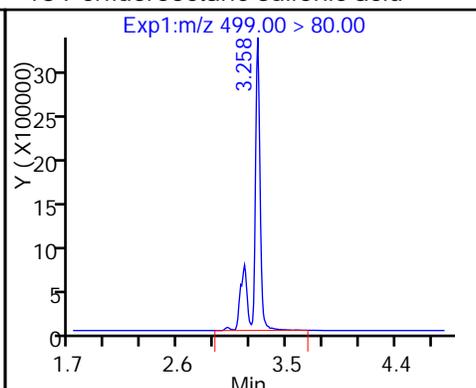
15 Perfluorooctanoic acid



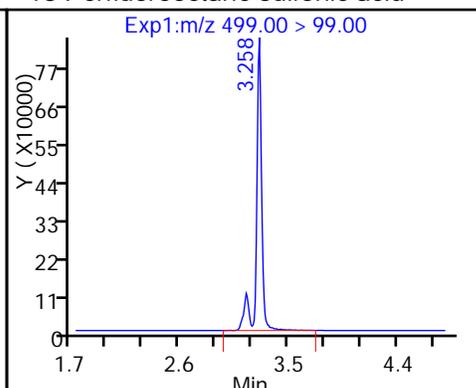
13 Perfluoroheptanesulfonic Acid



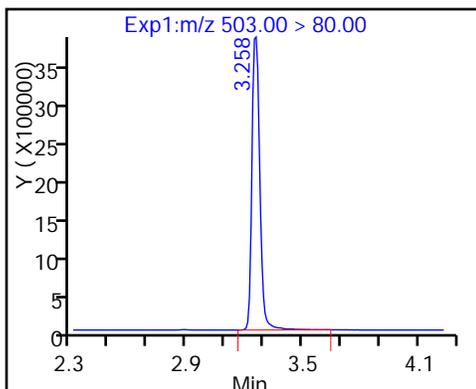
18 Perfluorooctane sulfonic acid



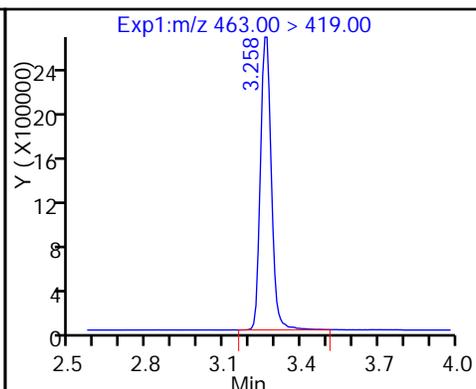
18 Perfluorooctane sulfonic acid



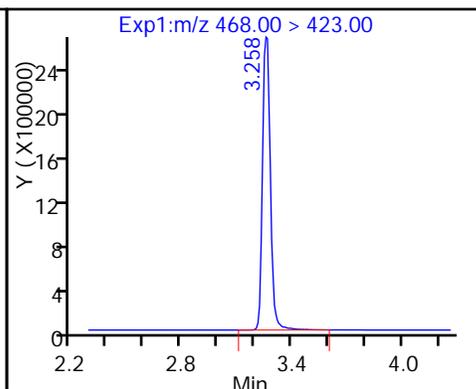
D 17 13C4 PFOS



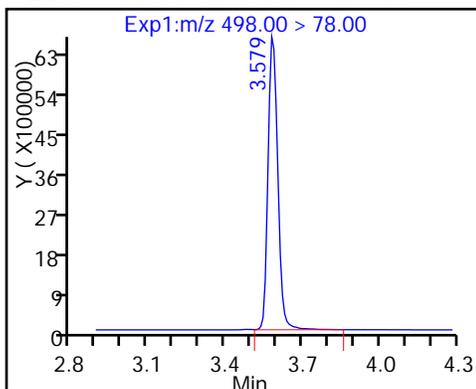
20 Perfluorononanoic acid



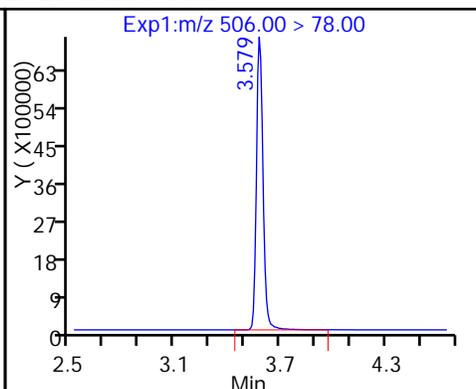
D 19 13C5 PFNA



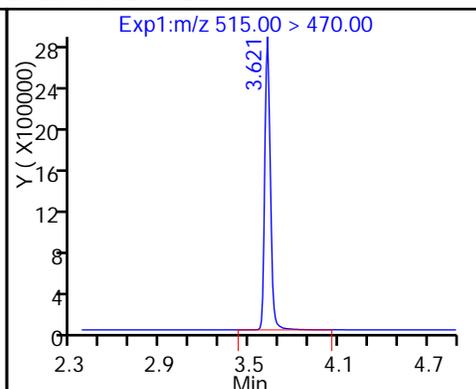
22 Perfluorooctane Sulfonamide

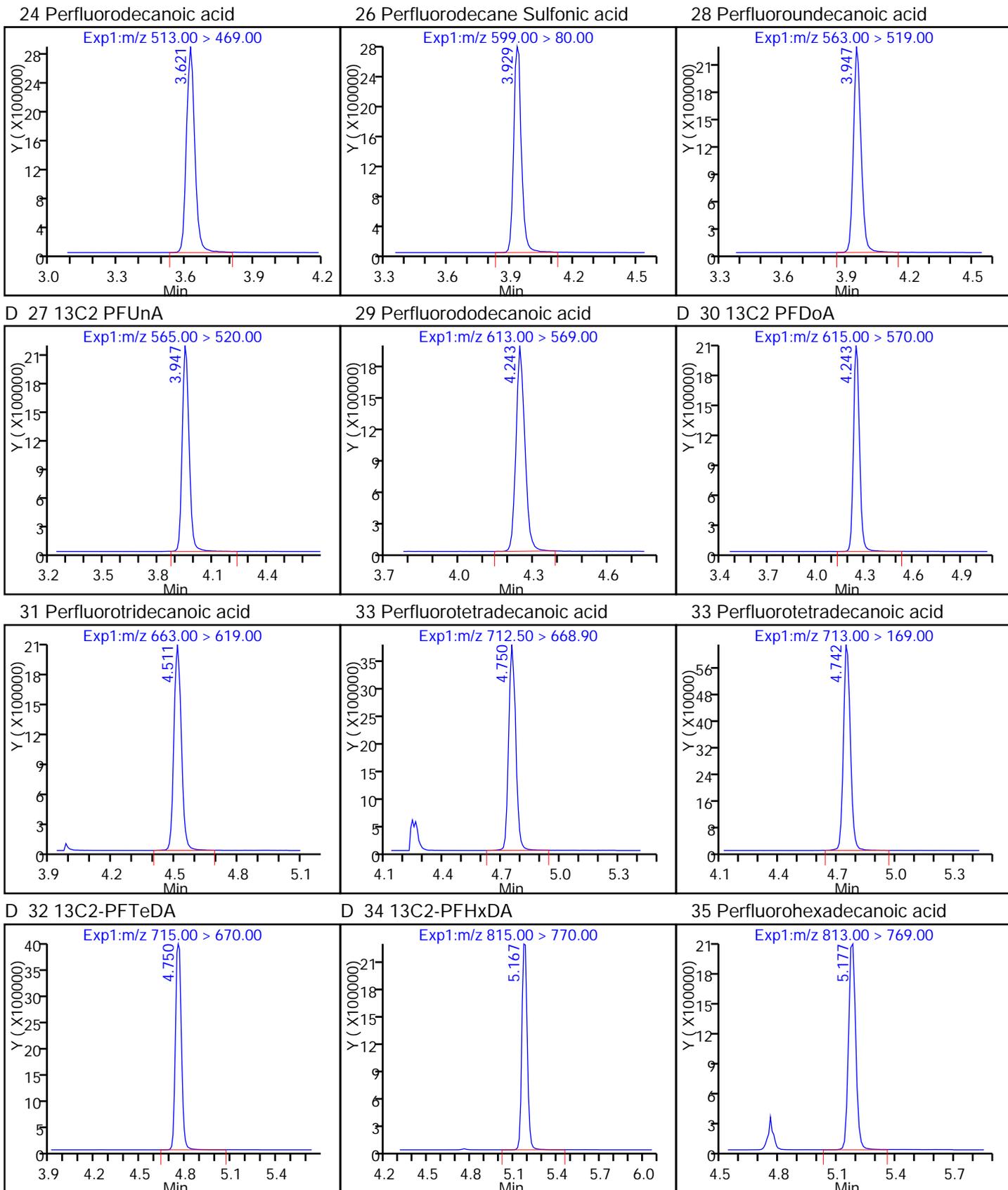


D 21 13C8 FOSA

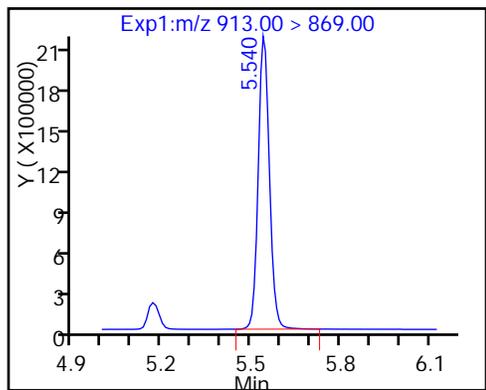


D 23 13C2 PFDA





36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: ICV 320-140564/11 Calibration Date: 12/03/2016 14:41
 Instrument ID: A8_N Calib Start Date: 12/03/2016 13:48
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/03/2016 15:33
 Lab File ID: 03DEC2016A_011.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.596	1.545		42.8	44.3	-3.2	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.072	0.9942		46.4	50.0	-7.2	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.085	0.9006		39.6	47.8	-17.0	25.0
13C4 PFBA	Ave	335829	312410		46.5	50.0	-7.0	50.0
13C5-PFPeA	Ave	264545	243307		46.0	50.0	-8.0	50.0
13C2 PFHxA	Ave	237486	229267		48.3	50.0	-3.5	50.0
13C4-PFHpA	Ave	207413	195333		47.1	50.0	-5.8	50.0
18O2 PFHxS	Ave	312342	294291		44.6	47.3	-5.8	50.0
13C4 PFOA	Ave	219258	198224		45.2	50.0	-9.6	50.0
13C4 PFOS	Ave	246009	225320		43.8	47.8	-8.4	50.0
13C5 PFNA	Ave	166415	156724		47.1	50.0	-5.8	50.0
13C8 FOSA	Ave	402279	382595		47.6	50.0	-4.9	50.0
13C2 PFDA	Ave	157817	146344		46.4	50.0	-7.3	50.0
13C2 PFUnA	Ave	118762	109554		46.1	50.0	-7.8	50.0
13C2 PFDoA	Ave	112084	109278		48.7	50.0	-2.5	50.0
13C2-PFTeDA	Ave	231173	214390		46.4	50.0	-7.3	50.0
13C2-PFHxDA	Ave	129725	121978		47.0	50.0	-6.0	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_011.d
 Lims ID: ICV
 Client ID:
 Sample Type: ICV
 Inject. Date: 03-Dec-2016 14:41:13 ALS Bottle#: 44 Worklist Smp#: 11
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: ICV_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist:

Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:26:23 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:47:22

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.566	1.574	-0.008	15620513	46.5		93.0	1612048	
1 Perfluorobutyric acid	212.90 > 169.00	1.566	1.577	-0.011	13190679	48.3			70053	
3 Perfluoropentanoic acid	262.90 > 219.00	1.848	1.861	-0.013	11502032	46.6			76938	
D 4 13C5-PFPeA	267.90 > 223.00	1.848	1.861	-0.013	12165345	46.0		92.0	911947	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.887	1.900	-0.013	20121282	42.8				
	298.90 > 99.00	1.887	1.900	-0.013	9272505		2.17(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.158	2.164	-0.006	10017573	45.8			244705	
D 6 13C2 PFHxA	315.00 > 270.00	2.158	2.164	-0.006	11463370	48.3		96.5	672140	
D 11 13C4-PFHpA	367.00 > 322.00	2.499	2.511	-0.012	9766668	47.1		94.2	852072	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.499	2.512	-0.013	9294490	46.3			100010	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.514	2.531	-0.017	13456842	41.7				
D 10 18O2 PFHxS	403.00 > 84.00	2.514	2.531	-0.017	13919947	44.6		94.2	810316	
D 14 13C4 PFOA	417.00 > 372.00	2.870	2.880	-0.010	9911203	45.2		90.4	548349	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.870	2.887	-0.017	1.000	9853753	46.4			123535	
413.00 > 169.00	2.862	2.887	-0.025	0.997	6029090		1.63(0.90-1.10)		244064	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.870	2.888	-0.018	1.000	13003397	49.0				
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.244	3.258	-0.014	1.000	9689735	39.6			1641854	
499.00 > 99.00	3.235	3.258	-0.023	0.997	2361166		4.10(0.90-1.10)		256440	
D 17 13C4 PFOS										
503.00 > 80.00	3.235	3.259	-0.024		10770309	43.8		91.6	336185	
D 19 13C5 PFNA										
468.00 > 423.00	3.244	3.263	-0.019		7836220	47.1		94.2	638662	
20 Perfluorononanoic acid										
463.00 > 419.00	3.244	3.263	-0.019	1.000	7315987	46.9			101636	
D 21 13C8 FOSA										
506.00 > 78.00	3.557	3.571	-0.014		19129770	47.6		95.1	889404	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.565	3.574	-0.009	1.000	16827146	47.1			307442	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.607	3.623	-0.016	1.000	6727679	47.9			161083	
D 23 13C2 PFDA										
515.00 > 470.00	3.607	3.626	-0.019		7317194	46.4		92.7	293278	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.912	3.936	-0.024	1.000	6797886	47.2				
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.938	3.955	-0.017	1.000	5226274	44.8			120213	
D 27 13C2 PFUnA										
565.00 > 520.00	3.938	3.958	-0.020		5477711	46.1		92.2	312875	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.224	4.250	-0.026	1.000	4931801	47.6			90828	
D 30 13C2 PFDoA										
615.00 > 570.00	4.232	4.251	-0.019		5463889	48.7		97.5	309492	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.503	4.518	-0.015	1.000	4773745	46.0			36970	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.732	4.759	-0.027		10719523	46.4		92.7	360593	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.740	4.761	-0.021	1.000	9370597	46.3			4744	
713.00 > 169.00	4.732	4.761	-0.029	0.998	1459995		6.42(0.00-0.00)		102221	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.158	5.186	-0.028		6098899	47.0		94.0	113553	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.158	5.186	-0.028	1.000	5084325	46.1			4219	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.523	5.559	-0.036	1.000	5482963	50.5			9364	

Reagents:

LCPFCIC_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_011.d

Injection Date: 03-Dec-2016 14:41:13

Instrument ID: A8_N

Lims ID: ICV

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 44

Worklist Smp#: 11

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

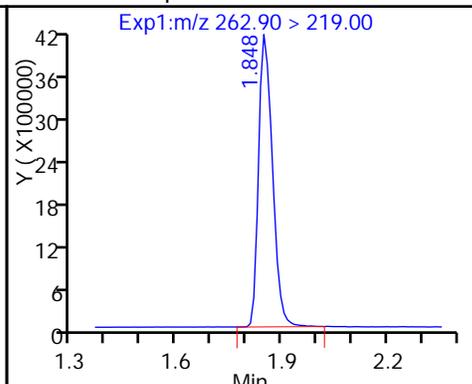
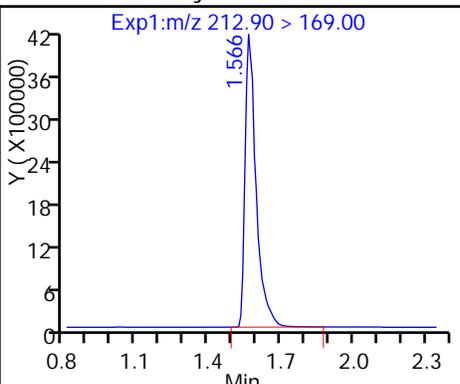
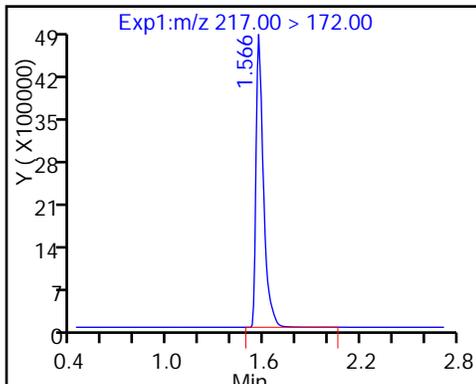
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

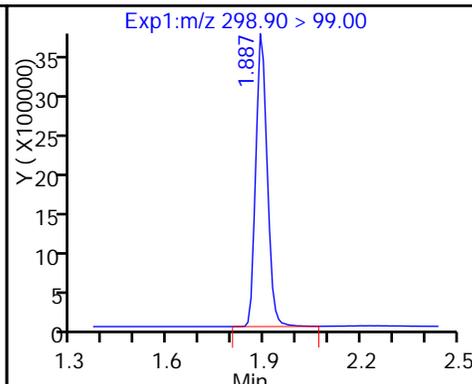
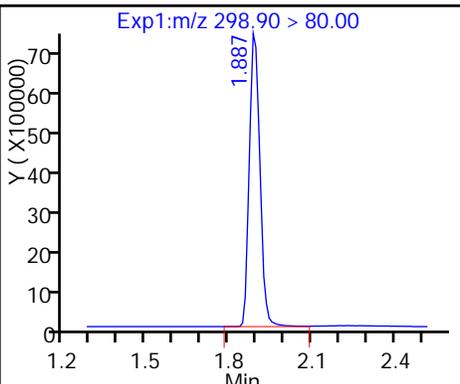
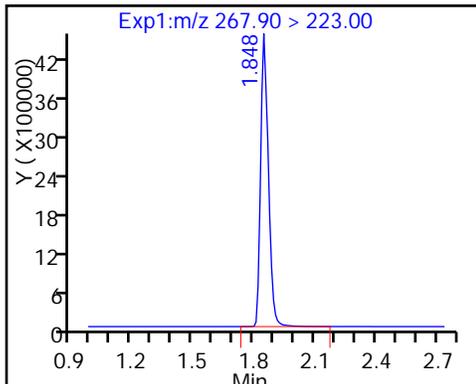
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

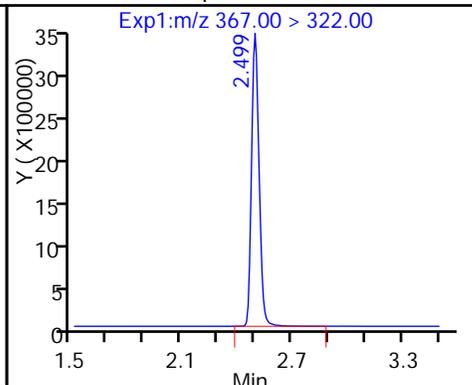
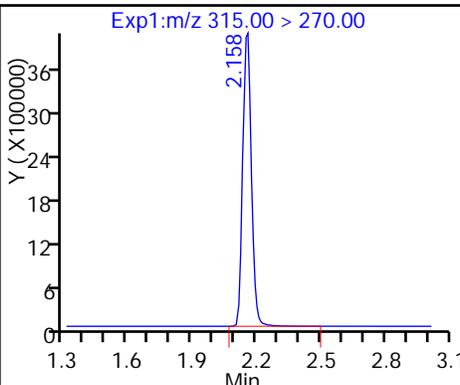
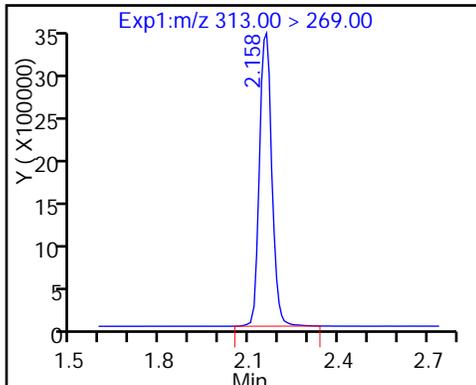
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

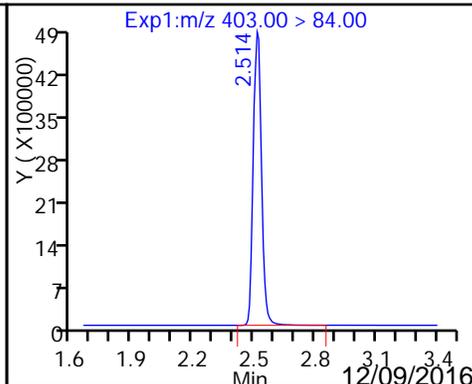
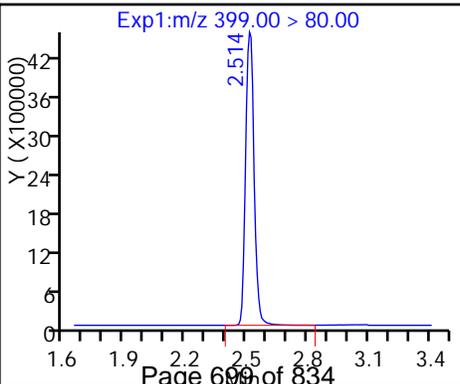
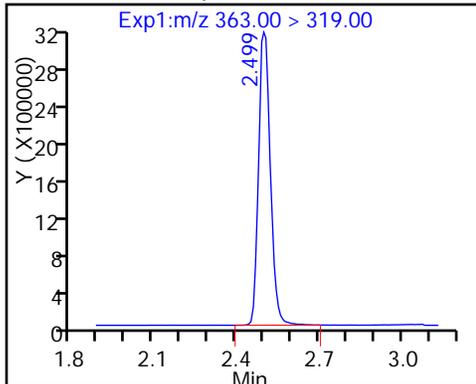
D 11 13C4-PFHpA



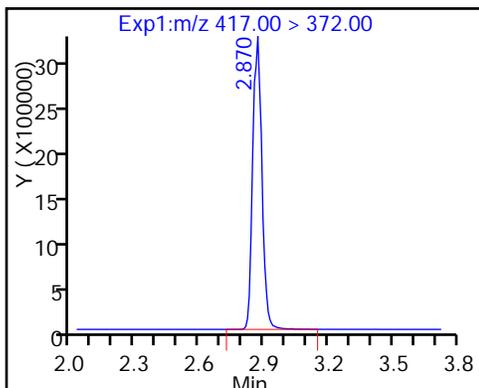
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

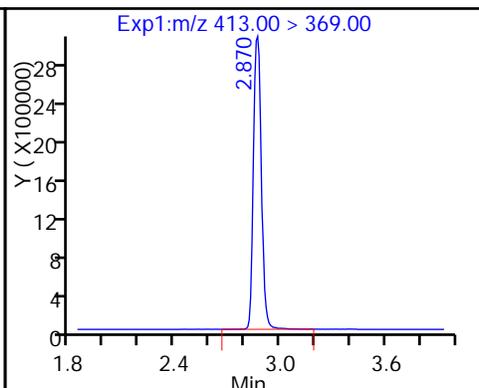
D 10 18O2 PFHxS



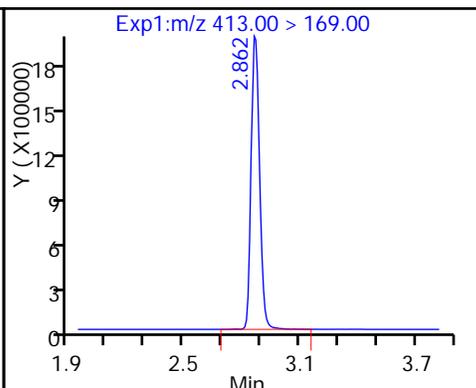
D 14 13C4 PFOA



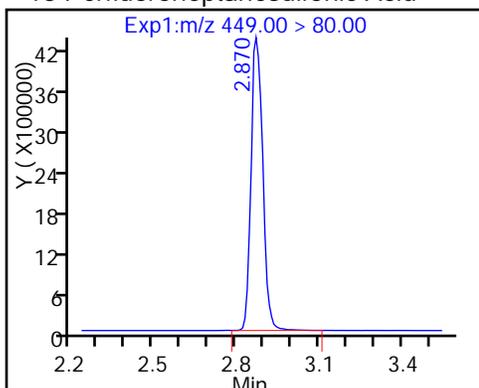
15 Perfluorooctanoic acid



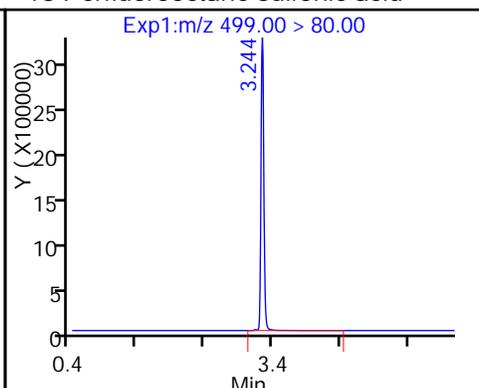
15 Perfluorooctanoic acid



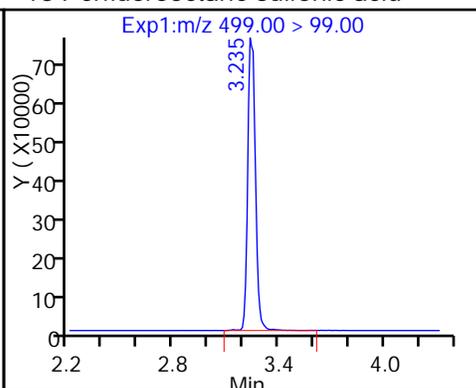
13 Perfluoroheptanesulfonic Acid



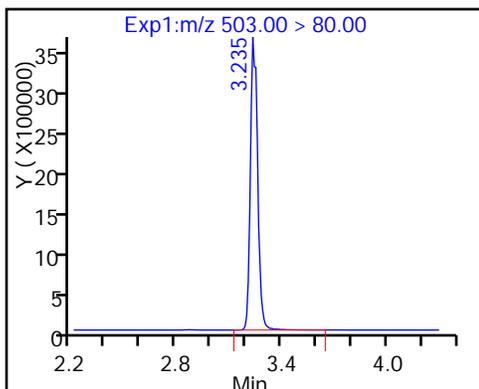
18 Perfluorooctane sulfonic acid



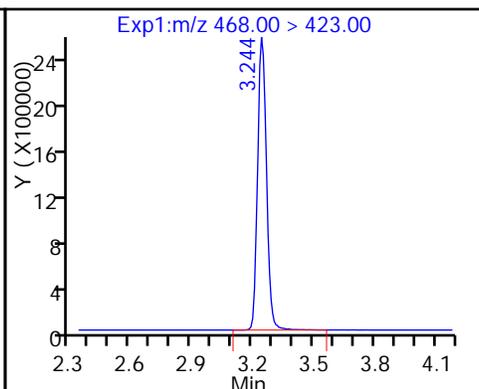
18 Perfluorooctane sulfonic acid



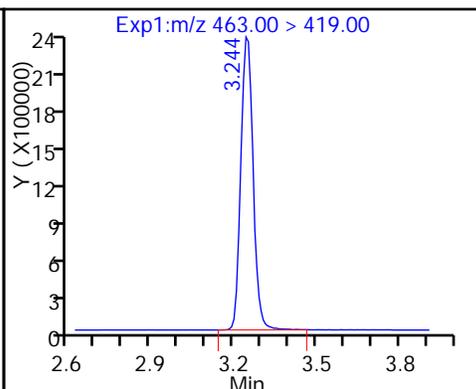
D 17 13C4 PFOS



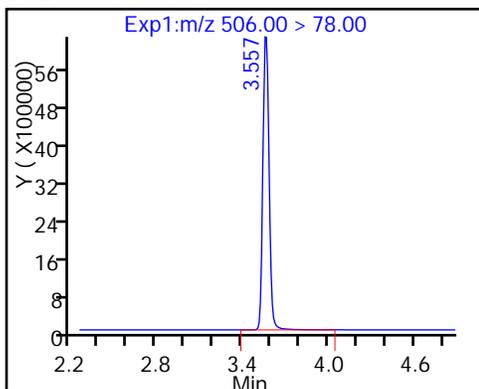
D 19 13C5 PFNA



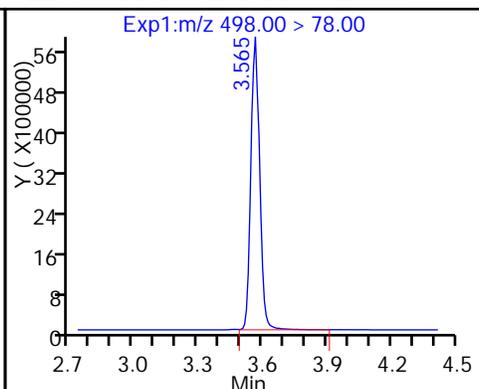
20 Perfluorononanoic acid



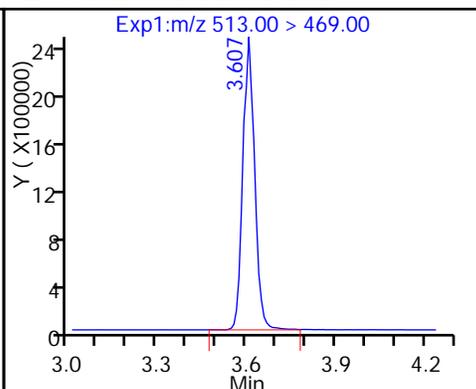
D 21 13C8 FOSA



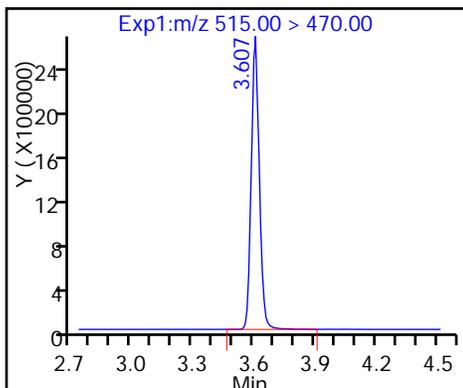
22 Perfluorooctane Sulfonamide



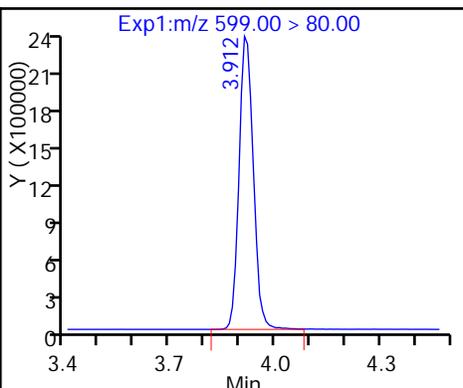
24 Perfluorodecanoic acid



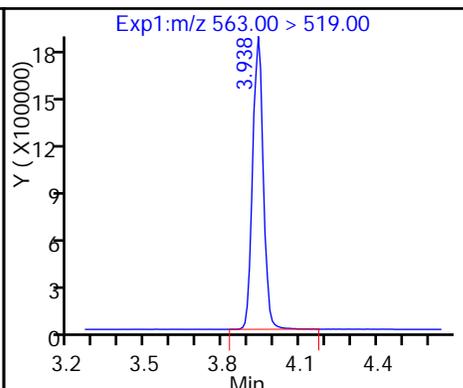
D 23 13C2 PFDA



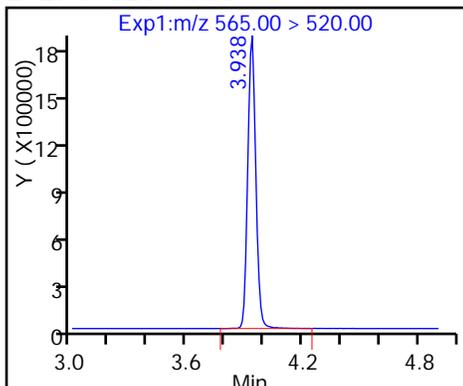
26 Perfluorodecane Sulfonic acid



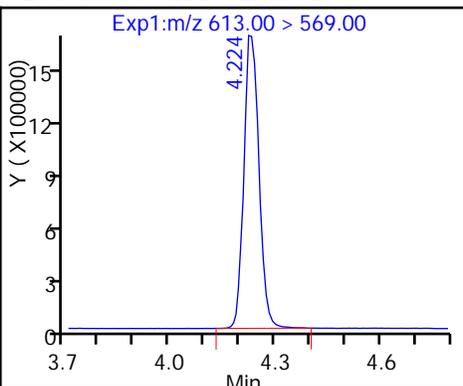
28 Perfluoroundecanoic acid



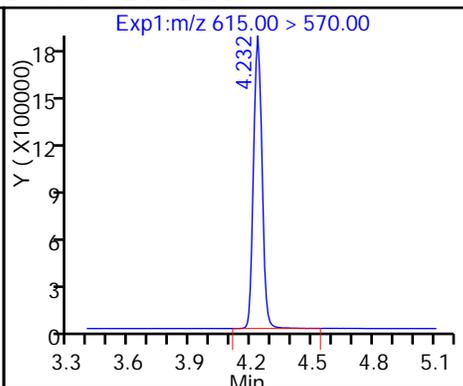
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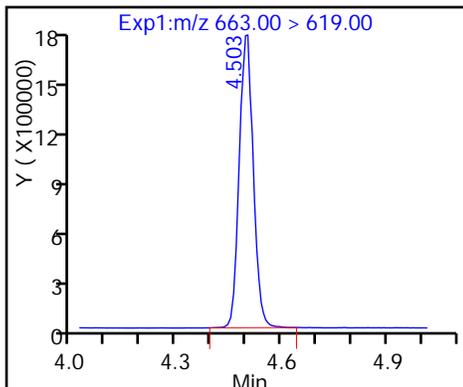
29 Perfluorododecanoic acid



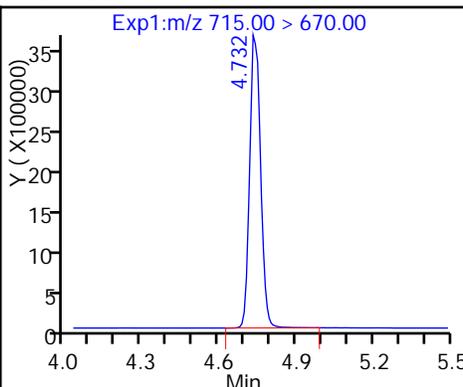
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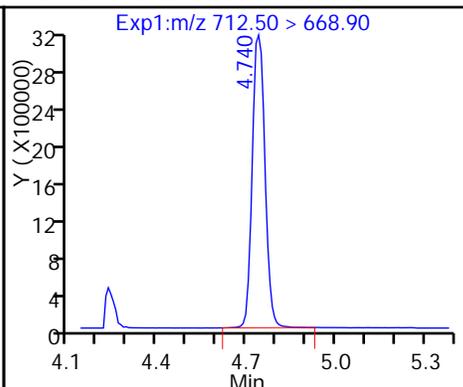
31 Perfluorotridecanoic acid



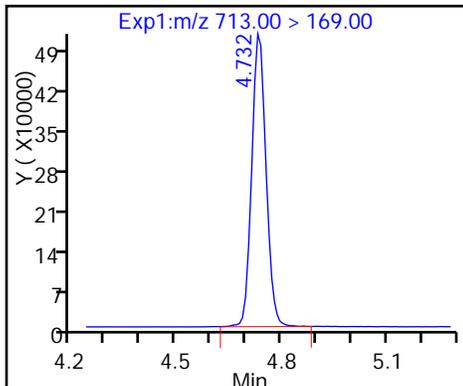
D 32 13C2-PFTeDA



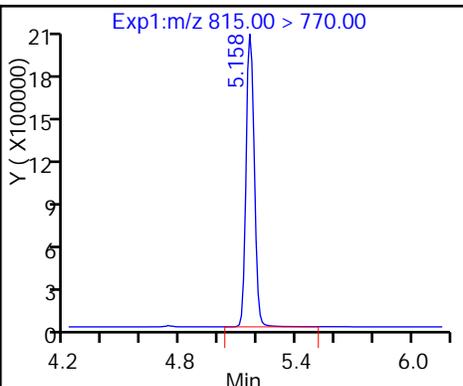
33 Perfluorotetradecanoic acid



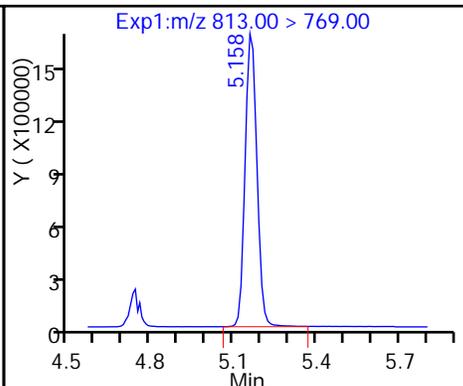
33 Perfluorotetradecanoic acid



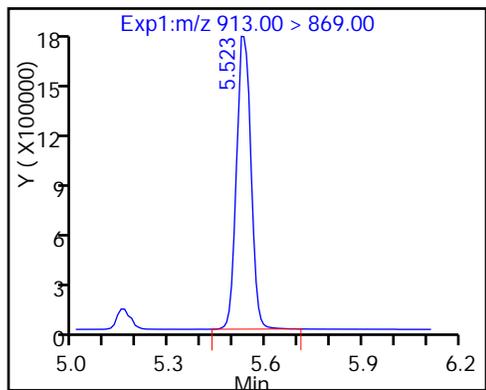
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-140564/30 Calibration Date: 12/03/2016 17:03
 Instrument ID: A8_N Calib Start Date: 12/03/2016 13:48
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/03/2016 15:33
 Lab File ID: 03DEC2016B_009.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.596	1.566		43.4	44.2	-1.9	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.072	1.016		47.4	50.0	-5.2	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.085	1.059		45.3	46.4	-2.4	25.0
13C4 PFBA	Ave	335829	313361		46.7	50.0	-6.7	50.0
13C5-PFPeA	Ave	264545	244850		46.3	50.0	-7.4	50.0
13C2 PFHxA	Ave	237486	217160		45.7	50.0	-8.6	50.0
13C4-PFHpA	Ave	207413	186664		45.0	50.0	-10.0	50.0
18O2 PFHxS	Ave	312342	300362		45.5	47.3	-3.8	50.0
13C4 PFOA	Ave	219258	202399		46.2	50.0	-7.7	50.0
13C4 PFOS	Ave	246009	229171		44.5	47.8	-6.8	50.0
13C5 PFNA	Ave	166415	153229		46.0	50.0	-7.9	50.0
13C8 FOSA	Ave	402279	368306		45.8	50.0	-8.4	50.0
13C2 PFDA	Ave	157817	144055		45.6	50.0	-8.7	50.0
13C2 PFUnA	Ave	118762	107655		45.3	50.0	-9.4	50.0
13C2 PFDoA	Ave	112084	107351		47.9	50.0	-4.2	50.0
13C2-PFTeDA	Ave	231173	206686		44.7	50.0	-10.6	50.0
13C2-PFHxDA	Ave	129725	113111		43.6	50.0	-12.8	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_009.d
 Lims ID: CCV L5
 Client ID:
 Sample Type: CCV
 Inject. Date: 03-Dec-2016 17:03:44 ALS Bottle#: 41 Worklist Smp#: 30
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L5
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:37:10 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:37:10

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.566	1.574	-0.008	15668067	46.7		93.3	1212976	
1 Perfluorobutyric acid	212.90 > 169.00	1.566	1.577	-0.011	13737697	50.2		100	89841	
3 Perfluoropentanoic acid	262.90 > 219.00	1.849	1.861	-0.012	12128535	48.8		97.6	97785	
D 4 13C5-PFPeA	267.90 > 223.00	1.849	1.861	-0.012	12242489	46.3		92.6	1494339	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.887	1.900	-0.013	20785751	43.4		98.1		
	298.90 > 99.00	1.887	1.900	-0.013	9458185		2.20(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.143	2.164	-0.021	10257258	49.6		99.1	206256	
D 6 13C2 PFHxA	315.00 > 270.00	2.143	2.164	-0.021	10858019	45.7		91.4	619877	
D 11 13C4-PFHpA	367.00 > 322.00	2.489	2.511	-0.022	9333205	45.0		90.0	537737	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.489	2.512	-0.023	9428316	49.2		98.4	112468	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.497	2.531	-0.034	14202827	43.1		94.7		
D 10 18O2 PFHxS	403.00 > 84.00	2.497	2.531	-0.034	14207124	45.5		96.2	1003243	
D 14 13C4 PFOA	417.00 > 372.00	2.854	2.880	-0.026	10119935	46.2		92.3	562804	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.854	2.887	-0.033	1.000	10284066	47.4		94.8	202666	
413.00 > 169.00	2.846	2.887	-0.041	0.997	6248942		1.65(0.90-1.10)		339460	
13 Perfluoroheptanesulfonic Acid										
449.00 > 80.00	2.862	2.888	-0.026	1.000	12744705	47.2		99.3		
18 Perfluorooctane sulfonic acid										
499.00 > 80.00	3.226	3.258	-0.032	1.000	11265789	45.3		97.6	979499	
499.00 > 99.00	3.226	3.258	-0.032	1.000	2568331		4.39(0.90-1.10)		255539	
D 17 13C4 PFOS										
503.00 > 80.00	3.217	3.259	-0.042		10954381	44.5		93.2	236210	
D 19 13C5 PFNA										
468.00 > 423.00	3.226	3.263	-0.037		7661469	46.0		92.1	827542	
20 Perfluorononanoic acid										
463.00 > 419.00	3.226	3.263	-0.037	1.000	7595586	49.8		99.5	106584	
D 21 13C8 FOSA										
506.00 > 78.00	3.548	3.571	-0.023		18415295	45.8		91.6	652368	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.548	3.574	-0.026	1.000	16711764	48.6		97.2	326893	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.582	3.623	-0.041	1.000	6874163	49.7		99.4	240558	
D 23 13C2 PFDA										
515.00 > 470.00	3.582	3.626	-0.044		7202726	45.6		91.3	250402	
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.893	3.936	-0.043	1.000	6912445	47.1		97.8		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.919	3.955	-0.036	1.000	5278542	46.0		92.0	157895	
D 27 13C2 PFUnA										
565.00 > 520.00	3.919	3.958	-0.039		5382740	45.3		90.6	645693	
29 Perfluorododecanoic acid										
613.00 > 569.00	4.212	4.250	-0.038	1.000	5011838	49.2		98.4	94279	
D 30 13C2 PFDaA										
615.00 > 570.00	4.212	4.251	-0.039		5367554	47.9		95.8	118003	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.484	4.518	-0.034	1.000	4658317	45.7		91.4	52356	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.714	4.759	-0.045		10334299	44.7		89.4	298039	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.714	4.761	-0.047	1.000	8894154	44.7		89.4	6477	
713.00 > 169.00	4.714	4.761	-0.047	1.000	1446397		6.15(0.00-0.00)		164686	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.132	5.186	-0.054		5655535	43.6		87.2	107822	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.132	5.186	-0.054	1.000	5063795	46.7		93.5	6291	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.494	5.559	-0.065	1.000	5010861	47.0		94.0	8065	

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_009.d

Injection Date: 03-Dec-2016 17:03:44

Instrument ID: A8_N

Lims ID: CCV L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 30

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

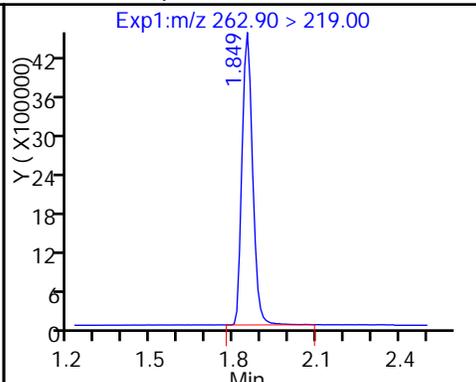
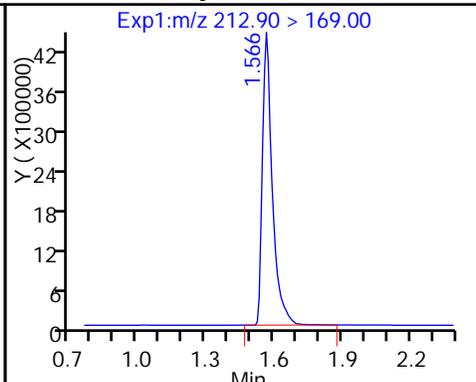
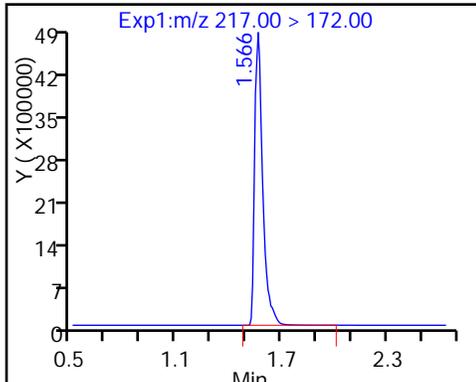
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

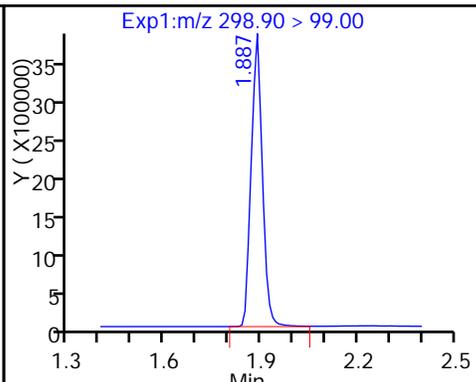
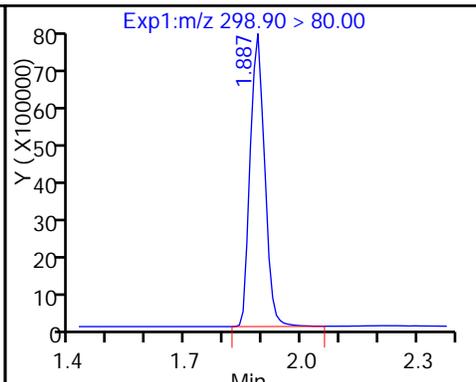
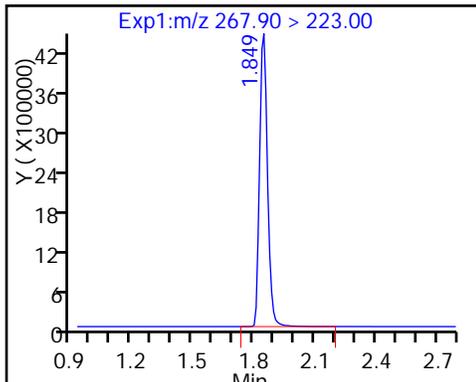
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

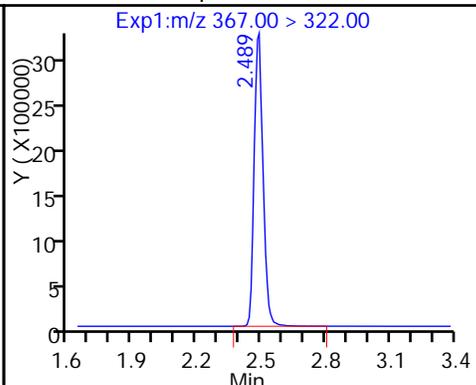
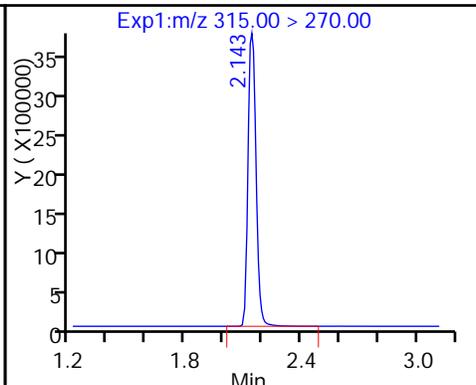
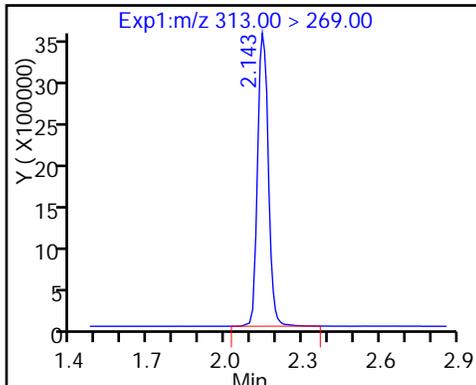
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

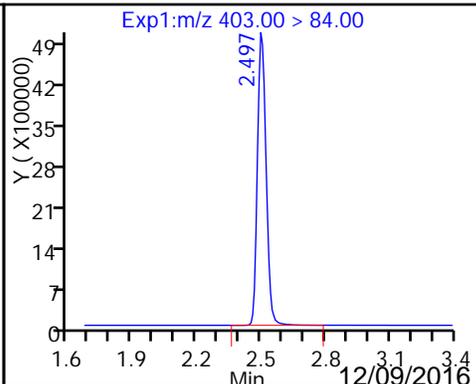
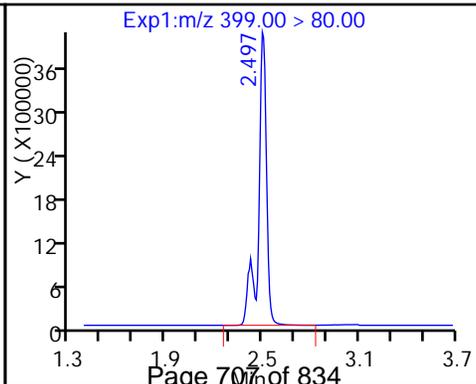
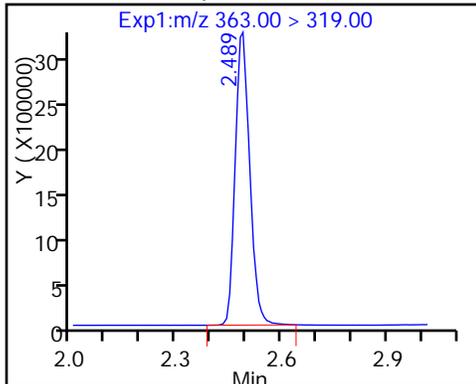
D 11 13C4-PFHpA



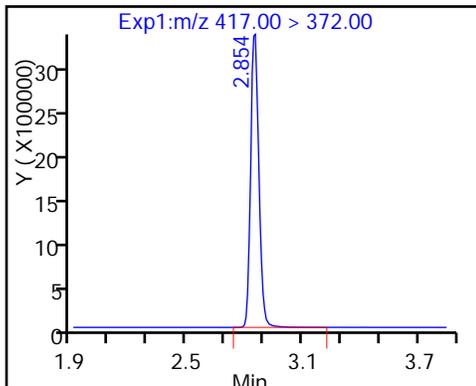
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

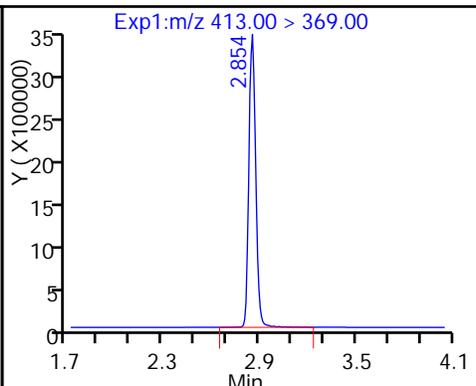
D 10 18O2 PFHxS



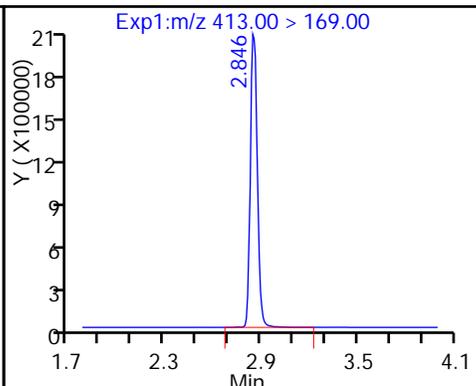
D 14 13C4 PFOA



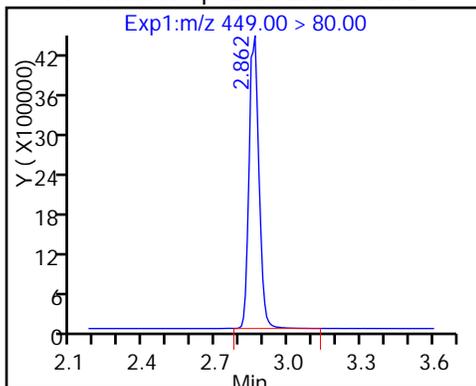
15 Perfluorooctanoic acid



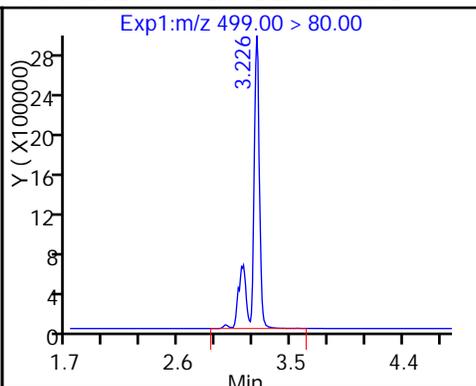
15 Perfluorooctanoic acid



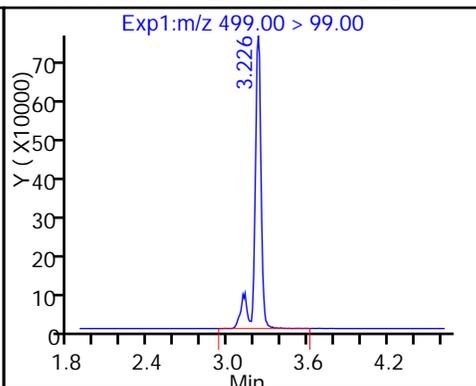
13 Perfluoroheptanesulfonic Acid



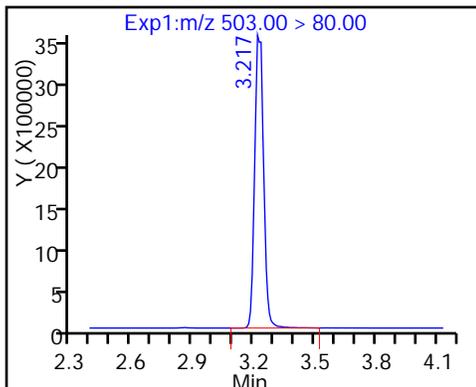
18 Perfluorooctane sulfonic acid



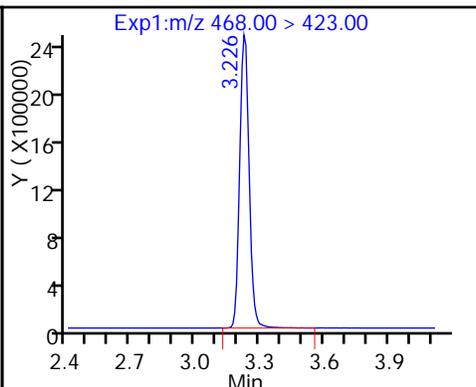
18 Perfluorooctane sulfonic acid



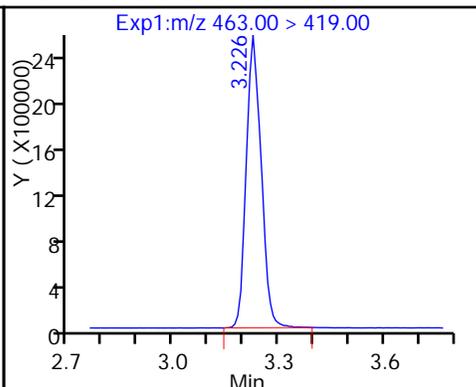
D 17 13C4 PFOS



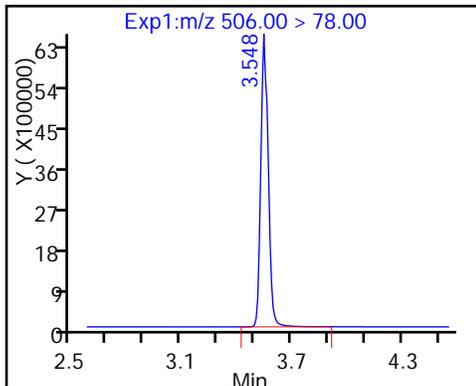
D 19 13C5 PFNA



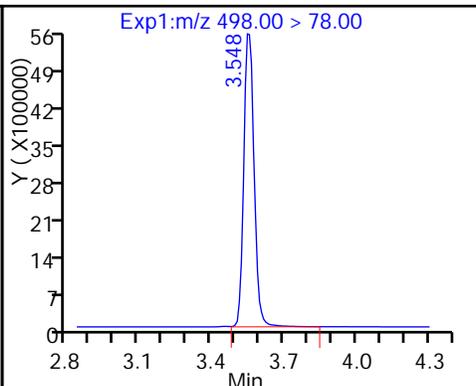
20 Perfluorononanoic acid



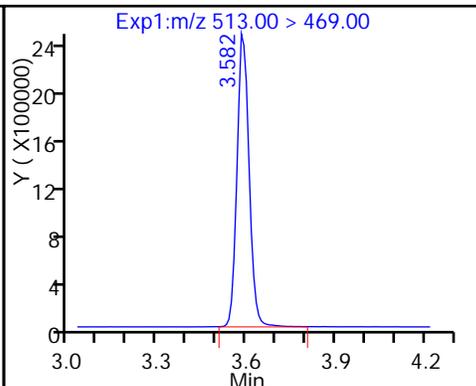
D 21 13C8 FOSA



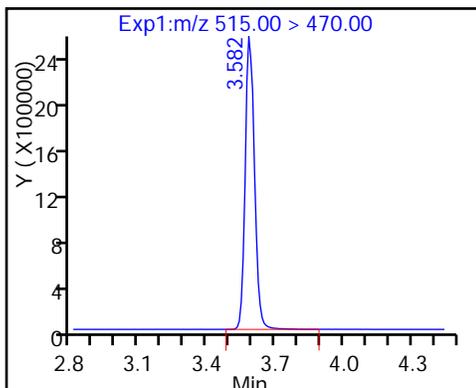
22 Perfluorooctane Sulfonamide



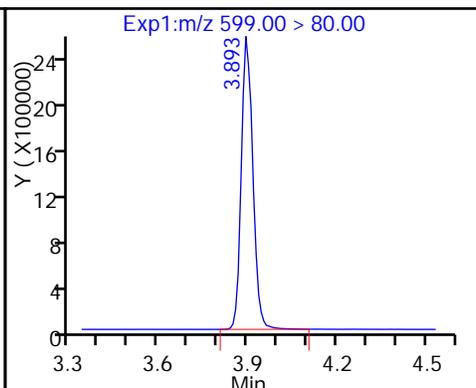
24 Perfluorodecanoic acid



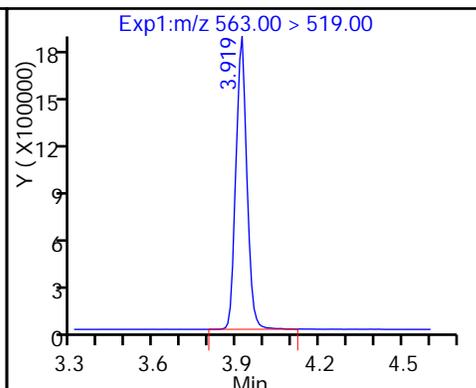
D 23 13C2 PFDA



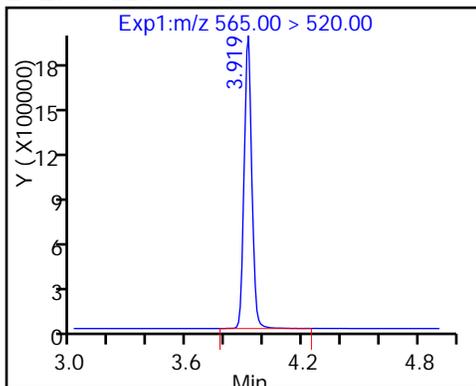
26 Perfluorodecane Sulfonic acid



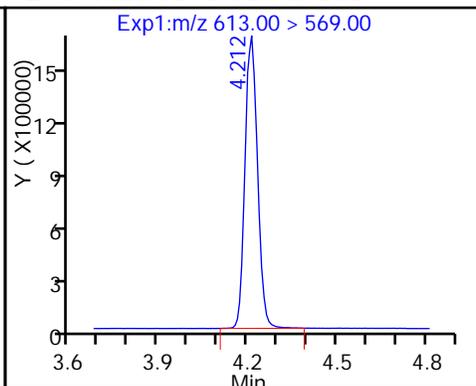
28 Perfluoroundecanoic acid



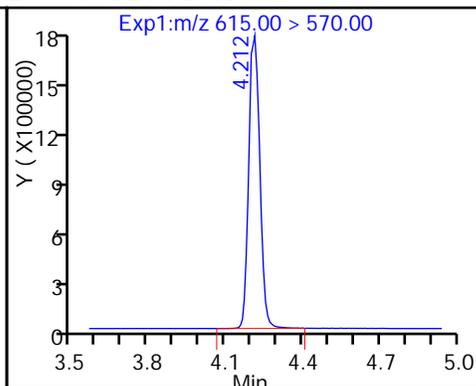
D 27 13C2 PFUa



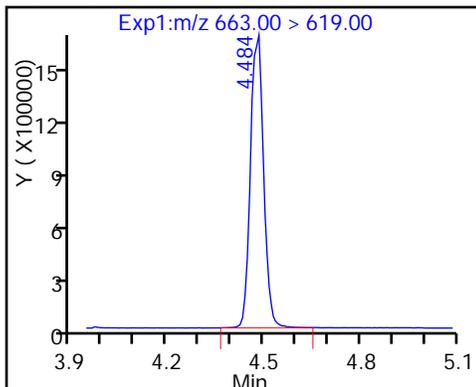
29 Perfluorododecanoic acid



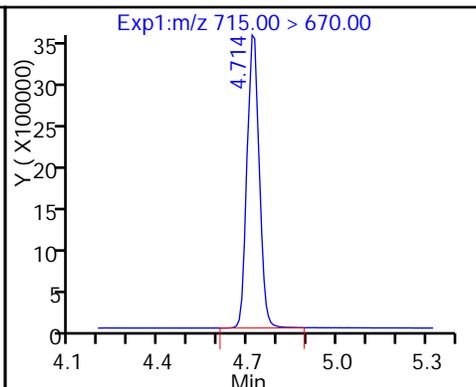
D 30 13C2 PFDa



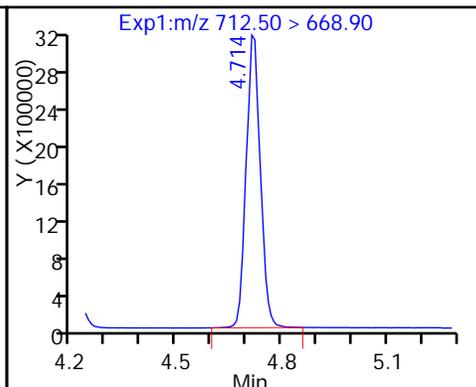
31 Perfluorotridecanoic acid



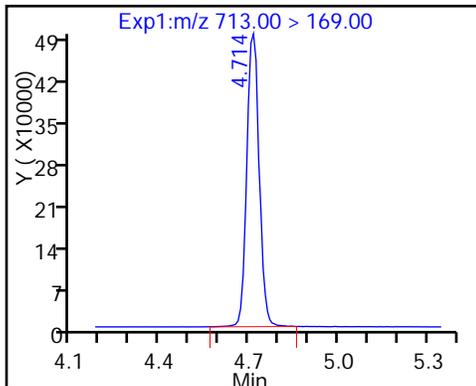
D 32 13C2-PFTeDA



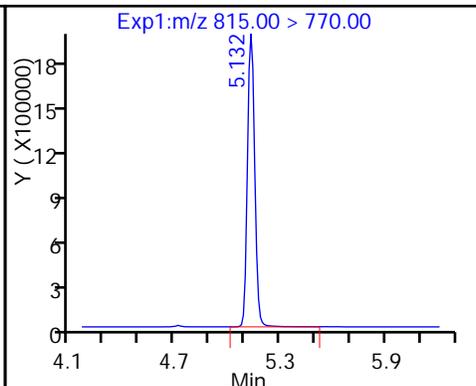
33 Perfluorotetradecanoic acid



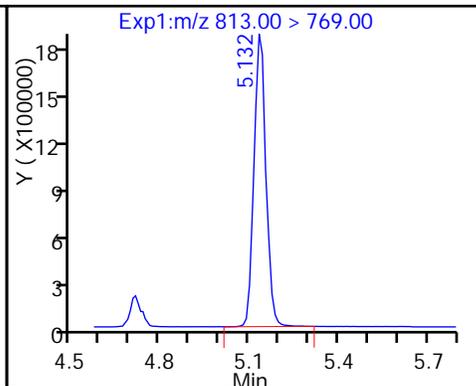
33 Perfluorotetradecanoic acid



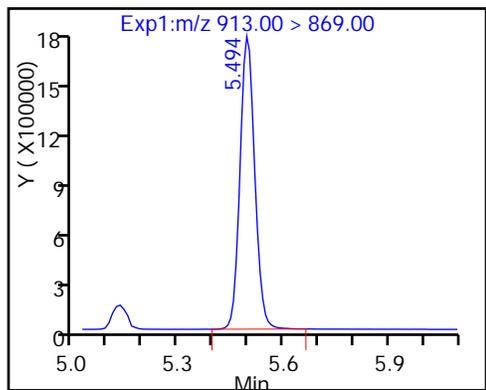
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-141054/5 Calibration Date: 12/07/2016 12:33
 Instrument ID: A8_N Calib Start Date: 12/03/2016 13:48
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/03/2016 15:33
 Lab File ID: 07DEC2016A_003.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.596	1.541		0.854	0.884	-3.4	50.0
Perfluorooctanoic acid (PFOA)	AveID	1.072	1.129		1.05	1.00	5.3	50.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.085	1.033		0.883	0.928	-4.8	50.0
13C4 PFBA	Ave	335829	347833		51.8	50.0	3.6	50.0
13C5-PFPeA	Ave	264545	268424		50.7	50.0	1.5	50.0
13C2 PFHxA	Ave	237486	238270		50.2	50.0	0.3	50.0
13C4-PFHpA	Ave	207413	222068		53.5	50.0	7.1	50.0
18O2 PFHxS	Ave	312342	329175		49.8	47.3	5.4	50.0
13C4 PFOA	Ave	219258	238888		54.5	50.0	9.0	50.0
13C4 PFOS	Ave	246009	260205		50.6	47.8	5.8	50.0
13C5 PFNA	Ave	166415	172520		51.8	50.0	3.7	50.0
13C8 FOSA	Ave	402279	413977		51.5	50.0	2.9	50.0
13C2 PFDA	Ave	157817	150832		47.8	50.0	-4.4	50.0
13C2 PFUnA	Ave	118762	116003		48.8	50.0	-2.3	50.0
13C2 PFDoA	Ave	112084	106415		47.5	50.0	-5.1	50.0
13C2-PFTeDA	Ave	231173	215301		46.6	50.0	-6.9	50.0
13C2-PFHxDA	Ave	129725	113138		43.6	50.0	-12.8	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37599.b\07DEC2016A_003.d
 Lims ID: CCV L2
 Client ID:
 Sample Type: CCVL
 Inject. Date: 07-Dec-2016 12:33:22 ALS Bottle#: 38 Worklist Smp#: 5
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: RB
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37599.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 07-Dec-2016 16:54:02 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK029

First Level Reviewer: westendorfc Date: 07-Dec-2016 13:15:35

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.549	1.545	0.004	17391644	51.8		104	1100126	
1 Perfluorobutyric acid	212.90 > 169.00	1.549	1.553	-0.004	306359	1.01		101	2645	
3 Perfluoropentanoic acid	262.90 > 219.00	1.819	1.823	-0.004	277648	1.02		102	2023	
D 4 13C5-PFPeA	267.90 > 223.00	1.829	1.823	0.006	13421195	50.7		101	972065	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.858	1.861	-0.003	448461	0.8536		96.6		
	298.90 > 99.00	1.858	1.861	-0.003	193951		2.31(0.00-0.00)			
D 6 13C2 PFHxA	315.00 > 270.00	2.114	2.114	0.0	11913489	50.2		100	718986	
7 Perfluorohexanoic acid	313.00 > 269.00	2.122	2.114	0.008	229334	1.01		101	8638	
D 11 13C4-PFHpA	367.00 > 322.00	2.460	2.458	0.002	11103394	53.5		107	1285501	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.460	2.458	0.002	226253	0.99		99.2	2493	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.400	2.476	-0.076	360442	1.00		110		
D 10 18O2 PFHxS	403.00 > 84.00	2.469	2.476	-0.007	15569961	49.8		105	806616	
15 Perfluorooctanoic acid	413.00 > 369.00	2.820	2.817	0.003	269704	1.05		105	3455	
	413.00 > 169.00	2.820	2.817	0.003	155122		1.74(0.90-1.10)		11443	
D 14 13C4 PFOA	417.00 > 372.00	2.820	2.817	0.003	11944378	54.5		109	834384	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.828	2.825	0.003	1.000	283824	0.9266	97.3	
D 17 13C4 PFOS	503.00	> 80.00	3.196	3.195	0.001		12437802	50.6	106	530071
D 19 13C5 PFNA	468.00	> 423.00	3.196	3.195	0.001		8625976	51.8	104	470795
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.196	3.195	0.001	1.000	249466	0.8834	95.2	22889
	499.00	> 99.00	3.188	3.195	-0.007	0.997	55540	4.49(0.90-1.10)		4147
20 Perfluorononanoic acid	463.00	> 419.00	3.196	3.195	0.001	1.000	170815	0.99	99.4	3586
D 21 13C8 FOSA	506.00	> 78.00	3.528	3.528	0.0		20698866	51.5	103	424878
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.528	3.528	0.0	1.000	394931	1.02	102	35277
24 Perfluorodecanoic acid	513.00	> 469.00	3.554	3.553	0.001	1.000	137775	0.9510	95.1	4027
D 23 13C2 PFDA	515.00	> 470.00	3.554	3.562	-0.008		7541616	47.8	95.6	223929
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.866	3.867	-0.001	1.000	148657	0.8930	92.6	
D 27 13C2 PFUnA	565.00	> 520.00	3.884	3.884	0.0		5800144	48.8	97.7	519081
28 Perfluoroundecanoic acid	563.00	> 519.00	3.884	3.884	0.0	1.000	120975	0.9785	97.9	3375
29 Perfluorododecanoic acid	613.00	> 569.00	4.178	4.173	0.005	1.000	98011	0.9705	97.1	2182
D 30 13C2 PFDaA	615.00	> 570.00	4.178	4.173	0.005		5320731	47.5	94.9	273006
31 Perfluorotridecanoic acid	663.00	> 619.00	4.443	4.452	-0.009	1.000	98135	0.9709	97.1	1794
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.681	4.684	-0.003	1.000	198704	1.01	101	759
	713.00	> 169.00	4.681	4.684	-0.003	1.000	32416	6.13(0.00-0.00)		11795
D 32 13C2-PFTeDA	715.00	> 670.00	4.690	4.693	-0.003		10765061	46.6	93.1	966268
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.102	5.104	-0.002	1.000	164498	0.8024	80.2	178
D 34 13C2-PFHxDA	815.00	> 770.00	5.102	5.104	-0.002		5656875	43.6	87.2	101566
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.460	5.462	-0.002	1.000	101370	0.9594	95.9	121

Reagents:

LCPFC-L2_00022

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37599.b\07DEC2016A_003.d

Injection Date: 07-Dec-2016 12:33:22

Instrument ID: A8_N

Lims ID: CCV L2

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 38

Worklist Smp#: 5

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

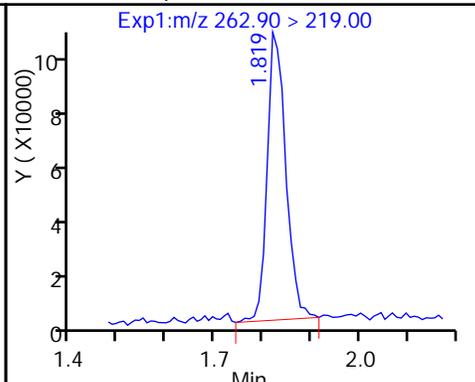
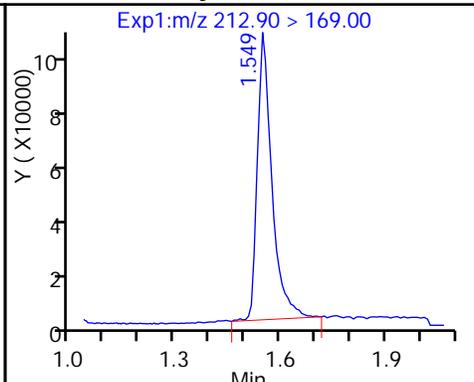
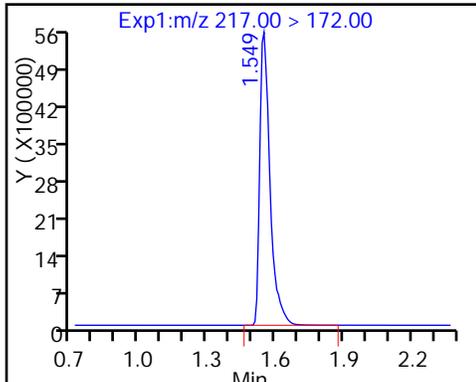
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

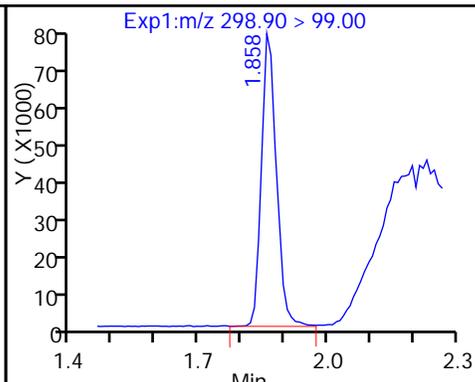
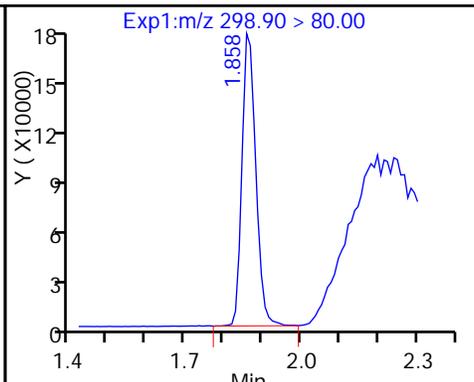
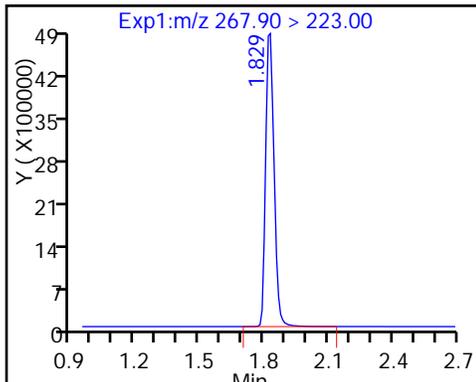
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

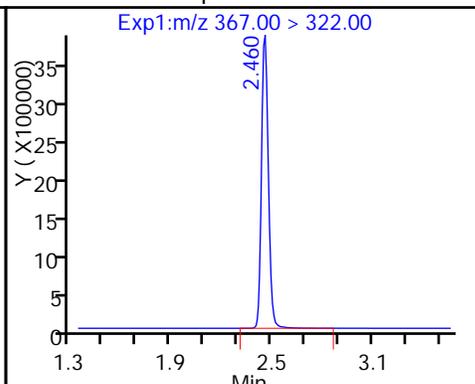
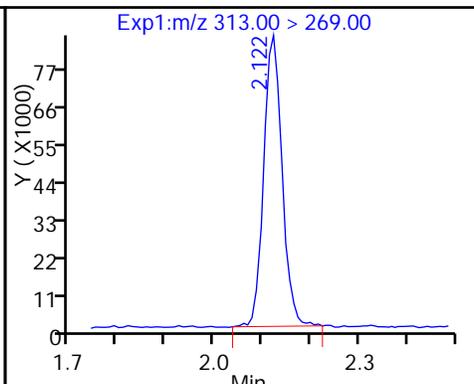
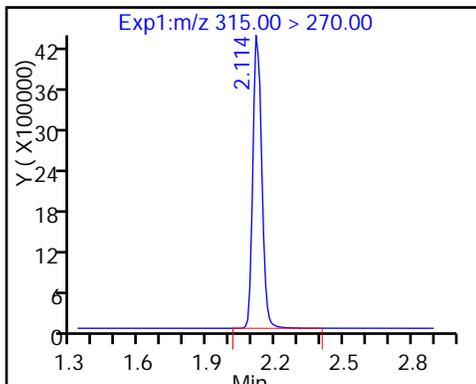
5 Perfluorobutanesulfonic acid



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

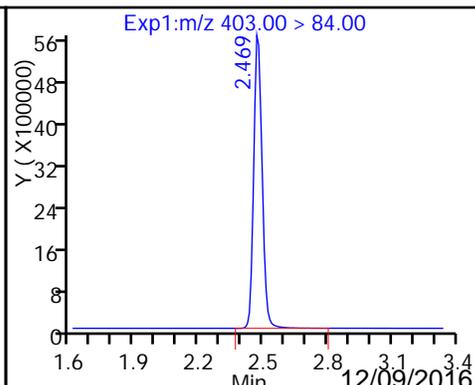
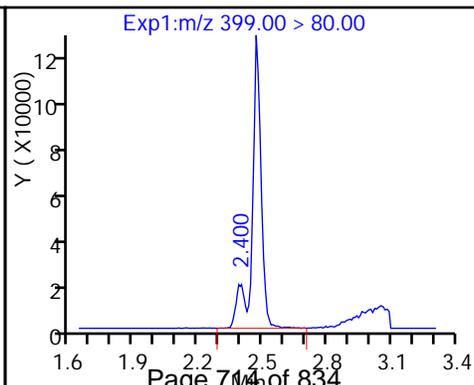
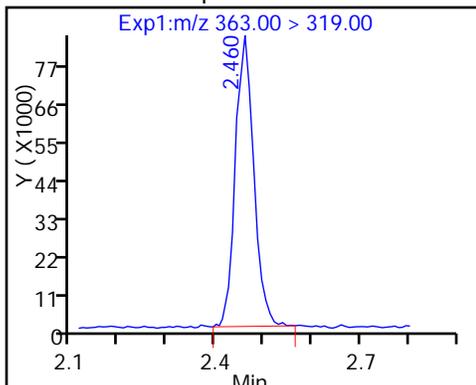
D 11 13C4-PFHpA

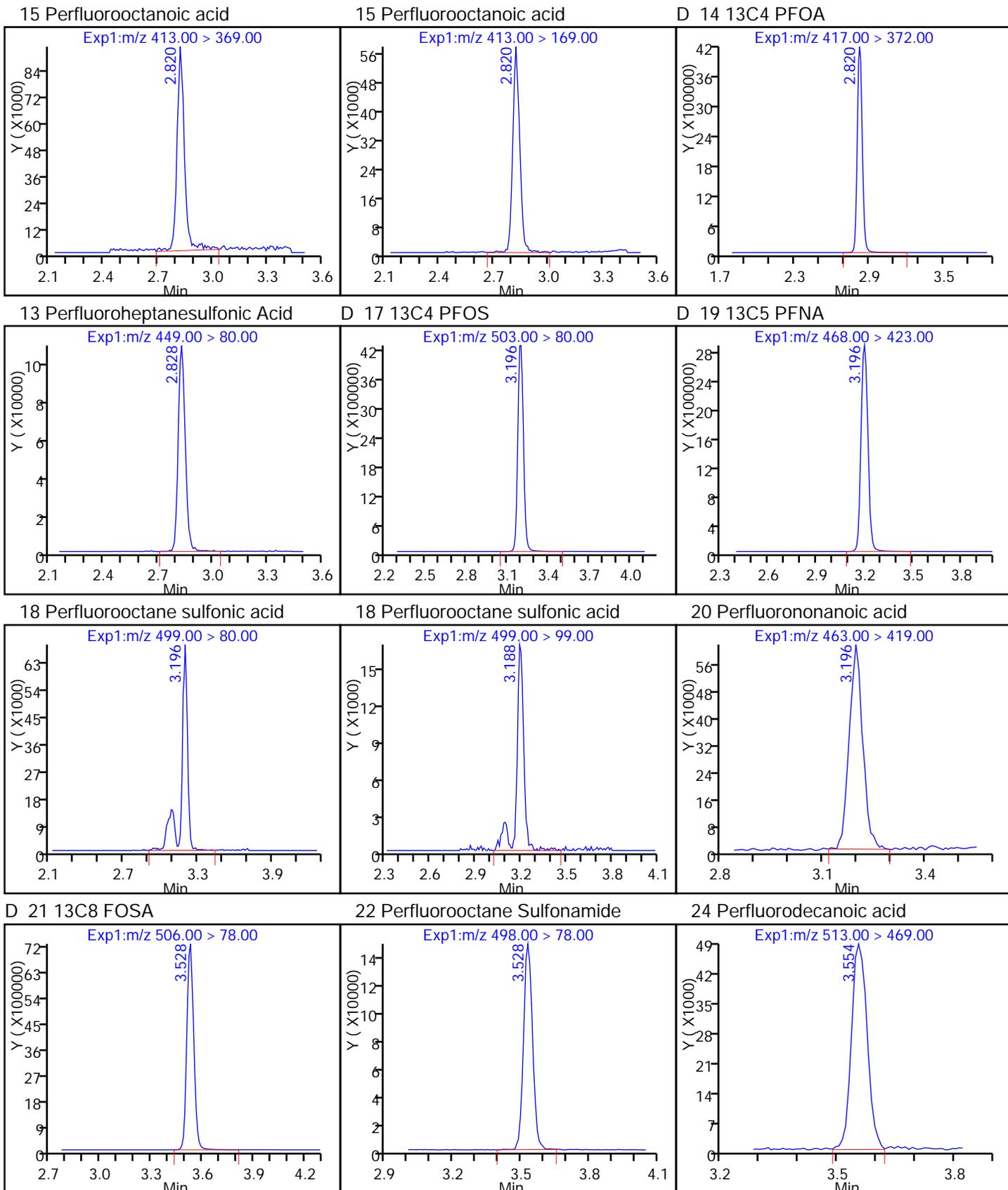


12 Perfluoroheptanoic acid

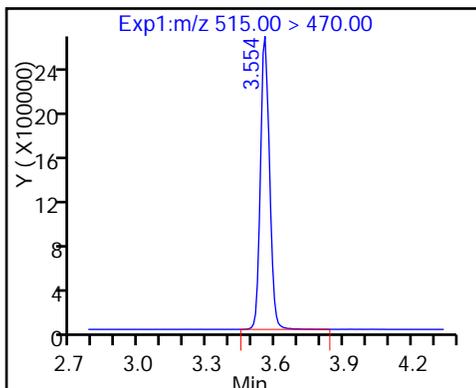
9 Perfluorohexanesulfonic acid

D 10 18O2 PFHxS

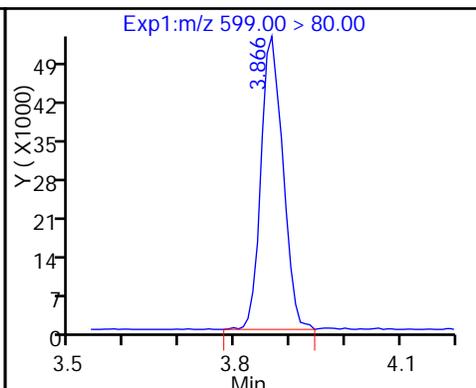




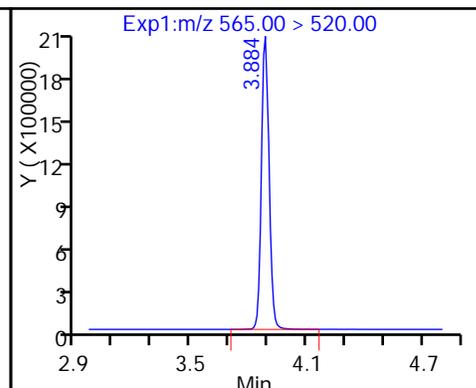
D 23 13C2 PFDA



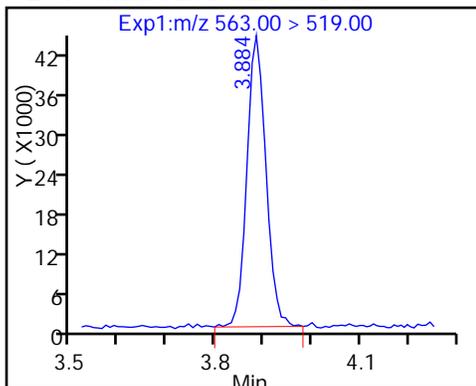
26 Perfluorodecane Sulfonic acid



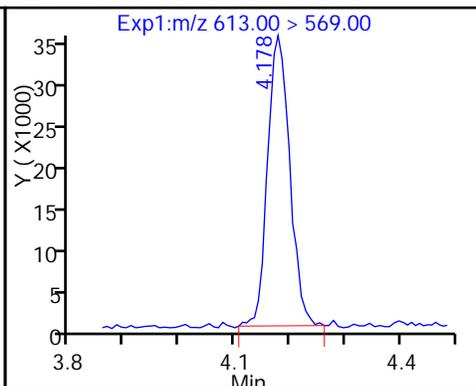
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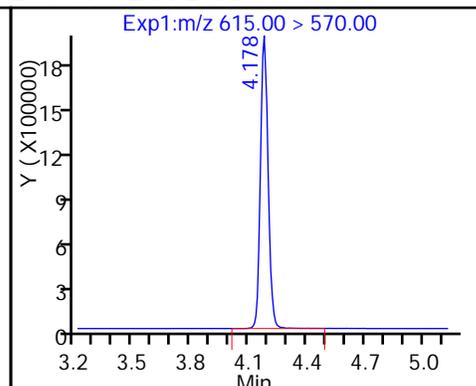
28 Perfluoroundecanoic acid



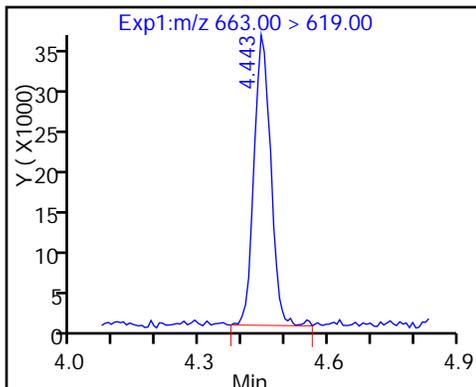
29 Perfluorododecanoic acid



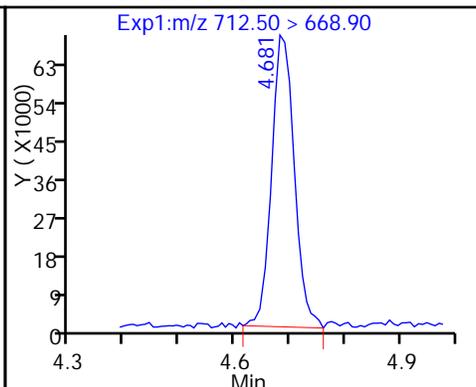
D 30 13C2 PFDaA



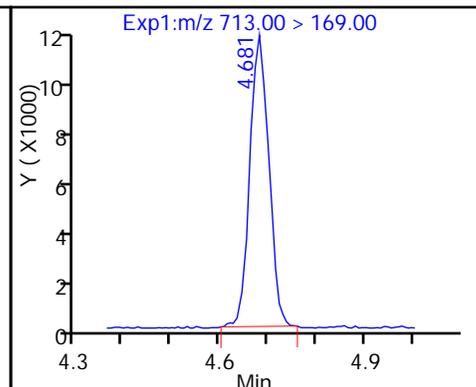
31 Perfluorotridecanoic acid



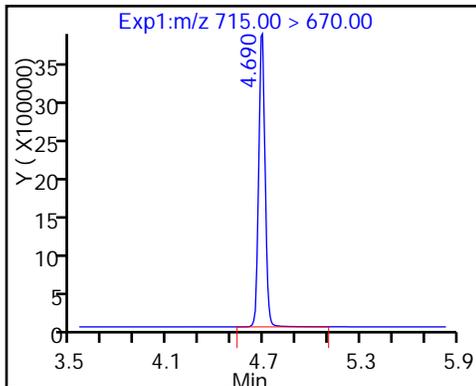
33 Perfluorotetradecanoic acid



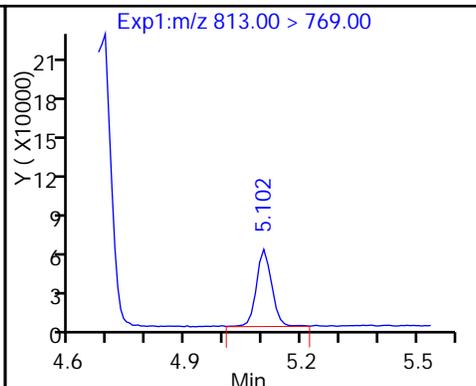
33 Perfluorotetradecanoic acid



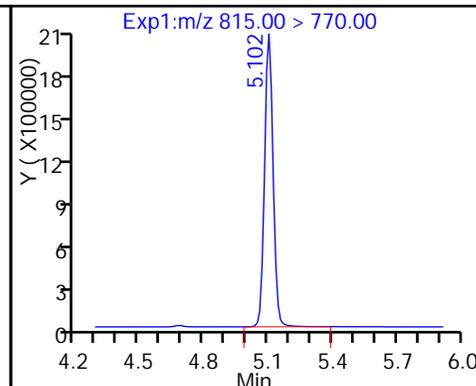
D 32 13C2-PFTeDA



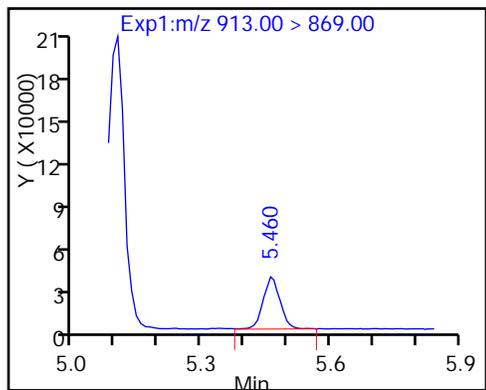
35 Perfluorohexadecanoic acid



D 34 13C2-PFHxDA



36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-141062/2 Calibration Date: 12/07/2016 16:40
 Instrument ID: A8_N Calib Start Date: 12/03/2016 13:48
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/03/2016 15:33
 Lab File ID: 07DEC2016B_002.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.596	1.604		44.4	44.2	0.5	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.072	1.030		48.1	50.0	-3.9	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.085	1.058		45.2	46.4	-2.5	25.0
13C4 PFBA	Ave	335829	312725		46.6	50.0	-6.9	50.0
13C5-PFPeA	Ave	264545	238010		45.0	50.0	-10.0	50.0
13C2 PFHxA	Ave	237486	220594		46.4	50.0	-7.1	50.0
13C4-PFHpA	Ave	207413	194634		46.9	50.0	-6.2	50.0
18O2 PFHxS	Ave	312342	293298		44.4	47.3	-6.1	50.0
13C4 PFOA	Ave	219258	195672		44.6	50.0	-10.8	50.0
13C4 PFOS	Ave	246009	231701		45.0	47.8	-5.8	50.0
13C5 PFNA	Ave	166415	147510		44.3	50.0	-11.4	50.0
13C8 FOSA	Ave	402279	362301		45.0	50.0	-9.9	50.0
13C2 PFDA	Ave	157817	134346		42.6	50.0	-14.9	50.0
13C2 PFUnA	Ave	118762	97918		41.2	50.0	-17.6	50.0
13C2 PFDoA	Ave	112084	94048		42.0	50.0	-16.1	50.0
13C2-PFTeDA	Ave	231173	194632		42.1	50.0	-15.8	50.0
13C2-PFHxDA	Ave	129725	101977		39.3	50.0	-21.4	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_002.d
 Lims ID: CCV L5
 Client ID:
 Sample Type: CCV
 Inject. Date: 07-Dec-2016 16:40:57 ALS Bottle#: 41 Worklist Smp#: 2
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L5
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:41:52 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d

Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:54:04

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.534	1.534	0.0	15636225	46.6		93.1	828932	
1 Perfluorobutyric acid	212.90 > 169.00	1.542	1.542	0.0	14059236	51.4		103	115968	
3 Perfluoropentanoic acid	262.90 > 219.00	1.820	1.820	0.0	11828155	49.0		97.9	115349	
D 4 13C5-PFPeA	267.90 > 223.00	1.820	1.820	0.0	11900494	45.0		90.0	856965	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.848	1.848	0.0	20799502	44.4		101		
	298.90 > 99.00	1.848	1.848	0.0	9409891		2.21(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.109	2.109	0.0	10356594	49.3		98.5	312643	
D 6 13C2 PFHxA	315.00 > 270.00	2.109	2.109	0.0	11029718	46.4		92.9	661333	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.460	2.460	0.0	13971121	43.4		95.4		
12 Perfluoroheptanoic acid	363.00 > 319.00	2.442	2.442	0.0	9773894	48.9		97.8	121575	
D 11 13C4-PFHpA	367.00 > 322.00	2.442	2.442	0.0	9731721	46.9		93.8	578610	
D 10 18O2 PFHxS	403.00 > 84.00	2.460	2.460	0.0	13872989	44.4		93.9	620057	
15 Perfluorooctanoic acid	413.00 > 369.00	2.802	2.802	0.0	10081606	48.1		96.1	359116	
	413.00 > 169.00	2.802	2.802	0.0	6185316		1.63(0.90-1.10)		319378	
D 14 13C4 PFOA	417.00 > 372.00	2.802	2.802	0.0	9783577	44.6		89.2	488275	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.810	2.810	0.0	1.000	12791414	46.9	98.5	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.061	3.061	0.0	1.000	11372952	45.2	97.5	335166
	499.00	> 99.00	3.169	3.061	0.108	1.035	2453380		4.64(0.90-1.10)	368490
D 17 13C4 PFOS	503.00	> 80.00	3.169	3.169	0.0		11075310	45.0	94.2	277945
D 19 13C5 PFNA	468.00	> 423.00	3.169	3.169	0.0		7375490	44.3	88.6	411723
20 Perfluorononanoic acid	463.00	> 419.00	3.177	3.177	0.0	1.000	7589604	51.6	103	141598
D 21 13C8 FOSA	506.00	> 78.00	3.510	3.510	0.0		18115044	45.0	90.1	375780
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.510	3.510	0.0	1.000	16706372	49.4	98.7	452673
D 23 13C2 PFDA	515.00	> 470.00	3.535	3.535	0.0		6717315	42.6	85.1	273157
24 Perfluorodecanoic acid	513.00	> 469.00	3.535	3.535	0.0	1.000	6363067	49.3	98.6	193170
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.846	3.846	0.0	1.000	6864650	46.3	96.1	
D 27 13C2 PFUnA	565.00	> 520.00	3.863	3.863	0.0		4895890	41.2	82.4	448575
28 Perfluoroundecanoic acid	563.00	> 519.00	3.863	3.863	0.0	1.000	4698388	45.0	90.0	113052
29 Perfluorododecanoic acid	613.00	> 569.00	4.153	4.153	0.0	1.000	4298857	48.2	96.3	84147
D 30 13C2 PFDaA	615.00	> 570.00	4.153	4.153	0.0		4702404	42.0	83.9	230400
31 Perfluorotridecanoic acid	663.00	> 619.00	4.418	4.418	0.0	1.000	4305616	48.2	96.4	70801
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.661	4.661	0.0	1.000	8283755	47.5	95.0	100455
	713.00	> 169.00	4.661	4.661	0.0	1.000	1285114		6.45(0.00-0.00)	117470
D 32 13C2-PFTeDA	715.00	> 670.00	4.661	4.661	0.0		9731605	42.1	84.2	838305
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.079	5.079	0.0	1.000	4650460	49.0	98.1	5313
D 34 13C2-PFHxDA	815.00	> 770.00	5.079	5.079	0.0		5098842	39.3	78.6	99009
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.436	5.436	0.0	1.000	4554958	48.8	97.6	6248

Reagents:

LCPFC-L5_00020

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_002.d

Injection Date: 07-Dec-2016 16:40:57

Instrument ID: A8_N

Lims ID: CCV L5

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 41

Worklist Smp#: 2

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

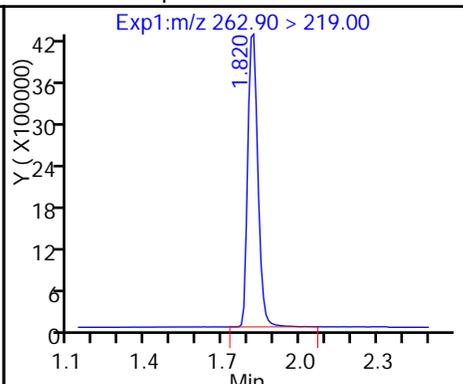
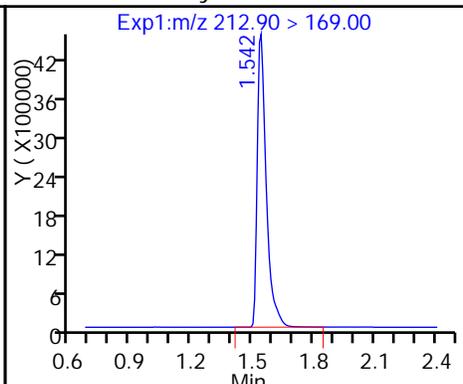
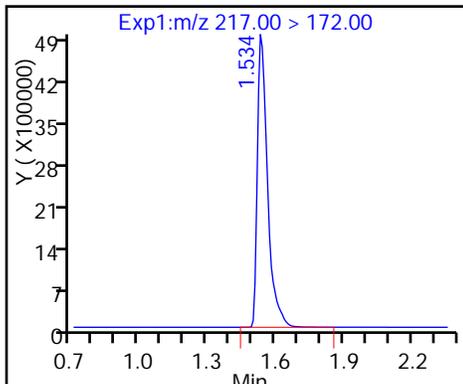
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

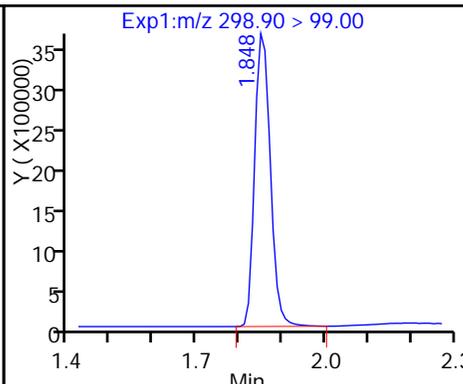
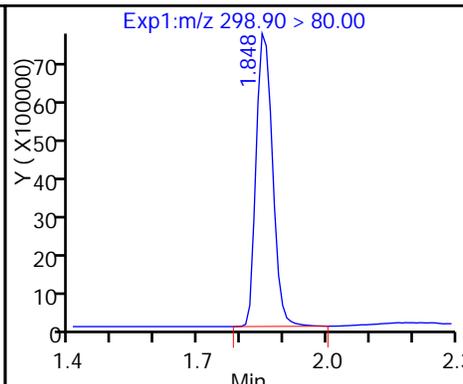
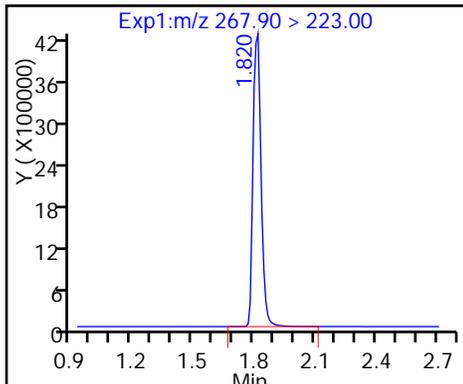
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

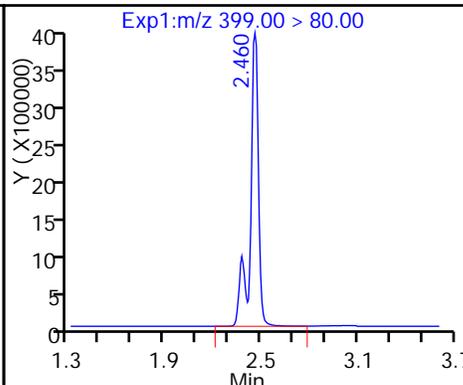
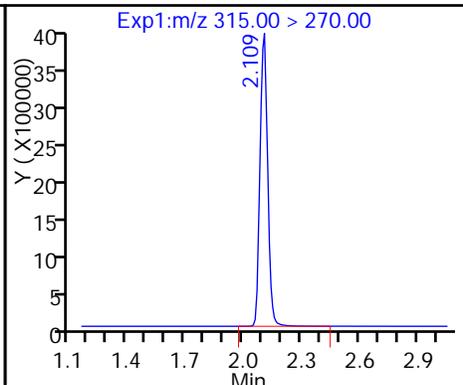
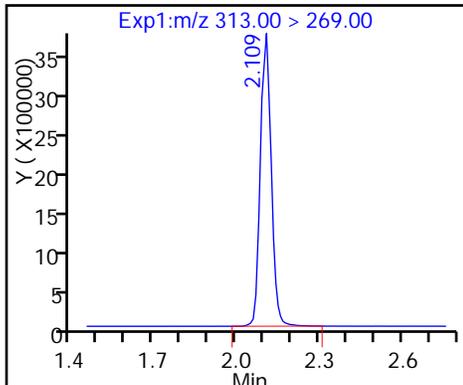
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

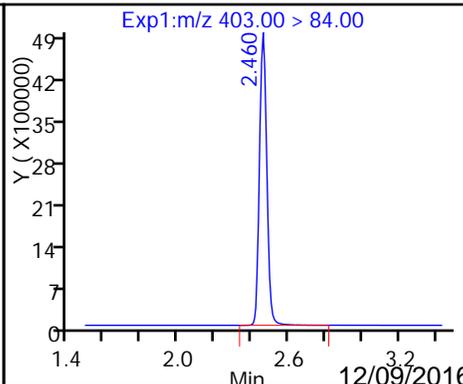
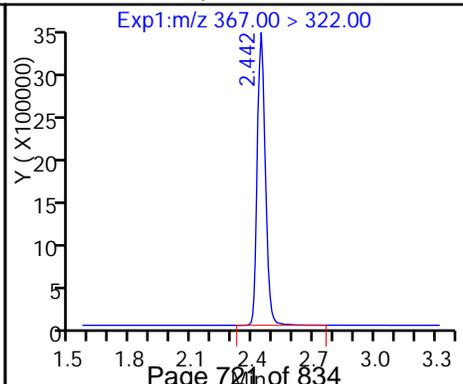
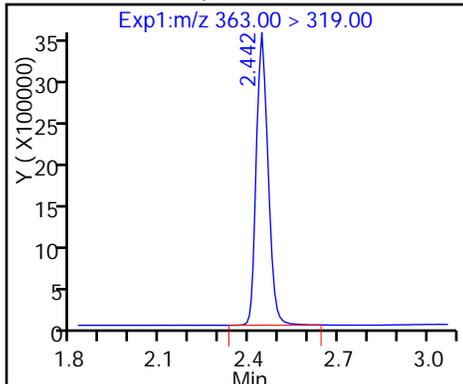
9 Perfluorohexanesulfonic acid

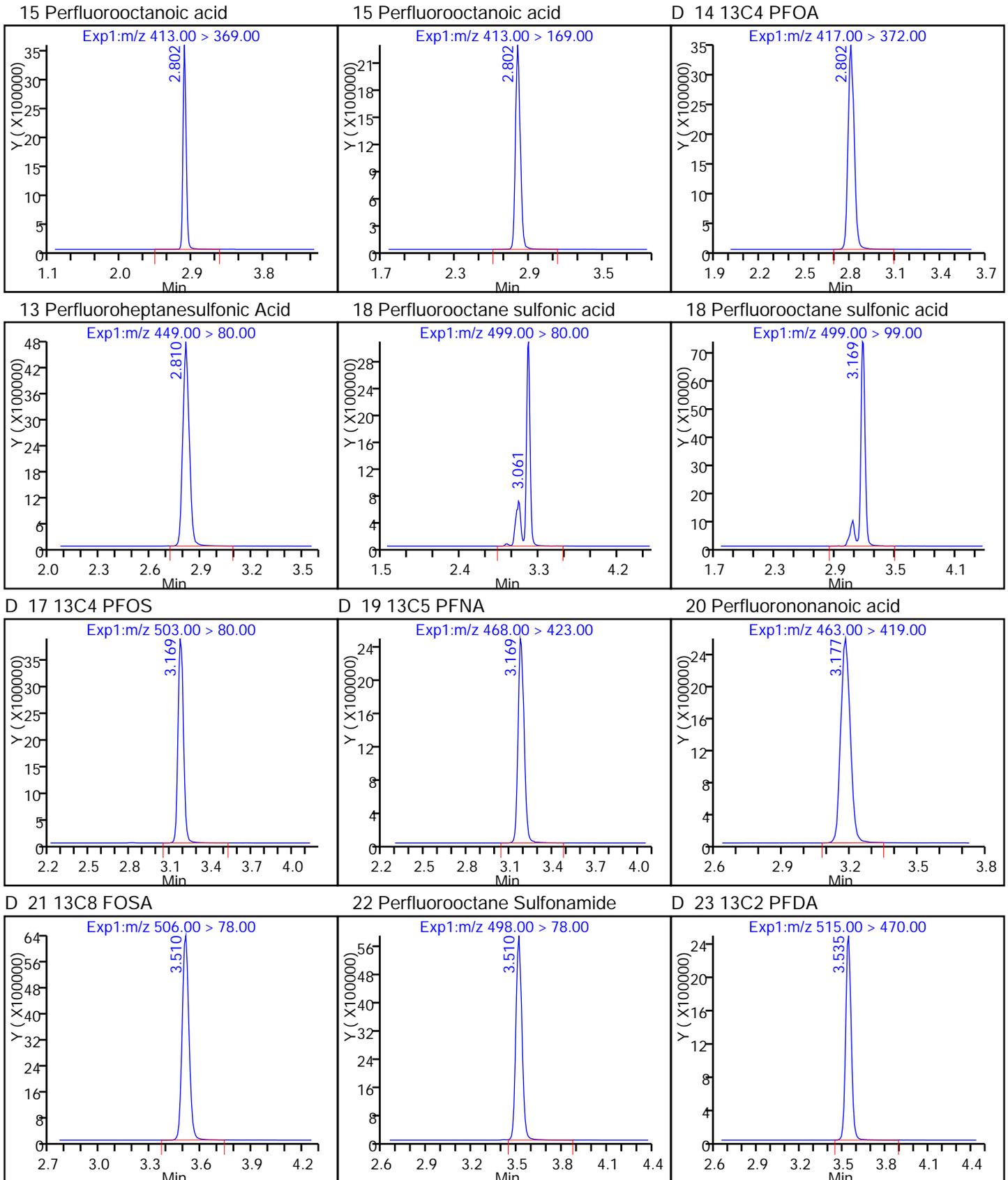


12 Perfluoroheptanoic acid

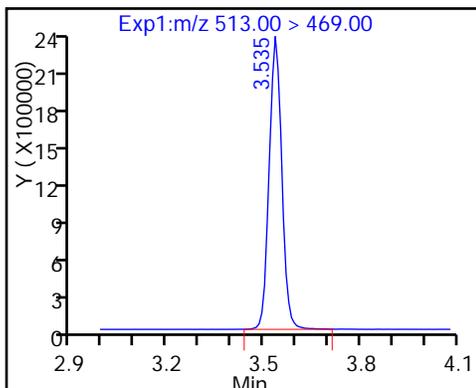
D 11 13C4-PFHpA

D 10 18O2 PFHxS

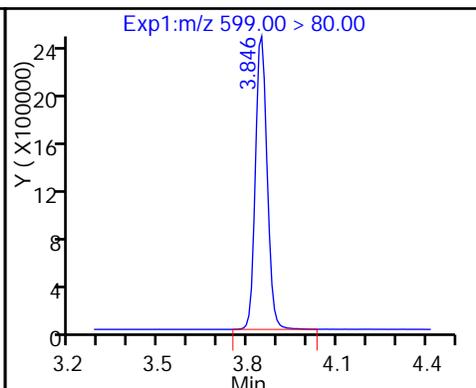




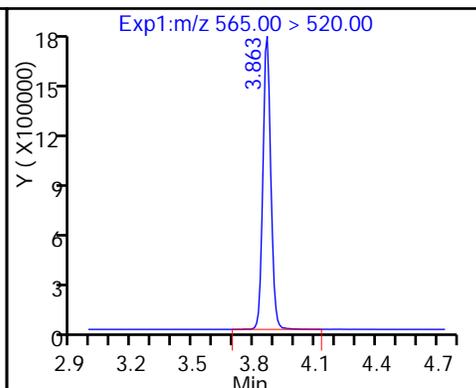
24 Perfluorodecanoic acid



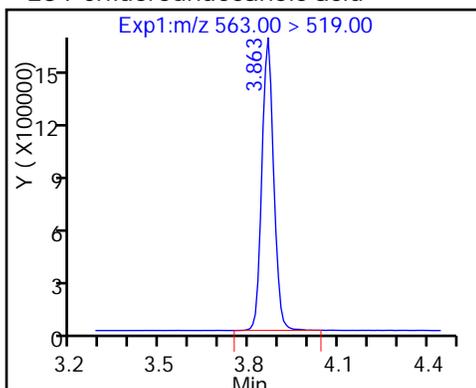
26 Perfluorodecane Sulfonic acid



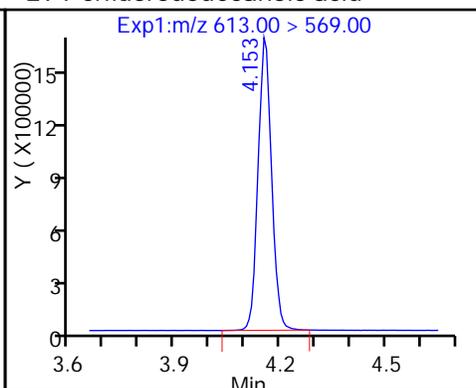
D 27 13C2 PFUnA



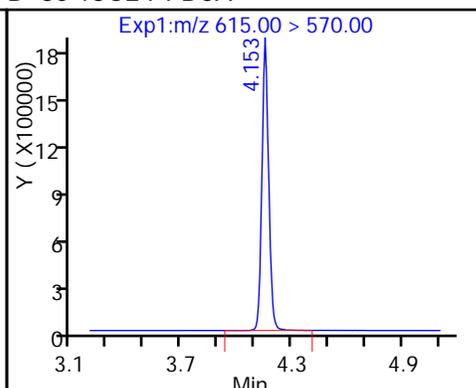
28 Perfluoroundecanoic acid



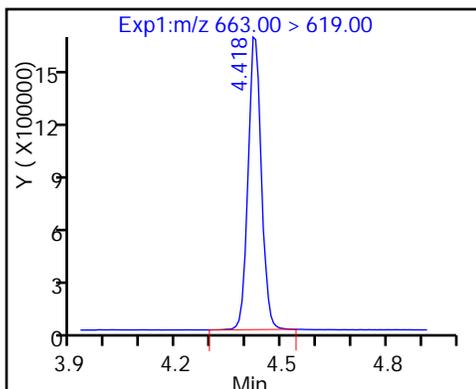
29 Perfluorododecanoic acid



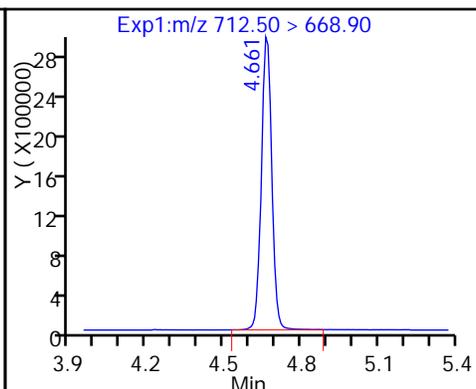
D 30 13C2 PFDa



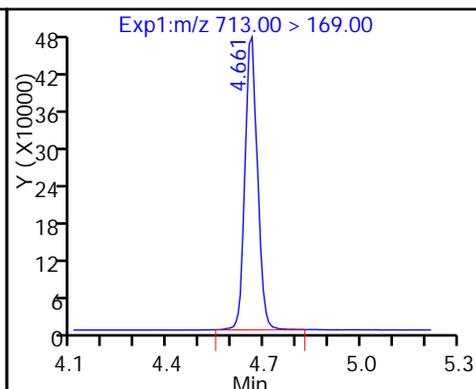
31 Perfluorotridecanoic acid



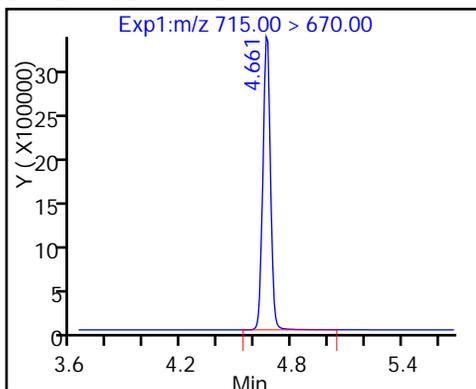
33 Perfluorotetradecanoic acid



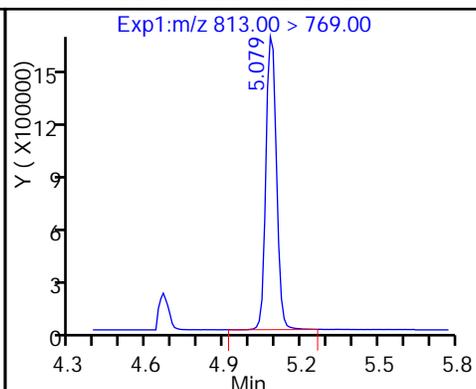
33 Perfluorotetradecanoic acid



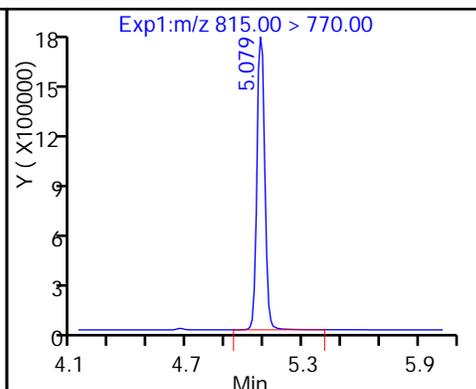
D 32 13C2-PFTeDA



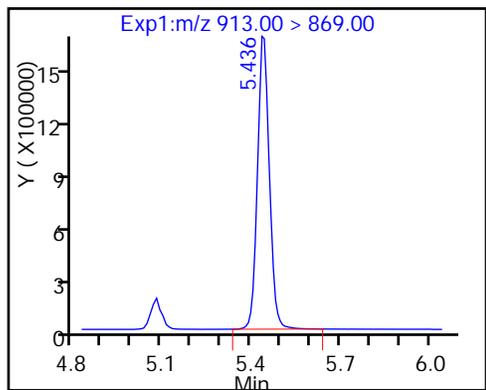
35 Perfluorohexadecanoic acid



D 34 13C2-PFHxDA



36 Perfluorooctadecanoic acid



FORM VII
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Lab Sample ID: CCV 320-141062/12 Calibration Date: 12/07/2016 17:55
 Instrument ID: A8_N Calib Start Date: 12/03/2016 13:48
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/03/2016 15:33
 Lab File ID: 07DEC2016B_012.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanesulfonic acid (PFBS)	AveID	1.596	1.698		18.8	17.7	6.4	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.072	1.112		20.7	20.0	3.7	25.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.085	1.115		19.1	18.6	2.7	25.0
13C4 PFBA	Ave	335829	345256		51.4	50.0	2.8	50.0
13C5-PFPeA	Ave	264545	263301		49.8	50.0	-0.5	50.0
13C2 PFHxA	Ave	237486	242078		51.0	50.0	1.9	50.0
13C4-PFHpA	Ave	207413	217275		52.4	50.0	4.8	50.0
18O2 PFHxS	Ave	312342	327905		49.7	47.3	5.0	50.0
13C4 PFOA	Ave	219258	222317		50.7	50.0	1.4	50.0
13C4 PFOS	Ave	246009	251697		48.9	47.8	2.3	50.0
13C5 PFNA	Ave	166415	166606		50.1	50.0	0.1	50.0
13C8 FOSA	Ave	402279	394496		49.0	50.0	-1.9	50.0
13C2 PFDA	Ave	157817	148417		47.0	50.0	-6.0	50.0
13C2 PFUnA	Ave	118762	108618		45.7	50.0	-8.5	50.0
13C2 PFDoA	Ave	112084	101693		45.4	50.0	-9.3	50.0
13C2-PFTeDA	Ave	231173	205294		44.4	50.0	-11.2	50.0
13C2-PFHxDA	Ave	129725	108113		41.7	50.0	-16.7	50.0

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_012.d
 Lims ID: CCV L4
 Client ID:
 Sample Type: CCV
 Inject. Date: 07-Dec-2016 17:55:57 ALS Bottle#: 40 Worklist Smp#: 12
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: CCV L4
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Sublist: chrom-A8_N*sub1
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 08-Dec-2016 09:42:07 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK033

First Level Reviewer: westendorfc Date: 08-Dec-2016 08:56:09

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.541	1.534	0.007	17262793	51.4		103	1403249	
1 Perfluorobutyric acid	212.90 > 169.00	1.541	1.542	-0.001	1.000	6671816	22.1	111	55460	
3 Perfluoropentanoic acid	262.90 > 219.00	1.819	1.820	-0.001	1.000	5599281	21.0	105	50089	
D 4 13C5-PFPeA	267.90 > 223.00	1.819	1.820	-0.001		13165063	49.8	99.5	1226213	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.848	1.848	0.0	1.000	9846174	18.8	106		
	298.90 > 99.00	1.848	1.848	0.0	1.000	4245854	2.32(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.103	2.109	-0.006	1.000	4846653	21.0	105	140191	
D 6 13C2 PFHxA	315.00 > 270.00	2.103	2.109	-0.006		12103899	51.0	102	880831	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.443	2.442	0.001	1.000	4714738	21.1	106	56032	
D 11 13C4-PFHpA	367.00 > 322.00	2.443	2.442	0.001		10863748	52.4	105	1271131	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.460	2.460	0.0	1.000	6438052	17.9	98.3		
D 10 18O2 PFHxS	403.00 > 84.00	2.451	2.460	-0.009		15509904	49.7	105	892994	
15 Perfluorooctanoic acid	413.00 > 369.00	2.800	2.802	-0.002	1.000	4943937	20.7	104	83959	
	413.00 > 169.00	2.800	2.802	-0.002	1.000	2964769	1.67(0.90-1.10)		130631	
D 14 13C4 PFOA	417.00 > 372.00	2.800	2.802	-0.002		11115843	50.7	101	984380	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags	
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.808	2.810	-0.002	1.000	6068100	20.5		108	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.169	3.061	0.108	1.000	5207884	19.1		103	3820
	499.00	> 99.00	3.177	3.061	0.116	1.003	1137641		4.58(0.90-1.10)		89754
D 17 13C4 PFOS	503.00	> 80.00	3.177	3.169	0.008		12031128	48.9		102	323326
D 19 13C5 PFNA	468.00	> 423.00	3.177	3.169	0.008		8330295	50.1		100	572933
20 Perfluorononanoic acid	463.00	> 419.00	3.169	3.177	-0.008	1.000	3514127	21.2		106	56920
D 21 13C8 FOSA	506.00	> 78.00	3.501	3.510	-0.009		19724818	49.0		98.1	836754
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.510	3.510	0.0	1.000	8004148	21.7		109	372606
D 23 13C2 PFDA	515.00	> 470.00	3.535	3.535	0.0		7420835	47.0		94.0	239982
24 Perfluorodecanoic acid	513.00	> 469.00	3.535	3.535	0.0	1.000	2986371	20.9		105	106578
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.846	3.846	0.0	1.000	3209681	19.9		103	
D 27 13C2 PFUnA	565.00	> 520.00	3.863	3.863	0.0		5430907	45.7		91.5	494854
28 Perfluoroundecanoic acid	563.00	> 519.00	3.863	3.863	0.0	1.000	2283019	19.7		98.6	64746
29 Perfluorododecanoic acid	613.00	> 569.00	4.153	4.153	0.0	1.000	1982079	20.5		103	54971
D 30 13C2 PFDaA	615.00	> 570.00	4.153	4.153	0.0		5084646	45.4		90.7	159279
31 Perfluorotridecanoic acid	663.00	> 619.00	4.418	4.418	0.0	1.000	1943594	20.1		101	37818
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.661	4.661	0.0	1.000	3732526	19.8		99.0	47957
	713.00	> 169.00	4.652	4.661	-0.009	0.998	582313		6.41(0.00-0.00)		72018
D 32 13C2-PFTeDA	715.00	> 670.00	4.661	4.661	0.0		10264724	44.4		88.8	553416
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.079	5.079	0.0	1.000	2098560	20.0		100	2398
D 34 13C2-PFHxDA	815.00	> 770.00	5.079	5.079	0.0		5405634	41.7		83.3	103268
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.437	5.436	0.001	1.000	1660459	16.4		82.2	2247

Reagents:

LCPFC-L4_00023

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161207-37602.b\07DEC2016B_012.d

Injection Date: 07-Dec-2016 17:55:57

Instrument ID: A8_N

Lims ID: CCV L4

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 40

Worklist Smp#: 12

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

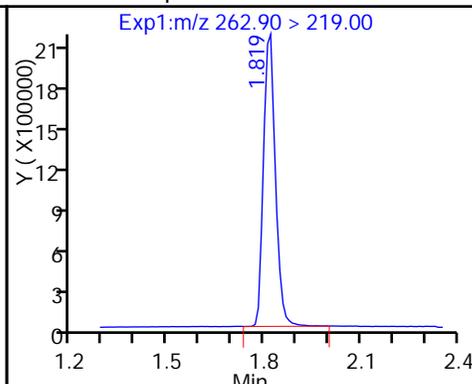
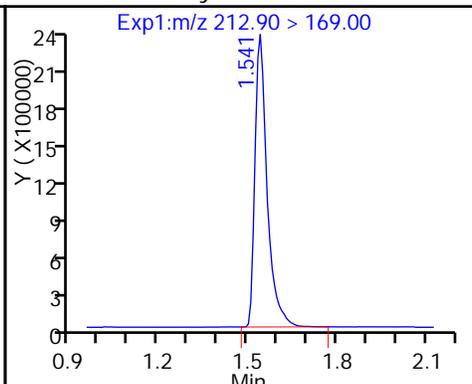
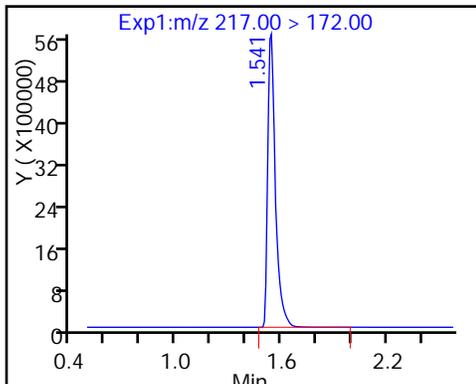
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

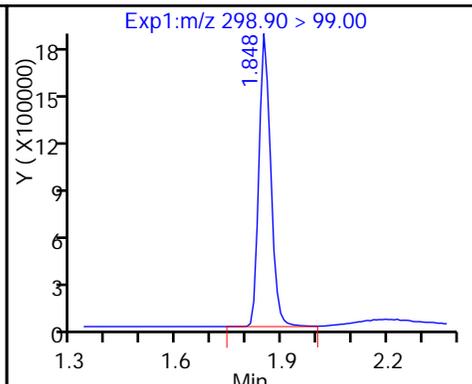
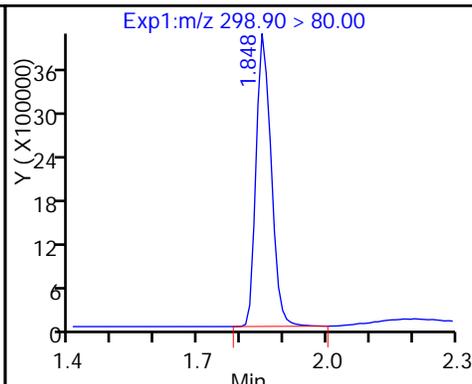
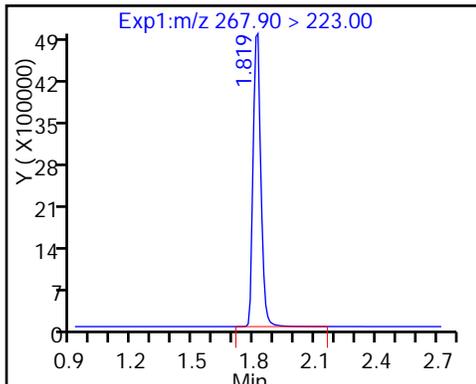
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

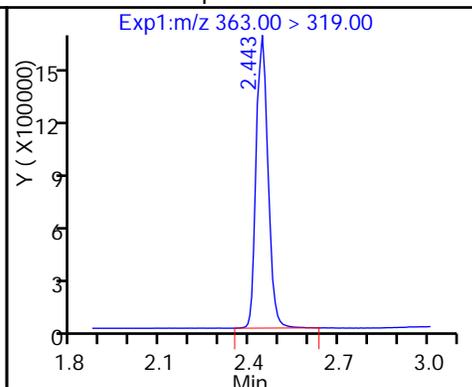
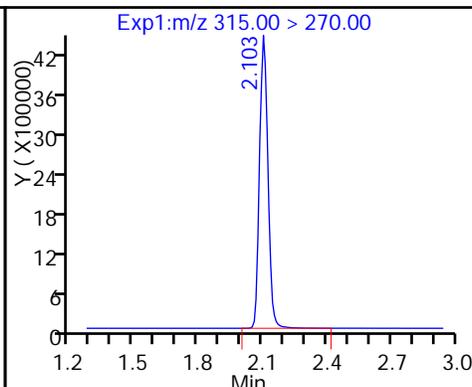
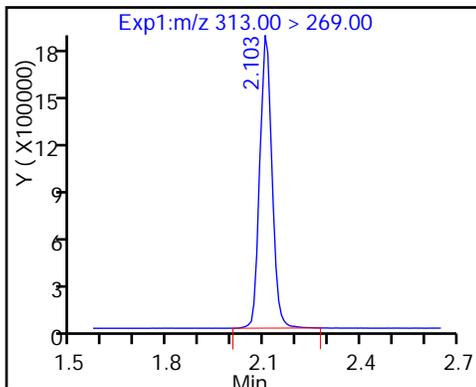
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

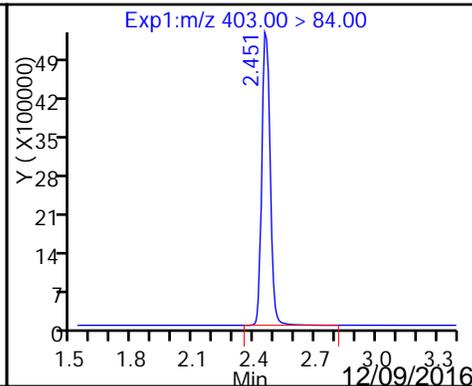
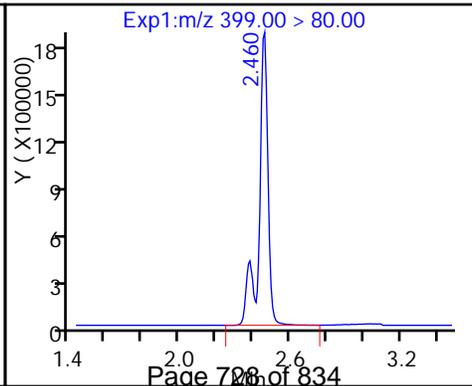
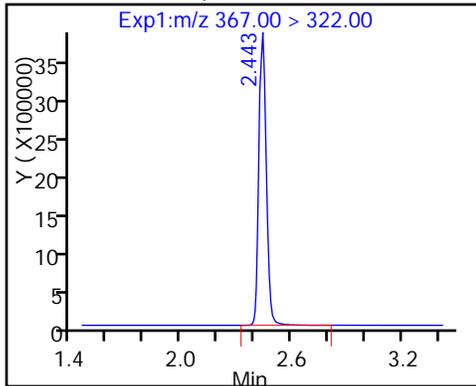
12 Perfluoroheptanoic acid

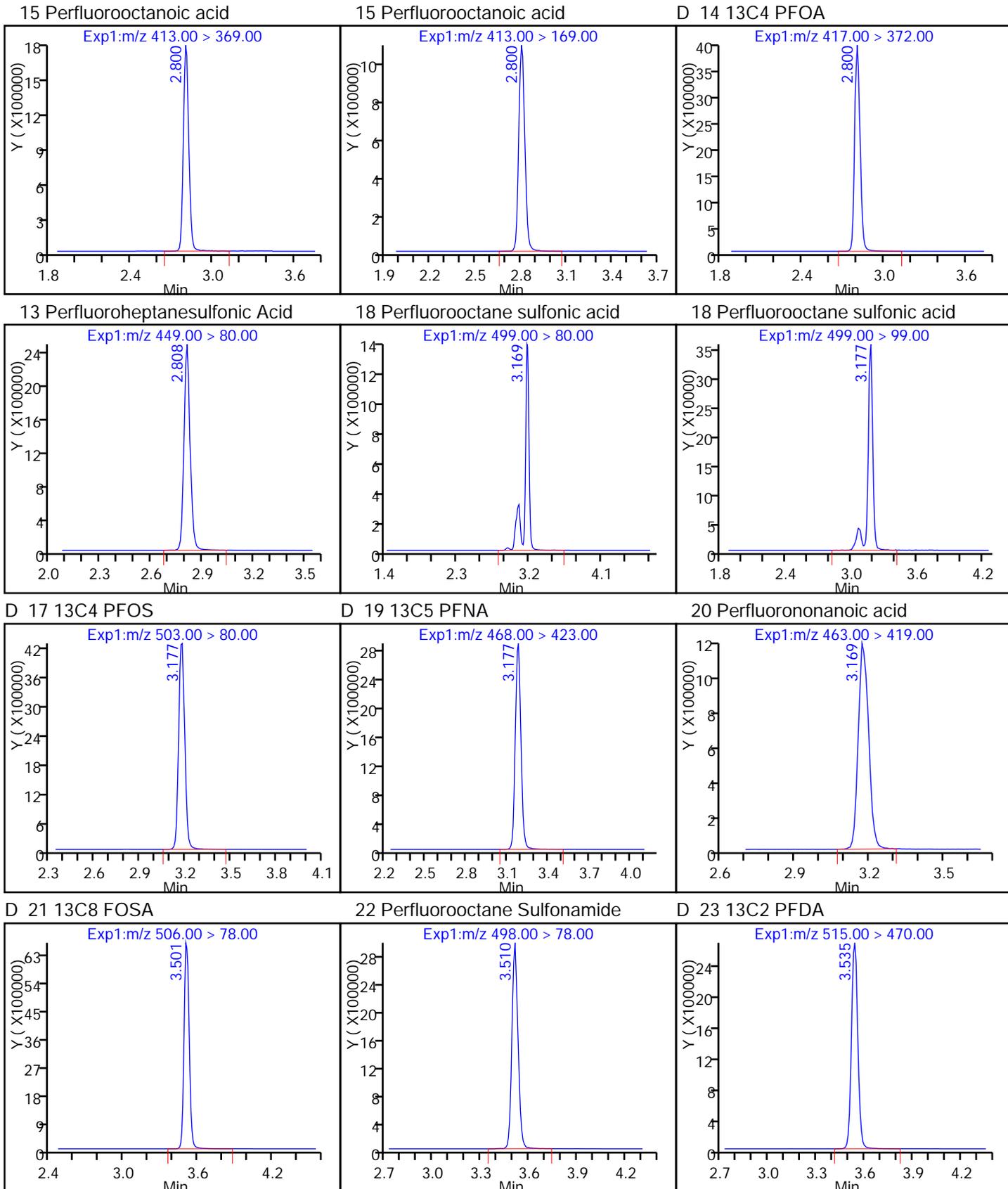


D 11 13C4-PFHpA

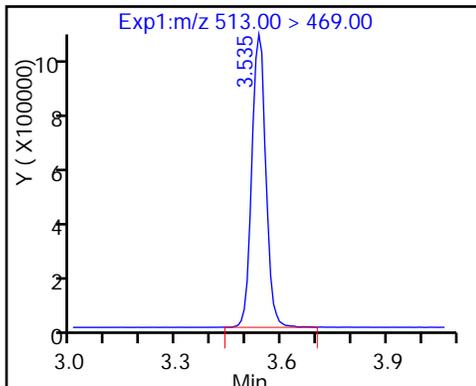
9 Perfluorohexanesulfonic acid

D 10 18O2 PFHxS

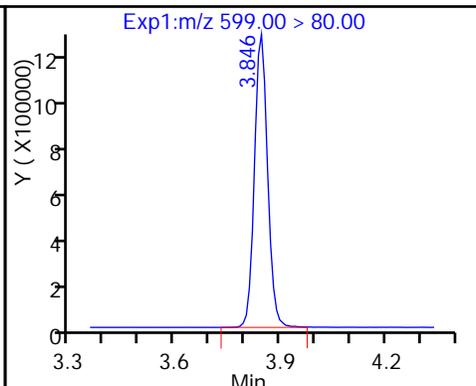




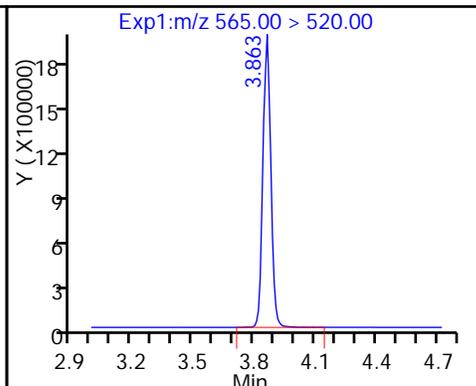
24 Perfluorodecanoic acid



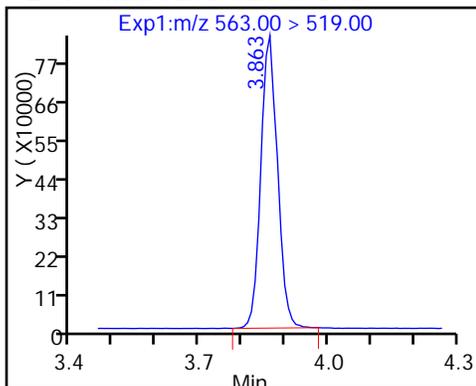
26 Perfluorodecane Sulfonic acid



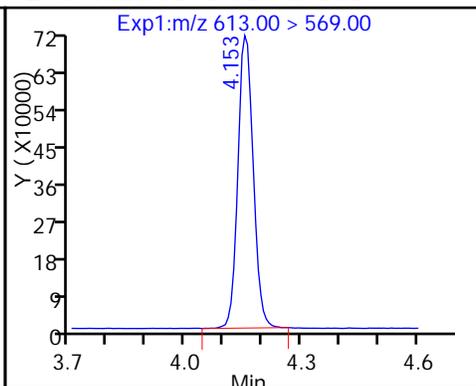
D 27 13C2 PFUnA



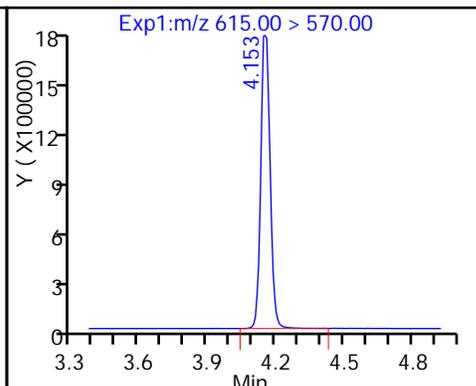
28 Perfluoroundecanoic acid



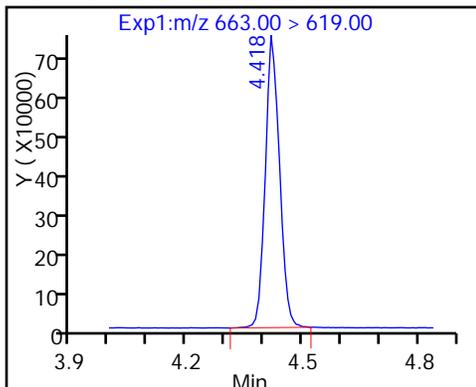
29 Perfluorododecanoic acid



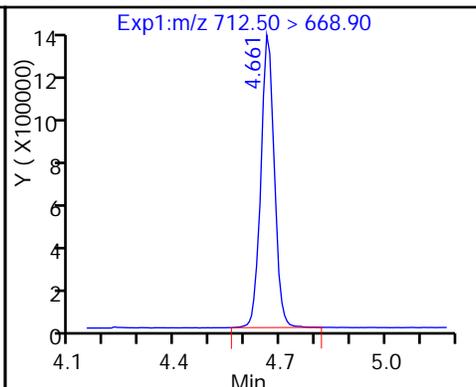
D 30 13C2 PFDaA



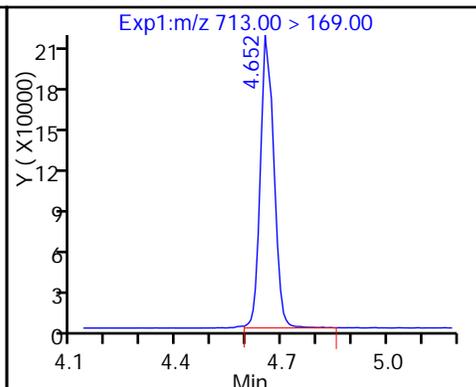
31 Perfluorotridecanoic acid



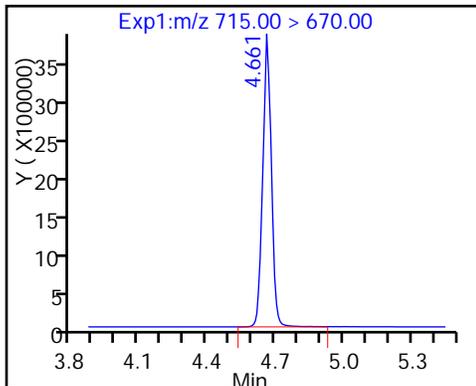
33 Perfluorotetradecanoic acid



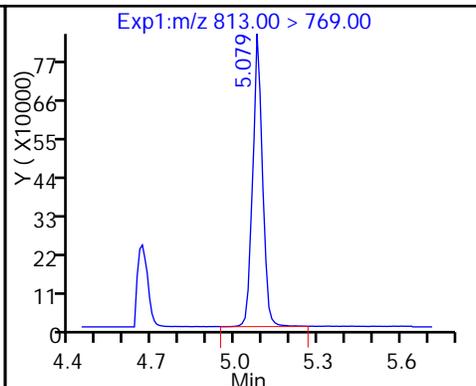
33 Perfluorotetradecanoic acid



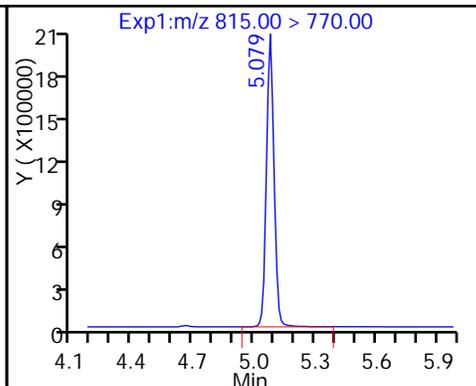
D 32 13C2-PFTeDA



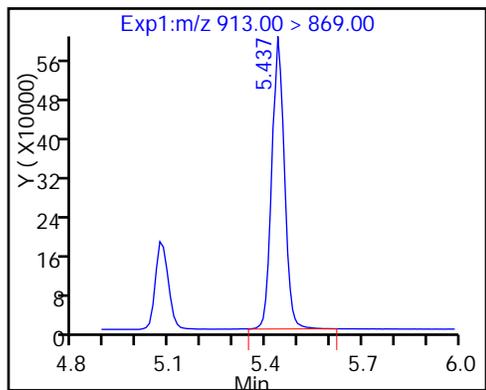
35 Perfluorohexadecanoic acid



D 34 13C2-PFHxDA



36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 320-139627/1-A
 Matrix: Solid Lab File ID: 02DEC2016C_001.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.00(g) Date Analyzed: 12/02/2016 14:59
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.30	U	0.50	0.10
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.30	U	0.50	0.13
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.30	U	0.40	0.10

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	108		25-150
STL00991	13C4 PFOS	94		25-150
STL00994	18O2 PFHxS	104		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_001.d
 Lims ID: MB 320-139627/1-A
 Client ID:
 Sample Type: MB
 Inject. Date: 02-Dec-2016 14:59:32 ALS Bottle#: 1 Worklist Smp#: 5
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: mb 320-139627/1-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:16:35 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:16:34

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.574	0.0	1.000	21795	0.0719		154	
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0		17111077	50.3	101	1388607	
3 Perfluoropentanoic acid	262.90 > 219.00	1.868	1.868	0.0	1.000	23535	0.0847		149	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.868	-0.010		13340497	49.4	98.7	1043506	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.897	1.907	-0.010	1.000	5255	0.0103			
	298.90 > 99.00	1.887	1.907	-0.020	0.995	2254	2.33(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.179	2.173	0.006	1.000	13669	0.0634		241	
D 6 13C2 PFHxA	315.00 > 270.00	2.162	2.173	-0.011		11285724	45.8	91.6	435771	
D 11 13C4-PFHpA	367.00 > 322.00	2.520	2.519	0.001		11132042	52.9	106	1111017	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.527	2.534	-0.007	1.000	54849	0.1642			
D 10 18O2 PFHxS	403.00 > 84.00	2.535	2.534	0.001		14913821	49.3	104	737756	
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.862	2.870	-0.008	1.000	3757	NR			
D 47 M2-6:2FTS	429.00 > 409.00	2.870	2.870	0.0		2973	0.0213	0.0		
D 14 13C4 PFOA	417.00 > 372.00	2.886	2.886	0.0		11840544	54.0	108	891757	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.902	2.894	0.008	1.000	58370	0.2416			382	
413.00 > 169.00	2.902	2.894	0.008	1.000	22471		2.60(0.90-1.10)		970	
D 17 13C4 PFOS										
503.00 > 80.00	3.268	3.267	0.001		11022180	45.1		94.3	717052	
D 19 13C5 PFNA										
468.00 > 423.00	3.268	3.275	-0.007		8809234	49.5		98.9	641246	
D 21 13C8 FOSA										
506.00 > 78.00	3.582	3.589	-0.007		12412970	30.0		59.9	447762	
43 Sodium 1H,1H,2H,2H-perfluorooctane										
527.00 > 507.00	3.632	3.624	0.008	1.005	1156	NR				
D 42 M2-8:2FTS										
529.00 > 509.00	3.616	3.624	-0.008		3435	0.0257		0.0		
D 23 13C2 PFDA										
515.00 > 470.00	3.632	3.631	0.001		7837443	47.2		94.4	389458	
24 Perfluorodecanoic acid										
513.00 > 469.00	3.632	3.631	0.001	1.000	4782	0.0315			163	
D 45 d3-NMeFOSAA										
573.00 > 419.00	3.787	3.790	-0.003		11498	0.1437		0.0		
44 N-methyl perfluorooctane sulfonami										
570.00 > 419.00	3.777	3.799	-0.022	0.997	2321	NR				
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.936	3.941	-0.005	1.000	808	0.005587				
D 46 d5-NEtFOSAA										
589.00 > 419.00	3.953	3.955	-0.002		16818	0.1888		0.0		
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.971	3.959	0.012	1.000	14456	0.1155			354	
49 N-ethyl perfluorooctane sulfonamid										
584.00 > 419.00	3.953	3.963	-0.010	1.000	3184	NR				
D 27 13C2 PFUnA										
565.00 > 520.00	3.962	3.967	-0.005		5905793	47.0		93.9	315207	
D 52 d-N-MeFOSA-M										
515.00 > 169.00	4.052	4.079	-0.027		5561	0.0523		0.0		
54 MeFOSA										
512.00 > 169.00	4.085	4.087	-0.002	1.000	998	NR				
D 30 13C2 PFDaA										
615.00 > 570.00	4.251	4.259	-0.008		4628188	39.6		79.2	143761	
D 51 d-N-EtFOSA-M										
531.00 > 169.00	4.251	4.261	-0.010		6878	0.0675		0.0		
53 N-ethylperfluoro-1-octanesulfonami										
526.00 > 169.00	4.285	4.271	0.014	1.000	842	NR				
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.743	4.772	-0.030	1.000	30983	0.1732			15.8	
713.00 > 169.00	4.759	4.772	-0.013	1.003	2876		10.77(0.00-0.00)		1371	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.759	4.772	-0.013		7246692	30.0		59.9	383039	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.178	5.184	-0.006		2329899	17.9		35.7	106732	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
35 Perfluorohexadecanoic acid	813.00 > 769.00	5.178	5.194	-0.016	1.000	34261	-0.4475		71.7	
36 Perfluorooctadecanoic acid	913.00 > 869.00	5.566	5.562	0.004	1.000	2419	0.0332		11.2	

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_001.d

Injection Date: 02-Dec-2016 14:59:32

Instrument ID: A8_N

Lims ID: MB 320-139627/1-A

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 1

Worklist Smp#: 5

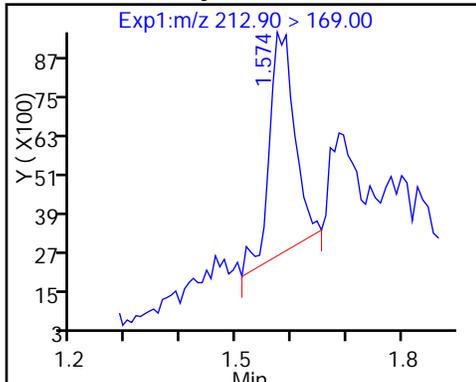
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

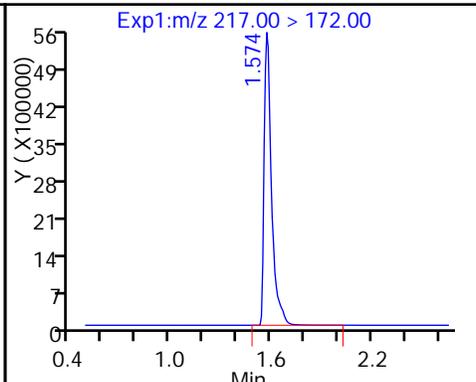
Method: A8_N

Limit Group: LC PFC_DOD ICAL

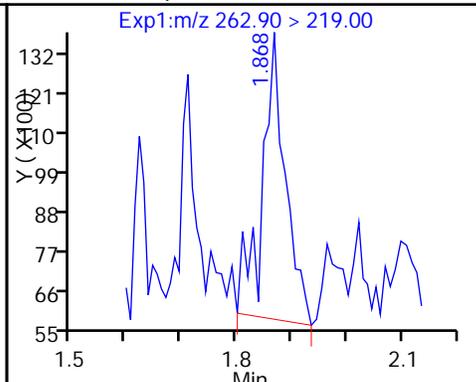
1 Perfluorobutyric acid



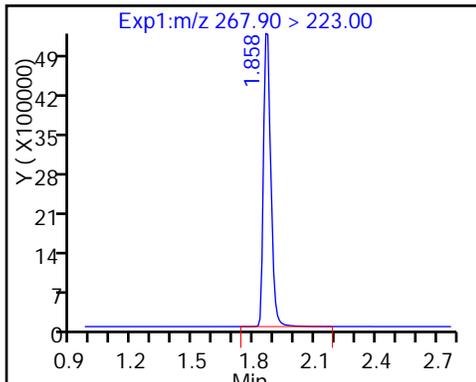
D 2 13C4 PFBA



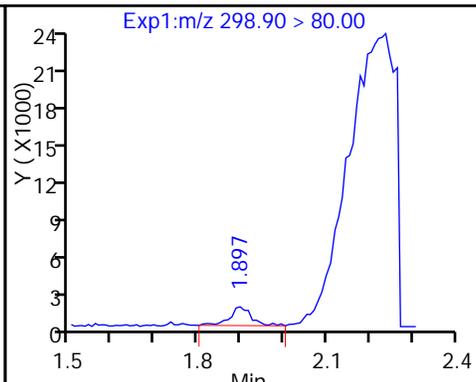
3 Perfluoropentanoic acid



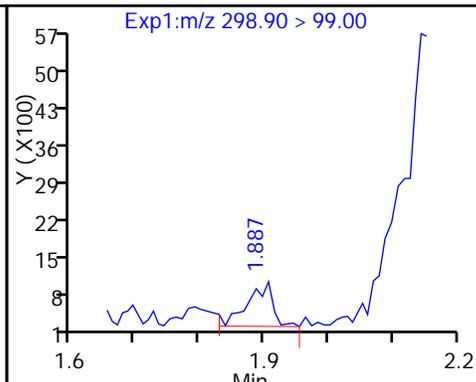
D 4 13C5-PFPeA



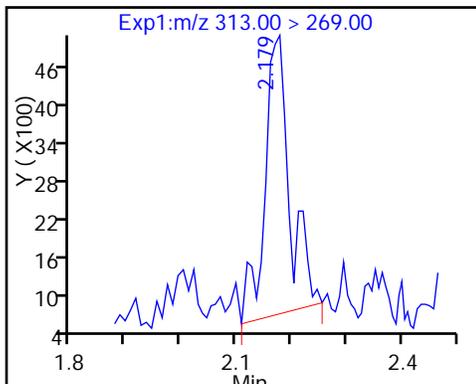
5 Perfluorobutanesulfonic acid



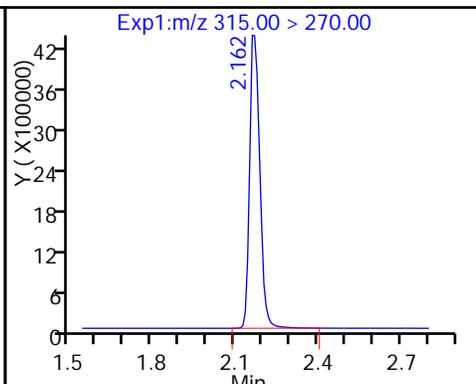
5 Perfluorobutanesulfonic acid



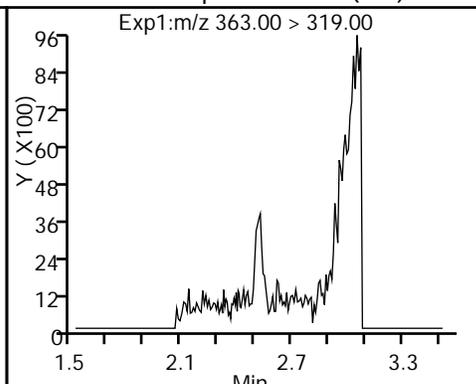
7 Perfluorohexanoic acid



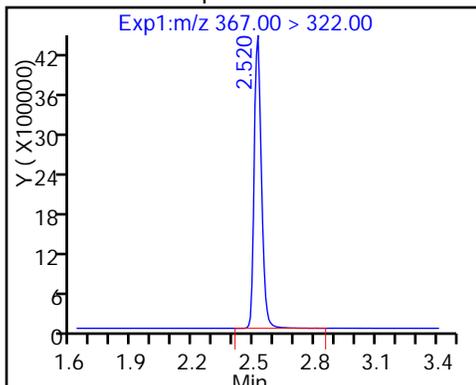
D 6 13C2 PFHxA



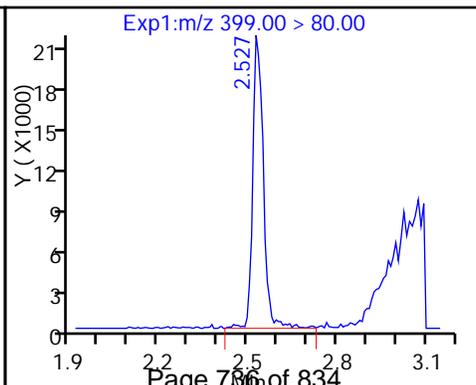
12 Perfluoroheptanoic acid (ND)



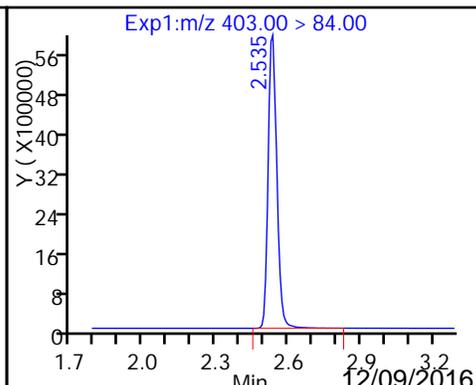
D 11 13C4-PFHpA



9 Perfluorohexanesulfonic acid



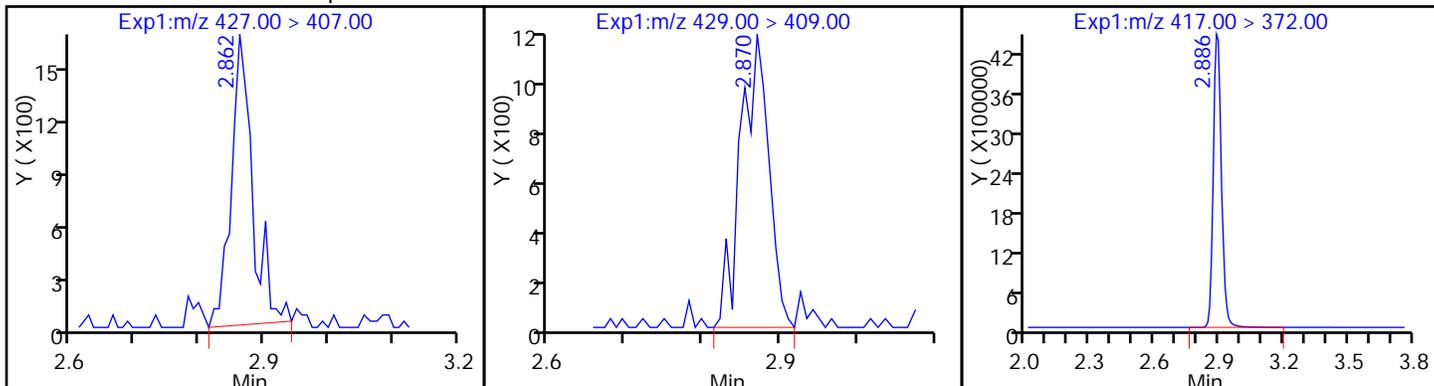
D 10 18O2 PFHxS



48 Sodium 1H,1H,2H,2H-perfluorooctanoate

D 47 M2-6:2FTS

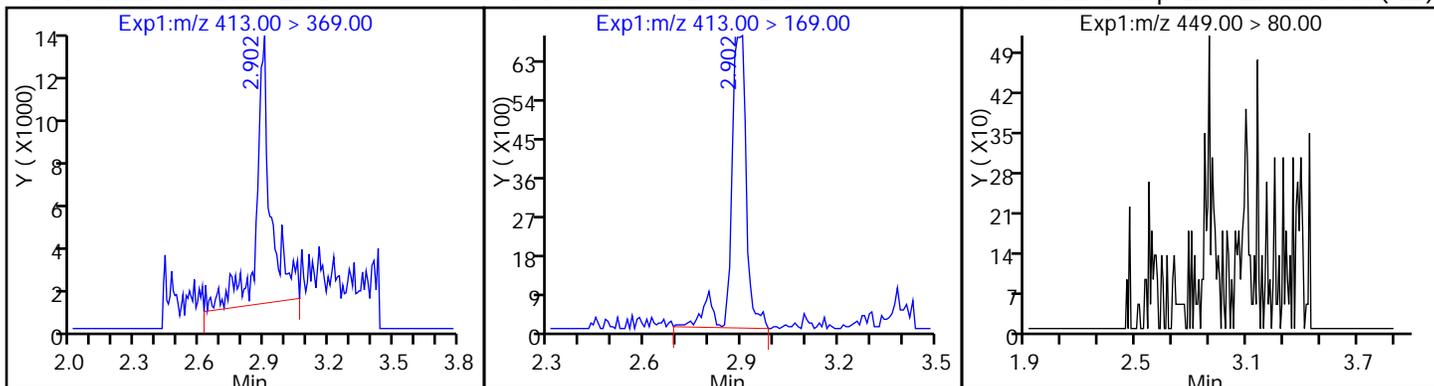
D 14 13C4 PFOA



15 Perfluorooctanoic acid

15 Perfluorooctanoic acid

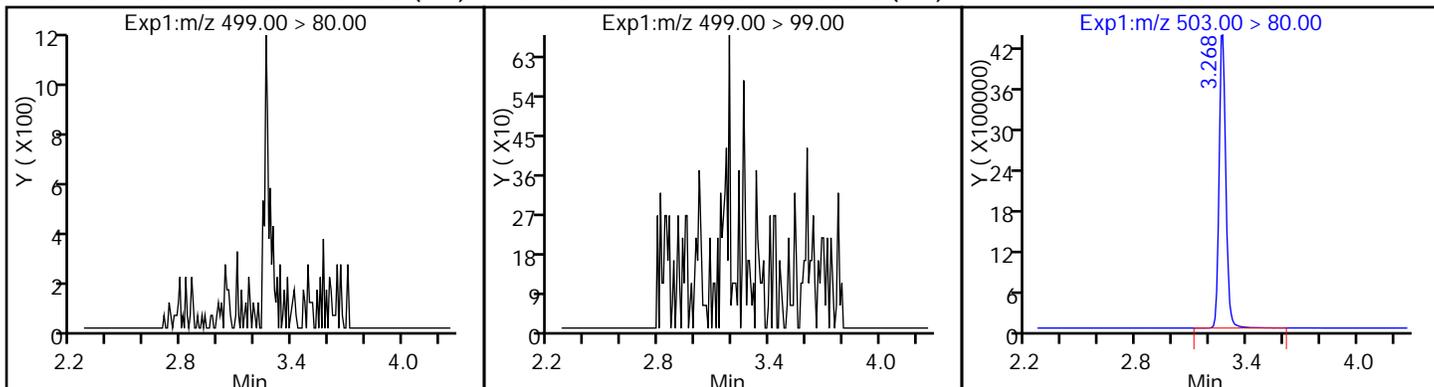
13 Perfluoroheptanesulfonic Acid (ND)



18 Perfluorooctane sulfonic acid (ND)

18 Perfluorooctane sulfonic acid (ND)

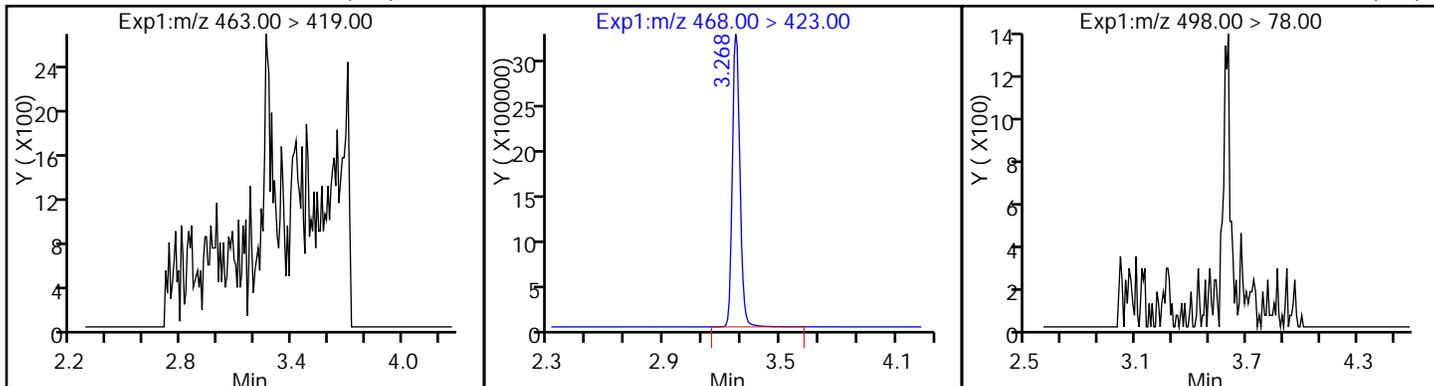
D 17 13C4 PFOS



20 Perfluorononanoic acid (ND)

D 19 13C5 PFNA

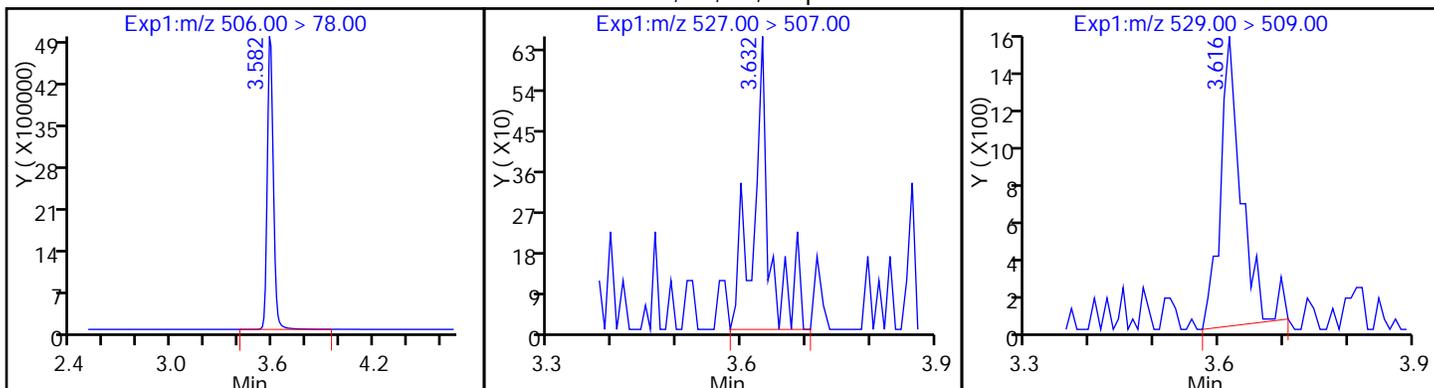
22 Perfluorooctane Sulfonamide (ND)



D 21 13C8 FOSA

43 Sodium 1H,1H,2H,2H-perfluorooctane

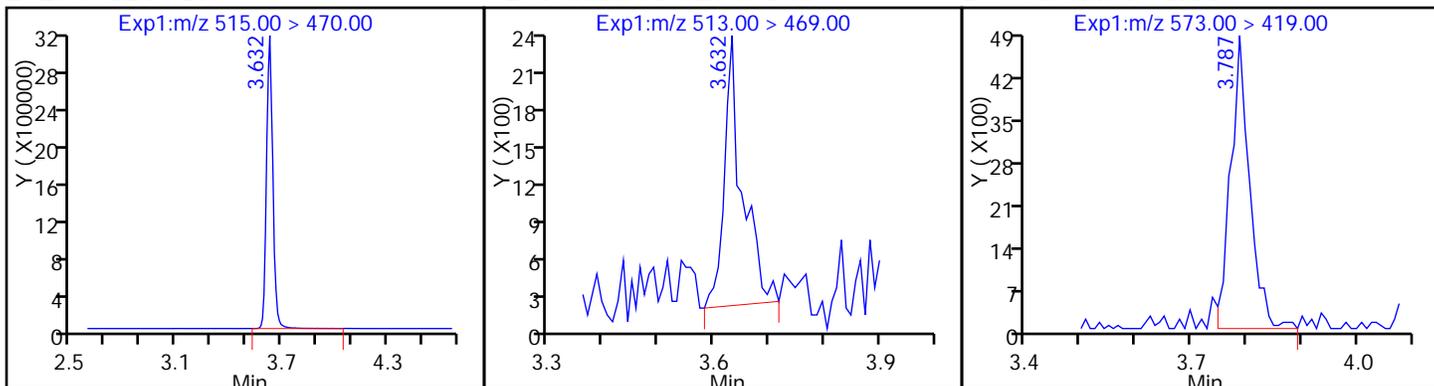
De 42 M2-8:2FTS



D 23 13C2 PFDA

24 Perfluorodecanoic acid

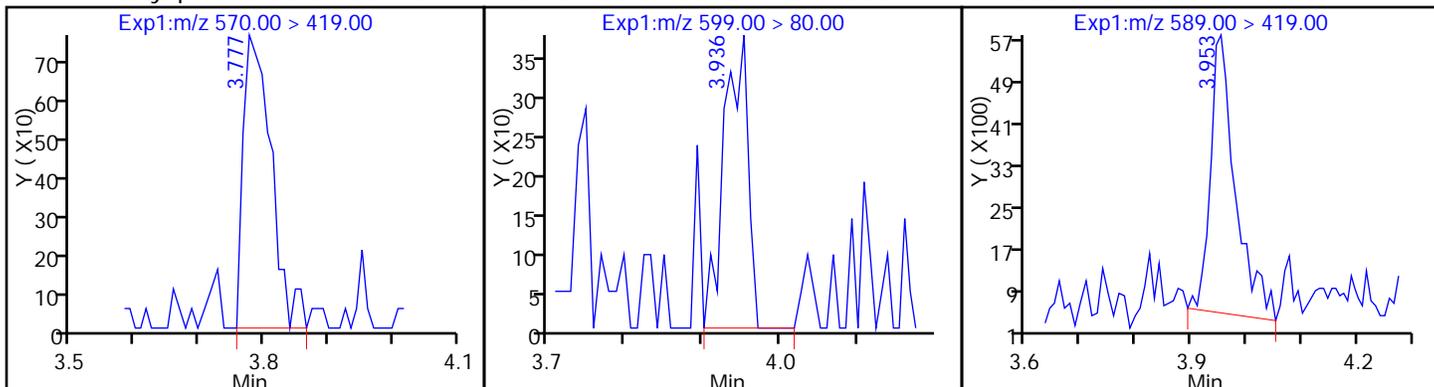
D 45 d3-NMeFOSAA



44 N-methyl perfluorooctane sulfonami

26 Perfluorodecane Sulfonic acid

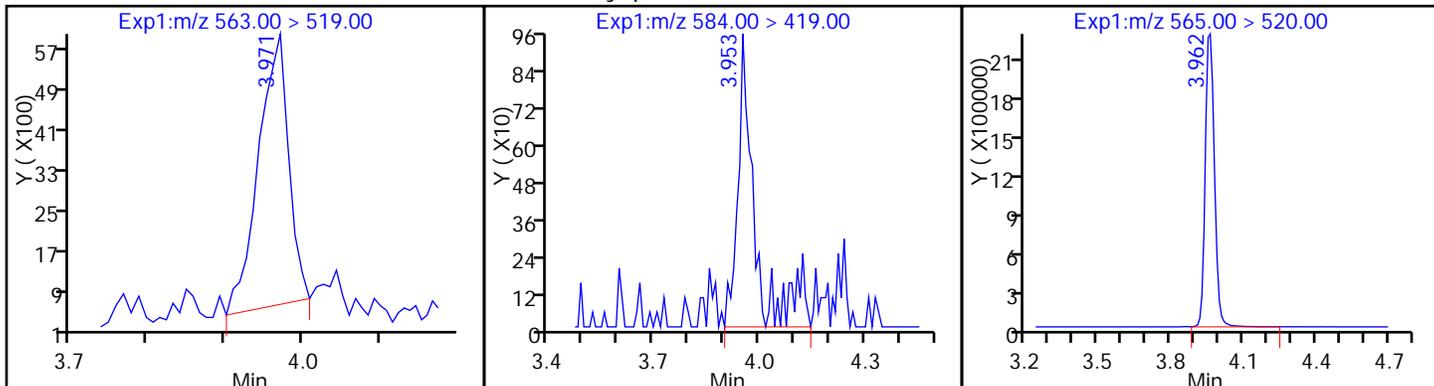
D 46 d5-NEtFOSAA



28 Perfluoroundecanoic acid

49 N-ethyl perfluorooctane sulfonamid

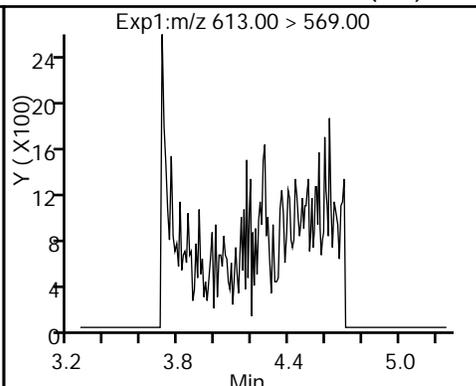
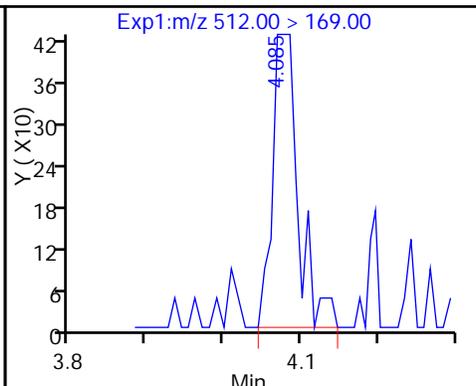
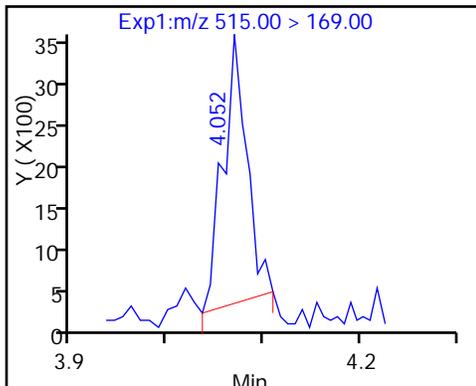
D 27 13C2 PFUnA



D 52 d-N-MeFOSA-M

54 MeFOSA

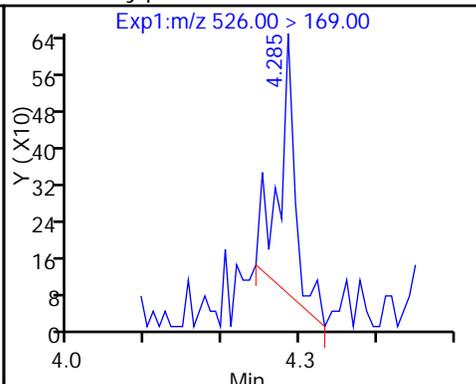
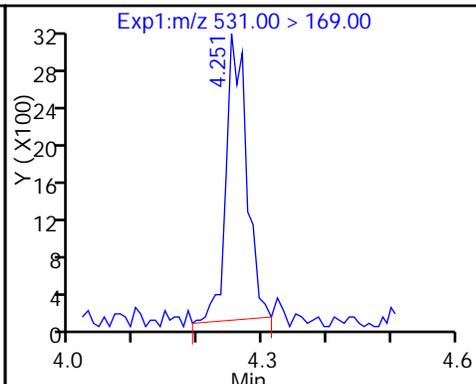
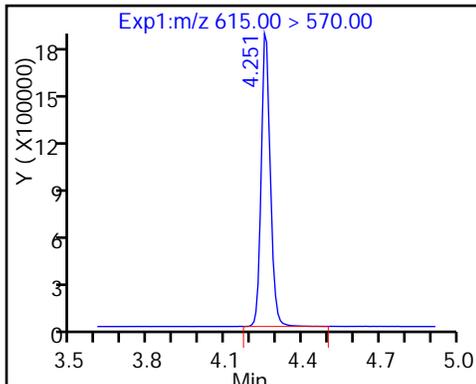
29 Perfluorododecanoic acid (ND)



D 30 13C2 PFDaA

D 51 d-N-EtFOSA-M

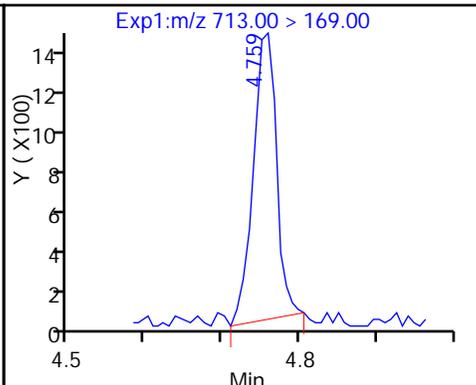
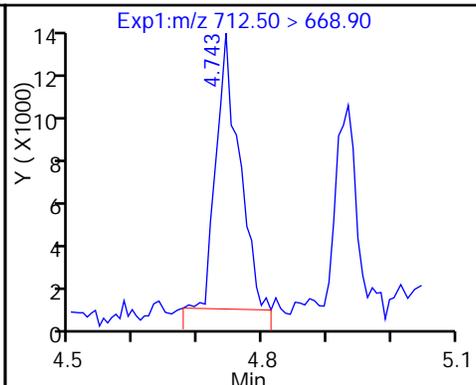
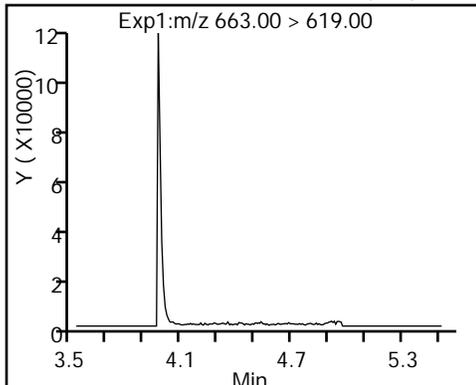
53 N-ethylperfluoro-1-octanesulfonami



31 Perfluorotridecanoic acid (ND)

33 Perfluorotetradecanoic acid

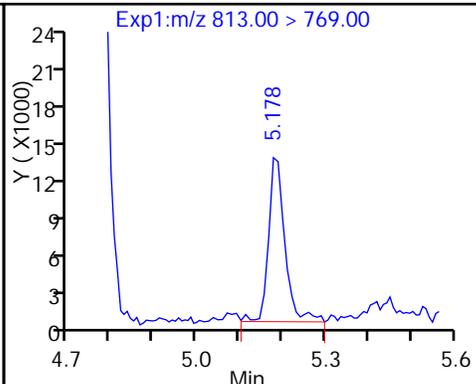
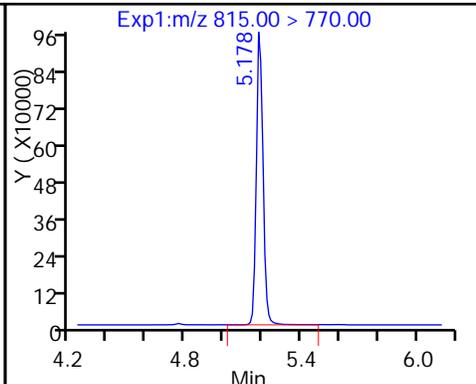
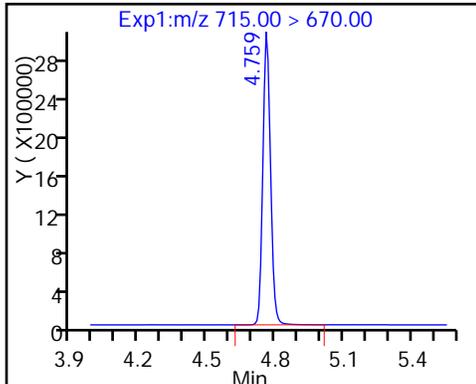
33 Perfluorotetradecanoic acid



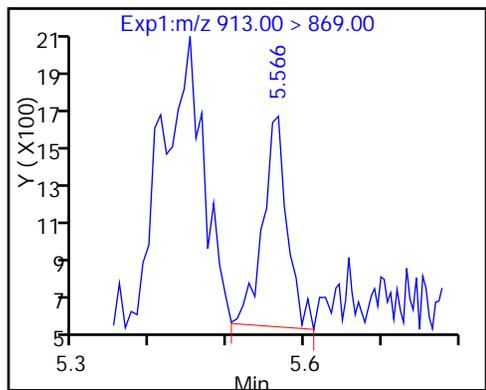
D 32 13C2-PFTeDA

D 34 13C2-PFHxDA

35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: MB 320-139749/1-A
 Matrix: Water Lab File ID: 03DEC2016B_001.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 250 (mL) Date Analyzed: 12/03/2016 16:03
 Con. Extract Vol.: 0.5 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.0030	U	0.0040	0.0013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	93		25-150
STL00991	13C4 PFOS	85		25-150
STL00994	18O2 PFHxS	86		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_001.d
 Lims ID: MB 320-139749/1-A
 Client ID:
 Sample Type: MB
 Inject. Date: 03-Dec-2016 16:03:39 ALS Bottle#: 32 Worklist Smp#: 22
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: mb 320-139749/1-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:30:02 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:24:04

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.566	1.574	-0.008	15226200	45.3		90.7	1569873	
1 Perfluorobutyric acid	212.90 > 169.00	1.598	1.577	0.021	1.000	26008	0.0977		100	
3 Perfluoropentanoic acid	262.90 > 219.00	1.848	1.861	-0.013	1.000	21886	0.0881		156	
D 4 13C5-PFPeA	267.90 > 223.00	1.848	1.861	-0.013		12242552	46.3	92.6	1166805	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.887	1.900	-0.013	1.000	5587	0.0130			
	298.90 > 99.00	1.868	1.900	-0.032	0.990	2503	2.23(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.158	2.164	-0.006	1.000	10301	0.0520		268	
D 6 13C2 PFHxA	315.00 > 270.00	2.150	2.164	-0.014		10398103	43.8	87.6	651663	
D 11 13C4-PFHpA	367.00 > 322.00	2.484	2.511	-0.027		9778883	47.1	94.3	1122402	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.492	2.512	-0.020	1.000	5152	0.0256		98.2	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.507	2.531	-0.024	1.000	43831	0.1481			
D 10 18O2 PFHxS	403.00 > 84.00	2.507	2.531	-0.024		12757089	40.8	86.3	681764	
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.830	2.836	-0.006	1.000	5869	NR			
D 47 M2-6:2FTS	429.00 > 409.00	2.839	2.836	0.003		2424	0.0230	0.0		
D 14 13C4 PFOA	417.00 > 372.00	2.855	2.880	-0.025		10158200	46.3	92.7	723151	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.855	2.887	-0.032	1.000	62196	0.2856			609	
413.00 > 169.00	2.871	2.887	-0.016	1.006	23771		2.62(0.90-1.10)		840	
D 17 13C4 PFOS										
503.00 > 80.00	3.235	3.259	-0.024		10026681	40.8		85.3	575603	
D 19 13C5 PFNA										
468.00 > 423.00	3.235	3.263	-0.028		7793773	46.8		93.7	2639754	
D 21 13C8 FOSA										
506.00 > 78.00	3.549	3.571	-0.022		2561806	6.37		12.7	217630	
43 Sodium 1H,1H,2H,2H-perfluorooctane										
527.00 > 507.00	3.583	3.588	-0.005	0.995	686	NR				
D 42 M2-8:2FTS										
529.00 > 509.00	3.599	3.591	0.008		1294	0.0133		0.0		
24 Perfluorodecanoic acid										
513.00 > 469.00	3.650	3.623	0.027	1.000	472	0.003505			31.5	
D 23 13C2 PFDA										
515.00 > 470.00	3.591	3.626	-0.035		7009564	44.4		88.8	202735	
D 45 d3-NMeFOSAA										
573.00 > 419.00	3.738	3.756	-0.018		8153	0.1124		0.0		
44 N-methyl perfluorooctane sulfonami										
570.00 > 419.00	3.748	3.759	-0.011	1.003	2600	NR				
D 46 d5-NEtFOSAA										
589.00 > 419.00	3.917	3.922	-0.005		12565	0.1577		0.0		
49 N-ethyl perfluorooctane sulfonamid										
584.00 > 419.00	3.917	3.929	-0.012	1.000	5719	NR				
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.908	3.936	-0.028	1.000	828	0.006170				
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.926	3.955	-0.029	1.000	16456	0.1458			370	
D 27 13C2 PFUnA										
565.00 > 520.00	3.926	3.958	-0.032		5294647	44.6		89.2	392201	
D 52 d-N-MeFOSA-M										
515.00 > 169.00	4.040	4.051	-0.011		3240	0.0308		0.0		
54 MeFOSA										
512.00 > 169.00	4.064	4.057	0.007	1.000	467	NR				
D 51 d-N-EtFOSA-M										
531.00 > 169.00	4.225	4.239	-0.014		3131	0.0316		0.0		
53 N-ethylperfluoro-1-octanesulfonami										
526.00 > 169.00	4.450	4.245	0.205	1.000	408	NR				
D 30 13C2 PFDoA										
615.00 > 570.00	4.216	4.251	-0.035		4946893	44.1		88.3	145179	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.723	4.759	-0.036		11899994	51.5		103	1138196	
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.723	4.761	-0.038	1.000	38242	0.2085			25.3	
713.00 > 169.00	4.715	4.761	-0.046	0.998	5325		7.18(0.00-0.00)		2143	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.148	5.186	-0.038		5517123	42.5		85.1	232893	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
35 Perfluorohexadecanoic acid	813.00 > 769.00	5.148	5.186	-0.038	1.000	72544	-0.0159		188	
36 Perfluorooctadecanoic acid	913.00 > 869.00	5.694	5.559	0.135	1.000	273	0.002779		1.6	

QC Flag Legend

Processing Flags

NR - Missing Quant Standard

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_001.d

Injection Date: 03-Dec-2016 16:03:39

Instrument ID: A8_N

Lims ID: MB 320-139749/1-A

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 32

Worklist Smp#: 22

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

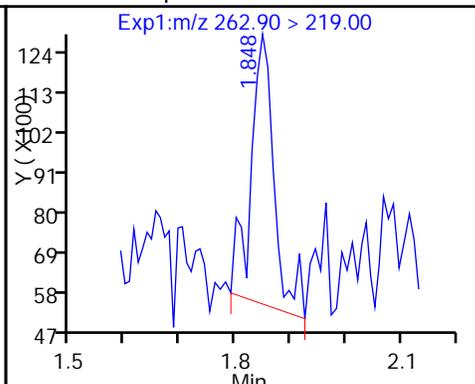
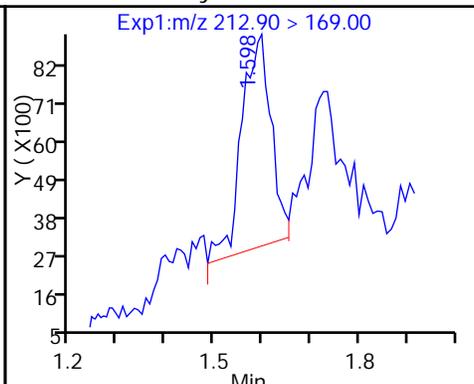
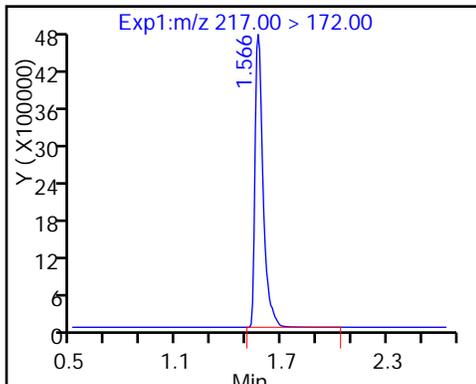
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

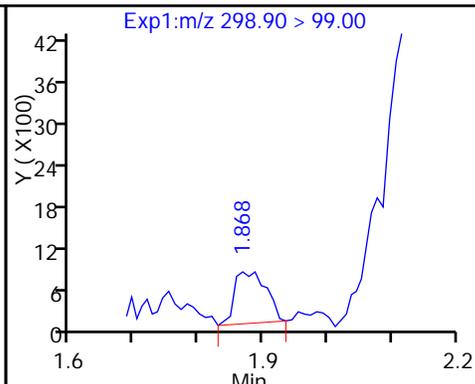
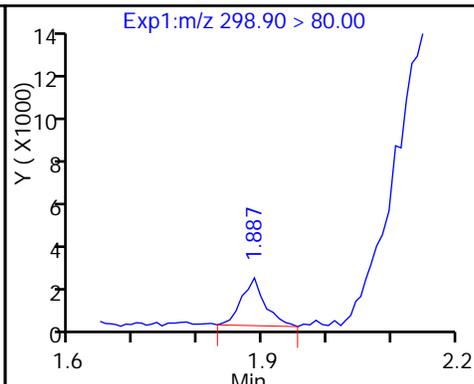
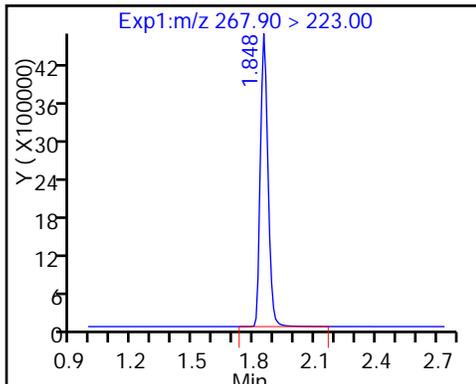
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

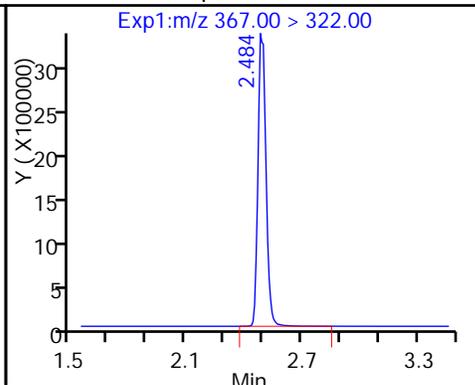
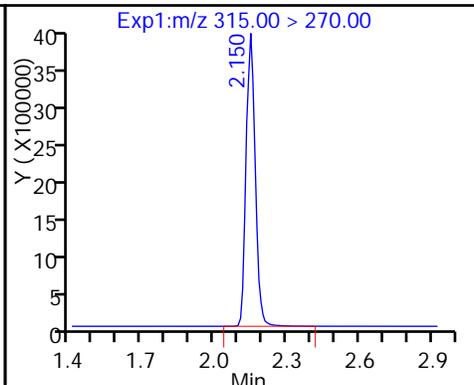
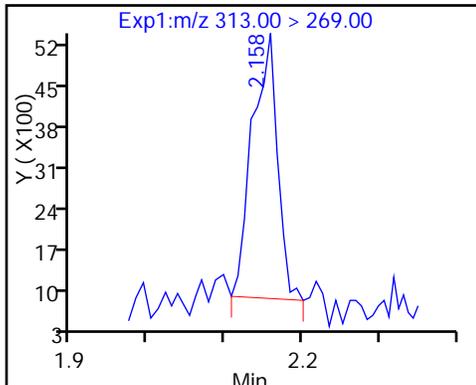
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

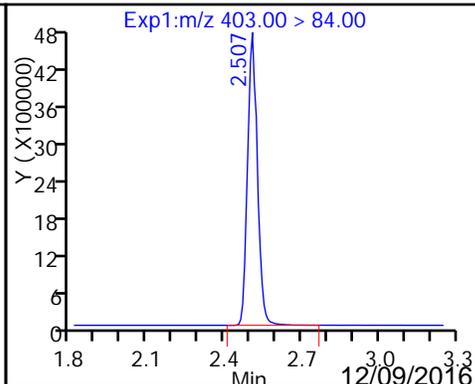
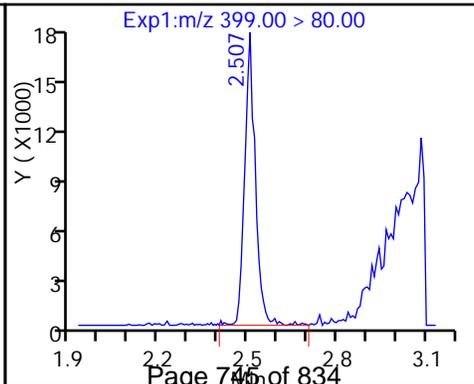
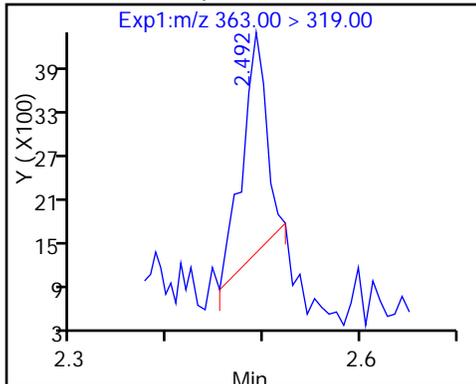
D 11 13C4-PFHpA



12 Perfluoroheptanoic acid

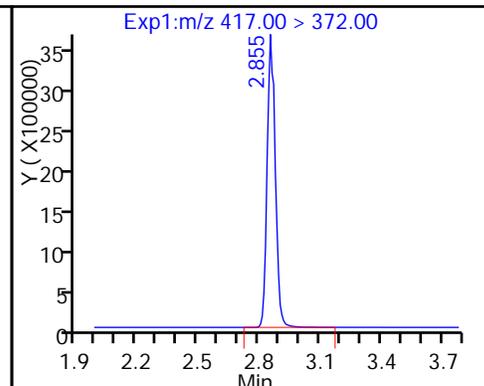
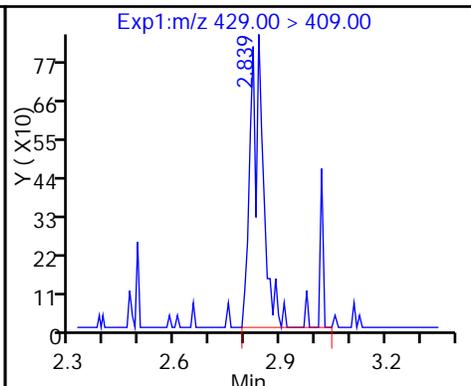
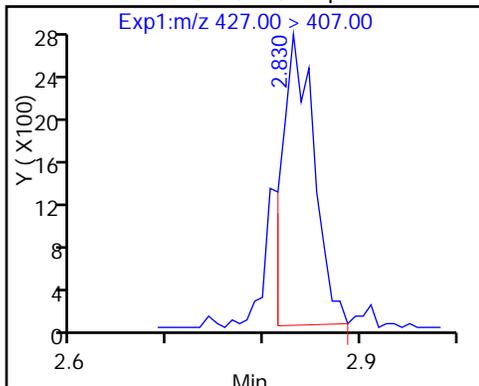
9 Perfluorohexanesulfonic acid

D 10 18O2 PFHxS



48 Sodium 1H,1H,2H,2H-perfluorooctanoate

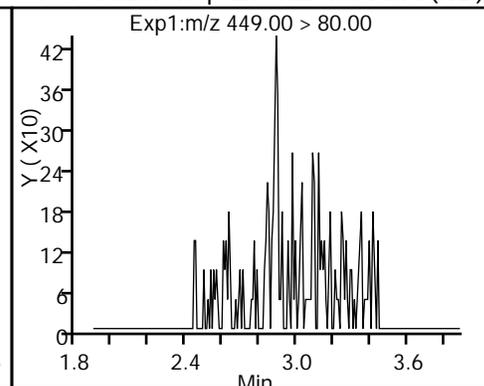
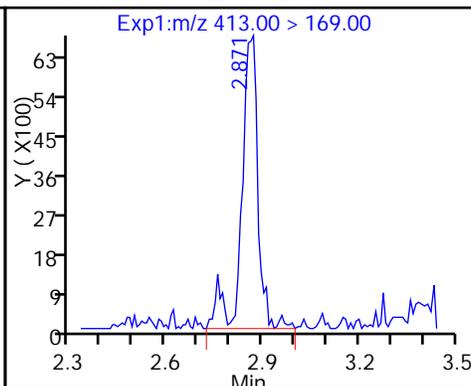
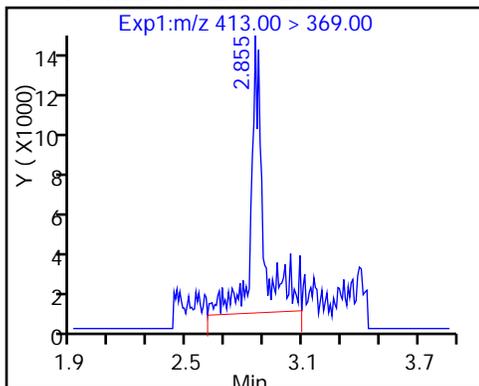
D 14 13C4 PFOA



15 Perfluorooctanoic acid

15 Perfluorooctanoic acid

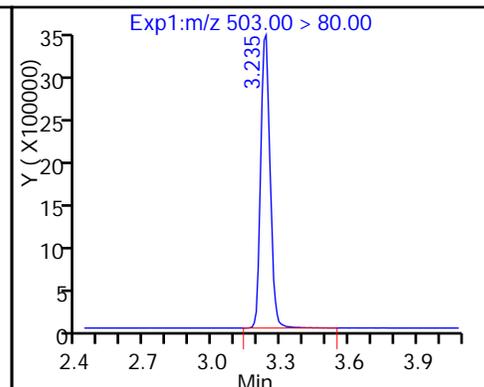
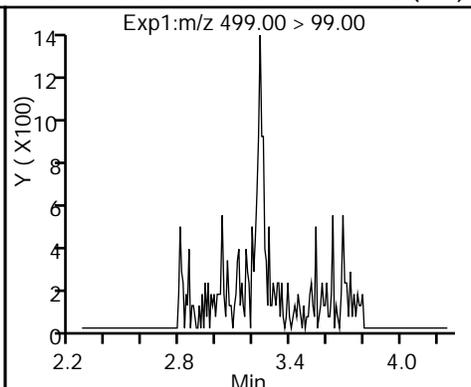
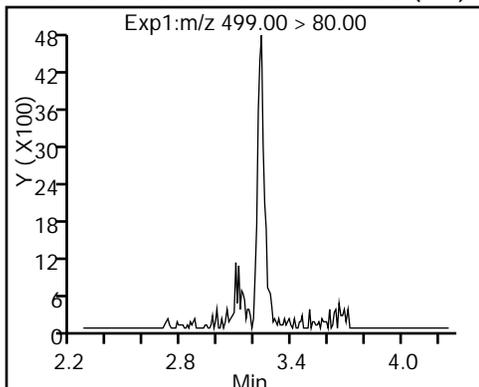
13 Perfluoroheptanesulfonic Acid (ND)



18 Perfluorooctane sulfonic acid (ND)

18 Perfluorooctane sulfonic acid (ND)

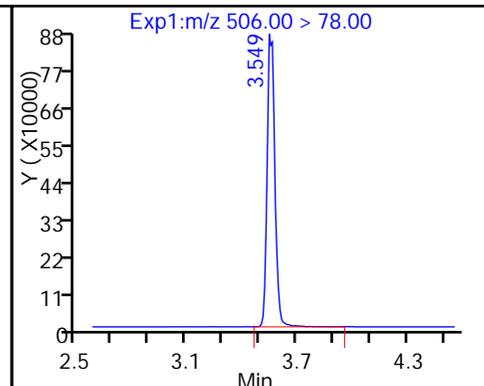
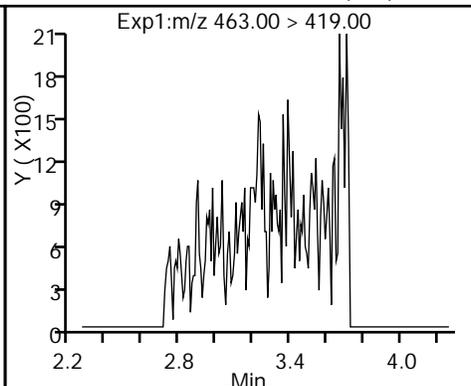
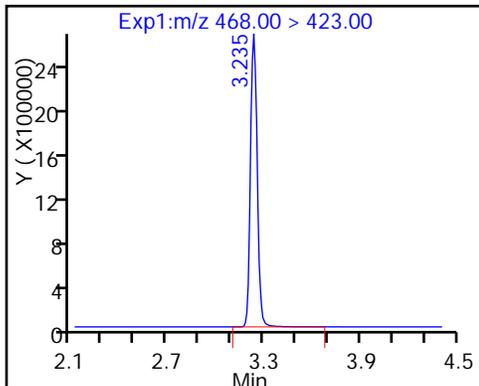
D 17 13C4 PFOS



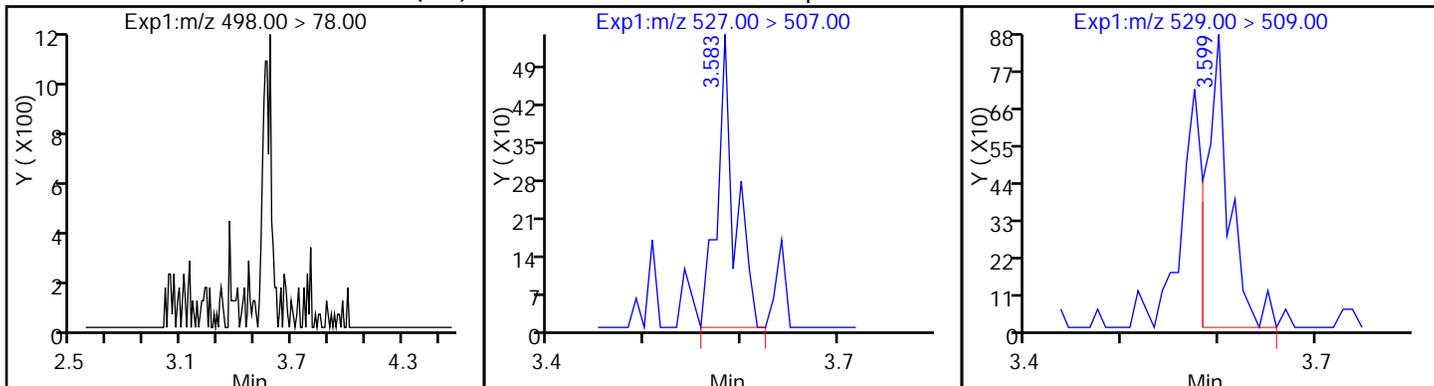
D 19 13C5 PFNA

20 Perfluorononanoic acid (ND)

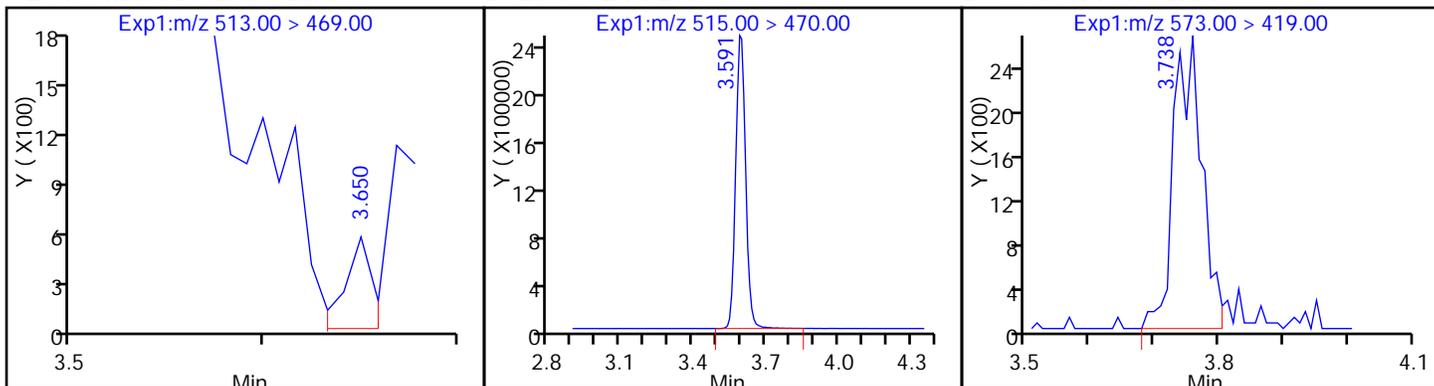
D 21 13C8 FOSA



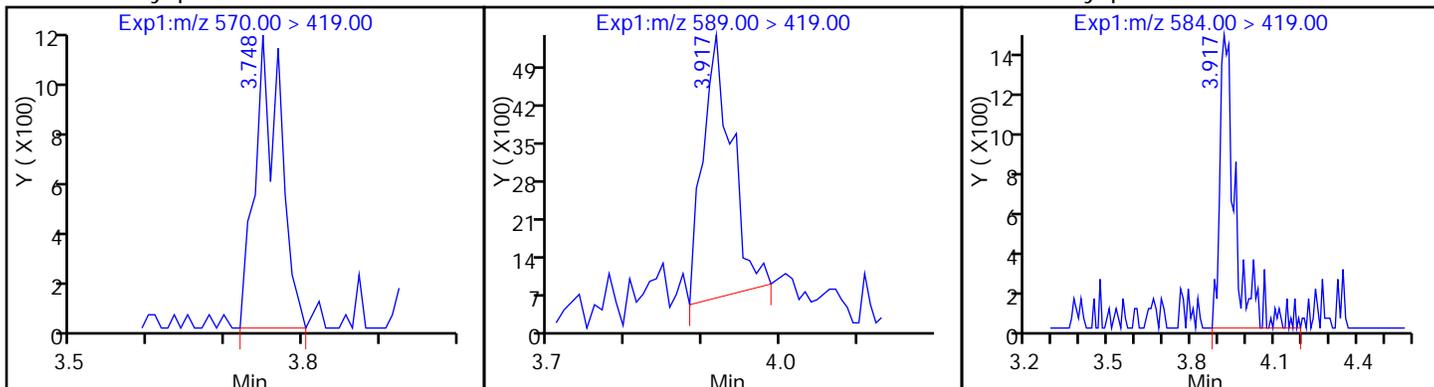
22 Perfluorooctane Sulfonamide (ND) 43 Sodium 1H,1H,2H,2H-perfluorooctane Sulfonamide 42 M2-8:2FTS



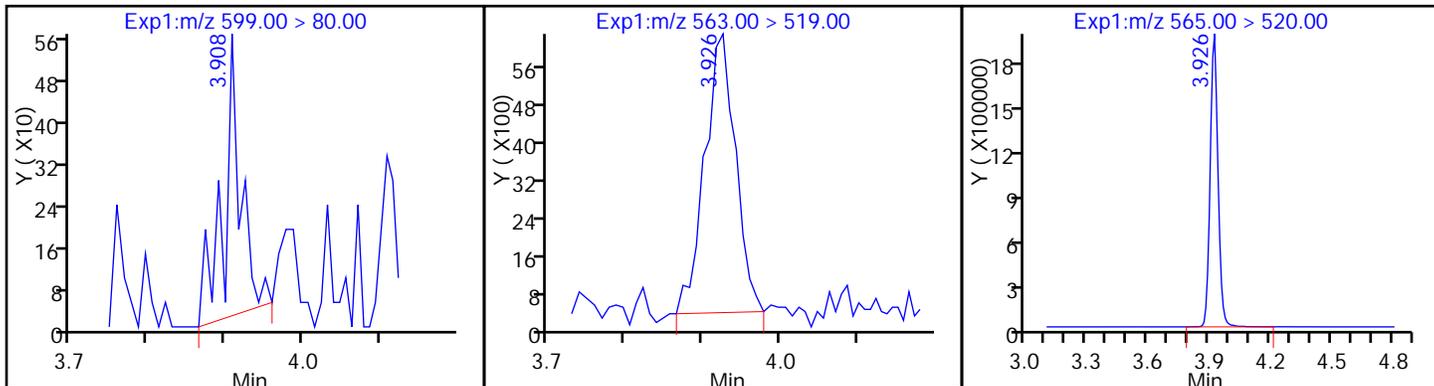
24 Perfluorodecanoic acid D 23 13C2 PFDA D 45 d3-NMeFOSAA



44 N-methyl perfluorooctane sulfonamid D 46 d5-NEtFOSAA 49 N-ethyl perfluorooctane sulfonamid



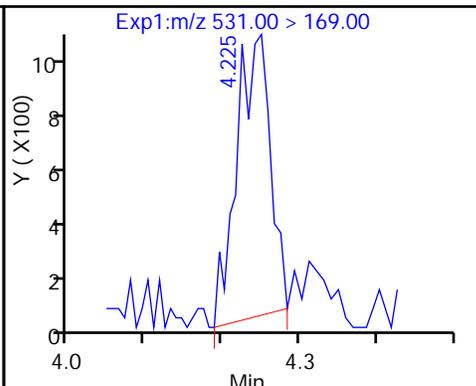
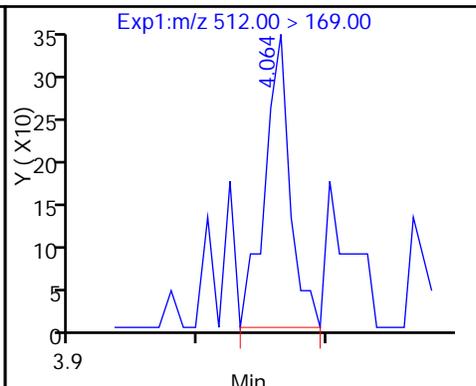
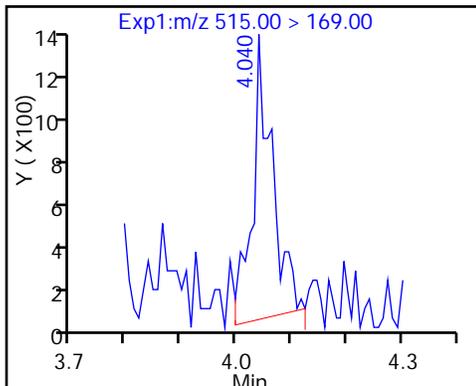
26 Perfluorodecane Sulfonic acid 28 Perfluoroundecanoic acid D 27 13C2 PFUnA



D 52 d-N-MeFOSA-M

54 MeFOSA

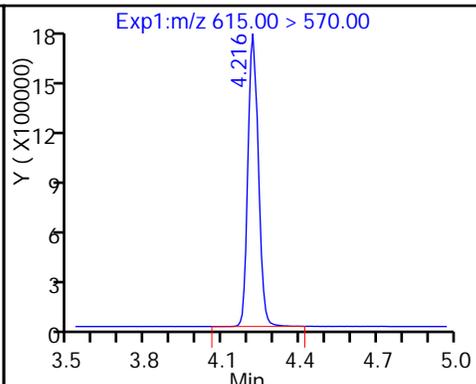
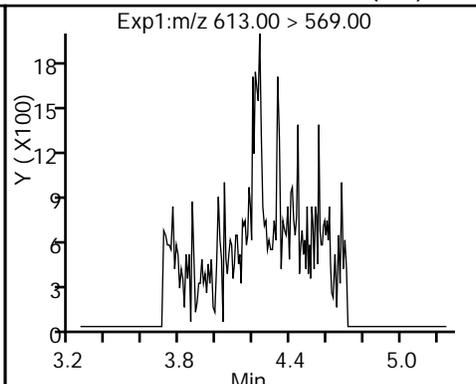
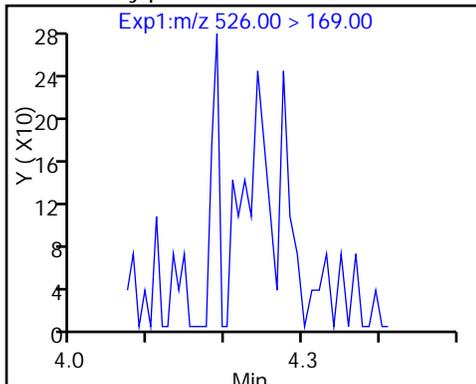
D 51 d-N-EtFOSA-M



53 N-ethylperfluoro-1-octanesulfonami

29 Perfluorododecanoic acid (ND)

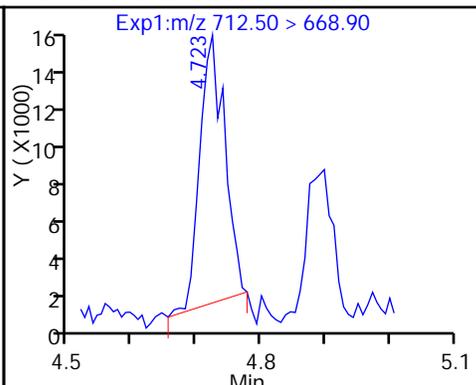
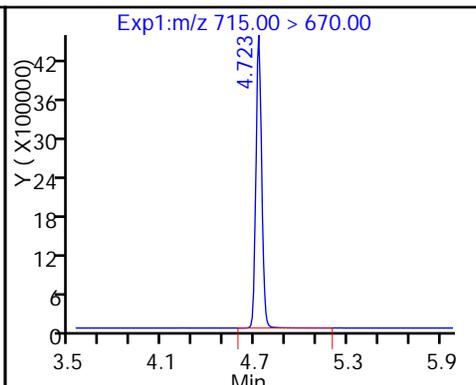
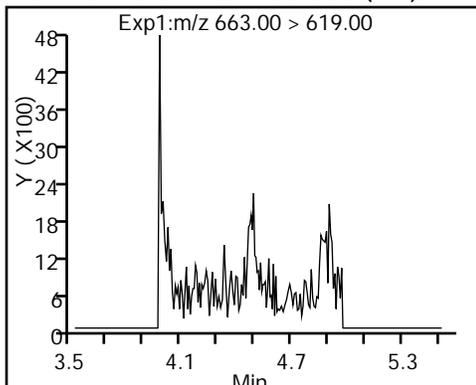
D 30 13C2 PFDaA



31 Perfluorotridecanoic acid (ND)

D 32 13C2-PFTeDA

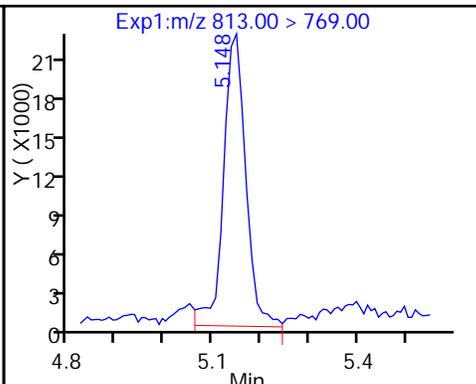
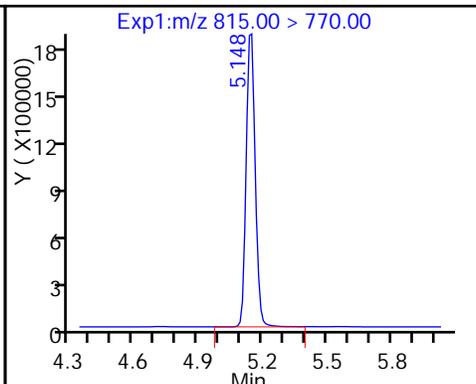
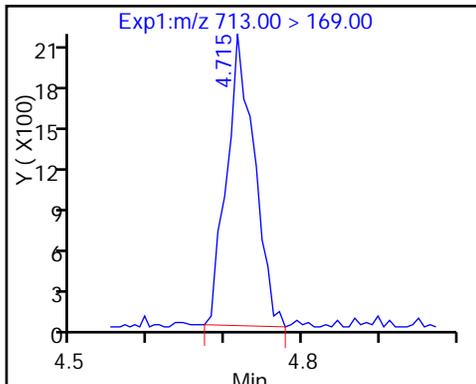
33 Perfluorotetradecanoic acid



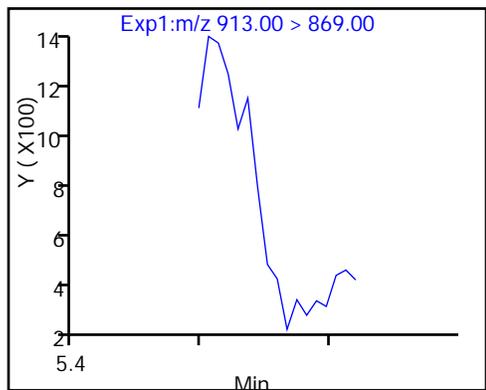
33 Perfluorotetradecanoic acid

D 34 13C2-PFHxDA

35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: ICB 320-140564/10
 Matrix: Water Lab File ID: 03DEC2016A_010.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: _____ Date Extracted: _____
 Sample wt/vol: 1(mL) Date Analyzed: 12/03/2016 14:33
 Con. Extract Vol.: _____ Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ng/mL

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.204	J	0.50	0.10
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.30	U	0.50	0.13
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.30	U	0.40	0.10

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	102		25-150
STL00991	13C4 PFOS	97		25-150
STL00994	18O2 PFHxS	102		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_010.d
 Lims ID: ICB
 Client ID:
 Sample Type: ICB
 Inject. Date: 03-Dec-2016 14:33:43 ALS Bottle#: 54 Worklist Smp#: 10
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: ICB_b
 Misc. Info.: Plate: 1 Rack: 1
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:26:19 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 09:46:44

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	16096656	47.9		95.9	1325245	
3 Perfluoropentanoic acid	262.90 > 219.00	1.849	1.861	-0.013	1.000	26187	0.1009		206	
D 4 13C5-PFPeA	267.90 > 223.00	1.858	1.861	-0.003	12785749	48.3		96.7	900580	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.897	1.900	-0.003	1.000	11990	0.0236			
	298.90 > 99.00	1.887	1.900	-0.013	0.995	5171	2.32(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.161	2.164	-0.003	1.000	14358	0.0642		402	
D 6 13C2 PFHxA	315.00 > 270.00	2.161	2.164	-0.003	11724024	49.4		98.7	702569	
D 11 13C4-PFHpA	367.00 > 322.00	2.504	2.511	-0.007	10534496	50.8		102	767171	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.504	2.512	-0.008	1.000	46404	0.2144		599	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.520	2.531	-0.011	1.000	57809	0.1658			
D 10 18O2 PFHxS	403.00 > 84.00	2.520	2.531	-0.011	15029673	48.1		102	1415639	
48 Sodium 1H,1H,2H,2H-perfluorooctane	427.00 > 407.00	2.837	2.836	0.001	1.000	14959	0.1994			
D 47 M2-6:2FTS	429.00 > 409.00	2.846	2.836	0.010	4240985	40.2		84.6		
D 14 13C4 PFOA	417.00 > 372.00	2.870	2.880	-0.010	11233082	51.2		102	944277	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
15 Perfluorooctanoic acid										
413.00 > 369.00	2.886	2.887	-0.001	1.000	49093	0.2039			305	
413.00 > 169.00	2.878	2.887	-0.009	0.997	17821		2.75(0.90-1.10)		1369	
D 17 13C4 PFOS										
503.00 > 80.00	3.251	3.259	-0.008		11404796	46.4		97.0	680924	
D 19 13C5 PFNA										
468.00 > 423.00	3.251	3.263	-0.012		8205328	49.3		98.6	665084	
D 21 13C8 FOSA										
506.00 > 78.00	3.563	3.571	-0.008		19326715	48.0		96.1	590557	
22 Perfluorooctane Sulfonamide										
498.00 > 78.00	3.563	3.574	-0.011	1.000	33109	0.0917			856	
43 Sodium 1H,1H,2H,2H-perfluorooctane										
527.00 > 507.00	3.588	3.588	0.0	0.998	1181	0.0176				
D 42 M2-8:2FTS										
529.00 > 509.00	3.596	3.591	0.005		3746326	38.6		80.5		
24 Perfluorodecanoic acid										
513.00 > 469.00	3.621	3.623	-0.002	1.000	7365	0.0500			262	
D 23 13C2 PFDA										
515.00 > 470.00	3.613	3.626	-0.013		7674974	48.6		97.3	318943	
D 45 d3-NMeFOSAA										
573.00 > 419.00	3.764	3.756	0.008		3213300	44.3		88.6		
44 N-methyl perfluorooctane sulfonami										
570.00 > 419.00	3.802	3.759	0.043	1.010	251	0.004550				
D 46 d5-NEtFOSAA										
589.00 > 419.00	3.932	3.922	0.010		3437758	43.2		86.3		
49 N-ethyl perfluorooctane sulfonamid										
584.00 > 419.00	3.950	3.929	0.021	1.004	1568	0.0298				
26 Perfluorodecane Sulfonic acid										
599.00 > 80.00	3.906	3.936	-0.030	1.000	4850	0.0318				
28 Perfluoroundecanoic acid										
563.00 > 519.00	3.941	3.955	-0.014	1.000	19419	0.1546			520	
D 27 13C2 PFUnA										
565.00 > 520.00	3.941	3.958	-0.017		5892898	49.6		99.2	340109	
D 52 d-N-MeFOSA-M										
515.00 > 169.00	4.058	4.051	0.007		4941569	47.0		94.0		
54 MeFOSA										
512.00 > 169.00	4.058	4.057	0.001	1.000	247	0.003093				
D 51 d-N-EtFOSA-M										
531.00 > 169.00	4.237	4.239	-0.002		4774609	48.1		96.3		
53 N-ethylperfluoro-1-octanesulfonami										
526.00 > 169.00	4.246	4.245	0.001	1.000	660	0.008032				
D 30 13C2 PFDoA										
615.00 > 570.00	4.237	4.251	-0.014		5399581	48.2		96.3	213278	
31 Perfluorotridecanoic acid										
663.00 > 619.00	4.502	4.518	-0.016	1.000	14415	0.1405			97.0	
D 32 13C2-PFTeDA										
715.00 > 670.00	4.740	4.759	-0.019		11538350	49.9		99.8	814530	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
33 Perfluorotetradecanoic acid										
712.50 > 668.90	4.749	4.761	-0.012	1.000	45349	0.2265			25.0	
713.00 > 169.00	4.740	4.761	-0.021	0.998	10278		4.41(0.00-0.00)		1682	
D 34 13C2-PFHxDA										
815.00 > 770.00	5.165	5.186	-0.021		6026380	46.5		92.9	130593	
35 Perfluorohexadecanoic acid										
813.00 > 769.00	5.165	5.186	-0.021	1.000	112545	0.2952			116	
36 Perfluorooctadecanoic acid										
913.00 > 869.00	5.534	5.559	-0.025	1.000	65863	0.6143			98.5	

Reagents:

LCPFC-L0_00002

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_010.d

Injection Date: 03-Dec-2016 14:33:43

Instrument ID: A8_N

Lims ID: ICB

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 54

Worklist Smp#: 10

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

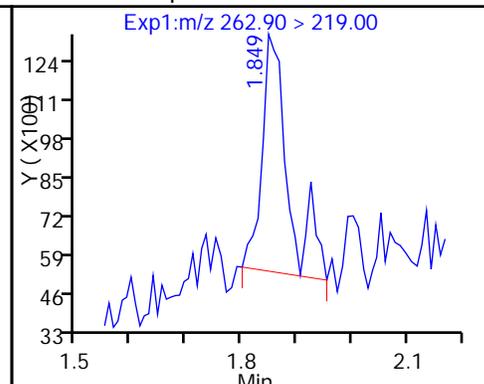
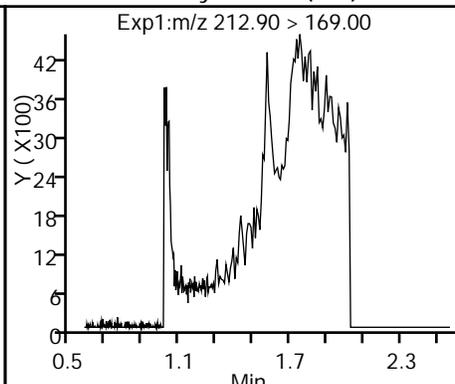
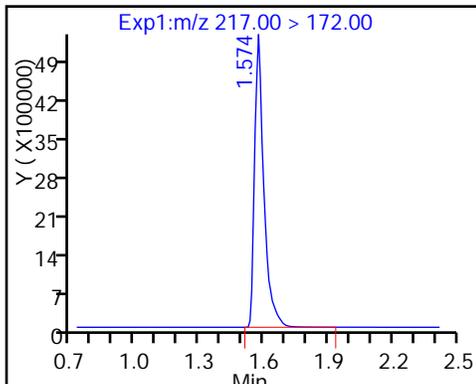
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid (ND)

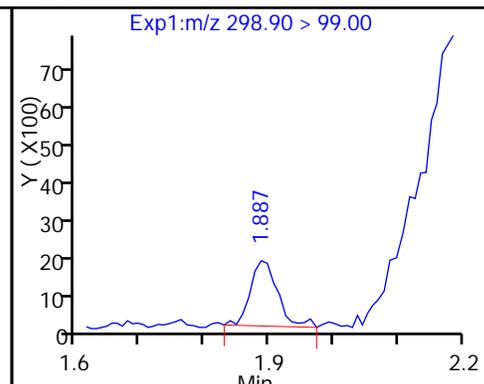
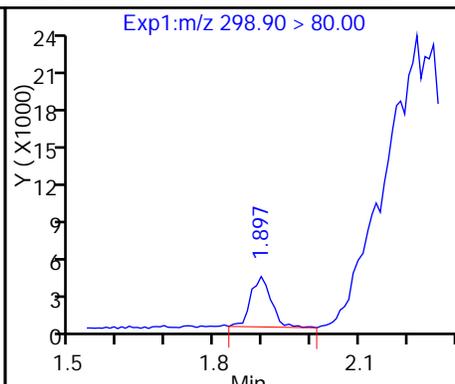
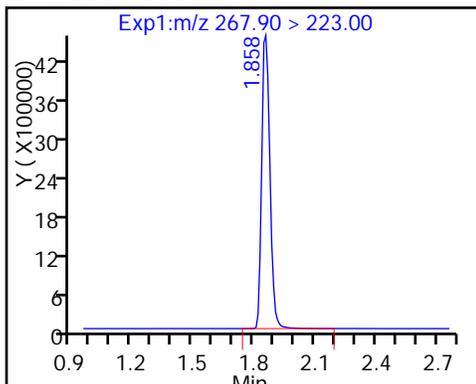
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

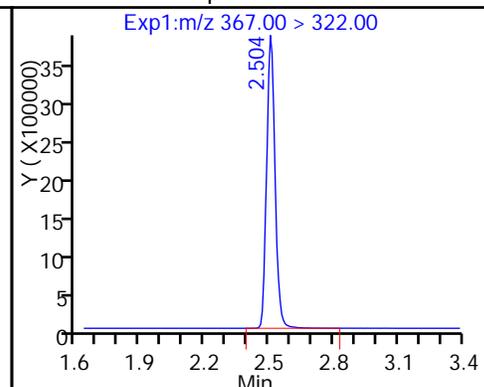
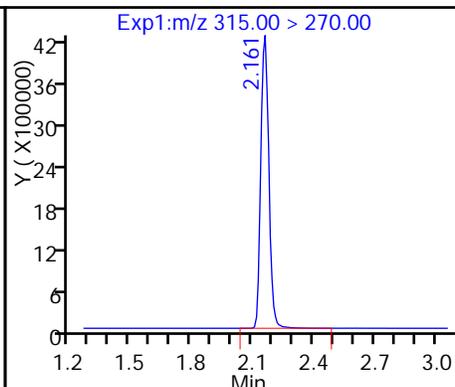
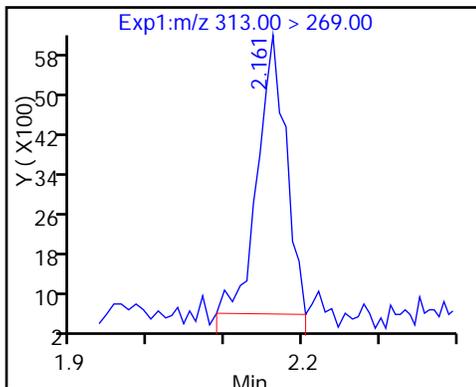
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

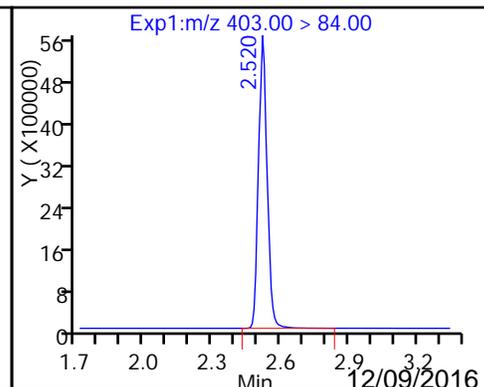
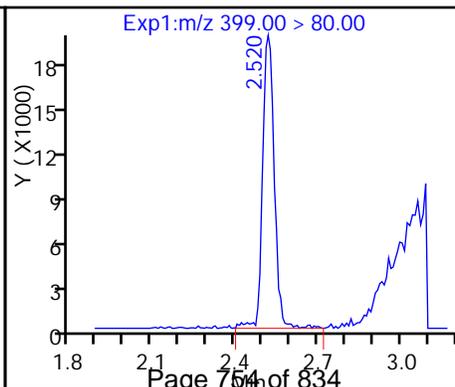
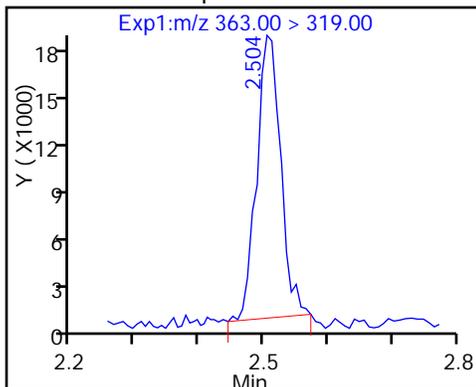
D 11 13C4-PFHpA



12 Perfluoroheptanoic acid

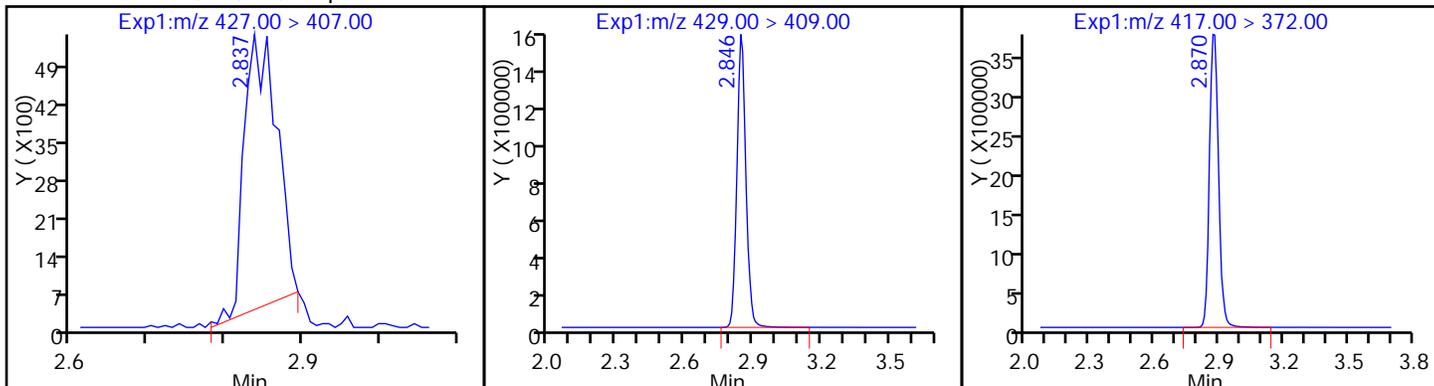
9 Perfluorohexanesulfonic acid

D 10 18O2 PFHxS



48 Sodium 1H,1H,2H,2H-perfluorooctanoate

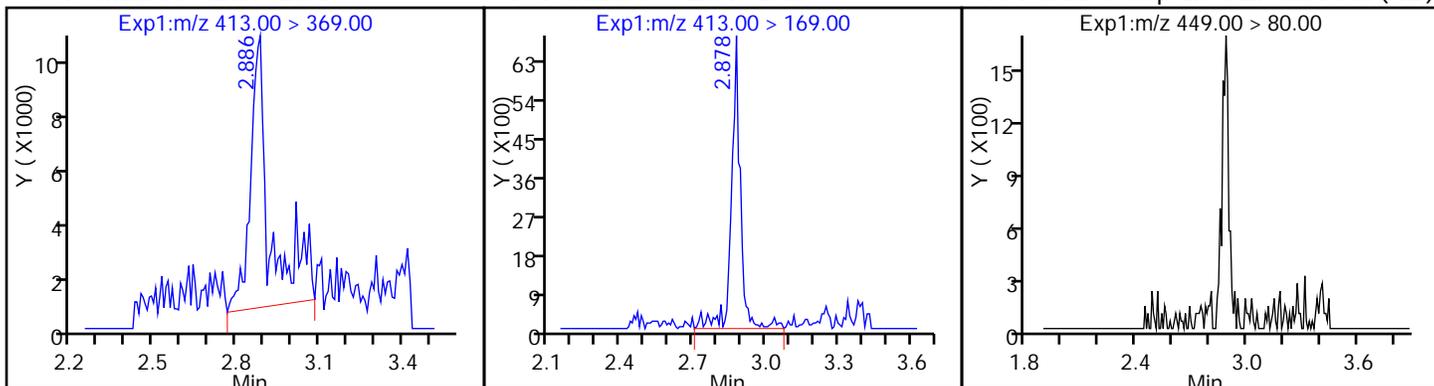
D 14 13C4 PFOA



15 Perfluorooctanoic acid

15 Perfluorooctanoic acid

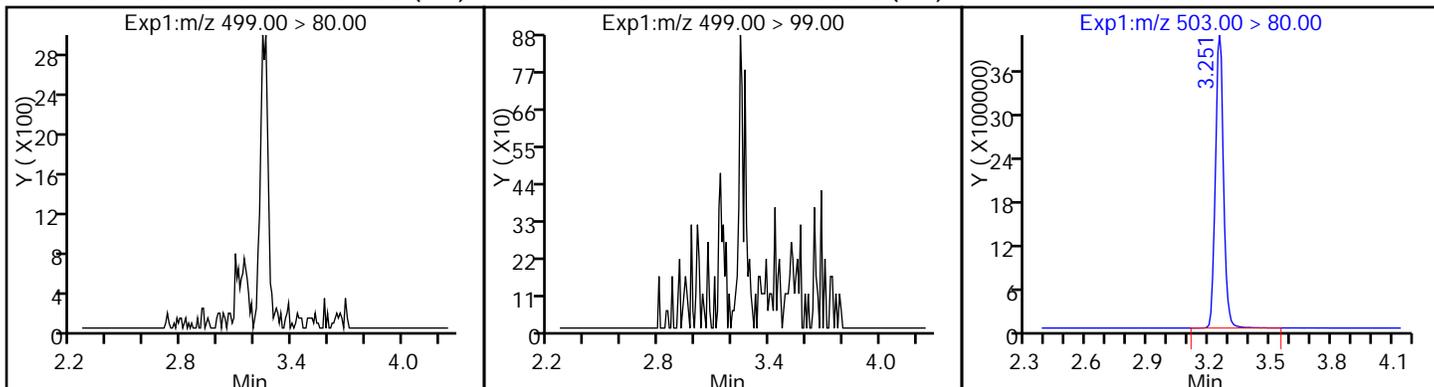
13 Perfluoroheptanesulfonic Acid (ND)



18 Perfluorooctane sulfonic acid (ND)

18 Perfluorooctane sulfonic acid (ND)

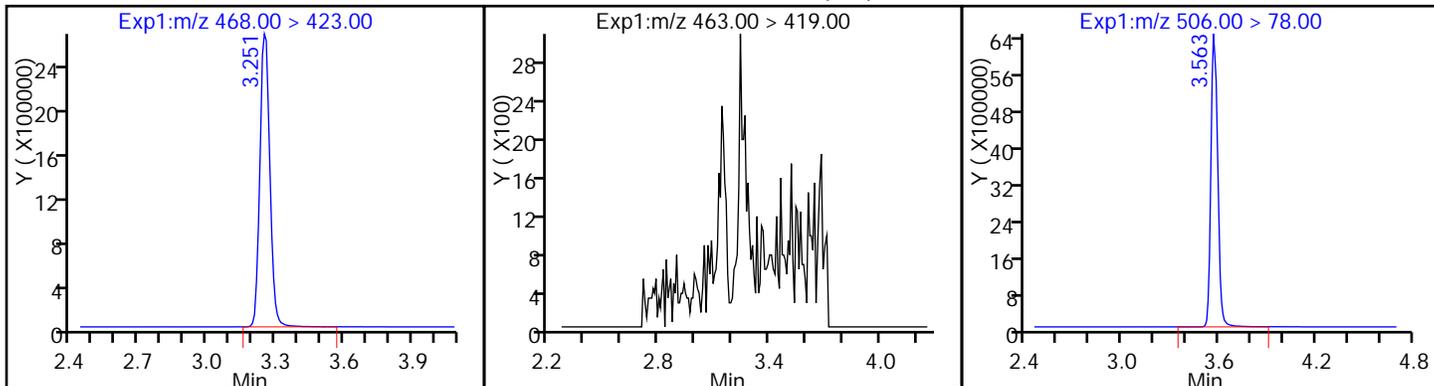
D 17 13C4 PFOS



D 19 13C5 PFNA

20 Perfluorononanoic acid (ND)

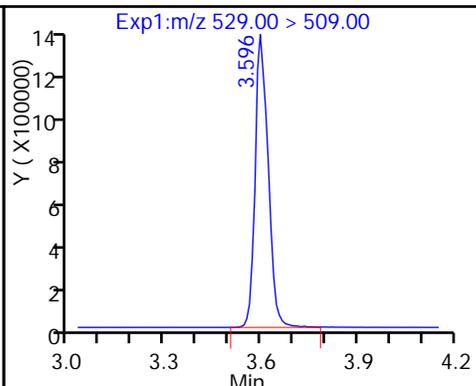
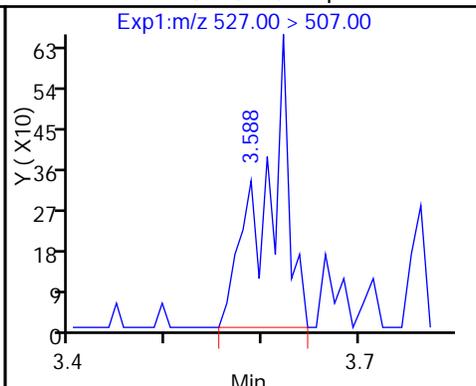
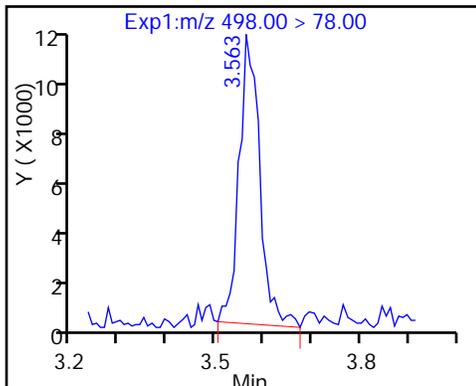
D 21 13C8 FOSA



22 Perfluorooctane Sulfonamide

43 Sodium 1H,1H,2H,2H-perfluorooctane

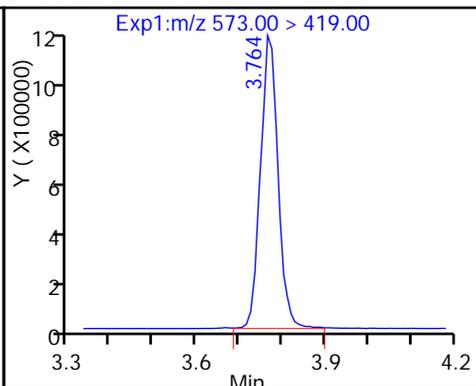
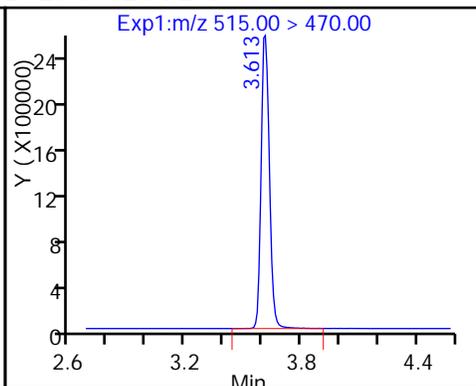
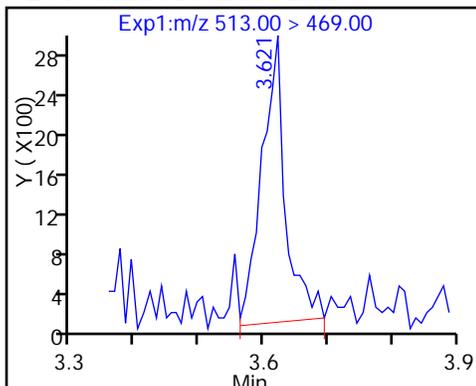
D 42 M2-8:2FTS



24 Perfluorodecanoic acid

D 23 13C2 PFDA

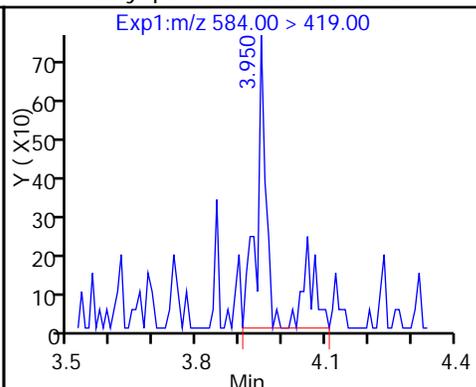
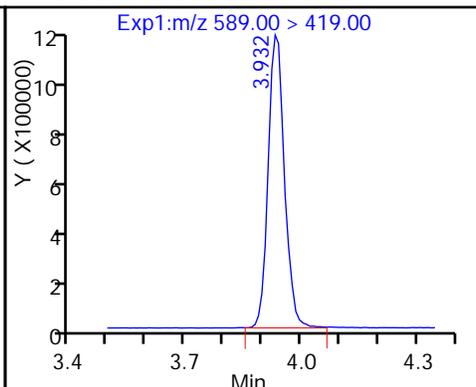
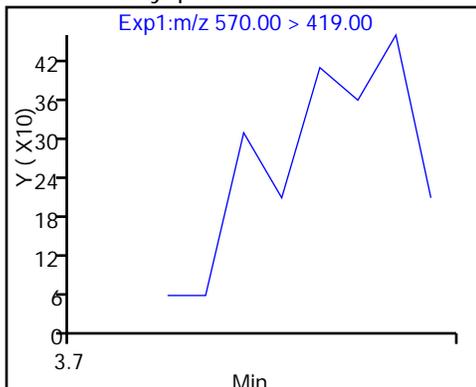
D 45 d3-NMeFOSAA



44 N-methyl perfluorooctane sulfonamid

D 46 d5-NEtFOSAA

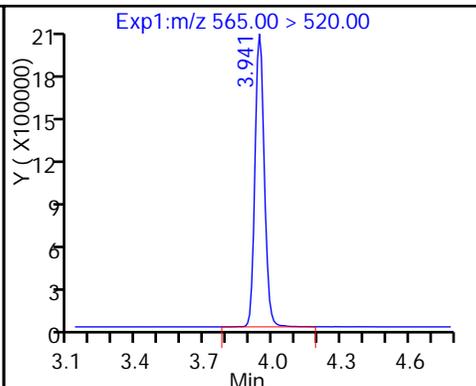
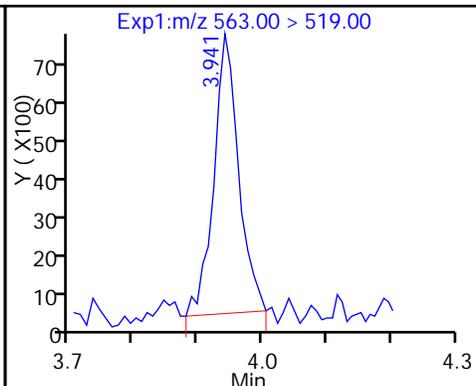
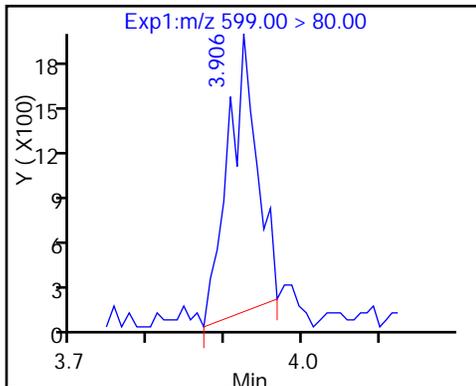
49 N-ethyl perfluorooctane sulfonamid



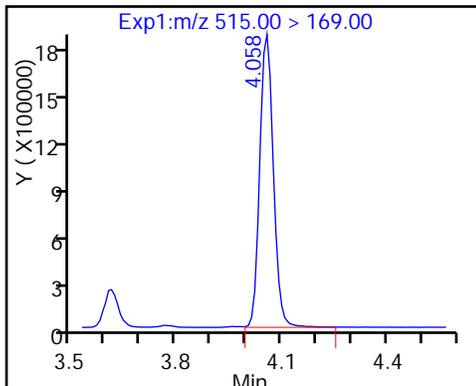
26 Perfluorodecane Sulfonic acid

28 Perfluoroundecanoic acid

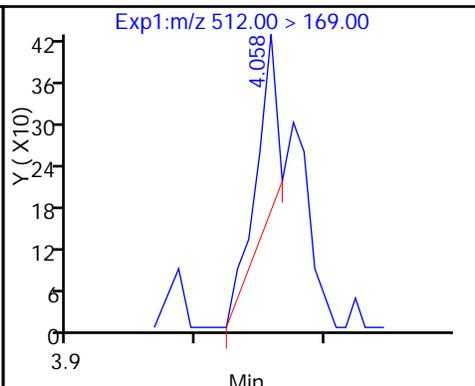
D 27 13C2 PFUnA



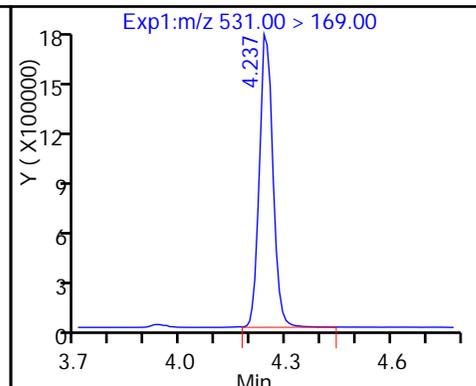
D 52 d-N-MeFOSA-M



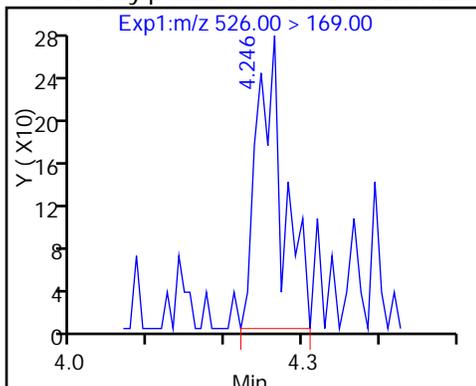
54 MeFOSA



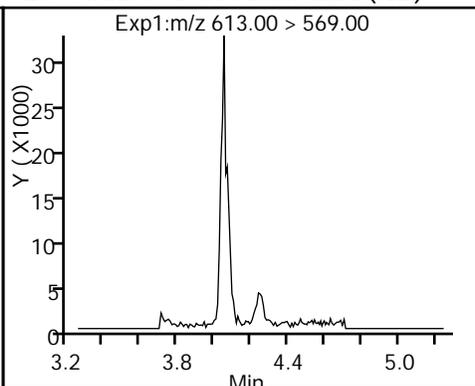
D 51 d-N-EtFOSA-M



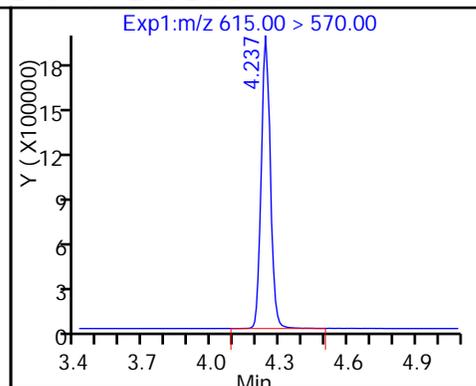
53 N-ethylperfluoro-1-octanesulfonami



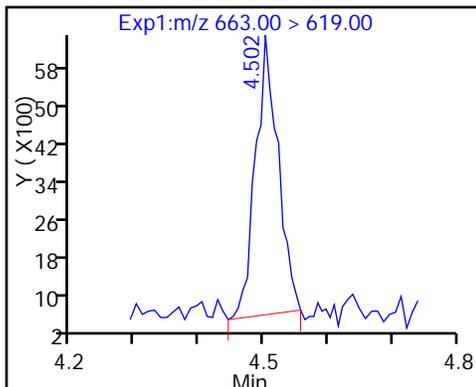
29 Perfluorododecanoic acid (ND)



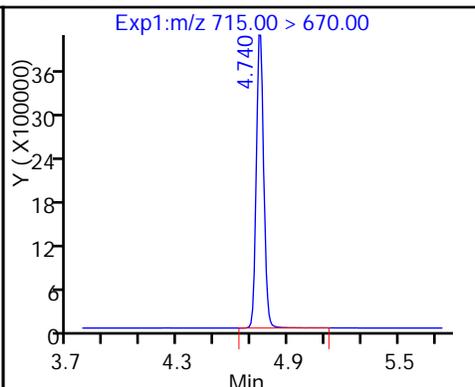
D 30 13C2 PFDaA



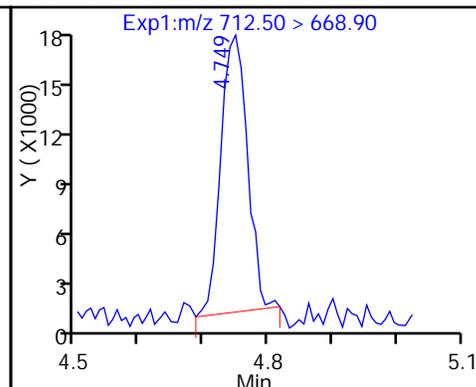
31 Perfluorotridecanoic acid



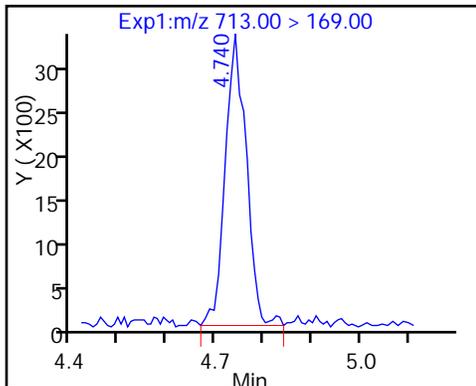
D 32 13C2-PFTeDA



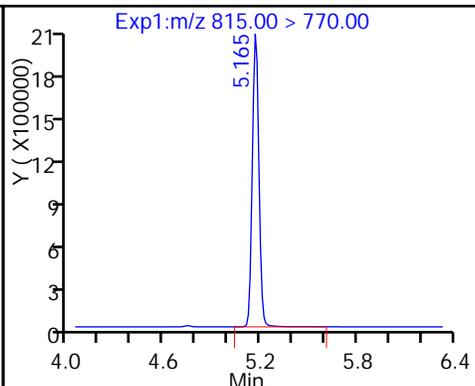
33 Perfluorotetradecanoic acid



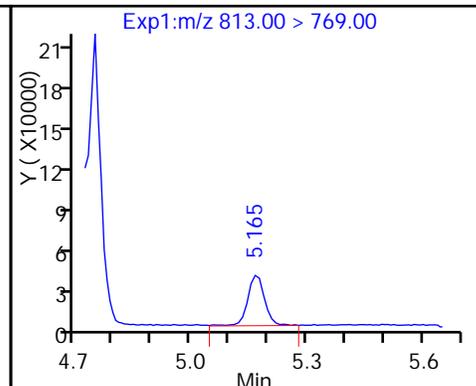
33 Perfluorotetradecanoic acid



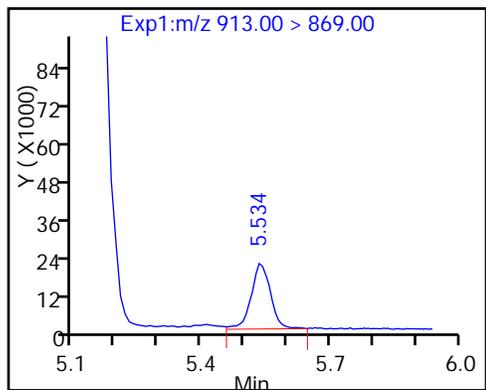
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 320-139627/2-A
 Matrix: Solid Lab File ID: 02DEC2016C_002.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: SHAKE Date Extracted: 11/28/2016 11:03
 Sample wt/vol: 5.00(g) Date Analyzed: 12/02/2016 15:07
 Con. Extract Vol.: 1.00(mL) Dilution Factor: 1
 Injection Volume: 2(uL) GC Column: Acquity ID: 2.1(mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140429 Units: ug/Kg

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	4.25		0.50	0.10
1763-23-1	Perfluorooctane Sulfonate (PFOS)	3.66		0.50	0.13
375-73-5	Perfluorobutanesulfonic acid (PFBS)	4.03		0.40	0.10

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	119		25-150
STL00991	13C4 PFOS	112		25-150
STL00994	18O2 PFHxS	120		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_002.d
 Lims ID: LCS 320-139627/2-A
 Client ID:
 Sample Type: LCS
 Inject. Date: 02-Dec-2016 15:07:03 ALS Bottle#: 2 Worklist Smp#: 6
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: lcs 320-139627/2-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 04-Dec-2016 18:17:14 Calib Date: 02-Dec-2016 12:29:29
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161202-37450.b\02DEC2016A_020.d
 Column 1 : Det: EXP1
 Process Host: XAWRK025

First Level Reviewer: chandrasenas Date: 04-Dec-2016 18:17:14

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
1 Perfluorobutyric acid										
212.90 > 169.00	1.582	1.574	0.008	1.000	7714134	21.3		106	70812	
D 2 13C4 PFBA										
217.00 > 172.00	1.574	1.574	0.0		20481402	60.2		120	1265290	
3 Perfluoropentanoic acid										
262.90 > 219.00	1.868	1.868	0.0	1.000	6535777	19.3		96.7	62609	
D 4 13C5-PFPeA										
267.90 > 223.00	1.868	1.868	0.0		16242034	60.1		120	1094905	
5 Perfluorobutanesulfonic acid										
298.90 > 80.00	1.906	1.907	-0.001	1.000	11783909	20.1		114		
298.90 > 99.00	1.906	1.907	-0.001	1.000	5085170		2.32(0.00-0.00)			
7 Perfluorohexanoic acid										
313.00 > 269.00	2.179	2.173	0.006	1.000	5304421	20.8		104	102430	
D 6 13C2 PFHxA										
315.00 > 270.00	2.179	2.173	0.006		13349831	54.2		108	1044992	
12 Perfluoroheptanoic acid										
363.00 > 319.00	2.520	2.519	0.001	1.000	5325594	20.3		101	84952	
D 11 13C4-PFHpA										
367.00 > 322.00	2.520	2.519	0.001		12690645	60.3		121	811314	
9 Perfluorohexanesulfonic acid										
399.00 > 80.00	2.536	2.534	0.002	1.000	6956903	18.1		99.4		
D 10 18O2 PFHxS										
403.00 > 84.00	2.536	2.534	0.002		17161737	56.7		120	2215913	
D 14 13C4 PFOA										
417.00 > 372.00	2.895	2.886	0.009		13102595	59.7		119	818355	
15 Perfluorooctanoic acid										
413.00 > 369.00	2.895	2.894	0.001	1.000	5685652	21.3		106	81058	
413.00 > 169.00	2.895	2.894	0.001	1.000	3320781		1.71(0.90-1.10)		187412	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.903	2.902	0.001	1.000	6325837	19.8		104
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.269	3.267	0.002	1.000	5437362	18.3		98.5
	499.00	> 99.00	3.276	3.267	0.009	1.002	1198769		4.54(0.90-1.10)	1603544 67382
D 17 13C4 PFOS	503.00	> 80.00	3.269	3.267	0.002		13071743	53.4		112
20 Perfluorononanoic acid	463.00	> 419.00	3.276	3.275	0.001	1.000	4107662	19.9		99.3
D 19 13C5 PFNA	468.00	> 423.00	3.276	3.275	0.001		10430504	58.6		117
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.591	3.589	0.002	1.000	6225860	20.9		104
D 21 13C8 FOSA	506.00	> 78.00	3.591	3.589	0.002		15929131	38.5		76.9
D 23 13C2 PFDA	515.00	> 470.00	3.633	3.631	0.002		9554899	57.5		115
24 Perfluorodecanoic acid	513.00	> 469.00	3.633	3.631	0.002	1.000	3686742	19.9		99.5
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.943	3.941	0.002	1.000	2787512	16.3		84.3
28 Perfluoroundecanoic acid	563.00	> 519.00	3.969	3.959	0.010	1.000	2452827	17.5		87.3
D 27 13C2 PFUnA	565.00	> 520.00	3.969	3.967	0.002		6630265	52.7		105
29 Perfluorododecanoic acid	613.00	> 569.00	4.260	4.259	0.001	1.000	1878687	19.1		95.7
D 30 13C2 PFDaA	615.00	> 570.00	4.260	4.259	0.001		5187445	44.4		88.7
31 Perfluorotridecanoic acid	663.00	> 619.00	4.530	4.520	0.010	1.000	1368885	12.7		63.4
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.763	4.772	-0.009	1.000	2525010	12.6		63.0
	713.00	> 169.00	4.763	4.772	-0.009	1.000	405389		6.23(0.00-0.00)	1335 85374
D 32 13C2-PFTeDA	715.00	> 670.00	4.763	4.772	-0.009		7044061	29.1		58.2
D 34 13C2-PFHxDA	815.00	> 770.00	5.186	5.184	0.002		2287334	17.5		35.1
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.186	5.194	-0.008	1.000	788728	6.89		34.5
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.569	5.562	0.007	1.000	812724	9.96		49.8

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b\02DEC2016C_002.d

Injection Date: 02-Dec-2016 15:07:03

Instrument ID: A8_N

Lims ID: LCS 320-139627/2-A

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 2

Worklist Smp#: 6

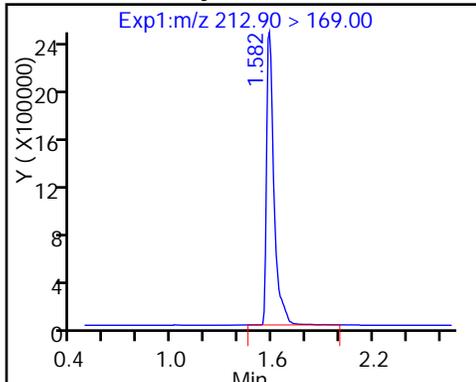
Injection Vol: 2.0 ul

Dil. Factor: 1.0000

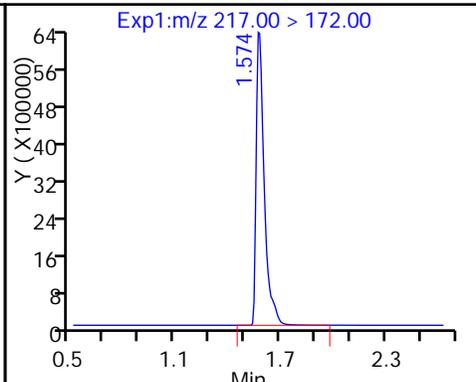
Method: A8_N

Limit Group: LC PFC_DOD ICAL

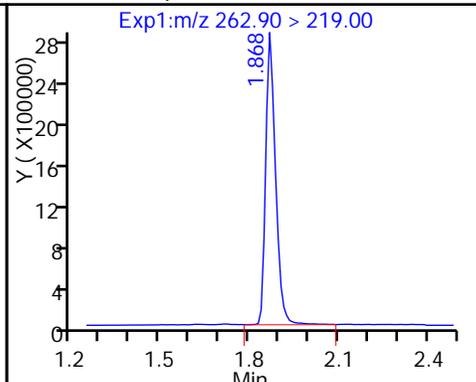
1 Perfluorobutyric acid



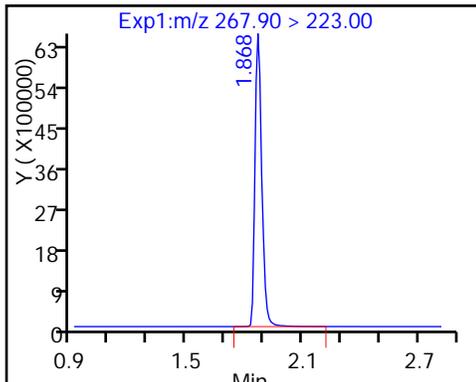
D 2 13C4 PFBA



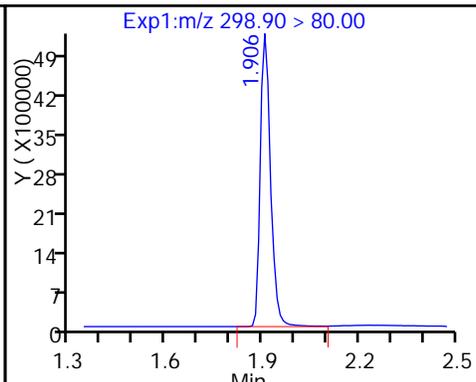
3 Perfluoropentanoic acid



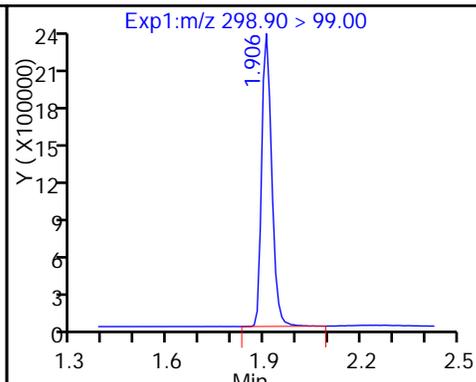
D 4 13C5-PFPeA



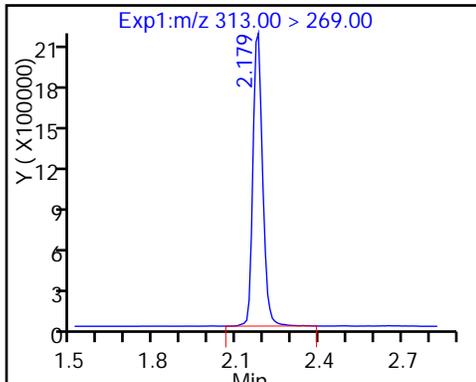
5 Perfluorobutanesulfonic acid



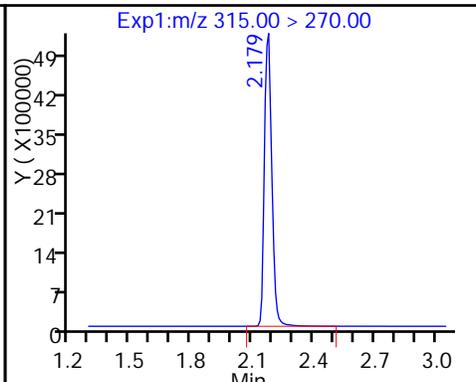
5 Perfluorobutanesulfonic acid



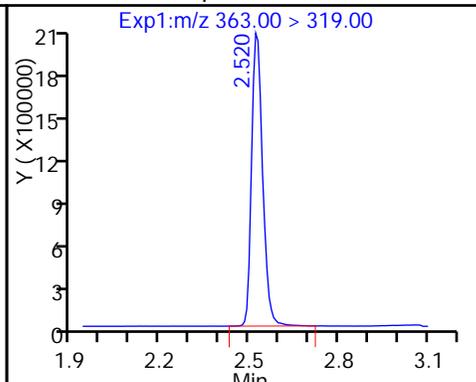
7 Perfluorohexanoic acid



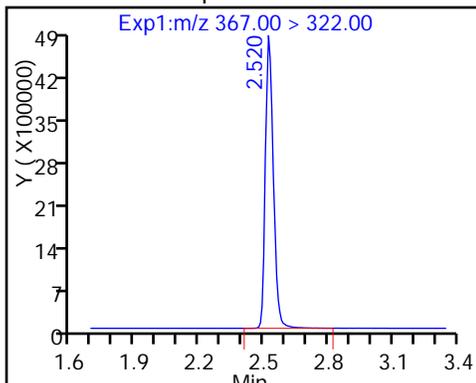
D 6 13C2 PFHxA



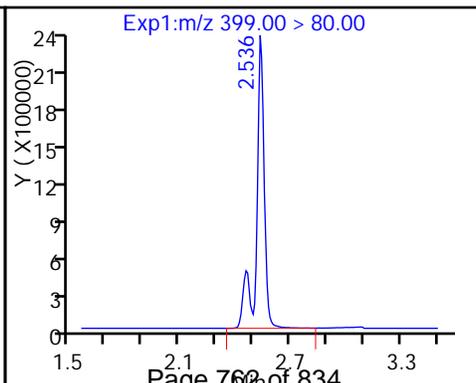
12 Perfluoroheptanoic acid



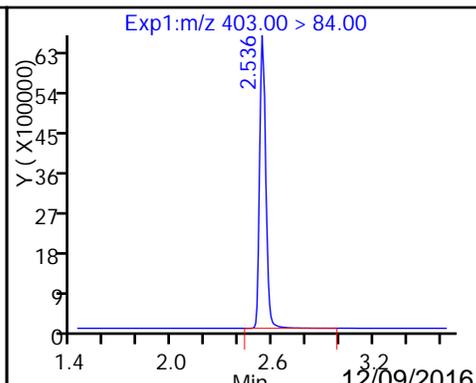
D 11 13C4-PFHpA



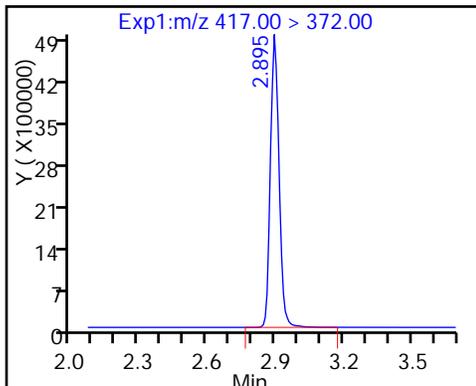
9 Perfluorohexanesulfonic acid



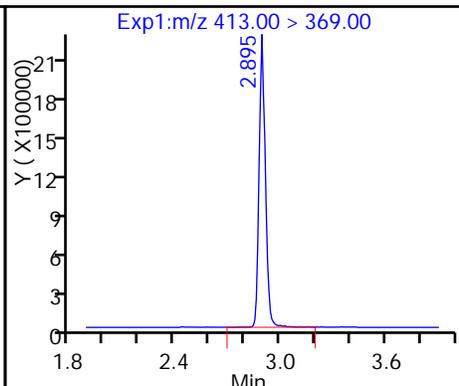
D 10 18O2 PFHxS



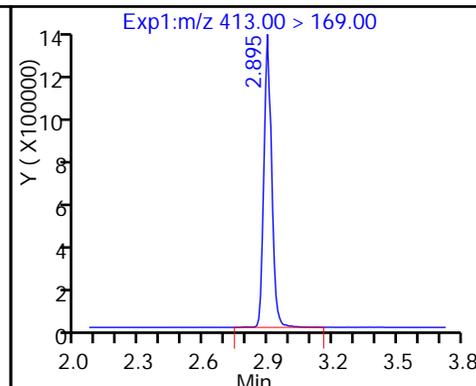
D 14 13C4 PFOA



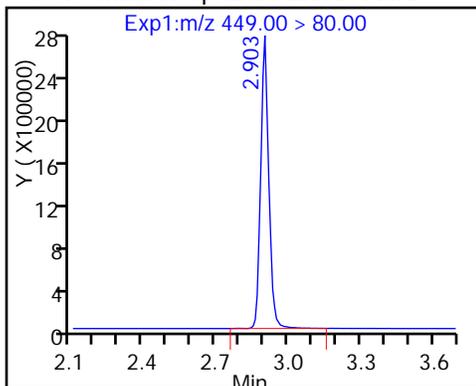
15 Perfluorooctanoic acid



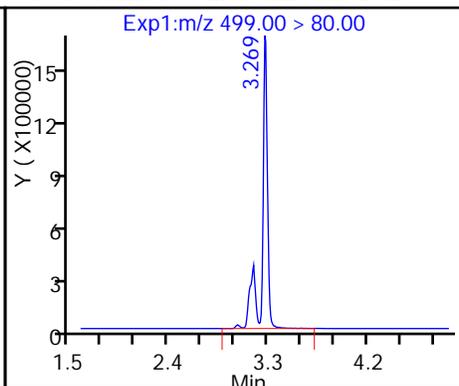
15 Perfluorooctanoic acid



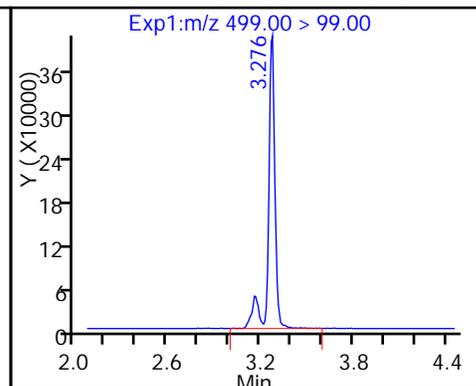
13 Perfluoroheptanesulfonic Acid



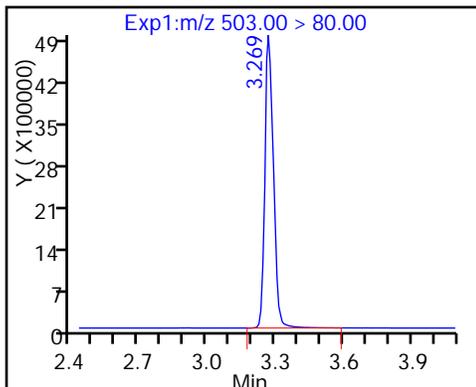
18 Perfluorooctane sulfonic acid



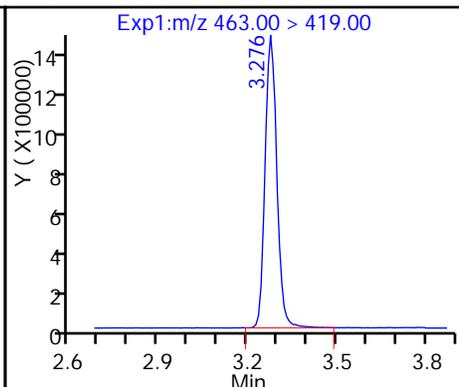
18 Perfluorooctane sulfonic acid



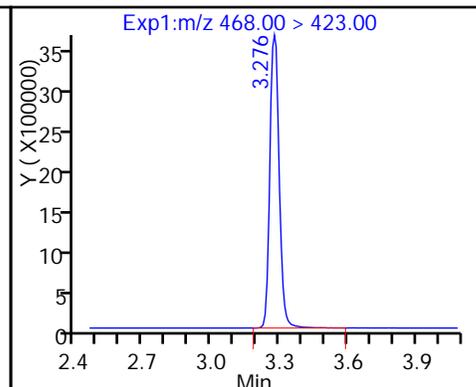
D 17 13C4 PFOS



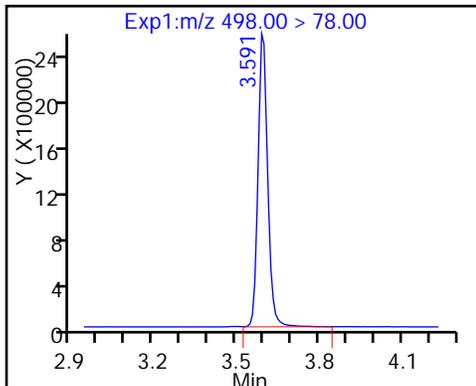
20 Perfluorononanoic acid



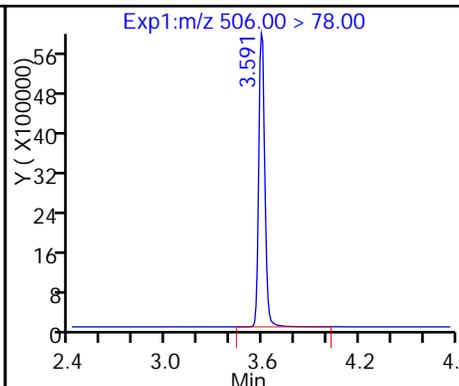
D 19 13C5 PFNA



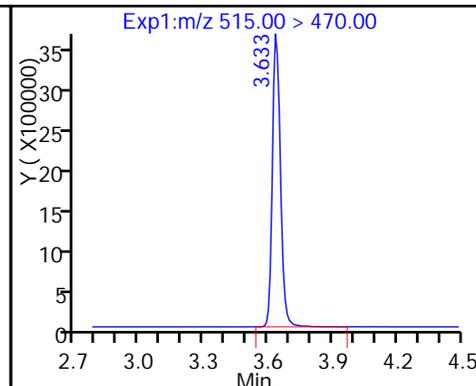
22 Perfluorooctane Sulfonamide

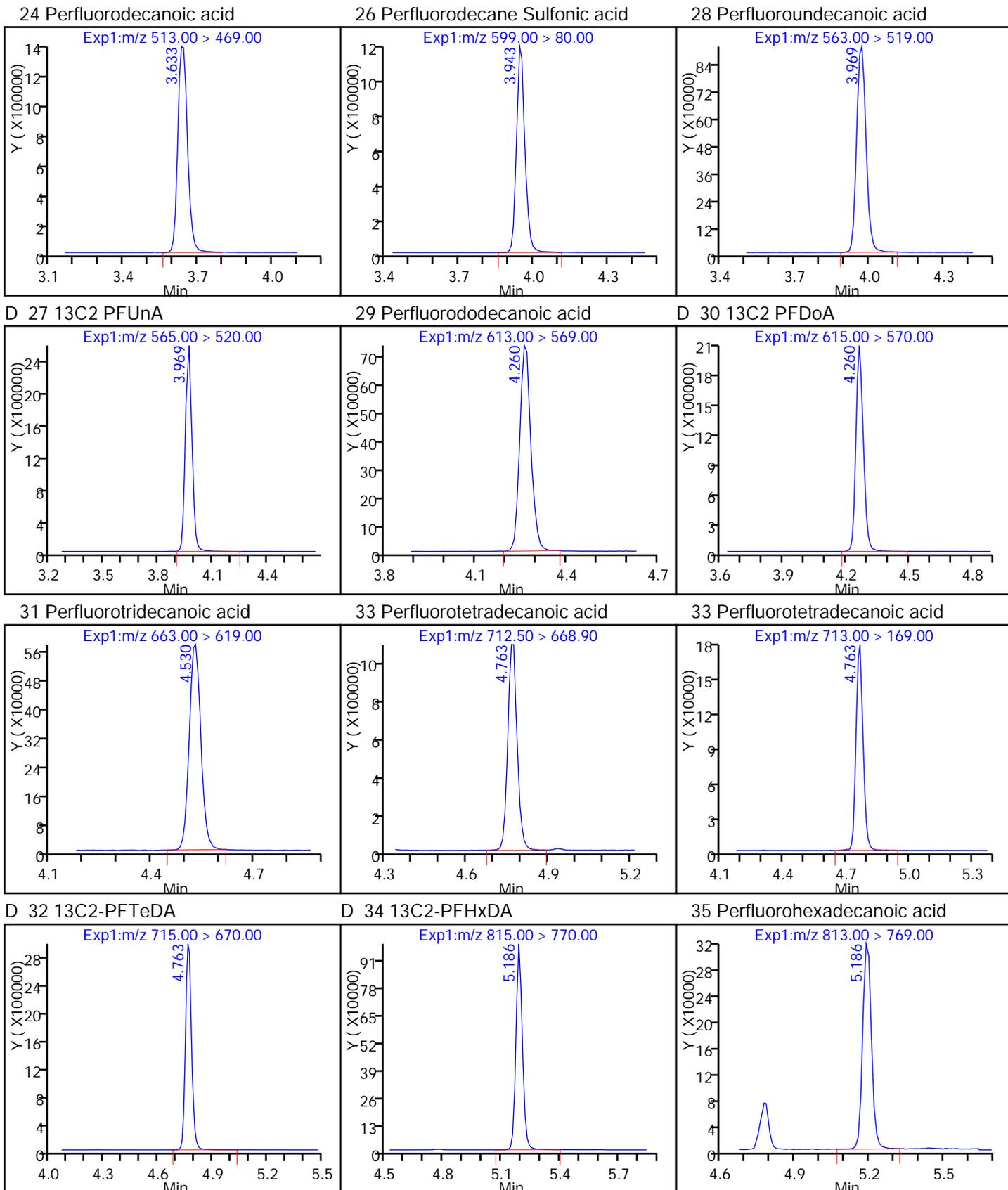


D 21 13C8 FOSA

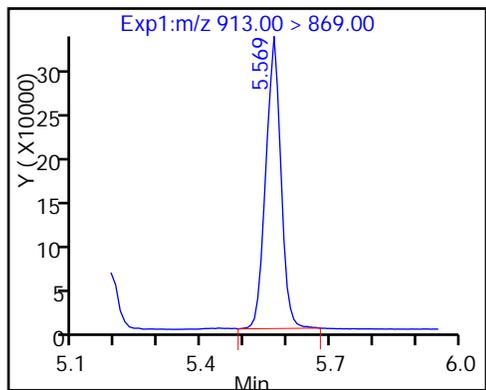


D 23 13C2 PFDA





36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCS 320-139749/2-A
 Matrix: Water Lab File ID: 03DEC2016B_002.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 250 (mL) Date Analyzed: 12/03/2016 16:11
 Con. Extract Vol.: 0.5 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.0296		0.0025	0.00075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.0261		0.0040	0.0013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0289		0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	108		25-150
STL00991	13C4 PFOS	104		25-150
STL00994	18O2 PFHxS	105		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_002.d
 Lims ID: LCS 320-139749/2-A
 Client ID:
 Sample Type: LCS
 Inject. Date: 03-Dec-2016 16:11:09 ALS Bottle#: 33 Worklist Smp#: 23
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: lcs 320-139749/2-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:30:02 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:27:02

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.558	1.574	-0.016	17560233	52.3		105	1375246	
1 Perfluorobutyric acid	212.90 > 169.00	1.558	1.577	-0.019	1.000	4857268	15.8	79.1	25509	
3 Perfluoropentanoic acid	262.90 > 219.00	1.849	1.861	-0.012	1.000	4268708	14.4	72.1	31016	
D 4 13C5-PFPeA	267.90 > 223.00	1.849	1.861	-0.012		14590160	55.2	110	741230	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.887	1.900	-0.013	1.000	7584330	14.5	81.8		
	298.90 > 99.00	1.887	1.900	-0.013	1.000	3290381	2.31(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.146	2.164	-0.018	1.000	3467656	14.9	74.6	82066	
D 6 13C2 PFHxA	315.00 > 270.00	2.146	2.164	-0.018		12196862	51.4	103	1065490	
D 11 13C4-PFHpA	367.00 > 322.00	2.485	2.511	-0.026		11238097	54.2	108	755063	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.485	2.512	-0.027	1.000	3523219	15.3	76.3	41059	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.507	2.531	-0.024	1.000	4525408	12.5	68.9		
D 10 18O2 PFHxS	403.00 > 84.00	2.507	2.531	-0.024		15545077	49.8	105	872611	
D 14 13C4 PFOA	417.00 > 372.00	2.847	2.880	-0.033		11822855	53.9	108	938614	
15 Perfluorooctanoic acid	413.00 > 369.00	2.847	2.887	-0.040	1.000	3754301	14.8	74.1	66200	
	413.00 > 169.00	2.855	2.887	-0.032	1.003	2189699	1.71(0.90-1.10)		79418	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.863	2.888	-0.025	1.000	4174557	13.8	72.7	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.227	3.258	-0.031	1.000	3625231	13.0	70.3	194917
	499.00	> 99.00	3.227	3.258	-0.031	1.000	816437		4.44(0.90-1.10)	61899
D 17 13C4 PFOS	503.00	> 80.00	3.227	3.259	-0.032		12244746	49.8	104	475079
D 19 13C5 PFNA	468.00	> 423.00	3.227	3.263	-0.036		9311847	56.0	112	652523
20 Perfluorononanoic acid	463.00	> 419.00	3.227	3.263	-0.036	1.000	2697531	14.5	72.7	55697
D 21 13C8 FOSA	506.00	> 78.00	3.549	3.571	-0.022		9046557	22.5	45.0	398080
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.558	3.574	-0.016	1.000	2588905	15.3	76.6	139330
24 Perfluorodecanoic acid	513.00	> 469.00	3.592	3.623	-0.031	1.000	2441183	14.7	73.4	68408
D 23 13C2 PFDA	515.00	> 470.00	3.592	3.626	-0.034		8656247	54.8	110	217190
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.900	3.936	-0.036	1.000	2203517	13.4	69.7	
28 Perfluoroundecanoic acid	563.00	> 519.00	3.917	3.955	-0.038	1.000	1723441	12.5	62.5	44705
D 27 13C2 PFUnA	565.00	> 520.00	3.917	3.958	-0.041		6471442	54.5	109	456235
29 Perfluorododecanoic acid	613.00	> 569.00	4.217	4.250	-0.033	1.000	1628778	14.2	70.8	30811
D 30 13C2 PFDoA	615.00	> 570.00	4.210	4.251	-0.041		6057552	54.0	108	187381
31 Perfluorotridecanoic acid	663.00	> 619.00	4.478	4.518	-0.040	1.000	1741016	15.1	75.6	30701
D 32 13C2-PFTeDA	715.00	> 670.00	4.715	4.759	-0.044		15011428	64.9	130	769401
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.715	4.761	-0.046	1.000	3709548	16.5	82.6	3985
	713.00	> 169.00	4.715	4.761	-0.046	1.000	633060		5.86(0.00-0.00)	121080
D 34 13C2-PFHxDA	815.00	> 770.00	5.137	5.186	-0.049		6159598	47.5	95.0	182215
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.137	5.186	-0.049	1.000	1622525	12.7	63.7	4013
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.505	5.559	-0.054	1.000	1681305	14.0	69.9	2156

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_002.d

Injection Date: 03-Dec-2016 16:11:09

Instrument ID: A8_N

Lims ID: LCS 320-139749/2-A

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 33

Worklist Smp#: 23

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

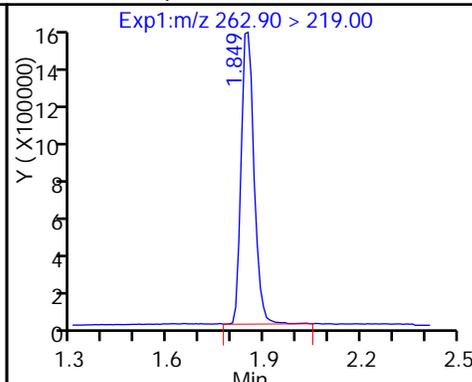
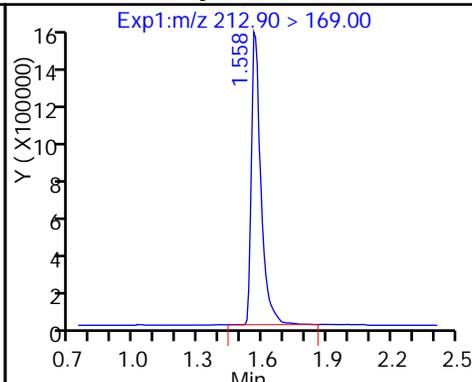
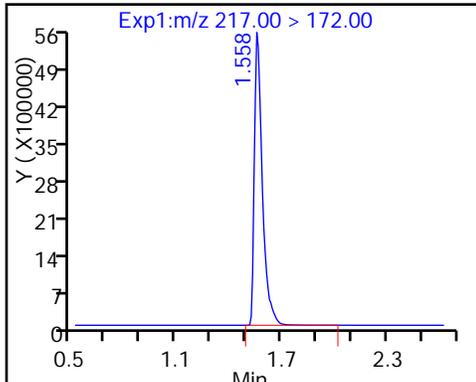
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

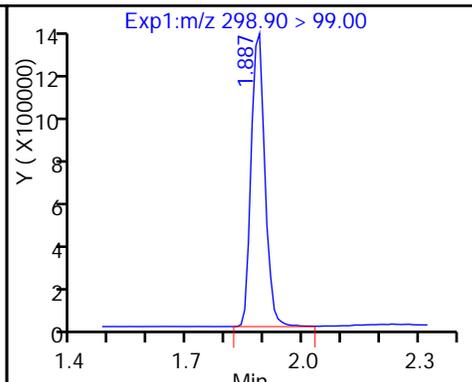
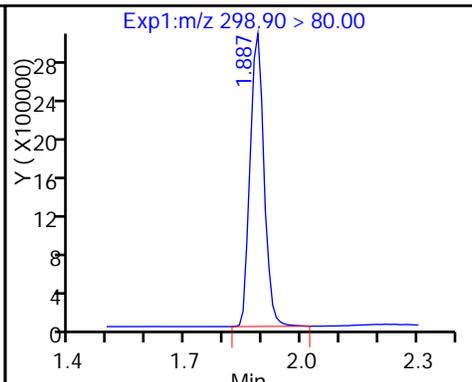
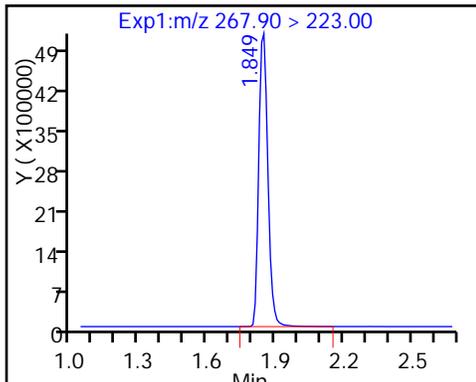
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

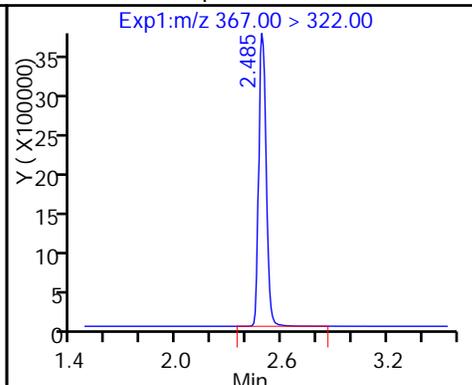
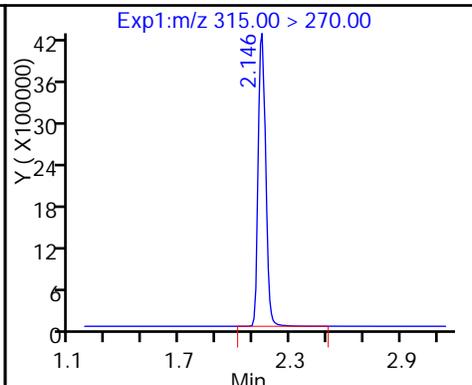
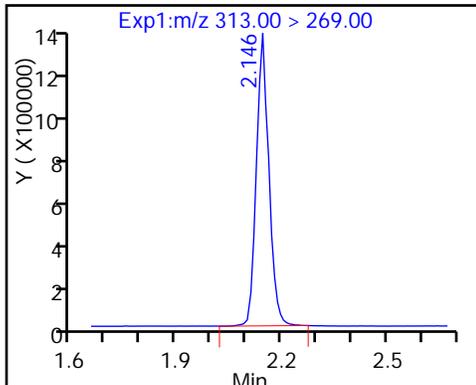
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

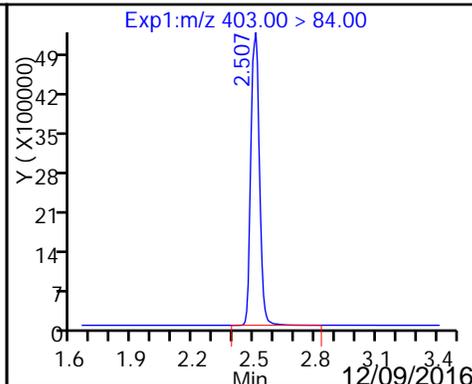
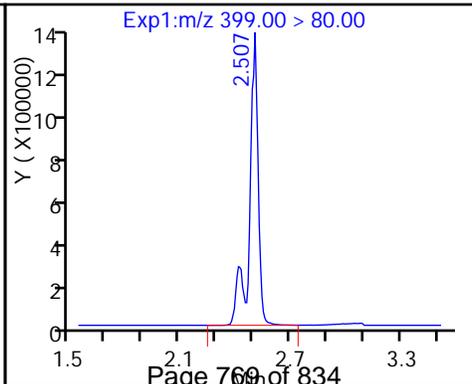
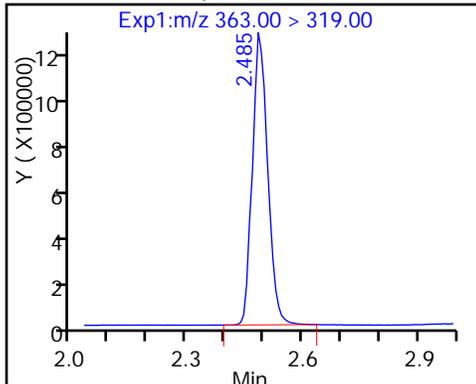
D 11 13C4-PFHpA



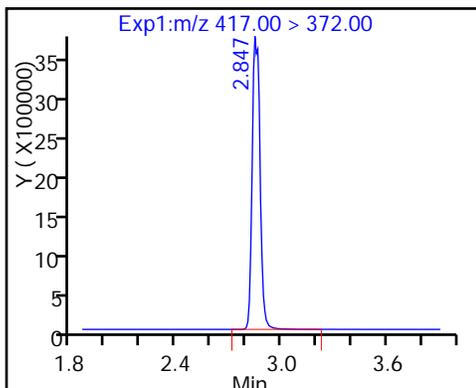
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

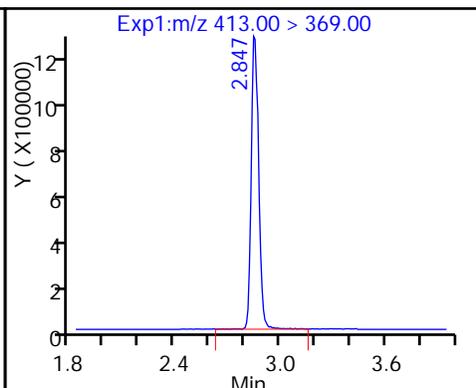
D 10 18O2 PFHxS



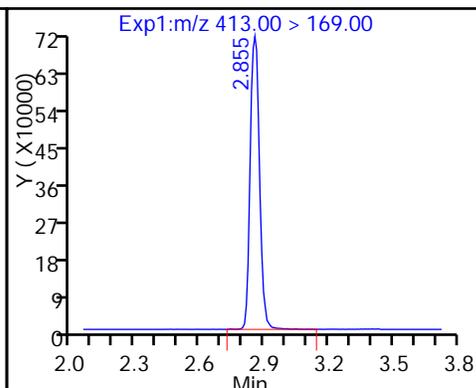
D 14 13C4 PFOA



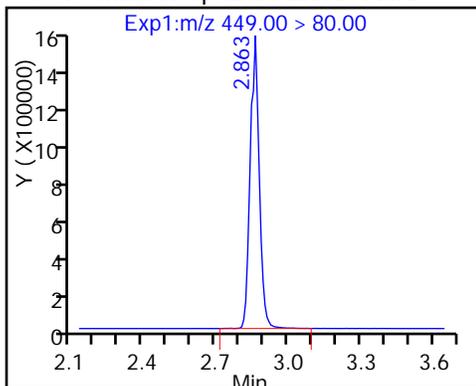
15 Perfluorooctanoic acid



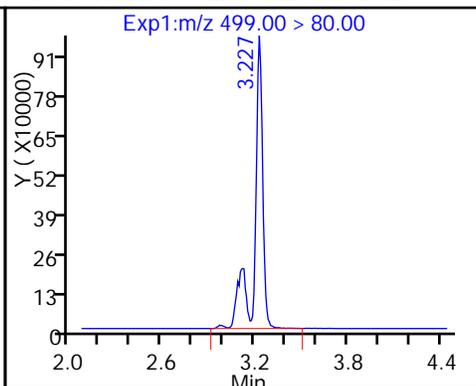
15 Perfluorooctanoic acid



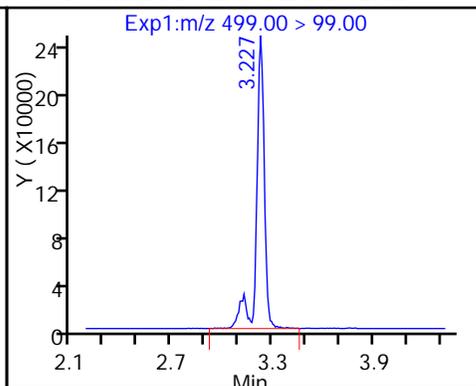
13 Perfluoroheptanesulfonic Acid



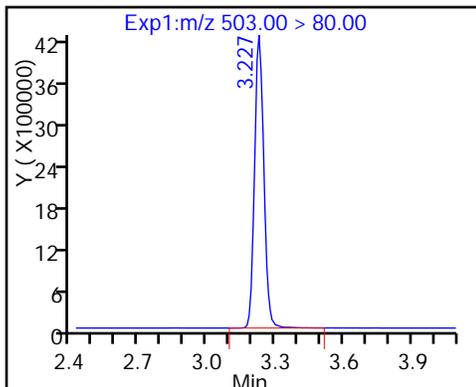
18 Perfluorooctane sulfonic acid



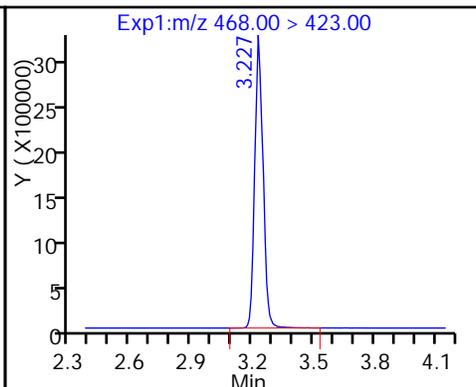
18 Perfluorooctane sulfonic acid



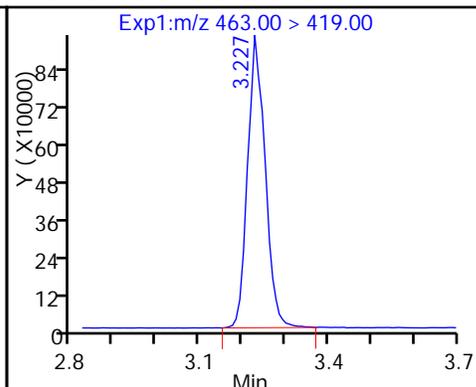
D 17 13C4 PFOS



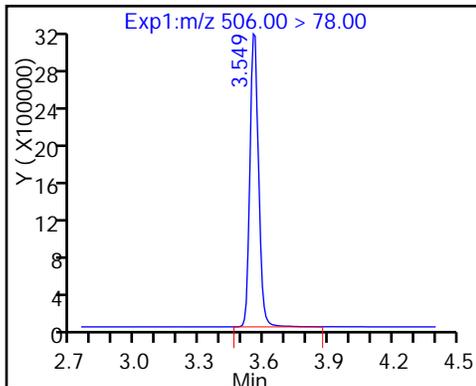
D 19 13C5 PFNA



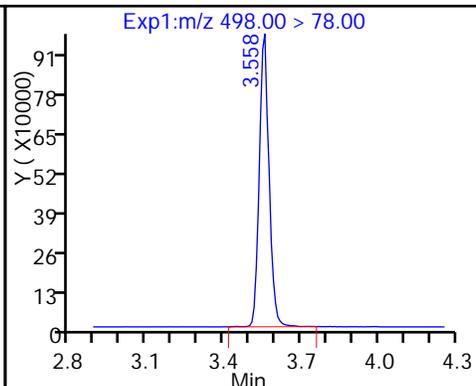
20 Perfluorononanoic acid



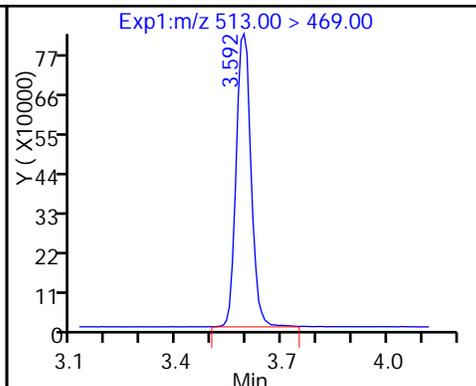
D 21 13C8 FOSA



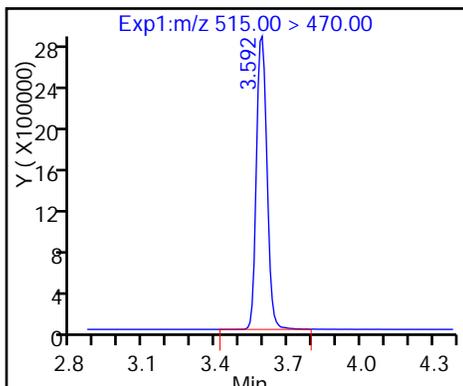
22 Perfluorooctane Sulfonamide



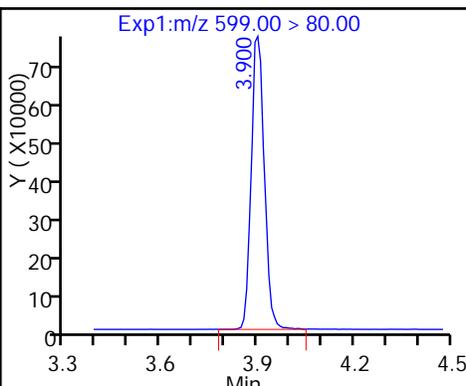
24 Perfluorodecanoic acid



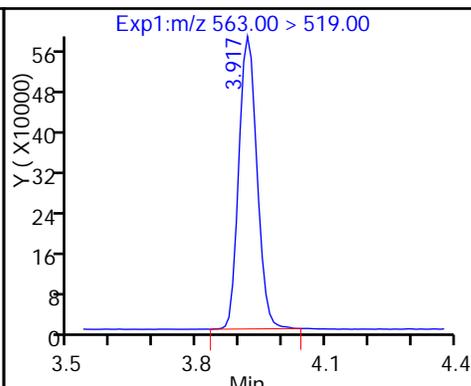
D 23 13C2 PFDA



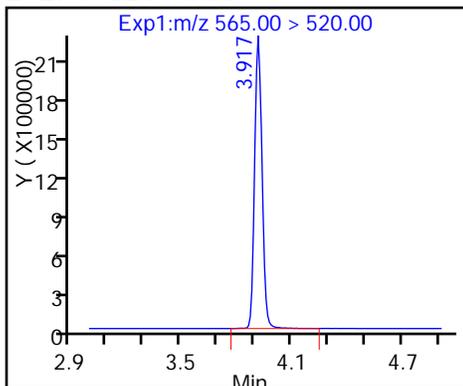
26 Perfluorodecane Sulfonic acid



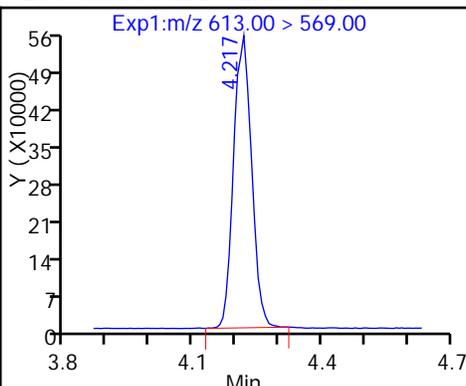
28 Perfluoroundecanoic acid



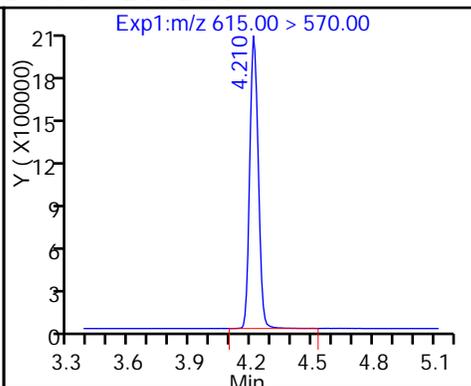
D 27 13C2 PFUa



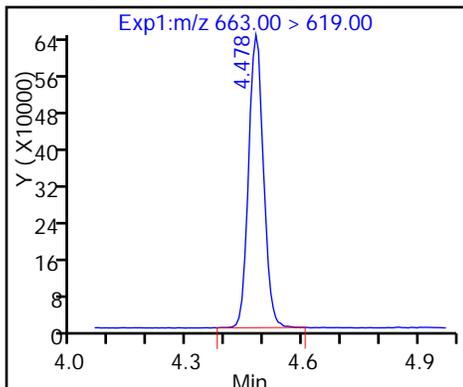
29 Perfluorododecanoic acid



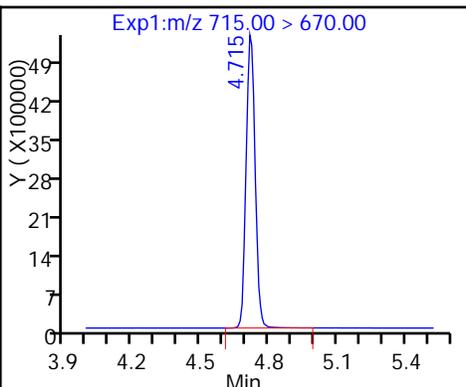
D 30 13C2 PFDa



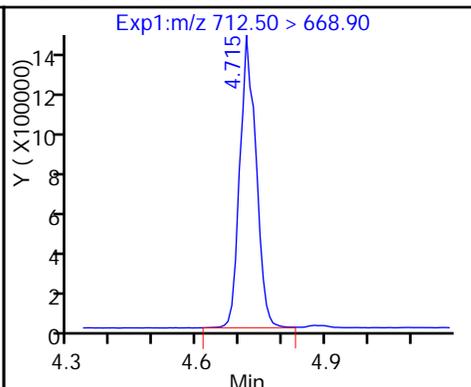
31 Perfluorotridecanoic acid



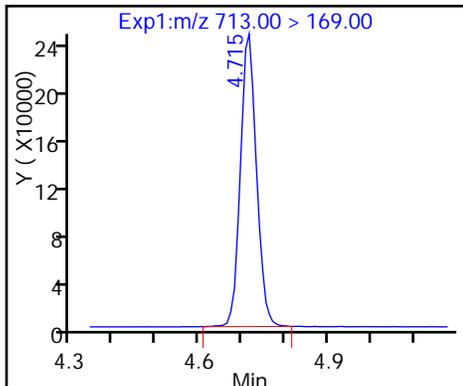
D 32 13C2-PFTeDA



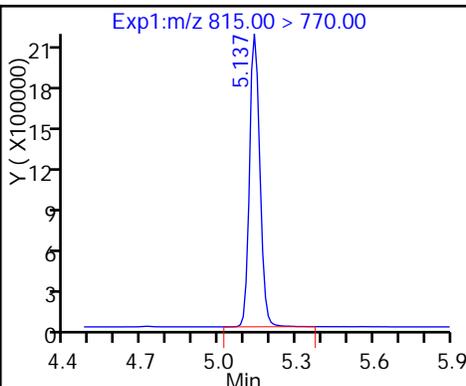
33 Perfluorotetradecanoic acid



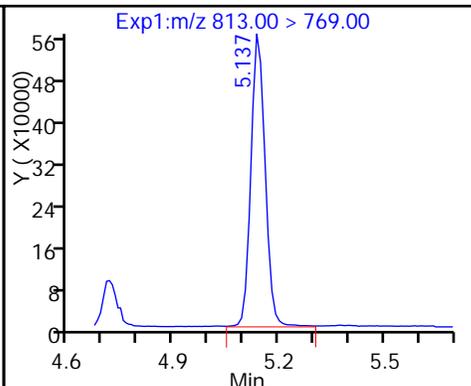
33 Perfluorotetradecanoic acid



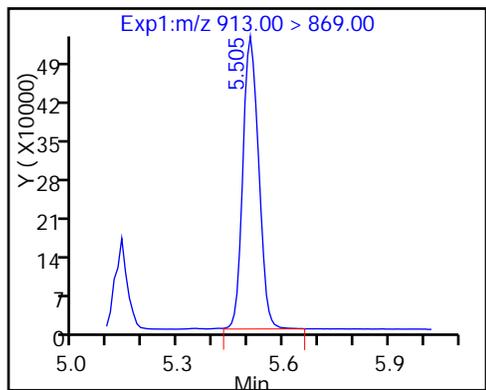
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



FORM I
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1
 SDG No.: _____
 Client Sample ID: _____ Lab Sample ID: LCSD 320-139749/3-A
 Matrix: Water Lab File ID: 03DEC2016B_003.d
 Analysis Method: 537 (Modified) Date Collected: _____
 Extraction Method: 3535 Date Extracted: 11/28/2016 16:55
 Sample wt/vol: 250 (mL) Date Analyzed: 12/03/2016 16:18
 Con. Extract Vol.: 0.5 (mL) Dilution Factor: 1
 Injection Volume: 2 (uL) GC Column: Acquity ID: 2.1 (mm)
 % Moisture: _____ GPC Cleanup: (Y/N) N
 Analysis Batch No.: 140564 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
335-67-1	Perfluorooctanoic acid (PFOA)	0.0314		0.0025	0.00075
1763-23-1	Perfluorooctane Sulfonate (PFOS)	0.0289		0.0040	0.0013
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0325		0.0025	0.00092

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00990	13C4 PFOA	116		25-150
STL00991	13C4 PFOS	104		25-150
STL00994	18O2 PFHxS	104		25-150

TestAmerica Sacramento
Target Compound Quantitation Report

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_003.d
 Lims ID: LCSD 320-139749/3-A
 Client ID:
 Sample Type: LCSD
 Inject. Date: 03-Dec-2016 16:18:41 ALS Bottle#: 34 Worklist Smp#: 24
 Injection Vol: 2.0 ul Dil. Factor: 1.0000
 Sample Info: lcsd 320-139749/3-a
 Misc. Info.: Plate: 1 Rack: 4
 Operator ID: A8-PC\A8 Instrument ID: A8_N
 Method: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\A8_N.m
 Limit Group: LC PFC_DOD ICAL
 Last Update: 05-Dec-2016 10:30:02 Calib Date: 03-Dec-2016 15:33:40
 Integrator: Picker
 Quant Method: Isotopic Dilution Quant By: Initial Calibration
 Last ICal File: \\ChromNA\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016A_018.d
 Column 1 : Det: EXP1
 Process Host: XAWRK020

First Level Reviewer: chandrasenas Date: 05-Dec-2016 10:29:05

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 2 13C4 PFBA	217.00 > 172.00	1.574	1.574	0.0	17866151	53.2		106	1821301	
1 Perfluorobutyric acid	212.90 > 169.00	1.574	1.577	-0.003	1.000	5293479	16.9	84.7	31833	
3 Perfluoropentanoic acid	262.90 > 219.00	1.849	1.861	-0.012	1.000	4737227	16.0	79.8	31580	
D 4 13C5-PFPeA	267.90 > 223.00	1.849	1.861	-0.012		14614763	55.2	110	1048466	
5 Perfluorobutanesulfonic acid	298.90 > 80.00	1.887	1.900	-0.013	1.000	8414372	16.3	92.0		
	298.90 > 99.00	1.887	1.900	-0.013	1.000	3551354	2.37(0.00-0.00)			
7 Perfluorohexanoic acid	313.00 > 269.00	2.157	2.164	-0.007	1.000	3907622	16.5	82.7	93403	
D 6 13C2 PFHxA	315.00 > 270.00	2.157	2.164	-0.007		12390841	52.2	104	587374	
D 11 13C4-PFHpA	367.00 > 322.00	2.492	2.511	-0.019		11963765	57.7	115	1060650	
12 Perfluoroheptanoic acid	363.00 > 319.00	2.492	2.512	-0.020	1.000	3976171	16.2	80.9	54713	
9 Perfluorohexanesulfonic acid	399.00 > 80.00	2.507	2.531	-0.024	1.000	5147254	14.5	79.5		
D 10 18O2 PFHxS	403.00 > 84.00	2.507	2.531	-0.024		15325146	49.1	104	1813532	
D 14 13C4 PFOA	417.00 > 372.00	2.855	2.880	-0.025		12757885	58.2	116	874253	
15 Perfluorooctanoic acid	413.00 > 369.00	2.855	2.887	-0.032	1.000	4297812	15.7	78.6	76631	
	413.00 > 169.00	2.855	2.887	-0.032	1.000	2552106	1.68(0.90-1.10)		113223	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluoroheptanesulfonic Acid	449.00	> 80.00	2.871	2.888	-0.017	1.000	4626287	15.4	81.0	
18 Perfluorooctane sulfonic acid	499.00	> 80.00	3.235	3.258	-0.023	1.000	3998252	14.5	77.9	251636
	499.00	> 99.00	3.235	3.258	-0.023	1.000	884783		4.52(0.90-1.10)	44320
D 17 13C4 PFOS	503.00	> 80.00	3.235	3.259	-0.024		12179746	49.5	104	383271
D 19 13C5 PFNA	468.00	> 423.00	3.235	3.263	-0.028		9401849	56.5	113	433505
20 Perfluorononanoic acid	463.00	> 419.00	3.235	3.263	-0.028	1.000	3047600	16.3	81.3	57192
D 21 13C8 FOSA	506.00	> 78.00	3.557	3.571	-0.014		5188852	12.9	25.8	208610
22 Perfluorooctane Sulfonamide	498.00	> 78.00	3.566	3.574	-0.008	1.000	1597785	16.5	82.4	70317
24 Perfluorodecanoic acid	513.00	> 469.00	3.599	3.623	-0.024	1.000	2761173	16.8	84.0	94373
D 23 13C2 PFDA	515.00	> 470.00	3.599	3.626	-0.027		8558073	54.2	108	483536
26 Perfluorodecane Sulfonic acid	599.00	> 80.00	3.909	3.936	-0.027	1.000	2430460	14.9	77.3	
28 Perfluoroundecanoic acid	563.00	> 519.00	3.927	3.955	-0.028	1.000	1914632	13.7	68.6	42765
D 27 13C2 PFUnA	565.00	> 520.00	3.918	3.958	-0.040		6546548	55.1	110	363893
29 Perfluorododecanoic acid	613.00	> 569.00	4.222	4.250	-0.028	1.000	1846976	15.7	78.6	38032
D 30 13C2 PFDoA	615.00	> 570.00	4.222	4.251	-0.029		6190705	55.2	110	193387
31 Perfluorotridecanoic acid	663.00	> 619.00	4.483	4.518	-0.035	1.000	1866805	15.9	79.4	22276
D 32 13C2-PFTeDA	715.00	> 670.00	4.720	4.759	-0.039		14267910	61.7	123	995101
33 Perfluorotetradecanoic acid	712.50	> 668.90	4.728	4.761	-0.033	1.000	4138056	18.0	90.2	2484
	713.00	> 169.00	4.720	4.761	-0.041	0.998	675300		6.13(0.00-0.00)	123708
D 34 13C2-PFHxDA	815.00	> 770.00	5.145	5.186	-0.041		6470420	49.9	99.8	167496
35 Perfluorohexadecanoic acid	813.00	> 769.00	5.145	5.186	-0.041	1.000	1855833	14.3	71.7	3647
36 Perfluorooctadecanoic acid	913.00	> 869.00	5.512	5.559	-0.047	1.000	1891606	15.4	76.9	3028

TestAmerica Sacramento

Data File: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b\03DEC2016B_003.d

Injection Date: 03-Dec-2016 16:18:41

Instrument ID: A8_N

Lims ID: LCSD 320-139749/3-A

Client ID:

Operator ID: A8-PC\A8

ALS Bottle#: 34

Worklist Smp#: 24

Injection Vol: 2.0 ul

Dil. Factor: 1.0000

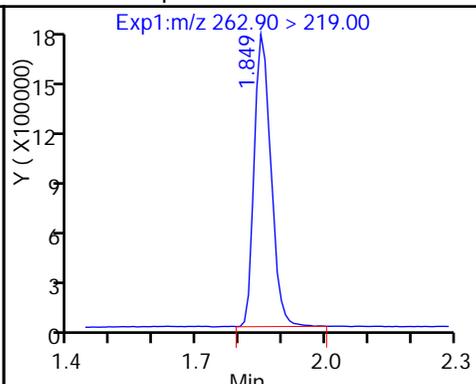
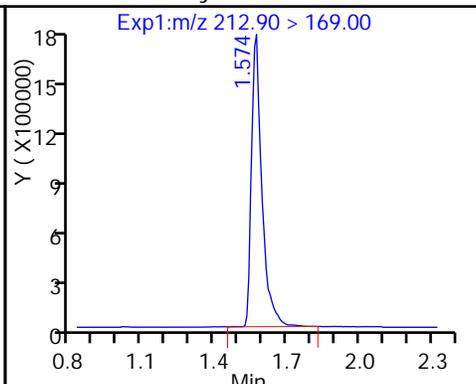
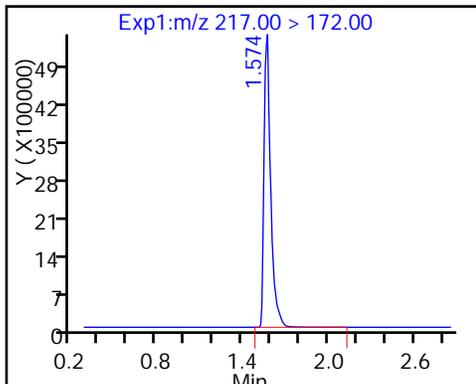
Method: A8_N

Limit Group: LC PFC_DOD ICAL

D 2 13C4 PFBA

1 Perfluorobutyric acid

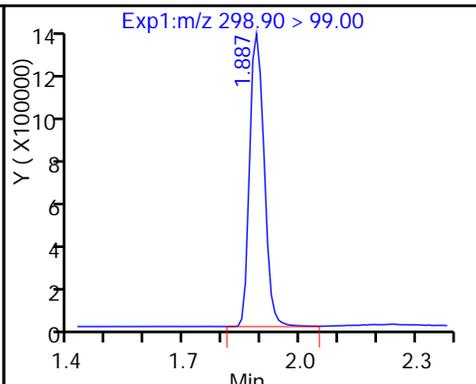
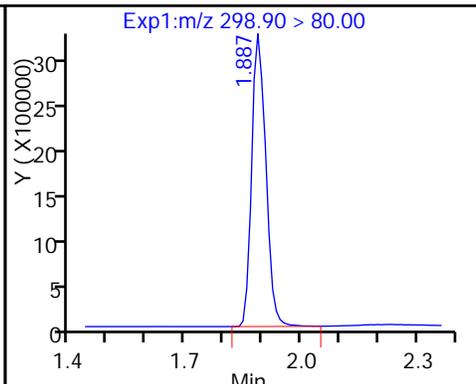
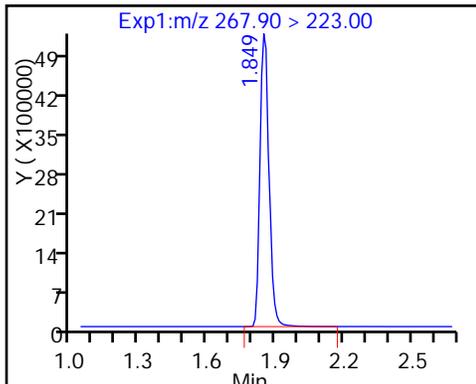
3 Perfluoropentanoic acid



D 4 13C5-PFPeA

5 Perfluorobutanesulfonic acid

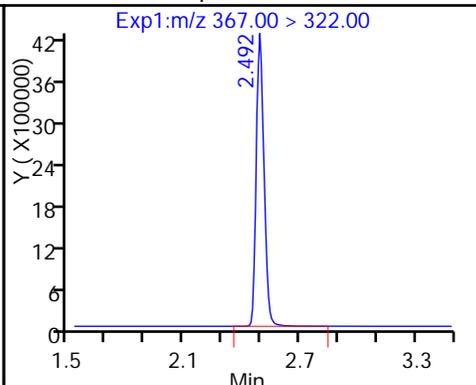
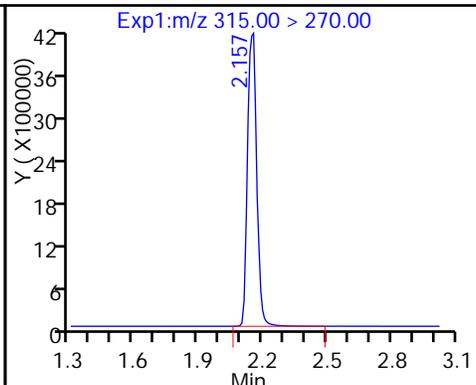
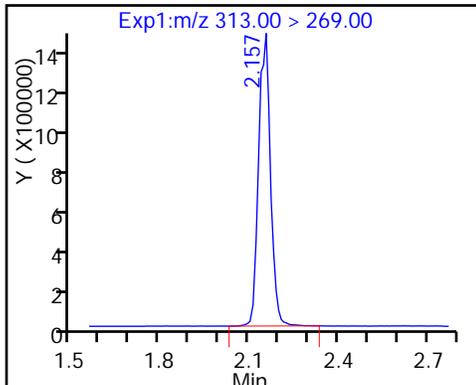
5 Perfluorobutanesulfonic acid



7 Perfluorohexanoic acid

D 6 13C2 PFHxA

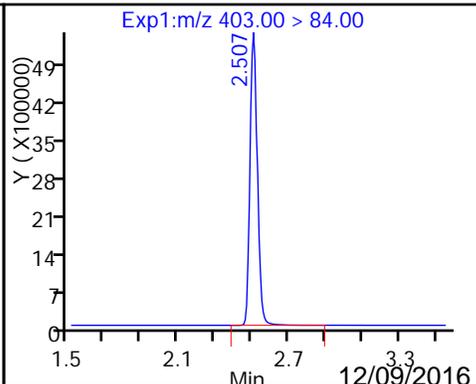
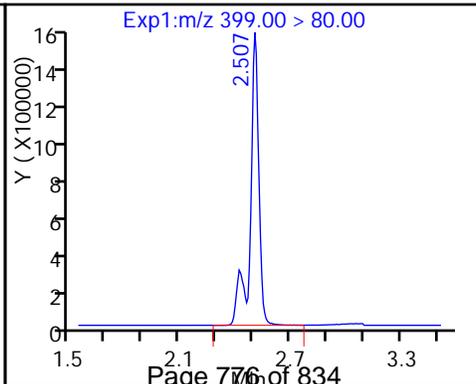
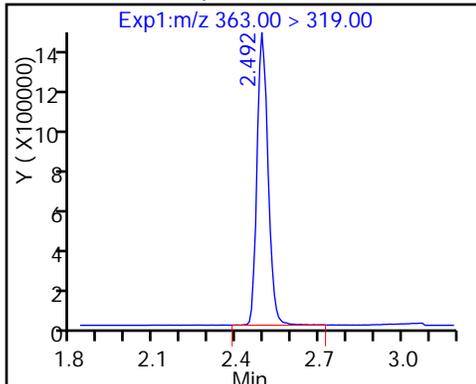
D 11 13C4-PFHpA



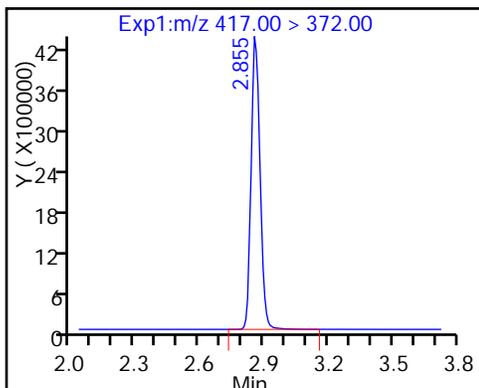
12 Perfluoroheptanoic acid

9 Perfluorohexanesulfonic acid

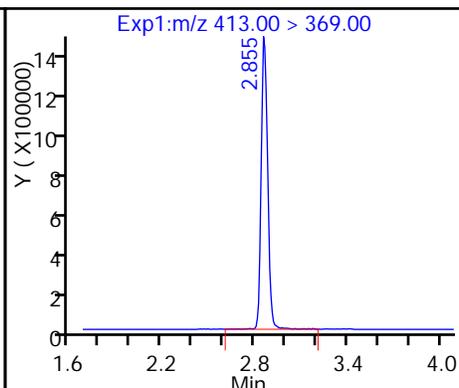
D 10 18O2 PFHxS



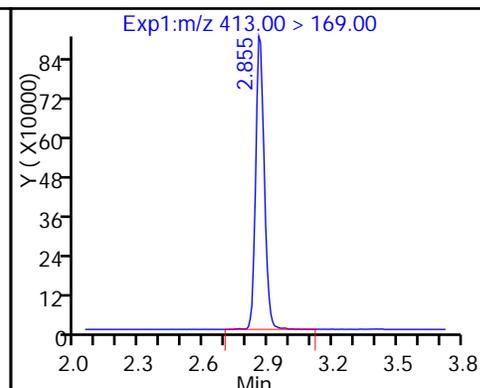
D 14 13C4 PFOA



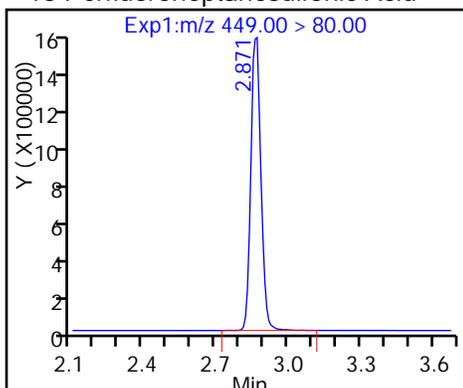
15 Perfluorooctanoic acid



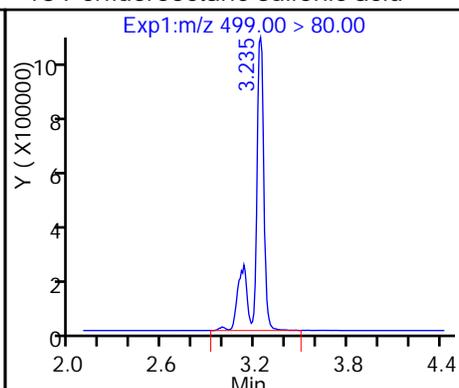
15 Perfluorooctanoic acid



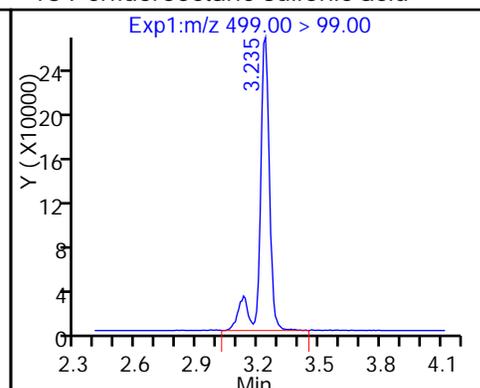
13 Perfluoroheptanesulfonic Acid



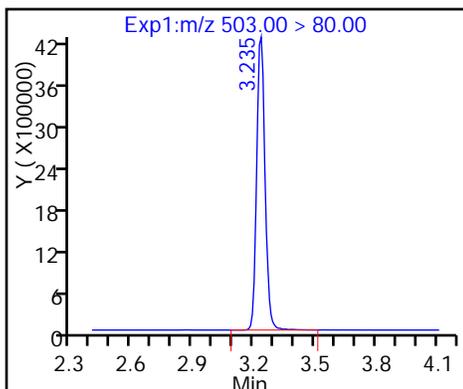
18 Perfluorooctane sulfonic acid



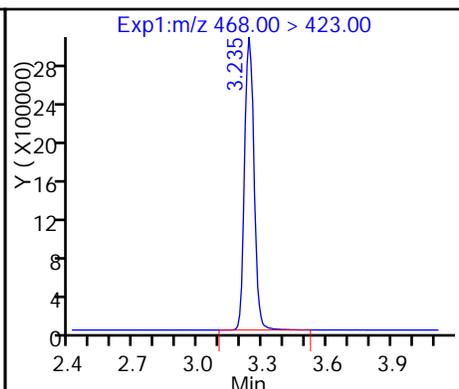
18 Perfluorooctane sulfonic acid



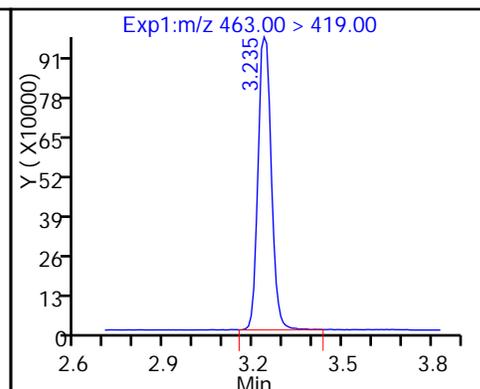
D 17 13C4 PFOS



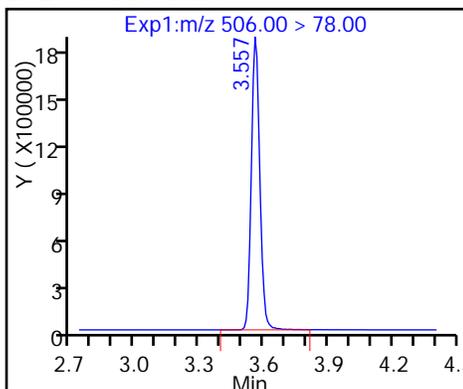
D 19 13C5 PFNA



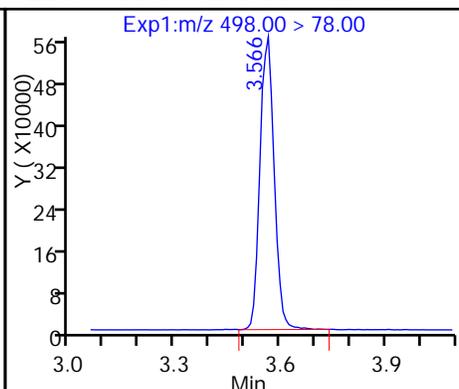
20 Perfluorononanoic acid



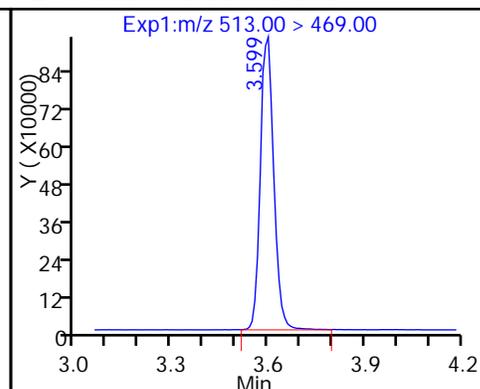
D 21 13C8 FOSA



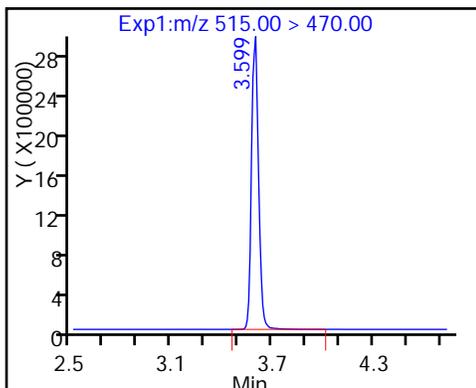
22 Perfluorooctane Sulfonamide



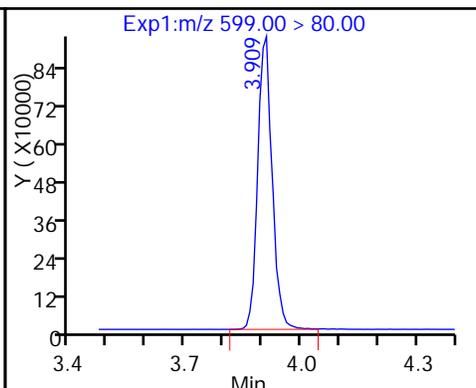
24 Perfluorodecanoic acid



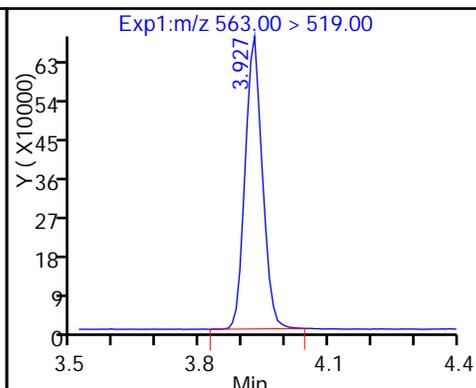
D 23 13C2 PFDA



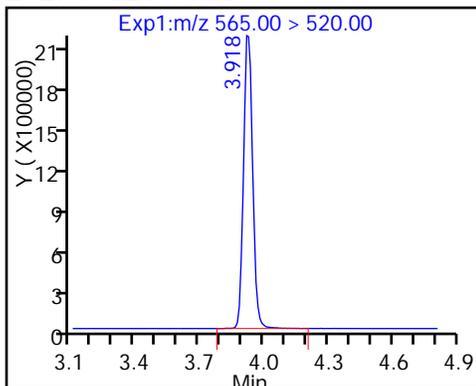
26 Perfluorodecane Sulfonic acid



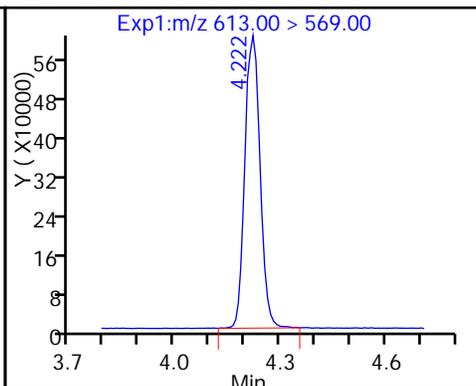
28 Perfluoroundecanoic acid



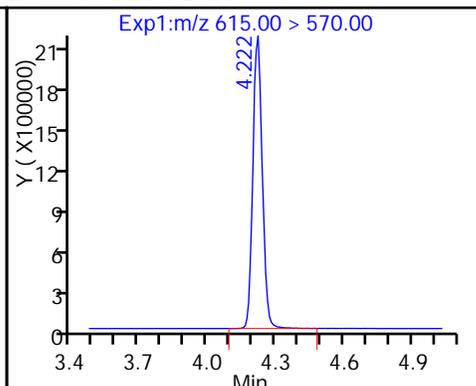
D 27 13C2 PFUa



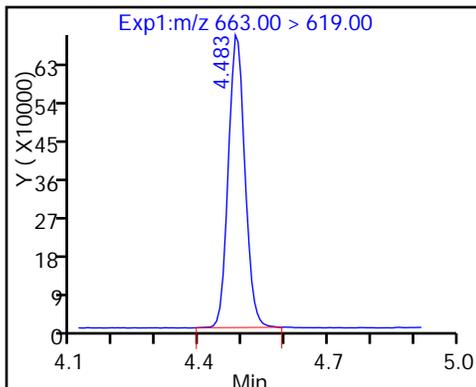
29 Perfluorododecanoic acid



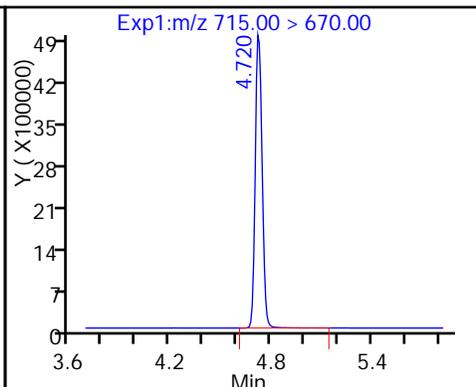
D 30 13C2 PFDa



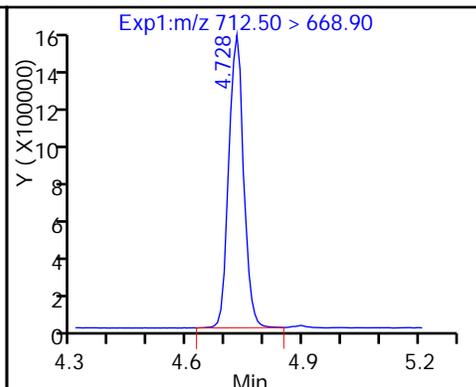
31 Perfluorotridecanoic acid



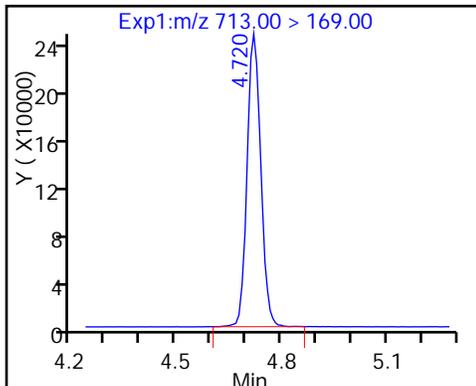
D 32 13C2-PFTeDA



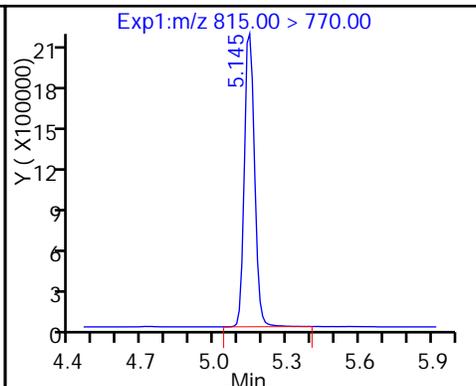
33 Perfluorotetradecanoic acid



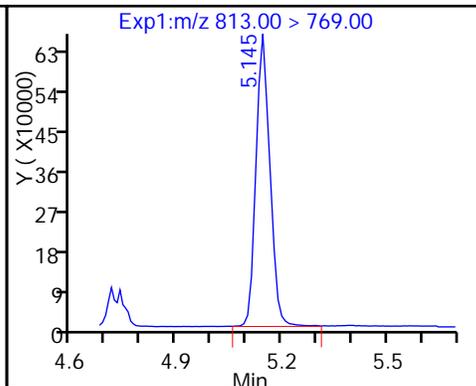
33 Perfluorotetradecanoic acid



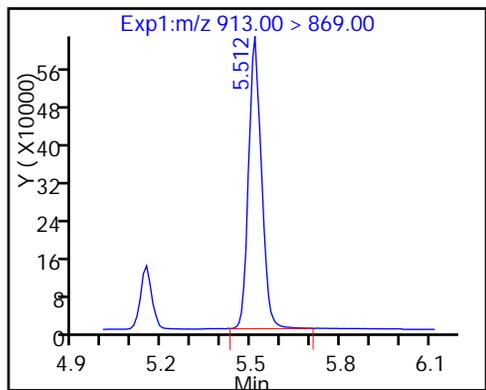
D 34 13C2-PFHxDA



35 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid



LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Start Date: 12/02/2016 10:06

Analysis Batch Number: 140382 End Date: 12/02/2016 14:22

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
RB 320-140382/1 CCB		12/02/2016 10:06	1		Acquity 2.1 (mm)
RB 320-140382/2 CCB		12/02/2016 10:14	1		Acquity 2.1 (mm)
RB 320-140382/3 CCB		12/02/2016 10:22	1		Acquity 2.1 (mm)
IC 320-140382/4		12/02/2016 10:29	1	02DEC2016A_004.d	Acquity 2.1 (mm)
IC 320-140382/5		12/02/2016 10:37	1	02DEC2016A_005.d	Acquity 2.1 (mm)
IC 320-140382/6		12/02/2016 10:44	1	02DEC2016A_006.d	Acquity 2.1 (mm)
IC 320-140382/7		12/02/2016 10:52	1	02DEC2016A_007.d	Acquity 2.1 (mm)
IC 320-140382/8		12/02/2016 10:59	1	02DEC2016A_008.d	Acquity 2.1 (mm)
IC 320-140382/9		12/02/2016 11:07	1	02DEC2016A_009.d	Acquity 2.1 (mm)
IC 320-140382/10		12/02/2016 11:14	1	02DEC2016A_010.d	Acquity 2.1 (mm)
ICB 320-140382/11		12/02/2016 11:22	1		Acquity 2.1 (mm)
ICV 320-140382/12		12/02/2016 11:29	1	02DEC2016A_012.d	Acquity 2.1 (mm)
RB 320-140382/13 CCB		12/02/2016 11:37	1		Acquity 2.1 (mm)
IC 320-140382/14		12/02/2016 11:44	1	02DEC2016A_014.d	Acquity 2.1 (mm)
IC 320-140382/15		12/02/2016 11:52	1	02DEC2016A_015.d	Acquity 2.1 (mm)
IC 320-140382/16		12/02/2016 11:59	1	02DEC2016A_016.d	Acquity 2.1 (mm)
IC 320-140382/17		12/02/2016 12:07	1	02DEC2016A_017.d	Acquity 2.1 (mm)
IC 320-140382/18		12/02/2016 12:14	1	02DEC2016A_018.d	Acquity 2.1 (mm)
IC 320-140382/19		12/02/2016 12:22	1	02DEC2016A_019.d	Acquity 2.1 (mm)
IC 320-140382/20		12/02/2016 12:29	1	02DEC2016A_020.d	Acquity 2.1 (mm)
ICB 320-140382/21		12/02/2016 12:36	1		Acquity 2.1 (mm)
ICV 320-140382/22		12/02/2016 12:44	1		Acquity 2.1 (mm)
RB 320-140382/23 CCB		12/02/2016 12:51	1		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 12:59	100		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:06	100		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:14	100		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:22	10		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:29	10		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:37	10		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:44	10		Acquity 2.1 (mm)
ZZZZZ		12/02/2016 13:52	10		Acquity 2.1 (mm)
RB 320-140382/32 CCB		12/02/2016 13:59	1		Acquity 2.1 (mm)
CCV 320-140382/33		12/02/2016 14:07	1		Acquity 2.1 (mm)
CCV 320-140382/34		12/02/2016 14:14	1		Acquity 2.1 (mm)
RB 320-140382/35 CCB		12/02/2016 14:22	1		Acquity 2.1 (mm)

LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Start Date: 12/02/2016 14:29

Analysis Batch Number: 140429 End Date: 12/02/2016 17:52

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
RB 320-140429/1 CCB		12/02/2016 14:29	1		Acquity 2.1(mm)
CCV 320-140429/2		12/02/2016 14:37	1	02DEC2016B_014.d	Acquity 2.1(mm)
CCV 320-140429/3		12/02/2016 14:44	1		Acquity 2.1(mm)
RB 320-140429/4 CCB		12/02/2016 14:52	1		Acquity 2.1(mm)
MB 320-139627/1-A		12/02/2016 14:59	1	02DEC2016C_001.d	Acquity 2.1(mm)
LCS 320-139627/2-A		12/02/2016 15:07	1	02DEC2016C_002.d	Acquity 2.1(mm)
ZZZZZ		12/02/2016 15:14	1		Acquity 2.1(mm)
ZZZZZ		12/02/2016 15:22	1		Acquity 2.1(mm)
ZZZZZ		12/02/2016 15:29	1		Acquity 2.1(mm)
ZZZZZ		12/02/2016 15:37	1		Acquity 2.1(mm)
ZZZZZ		12/02/2016 15:44	1		Acquity 2.1(mm)
320-23783-1		12/02/2016 15:52	1	02DEC2016C_008.d	Acquity 2.1(mm)
320-23783-2		12/02/2016 15:59	1	02DEC2016C_009.d	Acquity 2.1(mm)
320-23783-5		12/02/2016 16:07	1	02DEC2016C_010.d	Acquity 2.1(mm)
ZZZZZ		12/02/2016 16:14	10		Acquity 2.1(mm)
CCV 320-140429/16		12/02/2016 16:22	1	02DEC2016C_012.d	Acquity 2.1(mm)
CCV 320-140429/17		12/02/2016 16:29	1		Acquity 2.1(mm)
RB 320-140429/18 CCB		12/02/2016 16:37	1		Acquity 2.1(mm)
320-23783-6		12/02/2016 16:44	1	02DEC2016C_015.d	Acquity 2.1(mm)
320-23783-7		12/02/2016 16:52	1	02DEC2016C_016.d	Acquity 2.1(mm)
320-23783-8		12/02/2016 16:59	1	02DEC2016C_017.d	Acquity 2.1(mm)
320-23783-9		12/02/2016 17:07	1	02DEC2016C_018.d	Acquity 2.1(mm)
320-23783-10		12/02/2016 17:14	1	02DEC2016C_019.d	Acquity 2.1(mm)
320-23783-11		12/02/2016 17:22	1	02DEC2016C_020.d	Acquity 2.1(mm)
RB 320-140429/25 CCB		12/02/2016 17:29	1		Acquity 2.1(mm)
CCV 320-140429/26		12/02/2016 17:37	1	02DEC2016C_022.d	Acquity 2.1(mm)
CCV 320-140429/27		12/02/2016 17:44	1		Acquity 2.1(mm)
RB 320-140429/28 CCB		12/02/2016 17:52	1		Acquity 2.1(mm)

LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Start Date: 12/03/2016 13:26

Analysis Batch Number: 140564 End Date: 12/03/2016 18:33

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
RB 320-140564/1 CCB		12/03/2016 13:26	1		Acquity 2.1(mm)
RB 320-140564/2 CCB		12/03/2016 13:33	1		Acquity 2.1(mm)
RB 320-140564/3 CCB		12/03/2016 13:41	1		Acquity 2.1(mm)
IC 320-140564/4		12/03/2016 13:48	1	03DEC2016A_004.d	Acquity 2.1(mm)
IC 320-140564/5		12/03/2016 13:56	1	03DEC2016A_005.d	Acquity 2.1(mm)
IC 320-140564/6		12/03/2016 14:03	1	03DEC2016A_006.d	Acquity 2.1(mm)
IC 320-140564/7		12/03/2016 14:11	1	03DEC2016A_007.d	Acquity 2.1(mm)
IC 320-140564/8		12/03/2016 14:18	1	03DEC2016A_008.d	Acquity 2.1(mm)
IC 320-140564/9		12/03/2016 14:26	1	03DEC2016A_009.d	Acquity 2.1(mm)
ICB 320-140564/10		12/03/2016 14:33	1	03DEC2016A_010.d	Acquity 2.1(mm)
ICV 320-140564/11		12/03/2016 14:41	1	03DEC2016A_011.d	Acquity 2.1(mm)
RB 320-140564/12 CCB		12/03/2016 14:48	1		Acquity 2.1(mm)
IC 320-140564/13		12/03/2016 14:56	1	03DEC2016A_013.d	Acquity 2.1(mm)
IC 320-140564/14		12/03/2016 15:03	1	03DEC2016A_014.d	Acquity 2.1(mm)
IC 320-140564/15		12/03/2016 15:11	1	03DEC2016A_015.d	Acquity 2.1(mm)
IC 320-140564/16		12/03/2016 15:18	1	03DEC2016A_016.d	Acquity 2.1(mm)
IC 320-140564/17		12/03/2016 15:26	1	03DEC2016A_017.d	Acquity 2.1(mm)
IC 320-140564/18		12/03/2016 15:33	1	03DEC2016A_018.d	Acquity 2.1(mm)
ICB 320-140564/19		12/03/2016 15:41	1		Acquity 2.1(mm)
ICV 320-140564/20		12/03/2016 15:48	1		Acquity 2.1(mm)
RB 320-140564/21 CCB		12/03/2016 15:56	1		Acquity 2.1(mm)
MB 320-139749/1-A		12/03/2016 16:03	1	03DEC2016B_001.d	Acquity 2.1(mm)
LCS 320-139749/2-A		12/03/2016 16:11	1	03DEC2016B_002.d	Acquity 2.1(mm)
LCSD 320-139749/3-A		12/03/2016 16:18	1	03DEC2016B_003.d	Acquity 2.1(mm)
320-23783-3		12/03/2016 16:26	1	03DEC2016B_004.d	Acquity 2.1(mm)
320-23783-4		12/03/2016 16:33	1	03DEC2016B_005.d	Acquity 2.1(mm)
320-23783-12		12/03/2016 16:41	1	03DEC2016B_006.d	Acquity 2.1(mm)
320-23783-13		12/03/2016 16:48	1	03DEC2016B_007.d	Acquity 2.1(mm)
RB 320-140564/29 CCB		12/03/2016 16:56	1		Acquity 2.1(mm)
CCV 320-140564/30		12/03/2016 17:03	1	03DEC2016B_009.d	Acquity 2.1(mm)
CCV 320-140564/31		12/03/2016 17:11	1		Acquity 2.1(mm)
RB 320-140564/32 CCB		12/03/2016 17:18	1		Acquity 2.1(mm)
RB 320-140564/39 CCB		12/03/2016 18:11	1		Acquity 2.1(mm)
CCV 320-140564/40		12/03/2016 18:18	1		Acquity 2.1(mm)
CCV 320-140564/41		12/03/2016 18:26	1		Acquity 2.1(mm)
RB 320-140564/42 CCB		12/03/2016 18:33	1		Acquity 2.1(mm)

LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Start Date: 12/07/2016 12:18

Analysis Batch Number: 141054 End Date: 12/07/2016 16:25

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
RB 320-141054/3 CCB		12/07/2016 12:18	1		Acquity 2.1(mm)
RB 320-141054/4 CCB		12/07/2016 12:25	1		Acquity 2.1(mm)
CCV 320-141054/5 CCVL		12/07/2016 12:33	1	07DEC2016A_003. d	Acquity 2.1(mm)
CCV 320-141054/6 CCVL		12/07/2016 12:40	1		Acquity 2.1(mm)
CCV 320-141054/7		12/07/2016 12:48	1		Acquity 2.1(mm)
CCV 320-141054/8		12/07/2016 12:55	1		Acquity 2.1(mm)
RB 320-141054/9 CCB		12/07/2016 13:03	1		Acquity 2.1(mm)
RB 320-141054/20 CCB		12/07/2016 14:26	1		Acquity 2.1(mm)
CCV 320-141054/21		12/07/2016 14:33	1		Acquity 2.1(mm)
CCV 320-141054/22		12/07/2016 14:41	1		Acquity 2.1(mm)
RB 320-141054/23 CCB		12/07/2016 14:48	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 14:55	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:03	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:11	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:18	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:26	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:33	1		Acquity 2.1(mm)
ZZZZZ		12/07/2016 15:41	1		Acquity 2.1(mm)
RB 320-141054/31 CCB		12/07/2016 15:48	1		Acquity 2.1(mm)
RB 320-141054/32 CCB		12/07/2016 15:55	1		Acquity 2.1(mm)
RB 320-141054/33 CCB		12/07/2016 16:03	1		Acquity 2.1(mm)
CCV 320-141054/34		12/07/2016 16:10	1		Acquity 2.1(mm)
CCV 320-141054/35		12/07/2016 16:18	1		Acquity 2.1(mm)
RB 320-141054/36 CCB		12/07/2016 16:25	1		Acquity 2.1(mm)

LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: A8_N Start Date: 12/07/2016 16:40

Analysis Batch Number: 141062 End Date: 12/07/2016 21:10

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
CCV 320-141062/2		12/07/2016 16:40	1	07DEC2016B_002.d	Acquity 2.1(mm)
CCV 320-141062/3		12/07/2016 16:48	1		Acquity 2.1(mm)
RB 320-141062/4 CCB		12/07/2016 16:55	1		Acquity 2.1(mm)
320-23783-12 DL		12/07/2016 17:03	100	07DEC2016B_005.d	Acquity 2.1(mm)
320-23783-3 DL		12/07/2016 17:10	10	07DEC2016B_006.d	Acquity 2.1(mm)
320-23783-4 DL		12/07/2016 17:18	10	07DEC2016B_007.d	Acquity 2.1(mm)
320-23783-13 DL		12/07/2016 17:25	10	07DEC2016B_008.d	Acquity 2.1(mm)
320-23783-5 DL		12/07/2016 17:33	10	07DEC2016B_009.d	Acquity 2.1(mm)
ZZZZZ		12/07/2016 17:40	1		Acquity 2.1(mm)
RB 320-141062/11 CCB		12/07/2016 17:48	1		Acquity 2.1(mm)
CCV 320-141062/12		12/07/2016 17:55	1	07DEC2016B_012.d	Acquity 2.1(mm)
CCV 320-141062/13		12/07/2016 18:03	1		Acquity 2.1(mm)
RB 320-141062/14 CCB		12/07/2016 18:10	1		Acquity 2.1(mm)
RB 320-141062/20 CCB		12/07/2016 18:55	1		Acquity 2.1(mm)
RB 320-141062/21 CCB		12/07/2016 19:03	1		Acquity 2.1(mm)
CCV 320-141062/22		12/07/2016 19:10	1		Acquity 2.1(mm)
CCV 320-141062/23		12/07/2016 19:18	1		Acquity 2.1(mm)
RB 320-141062/24 CCB		12/07/2016 19:25	1		Acquity 2.1(mm)
RB 320-141062/36 CCB		12/07/2016 20:55	1		Acquity 2.1(mm)
CCV 320-141062/37		12/07/2016 21:03	1		Acquity 2.1(mm)
CCV 320-141062/38		12/07/2016 21:10	1		Acquity 2.1(mm)

LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 139627 Batch Start Date: 11/28/16 11:03 Batch Analyst: Arauz, Horacio J

Batch Method: SHAKE Batch End Date: 11/30/16 16:30

Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount	LCMPFCSU 00046	LCPFCSU 00070		
MB 320-139627/1		SHAKE, 537 (Modified)		5.00 g	1.00 mL	50 uL			
LCS 320-139627/2		SHAKE, 537 (Modified)		5.00 g	1.00 mL	50 uL	40 uL		
320-23783-A-1	DPT-16-35-SO-00-01	SHAKE, 537 (Modified)	T	5.01 g	1.00 mL	50 uL			
320-23783-A-2	DPT-16-35-SO-19-20	SHAKE, 537 (Modified)	T	5.07 g	1.00 mL	50 uL			
320-23783-A-5	DPT-16-19-SO-00-01	SHAKE, 537 (Modified)	T	5.02 g	1.00 mL	50 uL			
320-23783-A-6	DPT-16-19-SO-17-18	SHAKE, 537 (Modified)	T	5.06 g	1.00 mL	50 uL			
320-23783-A-7	DPT-16-32-SO-14-15	SHAKE, 537 (Modified)	T	5.03 g	1.00 mL	50 uL			
320-23783-A-8	DPT-16-31-SO-14-15	SHAKE, 537 (Modified)	T	5.00 g	1.00 mL	50 uL			
320-23783-A-9	DPT-16-30-SO-14-15	SHAKE, 537 (Modified)	T	5.02 g	1.00 mL	50 uL			
320-23783-A-10	DPT-16-28-SO-14-15	SHAKE, 537 (Modified)	T	5.03 g	1.00 mL	50 uL			
320-23783-A-11	DPT-16-15-SO-14-15	SHAKE, 537 (Modified)	T	5.05 g	1.00 mL	50 uL			

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 139627 Batch Start Date: 11/28/16 11:03 Batch Analyst: Arauz, Horacio J

Batch Method: SHAKE Batch End Date: 11/30/16 16:30

Batch Notes	
Acetic Acid ID	429065
Balance ID	QA-070
Batch Comment	Pipette MD05306
Hexane ID	J.T.BAKER 0000146278
Manifold ID	5,9
Methanol ID	789824
Methanolic Potassium Hydroxide ID	757037
Millipore Water Dispense Date	11-29-16
Sodium Hydroxide ID	758321
Ammonium Hydroxide/MeOH ID	794501
Analyst ID - Reagent Drop Witness	NSH
Blank Sand Lot #	Fisher 156690
SPE Cartridge ID	016636277A
SPE Cartridge Type	WAX 150mg

Basis	Basis Description
T	Total/NA

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 139749 Batch Start Date: 11/28/16 16:55 Batch Analyst: Marchenko, Veronika P

Batch Method: 3535 Batch End Date: 11/29/16 14:06

Lab Sample ID	Client Sample ID	Method Chain	Basis	GrossWeight	TareWeight	InitialAmount	FinalAmount	LCMPFCSU 00046	LCPFCSU 00070
MB 320-139749/1		3535, 537 (Modified)			25.53 g	250 mL	0.5 mL	25 uL	
LCS 320-139749/2		3535, 537 (Modified)			25.67 g	250 mL	0.5 mL	25 uL	20 uL
LCSD 320-139749/3		3535, 537 (Modified)			25.81 g	250 mL	0.5 mL	25 uL	20 uL
320-23783-A-3	DPT-16-35-GW-21-25	3535, 537 (Modified)	T	276.82 g	26.59 g	250.2 mL	0.5 mL	25 uL	
320-23783-A-4	DPT-16-35-GW-31-35	3535, 537 (Modified)	T	275.75 g	26.19 g	249.6 mL	0.5 mL	25 uL	
320-23783-A-12	DPT-16-15-GW-18-22	3535, 537 (Modified)	T	275.60 g	25.93 g	249.7 mL	0.5 mL	25 uL	
320-23783-A-13	DPT-16-15-GW-31-35	3535, 537 (Modified)	T	272.23 g	26.37 g	245.9 mL	0.5 mL	25 uL	

Lab Sample ID	Client Sample ID	Method Chain	Basis	AnalysisComment					
MB 320-139749/1		3535, 537 (Modified)							
LCS 320-139749/2		3535, 537 (Modified)							
LCSD 320-139749/3		3535, 537 (Modified)		See NCM regarding these samples, contains sediment.					
320-23783-A-3	DPT-16-35-GW-21-25	3535, 537 (Modified)	T						
320-23783-A-4	DPT-16-35-GW-31-35	3535, 537 (Modified)	T						
320-23783-A-12	DPT-16-15-GW-18-22	3535, 537 (Modified)	T						
320-23783-A-13	DPT-16-15-GW-31-35	3535, 537 (Modified)	T						

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 139749 Batch Start Date: 11/28/16 16:55

Batch Analyst: Marchenko, Veronika P

Batch Method: 3535 Batch End Date: 11/29/16 14:06

Batch Notes	
Balance ID	QA-070
Batch Comment	0.1N NaOH/H2O: 794892
H2O ID	11/28/16
Hexane ID	0000146278
Manifold ID	1
Methanol ID	769613
Pipette ID	MD05306
Analyst ID - Reagent Drop	VPM
Analyst ID - SU Reagent Drop	VPM
Analyst ID - SU Reagent Drop Witness	KMK
Solvent Lot #	794501
Solvent Name	0.3% NH4OH/MeOH
SOP Number	WS-LC-0025
SPE Cartridge Type	WAX 500mg
Solid Phase Extraction Disk ID	002836112A

Basis	Basis Description
T	Total/NA

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

HPLC/LCMS Data Review Checklist

Job Number(s): 23542; 23651; 23718; 23783 Work List ID(s): 37450; 37461
 Extraction Batch: 138217; 138291; 139627 Analysis Batch(es): 140382; 140429
 Delivery Rank: 4 Due Date: 11/19/16; 11/25/16

A. Calibration/Instrument Run QC	1 st Level	2 nd Level	N/A
1. ICAL locked in Chrom and TALS? ICAL Batch# <u>140382;</u>	✓	✓	
2. ICAL, CCV Frequency & Criteria met.	✓	✓	
• RF _{average} criteria appropriate for the method.	✓	✓	
• Linear Regression criteria appropriate if required ($r \geq 0.995$).	✓	✓	
• Quadratic fit criteria appropriate if required ($r^2 \geq 0.990$).			✓
• For Linear Regression and Quadratic fit – Does the y-intercept support ½ the reporting limit as described in CA-Q-S-005?	✓	✓	
• All curve points show calculated concentrations.	✓	✓	
3. Peaks correctly ID'd by data system.	✓	✓	
5. Tune check frequency & criteria met and Tune check report attached.	✓	✓	
B. QA/QC			
1. Are all QC samples properly linked in TALS?	✓	✓	
2. Method blank, LCS/LCSD and MS/SD frequencies met.	✓	✓	
3. LCS/LCSD and MB data are within control limits. If not, NCM is present.	✓	✓	
4. Are MS/MSD recoveries and RPD within control limits?	✓	✓	
5. Holding Times were met for prep and analytical.	✓	✓	
6. IS/Surrogate recoveries meet criteria or properly noted.	✓	✓	
C. Sample Analysis			
1. Was correct analysis performed and were project instructions followed?	✓	✓	
2. If required, are compounds within RT windows?	✓	✓	
3. If required, are positive hits confirmed and >40% RPD flagged?			✓
4. Manual Integrations reviewed and appropriate.	✓	✓	
5. All analytes correctly reported. (Primary, secondary, acceptable status)	✓	✓	
6. Correct reporting limits used. (based on client request, prep factors, and dilutions)	✓	✓	
D. Documentation			
1. Are all non-conformances documented/attached? NCM# <u>71898; 71899</u>	✓	✓	
2. Do results make sense (e.g. dilutions, etc.)?	✓	✓	
3. Have all flags been reviewed for appropriateness?	✓	✓	
4. For level 3 and 4 reports, have forms and raw data been reviewed?	✓	✓	
5. Was QC Checker run for this job?	✓	✓	

*Upon completion of this checklist, the reviewer must scan and attach the checklist to the TALS job.

1st Level (Analyst):  Date: 12/4/16

2nd Level Reviewer:  Date: 12/5/16

NCMS: 71898; 71899; 71937; 71940

TestAmerica Laboratories
Worklist QC Batch Report

Worklist Name: 02DEC2016A_PFC Worklist Number: 37450
 Instrument Name: A8_N Chrom Method: A8_N
 Data Directory: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37450.b
 QC Batching: Disabled Limit Group Batching: Enabled

QC Batch: 1	LC PFC_DOD ICAL Raw Batch: 140382	LC PFC ICAL Raw Batch: 140383	LC PFAS ICAL Raw Batch: 140384
# 1 RB	# 1 RB		
# 2 RB	# 2 RB		
# 3 RB	# 3 RB		
# 4 IC L1	# 4 IC L1	# 4 IC L1	# 4 IC L1
# 5 IC L2	# 5 IC L2	# 5 IC L2	# 5 IC L2
# 6 IC L3	# 6 IC L3	# 6 IC L3	# 6 IC L3
# 7 IC L4	# 7 IC L4	# 7 IC L4	# 7 IC L4
# 8 IC L5	# 8 IC L5	# 8 IC L5	# 8 IC L5
# 9 IC L6	# 9 IC L6	# 9 IC L6	# 9 IC L6
#10 IC L7	#10 IC L7	#10 IC L7	#10 IC L7
#11 ICB	#11 ICB	#11 ICB	#11 ICB
#12 ICV	#12 ICV <i>Tune NCM</i>	#12 ICV	
#13 RB	#13 RB <i>71898</i>	#13 RB	#13 RB
#14 IC L1 Add-on	#14 IC L1 Add-on	#14 IC L1 Add-on	#14 IC L1 Add-on
#15 IC L2 Add-on	#15 IC L2 Add-on	#15 IC L2 Add-on	#15 IC L2 Add-on
#16 IC L3 Add-on	#16 IC L3 Add-on	#16 IC L3 Add-on	#16 IC L3 Add-on
#17 IC L4 Add-on	#17 IC L4 Add-on	#17 IC L4 Add-on	#17 IC L4 Add-on
#18 IC L5 Add-on	#18 IC L5 Add-on	#18 IC L5 Add-on	#18 IC L5 Add-on
#19 IC L6 Add-on	#19 IC L6 Add-on	#19 IC L6 Add-on	#19 IC L6 Add-on
#20 IC L7 Add-on	#20 IC L7 Add-on	#20 IC L7 Add-on	#20 IC L7 Add-on
#21 ICB	#21 ICB	#21 ICB	#21 ICB
#22 ICV Add-on	#22 ICV Add-on	#22 ICV Add-on	
#23 RB	#23 RB	#23 RB	#23 RB
#24 320-23542-A-2-A	#24 320-23542-A-2-A		
#25 320-23542-A-2-B MS	#25 320-23542-A-2-B MS		
#26 320-23542-A-2-C MSD	#26 320-23542-A-2-C MSD		
#27 320-23542-A-1-A	#27 320-23542-A-1-A		
#28 320-23542-A-4-A	#28 320-23542-A-4-A <i>Misinject</i>		
#29 320-23542-A-3-A	#29 320-23542-A-3-A		
#30 320-23542-A-3-B MS	#30 320-23542-A-3-B MS		
#31 320-23542-A-3-C MSD	#31 320-23542-A-3-C MSD		
#32 RB	#32 RB	#32 RB	#32 RB
#33 CCV L5	#33 CCV L5	#33 CCV L5	#33 CCV L5
#34 CCV L5 Add-on	#34 CCV L5 Add-on	#34 CCV L5 Add-on	#34 CCV L5 Add-on
#35 RB	#35 RB	#35 RB	#35 RB

MS/MSD high targets NCM 71899

TestAmerica Laboratories
Worklist QC Batch Report

Worklist Name: 02DEC2016B_PFC Worklist Number: 37461
 Instrument Name: A8_N Chrom Method: A8_N
 Data Directory: \\ChromNa\Sacramento\ChromData\A8_N\20161202-37461.b
 QC Batching: Disabled Limit Group Batching: Enabled

QC Batch: 1	LC PFC_DOD ICAL Raw Batch: 140429	LC PFC ICAL Raw Batch: 140430	LC PFAS ICAL Raw Batch: 140431
# 1 RB	# 1 RB	# 1 RB	# 1 RB
# 2 CCV L5	# 2 CCV L5	# 2 CCV L5	# 2 CCV L5
# 3 CCV L5 Add-on	# 3 CCV L5 Add-on	# 3 CCV L5 Add-on	# 3 CCV L5 Add-on
# 4 RB	# 4 RB	# 4 RB	# 4 RB
# 5 MB 320-139627/1-A	# 5 MB 320-139627/1-A		
# 6 LCS 320-139627/2-A	# 6 LCS 320-139627/2-A		
# 7 320-23651-A-2-A	# 7 320-23651-A-2-A		
# 8 320-23651-A-3-A	# 8 320-23651-A-3-A		
# 9 320-23718-A-3-A	# 9 320-23718-A-3-A		
#10 320-23718-A-3-B MS	#10 320-23718-A-3-B MS		
#11 320-23718-A-3-C MSD	#11 320-23718-A-3-C MSD		
#12 320-23783-A-1-A	#12 320-23783-A-1-A		
#13 320-23783-A-2-A	#13 320-23783-A-2-A		
#14 320-23783-A-5-A	#14 320-23783-A-5-A		
#15 320-23542-A-4-A	#15 320-23542-A-4-A		
#16 CCV L5	#16 CCV L5	#16 CCV L5	#16 CCV L5
#17 CCV L5 Add-on	#17 CCV L5 Add-on	#17 CCV L5 Add-on	#17 CCV L5 Add-on
#18 RB	#18 RB	#18 RB	#18 RB
#19 320-23783-A-6-A	#19 320-23783-A-6-A		
#20 320-23783-A-7-A	#20 320-23783-A-7-A		
#21 320-23783-A-8-A	#21 320-23783-A-8-A		
#22 320-23783-A-9-A	#22 320-23783-A-9-A		
#23 320-23783-A-10-A	#23 320-23783-A-10-A		
#24 320-23783-A-11-A	#24 320-23783-A-11-A		
#25 RB	#25 RB	#25 RB	#25 RB
#26 CCV L5	#26 CCV L5	#26 CCV L5	#26 CCV L5
#27 CCV L5 Add-on	#27 CCV L5 Add-on	#27 CCV L5 Add-on	#27 CCV L5 Add-on
#28 RB	#28 RB	#28 RB	#28 RB

E 10x PFOs. E flag NCM 71937

ICV 140382

Tune + SBC 12/4/16

Tune NCM 71998

AS

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

AS 11/20/16
AS 12/2/16

Batch Number: 320-138217
Method Code: 320-3535_IWWT-320

Analyst: Arauz, Horacio J

Batch Open: 11/17/2016 8:49:00AM
Batch End: 11/18/2016 2:20:00PM

Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	InitAmt		PHs Adj1 Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
			FltAmt	FltAmt						
1 MB-320-138217/1 N/A	N/A		250 mL	0.5 mL		N/A	N/A	N/A		
2 LCS-320-138217/2 N/A	N/A		250 mL	0.5 mL		N/A	N/A	N/A		
3 320-23542-A-1 (PFC_IDA_DOD5)	N/A (320-23542-1)	266.74 g 26.14 g	240.6 mL	0.5 mL		11/19/16	16_Days	4	10x PFOS ✓	
4 320-23542-A-2 (PFC_IDA_DOD5)	N/A (320-23542-1)	284.91 g 26.35 g	258.6 mL	0.5 mL		11/19/16	16_Days	4	100x PFOS ✓	
5 320-23542-A-2-MS (PFC_IDA_DOD5)	N/A (320-23542-1)	277.51 g 26.23 g	251.3 mL	0.5 mL		11/19/16	16_Days	4		
6 320-23542-A-2-MSD (PFC_IDA_DOD5)	N/A (320-23542-1)	273.81 g 26.03 g	247.8 mL	0.5 mL		11/19/16	16_Days	4		
7 320-23542-A-4 (PFC_IDA_DOD5)	N/A (320-23542-1)	280.63 g 26.42 g	254.2 mL	0.5 mL		11/19/16	16_Days	4	10x PFOA PI ✓ PFOS	

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Number: 320-138217

Method Code: 320-3535_IVWT-320

Batch Open: 11/17/2016 8:49:00AM

Batch End: 11-6-16 19:20

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Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	Int&Amnt Fin&Amnt	PHs Rcvd Adj1 Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
1 MB-320-1382171 N/A	N/A		250 mL 0.5 mL		N/A	N/A	N/A		
2 LGS-320-1382172 N/A	N/A		250 mL 0.5 mL		N/A	N/A	N/A		
3 320-23542-A-1 (PFC_IDA_DOD5)	N/A (320-23542-1)	266.74 g	0.5 mL		11/19/16	16_Days	4		
4 320-23542-A-2 (PFC_IDA_DOD5)	N/A (320-23542-1)	284.91 g	0.5 mL		11/19/16	16_Days	4		
5 320-23542-A-2-MS (PFC_IDA_DOD5)	N/A (320-23542-1)	277.51 g	0.5 mL		11/19/16	16_Days	4		
6 320-23542-A-2-MSD (PFC_IDA_DOD5)	N/A (320-23542-1)	273.81 g	0.5 mL		11/19/16	16_Days	4		
7 320-23542-A-4 (PFC_IDA_DOD5)	N/A (320-23542-1)	280.63 g	0.5 mL		11/19/16	16_Days	4		

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Open: 11/17/2016 8:49:00AM

Batch End:

Batch Number: 320-138217

Method Code: 320-3535_IVWT-320

Batch Notes

Manifold ID 5

Methanol ID 769617

Hexane ID 000146278

Sodium Hypochlorite ID NA

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type WAX 500mg

Solid Phase Extraction Disk ID 002836112A

H2O ID 11-14-16

Pipette ID MD05306

Solvent Name 0.3% NH4OH-Me

Solvent Lot # 776672

Analyst ID - Reagent Drop HJA

Analyst ID - SU Reagent Drop HJA

Analyst ID - SU Reagent Drop Witness *MSK*

Acid Name NA

Acid ID NA

Reagent ID NA

Reagent Lot Number NA

NaCl ID NA

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Open: 11/17/2016 8:49:00AM

Batch End:

Batch Number: 320-138217
Method Code: 320-3535_IVWT-320

SOP Number WS-LC-0025

Batch Comment 0.1N NaOH/H2O_00026

Comments

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-138217
 Method Code: 320-3535_JVWT-320

Analyst: Arauz, Horacio J

Batch Open: 11/17/2016 8:49:00AM
 Batch End:

Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-138217/1	LCMPFCSU_00046	25 uL	0.5 mL	HSA 11-17-16	New 11-17-16
LCS 320-138217/2	LCMPFCSU_00046	25 uL	0.5 mL		
LCS 320-138217/2	LCPFCSU_00066	20 uL	0.5 mL		
320-23542-A-1	LCMPFCSU_00046	25 uL	0.5 mL		
320-23542-A-2	LCMPFCSU_00046	25 uL	0.5 mL		
320-23542-A-2 MS	LCMPFCSU_00046	25 uL	0.5 mL		
320-23542-A-2 MS	LCPFCSU_00066	20 uL	0.5 mL		
320-23542-A-2 MSD	LCMPFCSU_00046	25 uL	0.5 mL		
320-23542-A-2 MSD	LCPFCSU_00066	20 uL	0.5 mL		
320-23542-A-4	LCMPFCSU_00046	25 uL	0.5 mL		

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Open: 11/17/2016 8:49:00AM

Batch End:

Batch Number: 320-138217

Method Code: 320-3535_JVWT-320

Reagent	Other Reagents:	Amount/Units	Lot#:

Preparation Batch Number(s): 320-138217 Test: PFC-L

Earliest Holding Time: 11-21-16

Sample List Tab		1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method		/	/
All necessary NCMs filed (including holding time)		/	/
Method/sample/login/QAS checked and correct		/	/
Worksheet Tab		1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved		NA	NA
Weights in anticipated range and not targeted		/	/
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and CI Check)		/	/
The pH is transcribed correctly in TALS		NA	NA
All additional information transcribed into TALS is correct and raw data is attached		/	/
Comments are transcribed correctly in TALS		/	/
Reagents Tab		1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS		/	/
All spike amounts correct and added to necessary samples and QC		/	/
Batch Information		1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly		/	/
All necessary 'batch information' complete and entered into TALS correctly		/	/

1st Level Reviewer: HSA

Date: 11-18-16

2nd Level Reviewer: VPM

Date: 11/18/16

Comments: _____

Method ID PFC-IDA-DODS

Job # 23542

Analyst (Print Name) Shykhara Chandrasena Analyst Initials SBC

Date 12/2/16

Sample#	Original F.V. (uL)	Aliquot (uL)	Dilution F.V. (uL)	Dilution Factor
1	500	100	1000	10x
2	↓	↓	↓	↓
2MS	↓	↓	↓	↓
2MSD	↓	↓	↓	↓
3	↓	↓	↓	↓
3MS	↓	↓	↓	↓
3MSD	↓	↓	↓	↓
4	↓	↓	↓	↓
2+	↓	↓	↓	100x
2-MS	↓	↓	↓	100x
2MSD	↓	↓	↓	100x

Comments:

Solid SW-846-3500 Analysis Sheet
 (To Accompany Samples to Instruments)

Batch Number: 320-138291
 Method Code: 320-Shake_Bath_14D-320

Analyst: Sharifi, Nooshin

Batch Open: 11/17/2016 12:46:00PM

Batch End: 11-19-16 13:18pm

AD 11/20/16
 AB 12/2/16

Shake Extraction with Ultrasonic Bath Extraction

Input Sample Lab ID (Analytical Method)	SDG (Job #)	Initial Amount	Final Amount	Due Date	Analytical TAT	DIV Rank	Comments	Output Sample Lab ID
1 N/A MB-320-138291/1	N/A	4.98 g	1.0 mL	N/A	N/A	N/A		
2 N/A LCS-320-138291/2	N/A	5.03 g	1.0 mL	N/A	N/A	N/A		
3 320-23542-A-3 (PFC_IDA_DOD5)	N/A (320-23542-1)	5.05 g	1.0 mL	11/19/16	16_Days	4	10x PPOS ✓	
4 320-23542-A-3-MS (PFC_IDA_DOD5)	N/A (320-23542-1)	4.99 g	1.0 mL	11/19/16	16_Days	4	10x PPOS ✓	
5 320-23542-A-3-MSD (PFC_IDA_DOD5)	N/A (320-23542-1)	5.03 g	1.0 mL	11/19/16	16_Days	4	10x PPOS ✓	

100

Solid SW-846-3500 Analysis Sheet
(To Accompany Samples to Instruments)

Batch Number: 320-138291

Method Code: 320-Shake_Bath_14D-320

Analyst: Sharifi, Nooshin

Batch Open: 11/17/2016 12:46:00PM

Batch End:

Batch Notes

Balance ID QA-070

Blank Sand Lot # 156690

Filter ID NA

Millipore Water Dispense Date

Analyst ID - Reagent Drop Witness NSH

SPE Cartridge ID 016336091A

SPE Cartridge Type WAX150mg

Hexane ID 0000146278

Methanol ID 769617

Ammonium Hydroxide/MeOH ID 776672

Sodium Hydroxide ID 758321

Methanolic Potassium Hydroxide ID 757037

Manifold ID

Interference check solution ID NA

Acetic Acid ID 429065

Batch Comment Pipette ID: MD05306, MG05455

Comments

Solid SW-846-3500 Analysis Sheet
 (To Accompany Samples to Instruments)

Batch Number: 320-138291
 Method Code: 320-Shake_Bath_14D-320

Analyst: Sharifi, Nooshin

Batch Open: 11/17/2016 12:46:00PM
 Batch End:

Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-138291/1	LCMPFCSU_00046	50 uL	1.0 mL	NSH 11-17-16	HSA 11-17-16
LCS 320-138291/2	LCMPFCSU_00046	50 uL	1.0 mL		
LCS 320-138291/2	LCPFCSU_00064	40 uL	1.0 mL		
320-23542-A-3	LCMPFCSU_00046	50 uL	1.0 mL		
320-23542-A-3 MS	LCMPFCSU_00046	50 uL	1.0 mL		
320-23542-A-3 MS	LCPFCSU_00064	40 uL	1.0 mL		
320-23542-A-3 MSD	LCMPFCSU_00046	50 uL	1.0 mL		
320-23542-A-3 MSD	LCPFCSU_00064	40 uL	1.0 mL		

Other Reagents:	Amount/Units	Lot#:

Preparation Batch Number(s): 138291 Test: PFC-S
 Earliest Holding Time: 11-28-16

Sample List Tab		1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method		/	/
All necessary NCMs filed (including holding time)		NA	NA
Method/sample/login/QAS checked and correct		/	/
Worksheet Tab		1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved		NA	NA
Weights in anticipated range and not targeted		/	/
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and CI Check)		/	/
The pH is transcribed correctly in TALS		NA	NA
All additional information transcribed into TALS is correct and raw data is attached		/	/
Comments are transcribed correctly in TALS		/	/
Reagents Tab		1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS		/	/
All spike amounts correct and added to necessary samples and QC		/	/
Batch Information		1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly		/	/
All necessary 'batch information' complete and entered into TALS correctly		/	/

1st Level Reviewer: NSH Date: 11-19-16
 2nd Level Reviewer: HSA Date: 11-19-16
 Comments: _____

Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

AB 12/2/16

Batch Number: 320-139627

Analyst: Arauz, Horacio J

Batch Open: 11/28/2016 11:03:00AM

Method Code: 320-Shake_Bath_14D-320

Batch End: 11-30-16 16:30

Shake Extraction with Ultrasonic Bath Extraction

Input Sample Lab ID (Analytical Method)	SDG (Job #)	Initial Amount	Final Amount	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
MB-320-139627/1 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		MB 320-139627/1-A
LCS-320-139627/2 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		LCS 320-139627/2-A
320-23651-A-2 (PFC_IDA_DOD5)	N/A	5.07 g	1.00 mL	11/21/16	16_Days	4		LCS 320-23651-A-2-A
320-23651-A-3 (PFC_IDA_DOD5)	(320-23651-1)	5.05 g	1.00 mL	11/21/16	16_Days	4		320-23651-A-2-A
320-23718-A-3 (PFC_IDA_DOD5)	N/A	5.00 g	1.00 mL	11/25/16	16_Days	4		320-23718-A-3-A
320-23718-A-3-MS (PFC_IDA_DOD5)	(320-23718-1)	5.01 g	1.00 mL	11/25/16	16_Days	4		320-23718-A-3-A
320-23718-A-3-MSD (PFC_IDA_DOD5)	N/A	5.08 g	1.00 mL	11/25/16	16_Days	4		320-23718-A-3-B
320-23783-A-1 (PFC_IDA_DOD5)	(320-23783-1)	5.01 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-3-C
320-23783-A-2 (PFC_IDA_DOD5)	N/A	5.07 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-1-A
320-23783-A-5 (PFC_IDA_DOD5)	(320-23783-1)	5.02 g	1.00 mL	11/29/16	11_Day_Rush	4	10X PFOS 681.79	320-23783-A-2-A
320-23783-A-6 (PFC_IDA_DOD5)	N/A	5.06 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-3-A
320-23783-A-7 (PFC_IDA_DOD5)	N/A	5.03 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-6-A
320-23783-A-8 (PFC_IDA_DOD5)	(320-23783-1)	5.00 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-7-A
320-23783-A-9 (PFC_IDA_DOD5)	N/A	5.02 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-8-A
320-23783-A-10 (PFC_IDA_DOD5)	(320-23783-1)	5.03 g	1.00 mL	11/29/16	11_Day_Rush	4		320-23783-A-9-A

Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Open: 11/28/2016 11:03:00AM
Batch End:

Batch Number: 320-139627

Method Code: 320-Shake_Bath_14D-320

320-23783-A-11 (PFC_IDA_DOD5)	N/A (320-23783-1)	5.05 g	1.00 mL	11/29/16	11_Day_Rush	4
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Batch Notes

Balance ID QA-070

Blank Sand Lot # Fisher 156690

Filter ID NA

Millipore Water Dispense Date 11-29-16

Analyst ID - Reagent Drop Witness MSH

SPE Cartridge ID 016636277A

SPE Cartridge Type WAX 150mg

Hexane ID J.T.BAKER 0000146278

Methanol ID 789824

Ammonium Hydroxide/MeOH ID 794501

Sodium Hydroxide ID 758321

Methanolic Potassium Hydroxide ID 757037

Manifold ID 5,9

Interference check solution ID NA

Acetic Acid ID 429065

Batch Comment Pipette MD05306

Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139627

Analyst: Arauz, Horacio J

Batch Open: 11/28/2016 11:03:00AM

Method Code: 320-Shake_Bath_14D-320

Batch End:

Comments

Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139627

Analyst: Arauz, Horacio J

Batch Open: 11/28/2016 11:03:00AM

Method Code: 320-Shake_Bath_14D-320

Batch End:

Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-139627/1	LCMPFCSU_00046	50 uL	1.00 mL	HSA 11-28-16	NSH 11-28-16
LCS 320-139627/2	LCMPFCSU_00046	50 uL	1.00 mL		
LCS 320-139627/2	LCPFCSU_00070	40 uL	1.00 mL		
320-23651-A-2	LCMPFCSU_00046	50 uL	1.00 mL		
320-23651-A-3	LCMPFCSU_00046	50 uL	1.00 mL		
320-23718-A-3	LCMPFCSU_00046	50 uL	1.00 mL		
320-23718-A-3 MS	LCMPFCSU_00046	50 uL	1.00 mL		
320-23718-A-3 MS	LCPFCSU_00070	40 uL	1.00 mL		
320-23718-A-3 MSD	LCMPFCSU_00046	50 uL	1.00 mL		
320-23718-A-3 MSD	LCPFCSU_00070	40 uL	1.00 mL		
320-23783-A-1	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-2	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-5	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-6	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-7	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-8	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-9	LCMPFCSU_00046	50 uL	1.00 mL		
320-23783-A-10	LCMPFCSU_00046	50 uL	1.00 mL		

Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Arauz, Horacio J

Batch Open: 11/28/2016 11:03:00AM

Batch End:

Batch Number: 320-139627

Method Code: 320-Shake_Bath_14D-320

320-23783-A-11	LCMPFCSU_00046	50 uL	1.00 mL	HSA 11-28-16	NSH 11-28-16
----------------	----------------	-------	---------	--------------	--------------

Other Reagents:

Reagent

Amount/Units

Lot#:

Preparation Batch Number(s): 320-139627 Test: PEC-1DA-0005

Earliest Holding Time: 11-29-16

Sample List Tab		
	1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method	/	/
All necessary NCMs filed (including holding time)	NA	NA
Method/sample/login/QAS checked and correct	/	/
Worksheet Tab		
	1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved	/	/
Weights in anticipated range and not targeted	NA	NA
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and CI Check)	/	/
The pH is transcribed correctly in TALS	NA	NA
All additional information transcribed into TALS is correct and raw data is attached	/	/
Comments are transcribed correctly in TALS	/	/
Reagents Tab		
	1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS	/	/
All spike amounts correct and added to necessary samples and QC	/	/
Batch Information		
	1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly	/	/
All necessary 'batch information' complete and entered into TALS correctly	/	/

1st Level Reviewer: HSA

Date: 11-30-16

2nd Level Reviewer: VPM

Date: 11-30-16

Comments: _____

HPLC/LCMS Data Review Checklist

Job Number(s): 23793; 23770; 23801

Work List ID(s): 37491

Extraction Batch: 139749; 139834

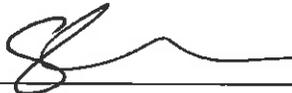
Analysis Batch(es): 140564; 140565

Delivery Rank 4; 2

Due Date: 11/29/16; 11/30/16

A. Calibration/Instrument Run QC	1 st Level	2 nd Level	N/A
1. ICAL locked in Chrom and TALS? ICAL Batch#	✓	✓	
2. ICAL, CCV Frequency & Criteria met.	✓	✓	
• RF _{average} criteria appropriate for the method.	✓	✓	
• Linear Regression criteria appropriate if required ($r \geq 0.995$).	✓	✓	
• Quadratic fit criteria appropriate if required ($r^2 \geq 0.990$).			✓
• For Linear Regression and Quadratic fit – Does the y-intercept support ½ the reporting limit as described in CA-Q-S-005?	✓	✓	
• All curve points show calculated concentrations.	✓	✓	
3. Peaks correctly ID'd by data system.	✓	✓	
5. Tune check frequency & criteria met and Tune check report attached.	✓	✓	
B. QA/QC			
1. Are all QC samples properly linked in TALS?	✓	✓	
2. Method blank, LCS/LCSD and MS/SD frequencies met.	✓	✓	
3. LCS/LCSD and MB data are within control limits. If not, NCM is present.	✓	✓	
4. Are MS/MSD recoveries and RPD within control limits? <i>SBC 12/5/16</i>	✓	✓	✓
5. Holding Times were met for prep and analytical.	✓	✓	
6. IS/Surrogate recoveries meet criteria or properly noted.	✓	✓	
C. Sample Analysis			
1. Was correct analysis performed and were project instructions followed?	✓	✓	
2. If required, are compounds within RT windows?	✓	✓	
3. If required, are positive hits confirmed and >40% RPD flagged?			✓
4. Manual Integrations reviewed and appropriate.	✓	✓	
5. All analytes correctly reported. (Primary, secondary, acceptable status)	✓	✓	
6. Correct reporting limits used. (based on client request, prep factors, and dilutions)	✓	✓	
D. Documentation			
1. Are all non-conformances documented/attached? NCM#	✓	✓	
2. Do results make sense (e.g. dilutions, etc.)?	✓	✓	
3. Have all flags been reviewed for appropriateness?	✓	✓	
4. For level 3 and 4 reports, have forms and raw data been reviewed?	✓	✓	
5. Was QC Checker run for this job?	✓	✓	

*Upon completion of this checklist, the reviewer must scan and attach the checklist to the TALS job.

1st Level (Analyst): 

Date: 12/5/16

2nd Level Reviewer: 

Date: 12/6/16

NCMs: 71975; 71977; 71973; 71979

TestAmerica Laboratories
Worklist QC Batch Report

Worklist Name: 03DEC2016A_PFC Worklist Number: 37491
 Instrument Name: A8_N Chrom Method: A8_N
 Data Directory: \\ChromNa\Sacramento\ChromData\A8_N\20161205-37491.b
 QC Batching: Disabled Limit Group Batching: Enabled

QC Batch: 1	LC PFC_DOD ICAL Raw Batch: 140564	LC PFC ICAL Raw Batch: 140565	LC PFAS ICAL Raw Batch: 140566
# 1 RB	# 1 RB		
# 2 RB	# 2 RB		
# 3 RB	# 3 RB		
# 4 IC L1	# 4 IC L1	# 4 IC L1	# 4 IC L1
# 5 IC L2	# 5 IC L2	# 5 IC L2	# 5 IC L2
# 6 IC L3	# 6 IC L3	# 6 IC L3	# 6 IC L3
# 7 IC L4	# 7 IC L4	# 7 IC L4	# 7 IC L4
# 8 IC L5	# 8 IC L5	# 8 IC L5	# 8 IC L5
# 9 IC L6	# 9 IC L6	# 9 IC L6	# 9 IC L6
#10 ICB	#10 ICB	#10 ICB	#10 ICB
#11 ICV	#11 ICV	#11 ICV	
#12 RB	#12 RB	#12 RB	#12 RB
#13 IC L1 Add-on	#13 IC L1 Add-on	#13 IC L1 Add-on	#13 IC L1 Add-on
#14 IC L2 Add-on	#14 IC L2 Add-on	#14 IC L2 Add-on	#14 IC L2 Add-on
#15 IC L3 Add-on	#15 IC L3 Add-on	#15 IC L3 Add-on	#15 IC L3 Add-on
#16 IC L4 Add-on	#16 IC L4 Add-on	#16 IC L4 Add-on	#16 IC L4 Add-on
#17 IC L5 Add-on	#17 IC L5 Add-on	#17 IC L5 Add-on	#17 IC L5 Add-on
#18 IC L6 Add-on	#18 IC L6 Add-on	#18 IC L6 Add-on	#18 IC L6 Add-on
#19 ICB	#19 ICB	#19 ICB	#19 ICB
#20 ICV Add-on	#20 ICV Add-on	#20 ICV Add-on	
#21 RB	#21 RB	#21 RB	#21 RB
#22 MB 320-139749/1-A	#22 MB 320-139749/1-A		
#23 LCS 320-139749/2-A	#23 LCS 320-139749/2-A		
#24 LCSD 320-139749/3-A	#24 LCSD 320-139749/3-A		
#25 320-23783-A-3-A	#25 320-23783-A-3-A 10x		
#26 320-23783-A-4-A	#26 320-23783-A-4-A 10x		
#27 320-23783-A-12-A	#27 320-23783-A-12-A IDA low NCM 71977 100x		
#28 320-23783-A-13-A	#28 320-23783-A-13-A RI, 10x		
#29 RB	#29 RB	#29 RB	#29 RB
#30 CCV L5	#30 CCV L5	#30 CCV L5	#30 CCV L5
#31 CCV L5 Add-on	#31 CCV L5 Add-on	#31 CCV L5 Add-on	#31 CCV L5 Add-on
#32 RB	#32 RB	#32 RB	#32 RB
#33 MB 320-139834/1-A		#33 MB 320-139834/1-A	#33 MB 320-139834/1-A
#34 LCS 320-139834/2-A		#34 LCS 320-139834/2-A	#34 LCS 320-139834/2-A
#35 LCSD 320-139834/3-A		#35 LCSD 320-139834/3-A	#35 LCSD 320-139834/3-A
#36 320-23770-A-1-A		#36 320-23770-A-1-A	#36 320-23770-A-1-A
#37 320-23801-A-1-A		#37 320-23801-A-1-A	#37 320-23801-A-1-A
#38 320-23801-A-2-A		#38 320-23801-A-2-A	#38 320-23801-A-2-A
#39 RB	#39 RB	#39 RB	#39 RB
#40 CCV L5	#40 CCV L5	#40 CCV L5	#40 CCV L5
#41 CCV L5 Add-on	#41 CCV L5 Add-on	#41 CCV L5 Add-on	#41 CCV L5 Add-on
#42 RB	#42 RB	#42 RB	#42 RB

E flag NCM 71977
IDA low NCM 71977 100x

Tune NCM 71975

FOSA IDA low NCM 71973

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139834

Analyst: Marchenko, Veronika P

Batch Open: 11/29/2016 12:27:00PM

Method Code: 320-3535_I/VWT-320

Batch End: 11-29-2016 9:33 PM

A8 11/30/16
12/2/16

Solid-Phase Extraction (SPE)

12/3/16

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	InitAmnt		PHs Rcvd Adj1 Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
			250 mL	0.5 mL						
1 MB-320-139834/1 N/A	N/A		250 mL	0.5 mL		N/A	N/A	N/A		MB 320-139834/1-A
2 LCS-320-139834/2 N/A	N/A	26.51 g	250 mL	0.5 mL		N/A	N/A	N/A		LCS 320-139834/2-A
3 LCSD-320-139834/3 N/A	N/A	25.54 g	250 mL	0.5 mL		N/A	N/A	N/A		LCSD 320-139834/3-A
4 320-23770-A-1 (PFC_IDA)	N/A (320-23770-1)	26.49 g 275.89 g	250 mL	0.5 mL		12/7/16	10_Days	2		320-23770-A-1-A
5 320-23801-A-1 (PFC_IDA)	N/A (320-23801-1)	24.16 g 275.99 g	250 mL	0.5 mL		11/30/16	8_Days	2		320-23801-A-1-A
6 320-23801-A-2 (PFC_IDA)	N/A (320-23801-1)	26.45 g 280.68 g	250 mL	0.5 mL		11/30/16	8_Days	2		320-23801-A-2-A

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139834

Analyst: Marchenko, Veronika P

Batch Open: 11/29/2016 12:27:00PM

Method Code: 320-3535_IVWT-320

Batch End:

Batch Notes

Manifold ID 1
Methanol ID 769613
Hexane ID 0000146278
Sodium Hypochlorite ID NA
First Start time NA
First End time NA
Balance ID QA-070
SPE Cartridge Type WAX 500mg
Solid Phase Extraction Disk ID 002836112A
H2O ID 11/29/16
Pipette ID MD05306
Solvent Name 0.3% NH4OH/MeOH
Solvent Lot # 758321
Analyst ID - Reagent Drop VPM
Analyst ID - SU Reagent Drop VPM
Analyst ID - SU Reagent Drop Witness ERW
Acid Name NA
Acid ID NA
Reagent ID NA
Reagent Lot Number NA
NaCl ID NA

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139834

Analyst: Marchenko, Veronika P

Batch Open: 11/29/2016 12:27:00PM

Method Code: 320-3535_IVWT-320

Batch End:

SOP Number WS-LC-0025

Batch Comment 0.1N NaOH/H2O: 758321

Comments

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Marchenko, Veronika P

Batch Open: 11/29/2016 12:27:00PM

Batch End:

Batch Number: 320-139834

Method Code: 320-3535_IVWT-320

Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-139834/1	LCMPFCSU_00046	25 uL	0.5 mL	VPM 11-29-16	ERW 11/29/16
LCS 320-139834/2	LCMPFCSU_00046	25 uL	0.5 mL		
LCS 320-139834/2	LCPFCSU_00070	20 uL	0.5 mL		
LCSD 320-139834/3	LCMPFCSU_00046	25 uL	0.5 mL		
LCSD 320-139834/3	LCPFCSU_00070	20 uL	0.5 mL		
320-23770-A-1	LCMPFCSU_00046	25 uL	0.5 mL		
320-23801-A-1	LCMPFCSU_00046	25 uL	0.5 mL		
320-23801-A-2	LCMPFCSU_00046	25 uL	0.5 mL		

Other Reagents:

Reagent	Amount/Units	Lot#:

Preparation Batch Number(s): 139834 Test: PFC-10A
 Earliest Holding Time: 11/29/16

Sample List Tab		1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method		/	✓
All necessary NCMs filed (including holding time)		/	✓
Method/sample/login/QAS checked and correct		/	✓
Worksheet Tab		1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved		NA	mt
Weights in anticipated range and not targeted		/	✓
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and Ci Check)		/	✓
The pH is transcribed correctly in TALS		NA	NA
All additional information transcribed into TALS is correct and raw data is attached		/	✓
Comments are transcribed correctly in TALS		/	✓
Reagents Tab		1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS		/	U
All spike amounts correct and added to necessary samples and QC		/	✓
Batch Information		1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly		/	U
All necessary 'batch information' complete and entered into TALS correctly		/	U

1st Level Reviewer: NSH

Date: 11-29-16

2nd Level Reviewer: [Signature]

Date: 11/29/16

Comments: _____

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Marchenko, Veronika P

Batch Number: 320-139749
 Method Code: 320-3535_IVWT-320

AG 11/30/16

12/2/16

Batch Open: 11/28/2016 4:55:00PM

Batch End: 11/29/2016 2:06:00PM

12/3/16

Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	InitAmt		Rcvd	PHs		Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
			Flt	Amnt		Adj1	Adj2					
1 MB-320-139749/1 N/A	N/A		250 mL					N/A	N/A	N/A		
		25.53 g	0.5 mL									
2 LCS-320-139749/2 N/A	N/A		250 mL					N/A	N/A	N/A		
		25.67 g	0.5 mL									
3 LCS-320-139749/3 N/A	N/A		250 mL					N/A	N/A	N/A	See NCM regarding these samples, contains sediment.	
		25.81 g	0.5 mL									
4 320-23783-A-3 (PFC_IDA_DOD5)	N/A (320-23783-1)	276.82 g	250.2 mL					11/29/16	11_Day_Rush	4	10x PFOA, PFOS	
		26.59 g	0.5 mL									
5 320-23783-A-4 (PFC_IDA_DOD5)	N/A (320-23783-1)	275.75 g	249.6 mL					11/29/16	11_Day_Rush	4	10x PFBS, PFOA, PFOS	
		26.19 g	0.5 mL									
6 320-23783-A-12 (PFC_IDA_DOD5)	N/A (320-23783-1)	275.60 g	249.7 mL					11/29/16	11_Day_Rush	4	100x PFBS, PFOA, PFOS	
		25.93 g	0.5 mL									
7 320-23783-A-13 (PFC_IDA_DOD5)	N/A (320-23783-1)	272.23 g	245.9 mL					11/29/16	11_Day_Rush	4	RT 1x possible PFBS, PFOA 10x PFOS	
		26.37 g	0.5 mL									

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139749

Analyst: Marchenko, Veronika P

Batch Open: 11/28/2016 4:55:00PM

Method Code: 320-3535_IVWT-320

Batch End:

Batch Notes

Manifold ID 1

Methanol ID 769613

Hexane ID 0000146278

Sodium Hypochlorite ID NA

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type WAX 500mg

Solid Phase Extraction Disk ID 002836112A

H2O ID 11/28/16

Pipette ID MD05306

Solvent Name 0.3% NH4OH/MeOH

Solvent Lot # 794501

Analyst ID - Reagent Drop VPM

Analyst ID - SU Reagent Drop VPM

Analyst ID - SU Reagent Drop Witness KMK 11-28-16

Acid Name NA

Acid ID NA

Reagent ID NA

Reagent Lot Number NA

NaCl ID NA

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139749

Analyst: Marchenko, Veronika P

Batch Open: 11/28/2016 4:55:00PM

Method Code: 320-3535_IVWT-320

Batch End:

SOP Number WS-LC-0025

Batch Comment 0.1N NaOH/H2O: 794892

Comments

Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-139749

Analyst: Marchenko, Veronika P

Batch Open: 11/28/2016 4:55:00PM

Method Code: 320-3535_IVWT-320

Batch End:

Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-139749/1	LCMPFCSU_00046	25 uL	0.5 mL	VPM 11/28/16	KMK
LCS 320-139749/2	LCMPFCSU_00046	25 uL	0.5 mL		
LCS 320-139749/2	LCPFCSU_00070	20 uL	0.5 mL		
LCSD 320-139749/3	LCMPFCSU_00046	25 uL	0.5 mL		
LCSD 320-139749/3	LCPFCSU_00070	20 uL	0.5 mL		
320-23783-A-3	LCMPFCSU_00046	25 uL	0.5 mL		
320-23783-A-4	LCMPFCSU_00046	25 uL	0.5 mL		
320-23783-A-12	LCMPFCSU_00046	25 uL	0.5 mL		
320-23783-A-13	LCMPFCSU_00046	25 uL	0.5 mL		

Other Reagents:

Amount/Units

Lot#:

Reagent

Preparation Batch Number(s): 139749 Test: PFC-10A-0005 (L)
 Earliest Holding Time: 11/28/16

Sample List Tab		1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method		/	/
All necessary NCMs filed (including holding time)		/	/
Method/sample/login/QAS checked and correct		/	/
Worksheet Tab		1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved		NA	NA
Weights in anticipated range and not targeted		/	/
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and CI Check)		/	/
The pH is transcribed correctly in TALS		NA	NA
All additional information transcribed into TALS is correct and raw data is attached		/	/
Comments are transcribed correctly in TALS		/	/
Reagents Tab		1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS		/	/
All spike amounts correct and added to necessary samples and QC		/	/
Batch Information		1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly		/	/
All necessary 'batch information' complete and entered into TALS correctly		/	/

1st Level Reviewer: VPM

Date: 11/29/16

2nd Level Reviewer: HSA

Date: 11-29-16

Comments: _____

GENERAL CHEMISTRY

COVER PAGE
GENERAL CHEMISTRY

Lab Name: TestAmerica Sacramento

Job Number: 320-23783-1

SDG No.: _____

Project: Former Bay Head Road Annex (RFP JU06-01)

Client Sample ID	Lab Sample ID
<u>DPT-16-35-SO-00-01</u>	<u>320-23783-1</u>
<u>DPT-16-35-SO-19-20</u>	<u>320-23783-2</u>
<u>DPT-16-19-SO-00-01</u>	<u>320-23783-5</u>
<u>DPT-16-19-SO-17-18</u>	<u>320-23783-6</u>
<u>DPT-16-32-SO-14-15</u>	<u>320-23783-7</u>
<u>DPT-16-31-SO-14-15</u>	<u>320-23783-8</u>
<u>DPT-16-30-SO-14-15</u>	<u>320-23783-9</u>
<u>DPT-16-28-SO-14-15</u>	<u>320-23783-10</u>
<u>DPT-16-15-SO-14-15</u>	<u>320-23783-11</u>

Comments:

9-IN
DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Sacramento

Job Number: 320-23783-1

SDG Number: _____

Matrix: Solid

Instrument ID: NOEQUIP

Method: D 2216

RL Date: 01/01/2012 08:18

Analyte	Wavelength/ Mass	RL (%)	
Percent Moisture		0.1	
Percent Solids		0.1	

9-IN
CALIBRATION BLANK DETECTION LIMITS
GENERAL CHEMISTRY

Lab Name: TestAmerica Sacramento Job Number: 320-23783-1
SDG Number: _____
Matrix: Solid Instrument ID: NOEQUIP
Method: D 2216 XRL Date: 01/01/2012 08:19

Analyte	Wavelength/ Mass	XRL (%)	
Percent Moisture		0.1	
Percent Solids		0.1	

13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: NOEQUIP Analysis Method: D 2216

Start Date: 12/01/2016 09:31 End Date: 12/01/2016 09:31

Lab Sample Id	D/F	Type	Time	Analytes																											
				% S	M o i s t																										
ZZZZZZ			09:31																												
ZZZZZZ			09:31																												
ZZZZZZ			09:31																												
ZZZZZZ			09:31																												
ZZZZZZ			09:31																												
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ZZZZZZ			09:31																												
ZZZZZZ			09:31																												
320-23783-1		1 T	09:31	X	X																										
320-23783-2		1 T	09:31	X	X																										
320-23783-5		1 T	09:31	X	X																										
320-23783-6		1 T	09:31	X	X																										

Prep Types: _____
T = Total/NA

13-IN
ANALYSIS RUN LOG
GENERAL CHEMISTRY

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Instrument ID: NOEQUIP Analysis Method: D 2216

Start Date: 12/01/2016 11:23 End Date: 12/01/2016 11:23

Lab Sample Id	D/F	Type	Time	Analytes																											
				% S	M																										
320-23783-7	1	T	11:23	X	X																										
320-23783-11	1	T	11:23	X	X																										
320-23783-8	1	T	11:23	X	X																										
320-23783-9	1	T	11:23	X	X																										
320-23783-10	1	T	11:23	X	X																										
ZZZZZZ			11:23																												
ZZZZZZ			11:23																												
ZZZZZZ			11:23																												
ZZZZZZ			11:23																												
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ZZZZZZ			11:23																												
ZZZZZZ			11:23																												

Prep Types: _____
T = Total/NA

GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 140129 Batch Start Date: 12/01/16 09:31 Batch Analyst: Martin, Christine

Batch Method: D 2216 Batch End Date: 12/02/16 09:43

Lab Sample ID	Client Sample ID	Method Chain	Basis	DISH#	DishWeight	SampleMassWet	SampleMassDry		
320-23783-A-1	DPT-16-35-SO-00-01	D 2216	T	17	0.99 g	10.88 g	9.54 g		
320-23783-A-2	DPT-16-35-SO-19-20	D 2216	T	18	0.99 g	13.25 g	11.14 g		
320-23783-A-5	DPT-16-19-SO-00-01	D 2216	T	19	1.01 g	12.45 g	10.93 g		
320-23783-A-6	DPT-16-19-SO-17-18	D 2216	T	20	1.00 g	10.78 g	8.79 g		

Batch Notes	
Balance ID	QA-069 No Unit
Date and Time Samples in Desiccator	12/02/2016 @ 09:05
Date and Time Samples out of Desiccator	12/02/2016 @ 09:43
Date samples were placed in the oven	12/01/2016
Oven Temp In	108 Degrees C
Time samples were place in the oven	10:16
Date samples were removed from oven	12/02/2016
Oven Temp Out	108 Degrees C
Time Samples were removed from oven	09:05
Oven ID	Soil Prep #1
Thermometer ID	151969626

Basis	Basis Description
T	Total/NA

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

GENERAL CHEMISTRY BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-23783-1

SDG No.: _____

Batch Number: 140181 Batch Start Date: 12/01/16 11:23 Batch Analyst: Lo, KitSeong S

Batch Method: D 2216 Batch End Date: 12/02/16 09:43

Lab Sample ID	Client Sample ID	Method Chain	Basis	DISH#	DishWeight	SampleMassWet	SampleMassDry		
320-23783-A-7	DPT-16-32-SO-14-15	D 2216	T	1	1.02 g	6.39 g	5.44 g		
320-23783-A-11	DPT-16-15-SO-14-15	D 2216	T	2	1.02 g	6.38 g	5.34 g		
320-23783-A-8	DPT-16-31-SO-14-15	D 2216	T	3	1.06 g	8.15 g	7.00 g		
320-23783-A-9	DPT-16-30-SO-14-15	D 2216	T	4	1.04 g	6.64 g	5.73 g		
320-23783-A-10	DPT-16-28-SO-14-15	D 2216	T	5	1.02 g	7.58 g	6.61 g		

Batch Notes	
Balance ID	QA-068 No Unit
Date and Time Samples in Desiccator	12/02/2016 @ 09:05
Date and Time Samples out of Desiccator	12/02/2016 @ 09:43
Date samples were placed in the oven	12/01/2016
Oven Temp In	107 Degrees C
Time samples were place in the oven	12:00
Date samples were removed from oven	12/02/2016
Oven Temp Out	108 Degrees C
Time Samples were removed from oven	09:05
Oven ID	soil prep #1
Thermometer ID	151969626

Basis	Basis Description
T	Total/NA

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

Preparation Batch Number(s): 140181

Test: 7. MOISTURE

Earliest Holding Time: N/A

Sample List Tab		1 st Level Reviewer	2 nd Level Reviewer
Samples identified to the correct method		/	/
All necessary NCMs filed (including holding time)		/	/
Method/sample/login/QAS checked and correct		/	/
Worksheet Tab		1 st Level Reviewer	2 nd Level Reviewer
All samples properly preserved		/	/
Weights in anticipated range and not targeted		/	/
All additional test requirements performed, documented, and uploaded to TALS correctly (e.g. final amount, initial amount, turbidity, and CI Check)		N/A	/
The pH is transcribed correctly in TALS		N/A	/
All additional information transcribed into TALS is correct and raw data is attached		/	/
Comments are transcribed correctly in TALS		/	/
Reagents Tab		1 st Level Reviewer	2 nd Level Reviewer
All necessary reagents not expired and entered into TALS		N/A	/
All spike amounts correct and added to necessary samples and QC		N/A	/
Batch Information		1 st Level Reviewer	2 nd Level Reviewer
Date and time accurate and entered into TALS correctly		/	/
All necessary 'batch information' complete and entered into TALS correctly		/	/

1st Level Reviewer: UMD

Date: 12/02/16

2nd Level Reviewer: [Signature]

Date: 12/2/16

Comments: _____

Shipping and Receiving Documents

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Temperature on Receipt _____
 Drinking Water? Yes No

Chain of Custody Record

TAL-4124 (1007)

Client: **AECOM** Project Manager: **Kurt VanGelder** Chain of Custody Number: **299019**
 Address: **3101 WILSON BLVD** Telephone Number (Area Code)/Fax Number: **(703) 682-9041** Page **1** of **2**
 City: **ARLINGTON VA** Zip Code: **22201** Site Contact: **MIKE LLINSKI** Lab Contact: **Jill Kellerman**

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix					Containers & Preservatives					Analysis (Attach list if more space is needed)	Special Instructions/ Conditions of Receipt	
			Air	Aqueous	Sed	Soil	Unpres	H2SO4	HNO3	HCl	NaOH	ZnAc/NaOH			
DPT-16-35-50-00-01	11/22/16	0924			✓	✓	1							✓ PFOA	
DPT-16-35-50-19-20		0936			✓	✓	1							✓ PFOA	
DPT-16-35-GW-21-25		1010		✓			2							✓ PFOA	
DPT-16-35-GW-31-35		1030		✓			2							✓ PFOA	
DPT-16-19-50-00-01		1120			✓	✓	1							✓ PFOA	
DPT-16-19-50-17-18		1130			✓	✓	1							✓ PFOA	
DPT-16-32-50-14-15	11/21/16	0936			✓	✓	1							✓ PFOA	
DPT-16-31-50-14-15		1012			✓	✓	1							✓ PFOA	
DPT-16-30-50-14-15		1046			✓	✓	1							✓ PFOA	
DPT-16-28-50-14-15		1154			✓	✓	1							✓ PFOA	
DPT-16-15-50-14-15		1250			✓	✓	1							✓ PFOA	
DPT-16-15-GW-18-22		1330			✓	✓	2							✓ PFOA	



320-23783 Chain of Custody

Possible Hazard Identification: Non-Hazard Flammable Skin Irritant Poison B Unknown Return To Client Disposal By Lab Archive For _____ Months
 Sample Disposal: Return To Client Disposal By Lab Archive For _____ Months
 Turn Around Time Required: 24 Hours 48 Hours 7 Days 14 Days 21 Days Other _____
 (A fee may be assessed if samples are retained longer than 1 month)

Relinquished By	Date	Time	1. Received By	Date	Time
<i>[Signature]</i>	11/22/16	1400	TRANSFER TO FEDEX	11/22/16	1400
			2. Received By	Date	Time
			<i>[Signature]</i> Troy Gr. Turpen THS	11/25/16	09:30
			3. Received By	Date	Time

Comments: **1.5" ice** * Sample labeled states "DPT-16-22-50-14-15" that 11/25/16

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Chain of Custody Record

Temperature on Receipt _____

Drinking Water? Yes No

TAL-4124 (1007)

Client: **AECOM** Project Manager: **Kurt Van Gelder** Chain of Custody Number: **295762**

Address: **3101 WILSON BLVD** Telephone Number (Area Code)/Fax Number: **11/22/16** Page **2** of **2**

City: **ARLINGTON VA** State: **VA** Zip Code: **22201** Site Contact: **MIKE GLINSKI** Lab Contact: **JILL K.**

Project Name and Location (State): **Former 13km Hend Rd Annex** Carrier/Vehicle Number: **FEDEX**

Contract/Purchase Order/Quote No.: **32007449 / 60444465**

Sample I.D. No. and Description (Containers for each sample may be combined on one line): **DPT-16-15-GW-31-35** Date: **11/21/16** Time: **1400**

Containers & Preservatives: **2**

Containers & Preservatives	Air	Aqueous	Sed.	Soil	Unpres.	H2SO4	HNO3	HCl	NaOH	ZnAc/NaOH
		<input checked="" type="checkbox"/>			2					

Analysis (Attach list if more space is needed): **✓ PFAS ✓ PFOS ✓ PFBs**

Special Instructions/Conditions of Receipt:

Sample Disposal: Return To Client Disposal By Lab Archive For _____ Months (A fee may be assessed if samples are retained longer than 1 month)

Turn Around Time Required: 24 Hours 48 Hours 7 Days 14 Days 21 Days Other: _____

1. Relinquished By: **MW Adams / AECOM** Date: **11/22/16** Time: **1400**

2. Relinquished By: **MW Adams / AECOM** Date: **11/22/16** Time: **1400**

3. Relinquished By: **Tracy Gump** Date: **11/25/16** Time: **09:30**

Comments: **1.5°C re**

Login Sample Receipt Checklist

Client: AECOM Technical Services Inc.

Job Number: 320-23783-1

Login Number: 23783
List Number: 1
Creator: Nelson, Kym D

List Source: TestAmerica Sacramento

Question	Answer	Comment
Radioactivity wasn't checked or is <=/ background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	CUSTODY SEALS
Sample custody seals, if present, are intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	N/A	
There are no discrepancies between the containers received and the COC.	False	IDs on containers do not match the COC. Logged in per COC.
Samples are received within Holding Time (excluding tests with immediate HTs)	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	N/A	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	

"DPT-16-15-GW-18-22", "537", "12/03/16", "16:41", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "5.8", "", "TRG", "No", "Y", "EM", "Y", "0.00075", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-18-22", "537", "12/03/16", "16:41", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "1.2", "", "TRG", "No", "Y", "E", "Y", "0.00092", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-18-22", "537", "12/03/16", "16:41", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "6.5", "", "TRG", "No", "Y", "EM", "Y", "0.0013", "0.0030", "0.0040", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-18-22", "537", "12/07/16", "17:03", "T", "NA", "DILUTION1", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "12", "", "TRG", "Yes", "Y", "D", "Y", "0.13", "0.30", "0.40", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-18-22", "537", "12/07/16", "17:03", "T", "NA", "DILUTION1", "335-67-1", "Perfluorooctanoic acid (PFOA)", "15", "", "TRG", "Yes", "Y", "D", "Y", "0.075", "0.20", "0.25", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-18-22", "537", "12/07/16", "17:03", "T", "NA", "DILUTION1", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.54", "", "TRG", "Yes", "Y", "D", "Y", "0.092", "0.20", "0.25", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/03/16", "16:48", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "0.46", "", "TRG", "No", "Y", "E", "Y", "0.0013", "0.0031", "0.0041", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/03/16", "16:48", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.27", "", "TRG", "Yes", "Y", "M", "Y", "0.00076", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/03/16", "16:48", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.048", "", "TRG", "Yes", "Y", "Y", "Y", "0.00093", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/07/16", "17:25", "T", "NA", "DILUTION1", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "0.45", "", "TRG", "Yes", "Y", "D", "Y", "0.013", "0.031", "0.041", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/07/16", "17:25", "T", "NA", "DILUTION1", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.035", "", "TRG", "No", "Y", "D", "Y", "0.0093", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-GW-31-35", "537", "12/07/16", "17:25", "T", "NA", "DILUTION1", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.27", "", "TRG", "No", "Y", "D", "Y", "0.0076", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-SO-14-15", "537", "12/02/16", "17:22", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "11", "", "TRG", "Yes", "Y", "M", "Y", "0.15", "0.37", "0.61", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-SO-14-15", "537", "12/02/16", "17:22", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "", "", "TRG", "Yes", "N", "U", "Y", "0.13", "0.37", "0.49", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-15-SO-14-15", "537", "12/02/16", "17:22", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.51", "", "TRG", "Yes", "Y", "J", "Y", "0.13", "0.37", "0.61", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-19-SO-00-01", "537", "12/02/16", "16:07", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "3.8", "", "TRG", "Yes", "Y", "Y", "Y", "0.12", "0.34", "0.57", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-19-SO-00-01", "537", "12/02/16", "16:07", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "130", "", "TRG", "No", "Y", "ME", "Y", "0.14", "0.34", "0.57", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-19-SO-00-01", "537", "12/02/16", "16:07", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.18", "", "TRG", "Yes", "Y", "J", "Y", "0.12", "0.34", "0.46", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

","","","","","","","","","","","","",""
"DPT-16-19-SO-00-01","537","12/07/16","17:33","T","NA","DILUTION1","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","No","N","U","Y","1.2","3.4","4.6","UG_KG","UG_KG","","","0","","","","0","","","","","",""
","","","","",""
"DPT-16-19-SO-00-01","537","12/07/16","17:33","T","NA","DILUTION1","1763-23-1","Perfluorooctane Sulfonate (PFOS)","170","","TRG","Yes","Y","D","Y","1.4","3.4","5.7","UG_KG","UG_KG","","","0","","","","0","","","","",""
","","","","",""
"DPT-16-19-SO-00-01","537","12/07/16","17:33","T","NA","DILUTION1","335-67-1","Perfluorooctanoic acid (PFOA)","3.5","","TRG","No","Y","JDM","Y","1.2","3.4","5.7","UG_KG","UG_KG","","","0","","","","0","","","",""
","","","","",""
"DPT-16-19-SO-17-18","537","12/02/16","16:44","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","6.5","","TRG","Yes","Y","M","Y","0.16","0.37","0.62","UG_KG","UG_KG","","","0","","","","0","","","",""
","","","","",""
"DPT-16-19-SO-17-18","537","12/02/16","16:44","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.13","0.37","0.50","UG_KG","UG_KG","","","0","","","","0","","","",""
","","","","",""
"DPT-16-19-SO-17-18","537","12/02/16","16:44","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","0.37","","TRG","Yes","Y","J","Y","0.13","0.37","0.62","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-28-SO-14-15","537","12/02/16","17:14","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","0.45","","TRG","Yes","Y","J","Y","0.12","0.35","0.58","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-28-SO-14-15","537","12/02/16","17:14","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.12","0.35","0.47","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-28-SO-14-15","537","12/02/16","17:14","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","10","","TRG","Yes","Y","M","Y","0.15","0.35","0.58","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-30-SO-14-15","537","12/02/16","17:07","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.12","0.36","0.48","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-30-SO-14-15","537","12/02/16","17:07","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","57","","TRG","Yes","Y","M","Y","0.15","0.36","0.59","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-30-SO-14-15","537","12/02/16","17:07","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","5.5","","TRG","Yes","Y","","Y","0.12","0.36","0.59","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-31-SO-14-15","537","12/02/16","16:59","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.12","0.36","0.48","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-31-SO-14-15","537","12/02/16","16:59","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","0.27","","TRG","Yes","Y","J","Y","0.12","0.36","0.60","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-31-SO-14-15","537","12/02/16","16:59","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","11","","TRG","Yes","Y","M","Y","0.15","0.36","0.60","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-32-SO-14-15","537","12/02/16","16:52","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","46","","TRG","Yes","Y","M","Y","0.15","0.36","0.60","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-32-SO-14-15","537","12/02/16","16:52","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.12","0.36","0.48","UG_KG","UG_KG","","","0","","","","0","","",""
","","","","",""
"DPT-16-32-SO-14-15","537","12/02/16","16:52","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","1.1","","TRG","Yes","Y","","Y","0.12","0.36","0.60","UG_KG","UG_KG","","","0","","","","0","","",""

"DPT-16-35-GW-21-25", "537", "12/03/16", "16:26", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.60", "", "TRG", "No", "Y", "EM", "Y", "0.00075", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-21-25", "537", "12/03/16", "16:26", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.28", "", "TRG", "Yes", "Y", "", "Y", "0.00092", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-21-25", "537", "12/03/16", "16:26", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "1.9", "", "TRG", "No", "Y", "EM", "Y", "0.0013", "0.0030", "0.0040", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-21-25", "537", "12/07/16", "17:10", "T", "NA", "DILUTION1", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.19", "", "TRG", "No", "Y", "D", "Y", "0.0092", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-21-25", "537", "12/07/16", "17:10", "T", "NA", "DILUTION1", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.76", "", "TRG", "Yes", "Y", "D", "Y", "0.0075", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-21-25", "537", "12/07/16", "17:10", "T", "NA", "DILUTION1", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "2.4", "", "TRG", "Yes", "Y", "D", "Y", "0.013", "0.030", "0.040", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/03/16", "16:33", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "0.72", "", "TRG", "No", "Y", "E", "Y", "0.0013", "0.0030", "0.0040", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/03/16", "16:33", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "1.1", "", "TRG", "No", "Y", "EM", "Y", "0.00075", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/03/16", "16:33", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.58", "", "TRG", "No", "Y", "E", "Y", "0.00092", "0.0020", "0.0025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/07/16", "17:18", "T", "NA", "DILUTION1", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.38", "", "TRG", "Yes", "Y", "D", "Y", "0.0092", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/07/16", "17:18", "T", "NA", "DILUTION1", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "0.80", "", "TRG", "Yes", "Y", "D", "Y", "0.013", "0.030", "0.040", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-GW-31-35", "537", "12/07/16", "17:18", "T", "NA", "DILUTION1", "335-67-1", "Perfluorooctanoic acid (PFOA)", "1.4", "", "TRG", "Yes", "Y", "D", "Y", "0.0075", "0.020", "0.025", "UG_L", "UG_L", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-00-01", "537", "12/02/16", "15:52", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "0.21", "", "TRG", "Yes", "Y", "J", "Y", "0.12", "0.35", "0.46", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-00-01", "537", "12/02/16", "15:52", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "28", "", "TRG", "Yes", "Y", "", "Y", "0.15", "0.35", "0.58", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-00-01", "537", "12/02/16", "15:52", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "12", "", "TRG", "Yes", "Y", "M", "Y", "0.12", "0.35", "0.58", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-19-20", "537", "12/02/16", "15:59", "T", "NA", "Initial", "375-73-5", "Perfluorobutanesulfonic acid (PFBS)", "", "", "TRG", "Yes", "N", "U", "Y", "0.12", "0.36", "0.48", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-19-20", "537", "12/02/16", "15:59", "T", "NA", "Initial", "1763-23-1", "Perfluorooctane Sulfonate (PFOS)", "4.0", "", "TRG", "Yes", "Y", "M", "Y", "0.15", "0.36", "0.60", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

"DPT-16-35-SO-19-20", "537", "12/02/16", "15:59", "T", "NA", "Initial", "335-67-1", "Perfluorooctanoic acid (PFOA)", "0.56", "", "TRG", "Yes", "Y", "J", "Y", "0.12", "0.36", "0.60", "UG_KG", "UG_KG", "", "", "0", "", "", "", "0", "", "", ""

","","","","","","","","","","",""
"LCS 320-139627/2-A","537","12/02/16","15:07","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","4.03","","SC","Yes","Y","","Y","0.10","0.30","0.40","UG_KG","UG_KG","","","0","3.54","4.03","114","0",
","","","","50","150",
"LCS 320-139627/2-A","537","12/02/16","15:07","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","3.66","","SC","Yes","Y","","Y","0.13","0.30","0.50","UG_KG","UG_KG","","","0","3.71","3.66","99","0",
","","","60","140",
"LCS 320-139627/2-A","537","12/02/16","15:07","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","4.25","","SC","Yes","Y","","Y","0.10","0.30","0.50","UG_KG","UG_KG","","","0","4.00","4.25","106","0",
","","","60","140",
"LCS 320-139749/2-A","537","12/03/16","16:11","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","0.0289","","SC","Yes","Y","","Y","0.00092","0.0020","0.0025","UG_L","UG_L","","","0","0.0354","0.0289",
","82","0","50","150",
"LCS 320-139749/2-A","537","12/03/16","16:11","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","0.0261","","SC","Yes","Y","","Y","0.0013","0.0030","0.0040","UG_L","UG_L","","","0","0.0371","0.0261",
","70","0","60","140",
"LCS 320-139749/2-A","537","12/03/16","16:11","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","0.0296","","SC","Yes","Y","","Y","0.00075","0.0020","0.0025","UG_L","UG_L","","","0","0.0400","0.0296",
","74","0","60","140",
"LCSD 320-139749/3-A","537","12/03/16","16:18","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","0.0325","","SC","Yes","Y","","Y","0.00092","0.0020","0.0025","UG_L","UG_L","","","0","0.0354","0.0325",
","92","12","50","150","30",
"LCSD 320-139749/3-A","537","12/03/16","16:18","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","0.0289","","SC","Yes","Y","","Y","0.0013","0.0030","0.0040","UG_L","UG_L","","","0","0.0371",
","0.0289","78","10","60","140","30",
"LCSD 320-139749/3-A","537","12/03/16","16:18","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","0.0314","","SC","Yes","Y","","Y","0.00075","0.0020","0.0025","UG_L","UG_L","","","0","0.0400",
","0.0314","79","6","60","140","30",
"MB 320-139627/1-A","537","12/02/16","14:59","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.10","0.30","0.40","UG_KG","UG_KG","","","0",
","","","
"MB 320-139627/1-A","537","12/02/16","14:59","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","","","TRG","Yes","N","U","Y","0.13","0.30","0.50","UG_KG","UG_KG","","","0",
","","","
"MB 320-139627/1-A","537","12/02/16","14:59","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","","","TRG","Yes","N","U","Y","0.10","0.30","0.50","UG_KG","UG_KG","","","0",
","","","
"MB 320-139749/1-A","537","12/03/16","16:03","T","NA","Initial","375-73-5","Perfluorobutanesulfonic acid (PFBS)","","","TRG","Yes","N","U","Y","0.00092","0.0020","0.0025","UG_L","UG_L","","","0",
","","","
"MB 320-139749/1-A","537","12/03/16","16:03","T","NA","Initial","1763-23-1","Perfluorooctane Sulfonate (PFOS)","","","TRG","Yes","N","U","Y","0.0013","0.0030","0.0040","UG_L","UG_L","","","0",
","","","
"MB 320-139749/1-A","537","12/03/16","16:03","T","NA","Initial","335-67-1","Perfluorooctanoic acid (PFOA)","","","TRG","Yes","N","U","Y","0.00075","0.0020","0.0025","UG_L","UG_L","","","0",
","","","



Data Validation Report

Project:	Former Bay Head Road Annex- Annapolis, MD	
Laboratory:	TestAmerica-West Sacramento, CA	
Job Number:	320-23783-1	
Analyses/Method:	Perfluorinated Compounds (PFCs) in Water, Soils, Sediments and Tissues by Liquid Chromatography/Mass Spectrometry/Mass Spectrometry (LC/MS/MS)/ Revision 1.4 (August 2015)	
Validation Level:	Limited	
Resolution Consultants	60444465-DM.DE	
Project Number:		
Prepared by:	Paula DiMattei/Resolution Consultants	Completed on: 12/13/16
Reviewed by:	Robert Kennedy/Resolution Consultants	Completed on: 12/13/16
File Name:	J23783-1_PFC memo.docx	

SUMMARY

The samples listed below were collected by Resolution Consultants from the Former Bay Head Road Annex site in Annapolis, MD on November 21-22, 2016.

Sample ID	Matrix/Sample Type
DPT-16-15-GW-18-22	Groundwater
DPT-16-15-GW-31-35	Groundwater
DPT-16-15-SO-14-15	Soil
DPT-16-19-SO-00-01	Soil
DPT-16-19-SO-17-18	Soil
DPT-16-28-SO-14-15	Soil
DPT-16-30-SO-14-15	Soil
DPT-16-31-SO-14-15	Soil
DPT-16-32-SO-14-15	Soil
DPT-16-35-GW-21-25	Groundwater
DPT-16-35-GW-31-35	Groundwater
DPT-16-35-SO-00-01	Soil
DPT-16-35-SO-19-20	Soil

Data validation activities were conducted with reference to:

- TestAmerica-West Sacramento SOP: Perfluorinated Compounds (PFCs) in Water, Soils, Sediments and Tissues by Liquid Chromatography/Mass Spectrometry/Mass Spectrometry (LC/MS/MS)/Revision 1.4 (August 2015);

- USEPA Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review (September 2016);
- USEPA Contract Laboratory Program National Functional Guidelines for High Resolution Superfund Methods Data Review (April 2016)
- Quality Systems Manual (QSM) for Environmental Laboratories, Version 5.0 (DoD, July 2013); and
- the project-specific Sampling and Analysis Plan.

In the absence of method-specific information, laboratory quality control (QC) limits, project-specific requirements and/or professional judgment were used as appropriate.

REVIEW ELEMENTS

The data were evaluated based on the following review elements (where applicable to the method):

- X Data completeness (chain-of-custody (COC)/sample integrity)
- ✓ Holding times/sample preservation
- ✓ Initial calibration/initial and continuing calibration verification
- ✓ Laboratory method blanks/equipment blanks
- NA Matrix spike (MS) and/or matrix spike duplicate (MSD) results
- ✓ Laboratory control sample (LCS)/laboratory control sample duplicate (LCSD) results
- NA Field duplicate results
- X Labeled compound results
- ✓ Sample results/reporting issues

The symbol (✓) indicates that no validation qualifiers were applied based on this parameter. An “NA” indicates that the parameter was not included as part of this data set or was not applicable to this validation and therefore not reviewed. The symbol (X) indicates that a QC nonconformance resulted in the qualification of data. Any QC nonconformance that resulted in the qualification of data is discussed below. In addition, nonconformances or other issues that were noted during validation, but did not result in qualification of data, may be discussed for informational purposes only.

The data appear valid as reported and may be used for decision making purposes. Select data points were qualified as estimated due to nonconformances of certain QC criteria (see discussion below). Qualified sample results are presented in Table 1.

RESULTS

Data Completeness (chain-of-custody (COC)/Sample Integrity)

The data package was reviewed and found to meet acceptance criteria for completeness:

- The COCs were reviewed for completeness of information relevant to the samples and requested analyses, and for signatures indicating transfer of sample custody.
- The laboratory sample login sheet(s) were reviewed for issues potentially affecting sample integrity, including the condition of sample containers upon receipt at the laboratory.
- Completeness of analyses was verified by comparing the reported results to the COC requests.

The laboratory noted that the sample container label for sample 320-23783-10 identified the sample as DPT-16-22-SO-14-15 while the COC identified this sample as DPT-16-28-SO-14-15. The sample ID listed on the COC was used to identify this sample.

The laboratory noted in the case narrative that all groundwater samples were decanted to new bottles prior to spiking and extraction because of the presence of excessive amounts of sediment present in the sample bottles. In these cases, the sample bottles are not rinsed as required by the method. Consequently, professional judgment was applied to qualify the positive results for all target compounds in these samples as estimated (J-) indicating a potential loss of target compounds that may have remained in the original sample bottle. Qualified sample results are presented in Table 1.

Holding Times/Sample Preservation

Sample preservation and preparation/analysis holding times were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Initial Calibration/Initial and Continuing Calibration Verification

Calibration data were reviewed for conformance with the QC acceptance criteria to ensure that:

- the initial calibration (ICAL) percent relative standard deviation (%RSD) or correlation coefficient (r) or coefficient of determination (r^2) method acceptance criteria were met;
- the initial calibration verification standard (ICV) percent recovery acceptance criteria were met; and
- the continuing calibration verification standard (CCV) frequency and method percent difference or percent drift (%D) criteria were met.

All QC acceptance criteria were met or qualification of the data was not required.

Laboratory Method Blanks/Equipment Blanks

Laboratory method blanks and equipment blanks are evaluated as to whether there are contaminants detected above the detection limit (DL). Target compounds were not detected in the laboratory method blanks associated with the samples in this data set. An equipment blank was not submitted with the samples in this data set.

MS/MSD Results

MS/MSD analyses were not performed on a sample from this data set. The data were not qualified on this basis.

LCS/LCSD Results

The LCS/LCSD percent recoveries (%Rs) and relative percent differences (RPDs) were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met.

Field Duplicate Results

A field duplicate pair was not submitted with this data set. Data were not qualified on this basis.

Labeled Compound Results

The labeled compound results were reviewed for conformance with the QC acceptance criteria. All QC acceptance criteria were met except for the labeled compound results summarized below.

Sample ID	Labeled Compound	% Recovery	QC Limits
DPT-16-15-GW-31-35	13C ₄ -PFOA	22	25-150

Samples were qualified as follows:

Actions: (Based on NFG 2011)

Criteria	Actions	
	Detected	Not detected
%R > Upper Acceptance Limit	J	UJ
%R >10% but < Lower Acceptance Limit	J	UJ
%R <10%	See below	
<10% and S/N >10:1	J	R
<10% and S/N <10:1	R	R

Qualified sample results are presented in Table 1.

Sample Results/Reporting Issues

If applicable, compounds detected at concentrations less than the LOQ but greater than the DL are qualified by the laboratory as estimated (J). This "J" qualifier is retained during data validation.

QUALIFICATION ACTIONS

Sample results qualified as a result of validation actions are summarized in Table 1. All actions are described above.

ATTACHMENTS

Attachment A: Qualifier Codes and Explanations

Attachment B: Reason Codes and Explanations

Table 1 - Data Validation Summary of Qualified Data

Sample ID	Matrix	Compound	Result	LOD	LOQ	Units	Validation Qualifiers	Validation Reason
DPT-16-15-GW-18-22	WG	Perfluorooctanesulfonic Acid (PFOS)	12	0.30	0.40	µg/L	J-	si
DPT-16-15-GW-18-22	WG	Perfluorooctanoic Acid (PFOA)	15	0.20	0.25	µg/L	J-	si
DPT-16-15-GW-18-22	WG	Perfluorobutanesulfonic Acid (PFBS)	0.54	0.20	0.25	µg/L	J-	si
DPT-16-15-GW-31-35	WG	Perfluorooctanoic Acid (PFOA)	0.27	0.0020	0.0025	µg/L	J	lc,si
DPT-16-15-GW-31-35	WG	Perfluorobutanesulfonic Acid (PFBS)	0.048	0.0020	0.0025	µg/L	J-	si
DPT-16-15-GW-31-35	WG	Perfluorooctanesulfonic Acid (PFOS)	0.45	0.031	0.041	µg/L	J-	si
DPT-16-35-GW-21-25	WG	Perfluorobutanesulfonic Acid (PFBS)	0.28	0.0020	0.0025	µg/L	J-	si
DPT-16-35-GW-21-25	WG	Perfluorooctanesulfonic Acid (PFOS)	2.4	0.030	0.040	µg/L	J-	si
DPT-16-35-GW-21-25	WG	Perfluorooctanoic Acid (PFOA)	0.76	0.020	0.025	µg/L	J-	si
DPT-16-35-GW-31-35	WG	Perfluorooctanesulfonic Acid (PFOS)	0.80	0.030	0.040	µg/L	J-	si
DPT-16-35-GW-31-35	WG	Perfluorooctanoic Acid (PFOA)	1.4	0.020	0.025	µg/L	J-	si
DPT-16-35-GW-31-35	WG	Perfluorobutanesulfonic Acid (PFBS)	0.38	0.020	0.025	µg/L	J-	si

Attachment A

Qualifier Codes and Explanations

Qualifier	Explanation
J	The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
J+	The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample and is potentially biased high.
J-	The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample and is potentially biased low.
JN	The analyte was tentatively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
UJ	The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
R	The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Attachment B

Reason Codes and Explanations

Reason Code	Explanation
be	Equipment blank contamination
bf	Field blank contamination
bl	Laboratory blank contamination
c	Calibration issue
d	Reporting limit raised due to chromatographic interference
fd	Field duplicate RPDs
h	Holding times
i	Internal standard areas
k	Estimated Maximum Possible Concentration (EMPC)
l	LCS or OPR recoveries
lc	Labeled compound recovery
ld	Laboratory duplicate RPDs
lp	Laboratory control sample/laboratory control sample duplicate RPDs
m	Matrix spike recovery
md	Matrix spike/matrix spike duplicate RPDs
nb	Negative laboratory blank contamination
p	Chemical preservation issue
r	Dual column RPD
q	Quantitation issue
s	Surrogate recovery
si	Sample integrity issue
su	Ion suppression
t	Temperature preservation issue
x	Percent solids
y	Serial dilution results
z	ICS results

INSTALLATION_ID	SITE_NAME	LOCATION_NAME	LOCATION_TYPE_DESC	COORD_X	COORD_Y	SAMPLE_NAME	SAMPLE_MATRIX_DESC	COLLECT_DATE	ANALYTICAL_METHOD_GRP_DESC	SDG
DAVID_TAYLOR_RC	SITE 00003	DPT-16-15	Direct Push/Geoprobe	1474570	496367	DPT-16-15-GW-18-22	Ground water	11/21/2016	Perfluoroalkyl Compounds	320-23783-1
DAVID_TAYLOR_RC	SITE 00003	DPT-16-15	Direct Push/Geoprobe	1474570	496367	DPT-16-15-GW-31-35	Ground water	11/21/2016	Perfluoroalkyl Compounds	320-23783-1
DAVID_TAYLOR_RC	SITE 00003	DPT-16-35	Direct Push/Geoprobe	1474570	496312	DPT-16-35-GW-21-25	Ground water	11/22/2016	Perfluoroalkyl Compounds	320-23783-1
DAVID_TAYLOR_RC	SITE 00003	DPT-16-35	Direct Push/Geoprobe	1474570	496312	DPT-16-35-GW-31-35	Ground water	11/22/2016	Perfluoroalkyl Compounds	320-23783-1