



**Off-Base Drinking Water Sample Results,  
Level 2 Laboratory Report, Level 4 Laboratory Report,  
Electronic Data Deliverable, Data Validation Report,  
and the Sample Location Figure, SDG J17154-1**

*Naval Air Station Oceana  
Virginia Beach, Virginia*

July 2019

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

## ANALYTICAL REPORT

TestAmerica Laboratories, Inc.  
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TestAmerica Job ID: 320-17154-1  
TestAmerica SDG: CTO WE7G PFC Sampling  
Client Project/Site: CTO WE7G PFC Sampling  
Revision: 1

For:  
CH2M Hill, Inc.  
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Virginia Beach, Virginia 23462

Attn: Laurie George



Authorized for release by:  
3/22/2016 12:42:28 PM

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### LINKS

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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: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Qualifiers

### LCMS

Qualifier	Qualifier Description
U	Undetected at the Limit of Detection.
J	Estimated: The analyte was positively identified; the quantitation is an estimation

## 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: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Job ID: 320-17154-1**

**Laboratory: TestAmerica Sacramento**

**Narrative**

## CASE NARRATIVE

**Client: CH2M Hill, Inc.**

**Project: CTO WE7G PFC Sampling**

**Report Number: 320-17154-1**

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.

### **Revision**

This report was revised March 22, 2016 to include all method blank data for all analytes and to remove the B qualifiers from the LCS/LCSD in batch 100093. No other data has changed as a result of this revision.

### **RECEIPT**

The samples were received on 02/04/2016; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 0.3 C.

### **PFC**

Samples OF-FB09-0216 (320-17154-1), OF-FB37-0216 (320-17154-3), OF-FB11-0216 (320-17154-6), OF-FB28-0216 (320-17154-8) and OF-FB67-0216 (320-17154-10) were analyzed for PFC in accordance with PFC. The samples were prepared on 02/09/2016 and 02/10/2016 and analyzed on 02/20/2016.

# Case Narrative

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

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## Job ID: 320-17154-1 (Continued)

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### Laboratory: TestAmerica Sacramento (Continued)

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

#### Organic Prep

Method(s) 3535: Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with batch 320-100093

Method(s) 3535: Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with batch 320-100147.

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

#### Subcontract Work

**PFC:** This method was subcontracted to Maxxam Analytics Inc.. The subcontract laboratory certification is different from that of the facility issuing the final report. Any analytical or quality issues are noted in the subcontract portion of the report.

The DL/LOD/LOQ limits for Maxxam, along with information on the data qualifiers, are included in the narrative (Page 8 of 143 of the subcontract report; page 494 of 632 of the entire report).



## Detection Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB09-0216**

**Lab Sample ID: 320-17154-1**

No Detections.

**Client Sample ID: OF-RW09-0216**

**Lab Sample ID: 320-17154-2**

No Detections.

**Client Sample ID: OF-FB37-0216**

**Lab Sample ID: 320-17154-3**

No Detections.

**Client Sample ID: OF-RW37-0216**

**Lab Sample ID: 320-17154-4**

No Detections.

**Client Sample ID: OF-RW11-0216**

**Lab Sample ID: 320-17154-5**

No Detections.

**Client Sample ID: OF-FB11-0216**

**Lab Sample ID: 320-17154-6**

No Detections.

**Client Sample ID: OF-RW28-0216**

**Lab Sample ID: 320-17154-7**

No Detections.

**Client Sample ID: OF-FB28-0216**

**Lab Sample ID: 320-17154-8**

No Detections.

**Client Sample ID: OF-RW67-0216**

**Lab Sample ID: 320-17154-9**

No Detections.

**Client Sample ID: OF-FB67-0216**

**Lab Sample ID: 320-17154-10**

No Detections.

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

# Client Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB09-0216**

**Lab Sample ID: 320-17154-1**

Date Collected: 02/03/16 16:40

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00079	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00074	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00091	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00086	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0013	ug/L		02/09/16 10:23	02/20/16 10:16	1

Isotope Dilution	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
13C4-PFHpA	122		25 - 150	02/09/16 10:23	02/20/16 10:16	1
13C4 PFOA	126		25 - 150	02/09/16 10:23	02/20/16 10:16	1
13C5 PFNA	127		25 - 150	02/09/16 10:23	02/20/16 10:16	1
18O2 PFHxS	113		25 - 150	02/09/16 10:23	02/20/16 10:16	1
13C4 PFOS	107		25 - 150	02/09/16 10:23	02/20/16 10:16	1

**Client Sample ID: OF-FB37-0216**

**Lab Sample ID: 320-17154-3**

Date Collected: 02/03/16 17:20

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.00073	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.00068	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.00059	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.00083	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.00079	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0012	ug/L		02/09/16 10:23	02/20/16 10:37	1

Isotope Dilution	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
13C4-PFHpA	107		25 - 150	02/09/16 10:23	02/20/16 10:37	1
13C4 PFOA	111		25 - 150	02/09/16 10:23	02/20/16 10:37	1
13C5 PFNA	113		25 - 150	02/09/16 10:23	02/20/16 10:37	1
18O2 PFHxS	100		25 - 150	02/09/16 10:23	02/20/16 10:37	1
13C4 PFOS	95		25 - 150	02/09/16 10:23	02/20/16 10:37	1

**Client Sample ID: OF-FB11-0216**

**Lab Sample ID: 320-17154-6**

Date Collected: 02/03/16 16:10

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.00073	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.00068	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.00059	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.00083	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.00079	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0012	ug/L		02/09/16 10:23	02/20/16 10:58	1

Isotope Dilution	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
13C4-PFHpA	111		25 - 150	02/09/16 10:23	02/20/16 10:58	1
13C4 PFOA	113		25 - 150	02/09/16 10:23	02/20/16 10:58	1
13C5 PFNA	116		25 - 150	02/09/16 10:23	02/20/16 10:58	1
18O2 PFHxS	95		25 - 150	02/09/16 10:23	02/20/16 10:58	1

TestAmerica Sacramento

# Client Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB11-0216**

**Date Collected: 02/03/16 16:10**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-6**

**Matrix: Water**

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons (Continued)**

<i>Isotope Dilution</i>	<i>%Recovery</i>	<i>Qualifier</i>	<i>Limits</i>	<i>Prepared</i>	<i>Analyzed</i>	<i>Dil Fac</i>
13C4 PFOS	93		25 - 150	02/09/16 10:23	02/20/16 10:58	1

**Client Sample ID: OF-FB28-0216**

**Date Collected: 02/03/16 17:43**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-8**

**Matrix: Water**

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0024	0.00077	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorooctanoic acid (PFOA)	0.0019	U	0.0024	0.00072	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorononanoic acid (PFNA)	0.0019	U	0.0024	0.00063	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0024	0.00088	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0024	0.00083	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorooctanesulfonic acid (PFOS)	0.0029	U	0.0038	0.0012	ug/L		02/10/16 07:14	02/20/16 12:23	1

<i>Isotope Dilution</i>	<i>%Recovery</i>	<i>Qualifier</i>	<i>Limits</i>	<i>Prepared</i>	<i>Analyzed</i>	<i>Dil Fac</i>
13C4-PFHpA	124		25 - 150	02/10/16 07:14	02/20/16 12:23	1
13C4 PFOA	132		25 - 150	02/10/16 07:14	02/20/16 12:23	1
13C5 PFNA	137		25 - 150	02/10/16 07:14	02/20/16 12:23	1
18O2 PFHxS	113		25 - 150	02/10/16 07:14	02/20/16 12:23	1
13C4 PFOS	108		25 - 150	02/10/16 07:14	02/20/16 12:23	1

**Client Sample ID: OF-FB67-0216**

**Date Collected: 02/03/16 18:21**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-10**

**Matrix: Water**

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0023	0.00075	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorooctanoic acid (PFOA)	0.0019	U	0.0023	0.00070	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorononanoic acid (PFNA)	0.0019	U	0.0023	0.00061	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0023	0.00086	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0023	0.00081	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorooctanesulfonic acid (PFOS)	0.0028	U	0.0037	0.0012	ug/L		02/10/16 07:14	02/20/16 12:45	1

<i>Isotope Dilution</i>	<i>%Recovery</i>	<i>Qualifier</i>	<i>Limits</i>	<i>Prepared</i>	<i>Analyzed</i>	<i>Dil Fac</i>
13C4-PFHpA	126		25 - 150	02/10/16 07:14	02/20/16 12:45	1
13C4 PFOA	131		25 - 150	02/10/16 07:14	02/20/16 12:45	1
13C5 PFNA	130		25 - 150	02/10/16 07:14	02/20/16 12:45	1
18O2 PFHxS	113		25 - 150	02/10/16 07:14	02/20/16 12:45	1
13C4 PFOS	112		25 - 150	02/10/16 07:14	02/20/16 12:45	1

# Isotope Dilution Summary

Client: CH2M Hill, Inc.  
 Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
 SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - 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)				
		<sup>13</sup> C4-PFHp (25-150)	<sup>13</sup> C4 PFO (25-150)	<sup>13</sup> C5 PFNA (25-150)	<sup>18</sup> O2 PFHx (25-150)	<sup>13</sup> C4 PFOS (25-150)
320-17154-1	OF-FB09-0216	122	126	127	113	107
320-17154-3	OF-FB37-0216	107	111	113	100	95
320-17154-6	OF-FB11-0216	111	113	116	95	93
320-17154-8	OF-FB28-0216	124	132	137	113	108
320-17154-10	OF-FB67-0216	126	131	130	113	112
LCS 320-100093/2-A	Lab Control Sample	95	95	98	90	82
LCS 320-100147/2-A	Lab Control Sample	124	119	117	108	99
LCSD 320-100093/3-A	Lab Control Sample Dup	95	88	91	86	80
LCSD 320-100147/3-A	Lab Control Sample Dup	123	115	116	107	100
MB 320-100093/1-A	Method Blank				95	
MB 320-100093/1-A	Method Blank	103	101	103		92
MB 320-100147/1-A	Method Blank	118	114	118	97	98

### Surrogate Legend

- <sup>13</sup>C4-PFHpA = <sup>13</sup>C4-PFHpA
- <sup>13</sup>C4 PFOA = <sup>13</sup>C4 PFOA
- <sup>13</sup>C5 PFNA = <sup>13</sup>C5 PFNA
- <sup>18</sup>O2 PFHxS = <sup>18</sup>O2 PFHxS
- <sup>13</sup>C4 PFOS = <sup>13</sup>C4 PFOS

# QC Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons

**Lab Sample ID: MB 320-100093/1-A**  
**Matrix: Water**  
**Analysis Batch: 100305**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	MB Result	MB Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00087	ug/L		02/09/16 10:23	02/13/16 00:22	1
Isotope Dilution		MB %Recovery	MB Qualifier	Limits			Prepared	Analyzed	Dil Fac
18O2 PFHxS		95		25 - 150			02/09/16 10:23	02/13/16 00:22	1

**Lab Sample ID: MB 320-100093/1-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	MB Result	MB Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00080	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorooctanesulfonic acid (PFOS)	0.00143	J	0.0040	0.0013	ug/L		02/09/16 10:23	02/20/16 06:23	1
Isotope Dilution		MB %Recovery	MB Qualifier	Limits			Prepared	Analyzed	Dil Fac
13C4-PFHpA		103		25 - 150			02/09/16 10:23	02/20/16 06:23	1
13C4 PFOA		101		25 - 150			02/09/16 10:23	02/20/16 06:23	1
13C5 PFNA		103		25 - 150			02/09/16 10:23	02/20/16 06:23	1
13C4 PFOS		92		25 - 150			02/09/16 10:23	02/20/16 06:23	1

**Lab Sample ID: LCS 320-100093/2-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	Limits
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0513		ug/L		128	60 - 140
Perfluorooctanoic acid (PFOA)	0.0400	0.0485		ug/L		121	60 - 140
Perfluorononanoic acid (PFNA)	0.0400	0.0537		ug/L		134	60 - 140
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0459		ug/L		130	50 - 150
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0480		ug/L		127	60 - 140
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0515		ug/L		135	60 - 140
Isotope Dilution		LCS %Recovery	LCS Qualifier	Limits			
13C4-PFHpA		95		25 - 150			
13C4 PFOA		95		25 - 150			
13C5 PFNA		98		25 - 150			
18O2 PFHxS		90		25 - 150			
13C4 PFOS		82		25 - 150			

# QC Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons (Continued)

**Lab Sample ID: LCSD 320-100093/3-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	Spike Added	LCSD Result	LCSD Qualifier	Unit	D	%Rec	Limits	RPD	RPD Limit
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0506		ug/L		126	60 - 140	1	30
Perfluorooctanoic acid (PFOA)	0.0400	0.0493		ug/L		123	60 - 140	2	30
Perfluorononanoic acid (PFNA)	0.0400	0.0517		ug/L		129	60 - 140	4	30
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0472		ug/L		133	50 - 150	3	30
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0486		ug/L		128	60 - 140	1	30
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0452		ug/L		118	60 - 140	13	30

Isotope Dilution	LCSD %Recovery	LCSD Qualifier	Limits
13C4-PFHpa	95		25 - 150
13C4 PFOA	88		25 - 150
13C5 PFNA	91		25 - 150
18O2 PFHxS	86		25 - 150
13C4 PFOS	80		25 - 150

**Lab Sample ID: MB 320-100147/1-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

Analyte	MB Result	MB Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00080	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00087	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0013	ug/L		02/10/16 07:14	02/20/16 11:20	1

Isotope Dilution	MB %Recovery	MB Qualifier	Limits	Prepared	Analyzed	Dil Fac
13C4-PFHpa	118		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C4 PFOA	114		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C5 PFNA	118		25 - 150	02/10/16 07:14	02/20/16 11:20	1
18O2 PFHxS	97		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C4 PFOS	98		25 - 150	02/10/16 07:14	02/20/16 11:20	1

**Lab Sample ID: LCS 320-100147/2-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	%Rec	Limits
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0498		ug/L		124	60 - 140
Perfluorooctanoic acid (PFOA)	0.0400	0.0476		ug/L		119	60 - 140
Perfluorononanoic acid (PFNA)	0.0400	0.0528		ug/L		132	60 - 140
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0412		ug/L		116	50 - 150
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0503		ug/L		133	60 - 140
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0529		ug/L		138	60 - 140

TestAmerica Sacramento



# QC Sample Results

Client: CH2M Hill, Inc.  
 Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
 SDG: CTO WE7G PFC Sampling

<i>Isotope Dilution</i>	<i>LCS %Recovery</i>	<i>LCS Qualifier</i>	<i>Limits</i>
13C4-PFHpA	124		25 - 150
13C4 PFOA	119		25 - 150
13C5 PFNA	117		25 - 150
18O2 PFHxS	108		25 - 150
13C4 PFOS	99		25 - 150

**Lab Sample ID: LCSD 320-100147/3-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

<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>
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0461		ug/L		115	60 - 140	8	30
Perfluorooctanoic acid (PFOA)	0.0400	0.0488		ug/L		122	60 - 140	2	30
Perfluorononanoic acid (PFNA)	0.0400	0.0534		ug/L		133	60 - 140	1	30
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0461		ug/L		130	50 - 150	11	30
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0473		ug/L		125	60 - 140	6	30
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0502		ug/L		131	60 - 140	5	30

<i>Isotope Dilution</i>	<i>LCSD %Recovery</i>	<i>LCSD Qualifier</i>	<i>Limits</i>
13C4-PFHpA	123		25 - 150
13C4 PFOA	115		25 - 150
13C5 PFNA	116		25 - 150
18O2 PFHxS	107		25 - 150
13C4 PFOS	100		25 - 150

# QC Association Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## LCMS

### Prep Batch: 100093

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-1	OF-FB09-0216	Total/NA	Water	3535	
320-17154-3	OF-FB37-0216	Total/NA	Water	3535	
320-17154-6	OF-FB11-0216	Total/NA	Water	3535	
LCS 320-100093/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-100093/3-A	Lab Control Sample Dup	Total/NA	Water	3535	
MB 320-100093/1-A	Method Blank	Total/NA	Water	3535	

### Prep Batch: 100147

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-8	OF-FB28-0216	Total/NA	Water	3535	
320-17154-10	OF-FB67-0216	Total/NA	Water	3535	
LCS 320-100147/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-100147/3-A	Lab Control Sample Dup	Total/NA	Water	3535	
MB 320-100147/1-A	Method Blank	Total/NA	Water	3535	

### Analysis Batch: 100305

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
MB 320-100093/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100093

### Analysis Batch: 100906

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-1	OF-FB09-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-3	OF-FB37-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-6	OF-FB11-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-8	OF-FB28-0216	Total/NA	Water	WS-LC-0025	100147
320-17154-10	OF-FB67-0216	Total/NA	Water	WS-LC-0025	100147
LCS 320-100093/2-A	Lab Control Sample	Total/NA	Water	WS-LC-0025	100093
LCS 320-100147/2-A	Lab Control Sample	Total/NA	Water	WS-LC-0025	100147
LCSD 320-100093/3-A	Lab Control Sample Dup	Total/NA	Water	WS-LC-0025	100093
LCSD 320-100147/3-A	Lab Control Sample Dup	Total/NA	Water	WS-LC-0025	100147
MB 320-100093/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100093
MB 320-100147/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100147

# Lab Chronicle

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB09-0216**

**Date Collected: 02/03/16 16:40**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-1**

**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			504.9 mL	1.00 mL	100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	504.9 mL	1.00 mL	100906	02/20/16 10:16	JRB	TAL SAC

**Client Sample ID: OF-FB37-0216**

**Date Collected: 02/03/16 17:20**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-3**

**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			550.7 mL	1.00 mL	100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	550.7 mL	1.00 mL	100906	02/20/16 10:37	JRB	TAL SAC

**Client Sample ID: OF-FB11-0216**

**Date Collected: 02/03/16 16:10**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-6**

**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			551.4 mL	1.00 mL	100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	551.4 mL	1.00 mL	100906	02/20/16 10:58	JRB	TAL SAC

**Client Sample ID: OF-FB28-0216**

**Date Collected: 02/03/16 17:43**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-8**

**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			521.6 mL	1.00 mL	100147	02/10/16 07:14	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	521.6 mL	1.00 mL	100906	02/20/16 12:23	JRB	TAL SAC

**Client Sample ID: OF-FB67-0216**

**Date Collected: 02/03/16 18:21**

**Date Received: 02/04/16 09:45**

**Lab Sample ID: 320-17154-10**

**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			535.9 mL	1.00 mL	100147	02/10/16 07:14	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	535.9 mL	1.00 mL	100906	02/20/16 12:45	JRB	TAL SAC

**Laboratory References:**

Maxxam = Maxxam Analytics Inc., PO BOX 57437, Postal Station A, Toronto, Ontario M5W 5M5

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

# Certification Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Laboratory: TestAmerica Sacramento

The certifications listed below are applicable to this report.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2928-01	01-31-17
Oregon	NELAP	10	CA200005	01-29-17

The following analytes are included in this report, but are not certified under this certification:

Analysis Method	Prep Method	Matrix	Analyte
WS-LC-0025	3535	Water	Perfluoroheptanoic acid (PFHpA)
WS-LC-0025	3535	Water	Perfluorononanoic acid (PFNA)
WS-LC-0025	3535	Water	Perfluorooctanoic acid (PFOA)

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

Analysis Method	Prep Method	Matrix	Analyte
WS-LC-0025	3535	Water	Perfluorobutanesulfonic acid (PFBS)
WS-LC-0025	3535	Water	Perfluorohexanesulfonic acid (PFHxS)
WS-LC-0025	3535	Water	Perfluorooctanesulfonic acid (PFOS)

## Laboratory: TestAmerica Denver

The certifications listed below are applicable to this report.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2907.01	10-31-17
Oregon	NELAP	10	4025	01-09-17

# Method Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

Method	Method Description	Protocol	Laboratory
WS-LC-0025	Perfluorinated Hydrocarbons	TAL SOP	TAL SAC
PFOA - PFOA	PFOA	NONE	Maxxam

**Protocol References:**

NONE = NONE

TAL SOP = TestAmerica Laboratories, Standard Operating Procedure

**Laboratory References:**

Maxxam = Maxxam Analytics Inc., PO BOX 57437, Postal Station A, Toronto, Ontario M5W 5M5

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

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- 14
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- 16

# Sample Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-17154-1	OF-FB09-0216	Water	02/03/16 16:40	02/04/16 09:45
320-17154-2	OF-RW09-0216	Water	02/03/16 16:45	02/04/16 09:45
320-17154-3	OF-FB37-0216	Water	02/03/16 17:20	02/04/16 09:45
320-17154-4	OF-RW37-0216	Water	02/03/16 17:28	02/04/16 09:45
320-17154-5	OF-RW11-0216	Water	02/03/16 16:04	02/04/16 09:45
320-17154-6	OF-FB11-0216	Water	02/03/16 16:10	02/04/16 09:45
320-17154-7	OF-RW28-0216	Water	02/03/16 17:41	02/04/16 09:45
320-17154-8	OF-FB28-0216	Water	02/03/16 17:43	02/04/16 09:45
320-17154-9	OF-RW67-0216	Water	02/03/16 18:19	02/04/16 09:45
320-17154-10	OF-FB67-0216	Water	02/03/16 18:21	02/04/16 09:45





Your Project #: 320-17154  
Your C.O.C. #: 283602

**Attention:PFC Reporting Group**

TestAmerica  
Sacramento  
880 Riverside Parkway  
West Sacramento, CA  
USA 95605

**Report Date: 2016/02/26**  
Report #: R3909946  
Version: 2 - Revision

**CERTIFICATE OF ANALYSIS – REVISED REPORT**

**MAXXAM JOB #: B630797**  
**Received: 2016/02/13, 13:40**

Sample Matrix: Water  
# Samples Received: 5

Analyses	Date		Laboratory Method	Reference
	Quantity Extracted	Date Analyzed		
Low level PFOS and PFOA in water	5	2016/02/18	2016/02/19 CAM SOP-00894	EPA 537 m

Reference Method suffix "m" indicates test methods incorporate validated modifications from specific reference methods to improve performance.

\* RPDs calculated using raw data. The rounding of final results may result in the apparent difference.

U = Undetected at the limit of quantitation.

J = Estimated concentration between the EDL & RDL.

B = Blank Contamination.

Q = One or more quality control criteria failed.

E = Analyte concentration exceeds the maximum concentration level.

K = Estimated maximum possible concentration due to ion abundance ratio failure.

**Encryption Key**

Please direct all questions regarding this Certificate of Analysis to your Project Manager.

Melissa DiGrazia, Project Manager - ATUT

Email: MDiGrazia@maxxam.ca

Phone# (905) 817-5700

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Maxxam has procedures in place to guard against improper use of the electronic signature and have the required "signatories", as per section 5.10.2 of ISO/IEC 17025:2005(E), signing the reports. For Service Group specific validation please refer to the Validation Signature Page.

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**RESULTS OF ANALYSES OF WATER**

Maxxam ID		BVX826	BVX827	BVX828	BVX829	BVX830			
Sampling Date		2016/02/03 16:45	2016/02/03 17:28	2016/02/03 16:04	2016/02/03 17:41	2016/02/03 18:19			
COC Number		283602	283602	283602	283602	283602			
	UNITS	OF-RW09-0216	OF-RW37-0216	OF-RW11-0216	OF-RW28-0216	OF-RW67-0216	MDL	QC Batch	RDL
Perfluorobutane Sulfonate (PFBS)	ng/L	0.27 U	0.79 J	0.27 U	0.27 U	0.97 J	0.27	4386408	2.0
Perfluoroheptanoic Acid (PFHpA)	ng/L	0.39 U	0.46 J	0.39 U	0.39 U	0.39 U	0.39	4386408	2.0
Perfluorohexane Sulfonate (PFHxS)	ng/L	0.71 J	3.3	0.75 J	0.75 J	0.40 U	0.40	4386408	2.0
Perfluoro-n-Octanoic Acid (PFOA)	ng/L	0.41 J	7.2	0.39 U	0.39 U	0.39 U	0.39	4386408	2.0
Perfluorononanoic Acid (PFNA)	ng/L	0.33 U	0.33 U	0.33 U	0.37 J	0.33 U	0.33	4386408	2.0
Perfluorooctane Sulfonate (PFOS)	ng/L	0.32 J	0.81 J	0.87 J	0.30 U	0.30 U	0.30	4386408	2.0
Surrogate Recovery (%)									
13C4-Perfluoroheptanoic acid	%	54	52	55	54	48 (1)	N/A	4386408	N/A
13C4-Perfluorooctanesulfonate	%	59	64	54	59	51	N/A	4386408	N/A
13C4-Perfluorooctanoic acid	%	58	57	59	54	51	N/A	4386408	N/A
13C5-Perfluorononanoic acid	%	62	65	61	57	55	N/A	4386408	N/A
18O2-Perfluorohexanesulfonate	%	69	64	60	50	56	N/A	4386408	N/A
QC Batch = Quality Control Batch									
N/A = Not Applicable									
(1) Surrogate recovery was below the defined lower control limit (LCL). Laboratory spiked water resulted in satisfactory recovery of the surrogate. When considered together, these QC data suggest that matrix interferences may be biasing the data low. Because quantitation is performed using isotope dilution techniques, any losses of the native compound that may occur during any of the sample preparation, extraction, cleanup or determinative steps will be mirrored by a similar loss of the labeled standard, and as such can be accounted for and corrected. Therefore, the quantification of these target compounds is not affected by the low surrogate recovery.									



**TEST SUMMARY**

**Maxxam ID:** BVX826  
**Sample ID:** OF-RW09-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX827  
**Sample ID:** OF-RW37-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX828  
**Sample ID:** OF-RW11-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX829  
**Sample ID:** OF-RW28-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX830  
**Sample ID:** OF-RW67-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara



**GENERAL COMMENTS**

Perfluorinated Compounds (PFCs):

Perfluoroheptanoic acid (PFHpA) MDL = 0.39, LOD = 1.0, LOQ = 2.0  
Perfluorooctanoic acid (PFOA) MDL = 0.39, LOD = 1.0, LOQ = 2.0  
Perfluorononanoic acid (PFNA) MDL = 0.33, LOD = 1.0, LOQ = 2.0  
Perfluorobutane sulfonate (PFBS) MDL = 0.27, LOD = 1.0, LOQ = 2.0  
Perfluorohexane sulfonate (PFHxA) MDL = 0.40, LOD = 1.0, LOQ = 2.0  
Perfluorooctane sulfonate (PFOS) MDL = 0.30, LOD = 1.0, LOQ = 2.0  
All Units are in ng/L

**Results relate only to the items tested.**



**QUALITY ASSURANCE REPORT**

QA/QC	Batch	Init	QC Type	Parameter	Date Analyzed	Value	% Recovery	UNITS	QC Limits		
4386408	CM5	Matrix Spike	13C4-Perfluoroheptanoic acid	2016/02/19		50	%	50 - 130			
			13C4-Perfluorooctanesulfonate	2016/02/19		59	%	50 - 130			
			13C4-Perfluorooctanoic acid	2016/02/19		58	%	50 - 130			
			13C5-Perfluorononanoic acid	2016/02/19		61	%	50 - 130			
			18O2-Perfluorohexanesulfonate	2016/02/19		53	%	50 - 130			
			Perfluorobutane Sulfonate (PFBS)	2016/02/19		157 (1)	%	70 - 130			
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19		133 (1)	%	70 - 130			
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19		148 (1)	%	70 - 130			
			Perfluorononanoic Acid (PFNA)	2016/02/19		120	%	70 - 130			
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19		125	%	70 - 130			
			Perfluorooctane Sulfonate (PFOS)	2016/02/19		116	%	70 - 130			
			4386408	CM5	Matrix Spike DUP	13C4-Perfluoroheptanoic acid	2016/02/19		52	%	50 - 130
						13C4-Perfluorooctanesulfonate	2016/02/19		64	%	50 - 130
						13C4-Perfluorooctanoic acid	2016/02/19		66	%	50 - 130
						13C5-Perfluorononanoic acid	2016/02/19		70	%	50 - 130
						18O2-Perfluorohexanesulfonate	2016/02/19		58	%	50 - 130
						Perfluorobutane Sulfonate (PFBS)	2016/02/19		150 (1)	%	70 - 130
Perfluoroheptanoic Acid (PFHpA)	2016/02/19					125	%	70 - 130			
Perfluorohexane Sulfonate (PFHxS)	2016/02/19					133 (1)	%	70 - 130			
Perfluorononanoic Acid (PFNA)	2016/02/19					104	%	70 - 130			
Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19					114	%	70 - 130			
Perfluorooctane Sulfonate (PFOS)	2016/02/19					109	%	70 - 130			
4386408	CM5	MS/MSD RPD				Perfluorobutane Sulfonate (PFBS)	2016/02/19	4.2		%	30
						Perfluoroheptanoic Acid (PFHpA)	2016/02/19	6.2		%	30
						Perfluorohexane Sulfonate (PFHxS)	2016/02/19	11		%	30
						Perfluorononanoic Acid (PFNA)	2016/02/19	14		%	30
						Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19	8.5		%	30
						Perfluorooctane Sulfonate (PFOS)	2016/02/19	6.4		%	30
			4386408	CM5	Spiked Blank	13C4-Perfluoroheptanoic acid	2016/02/19		67	%	50 - 130
13C4-Perfluorooctanesulfonate	2016/02/19					77	%	50 - 130			
13C4-Perfluorooctanoic acid	2016/02/19					70	%	50 - 130			
13C5-Perfluorononanoic acid	2016/02/19					73	%	50 - 130			
18O2-Perfluorohexanesulfonate	2016/02/19					83	%	50 - 130			
Perfluorobutane Sulfonate (PFBS)	2016/02/19					107	%	70 - 130			
Perfluoroheptanoic Acid (PFHpA)	2016/02/19					121	%	70 - 130			
Perfluorohexane Sulfonate (PFHxS)	2016/02/19					115	%	70 - 130			
Perfluorononanoic Acid (PFNA)	2016/02/19					118	%	70 - 130			
Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19					124	%	70 - 130			
Perfluorooctane Sulfonate (PFOS)	2016/02/19					112	%	70 - 130			
4386408	CM5	Method Blank				13C4-Perfluoroheptanoic acid	2016/02/19		74	%	50 - 130
						13C4-Perfluorooctanesulfonate	2016/02/19		82	%	50 - 130
						13C4-Perfluorooctanoic acid	2016/02/19		74	%	50 - 130
						13C5-Perfluorononanoic acid	2016/02/19		76	%	50 - 130
						18O2-Perfluorohexanesulfonate	2016/02/19		86	%	50 - 130
						Perfluorobutane Sulfonate (PFBS)	2016/02/19	0.27 U, MDL=0.27		ng/L	
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19	0.39 U, MDL=0.39		ng/L				
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19	0.40 U, MDL=0.40		ng/L				
			Perfluorononanoic Acid (PFNA)	2016/02/19	0.33 U, MDL=0.33		ng/L				
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19	0.39 U, MDL=0.39		ng/L				

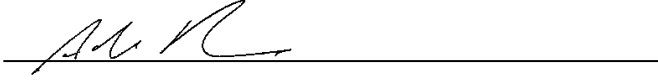
**QUALITY ASSURANCE REPORT(CONT'D)**

QA/QC				Date		%		
Batch	Init	QC Type	Parameter	Analyzed	Value	Recovery	UNITS	QC Limits
			Perfluorooctane Sulfonate (PFOS)	2016/02/19	0.30 U, MDL=0.30		ng/L	
<p>Matrix Spike: A sample to which a known amount of the analyte of interest has been added. Used to evaluate sample matrix interference.</p> <p>Spiked Blank: A blank matrix sample to which a known amount of the analyte, usually from a second source, has been added. Used to evaluate method accuracy.</p> <p>Method Blank: A blank matrix containing all reagents used in the analytical procedure. Used to identify laboratory contamination.</p> <p>Surrogate: A pure or isotopically labeled compound whose behavior mirrors the analytes of interest. Used to evaluate extraction efficiency.</p> <p>(1) Recovery or RPD for this parameter is outside control limits. The overall quality control for this analysis meets acceptability criteria.</p>								



**VALIDATION SIGNATURE PAGE**

The analytical data and all QC contained in this report were reviewed and validated by the following individual(s).



Adam Robinson, Supervisor, LC/MS/MS

---

Maxxam has procedures in place to guard against improper use of the electronic signature and have the required "signatories", as per section 5.10.2 of ISO/IEC 17025:2005(E), signing the reports. For Service Group specific validation please refer to the Validation Signature Page.

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CTD WE7G

320-17154 Chain of Custody

Temperature on Receipt \_\_\_\_\_  
Drinking Water? Yes  No

# Chain of Custody Record

TAL-4124 (1007)

Client: CH2M HILL Project Manager: Bill Friedman Chain of Custody Number: 283602  
 Address: 5701 Cleveland St, Suite 200 Telephone Number (Area Code)/Fax Number: 757-671-6223 Lab Number: 02/03/16  
 City: Virginia Beach State: VA Zip Code: 23462 Site Contact: \_\_\_\_\_ Lab Contact: \_\_\_\_\_ Page: 1 of 1  
 Project Name and Location (State): CTD WE7G PFC Sampling Carrier/Mailbox Number: FedEx  
 Contract/Purchase Order/Quote No.: PO# 938652

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix						Containers & Preservatives						Special Instructions/ Conditions of Receipt		
			Air	Soil	Urbres	H2SO4	HNO3	HCl	HNO3	ZnAc	H2O2	HOAc					
OF-FB09-0216	02/03/16	1646	X		X												
OF-RW09-0216		1645															
OF-FB37-0216		1720															
OF-RW37-0216		1728															
OF-RW11-0216		1604															
OF-FB11-0216		1610															
OF-RW28-0216		1741															
OF-FB28-0216		1743															
OF-RW67-0216		1819															
OF-FB67-0216		1821															

Possible Hazard Identification:  
 Non-Hazard  Flammable  Skin Irritant  Poison B  Poison A  Unknown  Return To Client  Archive For \_\_\_\_\_ Months (A fee may be assessed if samples are retained longer than 1 month)  
 Disposal By Lab  Sample Disposal

Turn Around Time Required:  
 24 Hours  48 Hours  7 Days  14 Days  21 Days  Other \_\_\_\_\_

QC Requirements (Specify): \_\_\_\_\_

1. Relinquished By: Y. Williams Date: 02/03/16 Time: 19:30  
 2. Relinquished By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 3. Relinquished By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Received By: J. Sale Date: 2/1/16 Time: 9:45  
 Received By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 Received By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Comments: 0.3



**Chain of Custody Record**

Due 2/22

Temperature on Receipt \_\_\_\_\_



320-17154 Chain of Custody

CTD WEF7G

Drinking Water? Yes  No

Project Manager [Redacted]	Date 02/03/16	Chain of Custody Number 283602
Telephone Number (Area Code)/Fax Number [Redacted]	Lab Number	Page 1 of 1
Site Contact	Lab Contact	

Carrier/Waybill Number  
FedEx

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						Select PCBs	Special Instructions/ Conditions of Receipt	
			Air	Aqueous	Solid	Soil	Uppres	HPDCA	AWC3	HCl	NiOH	ZnAc2			NiOH
<del>OF-RW09-0216</del>	<del>02/02/16</del>	<del>1646</del>	<del>X</del>												
OF-RW09-0216		1645													
<del>OF-RW37-0216</del>		1720													
OF-RW37-0216		1728													
<del>OF-RW11-0216</del>		1604													
OF-RW11-0216		1610													
<del>OF-RW28-0216</del>		1741													
OF-RW28-0216		1743													
<del>OF-RW67-0216</del>		1819													
OF-RW67-0216		1821													

13-Feb-16 13:40  
Hongmei Zhao (Grace)  
B630797

RGN ENV-951

Possible Hazard Identification  
 Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown

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 \_\_\_\_\_

QC Requirements (Specify)

1. Relinquished By Kathryn Smith	Date 02/03/16	Time 19:30	1. Received By [Signature]	Date 2/9/16	Time 9:45
2. Relinquished By	Date	Time	2. Received By [Signature]	Date 206/02/13	Time 13:40
3. Relinquished By	Date	Time	3. Received By	Date	Time

Comments

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

0.3<sup>v</sup> REFER TO ACTR

## Login Sample Receipt Checklist

Client: CH2M Hill, Inc.

Job Number: 320-17154-1  
SDG Number: CTO WE7G PFC Sampling

**Login Number: 17154**  
**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	
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.	True	
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.	N/A	
Sample Preservation Verified.	True	
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: 320-17154-1  
SDG Number: CTO WE7G PFC Sampling  
Job Description: CTO WE7G PFC Sampling

For:  
CH2M Hill, Inc.  
5701 Cleveland Street  
Suite 200  
Virginia Beach, VA 23462  
Attention: Laurie George



Approved for release.  
Laura Turpen  
Project Manager I  
3/22/2016 12:44 PM

---

Laura Turpen, Project Manager I  
880 Riverside Parkway, West Sacramento, CA, 95605  
(916)374-4414  
laura.turpen@testamericainc.com  
03/22/2016  
Revision: 1

The test results in this report relate only to the samples in this report and meet all requirements of NELAP, with any exceptions noted. Pursuant to NELAP, this report shall not be reproduced except in full, without the written approval of the laboratory. All questions regarding this report should be directed to the TestAmerica Denver Project Manager.

The Lab Certification ID# is 4025.

Reporting limits are adjusted for sample size used, dilutions and moisture content if applicable.

**TestAmerica Laboratories, Inc.**

TestAmerica Sacramento 880 Riverside Parkway, West Sacramento, CA 95605  
Tel (916) 373-5600 Fax (916) 372-1059 [www.testamericainc.com](http://www.testamericainc.com)



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# Definitions/Glossary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Qualifiers

### LCMS

Qualifier	Qualifier Description
U	Undetected at the Limit of Detection.
J	Estimated: The analyte was positively identified; the quantitation is an estimation

## 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: CH2M Hill, Inc.

Project: CTO WE7G PFC Sampling

Report Number: 320-17154-1

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.

### Revision

This report was revised March 22, 2016 to include all method blank data for all analytes and to remove the B qualifiers from the LCS/LCSD in batch 100093. No other data has changed as a result of this revision.

### RECEIPT

The samples were received on 02/04/2016; the samples arrived in good condition, properly preserved and on ice. The temperature of the coolers at receipt was 0.3 C.

### PFC

Samples OF-FB09-0216 (320-17154-1), OF-FB37-0216 (320-17154-3), OF-FB11-0216 (320-17154-6), OF-FB28-0216 (320-17154-8) and OF-FB67-0216 (320-17154-10) were analyzed for PFC in accordance with PFC. The samples were prepared on 02/09/2016 and 02/10/2016 and analyzed on 02/20/2016.

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

### Organic Prep

Method(s) 3535: Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with batch 320-100093

Method(s) 3535: Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate (MS/MSD) associated with batch 320-100147.

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

### Subcontract Work

**PFC:** This method was subcontracted to Maxxam Analytics Inc.. The subcontract laboratory certification is different from that of the facility issuing the final report. Any analytical or quality issues are noted in the subcontract portion of the report.

The DL/LOD/LOQ limits for Maxxam, along with information on the data qualifiers, are included in the narrative (Page 8 of 143 of the subcontract report; page 494 of 632 of the entire report).

# Detection Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB09-0216**

**Lab Sample ID: 320-17154-1**

No Detections.

**Client Sample ID: OF-RW09-0216**

**Lab Sample ID: 320-17154-2**

No Detections.

**Client Sample ID: OF-FB37-0216**

**Lab Sample ID: 320-17154-3**

No Detections.

**Client Sample ID: OF-RW37-0216**

**Lab Sample ID: 320-17154-4**

No Detections.

**Client Sample ID: OF-RW11-0216**

**Lab Sample ID: 320-17154-5**

No Detections.

**Client Sample ID: OF-FB11-0216**

**Lab Sample ID: 320-17154-6**

No Detections.

**Client Sample ID: OF-RW28-0216**

**Lab Sample ID: 320-17154-7**

No Detections.

**Client Sample ID: OF-FB28-0216**

**Lab Sample ID: 320-17154-8**

No Detections.

**Client Sample ID: OF-RW67-0216**

**Lab Sample ID: 320-17154-9**

No Detections.

**Client Sample ID: OF-FB67-0216**

**Lab Sample ID: 320-17154-10**

No Detections.

This Detection Summary does not include radiochemical test results.

# Client Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB09-0216**

**Lab Sample ID: 320-17154-1**

Date Collected: 02/03/16 16:40

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00079	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00074	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00091	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00086	ug/L		02/09/16 10:23	02/20/16 10:16	1
Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0013	ug/L		02/09/16 10:23	02/20/16 10:16	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4-PFHpA	122		25 - 150				02/09/16 10:23	02/20/16 10:16	1
13C4 PFOA	126		25 - 150				02/09/16 10:23	02/20/16 10:16	1
13C5 PFNA	127		25 - 150				02/09/16 10:23	02/20/16 10:16	1
18O2 PFHxS	113		25 - 150				02/09/16 10:23	02/20/16 10:16	1
13C4 PFOS	107		25 - 150				02/09/16 10:23	02/20/16 10:16	1

**Client Sample ID: OF-FB37-0216**

**Lab Sample ID: 320-17154-3**

Date Collected: 02/03/16 17:20

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.00073	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.00068	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.00059	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.00083	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.00079	ug/L		02/09/16 10:23	02/20/16 10:37	1
Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0012	ug/L		02/09/16 10:23	02/20/16 10:37	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4-PFHpA	107		25 - 150				02/09/16 10:23	02/20/16 10:37	1
13C4 PFOA	111		25 - 150				02/09/16 10:23	02/20/16 10:37	1
13C5 PFNA	113		25 - 150				02/09/16 10:23	02/20/16 10:37	1
18O2 PFHxS	100		25 - 150				02/09/16 10:23	02/20/16 10:37	1
13C4 PFOS	95		25 - 150				02/09/16 10:23	02/20/16 10:37	1

**Client Sample ID: OF-FB11-0216**

**Lab Sample ID: 320-17154-6**

Date Collected: 02/03/16 16:10

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.00073	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.00068	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.00059	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.00083	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.00079	ug/L		02/09/16 10:23	02/20/16 10:58	1
Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0012	ug/L		02/09/16 10:23	02/20/16 10:58	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4-PFHpA	111		25 - 150				02/09/16 10:23	02/20/16 10:58	1
13C4 PFOA	113		25 - 150				02/09/16 10:23	02/20/16 10:58	1
13C5 PFNA	116		25 - 150				02/09/16 10:23	02/20/16 10:58	1
18O2 PFHxS	95		25 - 150				02/09/16 10:23	02/20/16 10:58	1



# Client Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

**Client Sample ID: OF-FB11-0216**

**Lab Sample ID: 320-17154-6**

Date Collected: 02/03/16 16:10

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons (Continued)**

Isotope Dilution	%Recovery	Qualifier	Limits	Prepared	Analyzed	Dil Fac
<sup>13</sup> C4 PFOS	93		25 - 150	02/09/16 10:23	02/20/16 10:58	1

**Client Sample ID: OF-FB28-0216**

**Lab Sample ID: 320-17154-8**

Date Collected: 02/03/16 17:43

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0024	0.00077	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorooctanoic acid (PFOA)	0.0019	U	0.0024	0.00072	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorononanoic acid (PFNA)	0.0019	U	0.0024	0.00063	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0024	0.00088	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0024	0.00083	ug/L		02/10/16 07:14	02/20/16 12:23	1
Perfluorooctanesulfonic acid (PFOS)	0.0029	U	0.0038	0.0012	ug/L		02/10/16 07:14	02/20/16 12:23	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
<sup>13</sup> C4-PFHpA	124		25 - 150				02/10/16 07:14	02/20/16 12:23	1
<sup>13</sup> C4 PFOA	132		25 - 150				02/10/16 07:14	02/20/16 12:23	1
<sup>13</sup> C5 PFNA	137		25 - 150				02/10/16 07:14	02/20/16 12:23	1
<sup>18</sup> O2 PFHxS	113		25 - 150				02/10/16 07:14	02/20/16 12:23	1
<sup>13</sup> C4 PFOS	108		25 - 150				02/10/16 07:14	02/20/16 12:23	1

**Client Sample ID: OF-FB67-0216**

**Lab Sample ID: 320-17154-10**

Date Collected: 02/03/16 18:21

Matrix: Water

Date Received: 02/04/16 09:45

**Method: WS-LC-0025 - Perfluorinated Hydrocarbons**

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0023	0.00075	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorooctanoic acid (PFOA)	0.0019	U	0.0023	0.00070	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorononanoic acid (PFNA)	0.0019	U	0.0023	0.00061	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0023	0.00086	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0023	0.00081	ug/L		02/10/16 07:14	02/20/16 12:45	1
Perfluorooctanesulfonic acid (PFOS)	0.0028	U	0.0037	0.0012	ug/L		02/10/16 07:14	02/20/16 12:45	1
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
<sup>13</sup> C4-PFHpA	126		25 - 150				02/10/16 07:14	02/20/16 12:45	1
<sup>13</sup> C4 PFOA	131		25 - 150				02/10/16 07:14	02/20/16 12:45	1
<sup>13</sup> C5 PFNA	130		25 - 150				02/10/16 07:14	02/20/16 12:45	1
<sup>18</sup> O2 PFHxS	113		25 - 150				02/10/16 07:14	02/20/16 12:45	1
<sup>13</sup> C4 PFOS	112		25 - 150				02/10/16 07:14	02/20/16 12:45	1

# Default Detection Limits

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons

Analyte	LOQ	DL	Units	Method
Perfluorobutanesulfonic acid (PFBS)	0.0025	0.00092	ug/L	WS-LC-0025
Perfluoroheptanoic acid (PFHpA)	0.0025	0.00080	ug/L	WS-LC-0025
Perfluorohexanesulfonic acid (PFHxS)	0.0025	0.00087	ug/L	WS-LC-0025
Perfluorononanoic acid (PFNA)	0.0025	0.00065	ug/L	WS-LC-0025
Perfluorooctanesulfonic acid (PFOS)	0.0040	0.0013	ug/L	WS-LC-0025
Perfluorooctanoic acid (PFOA)	0.0025	0.00075	ug/L	WS-LC-0025

# Isotope Dilution Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons

Matrix: Water

Prep Type: Total/NA

Lab Sample ID	Client Sample ID	Percent Isotope Dilution Recovery (Acceptance Limits)				
		<sup>13</sup> C4-PFHp (25-150)	<sup>13</sup> C4 PFO (25-150)	<sup>13</sup> C5 PFN (25-150)	<sup>18</sup> O2 PFHx (25-150)	<sup>13</sup> C4 PFO (25-150)
320-17154-1	OF-FB09-0216	122	126	127	113	107
320-17154-3	OF-FB37-0216	107	111	113	100	95
320-17154-6	OF-FB11-0216	111	113	116	95	93
320-17154-8	OF-FB28-0216	124	132	137	113	108
320-17154-10	OF-FB67-0216	126	131	130	113	112
LCS 320-100093/2-A	Lab Control Sample	95	95	98	90	82
LCS 320-100147/2-A	Lab Control Sample	124	119	117	108	99
LCSD 320-100093/3-A	Lab Control Sample Dup	95	88	91	86	80
LCSD 320-100147/3-A	Lab Control Sample Dup	123	115	116	107	100
MB 320-100093/1-A	Method Blank				95	
MB 320-100093/1-A	Method Blank	103	101	103		92
MB 320-100147/1-A	Method Blank	118	114	118	97	98

### Surrogate Legend

<sup>13</sup>C4-PFHpA = <sup>13</sup>C4-PFHpA  
<sup>13</sup>C4 PFOA = <sup>13</sup>C4 PFOA  
<sup>13</sup>C5 PFNA = <sup>13</sup>C5 PFNA  
<sup>18</sup>O2 PFHxS = <sup>18</sup>O2 PFHxS  
<sup>13</sup>C4 PFOS = <sup>13</sup>C4 PFOS

# QC Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons

**Lab Sample ID: MB 320-100093/1-A**  
**Matrix: Water**  
**Analysis Batch: 100305**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	MB	MB	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier							
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00087	ug/L		02/09/16 10:23	02/13/16 00:22	1
<b>MB MB</b>									
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
18O2 PFHxS	95		25 - 150				02/09/16 10:23	02/13/16 00:22	1

**Lab Sample ID: MB 320-100093/1-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	MB	MB	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier							
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00080	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		02/09/16 10:23	02/20/16 06:23	1
Perfluorooctanesulfonic acid (PFOS)	0.00143	J	0.0040	0.0013	ug/L		02/09/16 10:23	02/20/16 06:23	1
<b>MB MB</b>									
Isotope Dilution	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4-PFHpA	103		25 - 150				02/09/16 10:23	02/20/16 06:23	1
13C4 PFOA	101		25 - 150				02/09/16 10:23	02/20/16 06:23	1
13C5 PFNA	103		25 - 150				02/09/16 10:23	02/20/16 06:23	1
13C4 PFOS	92		25 - 150				02/09/16 10:23	02/20/16 06:23	1

**Lab Sample ID: LCS 320-100093/2-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	Spike Added	LCS	LCS	Unit	D	%Rec	%Rec.	Limits
		Result	Qualifier					
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0513		ug/L		128		60 - 140
Perfluorooctanoic acid (PFOA)	0.0400	0.0485		ug/L		121		60 - 140
Perfluorononanoic acid (PFNA)	0.0400	0.0537		ug/L		134		60 - 140
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0459		ug/L		130		50 - 150
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0480		ug/L		127		60 - 140
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0515		ug/L		135		60 - 140
<b>LCS LCS</b>								
Isotope Dilution	%Recovery	Qualifier	Limits					
13C4-PFHpA	95		25 - 150					
13C4 PFOA	95		25 - 150					
13C5 PFNA	98		25 - 150					
18O2 PFHxS	90		25 - 150					
13C4 PFOS	82		25 - 150					

# QC Sample Results

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Method: WS-LC-0025 - Perfluorinated Hydrocarbons (Continued)

**Lab Sample ID: LCSD 320-100093/3-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 100093**

Analyte	Spike Added	LCS	LCS	Unit	D	%Rec	%Rec.		RPD	Limit
							Limits	RPD		
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0506		ug/L		126	60 - 140	1	30	
Perfluorooctanoic acid (PFOA)	0.0400	0.0493		ug/L		123	60 - 140	2	30	
Perfluorononanoic acid (PFNA)	0.0400	0.0517		ug/L		129	60 - 140	4	30	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0472		ug/L		133	50 - 150	3	30	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0486		ug/L		128	60 - 140	1	30	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0452		ug/L		118	60 - 140	13	30	

Isotope Dilution	LCS		Limits
	%Recovery	Qualifier	
13C4-PFHpA	95		25 - 150
13C4 PFOA	88		25 - 150
13C5 PFNA	91		25 - 150
18O2 PFHxS	86		25 - 150
13C4 PFOS	80		25 - 150

**Lab Sample ID: MB 320-100147/1-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

Analyte	MB		LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
	Result	Qualifier							
Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.00080	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.00075	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.00065	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.00092	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.00087	ug/L		02/10/16 07:14	02/20/16 11:20	1
Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0013	ug/L		02/10/16 07:14	02/20/16 11:20	1

Isotope Dilution	MB		Limits	Prepared	Analyzed	Dil Fac
	%Recovery	Qualifier				
13C4-PFHpA	118		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C4 PFOA	114		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C5 PFNA	118		25 - 150	02/10/16 07:14	02/20/16 11:20	1
18O2 PFHxS	97		25 - 150	02/10/16 07:14	02/20/16 11:20	1
13C4 PFOS	98		25 - 150	02/10/16 07:14	02/20/16 11:20	1

**Lab Sample ID: LCS 320-100147/2-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

Analyte	Spike Added	LCS	LCS	Unit	D	%Rec	%Rec.	
							Limits	RPD
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0498		ug/L		124	60 - 140	
Perfluorooctanoic acid (PFOA)	0.0400	0.0476		ug/L		119	60 - 140	
Perfluorononanoic acid (PFNA)	0.0400	0.0528		ug/L		132	60 - 140	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0412		ug/L		116	50 - 150	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0503		ug/L		133	60 - 140	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0529		ug/L		138	60 - 140	

# QC Sample Results

Client: CH2M Hill, Inc.  
 Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
 SDG: CTO WE7G PFC Sampling

<i>Isotope Dilution</i>	<i>LCS LCS</i>		<i>Limits</i>
	<i>%Recovery</i>	<i>Qualifier</i>	
13C4-PFHpA	124		25 - 150
13C4 PFOA	119		25 - 150
13C5 PFNA	117		25 - 150
18O2 PFHxS	108		25 - 150
13C4 PFOS	99		25 - 150

**Lab Sample ID: LCSD 320-100147/3-A**  
**Matrix: Water**  
**Analysis Batch: 100906**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 100147**

<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	0.0488		ug/L		122	60 - 140	2	30
Perfluorononanoic acid (PFNA)	0.0400	0.0534		ug/L		133	60 - 140	1	30
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0461		ug/L		130	50 - 150	11	30
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0473		ug/L		125	60 - 140	6	30
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0502		ug/L		131	60 - 140	5	30

<i>Isotope Dilution</i>	<i>LCSD LCSD</i>		<i>Limits</i>
	<i>%Recovery</i>	<i>Qualifier</i>	
13C4-PFHpA	123		25 - 150
13C4 PFOA	115		25 - 150
13C5 PFNA	116		25 - 150
18O2 PFHxS	107		25 - 150
13C4 PFOS	100		25 - 150

# QC Association Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## LCMS

### Prep Batch: 100093

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-1	OF-FB09-0216	Total/NA	Water	3535	
320-17154-3	OF-FB37-0216	Total/NA	Water	3535	
320-17154-6	OF-FB11-0216	Total/NA	Water	3535	
LCS 320-100093/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-100093/3-A	Lab Control Sample Dup	Total/NA	Water	3535	
MB 320-100093/1-A	Method Blank	Total/NA	Water	3535	

### Prep Batch: 100147

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-8	OF-FB28-0216	Total/NA	Water	3535	
320-17154-10	OF-FB67-0216	Total/NA	Water	3535	
LCS 320-100147/2-A	Lab Control Sample	Total/NA	Water	3535	
LCSD 320-100147/3-A	Lab Control Sample Dup	Total/NA	Water	3535	
MB 320-100147/1-A	Method Blank	Total/NA	Water	3535	

### Analysis Batch: 100305

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
MB 320-100093/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100093

### Analysis Batch: 100906

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-17154-1	OF-FB09-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-3	OF-FB37-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-6	OF-FB11-0216	Total/NA	Water	WS-LC-0025	100093
320-17154-8	OF-FB28-0216	Total/NA	Water	WS-LC-0025	100147
320-17154-10	OF-FB67-0216	Total/NA	Water	WS-LC-0025	100147
LCS 320-100093/2-A	Lab Control Sample	Total/NA	Water	WS-LC-0025	100093
LCS 320-100147/2-A	Lab Control Sample	Total/NA	Water	WS-LC-0025	100147
LCSD 320-100093/3-A	Lab Control Sample Dup	Total/NA	Water	WS-LC-0025	100093
LCSD 320-100147/3-A	Lab Control Sample Dup	Total/NA	Water	WS-LC-0025	100147
MB 320-100093/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100093
MB 320-100147/1-A	Method Blank	Total/NA	Water	WS-LC-0025	100147

# Lab Chronicle

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

## Client Sample ID: OF-FB09-0216

Date Collected: 02/03/16 16:40

Date Received: 02/04/16 09:45

## Lab Sample ID: 320-17154-1

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	100906	02/20/16 10:16	JRB	TAL SAC

## Client Sample ID: OF-FB37-0216

Date Collected: 02/03/16 17:20

Date Received: 02/04/16 09:45

## Lab Sample ID: 320-17154-3

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	100906	02/20/16 10:37	JRB	TAL SAC

## Client Sample ID: OF-FB11-0216

Date Collected: 02/03/16 16:10

Date Received: 02/04/16 09:45

## Lab Sample ID: 320-17154-6

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			100093	02/09/16 10:23	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	100906	02/20/16 10:58	JRB	TAL SAC

## Client Sample ID: OF-FB28-0216

Date Collected: 02/03/16 17:43

Date Received: 02/04/16 09:45

## Lab Sample ID: 320-17154-8

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			100147	02/10/16 07:14	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	100906	02/20/16 12:23	JRB	TAL SAC

## Client Sample ID: OF-FB67-0216

Date Collected: 02/03/16 18:21

Date Received: 02/04/16 09:45

## Lab Sample ID: 320-17154-10

Matrix: Water

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			100147	02/10/16 07:14	HJA	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	100906	02/20/16 12:45	JRB	TAL SAC

### Laboratory References:

Maxxam = Maxxam Analytics Inc., PO BOX 57437, Postal Station A, Toronto, Ontario M5W 5M5

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



# Certification Summary

Client: CH2M Hill, Inc.  
 Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
 SDG: CTO WE7G PFC Sampling

## Laboratory: TestAmerica Sacramento

The certifications listed below are applicable to this report.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2928-01	01-31-17
Oregon	NELAP	10	CA200005	01-29-17

The following analytes are included in this report, but are not certified under this certification:

Analysis Method	Prep Method	Matrix	Analyte
WS-LC-0025	3535	Water	Perfluoroheptanoic acid (PFHpA)
WS-LC-0025	3535	Water	Perfluorononanoic acid (PFNA)
WS-LC-0025	3535	Water	Perfluorooctanoic acid (PFOA)

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

Method	Prep Method	Matrix	Analyte
WS-LC-0025	3535	Water	Perfluorobutanesulfonic acid (PFBS)
WS-LC-0025	3535	Water	Perfluorohexanesulfonic acid (PFHxS)
WS-LC-0025	3535	Water	Perfluorooctanesulfonic acid (PFOS)

## Laboratory: TestAmerica Denver

The certifications listed below are applicable to this report.

Authority	Program	EPA Region	Certification ID	Expiration Date
A2LA	DoD ELAP		2907.01	10-31-17
Oregon	NELAP	10	4025	01-09-17

# Method Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

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<b>Method</b>	<b>Method Description</b>	<b>Protocol</b>	<b>Laboratory</b>
WS-LC-0025	Perfluorinated Hydrocarbons	TAL SOP	TAL SAC
PFOA - PFOA	PFOA	NONE	Maxxam

**Protocol References:**

NONE = NONE

TAL SOP = TestAmerica Laboratories, Standard Operating Procedure

**Laboratory References:**

Maxxam = Maxxam Analytics Inc., PO BOX 57437, Postal Station A, Toronto, Ontario M5W 5M5

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

# Sample Summary

Client: CH2M Hill, Inc.  
Project/Site: CTO WE7G PFC Sampling

TestAmerica Job ID: 320-17154-1  
SDG: CTO WE7G PFC Sampling

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-17154-1	OF-FB09-0216	Water	02/03/16 16:40	02/04/16 09:45
320-17154-2	OF-RW09-0216	Water	02/03/16 16:45	02/04/16 09:45
320-17154-3	OF-FB37-0216	Water	02/03/16 17:20	02/04/16 09:45
320-17154-4	OF-RW37-0216	Water	02/03/16 17:28	02/04/16 09:45
320-17154-5	OF-RW11-0216	Water	02/03/16 16:04	02/04/16 09:45
320-17154-6	OF-FB11-0216	Water	02/03/16 16:10	02/04/16 09:45
320-17154-7	OF-RW28-0216	Water	02/03/16 17:41	02/04/16 09:45
320-17154-8	OF-FB28-0216	Water	02/03/16 17:43	02/04/16 09:45
320-17154-9	OF-RW67-0216	Water	02/03/16 18:19	02/04/16 09:45
320-17154-10	OF-FB67-0216	Water	02/03/16 18:21	02/04/16 09:45

LCMS MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4 Analysis Batch Number: 100305

Lab Sample ID: STD 320-100305/3 IC Client Sample ID: \_\_\_\_\_

Date Analyzed: 02/12/16 10:56 Lab File ID: 12FEB2016A4A\_004.d GC Column: Acquity ID: 2.1(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Perfluorobutanoic acid (PFBA)	5.69	Baseline	westendor fc	02/15/16 08:14
Perfluorooctanesulfonic acid (PFOS)	11.58	Baseline	westendor fc	02/15/16 08:20
Perfluorononanoic acid (PFNA)	11.60	Baseline	westendor fc	02/15/16 08:20

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
LCMPFCSU_00025	07/29/16	01/29/16	Methanol, Lot Baker 115491	5 mL	LCM2PFHxDA_00003	0.1 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.1 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.1 mL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00004	0.1 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.1 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.1 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.1 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.1 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.1 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.1 mL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.1 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.1 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.1 mL	13C4 PFOS	0.956 ug/mL
LCMPFudA_00005	0.1 mL	13C2 PFUnA	1 ug/mL					
.LCM2PFHxDA_00003	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)	13C2-PFHxDA	50 ug/mL	
.LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)	13C2-PFTeDA	50 ug/mL	
.LCM4PFHPA_00003	05/22/20	Wellington Laboratories, Lot M4PFHpa0515			(Purchased Reagent)	13C4-PFHpa	50 ug/mL	
.LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)	13C5-PFPeA	50 ug/mL	
.LCM8FOSA_00006	12/15/16	Wellington Laboratories, Lot M8FOSA1214I			(Purchased Reagent)	13C8 FOSA	50 ug/mL	
.LCMPFBA_00004	10/31/19	Wellington Laboratories, Lot MPFBA1014			(Purchased Reagent)	13C4 PFBA	50 ug/mL	
.LCMPFDA_00004	04/13/19	Wellington Laboratories, Lot MPFDA0414			(Purchased Reagent)	13C2 PFDA	50 ug/mL	
.LCMPFDoA_00004	07/17/19	Wellington Laboratories, Lot MPFDoA0714			(Purchased Reagent)	13C2 PFDoA	50 ug/mL	
.LCMPFHxA_00005	04/13/19	Wellington Laboratories, Lot MPFHxA0414			(Purchased Reagent)	13C2 PFHxA	50 ug/mL	
.LCMPFHxS_00004	07/25/18	Wellington Laboratories, Lot MPFHxS0713			(Purchased Reagent)	1802 PFHxS	47.3 ug/mL	
.LCMPFNA_00003	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)	13C5 PFNA	50 ug/mL	
.LCMPFOA_00007	04/10/20	Wellington Laboratories, Lot MPFOA0415			(Purchased Reagent)	13C4 PFOA	50 ug/mL	
.LCMPFOS_00009	05/15/20	Wellington Laboratories, Lot MPFOS0515			(Purchased Reagent)	13C4 PFOS	47.8 ug/mL	
.LCMPFudA_00005	10/31/19	Wellington Laboratories, Lot MPFudA1014			(Purchased Reagent)	13C2 PFUnA	50 ug/mL	
LCPFC-L1_00018	06/29/16	12/30/15	MeOH/H2O, Lot 90285	5 mL	LCMPFCSU_00024	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
							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
					LCPFCSP_00040	25 uL	Perfluorobutyric acid	0.5 ng/mL
		Perfluorobutanesulfonic acid (PFBS)	0.442 ng/mL					
		Perfluorodecanoic acid	0.5 ng/mL					

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorododecanoic acid	0.5 ng/mL
							Perfluorodecane Sulfonic acid	0.482 ng/mL
							Perfluoroheptanoic acid (PFHpA)	0.5 ng/mL
							Perfluoroheptanesulfonic Acid	0.476 ng/mL
							Perfluorohexanoic acid	0.5 ng/mL
							Perfluorohexadecanoic acid	0.5 ng/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.473 ng/mL
							Perfluorononanoic acid (PFNA)	0.5 ng/mL
							Perfluorooctanoic acid (PFOA)	0.5 ng/mL
							Perfluorooctadecanoic acid	0.5 ng/mL
							Perfluorooctanesulfonic acid (PFOS)	0.478 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHFA_00003	0.2 mL	13C4-PFHFA	1 ug/mL
					LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.2 mL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00003	11/29/17		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00003	11/29/17		Wellington Laboratories, Lot M2PFTeDA1112		(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHFA_00003	05/22/20		Wellington Laboratories, Lot M4PFHFA0515		(Purchased Reagent)		13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00004	05/22/20		Wellington Laboratories, Lot M5PFPeA0515		(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00004	10/31/19		Wellington Laboratories, Lot MPFBA1014		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00004	04/13/19		Wellington Laboratories, Lot MPFDA0414		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00003	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00007	04/10/20		Wellington Laboratories, Lot MPFOA0415		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00009	05/15/20		Wellington Laboratories, Lot MPFOS0515		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00005	10/31/19		Wellington Laboratories, Lot MPFUdA1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCPFCSP_00040	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFCSP_00039	0.5 mL	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 (PFHpA)	0.1 ug/mL
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL
							Perfluorohexanoic acid	0.1 ug/mL
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.0946 ug/mL
							Perfluorononanoic acid (PFNA)	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctanesulfonic acid (PFOS)	0.0956 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_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL					
LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL					
LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL					
LCPFUdA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL					
...LCPFBA_00003	03/05/18	Wellington Laboratories, Lot PFBA0313			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFBSA_00001	10/09/19	Wellington Laboratories, Lot LPFBS1014			(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
...LCPFDA 00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDoA 00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFDSA 00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
...LCPFHpSA 00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA 00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
...LCPFNA 00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
...LCPFOA 00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFODA 00004	04/25/17		Wellington Laboratories, Lot PFOA0807		(Purchased Reagent)		Perfluorooctandecanoic acid	50 ug/mL
...LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
...LCPFOSA 00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA 00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA 00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA 00003	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
...LCPFUdA 00003	06/19/18		Wellington Laboratories, Lot PFUdA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L2_00018</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00024	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_00040	50 uL	Perfluorobutyric acid	1 ng/mL
							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 (PFHpA)	1 ng/mL
							Perfluoroheptanesulfonic Acid	0.952 ng/mL
							Perfluorohexanoic acid	1 ng/mL
							Perfluorohexadecanoic acid	1 ng/mL
Perfluorohexanesulfonic acid (PFHxS)	0.946 ng/mL							
Perfluorononanoic acid (PFNA)	1 ng/mL							



REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration	
					Reagent ID	Volume Added			
							Perfluorooctanoic acid (PFOA)	1 ng/mL	
							Perfluorooctadecanoic acid	1 ng/mL	
							Perfluorooctanesulfonic acid (PFOS)	0.956 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL	
					LCM2PFTeDA 00003	0.2 mL	13C2-PFTeDA	1 ug/mL	
					LCM4PFHFA 00003	0.2 mL	13C4-PFHFA	1 ug/mL	
					LCM5PFPEA 00004	0.2 mL	13C5-PFPeA	1 ug/mL	
					LCM8FOSA 00006	0.2 mL	13C8 FOSA	1 ug/mL	
					LCMPFBA 00004	0.2 mL	13C4 PFBA	1 ug/mL	
					LCMPFDA 00004	0.2 mL	13C2 PFDA	1 ug/mL	
					LCMPFDoA 00004	0.2 mL	13C2 PFDoA	1 ug/mL	
					LCMPFHxA 00005	0.2 mL	13C2 PFHxA	1 ug/mL	
					LCMPFHxS 00004	0.2 mL	18O2 PFHxS	0.946 ug/mL	
					LCMPFNA 00003	0.2 mL	13C5 PFNA	1 ug/mL	
					LCMPFOA 00007	0.2 mL	13C4 PFOA	1 ug/mL	
					LCMPFOS 00009	0.2 mL	13C4 PFOS	0.956 ug/mL	
					LCMPFUDa 00005	0.2 mL	13C2 PFUnA	1 ug/mL	
..LCM2PFHxDA 00003	11/29/17		Wellington Laboratories, Lot M2PFHxDA1112				(Purchased Reagent)	13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA 00003	11/29/17		Wellington Laboratories, Lot M2PFTeDA1112				(Purchased Reagent)	13C2-PFTeDA	50 ug/mL
..LCM4PFHFA 00003	05/22/20		Wellington Laboratories, Lot M4PFHFA0515				(Purchased Reagent)	13C4-PFHFA	50 ug/mL
..LCM5PFPEA 00004	05/22/20		Wellington Laboratories, Lot M5PFPeA0515				(Purchased Reagent)	13C5-PFPeA	50 ug/mL
..LCM8FOSA 00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I				(Purchased Reagent)	13C8 FOSA	50 ug/mL
..LCMPFBA 00004	10/31/19		Wellington Laboratories, Lot MPFBA1014				(Purchased Reagent)	13C4 PFBA	50 ug/mL
..LCMPFDA 00004	04/13/19		Wellington Laboratories, Lot MPFDA0414				(Purchased Reagent)	13C2 PFDA	50 ug/mL
..LCMPFDoA 00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714				(Purchased Reagent)	13C2 PFDoA	50 ug/mL
..LCMPFHxA 00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414				(Purchased Reagent)	13C2 PFHxA	50 ug/mL
..LCMPFHxS 00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713				(Purchased Reagent)	18O2 PFHxS	47.3 ug/mL
..LCMPFNA 00003	04/13/19		Wellington Laboratories, Lot MPFNA0414				(Purchased Reagent)	13C5 PFNA	50 ug/mL
..LCMPFOA 00007	04/10/20		Wellington Laboratories, Lot MPFOA0415				(Purchased Reagent)	13C4 PFOA	50 ug/mL
..LCMPFOS 00009	05/15/20		Wellington Laboratories, Lot MPFOS0515				(Purchased Reagent)	13C4 PFOS	47.8 ug/mL
..LCMPFUDa 00005	10/31/19		Wellington Laboratories, Lot MPFUDa1014				(Purchased Reagent)	13C2 PFUnA	50 ug/mL
.LCPFCSP_00040	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFCSP_00039	0.5 mL	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 (PFHpA)	0.1 ug/mL	
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL	
							Perfluorohexanoic acid	0.1 ug/mL	

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.0946 ug/mL
							Perfluorononanoic acid (PFNA)	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctanesulfonic acid (PFOS)	0.0956 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_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
...LCPFBA_00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDoA_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFDSA_00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
...LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA_00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
...LCPFNA 00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
...LCPFOA 00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFODA 00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
...LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
...LCPFOSA 00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA 00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA 00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA 00003	12/10/18		Wellington Laboratories, Lot PFTTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
...LCPFUDA 00003	06/19/18		Wellington Laboratories, Lot PFUDA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L2_00019</b>	06/29/16	01/08/16	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00024	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_00040	50 uL	Perfluorobutyric acid	1 ng/mL
							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 (PFHpA)	1 ng/mL
							Perfluoroheptanesulfonic Acid	0.952 ng/mL
							Perfluorohexanoic acid	1 ng/mL
							Perfluorohexadecanoic acid	1 ng/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.946 ng/mL
							Perfluorononanoic acid (PFNA)	1 ng/mL
							Perfluorooctanoic acid (PFOA)	1 ng/mL
Perfluorooctadecanoic acid	1 ng/mL							
Perfluorooctanesulfonic acid (PFOS)	0.956 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							

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCMPFCSU_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.2 mL	13C4-PFHpA	1 ug/mL
					LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.2 mL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.2 mL	13C4 PFOS	0.956 ug/mL
LCMPFUdA_00005	0.2 mL	13C2 PFUnA	1 ug/mL					
..LCM2PFHxDA_00003	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)	13C2-PFHxDA	50 ug/mL	
..LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)	13C2-PFTeDA	50 ug/mL	
..LCM4PFHPA_00003	05/22/20	Wellington Laboratories, Lot M4PFHpa0515			(Purchased Reagent)	13C4-PFHpA	50 ug/mL	
..LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)	13C5-PFPeA	50 ug/mL	
..LCM8FOSA_00006	12/15/16	Wellington Laboratories, Lot M8FOSA1214I			(Purchased Reagent)	13C8 FOSA	50 ug/mL	
..LCMPFBA_00004	10/31/19	Wellington Laboratories, Lot MPFBA1014			(Purchased Reagent)	13C4 PFBA	50 ug/mL	
..LCMPFDA_00004	04/13/19	Wellington Laboratories, Lot MPFDA0414			(Purchased Reagent)	13C2 PFDA	50 ug/mL	
..LCMPFDoA_00004	07/17/19	Wellington Laboratories, Lot MPFDoA0714			(Purchased Reagent)	13C2 PFDoA	50 ug/mL	
..LCMPFHxA_00005	04/13/19	Wellington Laboratories, Lot MPFHxA0414			(Purchased Reagent)	13C2 PFHxA	50 ug/mL	
..LCMPFHxS_00004	07/25/18	Wellington Laboratories, Lot MPFHxS0713			(Purchased Reagent)	1802 PFHxS	47.3 ug/mL	
..LCMPFNA_00003	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)	13C5 PFNA	50 ug/mL	
..LCMPFOA_00007	04/10/20	Wellington Laboratories, Lot MPFOA0415			(Purchased Reagent)	13C4 PFOA	50 ug/mL	
..LCMPFOS_00009	05/15/20	Wellington Laboratories, Lot MPFOS0515			(Purchased Reagent)	13C4 PFOS	47.8 ug/mL	
..LCMPFUdA_00005	10/31/19	Wellington Laboratories, Lot MPFUdA1014			(Purchased Reagent)	13C2 PFUnA	50 ug/mL	
.LCPFCSP_00040	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFCSP_00039	0.5 mL	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 (PFHpA)	0.1 ug/mL
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL
							Perfluorohexanoic acid	0.1 ug/mL
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.0946 ug/mL
							Perfluorononanoic acid (PFNA)	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctanesulfonic acid (PFOS)	0.0956 ug/mL
							Perfluorooctane Sulfonamide	0.1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							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_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctandecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUDA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
...LCPFBA_00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDoA_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFDSA_00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
...LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA_00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
...LCPFNA_00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
...LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFODA_00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctandecanoic acid	50 ug/mL
...LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
...LCPFOSA_00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA_00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA_00003	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration		
					Reagent ID	Volume Added				
...LCPFuDA_00003	06/19/18		Wellington Laboratories, Lot PFUdA0613			(Purchased Reagent)	Perfluoroundecanoic acid	50 ug/mL		
<b>LCPF3-L3_00016</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00024	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		
							LCPF3SP_00040	250 uL	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 (PFHpA)	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 (PFHxS)	4.73 ng/mL				
					Perfluorononanoic acid (PFNA)	5 ng/mL				
					Perfluorooctanoic acid (PFOA)	5 ng/mL				
					Perfluorooctadecanoic acid	5 ng/mL				
Perfluorooctanesulfonic acid (PFOS)	4.78 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL		
							LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
							LCM4PFHPA_00003	0.2 mL	13C4-PFHpA	1 ug/mL
							LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
							LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
							LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
							LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
							LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
							LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCMPFHxS 00004	0.2 mL	1802 PFHxS	0.946 ug/mL
					LCMPFNA 00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA 00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS 00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa 00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA 00003	11/29/17		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA 00003	11/29/17		Wellington Laboratories, Lot M2PFTeDA1112		(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA 00003	05/22/20		Wellington Laboratories, Lot M4PFHPA0515		(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA 00004	05/22/20		Wellington Laboratories, Lot M5PFPeA0515		(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA 00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA 00004	10/31/19		Wellington Laboratories, Lot MPFBA1014		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA 00004	04/13/19		Wellington Laboratories, Lot MPFDA0414		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA 00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA 00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS 00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		1802 PFHxS	47.3 ug/mL
..LCMPFNA 00003	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA 00007	04/10/20		Wellington Laboratories, Lot MPFOA0415		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS 00009	05/15/20		Wellington Laboratories, Lot MPFOS0515		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa 00005	10/31/19		Wellington Laboratories, Lot MPFUDa1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
..LCPFCSP_00040	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFCSP_00039	0.5 mL	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 (PFHpA)	0.1 ug/mL
							Perfluoroheptanesulfonic Acid	0.0952 ug/mL
							Perfluorohexanoic acid	0.1 ug/mL
							Perfluorohexadecanoic acid	0.1 ug/mL
							Perfluorohexanesulfonic acid (PFHxS)	0.0946 ug/mL
							Perfluorononanoic acid (PFNA)	0.1 ug/mL
							Perfluorooctanoic acid (PFOA)	0.1 ug/mL
							Perfluorooctadecanoic acid	0.1 ug/mL
							Perfluorooctanesulfonic acid (PFOS)	0.0956 ug/mL
							Perfluorooctane Sulfonylamide	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_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA 00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA 00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA 00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
...LCPFBA_00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
...LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
...LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
...LCPFDaA_00003	01/03/18		Wellington Laboratories, Lot PFDaA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
...LCPFDSA_00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
...LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
...LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHps1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
...LCPFHxA_00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
...LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
...LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
...LCPFNA_00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
...LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
...LCPFODA_00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
...LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
...LCPFOSA_00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
...LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
...LCPFTeDA_00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
...LCPFTrDA_00003	12/10/18		Wellington Laboratories, Lot PFTTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
...LCPFUdA_00003	06/19/18		Wellington Laboratories, Lot PFUdA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L4_00017</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00024	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



REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							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_00039	100 uL	Perfluorobutyric acid	20 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	17.68 ng/mL
							Perfluorodecanoic acid	20 ng/mL
							Perfluorododecanoic acid	20 ng/mL
							Perfluorodecane Sulfonic acid (PFHpA)	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 (PFHxS)	18.92 ng/mL
							Perfluorononanoic acid (PFNA)	20 ng/mL
							Perfluorooctanoic acid (PFOA)	20 ng/mL
							Perfluorooctandecanoic acid	20 ng/mL
		Perfluorooctanesulfonic acid (PFOS)	19.12 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.2 mL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.2 mL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00003	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00003	05/22/20	Wellington Laboratories, Lot M4PFHpA0515			(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCM8FOSA 00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA 00004	10/31/19		Wellington Laboratories, Lot MPFBA1014		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA 00004	04/13/19		Wellington Laboratories, Lot MPFDA0414		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA 00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA 00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS 00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA 00003	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA 00007	04/10/20		Wellington Laboratories, Lot MPFOA0415		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS 00009	05/15/20		Wellington Laboratories, Lot MPFOS0515		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa 00005	10/31/19		Wellington Laboratories, Lot MPFUDa1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA 00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA 00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA 00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA 00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpa_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA 00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA 00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA 00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA 00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA 00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA 00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA 00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA 00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA 00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA 00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUda 00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA 00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA 00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA 00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDSA 00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpa_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
..LCPFHpSA 00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA 00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
..LCPFNA 00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
..LCPFOA 00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA 00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
..LCPFOSA 00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA 00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA 00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA 00003	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUDA 00003	06/19/18		Wellington Laboratories, Lot PFUDA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L5_00016</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	5 mL	LCMPFCSU_00024	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_00039	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 (PFHpA)	50 ng/mL		
					Perfluoroheptanesulfonic Acid	47.6 ng/mL		
					Perfluorohexanoic acid	50 ng/mL		
					Perfluorohexadecanoic acid	50 ng/mL		
					Perfluorohexanesulfonic acid (PFHxS)	47.3 ng/mL		
					Perfluorononanoic acid (PFNA)	50 ng/mL		
					Perfluorooctanoic acid (PFOA)	50 ng/mL		
					Perfluorooctandecanoic acid	50 ng/mL		
Perfluorooctanesulfonic acid (PFOS)	47.8 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.2 mL	13C4-PFHpA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCM5PFPEA 00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA 00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA 00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA 00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA 00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA 00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS 00004	0.2 mL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA 00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA 00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS 00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA 00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA 00003	11/29/17		Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA 00003	11/29/17		Wellington Laboratories, Lot M2PFTeDA1112		(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA 00003	05/22/20		Wellington Laboratories, Lot M4PFHPA0515		(Purchased Reagent)		13C4-PFHpA	50 ug/mL
..LCM5PFPEA 00004	05/22/20		Wellington Laboratories, Lot M5PFPeA0515		(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA 00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA 00004	10/31/19		Wellington Laboratories, Lot MPFBA1014		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA 00004	04/13/19		Wellington Laboratories, Lot MPFDA0414		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA 00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA 00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS 00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA 00003	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA 00007	04/10/20		Wellington Laboratories, Lot MPFOA0415		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS 00009	05/15/20		Wellington Laboratories, Lot MPFOS0515		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA 00005	10/31/19		Wellington Laboratories, Lot MPFUdA1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
..LCPFCSP_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA 00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA 00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA 00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA 00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA 00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA 00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA 00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA 00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA 00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA 00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA 00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPPeA 00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPPTeDA 00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPPTrDA 00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA 00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFBA_00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDSA_00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
..LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA_00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
..LCPFNA_00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
..LCPFoA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFoDA_00004	04/25/17		Wellington Laboratories, Lot PFoDA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFoSA_00004	06/20/19		Wellington Laboratories, Lot LPFoSA0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFoSA)	47.8 ug/mL
..LCPFoSA_00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA_00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00003	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFuDA_00003	06/19/18		Wellington Laboratories, Lot PFuDA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L6_00015</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	2 mL	LCMPFCSU_00024	100 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_00039	400 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 (PFHpA)	200 ng/mL
Perfluoroheptanesulfonic Acid	190.4 ng/mL							
Perfluorohexanoic acid	200 ng/mL							
Perfluorohexadecanoic acid	200 ng/mL							

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluorohexanesulfonic acid (PFHxS)	189.2 ng/mL
							Perfluorononanoic acid (PFNA)	200 ng/mL
							Perfluorooctanoic acid (PFOA)	200 ng/mL
							Perfluorooctadecanoic acid	200 ng/mL
							Perfluorooctanesulfonic acid (PFOS)	191.2 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.2 mL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.2 mL	1802 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00003	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00003	05/22/20	Wellington Laboratories, Lot M4PFHPa0515			(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00006	12/15/16	Wellington Laboratories, Lot M8FOSA1214I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00004	10/31/19	Wellington Laboratories, Lot MPFBA1014			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00004	04/13/19	Wellington Laboratories, Lot MPFDA0414			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00004	07/17/19	Wellington Laboratories, Lot MPFDoA0714			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00005	04/13/19	Wellington Laboratories, Lot MPFHxA0414			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00004	07/25/18	Wellington Laboratories, Lot MPFHxS0713			(Purchased Reagent)		1802 PFHxS	47.3 ug/mL
..LCMPFNA_00003	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00007	04/10/20	Wellington Laboratories, Lot MPFOA0415			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00009	05/15/20	Wellington Laboratories, Lot MPFOS0515			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00005	10/31/19	Wellington Laboratories, Lot MPFUDa1014			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFHpA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA_00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDSA_00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpA_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
..LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHps1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA_00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
..LCPFNA_00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
..LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
..LCPFOSA_00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA_00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00003	12/10/18		Wellington Laboratories, Lot PFTTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00003	06/19/18		Wellington Laboratories, Lot PFUdA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L7_00015</b>	06/29/16	12/30/15	MeOH/H2O, Lot 090285	2 mL	LCMPFCSU_00024	100 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

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							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_00039	800 uL	Perfluorobutyric acid	400 ng/mL
							Perfluorobutanesulfonic acid (PFBS)	353.6 ng/mL
							Perfluorodecanoic acid	400 ng/mL
							Perfluorododecanoic acid	400 ng/mL
							Perfluorodecane Sulfonic acid (PFHpA)	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 (PFHxS)	378.4 ng/mL
							Perfluorononanoic acid (PFNA)	400 ng/mL
							Perfluorooctanoic acid (PFOA)	400 ng/mL
							Perfluorooctandecanoic acid	400 ng/mL
		Perfluorooctanesulfonic acid (PFOS)	382.4 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_00024	06/29/16	12/29/15	Methanol, Lot Baker 115491	10 mL	LCM2PFHxDA_00003	0.2 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.2 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.2 mL	13C4-PFHpa	1 ug/mL
					LCM5PFPEA_00004	0.2 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.2 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00004	0.2 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00004	0.2 mL	13C2 PFDoA	1 ug/mL
					LCMPFHxA_00005	0.2 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.2 mL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.2 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.2 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.2 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00005	0.2 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00003	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00003	05/22/20	Wellington Laboratories, Lot M4PFHpA0515			(Purchased Reagent)		13C4-PFHpa	50 ug/mL
..LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPeA0515			(Purchased Reagent)		13C5-PFPeA	50 ug/mL



REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCM8FOSA 00006	12/15/16		Wellington Laboratories, Lot M8FOSA1214I		(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA 00004	10/31/19		Wellington Laboratories, Lot MPFBA1014		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA 00004	04/13/19		Wellington Laboratories, Lot MPFDA0414		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA 00004	07/17/19		Wellington Laboratories, Lot MPFDoA0714		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA 00005	04/13/19		Wellington Laboratories, Lot MPFHxA0414		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS 00004	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA 00003	04/13/19		Wellington Laboratories, Lot MPFNA0414		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA 00007	04/10/20		Wellington Laboratories, Lot MPFOA0415		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS 00009	05/15/20		Wellington Laboratories, Lot MPFOS0515		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa 00005	10/31/19		Wellington Laboratories, Lot MPFUDa1014		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHpa_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUda_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA 00003	03/05/18		Wellington Laboratories, Lot PFBA0313		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBSA_00001	10/09/19		Wellington Laboratories, Lot LPFBS1014		(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
..LCPFDA 00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDoA 00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDSA 00001	09/13/18		Wellington Laboratories, Lot LPFDS0913		(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
..LCPFHpa_00004	05/09/19		Wellington Laboratories, Lot PFHpA0514		(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
..LCPFHpSA_00001	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
..LCPFHxA 00003	05/09/19		Wellington Laboratories, Lot PFHxA0514		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA 00004	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxSA_00001	05/09/19		Wellington Laboratories, Lot LPFHxS0514		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
..LCPFNA 00004	05/09/19		Wellington Laboratories, Lot PFNA0514		(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
..LCPFOA 00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA 00004	04/25/17		Wellington Laboratories, Lot PFODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFOS_00004	06/20/19		Wellington Laboratories, Lot LPFOS0614		(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
..LCPFOSA 00005	07/31/18		Wellington Laboratories, Lot FOSA0714I		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA 00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFTeDA 00003	06/19/18		Wellington Laboratories, Lot PFTeDA0613		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA 00003	12/10/18		Wellington Laboratories, Lot PFTrDA1213		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUDA 00003	06/19/18		Wellington Laboratories, Lot PFUDA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFCIC_00014</b>	03/15/16	09/15/15	MeOH/H2O, Lot 09285	5 mL	LCMPFCSU_00018	250 uL	18O2 PFHxS	47.3 ng/mL
					LCPFACMXB_00006	125 uL	Perfluorohexanesulfonic acid (PFHxS)	47.25 ng/mL
.LCMPFCSU_00018	03/15/16	09/15/15	Methanol, Lot Fisher 153635	10 mL	LCMPFHxS_00003	0.2 mL	18O2 PFHxS	0.946 ug/mL
..LCMPFHxS_00003	07/25/18		Wellington Laboratories, Lot MPFHxS0713		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
.LCPFACMXB_00006	01/08/18		Wellington Laboratories, Lot PFACMXB0312		(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	1.89 ug/mL
<b>LCPFCIC_00016</b>	06/16/16	12/22/15	MeOH/H2O, Lot 09285	5 mL	LCMPFCSU_00023	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
					LCPFACMXB_00008	125 uL	Perfluorobutanesulfonic acid (PFBS)	44.25 ng/mL
							Perfluoroheptanoic acid (PFHpA)	50 ng/mL
							Perfluorohexanesulfonic acid (PFHxS)	47.25 ng/mL
							Perfluorononanoic acid (PFNA)	50 ng/mL
							Perfluorooctanesulfonic acid (PFOS)	47.75 ng/mL
							Perfluorooctanoic acid (PFOA)	50 ng/mL
.LCMPFCSU_00023	06/21/16	12/21/15	Methanol, Lot Baker 115491	5 mL	LCM2PFHxDA_00002	0.1 mL	13C2-PFHxDA	1 ug/mL
					LCM2PFTeDA_00003	0.1 mL	13C2-PFTeDA	1 ug/mL
					LCM4PFHPA_00003	0.1 mL	13C4-PFHpA	1 ug/mL
					LCM5PFPEA_00004	0.1 mL	13C5-PFPeA	1 ug/mL
					LCM8FOSA_00006	0.1 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00004	0.1 mL	13C4 PFBA	1 ug/mL
					LCMPFDA_00005	0.1 mL	13C2 PFDA	1 ug/mL
					LCMPFDoA_00003	0.1 mL	13C2 PFDoA	1 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCMPFHxA_00006	0.1 mL	13C2 PFHxA	1 ug/mL
					LCMPFHxS_00004	0.1 mL	18O2 PFHxS	0.946 ug/mL
					LCMPFNA_00003	0.1 mL	13C5 PFNA	1 ug/mL
					LCMPFOA_00007	0.1 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00009	0.1 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00004	0.1 mL	13C2 PFUnA	1 ug/mL
..LCM2PFHxDA_00002	11/29/17	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00003	11/29/17	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHFA_00003	05/22/20	Wellington Laboratories, Lot M4PFHFA0515			(Purchased Reagent)		13C4-PFHFA	50 ug/mL
..LCM5PFPEA_00004	05/22/20	Wellington Laboratories, Lot M5PFPEA0515			(Purchased Reagent)		13C5-PFPEA	50 ug/mL
..LCM8FOSA_00006	12/15/16	Wellington Laboratories, Lot M8FOSA1214I			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00004	10/31/19	Wellington Laboratories, Lot MPFBA1014			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00005	04/13/19	Wellington Laboratories, Lot MPFDA0414			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDoA_00003	07/17/19	Wellington Laboratories, Lot MPFDoA0714			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00006	04/13/19	Wellington Laboratories, Lot MPFHxA0414			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00004	07/25/18	Wellington Laboratories, Lot MPFHxS0713			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00003	04/13/19	Wellington Laboratories, Lot MPFNA0414			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00007	04/10/20	Wellington Laboratories, Lot MPFOA0415			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00009	05/15/20	Wellington Laboratories, Lot MPFOS0515			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00004	10/31/19	Wellington Laboratories, Lot MPFUdA1014			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFACMXB_00008	06/20/19	Wellington Laboratories, Lot PFACMXB0614			(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	1.77 ug/mL
							Perfluoroheptanoic acid (PFHpA)	2 ug/mL
							Perfluorohexanesulfonic acid (PFHxS)	1.89 ug/mL
							Perfluorononanoic acid (PFNA)	2 ug/mL
							Perfluorooctanesulfonic acid (PFOS)	1.91 ug/mL
							Perfluorooctanoic acid (PFOA)	2 ug/mL
LCPFCSP_00039	06/30/16	12/30/15	Methanol, Lot 090285	5 mL	LCPFBA_00003	0.1 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00003	0.1 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFBSA_00001	0.1 mL	Perfluorobutanesulfonic acid (PFBS)	0.884 ug/mL
					LCPFDA_00003	0.1 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDoA_00003	0.1 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00003	0.1 mL	PFDoS (Perfluoro-1-dodecanesulfonate)	0.968 ug/mL
					LCPFDS_00003	0.1 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFDSA_00001	0.1 mL	Perfluorodecane Sulfonic acid	0.964 ug/mL
					LCPFHFA_00004	0.1 mL	Perfluoroheptanoic acid (PFHpA)	1 ug/mL
					LCPFHpS_00005	0.1 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHpSA_00001	0.1 mL	Perfluoroheptanesulfonic Acid	0.952 ug/mL
					LCPFHxA_00003	0.1 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00004	0.1 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00003	0.1 mL	Perfluorohexane Sulfonate	0.946 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFHxSA_00001	0.1 mL	Perfluorohexanesulfonic acid (PFHxS)	0.946 ug/mL
					LCPFNA_00004	0.1 mL	Perfluorononanoic acid (PFNA)	1 ug/mL
					LCPFNS_00002	0.1 mL	PFNS (Perfluoro-1-nonanesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.1 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00004	0.1 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00004	0.1 mL	Perfluorooctanesulfonic acid (PFOS)	0.956 ug/mL
					LCPFOSA_00005	0.1 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.1 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00002	0.1 mL	PFPeS (Perfluoro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00003	0.1 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00003	0.1 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00003	0.1 mL	Perfluoroundecanoic acid	1 ug/mL
.LCPFBA_00003	03/05/18	Wellington Laboratories, Lot PFBA0313			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
.LCPFBS_00003	10/09/19	Wellington Laboratories, Lot LPPFBS1014			(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
.LCPFBSA_00001	10/09/19	Wellington Laboratories, Lot LPPFBS1014			(Purchased Reagent)		Perfluorobutanesulfonic acid (PFBS)	44.2 ug/mL
.LCPFDA_00003	06/18/18	Wellington Laboratories, Lot PFDA0613			(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
.LCPFDoA_00003	01/03/18	Wellington Laboratories, Lot PFDoA0113			(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
.LCPFDoS_00003	10/06/16	Wellington Laboratories, Lot LPPFDoS1011			(Purchased Reagent)		PFDoS (Perfluoro-1-dodecanesulfonate)	48.4 ug/mL
.LCPFDS_00003	09/13/18	Wellington Laboratories, Lot LPPFDS0913			(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
.LCPFDSA_00001	09/13/18	Wellington Laboratories, Lot LPPFDS0913			(Purchased Reagent)		Perfluorodecane Sulfonic acid	48.2 ug/mL
.LCPFHpA_00004	05/09/19	Wellington Laboratories, Lot PFHpA0514			(Purchased Reagent)		Perfluoroheptanoic acid (PFHpA)	50 ug/mL
.LCPFHpS_00005	01/28/19	Wellington Laboratories, Lot LPPFHpS0114			(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
.LCPFHpSA_00001	11/21/17	Wellington Laboratories, Lot LPPFHpS1112			(Purchased Reagent)		Perfluoroheptanesulfonic Acid	47.6 ug/mL
.LCPFHxA_00003	05/09/19	Wellington Laboratories, Lot PFHxA0514			(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
.LCPFHxDA_00004	11/28/17	Wellington Laboratories, Lot PFHxDA0707			(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
.LCPFHxS_00003	05/09/19	Wellington Laboratories, Lot LPPFHxS0514			(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
.LCPFHxSA_00001	05/09/19	Wellington Laboratories, Lot LPPFHxS0514			(Purchased Reagent)		Perfluorohexanesulfonic acid (PFHxS)	47.3 ug/mL
.LCPFNA_00004	05/09/19	Wellington Laboratories, Lot PFNA0514			(Purchased Reagent)		Perfluorononanoic acid (PFNA)	50 ug/mL
.LCPFNS_00002	07/04/17	Wellington Laboratories, Lot LPPFNS0712			(Purchased Reagent)		PFNS (Perfluoro-1-nonanesulfonate)	48 ug/mL
.LCPFOA_00004	10/11/18	Wellington Laboratories, Lot PFOA1013			(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
.LCPFODA_00004	04/25/17	Wellington Laboratories, Lot PFODA0807			(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
.LCPFOS_00004	06/20/19	Wellington Laboratories, Lot LPPFOS0614			(Purchased Reagent)		Perfluorooctanesulfonic acid (PFOS)	47.8 ug/mL
.LCPFOSA_00005	07/31/18	Wellington Laboratories, Lot FOSA0714I			(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
.LCPFPeA_00003	01/03/18	Wellington Laboratories, Lot PFPeA0113			(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
.LCPFPeS_00002	07/04/17	Wellington Laboratories, Lot LPPPeS0712			(Purchased Reagent)		PFPeS (Perfluoro-1-pentanesulfonate)	46.9 ug/mL
.LCPFTeDA_00003	06/19/18	Wellington Laboratories, Lot PFTeDA0613			(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
.LCPFTrDA_00003	12/10/18	Wellington Laboratories, Lot PFTrDA1213			(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCPFUdA_00003	06/19/18		Wellington Laboratories, Lot PFUdA0613		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL

Reagent

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**LCM2PFHxDA\_00002**

Rec: 8/14/14 SKV

318141  
ID: LCM2PFHxDA\_00002  
Exp: 11/29/17 Prod: SKV  
13C2-PFHxDA at 50ug/ml

Scanned: 8/18/14 SKV

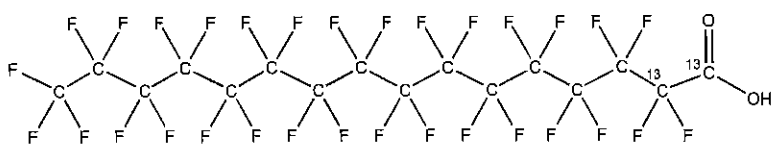


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** M2PFHxDA      **LOT NUMBER:** M2PFHxDA1112  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]hexadecanoic acid

**STRUCTURE:**      **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>14</sub>HF<sub>31</sub>O<sub>2</sub>      **MOLECULAR WEIGHT:** 816.11  
**CONCENTRATION:** 50 ± 2.5 µg/ml      **SOLVENT(S):** Methanol  
Water (<1%)  
**CHEMICAL PURITY:** >98%      **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
(1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 11/29/2012  
**EXPIRY DATE:** (mm/dd/yyyy) 11/29/2017  
**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:** 01/10/2013  
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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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:**

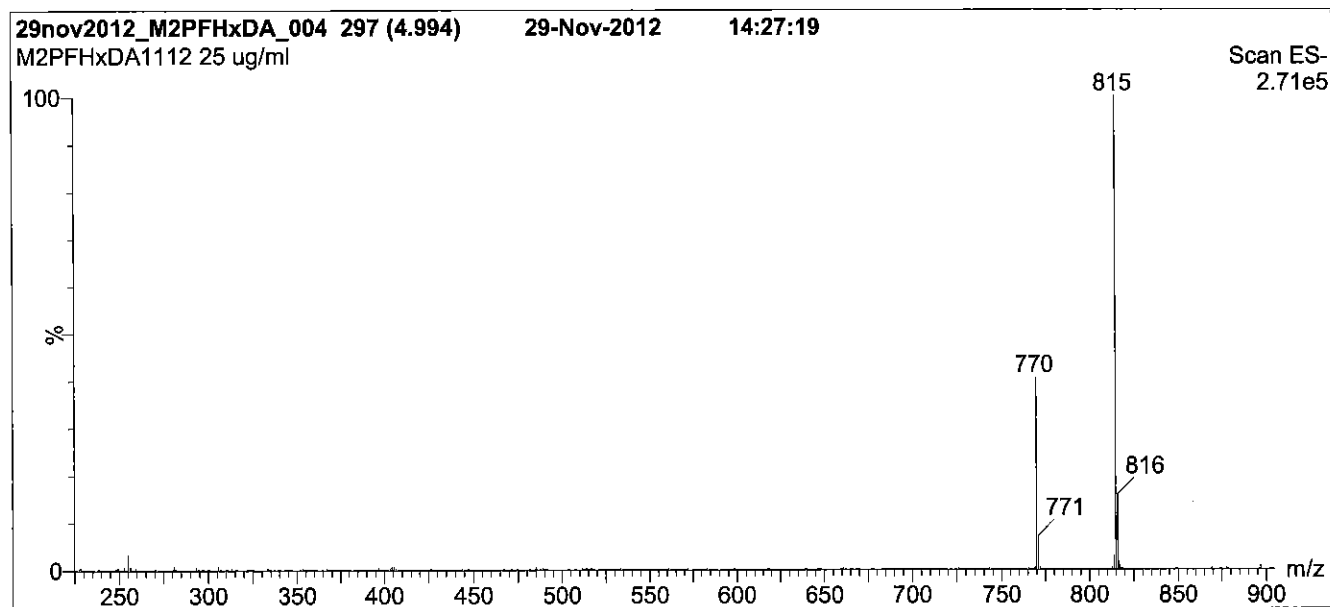
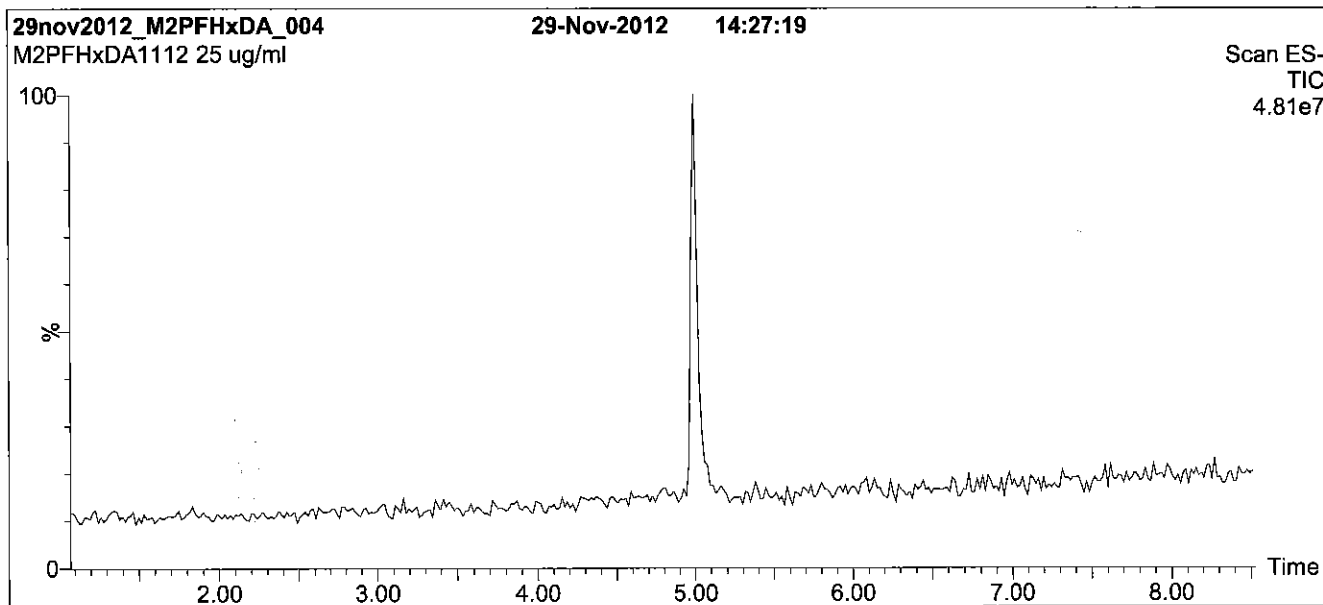
This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number AR-1523).



\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*



**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 RP<sub>18</sub>  
1.7 μm, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

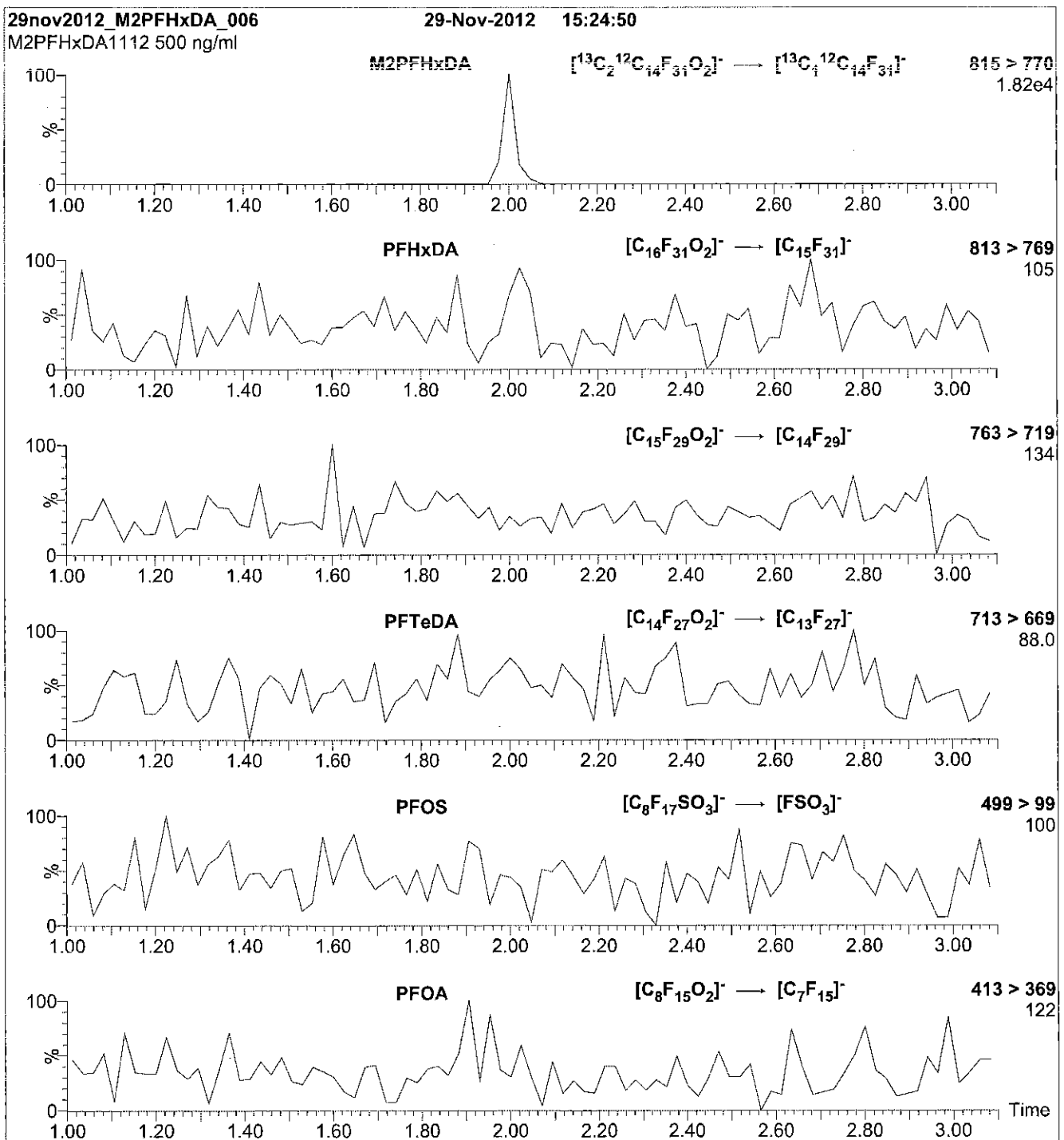
**Flow:** 300 μl/min

**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  $\mu\text{l}$  (500 ng/ml M2PFHxDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

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

Reagent

---

**LCM2PFHxDA\_00003**

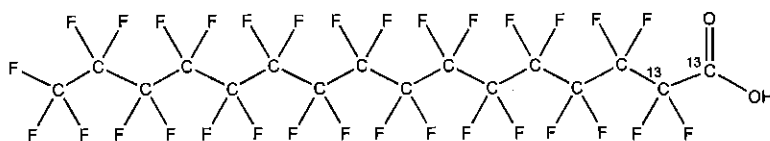


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** M2PFHxDA      **LOT NUMBER:** M2PFHxDA1112  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]hexadecanoic acid

**STRUCTURE:**      **CAS #:** Not available



<b>MOLECULAR FORMULA:</b>	<sup>13</sup> C <sub>2</sub> <sup>12</sup> C <sub>14</sub> HF <sub>31</sub> O <sub>2</sub>	<b>MOLECULAR WEIGHT:</b>	816.11
<b>CONCENTRATION:</b>	50 ± 2.5 µg/ml	<b>SOLVENT(S):</b>	Methanol Water (<1%)
<b>CHEMICAL PURITY:</b>	>98%	<b>ISOTOPIC PURITY:</b>	≥99% <sup>13</sup> C (1,2- <sup>13</sup> C <sub>2</sub> )
<b>LAST TESTED:</b> (mm/dd/yyyy)	11/29/2012		
<b>EXPIRY DATE:</b> (mm/dd/yyyy)	11/29/2017		
<b>RECOMMENDED STORAGE:</b>	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:** 04/01/2015  
(mm/dd/yyyy)  
 B.G. Chittim

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:

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

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.

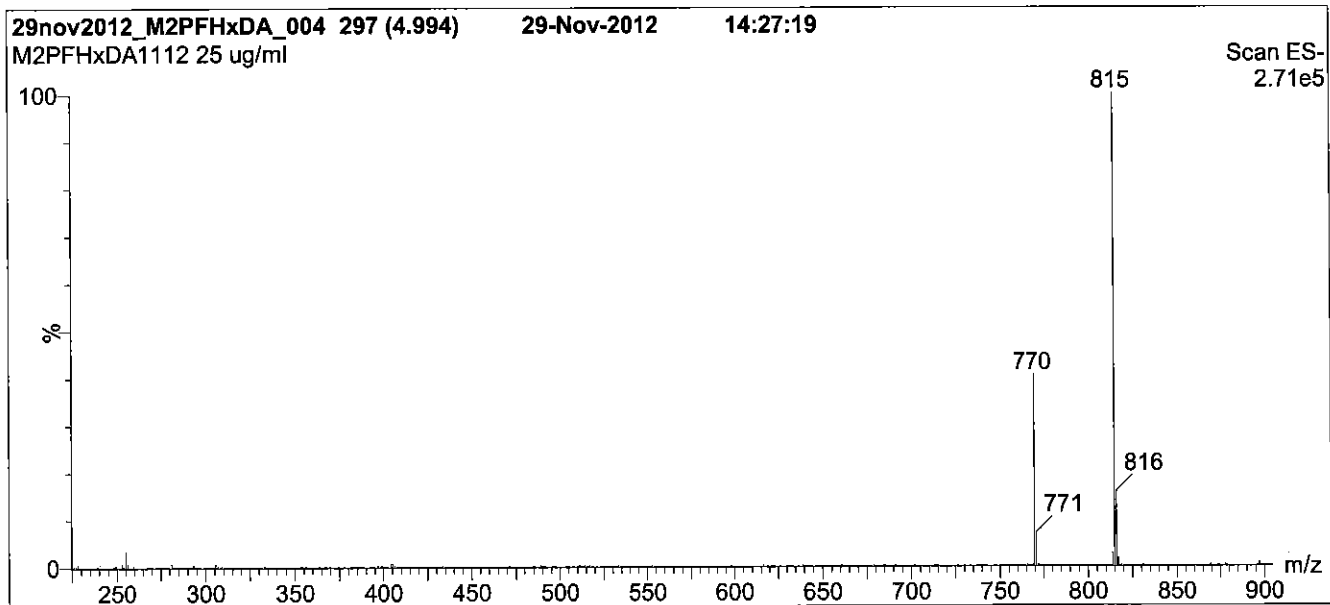
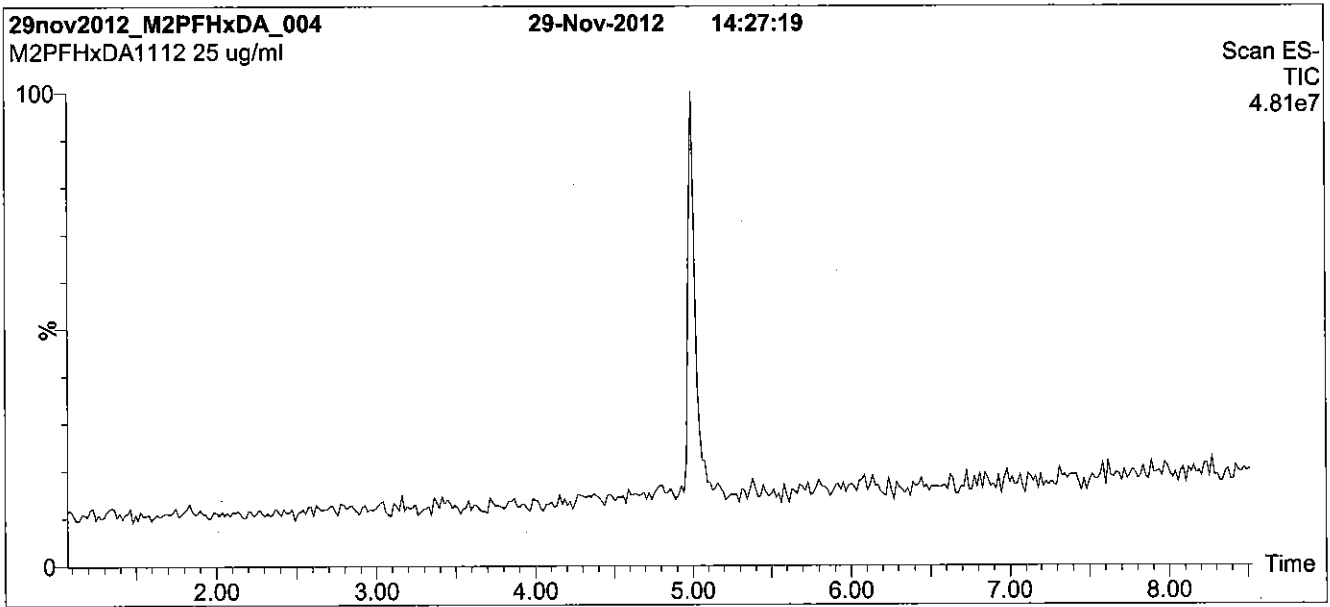
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

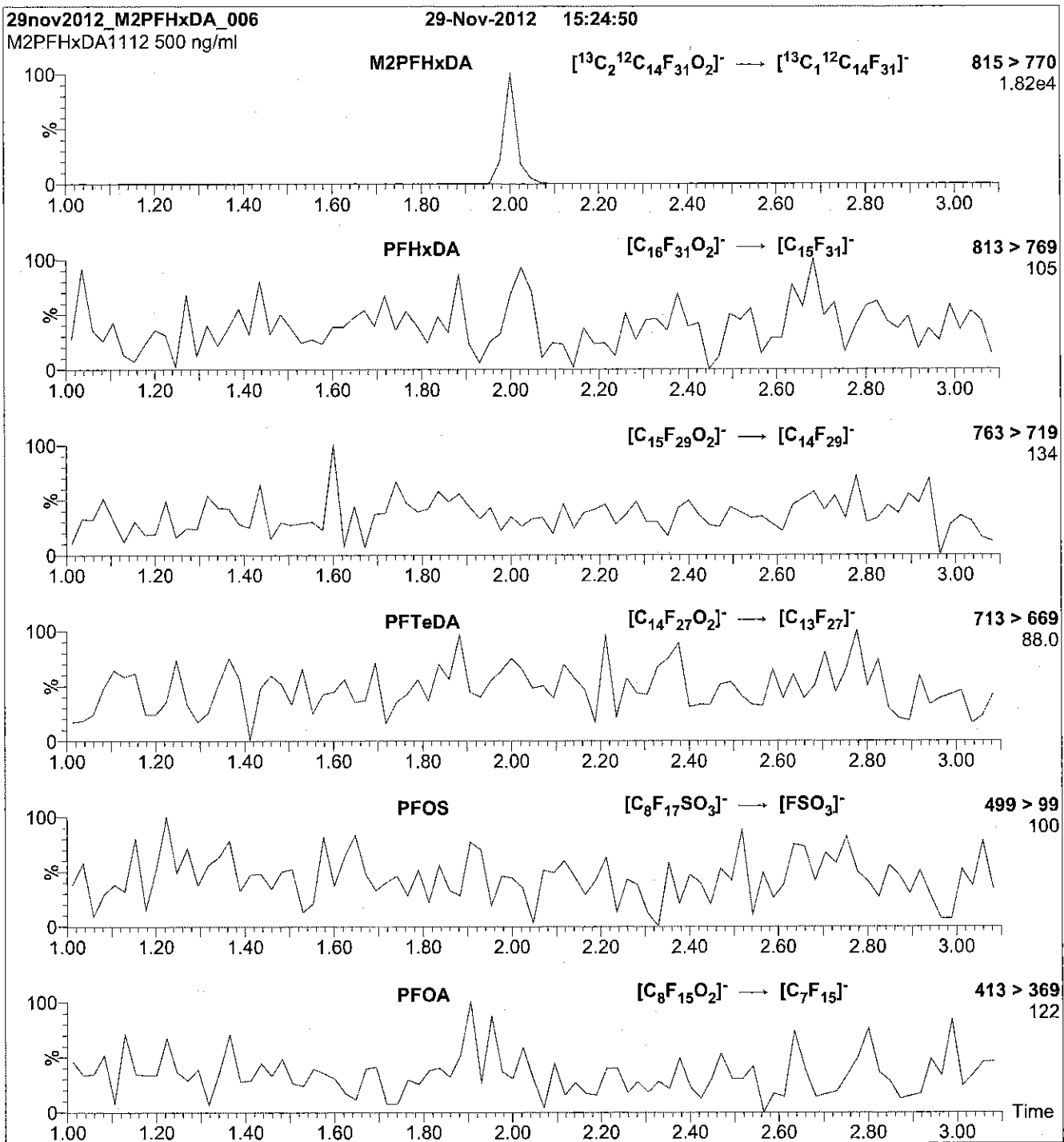
**Flow:** 300  $\mu$ l/min

**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  $\mu\text{l}$  (500 ng/ml M2PFHxDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

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

Reagent

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**LCM2PFTeDA\_00003**



r: 2/1/15 Stv

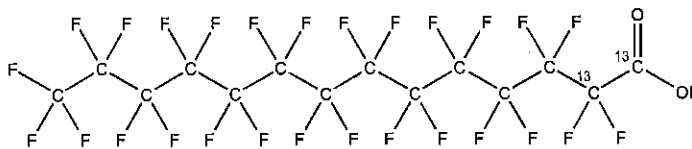


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** M2PFTeDA **LOT NUMBER:** M2PFTeDA1112  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]tetradecanoic acid

**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>12</sub>HF<sub>27</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 716.10  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
(1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 11/29/2012  
**EXPIRY DATE:** (mm/dd/yyyy) 11/29/2017  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

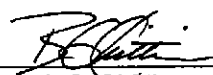
**DOCUMENTATION/ DATA ATTACHED:**

- 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:**

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### **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.

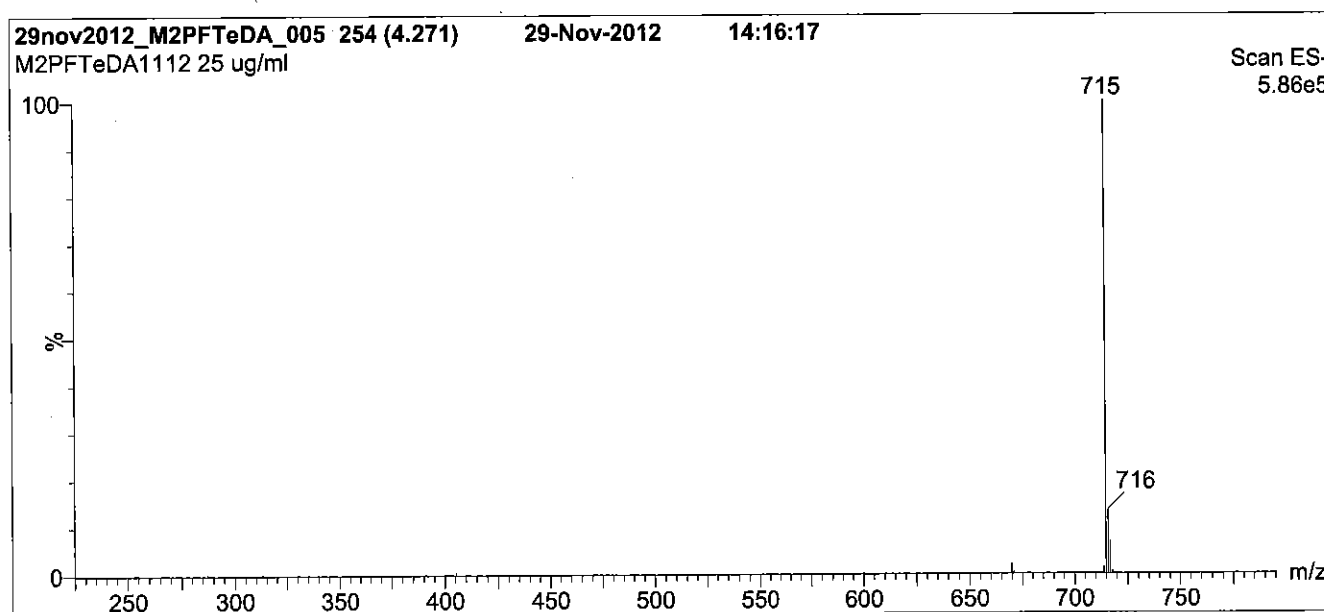
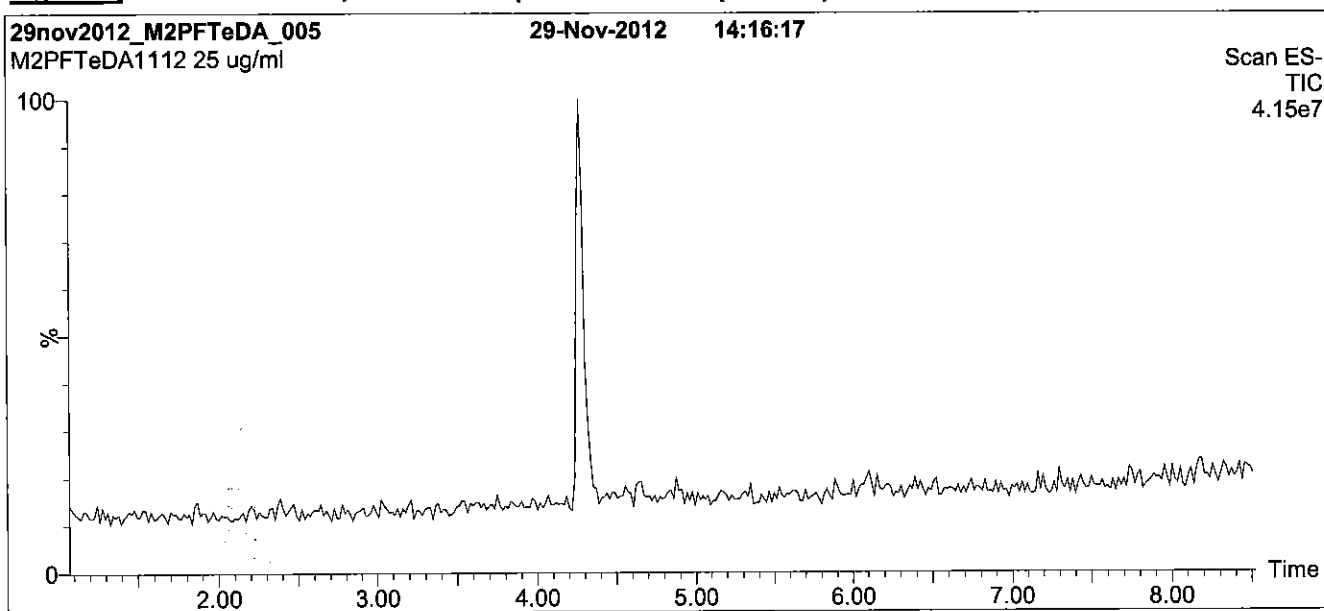
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

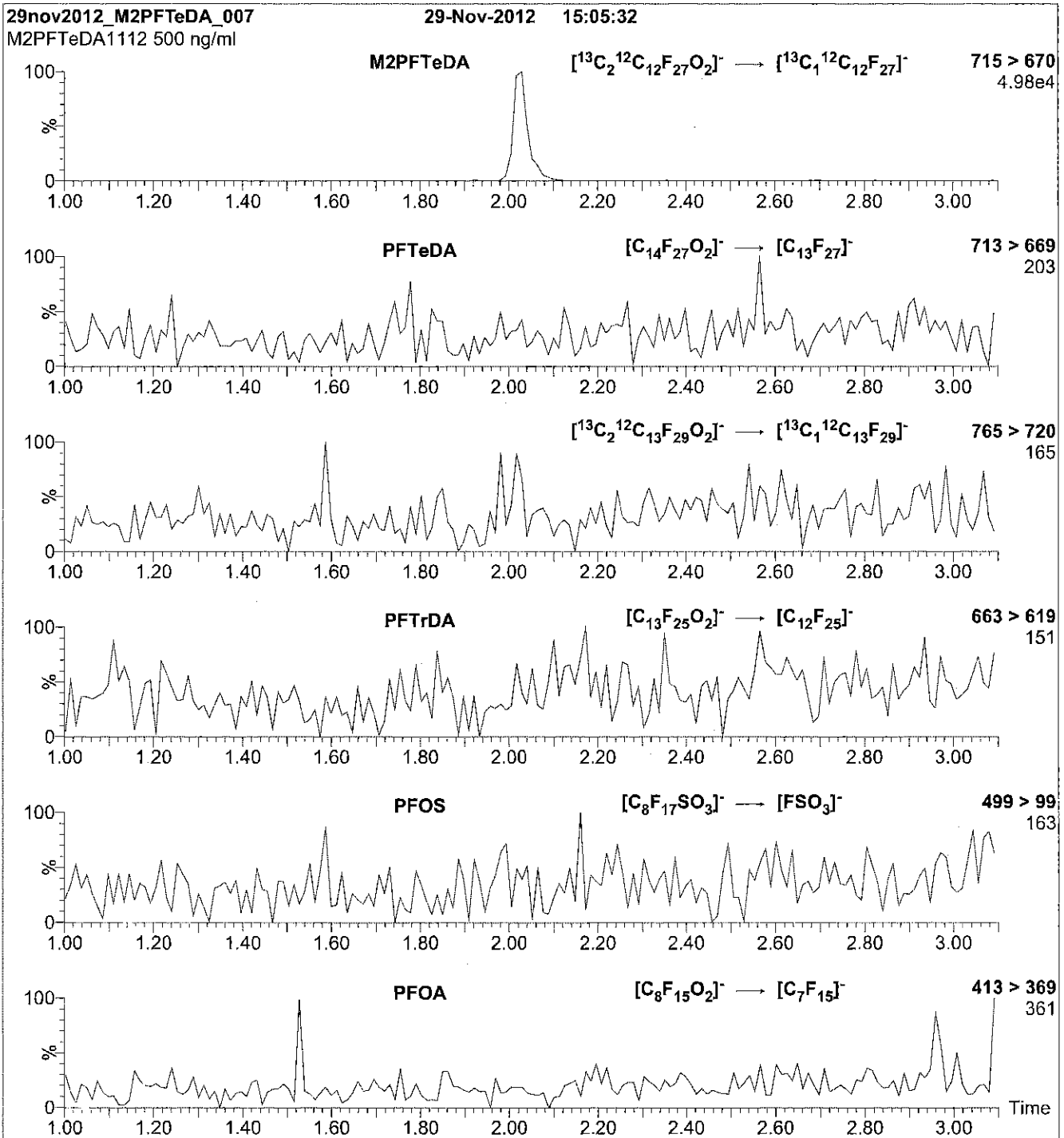
**Flow:** 300  $\mu$ l/min

**MS Parameters**

**Experiment:** Full Scan (225 - 1200 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: M2PFTeDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml M2PFTeDA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.66e-3  
Collision Energy (eV) = 14

Reagent

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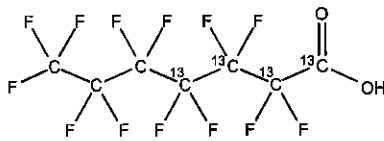
**LCM4PFHPA\_00003**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** M4PFHpA **LOT NUMBER:** M4PFHpA0515  
**COMPOUND:** Perfluoro-n-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]heptanoic acid  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>4</sub><sup>12</sup>C<sub>3</sub>HF<sub>13</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 368.03  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99%<sup>13</sup>C  
 (1,2,3,4-<sup>13</sup>C<sub>4</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 05/22/2015  
**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.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

Certified By: \_\_\_\_\_

B.G. Chittim

Date: 05/25/2015

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

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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:**

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.

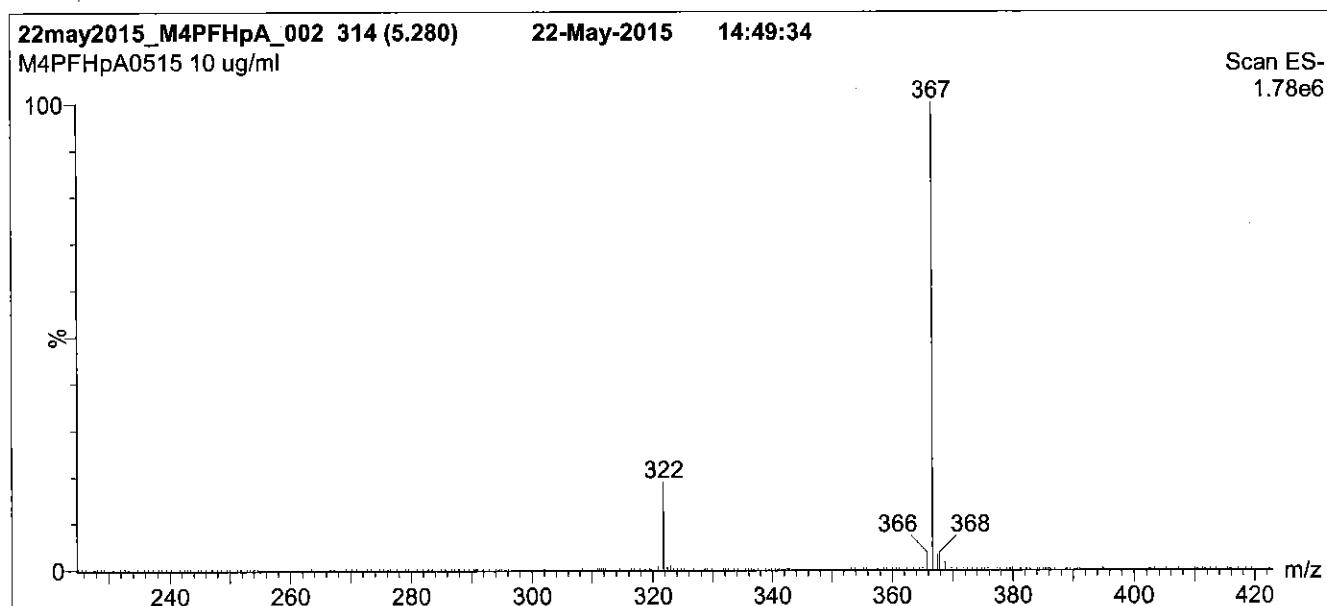
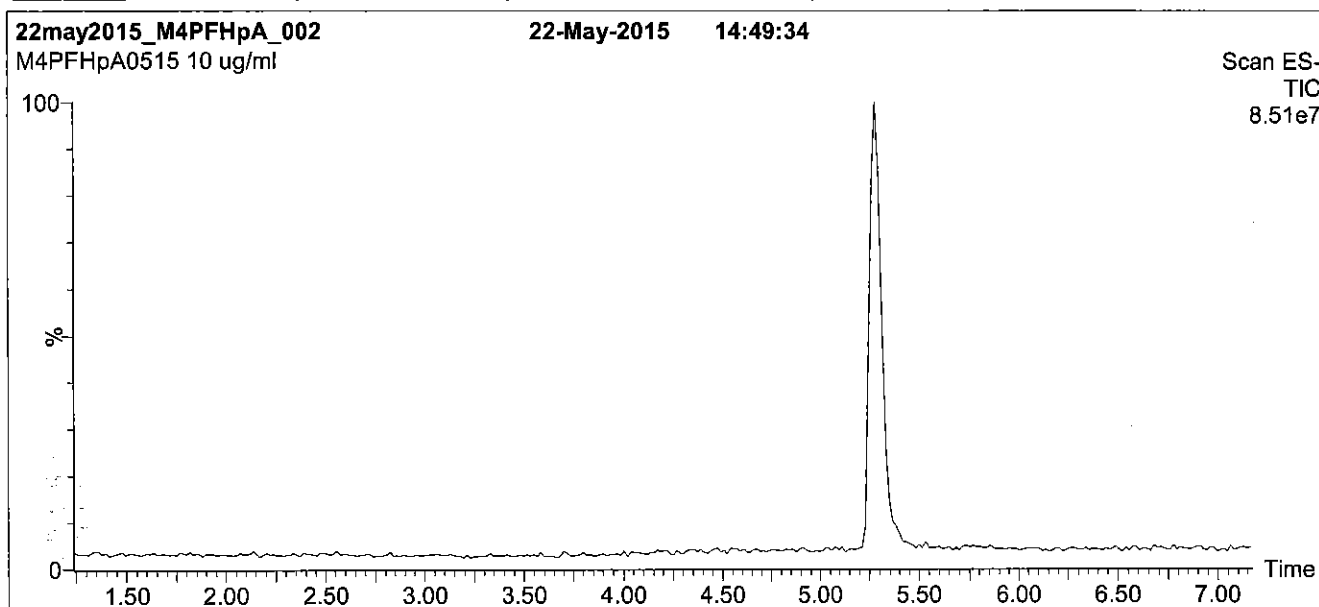
### **QUALITY MANAGEMENT:**

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: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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  $\mu$ l/min

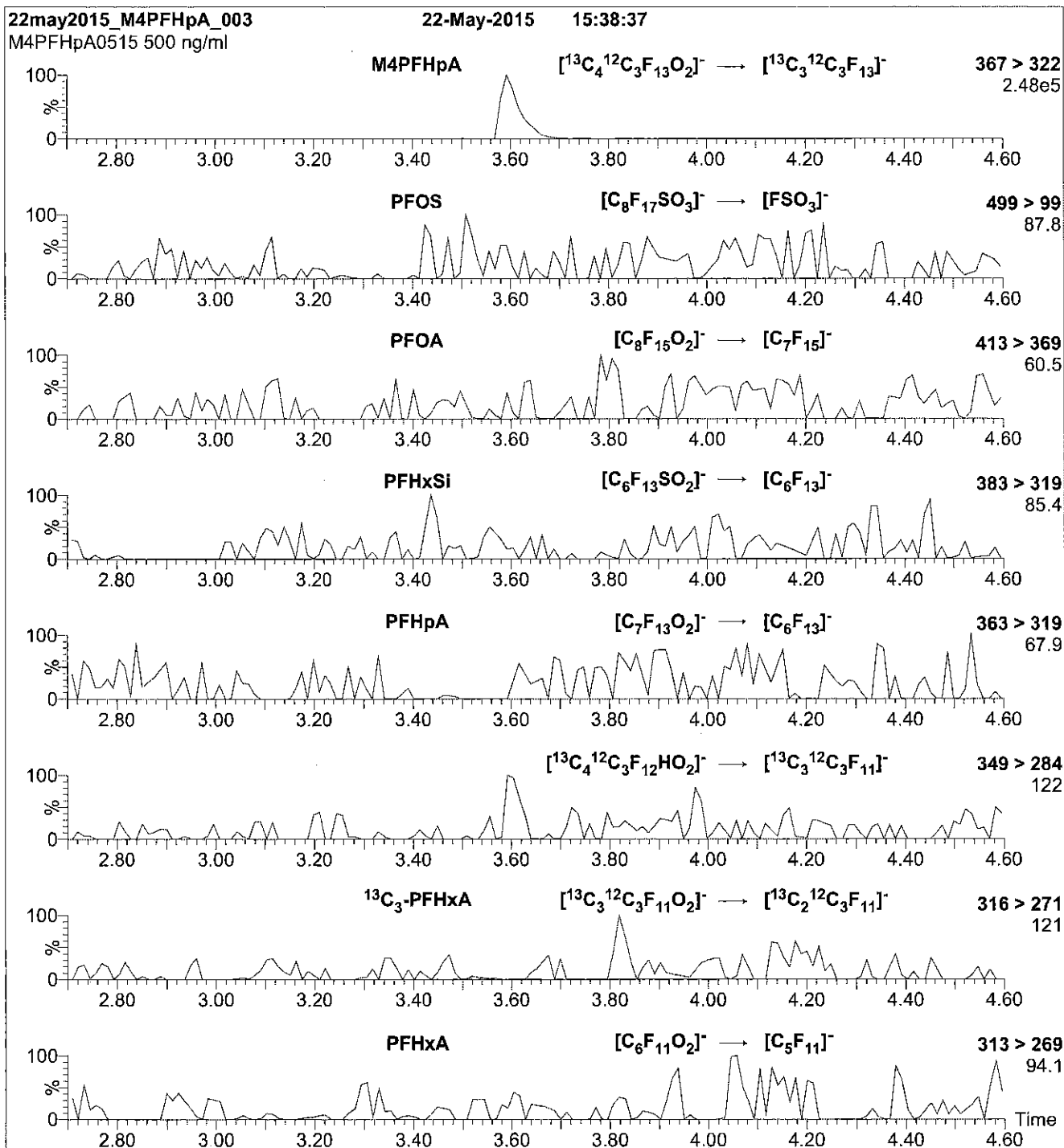
**MS Parameters**

**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: 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<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300 µl/min

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
 Collision Energy (eV) = 11

Reagent

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**LCM5PFPEA\_00004**



### **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.

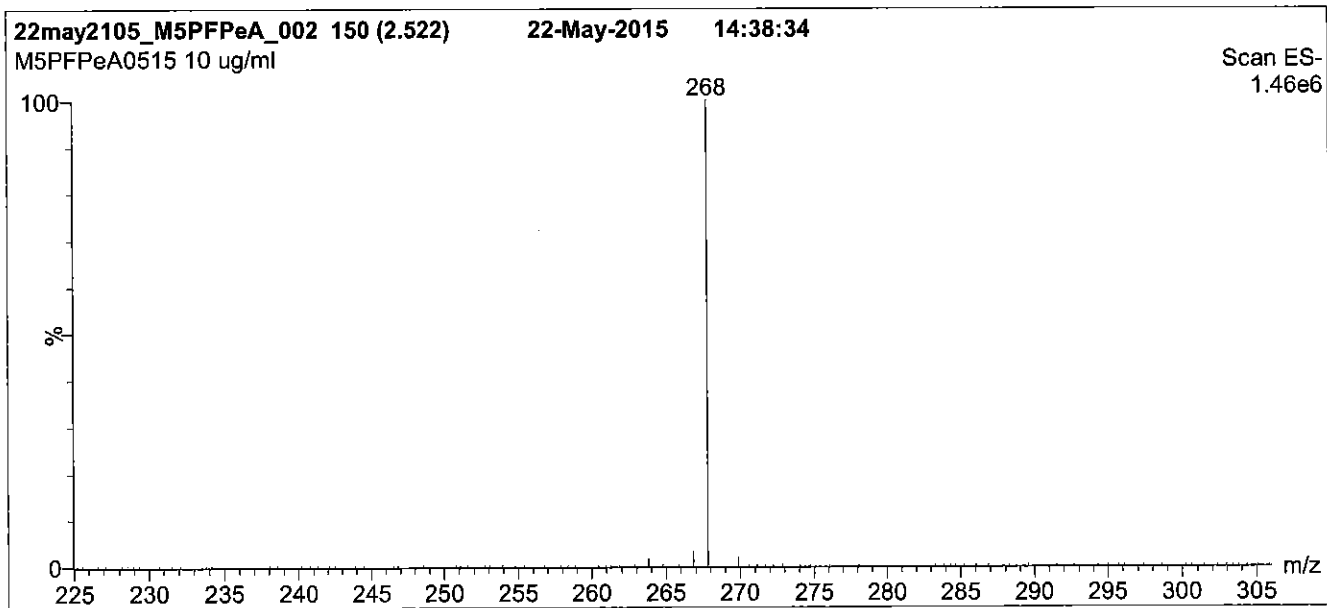
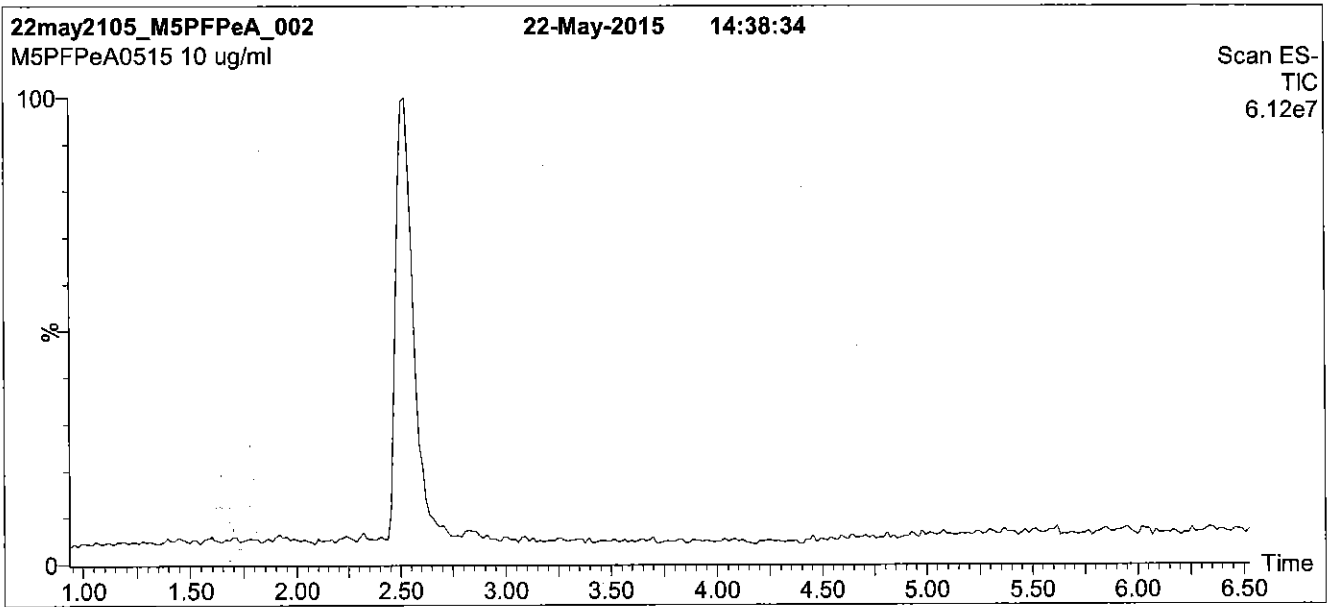
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto: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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

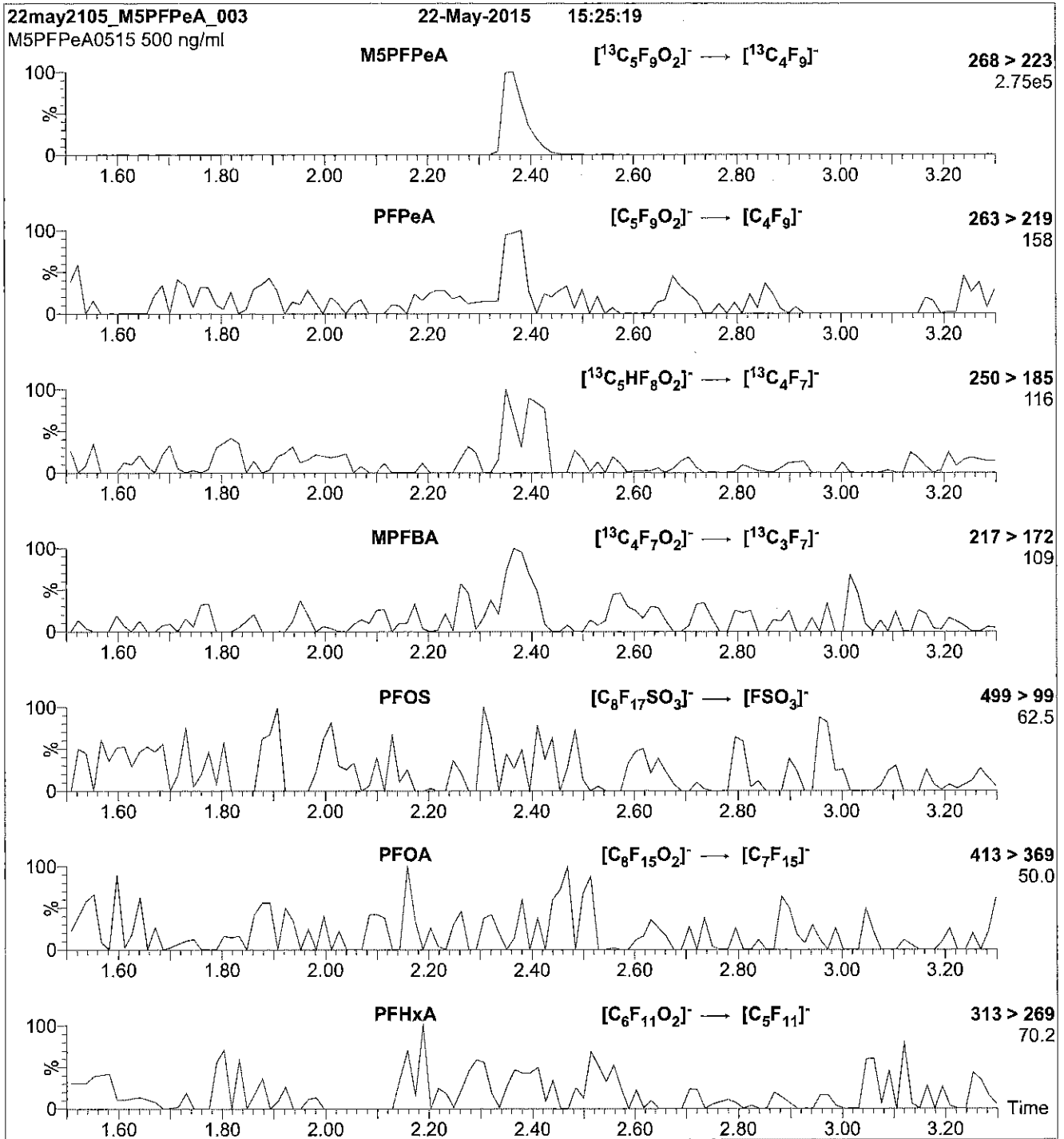
Flow: 300  $\mu$ l/min

**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  $\mu\text{l}$  (500 ng/ml M5PFPeA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 9

Reagent

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**LCM8FOSA\_00006**

rec: 9/15/15 sv



# WELLINGTON LABORATORIES

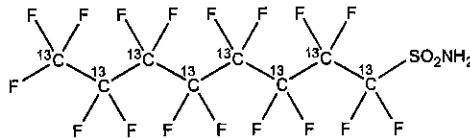
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** M8FOSA-I  
**COMPOUND:** Perfluoro-1-<sup>13</sup>C<sub>8</sub>octanesulfonamide

**LOT NUMBER:** M8FOSA1214I

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>8</sub>H<sub>2</sub>F<sub>17</sub>NO<sub>2</sub>S  
**CONCENTRATION:** 50 ± 2.5 µg/ml  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 12/15/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 12/15/2016  
**RECOMMENDED STORAGE:** Refrigerate ampoule

**MOLECULAR WEIGHT:** 507.09  
**SOLVENT(S):** Isopropanol  
**ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
(<sup>13</sup>C<sub>8</sub>)

**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:**

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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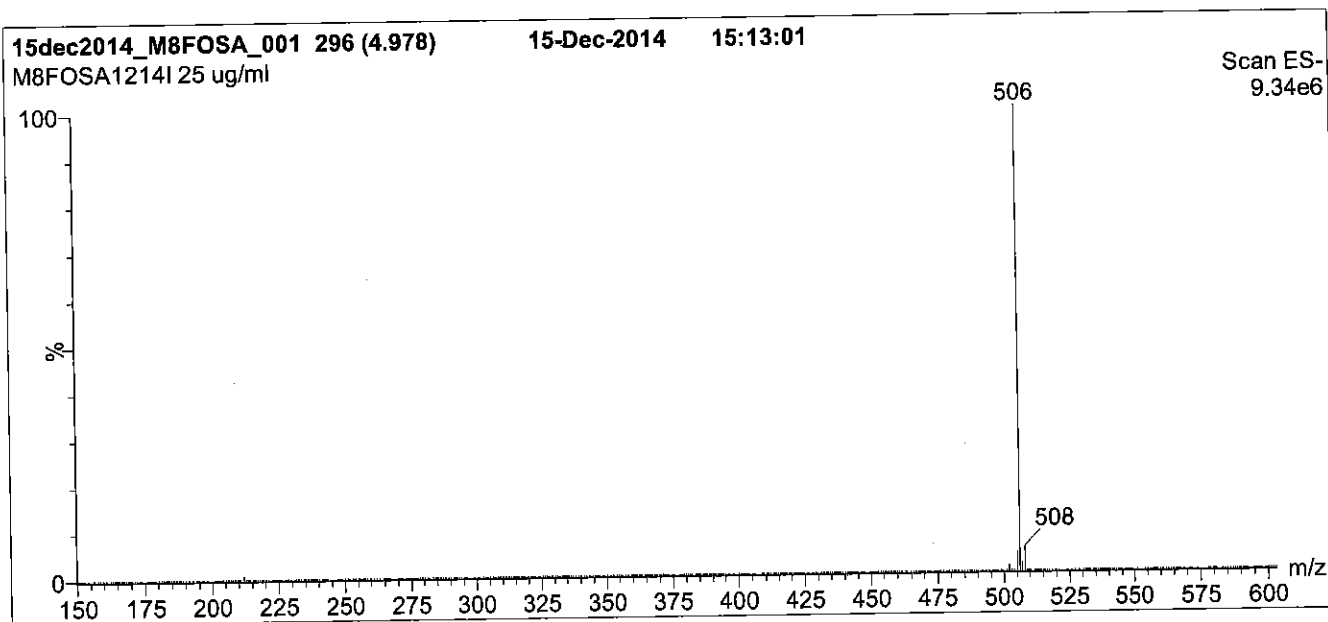
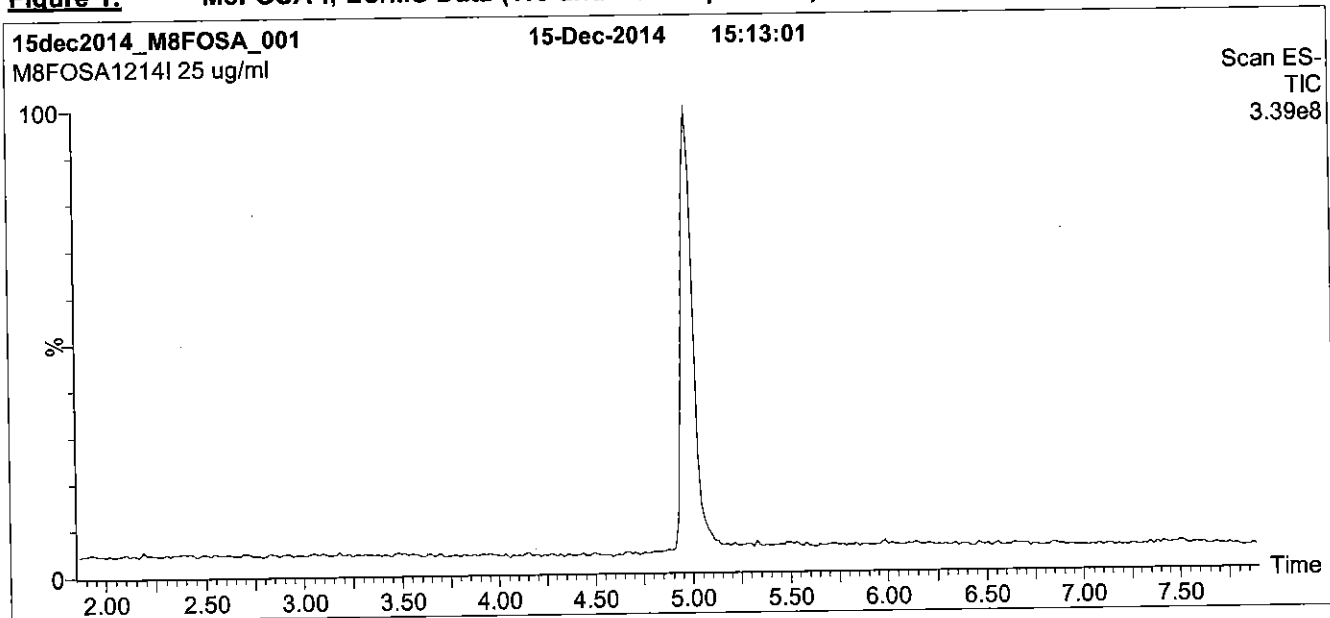
### **QUALITY MANAGEMENT:**

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: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

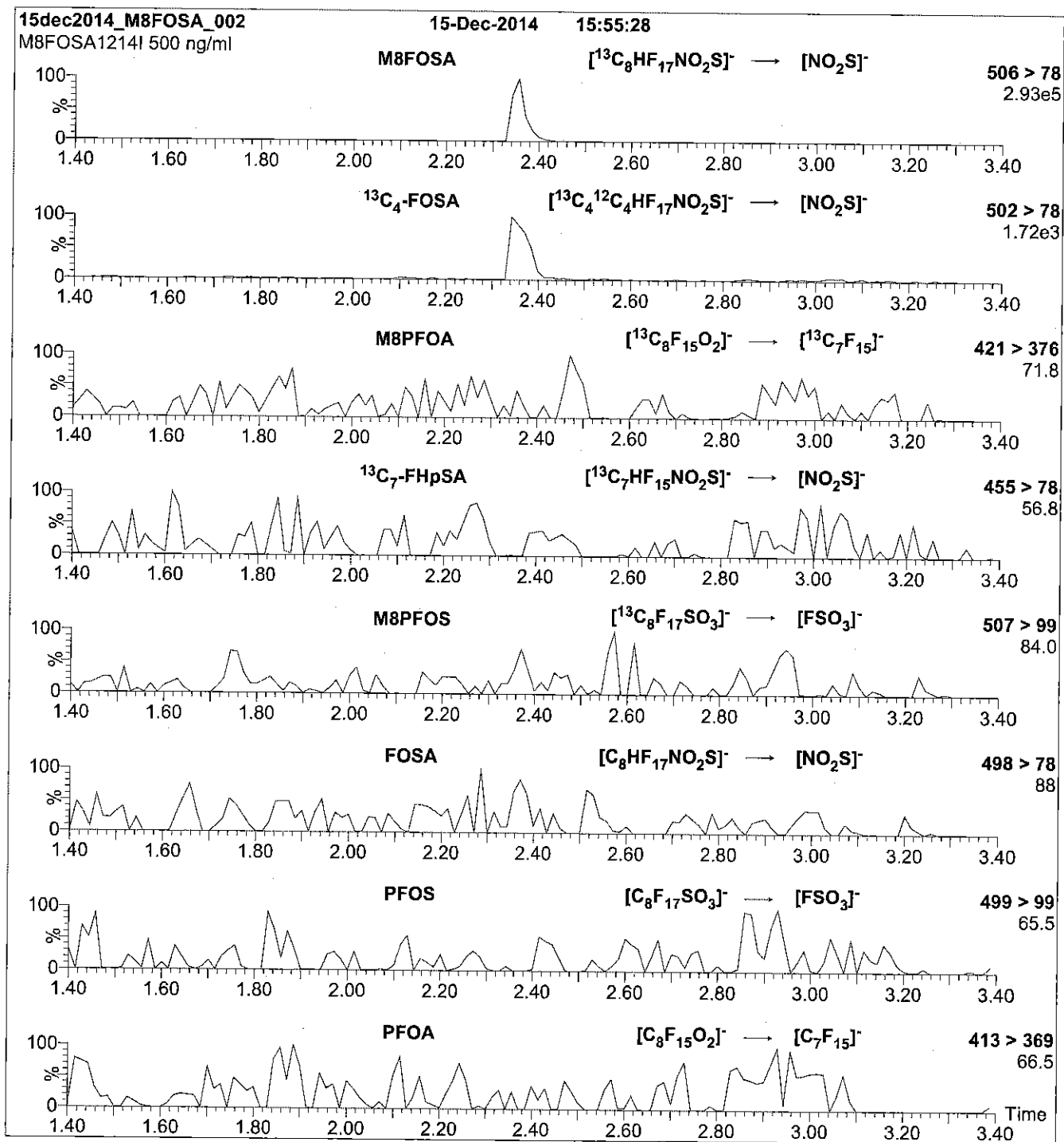
Flow: 300  $\mu$ l/min

**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<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300 µl/min

**MS Parameters**

Collision Gas (mbar) = 3.31e-3  
Collision Energy (eV) = 30

Reagent

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**LCMPFBA\_00004**



# WELLINGTON LABORATORIES

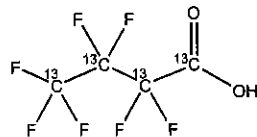
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFBA  
**COMPOUND:** Perfluoro-n-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]butanoic acid

**LOT NUMBER:** MPFBA1014

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>4</sub>HF<sub>7</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 218.01  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 10/31/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 10/31/2019  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**ISOTOPIC PURITY:** ≥99%<sup>13</sup>C  
(1,2,3,4-<sup>13</sup>C<sub>4</sub>)

**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:** 03/31/2015  
(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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### **EXPIRY DATE / PERIOD OF VALIDITY:**

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### **LIMITED WARRANTY:**

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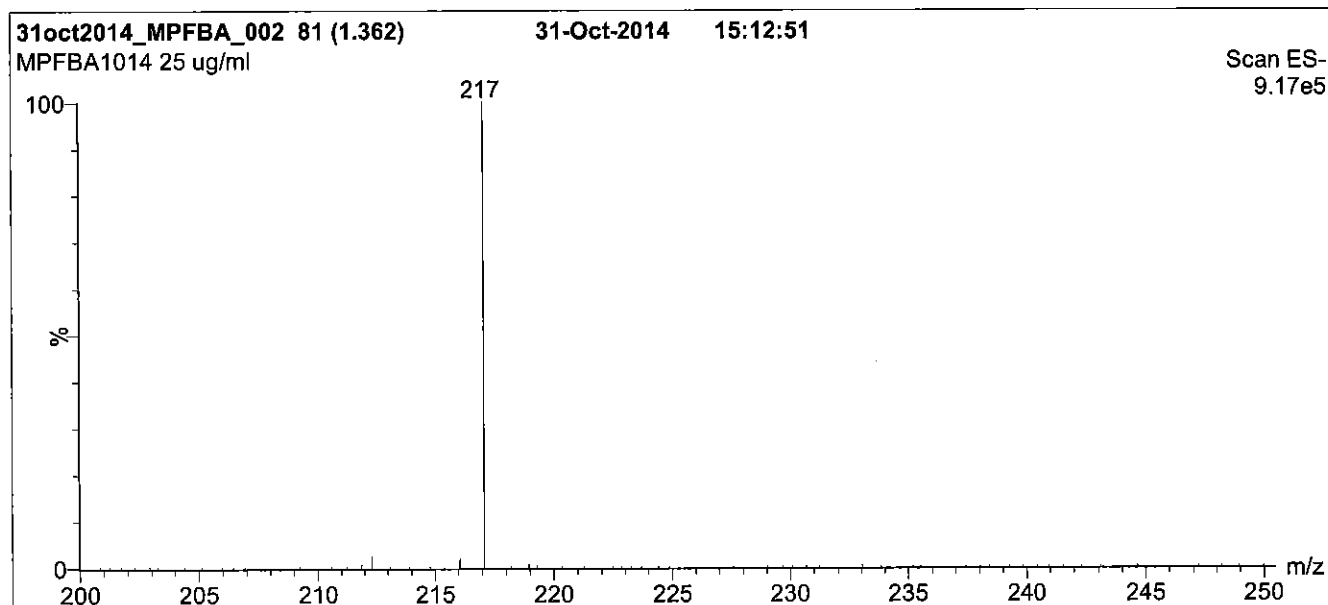
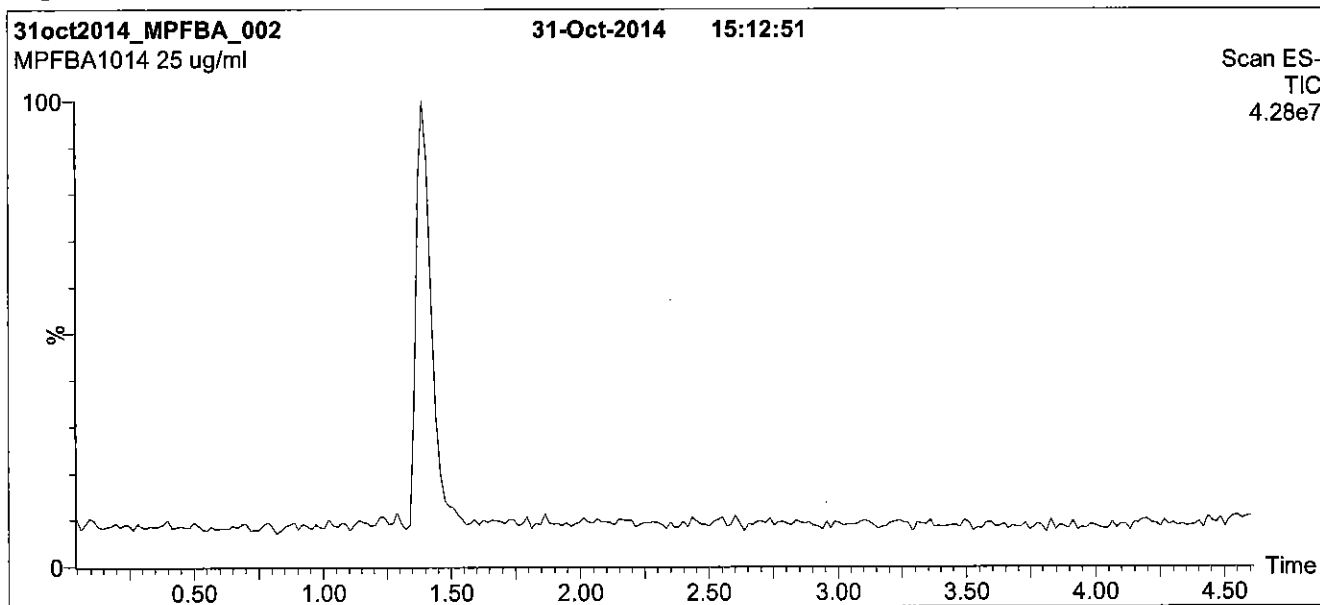
### **QUALITY MANAGEMENT:**

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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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 5 min and hold for 1.5 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min

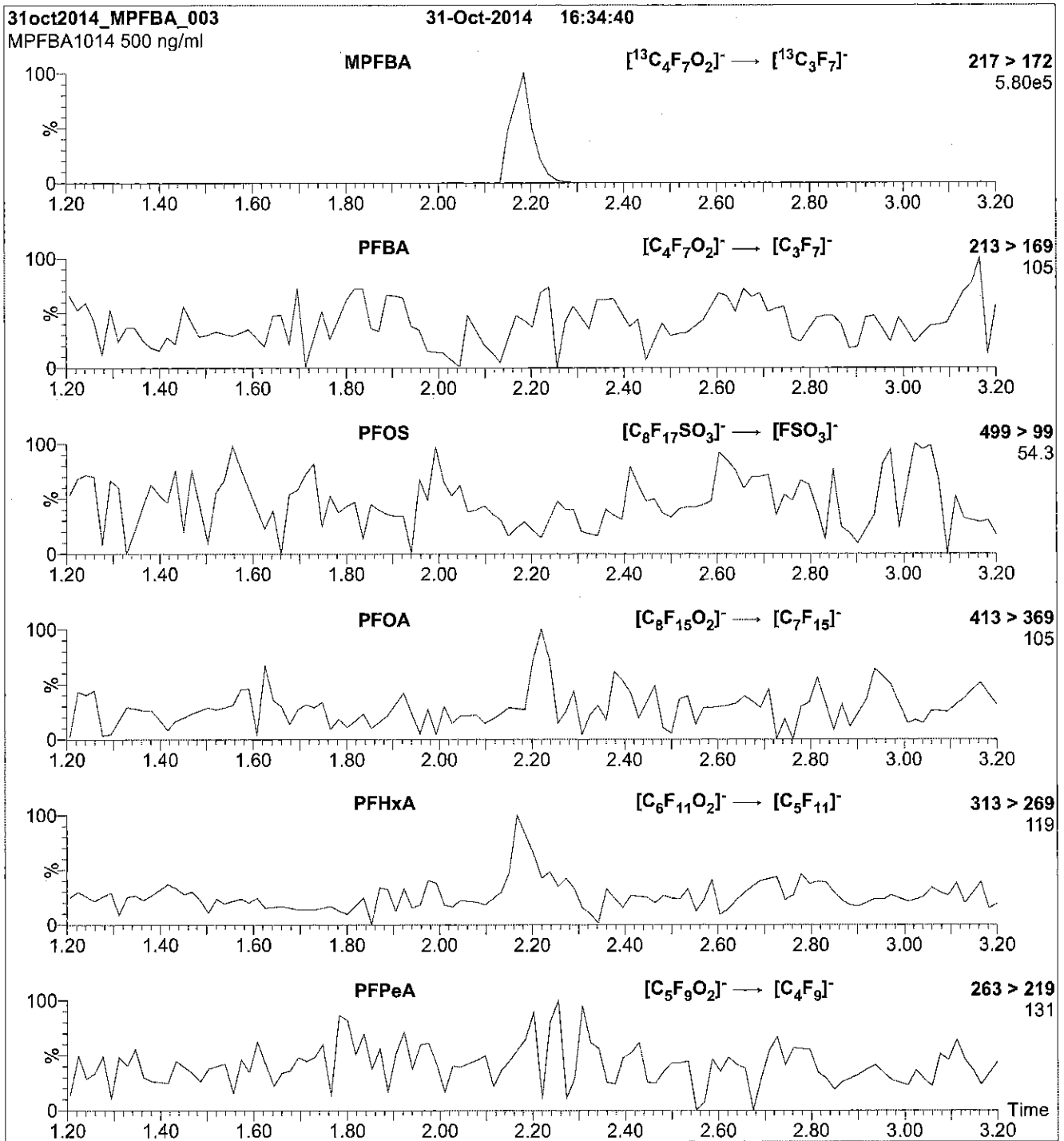
**Flow:** 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (200 - 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: MPFBA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml MPFBA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.28e-3  
Collision Energy (eV) = 10



Reagent

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**LCMPFDA\_00004**

R: 10-20-2011  
2011  
2011  
2011



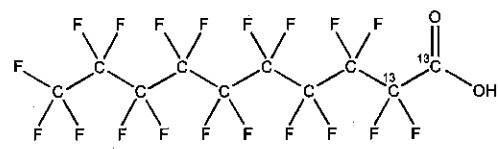
# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

12LCMS0262  
LCMPFDA-00001

**PRODUCT CODE:** MPFDA **LOT NUMBER:** MPFDA0411  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]decanoic acid

**STRUCTURE:** **CAS #** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>8</sub>HF<sub>19</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 516.07  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
(1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 04/07/2011  
**EXPIRY DATE:** (mm/dd/yyyy) 04/07/2014  
**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 <sup>13</sup>C<sub>1</sub>-PFNA.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:** **Date:** 04/19/2011  
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

Form#: 27, Issued 2004-11-10  
Revision#: 1, Revised 2010-07-26

MPFDA0411 (1 of 4)  
rev0

**INTENDED USE:**

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**HAZARDS:**

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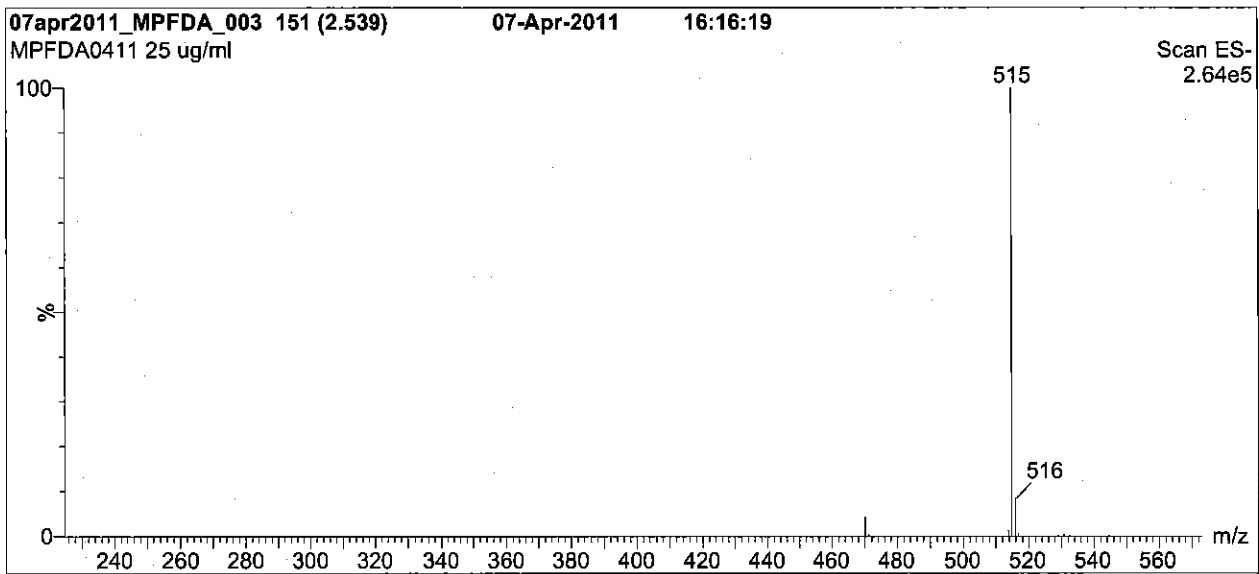
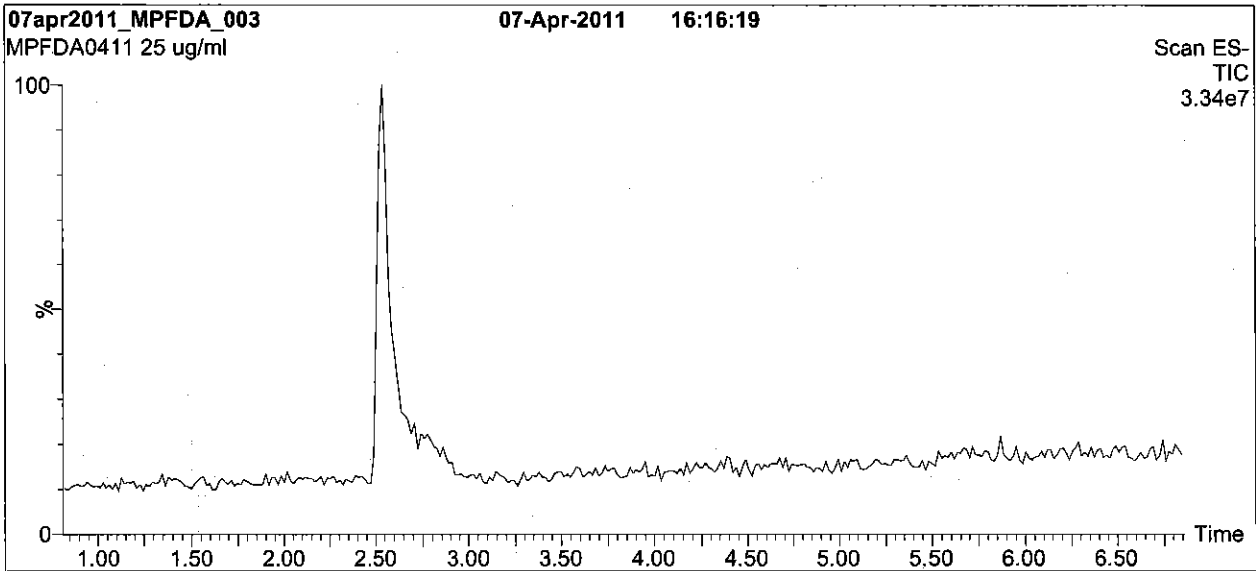
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**LIMITED WARRANTY:**

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.

\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

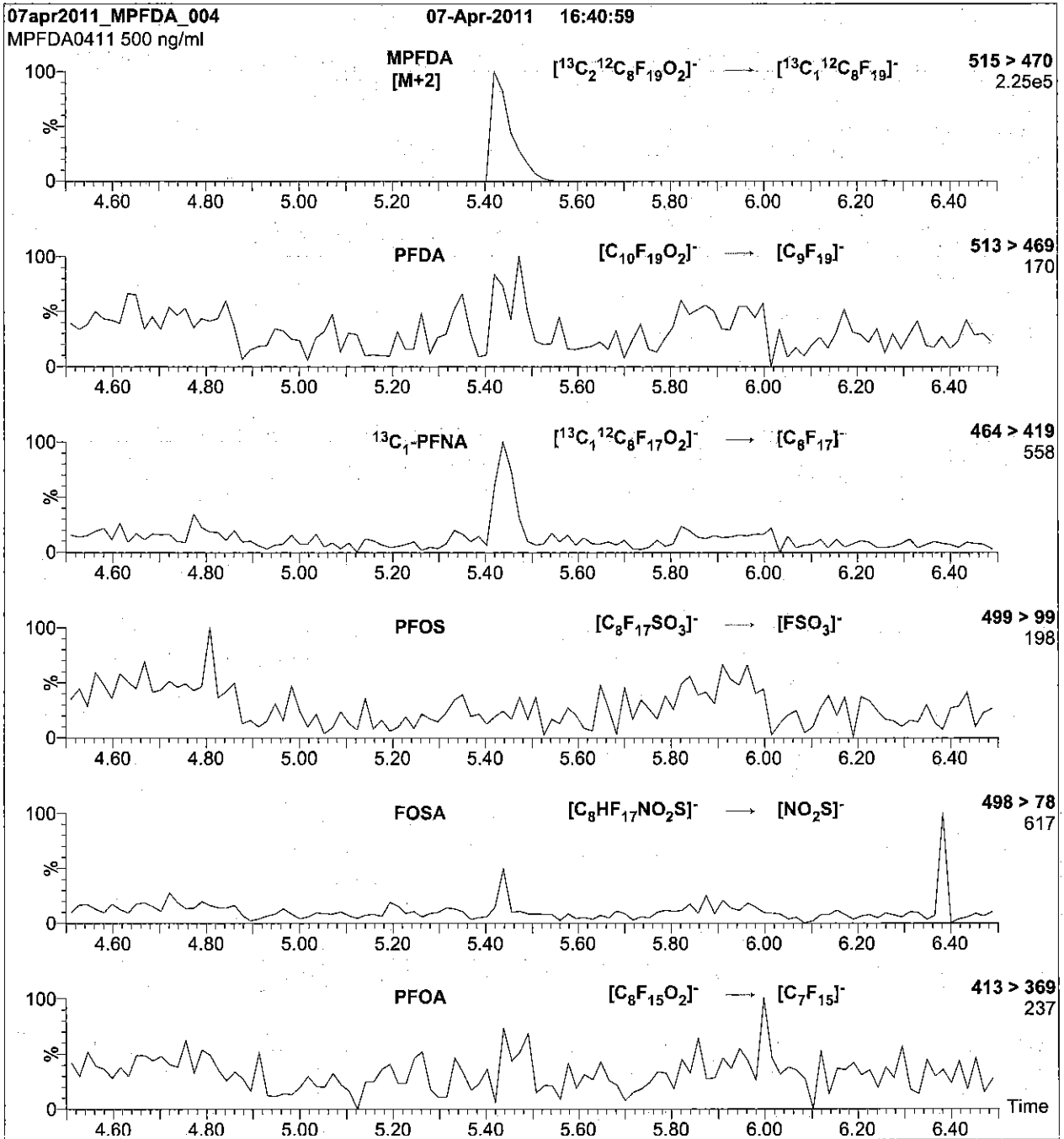
Mobile phase: Gradient  
Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min and hold for 1.5 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min

Flow: 300  $\mu$ 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: MPFDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml MPFDA)

**Mobile phase:** Isocratic 70% (80:20 MeOH:ACN) / 30%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

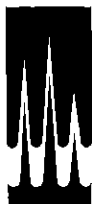
**MS Parameters**

Collision Gas (mbar) = 3.39e-3  
Collision Energy (eV) = 13

Reagent

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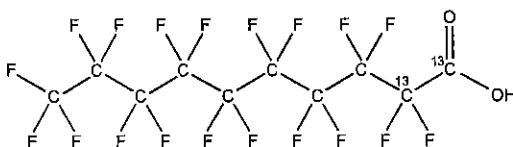
**LCMPFDA\_00005**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFDA **LOT NUMBER:** MPFDA0414  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]decanoic acid  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>8</sub>HF<sub>19</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 516.07  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
 (1,2-<sup>13</sup>C<sub>2</sub>)  
**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.
- Contains ~ 0.1% of <sup>13</sup>C<sub>1</sub>-PFNA.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

Certified By:

B.G. Chittim

Date: 04/15/2014

(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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:**

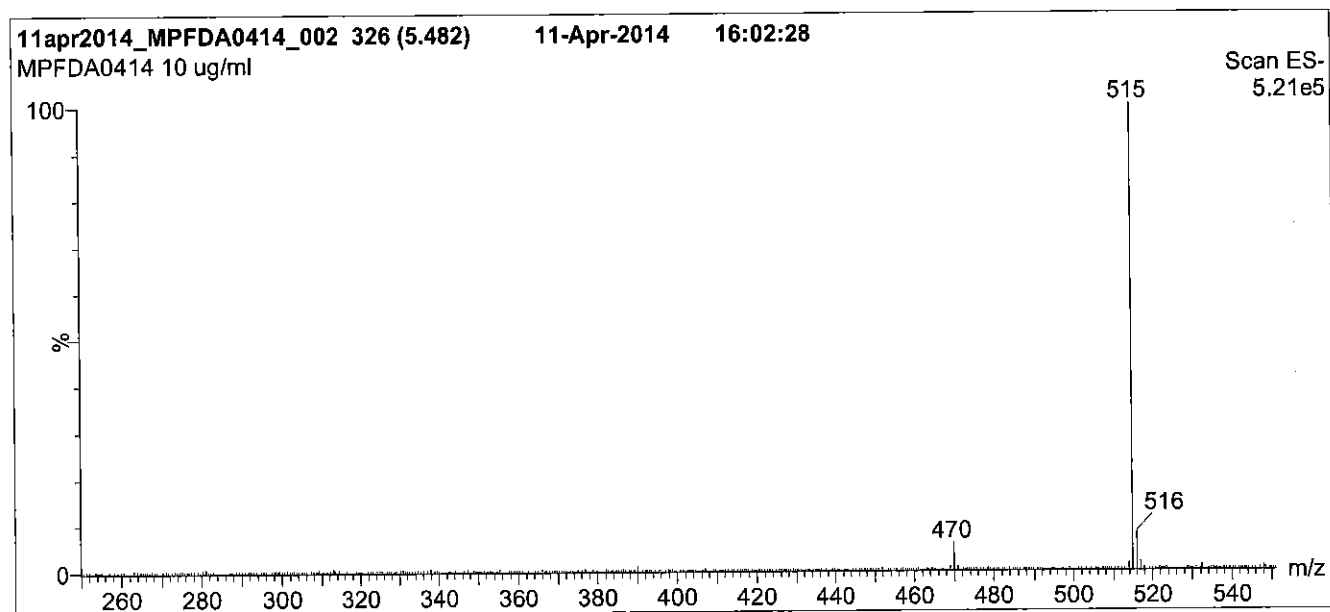
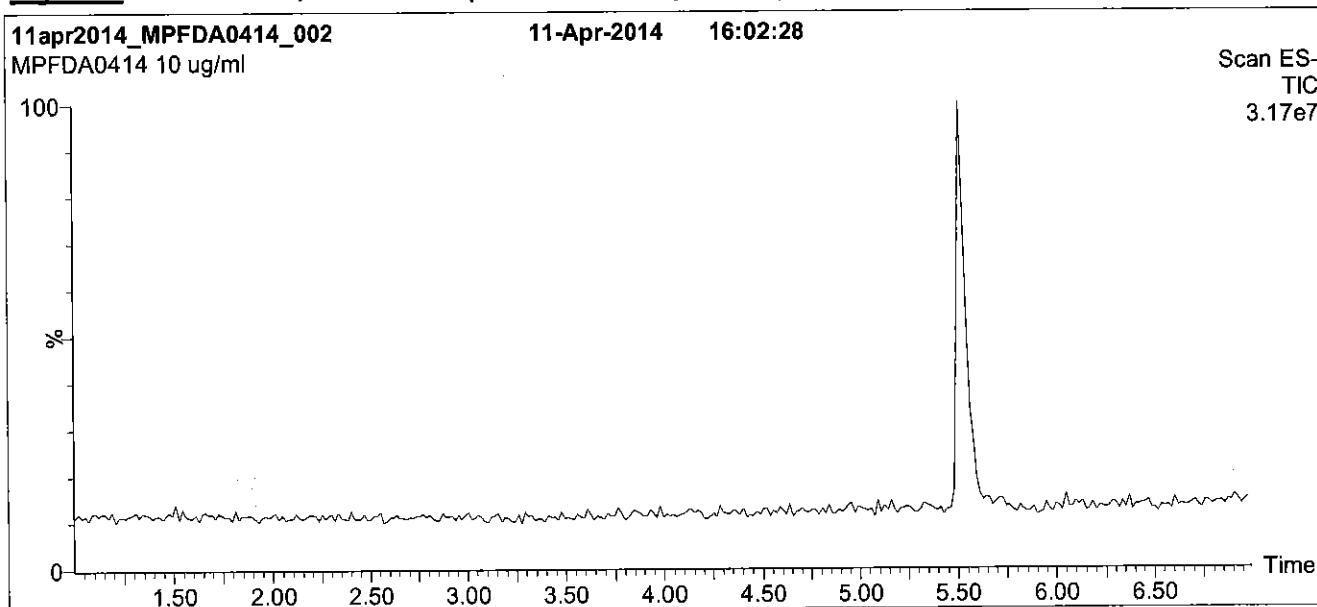
This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number AR-1523).



\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*



**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

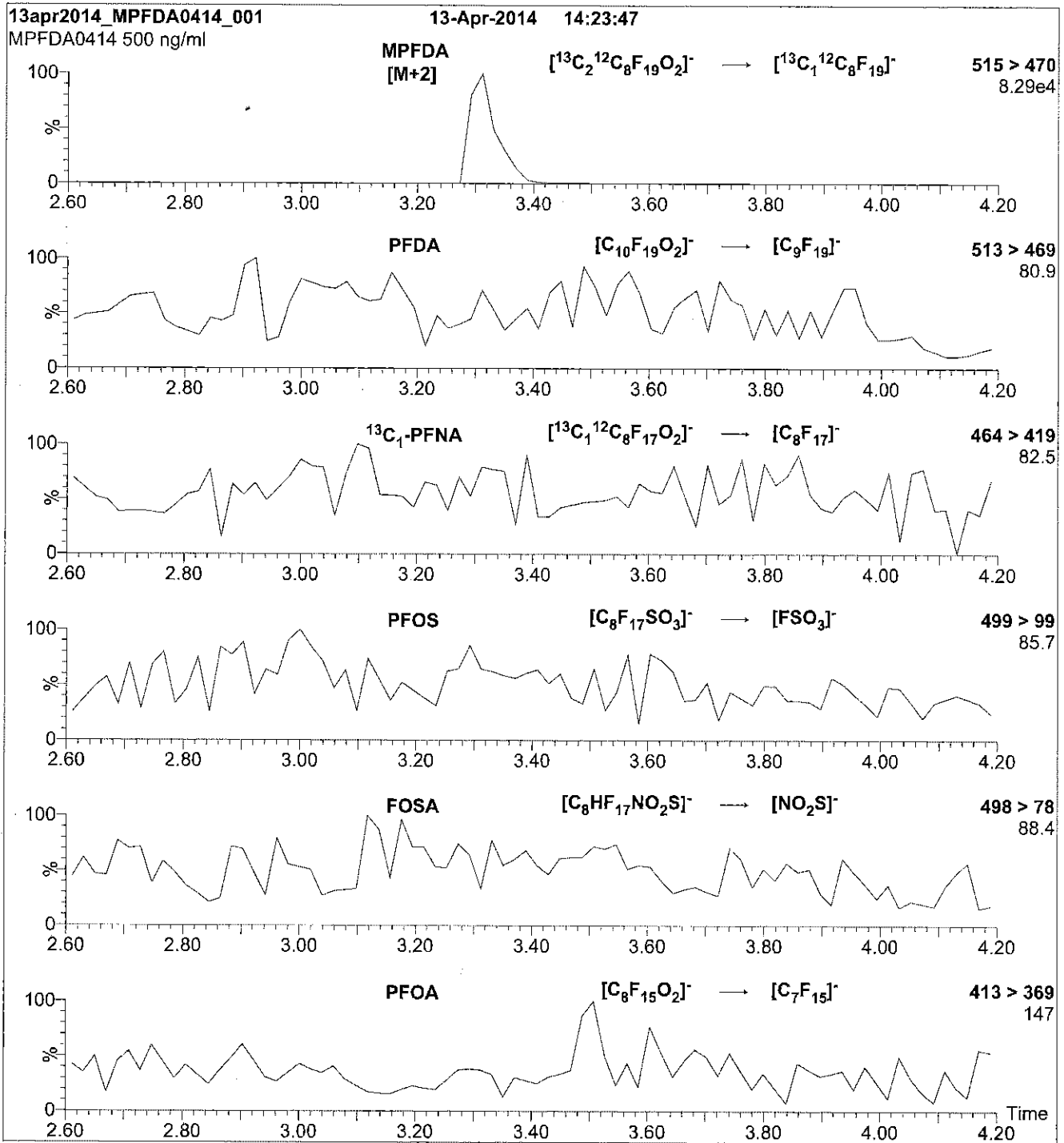
**Flow:** 300  $\mu$ l/min

**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: MPFDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml MPFDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.31e-3  
Collision Energy (eV) = 13

Reagent

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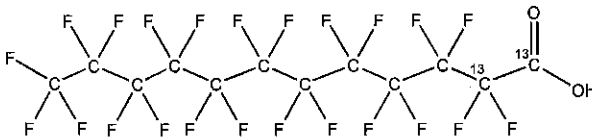
**LCMPFD<sub>o</sub>A\_00003**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFDoA **LOT NUMBER:** MPFDoA0714  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]dodecanoic acid  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub> <sup>12</sup>C<sub>10</sub> HF<sub>23</sub> O<sub>2</sub> **MOLECULAR WEIGHT:** 616.08  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
 (1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 07/17/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 07/17/2019  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

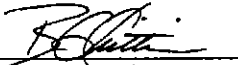
### DOCUMENTATION/ DATA ATTACHED:

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:** 07/21/2014  
 (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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

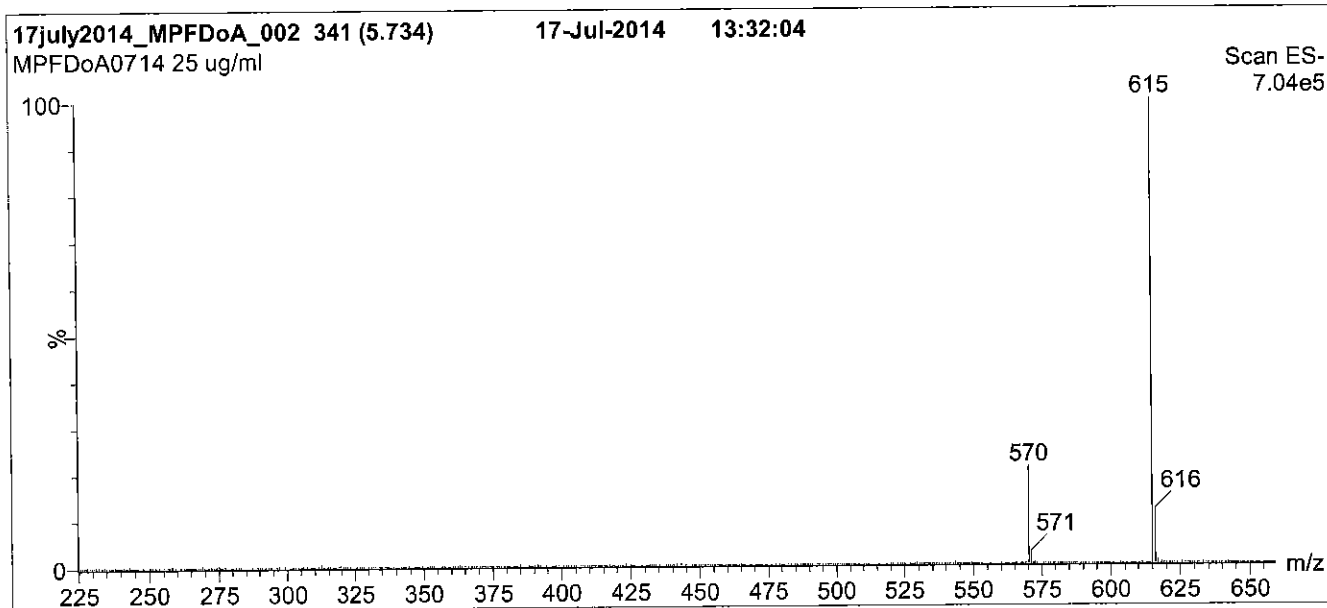
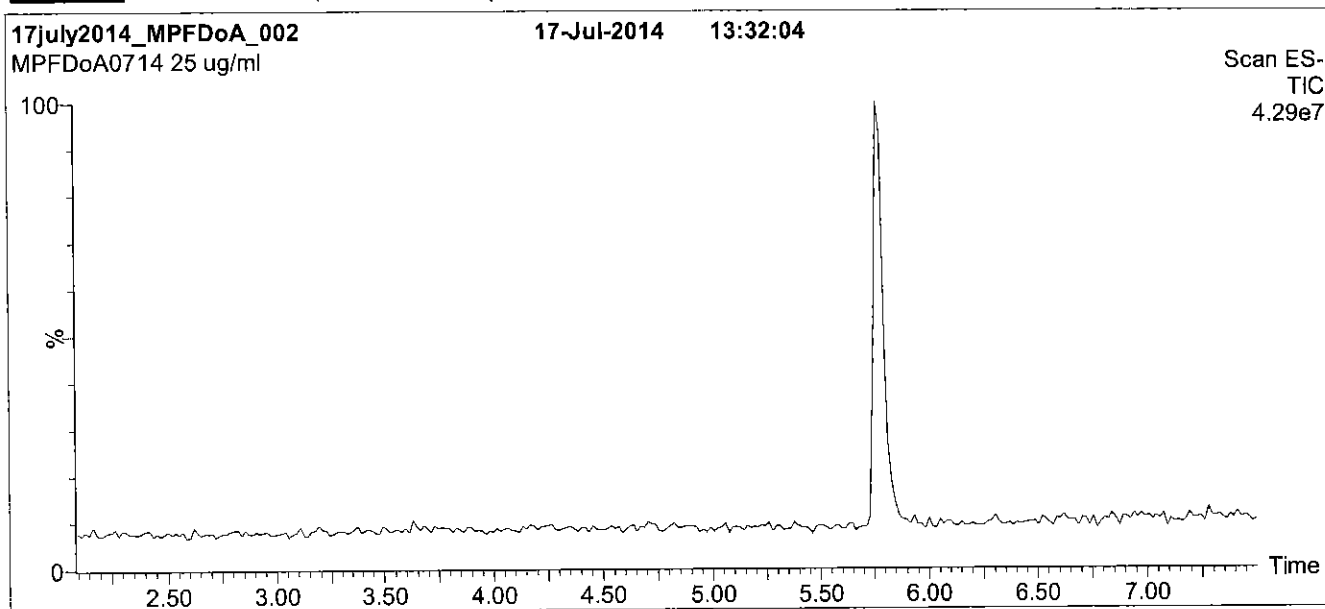
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number AR-1523).



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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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

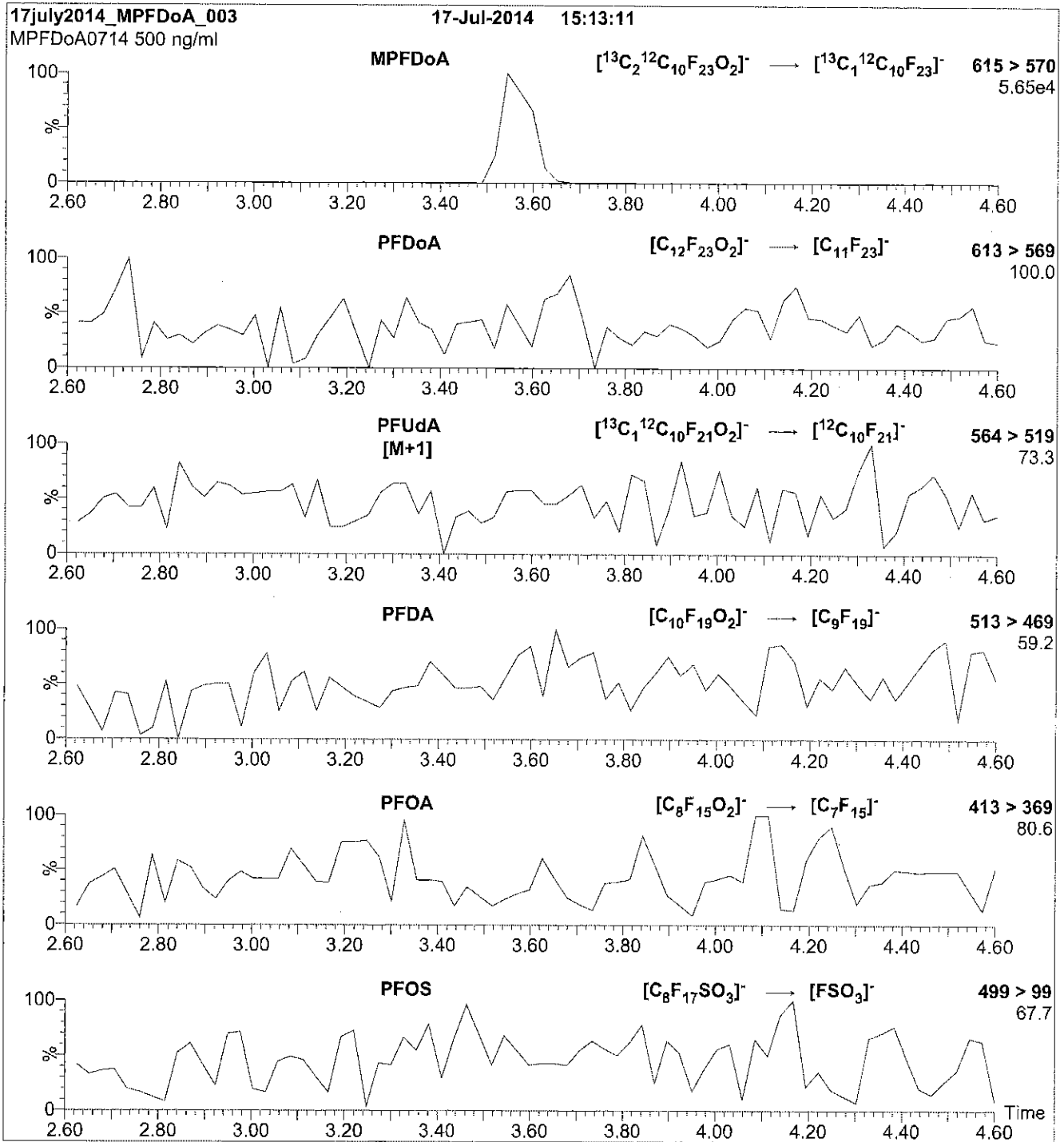
**Mobile phase:** Gradient  
 Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)  
 Ramp to 90% organic over 7 min and hold for 2 min  
 before returning to initial conditions in 0.5 min.  
 Time: 10 min

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (225 - 950 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  $\mu\text{l}$  (500 ng/ml MPFDoA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.43e-3  
Collision Energy (eV) = 13

Reagent

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**LCMPFD<sub>o</sub>A\_00004**





### **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.

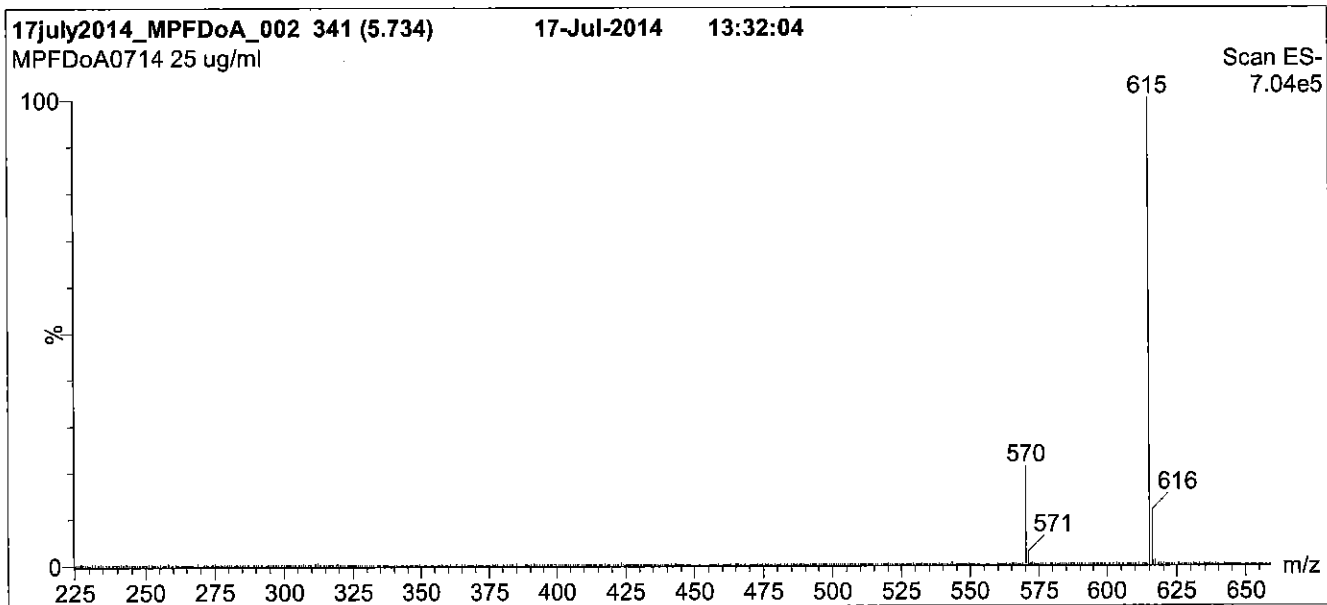
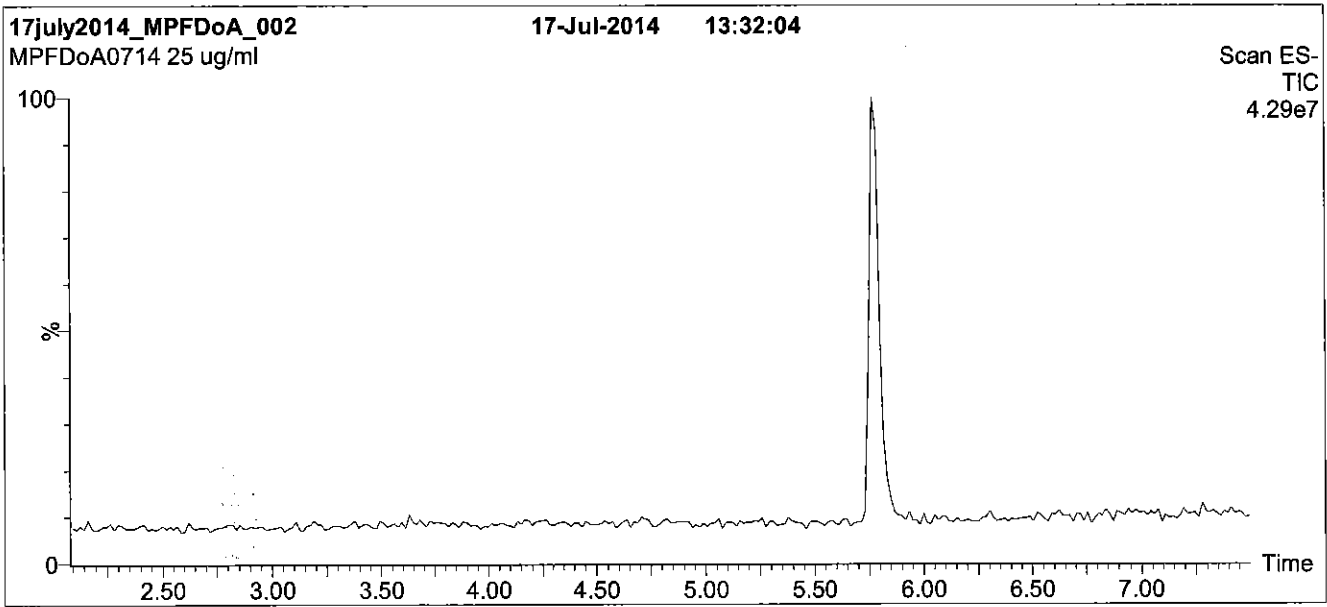
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>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

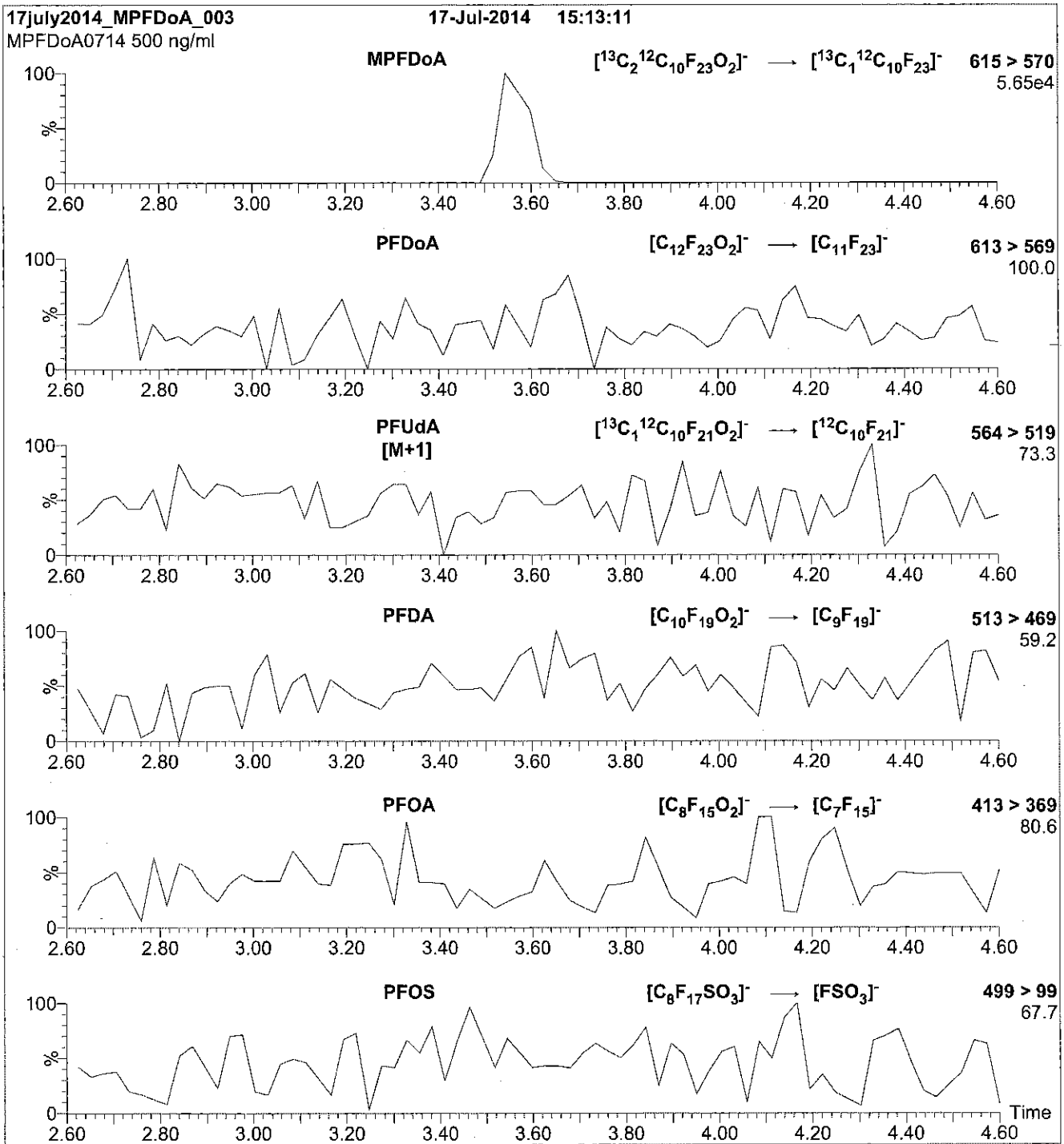
**Flow:** 300  $\mu$ l/min

**MS Parameters**

**Experiment:** Full Scan (225 - 950 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  $\mu\text{l}$  (500 ng/ml MPFDoA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.43e-3  
Collision Energy (eV) = 13

Reagent

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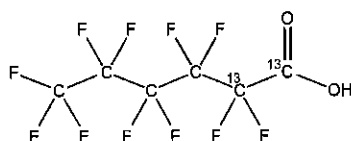
**LCMPFHxA\_00006**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFHxA  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]hexanoic acid  
**LOT NUMBER:** MPFHxA0414  
**STRUCTURE:**  
**CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>4</sub>HF<sub>11</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml  
**MOLECULAR WEIGHT:** 316.04  
**SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98%  
**ISOTOPIC PURITY:** ≥99%<sup>13</sup>C  
 (1,2-<sup>13</sup>C<sub>2</sub>)  
**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.
- 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: 04/15/2014

(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

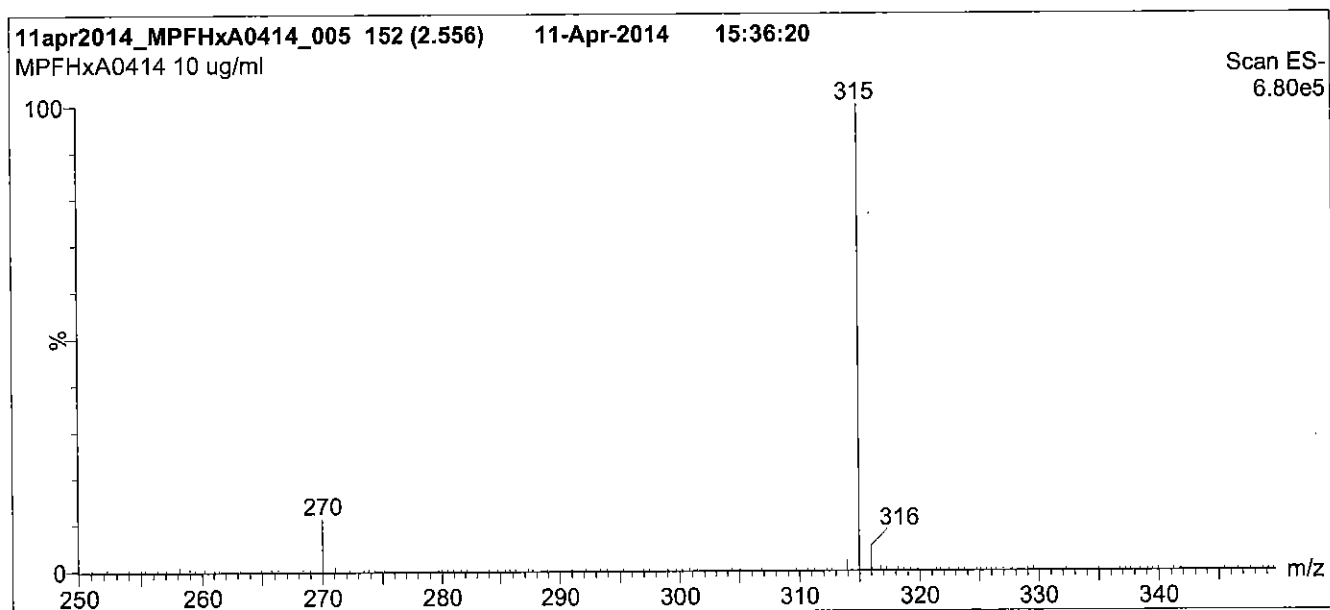
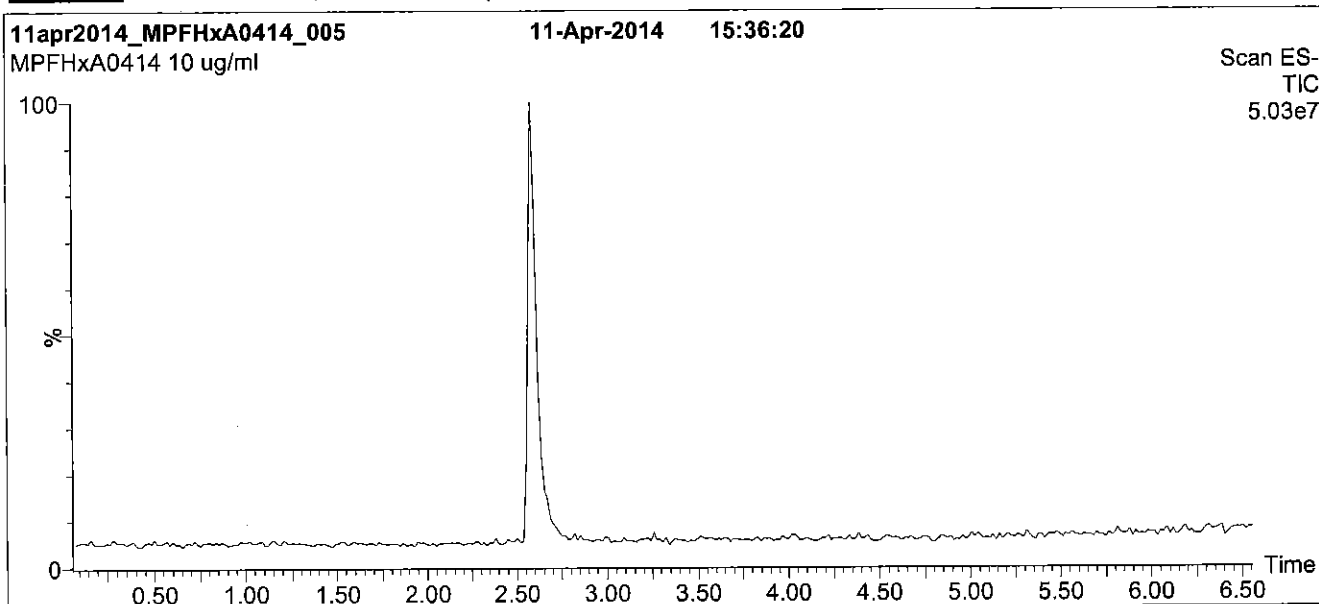
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**Figure 1: MPFHxA; 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<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min and hold for 2 min  
before returning to initial conditions over 0.5 min.  
Time: 10 min

**Flow:** 300  $\mu$ l/min

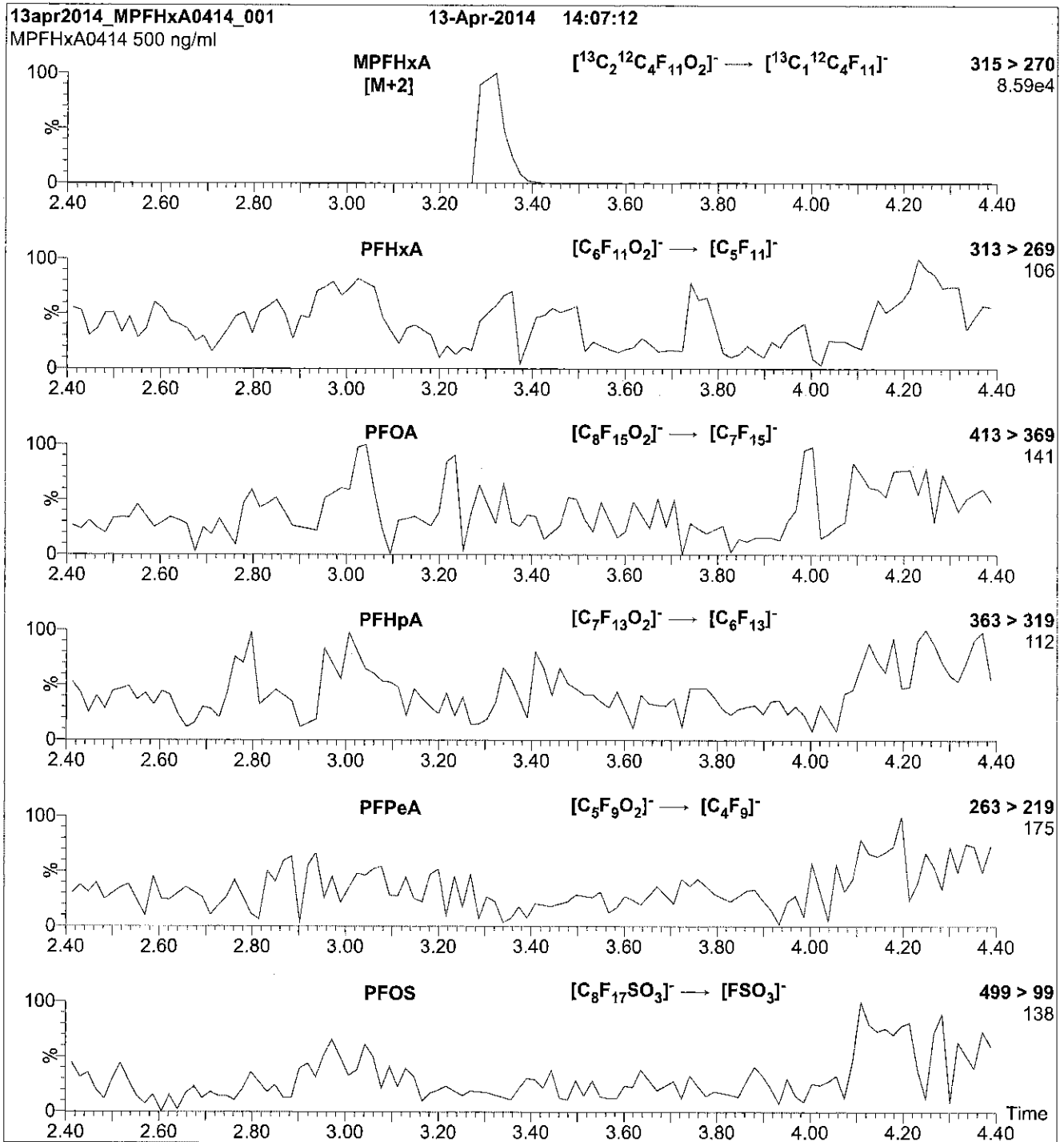
**MS Parameters**

**Experiment:** Full Scan (250 - 850 amu)

**Source:** Electrospray (negative)  
**Capillary Voltage (kV)** = 2.00  
**Cone Voltage (V)** = 15.00  
**Cone Gas Flow (l/hr)** = 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  $\mu$ l (500 ng/ml MPFHxA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.43e-3  
 Collision Energy (eV) = 10

Reagent

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**LCMPFHXS\_00003**

v. 21/15/12

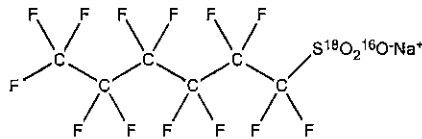


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFHxS **LOT NUMBER:** MPFHxS0713  
**COMPOUND:** Sodium perfluoro-1-hexane[<sup>18</sup>O<sub>2</sub>]sulfonate

**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>6</sub>F<sub>13</sub>S<sup>18</sup>O<sub>2</sub><sup>16</sup>O<sup>-</sup>Na **MOLECULAR WEIGHT:** 426.10  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol  
47.3 ± 2.4 µg/ml (MPFHxS anion)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** >94% (<sup>18</sup>O<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 07/25/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 07/25/2018  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

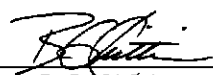
### DOCUMENTATION/ DATA ATTACHED:

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.
- The response factor for MPFHxS (C<sub>6</sub>F<sub>13</sub>S<sup>18</sup>O<sub>2</sub><sup>16</sup>O<sup>-</sup>) has been observed to be up to 10% lower than for PFHxS (C<sub>6</sub>F<sub>13</sub>S<sup>18</sup>O<sub>3</sub><sup>-</sup>) when both compounds are injected together. This difference may vary between instruments.
- Due to the isotopic purity of the starting material (<sup>18</sup>O<sub>2</sub> >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:** 04/15/2014  
(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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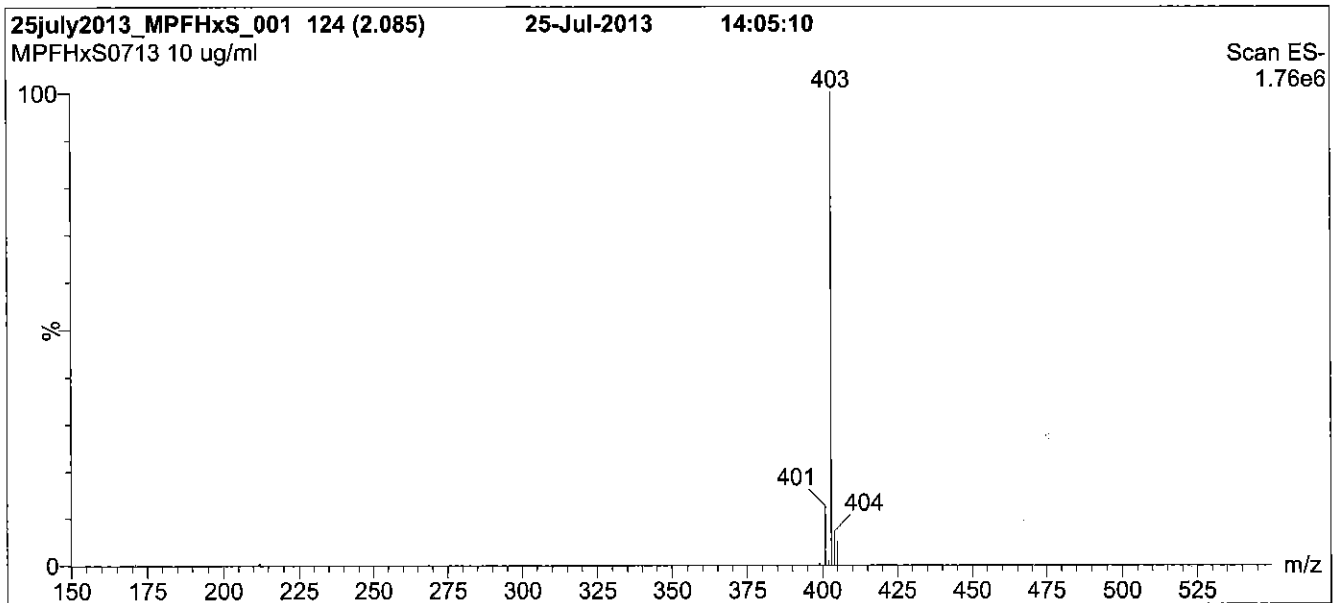
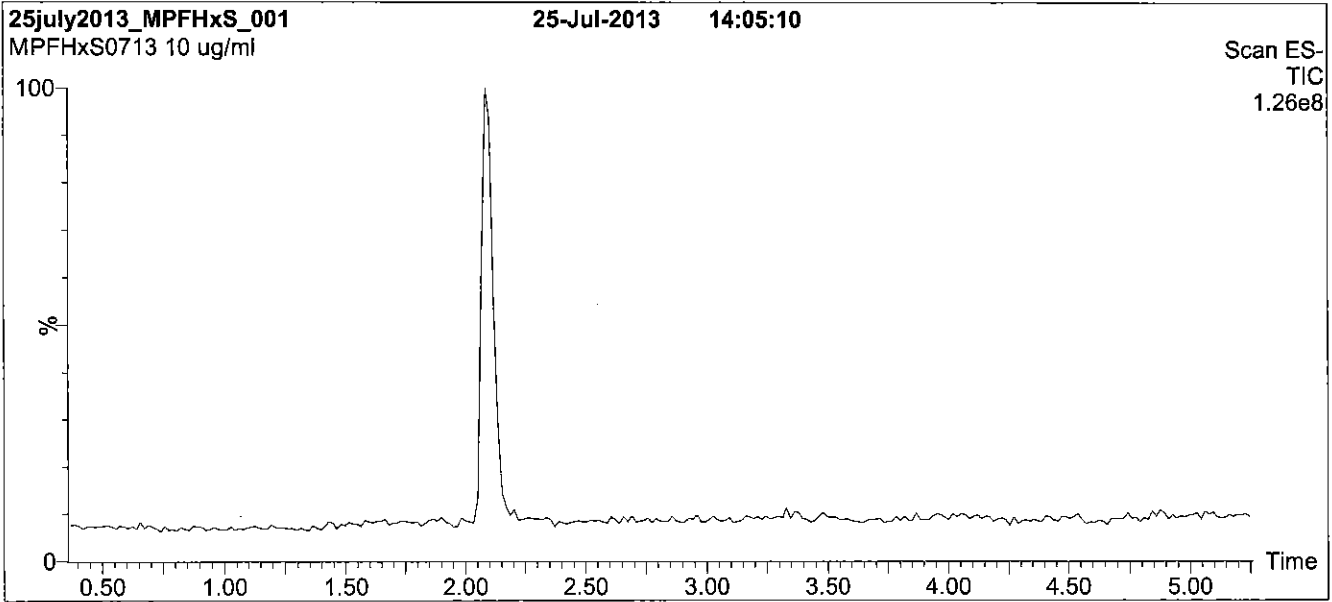
**QUALITY MANAGEMENT:**

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**Figure 1: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

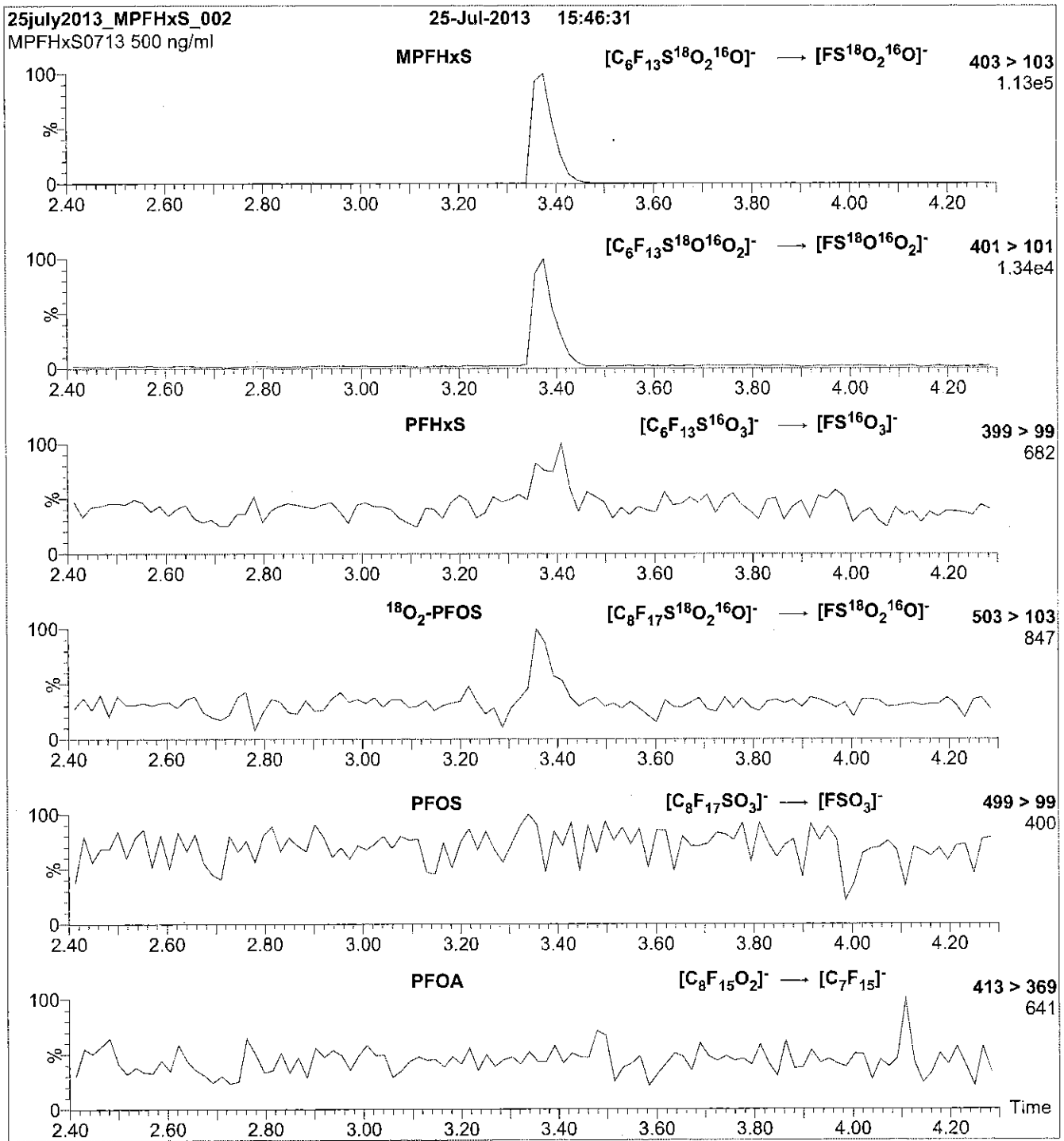
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
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  $\mu$ l (500 ng/ml MPFHxS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.58e-3  
 Collision Energy (eV) = 30

Reagent

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**LCMPFHXS\_00004**

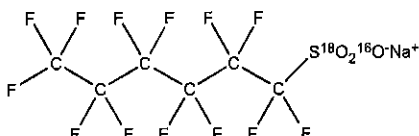


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFHxS **LOT NUMBER:** MPFHxS0713  
**COMPOUND:** Sodium perfluoro-1-hexane<sup>[18O<sub>2</sub>]</sup>sulfonate

**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>6</sub>F<sub>13</sub>S<sup>18</sup>O<sub>2</sub><sup>16</sup>O<sup>-</sup>Na<sup>+</sup> **MOLECULAR WEIGHT:** 426.10  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol  
 47.3 ± 2.4 µg/ml (MPFHxS anion)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** >94% (<sup>18</sup>O<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 07/25/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 07/25/2018  
**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.
- The response factor for MPFHxS (C<sub>6</sub>F<sub>13</sub>S<sup>18</sup>O<sub>2</sub><sup>16</sup>O<sup>-</sup>) has been observed to be up to 10% lower than for PFHxS (C<sub>6</sub>F<sub>13</sub>S<sup>16</sup>O<sub>3</sub><sup>-</sup>) when both compounds are injected together. This difference may vary between instruments.
- Due to the isotopic purity of the starting material (<sup>18</sup>O<sub>2</sub> >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: 03/30/2015  
(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.

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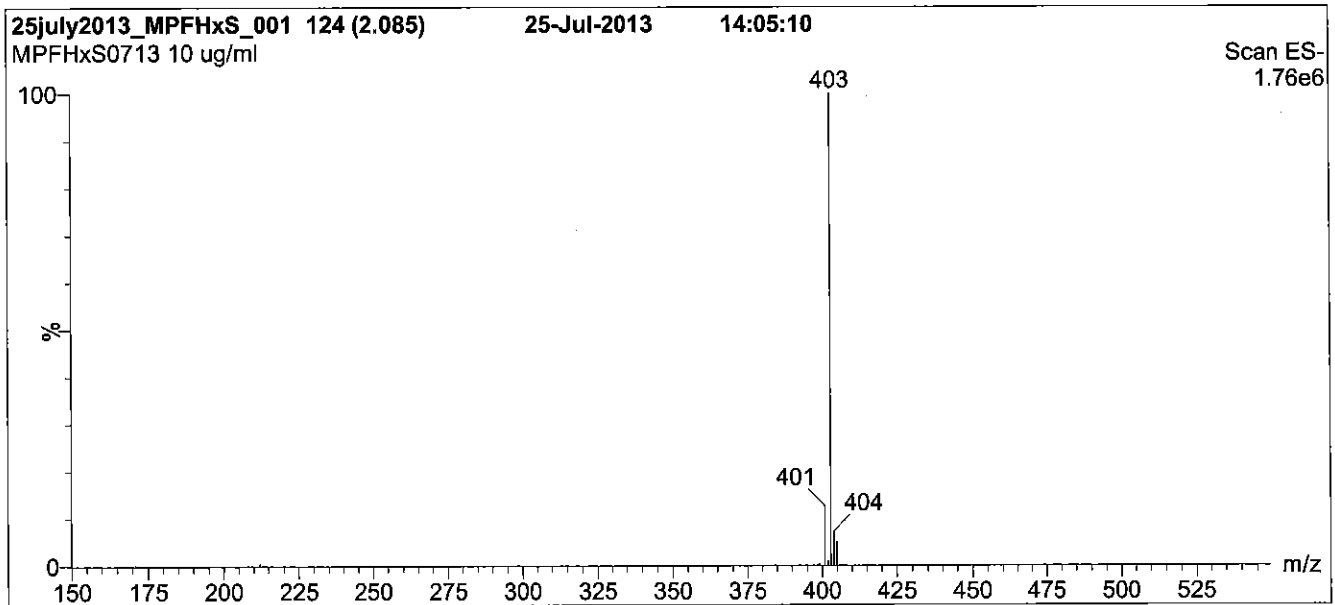
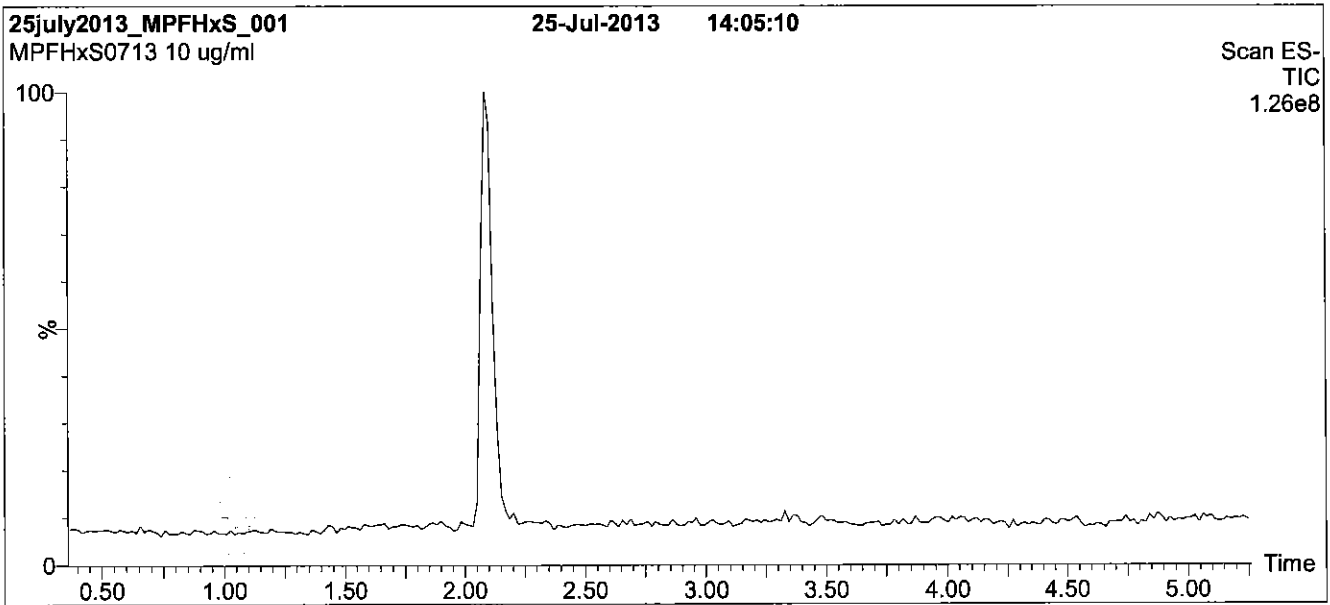
### **QUALITY MANAGEMENT:**

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**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

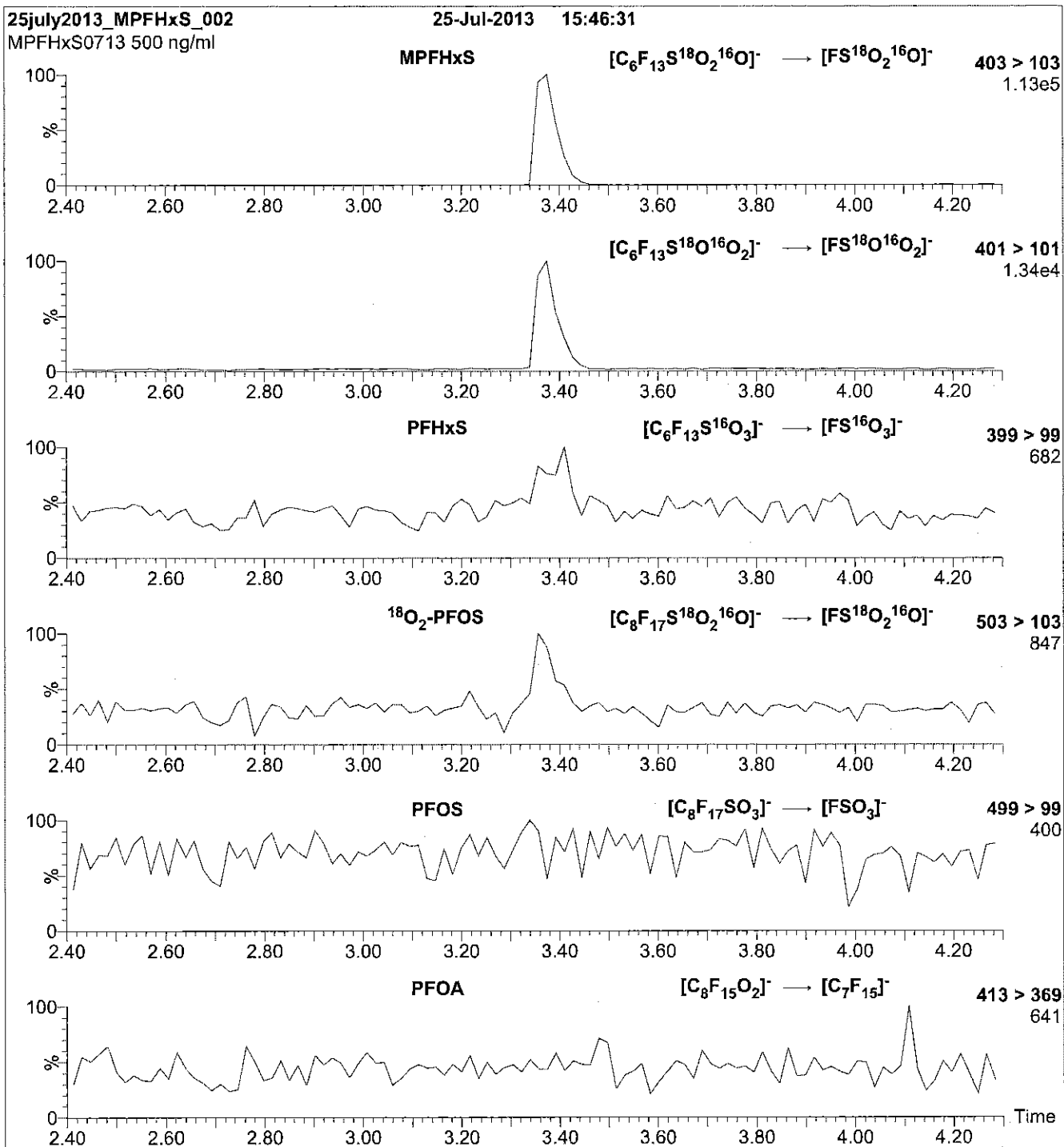
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
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  $\mu$ l (500 ng/ml MPFHxS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.58e-3  
Collision Energy (eV) = 30

Reagent

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**LCMPFNA\_00003**

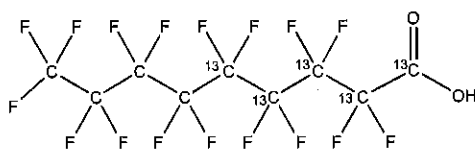


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFNA **LOT NUMBER:** MPFNA0414  
**COMPOUND:** Perfluoro-n-[1,2,3,4,5-<sup>13</sup>C<sub>5</sub>]nonanoic acid

**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:**  $^{13}\text{C}_5^{12}\text{C}_4\text{HF}_{17}\text{O}_2$  **MOLECULAR WEIGHT:** 469.04  
**CONCENTRATION:**  $50 \pm 2.5 \mu\text{g/ml}$  **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:**  $\geq 99\%^{13}\text{C}$   
**LAST TESTED:** (mm/dd/yyyy) 04/13/2014 (1,2,3,4,5-<sup>13</sup>C<sub>5</sub>)  
**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: 04/13/2014  
(mm/dd/yyyy)

Wellington Laboratories Inc., 345 Southgate Dr. Guelph ON N1G 3M5 CANADA  
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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 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 17025:2005 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 for the period of time specified by the 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.

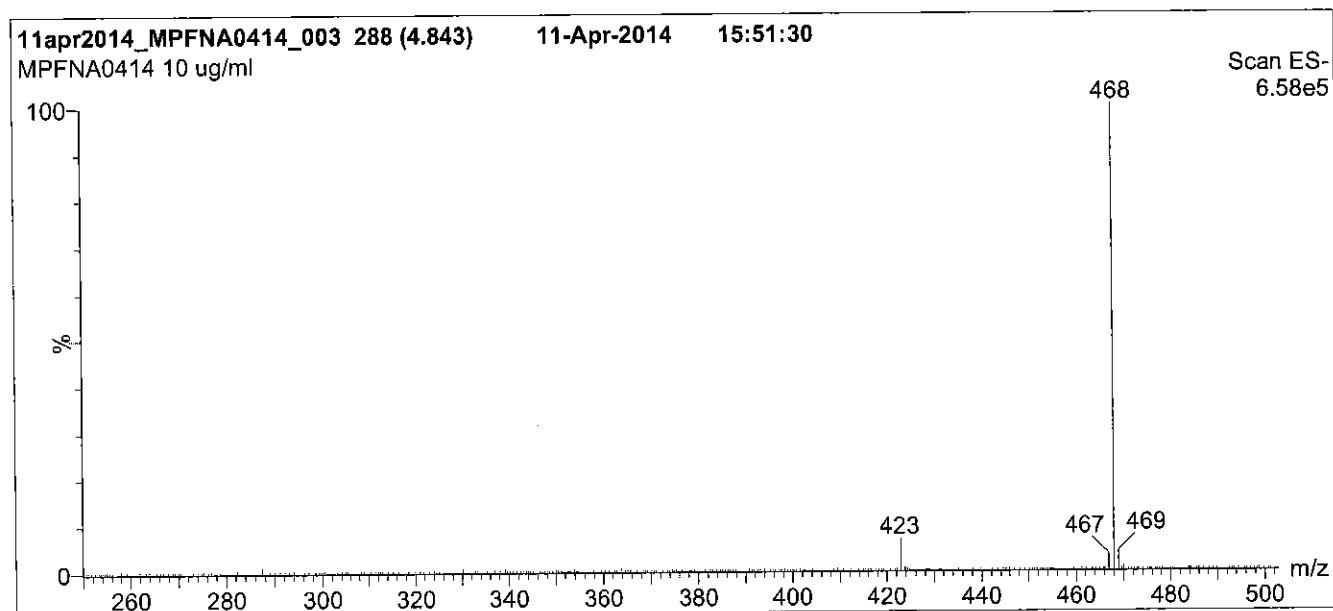
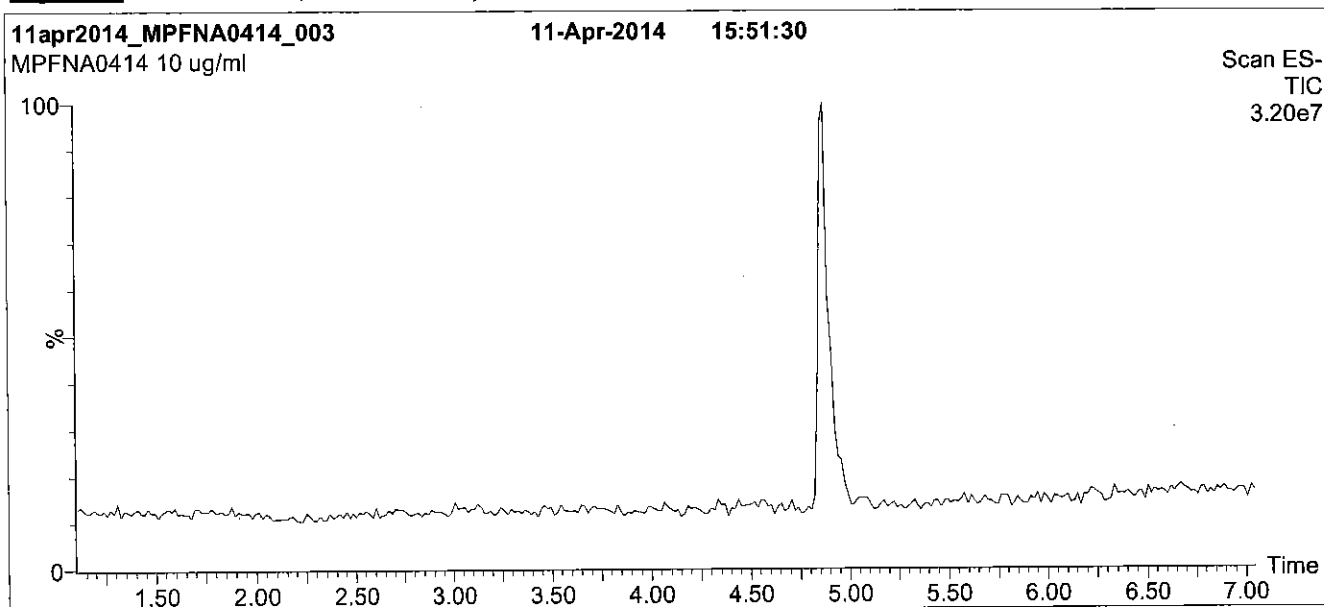
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto: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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

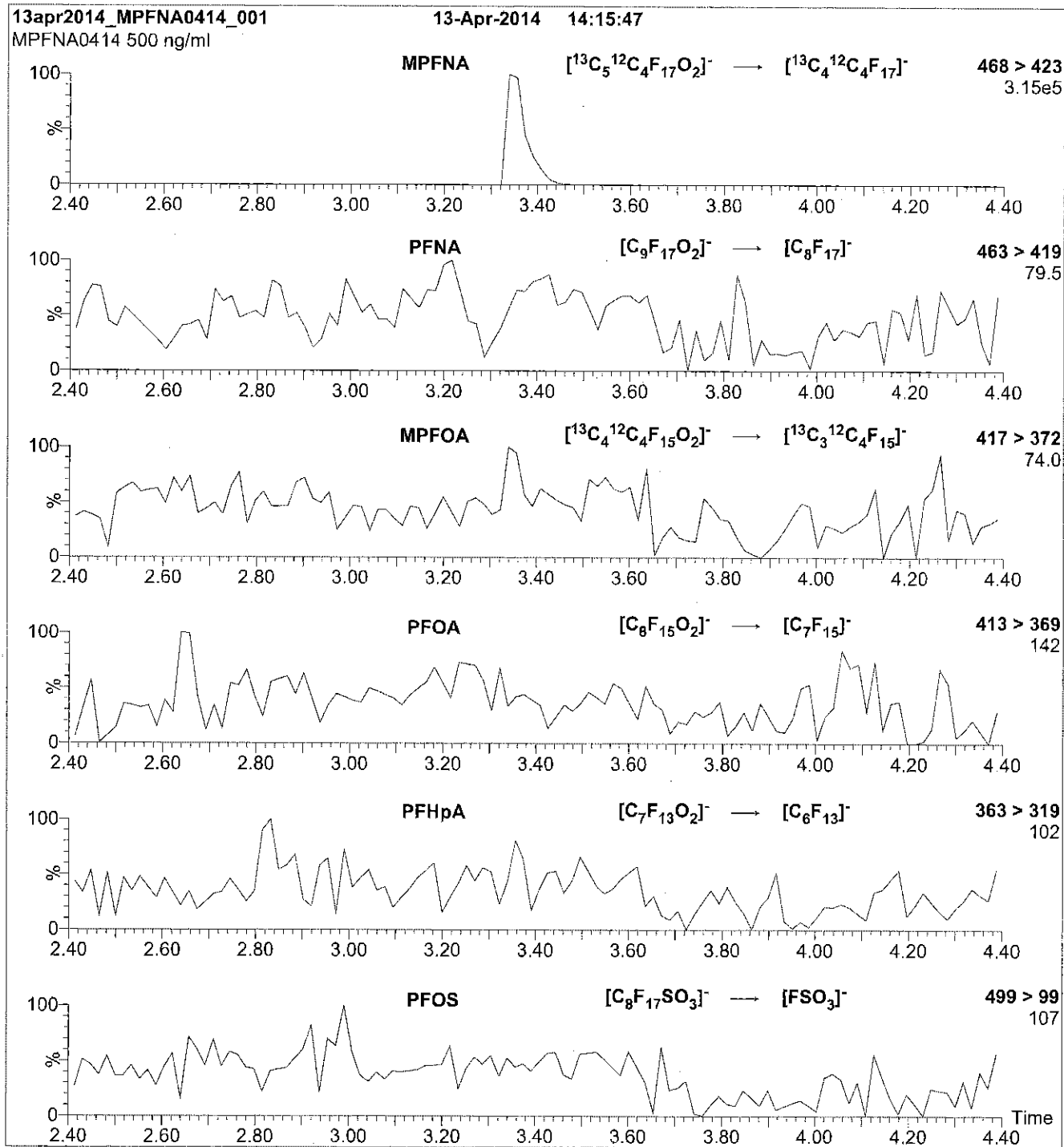
**Flow:** 300  $\mu$ l/min

**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  $\mu\text{l}$  (500 ng/ml MPFNA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.28e-3  
Collision Energy (eV) = 11



Reagent

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**LCMPFOA\_00007**

r: 9/5/15 sv



# WELLINGTON LABORATORIES

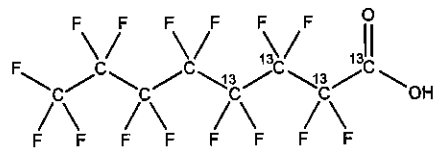
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFOA  
**COMPOUND:** Perfluoro-n-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]octanoic acid

**LOT NUMBER:** MPFOA0415

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>4</sub><sup>12</sup>C<sub>4</sub>HF<sub>16</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 418.04  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 04/10/2015  
**EXPIRY DATE:** (mm/dd/yyyy) 04/10/2020  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
(1,2,3,4-<sup>13</sup>C<sub>4</sub>)

**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:** 04/10/2015  
(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:**

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.

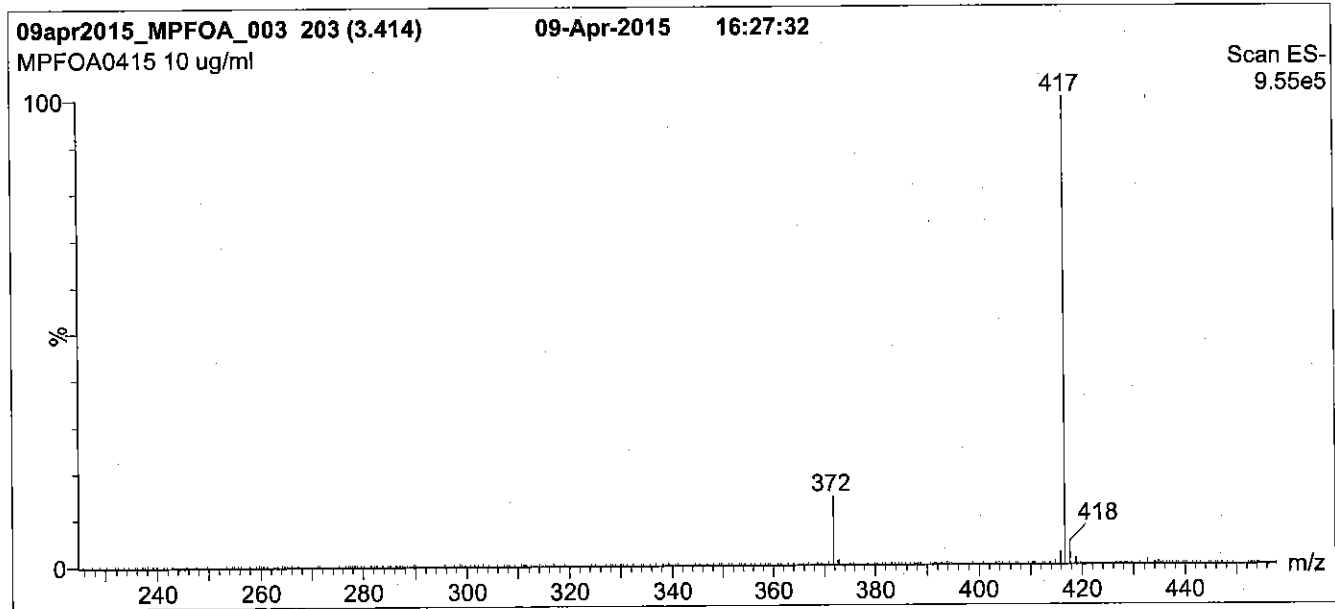
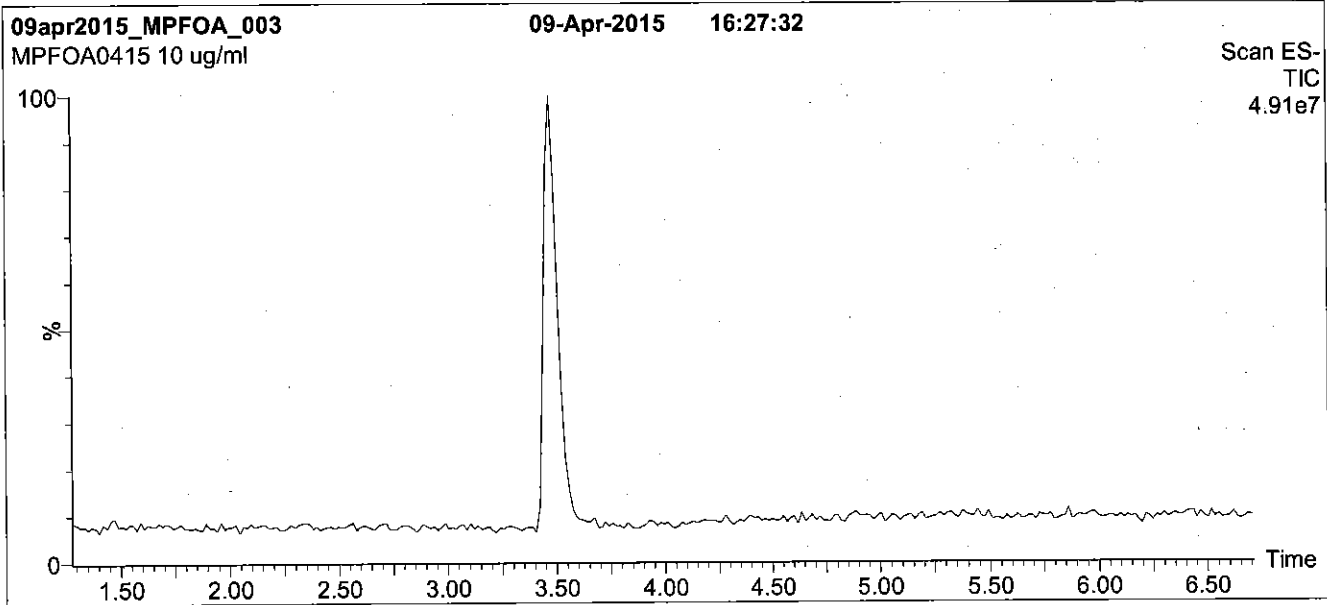
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**Figure 1: MPFOA; 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<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

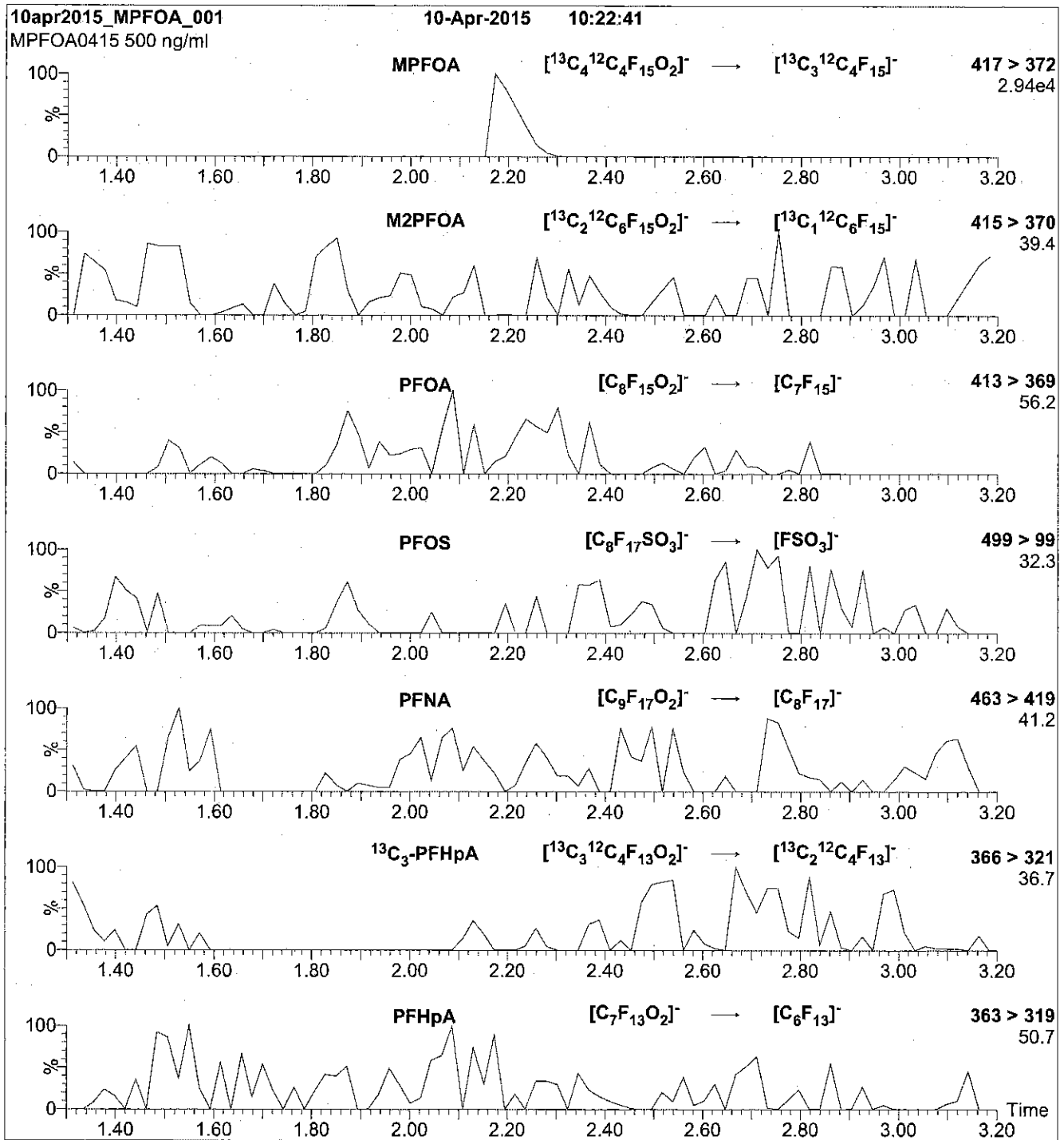
**Flow:** 300  $\mu$ l/min

**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) = 100  
Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

Injection: Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml MPFOA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.24e-3  
Collision Energy (eV) = 11

Reagent

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**LCMPFOS\_00009**

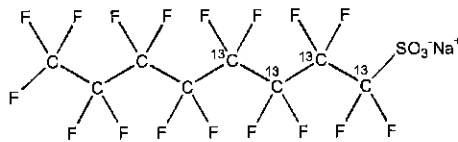
V: 91515 61



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFOS **LOT NUMBER:** MPFOS0515  
**COMPOUND:** Sodium perfluoro-1-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]octanesulfonate  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>4</sub><sup>12</sup>C<sub>4</sub>F<sub>17</sub>SO<sub>3</sub>Na **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% <sup>13</sup>C  
**LAST TESTED:** (mm/dd/yyyy) 05/15/2015 (1,2,3,4-<sup>13</sup>C<sub>4</sub>)  
**EXPIRY DATE:** (mm/dd/yyyy) 05/15/2020  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

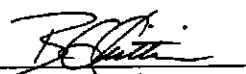
**DOCUMENTATION/ DATA ATTACHED:**

- 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 ~ 0.8% Sodium perfluoro-1-[1,2,3-<sup>13</sup>C<sub>3</sub>]heptanesulfonate.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
B.G. Chittim **Date:** 05/28/2015  
(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.

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### **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:**

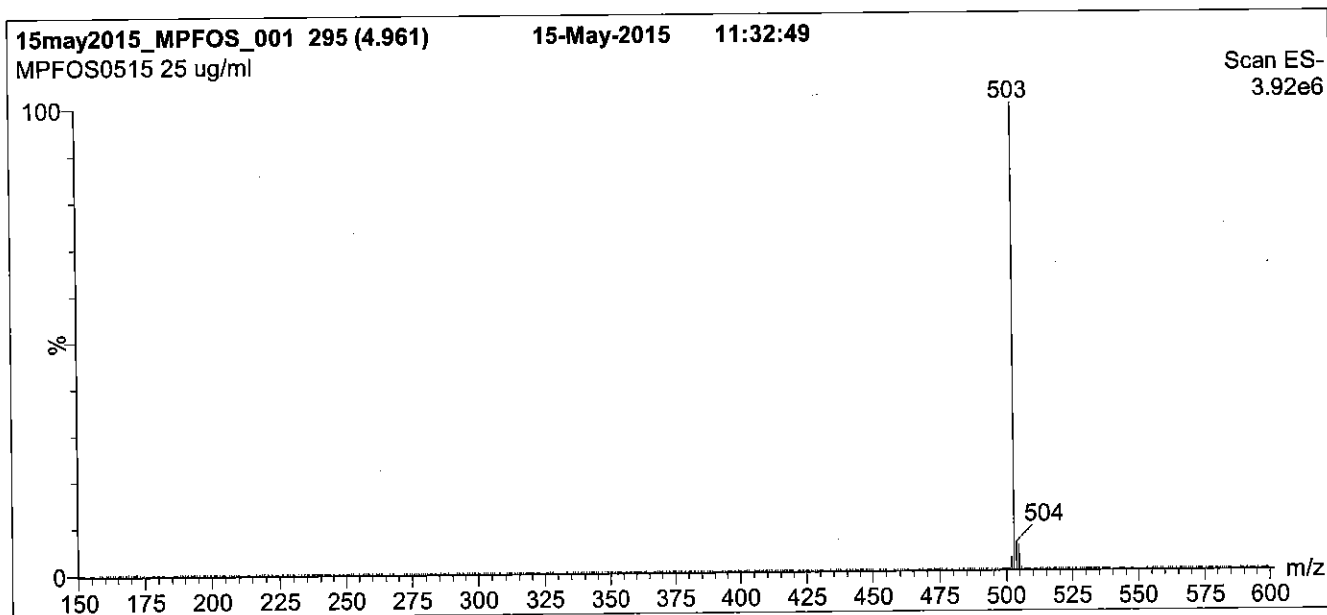
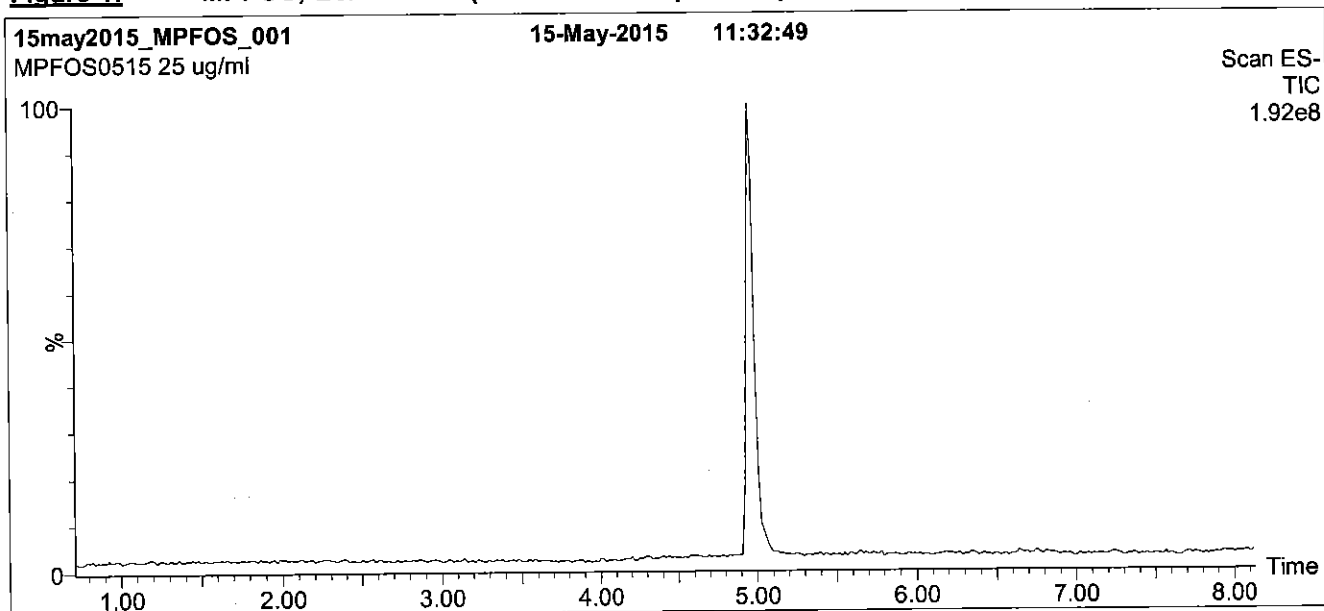
This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*



**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 45% (80:20 MeOH:ACN) / 55% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

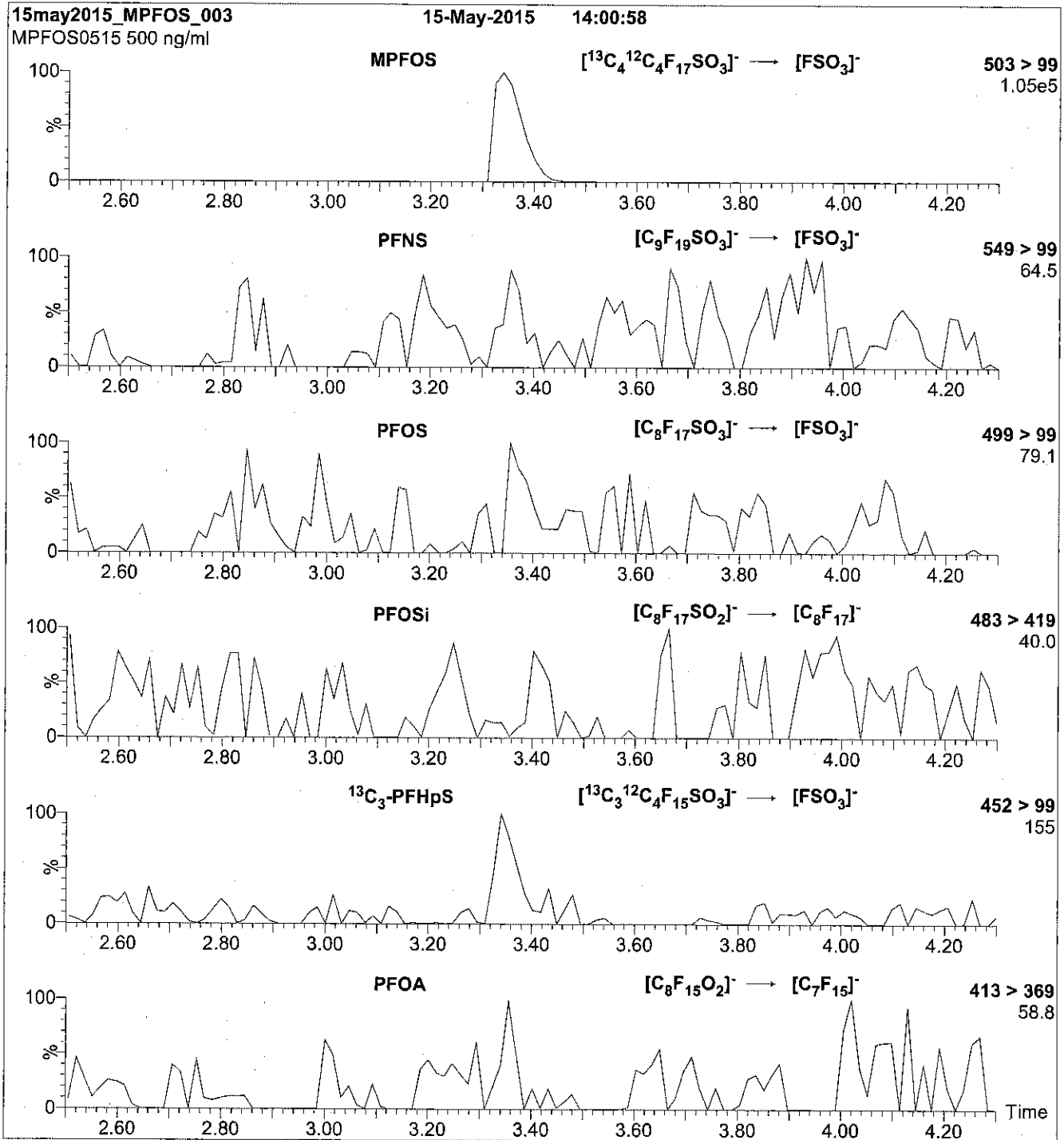
**Flow:** 300  $\mu$ l/min

**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: MPFOS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
10  $\mu\text{l}$  (500 ng/ml MPFOS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 40

Reagent

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**LCMPFUdA\_00004**

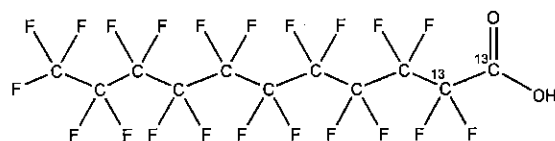
1:41/515 SKU



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFUdA **LOT NUMBER:** MPFUdA1014  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]undecanoic acid  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>9</sub>HF<sub>21</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 566.08  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
 (1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 10/31/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 10/31/2019  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

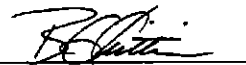
### DOCUMENTATION/ DATA ATTACHED:

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.
- Presence of 1-<sup>13</sup>C<sub>1</sub>-PFUdA (~1%; see Figure 2), 2-<sup>13</sup>C<sub>1</sub>-PFUdA (~1%), and PFUdA (~0.2%; see Figure 2) are due to the isotopic purity of the <sup>13</sup>C-precursor.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim **Date:** 11/03/2014  
(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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$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

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

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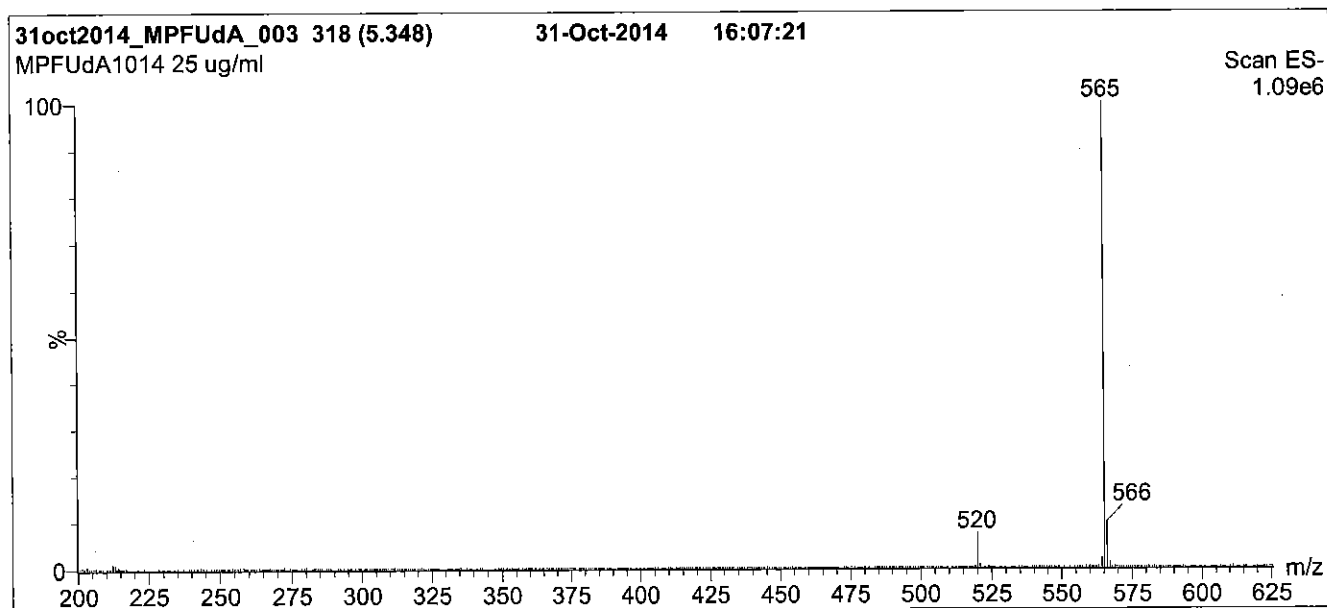
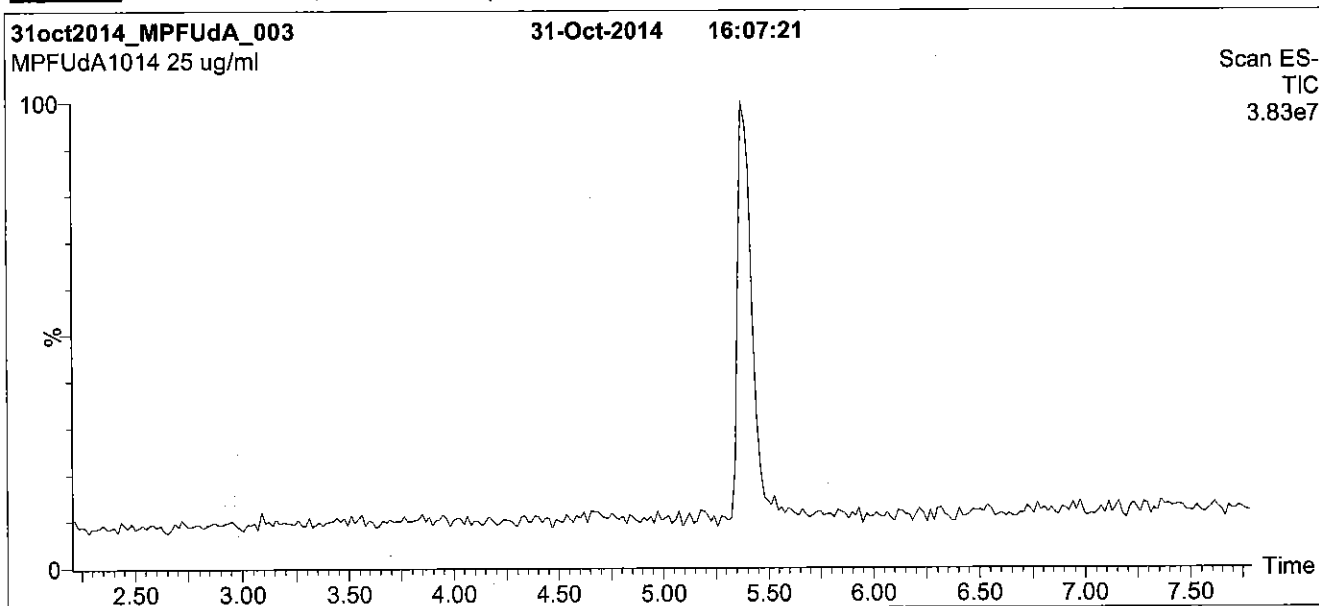
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

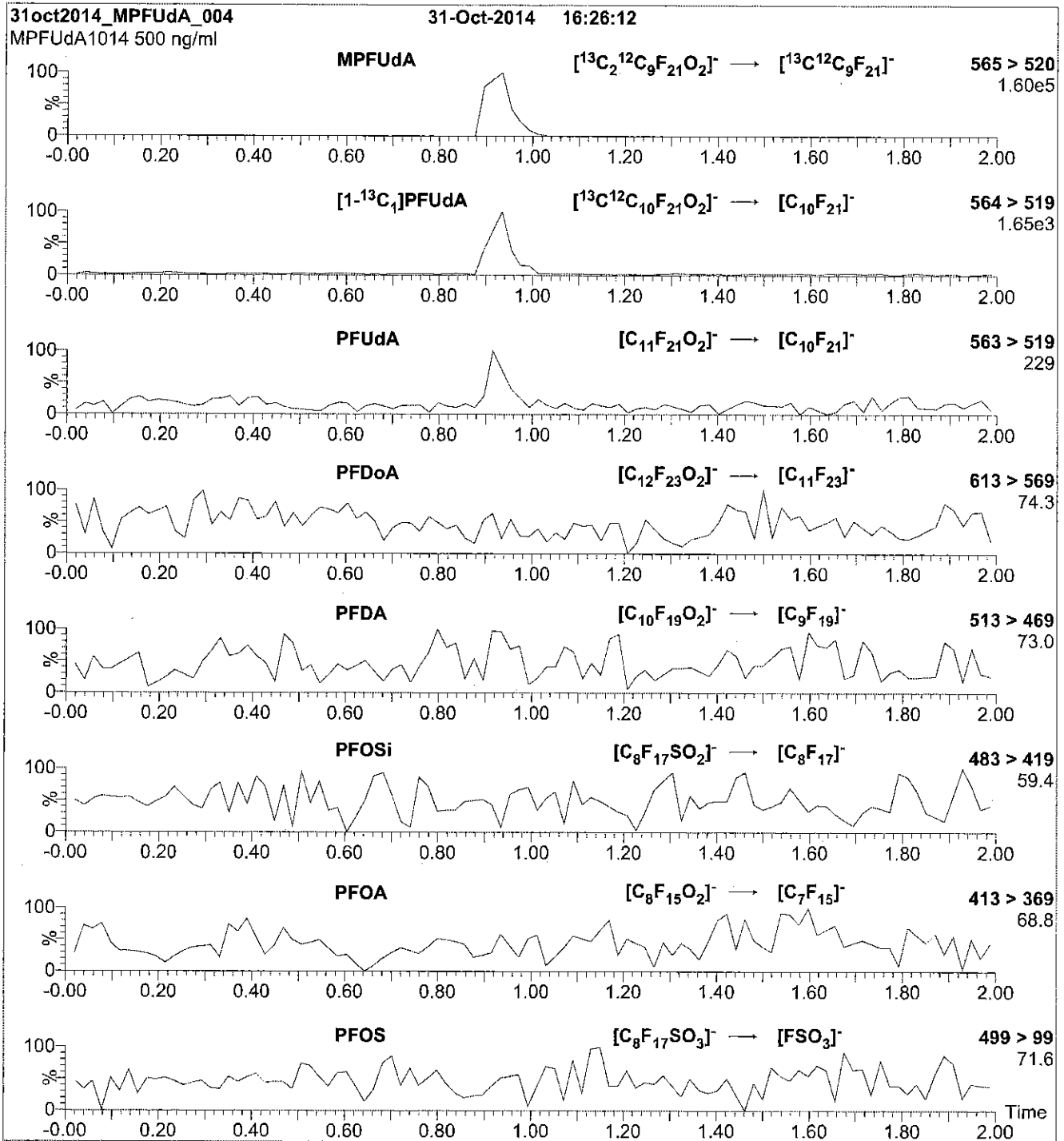
**Flow:** 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (200 - 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  $\mu\text{l}$  (500 ng/ml MPFUdA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.46e-3  
Collision Energy (eV) = 11

Reagent

---

**LCMPFUdA\_00005**

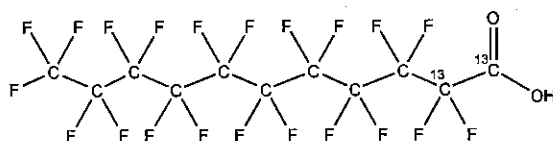




# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** MPFUdA **LOT NUMBER:** MPFUdA1014  
**COMPOUND:** Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]undecanoic acid  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** <sup>13</sup>C<sub>2</sub><sup>12</sup>C<sub>9</sub>HF<sub>21</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 566.08  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98% **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
 (1,2-<sup>13</sup>C<sub>2</sub>)  
**LAST TESTED:** (mm/dd/yyyy) 10/31/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 10/31/2019  
**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.
- Presence of 1-<sup>13</sup>C<sub>1</sub>-PFUdA (~1%; see Figure 2), 2-<sup>13</sup>C<sub>1</sub>-PFUdA (~1%), and PFUdA (~0.2%; see Figure 2) are due to the isotopic purity of the <sup>13</sup>C-precursor.

**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.

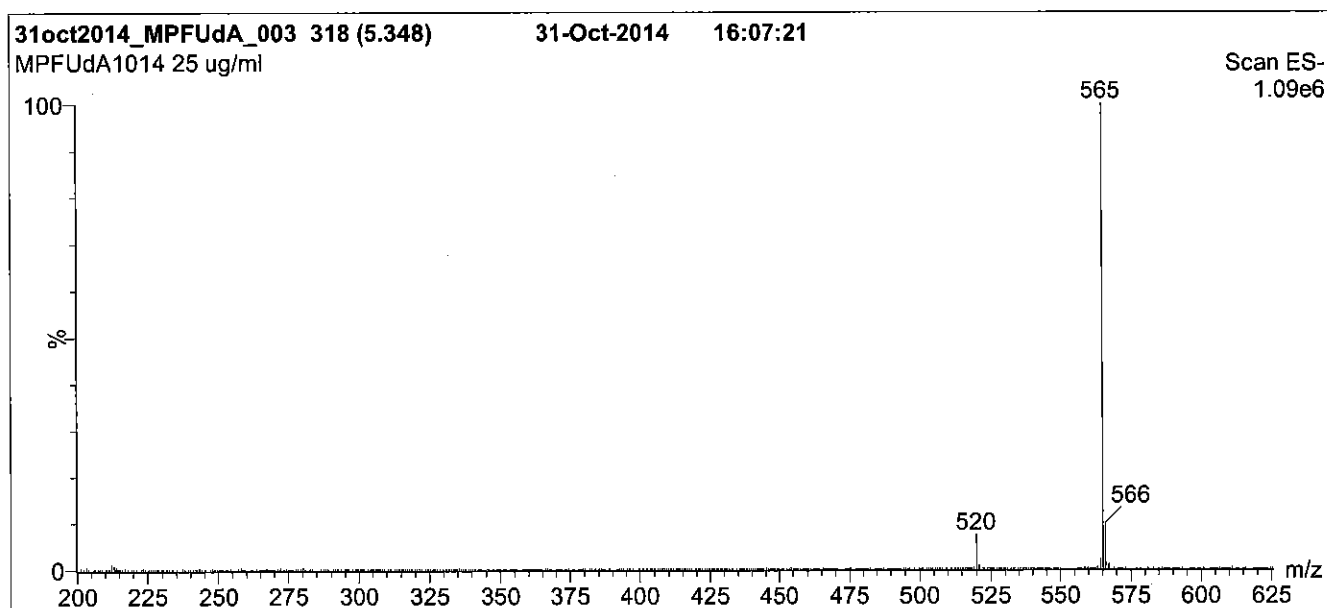
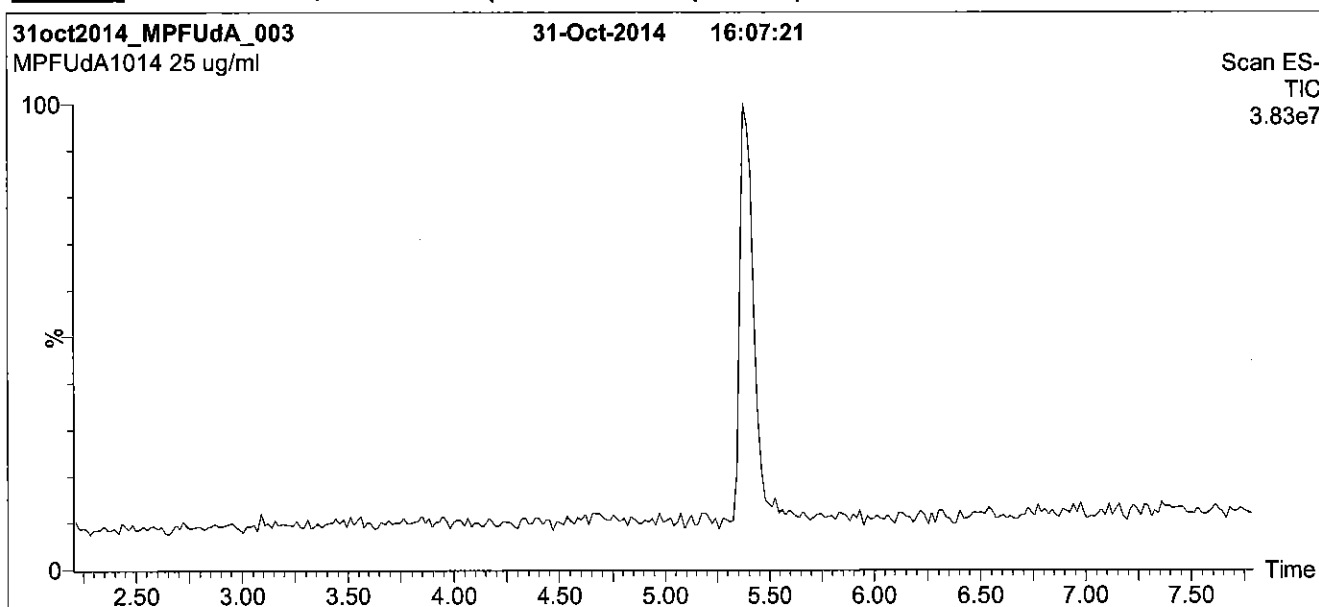
### **QUALITY MANAGEMENT:**

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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

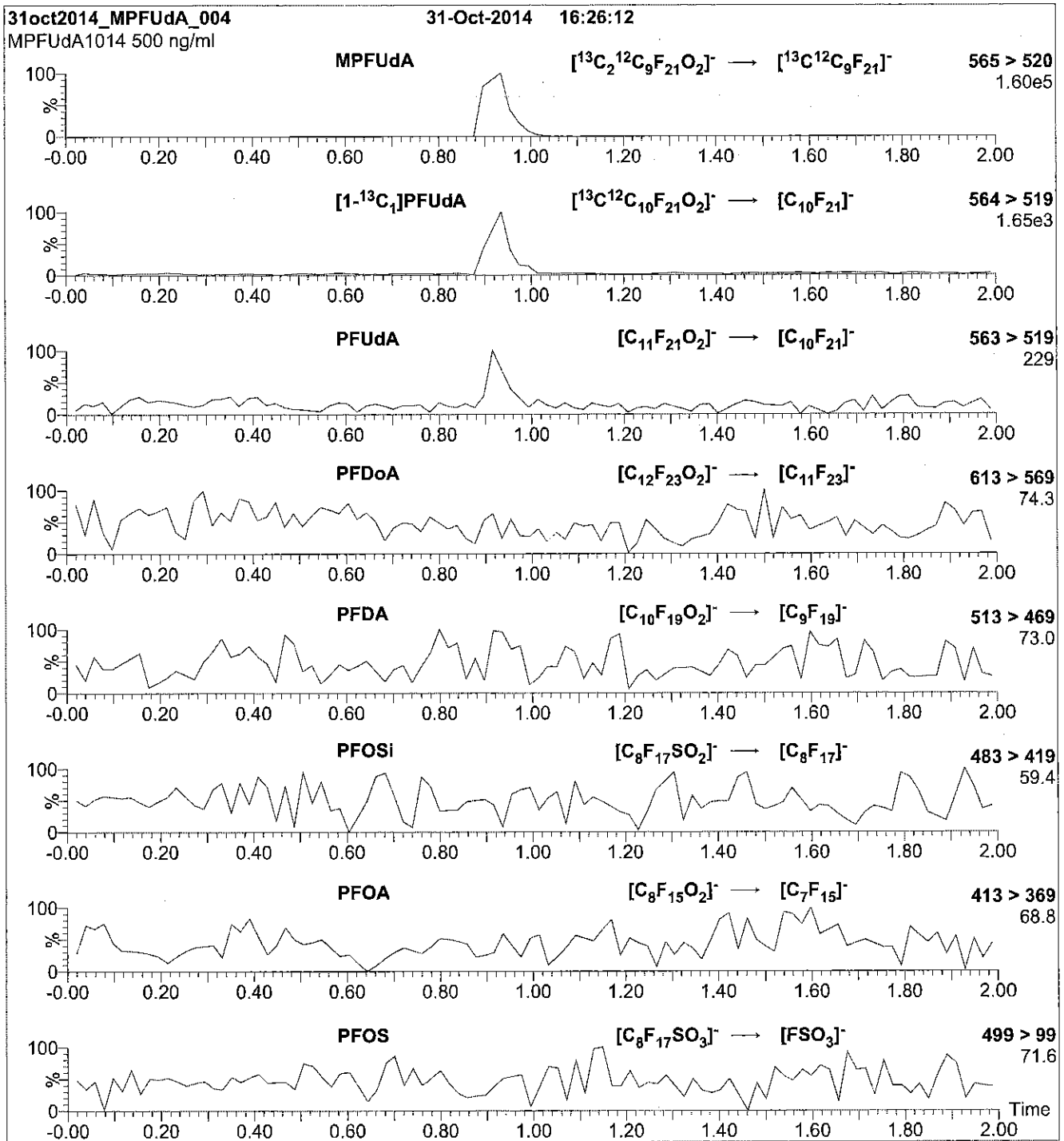
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (200 - 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  $\mu\text{l}$  (500 ng/ml MPFUdA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
 (both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

**MS Parameters**

Collision Gas (mbar) = 3.46e-3  
 Collision Energy (eV) = 11

Reagent

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**LCPFBA\_00003**

rec 7/15/14



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:**

PFBA

**LOT NUMBER:**

PFBA0313

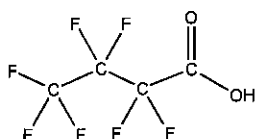
**COMPOUND:**

Perfluoro-n-butanoic acid

**STRUCTURE:**

**CAS #:**

375-22-4



**MOLECULAR FORMULA:**

C<sub>4</sub>HF<sub>7</sub>O<sub>2</sub>

**MOLECULAR WEIGHT:**

214.04

**CONCENTRATION:**

50 ± 2.5 µg/ml

**SOLVENT(S):**

Methanol  
Water (<1%)

**CHEMICAL PURITY:**

>98%

**LAST TESTED:** (mm/dd/yyyy)

03/05/2013

**EXPIRY DATE:** (mm/dd/yyyy)

03/05/2018

**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.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

Certified By:

B.G. Chittim

Date: 03/06/2013

(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

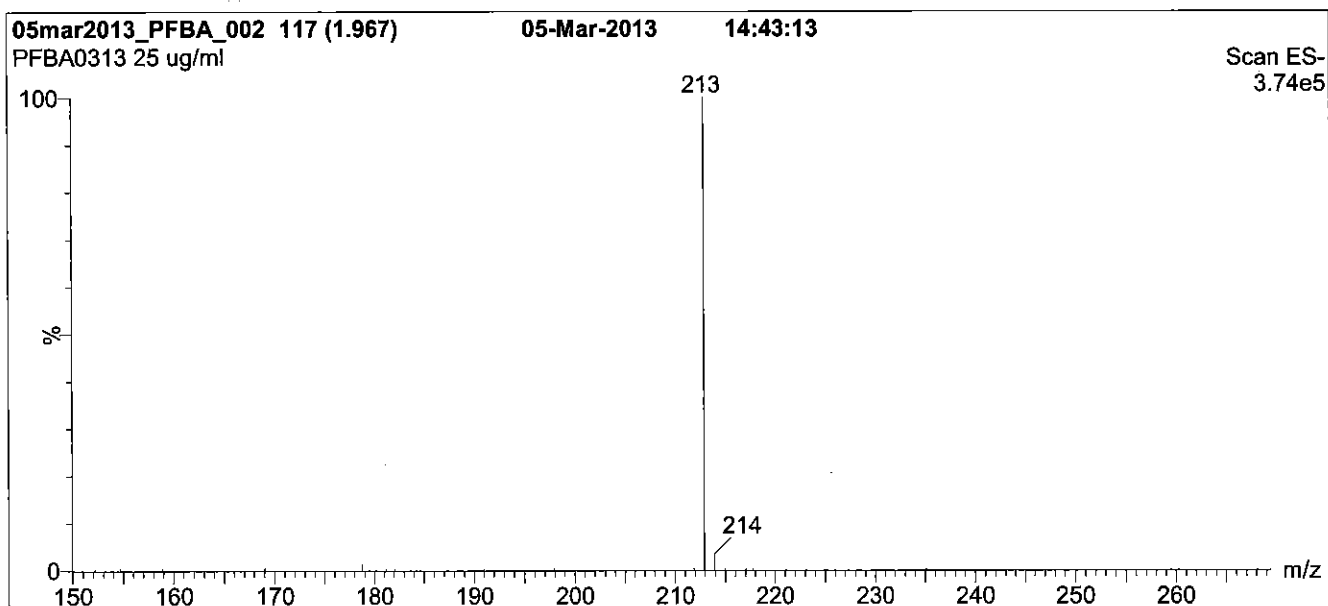
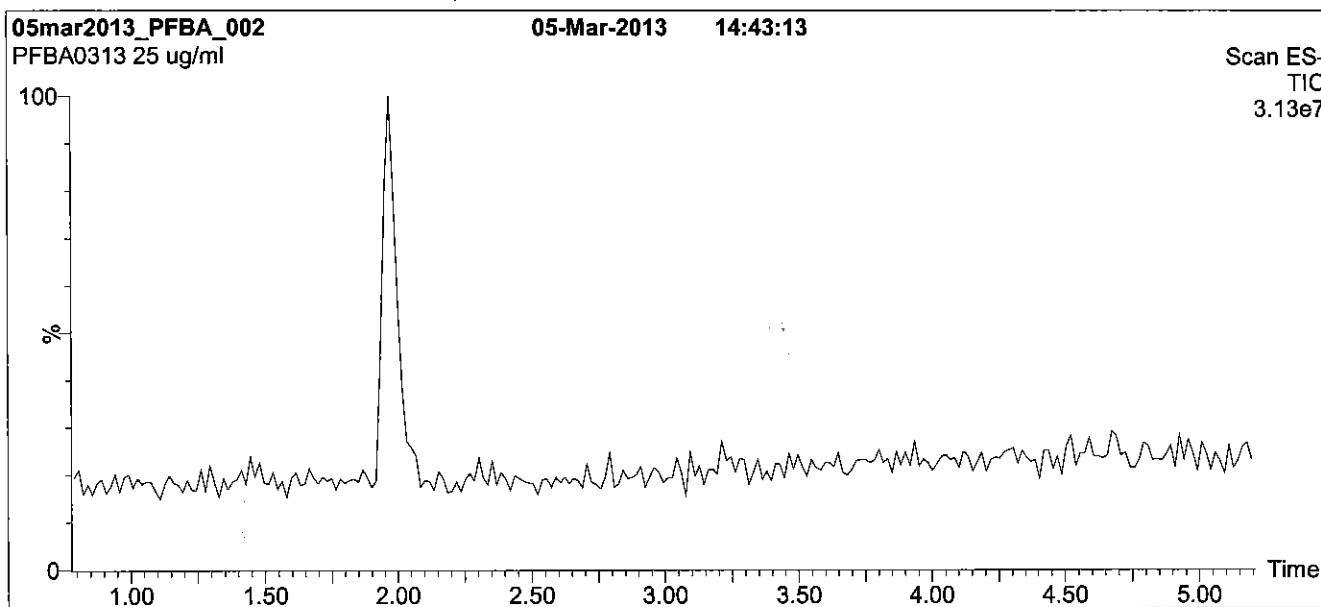
### **QUALITY MANAGEMENT:**

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**Figure 1: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 25% (80:20 MeOH:ACN) / 75% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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  $\mu$ l/min

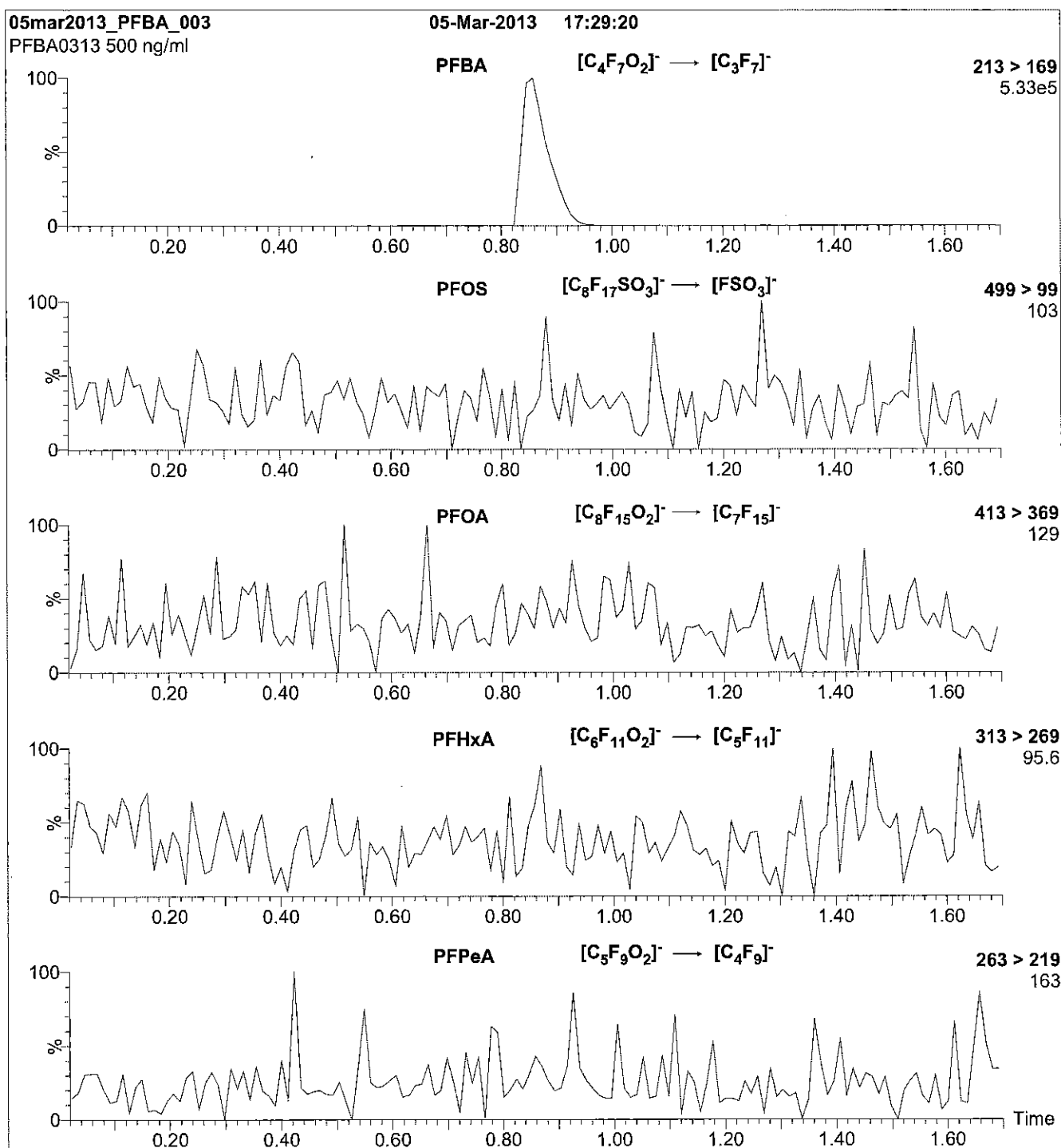
**MS Parameters**

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  $\mu$ l (500 ng/ml PFBA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

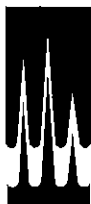
**MS Parameters**

Collision Gas (mbar) = 3.70e-3  
Collision Energy (eV) = 10

Reagent

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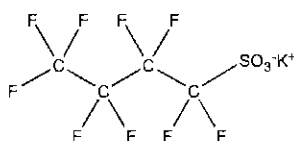
**LCPFBS\_00003**



# 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:** C<sub>4</sub>F<sub>9</sub>SO<sub>3</sub>K **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:** 10/17/2014  
 (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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**HAZARDS:**

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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.

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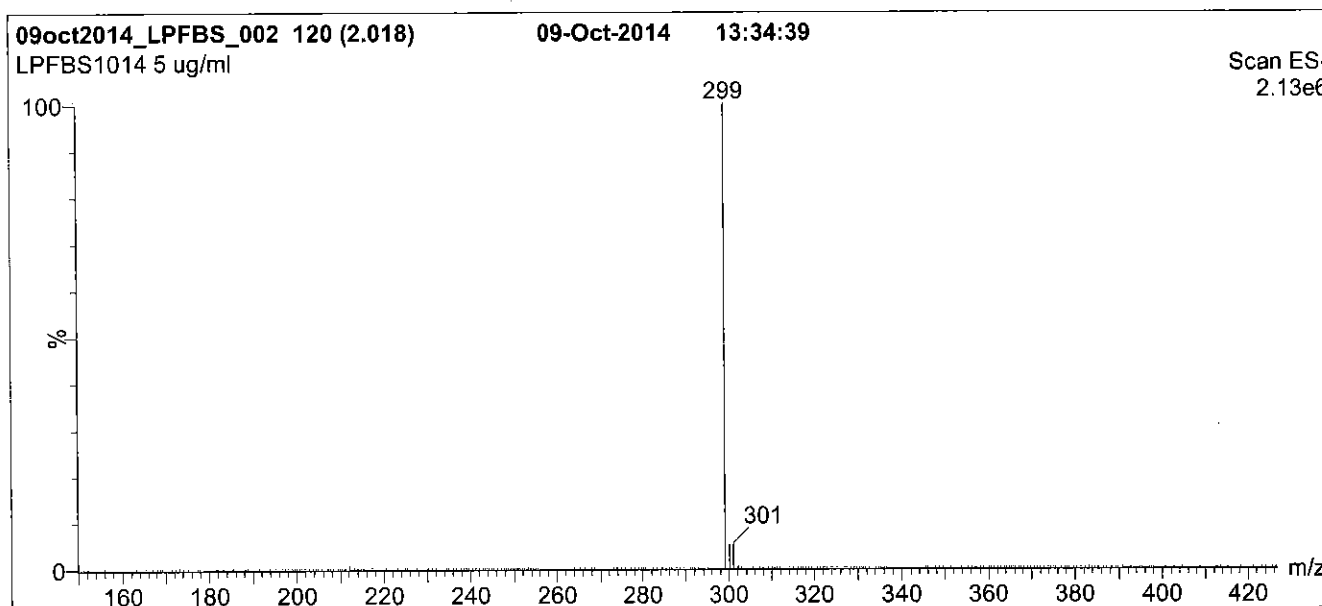
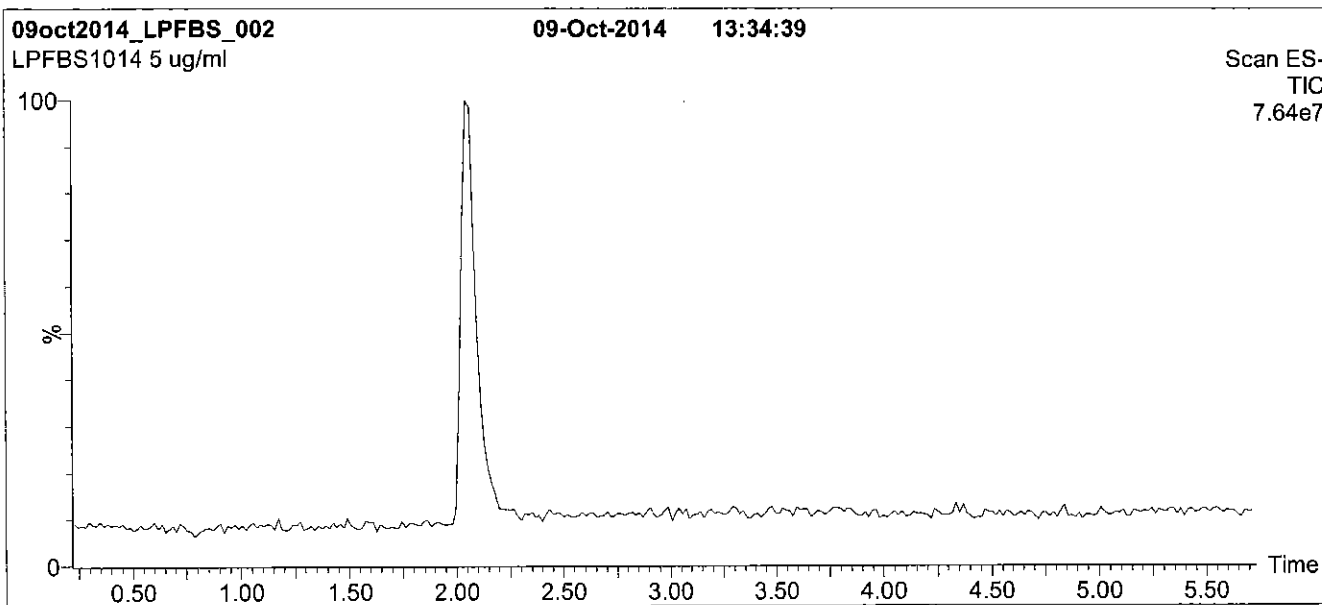
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**Figure 1: L-PFBS; LC/MS Data (TIC and Mass Spectrum)**



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

**Column:** Acquity UPLC BEH Shield RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

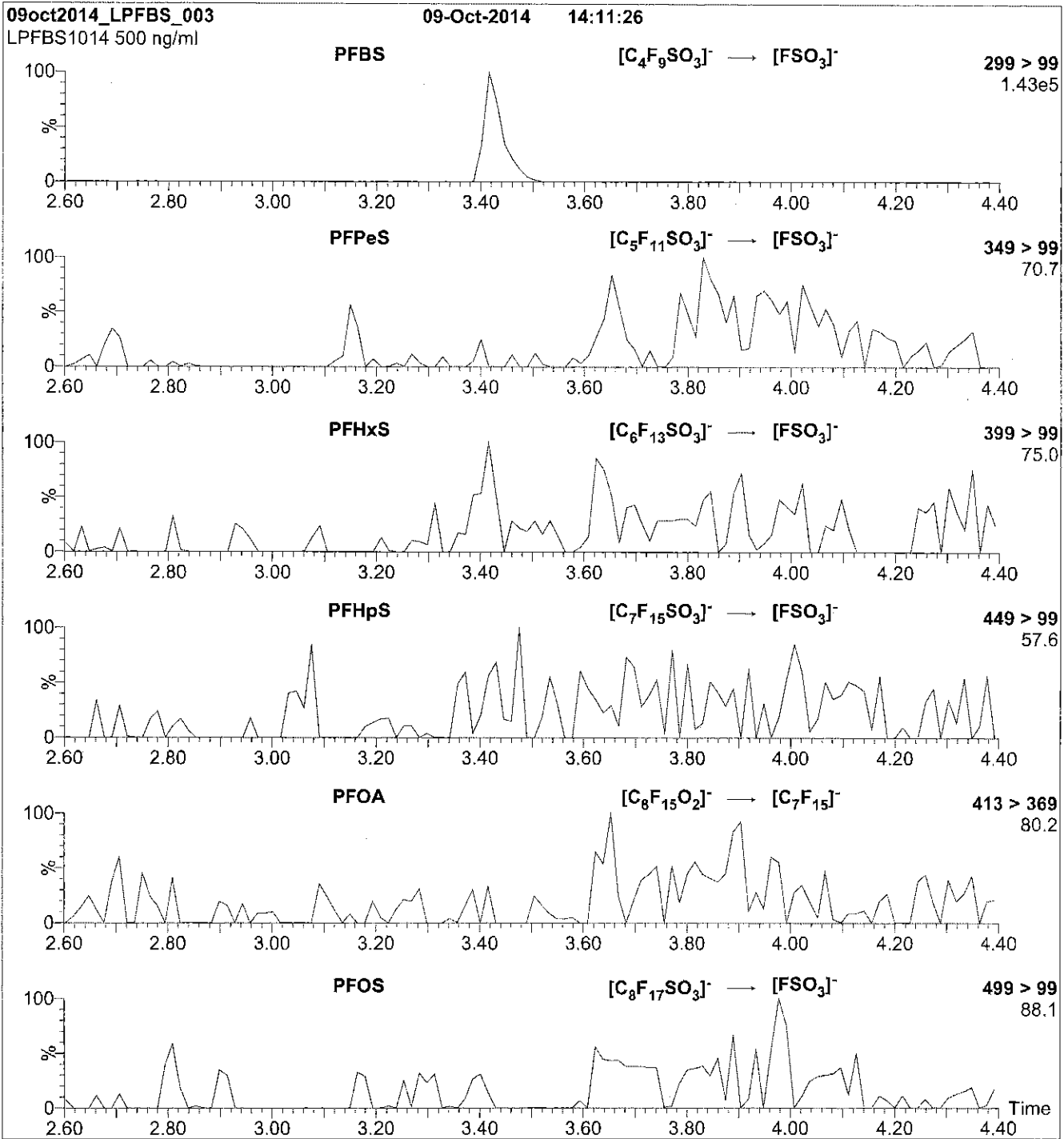
**Flow:** 300  $\mu$ l/min

**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  $\mu$ l (500 ng/ml L-PFBS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

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

Reagent

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**LCPFDA\_00003**

rec 7/16/14



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:**

PFDA

**LOT NUMBER:**

PFDA0613

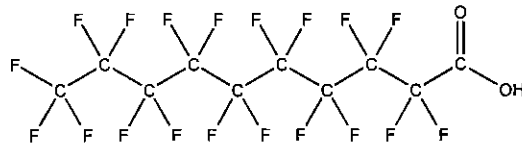
**COMPOUND:**

Perfluoro-n-decanoic acid

**STRUCTURE:**

**CAS #:**

335-76-2



**MOLECULAR FORMULA:**

C<sub>10</sub>H<sub>F<sub>19</sub></sub>O<sub>2</sub>

**MOLECULAR WEIGHT:**

514.08

**CONCENTRATION:**

50 ± 2.5 µg/ml

**SOLVENT(S):**

Methanol  
Water (<1%)

**CHEMICAL PURITY:**

>98%

**LAST TESTED:** (mm/dd/yyyy)

06/19/2013

**EXPIRY DATE:** (mm/dd/yyyy)

06/19/2018

**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 ~ 0.4% PFNA and ~ 0.1% PFOA.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

Certified By:

B.G. Chittim

Date: 07/03/2013

(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

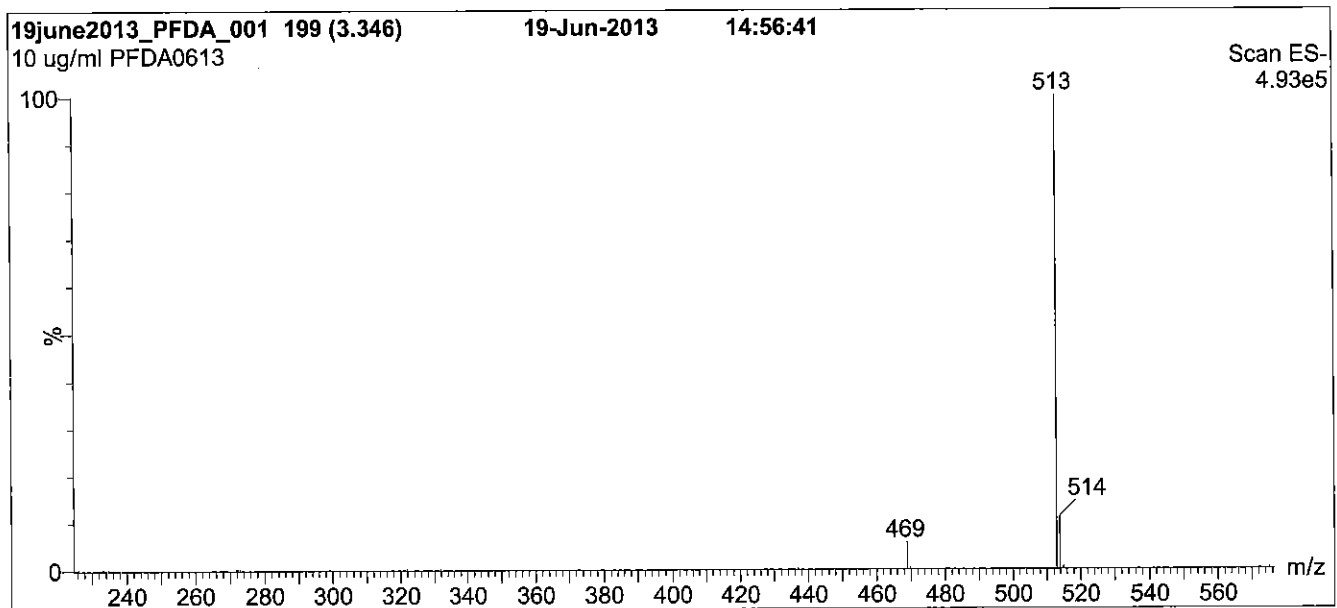
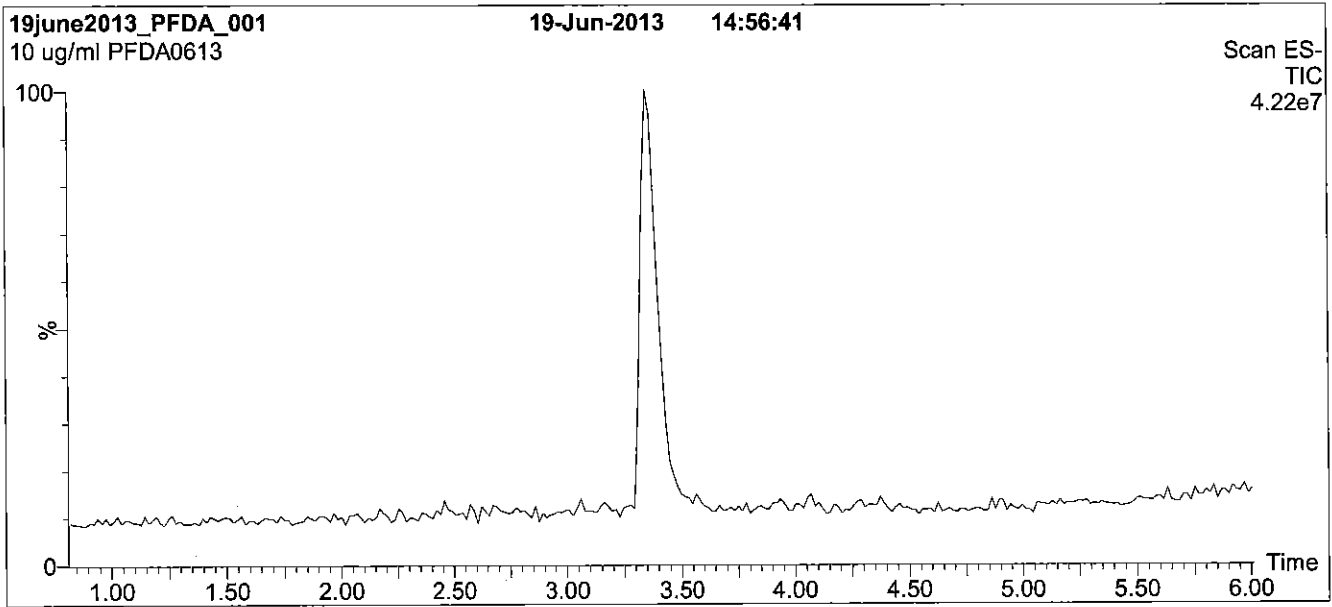
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

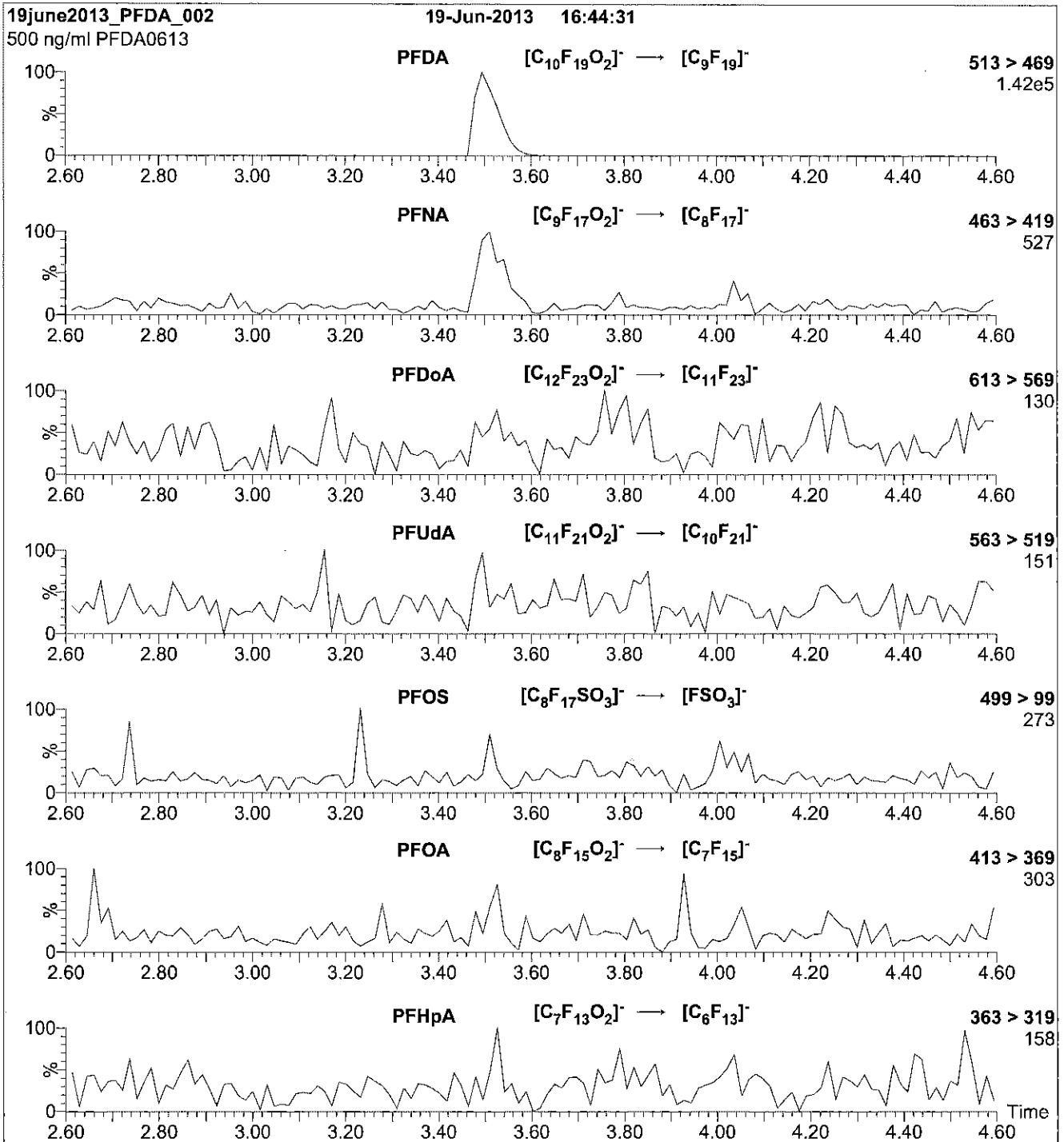
Flow: 300  $\mu$ l/min

**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: PFDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct loop injection  
10  $\mu$ l (500 ng/ml PFDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.58e-3  
Collision Energy (eV) = 13

Reagent

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**LCPFDoA\_00003**

Rec 7/15



# WELLINGTON LABORATORIES

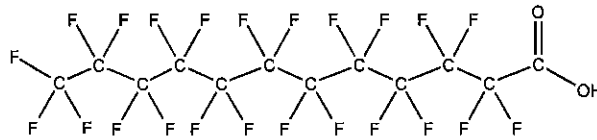
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFD0A  
**COMPOUND:** Perfluoro-n-dodecanoic acid

**LOT NUMBER:** PFD0A0113

**STRUCTURE:**

**CAS #:** 307-55-1



**MOLECULAR FORMULA:**  $C_{12}HF_{23}O_2$   
**CONCENTRATION:**  $50 \pm 2.5 \mu\text{g/ml}$

**MOLECULAR WEIGHT:** 614.10  
**SOLVENT(S):** Methanol  
 Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 01/03/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 01/03/2018  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

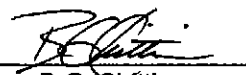
**DOCUMENTATION/ DATA ATTACHED:**

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

**ADDITIONAL INFORMATION:**

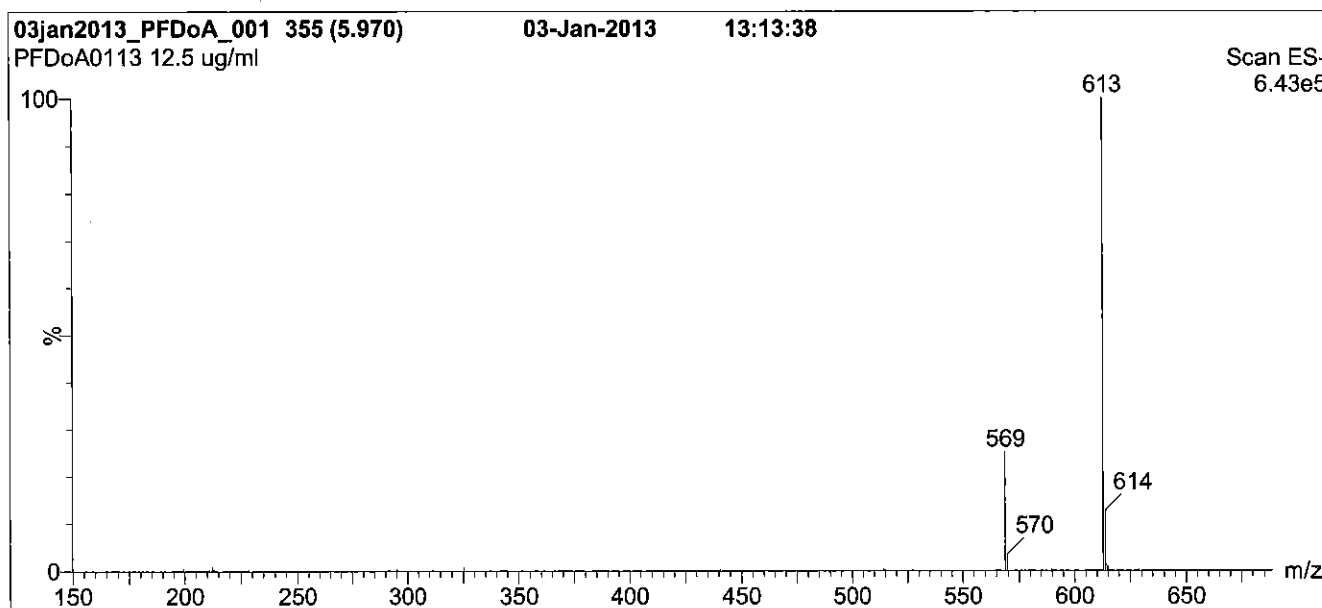
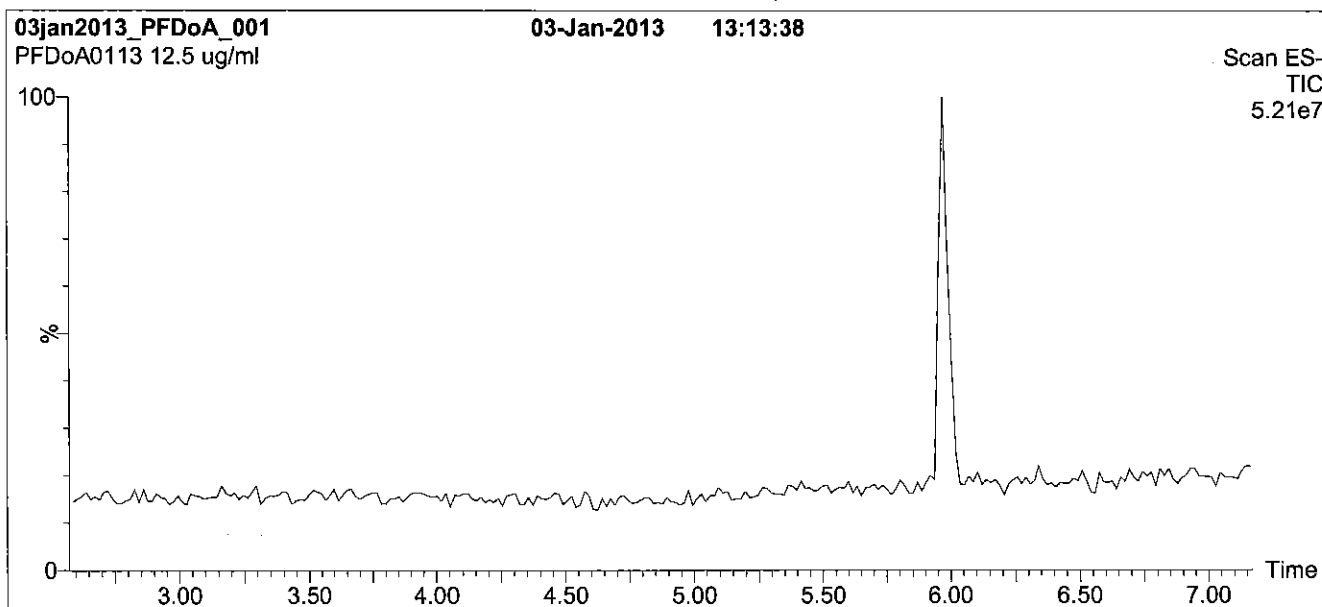
- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim  
**Date:** 02/01/2013  
 (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

**Figure 1: PFD<sub>o</sub>A; 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<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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  $\mu$ l/min

**MS Parameters**

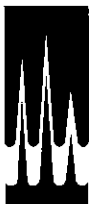
**Experiment:** Full Scan (150 - 850 amu)

**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

Reagent

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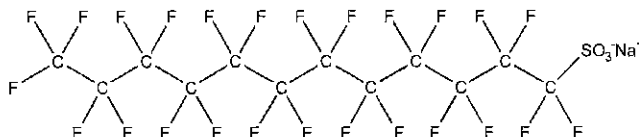
**LCPFDoS\_00003**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFDoS **LOT NUMBER:** LPFDoS1011  
**COMPOUND:** Sodium perfluoro-1-dodecanesulfonate  
**STRUCTURE:** **CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>12</sub>F<sub>25</sub>SO<sub>3</sub>Na **MOLECULAR WEIGHT:** 722.14  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol  
 48.4 ± 2.4 µg/ml (PFDoS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 10/06/2011  
**EXPIRY DATE:** (mm/dd/yyyy) 10/06/2016  
**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 ~ 0.3% of sodium perfluoro-1-tetradecanesulfonate and ~ 0.8% of perfluoro-n-dodecanoic acid (PFDoA).

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim **Date:** 01/15/2013  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

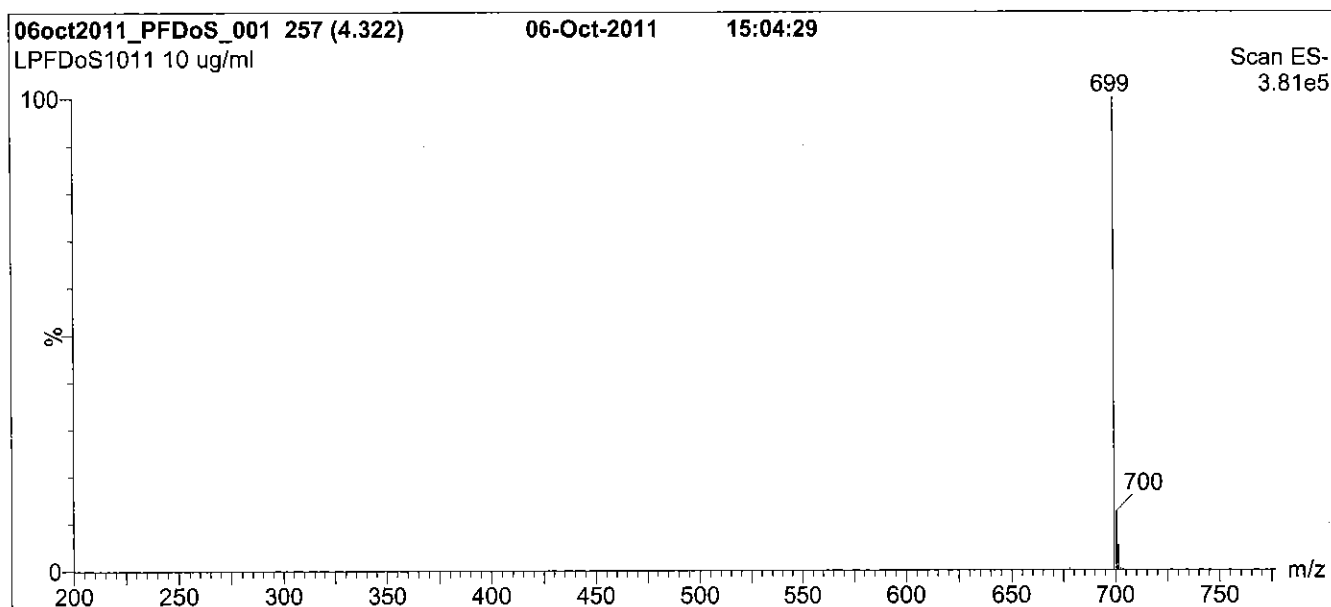
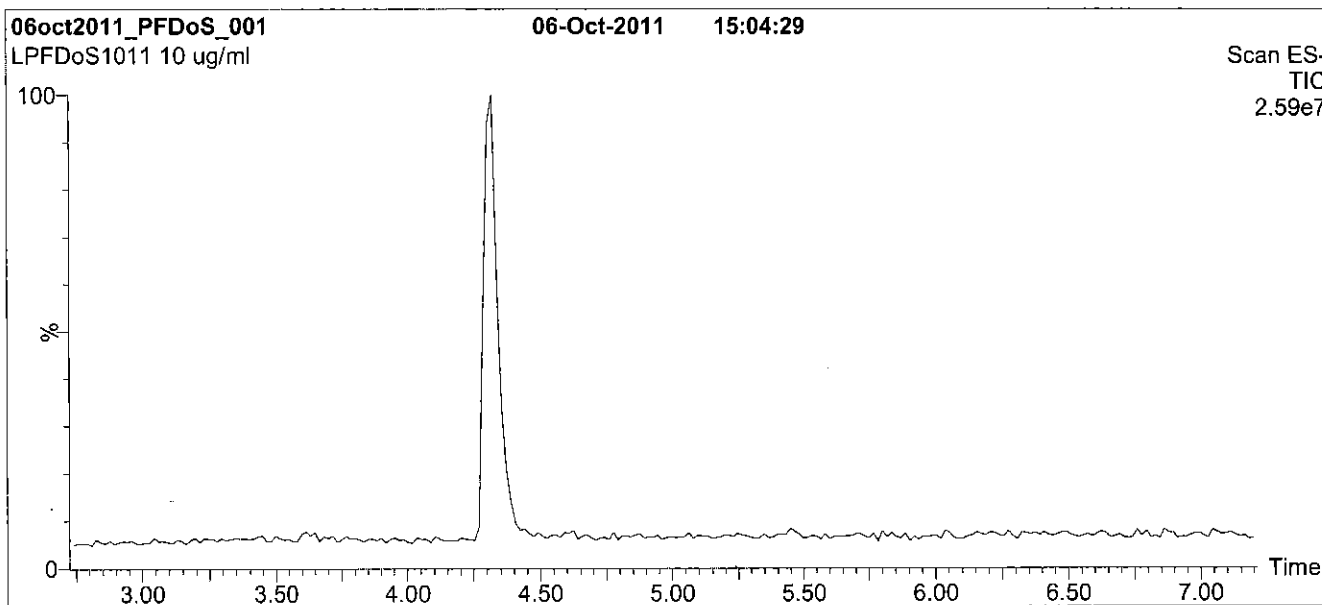
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number AR-1523).



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**Figure 1: L-PFDoS; 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<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 65% (80:20 MeOH:ACN) / 35% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

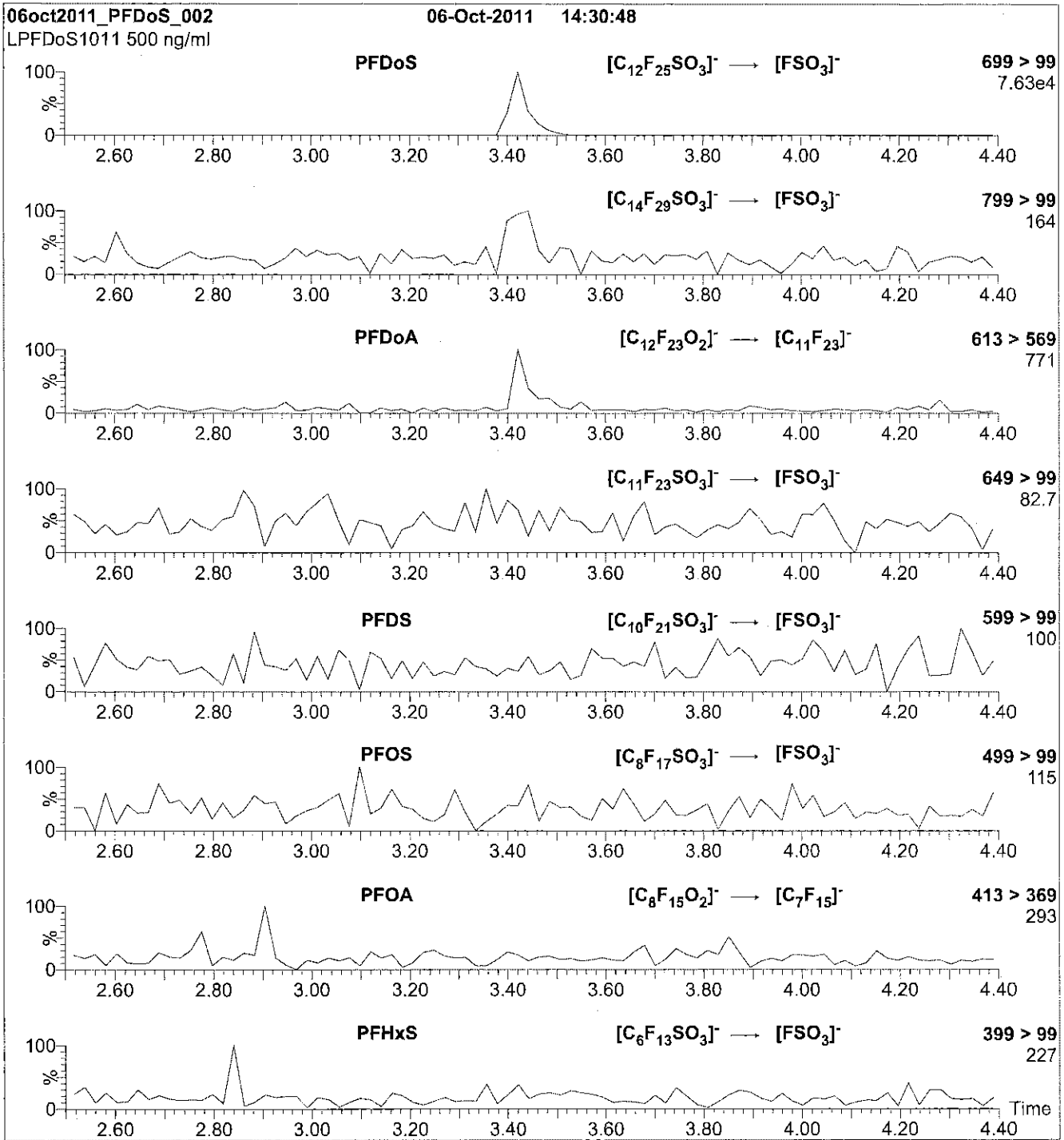
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (200 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 3.00  
Cone Voltage (V) = 80.00  
Cone Gas Flow (l/hr) = 50  
Desolvation Gas Flow (l/hr) = 750

**Figure 2: L-PFDoS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
10  $\mu$ l (500 ng/ml L-PFDoS)

**Mobile phase:** Isocratic 65% (80:20 MeOH:ACN) / 35% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

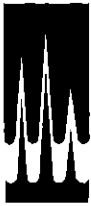
Collision Gas (mbar) = 3.54e-3  
Collision Energy (eV) = 50

Reagent

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**LCPFDS\_00003**

P: 2/11/15 SV



# WELLINGTON LABORATORIES

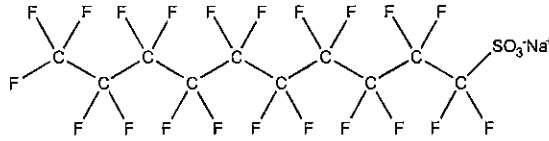
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFDS  
**COMPOUND:** Sodium perfluoro-1-decanesulfonate

**LOT NUMBER:** LPFDS0913

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>10</sub>F<sub>21</sub>SO<sub>3</sub>Na  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt)  
48.2 ± 2.4 µg/ml (PFDS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 09/13/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 09/13/2018  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**MOLECULAR WEIGHT:** 622.13  
**SOLVENT(S):** Methanol

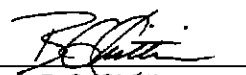
**DOCUMENTATION/ DATA ATTACHED:**

- 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:** 09/23/2013  
(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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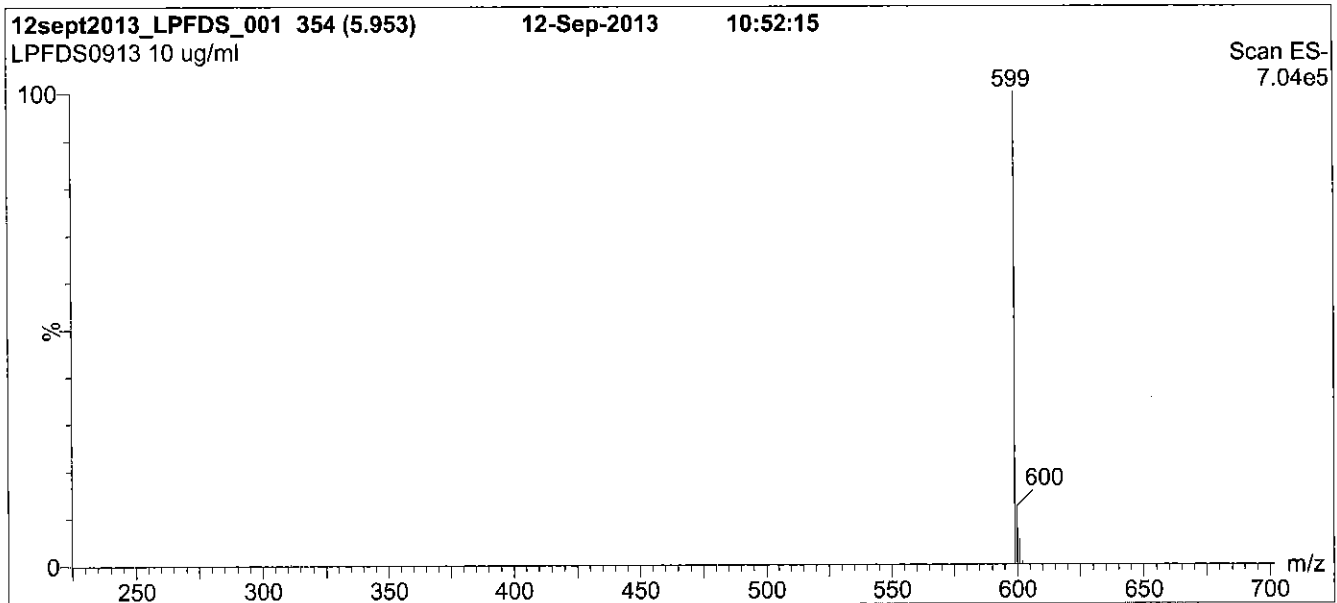
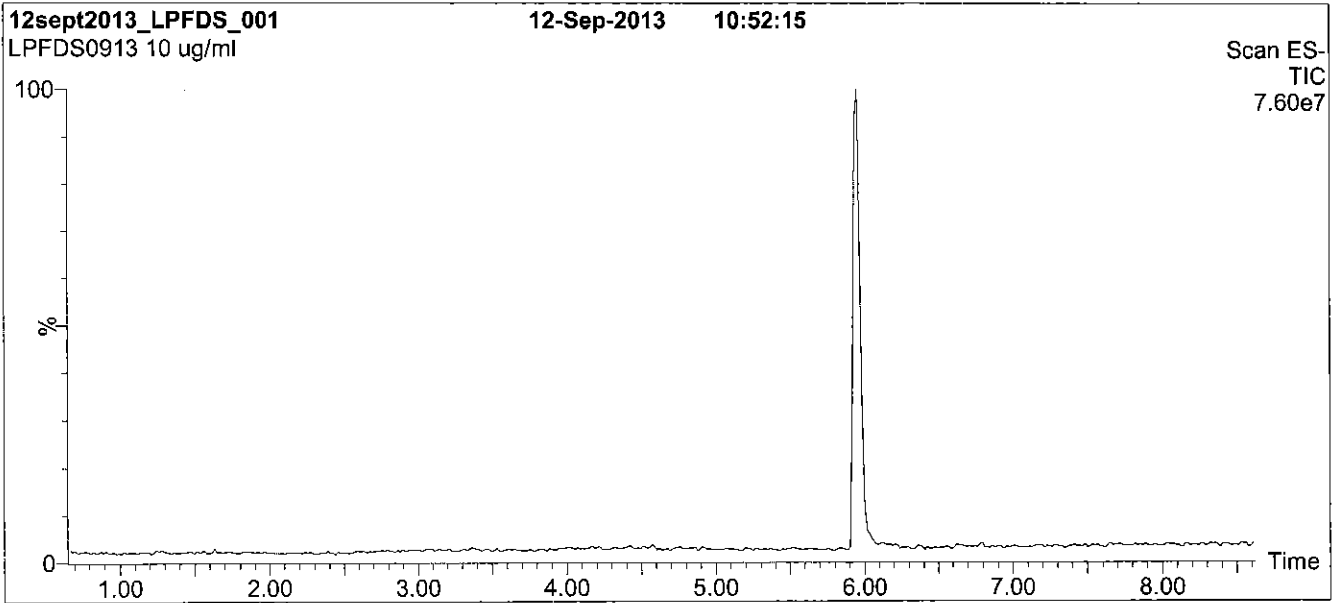
**QUALITY MANAGEMENT:**

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**Figure 1: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 45% (80:20 MeOH:ACN) / 55% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min and hold for  
1.5 min before returning to initial conditions in 0.5 min.  
Time: 11 min

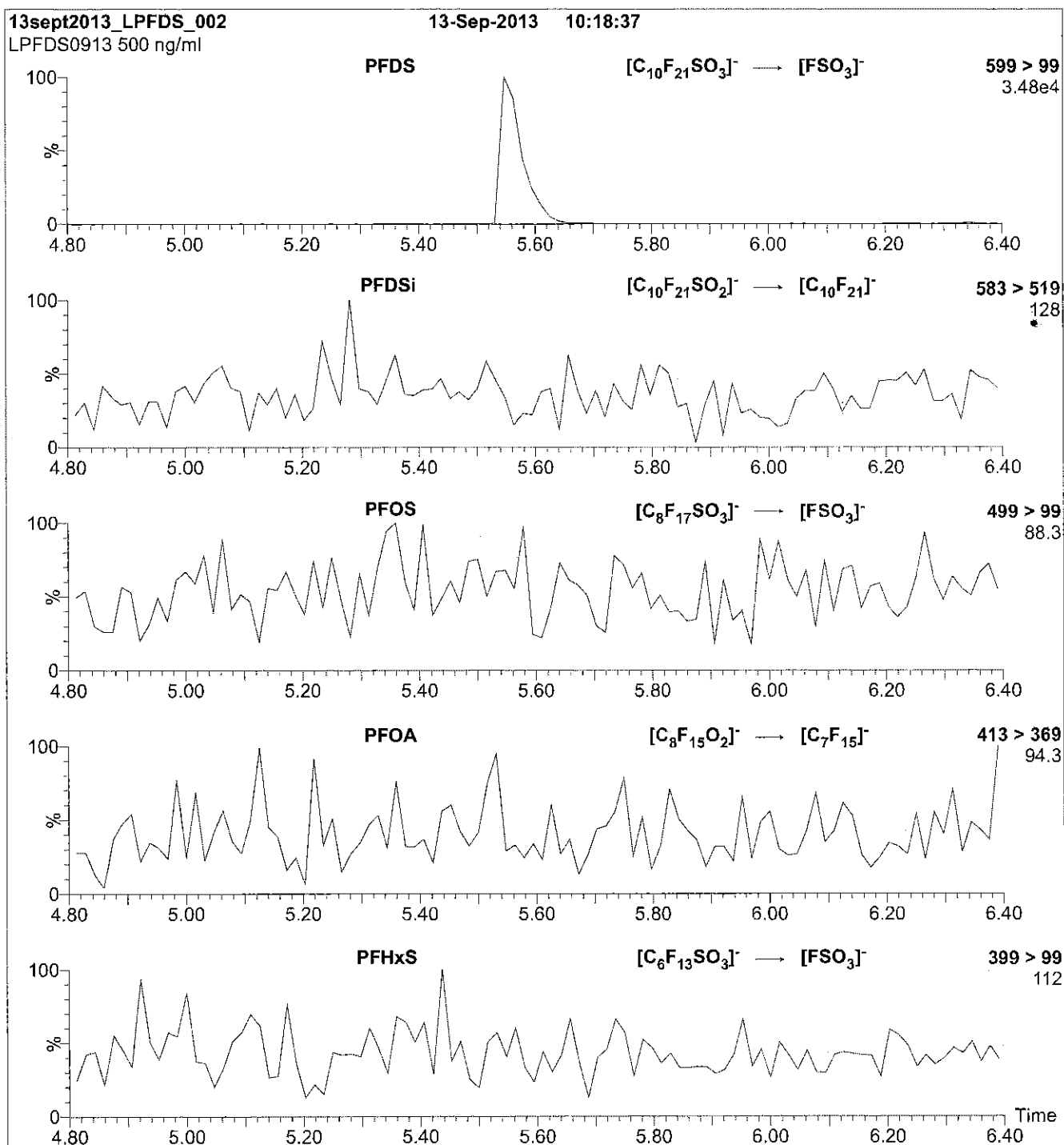
**Flow:** 300  $\mu$ l/min

**MS Parameters**

**Experiment:** Full Scan (225 - 850 amu)

**Source:** Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 70.00  
Cone Gas Flow (l/hr) = 60  
Desolvation Gas Flow (l/hr) = 650

**Figure 2: L-PFDS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct loop injection  
 10  $\mu$ l (500 ng/ml L-PFDS)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.78e-3  
 Collision Energy (eV) = 50



Reagent

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**LCPFHpA\_00004**



# WELLINGTON LABORATORIES

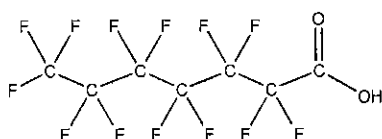
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFHpA  
**COMPOUND:** Perfluoro-n-heptanoic acid

**LOT NUMBER:** PFHpA0514

**STRUCTURE:**

**CAS #:** 375-85-9



**MOLECULAR FORMULA:** C<sub>7</sub>HF<sub>13</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 364.06  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 05/09/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 05/09/2019  
**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.
- 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: 05/22/2014  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

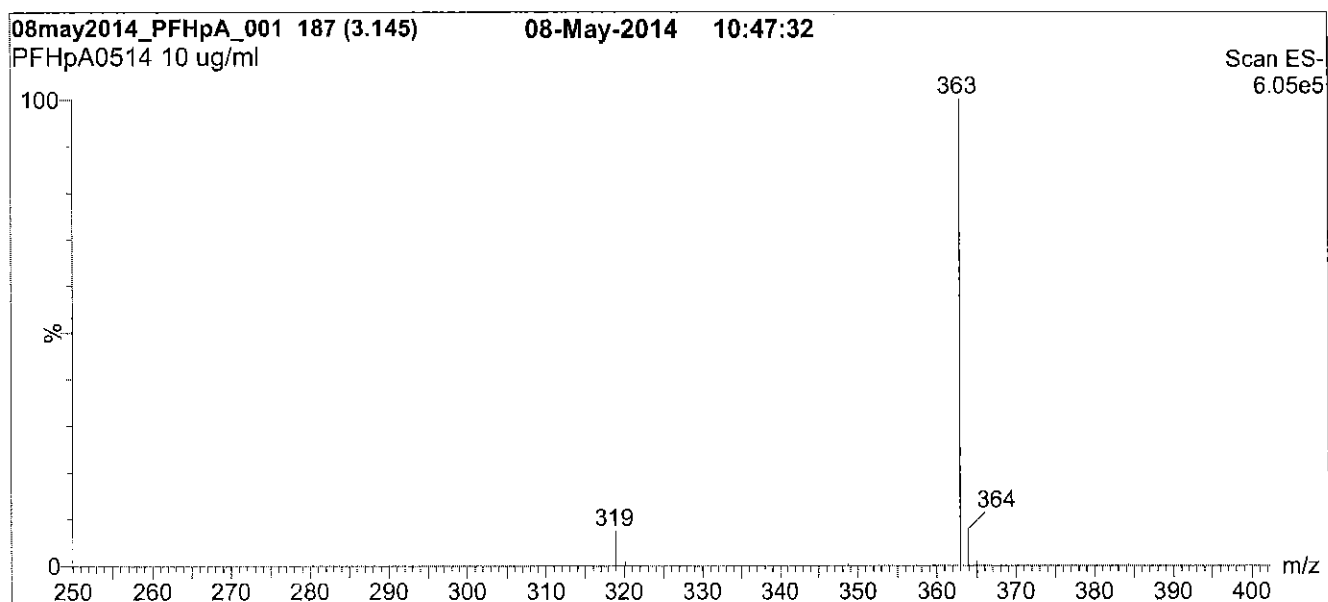
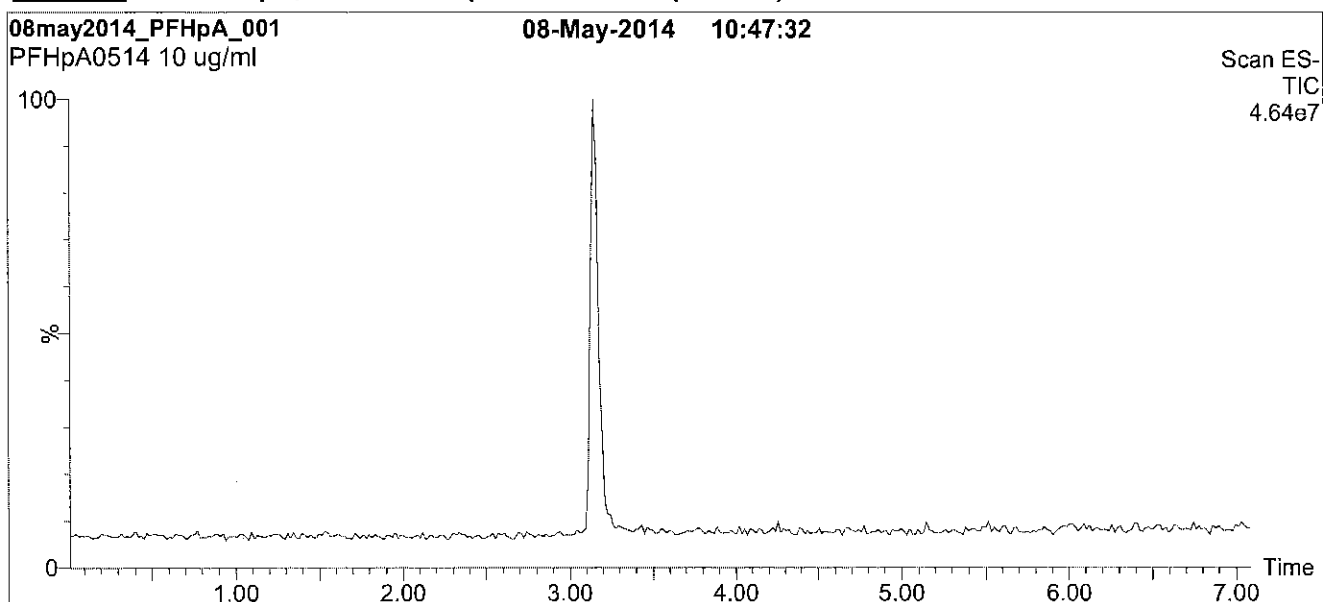
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto: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 C<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

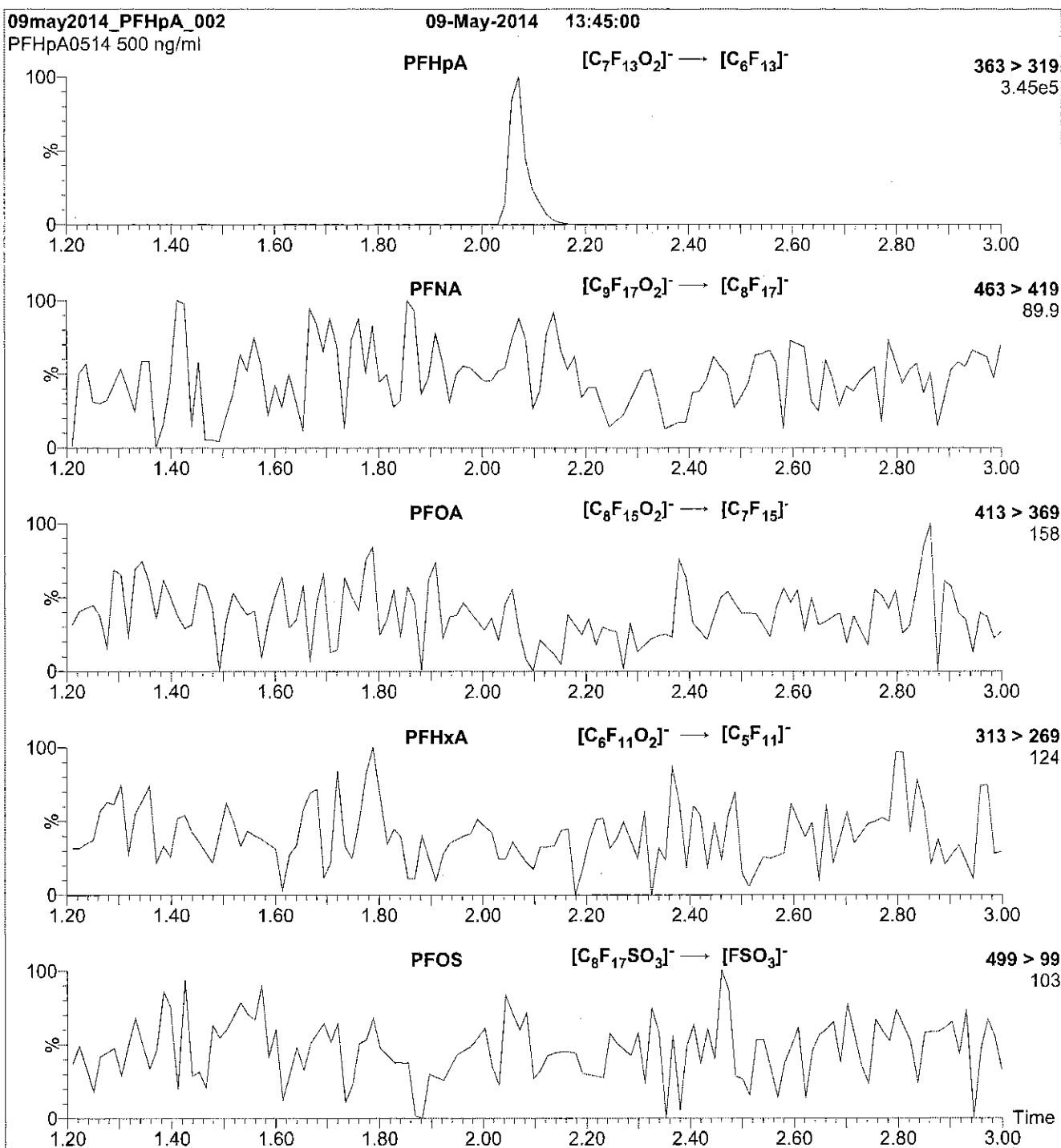
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (250 - 950 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  $\mu$ l (500 ng/ml PFHpA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.24e-3  
Collision Energy (eV) = 11

Reagent

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**LCPFHpS\_00005**

R: 4/15/15 SW



# WELLINGTON LABORATORIES

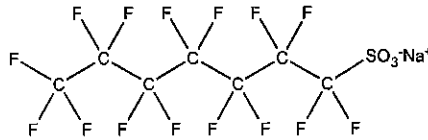
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFHpS  
**COMPOUND:** Sodium perfluoro-1-heptanesulfonate

**LOT NUMBER:** LPFHpS0114

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>7</sub>F<sub>15</sub>SO<sub>3</sub>Na  
**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) 01/28/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 01/28/2019  
**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 (C<sub>5</sub>F<sub>13</sub>SO<sub>3</sub>Na) and ~ 0.2% of L-PFOS (C<sub>8</sub>F<sub>17</sub>SO<sub>3</sub>Na).

**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:**

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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.

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### **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:**

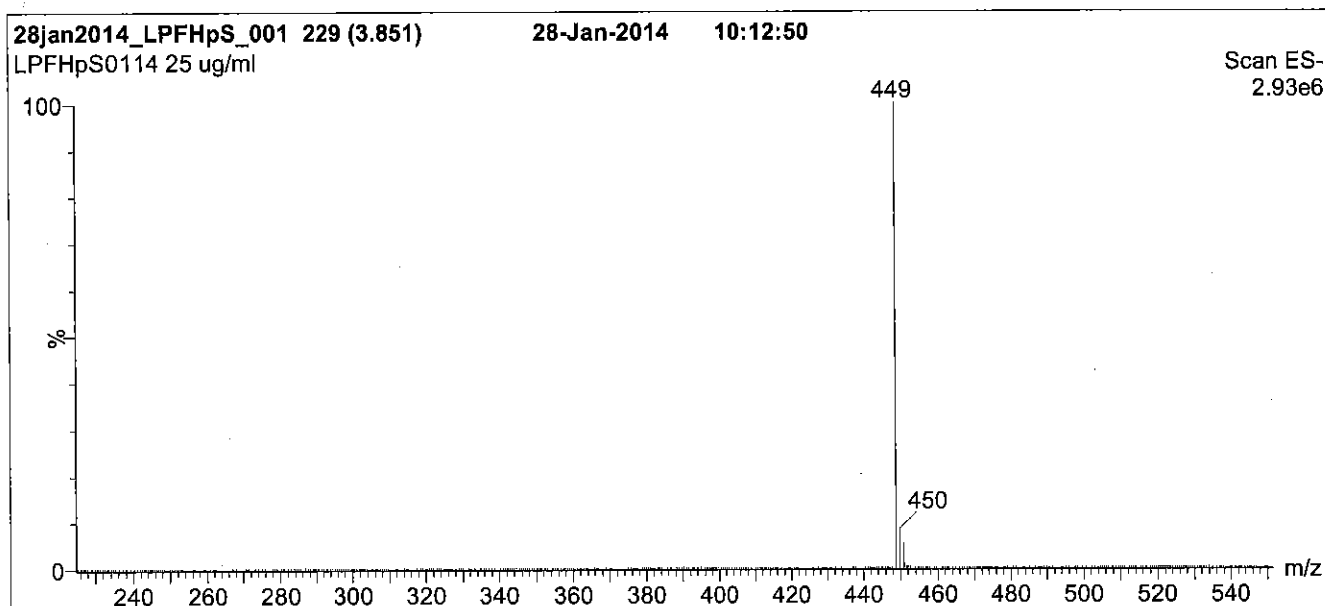
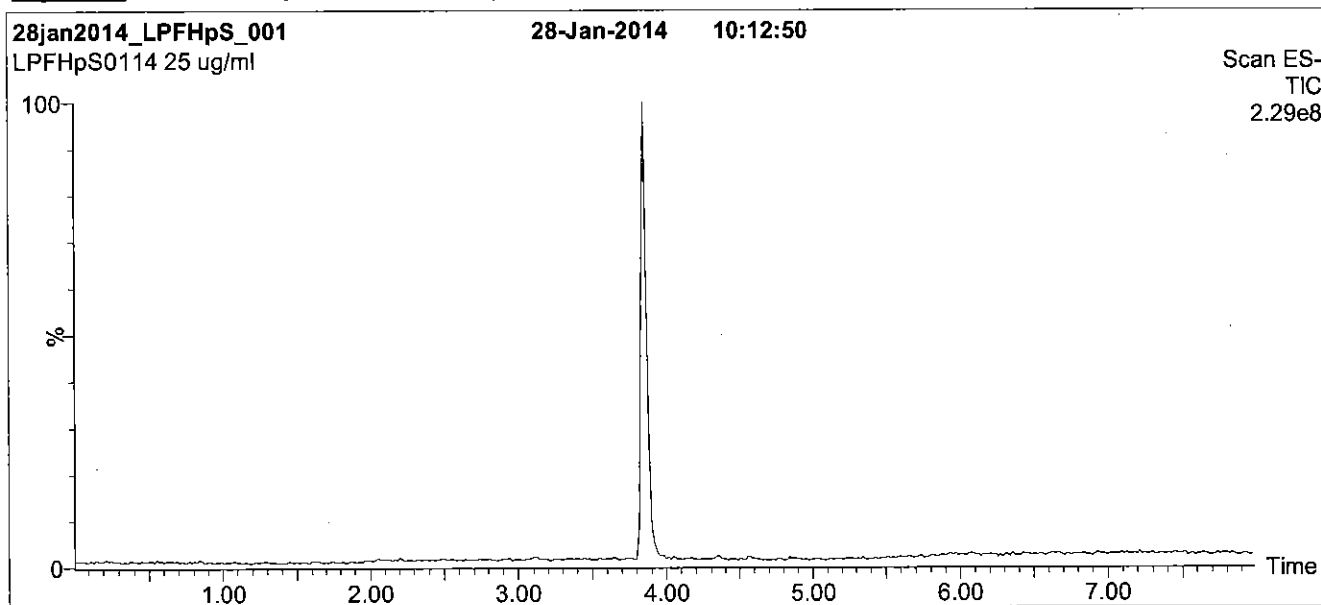
This product was produced using a Quality Management System registered to the latest versions of ISO 9001 by SAI Global, ISO/IEC 17025 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34 by ANSI-ASQ National Accreditation Board (ANAB; AR-1523).



\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*



**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

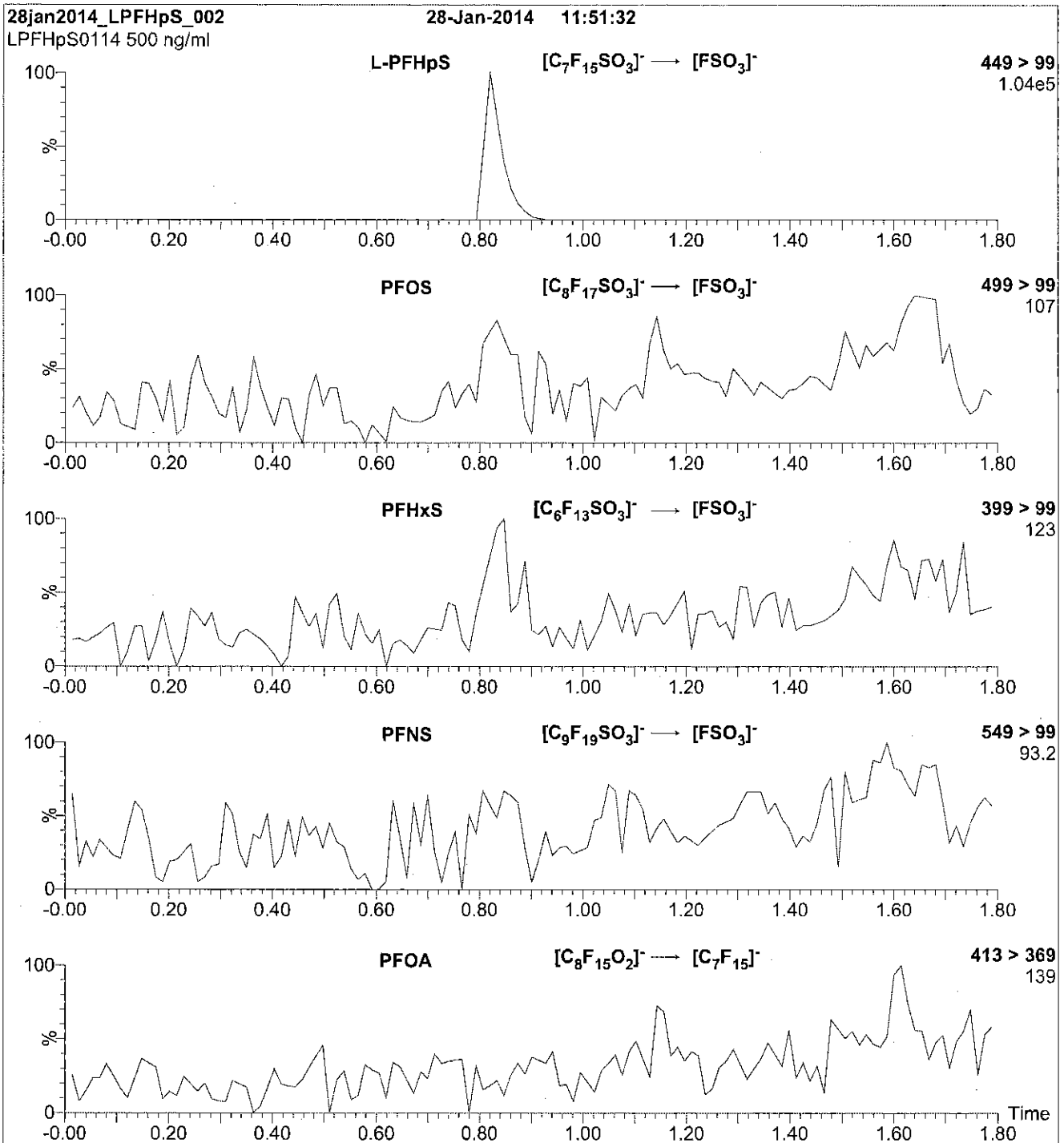
**Flow:** 300  $\mu$ l/min

**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  $\mu$ l (500 ng/ml L-PFHpS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.66e-3  
 Collision Energy (eV) = 35

Reagent

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**LCPFHxA\_00003**

V# 2/11/15 SW



# WELLINGTON LABORATORIES

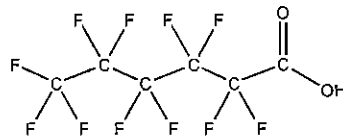
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFHxA  
**COMPOUND:** Perfluoro-n-hexanoic acid

**LOT NUMBER:** PFHxA0514

**STRUCTURE:**

**CAS #:** 307-24-4



**MOLECULAR FORMULA:** C<sub>6</sub>HF<sub>11</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 314.05  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 05/09/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 05/09/2019  
**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.
- 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: 05/22/2014  
(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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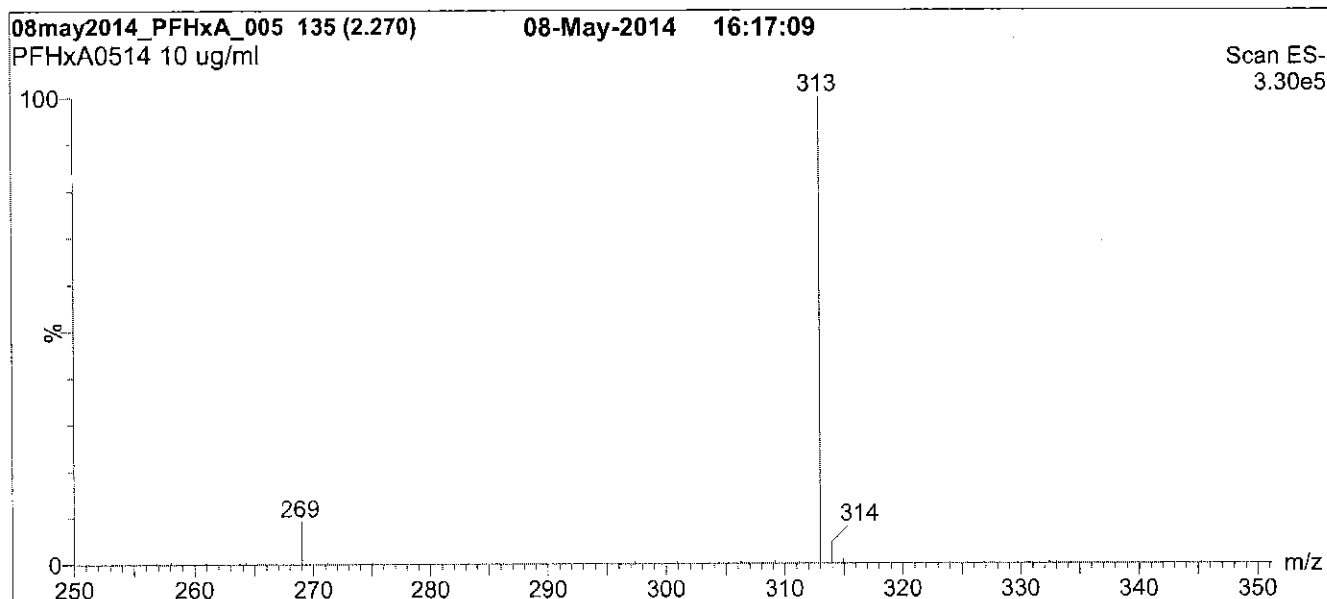
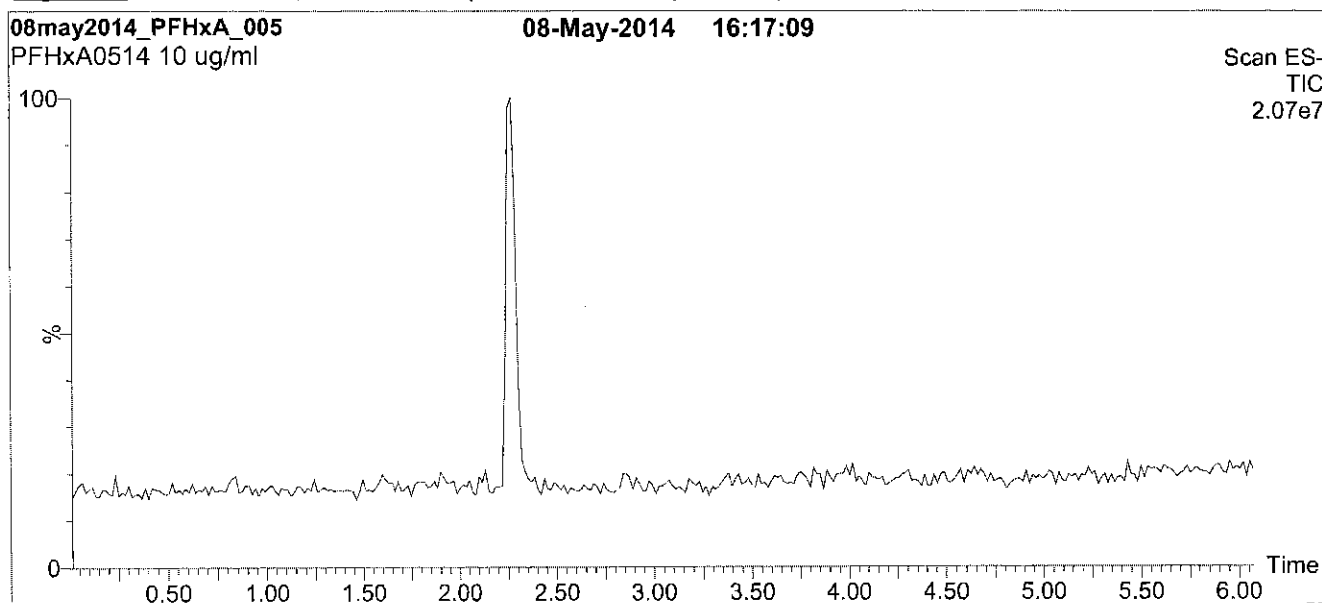
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**Figure 1: PFHxA; LC/MS Data (TIC and Mass Spectrum)**



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acquity UPLC BEH C<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

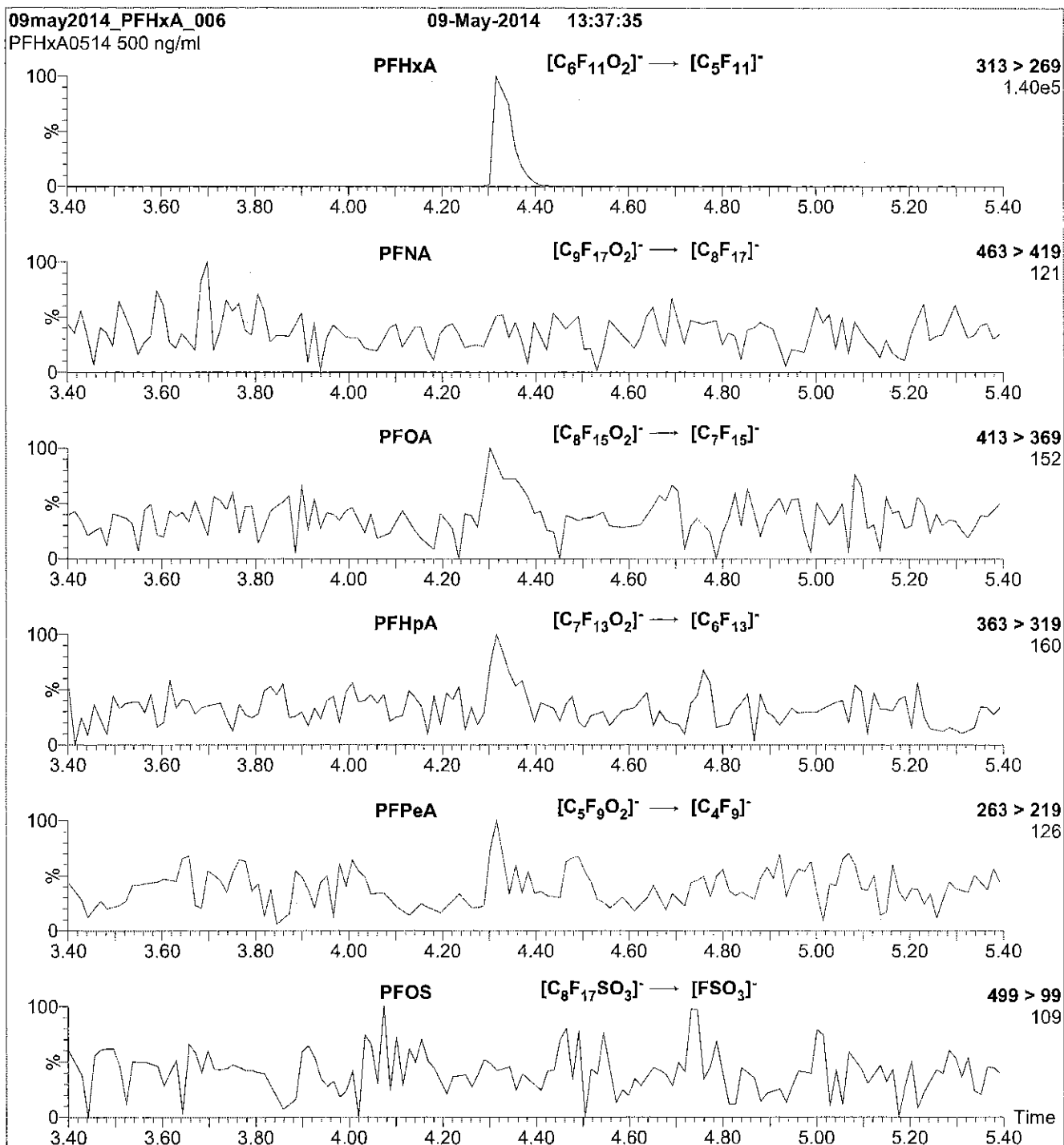
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (250 - 950 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  $\mu$ l (500 ng/ml PFHxA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.24e-3  
Collision Energy (eV) = 10

Reagent

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**LCPFHXS\_00003**





# WELLINGTON LABORATORIES

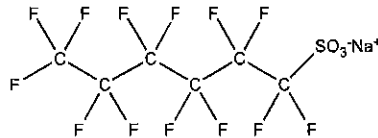
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFHxS  
**COMPOUND:** Sodium perfluoro-1-hexanesulfonate

**LOT NUMBER:** LPFHxS0514

**STRUCTURE:**

**CAS #:** 82382-12-5



**MOLECULAR FORMULA:** C<sub>6</sub>F<sub>13</sub>SO<sub>3</sub>Na  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt)  
 47.3 ± 2.4 µg/ml (PFHxS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 05/09/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 05/09/2019  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**MOLECULAR WEIGHT:** 422.10  
**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.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim  
**Date:** 05/16/2014  
(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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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:

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

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 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 17025:2005 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 for the period of time specified by the 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.

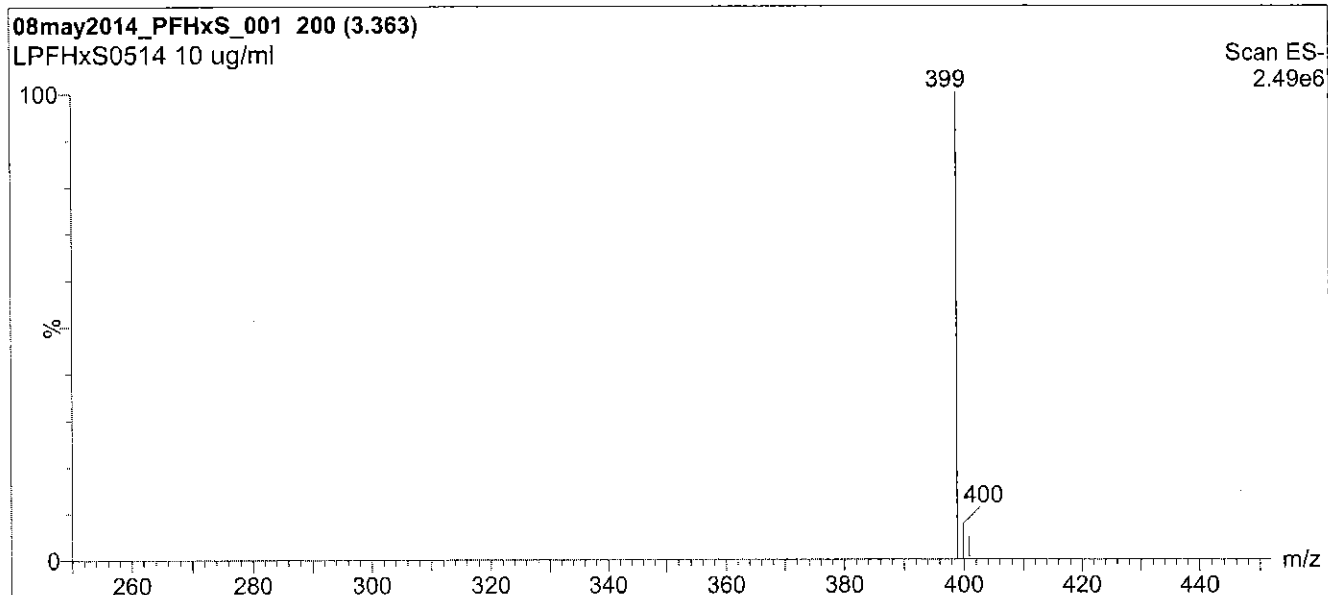
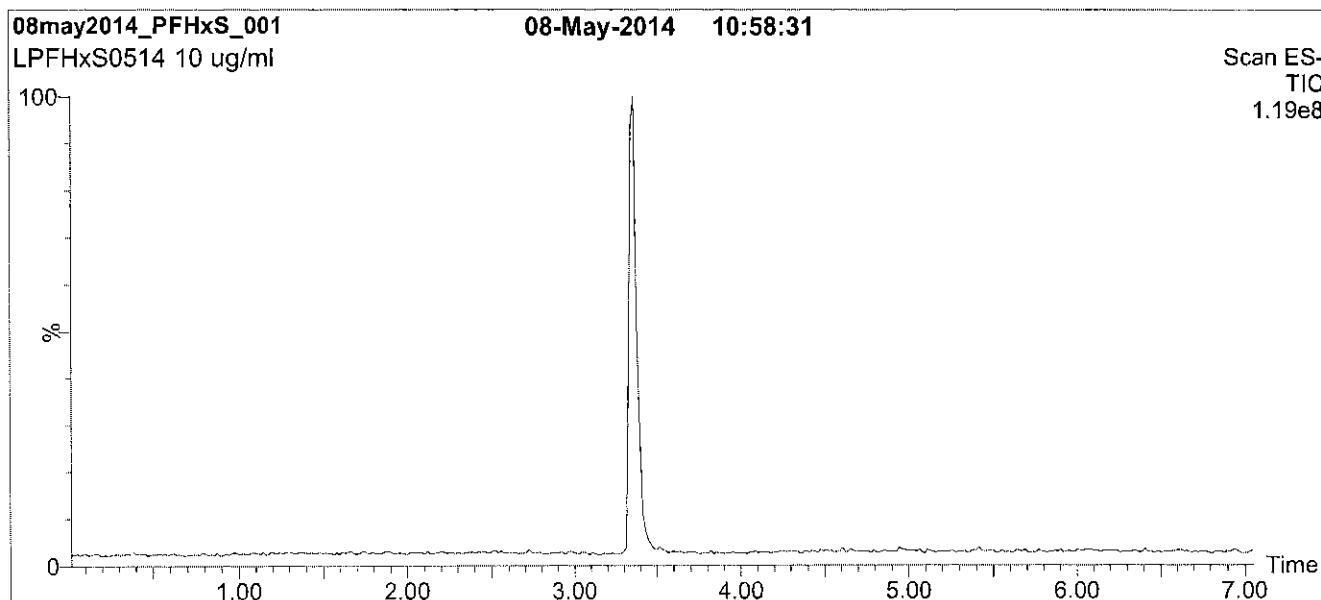
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**Figure 1: L-PFHxS; 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 C<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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

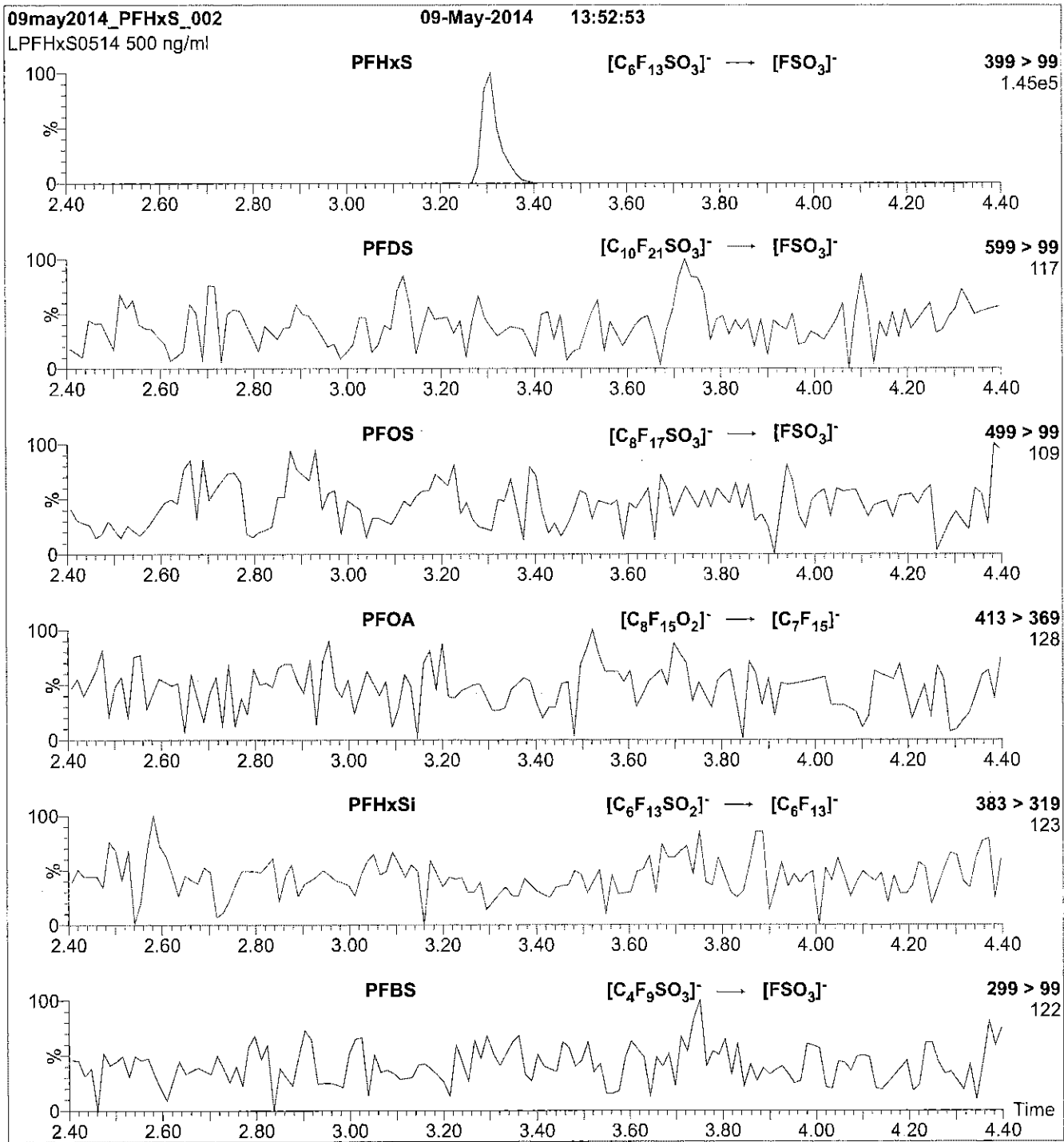
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (250 - 950 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 50.00  
Cone Gas Flow (l/hr) = 60  
Desolvation Gas Flow (l/hr) = 750

**Figure 2: L-PFHxS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
 10  $\mu$ l (500 ng/ml L-PFHxS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

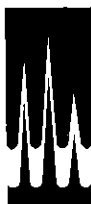
Collision Gas (mbar) = 3.17e-3  
 Collision Energy (eV) = 30

Reagent

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**LCPFNA\_00004**

r: 3/27/15 ✓  
s:



# WELLINGTON LABORATORIES

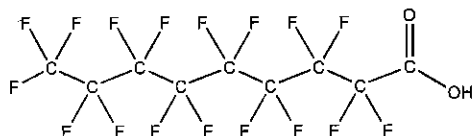
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFNA  
**COMPOUND:** Perfluoro-n-nonanoic acid

**LOT NUMBER:** PFNA0514

**STRUCTURE:**

**CAS #:** 375-95-1



**MOLECULAR FORMULA:** C<sub>9</sub>H<sub>17</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 464.08  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 05/09/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 05/09/2019  
**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.
- 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:** 05/22/2014  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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:

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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.

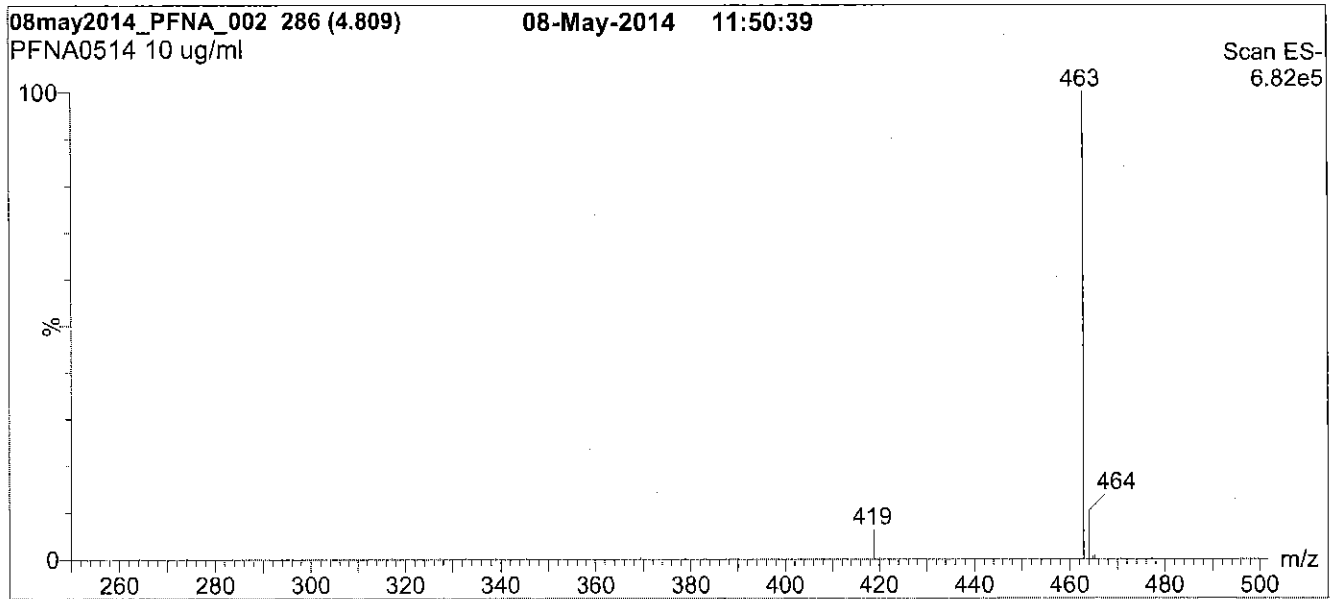
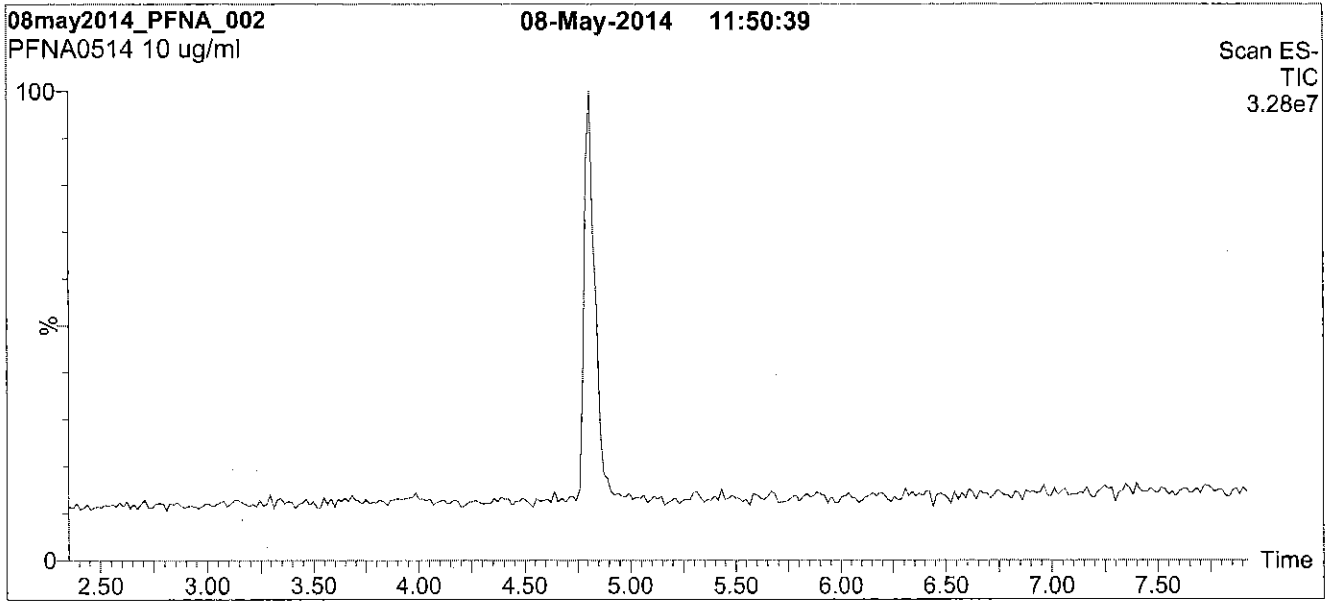
**QUALITY MANAGEMENT:**

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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 *micro* API MS

**Chromatographic Conditions**

Column: Acquity UPLC BEH C<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 50% (80:20 MeOH:ACN) / 50% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>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  $\mu$ l/min

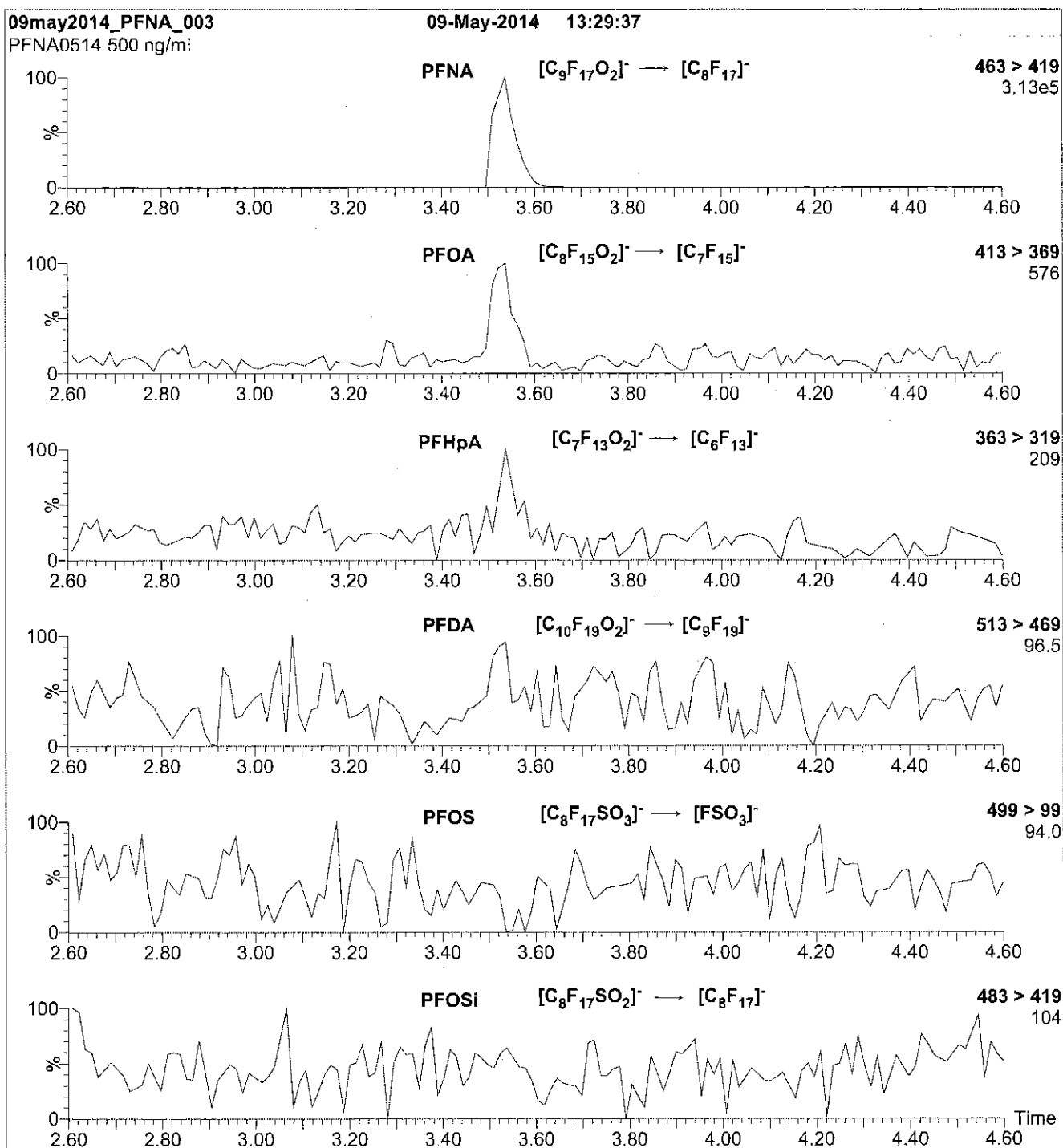
**MS Parameters**

Experiment: Full Scan (250 - 950 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  $\mu$ l (500 ng/ml PFNA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.17e-3  
Collision Energy (eV) = 11

Reagent

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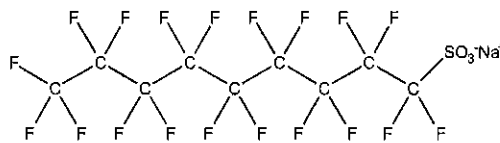
**LCPFNS\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFNS **LOT NUMBER:** LPFNS0712  
**COMPOUND:** Sodium perfluoro-1-nonanesulfonate  
**STRUCTURE:** **CAS #:** 98789-57-2



**MOLECULAR FORMULA:** C<sub>9</sub>F<sub>19</sub>SO<sub>3</sub>Na **MOLECULAR WEIGHT:** 572.12  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt) **SOLVENT(S):** Methanol  
 48.0 ± 2.4 µg/ml (PFNS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 07/04/2012  
**EXPIRY DATE:** (mm/dd/yyyy) 07/04/2017  
**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: 01/15/2013

(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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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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:

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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:**

Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the 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.

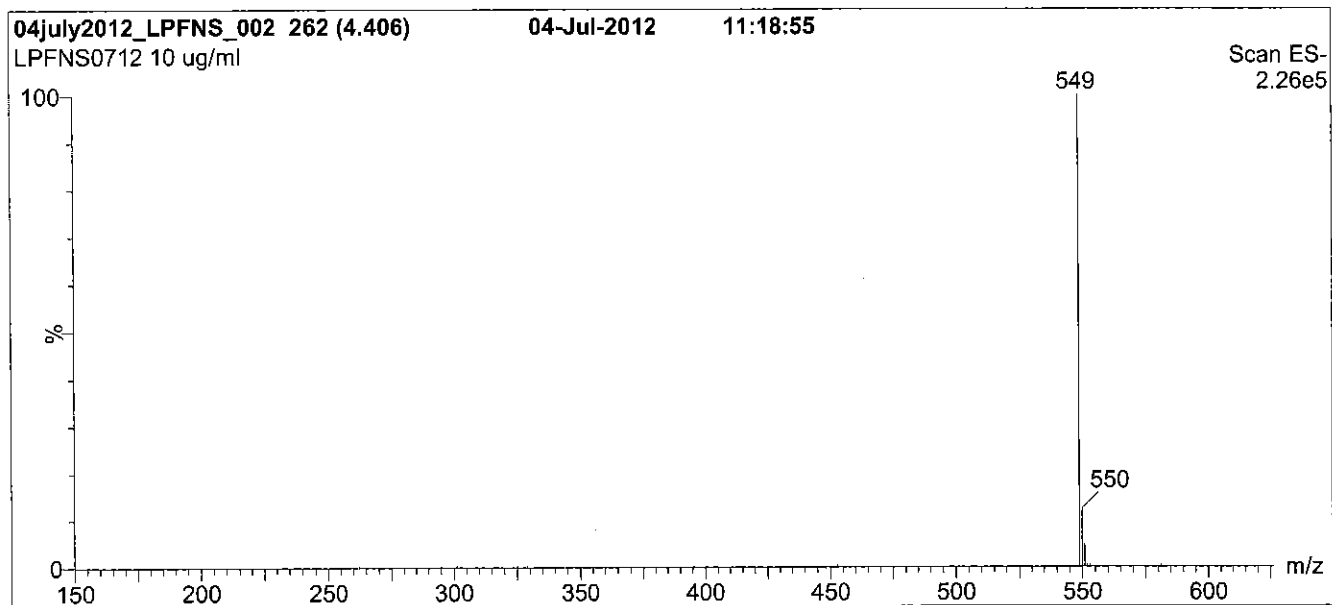
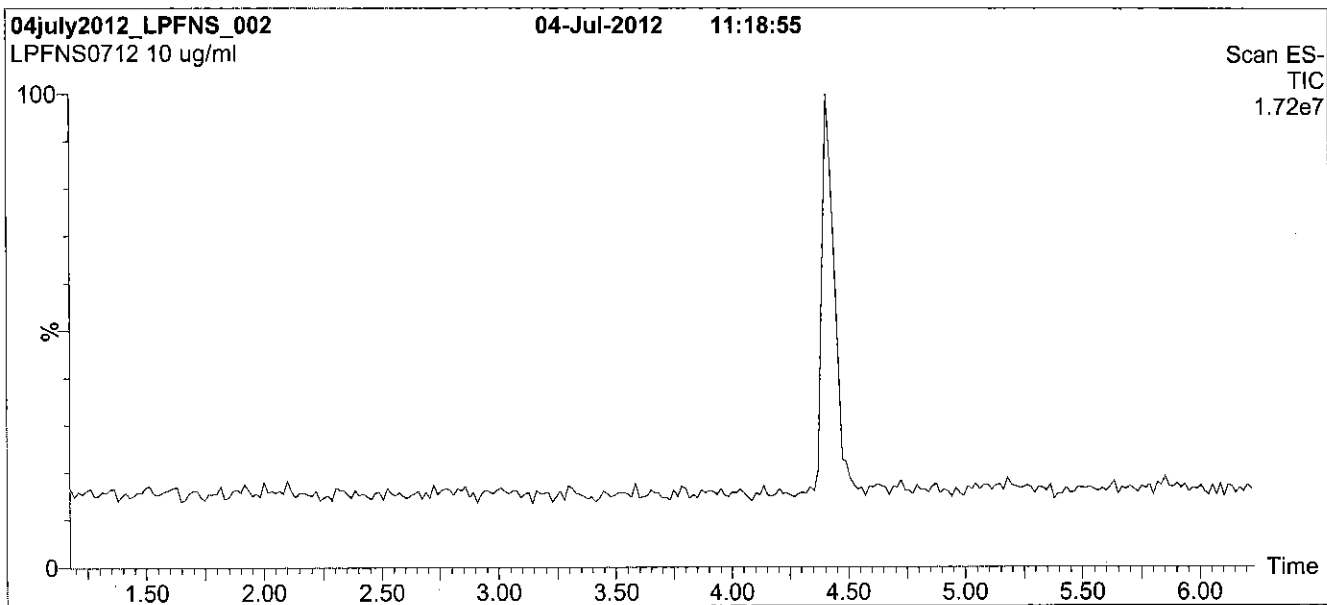
### **QUALITY MANAGEMENT:**

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**Figure 1: L-PFNS; 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<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

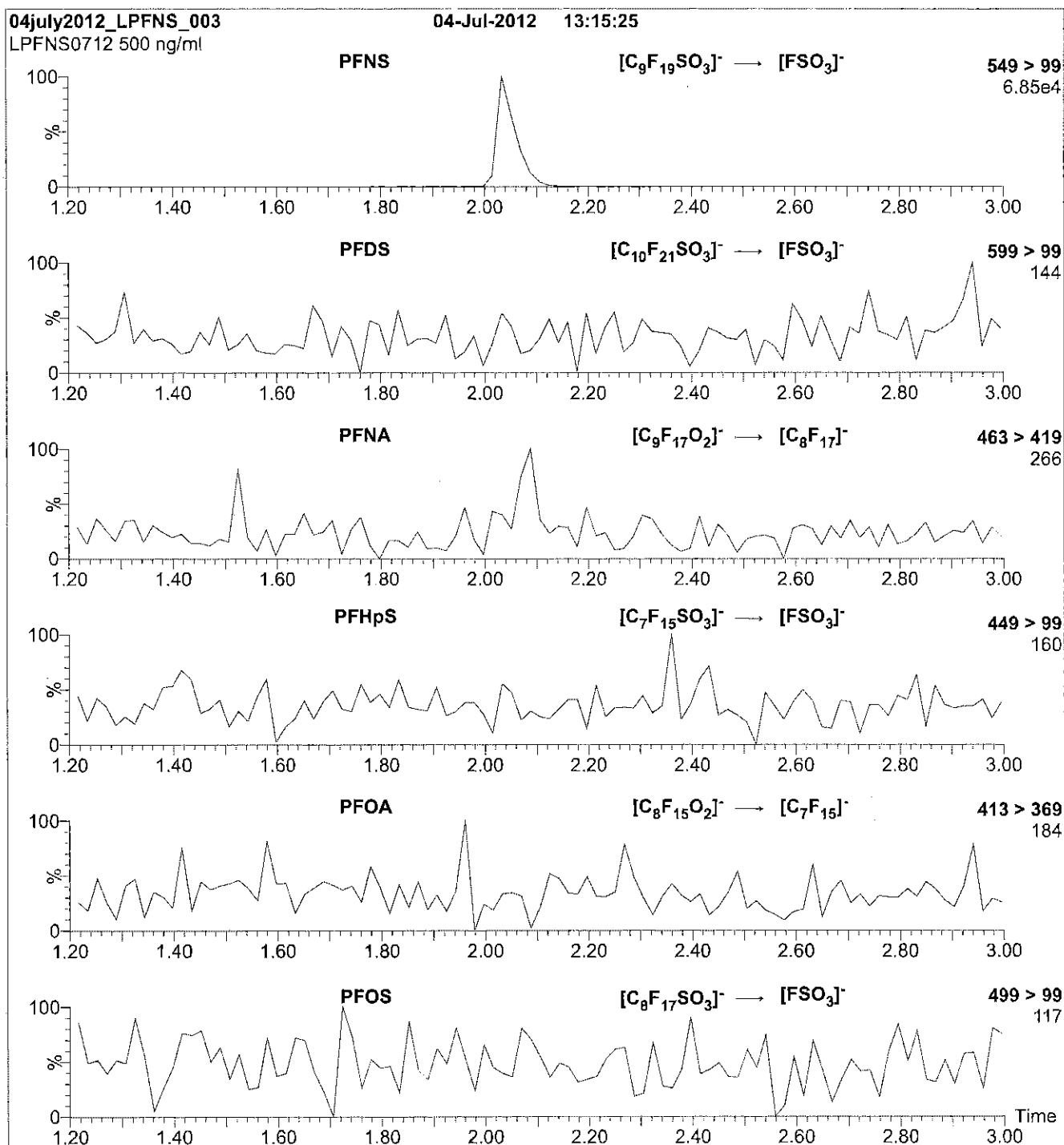
**Flow:** 300  $\mu$ l/min

**MS Parameters**

**Experiment:** Full Scan (150 - 850 amu)

**Source:** Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 65.00  
Cone Gas Flow (l/hr) = 50  
Desolvation Gas Flow (l/hr) = 750

**Figure 2: L-PFNS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
 10  $\mu$ l (500 ng/ml L-PFNS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.54e-3  
 Collision Energy (eV) = 45

Reagent

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**LCPFOA\_00004**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Rec 7/15/14

**PRODUCT CODE:**

PFOA

**LOT NUMBER:**

PFOA1013

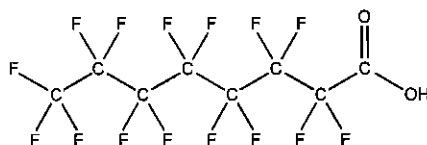
**COMPOUND:**

Perfluoro-n-octanoic acid

**STRUCTURE:**

**CAS #:**

335-67-1



**MOLECULAR FORMULA:**

$C_8H_{15}O_2$

**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)

10/11/2013

**EXPIRY DATE:** (mm/dd/yyyy)

10/11/2018

**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.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

Certified By:

B.G. Chittim

Date: 10/18/2013  
(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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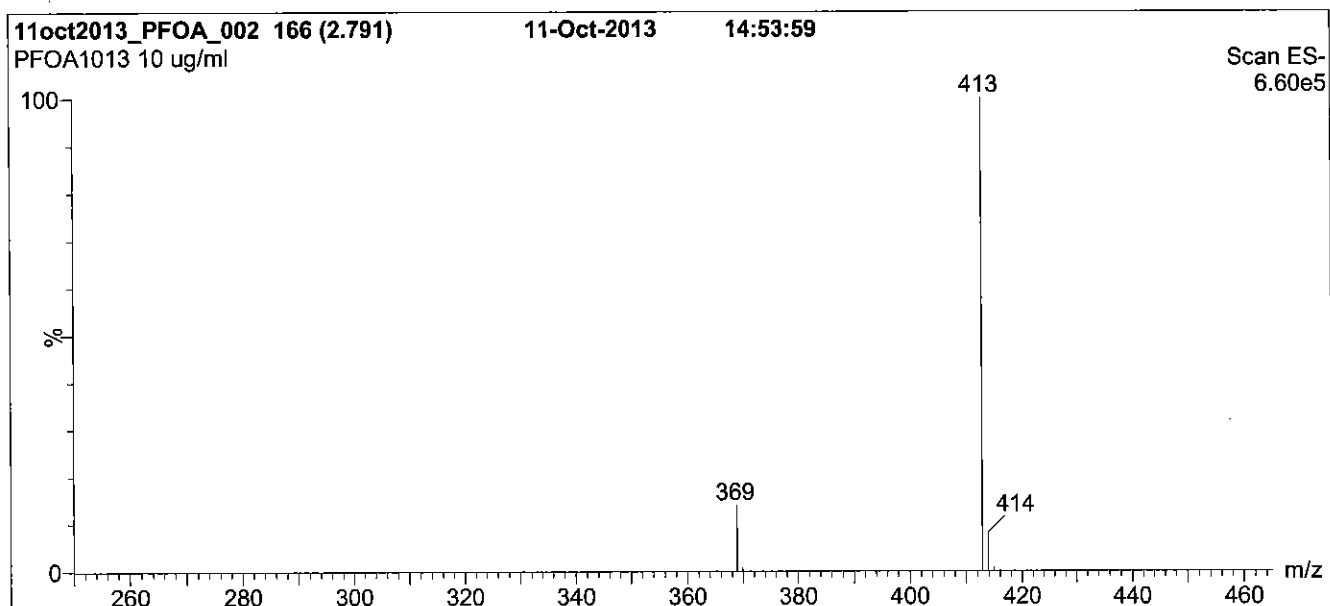
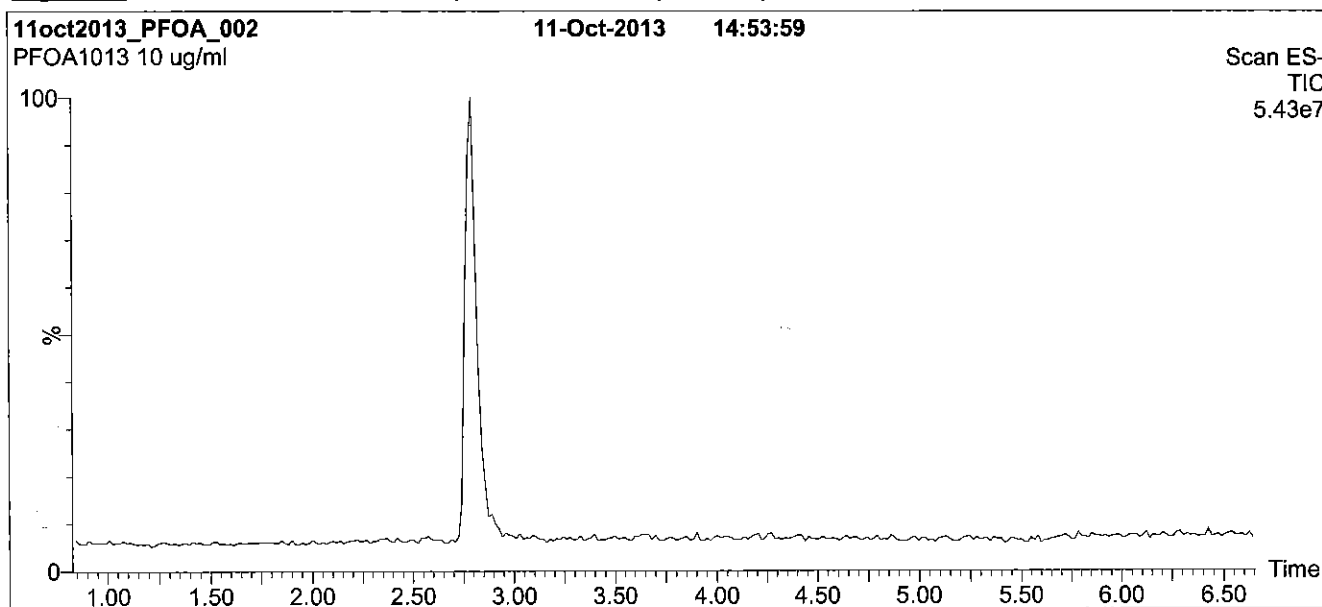
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**Figure 1: PFOA; LC/MS Data (TIC and Mass Spectrum)**



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

**Column:** Acquity UPLC BEH Shield RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

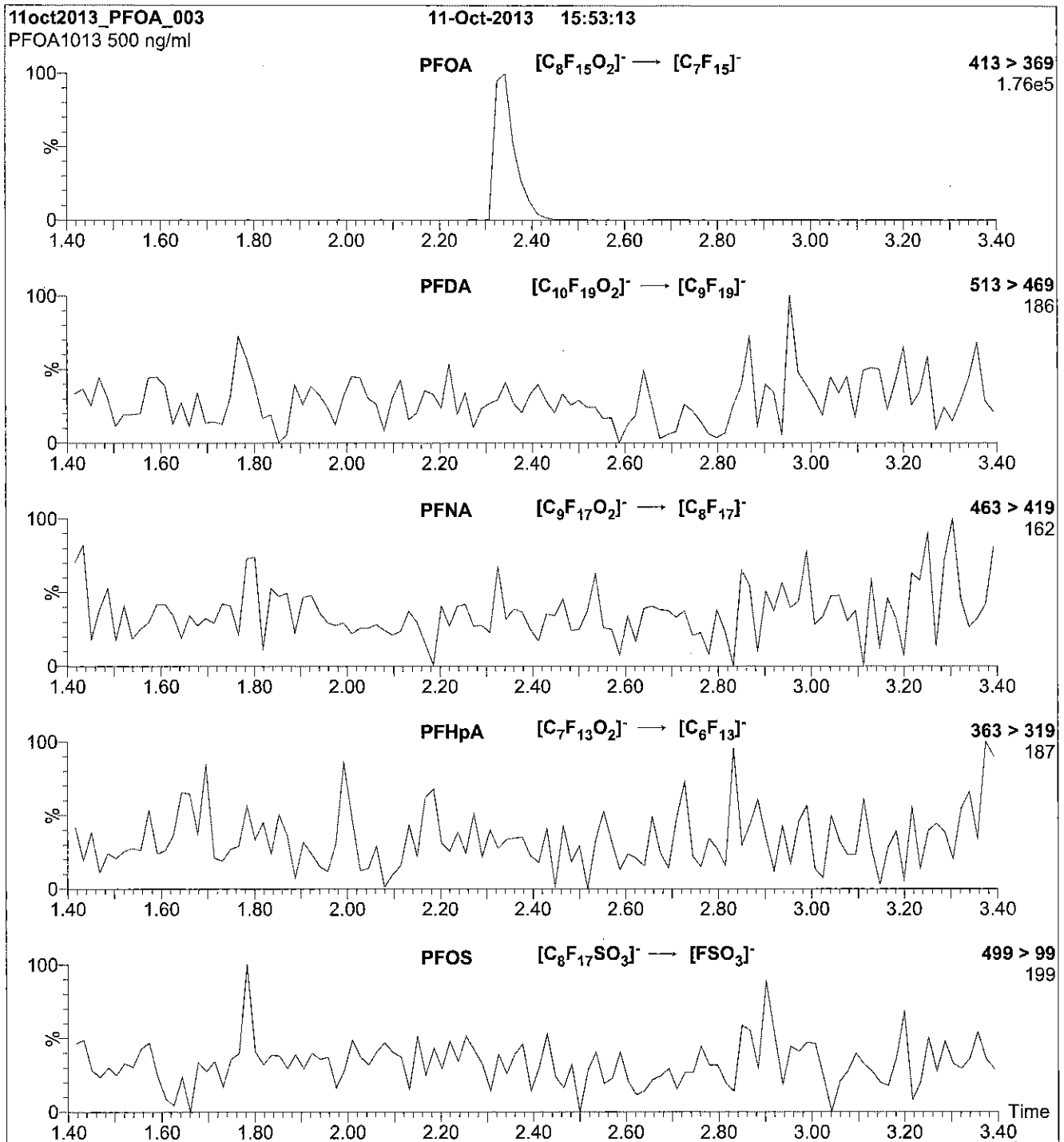
**Flow:** 300  $\mu$ l/min

**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) = 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  $\mu$ l (500 ng/ml PFOA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

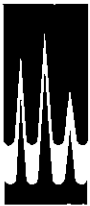
Collision Gas (mbar) = 3.28e-3  
Collision Energy (eV) = 11

Reagent

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**LCPFODA\_00004**

17 2/15 2

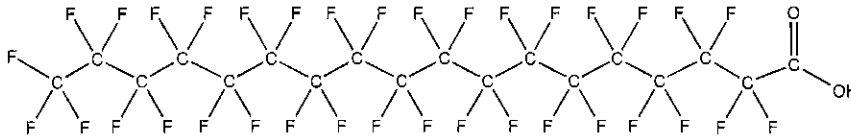


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFODA **LOT NUMBER:** PFODA0807  
**COMPOUND:** Perfluoro-n-octadecanoic acid

**STRUCTURE:** **CAS #:** 16517-11-6



**MOLECULAR FORMULA:** C<sub>18</sub>H<sub>35</sub>F<sub>35</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 914.15  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (4%)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 04/25/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 04/25/2017  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

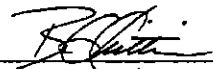
**DOCUMENTATION/ DATA ATTACHED:**

- 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:** 04/28/2014  
 (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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

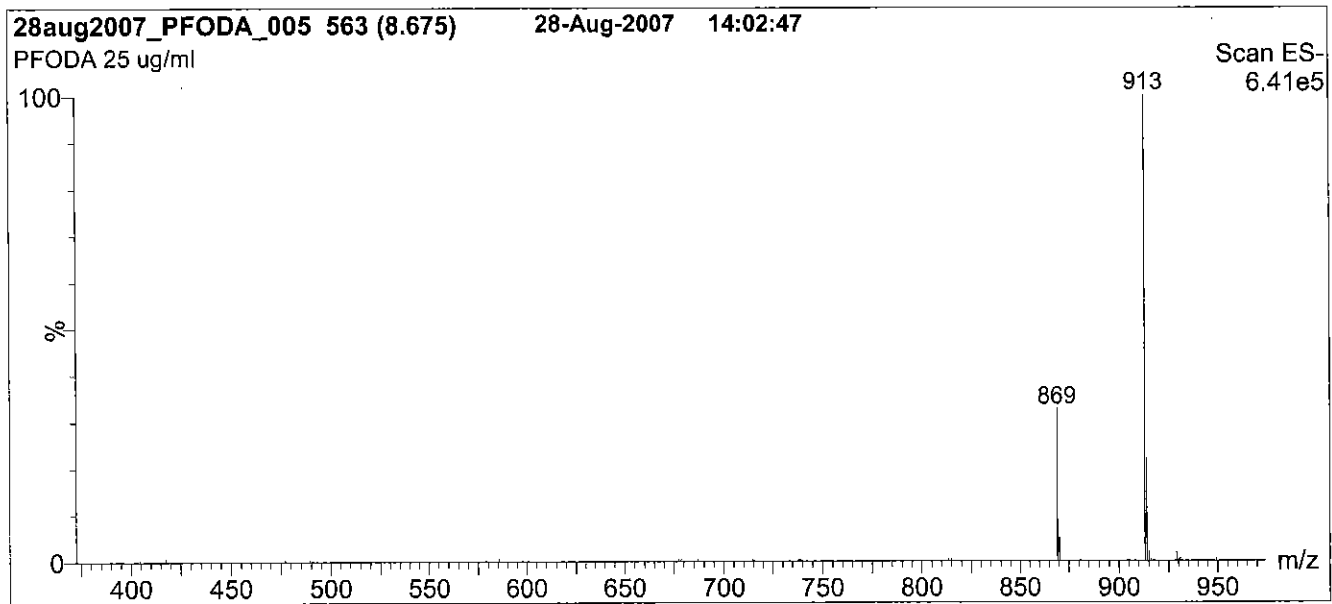
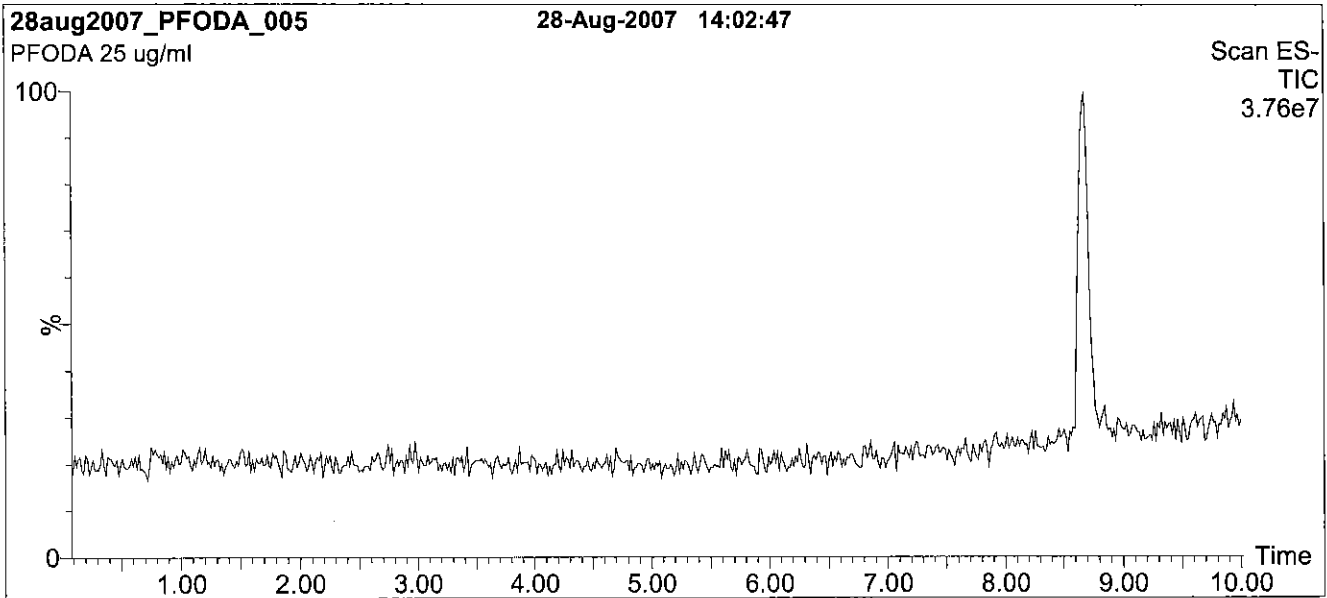
**QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 75% (80:20 MeOH:ACN) / 25% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)  
 Hold 5 min. Ramp to 100% organic over 6 min.  
 Hold 3 min before returning to initial conditions.  
 Time: 16 min

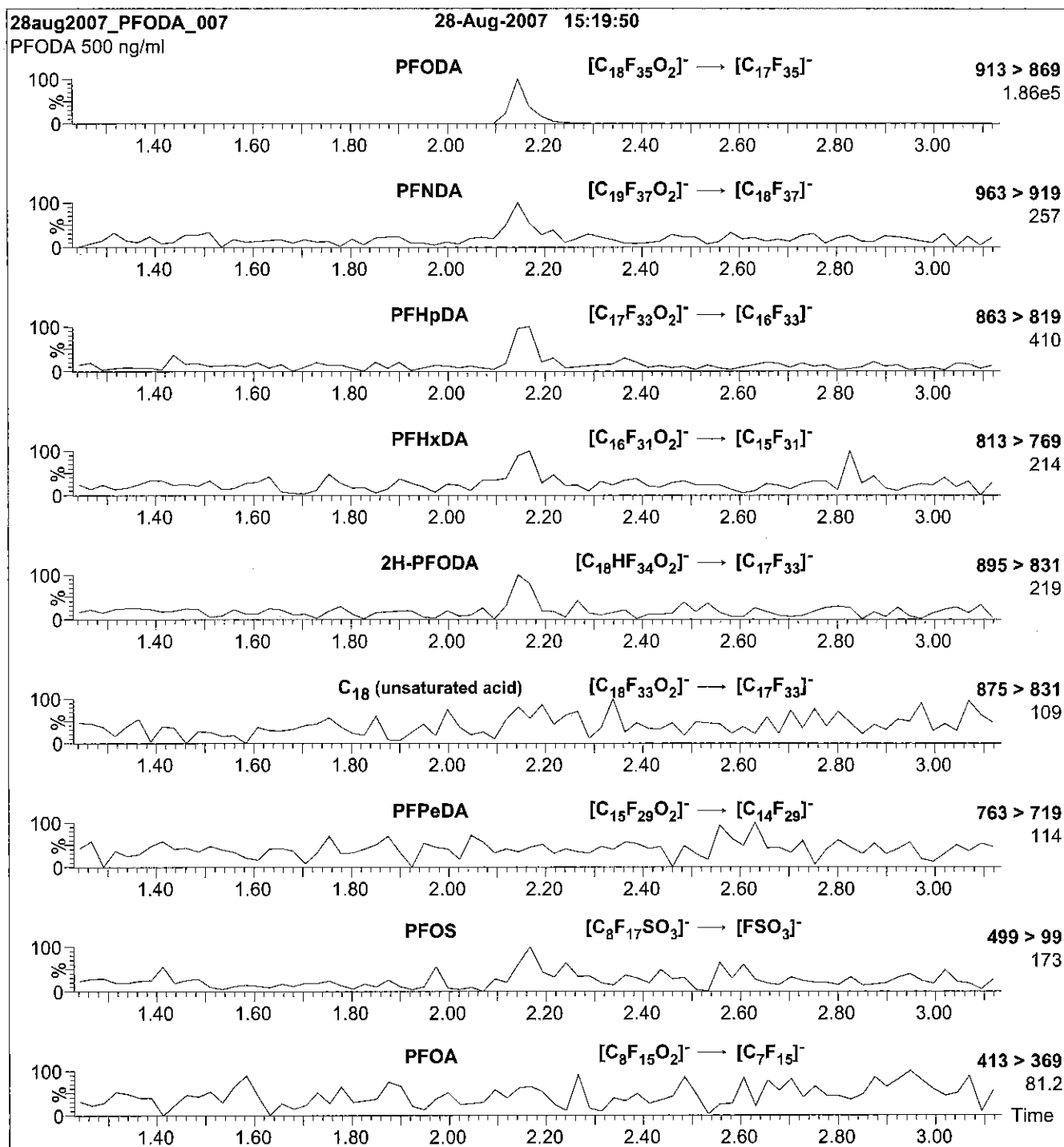
**Flow:** 300  $\mu$ l/min

**MS Parameters**

**Experiment:** Full Scan (225 - 1100 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) = 650

**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 75% (80:20 MeOH:ACN) / 25% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300 µl/min

**MS Parameters**

Collision Gas (mbar) = 3.58e-3  
Collision Energy (eV) = 15

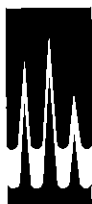


Reagent

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**LCPFOS\_00004**

3/17/15 SV



# WELLINGTON LABORATORIES

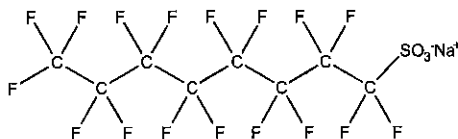
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFOS  
**COMPOUND:** Sodium perfluoro-1-octanesulfonate

**LOT NUMBER:** LPFOS0614

**STRUCTURE:**

**CAS #:** 4021-47-0



**MOLECULAR FORMULA:** C<sub>8</sub>F<sub>17</sub>SO<sub>3</sub>Na  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt)  
 47.8 ± 2.4 µg/ml (PFOS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 06/20/2014  
**EXPIRY DATE:** (mm/dd/yyyy) 06/20/2019  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**MOLECULAR WEIGHT:** 522.11  
**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:** 10/27/2014  
 (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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

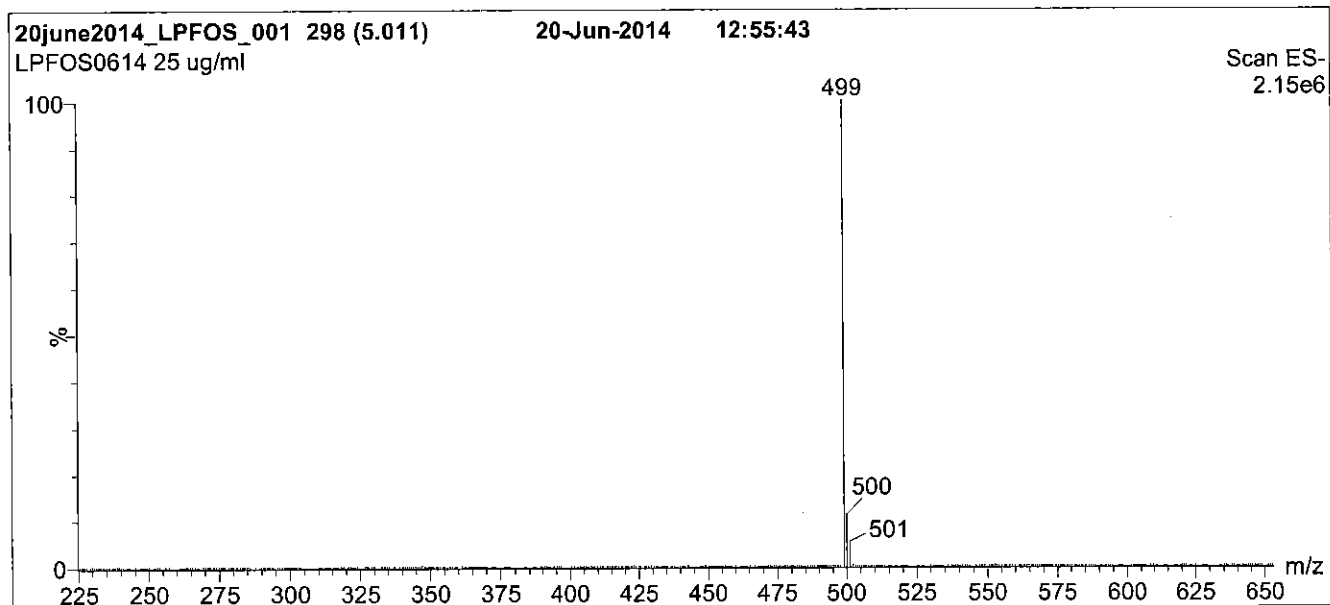
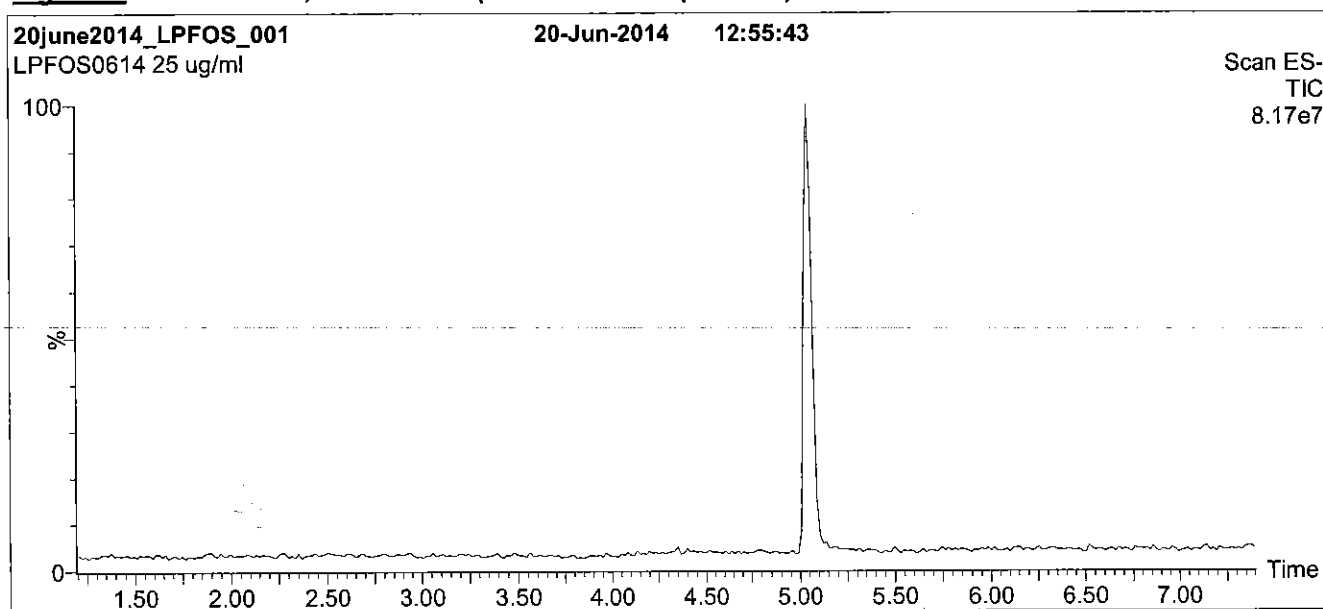
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**Figure 1: L-PFOS; 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<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 45% (80:20 MeOH:ACN) / 55% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

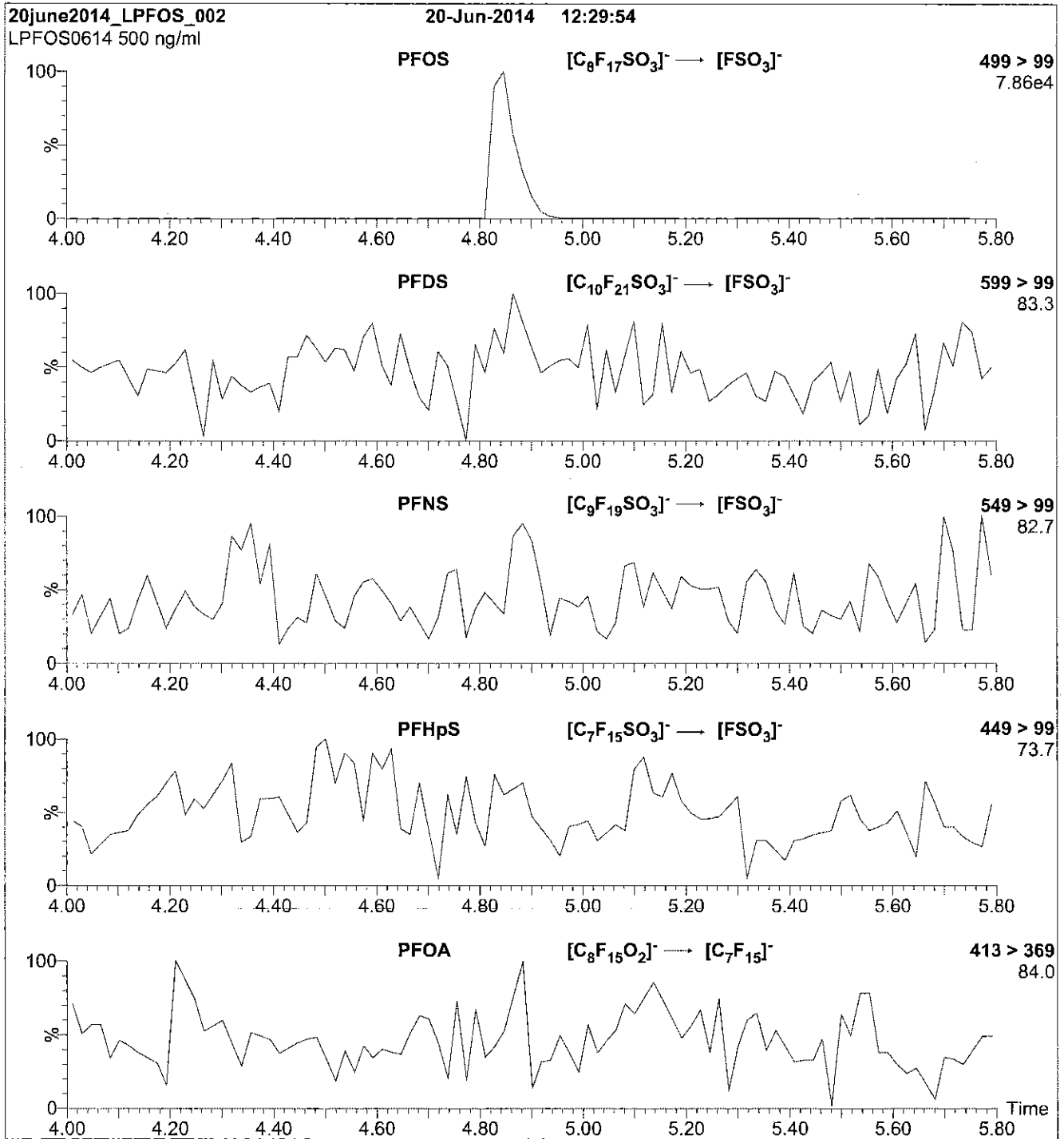
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (225 - 950 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: L-PFOS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
 10  $\mu$ l (500 ng/ml L-PFOS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.43e-3  
 Collision Energy (eV) = 40

Reagent

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**LCPFOSA\_00005**

ri 2/11/15 BV

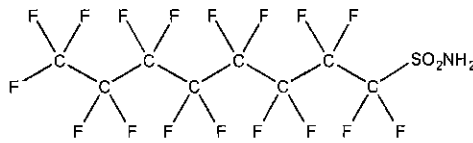


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** FOSA-I **LOT NUMBER:** FOSA0714I  
**COMPOUND:** Perfluoro-1-octanesulfonamide

**STRUCTURE:** **CAS #:** 754-91-6



**MOLECULAR FORMULA:** C<sub>8</sub>H<sub>2</sub>F<sub>17</sub>NO<sub>2</sub>S **MOLECULAR WEIGHT:** 499.14  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Isopropanol  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 07/31/2014  
**EXPIRY DATE:** (mm/dd/yyyy) Stability studies ongoing  
**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:** 08/05/2014  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 our products.

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**EXPIRY DATE / PERIOD OF VALIDITY:**

Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the 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:**

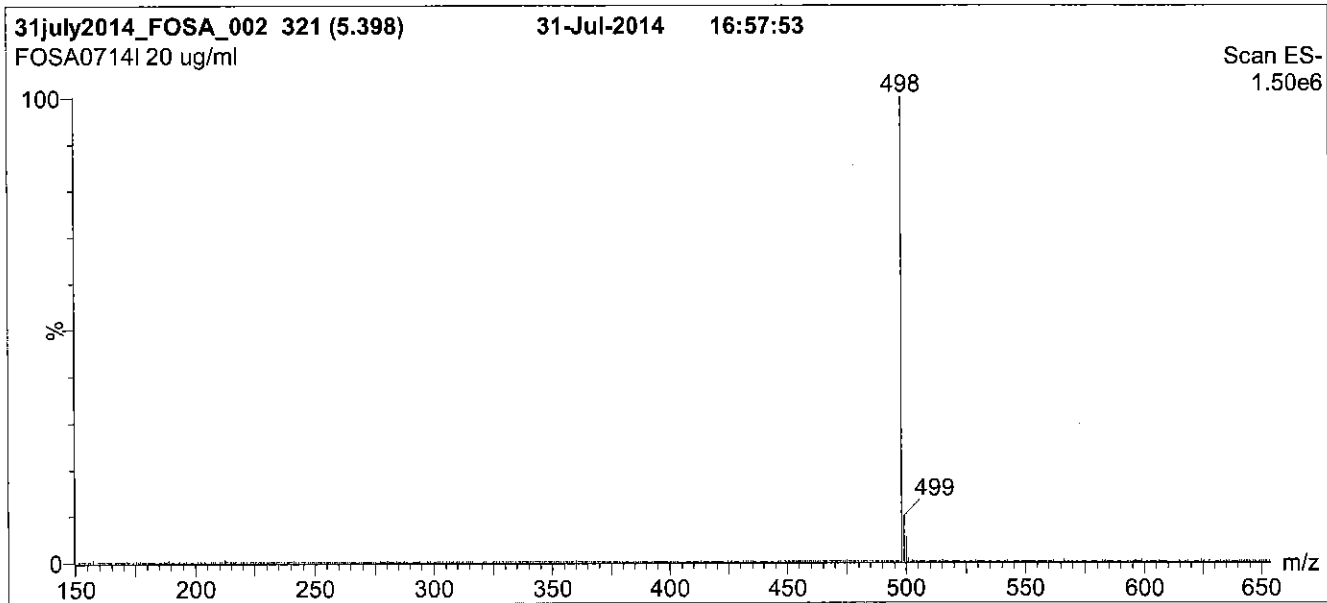
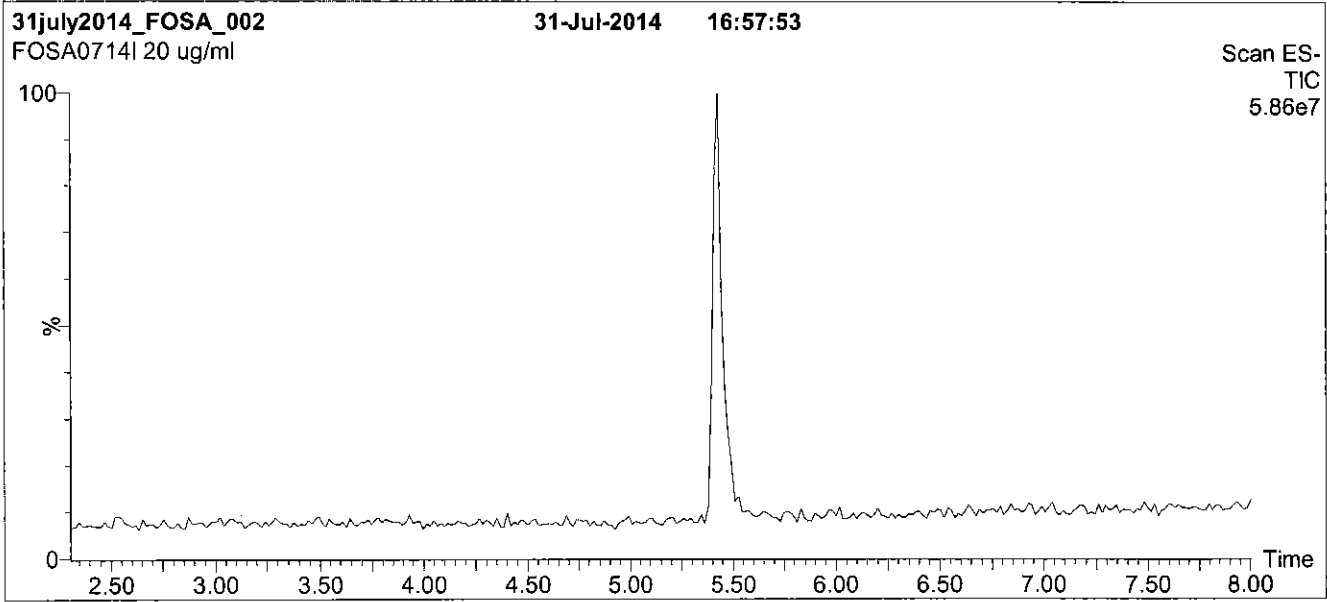
This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number AR-1523).



\*\*For additional information or assistance concerning this or any other products from Wellington Laboratories Inc., please visit our website at [www.well-labs.com](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto: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 C<sub>18</sub>  
 1.7 μm, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>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

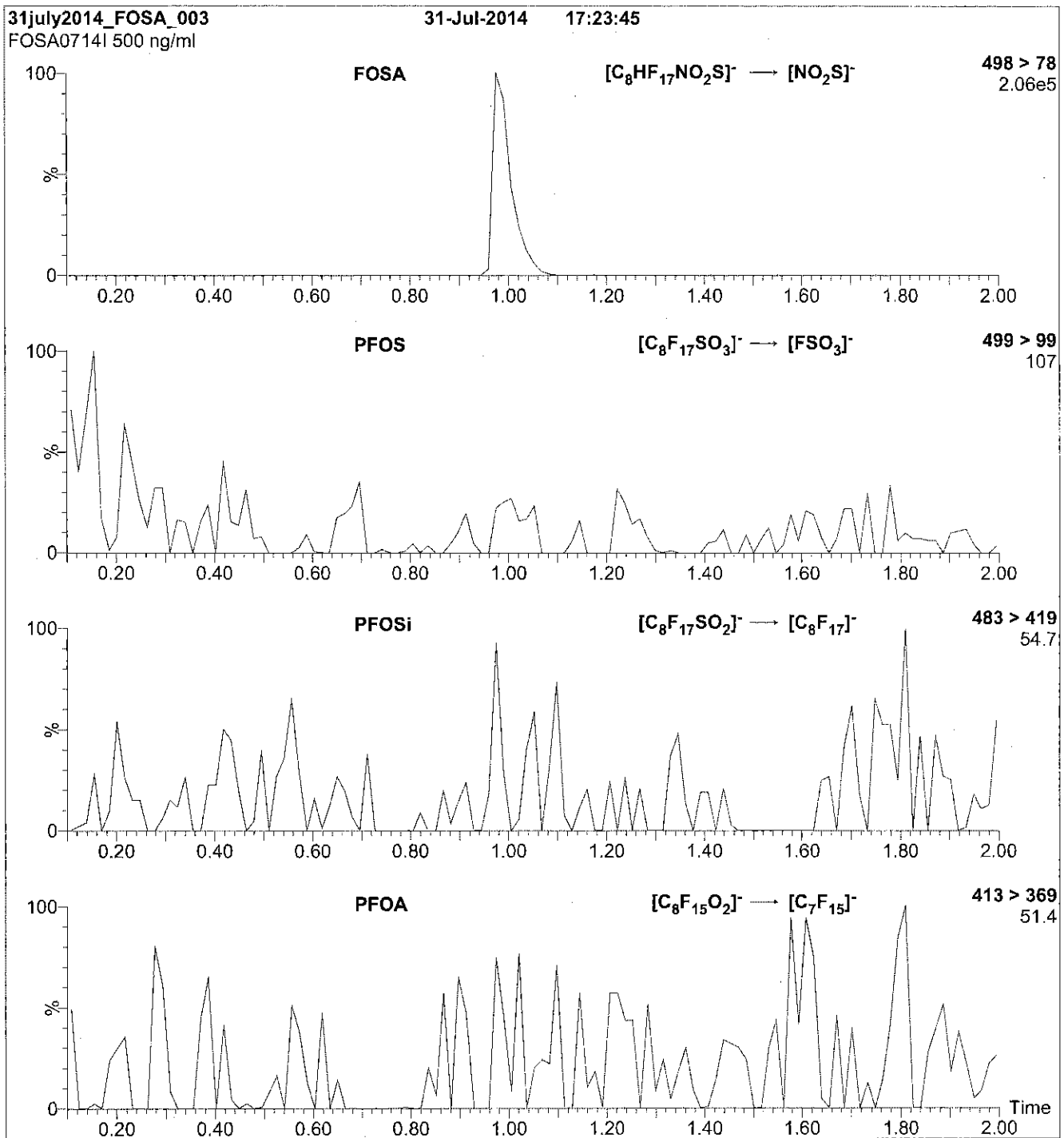
**Flow:** 300 μl/min

**MS Parameters**

**Experiment:** Full Scan (150 - 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: FOSA-I; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct loop injection  
10  $\mu$ l (500 ng/ml FOSA-I)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.58e-3  
Collision Energy (eV) = 30

Reagent

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**LCFPeA\_00003**

Rec 7/15/14



# WELLINGTON LABORATORIES

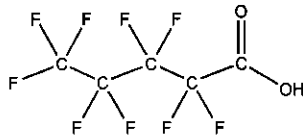
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFPeA  
**COMPOUND:** Perfluoro-n-pentanoic acid

**LOT NUMBER:** PFPeA0113

**STRUCTURE:**

**CAS #:** 2706-90-3



**MOLECULAR FORMULA:** C<sub>5</sub>HF<sub>9</sub>O<sub>2</sub>  
**CONCENTRATION:** 50 ± 2.5 µg/ml

**MOLECULAR WEIGHT:** 264.05  
**SOLVENT(S):** Methanol  
Water (<1%)

**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 01/03/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 01/03/2018  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

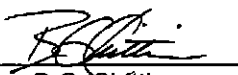
**DOCUMENTATION/ DATA ATTACHED:**

- 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 ~ 0.3% of Perfluoro-n-heptanoic acid (PFHpA) and ~ 0.2% of C<sub>5</sub>H<sub>2</sub>F<sub>8</sub>O<sub>2</sub> (hydrido - derivative) as measured by <sup>19</sup>F NMR.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
B.G. Chittim  
**Date:** 01/14/2013  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

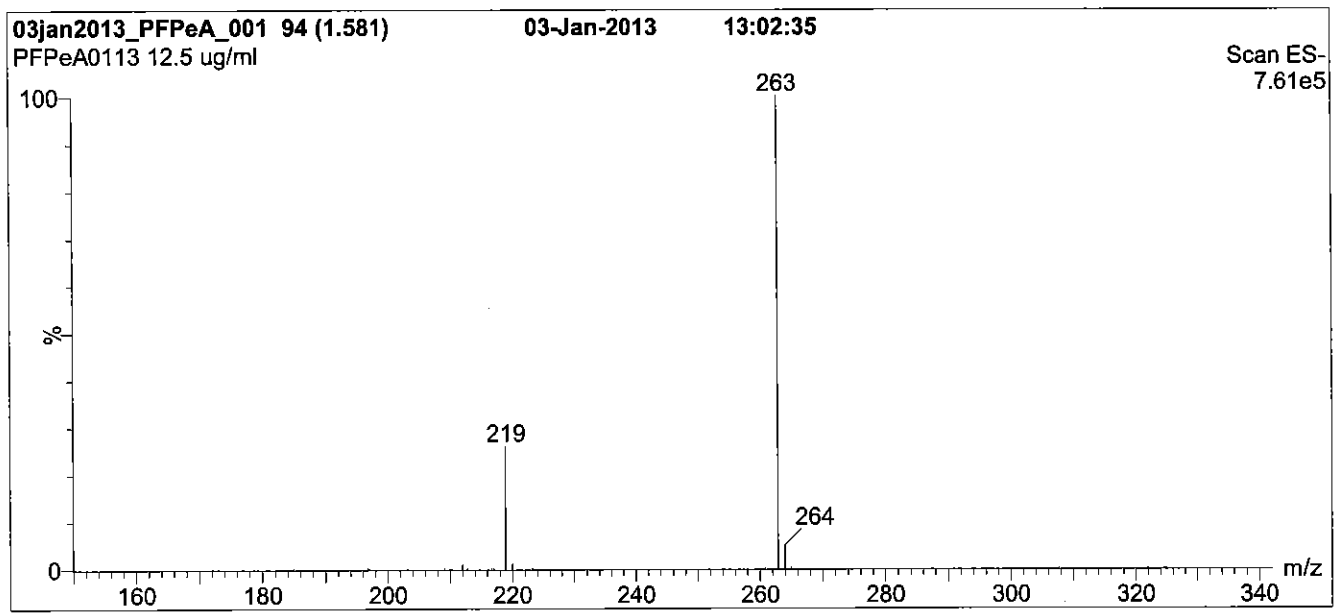
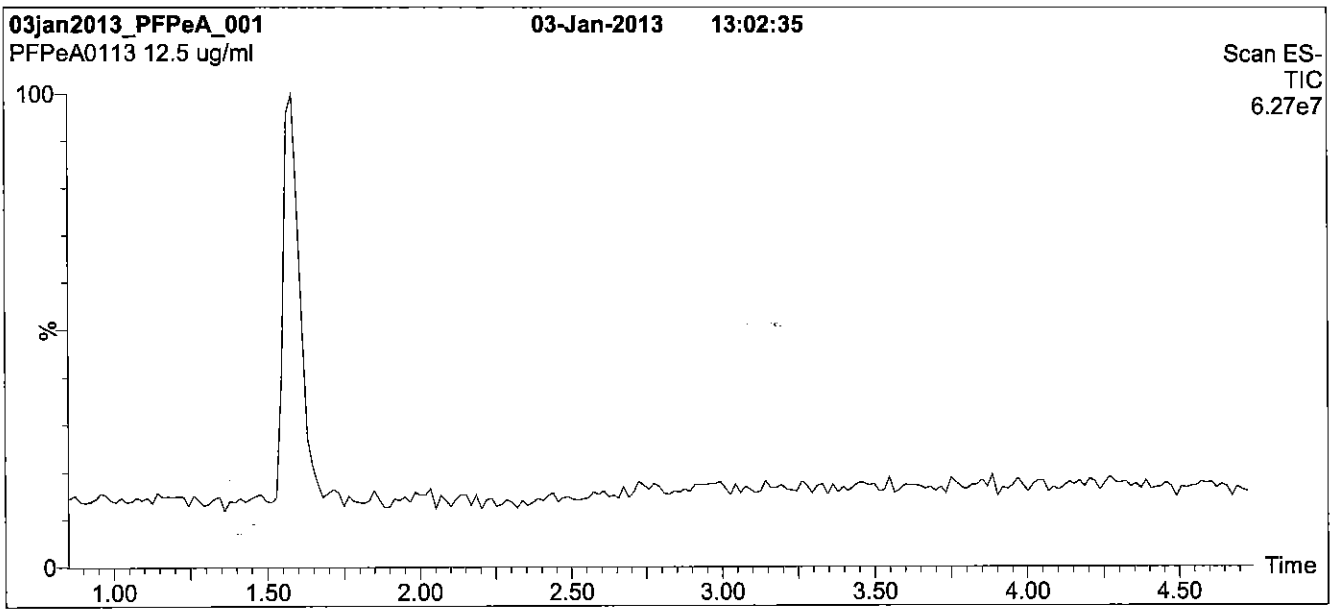
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto: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 RP<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
 Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>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

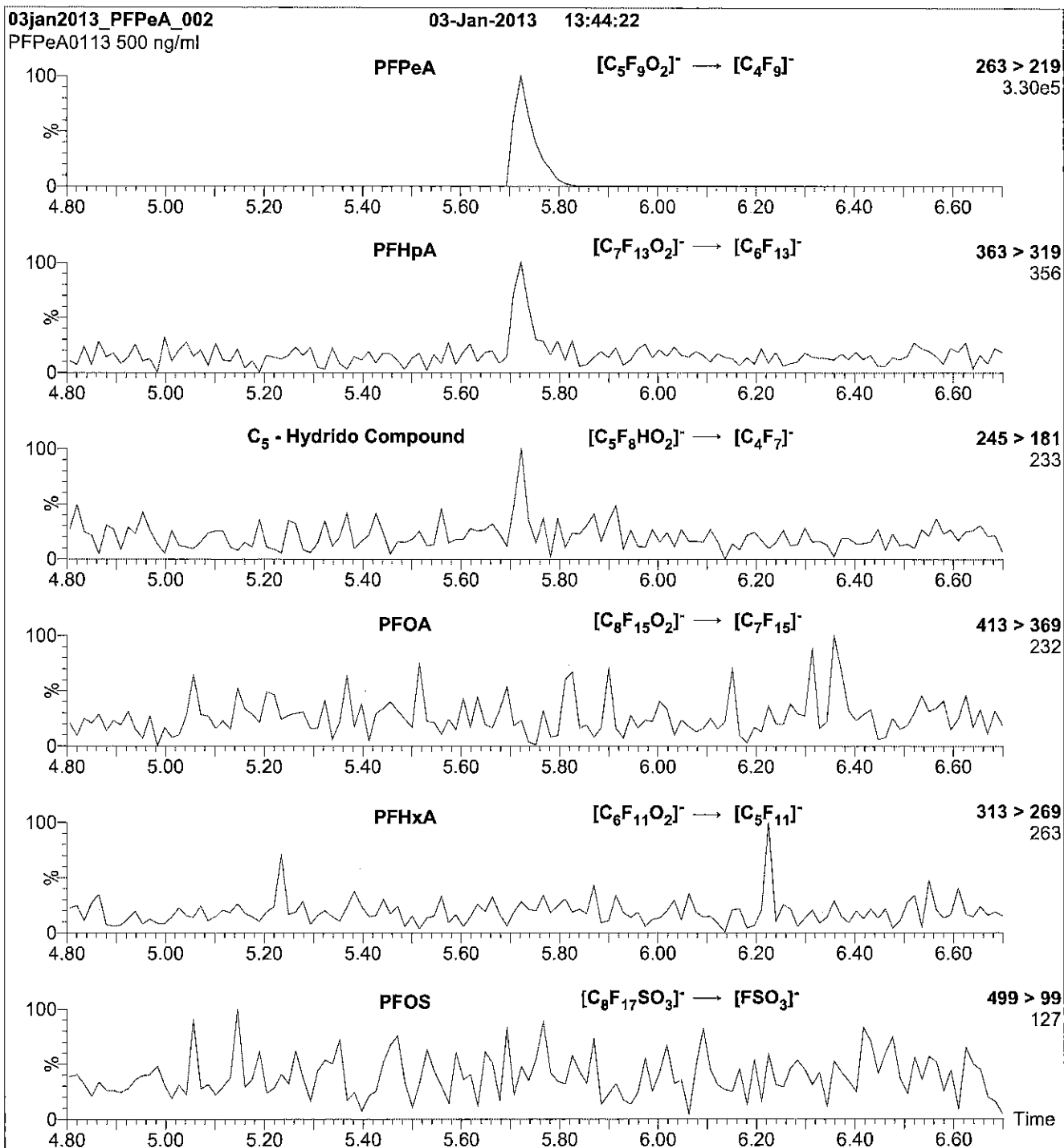
Flow: 300  $\mu$ l/min

**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  $\mu$ l (500 ng/ml PFPeA)

Mobile phase: Isocratic 70% (80:20 MeOH:ACN) / 30% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 9

Reagent

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**LCFPeS\_00002**



R 2445 2



# WELLINGTON LABORATORIES

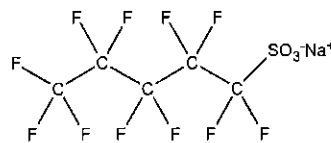
## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** L-PFPeS  
**COMPOUND:** Sodium perfluoro-1-pentanesulfonate

**LOT NUMBER:** LPFPeS0712

**STRUCTURE:**

**CAS #:** Not available



**MOLECULAR FORMULA:** C<sub>5</sub>F<sub>11</sub>SO<sub>3</sub>Na  
**CONCENTRATION:** 50.0 ± 2.5 µg/ml (Na salt)  
 46.9 ± 2.3 µg/ml (PFPeS anion)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 07/04/2012  
**EXPIRY DATE:** (mm/dd/yyyy) 07/04/2017  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

**MOLECULAR WEIGHT:** 372.09  
**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: 01/15/2013  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) are available upon request.

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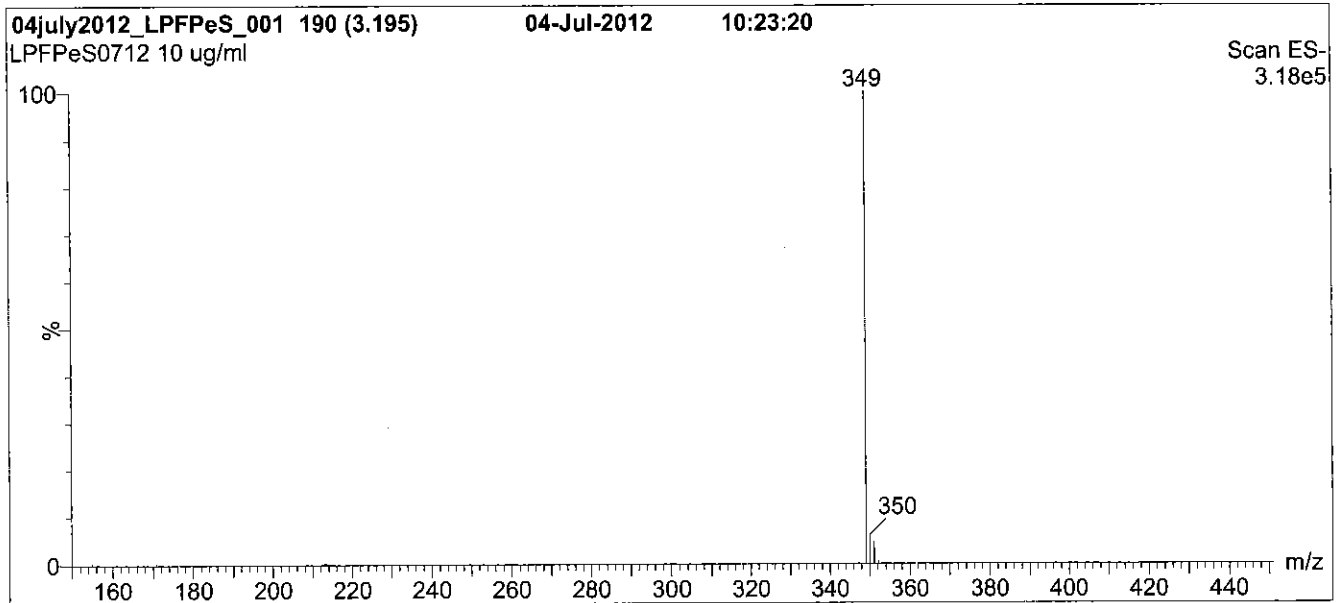
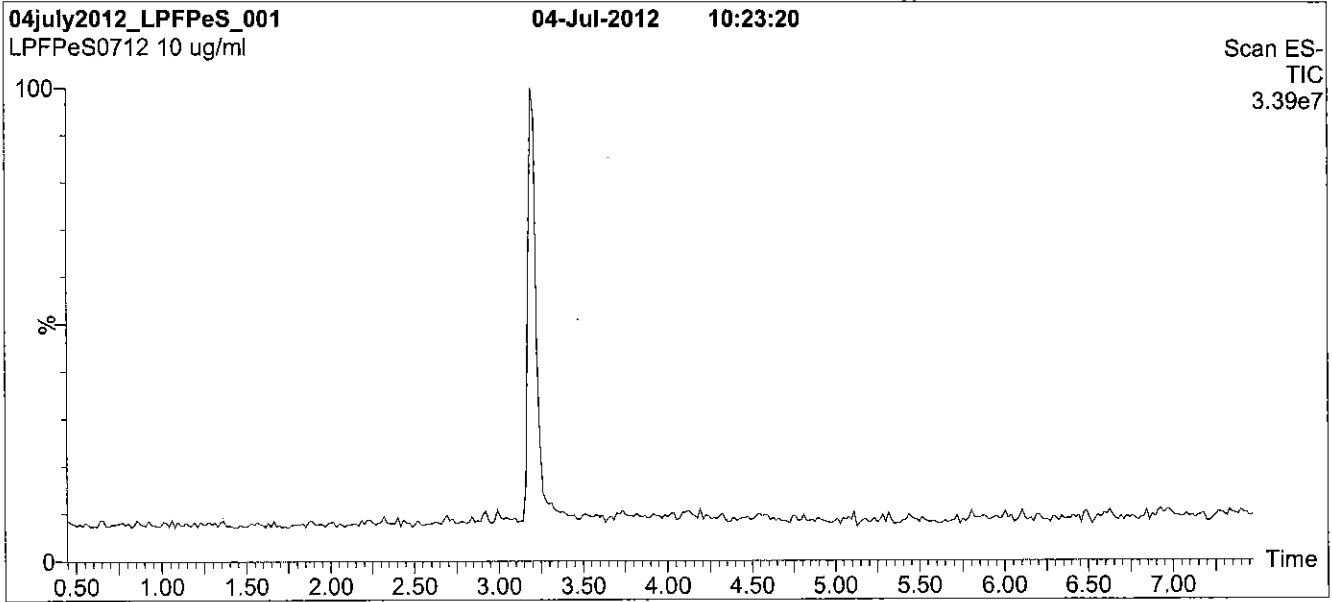
### **QUALITY MANAGEMENT:**

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**Figure 1: L-PFPeS; 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<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

Mobile phase: Gradient  
 Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)  
 Ramp to 90% organic over 7 min and hold for 1.5 min  
 before returning to initial conditions over 0.5 min.  
 Time: 10 min

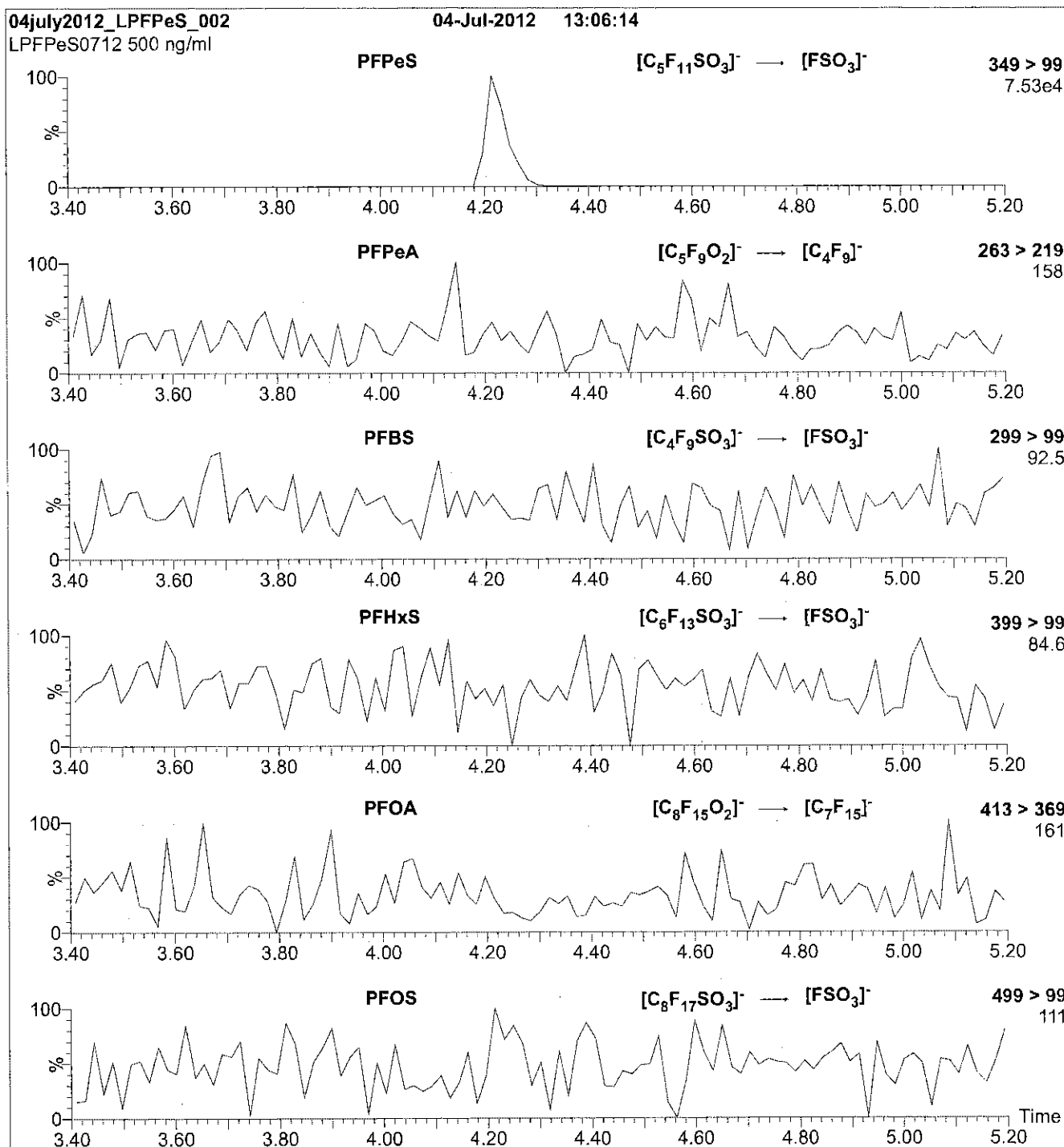
Flow: 300  $\mu$ l/min

**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: L-PFPeS; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

**Injection:** Direct loop injection  
 10  $\mu$ l (500 ng/ml L-PFPeS)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.66e-3  
 Collision Energy (eV) = 30

Reagent

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**LCPFTeDA\_00003**

v: 2/11/15 srw

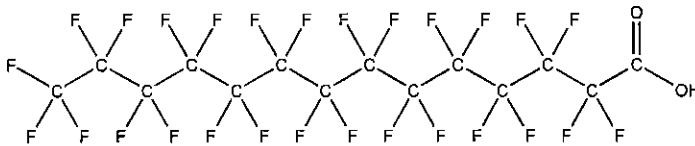


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFTeDA **LOT NUMBER:** PFTeDA0613  
**COMPOUND:** Perfluoro-n-tetradecanoic acid

**STRUCTURE:** **CAS #:** 376-06-7



**MOLECULAR FORMULA:** C<sub>14</sub>HF<sub>27</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 714.11  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
Water (<1%)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 06/19/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 06/19/2018  
**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 ~ 0.2% of PFDoA (C<sub>12</sub>HF<sub>23</sub>O<sub>2</sub>) and ~ 0.2% of PFPeDA (C<sub>15</sub>HF<sub>29</sub>O<sub>2</sub>).

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
B.G. Chittim **Date:** 07/17/2013  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

**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. Material Safety Data Sheets (MSDSs) are available upon request.

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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:

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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:**

Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the 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.

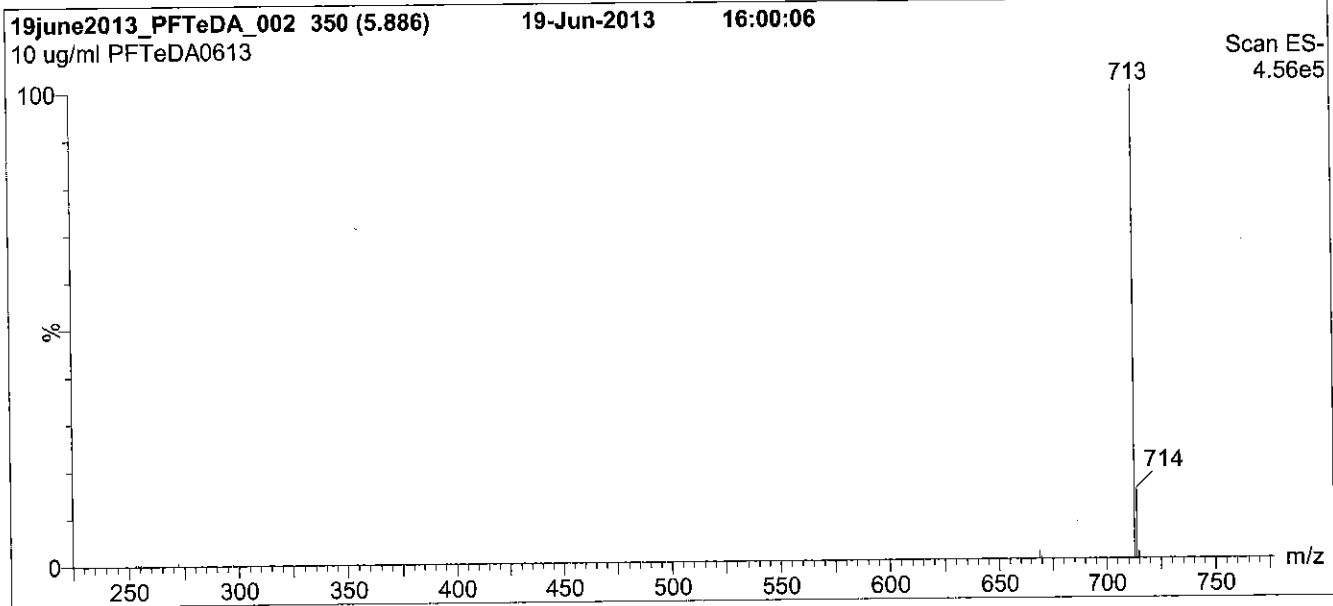
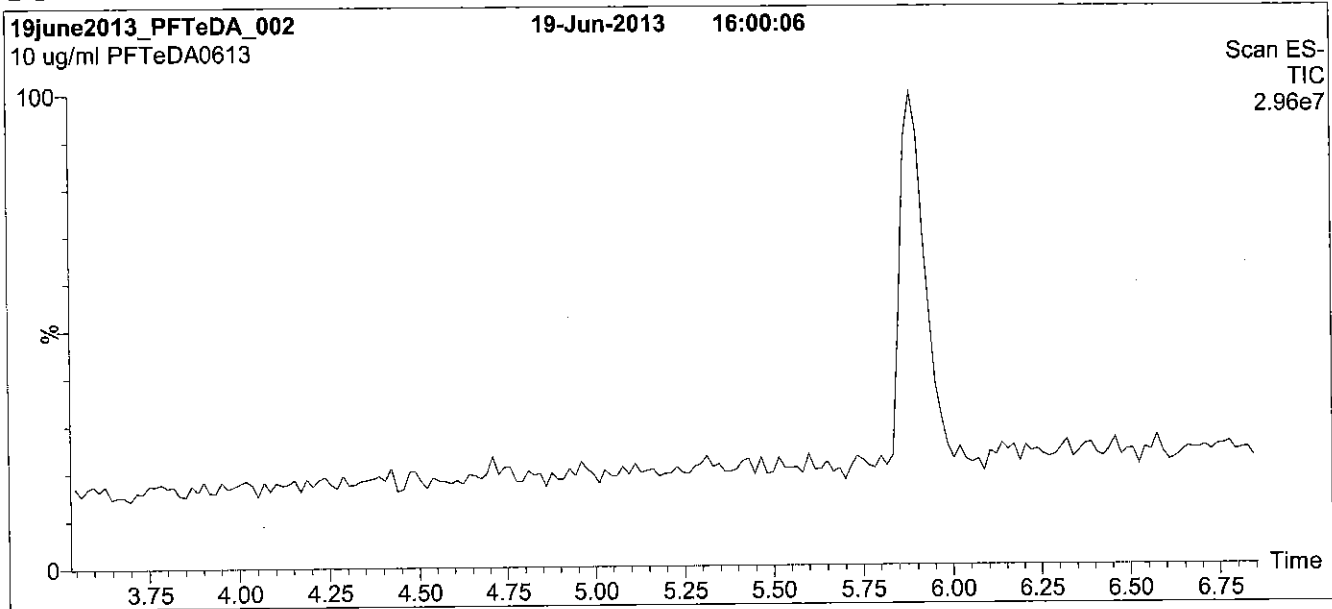
**QUALITY MANAGEMENT:**

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**Figure 1: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min and hold for 1.5 min  
before returning to initial conditions in 0.50 min.  
Time: 10 min

**Flow:** 300  $\mu$ l/min

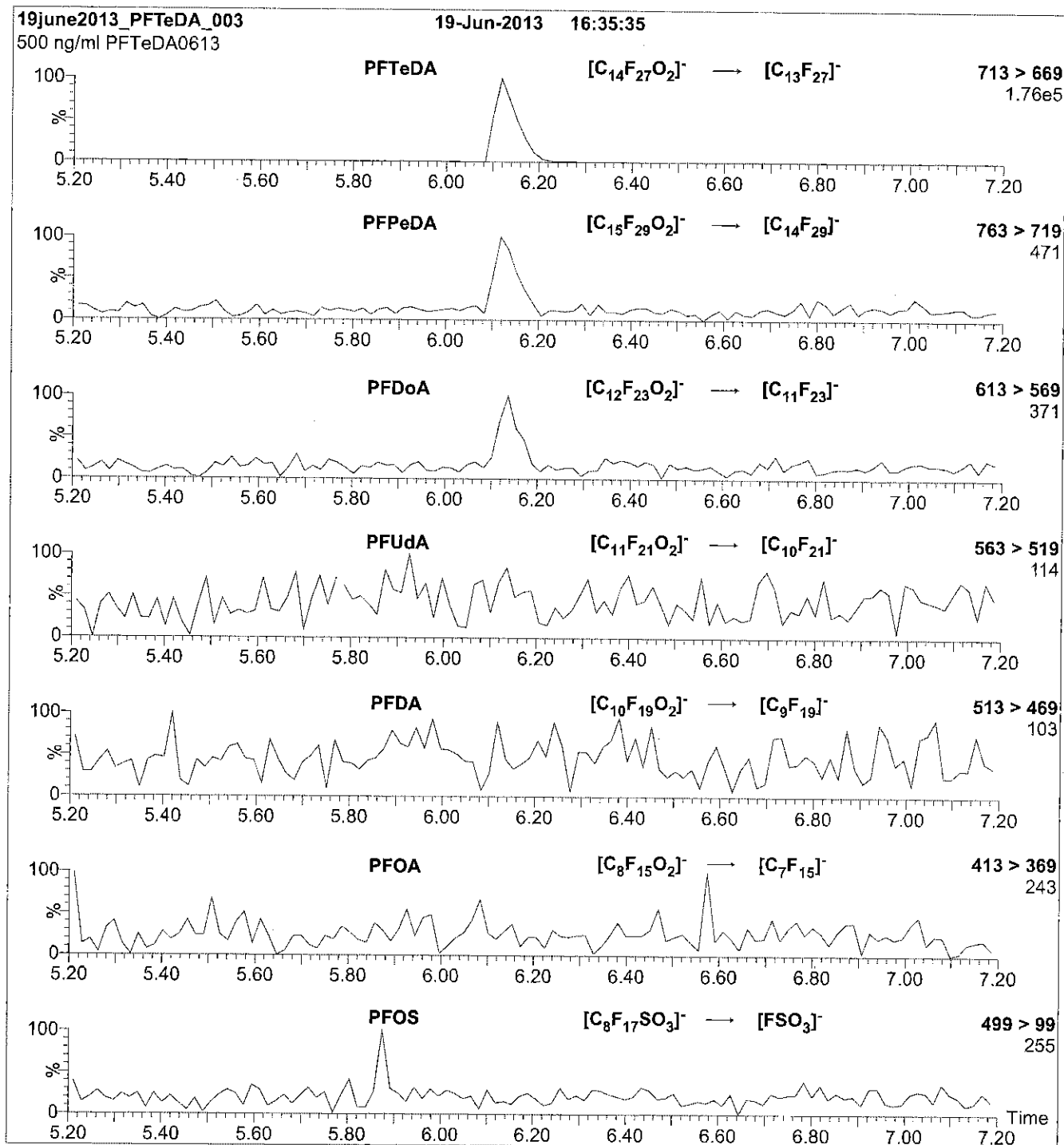
**MS Parameters**

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: PFTeDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct injection  
10  $\mu$ l (500 ng/ml PFTeDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

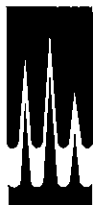
**MS Parameters**

Collision Gas (mbar) = 3.50e-3  
Collision Energy (eV) = 14

Reagent

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**LCPFT<sub>r</sub>DA\_00003**

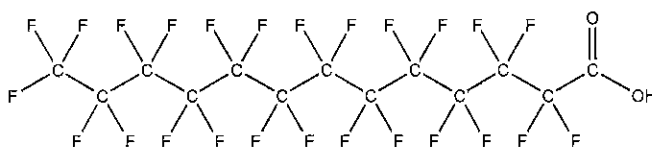


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFTrDA **LOT NUMBER:** PFTrDA1213  
**COMPOUND:** Perfluoro-n-tridecanoic acid

**STRUCTURE:** **CAS #:** 72629-94-8



**MOLECULAR FORMULA:**  $C_{13}HF_{26}O_2$  **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 ( $C_{11}HF_{21}O_2$ ), ~ 0.4% of PFDaA ( $C_{12}HF_{23}O_2$ ), and ~ 0.1% of PFTeDA ( $C_{14}HF_{27}O_2$ ).

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim **Date:** 12/11/2013  
 (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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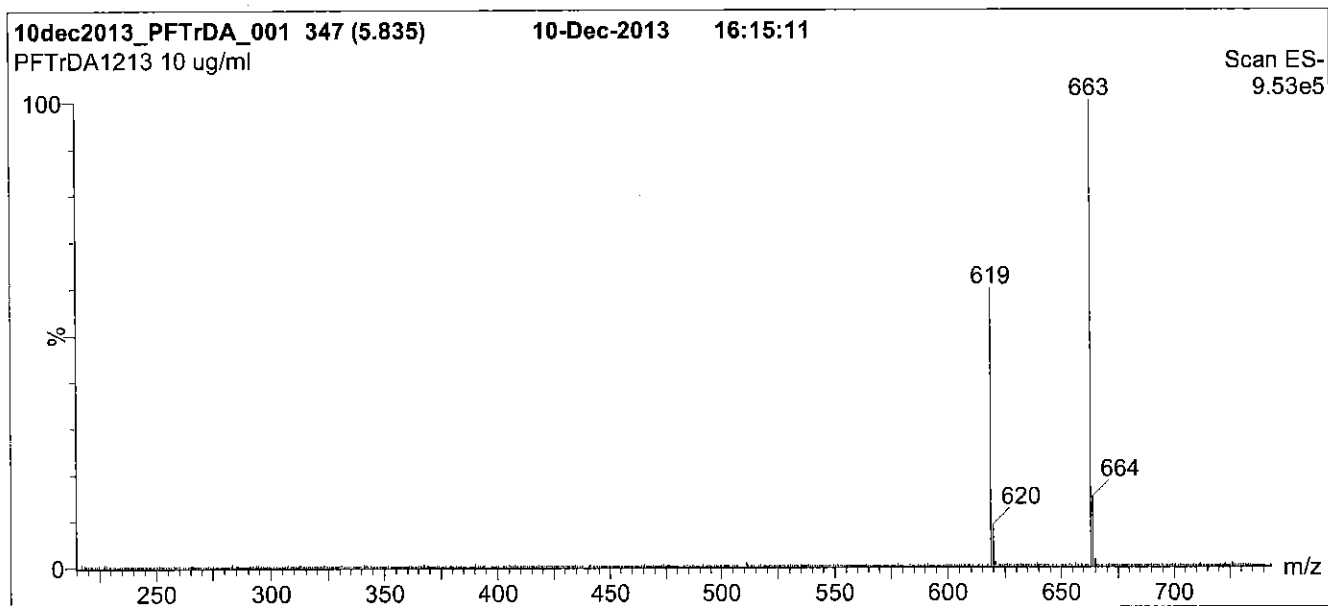
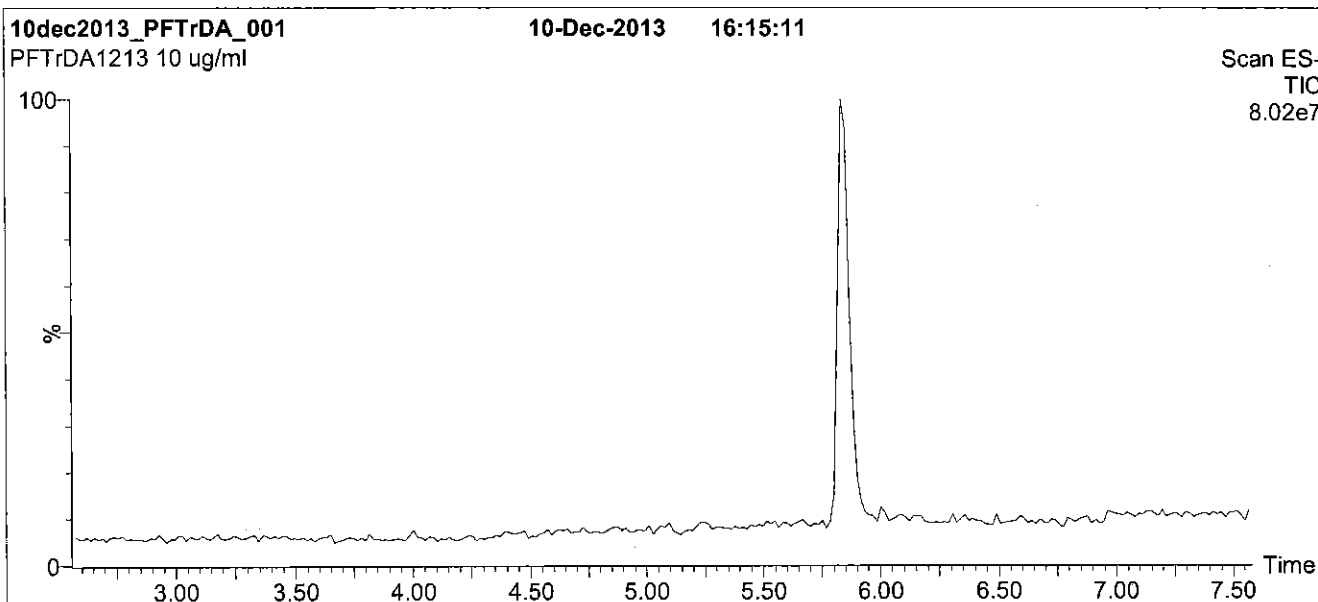
### **QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to ISO 9001:2008 by SAI Global, ISO/IEC 17025:2005 by the Canadian Association for Laboratory Accreditation Inc. (CALA; A 1226), and ISO GUIDE 34:2009 by ACLASS (certificate number 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](http://www.well-labs.com) or contact us directly at [info@well-labs.com](mailto:info@well-labs.com)\*\*

**Figure 1: PFTTrDA; 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<sub>18</sub>  
 1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
 Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc 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

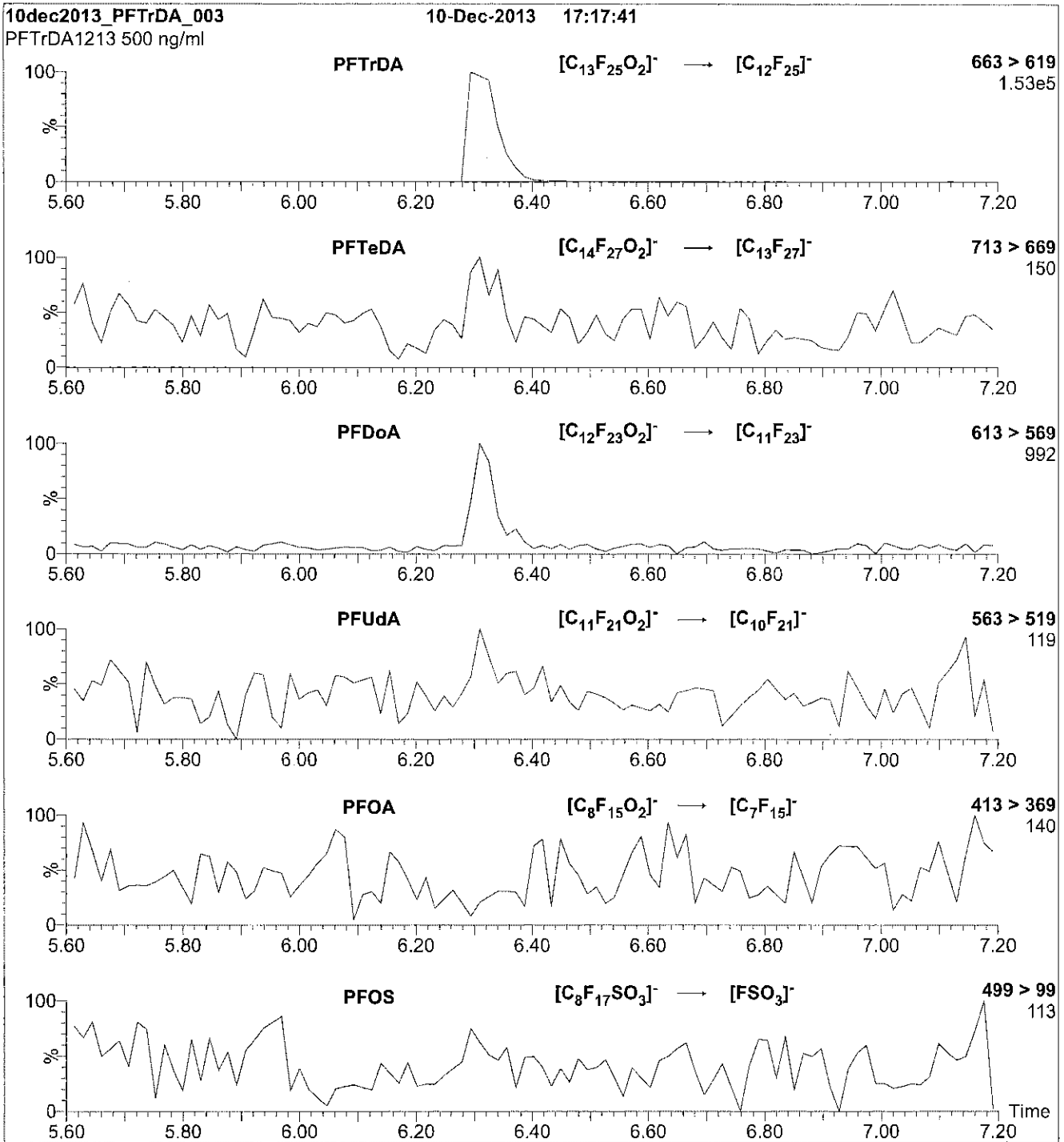
**Flow:** 300  $\mu$ l/min

**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: PFTTrDA; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figure 2:**

Injection: Direct loop injection  
 10  $\mu$ l (500 ng/ml PFTTrDA)

Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
 (both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.28e-3  
 Collision Energy (eV) = 15

Reagent

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**LCPFUdA\_00003**

PC 2/11/15 SFV

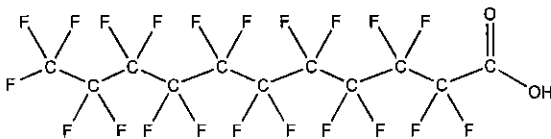


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:** PFUdA **LOT NUMBER:** PFUdA0613  
**COMPOUND:** Perfluoro-n-undecanoic acid

**STRUCTURE:** **CAS #:** 2058-94-8



**MOLECULAR FORMULA:** C<sub>11</sub>HF<sub>21</sub>O<sub>2</sub> **MOLECULAR WEIGHT:** 564.09  
**CONCENTRATION:** 50 ± 2.5 µg/ml **SOLVENT(S):** Methanol  
 Water (<1%)  
**CHEMICAL PURITY:** >98%  
**LAST TESTED:** (mm/dd/yyyy) 06/19/2013  
**EXPIRY DATE:** (mm/dd/yyyy) 06/19/2018  
**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place

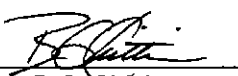
**DOCUMENTATION/ DATA ATTACHED:**

- 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.

**FOR LABORATORY USE ONLY: NOT FOR HUMAN OR DRUG USE**

**Certified By:**   
 B.G. Chittim **Date:** 07/03/2013  
(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. They are designed to be used as reference standards for the identification and/or quantification of specific chemical compound(s).

### **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. Material Safety Data Sheets (MSDSs) 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, 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 and/or LC/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 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 17025:2005 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 for the period of time specified by the 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.

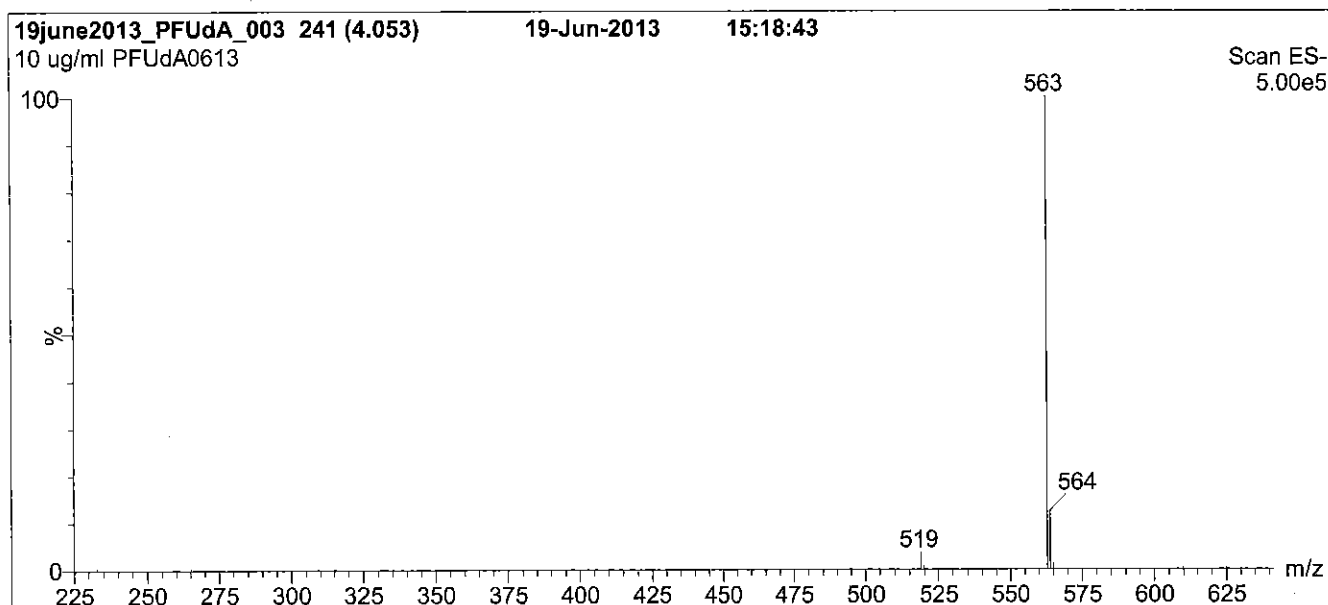
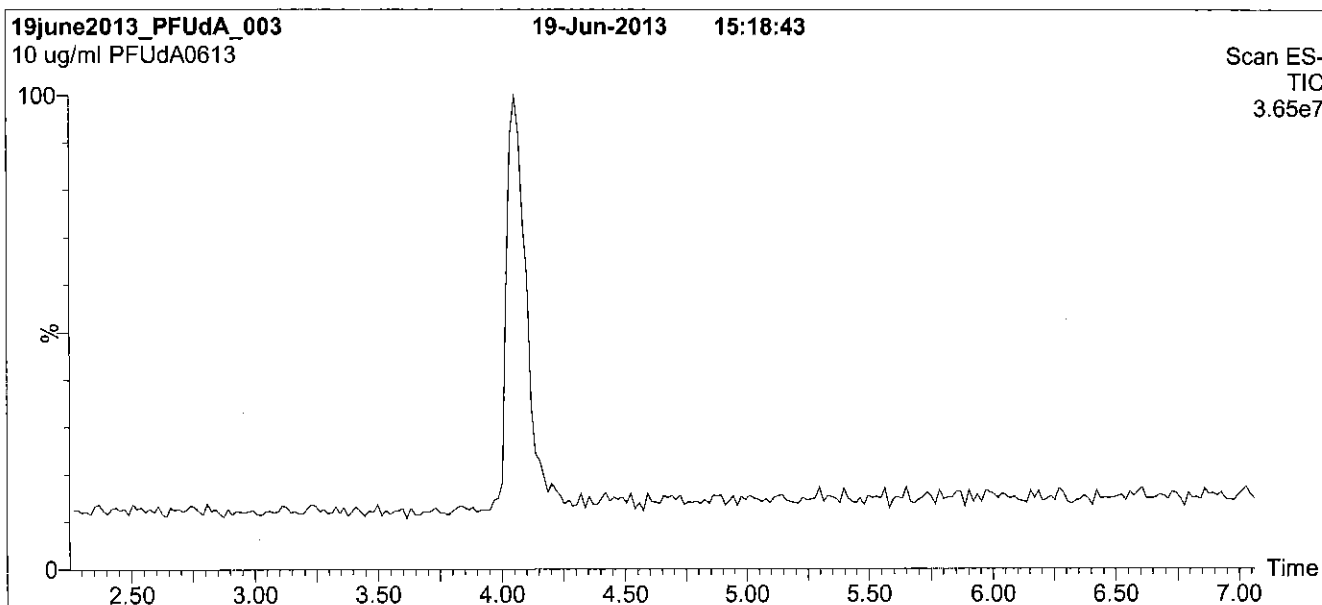
### **QUALITY MANAGEMENT:**

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**Figure 1: 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 RP<sub>18</sub>  
1.7  $\mu$ m, 2.1 x 100 mm

**Mobile phase:** Gradient  
Start: 60% (80:20 MeOH:ACN) / 40% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc 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

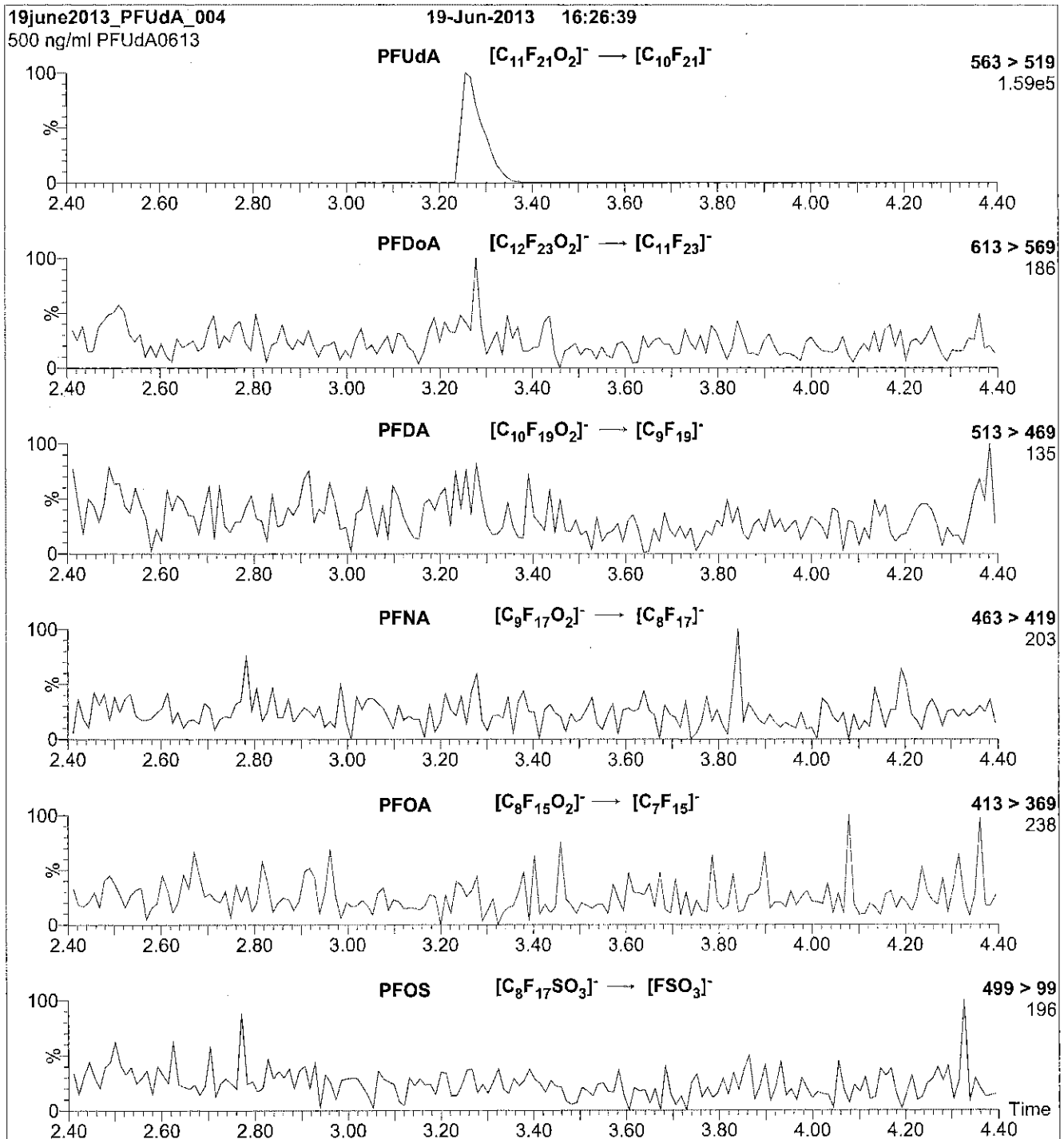
**Flow:** 300  $\mu$ l/min

**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)** = 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  $\mu$ l (500 ng/ml PFUdA)

**Mobile phase:** Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

**Flow:** 300  $\mu$ l/min

**MS Parameters**

Collision Gas (mbar) = 3.46e-3  
Collision Energy (eV) = 11

# Method PFC DOD

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Perfluronated Hydrocarbons (LC/MS)  
by Method PFC\_DOD

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Matrix: Water

Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	13CHpA #	PFHxS #	PFOA #	PFOS #	PFNA #
OF-FB09-0216	320-17154-1	122	113	126	107	127
OF-FB11-0216	320-17154-6	111	95	113	93	116
OF-FB28-0216	320-17154-8	124	113	132	108	137
OF-FB67-0216	320-17154-10	126	113	131	112	130
	MB 320-100147/1-A	118	97	114	98	118
	LCS 320-100093/2-A	95	90	95	82	98
	LCS 320-100147/2-A	124	108	119	99	117
	LCSD 320-100093/3-A	95	86	88	80	91
	LCSD 320-100147/3-A	123	107	115	100	116

13CHpA = 13C4-PFHpA  
 PFHxS = 1802 PFHxS  
 PFOA = 13C4 PFOA  
 PFOS = 13C4 PFOS  
 PFNA = 13C5 PFNA

QC LIMITS  
 25-150  
 25-150  
 25-150  
 25-150  
 25-150

# Column to be used to flag recovery values

FORM II WS-LC-0025

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Matrix: Water Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	13CHpA #	PFOA #	PFOS #	PFNA #
	MB 320-100093/1-A	103	101	92	103

13CHpA = 13C4-PFHpA  
PFOA = 13C4 PFOA  
PFOS = 13C4 PFOS  
PFNA = 13C5 PFNA

QC LIMITS  
25-150  
25-150  
25-150  
25-150

# Column to be used to flag recovery values

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Matrix: Water

Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFHxS #
	MB 320-100093/1-A	95

PFHxS = 1802 PFHxS

QC LIMITS  
25-150

# Column to be used to flag recovery values

FORM II WS-LC-0025

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Matrix: Water Level: Low  
 GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFHxS #	13CHpA #	PFOA #	PFOS #	PFNA #
OF-FB37-0216	320-17154-3	100	107	111	95	113

PFHxS = 1802 PFHxS  
 13CHpA = 13C4-PFHpA  
 PFOA = 13C4 PFOA  
 PFOS = 13C4 PFOS  
 PFNA = 13C5 PFNA

QC LIMITS  
 25-150  
 25-150  
 25-150  
 25-150  
 25-150

# Column to be used to flag recovery values

FORM II WS-LC-0025



FORM III  
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Matrix: Water Level: Low Lab File ID: 19FEB2016A6A\_048.d  
 Lab ID: LCS 320-100093/2-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0513	128	60-140	
Perfluorooctanoic acid (PFOA)	0.0400	0.0485	121	60-140	
Perfluorononanoic acid (PFNA)	0.0400	0.0537	134	60-140	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0459	130	50-150	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0480	127	60-140	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0515	135	60-140	
13C4-PFHpA	0.100	0.0955	95	25-150	
13C4 PFOA	0.100	0.0945	95	25-150	
13C5 PFNA	0.100	0.0975	98	25-150	
18O2 PFHxS	0.0946	0.0852	90	25-150	
13C4 PFOS	0.0956	0.0783	82	25-150	

# Column to be used to flag recovery and RPD values

FORM III  
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Matrix: Water Level: Low Lab File ID: 19FEB2016A6A\_062.d  
 Lab ID: LCS 320-100147/2-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0498	124	60-140	
Perfluorooctanoic acid (PFOA)	0.0400	0.0476	119	60-140	
Perfluorononanoic acid (PFNA)	0.0400	0.0528	132	60-140	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0412	116	50-150	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0503	133	60-140	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0529	138	60-140	
13C4-PFHpA	0.100	0.124	124	25-150	
13C4 PFOA	0.100	0.119	119	25-150	
13C5 PFNA	0.100	0.117	117	25-150	
18O2 PFHxS	0.0946	0.102	108	25-150	
13C4 PFOS	0.0956	0.0948	99	25-150	

# Column to be used to flag recovery and RPD values

FORM III  
LCMS LAB CONTROL SAMPLE DUPLICATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Matrix: Water Level: Low Lab File ID: 19FEB2016A6A\_049.d  
 Lab ID: LCSD 320-100093/3-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCSD CONCENTRATION (ug/L)	LCSD % REC	% RPD	QC LIMITS		#
					RPD	REC	
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0506	126	1	30	60-140	
Perfluorooctanoic acid (PFOA)	0.0400	0.0493	123	2	30	60-140	
Perfluorononanoic acid (PFNA)	0.0400	0.0517	129	4	30	60-140	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0472	133	3	30	50-150	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0486	128	1	30	60-140	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0452	118	13	30	60-140	
13C4-PFHpA	0.100	0.0955	95			25-150	
13C4 PFOA	0.100	0.0880	88			25-150	
13C5 PFNA	0.100	0.0909	91			25-150	
18O2 PFHxS	0.0946	0.0810	86			25-150	
13C4 PFOS	0.0956	0.0761	80			25-150	

# Column to be used to flag recovery and RPD values

FORM III  
LCMS LAB CONTROL SAMPLE DUPLICATE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Matrix: Water Level: Low Lab File ID: 19FEB2016A6A\_063.d  
 Lab ID: LCSD 320-100147/3-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ug/L)	LCSD CONCENTRATION (ug/L)	LCSD % REC	% RPD	QC LIMITS		#
					RPD	REC	
Perfluoroheptanoic acid (PFHpA)	0.0400	0.0461	115	8	30	60-140	
Perfluorooctanoic acid (PFOA)	0.0400	0.0488	122	2	30	60-140	
Perfluorononanoic acid (PFNA)	0.0400	0.0534	133	1	30	60-140	
Perfluorobutanesulfonic acid (PFBS)	0.0354	0.0461	130	11	30	50-150	
Perfluorohexanesulfonic acid (PFHxS)	0.0378	0.0473	125	6	30	60-140	
Perfluorooctanesulfonic acid (PFOS)	0.0382	0.0502	131	5	30	60-140	
13C4-PFHpA	0.100	0.123	123			25-150	
13C4 PFOA	0.100	0.115	115			25-150	
13C5 PFNA	0.100	0.116	116			25-150	
18O2 PFHxS	0.0946	0.101	107			25-150	
13C4 PFOS	0.0956	0.0953	100			25-150	

# Column to be used to flag recovery and RPD values

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab File ID: 12FEB2016A4A\_042.d Lab Sample ID: MB 320-100093/1-A  
 Matrix: Water Date Extracted: 02/09/2016 10:23  
 Instrument ID: A4 Date Analyzed: 02/13/2016 00:22  
 Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-100093/2-A	19FEB2016A6 A 048.d	02/20/2016 06:44
	LCSD 320-100093/3-A	19FEB2016A6 A 049.d	02/20/2016 07:05
OF-FB09-0216	320-17154-1	19FEB2016A6 A 058.d	02/20/2016 10:16
OF-FB37-0216	320-17154-3	19FEB2016A6 A 059.d	02/20/2016 10:37
OF-FB11-0216	320-17154-6	19FEB2016A6 A 060.d	02/20/2016 10:58

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab File ID: 19FEB2016A6A\_047.d Lab Sample ID: MB 320-100093/1-A  
 Matrix: Water Date Extracted: 02/09/2016 10:23  
 Instrument ID: A6 Date Analyzed: 02/20/2016 06:23  
 Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab File ID: 19FEB2016A6A\_061.d Lab Sample ID: MB 320-100147/1-A  
 Matrix: Water Date Extracted: 02/10/2016 07:14  
 Instrument ID: A6 Date Analyzed: 02/20/2016 11:20  
 Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-100147/2-A	19FEB2016A6 A 062.d	02/20/2016 11:41
	LCSD 320-100147/3-A	19FEB2016A6 A 063.d	02/20/2016 12:02
OF-FB28-0216	320-17154-8	19FEB2016A6 A 064.d	02/20/2016 12:23
OF-FB67-0216	320-17154-10	19FEB2016A6 A 065.d	02/20/2016 12:45

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-17154-1</u>
SDG No.: <u>CTO WE7G PFC Sampling</u>	
Client Sample ID: <u>OF-FB09-0216</u>	Lab Sample ID: <u>320-17154-1</u>
Matrix: <u>Water</u>	Lab File ID: <u>19FEB2016A6A_058.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>02/03/2016 16:40</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>02/09/2016 10:23</u>
Sample wt/vol: <u>504.9(mL)</u>	Date Analyzed: <u>02/20/2016 10:16</u>
Con. Extract Vol.: <u>1.00(mL)</u>	Dilution Factor: <u>1</u>
Injection Volume: <u>15(uL)</u>	GC Column: <u>Acquity</u> ID: <u>2.1(mm)</u>
% Moisture: _____	GPC Cleanup: (Y/N) <u>N</u>
Analysis Batch No.: <u>100906</u>	Units: <u>ug/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.0020	0.00079
335-67-1	Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.0020	0.00074
375-95-1	Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.0020	0.00091
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.0020	0.00086
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	122		25-150
STL00990	13C4 PFOA	126		25-150
STL00995	13C5 PFNA	127		25-150
STL00994	18O2 PFHxS	113		25-150
STL00991	13C4 PFOS	107		25-150



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_058.d  
 Lims ID: 320-17154-A-1-A Lab Sample ID: 320-17154-1  
 Client ID: OF-FB09-0216  
 Sample Type: Client  
 Inject. Date: 20-Feb-2016 10:16:31 ALS Bottle#: 11 Worklist Smp#: 56  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-17154-A-1-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:16:45

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 8 13C4-PFHpA	367.0 > 322.0	9.294	9.322	-0.028	1879807	61.0		122	7119	
D 11 18O2 PFHxS	403.0 > 84.0	9.329	9.358	-0.029	797748	53.3		113	3508	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.305	9.364	-0.059	1469	0.1290				
D 12 13C4 PFOA	417.0 > 372.0	10.419	10.450	-0.031	2152297	63.0		126	8027	
13 Perfluorooctanoic acid	413.0 > 369.0	10.426	10.450	-0.024	2132	0.0493			2.6	
D 16 13C4 PFOS	503.0 > 80.0	11.385	11.414	-0.029	930905	51.2		107	3639	
D 17 13C5 PFNA	468.0 > 423.0	11.408	11.434	-0.026	1792957	63.6		127	4359	

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_058.d

Injection Date: 20-Feb-2016 10:16:31

Instrument ID: A6

Lims ID: 320-17154-A-1-A

Lab Sample ID: 320-17154-1

Client ID: OF-FB09-0216

Operator ID: JRB

ALS Bottle#: 11

Worklist Smp#: 56

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

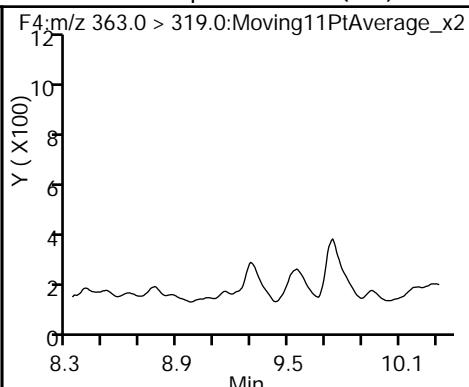
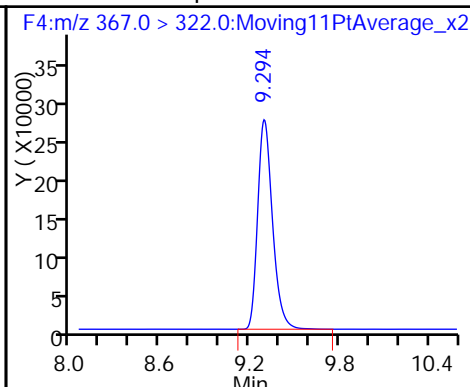
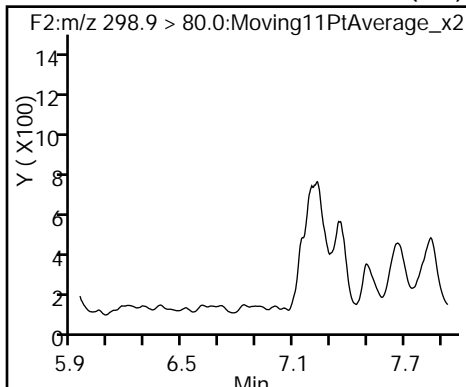
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

40 Perfluorobutanesulfonic acid (ND)

D 8 13C4-PFHpA

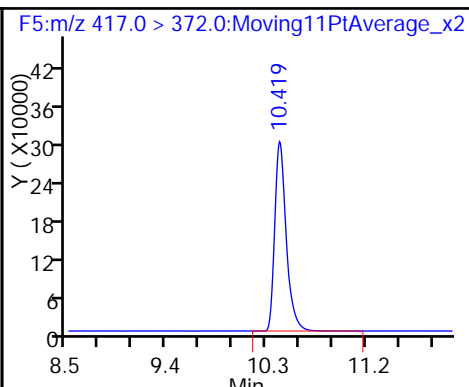
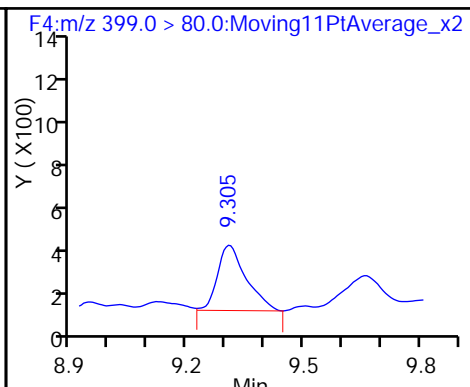
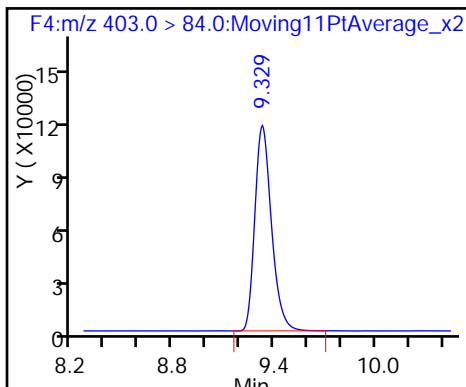
9 Perfluoroheptanoic acid (ND)



D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

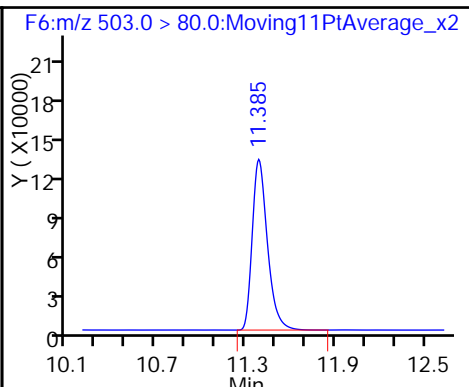
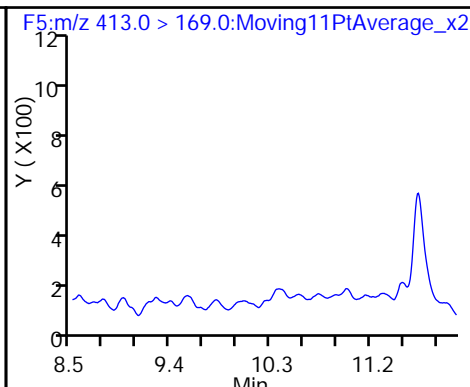
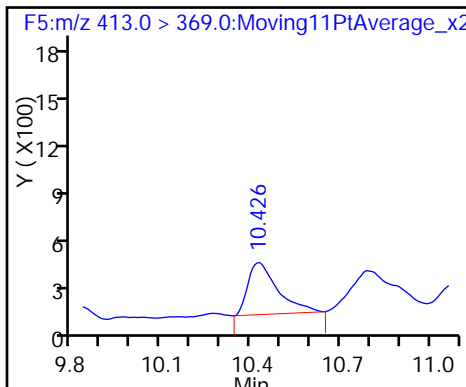
D 12 13C4 PFOA



13 Perfluorooctanoic acid

13 Perfluorooctanoic acid

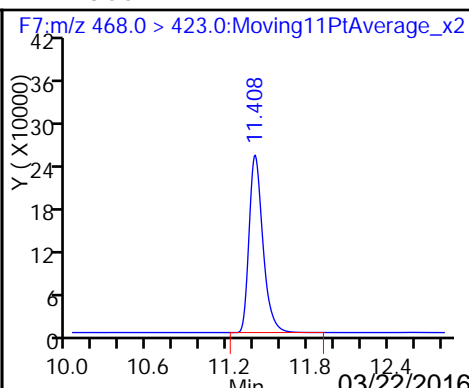
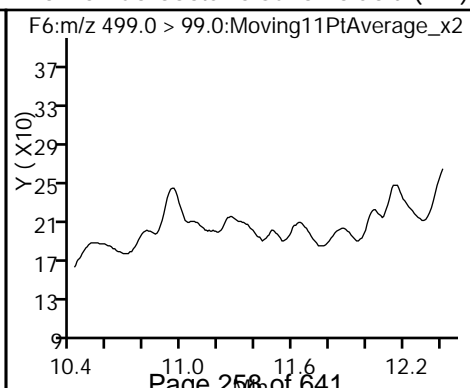
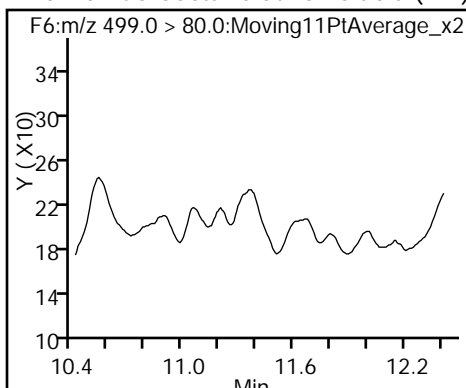
D 16 13C4 PFOS



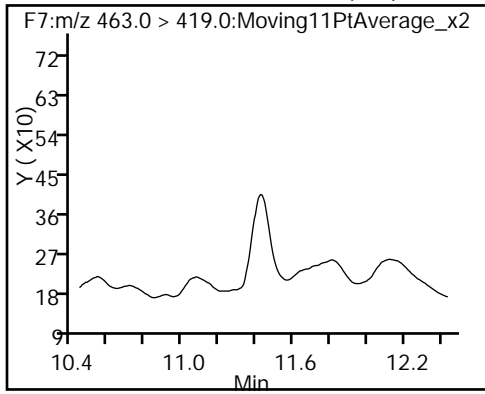
15 Perfluorooctane sulfonic acid (ND)

15 Perfluorooctane sulfonic acid (ND)

D 17 13C5 PFNA



18 Perfluorononanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-17154-1</u>
SDG No.: <u>CTO WE7G PFC Sampling</u>	
Client Sample ID: <u>OF-FB37-0216</u>	Lab Sample ID: <u>320-17154-3</u>
Matrix: <u>Water</u>	Lab File ID: <u>19FEB2016A6A_059.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>02/03/2016 17:20</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>02/09/2016 10:23</u>
Sample wt/vol: <u>550.7(mL)</u>	Date Analyzed: <u>02/20/2016 10:37</u>
Con. Extract Vol.: <u>1.00(mL)</u>	Dilution Factor: <u>1</u>
Injection Volume: <u>15(uL)</u>	GC Column: <u>Acquity</u> ID: <u>2.1(mm)</u>
% Moisture: _____	GPC Cleanup: (Y/N) <u>N</u>
Analysis Batch No.: <u>100906</u>	Units: <u>ug/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.0018	0.00073
335-67-1	Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.0018	0.00068
375-95-1	Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.0018	0.00059
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.0018	0.00083
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.0018	0.00079
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0027	0.0012

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	107		25-150
STL00990	13C4 PFOA	111		25-150
STL00995	13C5 PFNA	113		25-150
STL00994	18O2 PFHxS	100		25-150
STL00991	13C4 PFOS	95		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_059.d  
 Lims ID: 320-17154-A-3-A Lab Sample ID: 320-17154-3  
 Client ID: OF-FB37-0216  
 Sample Type: Client  
 Inject. Date: 20-Feb-2016 10:37:44 ALS Bottle#: 12 Worklist Smp#: 57  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-17154-A-3-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:17:16

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 8 13C4-PFHpA	367.0 > 322.0	9.300	9.322	-0.022	1644356	53.3		107	5755	
D 11 18O2 PFHxS	403.0 > 84.0	9.335	9.358	-0.023	707917	47.3		100	2549	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.364	9.364	0.0	905	0.0895				
D 12 13C4 PFOA	417.0 > 372.0	10.426	10.450	-0.024	1890597	55.4		111	7276	
D 16 13C4 PFOS	503.0 > 80.0	11.392	11.414	-0.022	825359	45.4		95.0	4506	
D 17 13C5 PFNA	468.0 > 423.0	11.408	11.434	-0.026	1597035	56.7		113	3808	

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_059.d

Injection Date: 20-Feb-2016 10:37:44

Instrument ID: A6

Lims ID: 320-17154-A-3-A

Lab Sample ID: 320-17154-3

Client ID: OF-FB37-0216

Operator ID: JRB

ALS Bottle#: 12

Worklist Smp#: 57

Injection Vol: 15.0 ul

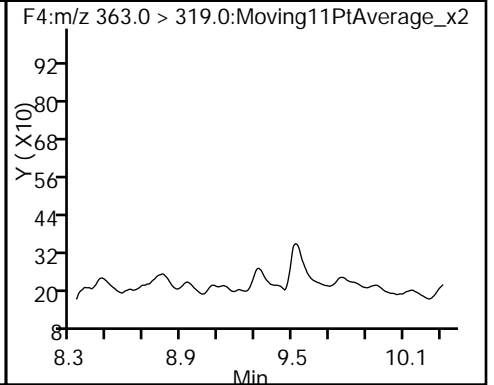
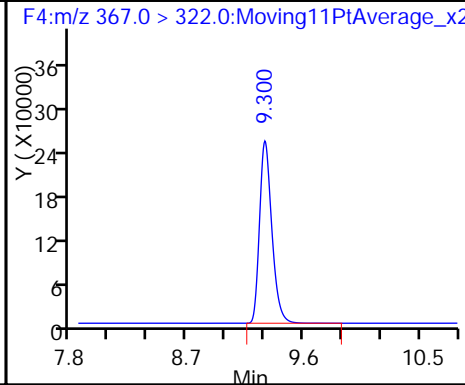
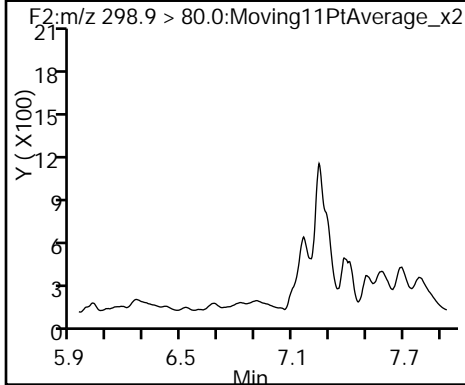
Dil. Factor: 1.0000

Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

40 Perfluorobutanesulfonic acid (ND) D 8 13C4-PFHpA

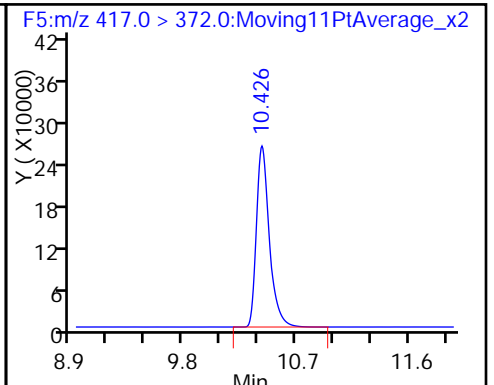
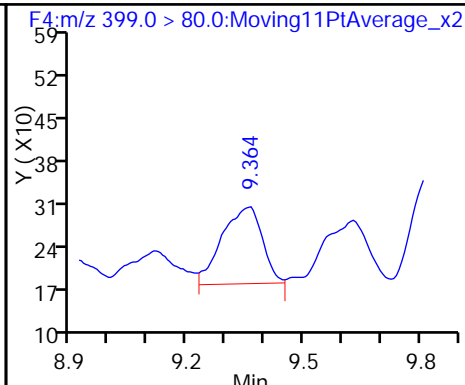
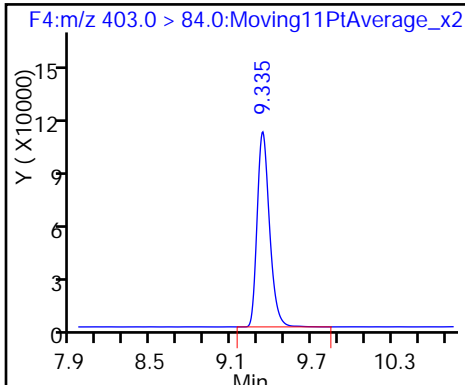
9 Perfluoroheptanoic acid (ND)



D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

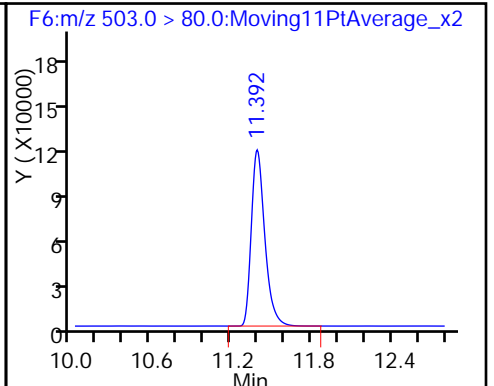
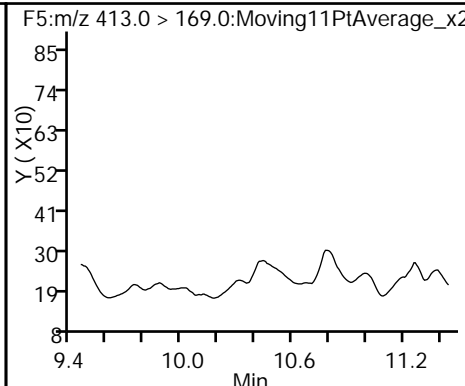
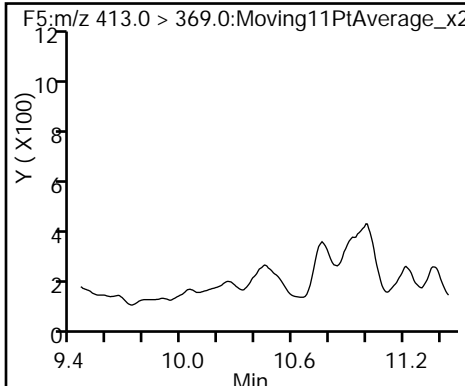
D 12 13C4 PFOA



13 Perfluorooctanoic acid (ND)

13 Perfluorooctanoic acid (ND)

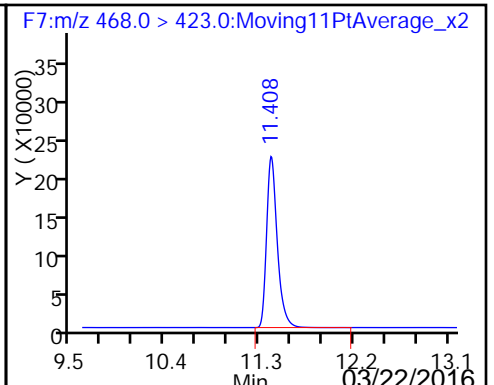
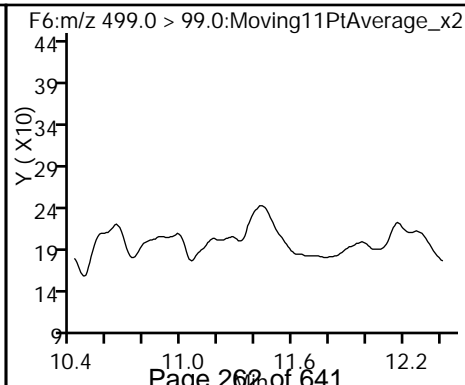
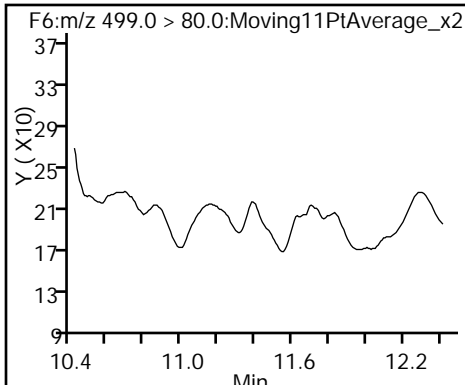
D 16 13C4 PFOS



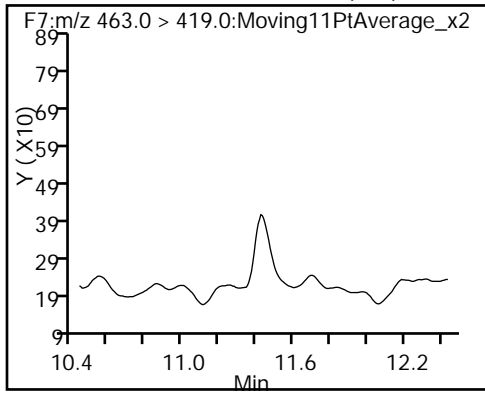
15 Perfluorooctane sulfonic acid (ND)

15 Perfluorooctane sulfonic acid (ND)

D 17 13C5 PFNA



18 Perfluorononanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-17154-1</u>
SDG No.: <u>CTO WE7G PFC Sampling</u>	
Client Sample ID: <u>OF-FB11-0216</u>	Lab Sample ID: <u>320-17154-6</u>
Matrix: <u>Water</u>	Lab File ID: <u>19FEB2016A6A_060.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>02/03/2016 16:10</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>02/09/2016 10:23</u>
Sample wt/vol: <u>551.4 (mL)</u>	Date Analyzed: <u>02/20/2016 10:58</u>
Con. Extract Vol.: <u>1.00 (mL)</u>	Dilution Factor: <u>1</u>
Injection Volume: <u>15 (uL)</u>	GC Column: <u>Acquity</u> ID: <u>2.1 (mm)</u>
% Moisture: _____	GPC Cleanup: (Y/N) <u>N</u>
Analysis Batch No.: <u>100906</u>	Units: <u>ug/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0018	U	0.0023	0.0018	0.00073
335-67-1	Perfluorooctanoic acid (PFOA)	0.0018	U	0.0023	0.0018	0.00068
375-95-1	Perfluorononanoic acid (PFNA)	0.0018	U	0.0023	0.0018	0.00059
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0018	U	0.0023	0.0018	0.00083
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0018	U	0.0023	0.0018	0.00079
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0027	U	0.0036	0.0027	0.0012

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	111		25-150
STL00990	13C4 PFOA	113		25-150
STL00995	13C5 PFNA	116		25-150
STL00994	18O2 PFHxS	95		25-150
STL00991	13C4 PFOS	93		25-150



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_060.d  
 Lims ID: 320-17154-A-6-A Lab Sample ID: 320-17154-6  
 Client ID: OF-FB11-0216  
 Sample Type: Client  
 Inject. Date: 20-Feb-2016 10:58:58 ALS Bottle#: 13 Worklist Smp#: 58  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-17154-A-6-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:17:44

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
40 Perfluorobutanesulfonic acid	298.9 > 80.0	7.130	6.939	0.191	1.000	704	0.0499			
D 8 13C4-PFHpA	367.0 > 322.0	9.299	9.322	-0.023		1708012	55.4	111	6858	
D 11 18O2 PFHxS	403.0 > 84.0	9.334	9.358	-0.024		671948	44.9	94.9	2363	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.346	9.364	-0.018	1.000	1160	0.1209			
D 12 13C4 PFOA	417.0 > 372.0	10.426	10.450	-0.024		1920771	56.3	113	8928	
D 16 13C4 PFOS	503.0 > 80.0	11.391	11.414	-0.023		808180	44.4	93.0	2623	
D 17 13C5 PFNA	468.0 > 423.0	11.414	11.434	-0.020		1630852	57.8	116	5881	

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_060.d

Injection Date: 20-Feb-2016 10:58:58

Instrument ID: A6

Lims ID: 320-17154-A-6-A

Lab Sample ID: 320-17154-6

Client ID: OF-FB11-0216

Operator ID: JRB

ALS Bottle#: 13

Worklist Smp#: 58

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

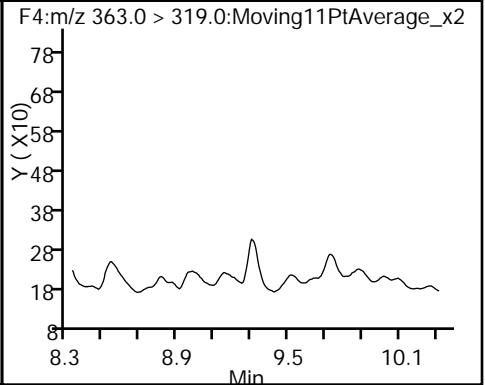
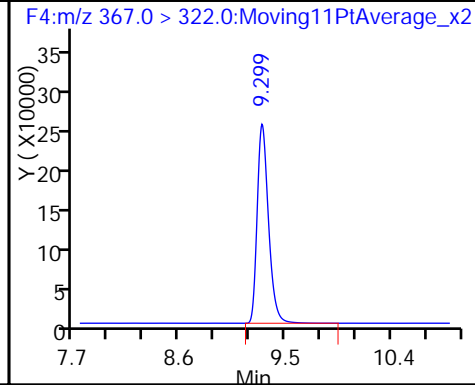
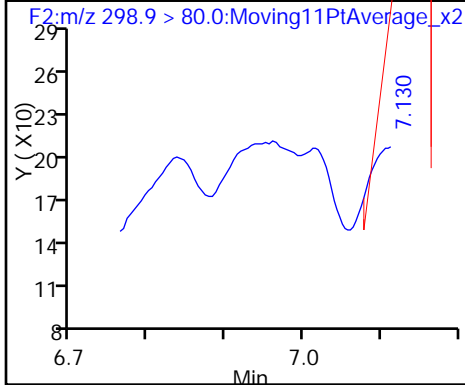
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

40 Perfluorobutanesulfonic acid

D 8 13C4-PFHpA

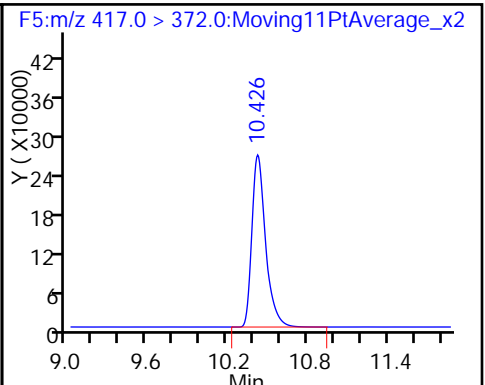
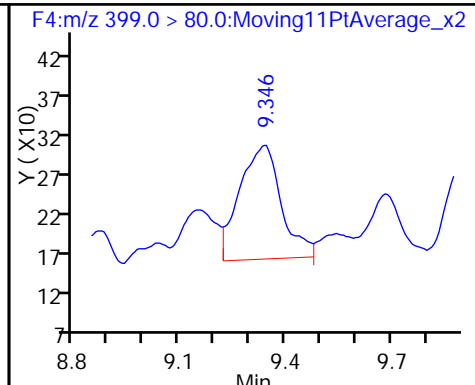
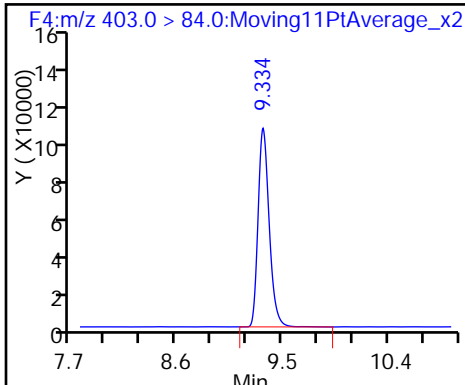
9 Perfluoroheptanoic acid (ND)



D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

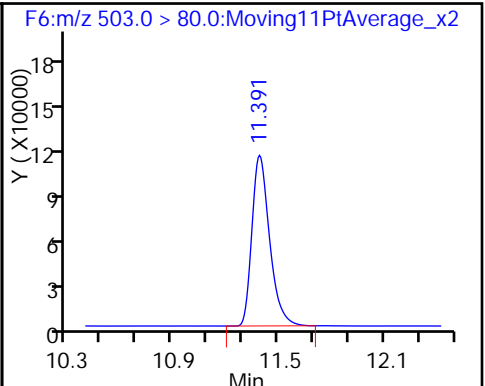
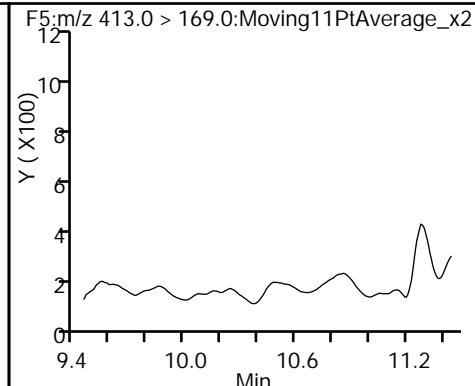
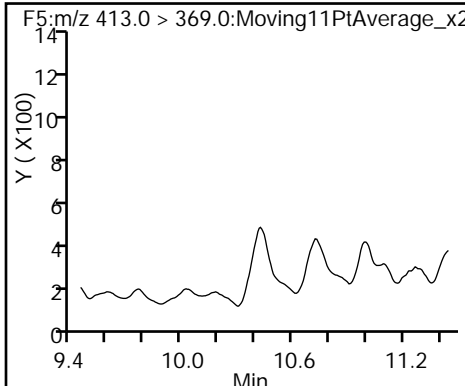
D 12 13C4 PFOA



13 Perfluorooctanoic acid (ND)

13 Perfluorooctanoic acid (ND)

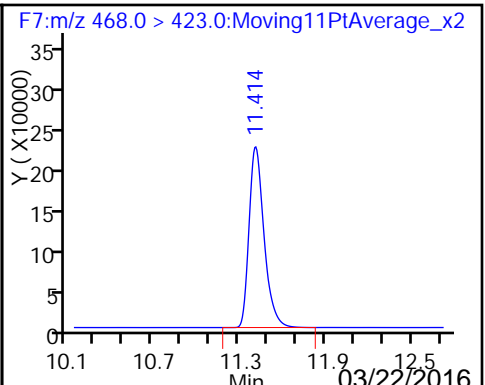
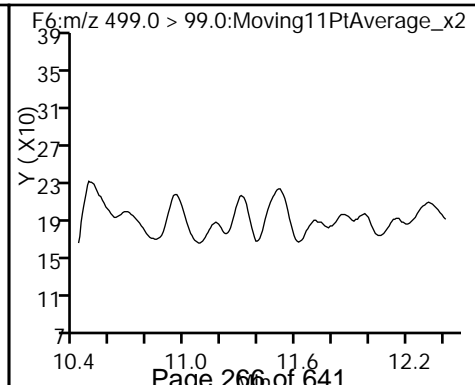
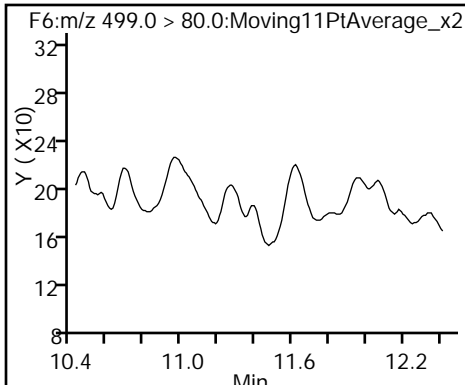
D 16 13C4 PFOS



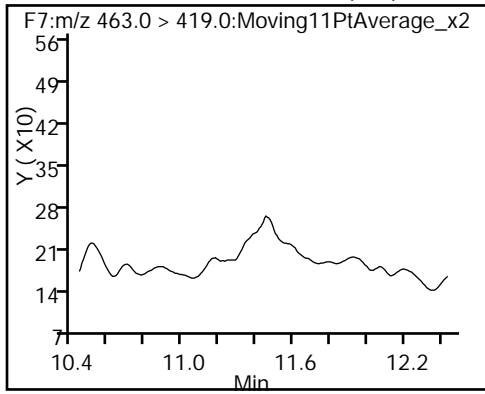
15 Perfluorooctane sulfonic acid (ND)

15 Perfluorooctane sulfonic acid (ND)

D 17 13C5 PFNA



18 Perfluorononanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-17154-1</u>
SDG No.: <u>CTO WE7G PFC Sampling</u>	
Client Sample ID: <u>OF-FB28-0216</u>	Lab Sample ID: <u>320-17154-8</u>
Matrix: <u>Water</u>	Lab File ID: <u>19FEB2016A6A_064.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>02/03/2016 17:43</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>02/10/2016 07:14</u>
Sample wt/vol: <u>521.6(mL)</u>	Date Analyzed: <u>02/20/2016 12:23</u>
Con. Extract Vol.: <u>1.00(mL)</u>	Dilution Factor: <u>1</u>
Injection Volume: <u>15(uL)</u>	GC Column: <u>Acquity</u> ID: <u>2.1(mm)</u>
% Moisture: _____	GPC Cleanup: (Y/N) <u>N</u>
Analysis Batch No.: <u>100906</u>	Units: <u>ug/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0024	0.0019	0.00077
335-67-1	Perfluorooctanoic acid (PFOA)	0.0019	U	0.0024	0.0019	0.00072
375-95-1	Perfluorononanoic acid (PFNA)	0.0019	U	0.0024	0.0019	0.00063
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0024	0.0019	0.00088
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0024	0.0019	0.00083
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0029	U	0.0038	0.0029	0.0012

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	124		25-150
STL00990	13C4 PFOA	132		25-150
STL00995	13C5 PFNA	137		25-150
STL00994	18O2 PFHxS	113		25-150
STL00991	13C4 PFOS	108		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_064.d  
 Lims ID: 320-17154-A-8-A Lab Sample ID: 320-17154-8  
 Client ID: OF-FB28-0216  
 Sample Type: Client  
 Inject. Date: 20-Feb-2016 12:23:54 ALS Bottle#: 17 Worklist Smp#: 62  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-17154-A-8-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:19:52

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
40 Perfluorobutanesulfonic acid	298.9 > 80.0	7.164	6.939	0.225	1.000	1108	0.0660			
D 8 13C4-PFHpA	367.0 > 322.0	9.297	9.322	-0.025		1912777	62.1	124	6834	
D 11 18O2 PFHxS	403.0 > 84.0	9.332	9.358	-0.026		799552	53.4	113	3404	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.315	9.364	-0.049	1.000	1461	0.1280			
D 12 13C4 PFOA	417.0 > 372.0	10.428	10.450	-0.022		2253482	66.0	132	10985	
13 Perfluorooctanoic acid	413.0 > 369.0	10.428	10.450	-0.022	1.000	2011	0.0445		3.1	
D 16 13C4 PFOS	503.0 > 80.0	11.393	11.414	-0.021		942004	51.8	108	3289	
D 17 13C5 PFNA	468.0 > 423.0	11.408	11.434	-0.026		1937075	68.7	137	6887	

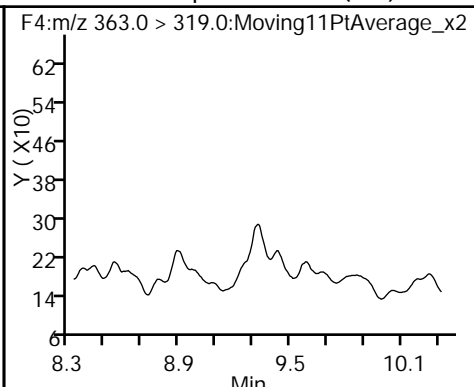
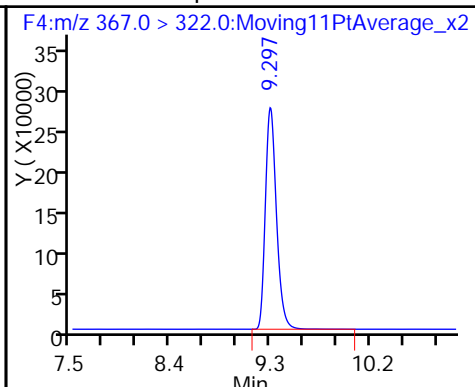
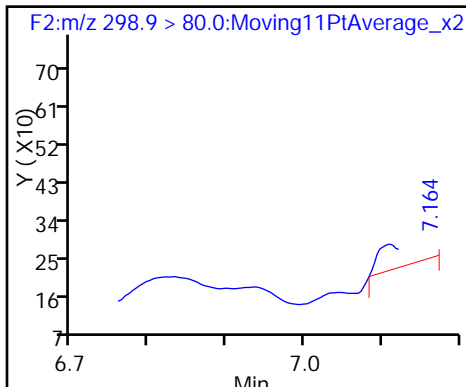
TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_064.d  
Injection Date: 20-Feb-2016 12:23:54 Instrument ID: A6  
Lims ID: 320-17154-A-8-A Lab Sample ID: 320-17154-8  
Client ID: OF-FB28-0216  
Operator ID: JRB ALS Bottle#: 17 Worklist Smp#: 62  
Injection Vol: 15.0 ul Dil. Factor: 1.0000  
Method: PFAC\_A6 Limit Group: LC PFC\_DOD ICAL

40 Perfluorobutanesulfonic acid

D 8 13C4-PFHpA

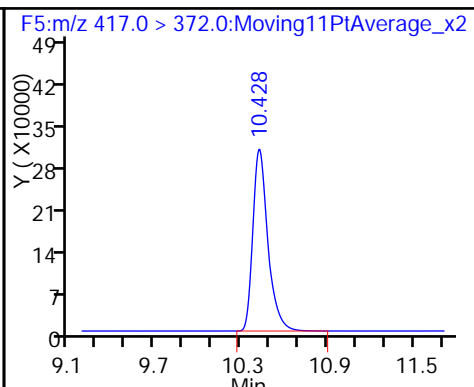
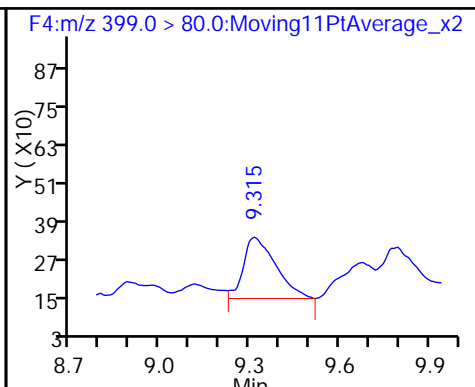
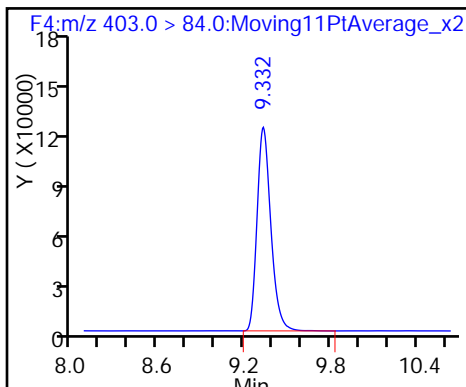
9 Perfluoroheptanoic acid (ND)



D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

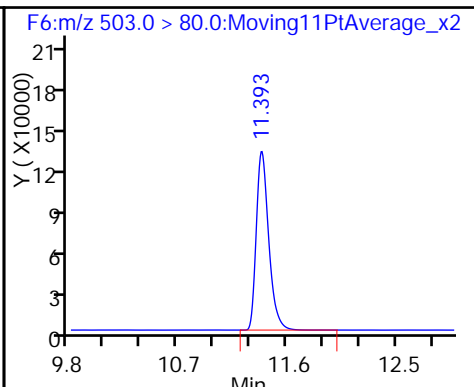
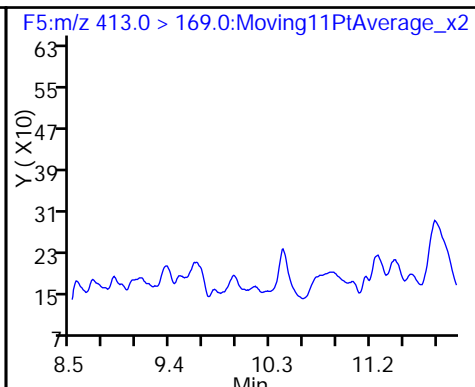
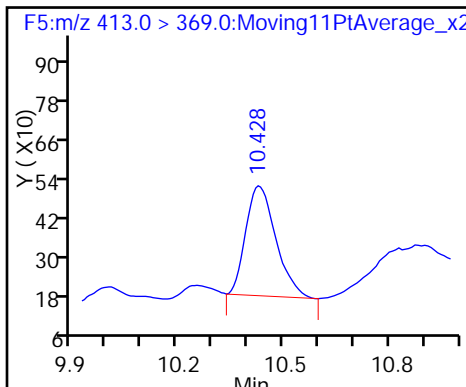
D 12 13C4 PFOA



13 Perfluorooctanoic acid

13 Perfluorooctanoic acid

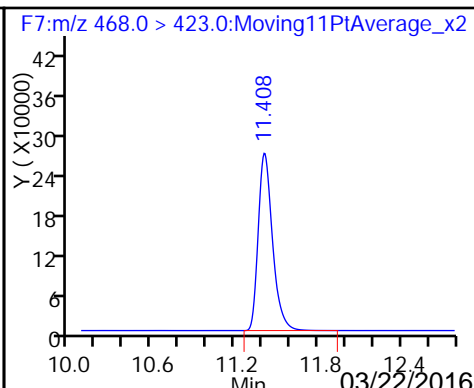
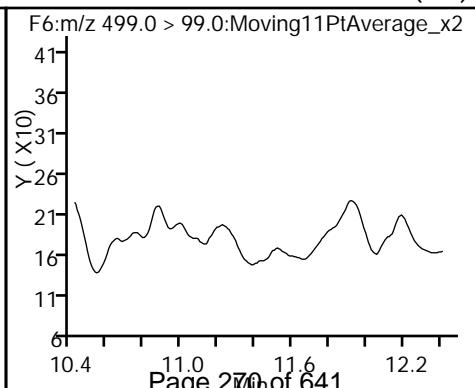
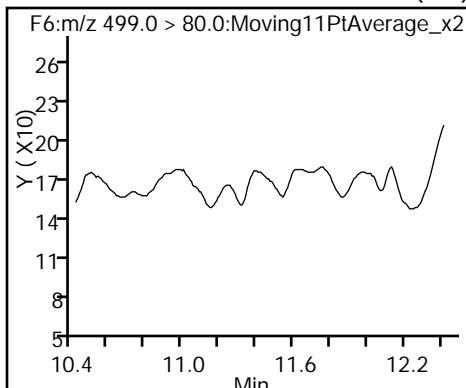
D 16 13C4 PFOS



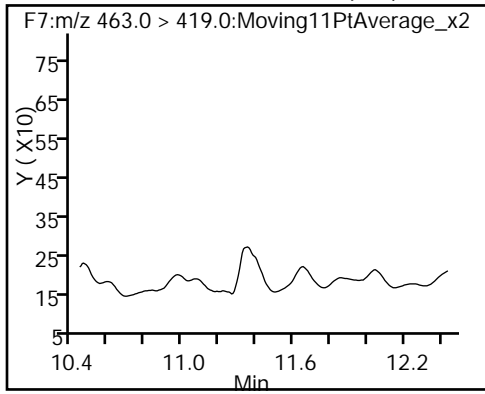
15 Perfluorooctane sulfonic acid (ND)

15 Perfluorooctane sulfonic acid (ND)

D 17 13C5 PFNA



18 Perfluorononanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-17154-1</u>
SDG No.: <u>CTO WE7G PFC Sampling</u>	
Client Sample ID: <u>OF-FB67-0216</u>	Lab Sample ID: <u>320-17154-10</u>
Matrix: <u>Water</u>	Lab File ID: <u>19FEB2016A6A_065.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>02/03/2016 18:21</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>02/10/2016 07:14</u>
Sample wt/vol: <u>535.9(mL)</u>	Date Analyzed: <u>02/20/2016 12:45</u>
Con. Extract Vol.: <u>1.00(mL)</u>	Dilution Factor: <u>1</u>
Injection Volume: <u>15(uL)</u>	GC Column: <u>Acquity</u> ID: <u>2.1(mm)</u>
% Moisture: _____	GPC Cleanup: (Y/N) <u>N</u>
Analysis Batch No.: <u>100906</u>	Units: <u>ug/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0019	U	0.0023	0.0019	0.00075
335-67-1	Perfluorooctanoic acid (PFOA)	0.0019	U	0.0023	0.0019	0.00070
375-95-1	Perfluorononanoic acid (PFNA)	0.0019	U	0.0023	0.0019	0.00061
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0019	U	0.0023	0.0019	0.00086
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0019	U	0.0023	0.0019	0.00081
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0028	U	0.0037	0.0028	0.0012

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	126		25-150
STL00990	13C4 PFOA	131		25-150
STL00995	13C5 PFNA	130		25-150
STL00994	18O2 PFHxS	113		25-150
STL00991	13C4 PFOS	112		25-150



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_065.d  
 Lims ID: 320-17154-A-10-A Lab Sample ID: 320-17154-10  
 Client ID: OF-FB67-0216  
 Sample Type: Client  
 Inject. Date: 20-Feb-2016 12:45:07 ALS Bottle#: 18 Worklist Smp#: 63  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-17154-A-10-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:21:03

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
40 Perfluorobutanesulfonic acid	298.9 > 80.0	7.167	6.939	0.228	1.000	3013	0.1791			
D 8 13C4-PFHpA	367.0 > 322.0	9.299	9.322	-0.023		1943094	63.0	126	7526	
D 11 18O2 PFHxS	403.0 > 84.0	9.334	9.358	-0.024		801416	53.5	113	4035	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.387	9.364	0.023	1.000	1955	0.1709			
D 12 13C4 PFOA	417.0 > 372.0	10.426	10.450	-0.024		2233440	65.4	131	8349	
13 Perfluorooctanoic acid	413.0 > 369.0	10.426	10.450	-0.024	1.000	2584	0.0576		4.0	
D 16 13C4 PFOS	503.0 > 80.0	11.392	11.414	-0.022		977166	53.7	112	3816	
D 17 13C5 PFNA	468.0 > 423.0	11.415	11.434	-0.019		1830448	64.9	130	10168	

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_065.d

Injection Date: 20-Feb-2016 12:45:07

Instrument ID: A6

Lims ID: 320-17154-A-10-A

Lab Sample ID: 320-17154-10

Client ID: OF-FB67-0216

Operator ID: JRB

ALS Bottle#: 18

Worklist Smp#: 63

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

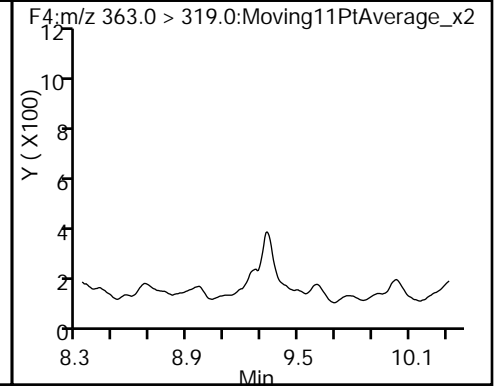
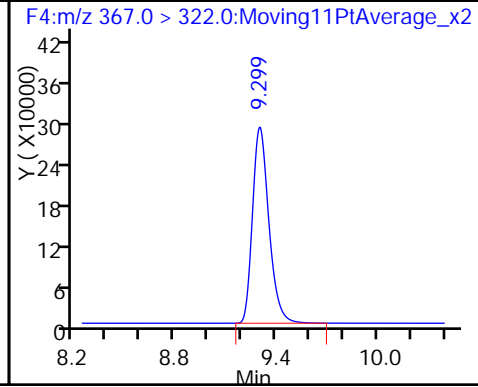
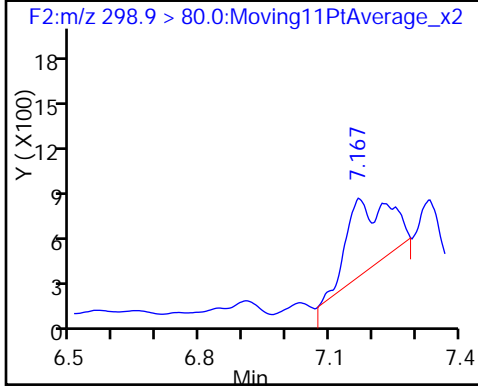
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

40 Perfluorobutanesulfonic acid

D 8 13C4-PFHpa

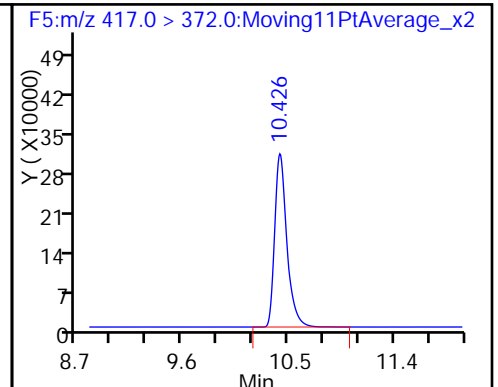
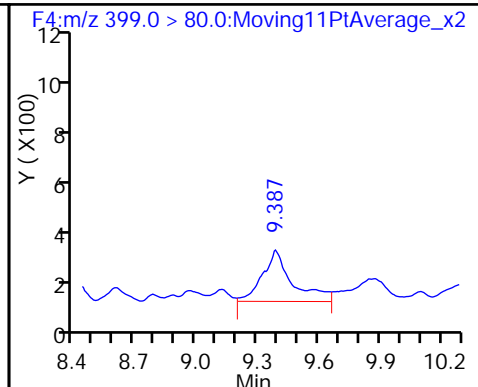
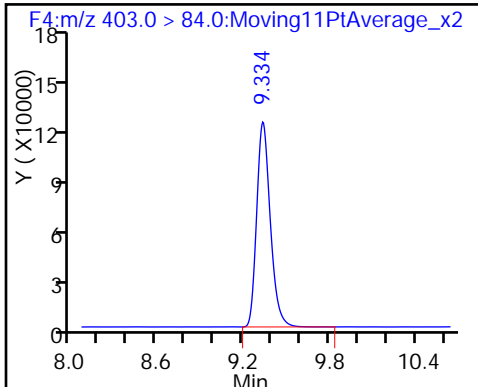
9 Perfluoroheptanoic acid (ND)



D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

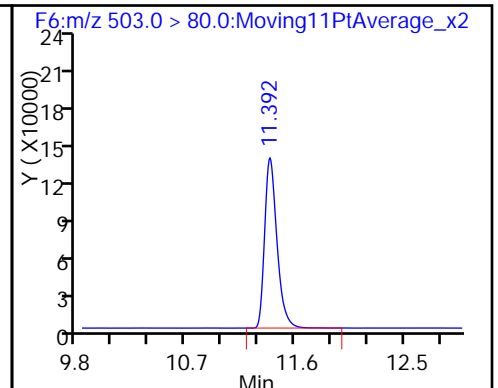
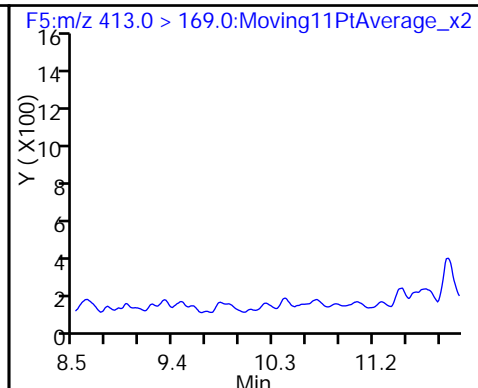
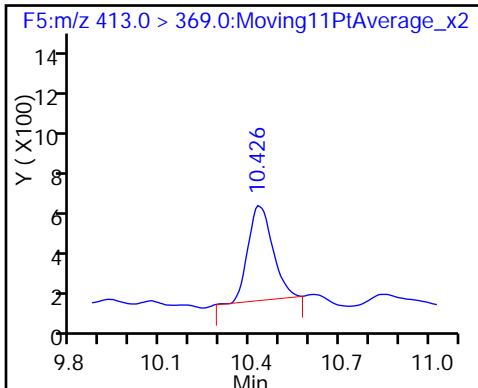
D 12 13C4 PFOA



13 Perfluorooctanoic acid

13 Perfluorooctanoic acid

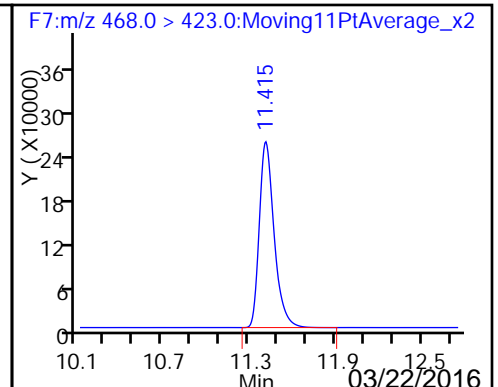
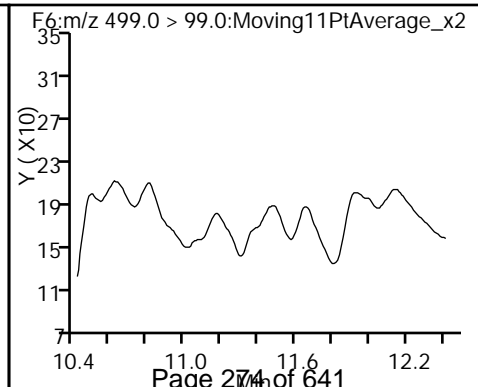
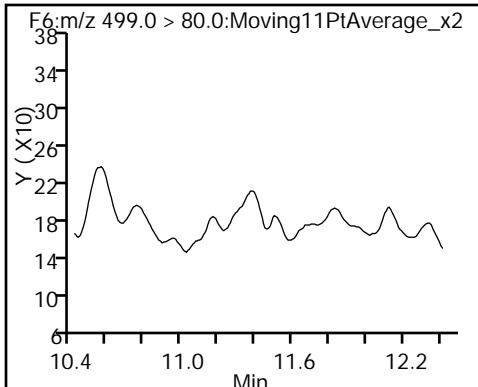
D 16 13C4 PFOS



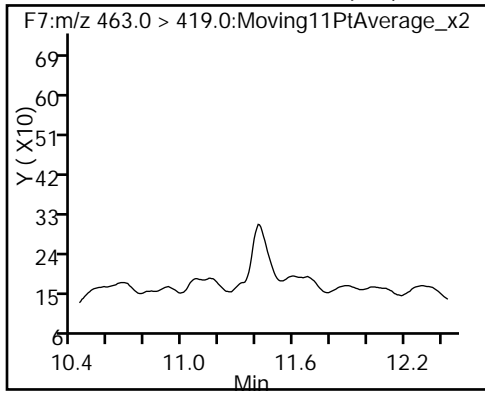
15 Perfluorooctane sulfonic acid (ND)

15 Perfluorooctane sulfonic acid (ND)

D 17 13C5 PFNA



18 Perfluorononanoic acid (ND)



FORM VI  
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA  
RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-17154-1

Analy Batch No.: 100305

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4

GC Column: Acquity

ID: 2.1(mm)

Heated Purge: (Y/N) N

Calibration Start Date: 02/12/2016 10:35

Calibration End Date: 02/12/2016 12:42

Calibration ID: 18968

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100305/2	12FEB2016A4A_003.d
Level 2	STD 320-100305/3	12FEB2016A4A_004.d
Level 3	STD 320-100305/4	12FEB2016A4A_005.d
Level 4	STD 320-100305/5	12FEB2016A4A_006.d
Level 5	STD 320-100305/6	12FEB2016A4A_007.d
Level 6	STD 320-100305/7	12FEB2016A4A_008.d
Level 7	STD 320-100305/8	12FEB2016A4A_009.d

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7				RT WINDOW	AVG RT
Perfluorobutanoic acid (PFBA)	5.653	5.692	5.655	5.753	5.735	5.750	5.747				5.403 - 5.903	5.712
Perfluoropentanoic acid (PFPeA)	6.907	6.943	6.897	7.058	7.015	7.043	7.039				6.657 - 7.157	6.986
Perfluorobutanesulfonic acid (PFBS)	++++	7.074	7.025	7.190	7.144	7.179	7.171				6.769 - 7.269	7.131
Perfluorohexanoic acid (PFHxA)	8.291	8.317	8.327	8.432	8.393	8.409	8.414				8.041 - 8.541	8.369
Perfluoroheptanoic acid (PFHpA)	++++	9.567	9.630	9.692	9.663	9.665	9.672				9.337 - 9.837	9.648
Perfluorohexanesulfonic acid (PFHxS)	++++	9.595	9.665	9.720	9.698	9.707	9.714				9.379 - 9.879	9.683
Perfluorooctanoic acid (PFOA)	10.715	10.670	10.753	10.770	10.760	10.762	10.771				10.465 - 10.965	10.743
Perfluoroheptanesulfonic Acid (PFHpS)	++++	10.661	10.753	10.770	10.760	10.762	10.771				10.455 - 10.955	10.746
Perfluorooctanesulfonic acid (PFOS)	++++	11.578	11.678	11.677	11.676	11.670	11.678				11.385 - 11.885	11.660
Perfluorononanoic acid (PFNA)	11.652	11.604	11.704	11.703	11.694	11.696	11.704				11.402 - 11.902	11.680
Perfluorodecanoic acid (PFDA)	12.445	12.404	12.508	12.496	12.493	12.487	12.498				12.195 - 12.695	12.476
Perfluorooctane Sulfonamide (FOSA)	12.963	12.912	13.021	13.013	13.010	13.004	13.012				12.713 - 13.213	12.991
Perfluorodecane Sulfonic acid	++++	13.040	13.128	13.111	13.116	13.102	++++				12.826 - 13.326	13.099
Perfluoroundecanoic acid (PFUnA)	++++	13.075	13.181	13.163	13.169	13.154	13.163				12.870 - 13.370	13.151
Perfluorododecanoic acid (PFDoA)	13.694	13.655	13.749	13.730	13.727	13.721	13.728				13.444 - 13.944	13.715
Perfluorotetradecanoic Acid (PFTriA)	14.174	14.135	14.220	14.205	14.210	14.197	14.211				13.924 - 14.424	14.193
Perfluorotetradecanoic acid (PFTeA)	++++	14.555	14.635	14.615	14.613	14.609	14.614				14.339 - 14.839	14.607
Perfluoro-n-hexadecanoic acid (PFHxDA)	++++	15.103	15.159	15.144	15.146	15.144	15.143				14.884 - 15.384	15.140
Perfluoro-n-octadecanoic acid (PFODA)	++++	15.414	15.461	15.446	15.448	15.441	15.444				15.191 - 15.691	15.442
13C4 PFBA	5.650	5.695	5.652	5.757	5.735	5.747	5.744				5.400 - 5.900	5.711
13C5-PFPeA	6.893	6.953	6.902	7.058	7.018	7.043	7.039				6.643 - 7.143	6.987
13C2 PFHxA	8.285	8.311	8.327	8.432	8.393	8.409	8.414				8.035 - 8.535	8.367
13C4-PFHpA	9.587	9.567	9.630	9.685	9.663	9.665	9.672				9.337 - 9.837	9.638
18O2 PFHxS	9.629	9.595	9.665	9.720	9.698	9.707	9.714				9.379 - 9.879	9.675
13C4 PFOA	10.705	10.661	10.762	10.779	10.760	10.762	10.771				10.455 - 10.955	10.743
13C4 PFOS	11.626	11.578	11.678	11.677	11.668	11.661	11.670				11.376 - 11.876	11.651
13C5 PFNA	11.652	11.604	11.704	11.703	11.694	11.687	11.704				11.402 - 11.902	11.678
13C2 PFDA	12.445	12.394	12.508	12.486	12.493	12.487	12.498				12.195 - 12.695	12.473
13C8 FOSA	12.963	12.912	13.021	13.013	13.010	13.004	13.012				12.713 - 13.213	12.991
13C2 PFUnA	13.120	13.075	13.181	13.163	13.169	13.154	13.163				12.870 - 13.370	13.146
13C2 PFDoA	13.694	13.655	13.739	13.730	13.727	13.721	13.728				13.444 - 13.944	13.713
13C2-PFTeDA	14.589	14.549	14.635	14.615	14.613	14.609	14.614				14.339 - 14.839	14.603
13C2-PFHxDA	15.134	15.103	15.159	15.144	15.146	15.139	15.143				14.884 - 15.384	15.138

FORM VI  
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1 Analy Batch No.: 100305

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/12/2016 10:35 Calibration End Date: 02/12/2016 12:42 Calibration ID: 18968

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100305/2	12FEB2016A4A_003.d
Level 2	STD 320-100305/3	12FEB2016A4A_004.d
Level 3	STD 320-100305/4	12FEB2016A4A_005.d
Level 4	STD 320-100305/5	12FEB2016A4A_006.d
Level 5	STD 320-100305/6	12FEB2016A4A_007.d
Level 6	STD 320-100305/7	12FEB2016A4A_008.d
Level 7	STD 320-100305/8	12FEB2016A4A_009.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R <sup>2</sup> OR COD	#	MIN R <sup>2</sup> OR COD
	LVL 1 LVL 5	LVL 2 LVL 6	LVL 3 LVL 7	LVL 4		B	M1	M2								
13C4 PFBA	43724 41504	46913 34216	43598 29068	45047	Ave		40581.3800			16.0		50.0				
13C5-PFPeA	34451 34059	36054 25700	36127 21962	35638	Ave		31998.7971			17.9		50.0				
13C2 PFHxA	42875 39681	45781 30298	44064 25927	45494	Ave		39160.1943			20.2		50.0				
13C4-PFHpA	45778 38237	47959 29360	43311 25753	42221	Ave		38945.7029			21.6		50.0				
18O2 PFHxS	23364 23478	27260 19068	26027 14366	25183	Ave		22677.9734			19.9		50.0				
13C4 PFOA	50863 41293	54619 29554	51061 24910	45734	Ave		42576.1771			26.8		50.0				
13C4 PFOS	6706.5 14179	9113.2 10098	10718 7313.8	12758	Ave		10126.7065			26.8		50.0				
13C5 PFNA	49362 39695	50456 33048	47231 27979	43449	Ave		41602.7543			20.5		50.0				
13C2 PFDA	71267 52112	78190 41272	71755 34061	62533	Ave		58741.3229			28.5		50.0				
13C8 FOSA	83903 96232	95349 77442	103205 64546	104593	Ave		89324.3886			16.4		50.0				
13C2 PFUnA	66421 51422	75174 41670	75316 36091	63423	Ave		58502.3943			26.9		50.0				
13C2 PFDoA	77834 69585	85784 59877	82255 48230	74292	Ave		71122.3257			18.6		50.0				
13C2-PFTeDA	60474 65785	71408 55680	71741 46948	71576	Ave		63373.0086			15.0		50.0				
13C2-PFHxDa	18269 36603	26650 30116	31998 27510	36475	Ave		29660.1057			21.5		50.0				

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

## CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-17154-1Analy Batch No.: 100305SDG No.: CTO WE7G PFC SamplingInstrument ID: A4GC Column: Acquity ID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 02/12/2016 10:35Calibration End Date: 02/12/2016 12:42Calibration ID: 18968

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R <sup>2</sup> OR COD	#	MIN R <sup>2</sup> OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7															
Perfluorobutanoic acid (PFBA)	21468 17932	17082 14917	23371	22403	20967	AveID		0.4901			11.8		35.0				
Perfluoropentanoic acid (PFPeA)	14714 12496	15934 10441	18319	17191	16202	AveID		0.4708			5.8		35.0				
Perfluorobutanesulfonic acid (PFBS)	++++ 12952	293560 9193.1	68810	28987	21233	L1ID	9.0196	0.6255						0.9990		0.9900	
Perfluorohexanoic acid (PFHxA)	18078 14787	16971 11751	17710	19721	17723	AveID		0.4308			8.8		35.0				
Perfluoroheptanoic acid (PFHpA)	++++ 15469	27094 12225	21661	21923	18594	AveID		0.5120			6.3		35.0				
Perfluorohexanesulfonic acid (PFHxS)	++++ 20390	40225 16108	33916	28150	25209	L2ID	0.3646	1.1114						0.9970		0.9900	
Perfluorooctanoic acid (PFOA)	25262 15958	27408 11088	26323	21407	19390	AveID		0.4910			6.6		35.0				
Perfluoroheptanesulfonic Acid (PFHpS)	++++ 20276	29776 15686	35251	33157	27743	AveID		2.5441			24.1		50.0				
Perfluorooctanesulfonic acid (PFOS)	++++ 35089	48115 28814	56672	52312	49745	AveID		4.2651			19.3		35.0				
Perfluorononanoic acid (PFNA)	50318 34992	35930 29202	46019	49227	41117	L1ID	-0.115	1.0504						1.0000		0.9900	
Perfluorodecanoic acid (PFDA)	70648 39007	63644 30089	65736	56732	49103	AveID		0.9142			6.1		35.0				
Perfluorooctane Sulfonamide (FOSA)	76012 79257	77919 63244	99217	99139	94248	AveID		0.9450			7.1		35.0				
Perfluorodecane Sulfonic acid	++++ 15557	20133 ++++	26070	24253	21983	AveID		1.9267			20.5		50.0				
Perfluoroundecanoic acid (PFUnA)	++++ 44481	76032 36825	77350	65907	54388	AveID		1.0372			2.1		35.0				
Perfluorododecanoic acid (PFDoA)	60746 50354	55500 40018	69542	62698	59128	AveID		0.8053			9.2		35.0				
Perfluorotridecanoic Acid (PFTriA)	67684 44810	64724 37595	75645	70540	57551	L2ID	0.0026	0.8342						0.9890	*	0.9900	
Perfluorotetradecanoic acid (PFTeA)	++++ 22265	40494 19239	30339	28687	26823	L2ID	0.0885	0.3784						0.9980		0.9900	
Perfluoro-n-hexadecanoic acid (PFHxDA)	++++ 61764	118084 53490	85190	80025	72186	L1ID	0.1544	1.0783						0.9990		0.9900	
Perfluoro-n-octadecanoic acid (PFODA)	++++ 57356	51667 50328	68278	73440	66807	L1ID	-0.544	1.0130						0.9980		0.9900	

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.: 320-17154-1 Analy Batch No.: 100305

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4 GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/12/2016 10:35 Calibration End Date: 02/12/2016 12:42 Calibration ID: 18968

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100305/2	12FEB2016A4A_003.d
Level 2	STD 320-100305/3	12FEB2016A4A_004.d
Level 3	STD 320-100305/4	12FEB2016A4A_005.d
Level 4	STD 320-100305/5	12FEB2016A4A_006.d
Level 5	STD 320-100305/6	12FEB2016A4A_007.d
Level 6	STD 320-100305/7	12FEB2016A4A_008.d
Level 7	STD 320-100305/8	12FEB2016A4A_009.d

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
13C4 PFBA	Ave	2186199 1710809	2345669 1453408	2179878	2252325	2075195	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C5-PFPeA	Ave	1722572 1285003	1802713 1098104	1806356	1781896	1702935	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFHxA	Ave	2143754 1514915	2289067 1296336	2203221	2274718	1984057	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C4-PFHpA	Ave	2288897 1467997	2397972 1287645	2165567	2111070	1911848	50.0 50.0	50.0 50.0	50.0	50.0	50.0
1802 PFHxS	Ave	1105120 901928	1289379 679535	1231093	1191134	1110488	47.3 47.3	47.3 47.3	47.3	47.3	47.3
13C4 PFOA	Ave	2543127 1477697	2730942 1245478	2553061	2286718	2064639	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C4 PFOS	Ave	320573 482668	435612 349599	512338	609839	677767	47.8 47.8	47.8 47.8	47.8	47.8	47.8
13C5 PFNA	Ave	2468093 1652409	2522779 1398930	2361526	2172455	1984772	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFDA	Ave	3563348 2063584	3909475 1703065	3587771	3126636	2605584	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C8 FOSA	Ave	4195173 3872124	4767443 3227314	5160232	5229659	4811591	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFUnA	Ave	3321059 2083502	3758689 1804563	3765798	3171146	2571081	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFDoA	Ave	3891686 2993831	4289188 2411510	4112735	3714597	3479267	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2-PFTeDA	Ave	3023681 2783983	3570399 2347410	3587033	3578785	3289262	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2-PFHxDA	Ave	913455 1505780	1332475 1375507	1599911	1823759	1830150	50.0 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.: 320-17154-1Analy Batch No.: 100305SDG No.: CTO WE7G PFC SamplingInstrument ID: A4GC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 02/12/2016 10:35Calibration End Date: 02/12/2016 12:42Calibration ID: 18968

## Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100305/2	12FEB2016A4A_003.d
Level 2	STD 320-100305/3	12FEB2016A4A_004.d
Level 3	STD 320-100305/4	12FEB2016A4A_005.d
Level 4	STD 320-100305/5	12FEB2016A4A_006.d
Level 5	STD 320-100305/6	12FEB2016A4A_007.d
Level 6	STD 320-100305/7	12FEB2016A4A_008.d
Level 7	STD 320-100305/8	12FEB2016A4A_009.d

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
Perfluorobutanoic acid (PFBA)		AveID	10734 3586362	17082 5966668	116853	448061	1048362	0.500 200	1.00 400	5.00	20.0	50.0
Perfluoropentanoic acid (PFPeA)		AveID	7357 2499117	15934 4176286	91597	343823	810081	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorobutanesulfonic acid (PFBS)		L1ID	++++ 2289993	259507 3250684	304142	512492	938494	++++ 177	0.884 354	4.42	17.7	44.2
Perfluorohexanoic acid (PFHxA)		AveID	9039 2957321	16971 4700473	88552	394416	886159	0.500 200	1.00 400	5.00	20.0	50.0
Perfluoroheptanoic acid (PFHpA)		AveID	++++ 3093796	27094 4889892	108307	438459	929700	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorohexanesulfonic acid (PFHxS)		L2ID	++++ 3857794	38053 6095086	160421	532601	1192387	++++ 189	0.946 378	4.73	18.9	47.3
Perfluorooctanoic acid (PFOA)		AveID	12631 3191541	27408 4435164	131613	428131	969508	0.500 200	1.00 400	5.00	20.0	50.0
Perfluoroheptanesulfonic Acid (PFHpS)		AveID	++++ 3860458	28347 5973185	167796	631312	1320551	++++ 190	0.952 381	4.76	19.0	47.6
Perfluorooctanesulfonic acid (PFOS)		AveID	++++ 6709046	45998 11018460	270891	1000210	2377814	++++ 191	0.956 382	4.78	19.1	47.8
Perfluorononanoic acid (PFNA)		L1ID	25159 6998423	35930 11680686	230097	984533	2055860	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorodecanoic acid (PFDA)		AveID	35324 7801447	63644 12035423	328682	1134637	2455156	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorooctane Sulfonamide (FOSA)		AveID	38006 15851344	77919 25297572	496084	1982787	4712375	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorodecane Sulfonic acid		AveID	++++ 2999356	19408 ++++	125658	467598	1059583	++++ 193	0.964 ++++	4.82	19.3	48.2
Perfluoroundecanoic acid (PFUnA)		AveID	++++ 8896158	76032 14730079	386752	1318141	2719401	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorododecanoic acid (PFDoA)		AveID	30373 10070858	55500 16007183	347709	1253957	2956390	0.500 200	1.00 400	5.00	20.0	50.0



## RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1 Analy Batch No.: 100305SDG No.: CTO WE7G PFC SamplingInstrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) NCalibration Start Date: 02/12/2016 10:35 Calibration End Date: 02/12/2016 12:42 Calibration ID: 18968

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
Perfluorotridecanoic Acid (PFTriA)		L2ID	33842 8962091	64724 15038082	378224	1410790	2877547	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorotetradecanoic acid (PFTeA)		L2ID	++++ 4452926	40494 7695775	151695	573742	1341162	++++ 200	1.00 400	5.00	20.0	50.0
Perfluoro-n-hexadecanoic acid (PFHxDA)		L1ID	++++ 12352725	118084 21395824	425951	1600494	3609294	++++ 200	1.00 400	5.00	20.0	50.0
Perfluoro-n-octadecanoic acid (PFODA)		L1ID	++++ 11471145	51667 20131167	341392	1468803	3340365	++++ 200	1.00 400	5.00	20.0	50.0

## Curve Type Legend:

AveID = Average isotope dilution
L1ID = Linear 1/conc IsoDil
L2ID = Linear 1/conc^2 IsoDil

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_003.d  
 Lims ID: Std L1  
 Client ID:  
 Sample Type: IC Calib Level: 1  
 Inject. Date: 12-Feb-2016 10:35:33 ALS Bottle#: 2 Worklist Smp#: 2  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L1  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:02:02 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 14:29:46

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
--------	----	--------	--------	--------	----------	--------------	---------------	------	-----	-------

D 1 13C4 PFBA	216.7 > 171.5	5.650	5.650	0.0	2186199	53.9		108	6396	
2 Perfluorobutyric acid	212.7 > 168.6	5.653	5.653	0.0	10734	0.5009		100	50.1	
D 3 13C5-PFPeA	267.6 > 222.7	6.893	6.893	0.0	1722572	53.8		108	2771	
4 Perfluoropentanoic acid	262.9 > 218.7	6.907	6.907	0.0	7357	0.4536		90.7	7.2	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.019	7.019	0.0	222434	NC			147	
	298.8 > 98.6	7.016	7.019	-0.003	135664		1.64(0.00-0.00)		111	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.019	7.019	0.0	222434	0.8005		181		
D 6 13C2 PFHxA	314.6 > 269.7	8.285	8.285	0.0	2143754	54.7		109	7491	
7 Perfluorohexanoic acid	312.9 > 268.7	8.291	8.291	0.0	9039	0.4894		97.9	41.0	
D 8 13C4-PFHpA	366.6 > 321.6	9.587	9.587	0.0	2288897	58.8		118	5094	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.587	9.587	0.0	22979	0.9804		196	70.4	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.629	9.629	0.0	23477	0.5761		122		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.629	9.629	0.0	23477	NC			99.2	
D 11 18O2 PFHxS	402.5 > 83.6	9.629	9.629	0.0	1105120	48.7		103	2974	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.705	10.705	0.0	1.000	17930	1.05	221		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.705	10.705	0.0	1.000	17930	NC		110	
D 12 13C4 PFOA	416.5 > 371.6	10.705	10.705	0.0		2543127	59.7	119	4035	
13 Perfluorooctanoic acid	412.8 > 368.8	10.715	10.715	0.0	1.000	12631	0.5058	101	36.0	
D 16 13C4 PFOS	502.4 > 79.7	11.626	11.626	0.0		320573	31.7	66.2	1287	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.635	11.635	0.0	1.000	32518	1.14	238	55.9	
D 17 13C5 PFNA	467.5 > 422.6	11.652	11.652	0.0		2468093	59.3	119	4243	
18 Perfluorononanoic acid	462.5 > 418.6	11.652	11.652	0.0	1.000	25159	0.5943	119	68.9	
D 19 13C2 PFDA	514.4 > 469.5	12.445	12.445	0.0		3563348	60.7	121	4269	
20 Perfluorodecanoic acid	512.5 > 468.5	12.445	12.445	0.0	1.000	35324	0.5422	108	88.0	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	12.963	12.963	0.0	1.000	38006	0.4793	95.9	123	
D 23 13C8 FOSA	505.4 > 77.6	12.963	12.963	0.0		4195173	47.0	93.9	2466	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.076	13.076	0.0	1.000	9585	NC		49.5	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.076	13.076	0.0	1.000	9585	0.7418	154		
D 26 13C2 PFUnA	564.3 > 519.5	13.120	13.120	0.0		3321059	56.8	114	5418	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.120	13.120	0.0	1.000	38320	0.5562	111	74.8	
D 28 13C2 PFDaA	614.4 > 569.4	13.694	13.694	0.0		3891686	54.7	109	4817	
29 Perfluorododecanoic acid	612.4 > 568.6	13.694	13.694	0.0	1.000	30373	0.4846	96.9	22.7	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.174	14.174	0.0	1.000	33842	0.5181	104	25.8	
D 33 13C2-PFTeDA	714.5 > 669.5	14.589	14.589	0.0		3023681	47.7	95.4	4248	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.589	14.589	0.0	1.000	22501	0.5301	106	21.1	
D 35 13C2-PFHxDA	814.8 > 769.6	15.134	15.134	0.0		913455	30.8	61.6	3550	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.134	15.134	0.0	1.000	89904	0.9280	186	133	
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.441	15.441	0.0	1.000	225283	110.8231	165	3207	

[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L1\_00018

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_003.d

Injection Date: 12-Feb-2016 10:35:33

Instrument ID: A4

Lims ID: Std L1

Client ID:

Operator ID: JRB

ALS Bottle#: 2

Worklist Smp#: 2

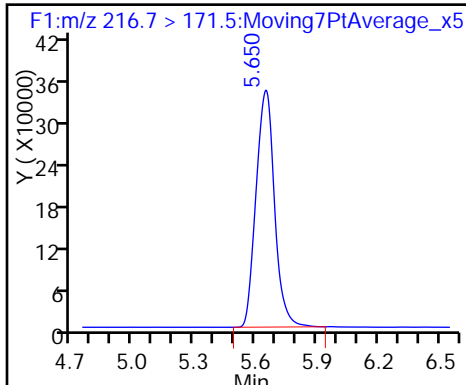
Injection Vol: 15.0 ul

Dil. Factor: 1.0000

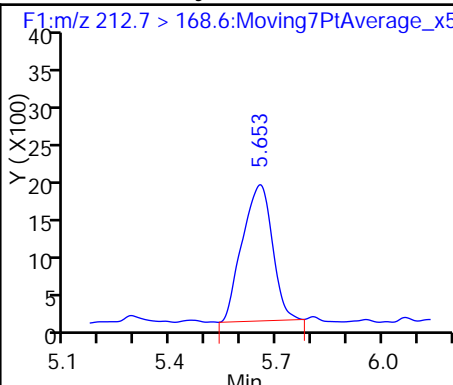
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

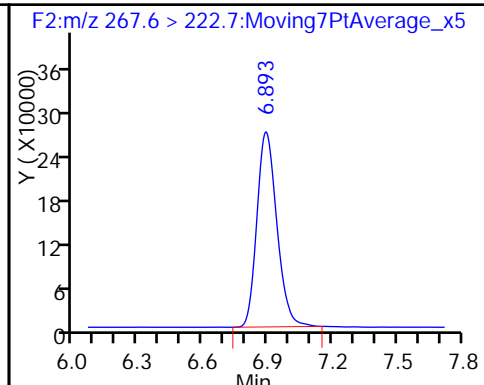
D 1 13C4 PFBA



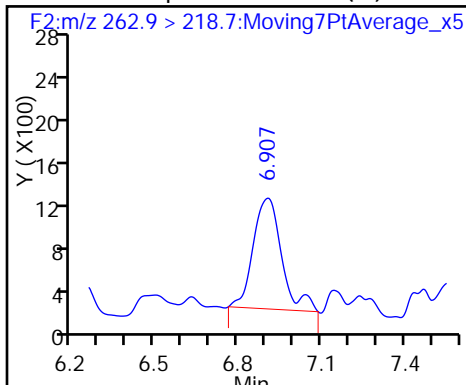
2 Perfluorobutyric acid



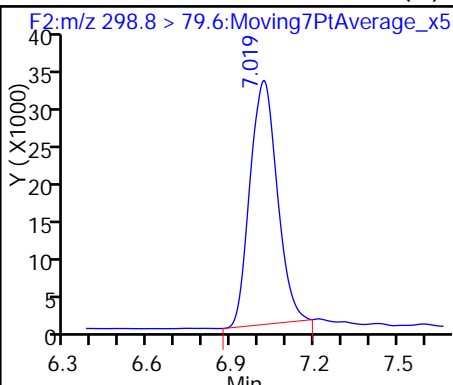
D 3 13C5-PFPeA (M)



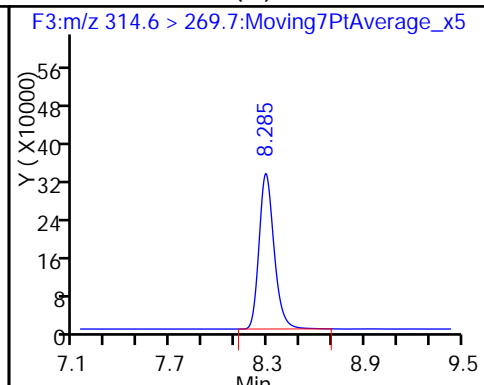
4 Perfluoropentanoic acid (M)



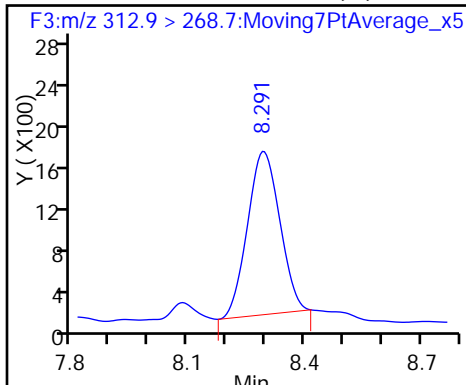
51 Perfluorobutanesulfonic acid (M)



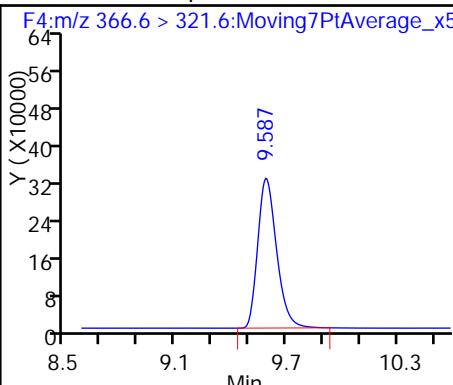
D 6 13C2 PFHxA (M)



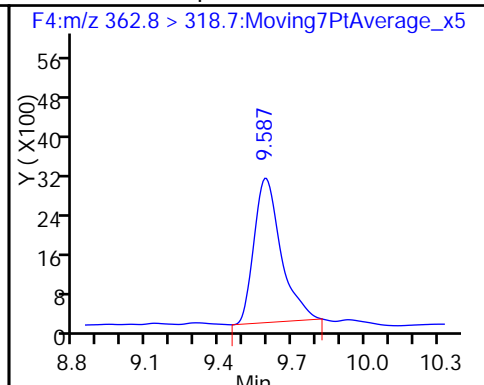
7 Perfluorohexanoic acid (M)



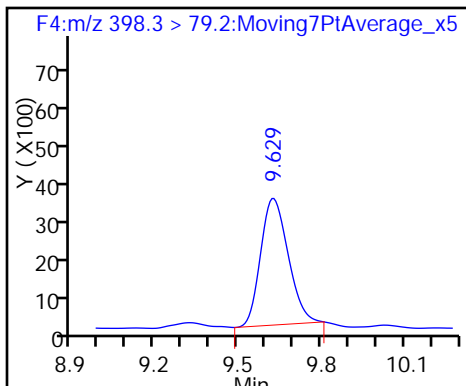
D 8 13C4-PFHpA



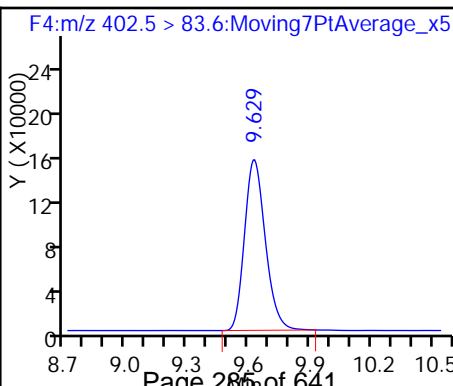
9 Perfluoroheptanoic acid



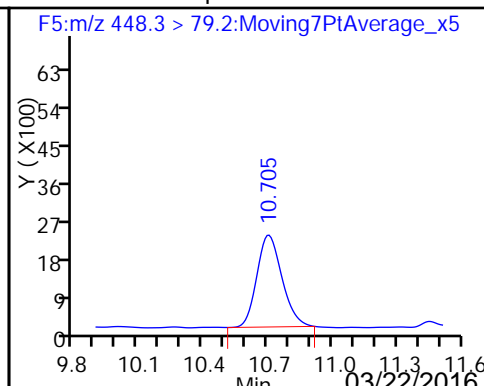
58 Perfluorohexanesulfonic acid



D 11 18O2 PFHxS



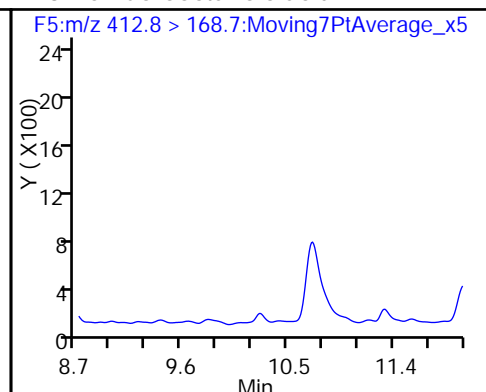
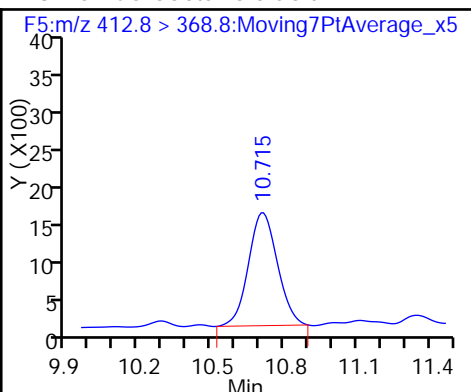
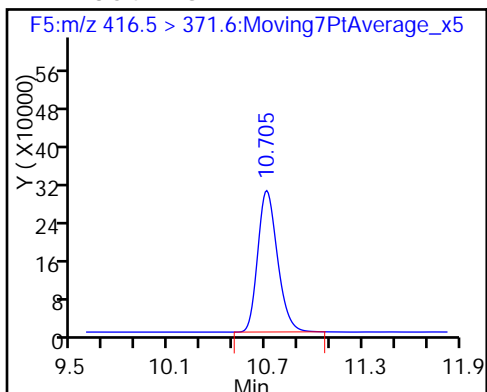
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

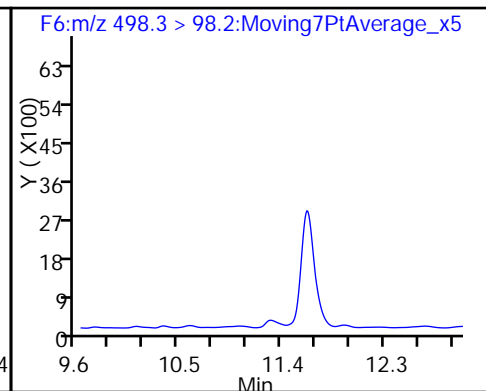
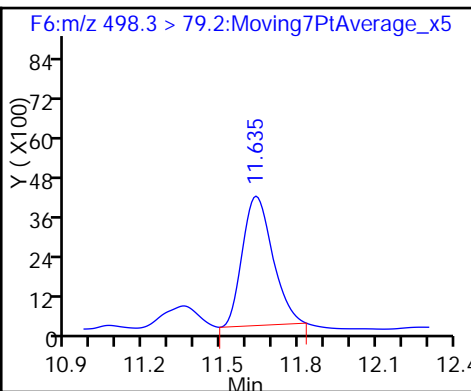
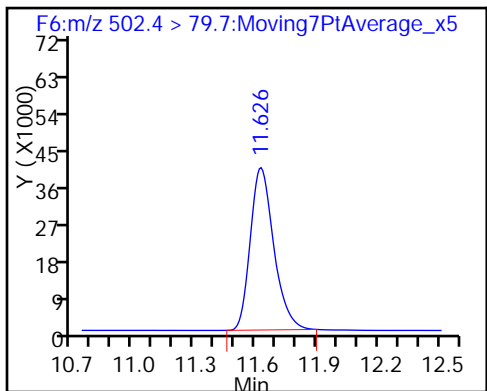
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid

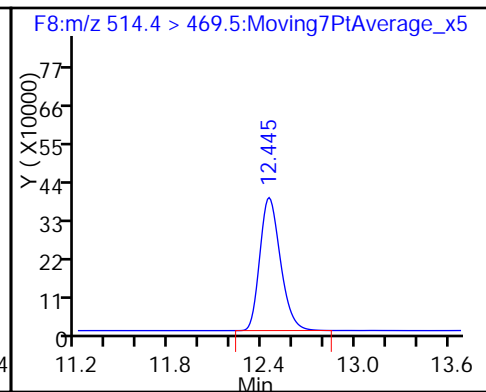
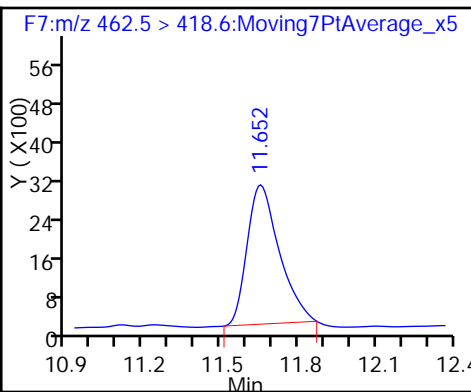
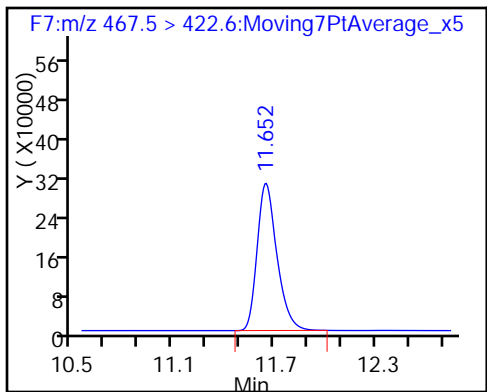
15 Perfluorooctane sulfonic acid



D 17 13C5 PFNA

18 Perfluorononanoic acid

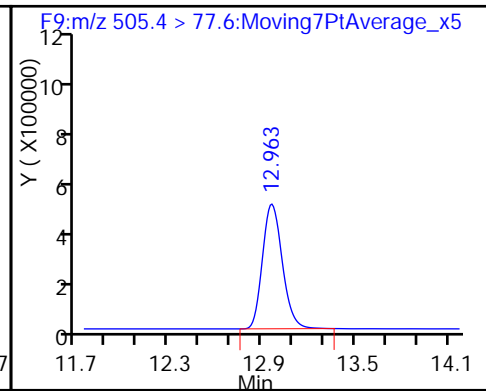
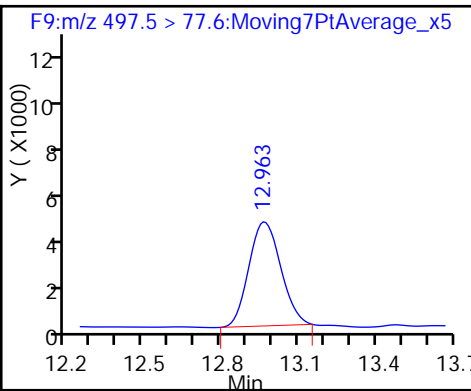
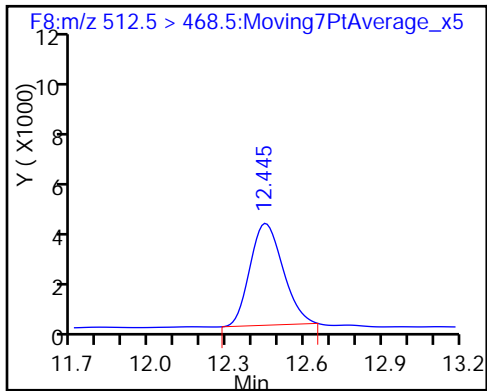
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

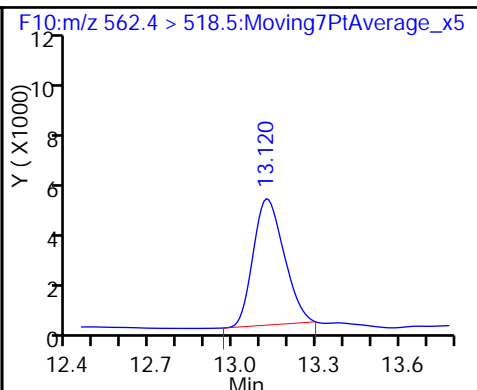
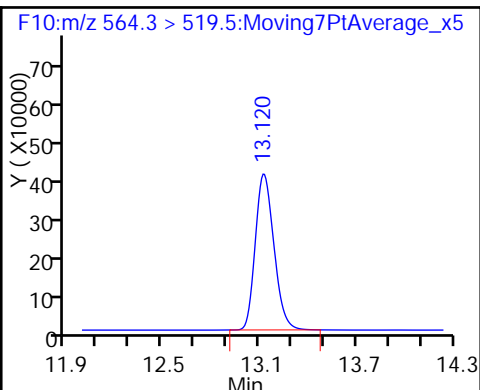
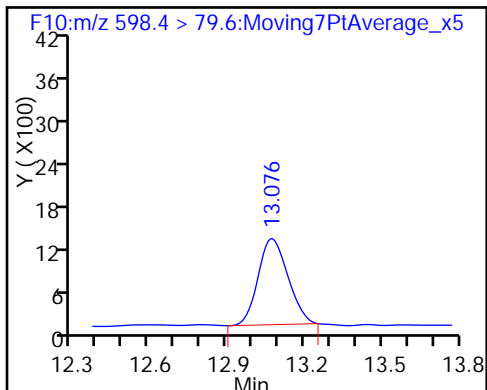
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

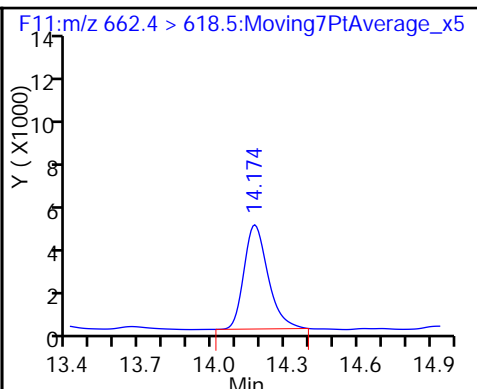
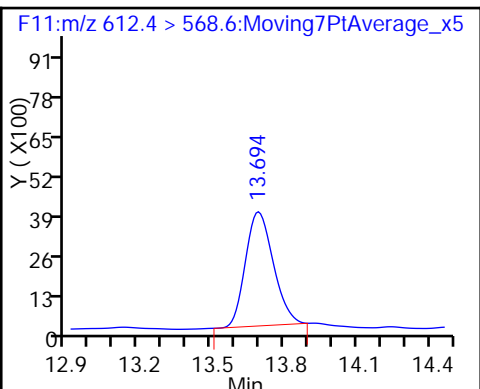
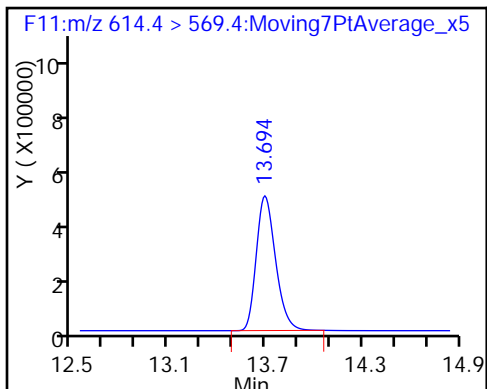
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

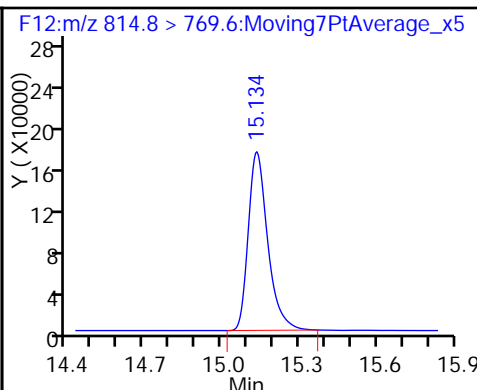
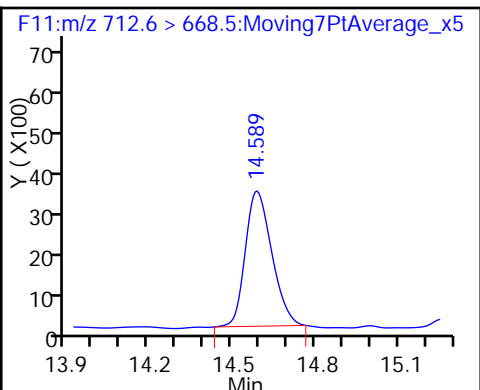
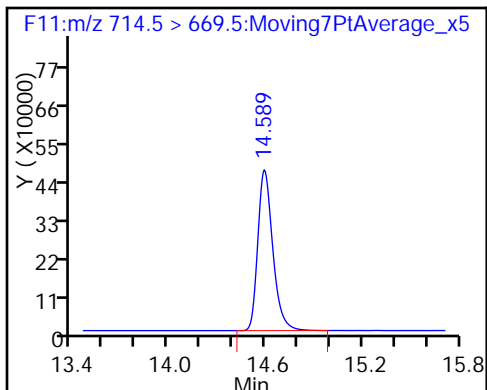
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

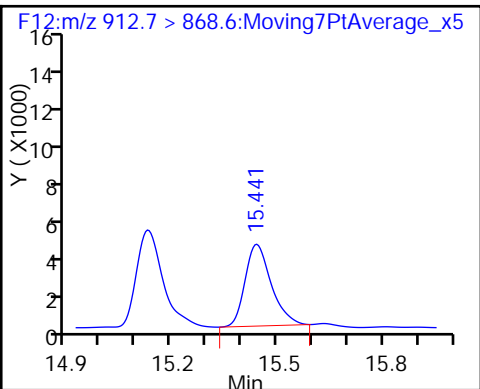
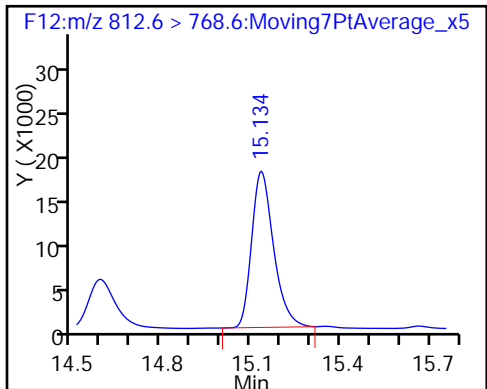
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d  
 Lims ID: Std L2  
 Client ID:  
 Sample Type: IC Calib Level: 2  
 Inject. Date: 12-Feb-2016 10:56:41 ALS Bottle#: 3 Worklist Smp#: 3  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L2  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:02:12 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 14:30:24

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	216.7 > 171.5	5.695	5.650	0.045	2345669	57.8		116	7846	
2 Perfluorobutyric acid	212.7 > 168.6	5.692	5.653	0.039	17082	0.7429		74.3	61.2	M
D 3 13C5-PFPeA	267.6 > 222.7	6.953	6.893	0.060	1802713	56.3		113	4423	
4 Perfluoropentanoic acid	262.9 > 218.7	6.943	6.907	0.036	15934	0.9387		93.9	13.6	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.074	7.019	0.055	259507	NC			152	
	298.8 > 98.6	7.077	7.019	0.058	164915		1.57(0.00-0.00)		141	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.074	7.019	0.055	259507	0.7997		90.5		
D 6 13C2 PFHxA	314.6 > 269.7	8.311	8.285	0.026	2289067	58.5		117	5220	
7 Perfluorohexanoic acid	312.9 > 268.7	8.317	8.291	0.026	16971	0.8605		86.0	101	
D 8 13C4-PFHpA	366.6 > 321.6	9.567	9.587	-0.020	2397972	61.6		123	4912	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.567	9.587	-0.020	27094	1.10		110	78.7	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.595	9.629	-0.034	38053	0.9280		98.1		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.595	9.629	-0.034	38053	NC			152	
D 11 18O2 PFHxS	402.5 > 83.6	9.595	9.629	-0.034	1289379	56.9		120	3861	



Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.661	10.705	-0.044	1.000	28347	1.22	128		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.661	10.705	-0.044	1.000	28347	NC		141	
D 12 13C4 PFOA	416.5 > 371.6	10.661	10.705	-0.044		2730942	64.1	128	4248	
13 Perfluorooctanoic acid	412.8 > 368.8	10.670	10.715	-0.045	1.000	27408	1.02	102	65.6	
	412.8 > 168.7	10.661	10.715	-0.054	0.999	9428	2.91(0.00-0.00)	102	37.6	
D 16 13C4 PFOS	502.4 > 79.7	11.578	11.626	-0.048		435612	43.0	90.0	1752	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.578	11.635	-0.057	1.000	45998	1.18	124	252	M
	498.3 > 98.2	11.578	11.635	-0.057	1.000	28574	1.61(0.00-0.00)	124	92.1	M
D 17 13C5 PFNA	467.5 > 422.6	11.604	11.652	-0.048		2522779	60.6	121	4314	
18 Perfluorononanoic acid	462.5 > 418.6	11.604	11.652	-0.048	1.000	35930	0.7871	78.7	77.9	M
D 19 13C2 PFDA	514.4 > 469.5	12.394	12.445	-0.051		3909475	66.6	133	4885	
20 Perfluorodecanoic acid	512.5 > 468.5	12.404	12.445	-0.041	1.000	63644	0.8904	89.0	147	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	12.912	12.963	-0.051	1.000	77919	0.8648	86.5	257	
D 23 13C8 FOSA	505.4 > 77.6	12.912	12.963	-0.051		4767443	53.4	107	2131	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.040	13.076	-0.036	1.000	19408	NC		118	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.040	13.076	-0.036	1.000	19408	1.11	115		
D 26 13C2 PFUnA	564.3 > 519.5	13.075	13.120	-0.045		3758689	64.2	128	4185	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.075	13.120	-0.045	1.000	76032	0.9752	97.5	138	
D 28 13C2 PFDoA	614.4 > 569.4	13.655	13.694	-0.039		4289188	60.3	121	4771	
29 Perfluorododecanoic acid	612.4 > 568.6	13.655	13.694	-0.039	1.000	55500	0.8034	80.3	34.0	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.135	14.174	-0.039	1.000	64724	0.9013	90.1	36.6	
D 33 13C2-PFTeDA	714.5 > 669.5	14.549	14.589	-0.040		3570399	56.3	113	4342	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.555	14.589	-0.034	1.000	40494	1.01	101	30.6	
D 35 13C2-PFHxDA	814.8 > 769.6	15.103	15.134	-0.031		1332475	44.9	89.8	3351	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.103	15.134	-0.031	1.000	118084	1.13	113	128	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.414	15.441	-0.027	1.000	51667	1.13	113	53.3	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

Review Flags

M - Manually Integrated

**Reagents:**

LCPFC-L2\_00018

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d

Injection Date: 12-Feb-2016 10:56:41

Instrument ID: A4

Lims ID: Std L2

Client ID:

Operator ID: JRB

ALS Bottle#: 3

Worklist Smp#: 3

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

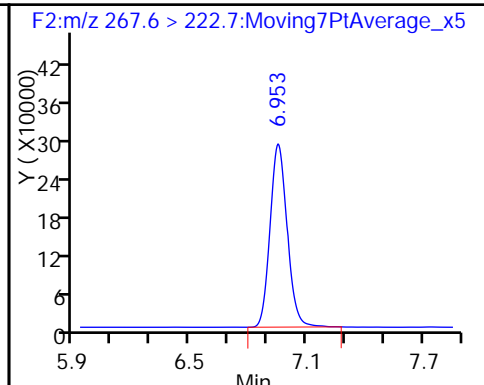
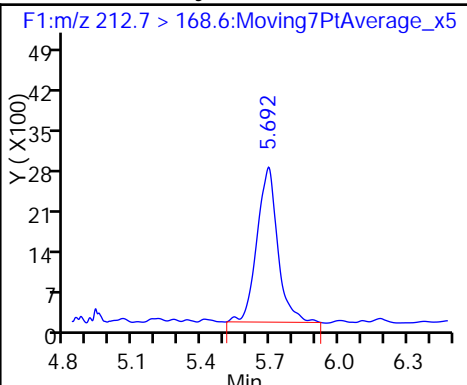
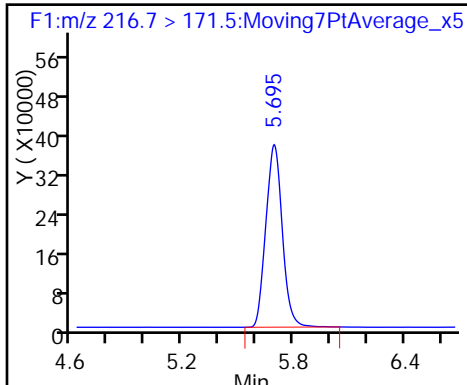
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid (M)

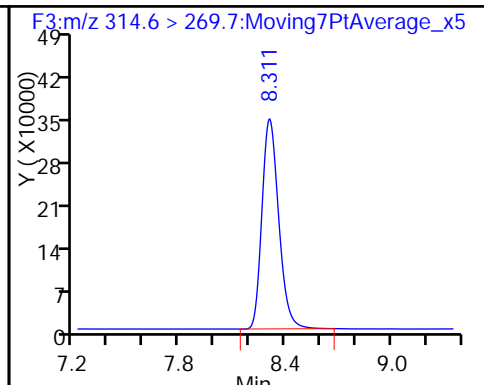
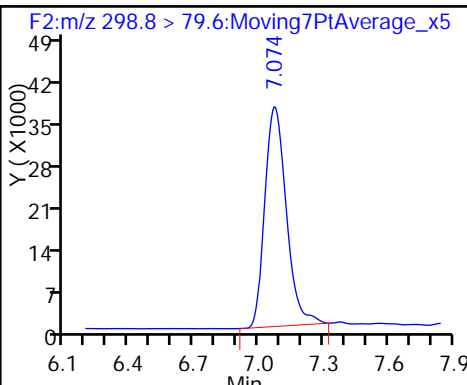
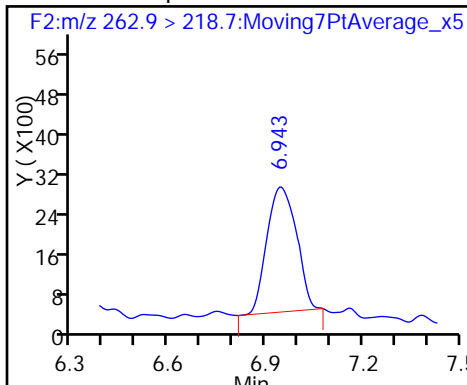
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

51 Perfluorobutanesulfonic acid

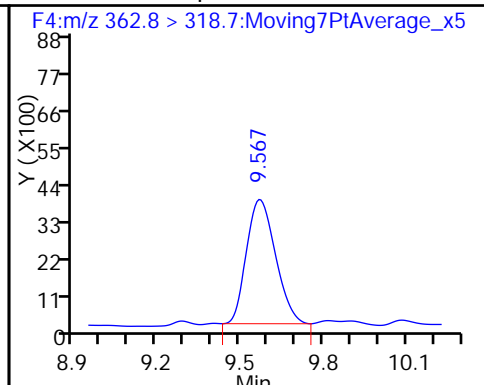
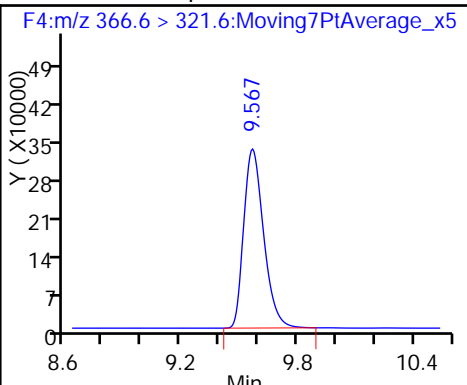
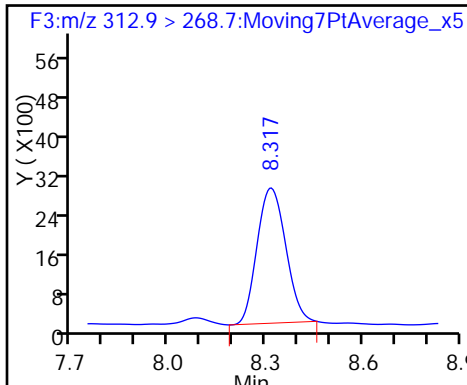
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

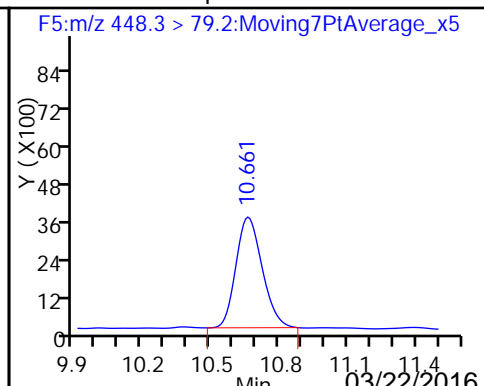
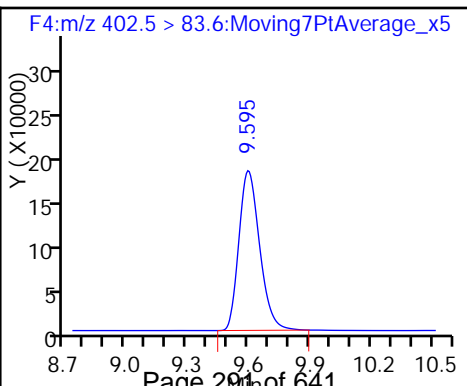
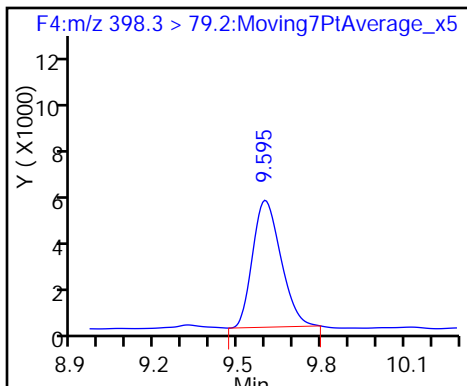
9 Perfluoroheptanoic acid



58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

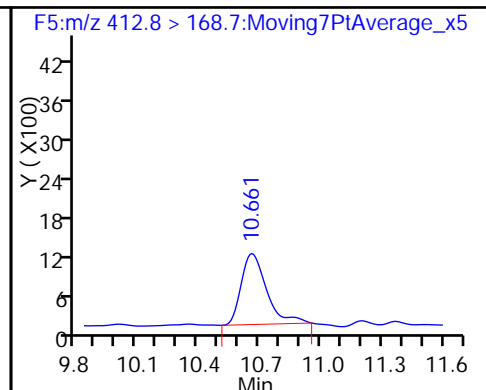
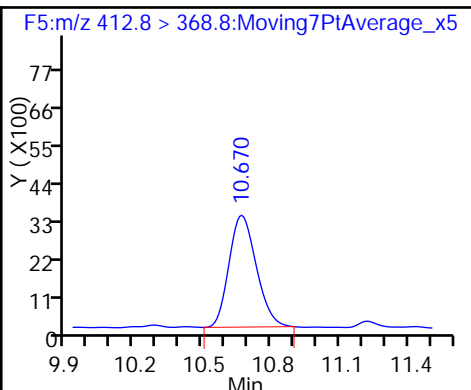
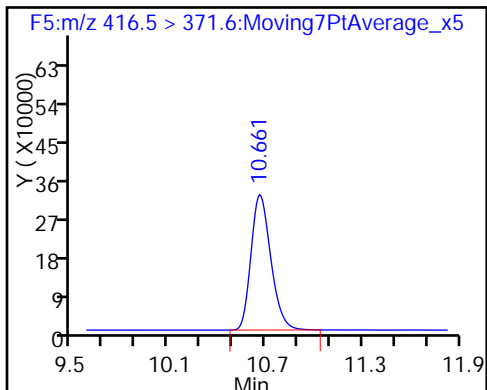
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

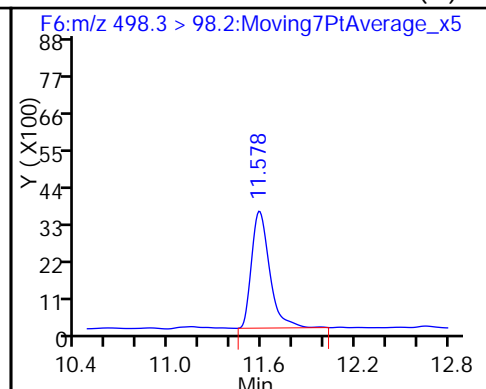
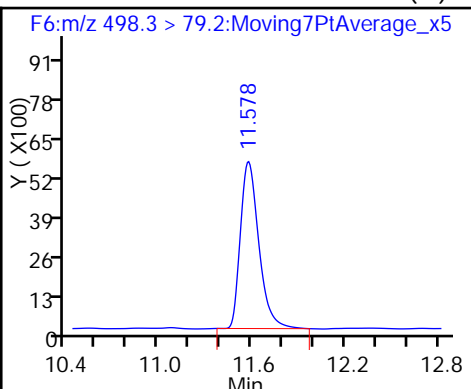
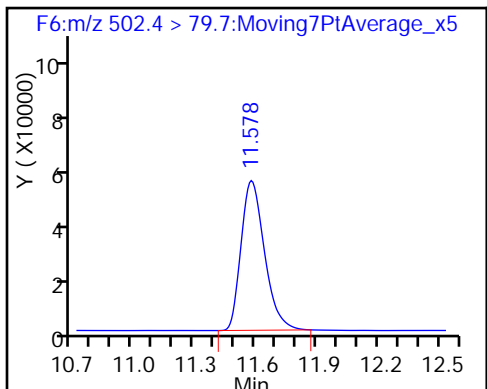
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid (M)

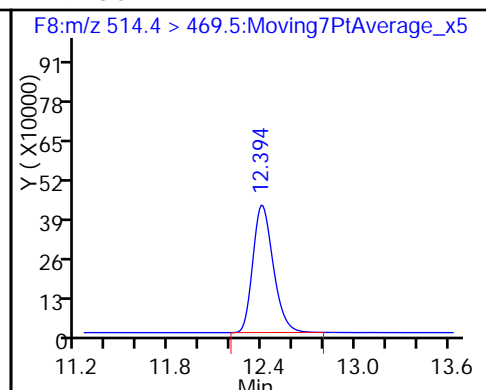
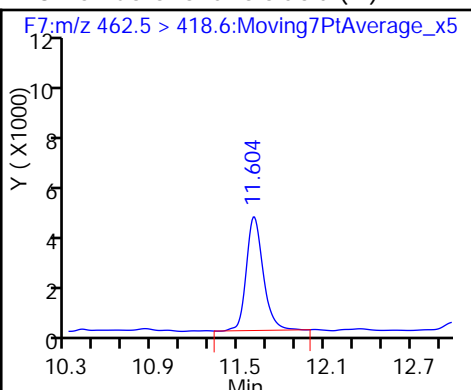
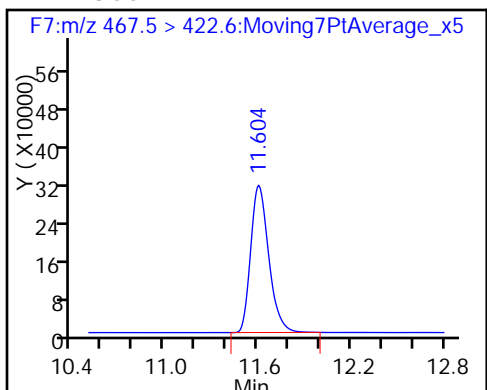
15 Perfluorooctane sulfonic acid (M)



D 17 13C5 PFNA

18 Perfluorononanoic acid (M)

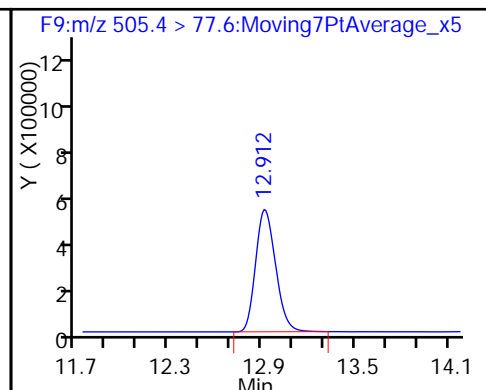
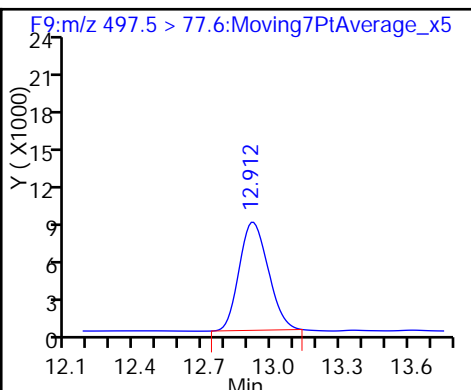
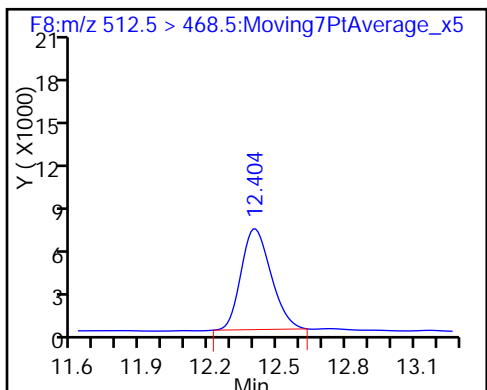
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

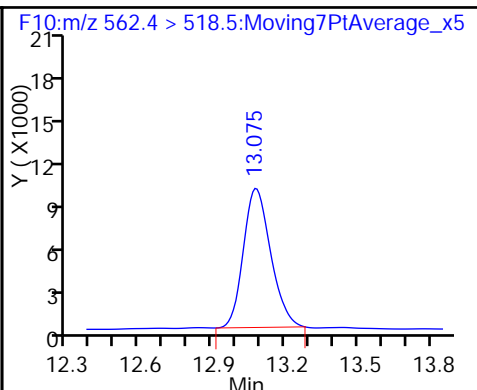
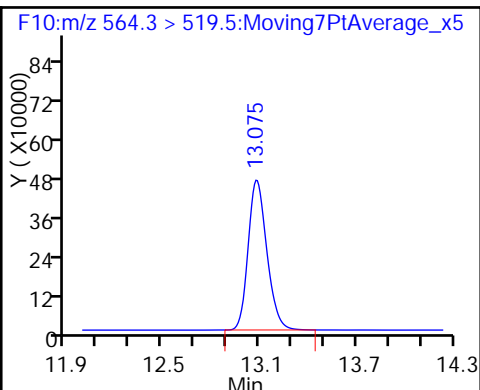
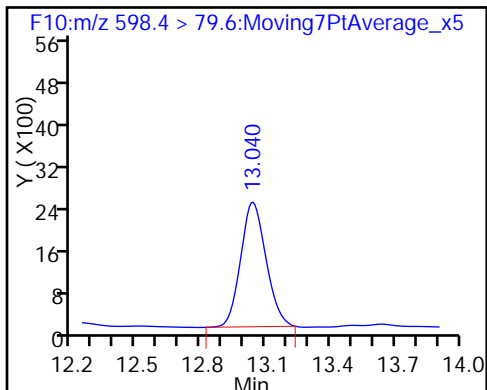
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

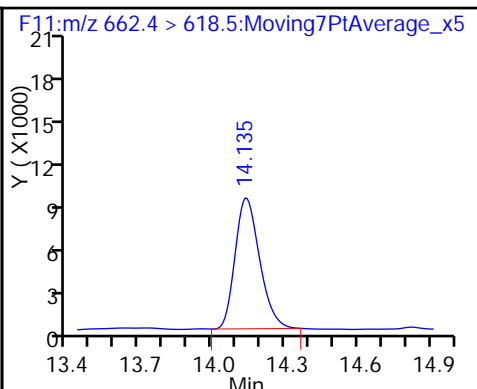
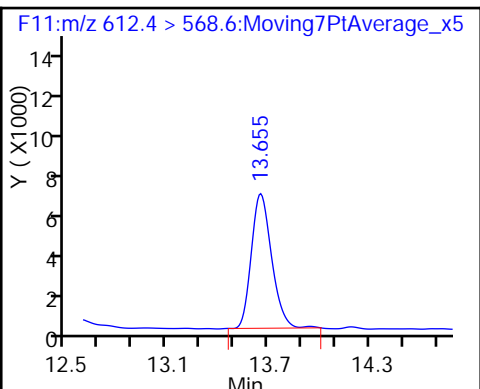
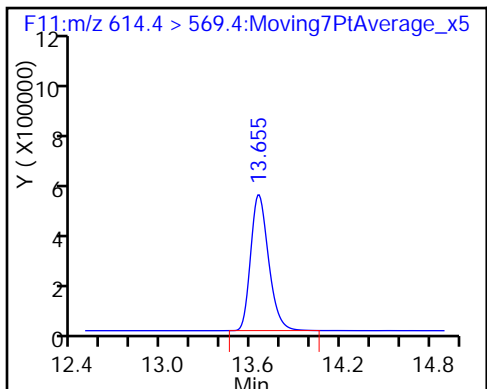
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

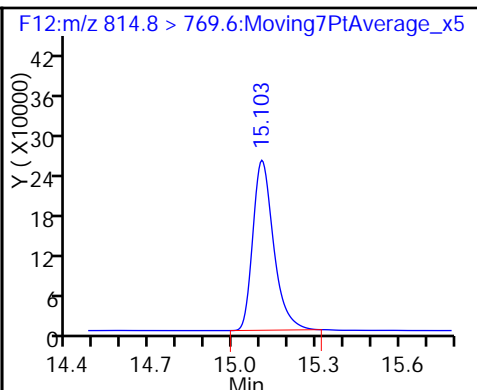
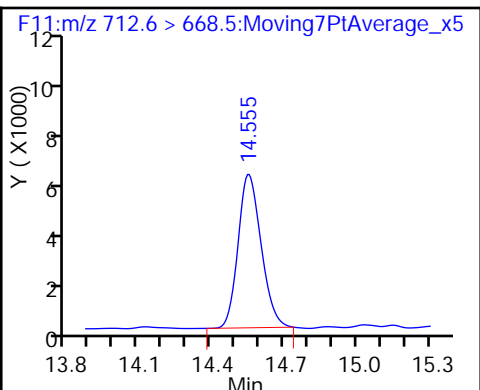
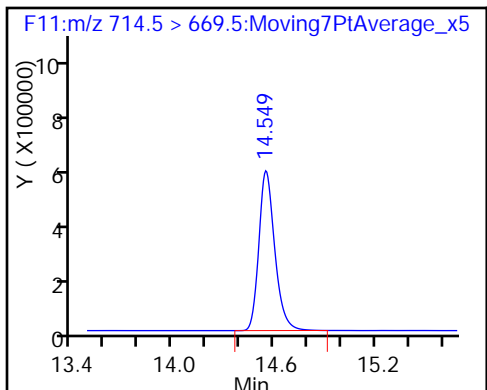
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

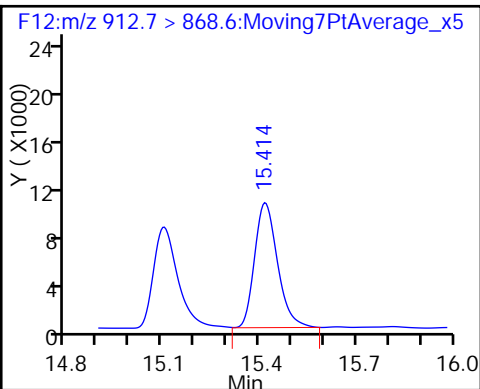
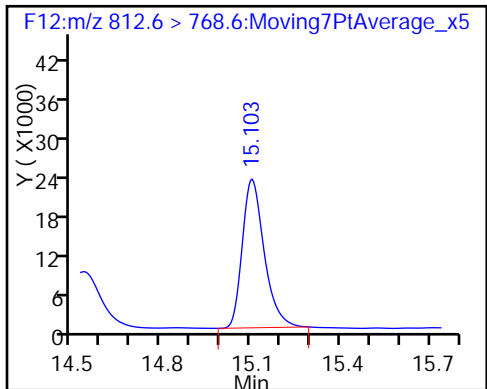
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxD A



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



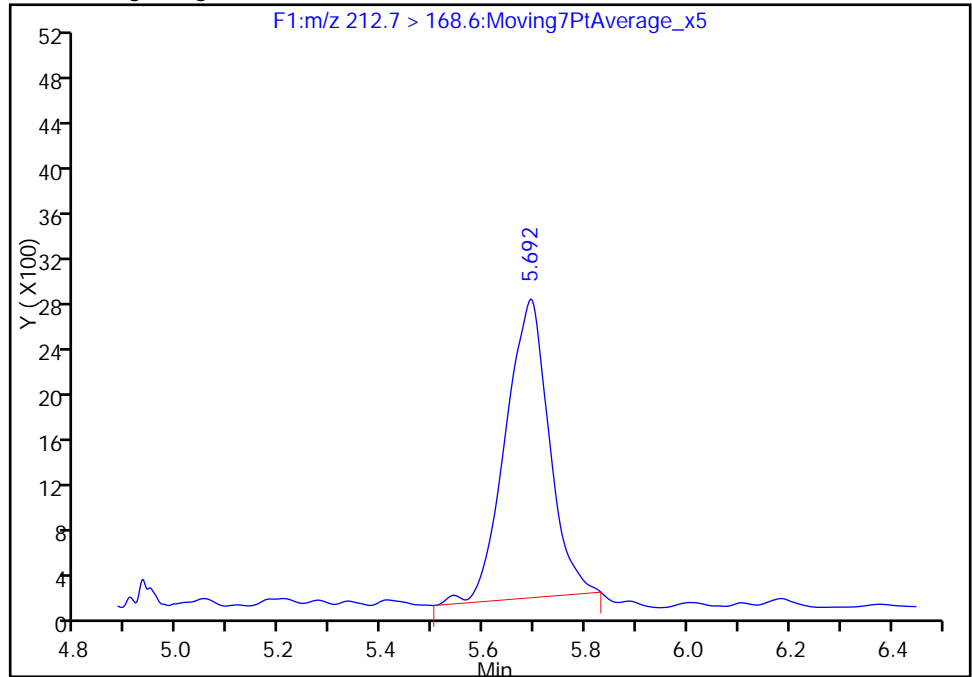
TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d  
Injection Date: 12-Feb-2016 10:56:41 Instrument ID: A4  
Lims ID: Std L2  
Client ID:  
Operator ID: JRB ALS Bottle#: 3 Worklist Smp#: 3  
Injection Vol: 15.0 ul Dil. Factor: 1.0000  
Method: PFAC\_A4 Limit Group: LC PFC\_DOD ICAL  
Column: Detector F1:M/RM

2 Perfluorobutyric acid, CAS: 375-22-4

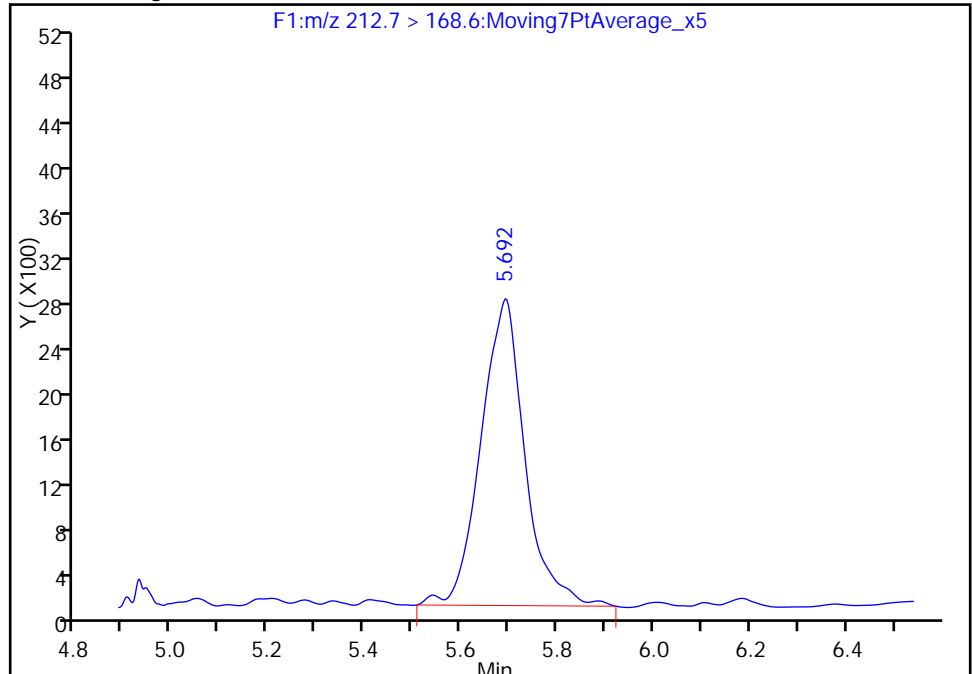
RT: 5.69  
Area: 15671  
Amount: 0.687562  
Amount Units: ng/ml

Processing Integration Results



RT: 5.69  
Area: 17082  
Amount: 0.742900  
Amount Units: ng/ml

Manual Integration Results



Reviewer: westendorfc, 15-Feb-2016 08:14:55  
Audit Action: Manually Integrated  
Audit Reason: Baseline

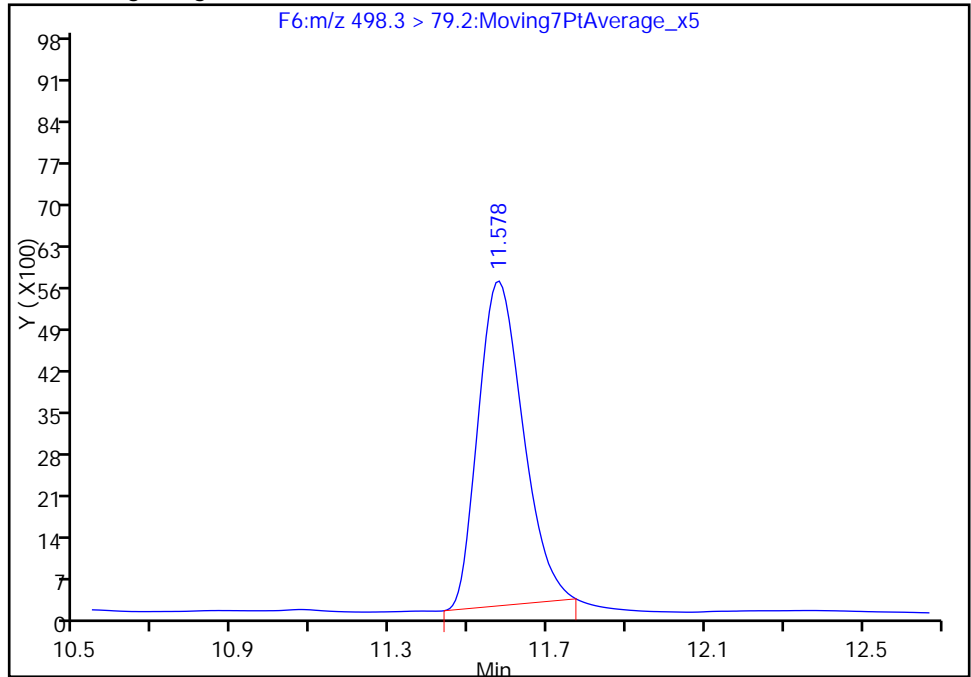
TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d  
Injection Date: 12-Feb-2016 10:56:41 Instrument ID: A4  
Lims ID: Std L2  
Client ID:  
Operator ID: JRB ALS Bottle#: 3 Worklist Smp#: 3  
Injection Vol: 15.0 ul Dil. Factor: 1.0000  
Method: PFAC\_A4 Limit Group: LC PFC\_DOD ICAL  
Column: Detector F6:M/RM

15 Perfluorooctane sulfonic acid, CAS: 1763-23-1

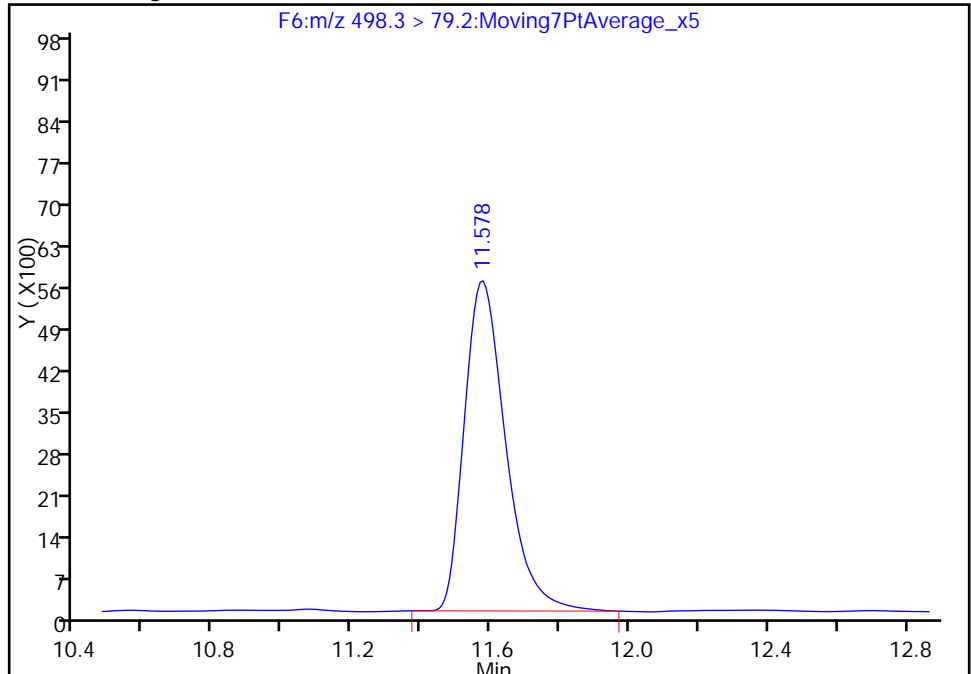
RT: 11.58  
Area: 43067  
Amount: 1.122782  
Amount Units: ng/ml

Processing Integration Results



RT: 11.58  
Area: 45998  
Amount: 1.183430  
Amount Units: ng/ml

Manual Integration Results



Reviewer: westendorfc, 15-Feb-2016 08:20:55  
Audit Action: Manually Integrated  
Audit Reason: Baseline

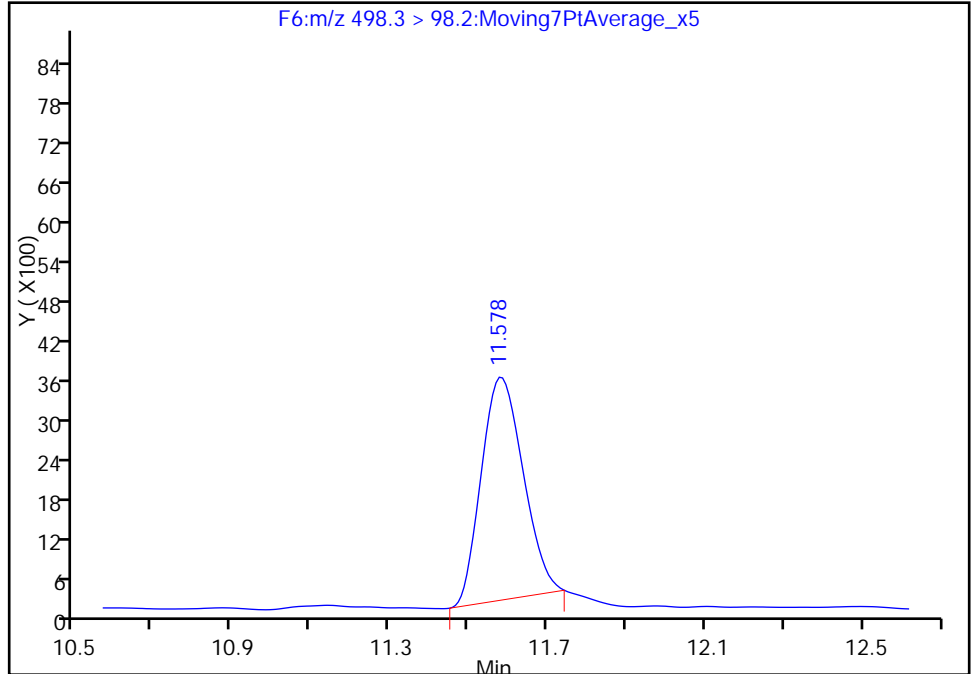
TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d  
Injection Date: 12-Feb-2016 10:56:41 Instrument ID: A4  
Lims ID: Std L2  
Client ID:  
Operator ID: JRB ALS Bottle#: 3 Worklist Smp#: 3  
Injection Vol: 15.0 ul Dil. Factor: 1.0000  
Method: PFAC\_A4 Limit Group: LC PFC\_DOD ICAL  
Column: Detector F6:MRM

15 Perfluorooctane sulfonic acid, CAS: 1763-23-1

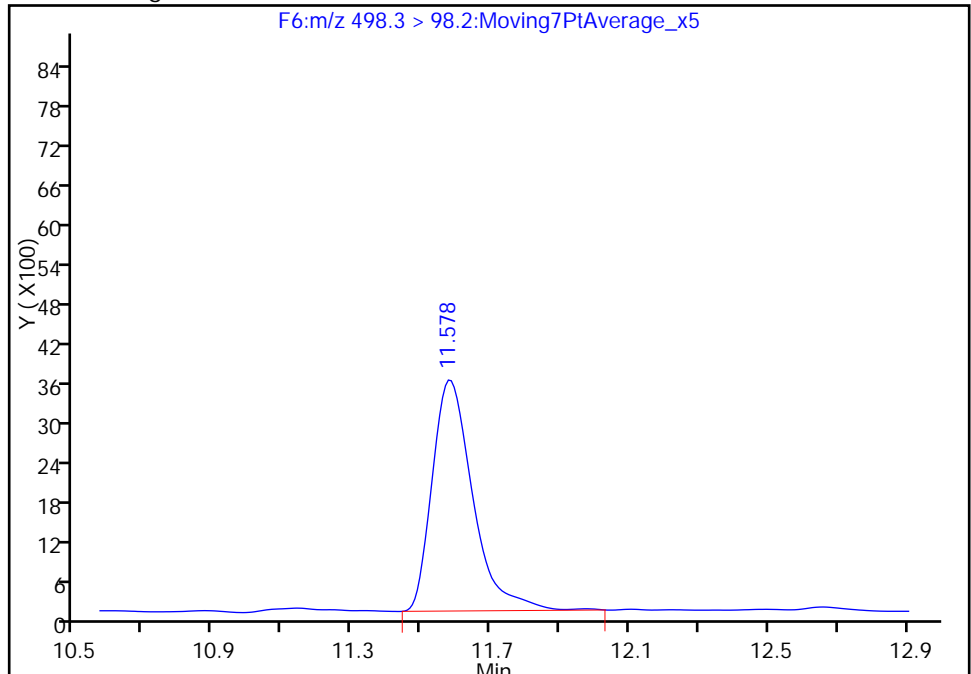
RT: 11.58  
Area: 24986  
Amount: 1.122782  
Amount Units: ng/ml

Processing Integration Results



RT: 11.58  
Area: 28574  
Amount: 1.183430  
Amount Units: ng/ml

Manual Integration Results



Reviewer: westendorfc, 15-Feb-2016 08:20:55  
Audit Action: Manually Integrated  
Audit Reason: Baseline



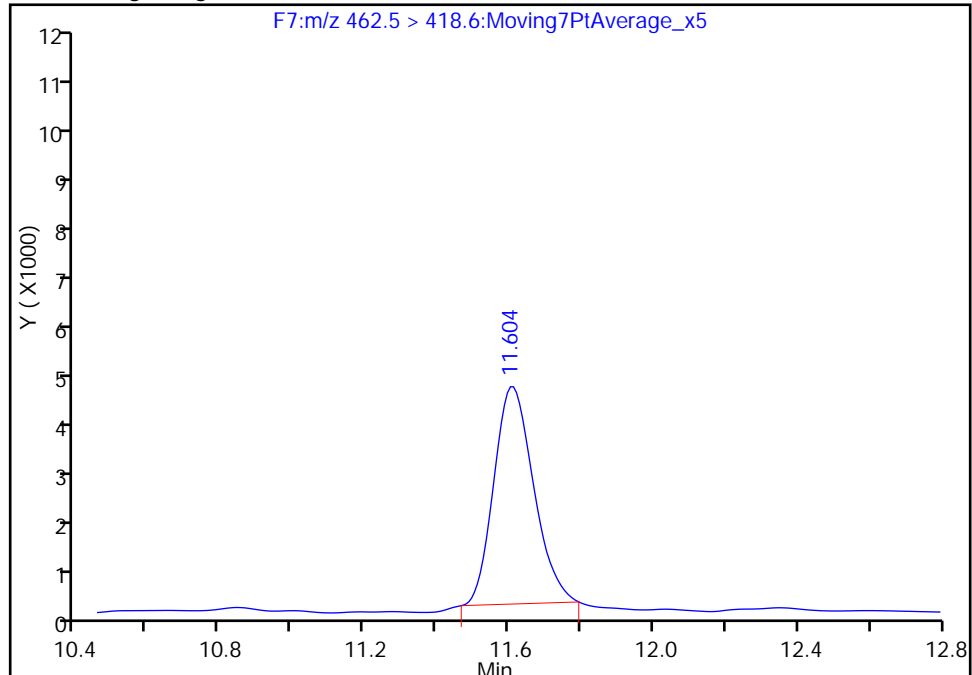
TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_004.d  
Injection Date: 12-Feb-2016 10:56:41 Instrument ID: A4  
Lims ID: Std L2  
Client ID:  
Operator ID: JRB ALS Bottle#: 3 Worklist Smp#: 3  
Injection Vol: 15.0 ul Dil. Factor: 1.0000  
Method: PFAC\_A4 Limit Group: LC PFC\_DOD ICAL  
Column: Detector F7:MRM

18 Perfluorononanoic acid, CAS: 375-95-1

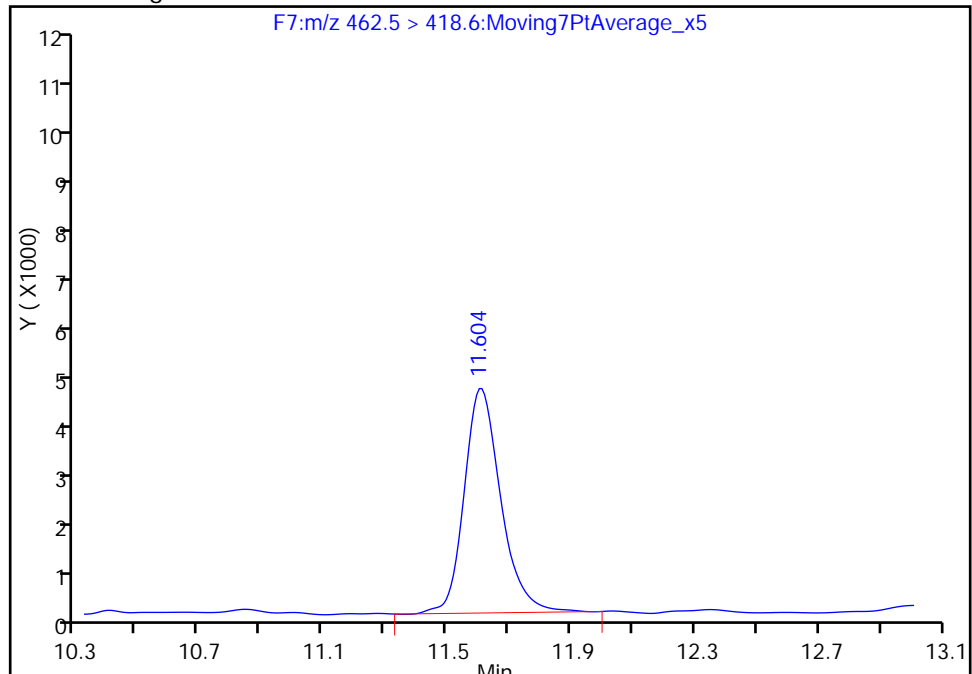
RT: 11.60  
Area: 32399  
Amount: 0.654364  
Amount Units: ng/ml

Processing Integration Results



RT: 11.60  
Area: 35930  
Amount: 0.787056  
Amount Units: ng/ml

Manual Integration Results



Reviewer: westendorfc, 15-Feb-2016 08:20:55  
Audit Action: Manually Integrated  
Audit Reason: Baseline

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_005.d  
 Lims ID: Std L3  
 Client ID:  
 Sample Type: IC Calib Level: 3  
 Inject. Date: 12-Feb-2016 11:17:51 ALS Bottle#: 4 Worklist Smp#: 4  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L3  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:02:20 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 14:30:58

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	216.7 > 171.5	5.652	5.673	-0.021	2179878	53.7		107	8464	
2 Perfluorobutyric acid	212.7 > 168.6	5.655	5.673	-0.018	116853	5.47		109	411	
D 3 13C5-PFPeA	267.6 > 222.7	6.902	6.923	-0.021	1806356	56.5		113	4928	
4 Perfluoropentanoic acid	262.9 > 218.7	6.897	6.925	-0.028	91597	5.38		108	86.2	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.025	7.047	-0.022	304142	NC			171	
	298.8 > 98.6	7.025	7.047	-0.022	183928		1.65(0.00-0.00)		131	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.025	7.047	-0.022	304142	4.26		96.4		
D 6 13C2 PFHxA	314.6 > 269.7	8.327	8.298	0.029	2203221	56.3		113	6734	
7 Perfluorohexanoic acid	312.9 > 268.7	8.327	8.304	0.023	88552	4.66		93.3	460	
D 8 13C4-PFHpA	366.6 > 321.6	9.630	9.577	0.053	2165567	55.6		111	4351	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.630	9.577	0.053	108307	4.88		97.7	279	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.665	9.612	0.053	160421	5.22		110		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.665	9.612	0.053	160421	NC			398	
D 11 18O2 PFHxS	402.5 > 83.6	9.665	9.612	0.053	1231093	54.3		115	3907	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.753	10.683	0.070	1.000	167796	6.15	129		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.753	10.683	0.070	1.000	167796	NC		913	
D 12 13C4 PFOA	416.5 > 371.6	10.762	10.683	0.079		2553061	60.0	120	4746	
13 Perfluorooctanoic acid	412.8 > 368.8	10.753	10.692	0.061	1.000	131613	5.25	105	261	
	412.8 > 168.7	10.762	10.692	0.070	1.001	41266	3.19(0.00-0.00)	105	123	
D 16 13C4 PFOS	502.4 > 79.7	11.678	11.602	0.076		512338	50.6	106	1864	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.678	11.606	0.072	1.000	270891	5.93	124	1015	
	498.3 > 98.2	11.678	11.606	0.072	1.000	170527	1.59(0.00-0.00)	124	450	
D 17 13C5 PFNA	467.5 > 422.6	11.704	11.628	0.076		2361526	56.8	114	3847	
18 Perfluorononanoic acid	462.5 > 418.6	11.704	11.628	0.076	1.000	230097	4.75	94.9	411	
D 19 13C2 PFDA	514.4 > 469.5	12.508	12.419	0.089		3587771	61.1	122	5014	
20 Perfluorodecanoic acid	512.5 > 468.5	12.508	12.424	0.084	1.000	328682	5.01	100	1356	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	13.021	12.937	0.084	1.000	496084	5.09	102	1219	
D 23 13C8 FOSA	505.4 > 77.6	13.021	12.937	0.084		5160232	57.8	116	2642	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.128	13.058	0.070	1.000	125658	NC		467	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.128	13.058	0.070	1.000	125658	6.08	126		
D 26 13C2 PFUnA	564.3 > 519.5	13.181	13.097	0.084		3765798	64.4	129	5102	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.181	13.097	0.084	1.000	386752	4.95	99.0	505	
D 28 13C2 PFDaA	614.4 > 569.4	13.739	13.674	0.065		4112735	57.8	116	5008	
29 Perfluorododecanoic acid	612.4 > 568.6	13.749	13.674	0.075	1.000	347709	5.25	105	222	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.220	14.155	0.065	1.000	378224	5.51	110	270	
D 33 13C2-PFTeDA	714.5 > 669.5	14.635	14.569	0.066		3587033	56.6	113	4316	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.635	14.572	0.063	1.000	151695	4.64	92.8	100	
D 35 13C2-PFHxDA	814.8 > 769.6	15.159	15.119	0.040		1599911	53.9	108	3273	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.159	15.119	0.040	1.000	425951	4.66	93.2	417	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7	> 868.6	15.461	15.428	0.033	1.000	341392	4.63	92.7	331

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L3\_00016

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_005.d

Injection Date: 12-Feb-2016 11:17:51

Instrument ID: A4

Lims ID: Std L3

Client ID:

Operator ID: JRB

ALS Bottle#: 4

Worklist Smp#: 4

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

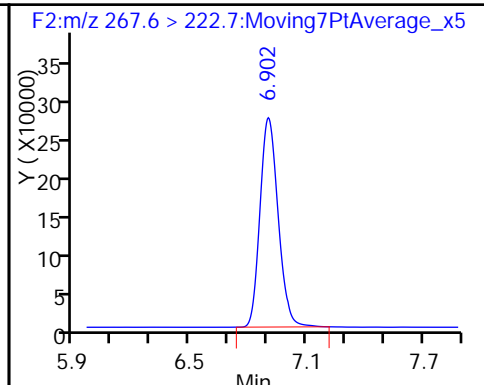
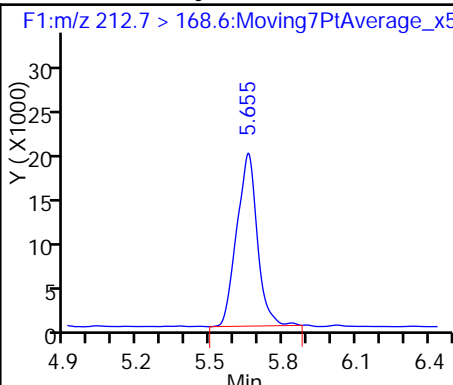
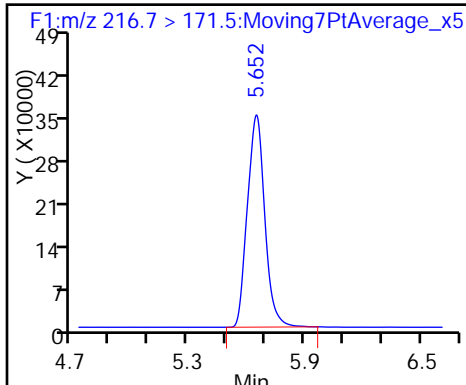
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

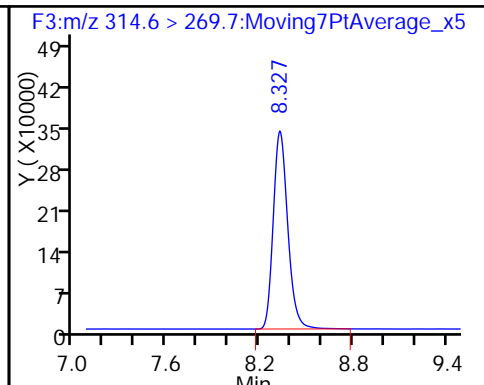
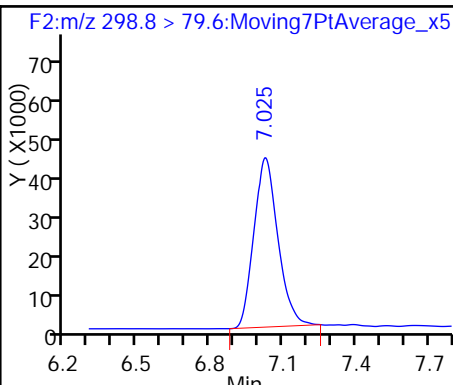
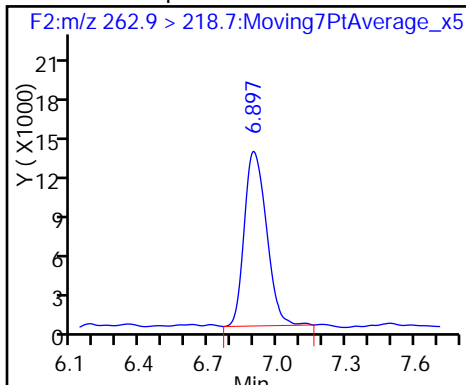
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

51 Perfluorobutanesulfonic acid

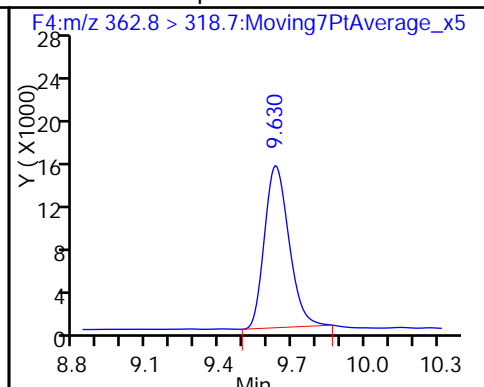
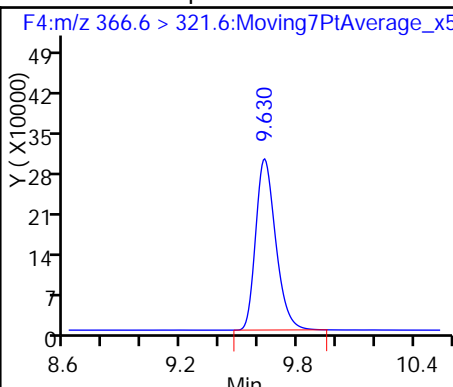
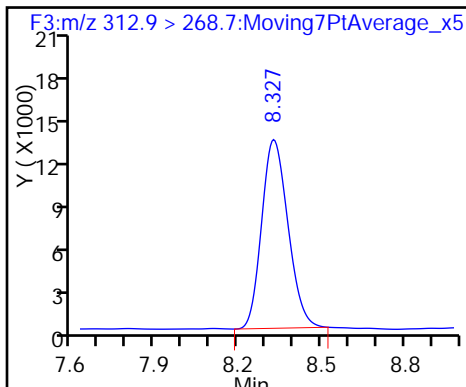
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

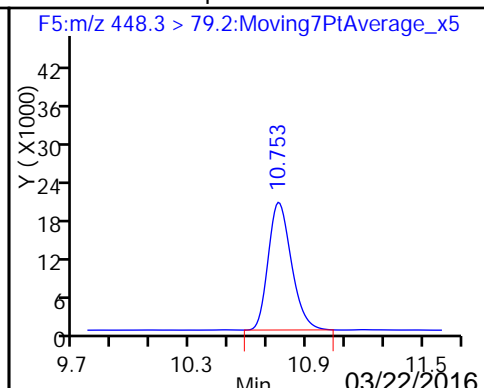
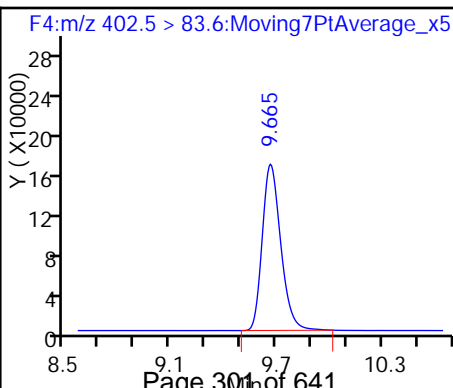
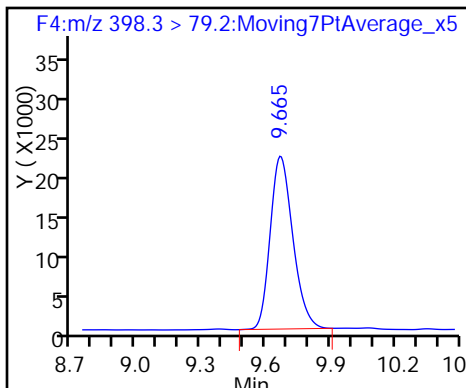
9 Perfluoroheptanoic acid



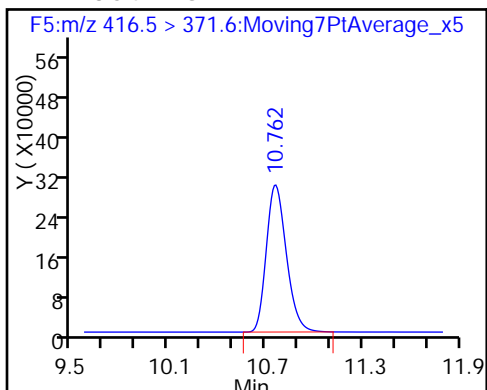
58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

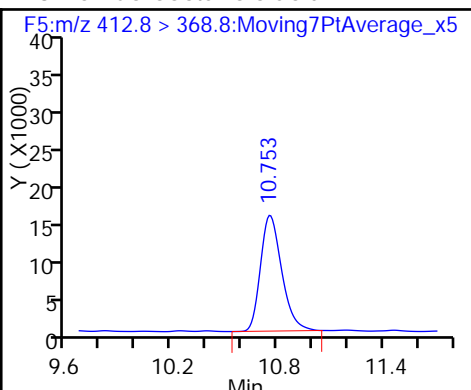
39 Perfluoroheptanesulfonic Acid



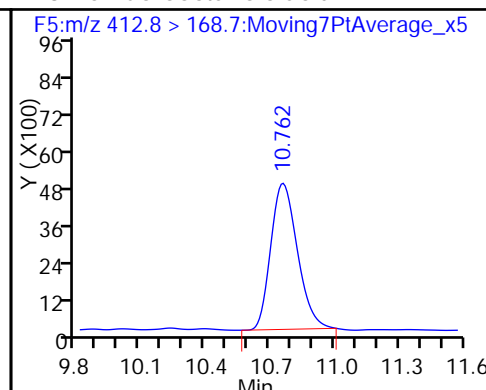
D 12 13C4 PFOA



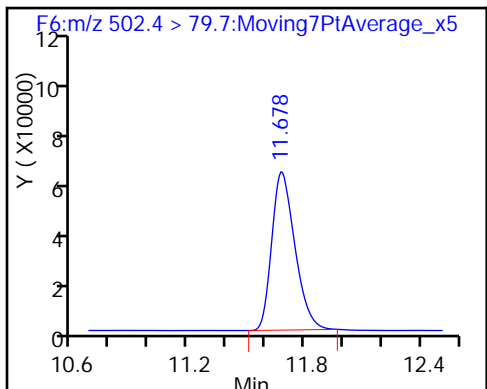
13 Perfluorooctanoic acid



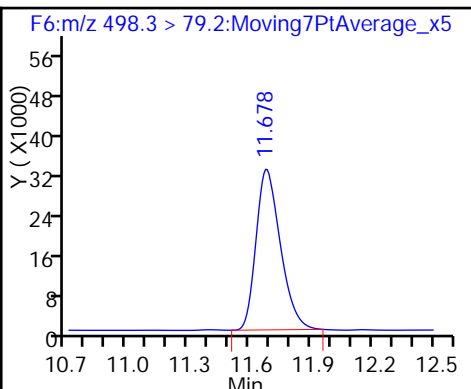
13 Perfluorooctanoic acid



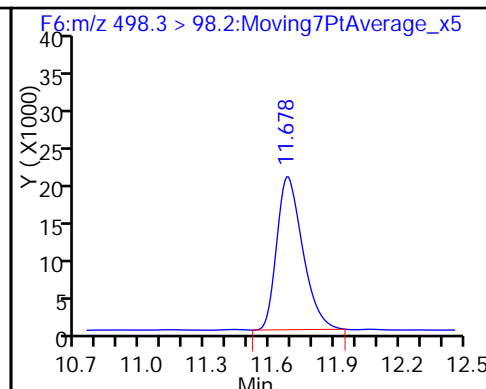
D 16 13C4 PFOS



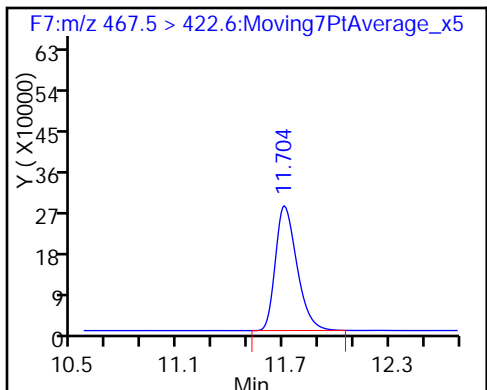
15 Perfluorooctane sulfonic acid



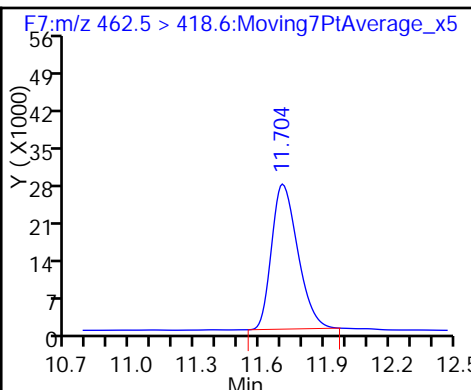
15 Perfluorooctane sulfonic acid



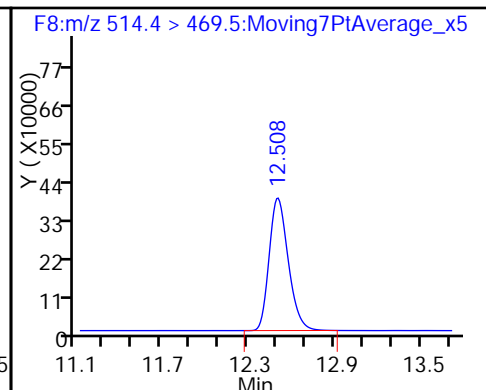
D 17 13C5 PFNA



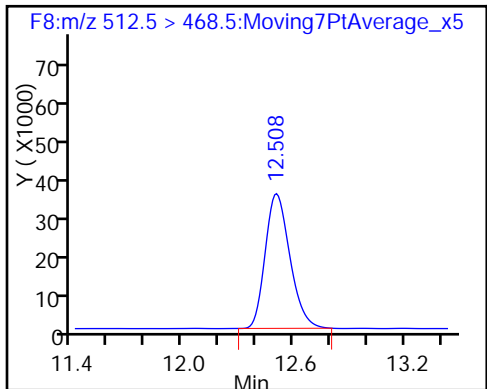
18 Perfluorononanoic acid



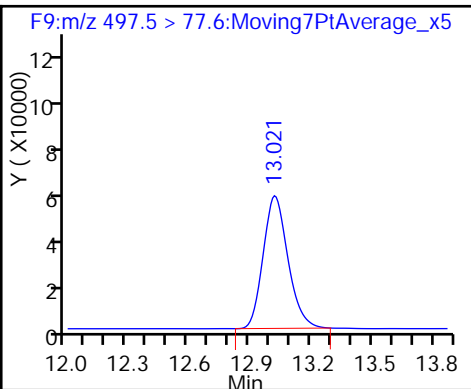
D 19 13C2 PFDA



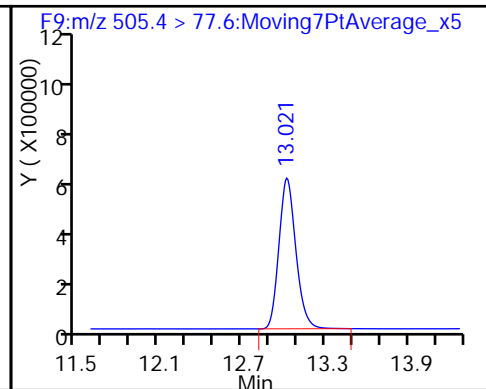
20 Perfluorodecanoic acid



24 Perfluorooctane Sulfonamide



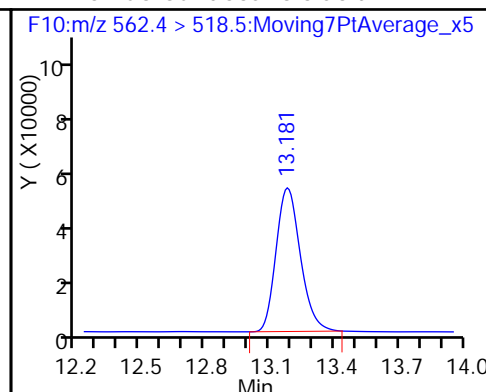
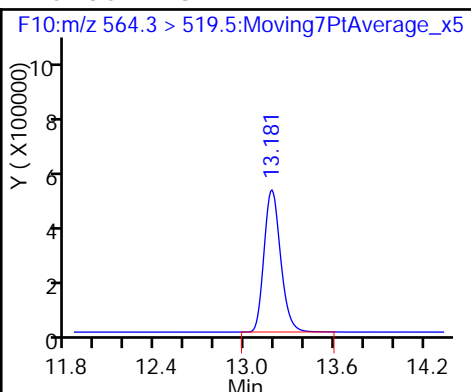
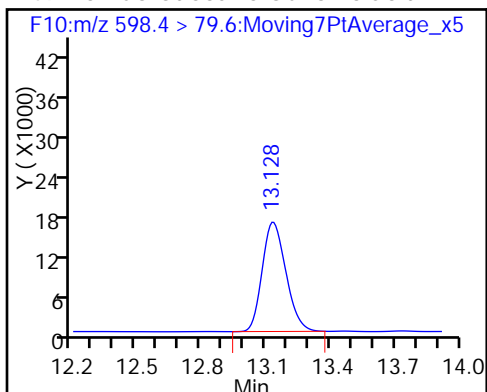
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

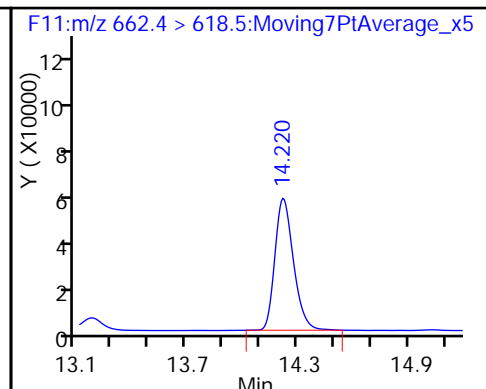
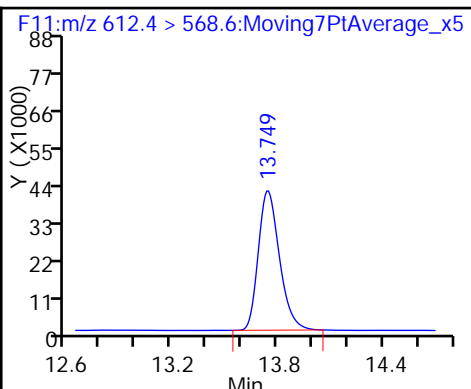
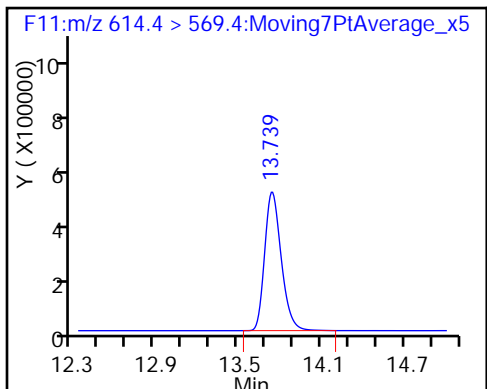
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

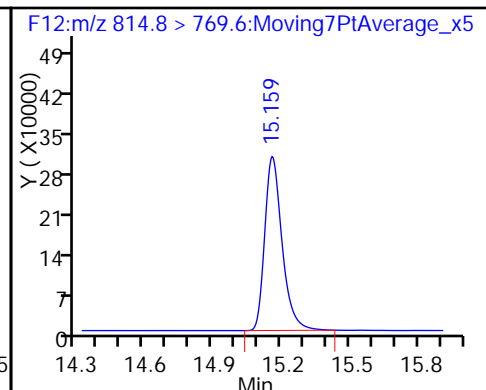
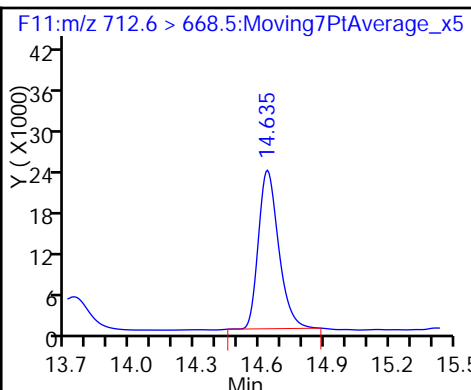
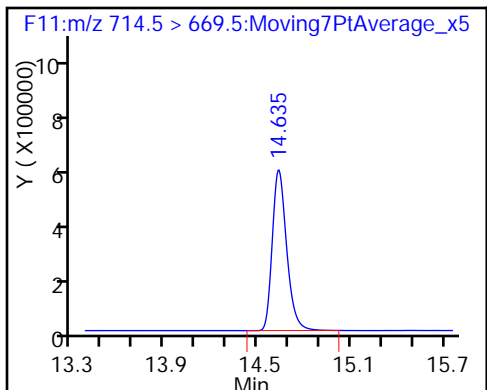
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

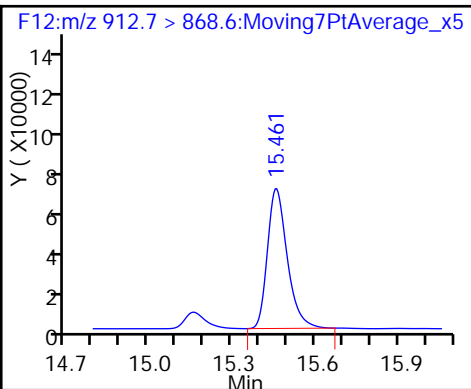
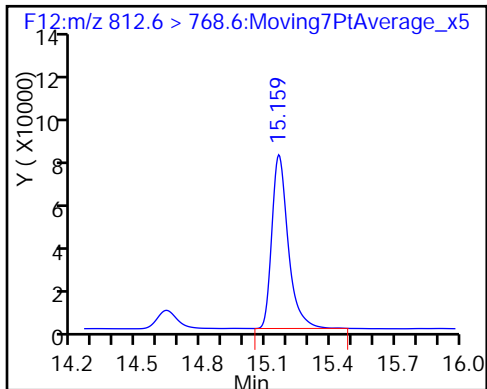
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_006.d  
 Lims ID: Std L4  
 Client ID:  
 Sample Type: IC Calib Level: 4  
 Inject. Date: 12-Feb-2016 11:39:03 ALS Bottle#: 5 Worklist Smp#: 5  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L4  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 11:03:53 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: westendorfc Date: 15-Feb-2016 11:03:53

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	216.7 > 171.5	5.757	5.712	0.045	2252325	55.5		111	6938	
2 Perfluorobutyric acid	212.7 > 168.6	5.753	5.712	0.041	448061	20.3		101	1325	
4 Perfluoropentanoic acid	262.9 > 218.7	7.058	6.986	0.072	343823	20.5		102	278	
D 3 13C5-PFPeA	267.6 > 222.7	7.058	6.986	0.072	1781896	55.7		111	4204	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.190	7.115	0.075	512492	NC			346	
	298.8 > 98.6	7.187	7.115	0.072	330932		1.55(0.00-0.00)		261	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.190	7.115	0.075	512492	18.1		102		
D 6 13C2 PFHxA	314.6 > 269.7	8.432	8.367	0.065	2274718	58.1		116	6607	
7 Perfluorohexanoic acid	312.9 > 268.7	8.432	8.369	0.063	394416	20.1		101	1381	
D 8 13C4-PFHpA	366.6 > 321.6	9.685	9.639	0.046	2111070	54.2		108	4381	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.692	9.640	0.052	438459	20.3		101	1619	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.720	9.675	0.045	532601	18.7		98.8		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.720	9.675	0.045	532601	NC			1420	
D 11 18O2 PFHxS	402.5 > 83.6	9.720	9.675	0.045	1191134	52.5		111	2897	



Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.770	10.740	0.030	1.000	631312	19.5	102		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.770	10.740	0.030	1.000	631312	NC		2265	
D 12 13C4 PFOA	416.5 > 371.6	10.779	10.743	0.036		2286718	53.7	107	3668	
13 Perfluorooctanoic acid	412.8 > 368.8	10.770	10.743	0.027	1.000	428131	19.1	95.3	775	
	412.8 > 168.7	10.779	10.743	0.036	1.001	164089	2.61(0.00-0.00)	95.3	523	
D 16 13C4 PFOS	502.4 > 79.7	11.677	11.651	0.026		609839	60.2	126	2302	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.677	11.656	0.021	1.000	1000210	18.4	96.1	2032	
	498.3 > 98.2	11.677	11.656	0.021	1.000	596288	1.68(0.00-0.00)	96.1	1477	
D 17 13C5 PFNA	467.5 > 422.6	11.703	11.678	0.025		2172455	52.2	104	3784	
18 Perfluorononanoic acid	462.5 > 418.6	11.703	11.680	0.023	1.000	984533	21.7	108	1232	
D 19 13C2 PFDA	514.4 > 469.5	12.486	12.473	0.013		3126636	53.2	106	4935	
20 Perfluorodecanoic acid	512.5 > 468.5	12.496	12.476	0.020	1.000	1134637	19.8	99.2	2598	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	13.013	12.991	0.022	1.000	1982787	20.1	100	2317	
D 23 13C8 FOSA	505.4 > 77.6	13.013	12.991	0.022		5229659	58.5	117	2282	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.111	13.098	0.013	1.000	467598	NC		1809	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.111	13.098	0.013	1.000	467598	19.0	98.7		
D 26 13C2 PFUnA	564.3 > 519.5	13.163	13.146	0.017		3171146	54.2	108	3316	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.163	13.146	0.017	1.000	1318141	20.0	100	1564	
D 28 13C2 PFDoA	614.4 > 569.4	13.730	13.713	0.017		3714597	52.2	104	3490	
29 Perfluorododecanoic acid	612.4 > 568.6	13.730	13.715	0.015	1.000	1253957	21.0	105	682	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.205	14.193	0.012	1.000	1410790	22.8	114	894	
D 33 13C2-PFTeDA	714.5 > 669.5	14.615	14.603	0.012		3578785	56.5	113	4100	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.615	14.604	0.011	1.000	573742	20.2	101	358	
D 35 13C2-PFHxDA	814.8 > 769.6	15.144	15.138	0.006		1823759	61.5	123	3523	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.144	15.139	0.005	1.000	1600494	19.8	99.2	1464	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.446	15.442	0.004	1.000	1468803	20.1	100	1170	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L4\_00017

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_006.d

Injection Date: 12-Feb-2016 11:39:03

Instrument ID: A4

Lims ID: Std L4

Client ID:

Operator ID: JRB

ALS Bottle#: 5

Worklist Smp#: 5

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

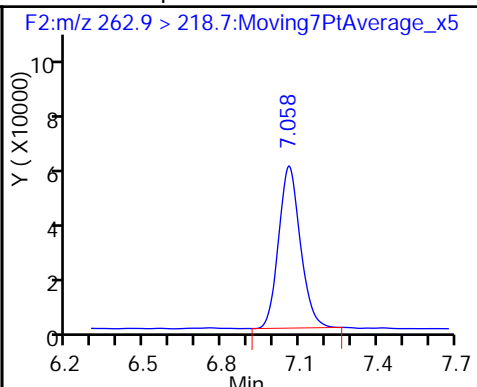
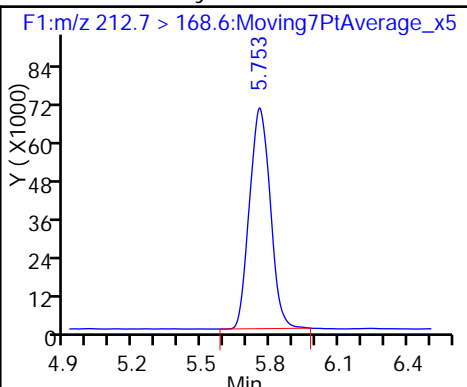
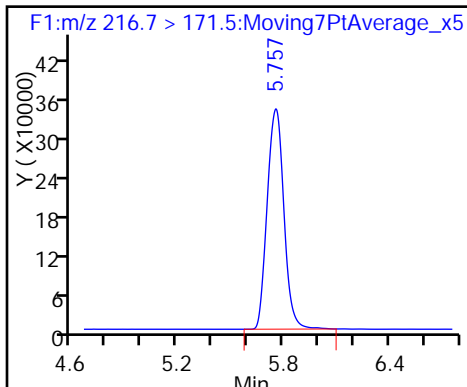
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

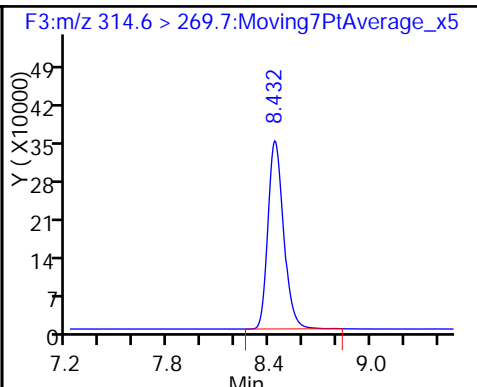
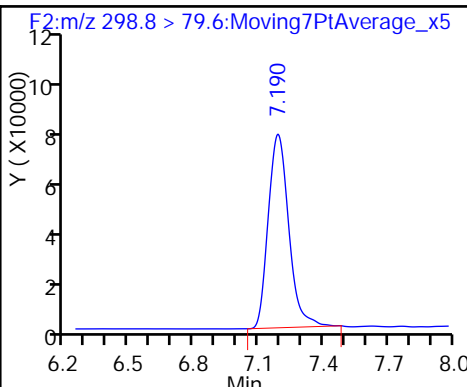
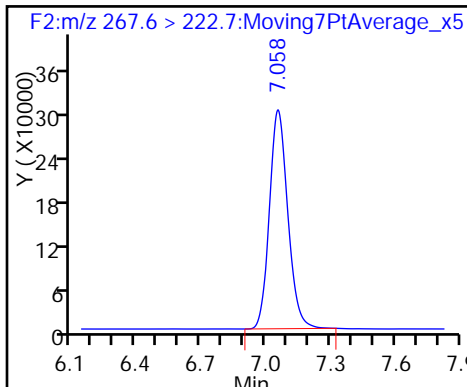
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

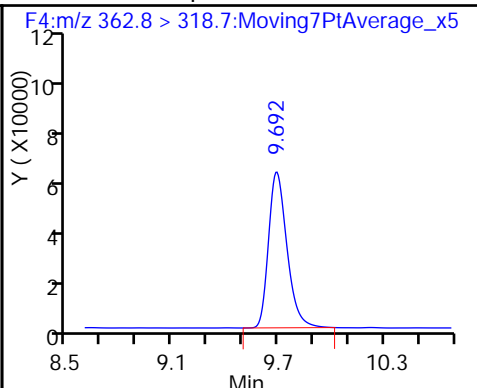
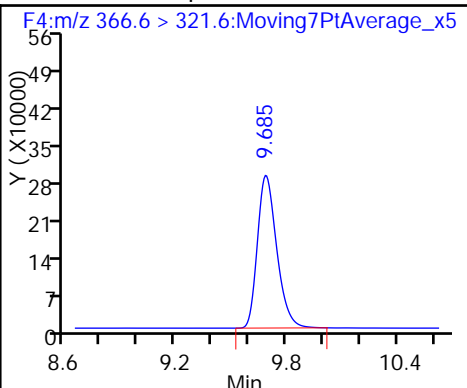
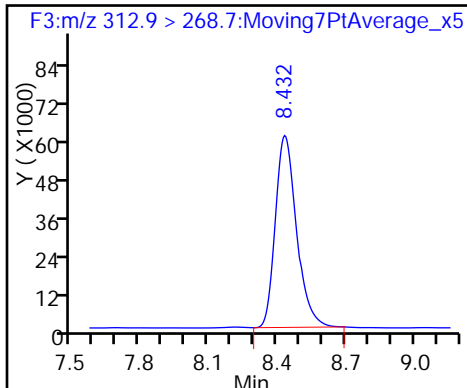
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

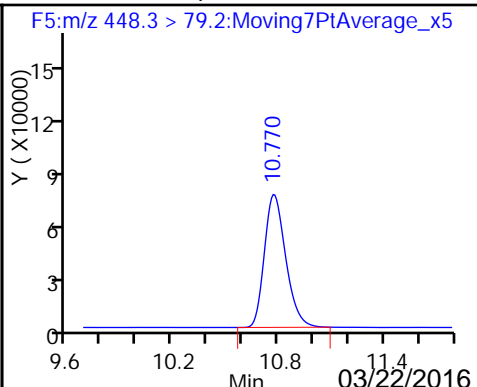
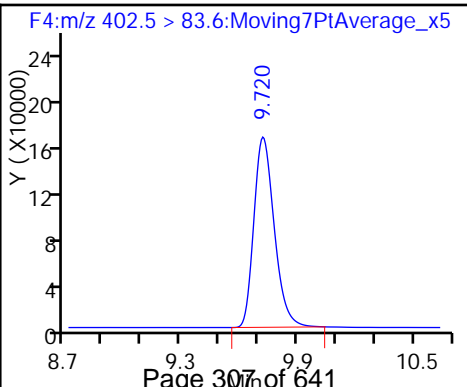
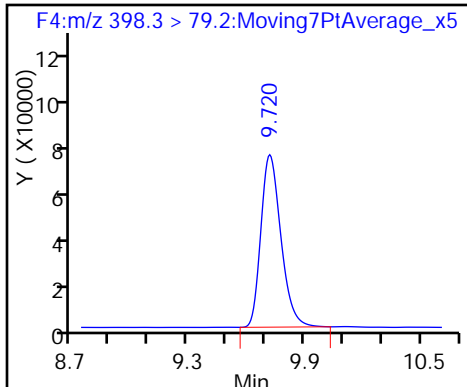
9 Perfluoroheptanoic acid



58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

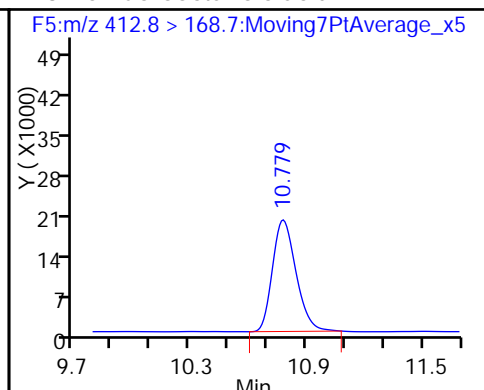
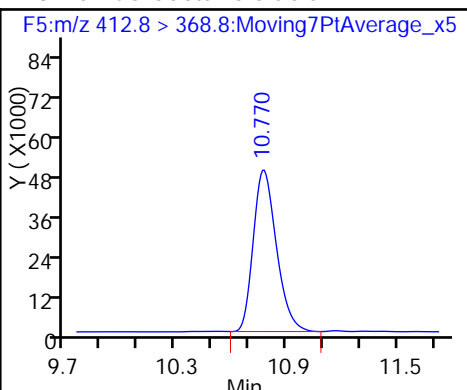
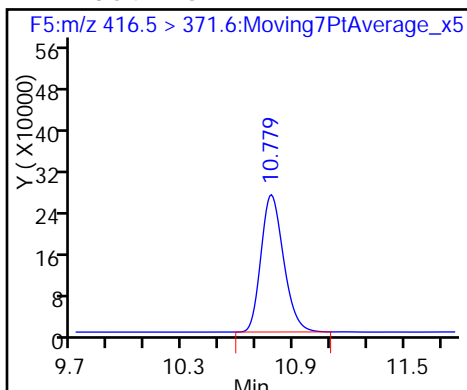
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

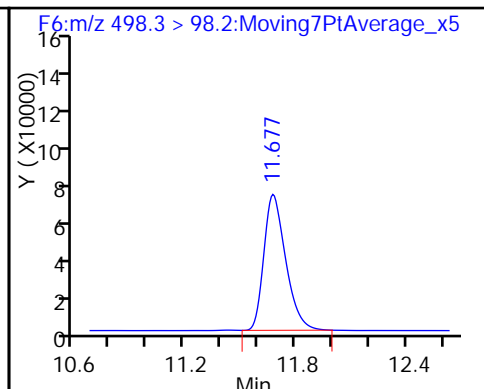
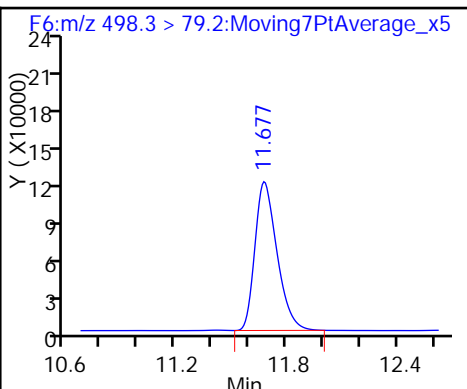
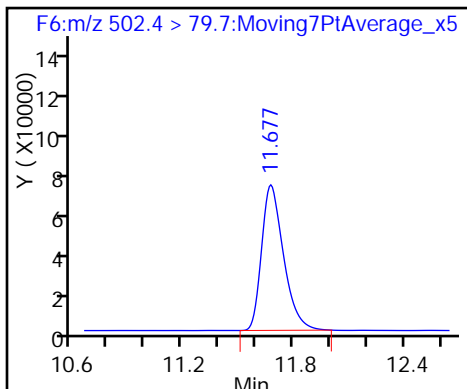
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid

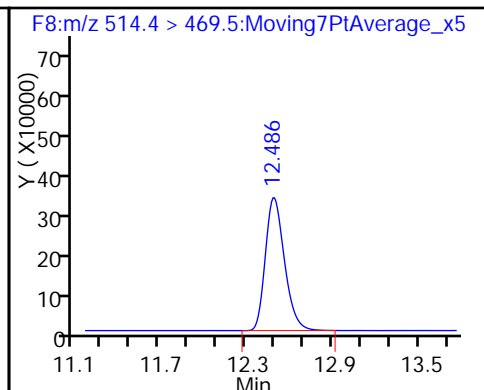
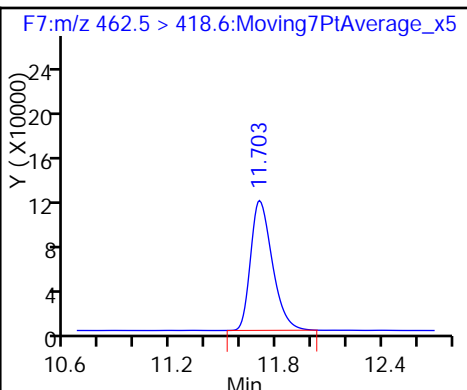
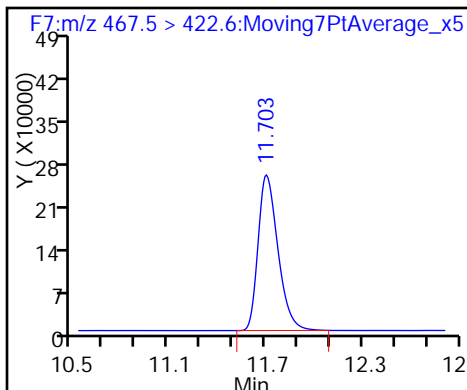
15 Perfluorooctane sulfonic acid



D 17 13C5 PFNA

18 Perfluorononanoic acid

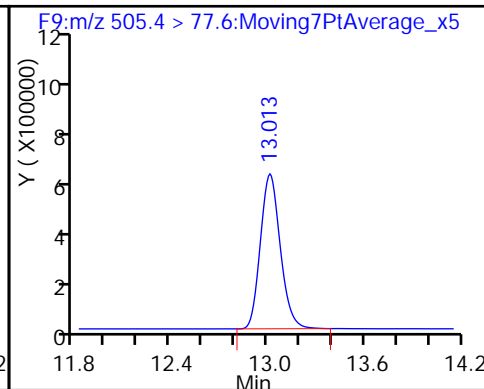
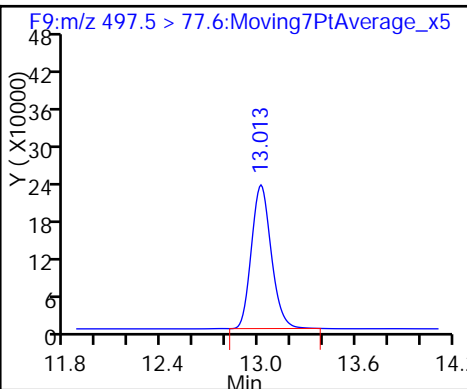
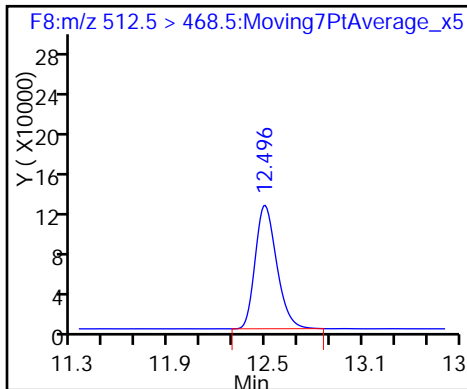
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

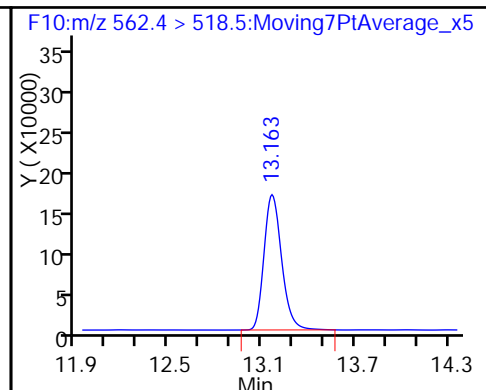
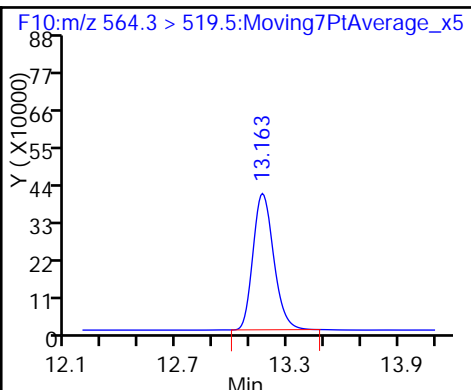
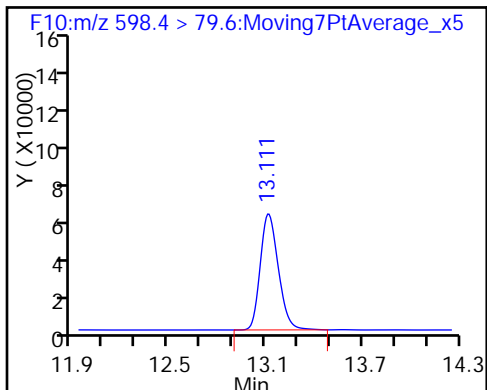
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

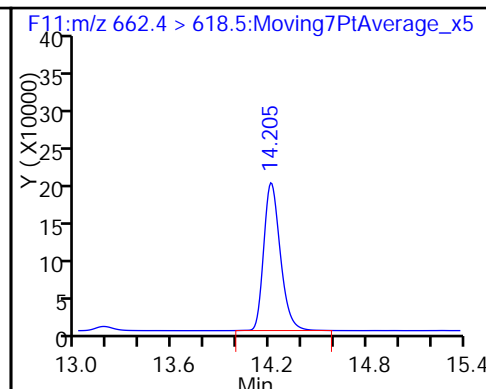
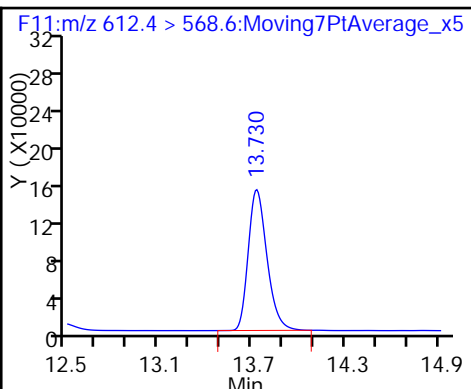
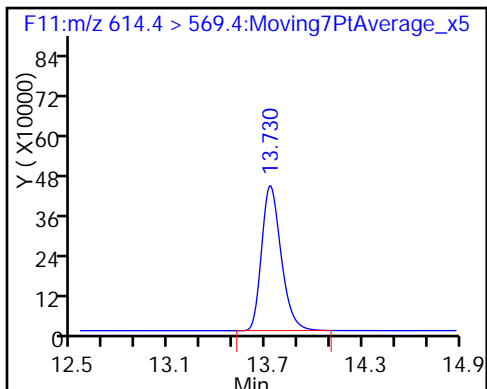
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

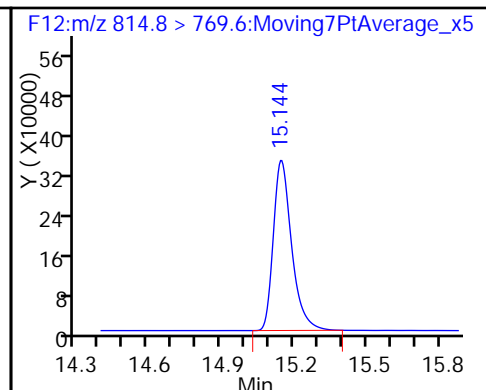
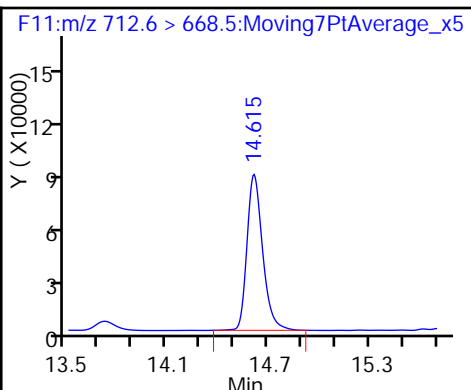
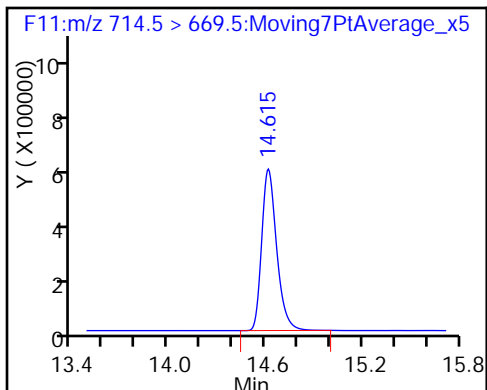
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

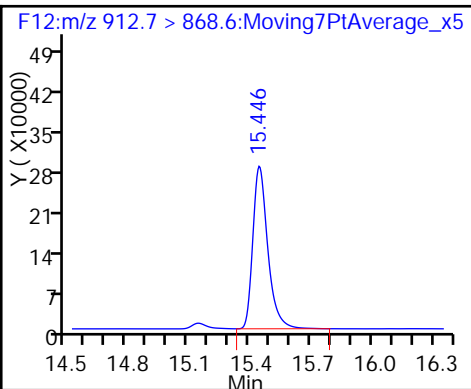
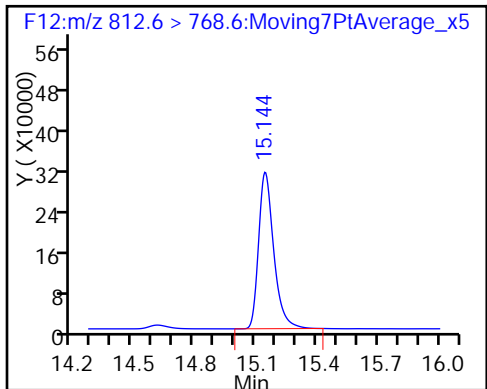
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_007.d  
 Lims ID: Std L5  
 Client ID:  
 Sample Type: IC Calib Level: 5  
 Inject. Date: 12-Feb-2016 12:00:14 ALS Bottle#: 6 Worklist Smp#: 6  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L5  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:31:48 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 14:27:49

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
--------	----	--------	--------	--------	----------	--------------	---------------	------	-----	-------

D 1 13C4 PFBA										
216.7 > 171.5	5.735	5.712	0.023		2075195	51.1		102	6828	
2 Perfluorobutyric acid										
212.7 > 168.6	5.735	5.712	0.023	1.000	1048362	51.5		103	2760	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.015	6.986	0.029	1.000	810081	50.5		101	925	
D 3 13C5-PFPeA										
267.6 > 222.7	7.018	6.986	0.032		1702935	53.2		106	4252	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.144	7.115	0.029	1.000	938494	NC			632	
298.8 > 98.6	7.147	7.115	0.032	1.000	594597		1.58(0.00-0.00)		483	
51 Perfluorobutanesulfonic acid										
298.8 > 79.6	7.144	7.115	0.029	1.000	938494	49.5		112		
D 6 13C2 PFHxA										
314.6 > 269.7	8.393	8.367	0.026		1984057	50.7		101	4591	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.393	8.369	0.024	1.000	886159	51.8		104	1218	
D 8 13C4-PFHpA										
366.6 > 321.6	9.663	9.639	0.024		1911848	49.1		98.2	3668	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.663	9.640	0.023	1.000	929700	47.5		95.0	1818	
58 Perfluorohexanesulfonic acid										
398.3 > 79.2	9.698	9.675	0.023	1.000	1192387	45.4		95.9		
10 Perfluorohexane Sulfonate										
398.3 > 79.2	9.698	9.675	0.023	1.000	1192387	NC			2528	
D 11 18O2 PFHxS										
402.5 > 83.6	9.698	9.675	0.023		1110488	49.0		104	3185	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.760	10.740	0.020	1.000	1320551	36.6	76.9		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.760	10.740	0.020	1.000	1320551	NC		3285	
D 12 13C4 PFOA	416.5 > 371.6	10.760	10.743	0.017		2064639	48.5	97.0	4189	
13 Perfluorooctanoic acid	412.8 > 368.8	10.760	10.743	0.017	1.000	969508	47.8	95.6	1437	
	412.8 > 168.7	10.770	10.743	0.027	1.001	375980	2.58(0.00-0.00)	95.6	1362	
D 16 13C4 PFOS	502.4 > 79.7	11.668	11.651	0.017		677767	66.9	140	1996	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.676	11.656	0.020	1.000	2377814	39.3	82.3	2245	
	498.3 > 98.2	11.676	11.656	0.020	1.000	1271737	1.87(0.00-0.00)	82.3	2060	
D 17 13C5 PFNA	467.5 > 422.6	11.694	11.678	0.016		1984772	47.7	95.4	2612	
18 Perfluorononanoic acid	462.5 > 418.6	11.694	11.680	0.014	1.000	2055860	49.4	98.8	1799	
D 19 13C2 PFDA	514.4 > 469.5	12.493	12.473	0.020		2605584	44.4	88.7	3301	
20 Perfluorodecanoic acid	512.5 > 468.5	12.493	12.476	0.017	1.000	2455156	51.5	103	4334	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	13.010	12.991	0.019	1.000	4712375	51.8	104	2004	
D 23 13C8 FOSA	505.4 > 77.6	13.010	12.991	0.019		4811591	53.9	108	2184	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.116	13.098	0.018	1.000	1059583	NC		2639	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.116	13.098	0.018	1.000	1059583	38.8	80.5		
D 26 13C2 PFUnA	564.3 > 519.5	13.169	13.146	0.023		2571081	43.9	87.9	3012	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.169	13.146	0.023	1.000	2719401	51.0	102	2174	
D 28 13C2 PFDaA	614.4 > 569.4	13.727	13.713	0.014		3479267	48.9	97.8	3001	
29 Perfluorododecanoic acid	612.4 > 568.6	13.727	13.715	0.012	1.000	2956390	52.8	106	1702	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.210	14.193	0.017	1.000	2877547	49.6	99.1	1541	
D 33 13C2-PFTeDA	714.5 > 669.5	14.613	14.603	0.010		3289262	51.9	104	4573	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.613	14.604	0.009	1.000	1341162	50.7	101	826	
D 35 13C2-PFHxDA	814.8 > 769.6	15.146	15.138	0.008		1830150	61.7	123	3757	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.146	15.139	0.007	1.000	3609294	48.0	95.9	2731	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7	> 868.6	15.448	15.442	0.006	1.000	3340365	47.9	95.9	2275

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L5\_00016

Amount Added: 1.00

Units: mL



TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_007.d

Injection Date: 12-Feb-2016 12:00:14

Instrument ID: A4

Lims ID: Std L5

Client ID:

Operator ID: JRB

ALS Bottle#: 6

Worklist Smp#: 6

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

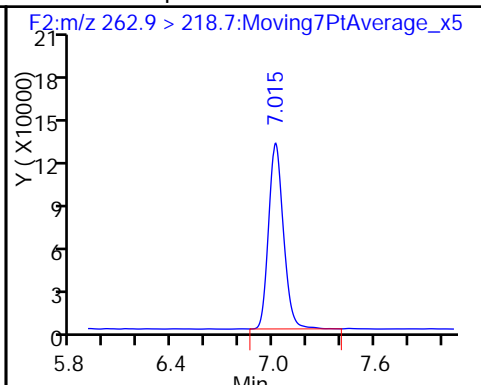
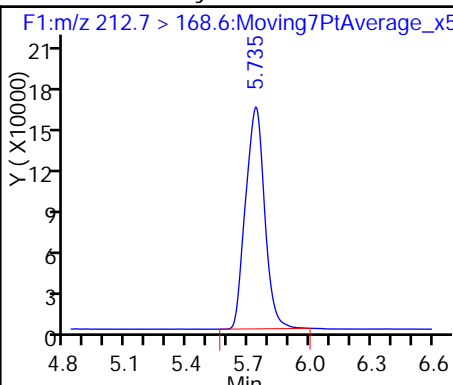
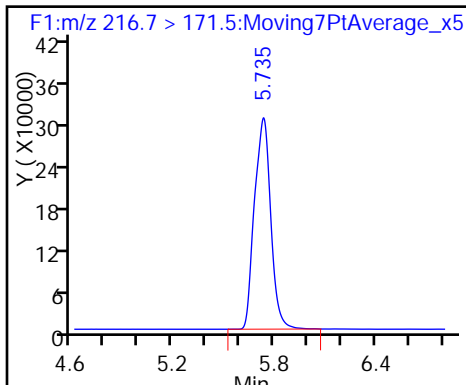
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

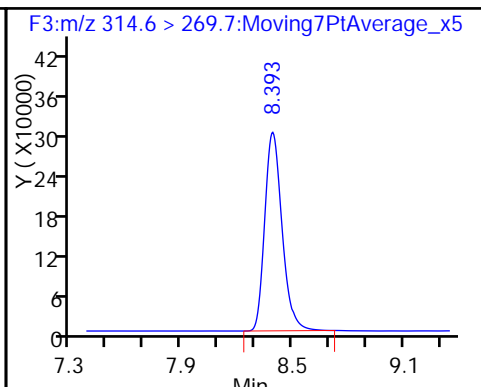
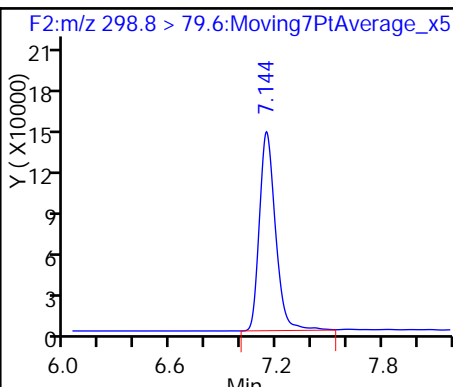
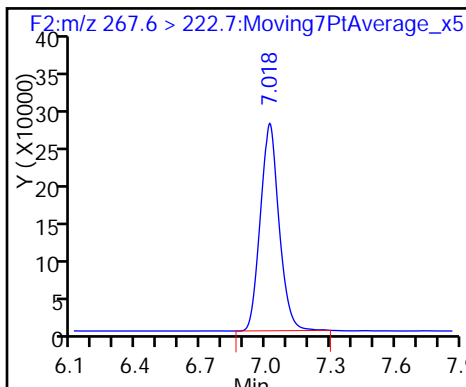
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

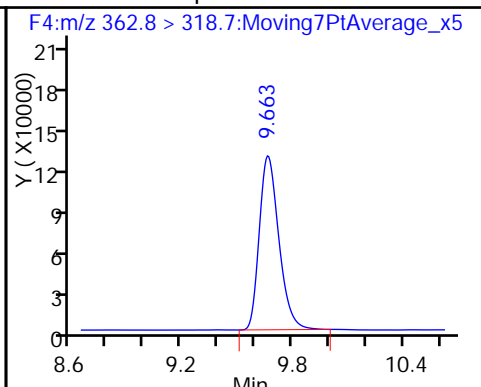
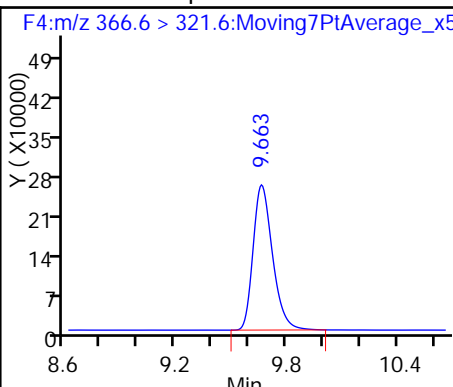
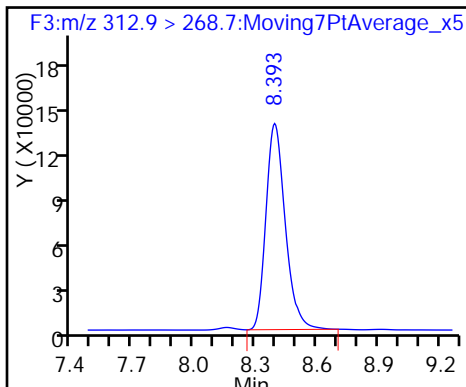
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

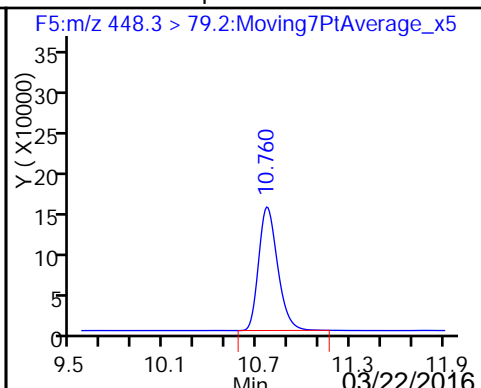
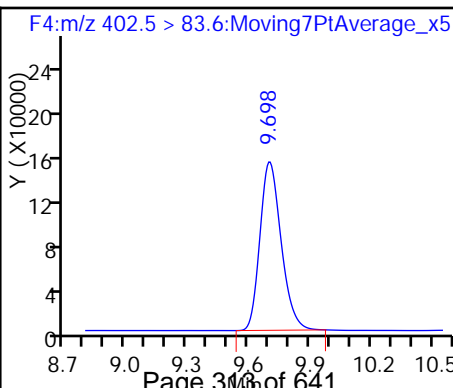
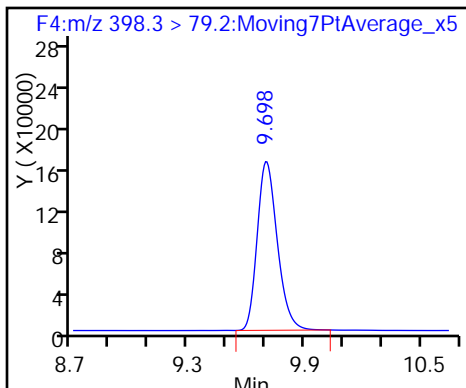
9 Perfluoroheptanoic acid



58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

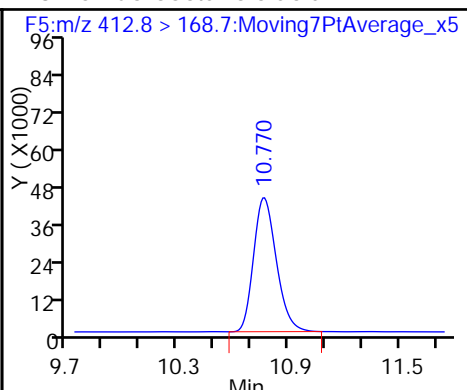
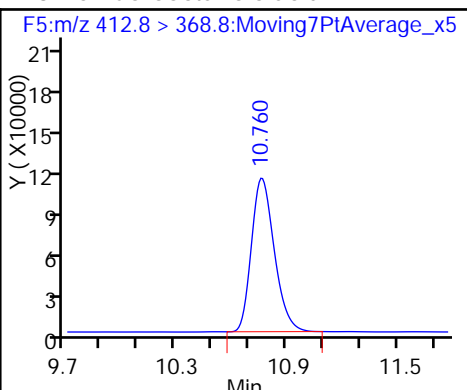
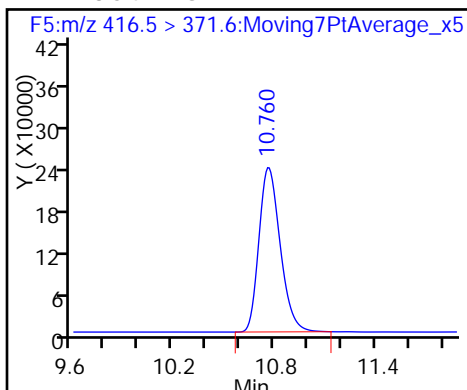
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

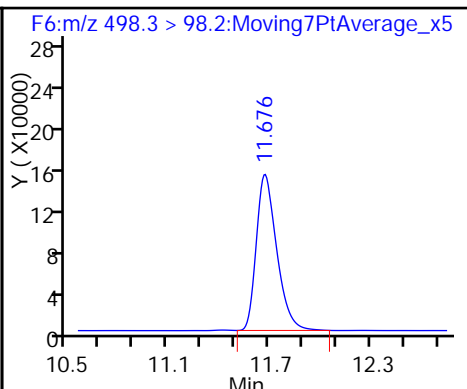
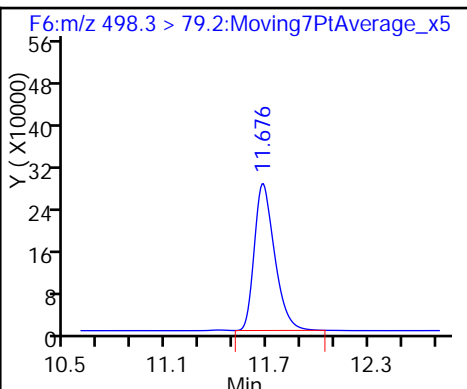
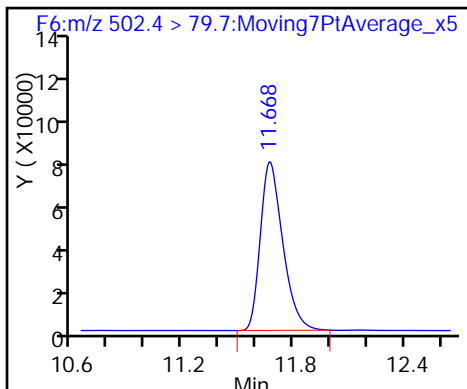
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid

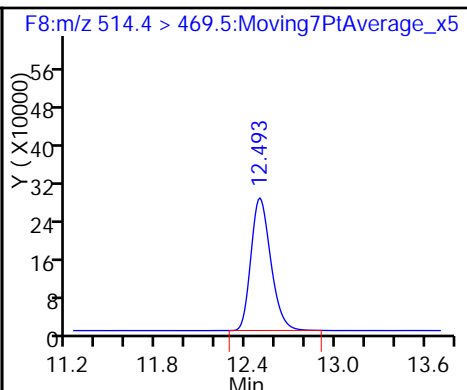
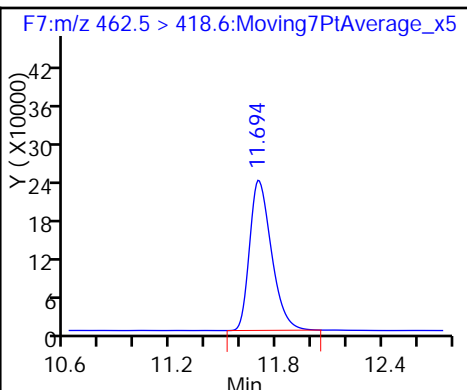
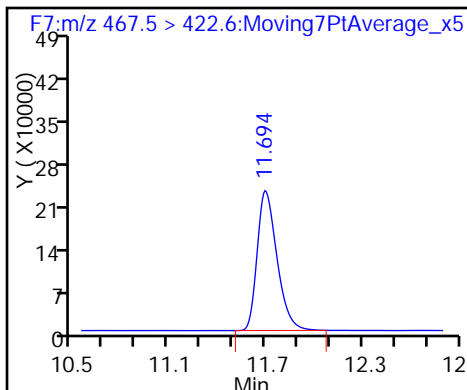
15 Perfluorooctane sulfonic acid



D 17 13C5 PFNA

18 Perfluorononanoic acid

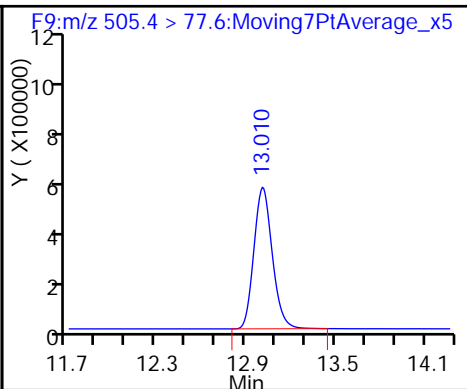
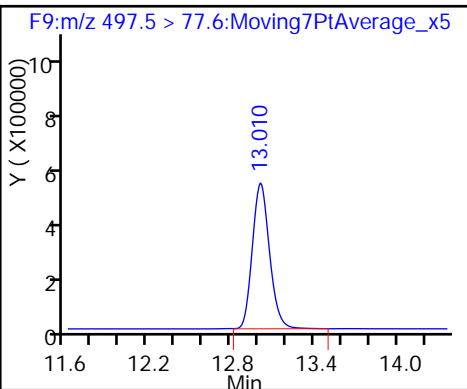
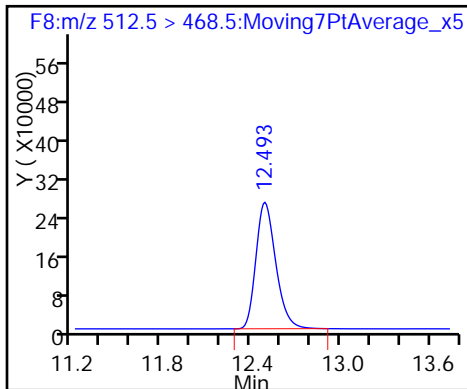
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

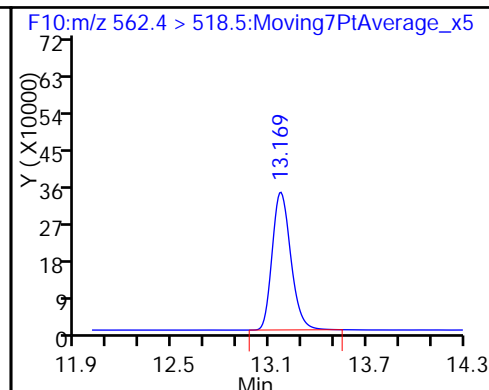
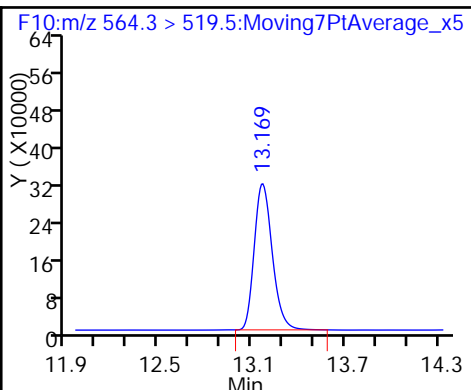
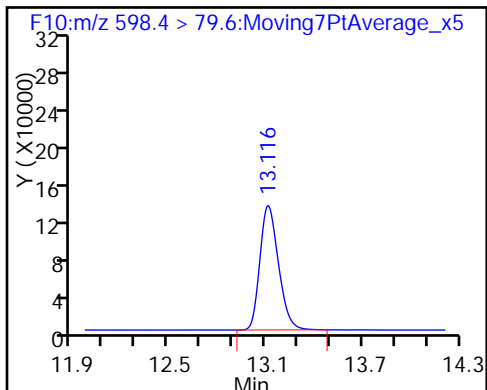
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

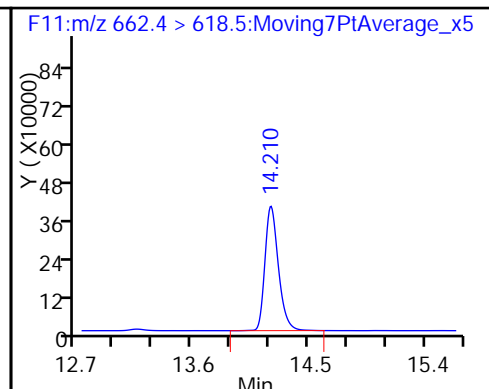
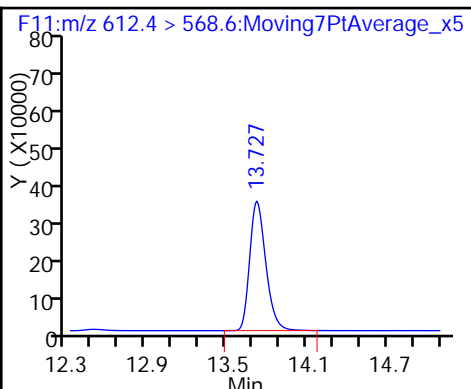
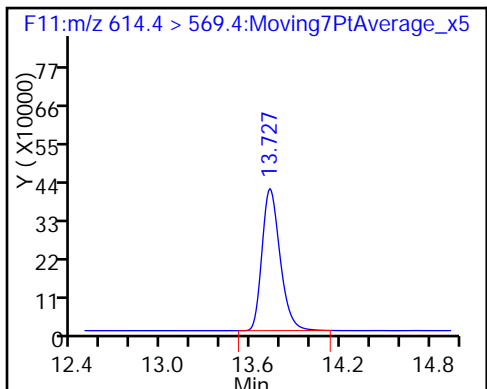
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

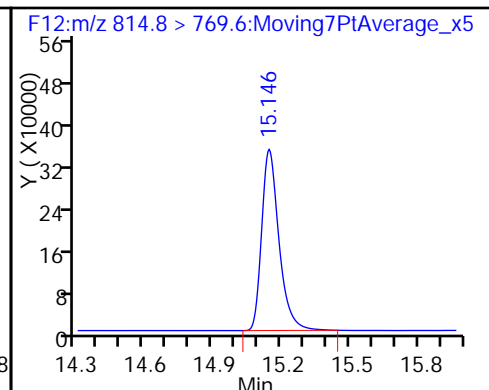
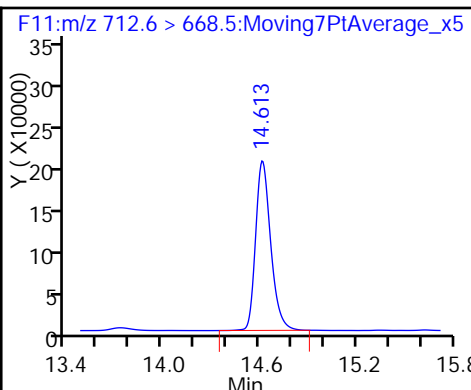
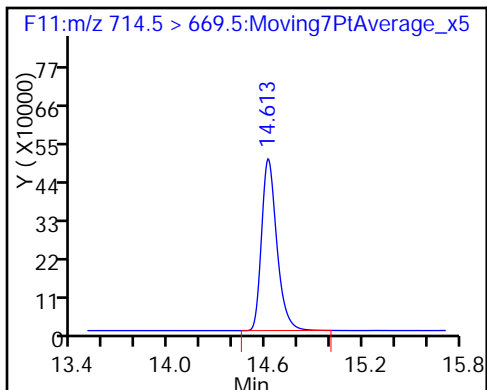
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

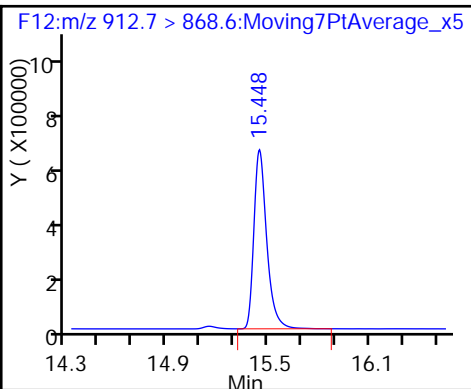
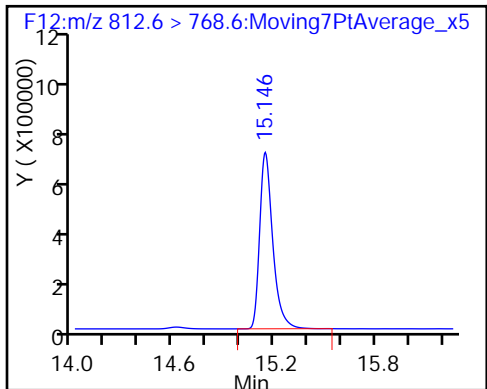
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_008.d  
 Lims ID: Std L6  
 Client ID:  
 Sample Type: IC Calib Level: 6  
 Inject. Date: 12-Feb-2016 12:21:25 ALS Bottle#: 7 Worklist Smp#: 7  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L6  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:02:34 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 14:28:17

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	216.7 > 171.5	5.747	5.698	0.049	1710809	42.2		84.3	4990	
2 Perfluorobutyric acid	212.7 > 168.6	5.750	5.698	0.052	1.000	3586362	213.9	107	8559	
4 Perfluoropentanoic acid	262.9 > 218.7	7.043	6.964	0.079	1.000	2499117	206.5	103	1697	
D 3 13C5-PFPeA	267.6 > 222.7	7.043	6.965	0.078	1285003	40.2		80.3	3117	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.179	7.090	0.089	1.000	2289993	NC		1422	
	298.8 > 98.6	7.175	7.090	0.085	0.999	1487976	1.54(0.00-0.00)		1019	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.179	7.090	0.089	1.000	2289993	177.6	100		
D 6 13C2 PFHxA	314.6 > 269.7	8.409	8.350	0.059	1514915	38.7		77.4	3765	
7 Perfluorohexanoic acid	312.9 > 268.7	8.409	8.352	0.057	1.000	2957321	226.6	113	1691	
D 8 13C4-PFHpA	366.6 > 321.6	9.665	9.627	0.038	1467997	37.7		75.4	3017	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.665	9.628	0.037	1.000	3093796	205.8	103	4736	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.707	9.661	0.046	1.000	3857794	181.7	96.0		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.707	9.661	0.046	1.000	3857794	NC		4285	
D 11 18O2 PFHxS	402.5 > 83.6	9.707	9.661	0.046	901928	39.8		84.1	2021	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.762	10.730	0.032	1.000	3860458		78.9		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.762	10.730	0.032	1.000	3860458	NC		6106	
D 12 13C4 PFOA	416.5 > 371.6	10.762	10.734	0.028		1477697		69.4	2215	
13 Perfluorooctanoic acid	412.8 > 368.8	10.762	10.734	0.028	1.000	3191541		110	3328	
	412.8 > 168.7	10.762	10.734	0.028	1.000	1130700	2.82(0.00-0.00)	110	1579	
D 16 13C4 PFOS	502.4 > 79.7	11.661	11.645	0.016		482668		99.7	1298	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.670	11.649	0.021	1.000	6709046		81.5	2708	
	498.3 > 98.2	11.670	11.649	0.021	1.000	4029707	1.66(0.00-0.00)	81.5	3419	
D 17 13C5 PFNA	467.5 > 422.6	11.687	11.671	0.016		1652409		79.4	2660	
18 Perfluorononanoic acid	462.5 > 418.6	11.696	11.671	0.025	1.000	6998423		101	3885	
D 19 13C2 PFDA	514.4 > 469.5	12.487	12.465	0.022		2063584		70.3	3127	
20 Perfluorodecanoic acid	512.5 > 468.5	12.487	12.469	0.018	1.000	7801447		103	6956	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	13.004	12.984	0.020	1.000	15851344		108	2457	
D 23 13C8 FOSA	505.4 > 77.6	13.004	12.984	0.020		3872124		86.7	1580	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.102	13.094	0.008	1.000	2999356	NC		3842	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.102	13.094	0.008	1.000	2999356		80.0		
D 26 13C2 PFUnA	564.3 > 519.5	13.154	13.142	0.012		2083502		71.2	3105	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.154	13.142	0.012	1.000	8896158		103	3584	
D 28 13C2 PFDoA	614.4 > 569.4	13.721	13.709	0.012		2993831		84.2	2833	
29 Perfluorododecanoic acid	612.4 > 568.6	13.721	13.711	0.010	1.000	10070858		104	3017	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.197	14.189	0.008	1.000	8962091		89.7	2919	
D 33 13C2-PFTeDA	714.5 > 669.5	14.609	14.600	0.009		2783983		87.9	2979	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.609	14.601	0.008	1.000	4452926		98.2	1728	
D 35 13C2-PFHxDA	814.8 > 769.6	15.139	15.137	0.002		1505780		102	2855	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.144	15.137	0.007	1.000	12352725		95.6	4106	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7	> 868.6	15.441	15.442	-0.001	1.000	11471145	189.7	94.8	4361

### QC Flag Legend

Processing Flags

NC - Not Calibrated

### Reagents:

LCPFC-L6\_00015

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_008.d

Injection Date: 12-Feb-2016 12:21:25

Instrument ID: A4

Lims ID: Std L6

Client ID:

Operator ID: JRB

ALS Bottle#: 7

Worklist Smp#: 7

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

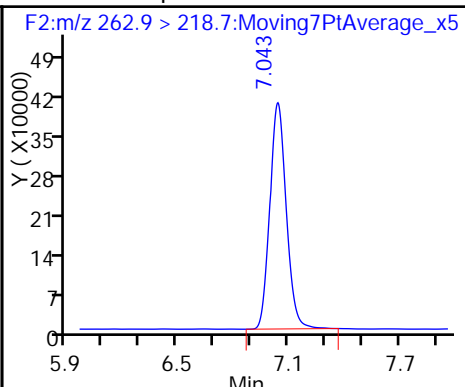
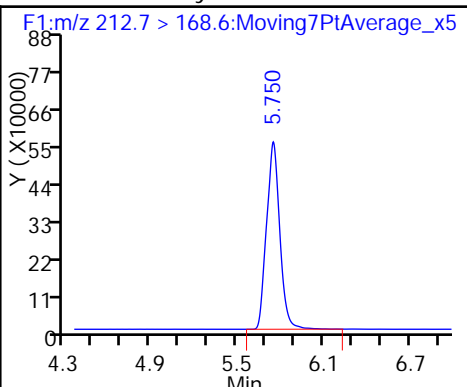
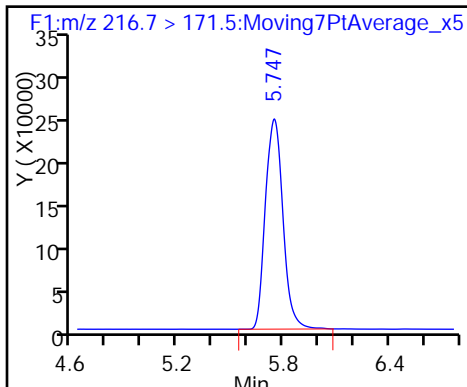
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

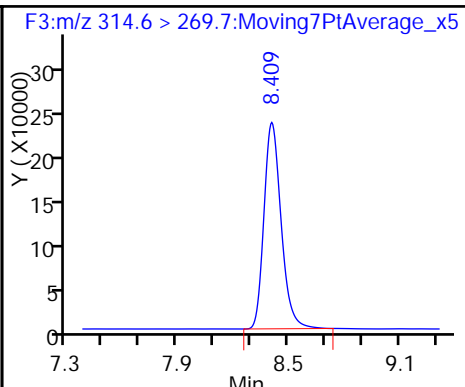
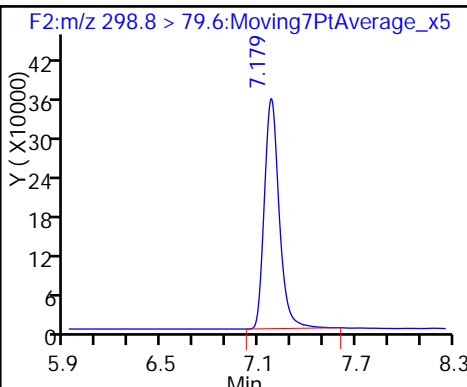
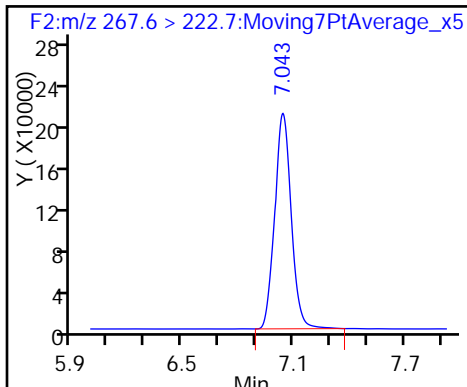
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

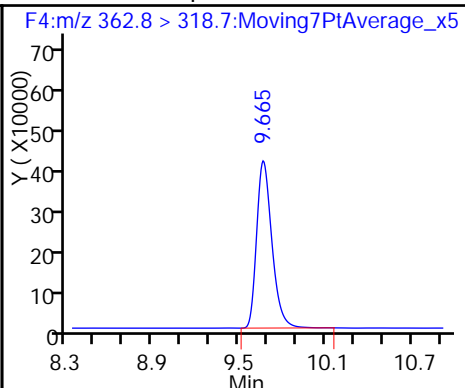
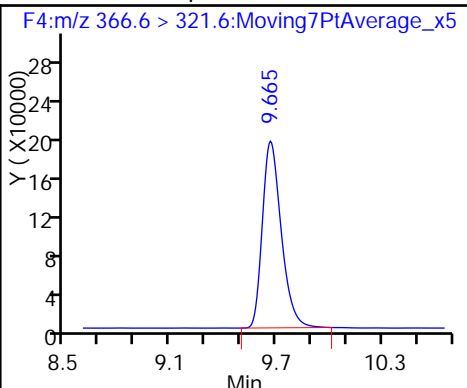
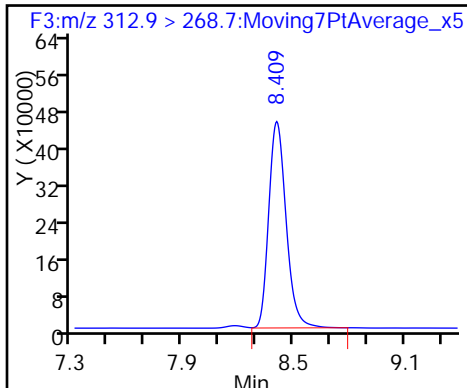
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

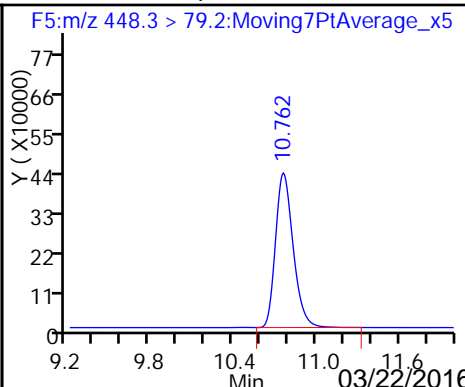
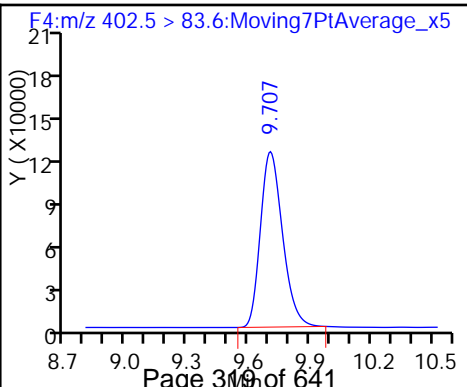
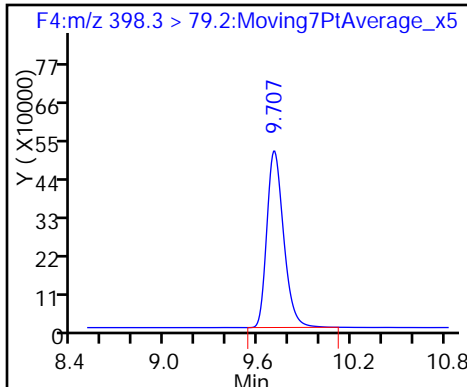
9 Perfluoroheptanoic acid



58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

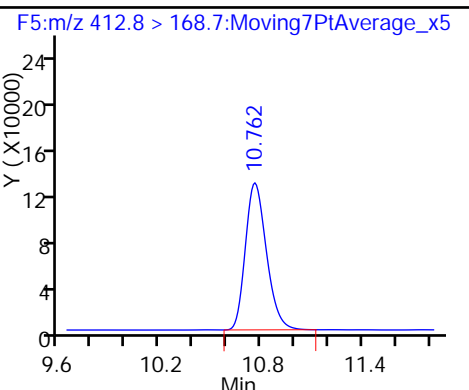
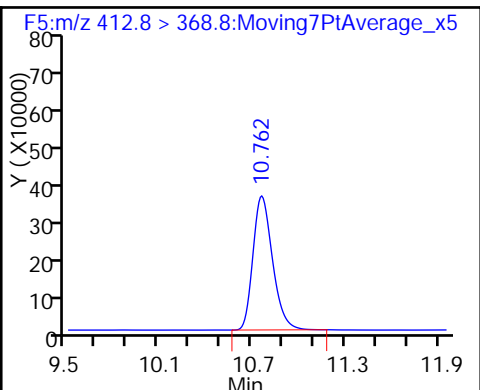
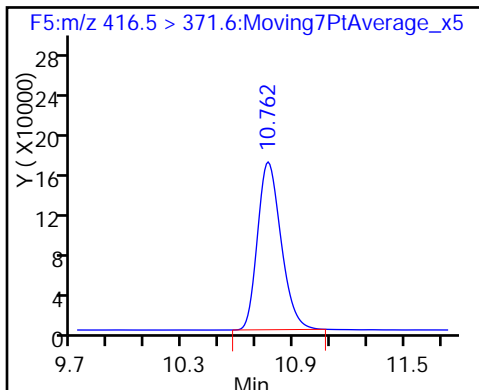
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

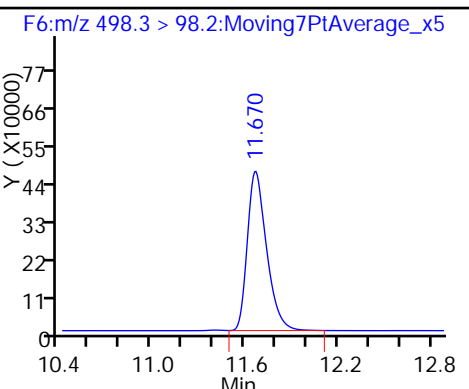
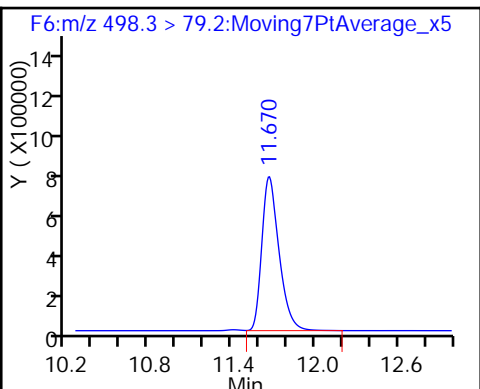
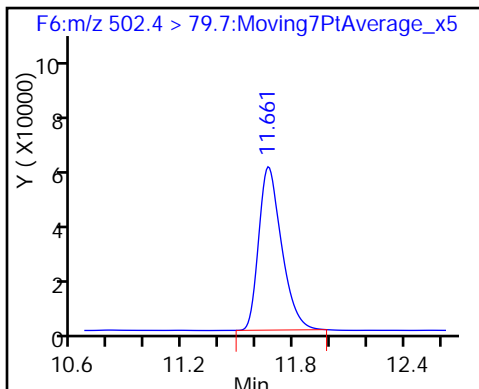
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid

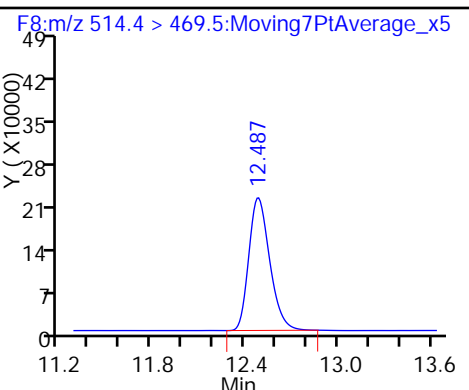
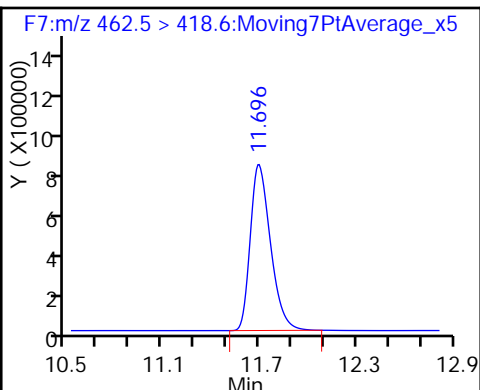
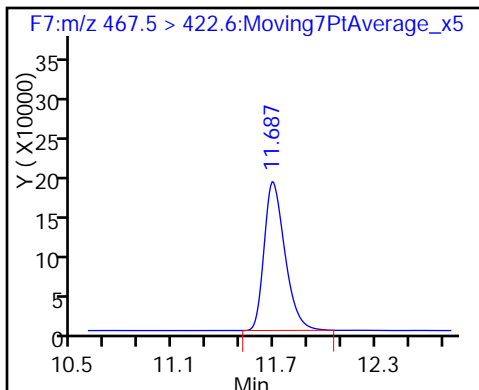
15 Perfluorooctane sulfonic acid



D 17 13C5 PFNA

18 Perfluorononanoic acid

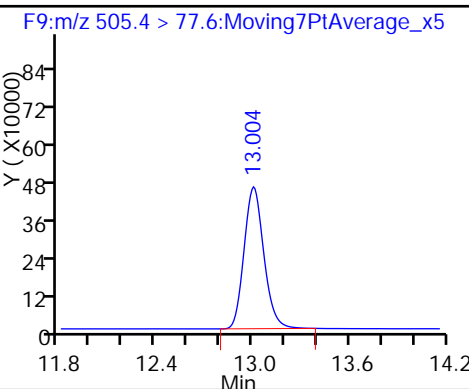
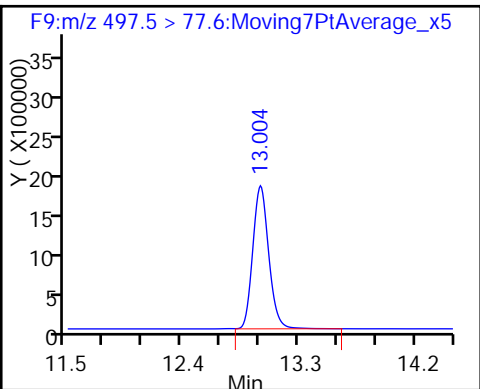
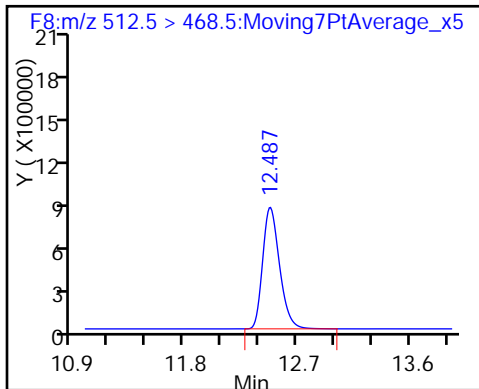
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

D 23 13C8 FOSA

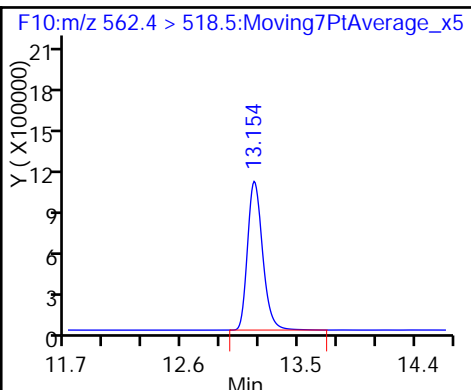
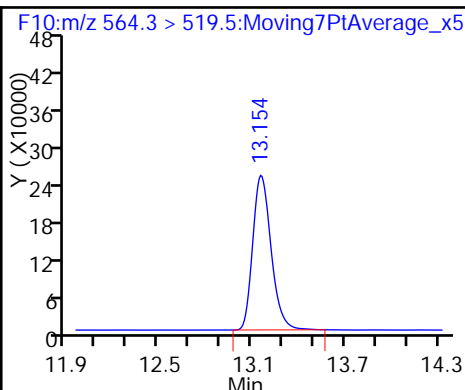
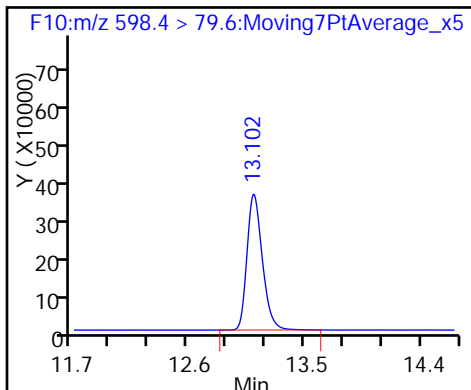




49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

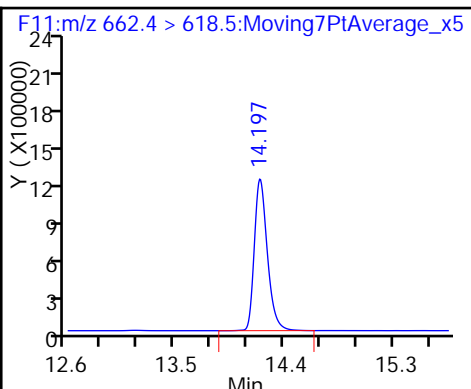
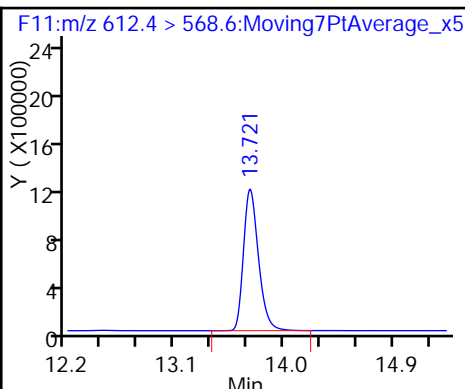
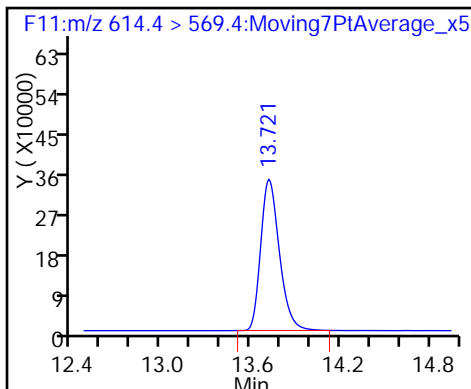
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

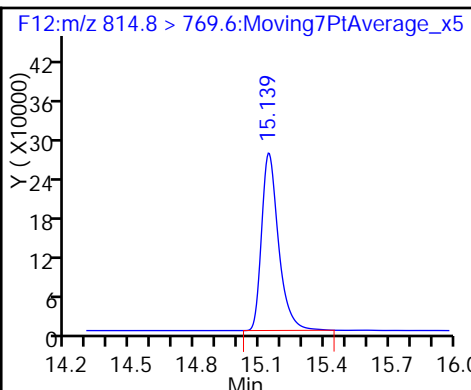
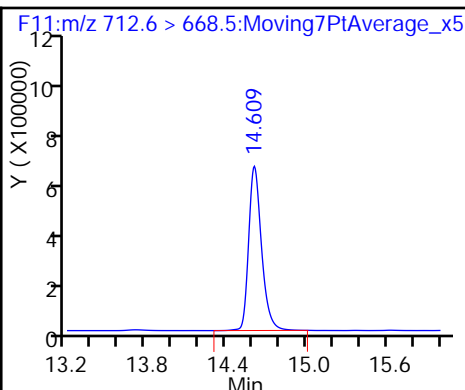
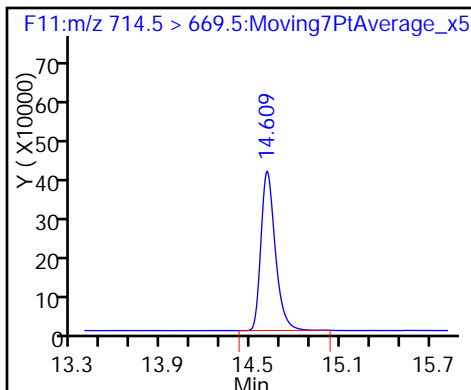
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

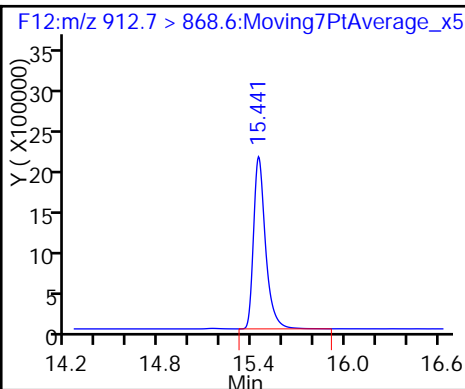
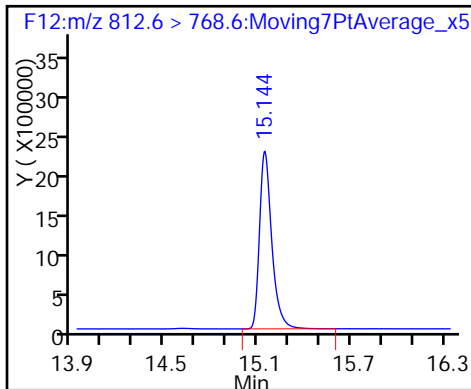
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d  
 Lims ID: Std L7  
 Client ID:  
 Sample Type: IC Calib Level: 7  
 Inject. Date: 12-Feb-2016 12:42:35 ALS Bottle#: 8 Worklist Smp#: 8  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L7  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12

Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:02:48 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 15:06:52

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	216.7 > 171.5	5.744	5.706	0.038	1453408	35.8		71.6	4360	
2 Perfluorobutyric acid	212.7 > 168.6	5.747	5.707	0.040	5966668	418.8		105	12387	
4 Perfluoropentanoic acid	262.9 > 218.7	7.039	6.977	0.062	4176286	403.9		101	2638	
D 3 13C5-PFPeA	267.6 > 222.7	7.039	6.978	0.061	1098104	34.3		68.6	1891	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.171	7.105	0.066	3250684	NC			2170	
	298.8 > 98.6	7.175	7.105	0.070	2095043		1.55(0.00-0.00)		1686	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.171	7.105	0.066	3250684	347.3		98.2		
D 6 13C2 PFHxA	314.6 > 269.7	8.414	8.360	0.054	1296336	33.1		66.2	2926	
7 Perfluorohexanoic acid	312.9 > 268.7	8.414	8.361	0.053	4700473	420.8		105	1630	
D 8 13C4-PFHpA	366.6 > 321.6	9.672	9.633	0.039	1287645	33.1		66.1	2298	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.672	9.634	0.038	4889892	370.8		92.7	6483	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.714	9.669	0.045	6095086	381.4		101		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.714	9.669	0.045	6095086	NC			10374	
D 11 18O2 PFHxS	402.5 > 83.6	9.714	9.669	0.045	679535	30.0		63.3	1556	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.771	10.735	0.036	1.000	5973185	321.0	84.3		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.771	10.735	0.036	1.000	5973185	NC		4795	
D 12 13C4 PFOA	416.5 > 371.6	10.771	10.738	0.033		1245478	29.3	58.5	1870	
13 Perfluorooctanoic acid	412.8 > 368.8	10.771	10.738	0.033	1.000	4435164	362.7	90.7	3425	
	412.8 > 168.7	10.771	10.738	0.033	1.000	1788026	2.48(0.00-0.00)	90.7	2827	
D 16 13C4 PFOS	502.4 > 79.7	11.670	11.648	0.022		349599	34.5	72.2	975	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.678	11.652	0.026	1.000	11018460	353.2	92.4	2325	
	498.3 > 98.2	11.678	11.652	0.026	1.000	6440011	1.71(0.00-0.00)	92.4	2904	
D 17 13C5 PFNA	467.5 > 422.6	11.704	11.674	0.030		1398930	33.6	67.3	2237	
18 Perfluorononanoic acid	462.5 > 418.6	11.704	11.676	0.028	1.000	11680686	397.6	99.4	5214	
D 19 13C2 PFDA	514.4 > 469.5	12.498	12.469	0.029		1703065	29.0	58.0	2597	
20 Perfluorodecanoic acid	512.5 > 468.5	12.498	12.472	0.026	1.000	12035423	386.5	96.6	4443	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	13.012	12.987	0.025	1.000	25297572	414.7	104	2260	
D 23 13C8 FOSA	505.4 > 77.6	13.012	12.987	0.025		3227314	36.1	72.3	1770	
25 Perfluorodecane Sulfonate	598.4 > 79.6	13.110	13.095	0.015	1.000	4169586	NC		3632	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	13.110	13.095	0.015	1.000	4169586	295.9	76.7		
D 26 13C2 PFUnA	564.3 > 519.5	13.163	13.144	0.019		1804563	30.8	61.7	2231	
27 Perfluoroundecanoic acid	562.4 > 518.5	13.163	13.144	0.019	1.000	14730079	393.5	98.4	3663	
D 28 13C2 PFDoA	614.4 > 569.4	13.728	13.711	0.017		2411510	33.9	67.8	2447	
29 Perfluorododecanoic acid	612.4 > 568.6	13.728	13.712	0.016	1.000	16007183	412.1	103	3036	
30 Perfluorotridecanoic acid	662.4 > 618.5	14.211	14.190	0.021	1.000	15038082	373.8	93.4	3203	
D 33 13C2-PFTeDA	714.5 > 669.5	14.614	14.602	0.012		2347410	37.0	74.1	2618	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.614	14.603	0.011	1.000	7695775	421.5	105	2631	
D 35 13C2-PFHxDA	814.8 > 769.6	15.143	15.138	0.005		1375507	46.4	92.8	3945	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.143	15.139	0.004	1.000	21395824	411.2	103	3796	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.444	15.442	0.002	1.000	20131167	412.6	103	4762	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L7\_00015

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d

Injection Date: 12-Feb-2016 12:42:35

Instrument ID: A4

Lims ID: Std L7

Client ID:

Operator ID: JRB

ALS Bottle#: 8

Worklist Smp#: 8

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

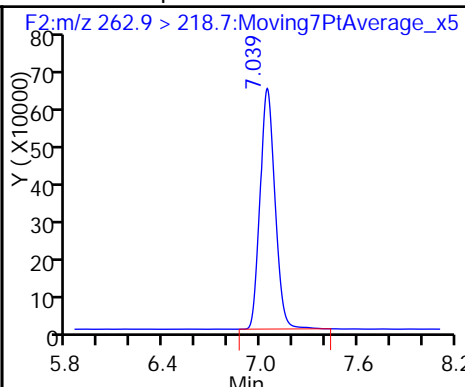
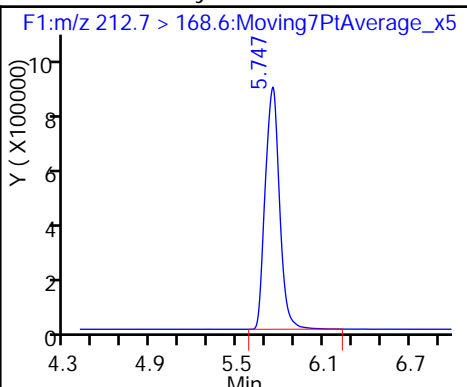
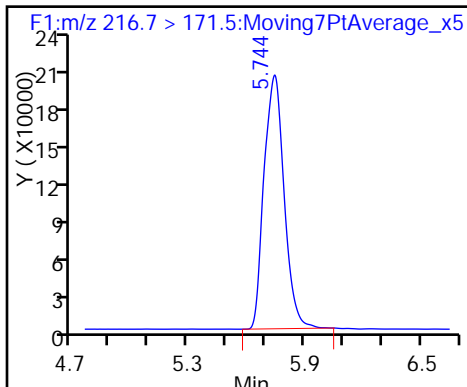
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

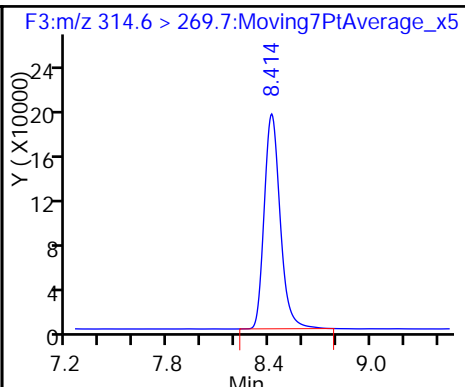
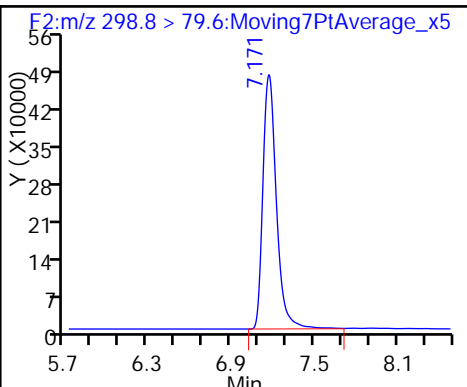
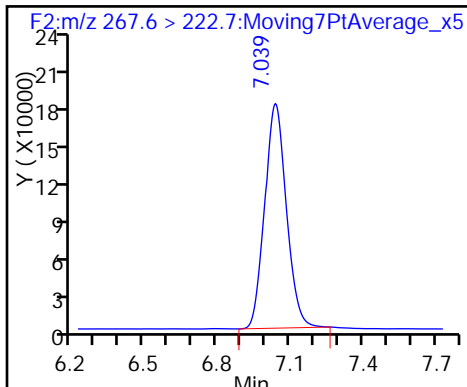
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

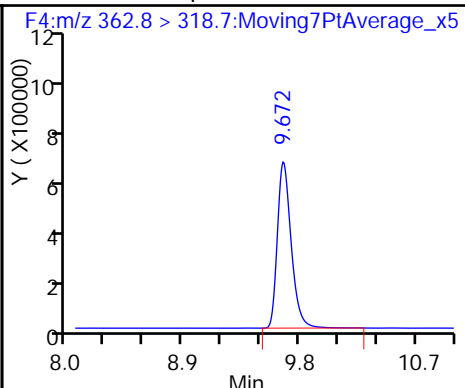
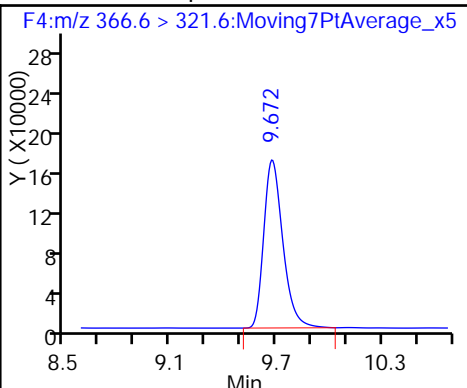
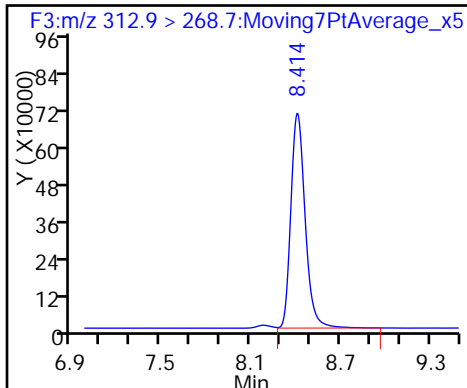
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

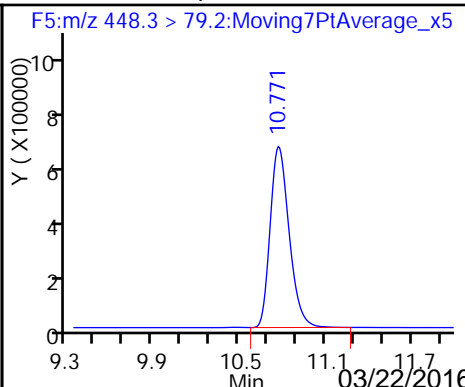
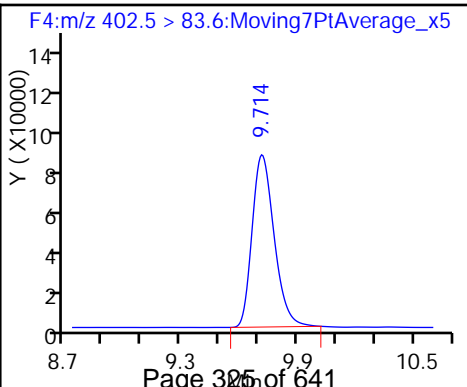
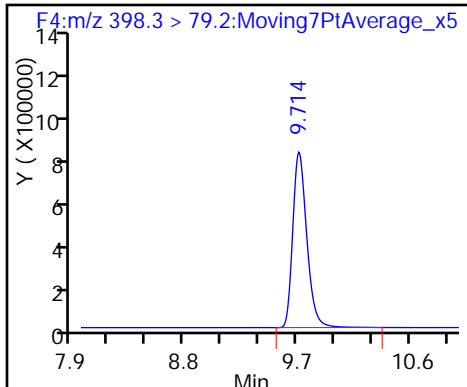
9 Perfluoroheptanoic acid



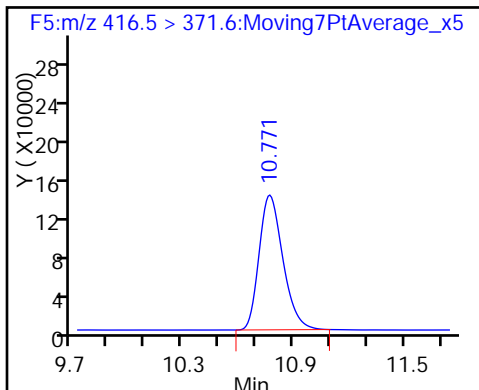
58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

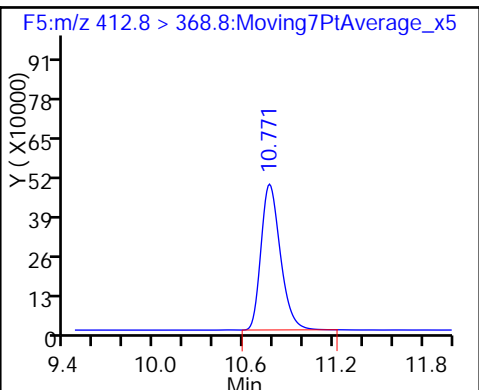
39 Perfluoroheptanesulfonic Acid



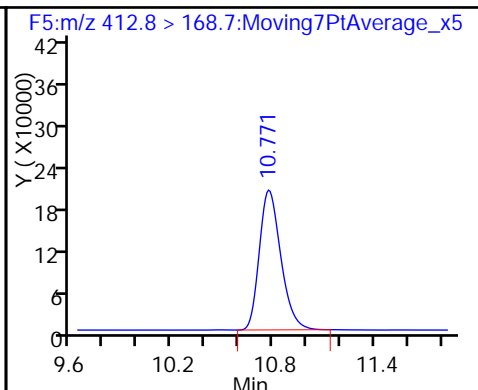
D 12 13C4 PFOA



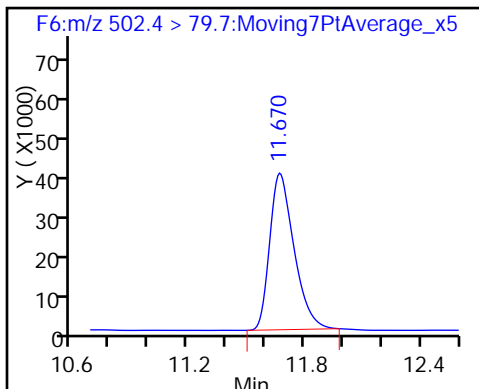
13 Perfluorooctanoic acid



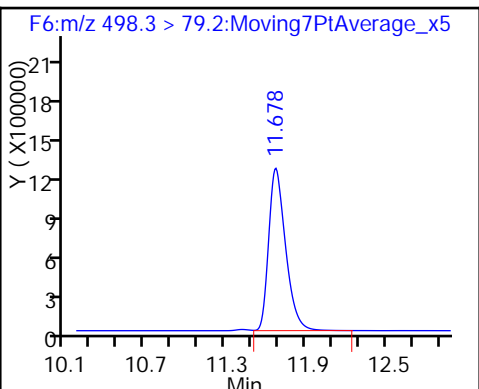
13 Perfluorooctanoic acid



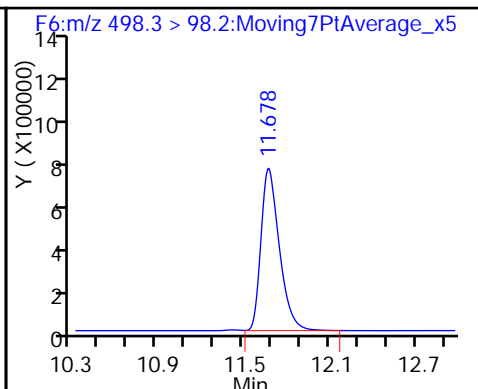
D 16 13C4 PFOS



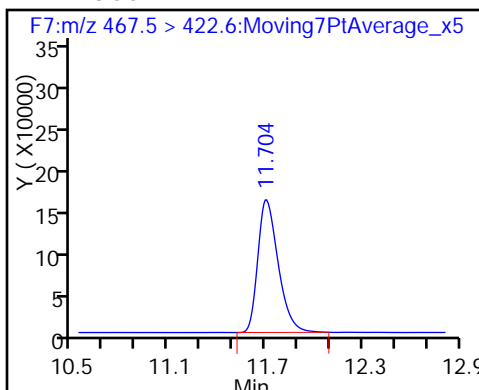
15 Perfluorooctane sulfonic acid



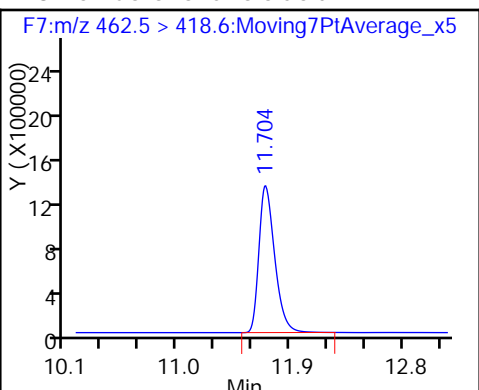
15 Perfluorooctane sulfonic acid



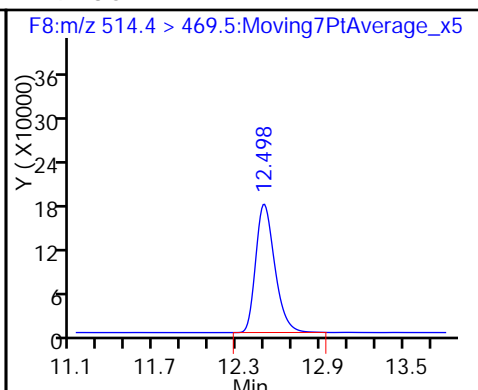
D 17 13C5 PFNA



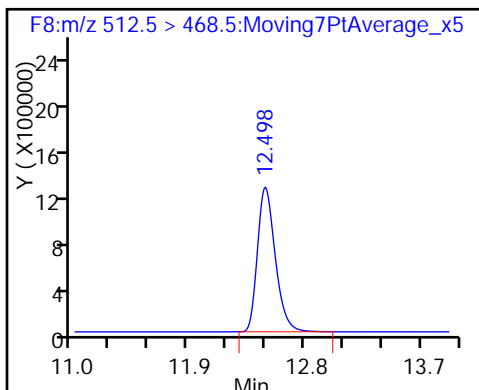
18 Perfluorononanoic acid



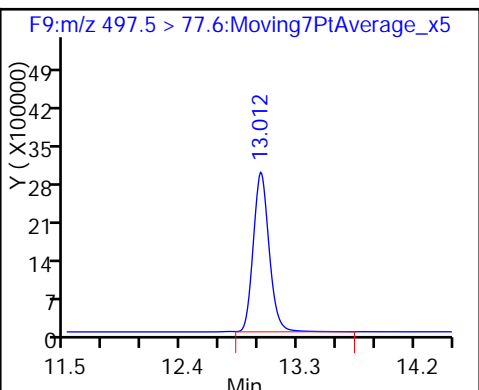
D 19 13C2 PFDA



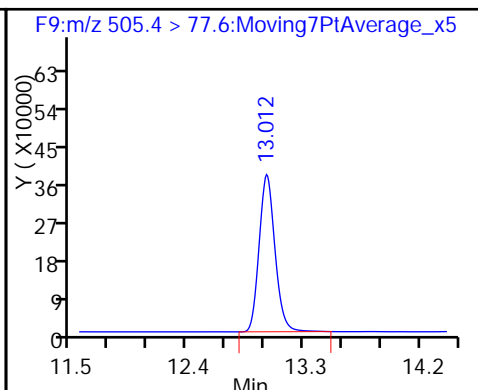
20 Perfluorodecanoic acid



24 Perfluorooctane Sulfonamide



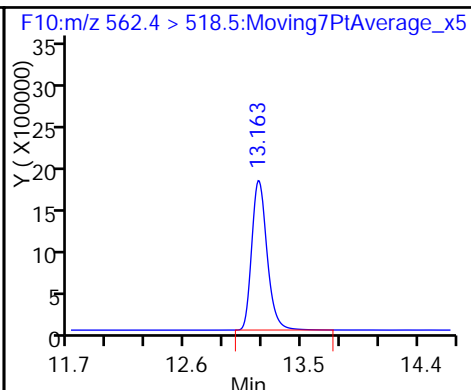
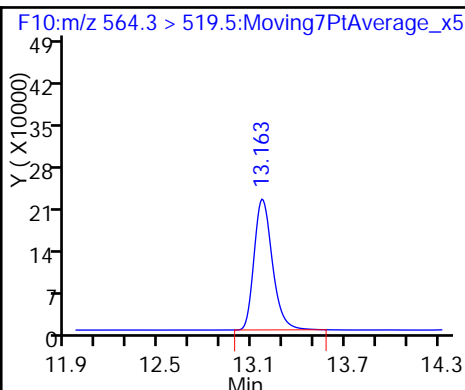
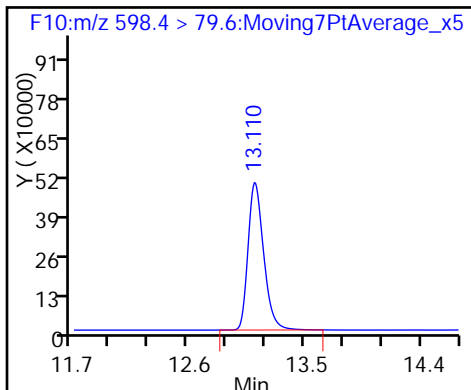
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

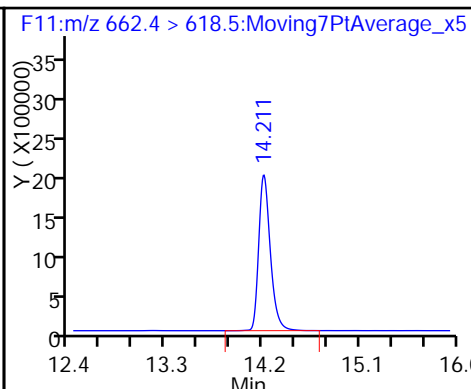
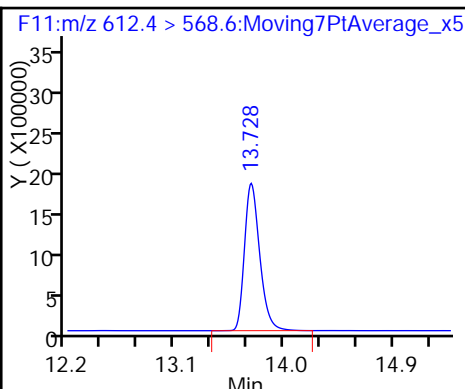
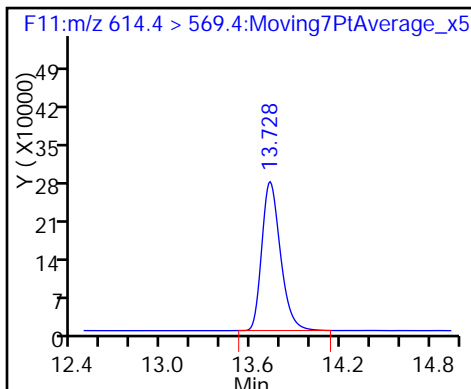
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

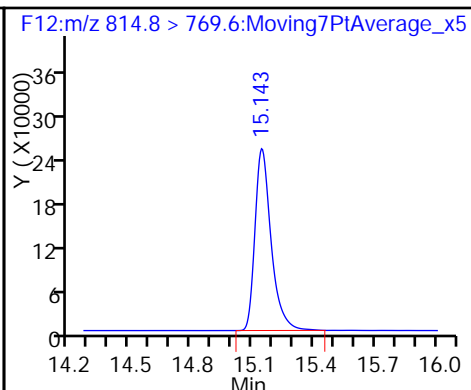
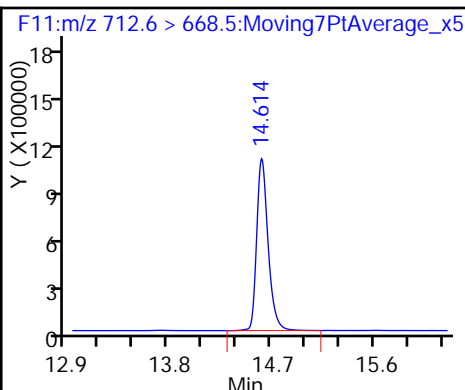
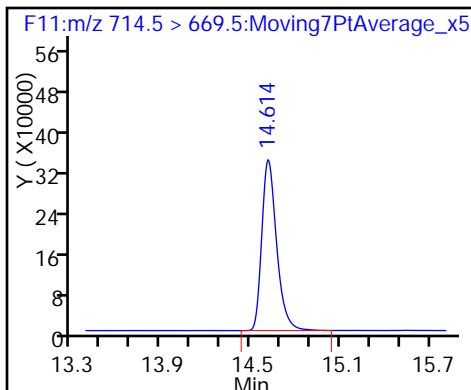
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

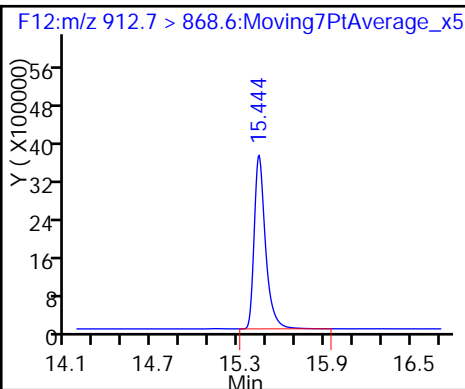
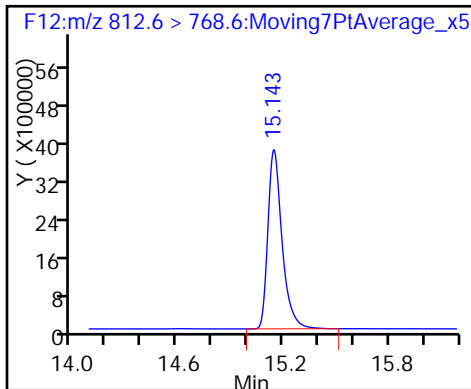
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VI  
 LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA  
 RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1 Analy Batch No.: 100906

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A6 GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/19/2016 15:08 Calibration End Date: 02/19/2016 17:15 Calibration ID: 19191

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100906/2	19FEB2016A6A_004.d
Level 2	STD 320-100906/3	19FEB2016A6A_005.d
Level 3	STD 320-100906/4	19FEB2016A6A_006.d
Level 4	STD 320-100906/5	19FEB2016A6A_007.d
Level 5	STD 320-100906/6	19FEB2016A6A_008.d
Level 6	STD 320-100906/7	19FEB2016A6A_009.d
Level 7	STD 320-100906/8	19FEB2016A6A_010.d

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7				RT WINDOW	AVG RT
Perfluorobutanoic acid (PFBA)	5.724	5.706	5.718	5.715	5.715	5.712	5.712				5.465 - 5.965	5.715
Perfluoropentanoic acid (PFPeA)	++++	6.837	6.842	6.828	6.828	6.819	6.818				6.578 - 7.078	6.829
Perfluorobutanesulfonic acid (PFBS)	++++	6.943	6.939	6.942	6.938	6.934	++++				6.689 - 7.189	6.939
Perfluorohexanoic acid (PFHxA)	++++	8.087	8.084	8.082	8.077	8.067	8.065				7.828 - 8.328	8.077
Perfluoroheptanoic acid (PFHpA)	++++	9.334	9.328	9.323	9.323	9.310	9.305				9.072 - 9.572	9.321
Perfluorohexanesulfonic acid (PFHxS)	++++	9.387	9.363	9.358	9.364	9.346	++++				9.114 - 9.614	9.364
Perfluorooctanoic acid (PFOA)	10.453	10.461	10.453	10.454	10.454	10.439	++++				10.200 - 10.700	10.452
Perfluoroheptanesulfonic Acid (PFHpS)	++++	10.475	10.461	10.461	10.461	10.439	++++				10.207 - 10.707	10.459
Perfluorooctanesulfonic acid (PFOS)	++++	11.436	11.420	11.421	11.422	11.399	++++				11.167 - 11.667	11.420
Perfluorononanoic acid (PFNA)	++++	11.451	11.443	11.437	11.437	11.421	11.422				11.189 - 11.689	11.435
Perfluorodecanoic acid (PFDA)	++++	12.296	12.286	12.276	12.280	12.254	++++				12.026 - 12.526	12.278
Perfluorooctane Sulfonamide (FOSA)	++++	12.824	12.814	12.814	12.818	12.793	12.793				12.562 - 13.062	12.809
Perfluorodecane Sulfonic acid	++++	12.970	12.960	12.950	12.954	12.929	++++				12.699 - 13.199	12.953
Perfluoroundecanoic acid (PFUnA)	13.010	13.010	13.010	13.002	13.005	12.981	++++				12.750 - 13.250	13.003
Perfluorododecanoic acid (PFDoA)	13.625	13.624	13.615	13.615	13.620	13.588	13.588				13.361 - 13.861	13.611
Perfluorotetradecanoic Acid (PFTriA)	14.147	14.139	14.139	14.132	14.136	14.109	14.109				13.880 - 14.380	14.130
Perfluorotetradecanoic acid (PFTeA)	14.585	14.585	14.578	14.572	14.576	14.552	14.552				14.321 - 14.821	14.571
Perfluoro-n-hexadecanoic acid (PFHxDA)	++++	15.218	15.213	15.213	15.210	15.193	15.193				14.958 - 15.458	15.207
Perfluoro-n-octandecanoic acid (PFODA)	15.557	15.549	15.544	15.544	15.546	15.529	15.529				15.293 - 15.793	15.543
13C4 PFBA	5.718	5.718	5.712	5.709	5.709	5.709	5.709				5.462 - 5.962	5.712
13C5-PFPeA	6.832	6.837	6.828	6.828	6.823	6.819	6.818				6.577 - 7.077	6.826
13C2 PFHxA	8.087	8.093	8.084	8.082	8.082	8.067	8.065				7.830 - 8.330	8.080
13C4-PFHpA	9.328	9.334	9.328	9.323	9.323	9.310	9.305				9.072 - 9.572	9.322
18O2 PFHxS	9.363	9.369	9.363	9.358	9.358	9.346	++++				9.108 - 9.608	9.360
13C4 PFOA	10.453	10.468	10.453	10.454	10.454	10.432	++++				10.200 - 10.700	10.452
13C4 PFOS	11.421	11.428	11.420	11.414	11.422	11.399	++++				11.164 - 11.664	11.417
13C5 PFNA	11.444	11.451	11.443	11.437	11.437	11.414	11.415				11.184 - 11.684	11.434
13C2 PFDA	12.286	12.286	12.286	12.276	12.280	12.254	++++				12.025 - 12.525	12.278
13C8 FOSA	12.814	12.824	12.814	12.814	12.818	12.793	12.793				12.560 - 13.060	12.810
13C2 PFUnA	13.010	13.010	13.010	13.002	13.005	12.981	++++				12.750 - 13.250	13.003
13C2 PFDoA	13.616	13.624	13.615	13.615	13.620	13.588	13.588				13.359 - 13.859	13.609
13C2-PFTeDA	14.579	14.585	14.578	14.576	14.576	14.552	14.552				14.321 - 14.821	14.572
13C2-PFHxDA	15.216	15.218	15.213	15.213	15.210	15.188	15.193				14.957 - 15.457	15.207



FORM VI  
LCMS BY EXTERNAL STANDARD - INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1 Analy Batch No.: 100906

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A6 GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/19/2016 15:08 Calibration End Date: 02/19/2016 17:15 Calibration ID: 19191

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100906/2	19FEB2016A6A_004.d
Level 2	STD 320-100906/3	19FEB2016A6A_005.d
Level 3	STD 320-100906/4	19FEB2016A6A_006.d
Level 4	STD 320-100906/5	19FEB2016A6A_007.d
Level 5	STD 320-100906/6	19FEB2016A6A_008.d
Level 6	STD 320-100906/7	19FEB2016A6A_009.d
Level 7	STD 320-100906/8	19FEB2016A6A_010.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R <sup>2</sup> OR COD	#	MIN R <sup>2</sup> OR COD
	LVL 1 LVL 5	LVL 2 LVL 6	LVL 3 LVL 7	LVL 4		B	M1	M2								
13C4 PFBA	17215 15573	18158 14290	17652 12637	16479	Ave		16000.4800			12.4		50.0				
13C5-PFPeA	33132 32092	37596 25883	31200 23439	31979	Ave		30760.1743			15.3		50.0				
13C2 PFHxA	31151 28249	32712 25751	29384 20807	30249	Ave		28328.9429			14.1		50.0				
13C4-PFHpA	33720 29961	37160 26279	32774 21986	33895	Ave		30825.1286			16.8		50.0				
18O2 PFHxS	16445 14414	15937 12069	15300 ++++	15633	Ave		14966.2297			10.5		50.0				
13C4 PFOA	36851 31471	39431 26220	35009 ++++	35877	Ave		34143.1967			13.7		50.0				
13C4 PFOS	19998 17191	20738 13980	18794 ++++	18393	Ave		18182.3013			13.2		50.0				
13C5 PFNA	32640 27720	35233 22861	29495 20201	29189	Ave		28191.1457			18.6		50.0				
13C2 PFDA	30433 26439	32141 20971	31311 ++++	30484	Ave		28629.7300			14.8		50.0				
13C8 FOSA	45384 42557	44125 37994	44477 31983	43988	Ave		41501.0343			11.7		50.0				
13C2 PFUnA	40172 34391	42368 28219	39609 ++++	39503	Ave		37377.2300			13.9		50.0				
13C2 PFDoA	43990 39123	44852 34183	40349 29026	42305	Ave		39118.3029			14.6		50.0				
13C2-PFTeDA	37943 36827	39228 31852	37299 30439	37717	Ave		35900.7657			9.3		50.0				
13C2-PFHxDA	45996 43682	46879 39154	44036 36320	45657	Ave		43103.4314			9.1		50.0				

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

## CURVE EVALUATION

Lab Name: TestAmerica SacramentoJob No.: 320-17154-1Analy Batch No.: 100906SDG No.: CTO WE7G PFC SamplingInstrument ID: A6GC Column: Acquity ID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 02/19/2016 15:08Calibration End Date: 02/19/2016 17:15Calibration ID: 19191

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R <sup>2</sup> OR COD	#	MIN R <sup>2</sup> OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
	LVL 6	LVL 7															
Perfluorobutanoic acid (PFBA)	17268 20409	18332 18204	24759	23146	22421	AveID		1.3041			15.6		35.0				
Perfluoropentanoic acid (PFPeA)	++++ 26519	41667 24011	33706	33699	31098	AveID		1.0434			4.7		35.0				
Perfluorobutanesulfonic acid (PFBS)	++++ 14042	11589 ++++	13996	16675	15750	AveID		0.9930			17.5		50.0				
Perfluorohexanoic acid (PFHxA)	++++ 28593	29048 23704	35504	30900	31979	AveID		1.0832			10.4		35.0				
Perfluoroheptanoic acid (PFHpA)	++++ 26755	35173 22997	33430	38937	32292	AveID		1.0429			6.5		35.0				
Perfluorohexanesulfonic acid (PFHxS)	++++ 8511.3	10001 ++++	10164	10841	9886.6	AveID		0.6753			4.5		35.0				
Perfluorooctanoic acid (PFOA)	41176 26427	32208 ++++	37796	35759	31611	AveID		1.0038			10.3		35.0				
Perfluoroheptanesulfonic Acid (PFHpS)	++++ 8507.0	9112.4 ++++	11112	11362	10932	L2ID	-0.177	0.6259						1.0000		0.9900	
Perfluorooctanesulfonic acid (PFOS)	++++ 13560	19579 ++++	18099	18656	16687	AveID		0.9724			2.6		35.0				
Perfluorononanoic acid (PFNA)	++++ 19794	23017 16627	24079	25918	22126	AveID		0.8075			10.2		35.0				
Perfluorodecanoic acid (PFDA)	++++ 20747	42554 ++++	32768	28572	26974	L2ID	0.3477	0.9748						0.9990		0.9900	
Perfluorooctane Sulfonamide (FOSA)	++++ 33984	42097 29306	40773	41641	39679	AveID		0.9268			2.4		35.0				
Perfluorodecane Sulfonic acid	++++ 7589.1	6534.2 ++++	10246	10522	10316	L1ID	-0.161	0.5576						0.9980		0.9900	
Perfluoroundecanoic acid (PFUnA)	76472 22856	51520 ++++	42693	34710	30224	L1ID	0.5568	0.8236						0.9980		0.9900	
Perfluorododecanoic acid (PFDoA)	28216 25891	23692 22968	32661	33950	30803	AveID		0.7311			14.6		35.0				
Perfluorotridecanoic Acid (PFTriA)	34658 29612	40040 23925	48646	43762	38156	AveID		0.9409			15.3		50.0				
Perfluorotetradecanoic acid (PFTeA)	49430 19999	42000 17283	27015	24942	23186	L2ID	0.2791	0.5968						0.9970		0.9900	
Perfluoro-n-hexadecanoic acid (PFHxDA)	++++ 37534	171017 32169	64991	49814	43976	L2ID	2.7370	1.0725						1.0000		0.9900	
Perfluoro-n-octadecanoic acid (PFODA)	49252 37638	39408 34132	49070	48056	44933	AveID		1.1108			9.8		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.: 320-17154-1 Analy Batch No.: 100906

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A6 GC Column: Acquity ID: 2.1(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/19/2016 15:08 Calibration End Date: 02/19/2016 17:15 Calibration ID: 19191

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100906/2	19FEB2016A6A_004.d
Level 2	STD 320-100906/3	19FEB2016A6A_005.d
Level 3	STD 320-100906/4	19FEB2016A6A_006.d
Level 4	STD 320-100906/5	19FEB2016A6A_007.d
Level 5	STD 320-100906/6	19FEB2016A6A_008.d
Level 6	STD 320-100906/7	19FEB2016A6A_009.d
Level 7	STD 320-100906/8	19FEB2016A6A_010.d

ANALYTE	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
		LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
13C4 PFBA	Ave	860754 714491	907904 631840	882584	823965	778630	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C5-PFPeA	Ave	1656599 1294135	1879822 1171928	1560021	1598957	1604599	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFHxA	Ave	1557529 1287540	1635595 1040351	1469185	1512461	1412469	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C4-PFHpA	Ave	1685988 1313950	1858023 1099321	1638704	1694741	1498068	50.0 50.0	50.0 50.0	50.0	50.0	50.0
1802 PFHxS	Ave	777834 570842	753810 ++++	723683	739449	681798	47.3 47.3	47.3 ++++	47.3	47.3	47.3
13C4 PFOA	Ave	1842532 1311023	1971561 ++++	1750451	1793859	1573533	50.0 50.0	50.0 ++++	50.0	50.0	50.0
13C4 PFOS	Ave	955905 668257	991283 ++++	898344	879178	821717	47.8 47.8	47.8 ++++	47.8	47.8	47.8
13C5 PFNA	Ave	1631991 1143055	1761627 1010032	1474756	1459436	1386004	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFDA	Ave	1521664 1048548	1607037 ++++	1565532	1524195	1321943	50.0 50.0	50.0 ++++	50.0	50.0	50.0
13C8 FOSA	Ave	2269224 1899691	2206225 1599127	2223851	2199385	2127859	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFUnA	Ave	2008621 1410933	2118412 ++++	1980471	1975170	1719562	50.0 50.0	50.0 ++++	50.0	50.0	50.0
13C2 PFDoA	Ave	2199484 1709160	2242621 1451296	2017427	2115244	1956174	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2-PFTeDA	Ave	1897170 1592616	1961399 1521958	1864931	1885833	1841361	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2-PFHxDA	Ave	2299786 1957713	2343949 1815999	2201810	2282833	2184111	50.0 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.: 320-17154-1Analy Batch No.: 100906SDG No.: CTO WE7G PFC SamplingInstrument ID: A6GC Column: AcquityID: 2.1(mm)Heated Purge: (Y/N) NCalibration Start Date: 02/19/2016 15:08Calibration End Date: 02/19/2016 17:15Calibration ID: 19191

## Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-100906/2	19FEB2016A6A_004.d
Level 2	STD 320-100906/3	19FEB2016A6A_005.d
Level 3	STD 320-100906/4	19FEB2016A6A_006.d
Level 4	STD 320-100906/5	19FEB2016A6A_007.d
Level 5	STD 320-100906/6	19FEB2016A6A_008.d
Level 6	STD 320-100906/7	19FEB2016A6A_009.d
Level 7	STD 320-100906/8	19FEB2016A6A_010.d

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
Perfluorobutanoic acid (PFBA)		AveID	8634 4081745	18332 7281727	123793	462922	1121027	0.500 200	1.00 400	5.00	20.0	50.0
Perfluoropentanoic acid (PFPeA)		AveID	++++ 5303760	41667 9604476	168531	673970	1554916	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorobutanesulfonic acid (PFBS)		AveID	++++ 2482631	10245 ++++	61862	294821	696169	++++ 177	0.884 ++++	4.42	17.7	44.2
Perfluorohexanoic acid (PFHxA)		AveID	++++ 5718698	29048 9481665	177522	618008	1598966	++++ 200	1.00 400	5.00	20.0	50.0
Perfluoroheptanoic acid (PFHpA)		AveID	++++ 5350969	35173 9198751	167149	778743	1614588	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorohexanesulfonic acid (PFHxS)		AveID	++++ 1610335	9461 ++++	48077	205104	467634	++++ 189	0.946 ++++	4.73	18.9	47.3
Perfluorooctanoic acid (PFOA)		AveID	20588 5285370	32208 ++++	188978	715180	1580565	0.500 200	1.00 ++++	5.00	20.0	50.0
Perfluoroheptanesulfonic Acid (PFHpS)		L2ID	++++ 1619742	8675 ++++	52893	216336	520355	++++ 190	0.952 ++++	4.76	19.0	47.6
Perfluorooctanesulfonic acid (PFOS)		AveID	++++ 2592636	18718 ++++	86511	356703	797626	++++ 191	0.956 ++++	4.78	19.1	47.8
Perfluorononanoic acid (PFNA)		AveID	++++ 3958780	23017 6650814	120395	518362	1106311	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorodecanoic acid (PFDA)		L2ID	++++ 4149384	42554 ++++	163841	571448	1348689	++++ 200	1.00 ++++	5.00	20.0	50.0
Perfluorooctane Sulfonamide (FOSA)		AveID	++++ 6796748	42097 11722467	203867	832823	1983962	++++ 200	1.00 400	5.00	20.0	50.0
Perfluorodecane Sulfonic acid		L1ID	++++ 1463184	6299 ++++	49387	202871	497252	++++ 193	0.964 ++++	4.82	19.3	48.2
Perfluoroundecanoic acid (PFUnA)		L1ID	38236 4571131	51520 ++++	213464	694202	1511190	0.500 200	1.00 ++++	5.00	20.0	50.0
Perfluorododecanoic acid (PFDoA)		AveID	14108 5178123	23692 9187053	163307	679009	1540137	0.500 200	1.00 400	5.00	20.0	50.0

## RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1 Analy Batch No.: 100906SDG No.: CTO WE7G PFC SamplingInstrument ID: A6 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) NCalibration Start Date: 02/19/2016 15:08 Calibration End Date: 02/19/2016 17:15 Calibration ID: 19191

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (NG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5
Perfluorotridecanoic Acid (PFTriA)		AveID	17329 5922305	40040 9569869	243228	875238	1907812	0.500 200	1.00 400	5.00	20.0	50.0
Perfluorotetradecanoic acid (PFTeA)		L2ID	24715 3999720	42000 6913347	135077	498848	1159290	0.500 200	1.00 400	5.00	20.0	50.0
Perfluoro-n-hexadecanoic acid (PFHxDA)		L2ID	++++ 7506851	171017 12867499	324953	996289	2198803	++++ 200	1.00 400	5.00	20.0	50.0
Perfluoro-n-octadecanoic acid (PFODA)		AveID	24626 7527521	39408 13652743	245350	961125	2246657	0.500 200	1.00 400	5.00	20.0	50.0

## Curve Type Legend:

AveID = Average isotope dilution L1ID = Linear 1/conc IsoDil L2ID = Linear 1/conc^2 IsoDil
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TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_004.d  
 Lims ID: Std L1  
 Client ID:  
 Sample Type: IC Calib Level: 1  
 Inject. Date: 19-Feb-2016 15:08:21 ALS Bottle#: 9 Worklist Smp#: 2  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L1  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5

Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:40 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d

Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA										
217.0 > 172.0	5.718	5.712	0.006		860754	53.8		108	3626	
2 Perfluorobutyric acid										
212.9 > 169.0	5.724	5.715	0.009	1.000	8634	0.3846		76.9	31.2	
D 3 13C5-PFPeA										
267.9 > 223.0	6.832	6.827	0.005		1656599	53.9		108	5445	
4 Perfluoropentanoic acid										
262.9 > 219.0	6.823	6.828	-0.005	1.000	24416	0.7063		141	4.9	
5 Perfluorobutane Sulfonate										
298.9 > 80.0	6.943	6.939	0.004	1.000	6549	NC			8.3	
298.9 > 99.0	6.947	6.939	0.008	1.001	2180		3.00(0.00-0.00)		6.8	
40 Perfluorobutanesulfonic acid										
298.9 > 80.0	6.943	6.939	0.004	1.000	6549	0.4011		90.7		
7 Perfluorohexanoic acid										
313.0 > 269.0	8.082	8.078	0.004	1.000	17559	0.5204		104	63.7	
D 6 13C2 PFHxA										
315.0 > 270.0	8.087	8.080	0.007		1557529	55.0		110	5408	
D 8 13C4-PFHpA										
367.0 > 322.0	9.328	9.322	0.006		1685988	54.7		109	6243	
9 Perfluoroheptanoic acid										
363.0 > 319.0	9.328	9.322	0.006	1.000	21816	0.6204		124	116	
D 11 18O2 PFHxS										
403.0 > 84.0	9.363	9.358	0.005		777834	52.0		110	2685	
10 Perfluorohexane Sulfonate										
399.0 > 80.0	9.381	9.364	0.017	1.000	6504	NC			21.0	
41 Perfluorohexanesulfonic acid										
399.0 > 80.0	9.381	9.364	0.017	1.000	6504	0.5857		124		
D 12 13C4 PFOA										
417.0 > 372.0	10.453	10.450	0.003		1842532	54.0		108	6072	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
413.0 > 369.0	10.453	10.450	0.003	1.000	20588	0.5566		111	21.4	
413.0 > 169.0	10.474	10.450	0.024	1.002	8191		2.51(0.00-0.00)	111	24.5	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.461	10.457	0.004	1.000	6248	NC			31.0	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.461	10.457	0.004	1.000	6248	0.7817		164		
D 16 13C4 PFOS										
503.0 > 80.0	11.421	11.414	0.007		955905	52.6		110	3294	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.421	11.417	0.004	1.000	10554	0.5427		114	37.9	
499.0 > 99.0	11.414	11.417	-0.003	0.999	3471		3.04(0.00-0.00)	114	16.8	
D 17 13C5 PFNA										
468.0 > 423.0	11.444	11.434	0.010		1631991	57.9		116	6715	
18 Perfluorononanoic acid										
463.0 > 419.0	11.458	11.439	0.019	1.000	11375	0.4316		86.3	38.1	
D 19 13C2 PFDA										
515.0 > 470.0	12.286	12.275	0.011		1521664	53.1		106	4154	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.286	12.276	0.010	1.000	28318	0.5978		120	92.0	
D 23 13C8 FOSA										
506.0 > 78.0	12.814	12.810	0.004		2269224	54.7		109	1690	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.825	12.812	0.013	1.000	16577	0.3941		78.8	56.9	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.950	12.949	0.001	1.000	2979	NC			5.8	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.950	12.949	0.001	1.000	2979	0.5554		115		
D 26 13C2 PFUnA										
565.0 > 520.0	13.010	13.000	0.010		2008621	53.7		107	5478	
27 Perfluoroundecanoic acid										
563.0 > 519.0	13.010	13.000	0.010	1.000	38236	0.4797		95.9	98.7	
D 28 13C2 PFDoA										
615.0 > 570.0	13.616	13.609	0.007		2199484	56.2		112	4878	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.625	13.611	0.014	1.000	14108	0.4387		87.7	18.9	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.147	14.130	0.017	1.000	17329	0.4187		83.7	16.0	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.585	14.571	0.014	1.000	24715	0.4738		94.8	13.2	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.579	14.571	0.008		1897170	52.8		106	6130	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.216	15.207	0.009		2299786	53.4		107	7006	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.216	15.208	0.008	1.000	178126	1.22		245	191	
36 Perfluorooctandecanoic acid										
913.0 > 869.0	15.557	15.543	0.014	1.000	24626	0.5040		101	32.9	

[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L1\_00018

Amount Added: 1.00

Units: mL



TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_004.d

Injection Date: 19-Feb-2016 15:08:21

Instrument ID: A6

Lims ID: Std L1

Client ID:

Operator ID: JRB

ALS Bottle#: 9

Worklist Smp#: 2

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

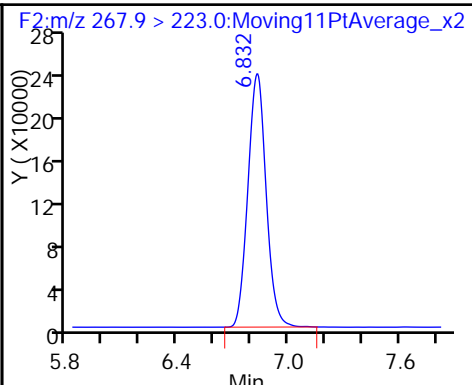
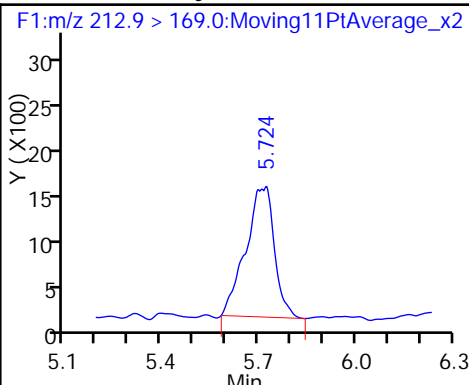
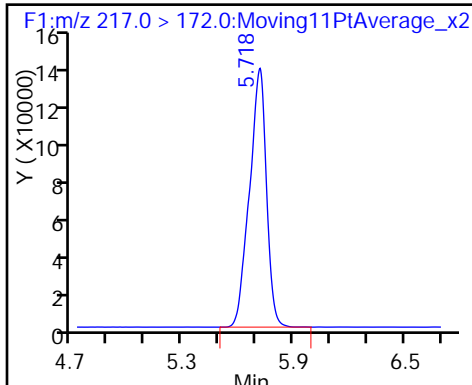
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

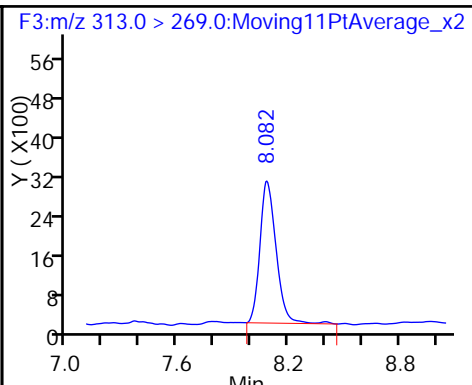
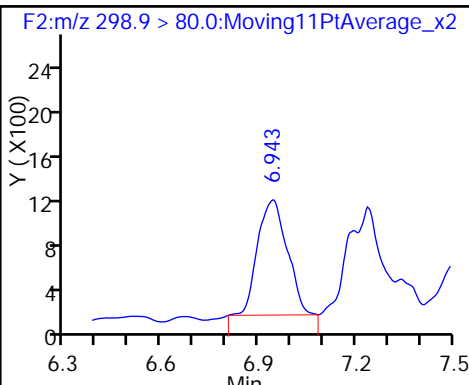
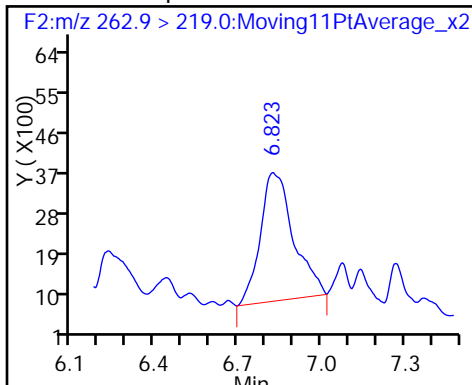
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

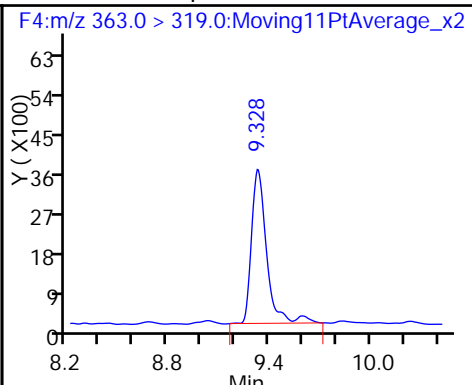
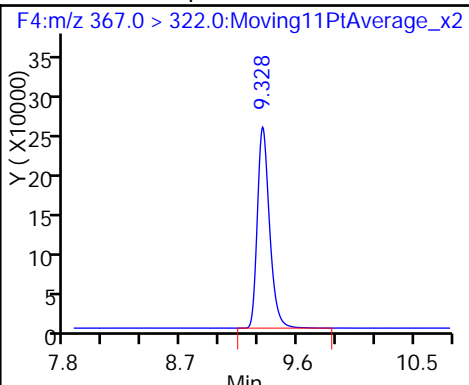
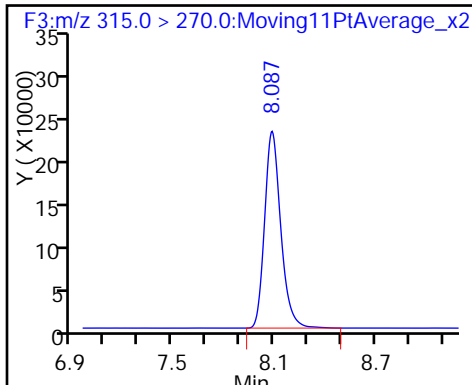
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

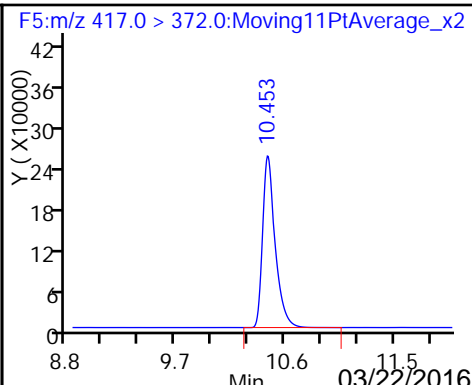
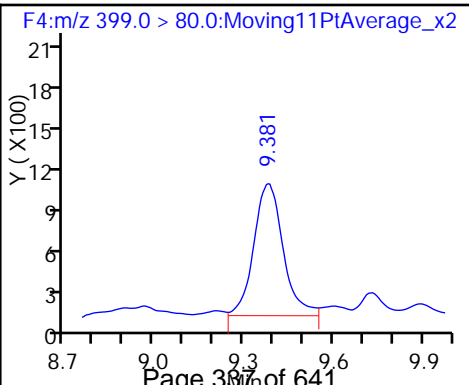
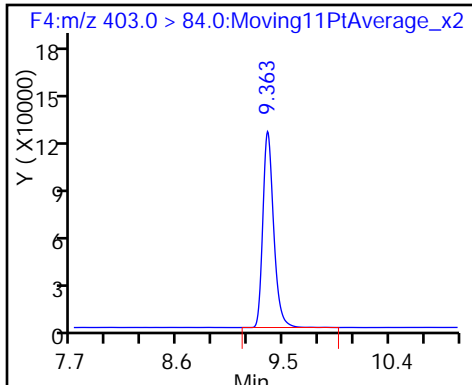
9 Perfluoroheptanoic acid

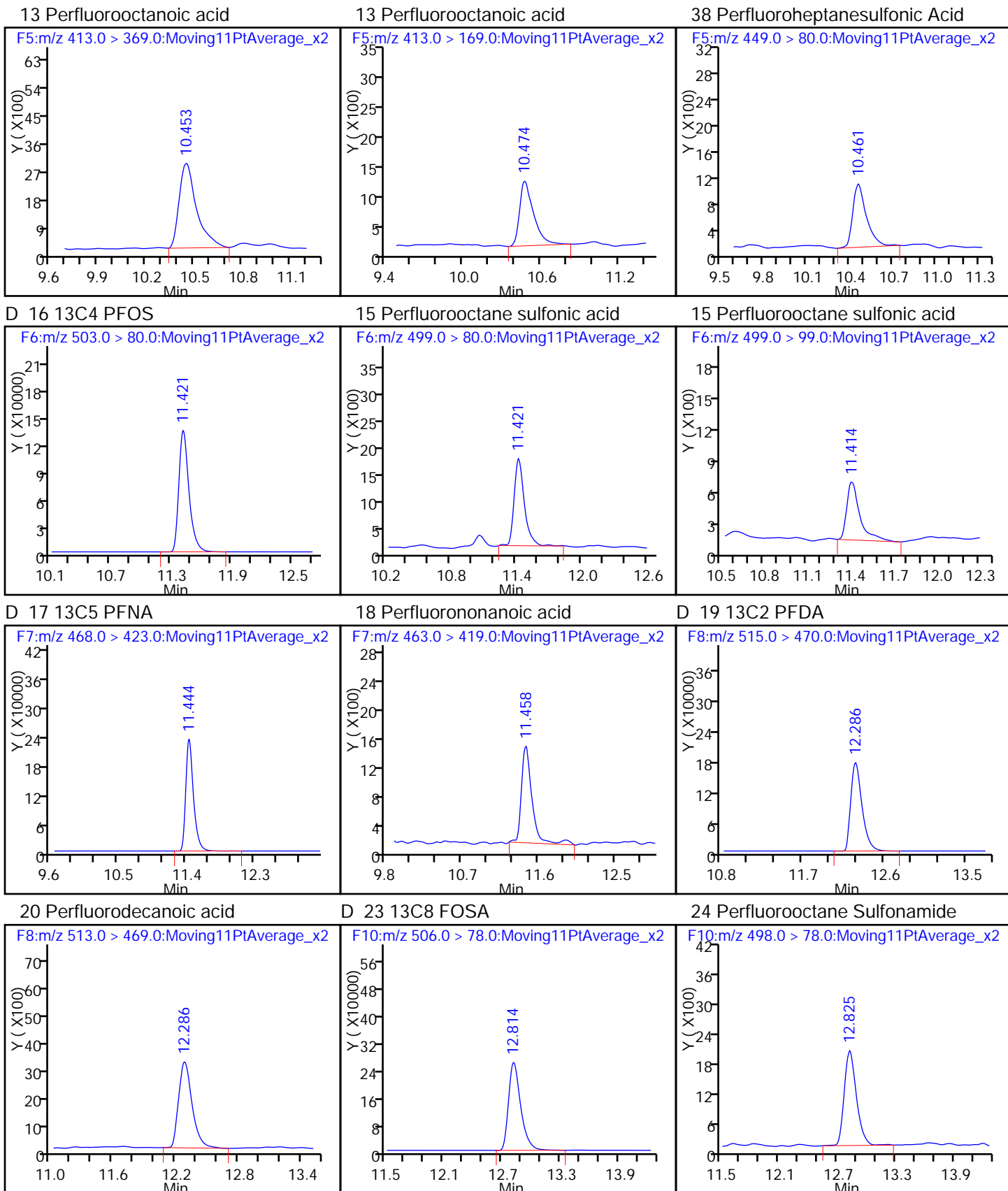


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

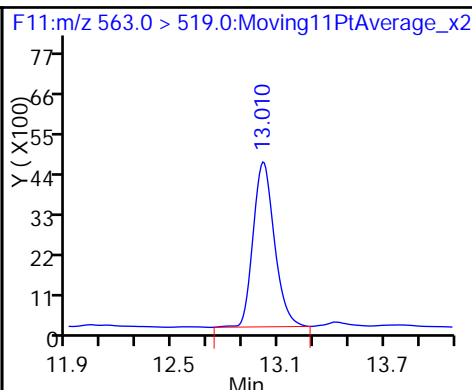
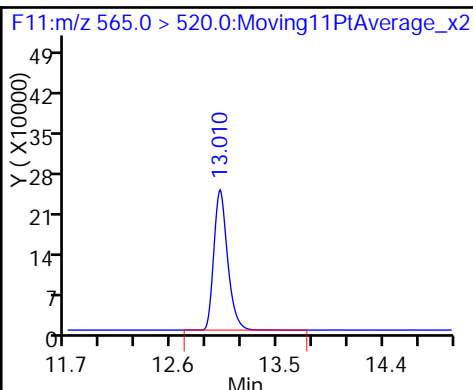
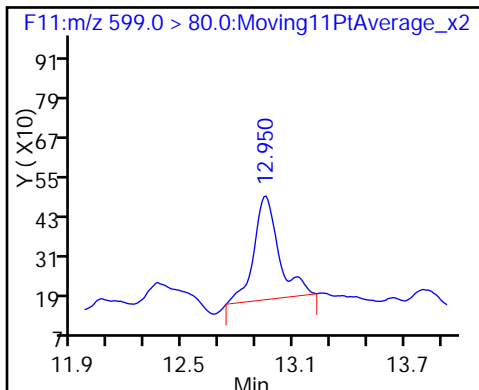




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

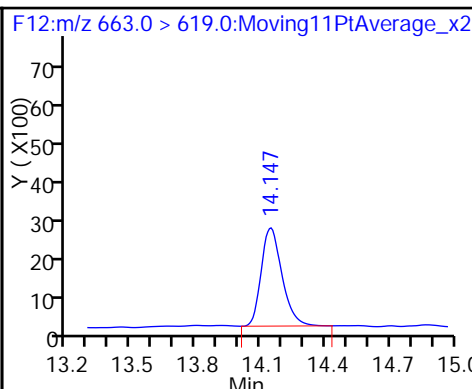
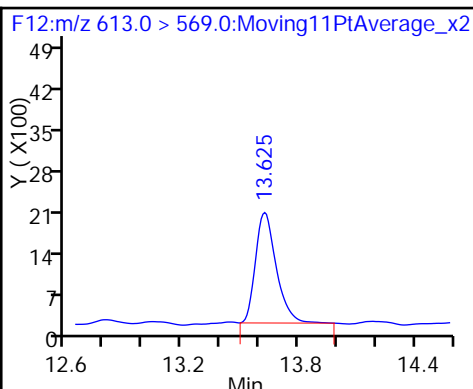
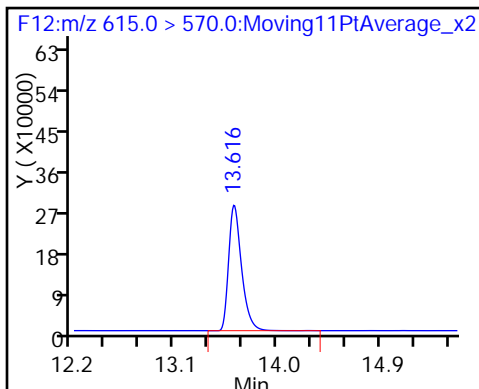
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

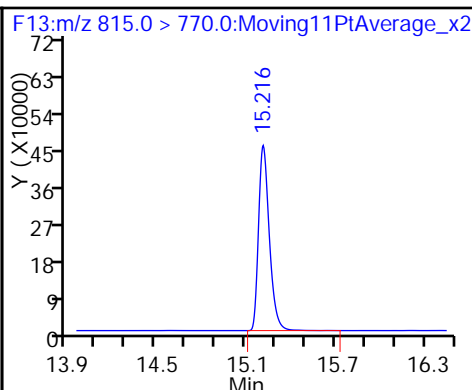
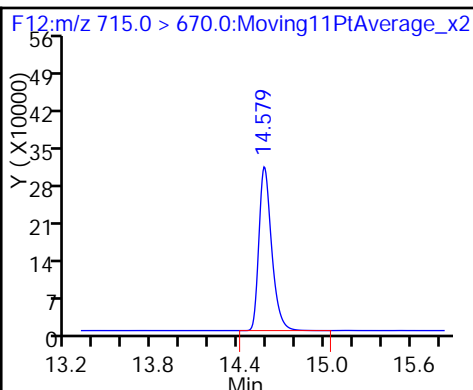
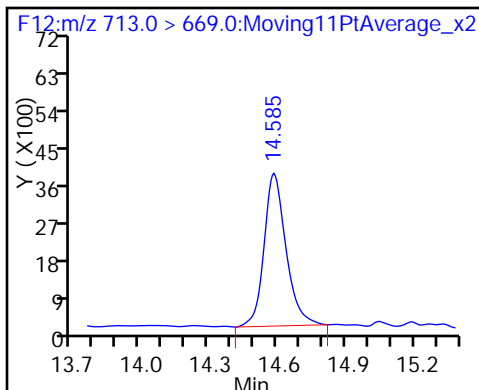
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

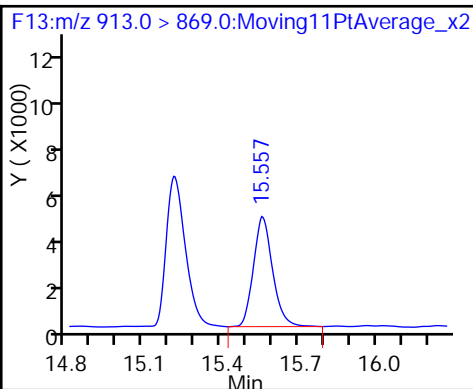
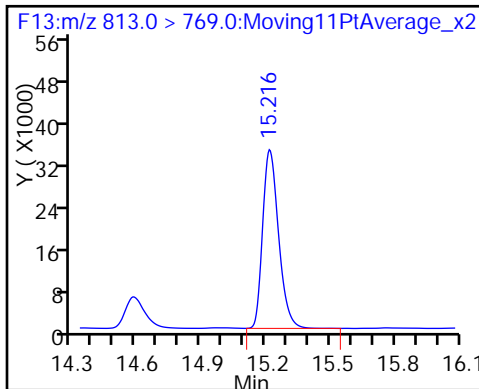
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_005.d  
 Lims ID: Std L2  
 Client ID:  
 Sample Type: IC Calib Level: 2  
 Inject. Date: 19-Feb-2016 15:29:34 ALS Bottle#: 10 Worklist Smp#: 3  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L2  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5

Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:43 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d

Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.718	5.712	0.006	907904	56.7		113	2336	
2 Perfluorobutyric acid	212.9 > 169.0	5.706	5.715	-0.009	18332	0.7742		77.4	48.8	
D 3 13C5-PFPeA	267.9 > 223.0	6.837	6.827	0.010	1879822	61.1		122	4021	
4 Perfluoropentanoic acid	262.9 > 219.0	6.837	6.828	0.009	41667	1.06		106	9.8	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.943	6.939	0.004	10245	NC			11.2	
	298.9 > 99.0	6.966	6.939	0.027	7619		1.34(0.00-0.00)		16.4	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.943	6.939	0.004	10245	0.6474		73.2		
7 Perfluorohexanoic acid	313.0 > 269.0	8.087	8.078	0.009	29048	0.8198		82.0	93.9	
D 6 13C2 PFHxA	315.0 > 270.0	8.093	8.080	0.013	1635595	57.7		115	3642	
D 8 13C4-PFHpA	367.0 > 322.0	9.334	9.322	0.012	1858023	60.3		121	6735	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.334	9.322	0.012	35173	0.9076		90.8	125	
D 11 18O2 PFHxS	403.0 > 84.0	9.369	9.358	0.011	753810	50.4		106	2618	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.387	9.364	0.023	9461	NC			32.6	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.387	9.364	0.023	9461	0.8791		92.9		
D 12 13C4 PFOA	417.0 > 372.0	10.468	10.450	0.018	1971561	57.7		115	5104	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
413.0 > 369.0	10.461	10.450	0.011	1.000	32208	0.8137		81.4	40.4	
413.0 > 169.0	10.468	10.450	0.018	1.001	15290		2.11(0.00-0.00)	81.4	38.9	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.475	10.457	0.018	1.000	8675	NC			20.9	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.475	10.457	0.018	1.000	8675	0.9509		99.9		
D 16 13C4 PFOS										
503.0 > 80.0	11.428	11.414	0.014		991283	54.5		114	3292	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.436	11.417	0.019	1.000	18718	0.9282		97.1	72.2	
499.0 > 99.0	11.428	11.417	0.011	0.999	4150		4.51(0.00-0.00)	97.1	19.5	
D 17 13C5 PFNA										
468.0 > 423.0	11.451	11.434	0.017		1761627	62.5		125	4150	
18 Perfluorononanoic acid										
463.0 > 419.0	11.451	11.439	0.012	1.000	23017	0.8091		80.9	58.4	
D 19 13C2 PFDA										
515.0 > 470.0	12.286	12.275	0.011		1607037	56.1		112	6838	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.296	12.276	0.020	1.000	42554	1.00		100	97.7	
D 23 13C8 FOSA										
506.0 > 78.0	12.824	12.810	0.014		2206225	53.2		106	1506	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.824	12.812	0.012	1.000	42097	1.03		103	188	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.970	12.949	0.021	1.000	6299	NC			16.0	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.970	12.949	0.021	1.000	6299	0.8330		86.4		
D 26 13C2 PFUnA										
565.0 > 520.0	13.010	13.000	0.010		2118412	56.7		113	6237	
27 Perfluoroundecanoic acid										
563.0 > 519.0	13.010	13.000	0.010	1.000	51520	0.8005		80.0	179	
D 28 13C2 PFDoA										
615.0 > 570.0	13.624	13.609	0.015		2242621	57.3		115	5907	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.624	13.611	0.013	1.000	23692	0.7225		72.3	32.2	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.139	14.130	0.009	1.000	40040	0.9488		94.9	60.9	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.585	14.571	0.014	1.000	42000	1.10		110	25.5	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.585	14.571	0.014		1961399	54.6		109	5975	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.218	15.207	0.011		2343949	54.4		109	6703	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.218	15.208	0.010	1.000	171017	1.00		100	211	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.549	15.543	0.006	1.000	39408	0.7910		79.1	62.1	

[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L2\_00019

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_005.d

Injection Date: 19-Feb-2016 15:29:34

Instrument ID: A6

Lims ID: Std L2

Client ID:

Operator ID: JRB

ALS Bottle#: 10

Worklist Smp#: 3

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

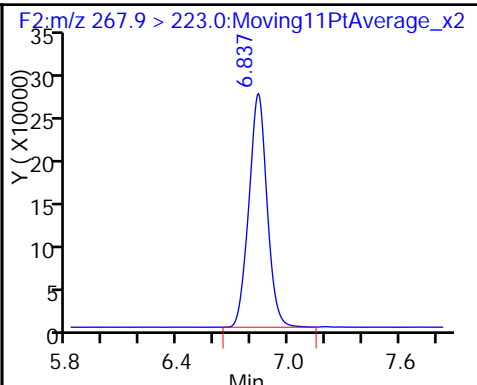
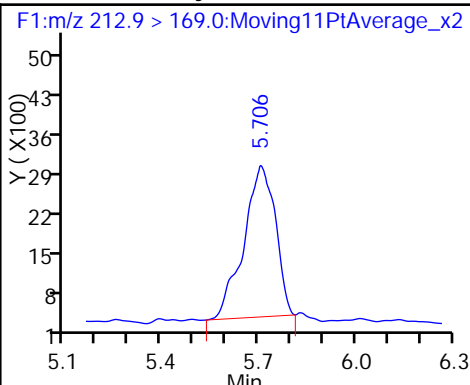
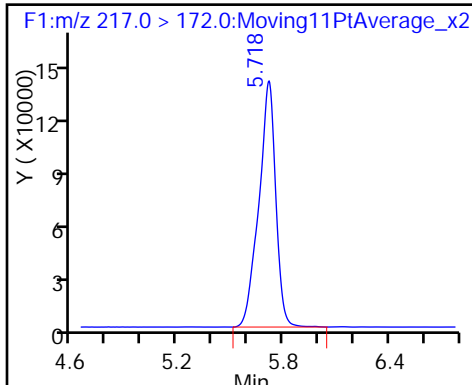
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

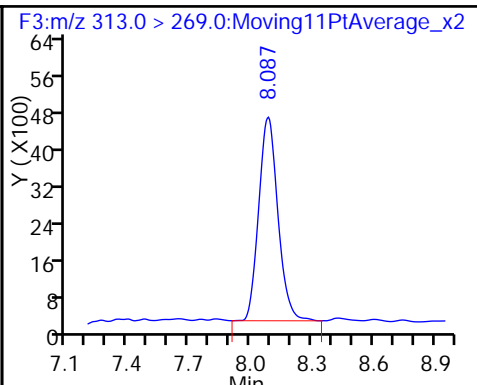
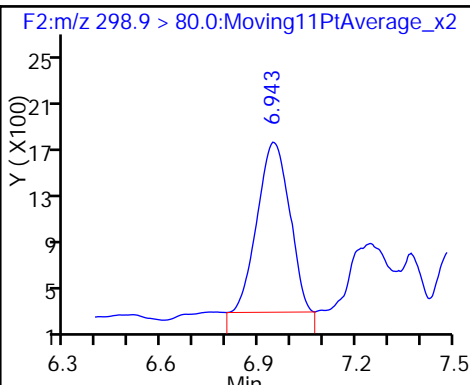
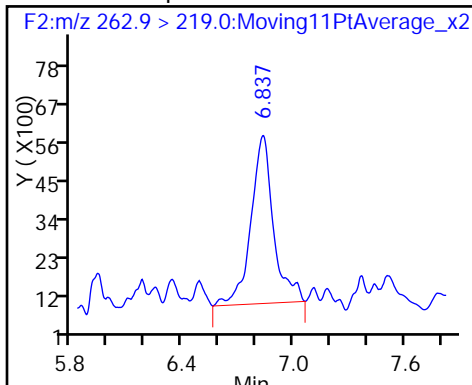
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

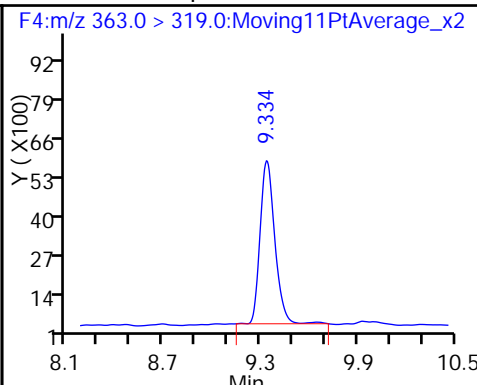
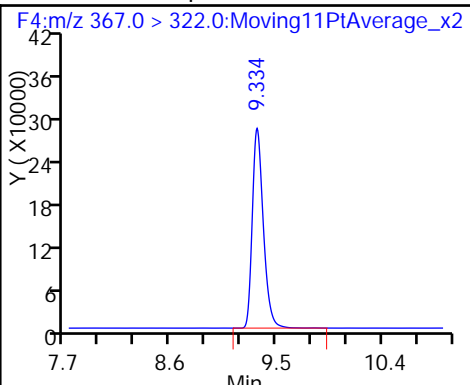
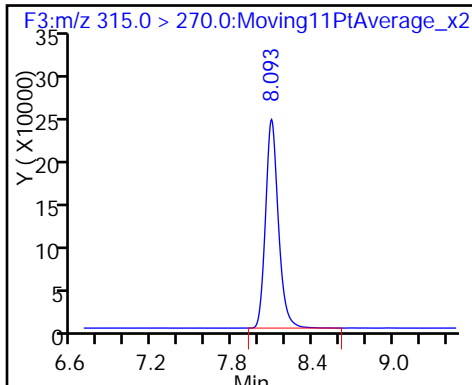
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

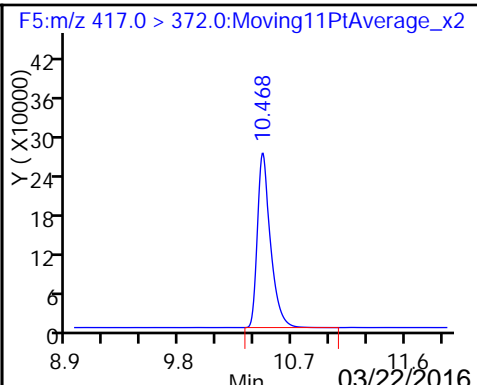
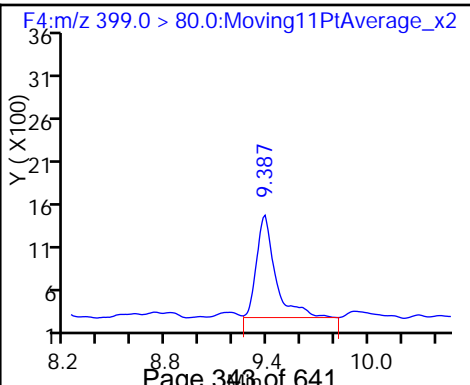
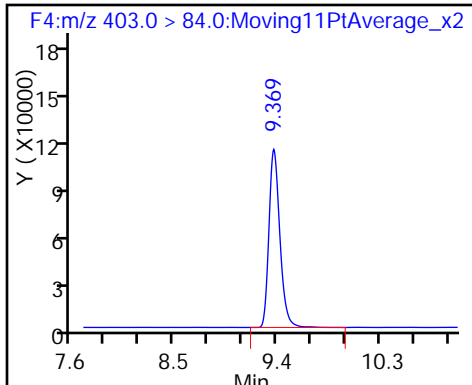
9 Perfluoroheptanoic acid

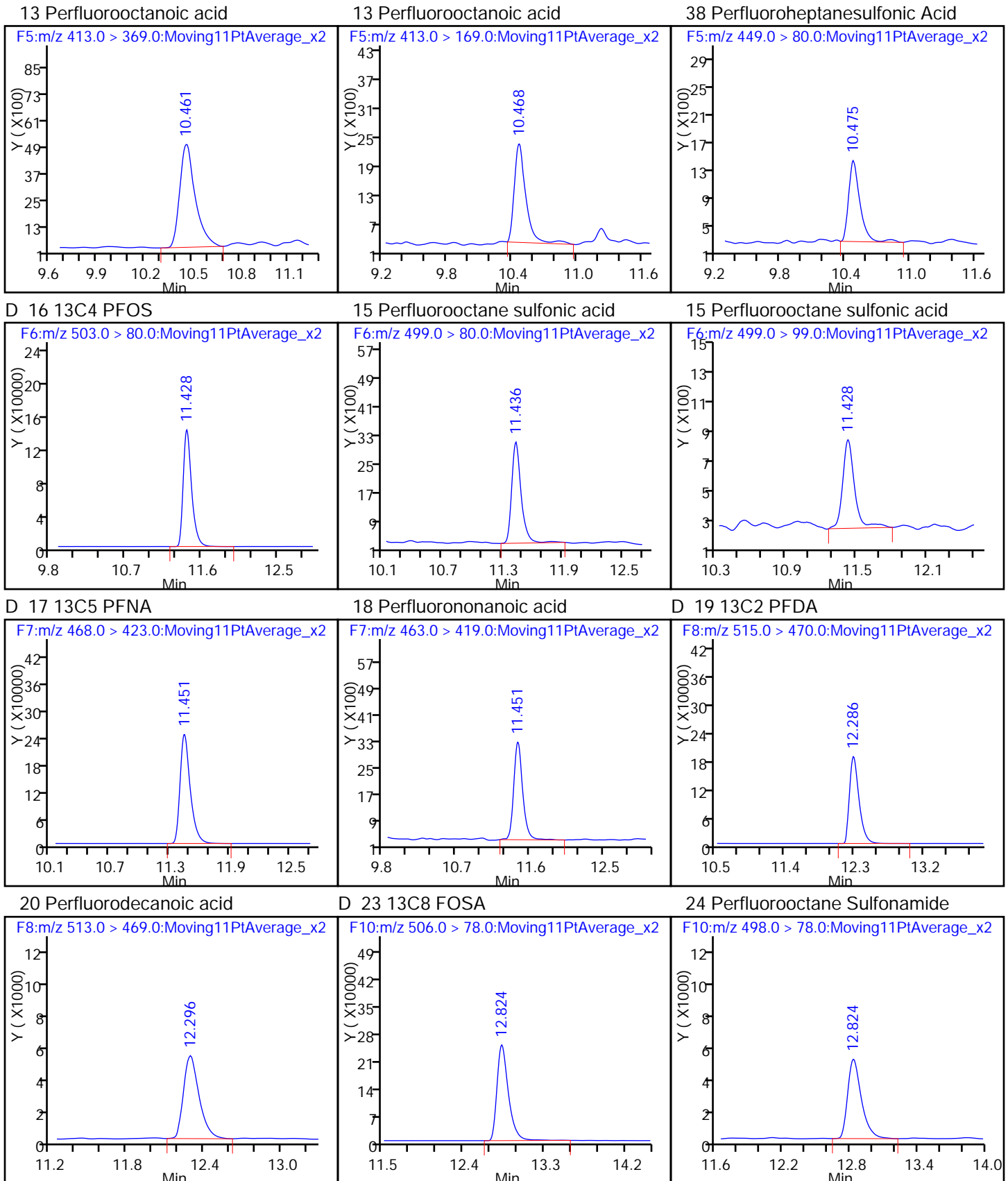


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA



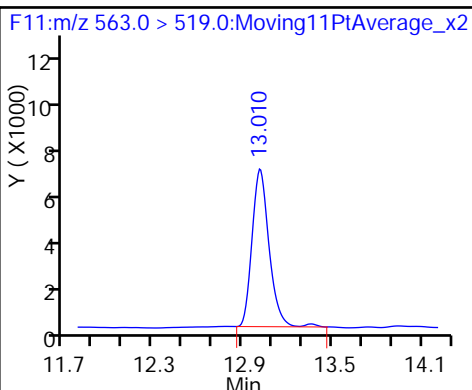
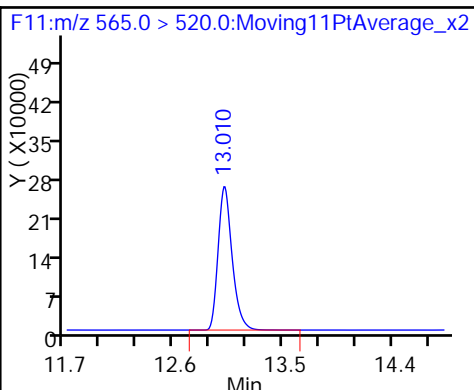
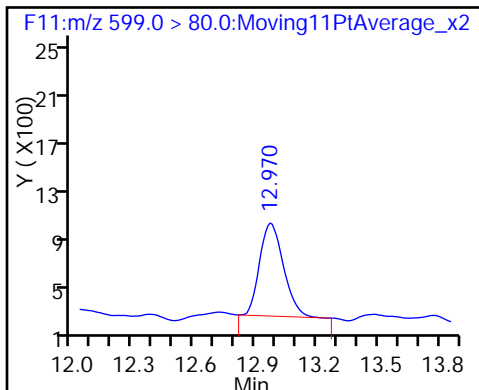




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

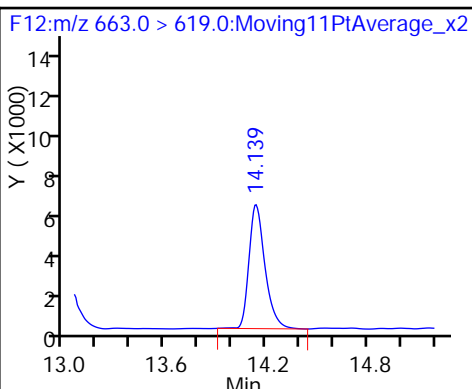
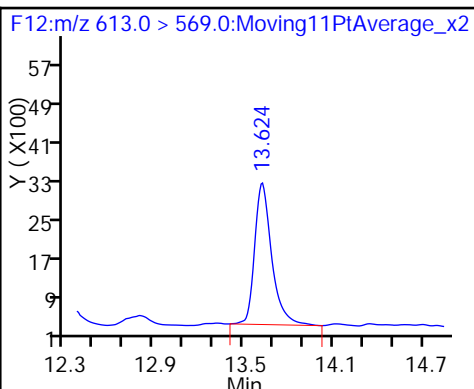
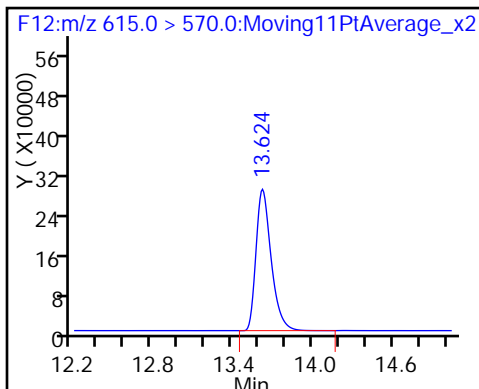
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

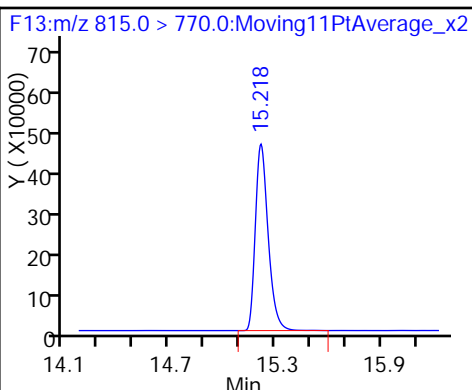
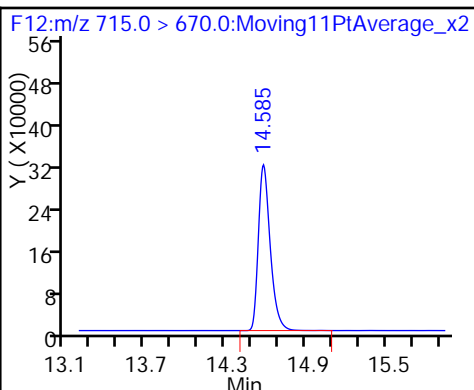
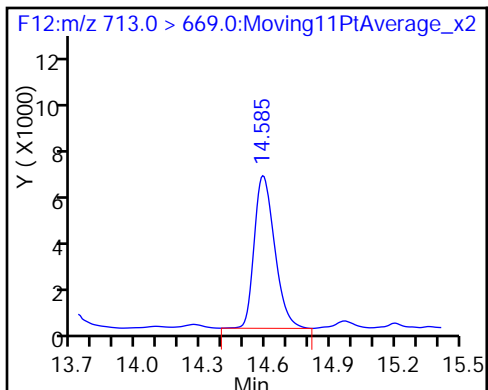
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

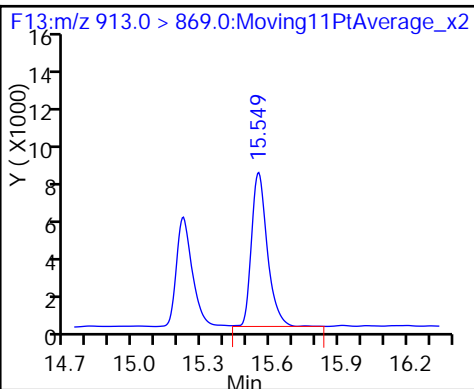
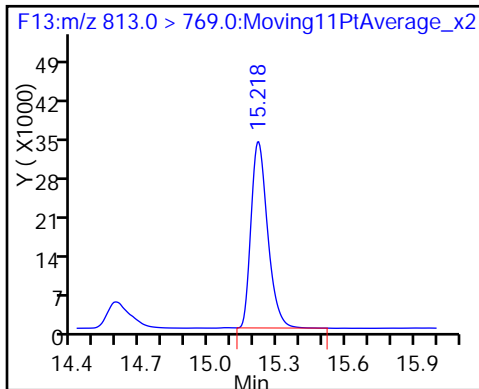
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_006.d  
 Lims ID: Std L3  
 Client ID:  
 Sample Type: IC Calib Level: 3  
 Inject. Date: 19-Feb-2016 15:50:47 ALS Bottle#: 11 Worklist Smp#: 4  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L3  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:45 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA										
217.0 > 172.0	5.712	5.712	0.0		882584	55.2		110	3246	
2 Perfluorobutyric acid										
212.9 > 169.0	5.718	5.715	0.003	1.000	123793	5.38		108	317	
D 3 13C5-PFPeA										
267.9 > 223.0	6.828	6.827	0.001		1560021	50.7		101	4030	
4 Perfluoropentanoic acid										
262.9 > 219.0	6.842	6.828	0.014	1.000	168531	5.18		104	40.4	
5 Perfluorobutane Sulfonate										
298.9 > 80.0	6.939	6.939	0.0	1.000	61862	NC			60.9	
298.9 > 99.0	6.948	6.939	0.009	1.001	38678		1.60(0.00-0.00)		67.4	
40 Perfluorobutanesulfonic acid										
298.9 > 80.0	6.939	6.939	0.0	1.000	61862	4.07		92.1		
7 Perfluorohexanoic acid										
313.0 > 269.0	8.084	8.078	0.006	1.000	177522	5.58		112	677	
D 6 13C2 PFHxA										
315.0 > 270.0	8.084	8.080	0.004		1469185	51.9		104	5297	
D 8 13C4-PFHpA										
367.0 > 322.0	9.328	9.322	0.006		1638704	53.2		106	5483	
9 Perfluoroheptanoic acid										
363.0 > 319.0	9.328	9.322	0.006	1.000	167149	4.89		97.8	599	
D 11 18O2 PFHxS										
403.0 > 84.0	9.363	9.358	0.005		723683	48.4		102	2722	
10 Perfluorohexane Sulfonate										
399.0 > 80.0	9.363	9.364	-0.001	1.000	48077	NC			141	
41 Perfluorohexanesulfonic acid										
399.0 > 80.0	9.363	9.364	-0.001	1.000	48077	4.65		98.4		
D 12 13C4 PFOA										
417.0 > 372.0	10.453	10.450	0.003		1750451	51.3		103	8337	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
413.0 > 369.0	10.453	10.450	0.003	1.000	188978	5.38		108	203	
413.0 > 169.0	10.461	10.450	0.011	1.001	67070		2.82(0.00-0.00)	108	145	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.461	10.457	0.004	1.000	52893	NC			204	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.461	10.457	0.004	1.000	52893	4.78		100		
D 16 13C4 PFOS										
503.0 > 80.0	11.420	11.414	0.006		898344	49.4		103	3427	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.420	11.417	0.003	1.000	86511	4.73		99.0	229	
499.0 > 99.0	11.420	11.417	0.003	1.000	39257		2.20(0.00-0.00)	99.0	114	
D 17 13C5 PFNA										
468.0 > 423.0	11.443	11.434	0.009		1474756	52.3		105	5146	
18 Perfluorononanoic acid										
463.0 > 419.0	11.443	11.439	0.004	1.000	120395	5.06		101	398	
D 19 13C2 PFDA										
515.0 > 470.0	12.286	12.275	0.011		1565532	54.7		109	4281	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.286	12.276	0.010	1.000	163841	5.01		100	357	
D 23 13C8 FOSA										
506.0 > 78.0	12.814	12.810	0.004		2223851	53.6		107	1602	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.814	12.812	0.002	1.000	203867	4.95		98.9	406	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.960	12.949	0.011	1.000	49387	NC			162	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.960	12.949	0.011	1.000	49387	5.00		104		
D 26 13C2 PFUnA										
565.0 > 520.0	13.010	13.000	0.010		1980471	53.0		106	5969	
27 Perfluoroundecanoic acid										
563.0 > 519.0	13.010	13.000	0.010	1.000	213464	5.87		117	515	
D 28 13C2 PFDoA										
615.0 > 570.0	13.615	13.609	0.006		2017427	51.6		103	3234	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.615	13.611	0.004	1.000	163307	5.54		111	163	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.139	14.130	0.009	1.000	243228	6.41		128	291	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.578	14.571	0.007	1.000	135077	5.14		103	90.0	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.578	14.571	0.007		1864931	51.9		104	5864	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.213	15.207	0.006		2201810	51.1		102	5730	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.213	15.208	0.005	1.000	324953	4.96		99.1	522	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.544	15.543	0.001	1.000	245350	5.47		109	345	

[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L3\_00016

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_006.d

Injection Date: 19-Feb-2016 15:50:47

Instrument ID: A6

Lims ID: Std L3

Client ID:

Operator ID: JRB

ALS Bottle#: 11

Worklist Smp#: 4

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

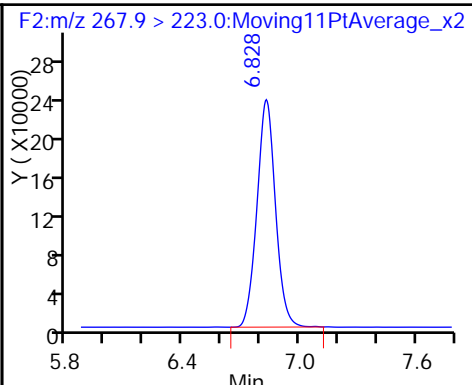
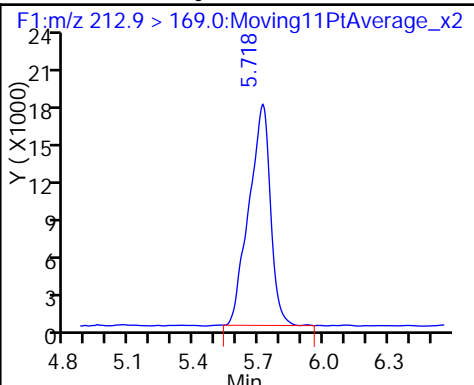
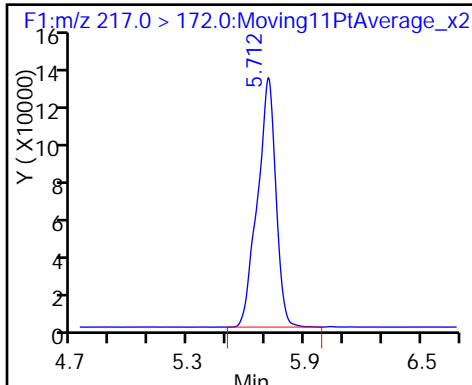
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

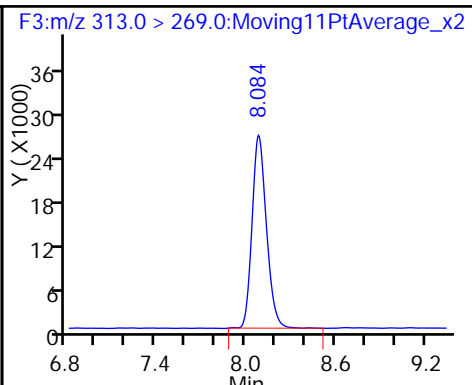
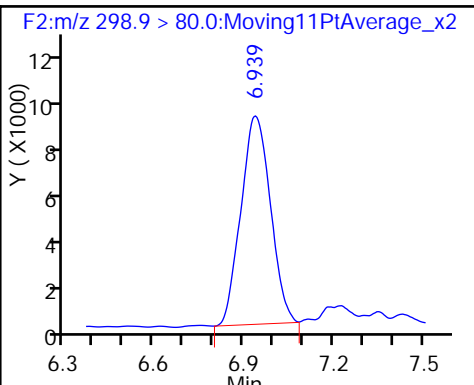
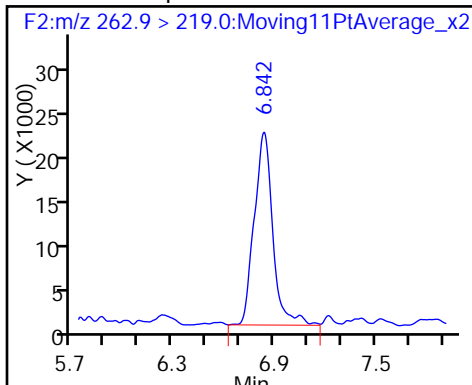
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

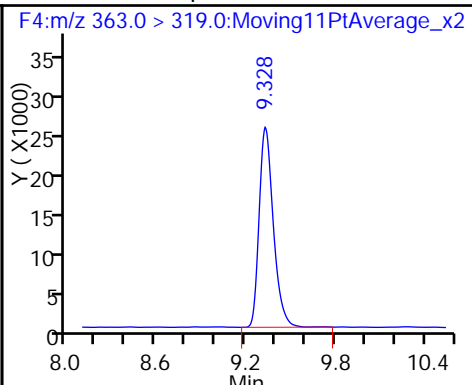
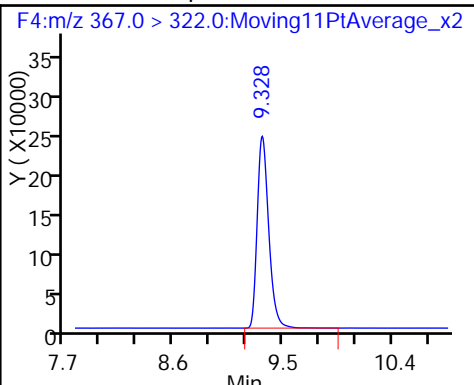
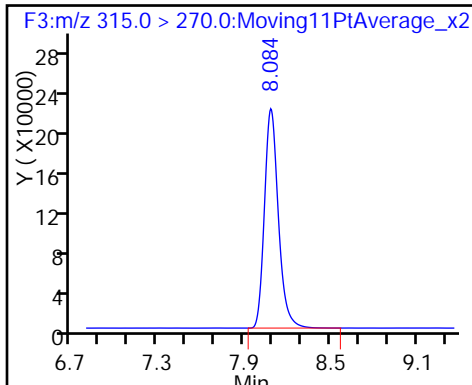
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

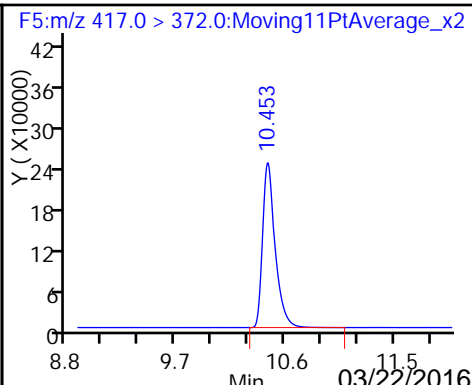
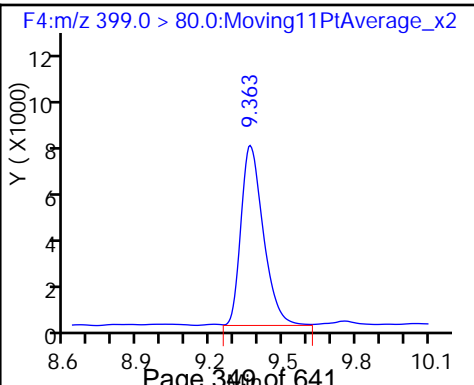
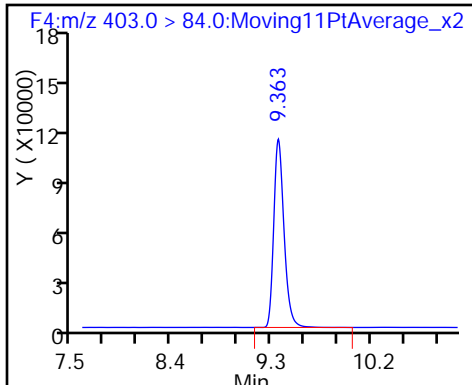
9 Perfluoroheptanoic acid

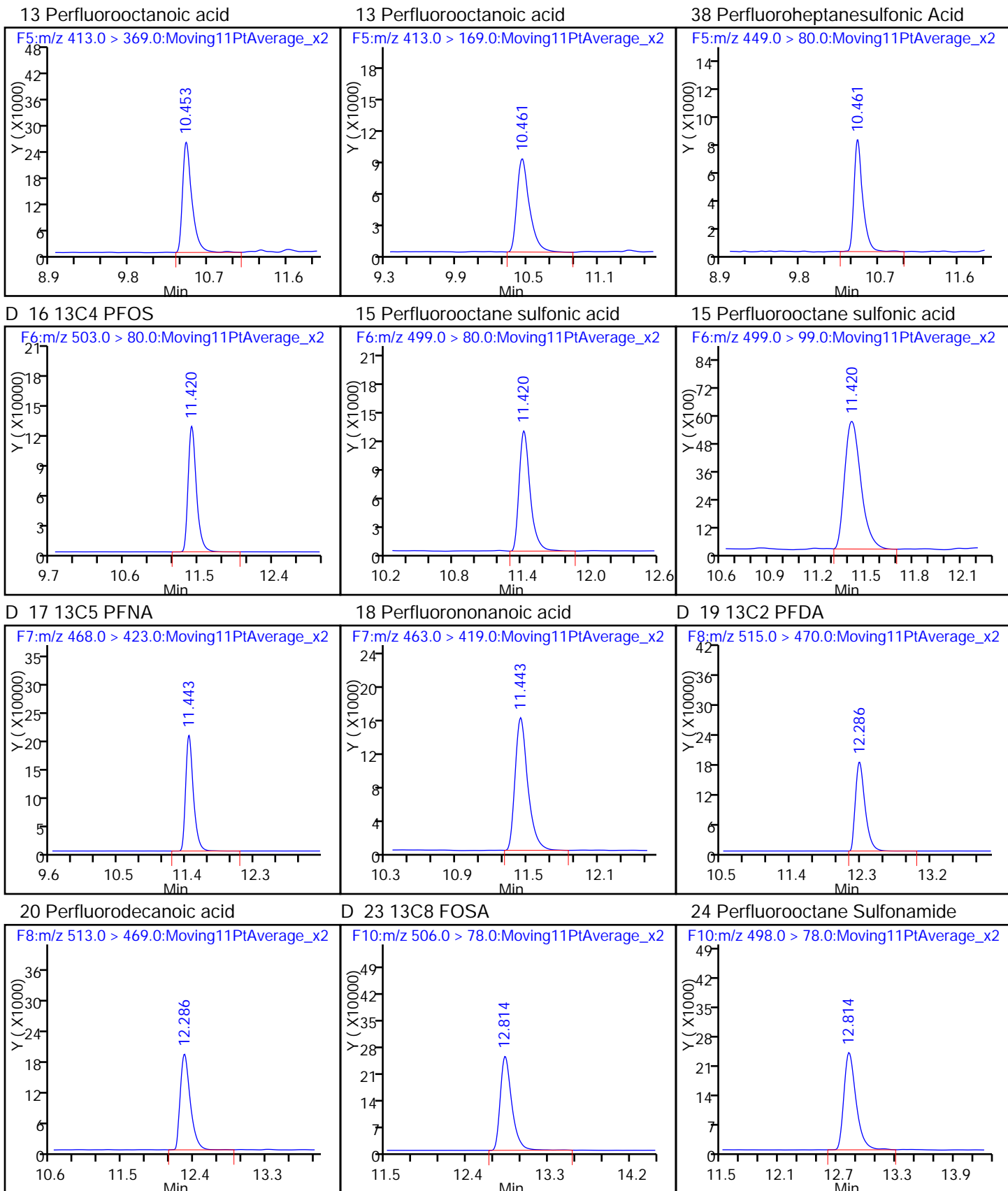


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

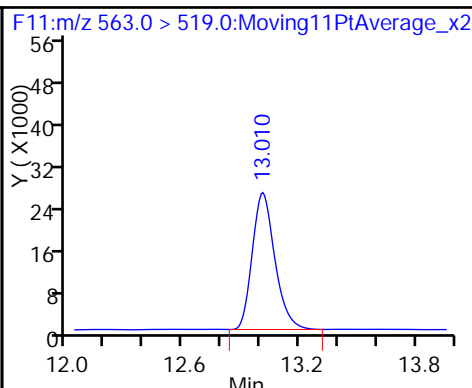
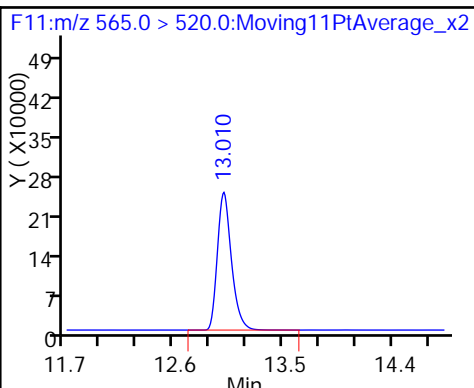
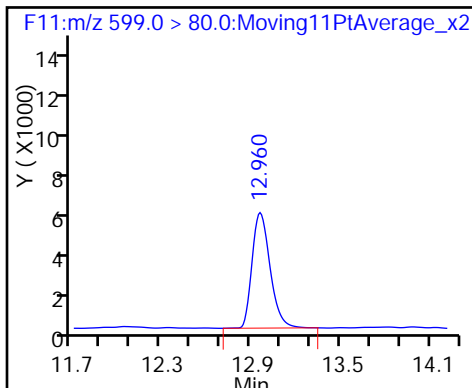




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

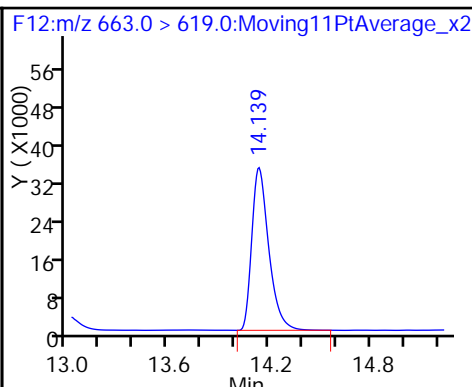
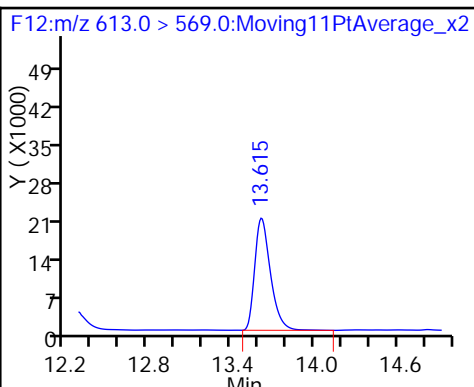
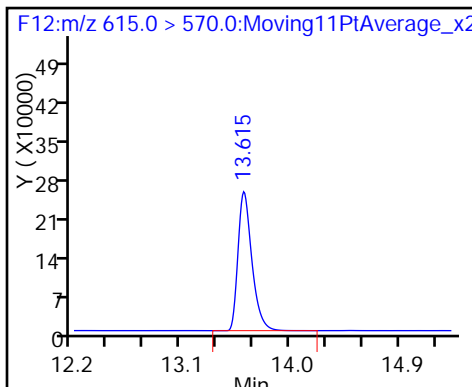
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

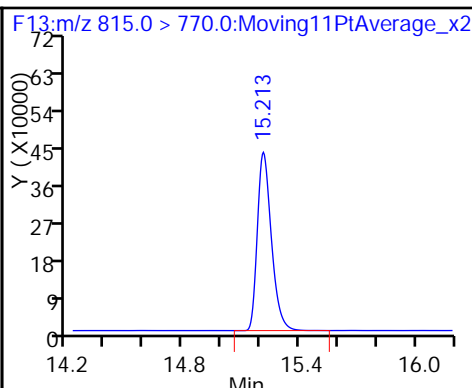
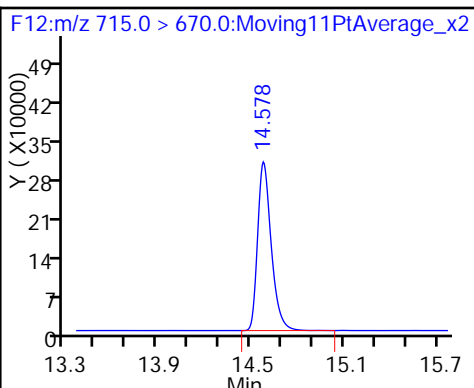
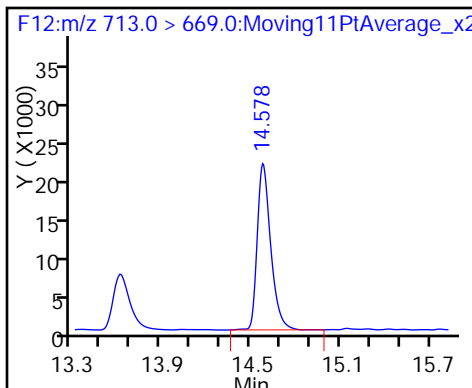
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

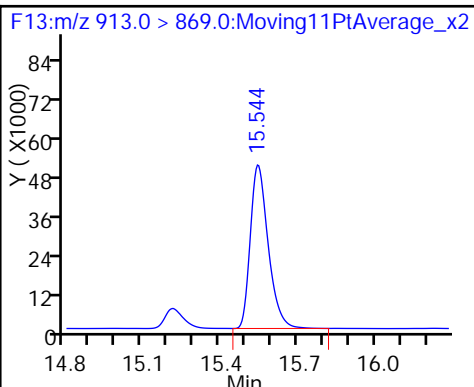
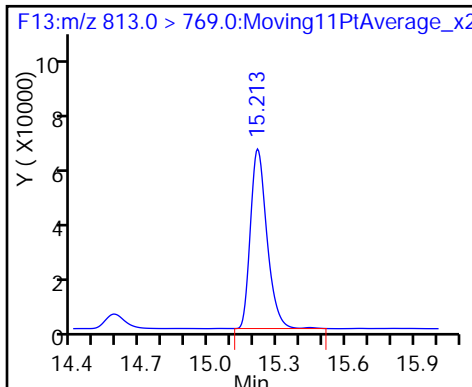
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_007.d  
 Lims ID: Std L4  
 Client ID:  
 Sample Type: IC Calib Level: 4  
 Inject. Date: 19-Feb-2016 16:11:58 ALS Bottle#: 12 Worklist Smp#: 5  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L4  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:48 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: westendorfc Date: 21-Feb-2016 15:43:25

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.709	5.712	-0.003	823965	51.5		103	1860	
2 Perfluorobutyric acid	212.9 > 169.0	5.715	5.715	0.0	462922	21.5		108	1792	
D 3 13C5-PFPeA	267.9 > 223.0	6.828	6.827	0.001	1598957	52.0		104	5454	
4 Perfluoropentanoic acid	262.9 > 219.0	6.828	6.828	0.0	673970	20.2		101	192	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.942	6.939	0.003	294821	NC			430	
	298.9 > 99.0	6.942	6.939	0.003	171078		1.72(0.00-0.00)		274	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.942	6.939	0.003	294821	19.0		107		
7 Perfluorohexanoic acid	313.0 > 269.0	8.082	8.078	0.004	618008	18.9		94.3	2035	
D 6 13C2 PFHxA	315.0 > 270.0	8.082	8.080	0.002	1512461	53.4		107	5873	
D 8 13C4-PFHpA	367.0 > 322.0	9.323	9.322	0.001	1694741	55.0		110	5704	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.323	9.322	0.001	778743	22.0		110	2475	
D 11 18O2 PFHxS	403.0 > 84.0	9.358	9.358	0.0	739449	49.4		104	2322	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.358	9.364	-0.006	205104	NC			759	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.358	9.364	-0.006	205104	19.4		103		



Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.454	10.450	0.004		1793859	52.5		105	5536	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.454	10.450	0.004	1.000	715180	19.9		99.3	762	
413.0 > 169.0	10.454	10.450	0.004	1.000	250759		2.85(0.00-0.00)	99.3	716	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.461	10.457	0.004	1.000	216336	NC			809	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.461	10.457	0.004	1.000	216336	19.1		100		
D 16 13C4 PFOS										
503.0 > 80.0	11.414	11.414	0.0		879178	48.4		101	3624	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.421	11.417	0.004	1.000	356703	19.9		104	413	
499.0 > 99.0	11.414	11.417	-0.003	0.999	193440		1.84(0.00-0.00)	104	731	
D 17 13C5 PFNA										
468.0 > 423.0	11.437	11.434	0.003		1459436	51.8		104	5635	
18 Perfluorononanoic acid										
463.0 > 419.0	11.437	11.439	-0.002	1.000	518362	22.0		110	1486	
D 19 13C2 PFDA										
515.0 > 470.0	12.276	12.275	0.001		1524195	53.2		106	4870	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.276	12.276	0.0	1.000	571448	18.9		94.4	1987	
D 23 13C8 FOSA										
506.0 > 78.0	12.814	12.810	0.004		2199385	53.0		106	2293	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.814	12.812	0.002	1.000	832823	20.4		102	833	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.950	12.949	0.001	1.000	202871	NC			616	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.950	12.949	0.001	1.000	202871	20.1		104		
D 26 13C2 PFUnA										
565.0 > 520.0	13.002	13.000	0.002		1975170	52.8		106	5266	
27 Perfluoroundecanoic acid										
563.0 > 519.0	13.002	13.000	0.002	1.000	694202	20.7		103	1673	
D 28 13C2 PFDaA										
615.0 > 570.0	13.615	13.609	0.006		2115244	54.1		108	6995	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.615	13.611	0.004	1.000	679009	22.0		110	827	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.132	14.130	0.002	1.000	875238	22.0		110	983	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.572	14.571	0.001	1.000	498848	19.3		96.5	333	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.579	14.571	0.008		1885833	52.5		105	5639	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.213	15.207	0.006		2282833	53.0		106	5499	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.213	15.208	0.005	1.000	996289	19.4		97.0	1274	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.544	15.543	0.001	1.000	961125	20.5	102	1418	

### QC Flag Legend

Processing Flags

NC - Not Calibrated

### Reagents:

LCPFC-L4\_00017

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_007.d

Injection Date: 19-Feb-2016 16:11:58

Instrument ID: A6

Lims ID: Std L4

Client ID:

Operator ID: JRB

ALS Bottle#: 12

Worklist Smp#: 5

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

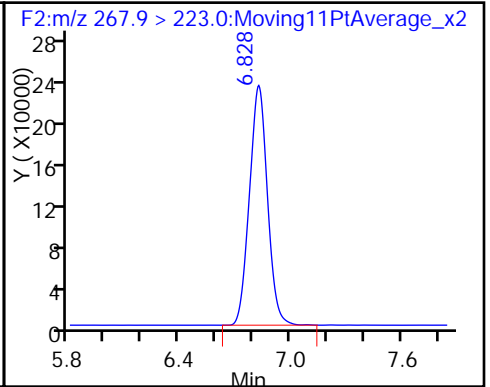
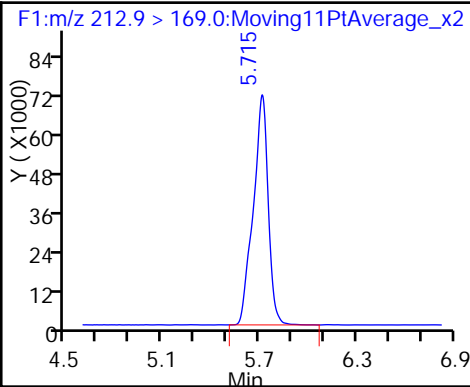
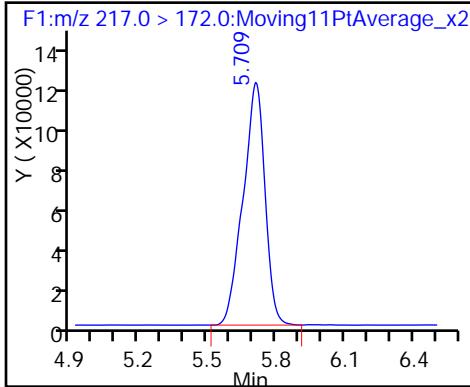
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

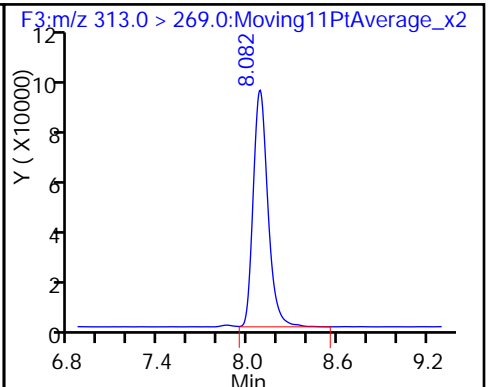
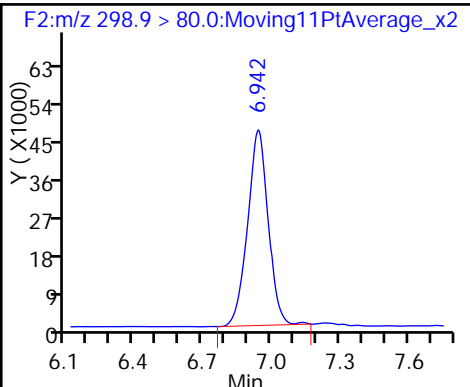
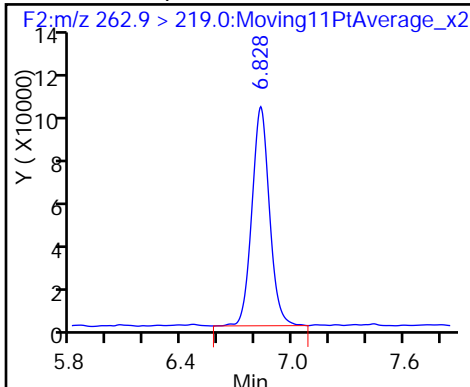
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

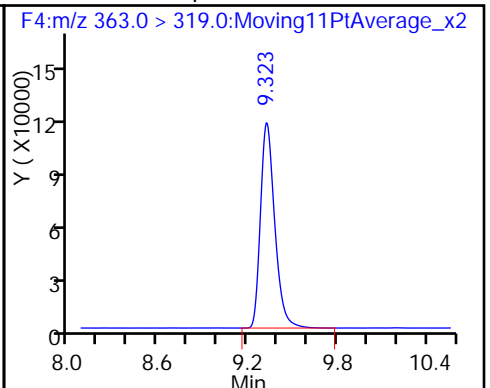
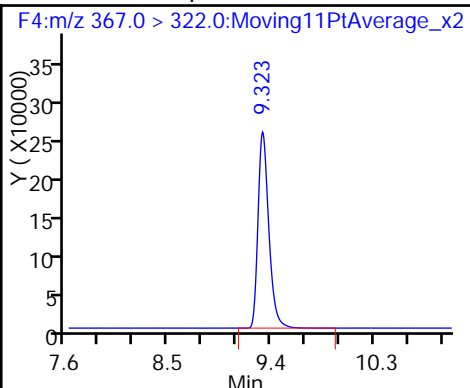
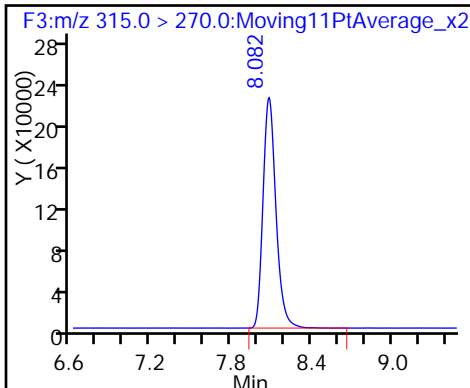
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

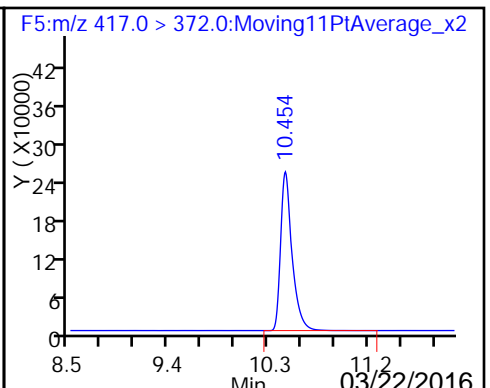
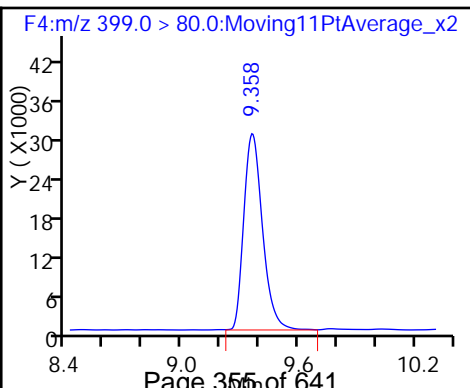
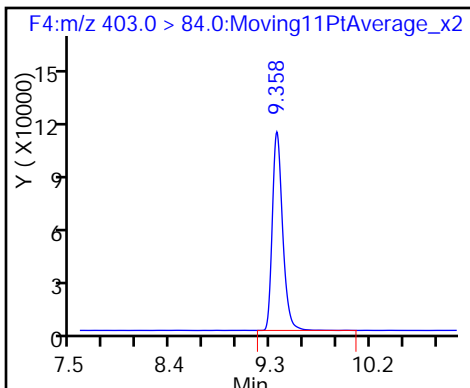
9 Perfluoroheptanoic acid

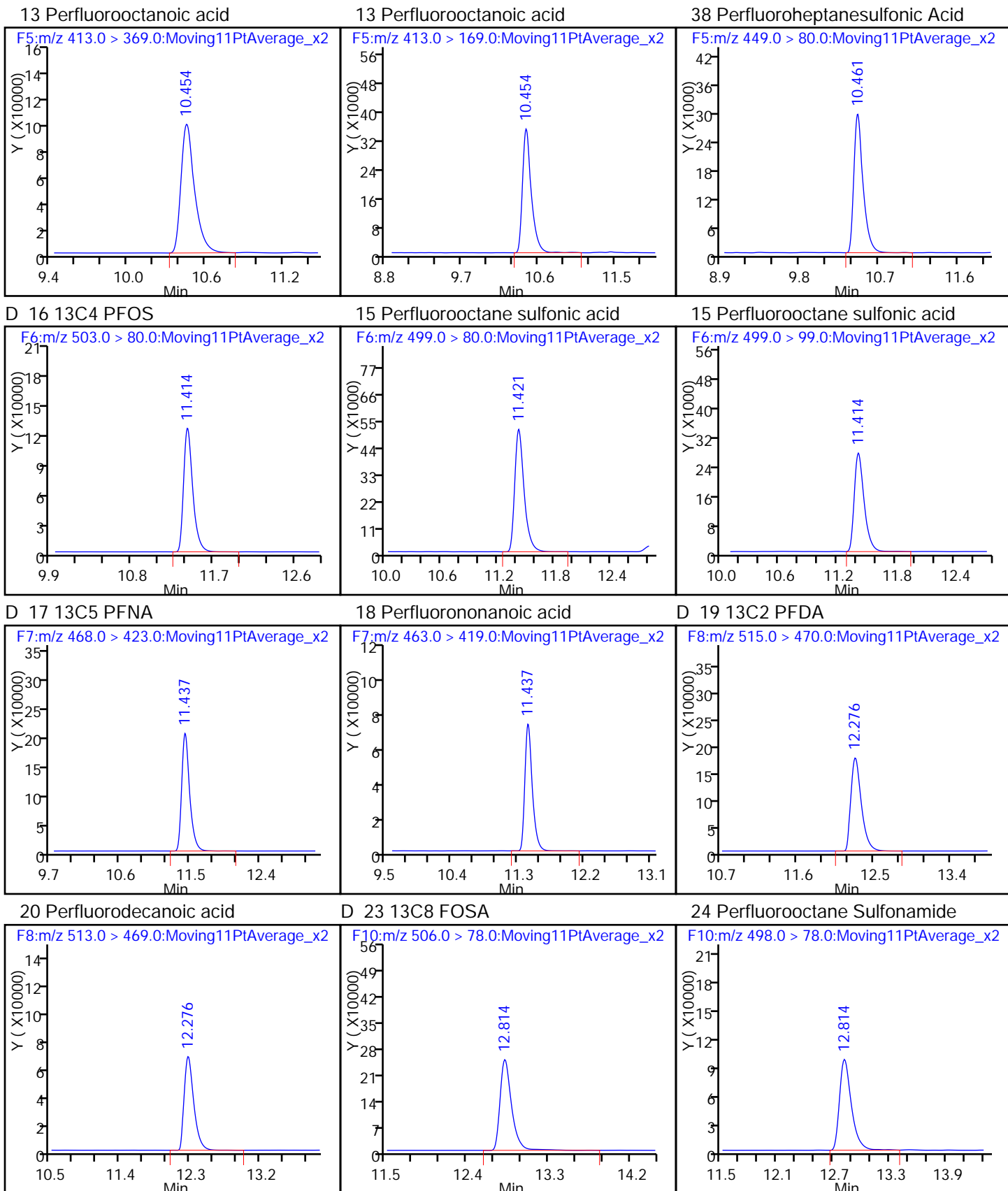


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

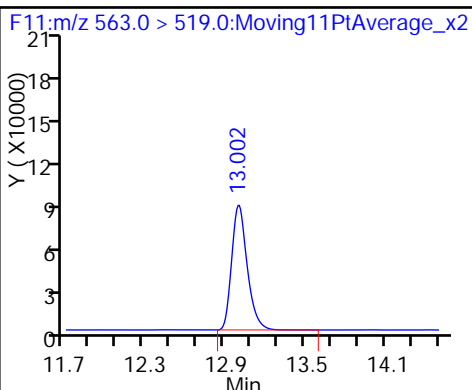
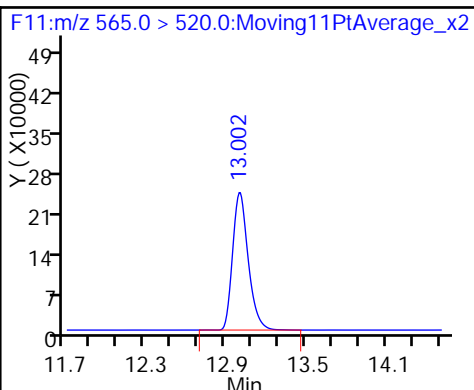
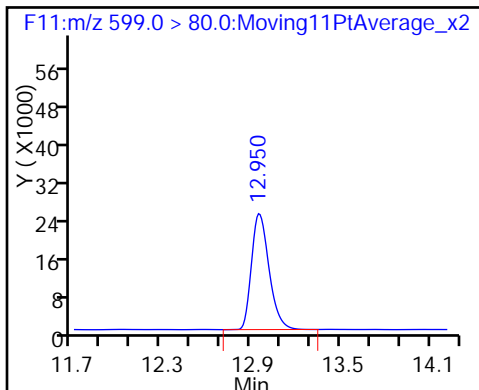




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

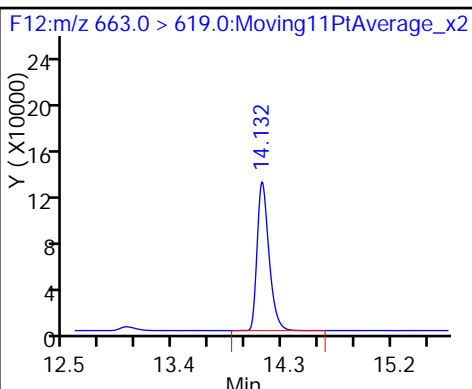
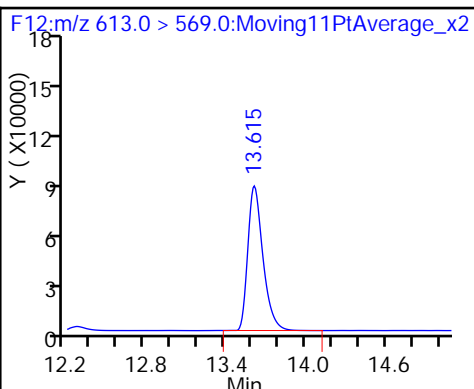
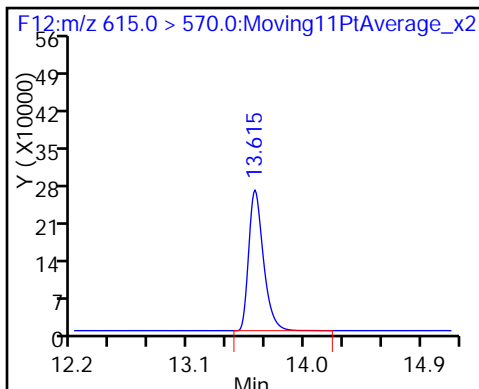
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

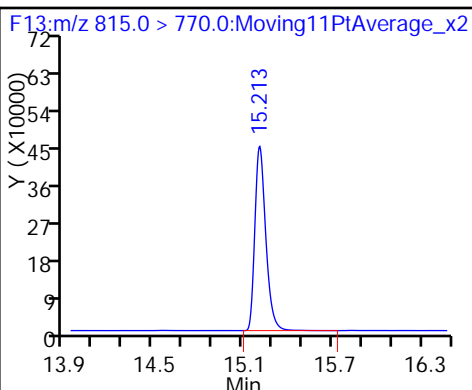
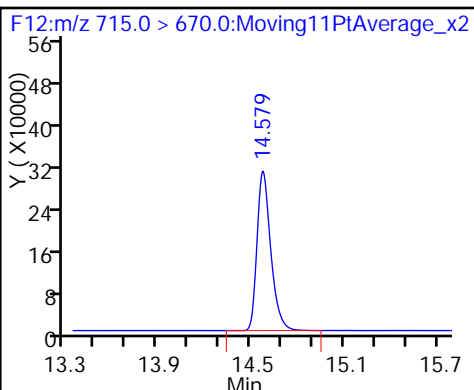
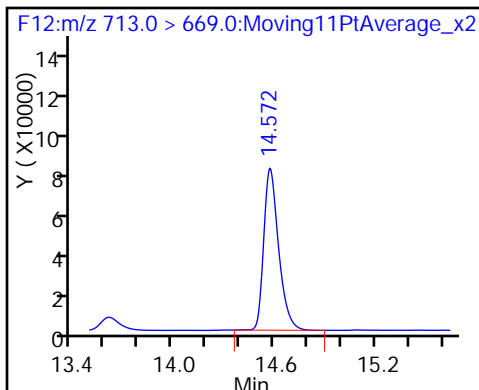
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

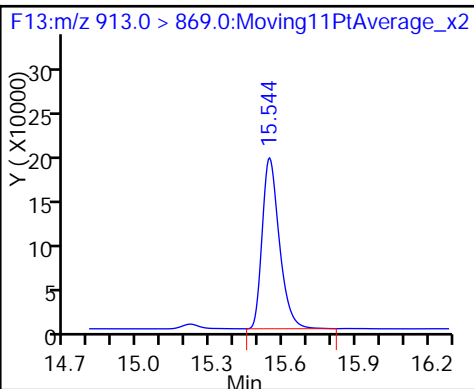
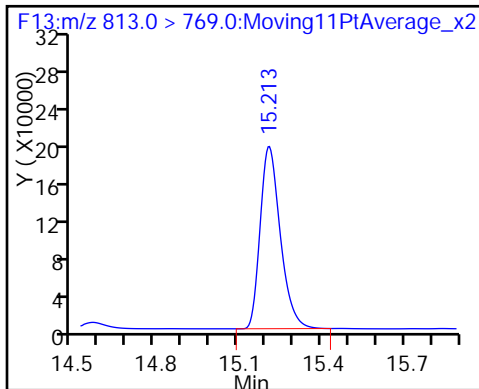
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_008.d  
 Lims ID: Std L5  
 Client ID:  
 Sample Type: IC Calib Level: 5  
 Inject. Date: 19-Feb-2016 16:33:12 ALS Bottle#: 13 Worklist Smp#: 6  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L5  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:51 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 19-Feb-2016 17:45:09

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.709	5.712	-0.003	778630	48.7		97.3	2140	
2 Perfluorobutyric acid	212.9 > 169.0	5.715	5.715	0.0	1121027	55.2		110	3299	
D 3 13C5-PFPeA	267.9 > 223.0	6.823	6.827	-0.004	1604599	52.2		104	5949	
4 Perfluoropentanoic acid	262.9 > 219.0	6.828	6.828	0.0	1554916	46.4		92.9	382	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.938	6.939	-0.001	696169	NC			934	
	298.9 > 99.0	6.943	6.939	0.004	378428		1.84(0.00-0.00)		526	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.938	6.939	-0.001	696169	48.6		110		
7 Perfluorohexanoic acid	313.0 > 269.0	8.077	8.078	-0.001	1598966	52.3		105	2420	
D 6 13C2 PFHxA	315.0 > 270.0	8.082	8.080	0.002	1412469	49.9		99.7	3689	
D 8 13C4-PFHpA	367.0 > 322.0	9.323	9.322	0.001	1498068	48.6		97.2	4589	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.323	9.322	0.001	1614588	51.7		103	4595	
D 11 18O2 PFHxS	403.0 > 84.0	9.358	9.358	0.0	681798	45.6		96.3	3254	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.364	9.364	0.0	467634	NC			1449	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.364	9.364	0.0	467634	48.0		102		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.454	10.450	0.004		1573533	46.1		92.2	4818	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.454	10.450	0.004	1.000	1580565	50.0		100	1517	
413.0 > 169.0	10.454	10.450	0.004	1.000	544646		2.90(0.00-0.00)	100	1443	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.461	10.457	0.004	1.000	520355	NC			2081	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.461	10.457	0.004	1.000	520355	48.6		102		
D 16 13C4 PFOS										
503.0 > 80.0	11.422	11.414	0.008		821717	45.2		94.5	2661	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.422	11.417	0.005	1.000	797626	47.7		99.8	421	
499.0 > 99.0	11.422	11.417	0.005	1.000	433010		1.84(0.00-0.00)	99.8	1470	
D 17 13C5 PFNA										
468.0 > 423.0	11.437	11.434	0.003		1386004	49.2		98.3	7596	
18 Perfluorononanoic acid										
463.0 > 419.0	11.437	11.439	-0.002	1.000	1106311	49.4		98.9	3403	
D 19 13C2 PFDA										
515.0 > 470.0	12.280	12.275	0.005		1321943	46.2		92.3	3965	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.280	12.276	0.004	1.000	1348689	52.0		104	2340	
D 23 13C8 FOSA										
506.0 > 78.0	12.818	12.810	0.008		2127859	51.3		103	1526	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.818	12.812	0.006	1.000	1983962	50.3		101	2804	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.954	12.949	0.005	1.000	497252	NC			1216	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.954	12.949	0.005	1.000	497252	52.2		108		
D 26 13C2 PFUnA										
565.0 > 520.0	13.005	13.000	0.005		1719562	46.0		92.0	3390	
27 Perfluoroundecanoic acid										
563.0 > 519.0	13.005	13.000	0.005	1.000	1511190	52.7		105	4312	
D 28 13C2 PFDaA										
615.0 > 570.0	13.620	13.609	0.011		1956174	50.0		100	4217	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.620	13.611	0.009	1.000	1540137	53.8		108	1765	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.136	14.130	0.006	1.000	1907812	51.8		104	1846	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.576	14.571	0.005	1.000	1159290	49.2		98.4	877	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.576	14.571	0.005		1841361	51.3		103	6184	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.210	15.207	0.003		2184111	50.7		101	5115	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.210	15.208	0.002	1.000	2198803	49.9		99.7	2557	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.546	15.543	0.003	1.000	2246657	51.7	103	2473	

### QC Flag Legend

Processing Flags

NC - Not Calibrated

### Reagents:

LCPFC-L5\_00016

Amount Added: 1.00

Units: mL



TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_008.d

Injection Date: 19-Feb-2016 16:33:12

Instrument ID: A6

Lims ID: Std L5

Client ID:

Operator ID: JRB

ALS Bottle#: 13

Worklist Smp#: 6

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

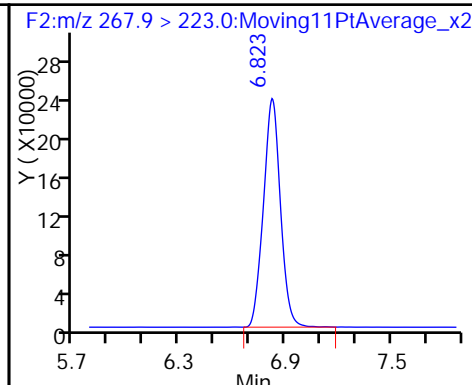
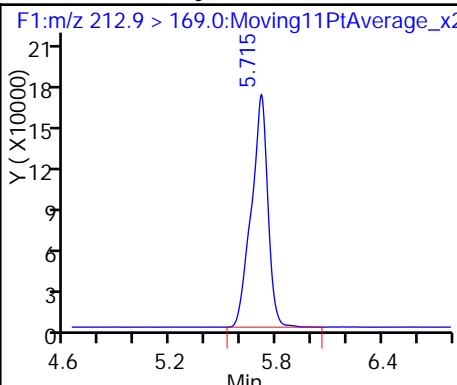
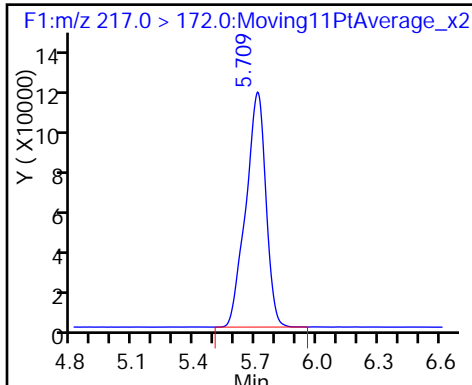
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

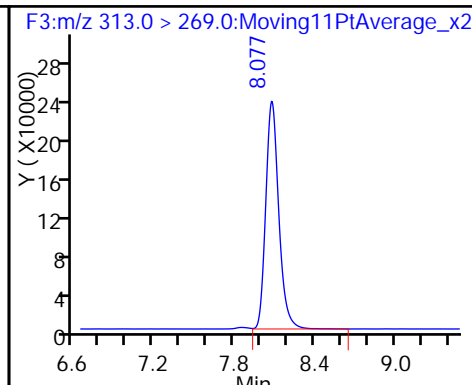
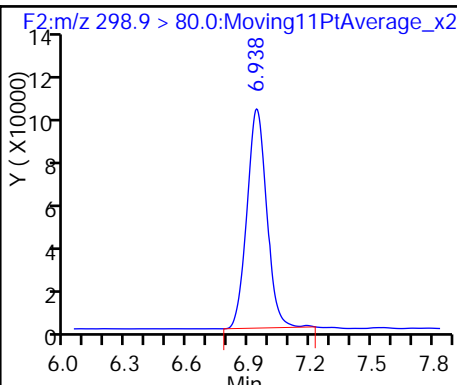
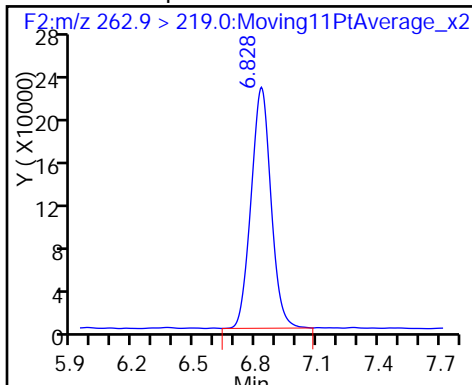
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

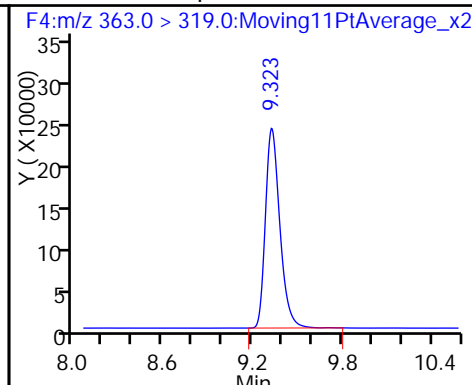
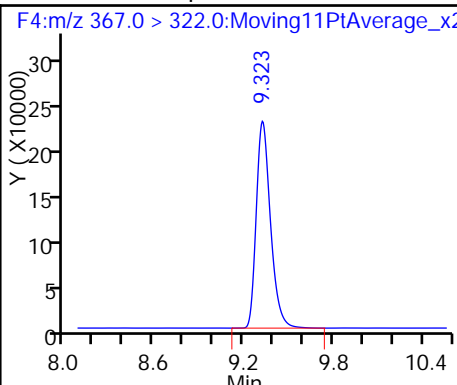
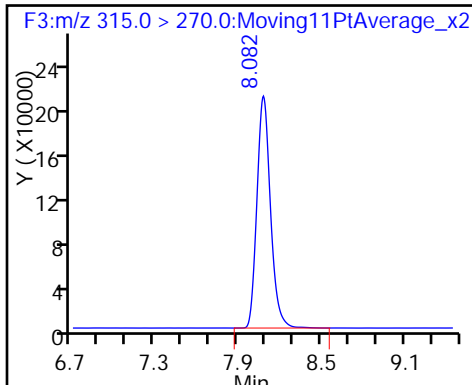
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

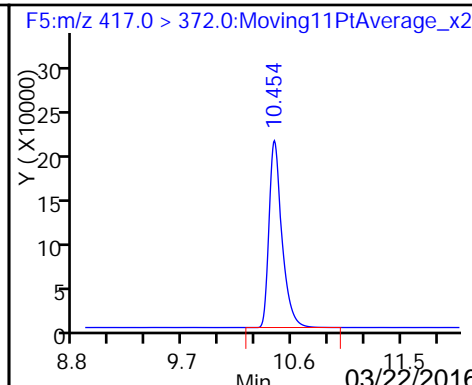
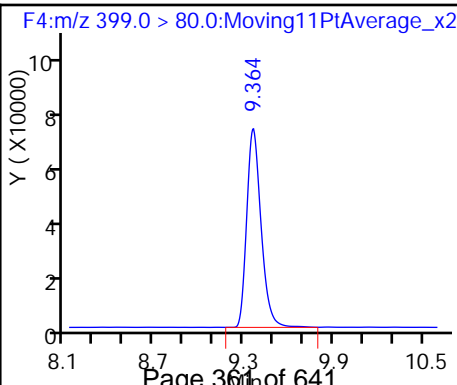
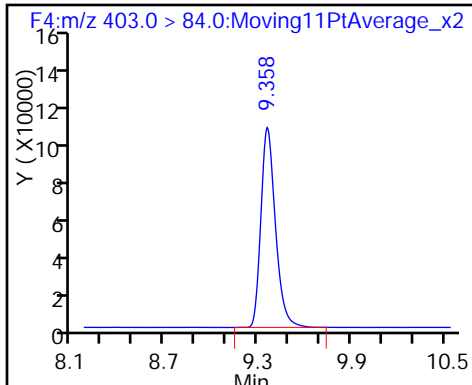
9 Perfluoroheptanoic acid

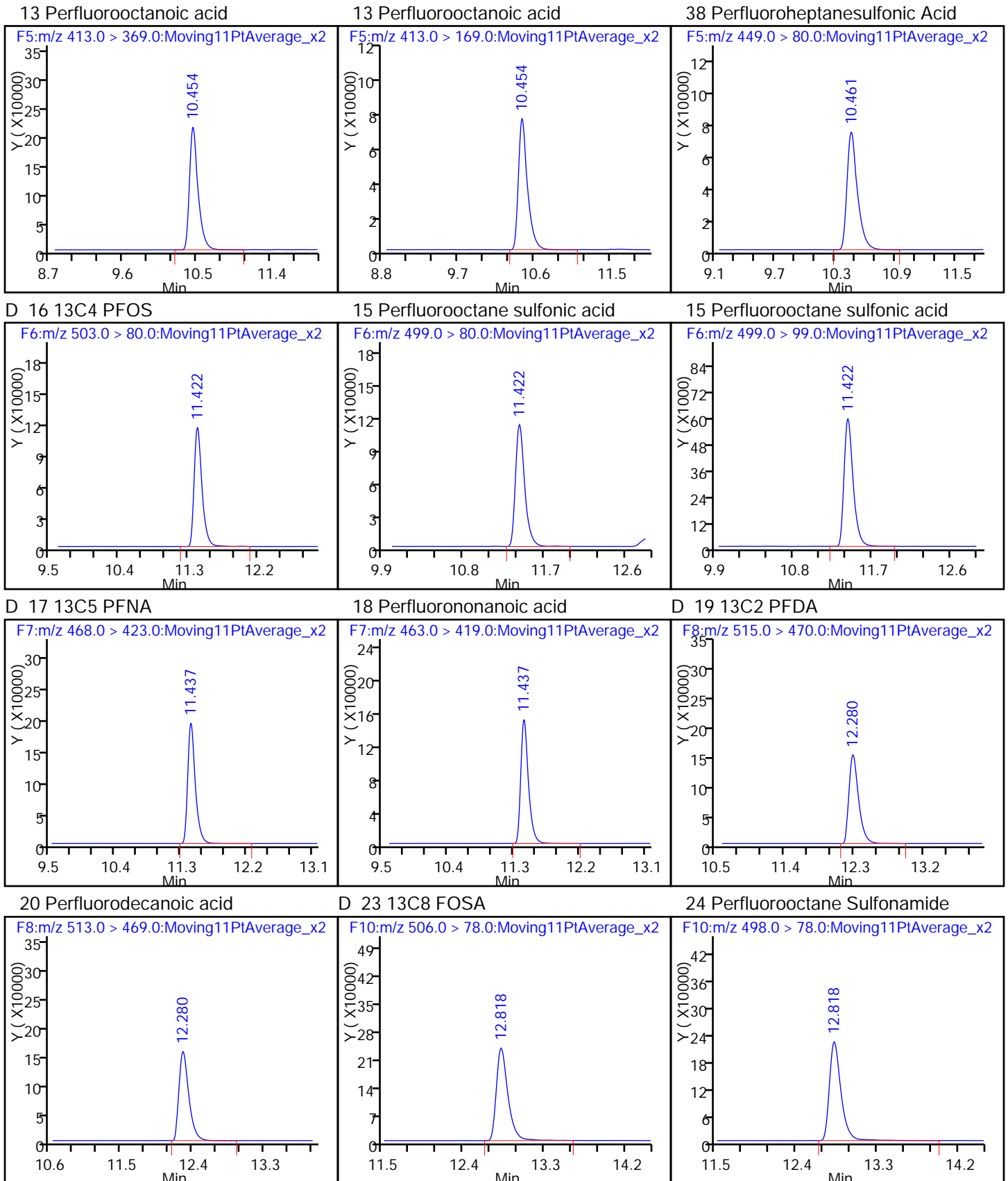


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

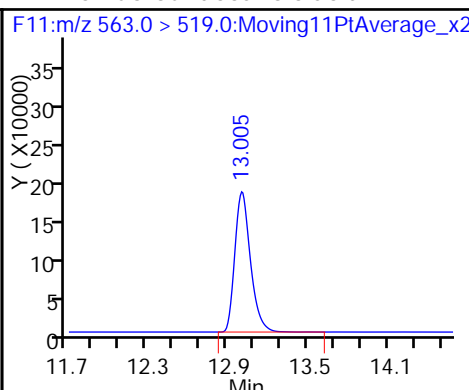
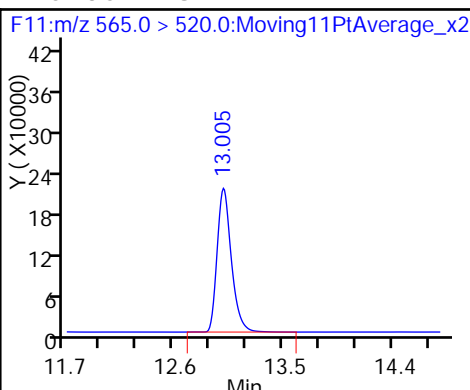
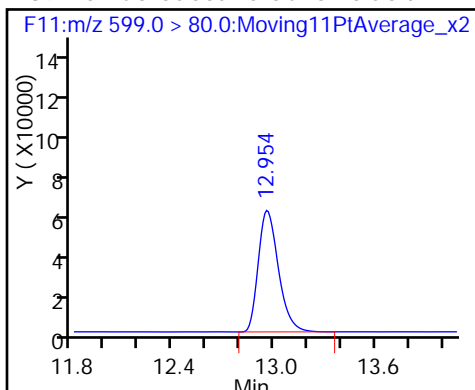




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

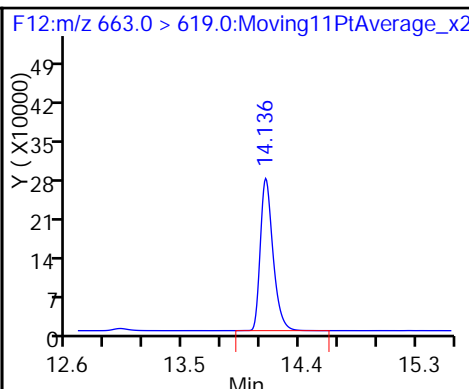
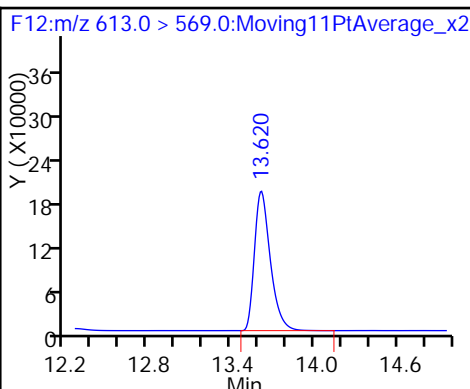
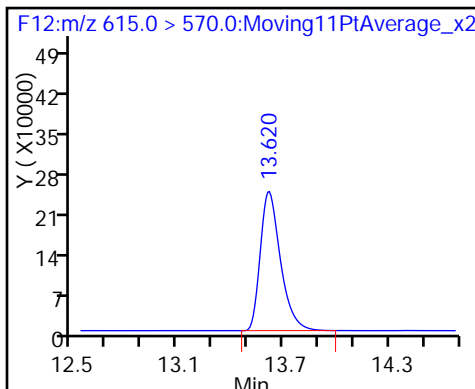
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

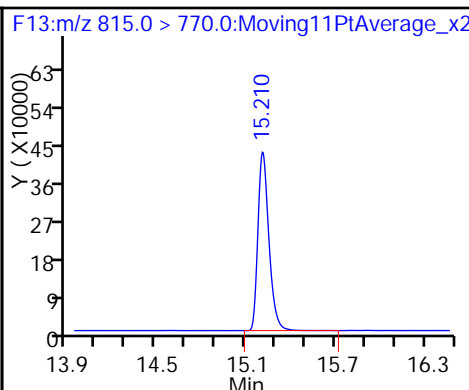
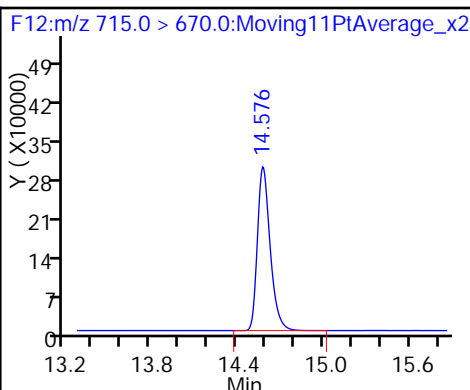
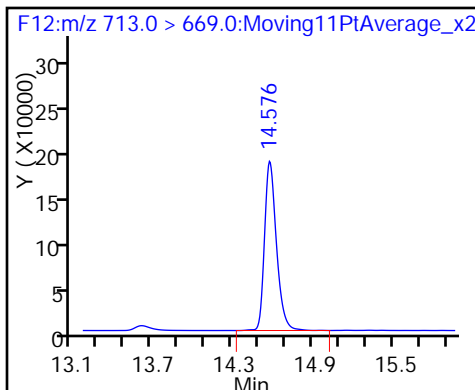
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

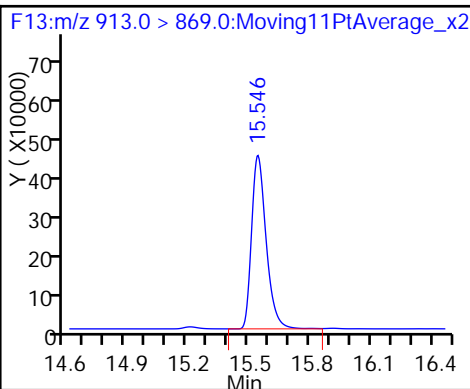
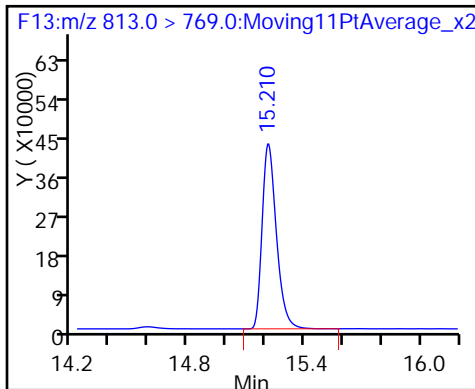
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_009.d  
 Lims ID: Std L6  
 Client ID:  
 Sample Type: IC Calib Level: 6  
 Inject. Date: 19-Feb-2016 16:54:25 ALS Bottle#: 14 Worklist Smp#: 7  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L6  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:53 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 19-Feb-2016 17:45:44

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.709	5.712	-0.003	714491	44.7		89.3	2310	
2 Perfluorobutyric acid	212.9 > 169.0	5.712	5.715	-0.003	4081745	219.0		110	10186	
D 3 13C5-PFPeA	267.9 > 223.0	6.819	6.827	-0.008	1294135	42.1		84.1	3701	
4 Perfluoropentanoic acid	262.9 > 219.0	6.819	6.828	-0.009	5303760	196.4		98.2	1105	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.934	6.939	-0.005	2482631	NC			3270	
	298.9 > 99.0	6.934	6.939	-0.005	1388611		1.79(0.00-0.00)		1814	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.934	6.939	-0.005	2482631	207.2		117		
7 Perfluorohexanoic acid	313.0 > 269.0	8.067	8.078	-0.011	5718698	205.0		103	2780	
D 6 13C2 PFHxA	315.0 > 270.0	8.067	8.080	-0.013	1287540	45.4		90.9	4384	
D 8 13C4-PFHpA	367.0 > 322.0	9.310	9.322	-0.012	1313950	42.6		85.3	4738	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.310	9.322	-0.012	5350969	195.3		97.6	8505	
D 11 18O2 PFHxS	403.0 > 84.0	9.346	9.358	-0.012	570842	38.1		80.6	2260	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.346	9.364	-0.018	1610335	NC			3760	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.346	9.364	-0.018	1610335	197.6		104		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.432	10.450	-0.018		1311023	38.4		76.8	4940	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.439	10.450	-0.011	1.000	5285370	200.8		100	4885	
413.0 > 169.0	10.439	10.450	-0.011	1.000	1833992		2.88(0.00-0.00)	100	3465	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.439	10.457	-0.018	1.000	1619742	NC			4928	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.439	10.457	-0.018	1.000	1619742	185.4		97.4		
D 16 13C4 PFOS										
503.0 > 80.0	11.399	11.414	-0.015		668257	36.8		76.9	2419	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.399	11.417	-0.018	1.000	2592636	190.7		99.7	325	
499.0 > 99.0	11.399	11.417	-0.018	1.000	1437770		1.80(0.00-0.00)	99.7	3069	
D 17 13C5 PFNA										
468.0 > 423.0	11.414	11.434	-0.020		1143055	40.5		81.1	4555	
18 Perfluorononanoic acid										
463.0 > 419.0	11.421	11.439	-0.018	1.000	3958780	214.5		107	10634	
D 19 13C2 PFDA										
515.0 > 470.0	12.254	12.275	-0.021		1048548	36.6		73.2	3638	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.254	12.276	-0.022	1.000	4149384	202.6		101	8461	
D 23 13C8 FOSA										
506.0 > 78.0	12.793	12.810	-0.017		1899691	45.8		91.5	1488	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.793	12.812	-0.019	1.000	6796748	193.0		96.5	1161	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.929	12.949	-0.020	1.000	1463184	NC			4274	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.929	12.949	-0.020	1.000	1463184	188.0		97.5		
D 26 13C2 PFUnA										
565.0 > 520.0	12.981	13.000	-0.019		1410933	37.7		75.5	4400	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.981	13.000	-0.019	1.000	4571131	196.0		98.0	6731	
D 28 13C2 PFDaA										
615.0 > 570.0	13.588	13.609	-0.021		1709160	43.7		87.4	3686	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.588	13.611	-0.023	1.000	5178123	207.2		104	5811	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.109	14.130	-0.021	1.000	5922305	184.1		92.1	4913	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.552	14.571	-0.019	1.000	3999720	195.6		97.8	2278	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.552	14.571	-0.019		1592616	44.4		88.7	6647	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.188	15.207	-0.019		1957713	45.4		90.8	4814	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.193	15.208	-0.015	1.000	7506851	202.2		101	4498	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.529	15.543	-0.014	1.000	7527521	198.2	99.1	4555	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L6\_00015

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_009.d

Injection Date: 19-Feb-2016 16:54:25

Instrument ID: A6

Lims ID: Std L6

Client ID:

Operator ID: JRB

ALS Bottle#: 14

Worklist Smp#: 7

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

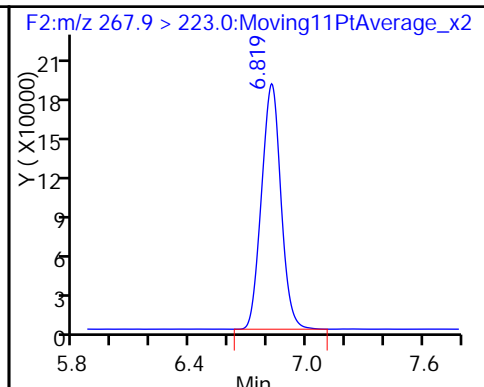
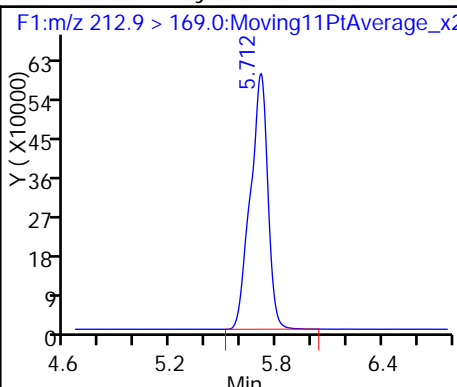
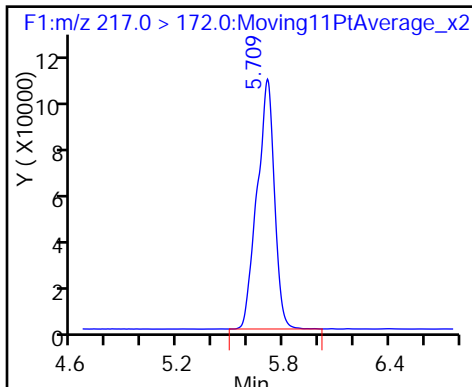
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

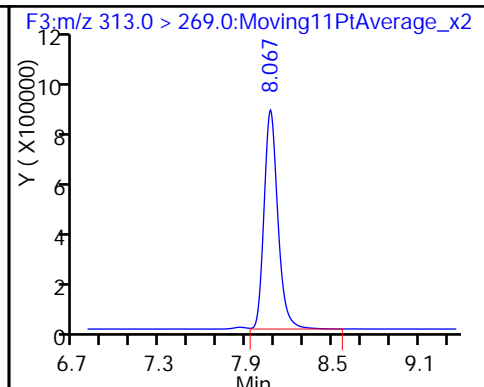
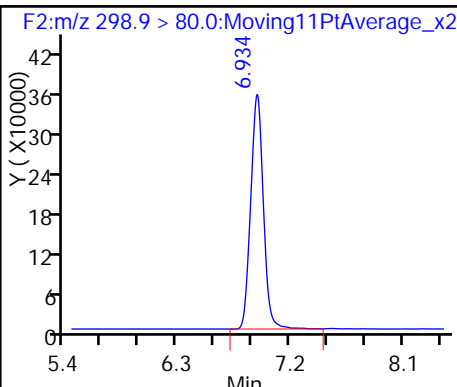
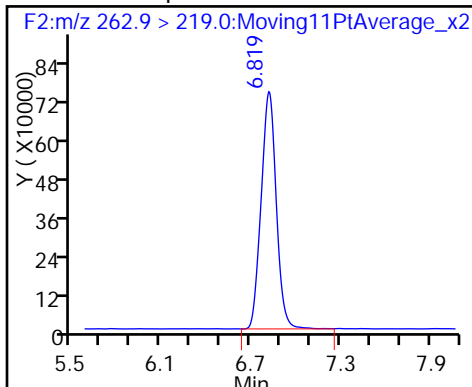
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

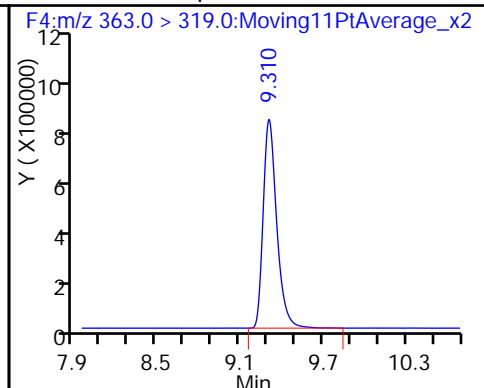
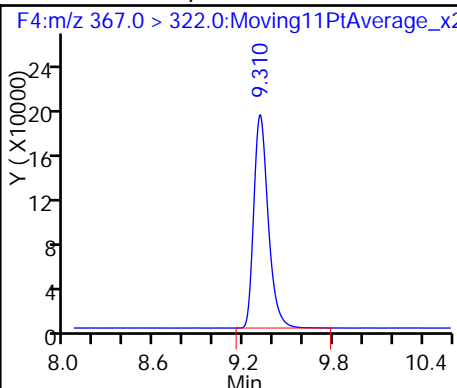
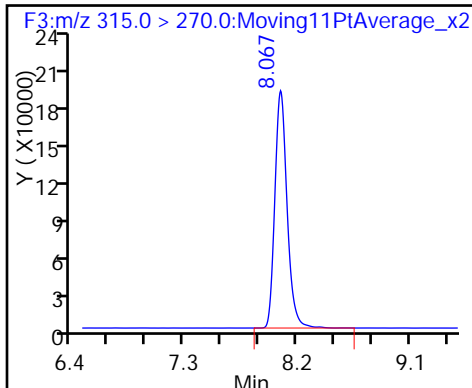
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

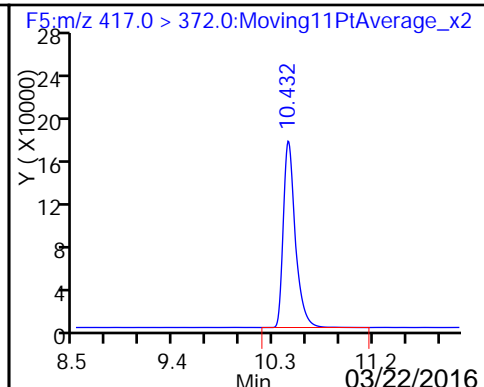
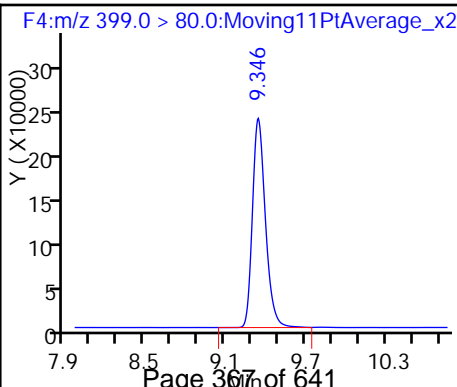
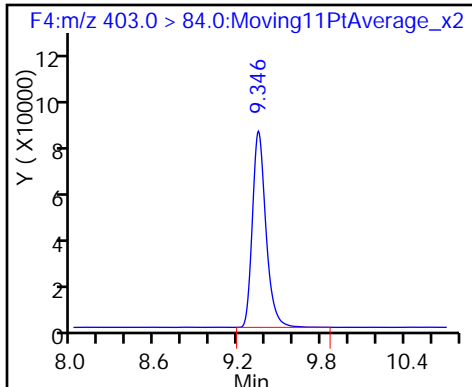
9 Perfluoroheptanoic acid

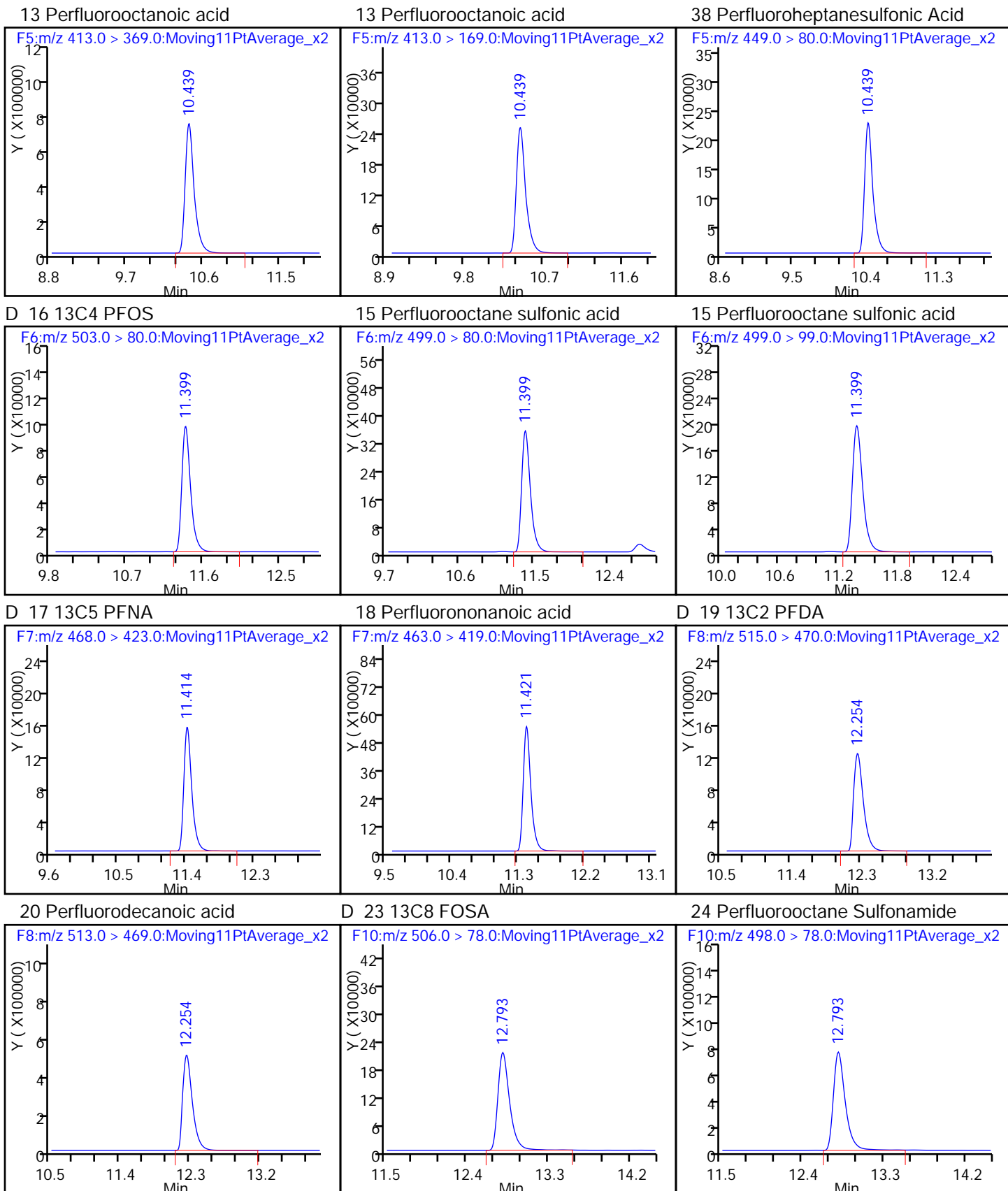


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA



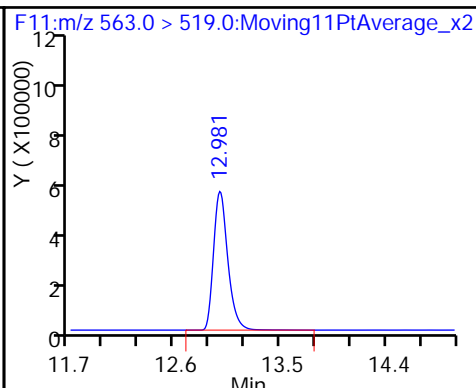
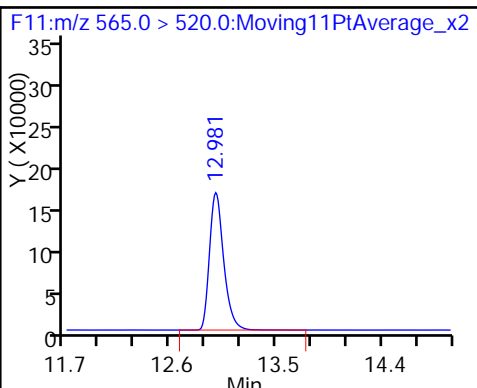
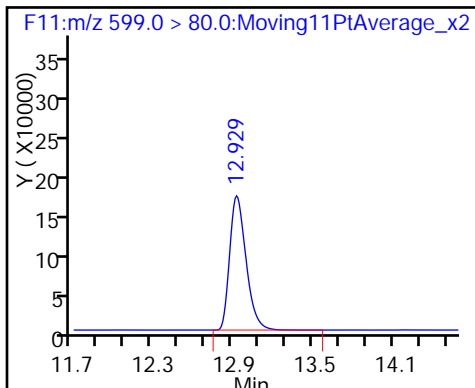




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

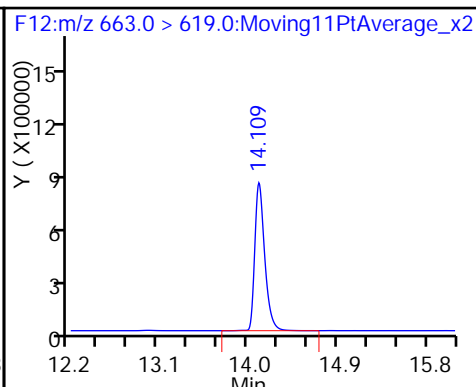
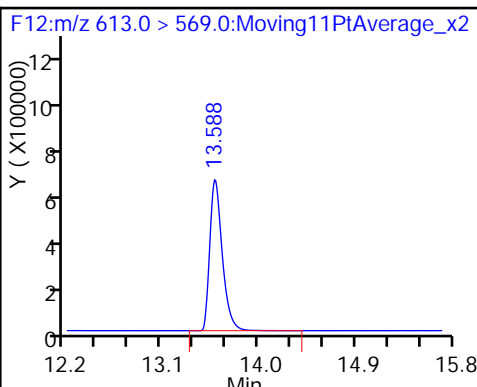
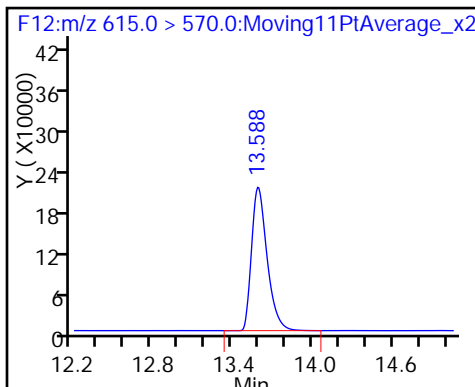
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

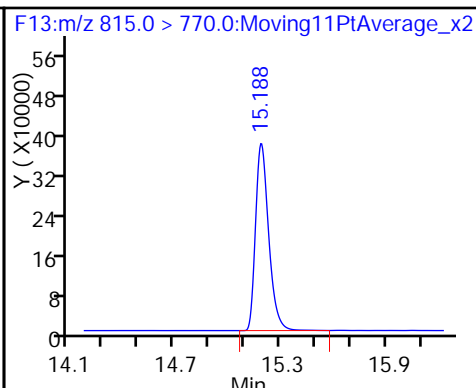
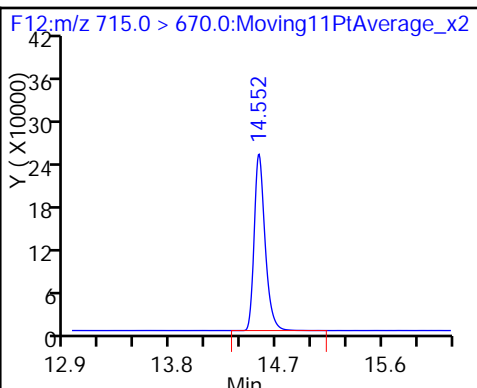
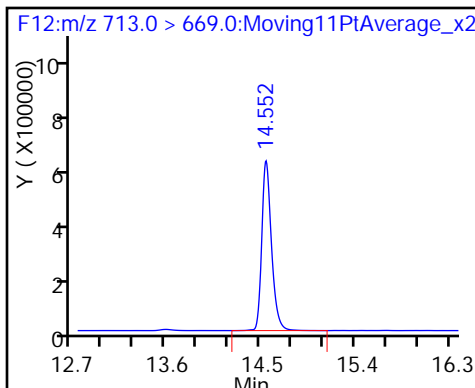
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

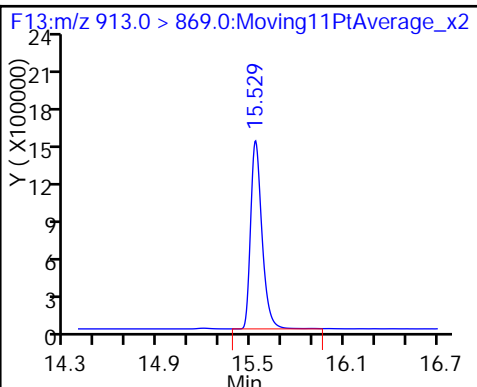
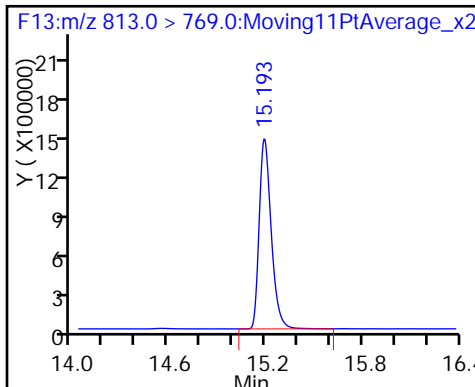
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Lims ID: Std L7  
 Client ID:  
 Sample Type: IC Calib Level: 7  
 Inject. Date: 19-Feb-2016 17:15:39 ALS Bottle#: 15 Worklist Smp#: 8  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: STD L7  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5

Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:58:55 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d

Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 19-Feb-2016 17:46:38

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.709	5.712	-0.003	631840	39.5		79.0	1919	
2 Perfluorobutyric acid	212.9 > 169.0	5.712	5.715	-0.003	7281727	441.9		110	11351	
D 3 13C5-PFPeA	267.9 > 223.0	6.818	6.827	-0.009	1171928	38.1		76.2	3691	
4 Perfluoropentanoic acid	262.9 > 219.0	6.818	6.828	-0.010	9604476	392.7		98.2	1997	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.933	6.939	-0.006	4163359	NC			2718	
	298.9 > 99.0	6.933	6.939	-0.006	2275031		1.83(0.00-0.00)		3395	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.933	6.939	-0.006	4163359	435.6		123		
7 Perfluorohexanoic acid	313.0 > 269.0	8.065	8.078	-0.013	9481665	420.7		105	2241	
D 6 13C2 PFHxA	315.0 > 270.0	8.065	8.080	-0.015	1040351	36.7		73.4	2818	
D 8 13C4-PFHpA	367.0 > 322.0	9.305	9.322	-0.017	1099321	35.7		71.3	3231	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.305	9.322	-0.017	9198751	401.2		100	11672	
D 11 18O2 PFHxS	403.0 > 84.0	9.346	9.358	-0.012	455229	30.4		64.3	1441	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.346	9.364	-0.018	2660458	NC			6698	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.346	9.364	-0.018	2660458	409.4		108		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.433	10.450	-0.017		1023740	30.0		60.0	3518	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.433	10.450	-0.017	1.000	8666978	421.7		105	4474	
413.0 > 169.0	10.433	10.450	-0.017	1.000	2909884		2.98(0.00-0.00)	105	3285	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.440	10.457	-0.017	1.000	2666346	NC			7584	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.440	10.457	-0.017	1.000	2666346	381.1		100		
D 16 13C4 PFOS										
503.0 > 80.0	11.392	11.414	-0.022		534710	29.4		61.5	2475	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.399	11.417	-0.018	1.000	4278056	393.3		103	332	
499.0 > 99.0	11.399	11.417	-0.018	1.000	2357374		1.81(0.00-0.00)	103	3975	
D 17 13C5 PFNA										
468.0 > 423.0	11.415	11.434	-0.019		1010032	35.8		71.7	2370	
18 Perfluorononanoic acid										
463.0 > 419.0	11.422	11.439	-0.017	1.000	6650814	407.7		102	13388	
D 19 13C2 PFDA										
515.0 > 470.0	12.255	12.275	-0.020		925036	32.3		64.6	1991	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.255	12.276	-0.021	1.000	6815253	377.5		94.4	4656	
D 23 13C8 FOSA										
506.0 > 78.0	12.793	12.810	-0.017		1599127	38.5		77.1	1713	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.793	12.812	-0.019	1.000	11722467	395.5		98.9	1678	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.929	12.949	-0.020	1.000	2170020	NC			3636	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.929	12.949	-0.020	1.000	2170020	348.2		90.3		
D 26 13C2 PFUnA										
565.0 > 520.0	12.981	13.000	-0.019		1172072	31.4		62.7	3708	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.981	13.000	-0.019	1.000	7799843	403.3		101	7175	
D 28 13C2 PFDaA										
615.0 > 570.0	13.588	13.609	-0.021		1451296	37.1		74.2	3521	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.588	13.611	-0.023	1.000	9187053	432.9		108	6144	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.109	14.130	-0.021	1.000	9569869	350.4		87.6	4611	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.552	14.571	-0.019	1.000	6913347	398.6		99.7	3248	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.552	14.571	-0.019		1521958	42.4		84.8	4569	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.193	15.207	-0.014		1815999	42.1		84.3	3968	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.193	15.208	-0.015	1.000	12867499	410.8		103	4108	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.529	15.543	-0.014	1.000	13652743	423.4	106	3366	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L7\_00015

Amount Added: 1.00

Units: mL

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d

Injection Date: 19-Feb-2016 17:15:39

Instrument ID: A6

Lims ID: Std L7

Client ID:

Operator ID: JRB

ALS Bottle#: 15

Worklist Smp#: 8

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

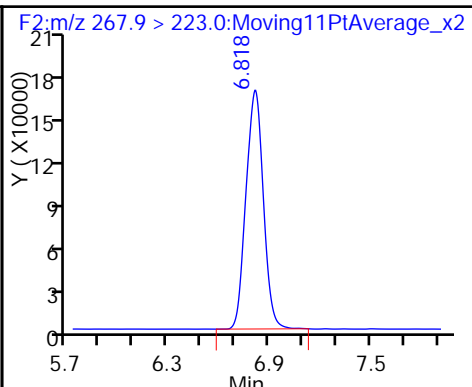
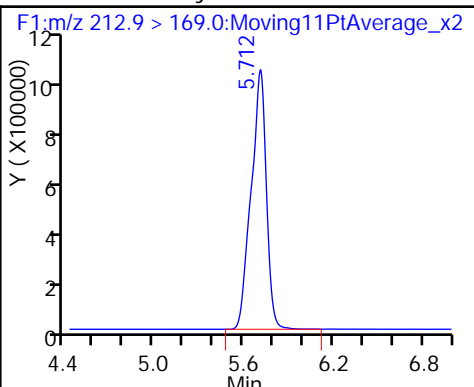
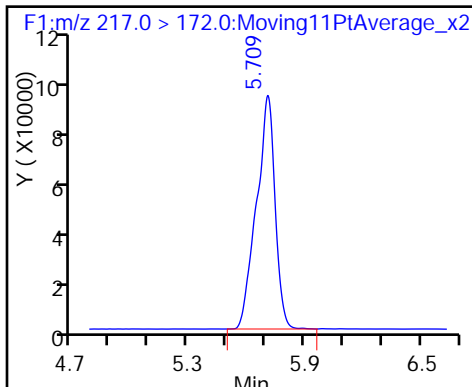
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

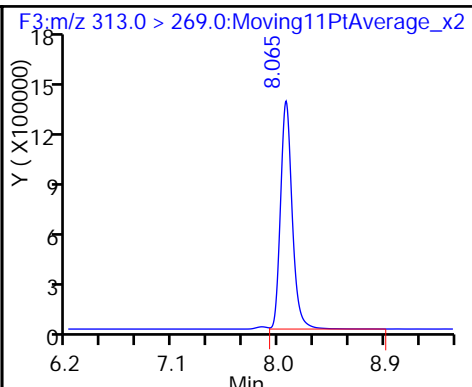
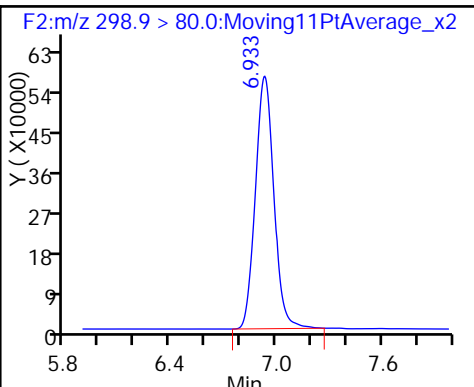
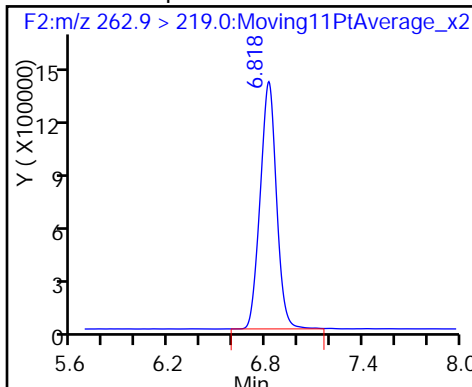
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

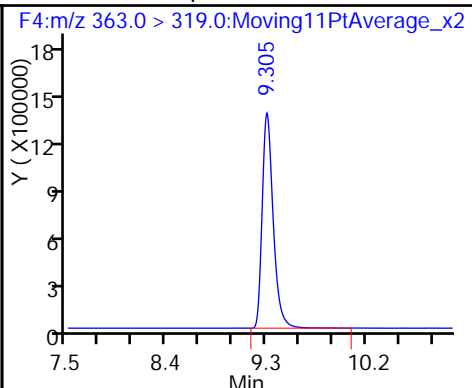
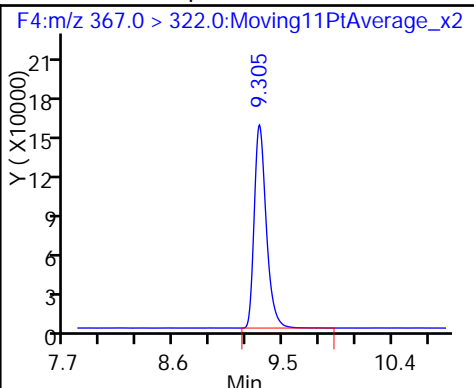
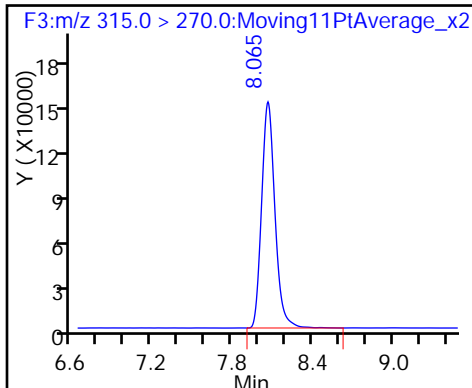
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

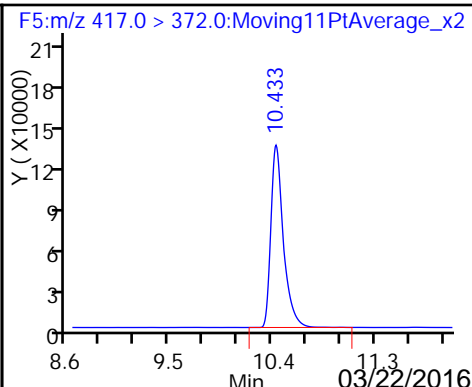
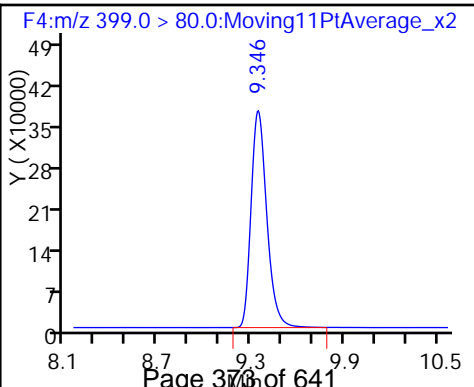
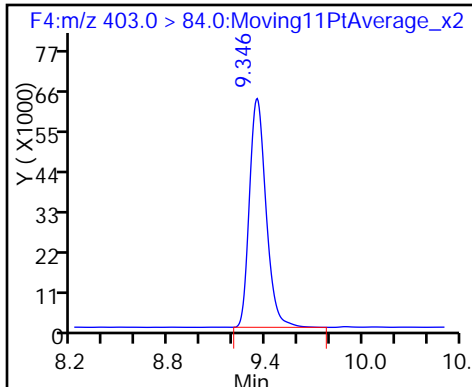
9 Perfluoroheptanoic acid

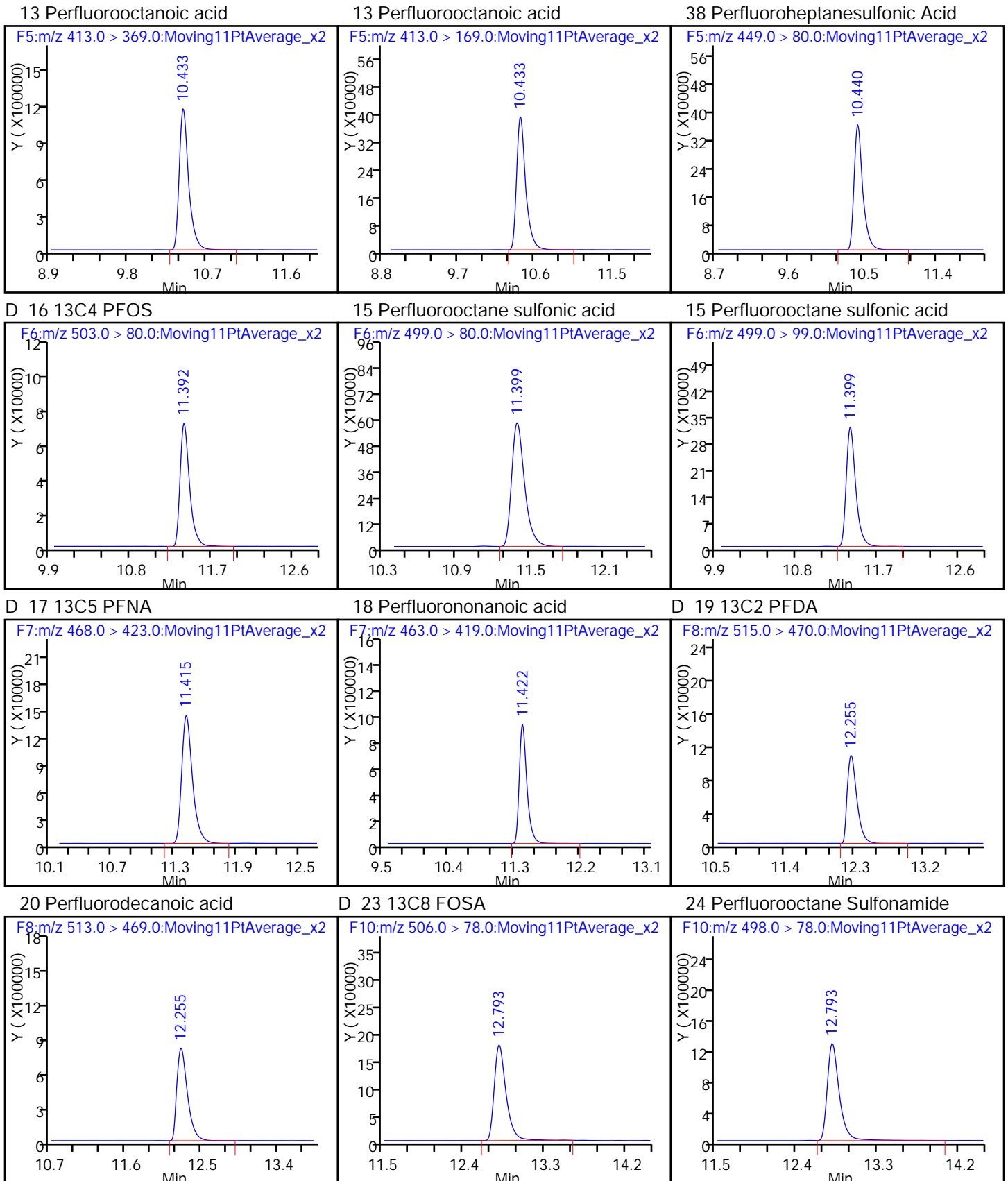


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

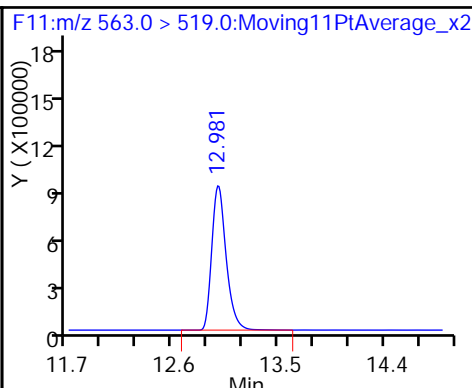
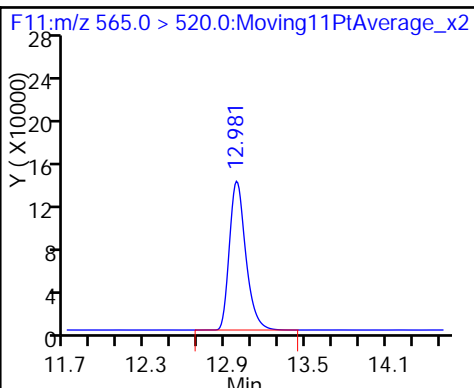
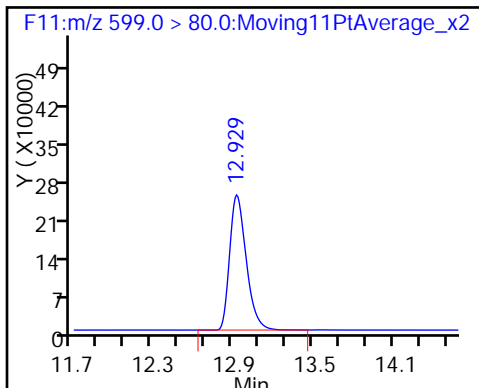




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

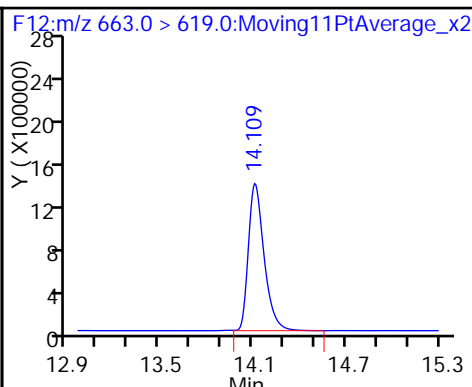
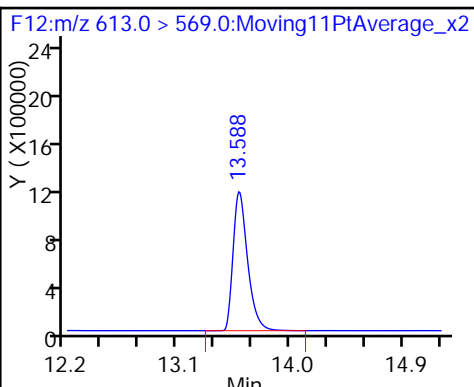
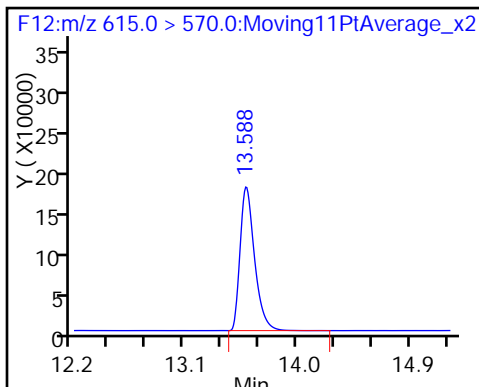
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

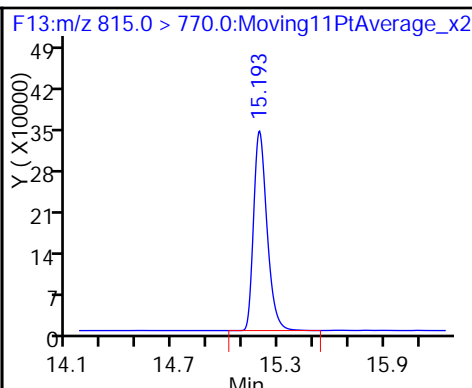
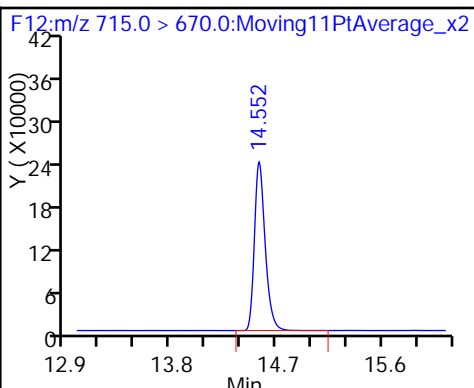
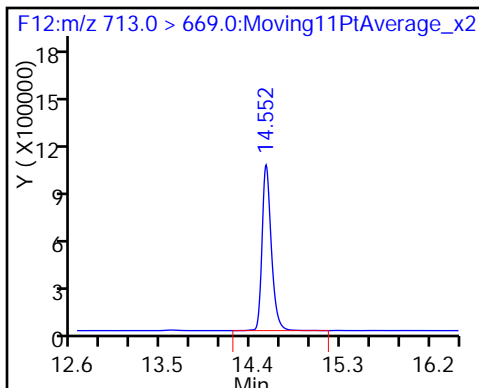
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

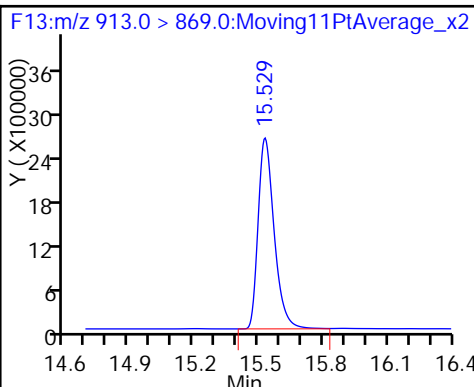
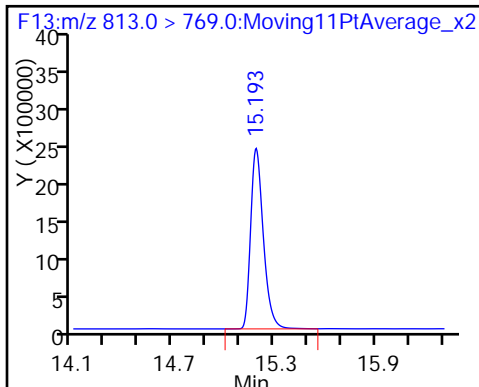
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: ICV 320-100305/10 Calibration Date: 02/12/2016 13:24  
 Instrument ID: A4 Calib Start Date: 02/12/2016 10:35  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/12/2016 12:42  
 Lab File ID: 12FEB2016A4A\_011.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	0.4901	0.5679		57.9	50.0	15.9	25.0
Perfluoropentanoic acid (PFPeA)	AveID	0.4708	0.5472		58.1	50.0	16.2	25.0
Perfluorobutanesulfonic acid (PFBS)	L1ID		0.8082		42.8	44.3	-3.4	25.0
Perfluorohexanoic acid (PFHxA)	AveID	0.4308	0.4788		55.6	50.0	11.2	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	0.5120	0.4775		46.6	50.0	-6.8	25.0
Perfluorohexanesulfonic acid (PFHxS)	L2ID		1.199		50.6	47.3	7.1	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	AveID	2.544	2.580		48.3	47.6	1.4	25.0
Perfluorooctanoic acid (PFOA)	AveID	0.4910	0.4815		49.0	50.0	-1.9	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	4.265	5.017		56.2	47.8	17.6	25.0
Perfluorononanoic acid (PFNA)	L1ID		1.133		54.0	50.0	8.0	25.0
Perfluorodecanoic acid (PFDA)	AveID	0.9142	0.9011		49.3	50.0	-1.4	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9450	1.016		53.7	50.0	7.5	25.0
Perfluorodecane Sulfonic acid	AveID	1.927	2.087		52.3	48.3	8.3	25.0
Perfluoroundecanoic acid (PFUnA)	AveID	1.037	1.039		50.1	50.0	0.1	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.8053	0.8703		54.0	50.0	8.1	25.0
Perfluorotridecanoic Acid (PFTriA)	L2ID		0.8591		51.5	50.0	3.0	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.3819		50.2	50.0	0.5	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L1ID		1.056		48.8	50.0	-2.3	25.0
Perfluoro-n-octadecanoic acid (PFODA)	L1ID		1.026		51.2	50.0	2.3	25.0



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_011.d  
 Lims ID: ICV  
 Client ID:  
 Sample Type: ICV  
 Inject. Date: 12-Feb-2016 13:24:55 ALS Bottle#: 9 Worklist Smp#: 10  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: ICV  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub6  
 Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:03:56 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d  
 Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: barnettj Date: 12-Feb-2016 17:40:03

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	216.7 > 171.5	5.729	5.712	0.017	1908642	47.0		94.1	6779	
2 Perfluorobutyric acid	212.7 > 168.6	5.732	5.712	0.020	1083841	57.9			3338	
4 Perfluoropentanoic acid	262.9 > 218.7	7.026	6.986	0.040	842323	58.1			647	
D 3 13C5-PFPeA	267.6 > 222.7	7.020	6.986	0.034	1539466	48.1		96.2	2904	
5 Perfluorobutane Sulfonate	298.8 > 79.6	7.155	7.115	0.040	800699	NC			536	
	298.8 > 98.6	7.155	7.115	0.040	553707		1.45(0.00-0.00)		494	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	7.155	7.115	0.040	800699	42.8				
D 6 13C2 PFHxA	314.6 > 269.7	8.383	8.367	0.016	1874527	47.9		95.7	4493	
7 Perfluorohexanoic acid	312.9 > 268.7	8.389	8.369	0.020	897607	55.6			1409	
22 PFPeS (Perflouro-1-pentanesulfonat	348.7 > 79.5	8.470	8.676	-0.206	982357	NC			2741	
D 8 13C4-PFHpA	366.6 > 321.6	9.643	9.639	0.004	1861616	47.8		95.6	4264	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.643	9.640	0.003	888852	46.6			1623	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.671	9.675	-0.004	1267917	50.6				
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.671	9.675	-0.004	1267917	NC			3702	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 11 18O2 PFHxS										
402.5 > 83.6	9.671	9.675	-0.004		1059018	46.7		98.7	2658	
39 Perfluoroheptanesulfonic Acid										
448.3 > 79.2	10.724	10.740	-0.016	1.000	1248257	48.3				
14 Perfluoroheptane Sulfonate										
448.3 > 79.2	10.724	10.740	-0.016	1.000	1248257	NC			3638	
D 12 13C4 PFOA										
416.5 > 371.6	10.724	10.743	-0.019		1999016	47.0		93.9	3452	
13 Perfluorooctanoic acid										
412.8 > 368.8	10.724	10.743	-0.019	1.000	962485	49.0			1436	
412.8 > 168.7	10.733	10.743	-0.010	1.001	366426		2.63(0.00-0.00)		845	
D 16 13C4 PFOS										
502.4 > 79.7	11.618	11.651	-0.033		485944	48.0		100	2096	
15 Perfluorooctane sulfonic acid										
498.3 > 79.2	11.626	11.656	-0.030	1.000	2435496	56.2			2710	
498.3 > 98.2	11.626	11.656	-0.030	1.000	1419627		1.72(0.00-0.00)		2054	
D 17 13C5 PFNA										
467.5 > 422.6	11.644	11.678	-0.034		1986876	47.8		95.5	3977	
18 Perfluorononanoic acid										
462.5 > 418.6	11.653	11.680	-0.027	1.000	2250334	54.0			2036	
D 19 13C2 PFDA										
514.4 > 469.5	12.434	12.473	-0.039		2782767	47.4		94.7	3747	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.434	12.476	-0.042	1.000	2507408	49.3			3108	
21 PFNS (Perflouro-1-nonanesulfonate)										
548.6 > 79.6	12.393	12.580	-0.187	1.000	762915	NC			2195	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	12.952	12.991	-0.039	1.000	4381313	53.7			1820	
D 23 13C8 FOSA										
505.4 > 77.6	12.952	12.991	-0.039		4313351	48.3		96.6	1996	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.058	13.098	-0.040	1.000	1023926	NC			3381	
49 Perfluorodecane Sulfonic acid										
598.4 > 79.6	13.058	13.098	-0.040	1.000	1023926	52.3				
D 26 13C2 PFUnA										
564.3 > 519.5	13.111	13.146	-0.035		2868368	49.0		98.1	3326	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.111	13.146	-0.035	1.000	2979313	50.1			2312	
D 28 13C2 PFDaA										
614.4 > 569.4	13.684	13.713	-0.029		3530689	49.6		99.3	3478	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.684	13.715	-0.031	1.000	3072566	54.0			1465	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.167	14.193	-0.026	1.000	3033306	51.5			1491	
31 PFDaS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.113	14.227	-0.114	1.000	315882	NC			1447	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.576	14.603	-0.027		3288418	51.9		104	3923	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.576	14.604	-0.028	1.000	1348229	50.2		890	
D 35 13C2-PFHxDA	814.8 > 769.6	15.124	15.138	-0.014		1776974	59.9	120	4070	
34 Perfluorohexadecanoic acid	812.6 > 768.6	15.124	15.139	-0.015	1.000	3730056	48.8		2582	
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.426	15.442	-0.016	1.000	3621991	51.2		2372	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFCIC\_00014

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_011.d

Injection Date: 12-Feb-2016 13:24:55

Instrument ID: A4

Lims ID: ICV

Client ID:

Operator ID: JRB

ALS Bottle#: 9

Worklist Smp#: 10

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

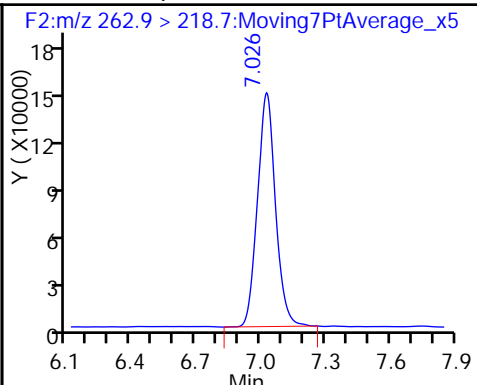
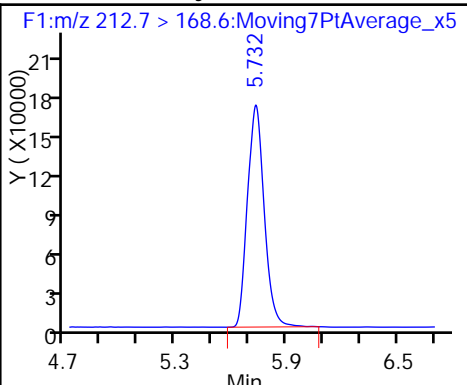
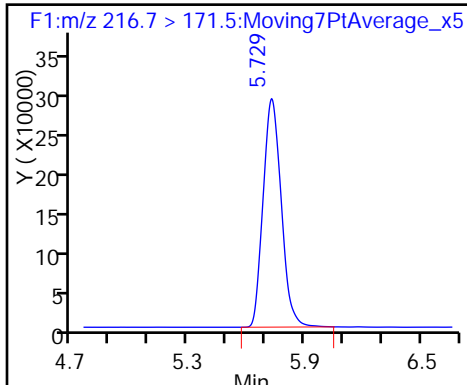
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

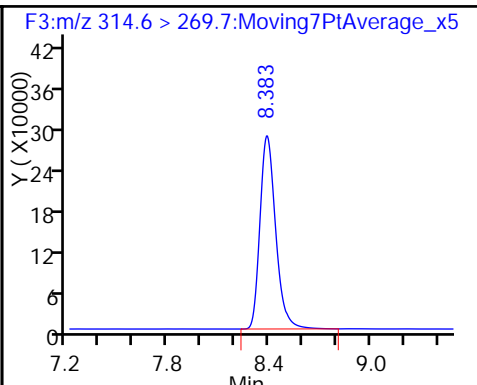
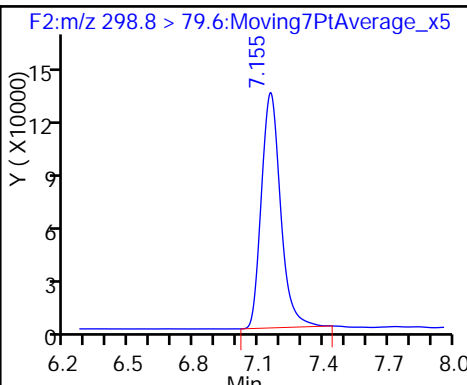
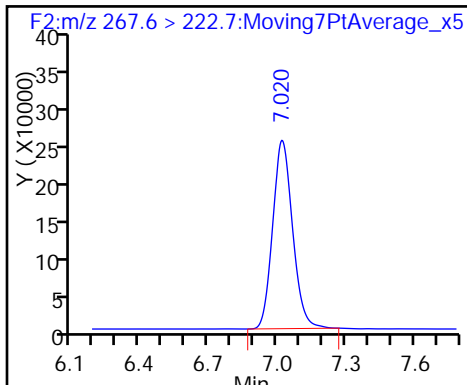
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

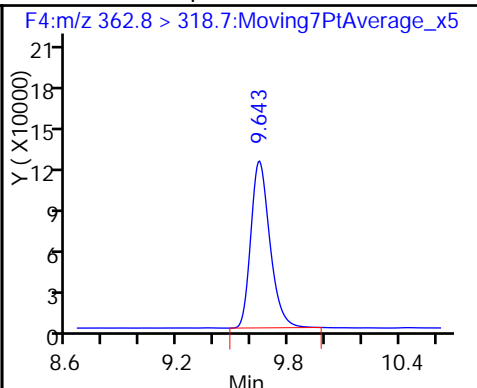
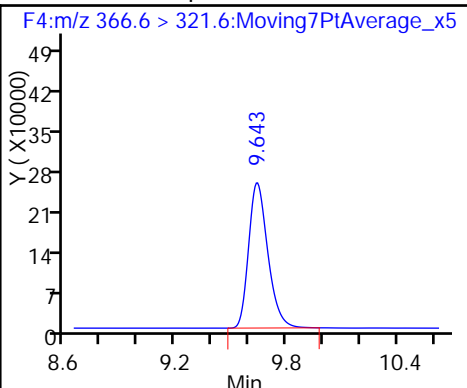
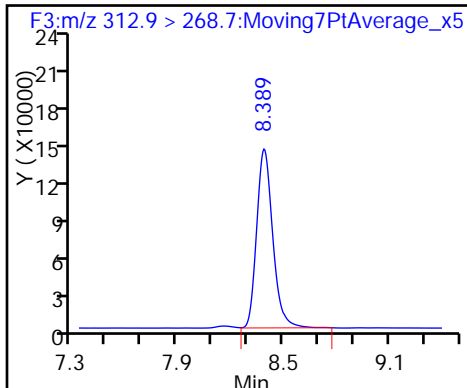
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

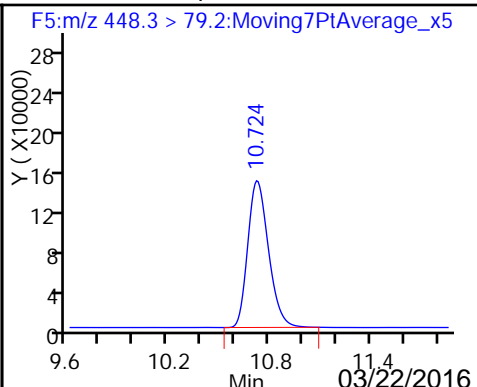
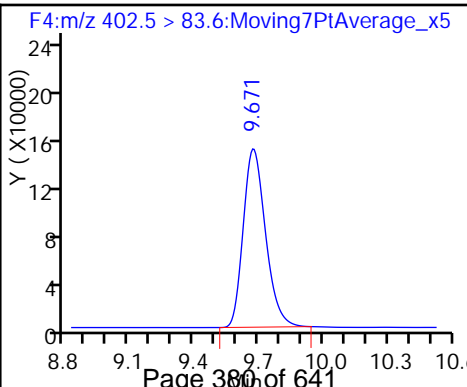
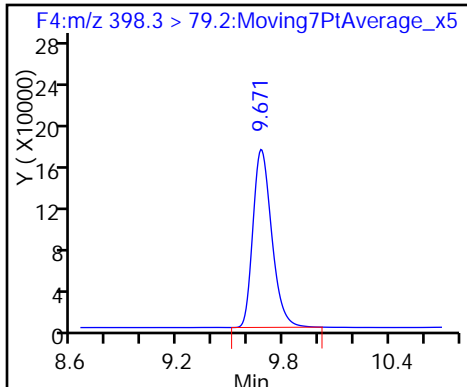
9 Perfluoroheptanoic acid



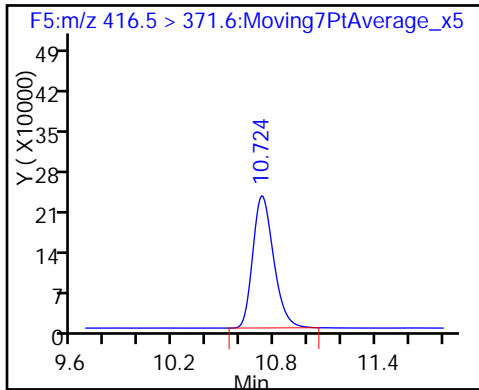
58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

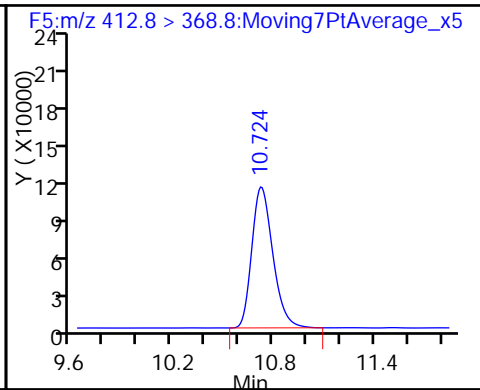
39 Perfluoroheptanesulfonic Acid



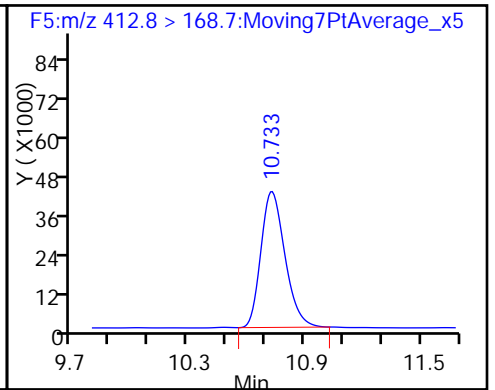
D 12 13C4 PFOA



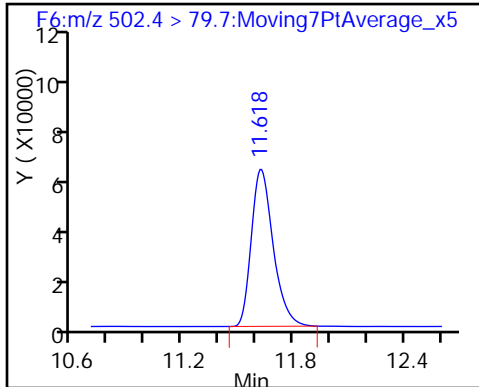
13 Perfluorooctanoic acid



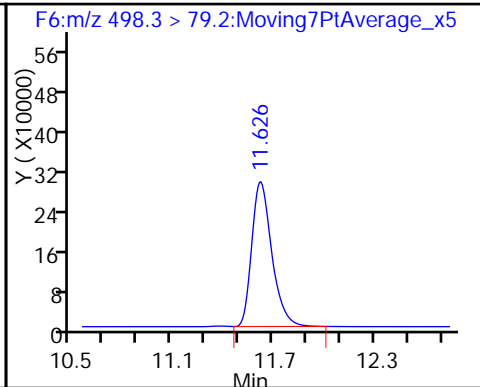
13 Perfluorooctanoic acid



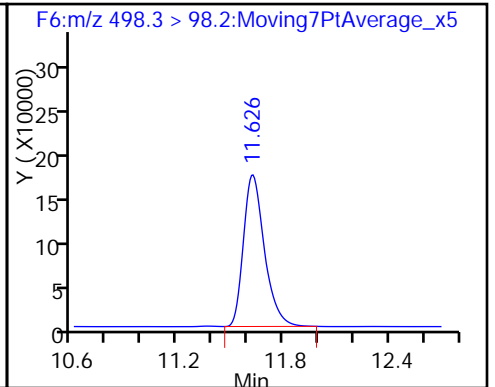
D 16 13C4 PFOS



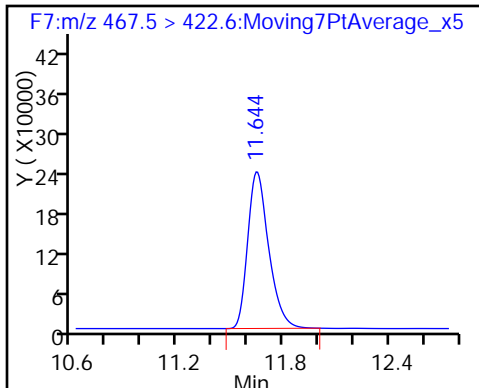
15 Perfluorooctane sulfonic acid



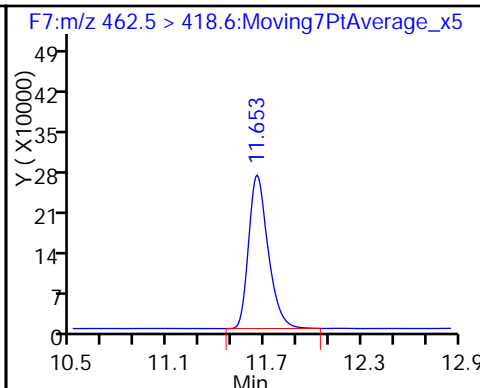
15 Perfluorooctane sulfonic acid



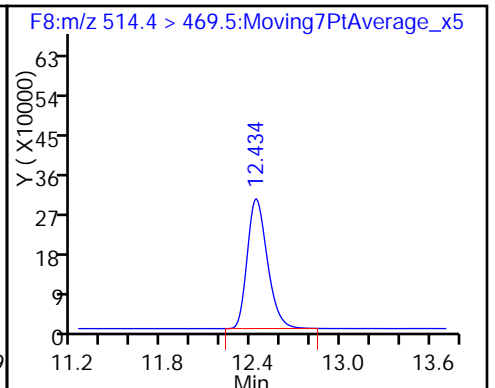
D 17 13C5 PFNA



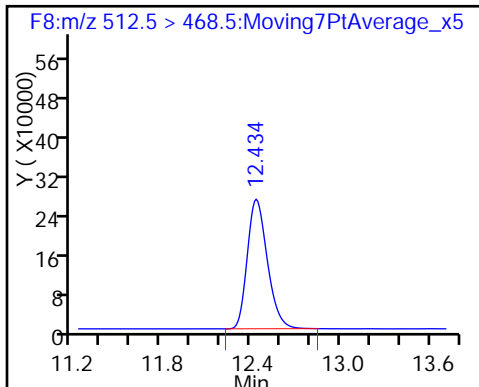
18 Perfluorononanoic acid



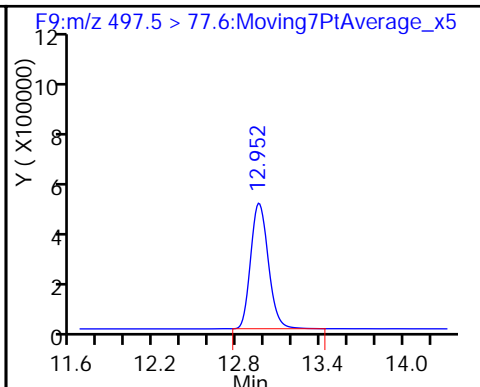
D 19 13C2 PFDA



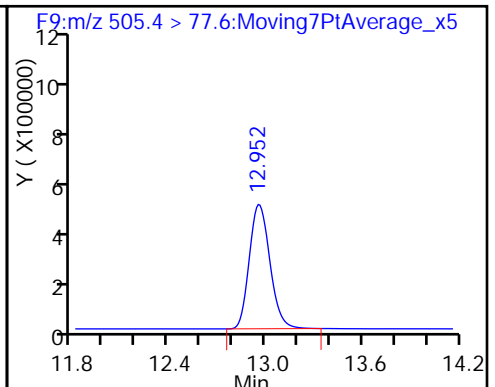
20 Perfluorodecanoic acid



24 Perfluorooctane Sulfonamide



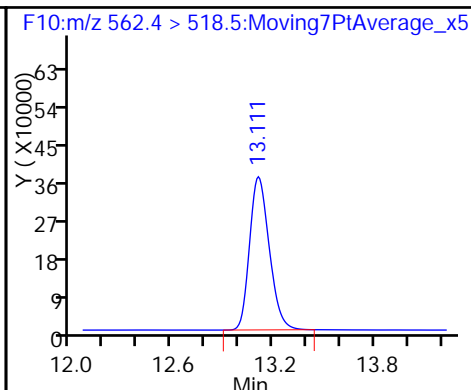
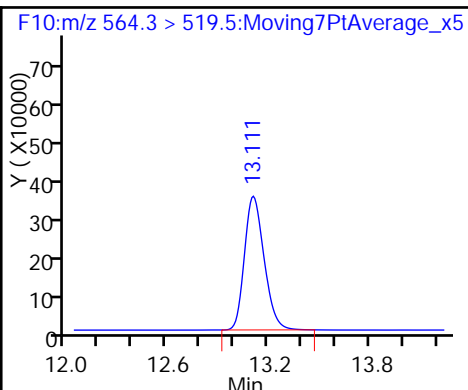
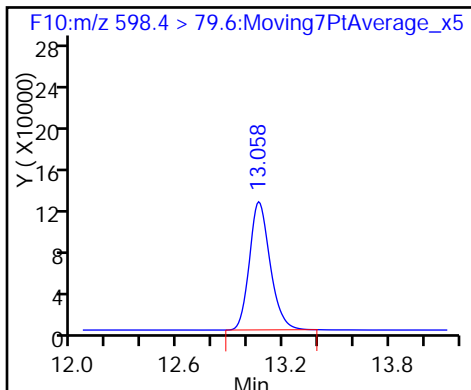
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

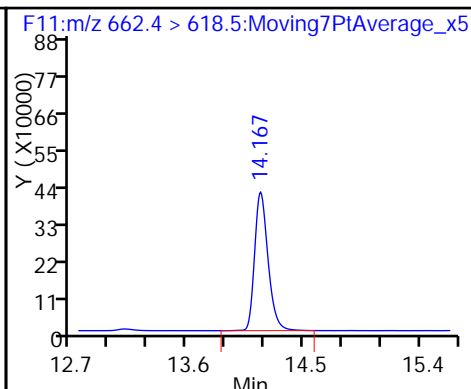
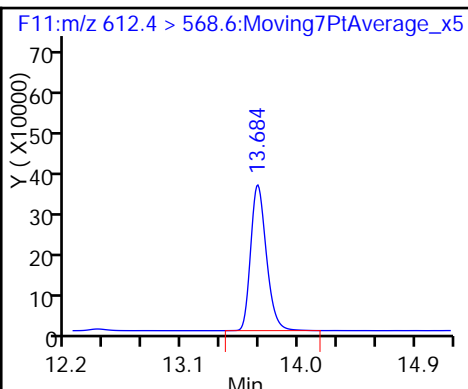
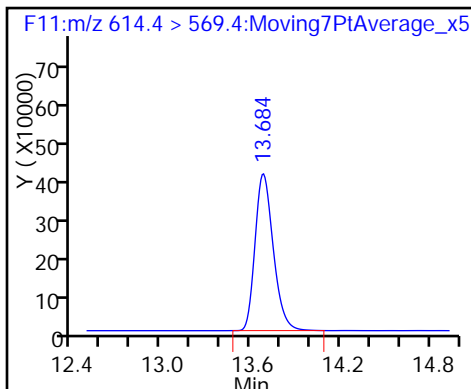
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

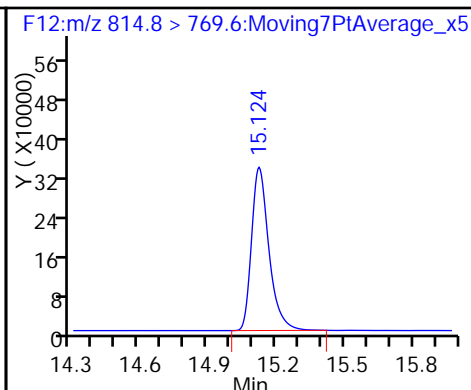
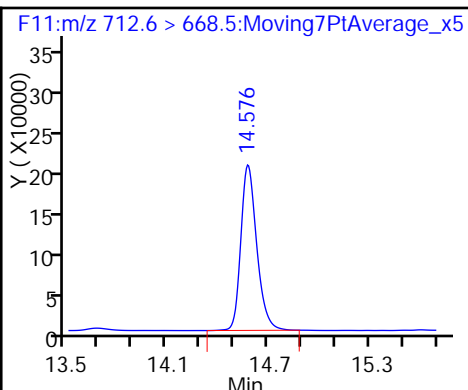
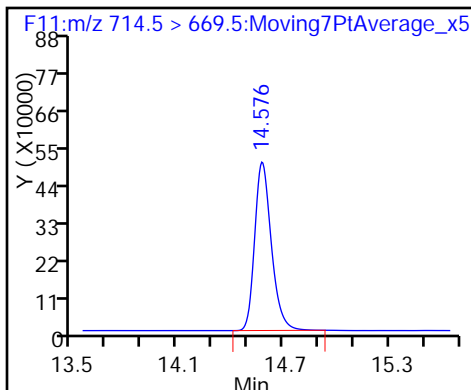
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

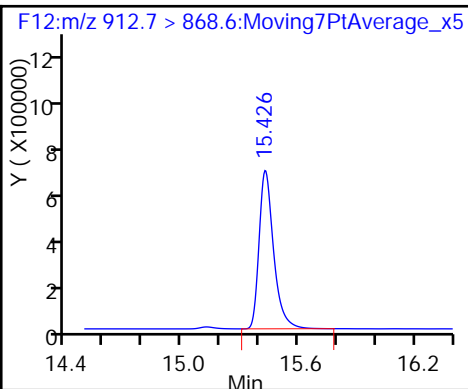
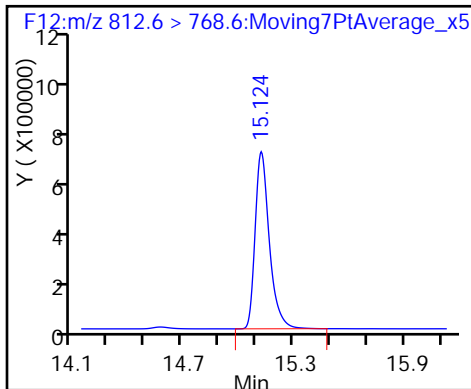
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: CCV 320-100305/40 Calibration Date: 02/13/2016 00:01  
 Instrument ID: A4 Calib Start Date: 02/12/2016 10:35  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/12/2016 12:42  
 Lab File ID: 12FEB2016A4A\_041.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	0.4901	0.5313		21.7	20.0	8.4	25.0
Perfluoropentanoic acid (PFPeA)	AveID	0.4708	0.4801		20.4	20.0	2.0	25.0
Perfluorobutanesulfonic acid (PFBS)	L1ID		1.029		14.7	17.7	-17.1	25.0
Perfluorohexanoic acid (PFHxA)	AveID	0.4308	0.4473		20.8	20.0	3.8	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	0.5120	0.5030		19.6	20.0	-1.8	25.0
Perfluorohexanesulfonic acid (PFHxS)	L2ID		1.130		18.9	18.9	-0.0	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	AveID	2.544	3.438		25.7	19.0	35.1*	25.0
Perfluorooctanoic acid (PFOA)	AveID	0.4910	0.4308		17.5	20.0	-12.3	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	4.265	5.395		24.2	19.1	26.5*	25.0
Perfluorononanoic acid (PFNA)	L1ID		1.160		22.2	20.0	11.0	25.0
Perfluorodecanoic acid (PFDA)	AveID	0.9142	0.8678		19.0	20.0	-5.1	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9450	1.005		21.3	20.0	6.4	25.0
Perfluorodecane Sulfonic acid	AveID	1.927	2.266		22.7	19.3	17.6	25.0
Perfluoroundecanoic acid (PFUnA)	AveID	1.037	0.9717		18.7	20.0	-6.3	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.8053	0.8602		21.4	20.0	6.8	25.0
Perfluorotridecanoic Acid (PFTriA)	L2ID		0.8802		21.1	20.0	5.5	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.3595		18.8	20.0	-6.2	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L1ID		0.8891		16.3	20.0	-18.3	25.0
Perfluoro-n-octadecanoic acid (PFODA)	L1ID		0.7773		15.9	20.0	-20.6	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_041.d  
 Lims ID: CCV L4  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 13-Feb-2016 00:01:23 ALS Bottle#: 47 Worklist Smp#: 40  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L4  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12  
 Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 10:07:12 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d  
 Column 1 : Det: F1:MRM  
 Process Host: XAWRK007

First Level Reviewer: westendorfc Date: 15-Feb-2016 09:21:48

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	216.7 > 171.5	5.570	5.712	-0.142	1932591	47.6		95.2	7956	
2 Perfluorobutyric acid	212.7 > 168.6	5.570	5.712	-0.142	410725	21.7		108	925	
4 Perfluoropentanoic acid	262.9 > 218.7	6.769	6.986	-0.217	283463	20.4		102	294	
D 3 13C5-PFPeA	267.6 > 222.7	6.764	6.986	-0.222	1476036	46.1		92.3	3796	
5 Perfluorobutane Sulfonate	298.8 > 79.6	6.888	7.115	-0.227	383963	NC			238	
	298.8 > 98.6	6.884	7.115	-0.231	255794		1.50(0.00-0.00)		197	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	6.888	7.115	-0.227	383963	14.7		82.9		
D 6 13C2 PFHxA	314.6 > 269.7	8.066	8.367	-0.301	1836068	46.9		93.8	4895	
7 Perfluorohexanoic acid	312.9 > 268.7	8.066	8.369	-0.303	328539	20.8		104	1083	
D 8 13C4-PFHpA	366.6 > 321.6	9.305	9.639	-0.334	1813051	46.6		93.1	4180	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.305	9.640	-0.335	364779	19.6		98.2	898	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.340	9.675	-0.335	451415	18.9		100.0		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.340	9.675	-0.335	451415	NC			1169	
D 11 18O2 PFHxS	402.5 > 83.6	9.340	9.675	-0.335	998501	44.0		93.1	2487	



Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.391	10.740	-0.349	1.000	549835	25.7	135		
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.391	10.740	-0.349	1.000	549835	NC		2176	
D 12 13C4 PFOA	416.5 > 371.6	10.391	10.743	-0.352		2224559	52.2	104	5414	
13 Perfluorooctanoic acid	412.8 > 368.8	10.391	10.743	-0.352	1.000	383302	17.5	87.7	981	
D 16 13C4 PFOS	502.4 > 79.7	11.313	11.651	-0.338		401507	39.6	82.9	1651	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.313	11.656	-0.343	1.000	866386	24.2	126	1548	
D 17 13C5 PFNA	467.5 > 422.6	11.336	11.678	-0.342		1963354	47.2	94.4	6100	
18 Perfluorononanoic acid	462.5 > 418.6	11.336	11.680	-0.344	1.000	911333	22.2	111	1384	
D 19 13C2 PFDA	514.4 > 469.5	12.144	12.473	-0.329		2887941	49.2	98.3	4099	
20 Perfluorodecanoic acid	512.5 > 468.5	12.144	12.476	-0.332	1.000	1002452	19.0	94.9	2524	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	12.673	12.991	-0.318	1.000	1499415	21.3	106	1343	
D 23 13C8 FOSA	505.4 > 77.6	12.673	12.991	-0.318		3729431	41.8	83.5	1371	
25 Perfluorodecane Sulfonate	598.4 > 79.6	12.789	13.098	-0.309	1.000	366949	NC		1105	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	12.789	13.098	-0.309	1.000	366949	22.7	118		
D 26 13C2 PFUnA	564.3 > 519.5	12.830	13.146	-0.316		2770335	47.4	94.7	2682	
27 Perfluoroundecanoic acid	562.4 > 518.5	12.830	13.146	-0.316	1.000	1076730	18.7	93.7	1203	
D 28 13C2 PFDaA	614.4 > 569.4	13.424	13.713	-0.289		3103590	43.6	87.3	3360	
29 Perfluorododecanoic acid	612.4 > 568.6	13.424	13.715	-0.291	1.000	1067827	21.4	107	1578	
30 Perfluorotridecanoic acid	662.4 > 618.5	13.914	14.193	-0.279	1.000	1092644	21.1	105	778	
D 33 13C2-PFTeDA	714.5 > 669.5	14.351	14.603	-0.252		2670650	42.1	84.3	3717	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.351	14.604	-0.253	1.000	446275	18.8	93.8	262	
D 35 13C2-PFHxDA	814.8 > 769.6	14.958	15.138	-0.180		1156686	39.0	78.0	2832	
34 Perfluorohexadecanoic acid	812.6 > 768.6	14.958	15.139	-0.181	1.000	1103795	16.3	81.7	1130	
36 Perfluorooctadecanoic acid	912.7 > 868.6	15.280	15.442	-0.162	1.000	264932	15.9	79.4	1068	

[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L4\_00017

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_041.d

Injection Date: 13-Feb-2016 00:01:23

Instrument ID: A4

Lims ID: CCV L4

Client ID:

Operator ID: JRB

ALS Bottle#: 47

Worklist Smp#: 40

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

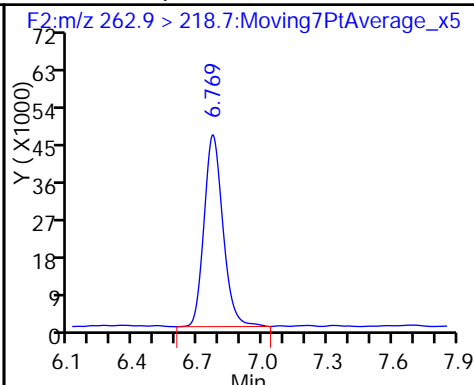
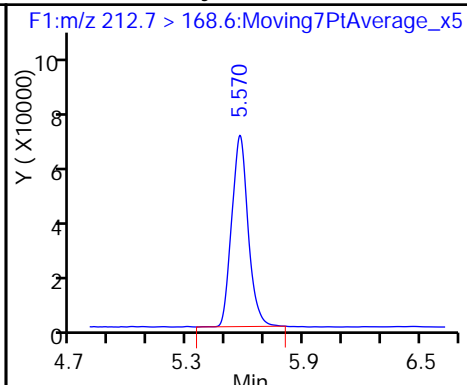
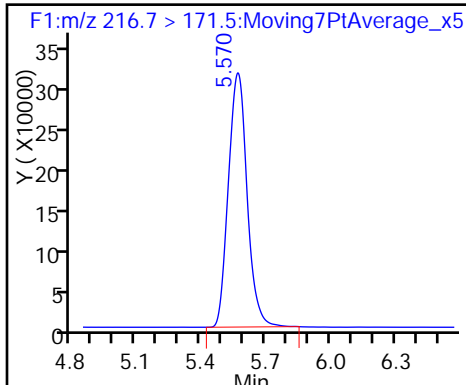
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

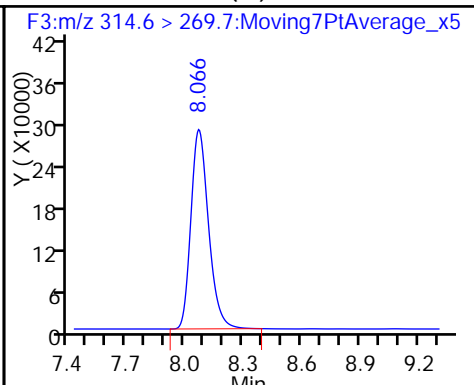
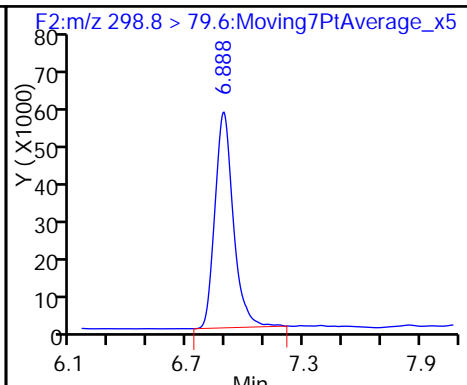
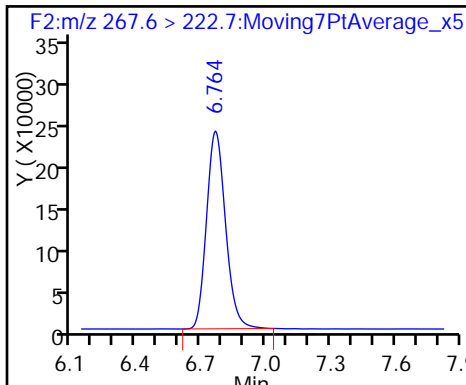
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

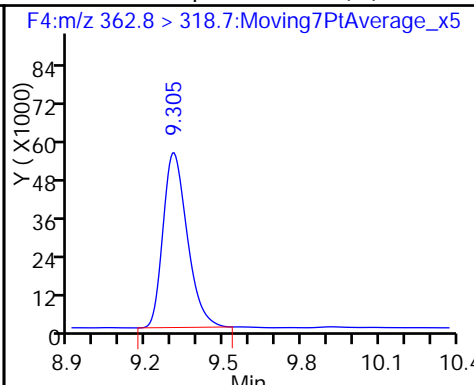
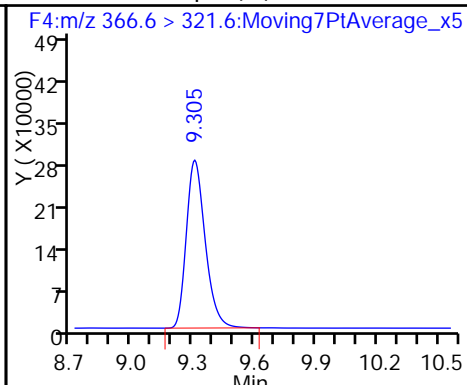
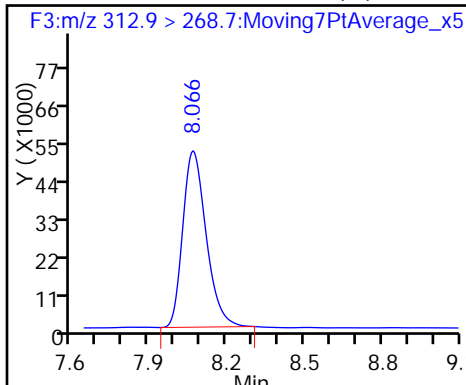
D 6 13C2 PFHxA (M)



7 Perfluorohexanoic acid (M)

D 8 13C4-PFHpA (M)

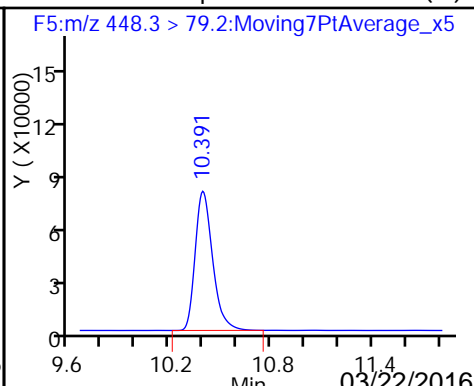
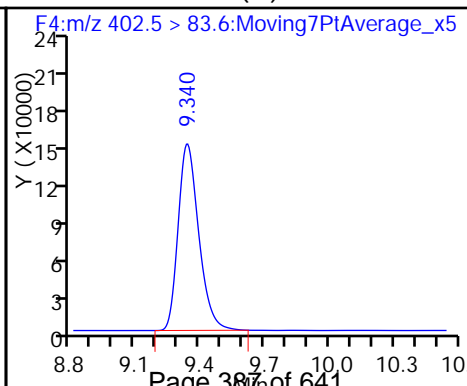
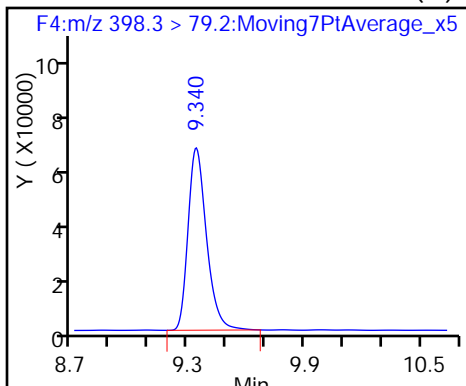
9 Perfluoroheptanoic acid (M)



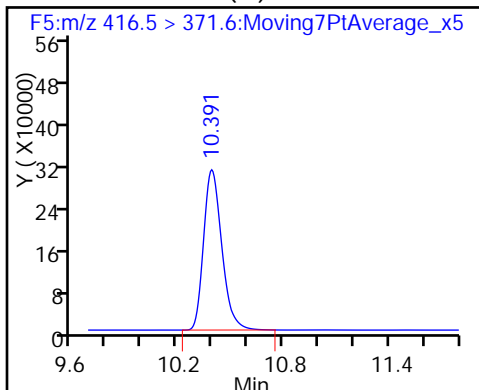
58 Perfluorohexanesulfonic acid (M)

D 11 18O2 PFHxS (M)

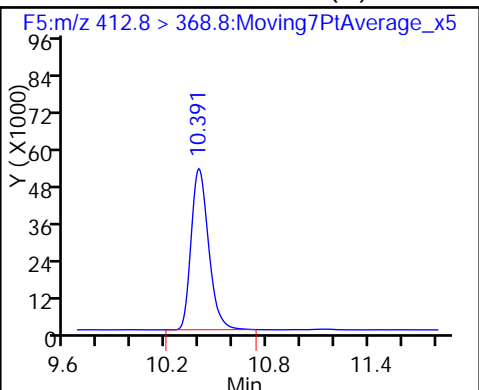
39 Perfluoroheptanesulfonic Acid (M)



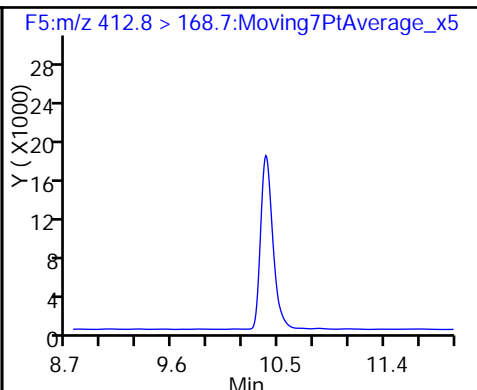
D 12 13C4 PFOA (M)



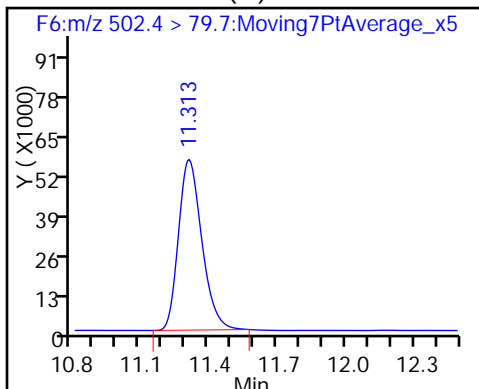
13 Perfluorooctanoic acid (M)



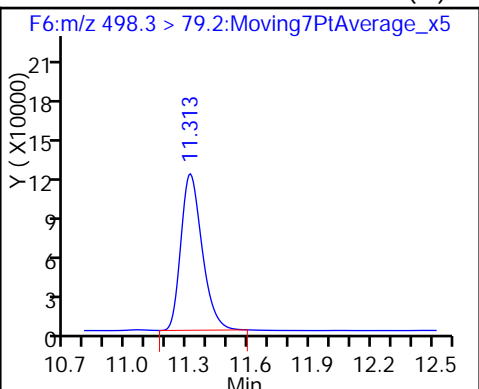
13 Perfluorooctanoic acid



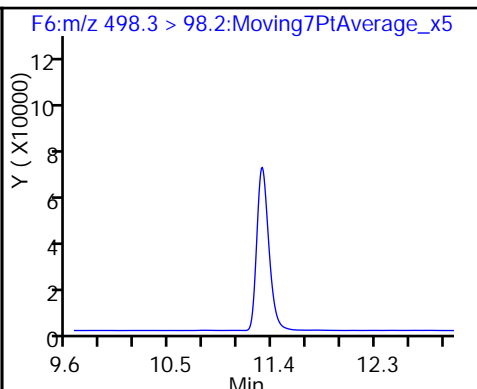
D 16 13C4 PFOS (M)



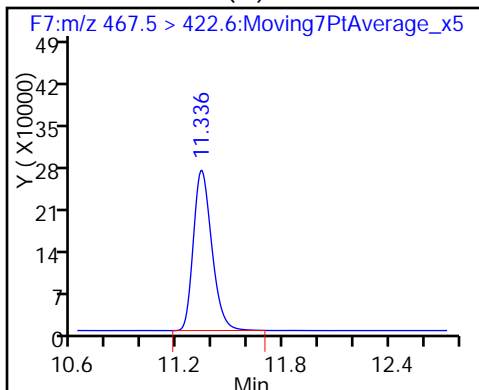
15 Perfluorooctane sulfonic acid (M)



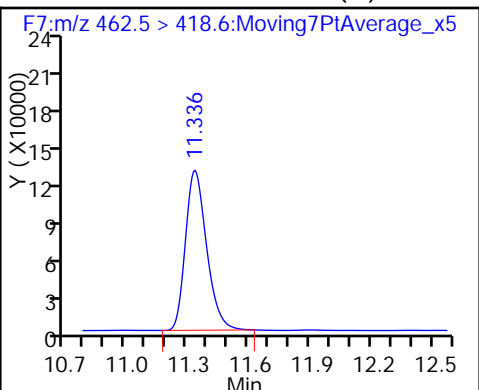
15 Perfluorooctane sulfonic acid



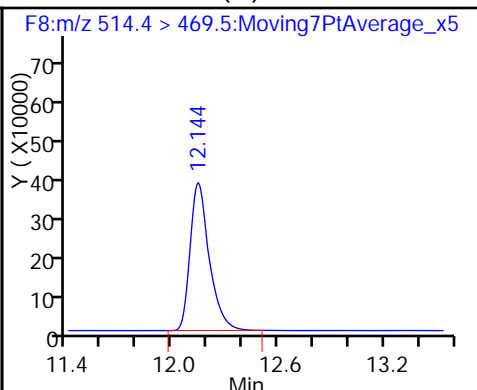
D 17 13C5 PFNA (M)



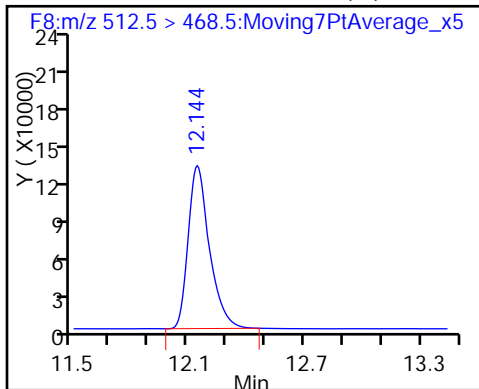
18 Perfluorononanoic acid (M)



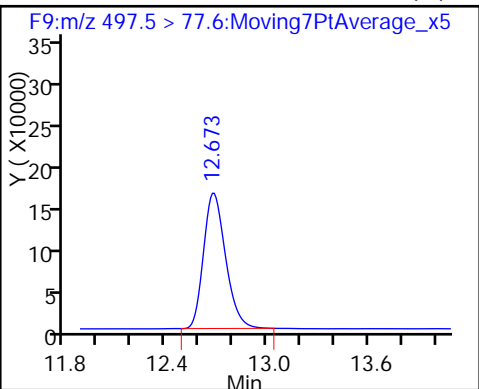
D 19 13C2 PFDA (M)



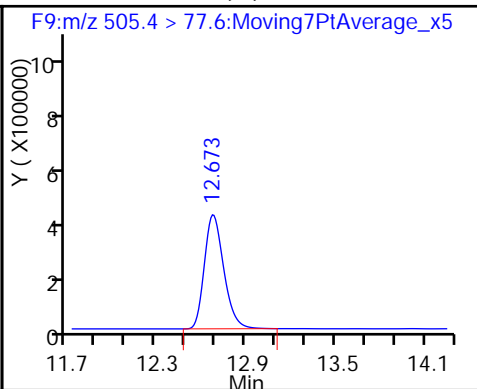
20 Perfluorodecanoic acid (M)



24 Perfluorooctane Sulfonamide (M)

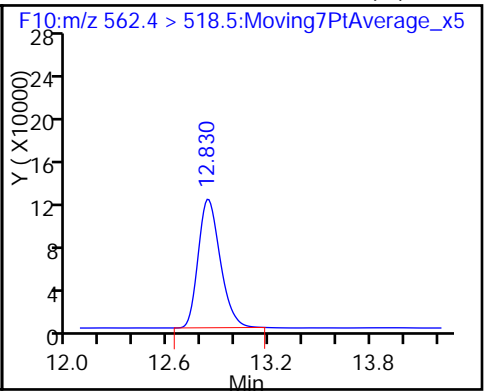
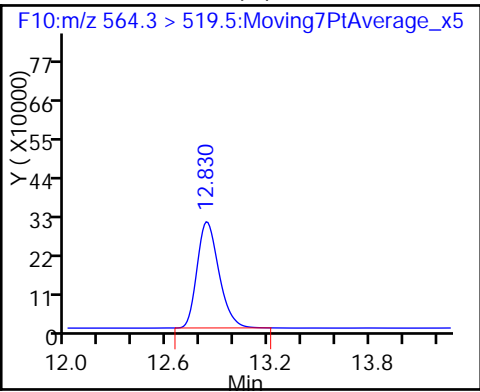
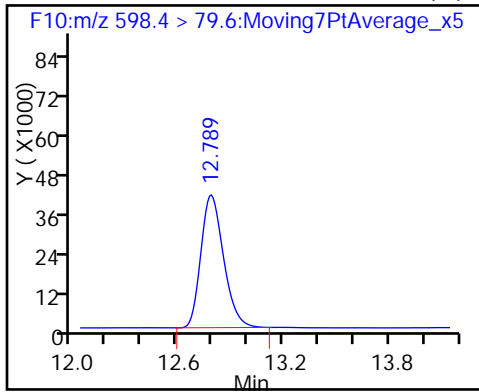


D 23 13C8 FOSA (M)



49 Perfluorodecane Sulfonic acid (M) D 26 13C2 PFUnA (M)

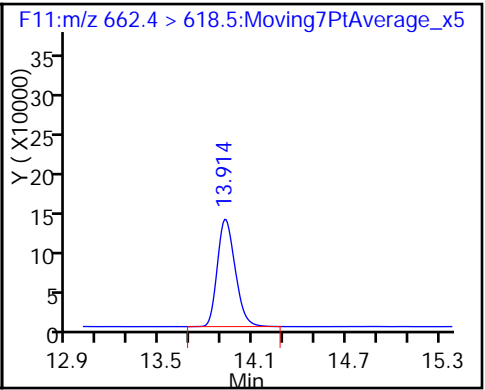
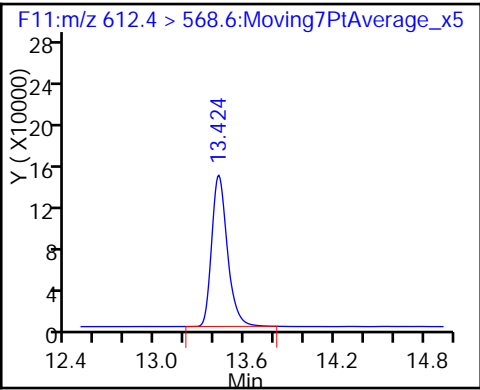
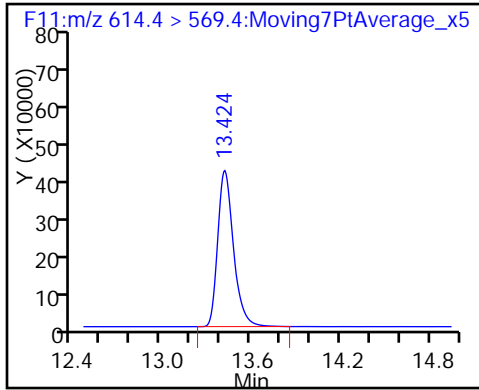
27 Perfluoroundecanoic acid (M)



D 28 13C2 PFDaA (M)

29 Perfluorododecanoic acid (M)

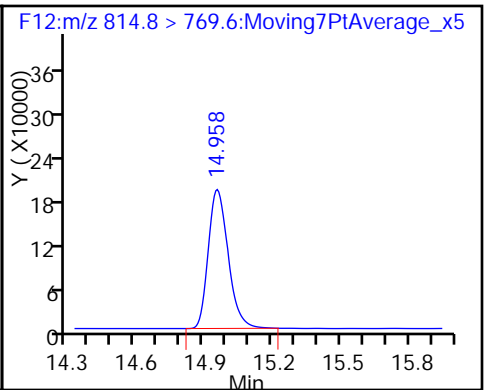
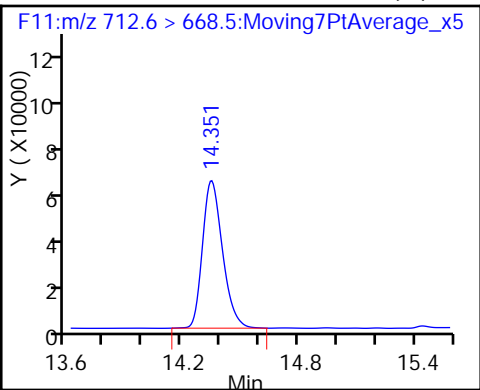
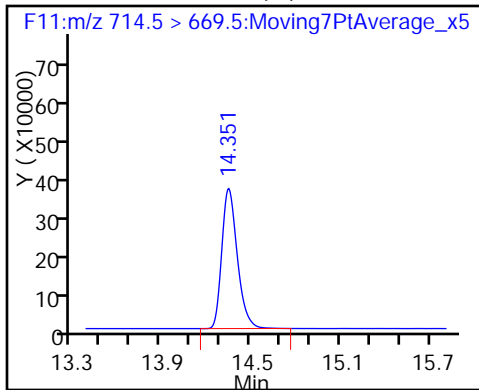
30 Perfluorotridecanoic acid (M)



D 33 13C2-PFTeDA (M)

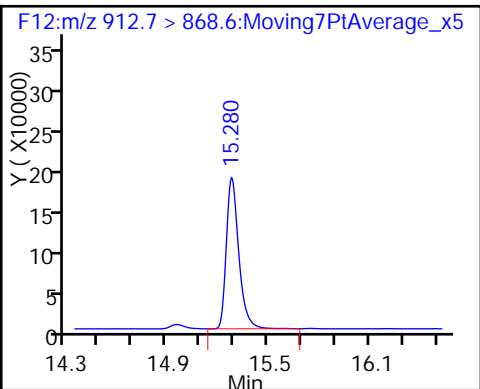
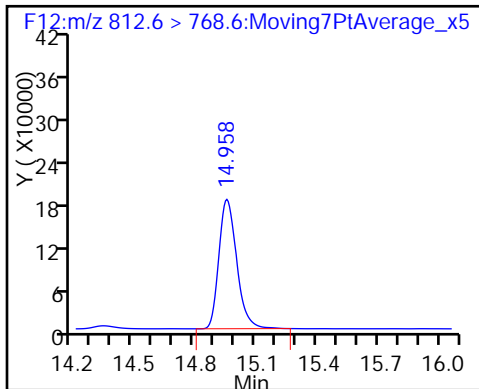
32 Perfluorotetradecanoic acid (M)

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: CCV 320-100305/51 Calibration Date: 02/13/2016 03:54  
 Instrument ID: A4 Calib Start Date: 02/12/2016 10:35  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/12/2016 12:42  
 Lab File ID: 12FEB2016A4A\_052.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	0.4901	0.5244		53.5	50.0	7.0	25.0
Perfluoropentanoic acid (PFPeA)	AveID	0.4708	0.4516		48.0	50.0	-4.1	25.0
Perfluorobutanesulfonic acid (PFBS)	L1ID		0.7015		35.2	44.2	-20.5	25.0
Perfluorohexanoic acid (PFHxA)	AveID	0.4308	0.3874		45.0	50.0	-10.1	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	0.5120	0.4541		44.3	50.0	-11.3	25.0
Perfluorohexanesulfonic acid (PFHxS)	L2ID		1.221		51.6	47.3	9.2	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	AveID	2.544	4.219		78.9	47.6	65.8*	25.0
Perfluorooctanoic acid (PFOA)	AveID	0.4910	0.4097		41.7	50.0	-16.5	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	4.265	7.089		79.4	47.8	66.2*	25.0
Perfluorononanoic acid (PFNA)	L1ID		0.9842		47.0	50.0	-6.1	25.0
Perfluorodecanoic acid (PFDA)	AveID	0.9142	0.8496		46.5	50.0	-7.1	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9450	0.9919		52.5	50.0	5.0	25.0
Perfluorodecane Sulfonic acid	AveID	1.927	2.658		66.5	48.2	37.9*	25.0
Perfluoroundecanoic acid (PFUnA)	AveID	1.037	1.005		48.5	50.0	-3.1	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.8053	0.8257		51.3	50.0	2.5	25.0
Perfluorotridecanoic Acid (PFTriA)	L2ID		0.7837		47.0	50.0	-6.1	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.3438		45.2	50.0	-9.6	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L1ID		0.8249		38.1	50.0	-23.8	25.0
Perfluoro-n-octadecanoic acid (PFODA)	L1ID		0.6924		34.7	50.0	-30.6*	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_052.d  
 Lims ID: CCV L5  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 13-Feb-2016 03:54:21 ALS Bottle#: 48 Worklist Smp#: 51  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L5  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub12  
 Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 16-Feb-2016 08:37:08 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d  
 Column 1 : Det: F1:MRM  
 Process Host: XAWRK014

First Level Reviewer: westendorfc Date: 15-Feb-2016 13:26:20

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	216.7 > 171.5	5.496	5.712	-0.216	1933902	47.7		95.3	6988	
2 Perfluorobutyric acid	212.7 > 168.6	5.502	5.712	-0.210	1014168	53.5		107	2752	
4 Perfluoropentanoic acid	262.9 > 218.7	6.668	6.986	-0.318	662247	48.0		95.9	686	
D 3 13C5-PFPeA	267.6 > 222.7	6.663	6.986	-0.323	1466604	45.8		91.7	3696	
5 Perfluorobutane Sulfonate	298.8 > 79.6	6.783	7.115	-0.332	547214	NC			354	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	6.783	7.115	-0.332	547214	35.2		79.5		
D 6 13C2 PFHxA	314.6 > 269.7	7.954	8.367	-0.413	1862377	47.6		95.1	5654	
7 Perfluorohexanoic acid	312.9 > 268.7	7.954	8.369	-0.415	721529	45.0		89.9	1439	
D 8 13C4-PFHpA	366.6 > 321.6	9.194	9.639	-0.445	1616208	41.5		83.0	4113	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.194	9.640	-0.446	733894	44.3		88.7	1564	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.229	9.675	-0.446	1019442	51.6		109		
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.229	9.675	-0.446	1019442	NC			1937	
D 11 18O2 PFHxS	402.5 > 83.6	9.229	9.675	-0.446	834789	36.8		77.8	2788	
39 Perfluoroheptanesulfonic Acid	448.3 > 79.2	10.308	10.740	-0.432	1225149	78.9		166		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
14 Perfluoroheptane Sulfonate	448.3 > 79.2	10.308	10.740	-0.432	1.000	1235149	NC		2954	
D 12 13C4 PFOA	416.5 > 371.6	10.308	10.743	-0.435		1917508	45.0	90.1	5722	
13 Perfluorooctanoic acid	412.8 > 368.8	10.308	10.743	-0.435	1.000	785613	41.7	83.5	1679	
D 16 13C4 PFOS	502.4 > 79.7	11.240	11.651	-0.411		294011	29.0	60.7	1240	
15 Perfluorooctane sulfonic acid	498.3 > 79.2	11.248	11.656	-0.408	1.000	2084119	79.4	166	2968	
D 17 13C5 PFNA	467.5 > 422.6	11.270	11.678	-0.408		1876713	45.1	90.2	4313	
18 Perfluorononanoic acid	462.5 > 418.6	11.270	11.680	-0.410	1.000	1846967	47.0	93.9	2443	
D 19 13C2 PFDA	514.4 > 469.5	12.091	12.473	-0.382		2662972	45.3	90.7	4929	
20 Perfluorodecanoic acid	512.5 > 468.5	12.099	12.476	-0.377	1.000	2262503	46.5	92.9	2657	
24 Perfluorooctane Sulfonamide	497.5 > 77.6	12.633	12.991	-0.358	1.000	3391522	52.5	105	1508	
D 23 13C8 FOSA	505.4 > 77.6	12.633	12.991	-0.358		3419233	38.3	76.6	1536	
25 Perfluorodecane Sulfonate	598.4 > 79.6	12.747	13.098	-0.351	1.000	787912	NC		2055	
49 Perfluorodecane Sulfonic acid	598.4 > 79.6	12.747	13.098	-0.351	1.000	787912	66.5	138		
D 26 13C2 PFUnA	564.3 > 519.5	12.799	13.146	-0.347		2548279	43.6	87.1	4647	
27 Perfluoroundecanoic acid	562.4 > 518.5	12.799	13.146	-0.347	1.000	2562150	48.5	96.9	2158	
D 28 13C2 PFDoA	614.4 > 569.4	13.393	13.713	-0.320		3091706	43.5	86.9	4335	
29 Perfluorododecanoic acid	612.4 > 568.6	13.393	13.715	-0.322	1.000	2552807	51.3	103	3289	
30 Perfluorotridecanoic acid	662.4 > 618.5	13.896	14.193	-0.297	1.000	2423005	47.0	93.9	1489	
D 33 13C2-PFTeDA	714.5 > 669.5	14.327	14.603	-0.276		2525070	39.8	79.7	3437	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.327	14.604	-0.277	1.000	1062771	45.2	90.4	513	
D 35 13C2-PFHxDA	814.8 > 769.6	14.944	15.138	-0.194		1032868	34.8	69.6	2340	
34 Perfluorohexadecanoic acid	812.6 > 768.6	14.944	15.139	-0.195	1.000	2550287	38.1	76.2	1971	
36 Perfluorooctandecanoic acid	912.7 > 868.6	15.265	15.442	-0.177	1.000	2140741	34.7	69.4	2051	



[QC Flag Legend](#)

Processing Flags

NC - Not Calibrated

[Reagents:](#)

LCPFC-L5\_00016

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_052.d

Injection Date: 13-Feb-2016 03:54:21

Instrument ID: A4

Lims ID: CCV L5

Client ID:

Operator ID: JRB

ALS Bottle#: 48

Worklist Smp#: 51

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

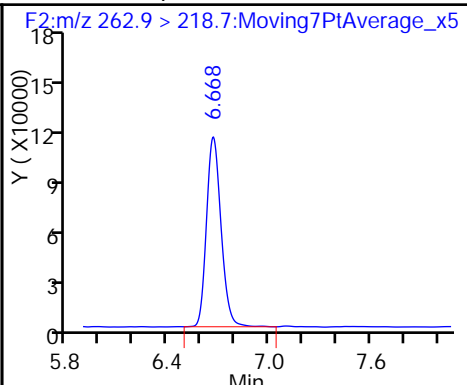
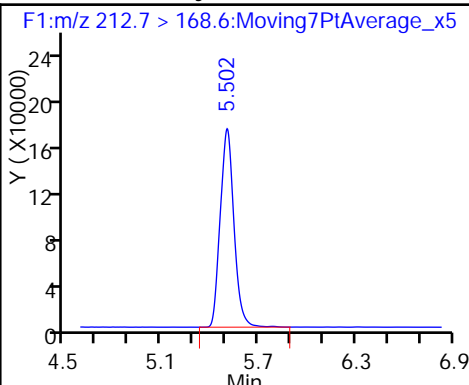
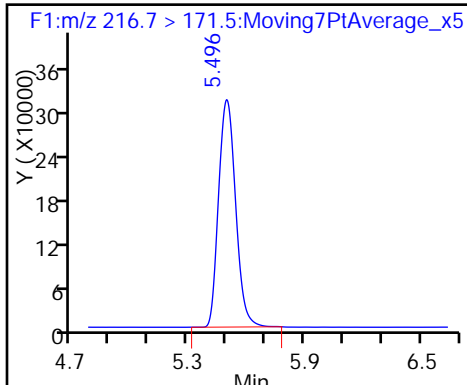
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Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

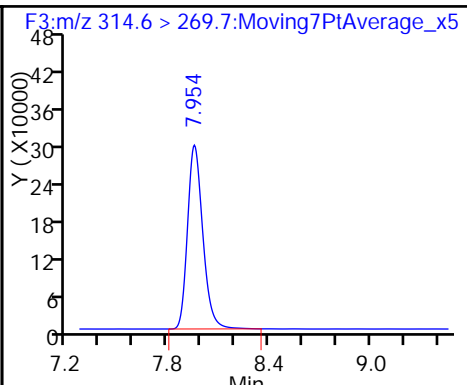
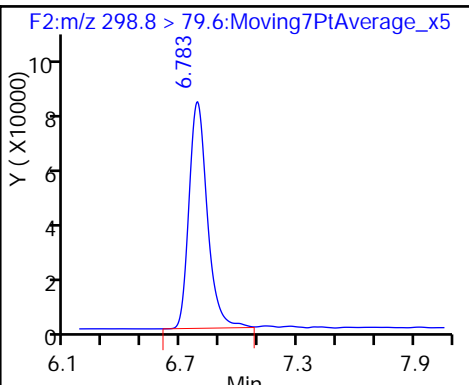
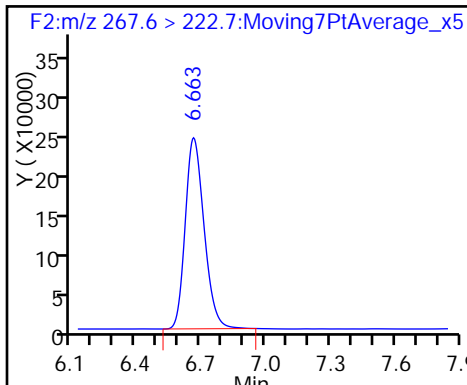
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

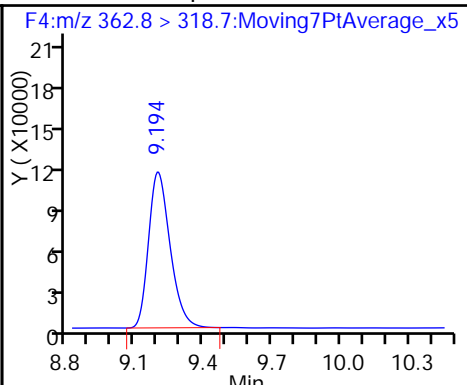
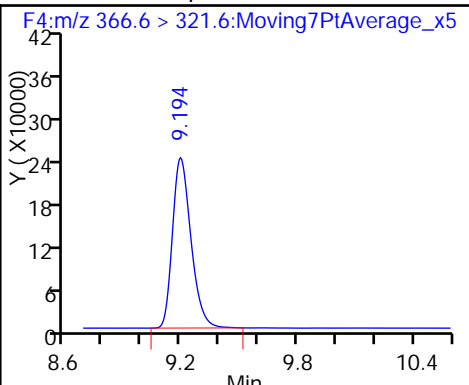
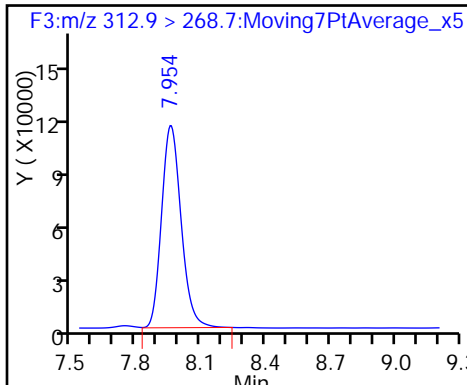
D 6 13C2 PFHxA



7 Perfluorohexanoic acid

D 8 13C4-PFHpA

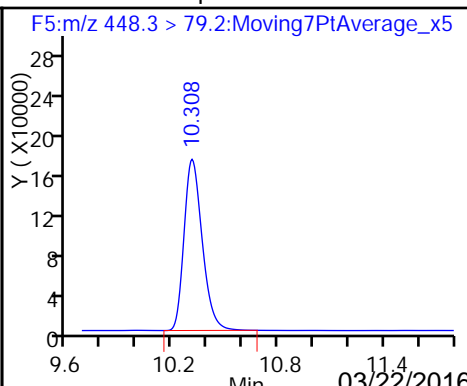
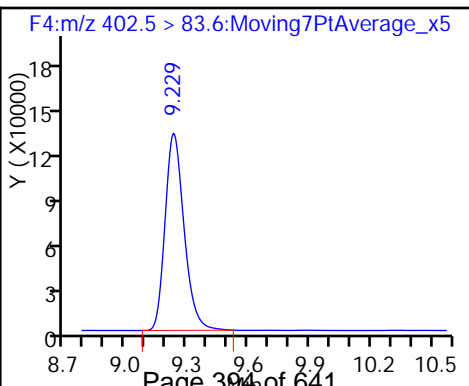
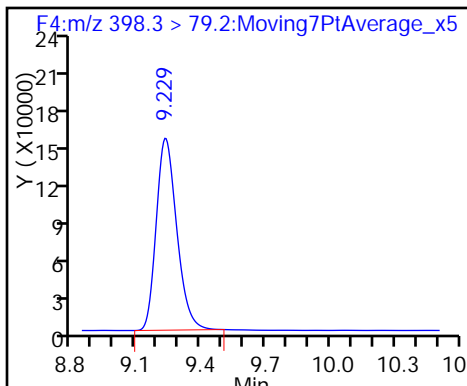
9 Perfluoroheptanoic acid



58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

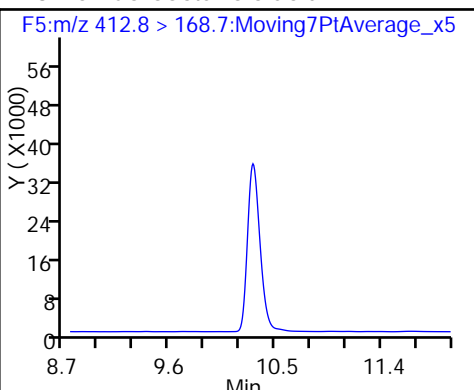
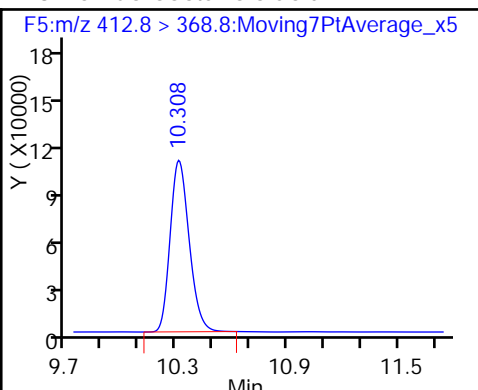
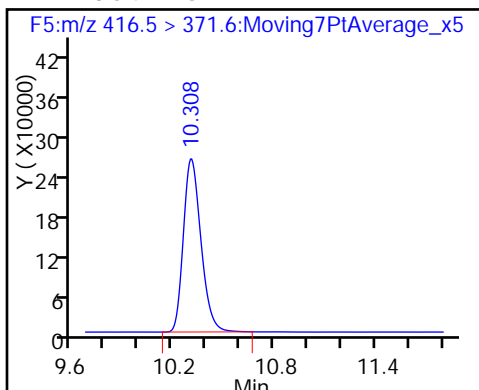
39 Perfluoroheptanesulfonic Acid



D 12 13C4 PFOA

13 Perfluorooctanoic acid

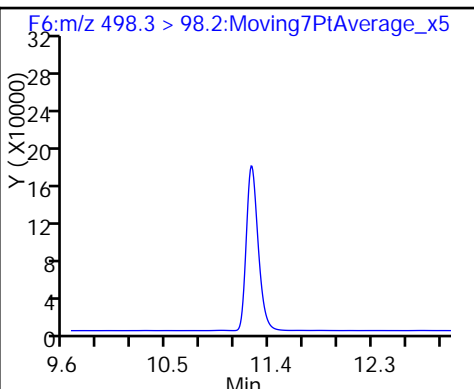
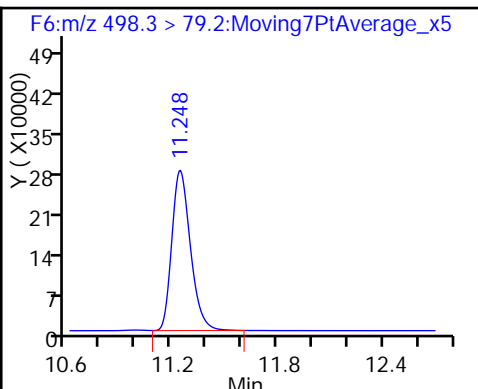
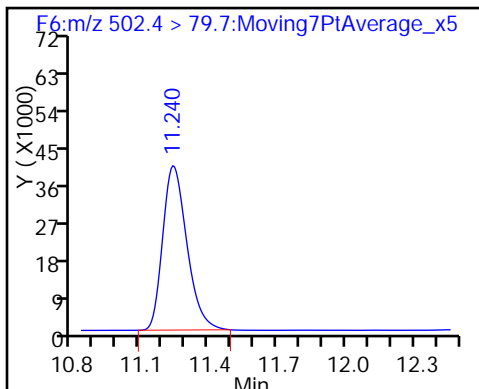
13 Perfluorooctanoic acid



D 16 13C4 PFOS

15 Perfluorooctane sulfonic acid

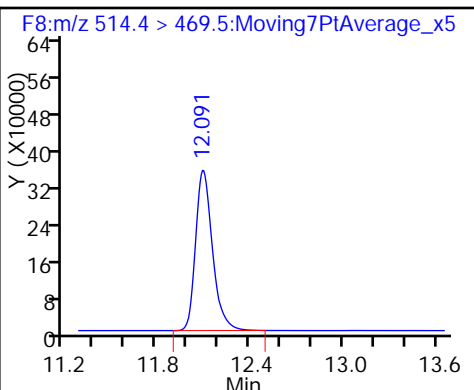
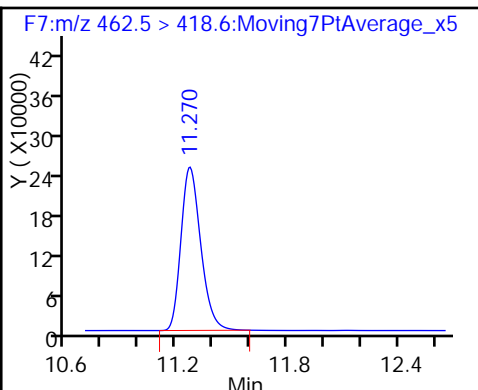
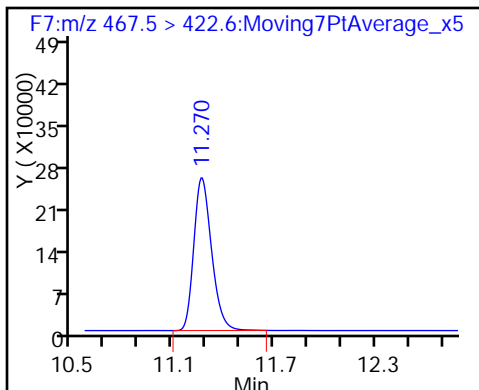
15 Perfluorooctane sulfonic acid



D 17 13C5 PFNA

18 Perfluorononanoic acid

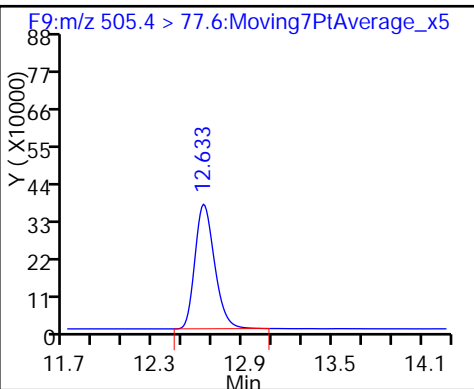
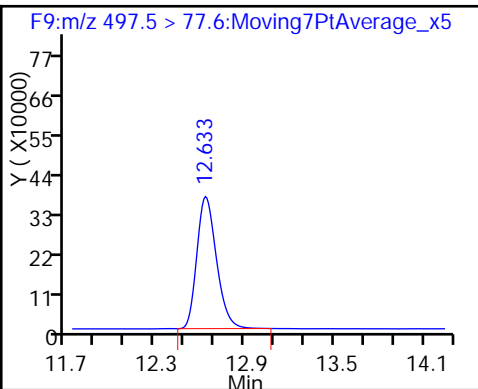
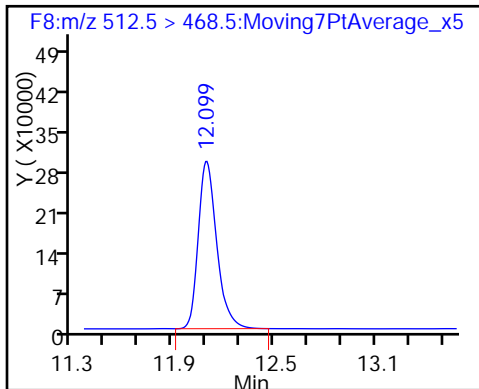
D 19 13C2 PFDA



20 Perfluorodecanoic acid

24 Perfluorooctane Sulfonamide

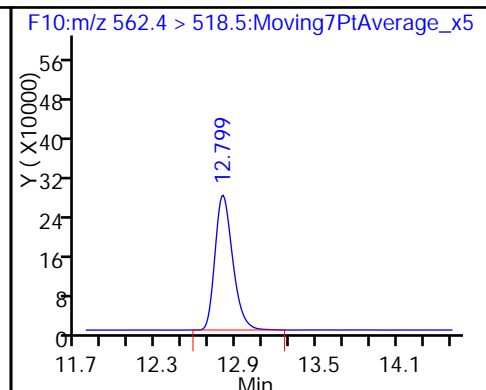
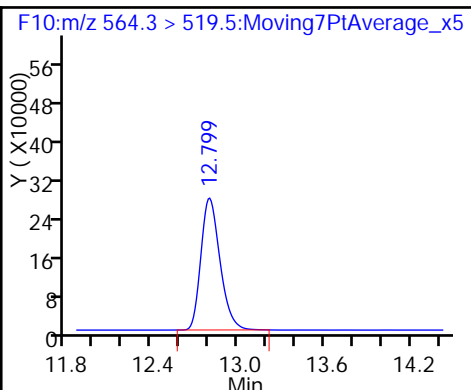
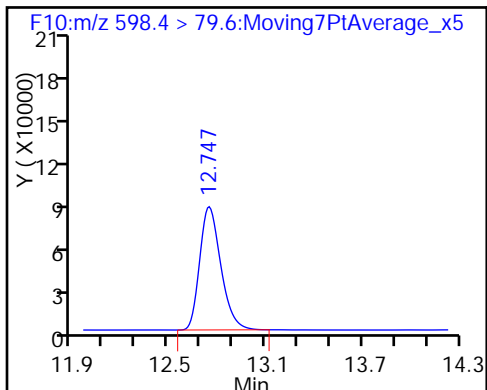
D 23 13C8 FOSA



49 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

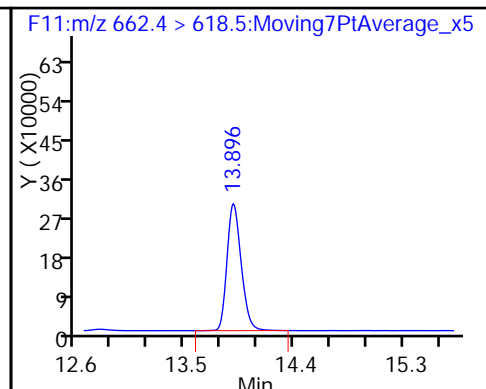
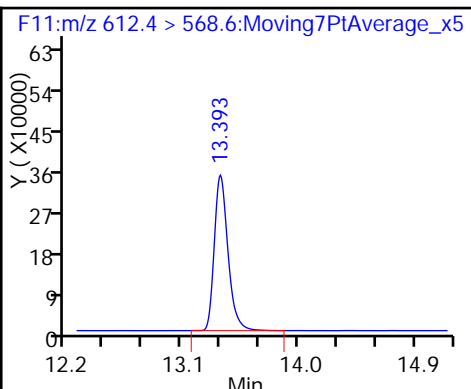
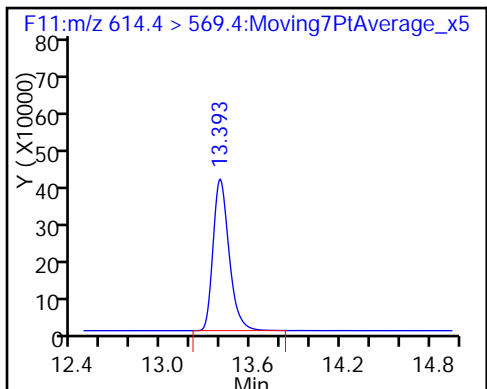
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

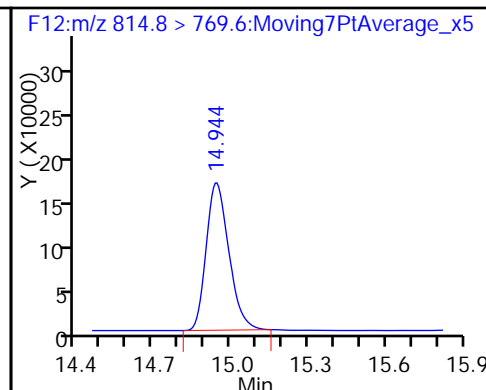
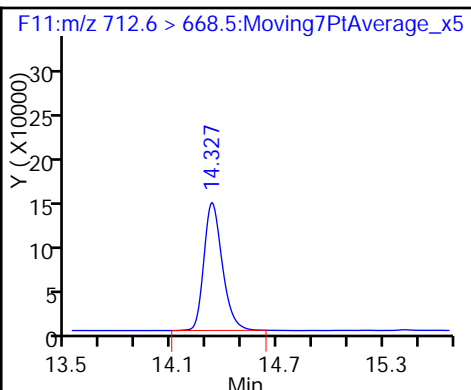
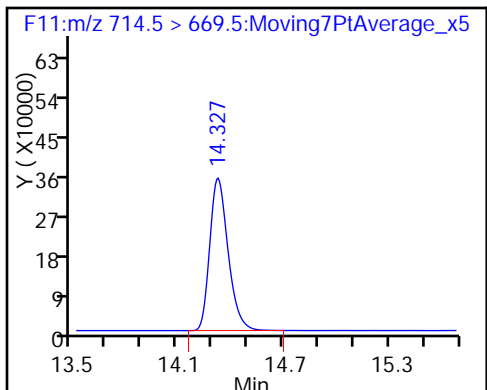
30 Perfluorotridecanoic acid



D 33 13C2-PFTeDA

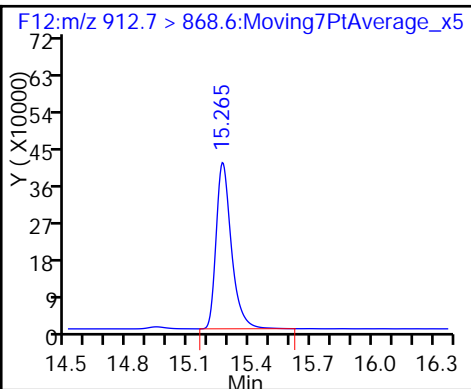
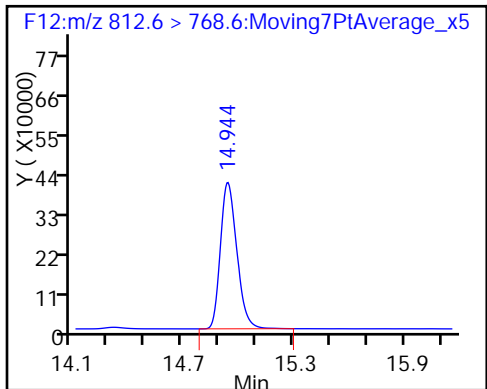
32 Perfluorotetradecanoic acid

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: ICV 320-100906/10 Calibration Date: 02/19/2016 17:58  
 Instrument ID: A6 Calib Start Date: 02/19/2016 15:08  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/19/2016 17:15  
 Lab File ID: 19FEB2016A6A\_012.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	1.304	1.537		58.9	50.0	17.9	25.0
Perfluoropentanoic acid (PFPeA)	AveID	1.043	1.090		52.2	50.0	4.5	25.0
Perfluorobutanesulfonic acid (PFBS)	AveID	0.9930	1.215		54.1	44.3	22.4	25.0
Perfluorohexanoic acid (PFHxA)	AveID	1.083	1.137		52.5	50.0	4.9	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.043	1.171		56.1	50.0	12.3	25.0
Perfluorohexanesulfonic acid (PFHxS)	AveID	0.6753	0.6993		48.9	47.3	3.6	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	L2ID		0.6445		49.3	47.6	3.6	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.004	1.026		51.1	50.0	2.2	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	0.9724	1.110		54.5	47.8	14.1	25.0
Perfluorononanoic acid (PFNA)	AveID	0.8075	0.8793		54.4	50.0	8.9	25.0
Perfluorodecanoic acid (PFDA)	L2ID		0.997		50.8	50.0	1.6	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9268	0.9404		50.7	50.0	1.5	25.0
Perfluorodecane Sulfonic acid	L1ID		0.6116		53.2	48.3	10.3	25.0
Perfluoroundecanoic acid (PFUnA)	L1ID		0.9581		57.5	50.0	15.0	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.7311	0.8388		57.4	50.0	14.7	25.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.9409	1.009		53.6	50.0	7.2	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.6279		52.1	50.0	4.3	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L2ID		1.181		52.5	50.0	5.1	25.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.111	1.148		51.7	50.0	3.4	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_012.d  
 Lims ID: ICV  
 Client ID:  
 Sample Type: ICV  
 Inject. Date: 19-Feb-2016 17:58:06 ALS Bottle#: 16 Worklist Smp#: 10  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: ICV  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A4\*sub6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:59:00 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.705	5.712	-0.007	693136	43.3		86.6	2632	
2 Perfluorobutyric acid	212.9 > 169.0	5.705	5.715	-0.010	1065572	58.9			3120	
D 3 13C5-PFPeA	267.9 > 223.0	6.816	6.827	-0.011	1409700	45.8		91.7	4713	
4 Perfluoropentanoic acid	262.9 > 219.0	6.816	6.828	-0.012	1536827	52.2			395	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.931	6.939	-0.008	654025	NC			569	
	298.9 > 99.0	6.931	6.939	-0.008	370497		1.77(0.00-0.00)		770	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.931	6.939	-0.008	654025	54.1				
7 Perfluorohexanoic acid	313.0 > 269.0	8.066	8.078	-0.012	1540103	52.5			1860	
D 6 13C2 PFHxA	315.0 > 270.0	8.066	8.080	-0.014	1354941	47.8		95.7	4025	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.142	8.158	-0.016	456683	NC			1959	
D 8 13C4-PFHpA	367.0 > 322.0	9.305	9.322	-0.017	1318603	42.8		85.6	4607	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.305	9.322	-0.017	1543763	56.1			4041	
D 11 18O2 PFHxS	403.0 > 84.0	9.341	9.358	-0.017	575394	38.4		81.3	2432	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.341	9.364	-0.023	401937	NC			1168	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.341	9.364	-0.023	401937	48.9				

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.433	10.450	-0.017		1436383	42.1		84.1	5380	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.433	10.450	-0.017	1.000	1474139	51.1			1374	
413.0 > 169.0	10.433	10.450	-0.017	1.000	511709		2.88(0.00-0.00)		1005	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.433	10.457	-0.024	1.000	444139	NC			1618	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.433	10.457	-0.024	1.000	444139	49.3				
D 16 13C4 PFOS										
503.0 > 80.0	11.393	11.414	-0.021		691986	38.1		79.6	3094	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.393	11.417	-0.024	1.000	767227	54.5			354	
499.0 > 99.0	11.393	11.417	-0.024	1.000	394536		1.94(0.00-0.00)		1323	
D 17 13C5 PFNA										
468.0 > 423.0	11.415	11.434	-0.019		1208840	42.9		85.8	3370	
18 Perfluorononanoic acid										
463.0 > 419.0	11.415	11.439	-0.024	1.000	1062919	54.4			3617	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.224	12.249	-0.025	1.000	380394	NC			1433	
D 19 13C2 PFDA										
515.0 > 470.0	12.256	12.275	-0.019		1189272	41.5		83.1	5143	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.256	12.276	-0.020	1.000	1186169	50.8			4224	
D 23 13C8 FOSA										
506.0 > 78.0	12.794	12.810	-0.016		1917001	46.2		92.4	1321	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.794	12.812	-0.018	1.000	1802677	50.7			1040	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.930	12.949	-0.019	1.000	427181	NC			1632	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.930	12.949	-0.019	1.000	427181	53.2				
D 26 13C2 PFUnA										
565.0 > 520.0	12.982	13.000	-0.018		1413914	37.8		75.7	4613	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.982	13.000	-0.018	1.000	1354687	57.5			3853	
D 28 13C2 PFDaA										
615.0 > 570.0	13.589	13.609	-0.020		1717701	43.9		87.8	4484	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.589	13.611	-0.022	1.000	1440856	57.4			2048	
31 PFDoS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.056	14.083	-0.027	1.000	481182	NC			1600	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.110	14.130	-0.020	1.000	1733242	53.6			2403	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.553	14.571	-0.018	1.000	1078544	52.1			883	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.546	14.571	-0.025		1649978	46.0		91.9	6367	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA										
815.0 > 770.0	15.191	15.207	-0.016		2024385	47.0		93.9	5896	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.191	15.208	-0.017	1.000	2029338	52.5			2611	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.527	15.543	-0.016	1.000	1972635	51.7			2318	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFCIC\_00016

Amount Added: 1.00

Units: mL



Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_012.d

Injection Date: 19-Feb-2016 17:58:06

Instrument ID: A6

Lims ID: ICV

Client ID:

Operator ID: JRB

ALS Bottle#: 16

Worklist Smp#: 10

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

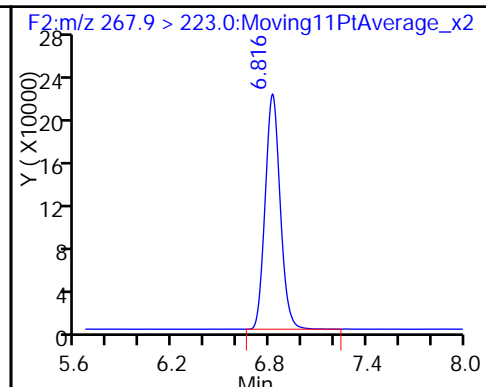
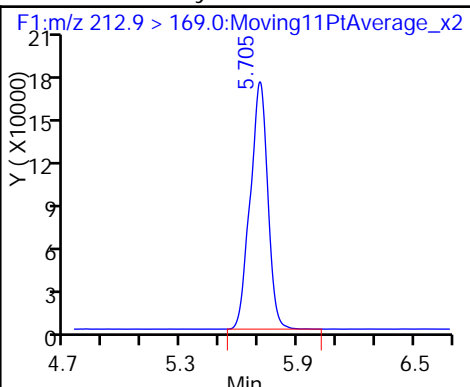
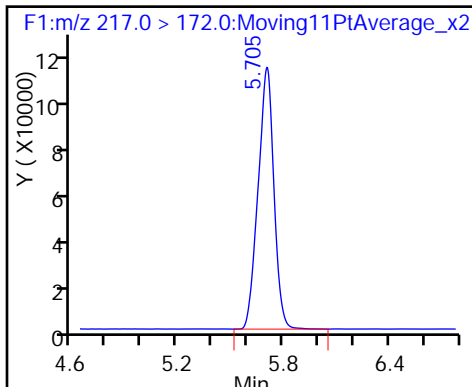
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

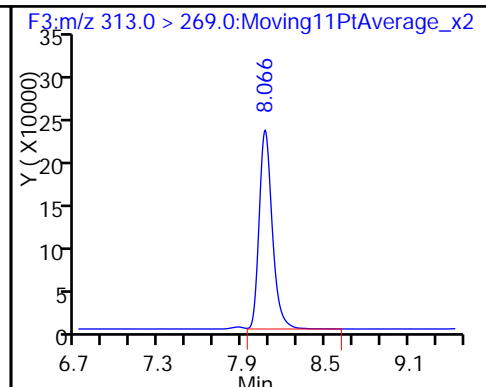
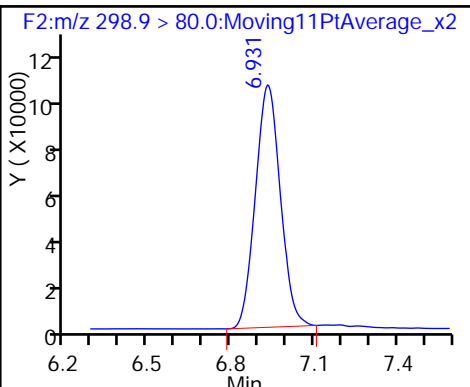
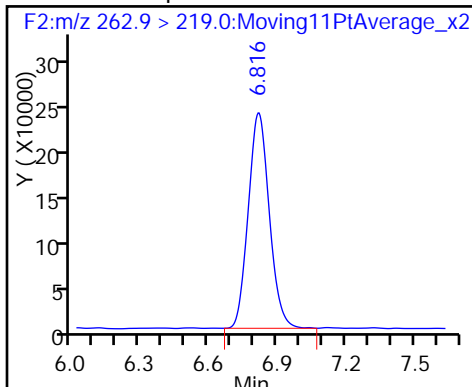
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

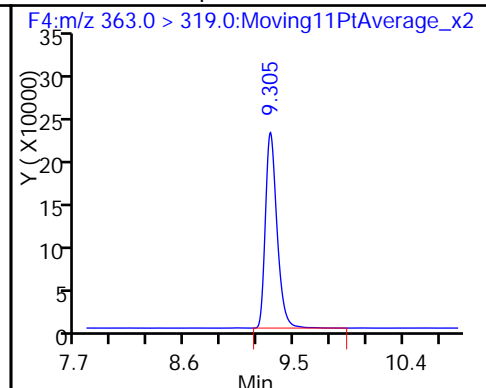
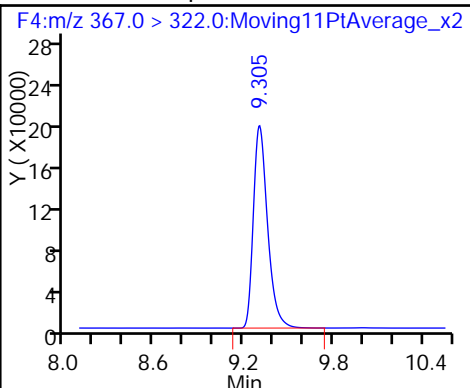
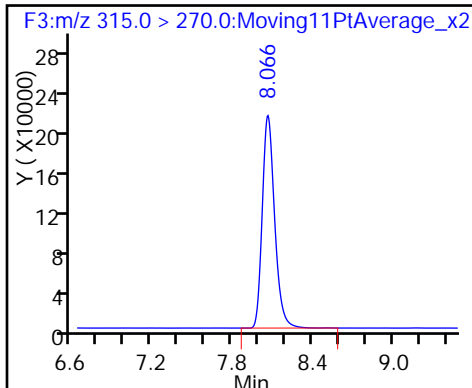
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

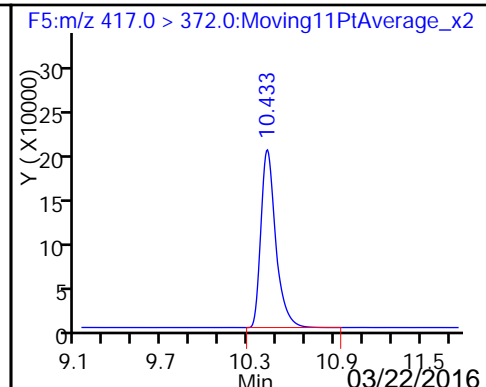
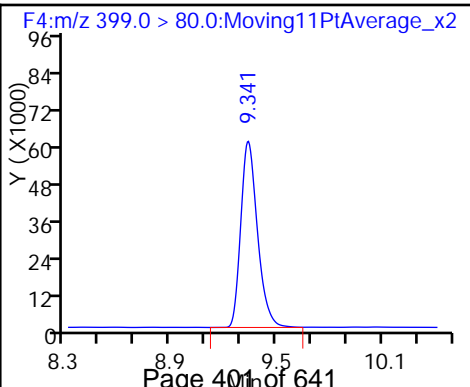
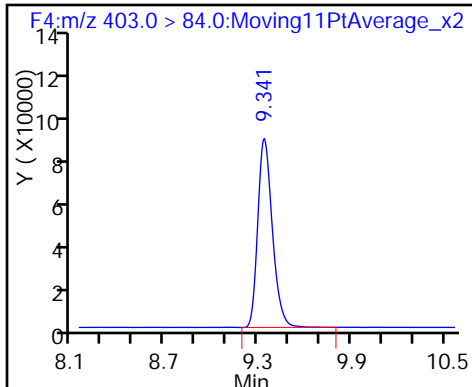
9 Perfluoroheptanoic acid

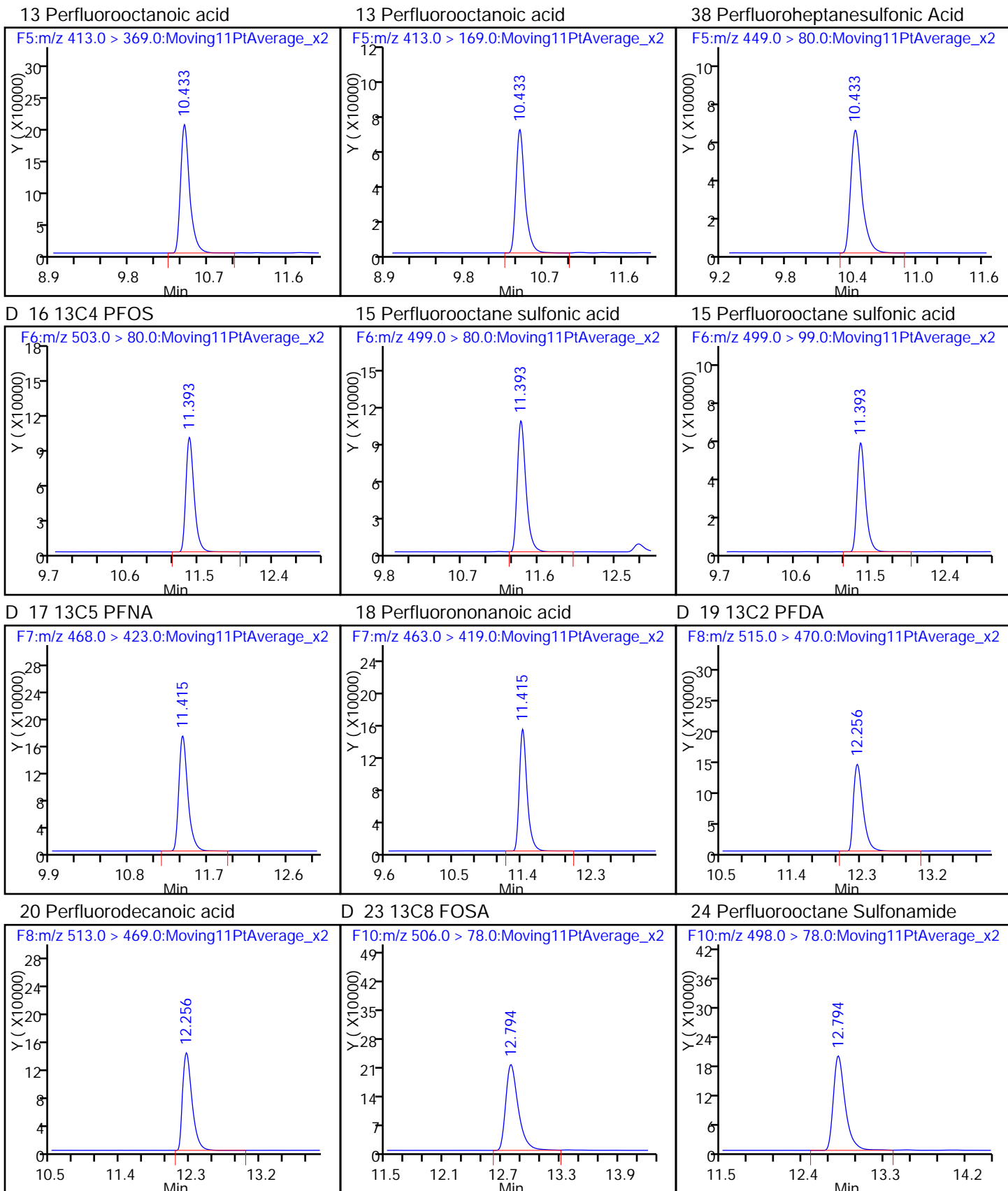


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

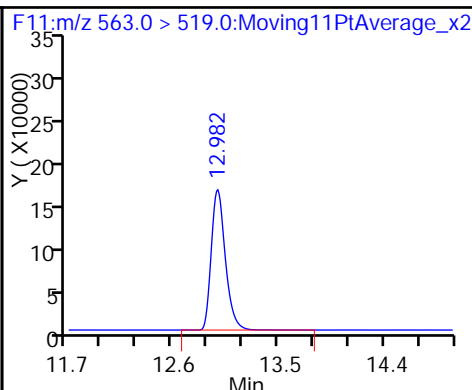
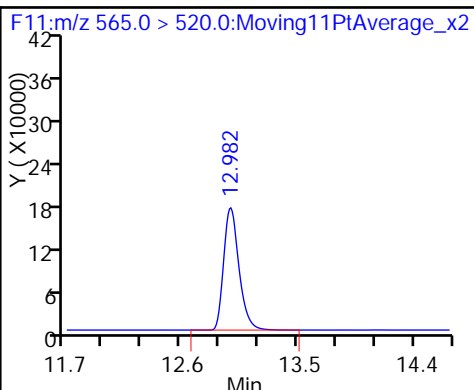
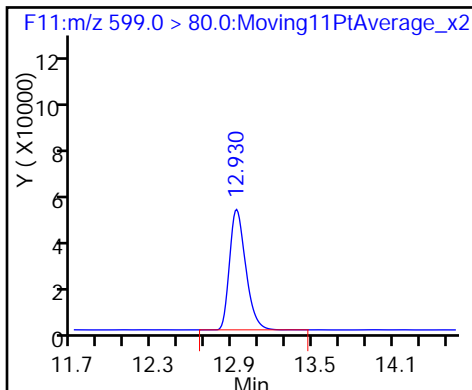




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

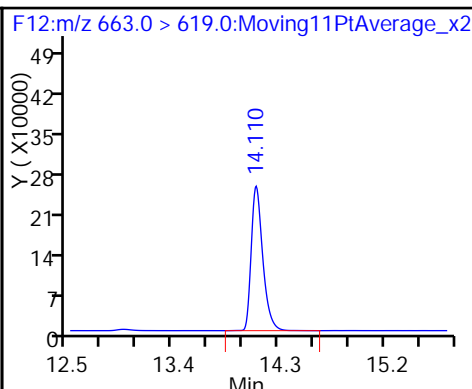
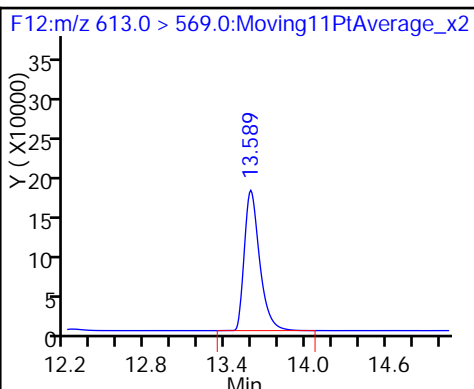
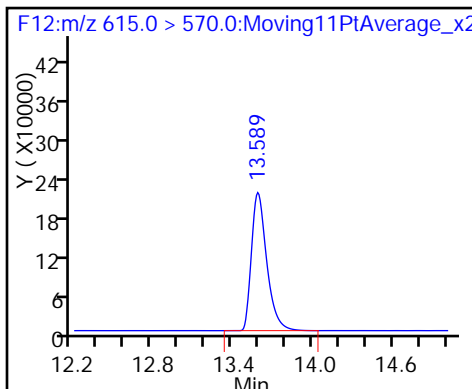
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

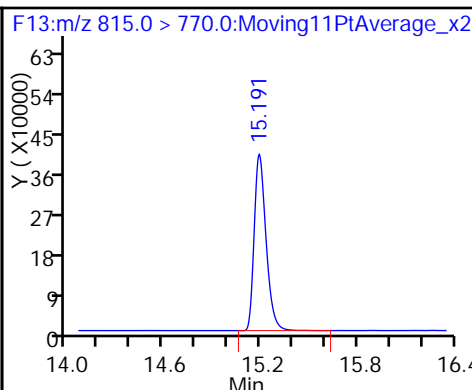
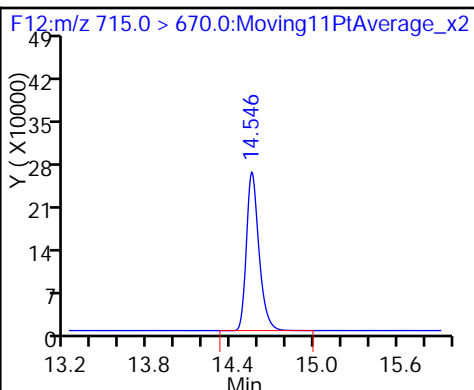
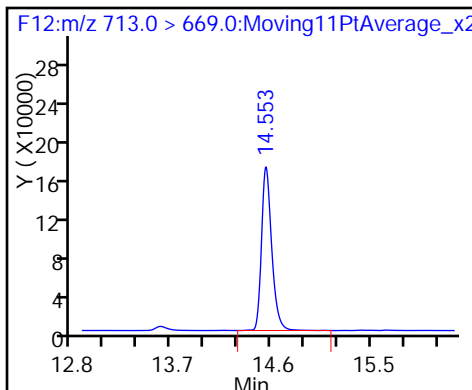
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

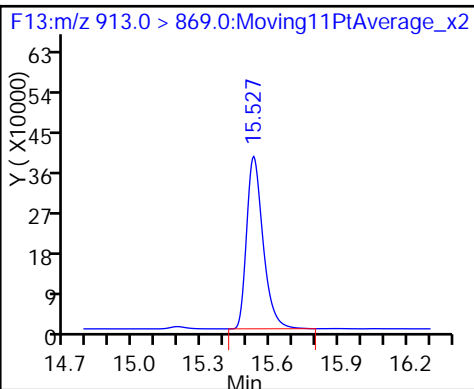
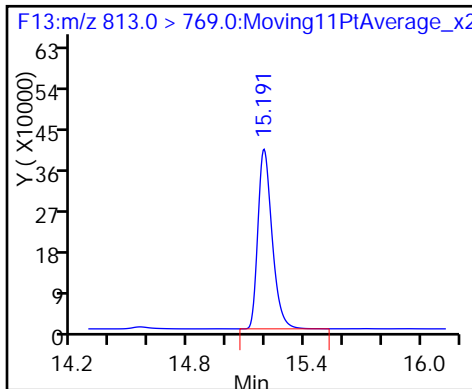
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: CCV 320-100906/43 Calibration Date: 02/20/2016 05:40  
 Instrument ID: A6 Calib Start Date: 02/19/2016 15:08  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/19/2016 17:15  
 Lab File ID: 19FEB2016A6A\_045.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	1.304	1.434		22.0	20.0	10.0	25.0
Perfluoropentanoic acid (PFPeA)	AveID	1.043	1.028		19.7	20.0	-1.5	25.0
Perfluorobutanesulfonic acid (PFBS)	AveID	0.9930	1.189		21.2	17.7	19.8	25.0
Perfluorohexanoic acid (PFHxA)	AveID	1.083	1.126		20.8	20.0	3.9	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.043	1.190		22.8	20.0	14.1	25.0
Perfluorohexanesulfonic acid (PFHxS)	AveID	0.6753	0.7464		20.9	18.9	10.5	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.004	1.028		20.5	20.0	2.4	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	L2ID		0.7230		22.3	19.0	17.0	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	0.9724	1.073		21.1	19.1	10.4	25.0
Perfluorononanoic acid (PFNA)	AveID	0.8075	0.7921		19.6	20.0	-1.9	25.0
Perfluorodecanoic acid (PFDA)	L2ID		0.9104		18.3	20.0	-8.4	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9268	0.8822		19.0	20.0	-4.8	25.0
Perfluorodecane Sulfonic acid	L1ID		0.6729		23.6	19.3	22.2	25.0
Perfluoroundecanoic acid (PFUnA)	L1ID		0.8914		21.0	20.0	4.9	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.7311	0.8023		21.9	20.0	9.7	25.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.9409	1.042		22.1	20.0	10.7	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.6163		20.2	20.0	0.9	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L2ID		1.126		18.4	20.0	-7.8	25.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.111	1.031		18.6	20.0	-7.1	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_045.d  
 Lims ID: CCV L4  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 20-Feb-2016 05:40:35 ALS Bottle#: 12 Worklist Smp#: 43  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L4  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5

Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 09:59:40 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d

Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: westendorfc

Date: 22-Feb-2016 08:19:42

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.700	5.712	-0.012	739750	46.2		92.5	3474	
2 Perfluorobutyric acid	212.9 > 169.0	5.697	5.715	-0.018	424275	22.0		110	916	
D 3 13C5-PFPeA	267.9 > 223.0	6.805	6.827	-0.022	1587614	51.6		103	4733	
4 Perfluoropentanoic acid	262.9 > 219.0	6.805	6.828	-0.023	652654	19.7		98.5	158	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.920	6.939	-0.019	269562	NC			336	
	298.9 > 99.0	6.920	6.939	-0.019	144355		1.87(0.00-0.00)		365	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.920	6.939	-0.019	269562	21.2		120		
7 Perfluorohexanoic acid	313.0 > 269.0	8.049	8.078	-0.029	603219	20.8		104	1219	
D 6 13C2 PFHxA	315.0 > 270.0	8.049	8.080	-0.031	1339541	47.3		94.6	6869	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.125	8.158	-0.033	203419	NC			674	
D 8 13C4-PFHpA	367.0 > 322.0	9.288	9.322	-0.034	1454678	47.2		94.4	6491	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.288	9.322	-0.034	692643	22.8		114	1501	
D 11 18O2 PFHxS	403.0 > 84.0	9.323	9.358	-0.035	606393	40.5		85.7	2799	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.317	9.364	-0.047	181033	NC			433	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.317	9.364	-0.047	1.000	181033	20.9	111		
D 12 13C4 PFOA	417.0 > 372.0	10.412	10.450	-0.038		1545984	45.3	90.6	5776	
13 Perfluorooctanoic acid	413.0 > 369.0	10.412	10.450	-0.038	1.000	635798	20.5	102	696	
	413.0 > 169.0	10.412	10.450	-0.038	1.000	210818		3.02(0.00-0.00)	628	
14 Perfluoroheptane Sulfonate	449.0 > 80.0	10.419	10.457	-0.038	1.000	200818	NC		691	
38 Perfluoroheptanesulfonic Acid	449.0 > 80.0	10.419	10.457	-0.038	1.000	200818	22.3	117		
D 16 13C4 PFOS	503.0 > 80.0	11.377	11.414	-0.037		697291	38.3	80.2	2392	
15 Perfluorooctane sulfonic acid	499.0 > 80.0	11.377	11.417	-0.040	1.000	299312	21.1	110	470	
	499.0 > 99.0	11.377	11.417	-0.040	1.000	141830		2.11(0.00-0.00)	479	
D 17 13C5 PFNA	468.0 > 423.0	11.393	11.434	-0.041		1345222	47.7	95.4	4165	
18 Perfluorononanoic acid	463.0 > 419.0	11.400	11.439	-0.039	1.000	426205	19.6	98.1	1868	
21 PFNS (Perflouro-1-nonanesulfonate)	549.0 > 80.0	12.203	12.249	-0.046	1.000	178226	NC		594	
D 19 13C2 PFDA	515.0 > 470.0	12.234	12.275	-0.041		1304985	45.6	91.2	4599	
20 Perfluorodecanoic acid	513.0 > 469.0	12.234	12.276	-0.042	1.000	475208	18.3	91.6	1260	
D 23 13C8 FOSA	506.0 > 78.0	12.783	12.810	-0.027		1591012	38.3	76.7	1051	
24 Perfluorooctane Sulfonamide	498.0 > 78.0	12.783	12.812	-0.029	1.000	561444	19.0	95.2	1156	
25 Perfluorodecane Sulfonate	599.0 > 80.0	12.918	12.949	-0.031	1.000	189262	NC		962	
39 Perfluorodecane Sulfonic acid	599.0 > 80.0	12.918	12.949	-0.031	1.000	189262	23.6	122		
D 26 13C2 PFUnA	565.0 > 520.0	12.960	13.000	-0.040		1676629	44.9	89.7	6651	
27 Perfluoroundecanoic acid	563.0 > 519.0	12.960	13.000	-0.040	1.000	597798	21.0	105	1584	
D 28 13C2 PFDaA	615.0 > 570.0	13.578	13.609	-0.031		1825247	46.7	93.3	7049	
29 Perfluorododecanoic acid	613.0 > 569.0	13.578	13.611	-0.033	1.000	585761	21.9	110	852	
31 PFDoS (Perflouro-1-dodecanesulfona	699.0 > 80.0	14.039	14.083	-0.044	1.000	191566	NC		863	
30 Perfluorotridecanoic acid	663.0 > 619.0	14.093	14.130	-0.037	1.000	760506	22.1	111	1212	
32 Perfluorotetradecanoic acid	713.0 > 669.0	14.539	14.571	-0.032	1.000	449980	20.2	101	341	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 33 13C2-PFTeDA										
715.0 > 670.0	14.539	14.571	-0.032		1556683	43.4		86.7	5943	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.177	15.207	-0.030		1775802	41.2		82.4	6748	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.177	15.208	-0.031	1.000	822211	18.4		92.2	1559	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.514	15.543	-0.029	1.000	753047	18.6		92.9	1332	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L4\_00017

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_045.d

Injection Date: 20-Feb-2016 05:40:35

Instrument ID: A6

Lims ID: CCV L4

Client ID:

Operator ID: JRB

ALS Bottle#: 12

Worklist Smp#: 43

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

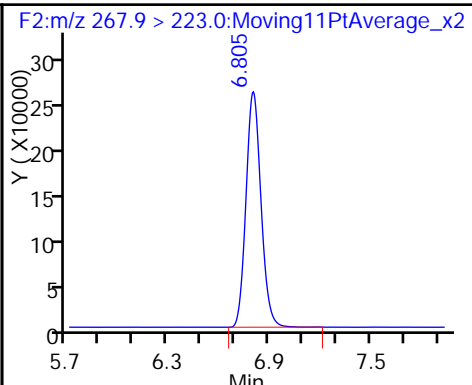
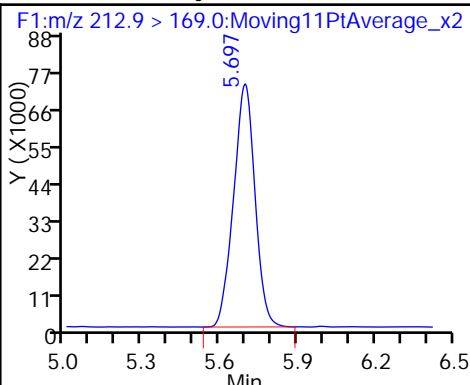
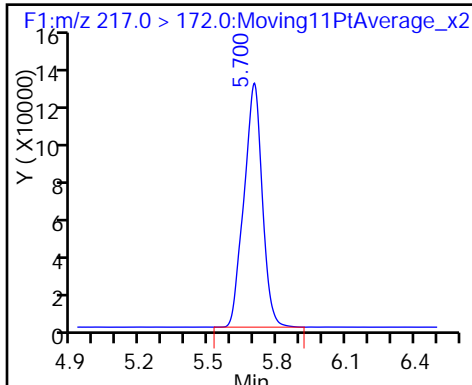
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

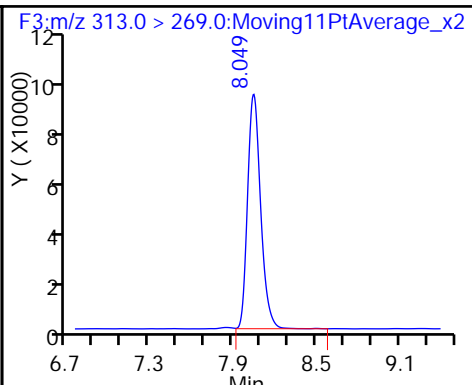
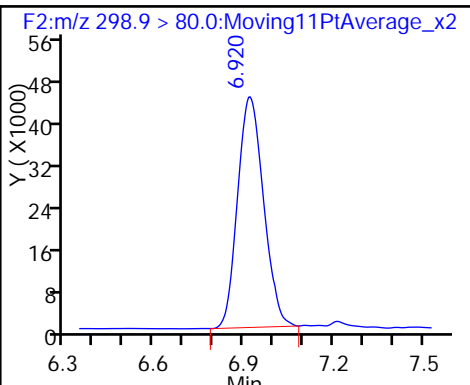
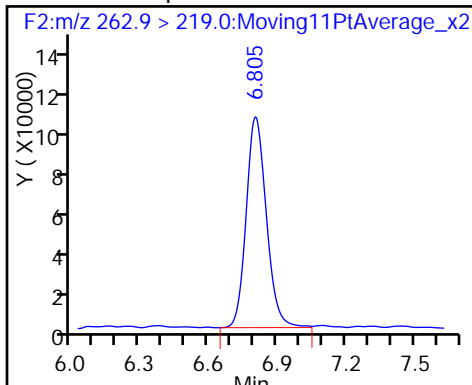
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

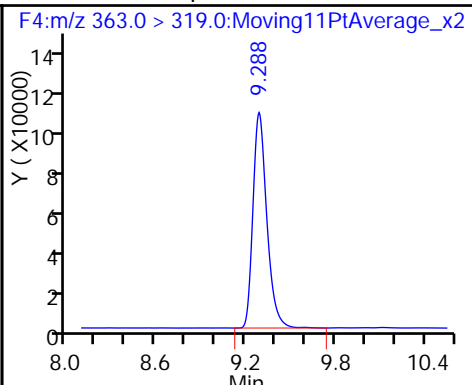
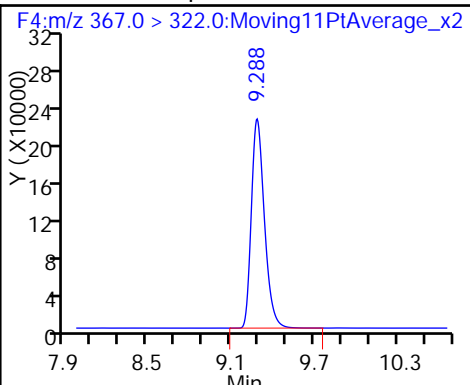
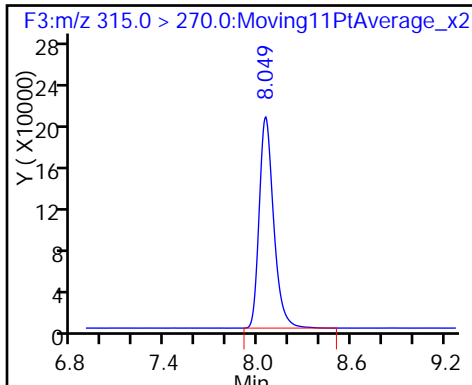
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

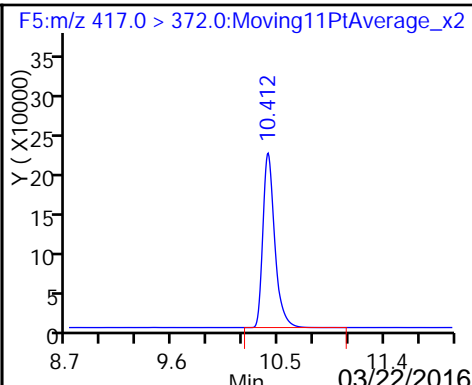
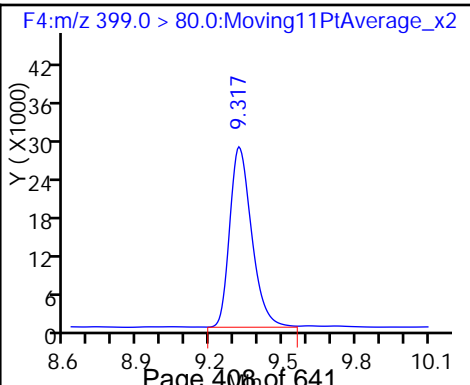
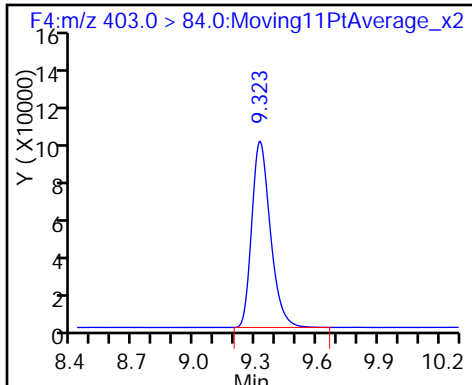
9 Perfluoroheptanoic acid



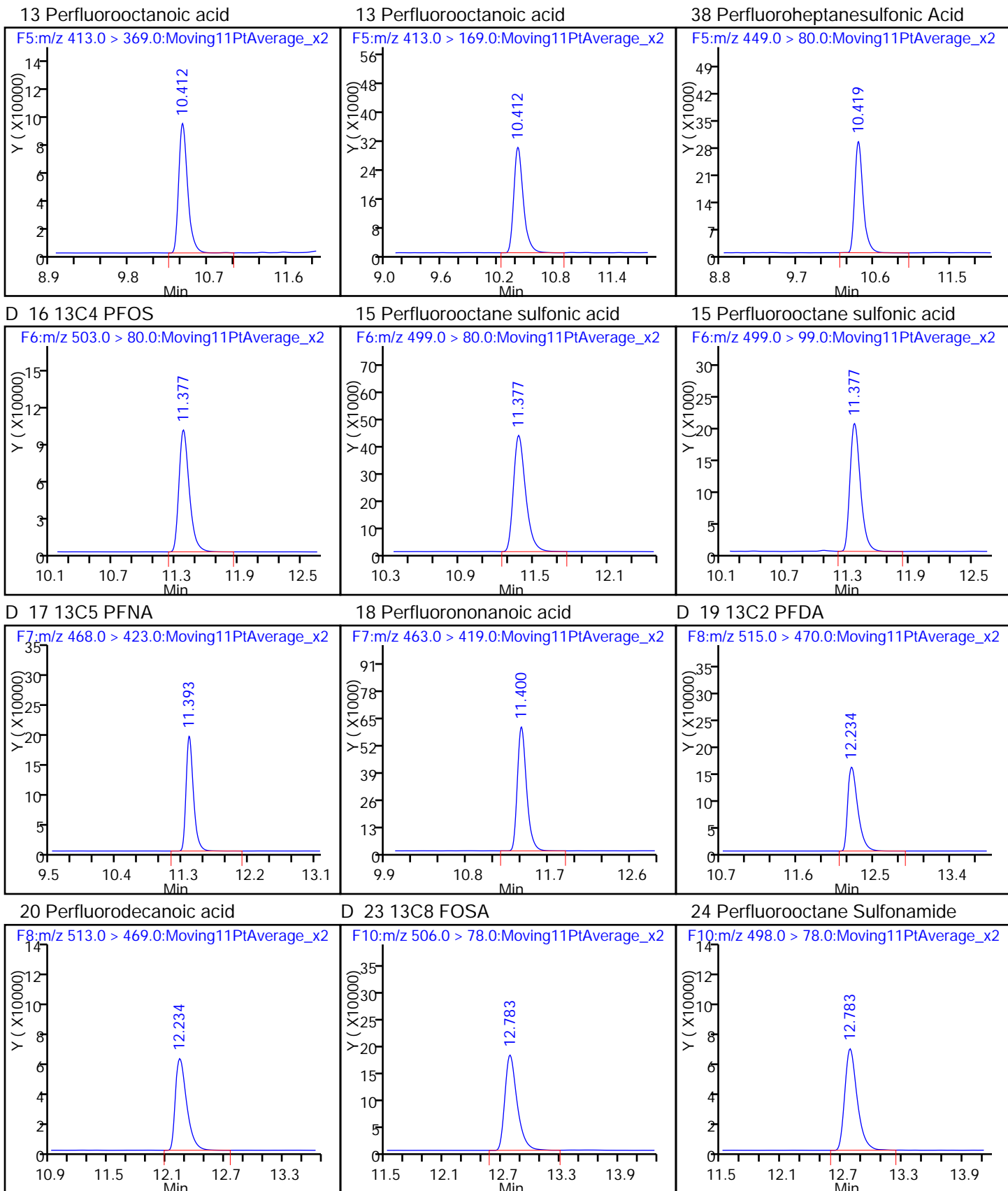
D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA



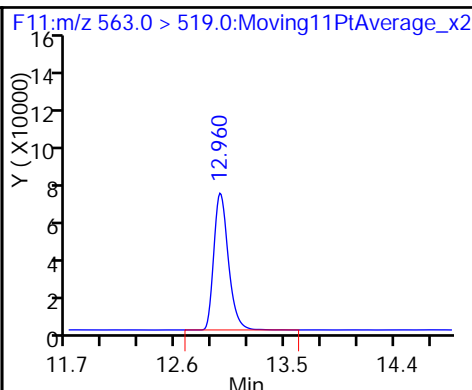
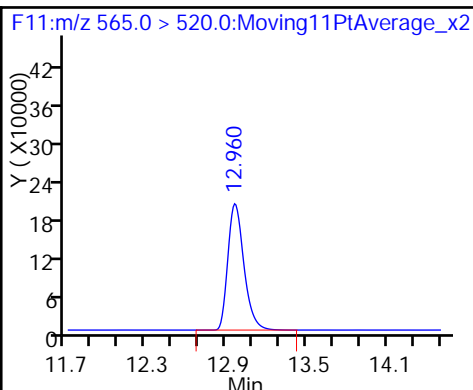
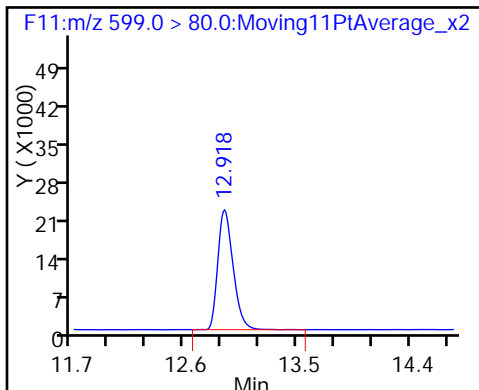




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

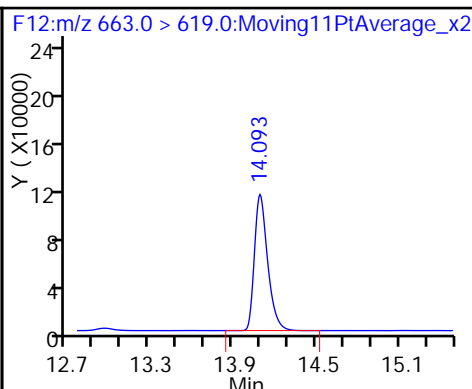
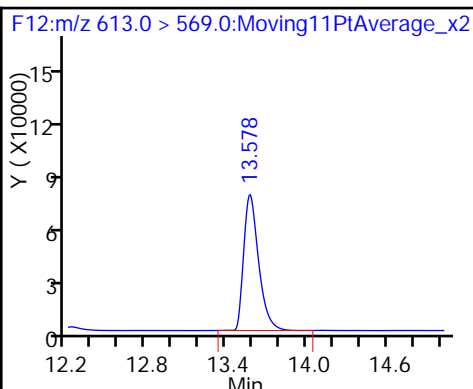
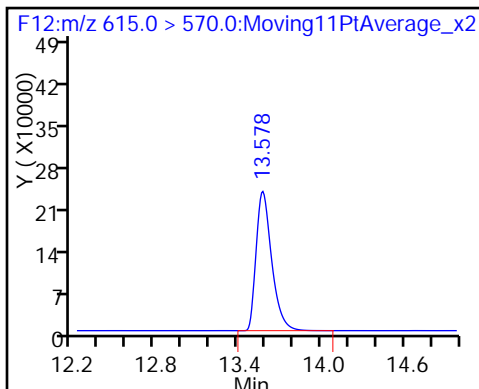
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

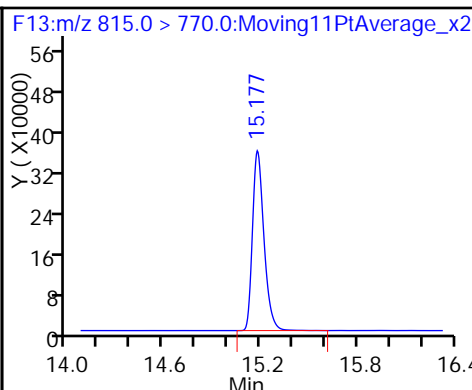
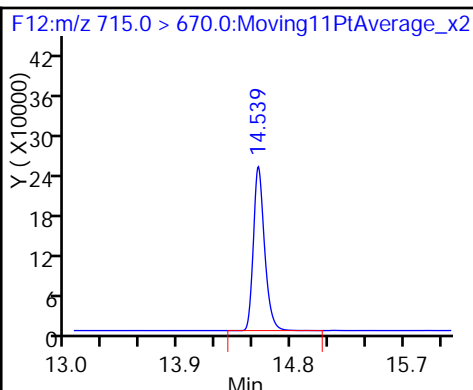
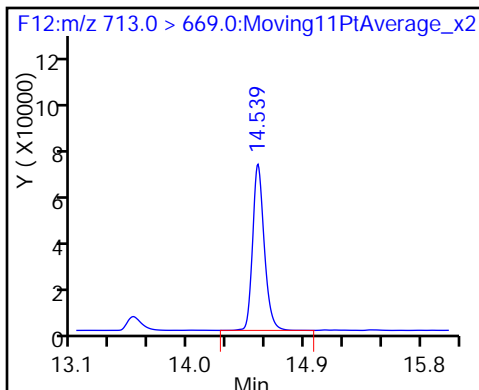
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

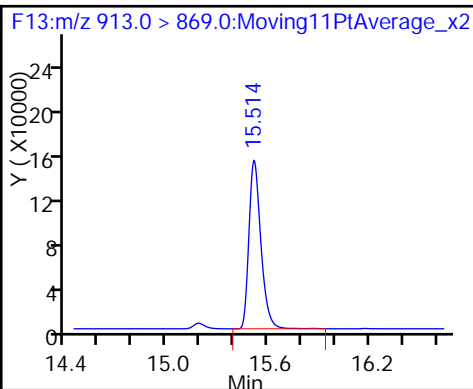
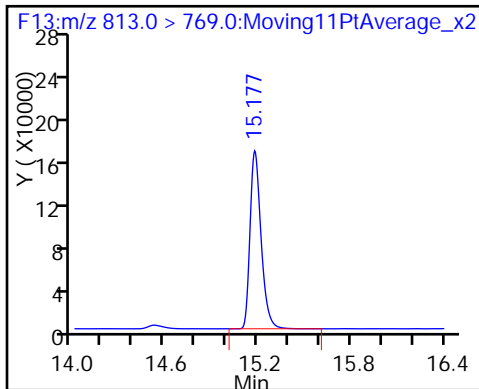
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: CCV 320-100906/54 Calibration Date: 02/20/2016 09:34  
 Instrument ID: A6 Calib Start Date: 02/19/2016 15:08  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/19/2016 17:15  
 Lab File ID: 19FEB2016A6A\_056.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	1.304	1.412		54.1	50.0	8.3	25.0
Perfluoropentanoic acid (PFPeA)	AveID	1.043	0.9755		46.7	50.0	-6.5	25.0
Perfluorobutanesulfonic acid (PFBS)	AveID	0.9930	1.104		49.1	44.2	11.1	25.0
Perfluorohexanoic acid (PFHxA)	AveID	1.083	1.144		52.8	50.0	5.6	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.043	1.018		48.8	50.0	-2.4	25.0
Perfluorohexanesulfonic acid (PFHxS)	AveID	0.6753	0.7199		50.4	47.3	6.6	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.004	1.000		49.8	50.0	-0.4	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	L2ID		0.6299		48.2	47.6	1.2	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	0.9724	1.010		49.6	47.8	3.8	25.0
Perfluorononanoic acid (PFNA)	AveID	0.8075	0.8407		52.1	50.0	4.1	25.0
Perfluorodecanoic acid (PFDA)	L2ID		1.010		51.5	50.0	2.9	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9268	0.8877		47.9	50.0	-4.2	25.0
Perfluorodecane Sulfonic acid	L1ID		0.6249		54.3	48.2	12.7	25.0
Perfluoroundecanoic acid (PFUnA)	L1ID		0.8572		51.4	50.0	2.7	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.7311	0.8069		55.2	50.0	10.4	25.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.9409	1.037		55.1	50.0	10.2	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.6323		52.5	50.0	5.0	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L2ID		1.064		47.0	50.0	-5.9	25.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.111	1.025		46.1	50.0	-7.7	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_056.d  
 Lims ID: CCV L5  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 20-Feb-2016 09:34:04 ALS Bottle#: 13 Worklist Smp#: 54  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L5  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 10:18:05 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.696	5.712	-0.016	735784	46.0		92.0	2562	
2 Perfluorobutyric acid	212.9 > 169.0	5.696	5.715	-0.019	1038956	54.1		108	3162	
D 3 13C5-PFPeA	267.9 > 223.0	6.800	6.827	-0.027	1596544	51.9		104	6136	
4 Perfluoropentanoic acid	262.9 > 219.0	6.805	6.828	-0.023	1557494	46.7		93.5	340	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.920	6.939	-0.019	634088	NC			805	
	298.9 > 99.0	6.915	6.939	-0.024	340139		1.86(0.00-0.00)		723	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.920	6.939	-0.019	634088	49.1		111		
7 Perfluorohexanoic acid	313.0 > 269.0	8.049	8.078	-0.029	1592823	52.8		106	3176	
D 6 13C2 PFHxA	315.0 > 270.0	8.049	8.080	-0.031	1392598	49.2		98.3	5464	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.131	8.158	-0.027	444944	NC			1556	
D 8 13C4-PFHpA	367.0 > 322.0	9.288	9.322	-0.034	1592041	51.6		103	6480	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.288	9.322	-0.034	1619972	48.8		97.6	3886	
D 11 18O2 PFHxS	403.0 > 84.0	9.323	9.358	-0.035	614918	41.1		86.9	2481	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.323	9.364	-0.041	442664	NC			1286	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.323	9.364	-0.041	442664	50.4		107		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.412	10.450	-0.038		1618742	47.4		94.8	5890	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.412	10.450	-0.038	1.000	1618258	49.8		99.6	2042	
413.0 > 169.0	10.419	10.450	-0.031	1.001	515853		3.14(0.00-0.00)		1384	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.419	10.457	-0.038	1.000	435358	NC			1348	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.419	10.457	-0.038	1.000	435358	48.2		101		
D 16 13C4 PFOS										
503.0 > 80.0	11.385	11.414	-0.029		694093	38.2		79.9	2463	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.385	11.417	-0.032	1.000	700885	49.6		104	445	
499.0 > 99.0	11.385	11.417	-0.032	1.000	365929		1.92(0.00-0.00)		1939	
D 17 13C5 PFNA										
468.0 > 423.0	11.400	11.434	-0.034		1363986	48.4		96.8	6799	
18 Perfluorononanoic acid										
463.0 > 419.0	11.400	11.439	-0.039	1.000	1146645	52.1		104	3094	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.204	12.249	-0.045	1.000	413024	NC			1920	
D 19 13C2 PFDA										
515.0 > 470.0	12.245	12.275	-0.030		1287906	45.0		90.0	3873	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.245	12.276	-0.031	1.000	1300889	51.5		103	4985	
D 23 13C8 FOSA										
506.0 > 78.0	12.783	12.810	-0.027		1825198	44.0		88.0	3566	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.783	12.812	-0.029	1.000	1620215	47.9		95.8	1429	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.919	12.949	-0.030	1.000	437365	NC			1584	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.919	12.949	-0.030	1.000	437365	54.3		113		
D 26 13C2 PFUnA										
565.0 > 520.0	12.971	13.000	-0.029		1701977	45.5		91.1	5975	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.971	13.000	-0.029	1.000	1458974	51.4		103	4237	
D 28 13C2 PFDaA										
615.0 > 570.0	13.578	13.609	-0.031		1913739	48.9		97.8	4707	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.578	13.611	-0.033	1.000	1544099	55.2		110	2102	
31 PFDoS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.048	14.083	-0.035	1.000	468581	NC			1591	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.102	14.130	-0.028	1.000	1985226	55.1		110	1838	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.546	14.571	-0.025	1.000	1209996	52.5		105	605	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.546	14.571	-0.025		1833393	51.1		102	6312	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA										
815.0 > 770.0	15.180	15.207	-0.027		2004655	46.5		93.0	6714	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.180	15.208	-0.028	1.000	2035836	47.0		94.1	2224	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.517	15.543	-0.026	1.000	1961521	46.1		92.3	1793	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L5\_00016

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_056.d

Injection Date: 20-Feb-2016 09:34:04

Instrument ID: A6

Lims ID: CCV L5

Client ID:

Operator ID: JRB

ALS Bottle#: 13

Worklist Smp#: 54

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

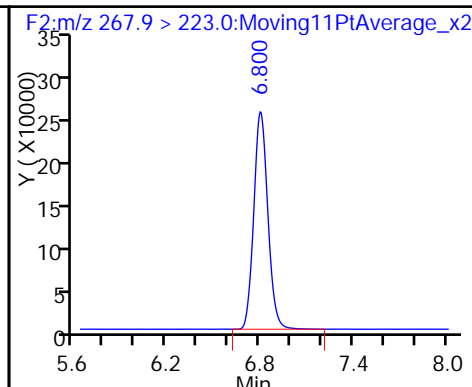
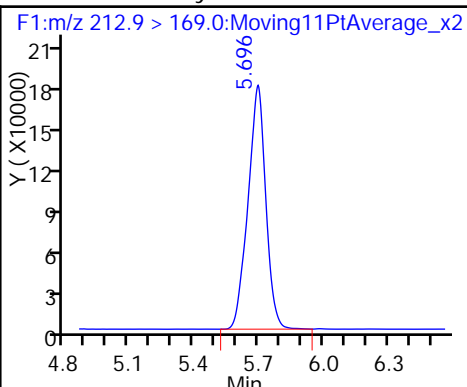
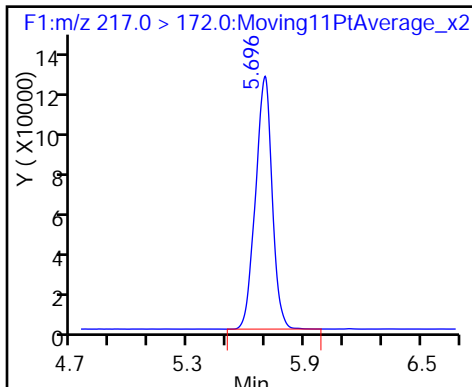
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

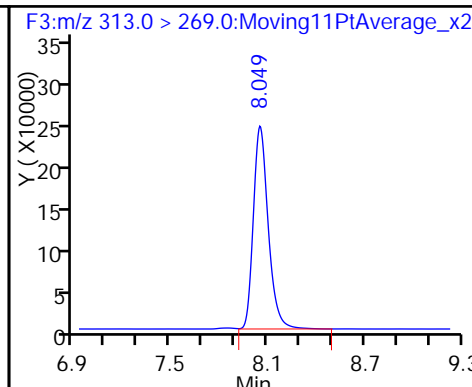
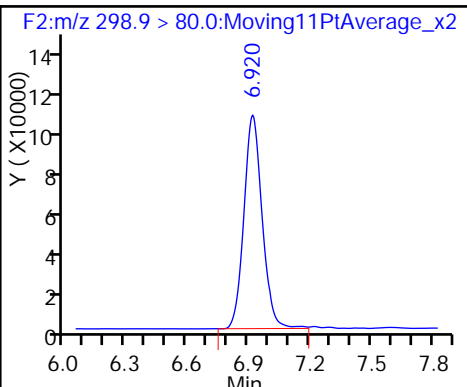
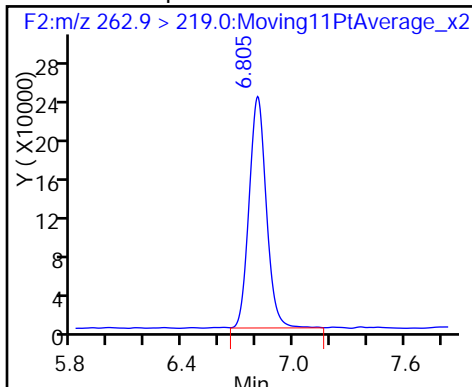
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

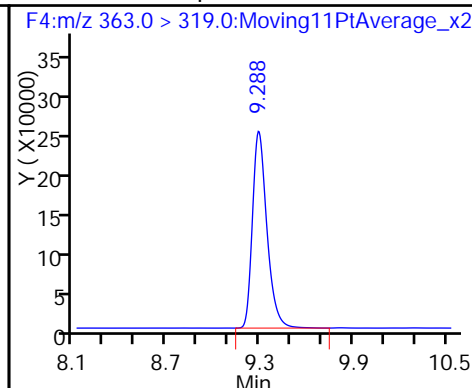
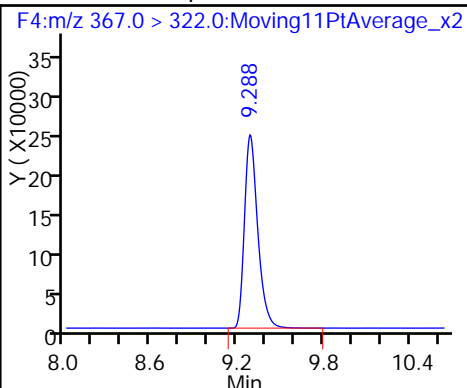
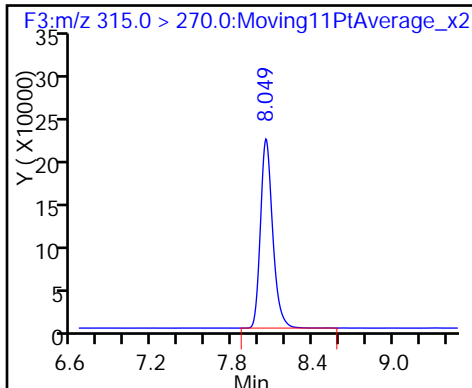
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

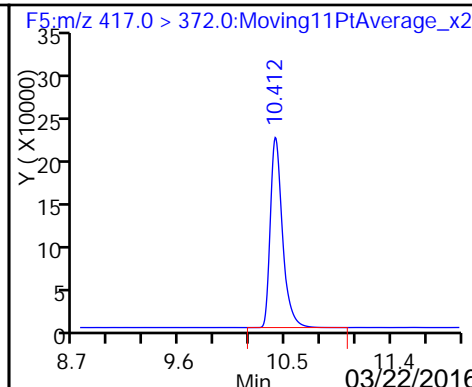
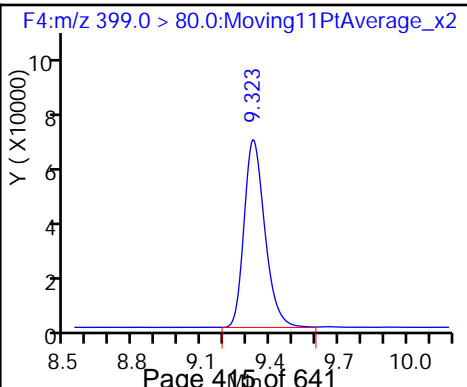
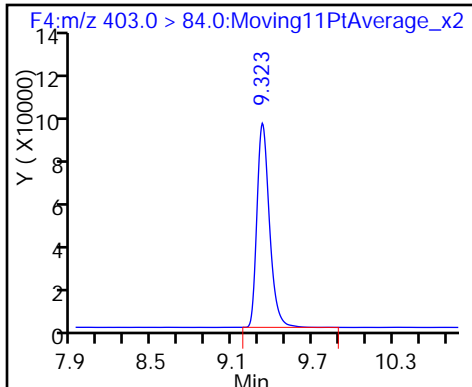
9 Perfluoroheptanoic acid

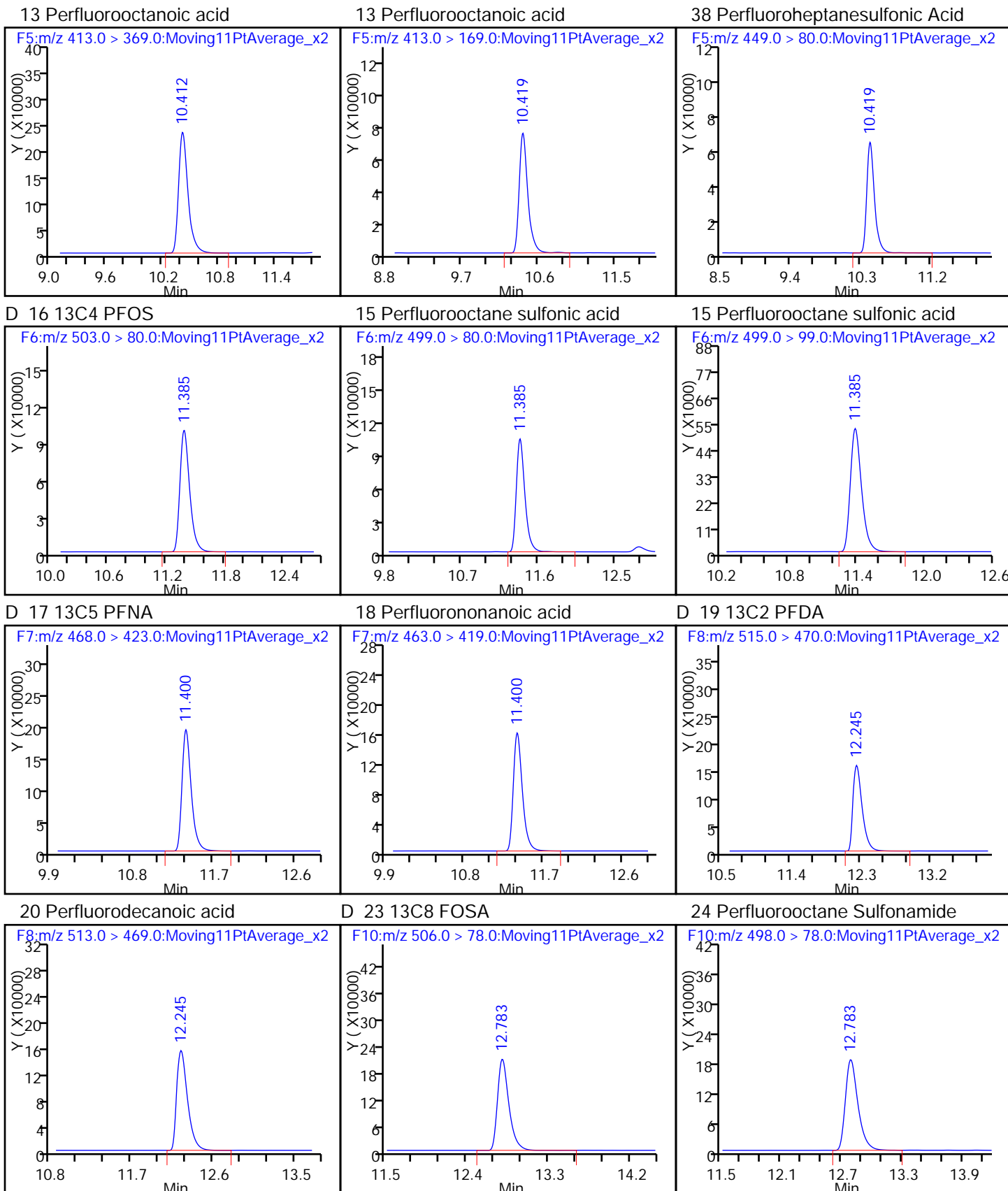


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA



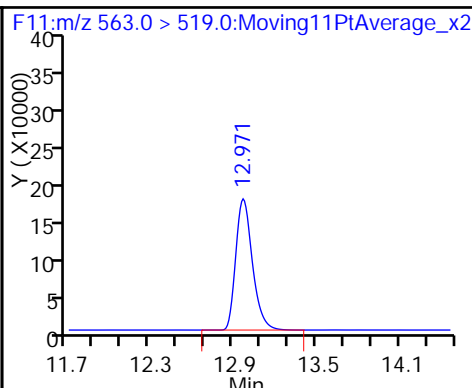
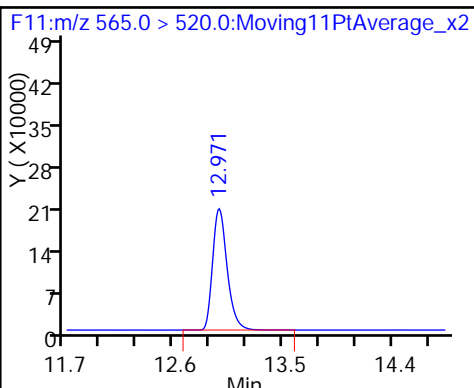
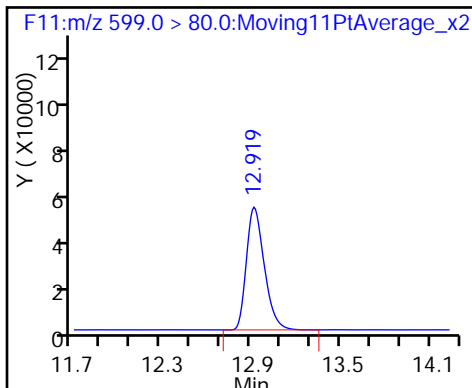




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

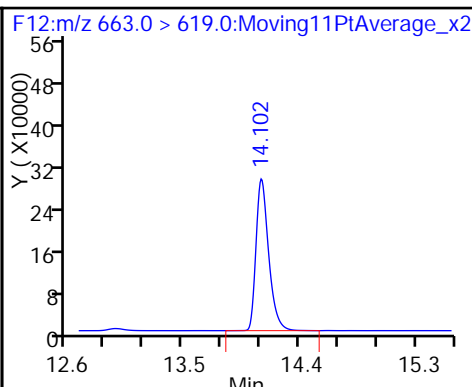
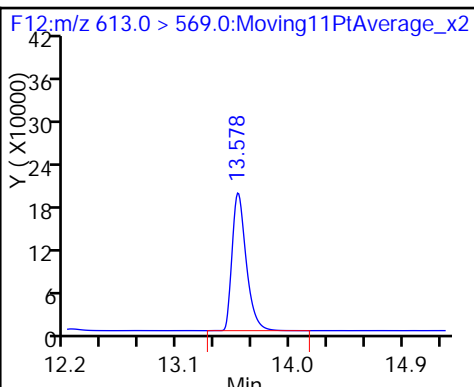
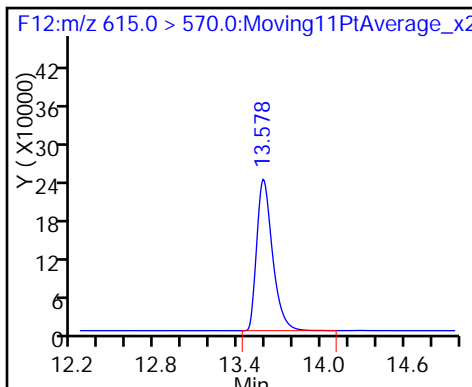
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

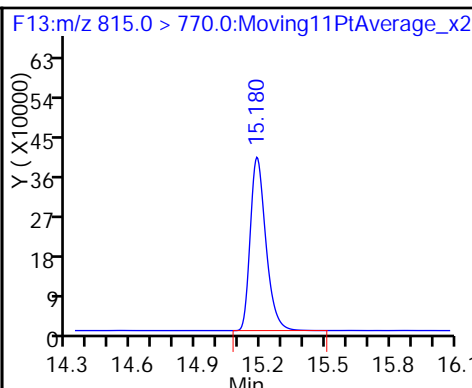
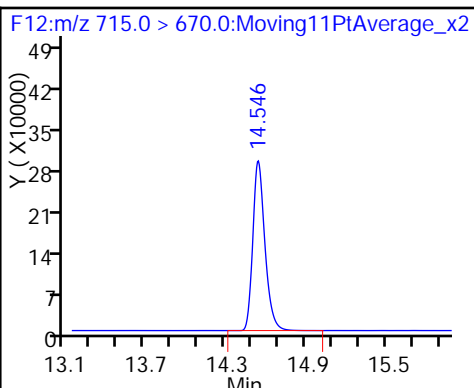
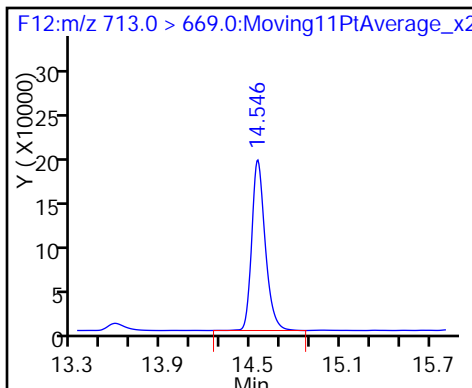
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

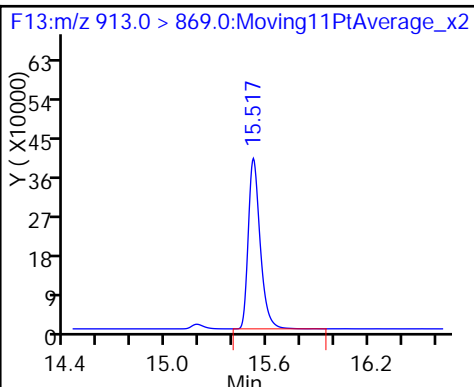
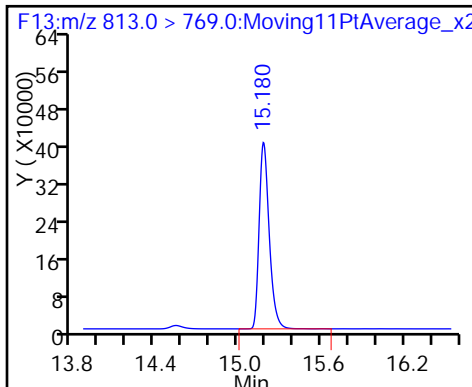
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Lab Sample ID: CCV 320-100906/65 Calibration Date: 02/20/2016 13:27  
 Instrument ID: A6 Calib Start Date: 02/19/2016 15:08  
 GC Column: Acquity ID: 2.10 (mm) Calib End Date: 02/19/2016 17:15  
 Lab File ID: 19FEB2016A6A\_067.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	1.304	1.413		21.7	20.0	8.4	25.0
Perfluoropentanoic acid (PFPeA)	AveID	1.043	1.005		19.3	20.0	-3.7	25.0
Perfluorobutanesulfonic acid (PFBS)	AveID	0.9930	0.9556		17.0	17.7	-3.8	25.0
Perfluorohexanoic acid (PFHxA)	AveID	1.083	1.152		21.3	20.0	6.3	25.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.043	1.149		22.0	20.0	10.2	25.0
Perfluorohexanesulfonic acid (PFHxS)	AveID	0.6753	0.6952		19.5	18.9	2.9	25.0
Perfluoroheptanesulfonic Acid (PFHpS)	L2ID		0.6011		18.6	19.0	-2.5	25.0
Perfluorooctanoic acid (PFOA)	AveID	1.004	0.9799		19.5	20.0	-2.4	25.0
Perfluorooctanesulfonic acid (PFOS)	AveID	0.9724	0.9873		19.4	19.1	1.5	25.0
Perfluorononanoic acid (PFNA)	AveID	0.8075	0.9268		23.0	20.0	14.8	25.0
Perfluorodecanoic acid (PFDA)	L2ID		0.9888		19.9	20.0	-0.4	25.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9268	0.9281		20.0	20.0	0.1	25.0
Perfluorodecane Sulfonic acid	L1ID		0.5449		19.1	19.3	-0.8	25.0
Perfluoroundecanoic acid (PFUnA)	L1ID		0.9003		21.2	20.0	5.9	25.0
Perfluorododecanoic acid (PFDoA)	AveID	0.7311	0.7269		19.9	20.0	-0.6	25.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.9409	0.9810		20.9	20.0	4.3	25.0
Perfluorotetradecanoic acid (PFTeA)	L2ID		0.5337		17.4	20.0	-12.9	25.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	L2ID		1.202		19.9	20.0	-0.7	25.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.111	0.9712		17.5	20.0	-12.6	25.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_067.d  
 Lims ID: CCV L4  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 20-Feb-2016 13:27:32 ALS Bottle#: 12 Worklist Smp#: 65  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L4  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Sublist: chrom-PFAC\_A6\*sub5  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 11:18:21 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: westendorfc Date: 21-Feb-2016 13:19:47

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.699	5.712	-0.013	844637	52.8		106	3042	
2 Perfluorobutyric acid	212.9 > 169.0	5.699	5.715	-0.016	477412	21.7		108	1440	
D 3 13C5-PFPeA	267.9 > 223.0	6.804	6.827	-0.023	1885401	61.3		123	5167	
4 Perfluoropentanoic acid	262.9 > 219.0	6.804	6.828	-0.024	757574	19.3		96.3	163	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.929	6.939	-0.010	277743	NC			326	
	298.9 > 99.0	6.920	6.939	-0.019	153377		1.81(0.00-0.00)		222	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.929	6.939	-0.010	277743	17.0		96.2		
7 Perfluorohexanoic acid	313.0 > 269.0	8.055	8.078	-0.023	730390	21.3		106	1455	
D 6 13C2 PFHxA	315.0 > 270.0	8.055	8.080	-0.025	1585078	56.0		112	5137	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.131	8.158	-0.027	217669	NC			1096	
D 8 13C4-PFHpA	367.0 > 322.0	9.288	9.322	-0.034	1821976	59.1		118	6569	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.294	9.322	-0.028	837547	22.0		110	3808	
D 11 18O2 PFHxS	403.0 > 84.0	9.329	9.358	-0.029	777616	52.0		110	3271	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.329	9.364	-0.035	216237	NC			848	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.329	9.364	-0.035	1.000	216237	19.5	103		
D 12 13C4 PFOA	417.0 > 372.0	10.419	10.450	-0.031		1951511	57.2	114	10200	
13 Perfluorooctanoic acid	413.0 > 369.0	10.419	10.450	-0.031	1.000	764939	19.5	97.6	803	
	413.0 > 169.0	10.419	10.450	-0.031	1.000	285474	2.68(0.00-0.00)		705	
14 Perfluoroheptane Sulfonate	449.0 > 80.0	10.419	10.457	-0.038	1.000	212630	NC		693	
38 Perfluoroheptanesulfonic Acid	449.0 > 80.0	10.419	10.457	-0.038	1.000	212630	18.6	97.5		
D 16 13C4 PFOS	503.0 > 80.0	11.385	11.414	-0.029		888097	48.8	102	3468	
15 Perfluorooctane sulfonic acid	499.0 > 80.0	11.385	11.417	-0.032	1.000	350731	19.4	102	295	
	499.0 > 99.0	11.385	11.417	-0.032	1.000	168205	2.09(0.00-0.00)		725	
D 17 13C5 PFNA	468.0 > 423.0	11.408	11.434	-0.026		1605407	56.9	114	7343	
18 Perfluorononanoic acid	463.0 > 419.0	11.408	11.439	-0.031	1.000	595129	23.0	115	1413	
21 PFNS (Perflouro-1-nonanesulfonate)	549.0 > 80.0	12.214	12.249	-0.035	1.000	216600	NC		797	
D 19 13C2 PFDA	515.0 > 470.0	12.245	12.275	-0.030		1606537	56.1	112	3436	
20 Perfluorodecanoic acid	513.0 > 469.0	12.245	12.276	-0.031	1.000	635386	19.9	99.6	1693	
D 23 13C8 FOSA	506.0 > 78.0	12.794	12.810	-0.016		2216095	53.4	107	2538	
24 Perfluorooctane Sulfonamide	498.0 > 78.0	12.794	12.812	-0.018	1.000	822680	20.0	100	1616	
25 Perfluorodecane Sulfonate	599.0 > 80.0	12.919	12.949	-0.030	1.000	195178	NC		580	
39 Perfluorodecane Sulfonic acid	599.0 > 80.0	12.919	12.949	-0.030	1.000	195178	19.1	99.2		
D 26 13C2 PFUnA	565.0 > 520.0	12.971	13.000	-0.029		2056267	55.0	110	5372	
27 Perfluoroundecanoic acid	563.0 > 519.0	12.971	13.000	-0.029	1.000	740493	21.2	106	2090	
D 28 13C2 PFDaA	615.0 > 570.0	13.579	13.609	-0.030		2309380	59.0	118	6770	
29 Perfluorododecanoic acid	613.0 > 569.0	13.579	13.611	-0.032	1.000	671444	19.9	99.4	904	
31 PFDaS (Perflouro-1-dodecanesulfona	699.0 > 80.0	14.048	14.083	-0.035	1.000	225190	NC		743	
30 Perfluorotridecanoic acid	663.0 > 619.0	14.102	14.130	-0.028	1.000	906215	20.9	104	935	
32 Perfluorotetradecanoic acid	713.0 > 669.0	14.546	14.571	-0.025	1.000	492963	17.4	87.1	248	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 33 13C2-PFTeDA										
715.0 > 670.0	14.546	14.571	-0.025		1919727	53.5		107	6759	
D 35 13C2-PFHxDA										
815.0 > 770.0	15.180	15.207	-0.027		2251795	52.2		104	5198	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.180	15.208	-0.028	1.000	1109943	19.9		99.3	1278	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.517	15.543	-0.026	1.000	897159	17.5		87.4	997	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

**Reagents:**

LCPFC-L4\_00017

Amount Added: 1.00

Units: mL

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_067.d

Injection Date: 20-Feb-2016 13:27:32

Instrument ID: A6

Lims ID: CCV L4

Client ID:

Operator ID: JRB

ALS Bottle#: 12

Worklist Smp#: 65

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

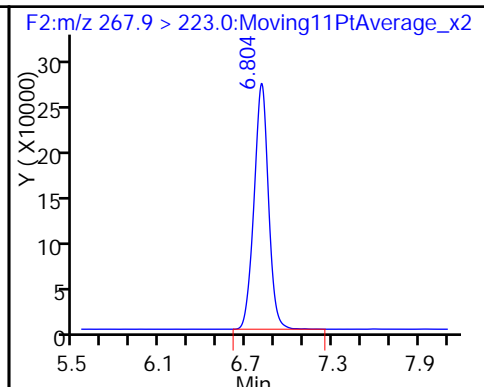
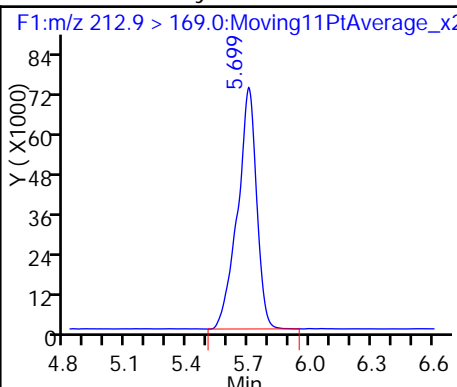
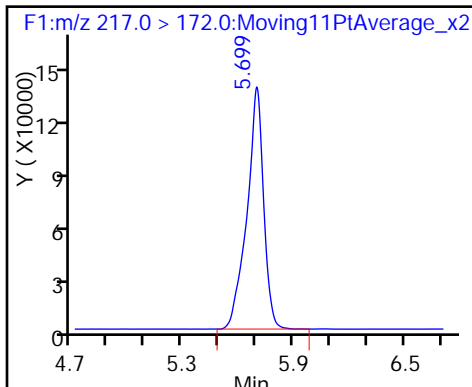
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

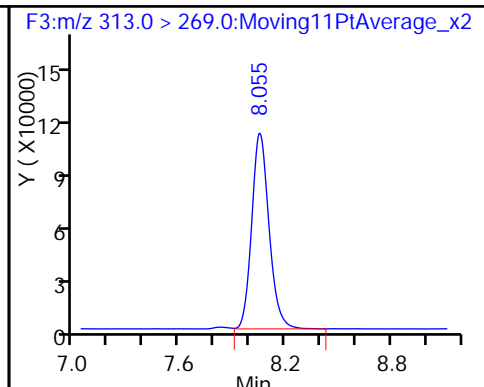
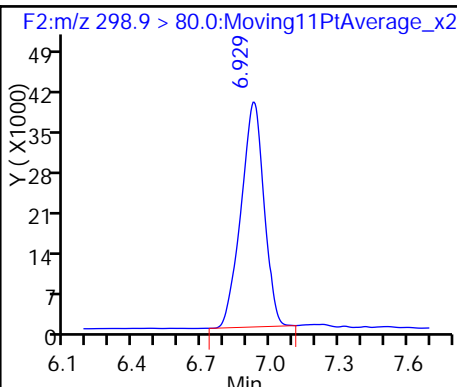
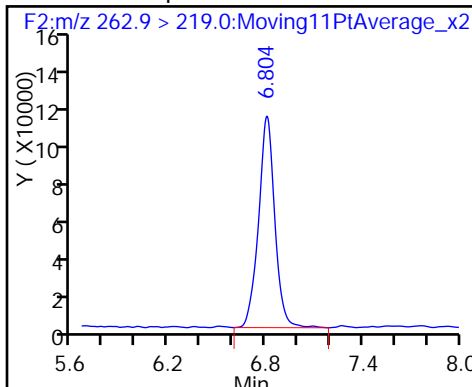
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

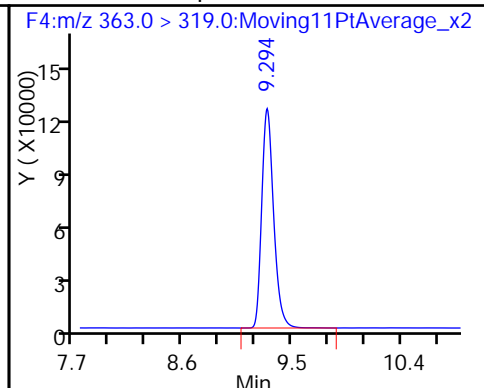
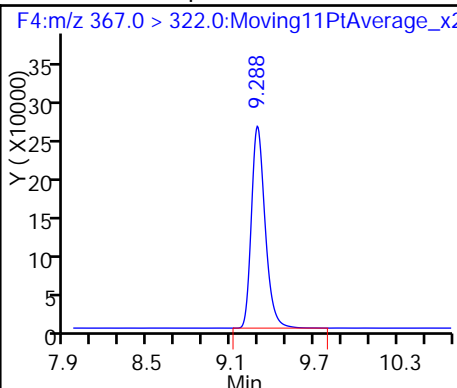
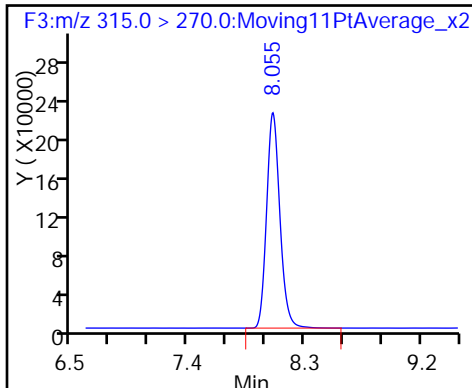
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

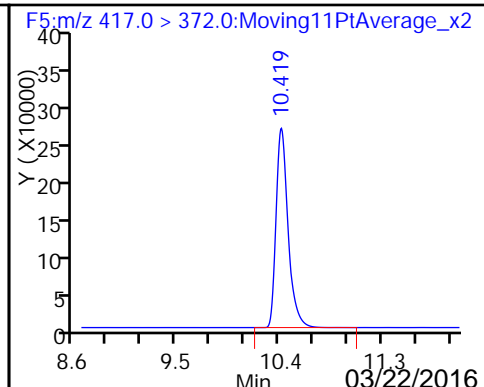
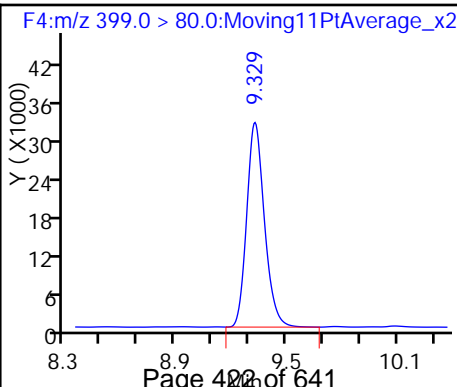
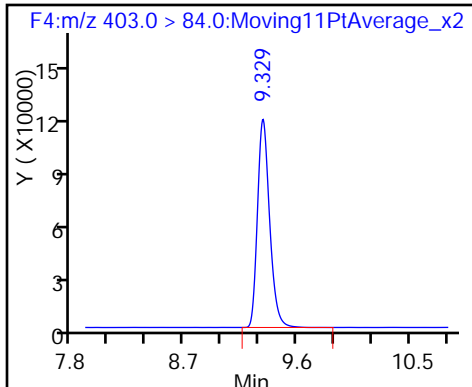
9 Perfluoroheptanoic acid

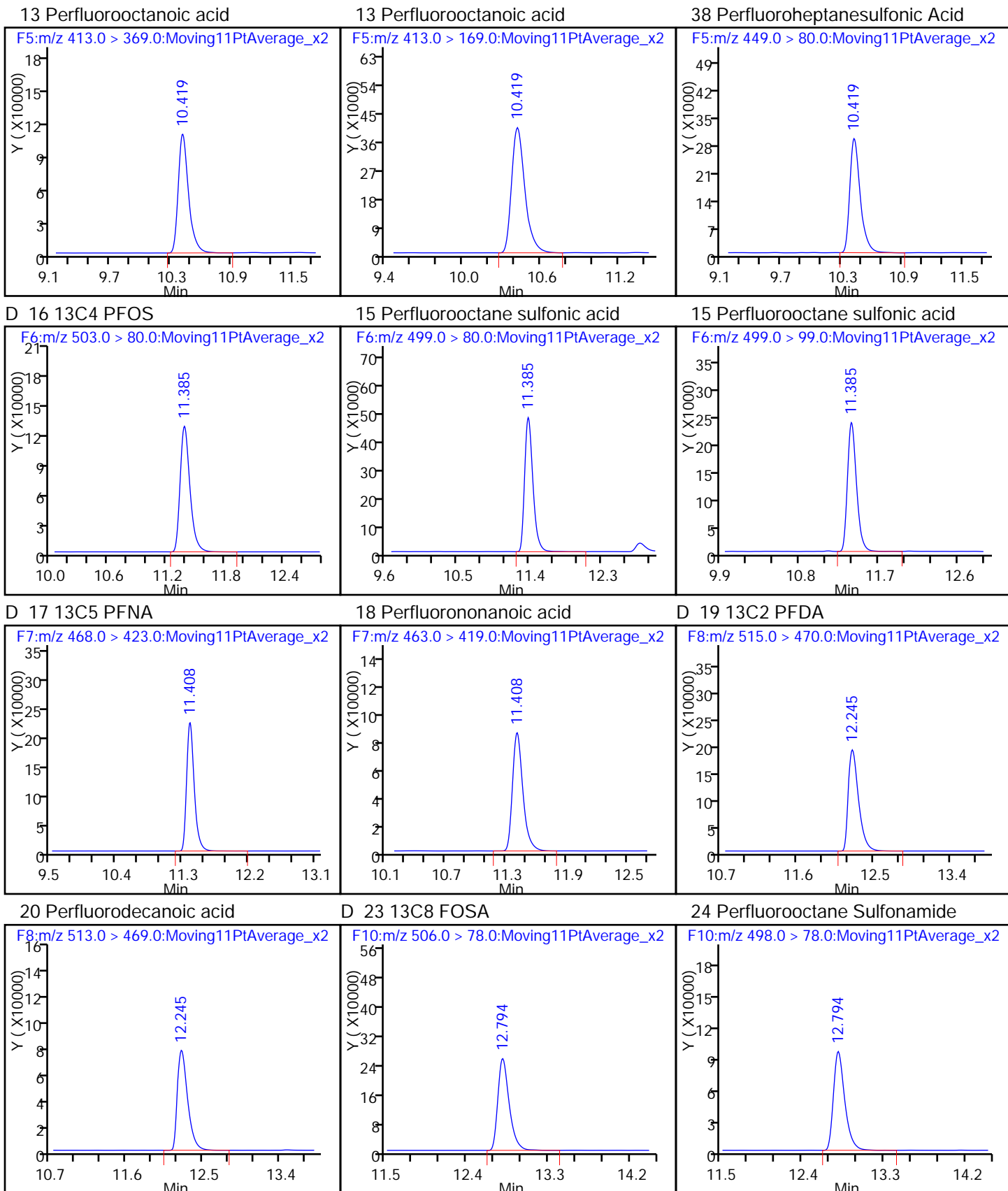


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

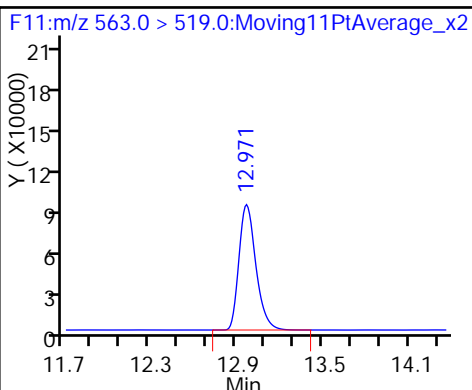
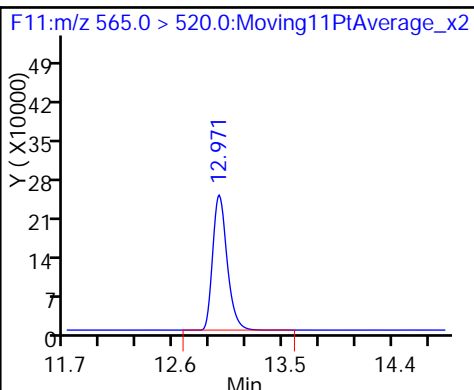
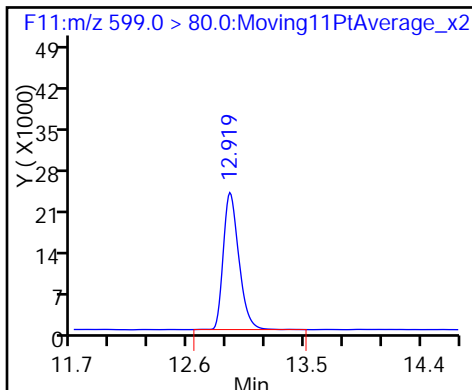




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUa

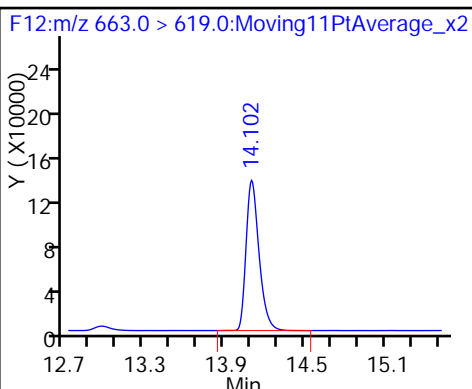
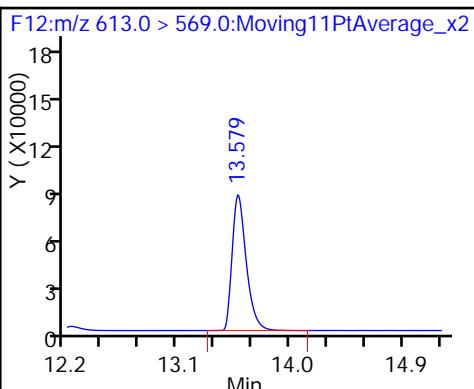
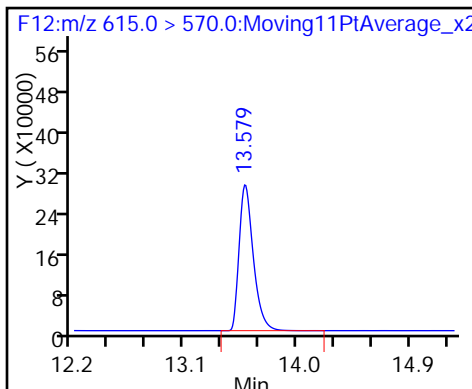
27 Perfluoroundecanoic acid



D 28 13C2 PFDa

29 Perfluorododecanoic acid

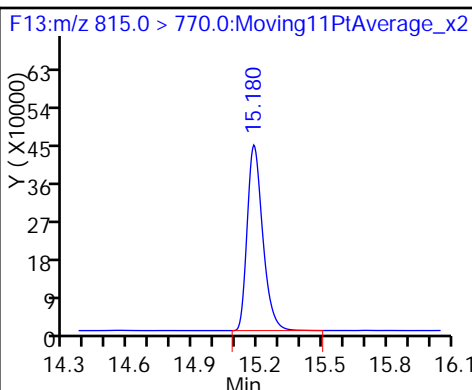
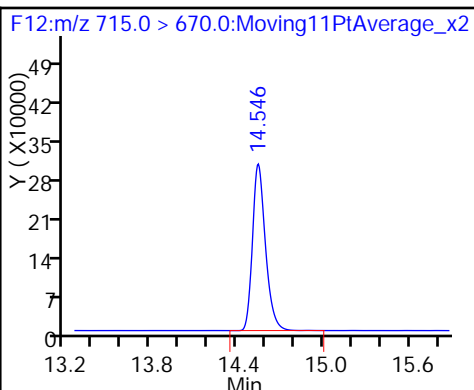
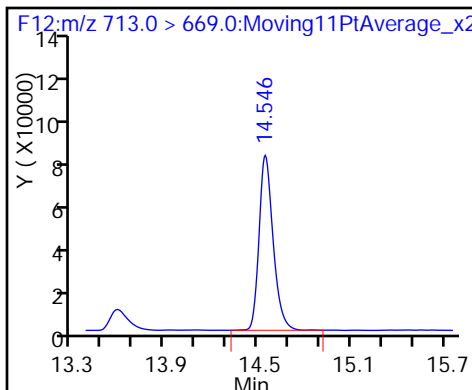
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

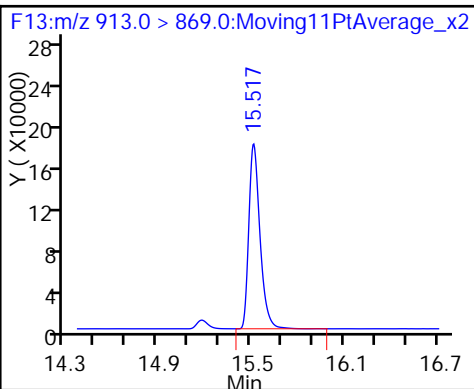
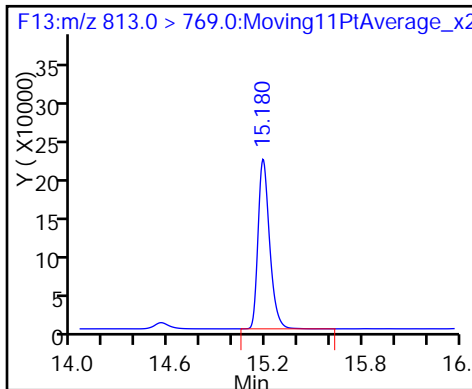
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid





FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 320-100093/1-A  
 Matrix: Water Lab File ID: 12FEB2016A4A\_042.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/09/2016 10:23  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/13/2016 00:22  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100305 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.0020	0.00087

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00994	18O2 PFHxS	95		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_042.d  
 Lims ID: MB 320-100093/1-A  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 13-Feb-2016 00:22:35 ALS Bottle#: 26 Worklist Smp#: 41  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: MB 320-100093/1-A box 18  
 Misc. Info.: Acquity BEH C18,1.7u, 3X150mm,T=35C  
 Operator ID: JRB Instrument ID: A4  
 Method: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\PFAC\_A4.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 15-Feb-2016 15:26:28 Calib Date: 12-Feb-2016 12:42:35  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_009.d  
 Column 1 : Det: F1:MRM  
 Process Host: XAWRK014

First Level Reviewer: westendorfc Date: 15-Feb-2016 09:32:44

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	216.7 > 171.5	5.502	5.712	-0.210	2107465	51.9		104	9071	
D 3 13C5-PFPeA	267.6 > 222.7	6.704	6.986	-0.282	1537026	48.0		96.1	4613	
5 Perfluorobutane Sulfonate	298.8 > 79.6	6.814	7.115	-0.301	224725	NC			128	
51 Perfluorobutanesulfonic acid	298.8 > 79.6	6.814	7.115	-0.301	224725	2.29				
D 6 13C2 PFHxA	314.6 > 269.7	8.017	8.367	-0.350	2045654	52.2		104	6042	
D 8 13C4-PFHpA	366.6 > 321.6	9.264	9.639	-0.375	1970849	50.6		101	6246	
9 Perfluoroheptanoic acid	362.8 > 318.7	9.264	9.640	-0.376	10783	0.5343			35.7	
58 Perfluorohexanesulfonic acid	398.3 > 79.2	9.293	9.675	-0.382	5224	-0.1095				
10 Perfluorohexane Sulfonate	398.3 > 79.2	9.293	9.675	-0.382	5224	NC			18.1	
D 11 18O2 PFHxS	402.5 > 83.6	9.299	9.675	-0.376	1017192	44.9		94.8	2451	
D 12 13C4 PFOA	416.5 > 371.6	10.370	10.743	-0.373	2259154	53.1		106	6800	
D 16 13C4 PFOS	502.4 > 79.7	11.313	11.651	-0.338	403051	39.8		83.3	1399	
D 17 13C5 PFNA	467.5 > 422.6	11.336	11.678	-0.342	1999657	48.1		96.1	5363	
D 19 13C2 PFDA	514.4 > 469.5	12.159	12.473	-0.314	2903115	49.4		98.8	6083	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 23 13C8 FOSA	505.4 > 77.6	12.701	12.991	-0.290		817248	9.15	18.3	792	
D 26 13C2 PFUnA	564.3 > 519.5	12.858	13.146	-0.288		2629742	45.0	89.9	3600	
D 28 13C2 PFDoA	614.4 > 569.4	13.460	13.713	-0.253		2970716	41.8	83.5	4140	
30 Perfluorotridecanoic acid	662.4 > 618.5	13.957	14.193	-0.236	1.000	2883	0.0550		2.9	
D 33 13C2-PFTeDA	714.5 > 669.5	14.379	14.603	-0.224		2359844	37.2	74.5	3807	
32 Perfluorotetradecanoic acid	712.6 > 668.5	14.386	14.604	-0.218	1.000	9243	0.1772		4.2	
D 35 13C2-PFHxDA	814.8 > 769.6	14.981	15.138	-0.157		1104662	37.2	74.5	2839	
34 Perfluorohexadecanoic acid	812.6 > 768.6	14.981	15.139	-0.158	1.000	43581	0.5370		58.9	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A4\20160212-28353.b\12FEB2016A4A\_042.d

Injection Date: 13-Feb-2016 00:22:35

Instrument ID: A4

Lims ID: MB 320-100093/1-A

Client ID:

Operator ID: JRB

ALS Bottle#: 26

Worklist Smp#: 41

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

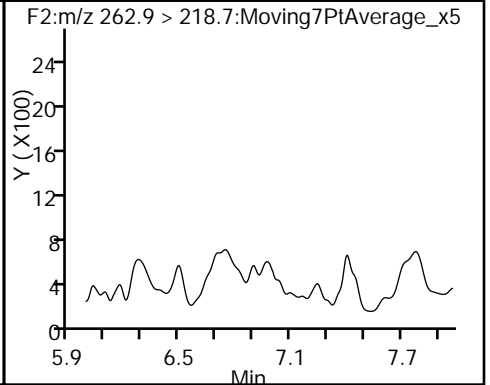
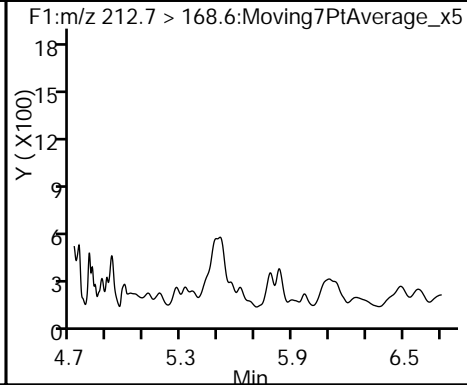
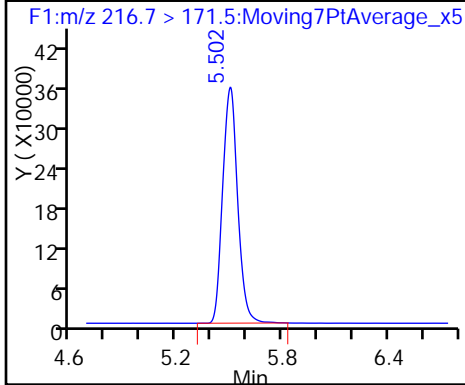
Method: PFAC\_A4

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid (ND)

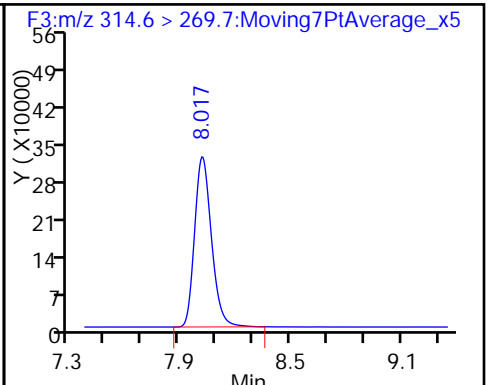
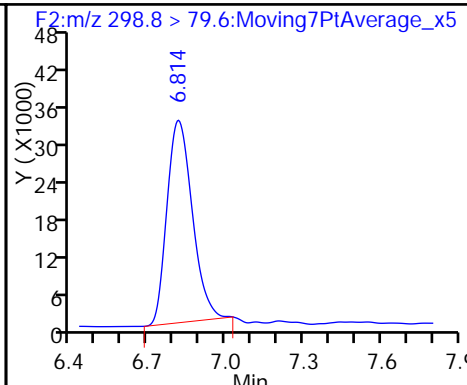
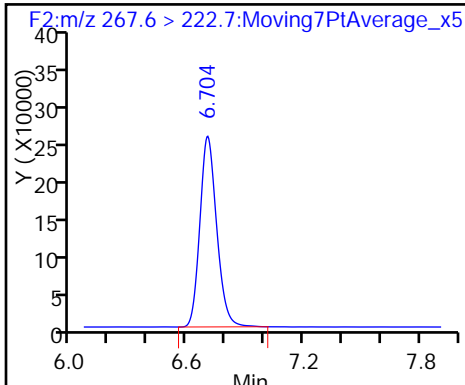
4 Perfluoropentanoic acid (ND)



D 3 13C5-PFPeA

51 Perfluorobutanesulfonic acid

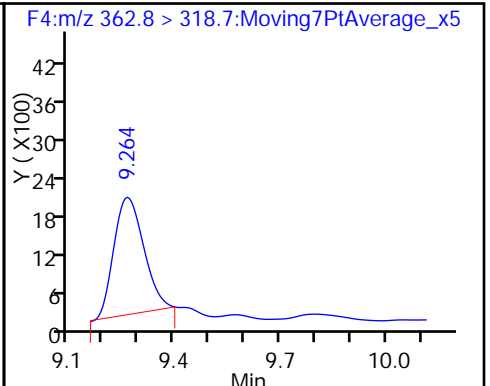
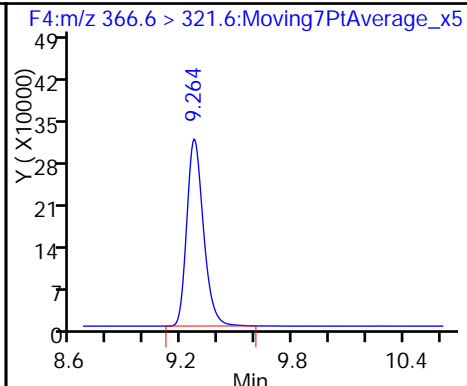
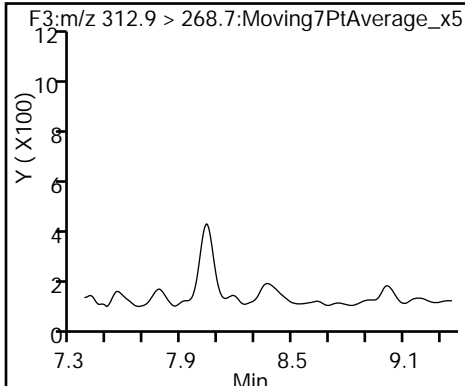
D 6 13C2 PFHxA



7 Perfluorohexanoic acid (ND)

D 8 13C4-PFHpA

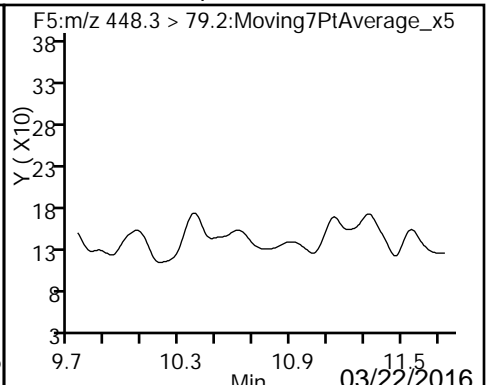
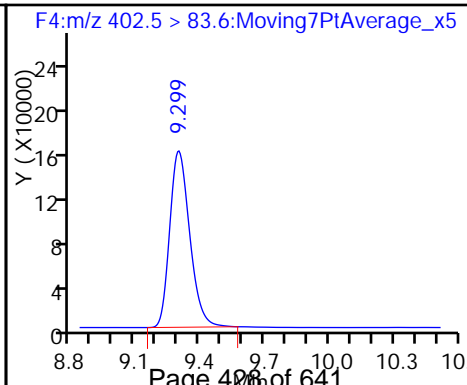
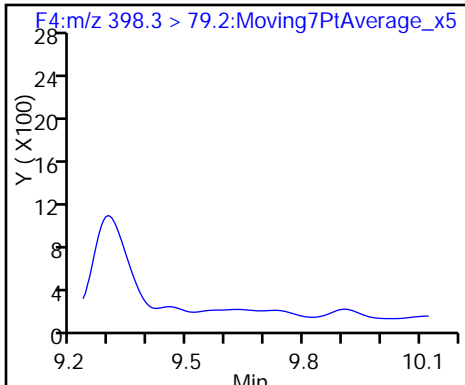
9 Perfluoroheptanoic acid



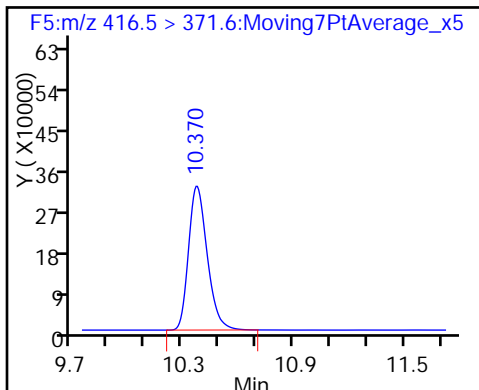
58 Perfluorohexanesulfonic acid

D 11 18O2 PFHxS

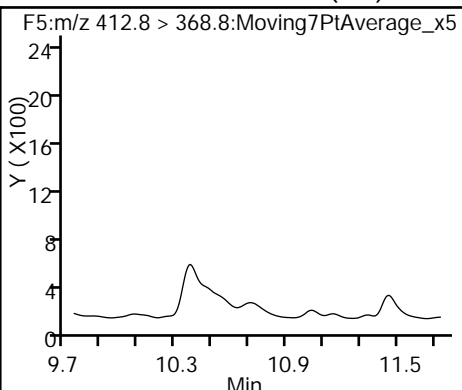
39 Perfluoroheptanesulfonic Acid (ND)



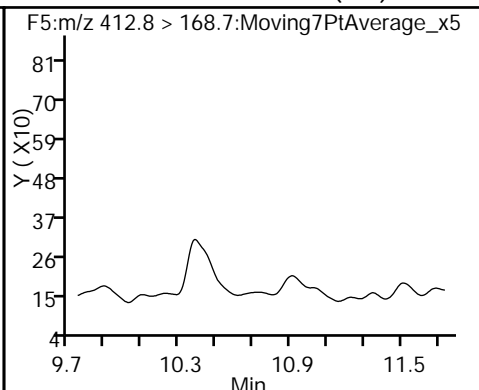
D 12 13C4 PFOA



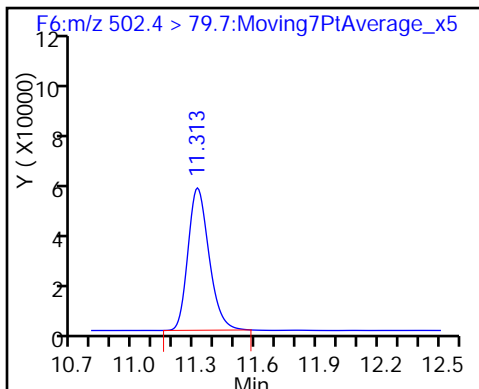
13 Perfluorooctanoic acid (ND)



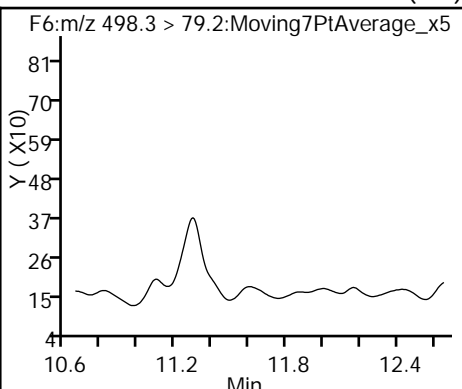
13 Perfluorooctanoic acid (ND)



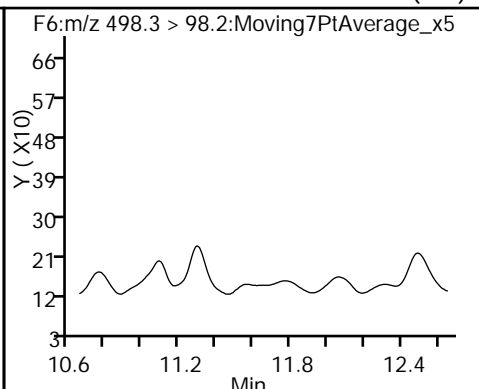
D 16 13C4 PFOS



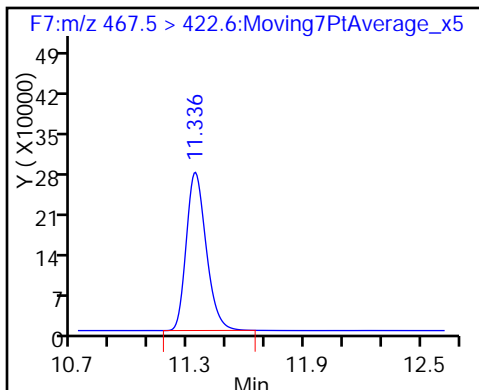
15 Perfluorooctane sulfonic acid (ND)



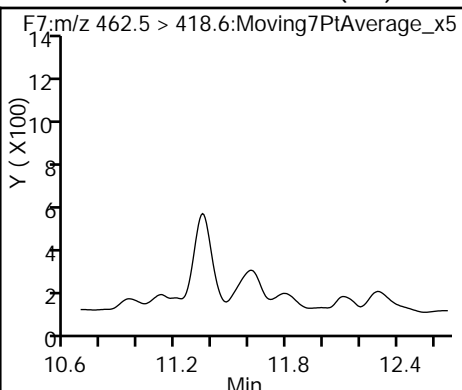
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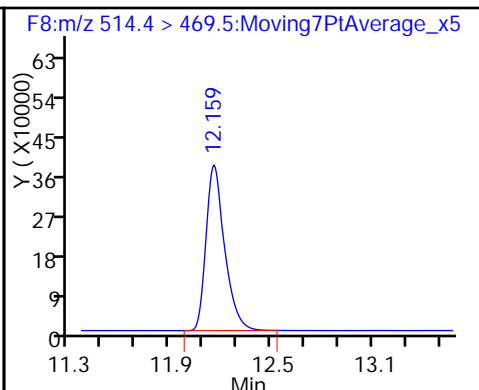
D 17 13C5 PFNA



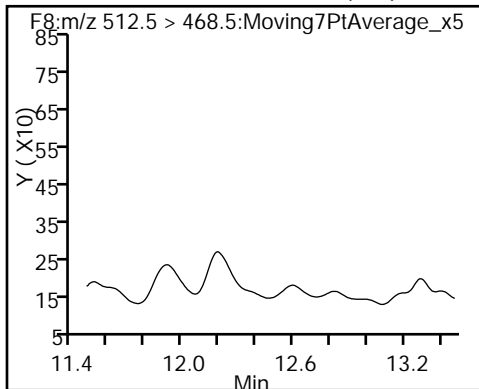
18 Perfluorononanoic acid (ND)



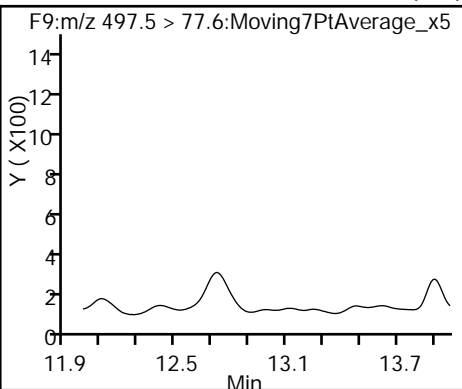
D 19 13C2 PFDA



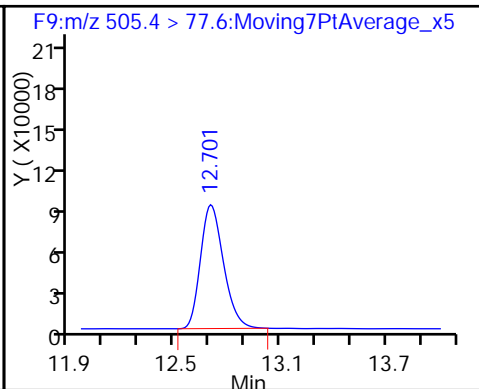
20 Perfluorodecanoic acid (ND)



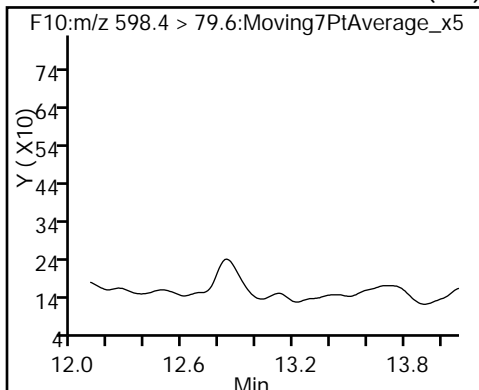
24 Perfluorooctane Sulfonamide (ND)



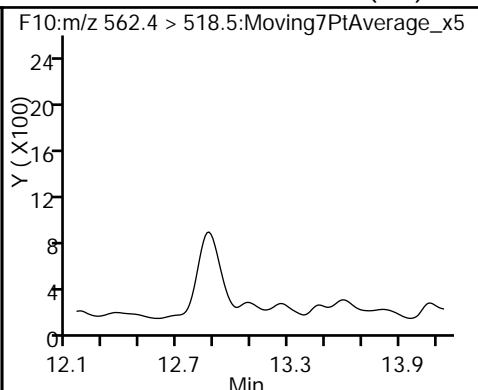
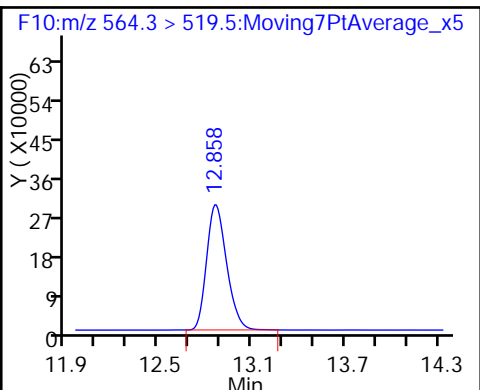
D 23 13C8 FOSA



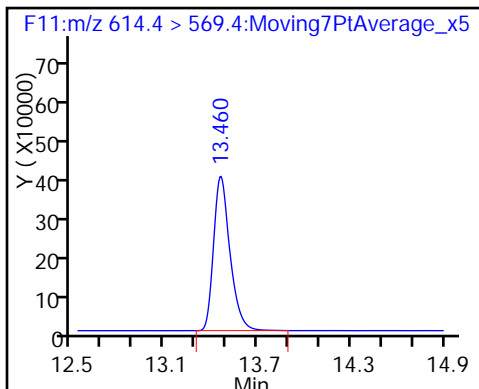
49 Perfluorodecane Sulfonic acid (ND) D 26 13C2 PFUnA



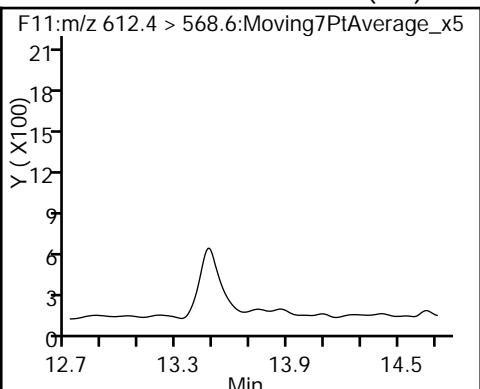
27 Perfluoroundecanoic acid (ND)



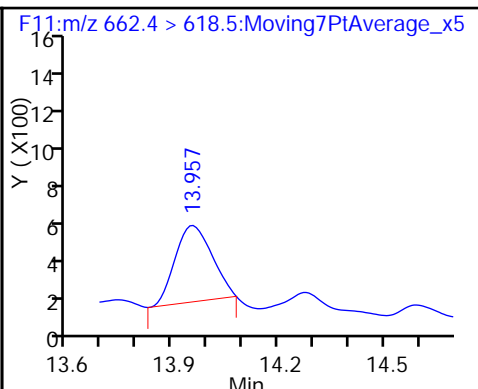
D 28 13C2 PFDaA



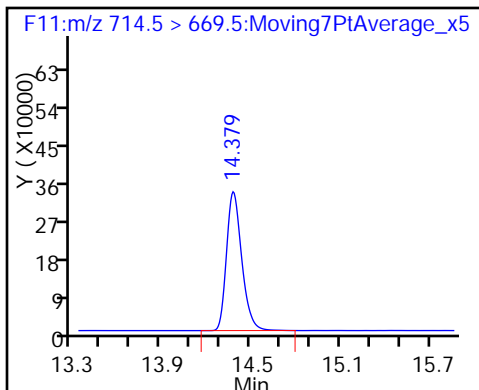
29 Perfluorododecanoic acid (ND)



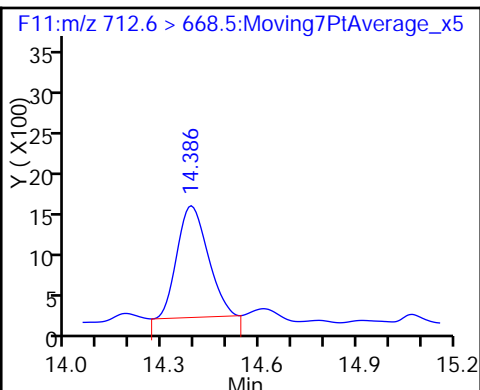
30 Perfluorotridecanoic acid



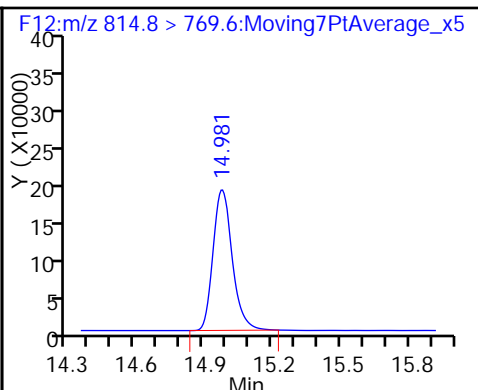
D 33 13C2-PFTeDA



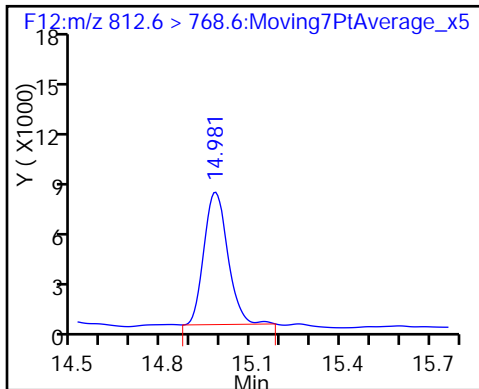
32 Perfluorotetradecanoic acid



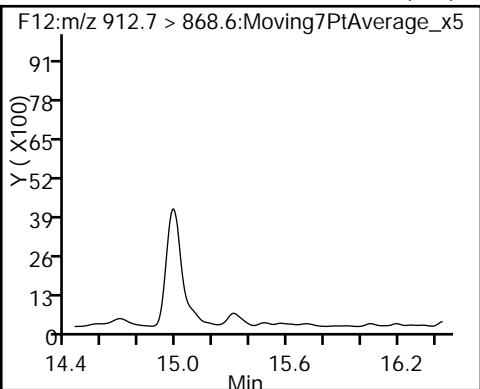
D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid



36 Perfluorooctadecanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 320-100093/1-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_047.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/09/2016 10:23  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 06:23  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.0020	0.00092
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.00143	J	0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	103		25-150
STL00990	13C4 PFOA	101		25-150
STL00995	13C5 PFNA	103		25-150
STL00991	13C4 PFOS	92		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_047.d  
 Lims ID: MB 320-100093/1-A  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 20-Feb-2016 06:23:01 ALS Bottle#: 1 Worklist Smp#: 45  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: MB 320-100093/1-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:29:40

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	217.0 > 172.0	5.690	5.712	-0.022	796044	49.8		99.5	3613	
2 Perfluorobutyric acid	212.9 > 169.0	5.635	5.715	-0.080	3850	0.1854			3.2	
D 3 13C5-PFPeA	267.9 > 223.0	6.800	6.827	-0.027	1689147	54.9		110	5481	
4 Perfluoropentanoic acid	262.9 > 219.0	6.823	6.828	-0.005	8535	0.2421			1.5	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.924	6.939	-0.015	2554	NC			5.0	
	298.9 > 99.0	6.938	6.939	-0.001	2255		1.13(0.00-0.00)		8.8	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.924	6.939	-0.015	2554	0.1761				
7 Perfluorohexanoic acid	313.0 > 269.0	8.047	8.078	-0.031	55187	1.78			69.6	
D 6 13C2 PFHxA	315.0 > 270.0	8.047	8.080	-0.033	1431721	50.5		101	4817	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.129	8.158	-0.029	1631	NC			9.9	
D 8 13C4-PFHpA	367.0 > 322.0	9.285	9.322	-0.037	1588868	51.5		103	7407	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.297	9.322	-0.025	1477	0.0446			5.8	
D 11 18O2 PFHxS	403.0 > 84.0	9.320	9.358	-0.038	690758	46.2		97.6	2532	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.320	9.364	-0.044	16976	NC			55.3	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.320	9.364	-0.044	16976	1.72				



Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA	417.0 > 372.0	10.407	10.450	-0.043		1722358	50.4	101	6977	
D 16 13C4 PFOS	503.0 > 80.0	11.379	11.414	-0.035		796221	43.8	91.6	2247	
15 Perfluorooctane sulfonic acid	499.0 > 80.0	11.365	11.417	-0.052	1.000	11575	0.7146		37.5	
	499.0 > 99.0	11.394	11.417	-0.023	1.003	6957		1.66(0.00-0.00)	21.9	
D 17 13C5 PFNA	468.0 > 423.0	11.395	11.434	-0.039		1454345	51.6	103	3476	
D 19 13C2 PFDA	515.0 > 470.0	12.238	12.275	-0.037		1396539	48.8	97.6	3223	
20 Perfluorodecanoic acid	513.0 > 469.0	12.228	12.276	-0.048	1.000	11670	0.0719		34.1	
D 23 13C8 FOSA	506.0 > 78.0	12.787	12.810	-0.023		425415	10.3	20.5	1512	
D 26 13C2 PFUnA	565.0 > 520.0	12.964	13.000	-0.036		1751587	46.9	93.7	4655	
27 Perfluoroundecanoic acid	563.0 > 519.0	12.954	13.000	-0.046	1.000	30855	0.3934		116	
D 28 13C2 PFDaA	615.0 > 570.0	13.581	13.609	-0.028		1880421	48.1	96.1	5575	
32 Perfluorotetradecanoic acid	713.0 > 669.0	14.548	14.571	-0.023	1.000	15510	0.2234		7.8	
D 33 13C2-PFTeDA	715.0 > 670.0	14.542	14.571	-0.029		1428562	39.8	79.6	5027	
D 35 13C2-PFHxDA	815.0 > 770.0	15.178	15.207	-0.029		1657995	38.5	76.9	4788	
34 Perfluorohexadecanoic acid	813.0 > 769.0	15.178	15.208	-0.030	1.000	134780	0.7895		194	
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.524	15.543	-0.019	1.000	1666	0.0399		4.1	

### QC Flag Legend

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_047.d

Injection Date: 20-Feb-2016 06:23:01

Instrument ID: A6

Lims ID: MB 320-100093/1-A

Client ID:

Operator ID: JRB

ALS Bottle#: 1

Worklist Smp#: 45

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

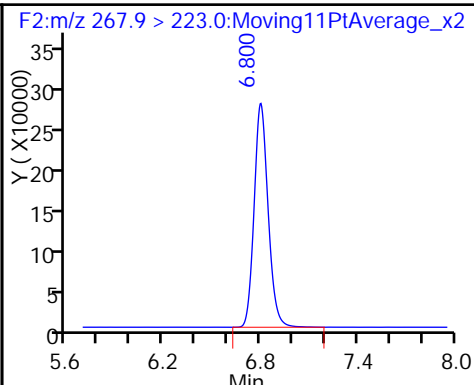
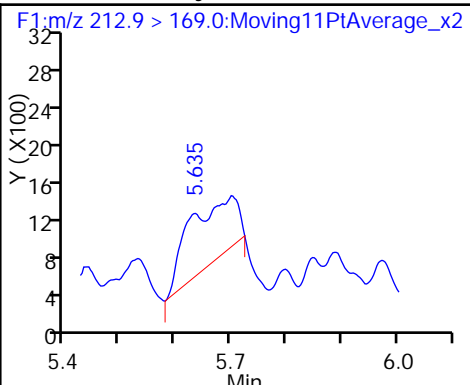
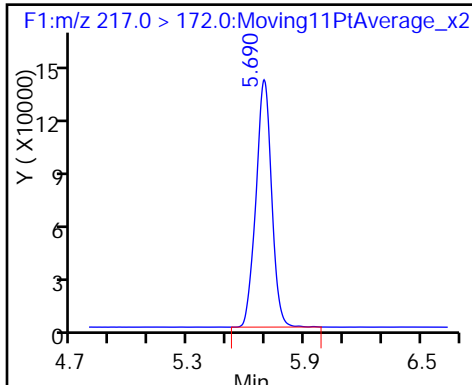
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

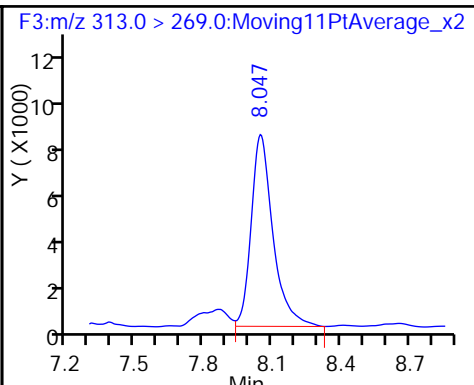
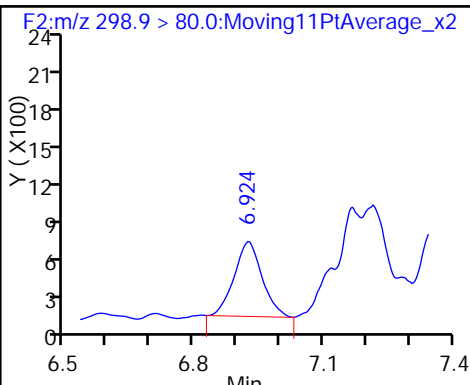
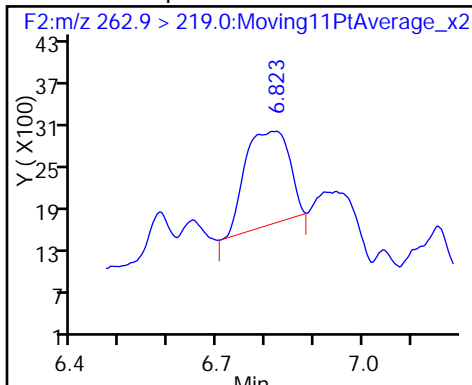
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

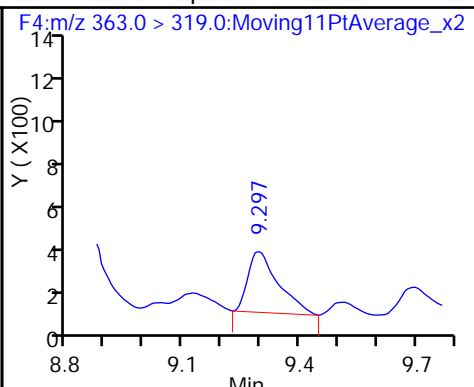
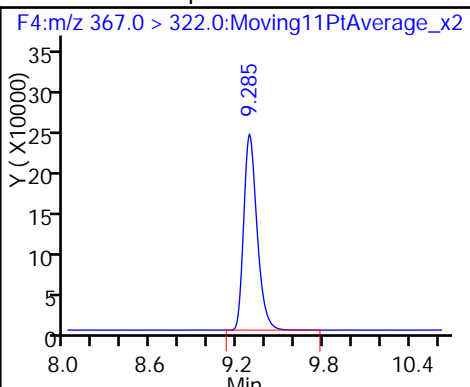
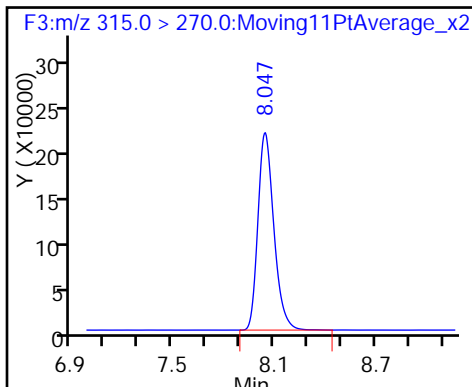
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

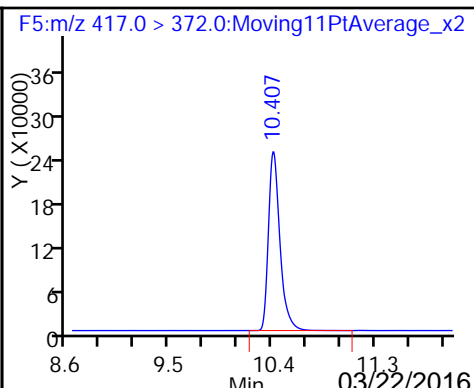
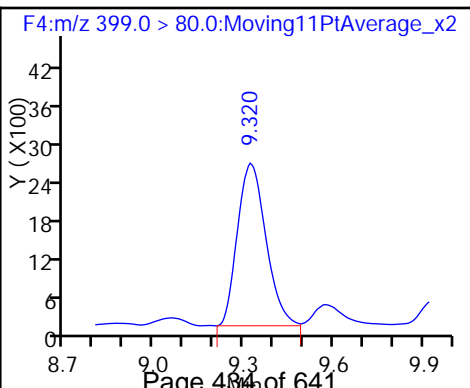
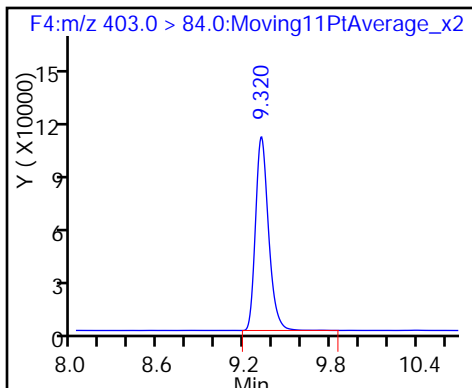
9 Perfluoroheptanoic acid

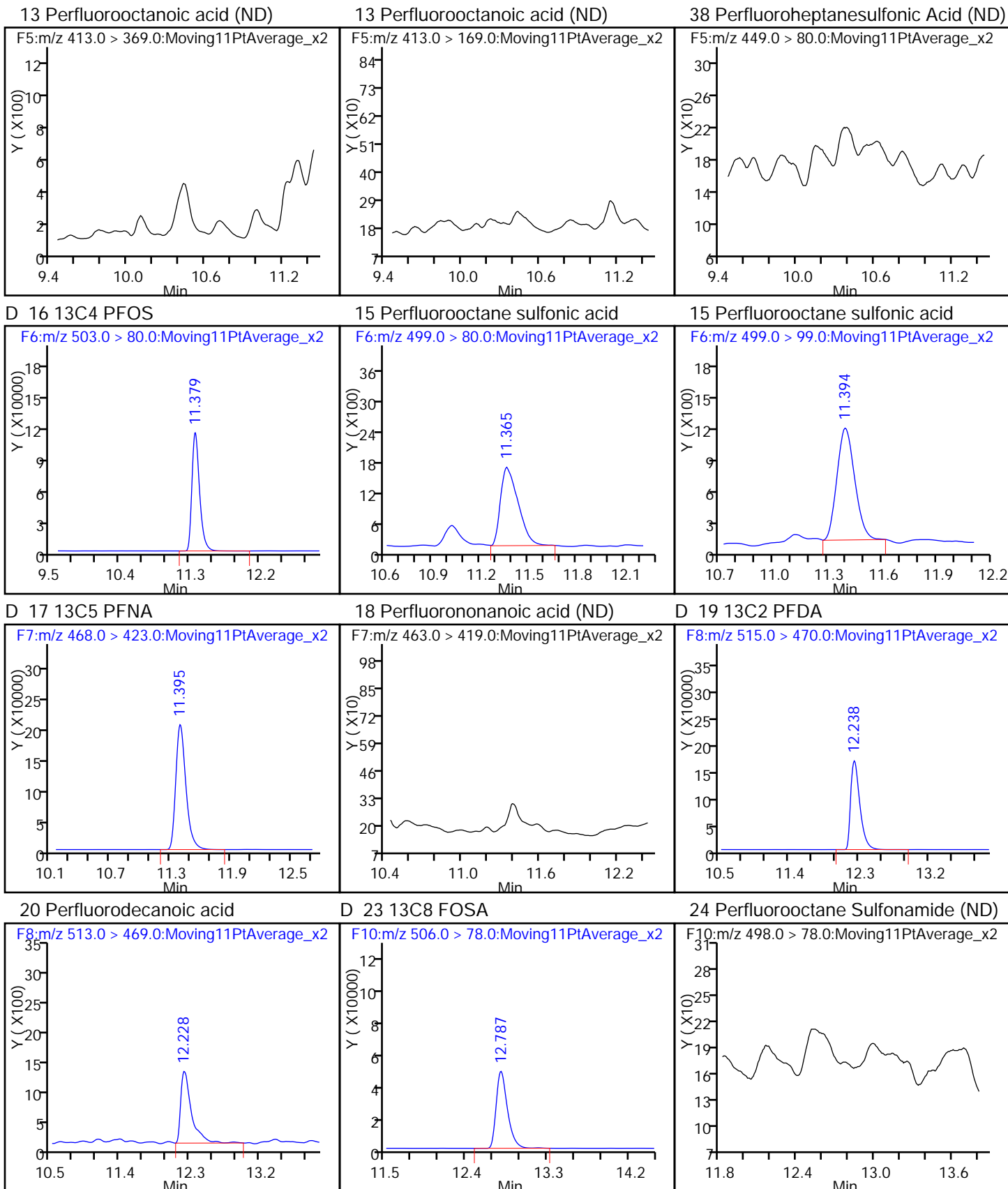


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

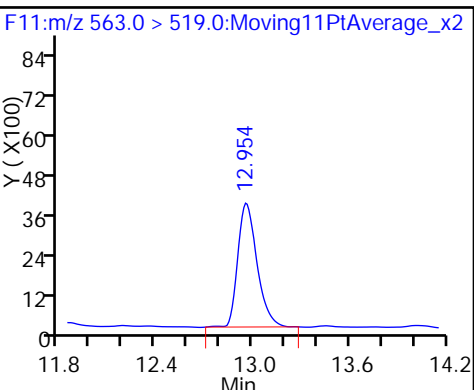
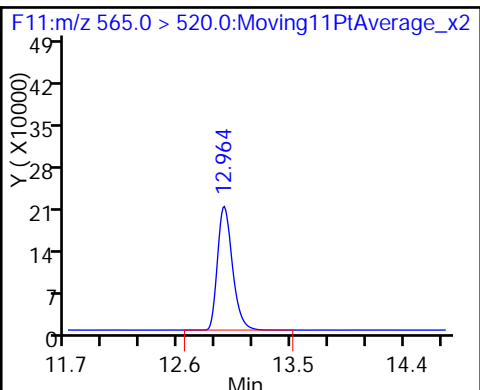
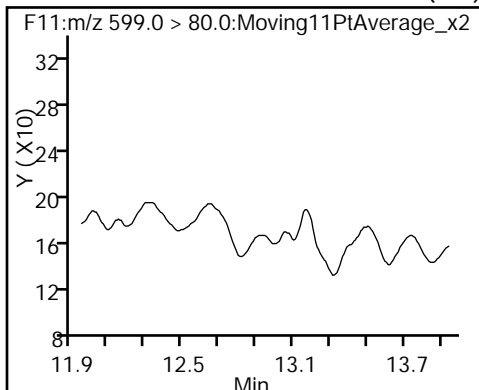
D 12 13C4 PFOA





39 Perfluorodecane Sulfonic acid (ND) D 26 13C2 PFUnA

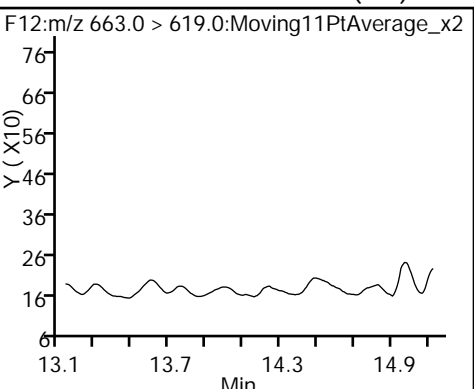
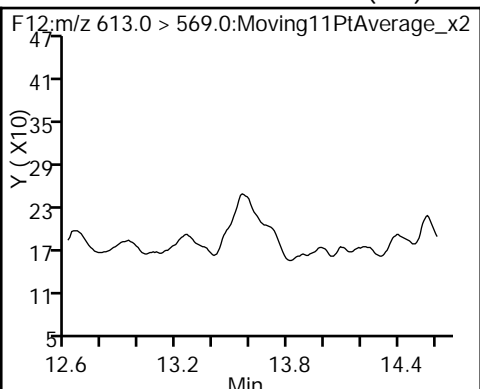
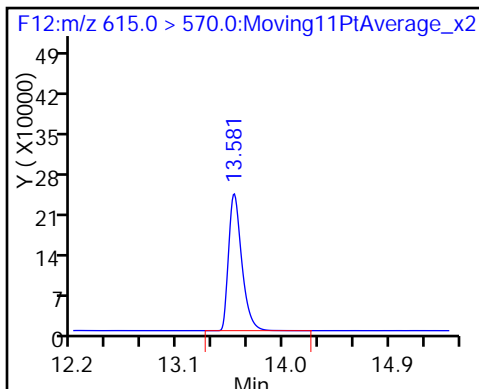
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid (ND)

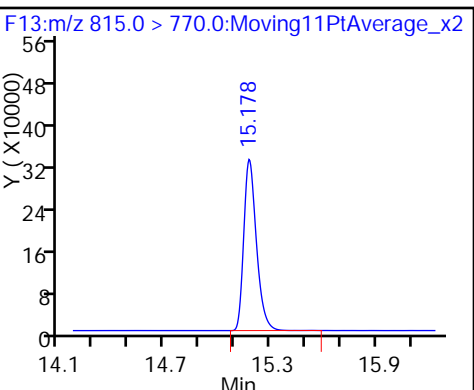
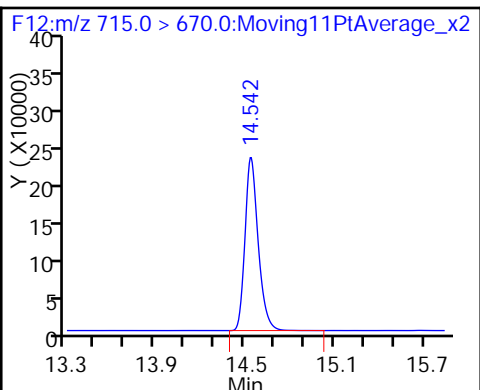
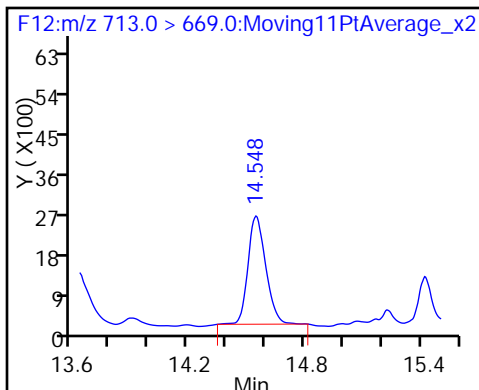
30 Perfluorotridecanoic acid (ND)



32 Perfluorotetradecanoic acid

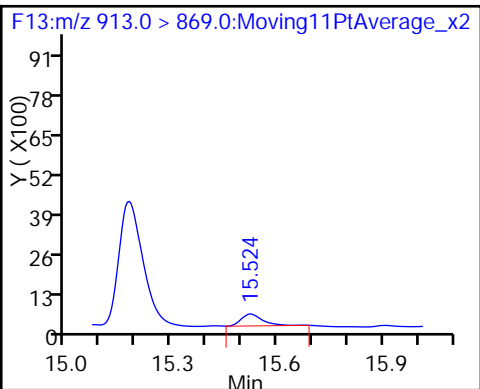
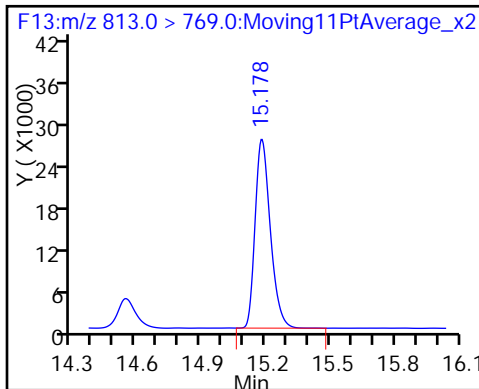
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 320-100147/1-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_061.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/10/2016 07:14  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 11:20  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0020	U	0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0020	U	0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0020	U	0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0020	U	0.0025	0.0020	0.00092
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0020	U	0.0025	0.0020	0.00087
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0030	U	0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	118		25-150
STL00990	13C4 PFOA	114		25-150
STL00995	13C5 PFNA	118		25-150
STL00994	18O2 PFHxS	97		25-150
STL00991	13C4 PFOS	98		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_061.d  
 Lims ID: MB 320-100147/1-A  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 20-Feb-2016 11:20:12 ALS Bottle#: 14 Worklist Smp#: 59  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: MB 320-100147/1-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:18:17

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.697	5.712	-0.015	822677	51.4		103	3372	
D 3 13C5-PFPeA	267.9 > 223.0	6.813	6.827	-0.014	1784267	58.0		116	6089	
4 Perfluoropentanoic acid	262.9 > 219.0	6.846	6.828	0.018	1205	0.0324			0.6	
5 Perfluorobutane Sulfonate	298.9 > 80.0	7.164	6.939	0.225	3607	NC			8.7	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	7.164	6.939	0.225	3607	0.2490				
7 Perfluorohexanoic acid	313.0 > 269.0	8.049	8.078	-0.029	7122	0.2144			28.7	
D 6 13C2 PFHxA	315.0 > 270.0	8.066	8.080	-0.014	1533227	54.1		108	4338	
D 8 13C4-PFHpA	367.0 > 322.0	9.299	9.322	-0.023	1824111	59.2		118	8633	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.288	9.322	-0.034	2422	0.0637			7.9	
D 11 18O2 PFHxS	403.0 > 84.0	9.335	9.358	-0.023	690169	46.1		97.5	2530	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.276	9.364	-0.088	300	NC			1.3	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.276	9.364	-0.088	300	0.0304				
D 12 13C4 PFOA	417.0 > 372.0	10.426	10.450	-0.024	1947965	57.1		114	6865	
D 16 13C4 PFOS	503.0 > 80.0	11.392	11.414	-0.022	852270	46.9		98.1	3706	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 17 13C5 PFNA	468.0 > 423.0	11.415	11.434	-0.019		1658130	58.8	118	6278	
D 19 13C2 PFDA	515.0 > 470.0	12.255	12.275	-0.020		1710616	59.7	119	4388	
20 Perfluorodecanoic acid	513.0 > 469.0	12.255	12.276	-0.021	1.000	12471	0.0172		44.1	
D 23 13C8 FOSA	506.0 > 78.0	12.804	12.810	-0.006		554432	13.4	26.7	1148	
D 26 13C2 PFUnA	565.0 > 520.0	12.980	13.000	-0.020		2089226	55.9	112	4731	
27 Perfluoroundecanoic acid	563.0 > 519.0	12.980	13.000	-0.020	1.000	25268	0.0582		60.4	
D 28 13C2 PFDoA	615.0 > 570.0	13.587	13.609	-0.022		2322857	59.4	119	5786	
32 Perfluorotetradecanoic acid	713.0 > 669.0	14.552	14.571	-0.019	1.000	16305	0.1205		6.6	
D 33 13C2-PFTeDA	715.0 > 670.0	14.552	14.571	-0.019		1865148	52.0	104	6265	
D 35 13C2-PFHxDA	815.0 > 770.0	15.182	15.207	-0.025		2228972	51.7	103	6252	
34 Perfluorohexadecanoic acid	813.0 > 769.0	15.182	15.208	-0.026	1.000	177273	1.01		251	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_061.d

Injection Date: 20-Feb-2016 11:20:12

Instrument ID: A6

Lims ID: MB 320-100147/1-A

Client ID:

Operator ID: JRB

ALS Bottle#: 14

Worklist Smp#: 59

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

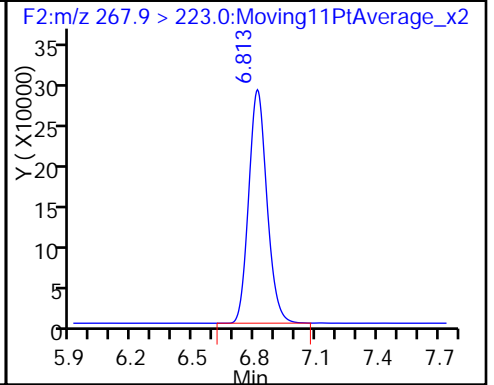
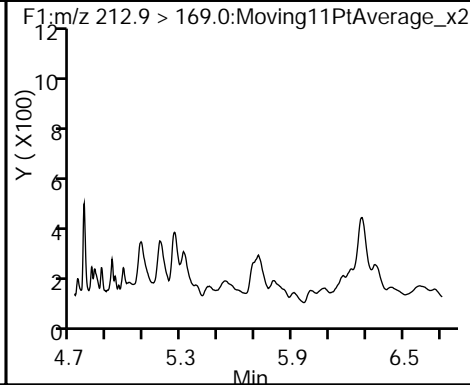
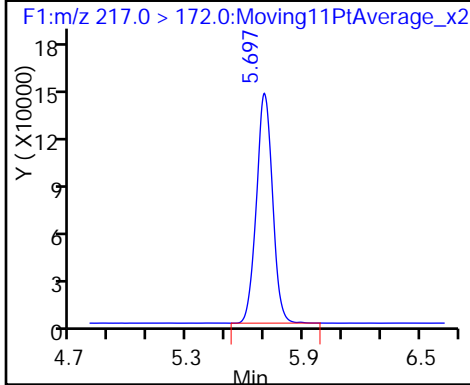
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid (ND)

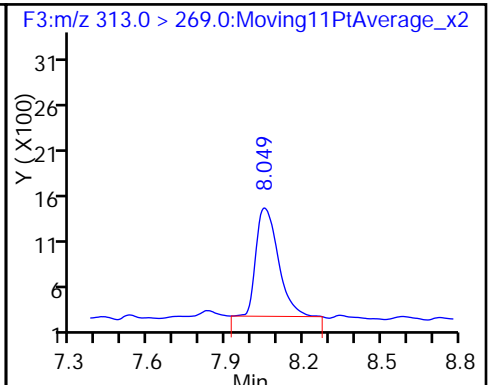
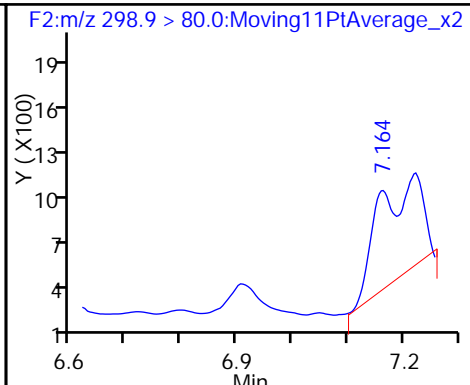
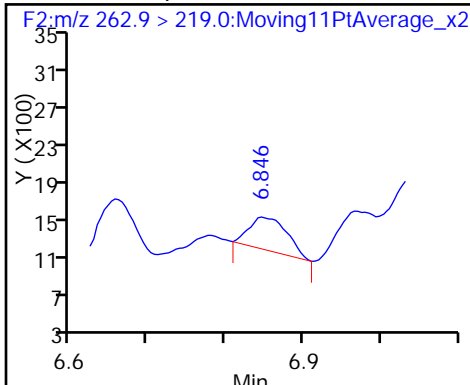
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

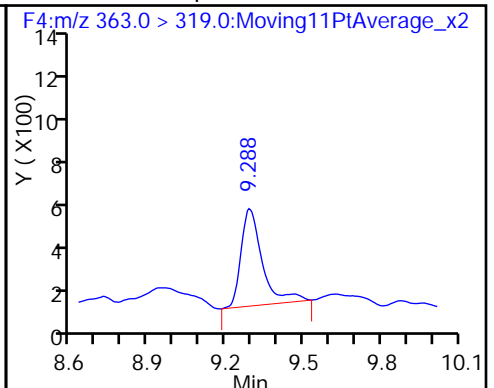
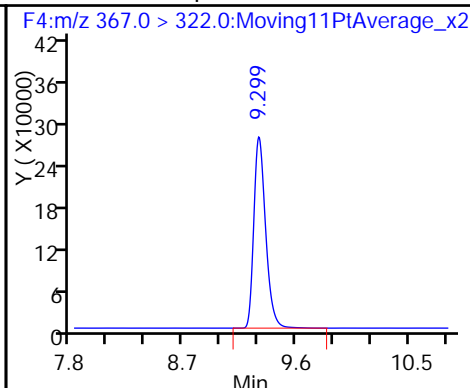
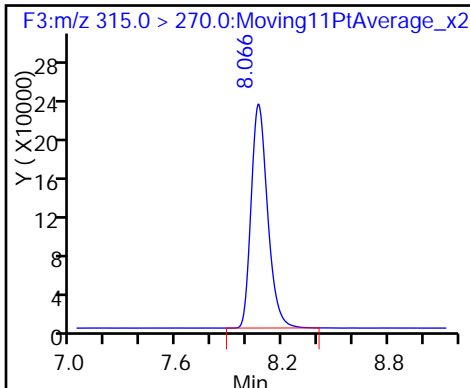
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

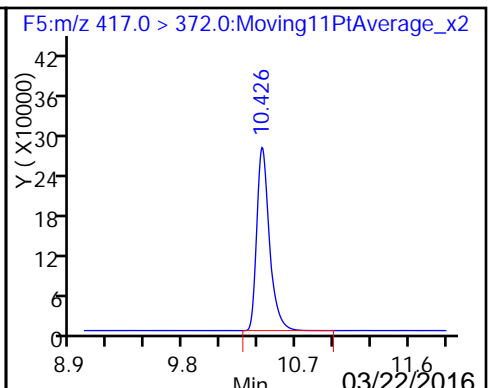
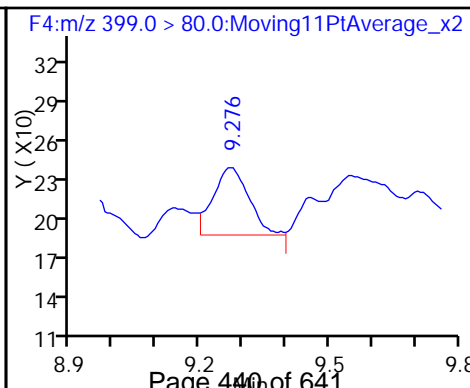
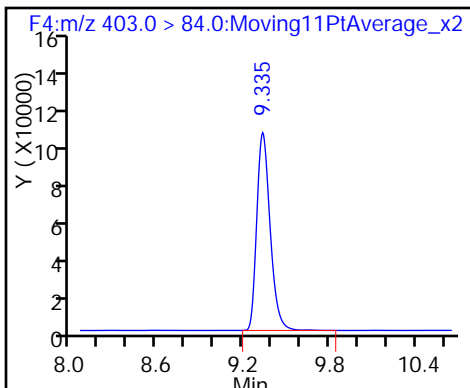
9 Perfluoroheptanoic acid



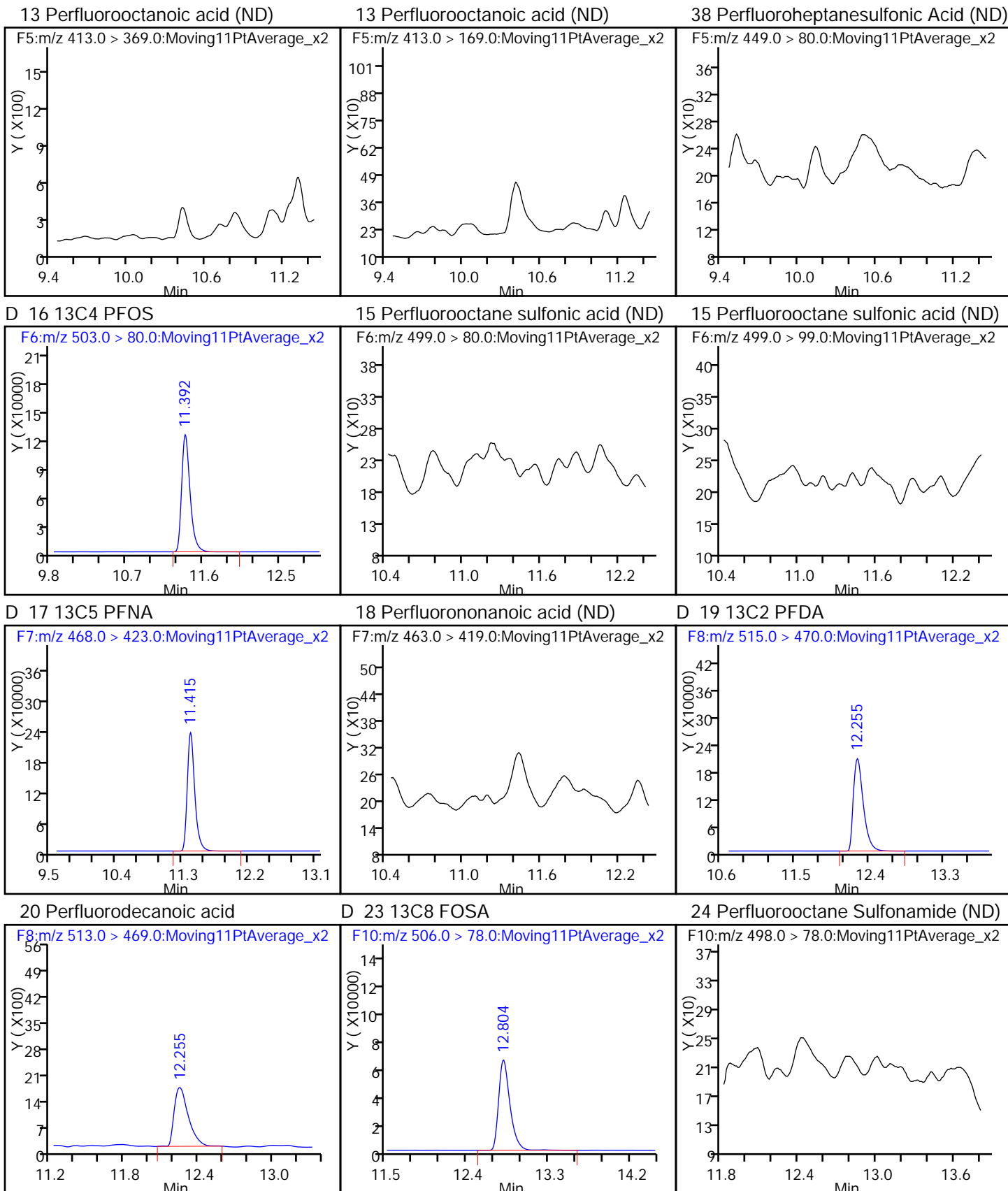
D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

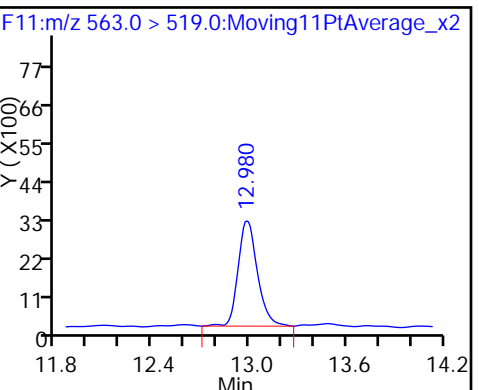
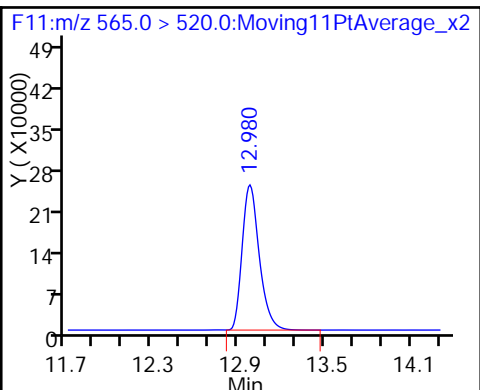
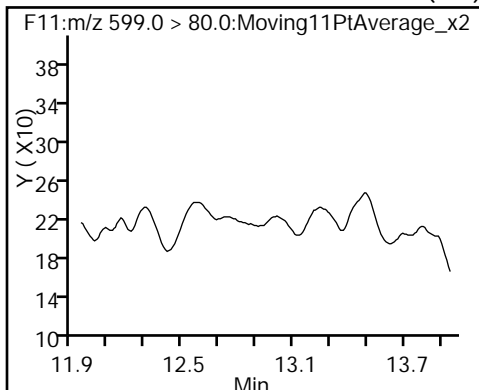






39 Perfluorodecane Sulfonic acid (ND) D 26 13C2 PFUnA

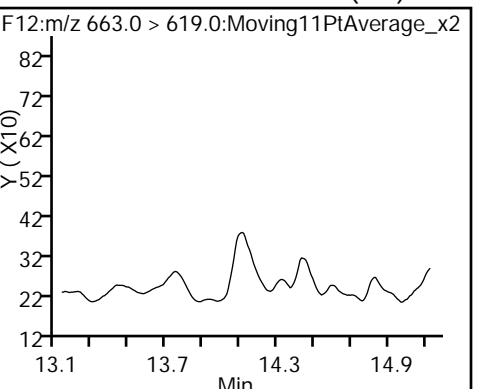
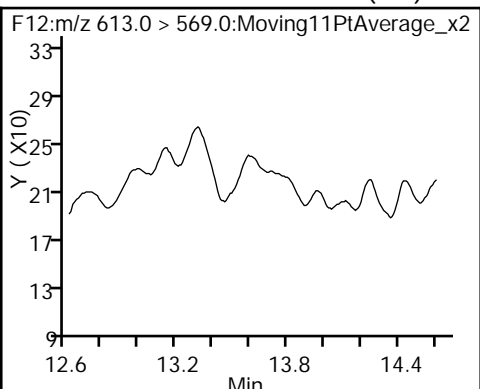
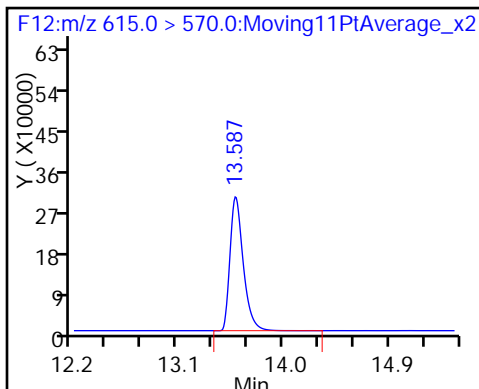
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid (ND)

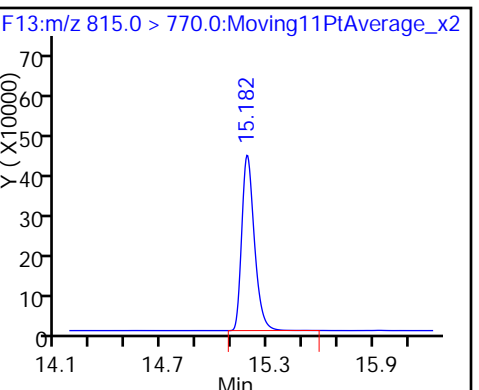
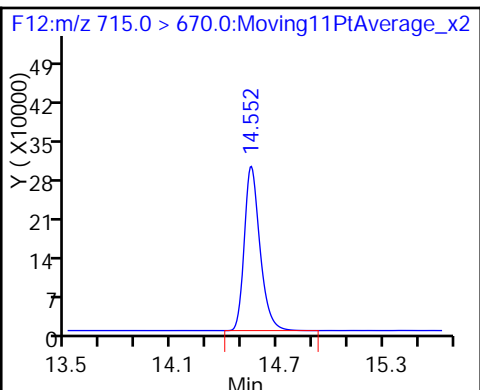
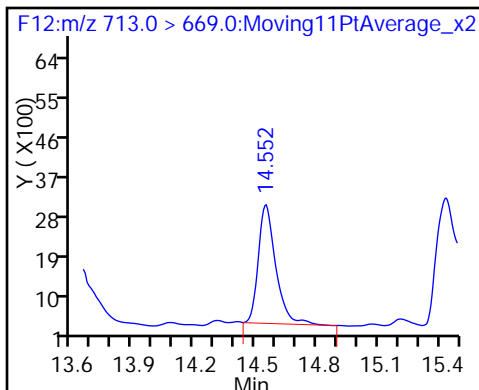
30 Perfluorotridecanoic acid (ND)



32 Perfluorotetradecanoic acid

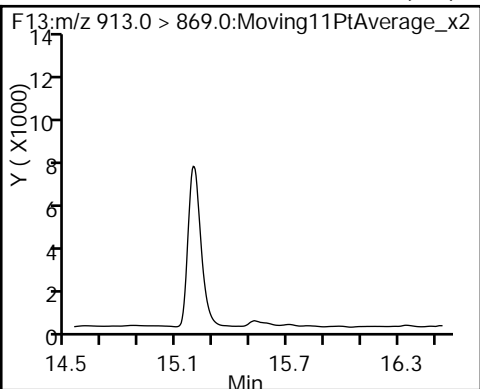
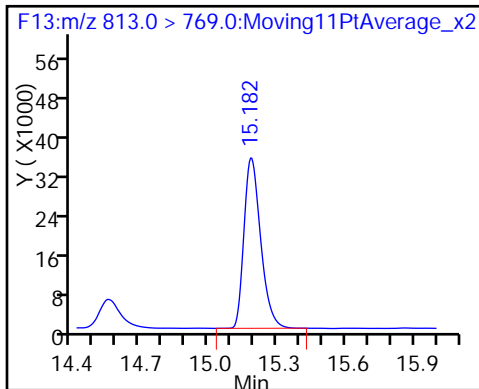
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid (ND)



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 320-100093/2-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_048.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/09/2016 10:23  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 06:44  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0513		0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0485		0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0537		0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0459		0.0025	0.0020	0.00092
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0480		0.0025	0.0020	0.00087
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0515		0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	95		25-150
STL00990	13C4 PFOA	95		25-150
STL00995	13C5 PFNA	98		25-150
STL00994	18O2 PFHxS	90		25-150
STL00991	13C4 PFOS	82		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_048.d  
 Lims ID: LCS 320-100093/2-A  
 Client ID:  
 Sample Type: LCS  
 Inject. Date: 20-Feb-2016 06:44:15 ALS Bottle#: 2 Worklist Smp#: 46  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCS 320-100093/2-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: westendorfc Date: 21-Feb-2016 15:38:20

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 13C4 PFBA	217.0 > 172.0	5.697	5.712	-0.015	758399	47.4		94.8	3128	
2 Perfluorobutyric acid	212.9 > 169.0	5.697	5.715	-0.018	480442	24.3		121	892	
D 3 13C5-PFPeA	267.9 > 223.0	6.805	6.827	-0.022	1590400	51.7		103	5164	
4 Perfluoropentanoic acid	262.9 > 219.0	6.809	6.828	-0.019	770505	23.2		116	224	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.924	6.939	-0.015	306961	NC			348	
	298.9 > 99.0	6.924	6.939	-0.015	166428		1.84(0.00-0.00)		290	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.924	6.939	-0.015	306961	22.9		130		
7 Perfluorohexanoic acid	313.0 > 269.0	8.055	8.078	-0.023	727454	25.3		127	1385	
D 6 13C2 PFHxA	315.0 > 270.0	8.055	8.080	-0.025	1326503	46.8		93.7	4748	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.131	8.158	-0.027	239493	NC			932	
D 8 13C4-PFHpA	367.0 > 322.0	9.288	9.322	-0.034	1471383	47.7		95.5	4291	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.294	9.322	-0.028	787432	25.7		128	3152	
D 11 18O2 PFHxS	403.0 > 84.0	9.323	9.358	-0.035	637656	42.6		90.1	3184	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.323	9.364	-0.041	218573	NC			829	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.323	9.364	-0.041	218573	24.0		127		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.419	10.450	-0.031		1613824	47.3		94.5	6332	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.419	10.450	-0.031	1.000	785558	24.2		121	1043	
413.0 > 169.0	10.419	10.450	-0.031	1.000	236834		3.32(0.00-0.00)		760	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.419	10.457	-0.038	1.000	252668	NC			1450	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.419	10.457	-0.038	1.000	252668	27.4		144		
D 16 13C4 PFOS										
503.0 > 80.0	11.385	11.414	-0.029		711939	39.2		81.9	3408	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.385	11.417	-0.032	1.000	373063	25.8		135	880	
499.0 > 99.0	11.385	11.417	-0.032	1.000	211135		1.77(0.00-0.00)		669	
D 17 13C5 PFNA										
468.0 > 423.0	11.401	11.434	-0.033		1374538	48.8		97.5	6797	
18 Perfluorononanoic acid										
463.0 > 419.0	11.408	11.439	-0.031	1.000	596415	26.9		134	1966	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.203	12.249	-0.046	1.000	212396	NC			888	
D 19 13C2 PFDA										
515.0 > 470.0	12.245	12.275	-0.030		1265366	44.2		88.4	5404	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.245	12.276	-0.031	1.000	649916	26.0		130	2595	
D 23 13C8 FOSA										
506.0 > 78.0	12.793	12.810	-0.017		165922	4.00		8.0	445	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.793	12.812	-0.019	1.000	58737	19.1		95.5	196	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.919	12.949	-0.030	1.000	211472	NC			820	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.919	12.949	-0.030	1.000	211472	25.8		134		
D 26 13C2 PFUnA										
565.0 > 520.0	12.970	13.000	-0.030		1646373	44.0		88.1	4319	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.970	13.000	-0.030	1.000	708085	25.4		127	2414	
D 28 13C2 PFDaA										
615.0 > 570.0	13.578	13.609	-0.031		1801826	46.1		92.1	5300	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.578	13.611	-0.033	1.000	663674	25.2		126	894	
31 PFDaS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.048	14.083	-0.035	1.000	215474	NC			709	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.101	14.130	-0.029	1.000	859635	25.4		127	1189	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.546	14.571	-0.025	1.000	471466	21.5		107	292	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.539	14.571	-0.032		1400557	39.0		78.0	3533	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA										
815.0 > 770.0	15.183	15.207	-0.024		1666171	38.7		77.3	4296	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.183	15.208	-0.025	1.000	958744	22.3		111	1477	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.514	15.543	-0.029	1.000	873096	21.8		109	1349	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_048.d

Injection Date: 20-Feb-2016 06:44:15

Instrument ID: A6

Lims ID: LCS 320-100093/2-A

Client ID:

Operator ID: JRB

ALS Bottle#: 2

Worklist Smp#: 46

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

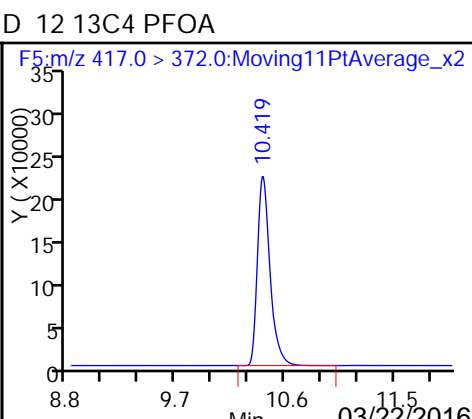
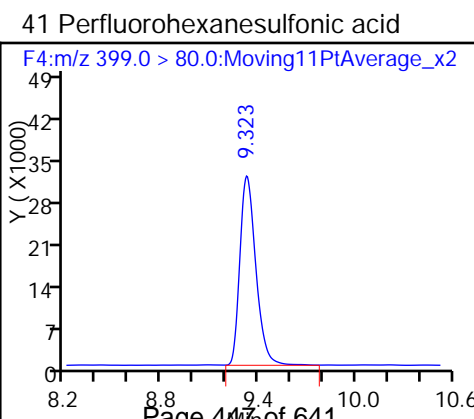
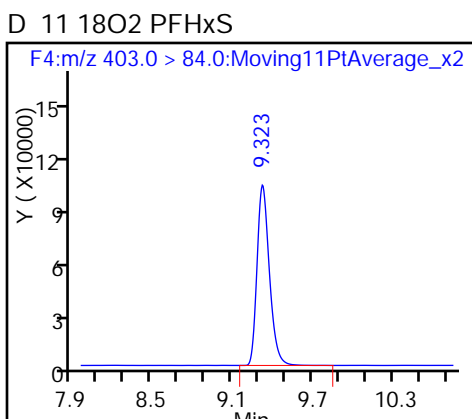
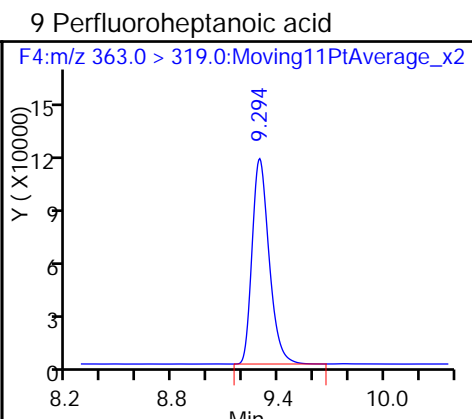
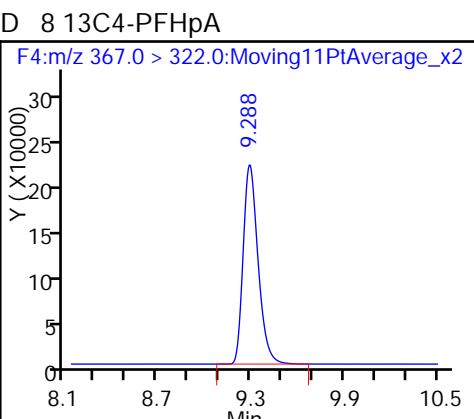
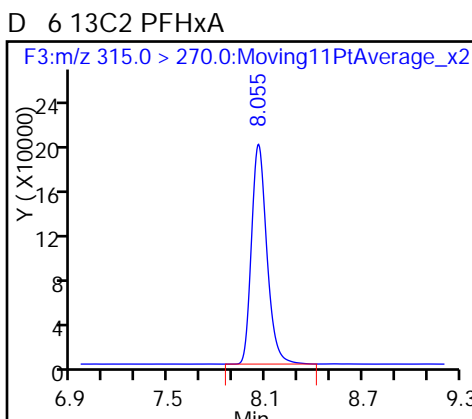
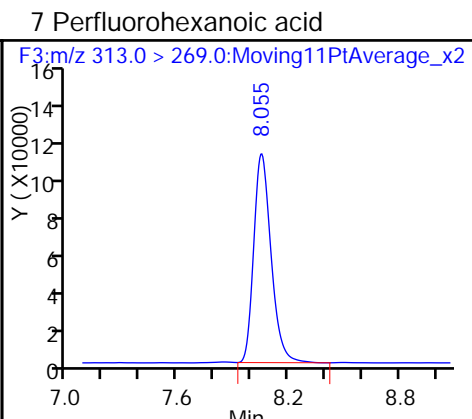
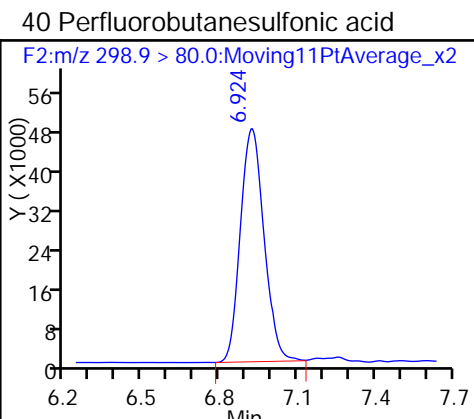
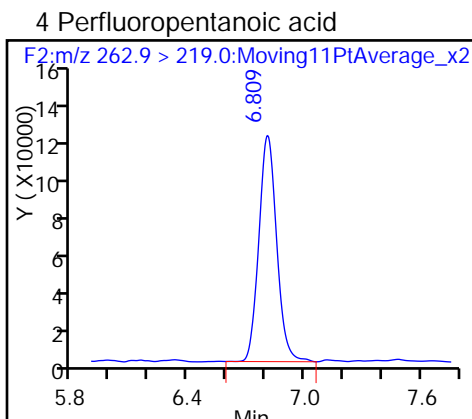
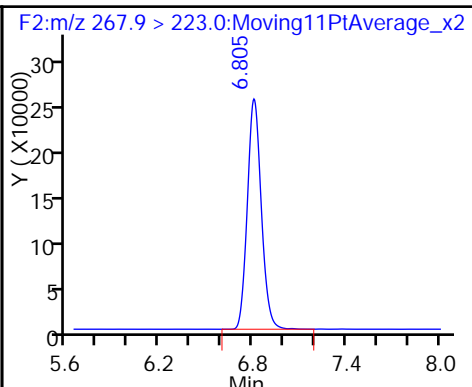
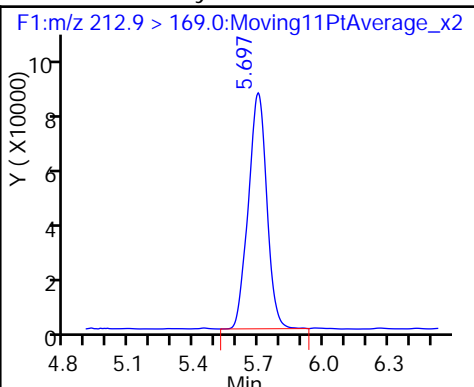
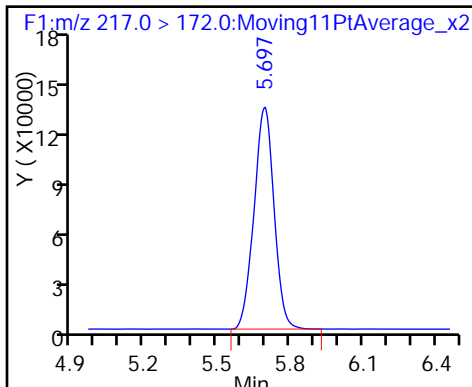
Method: PFAC\_A6

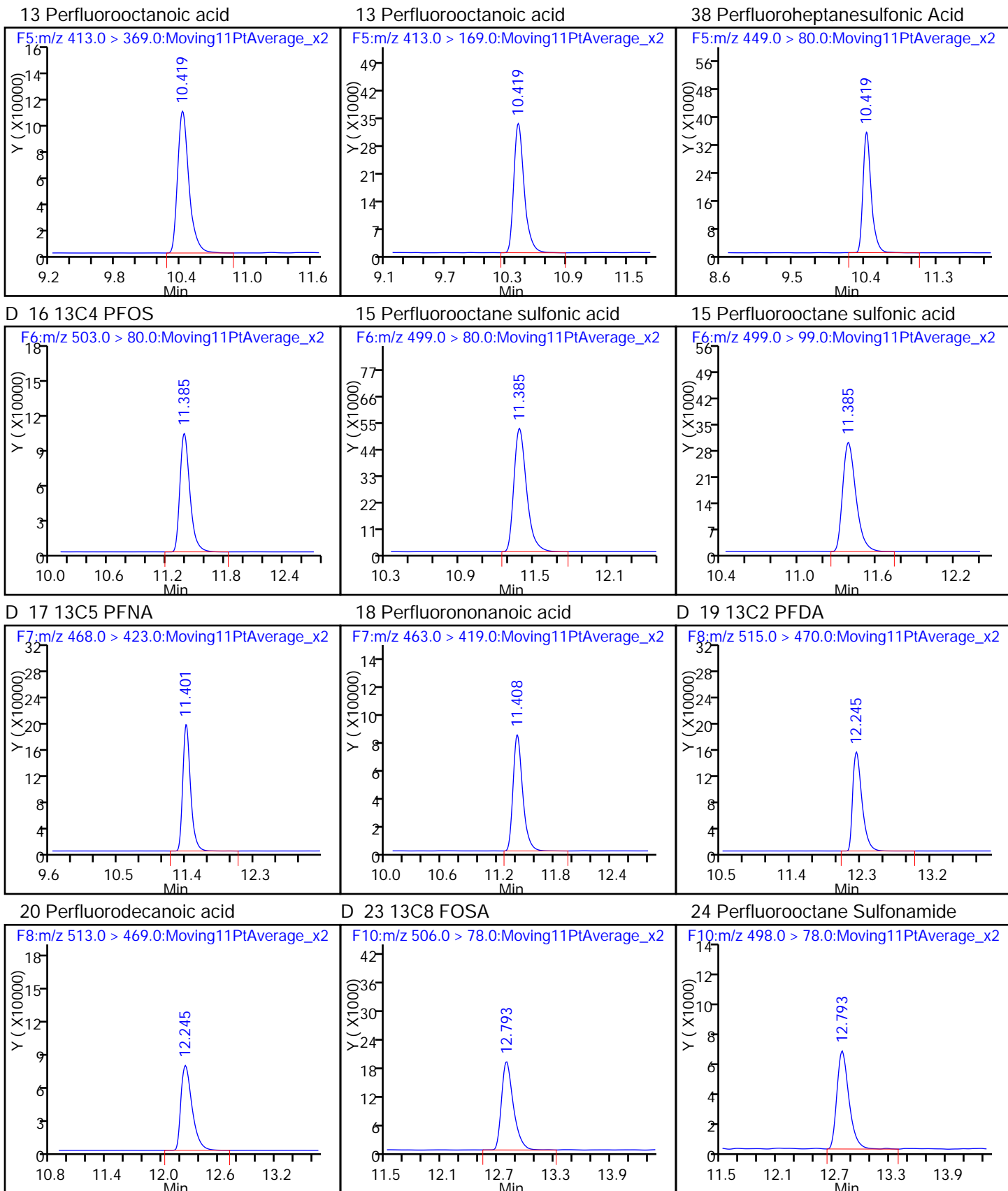
Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

D 3 13C5-PFPeA



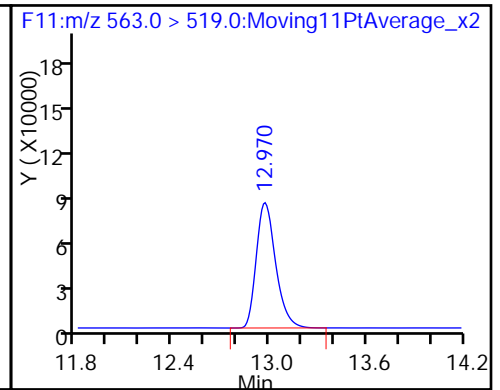
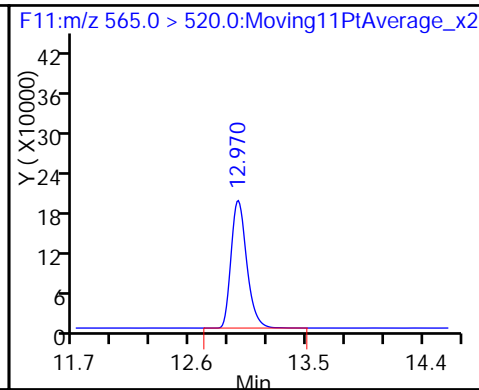
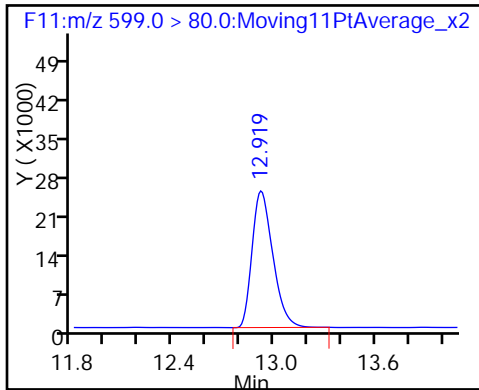




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

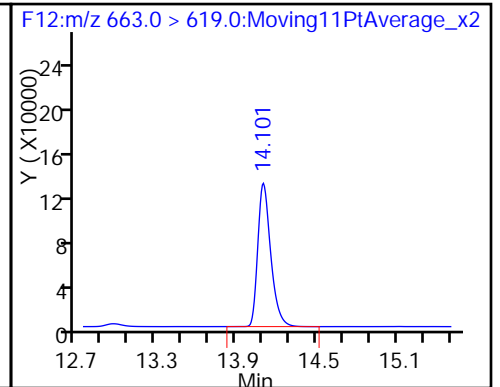
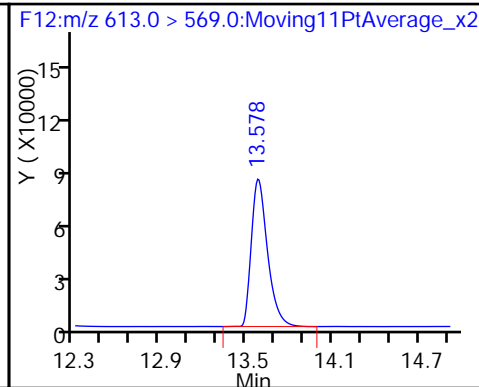
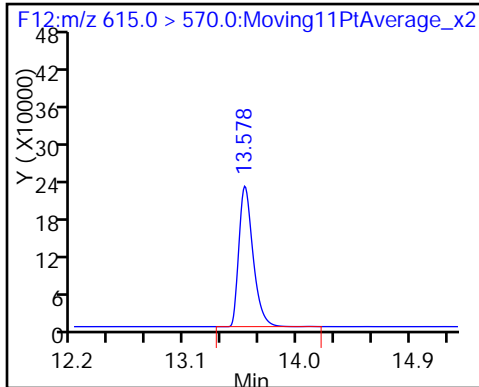
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

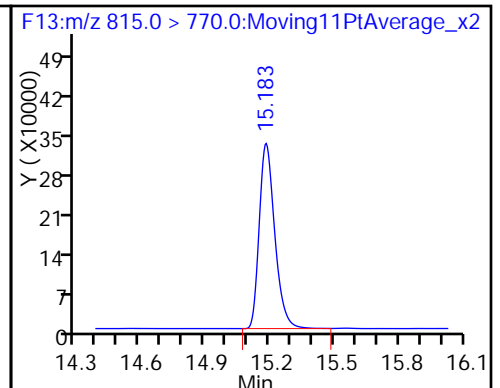
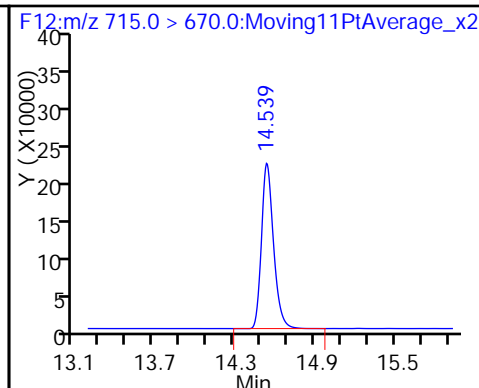
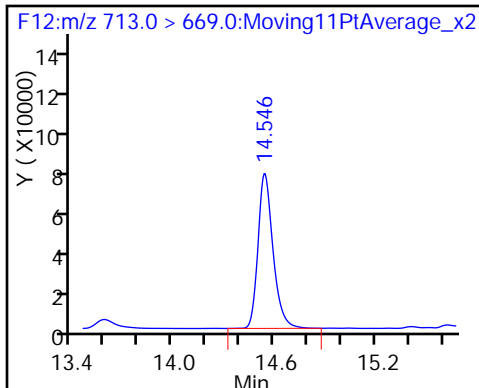
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

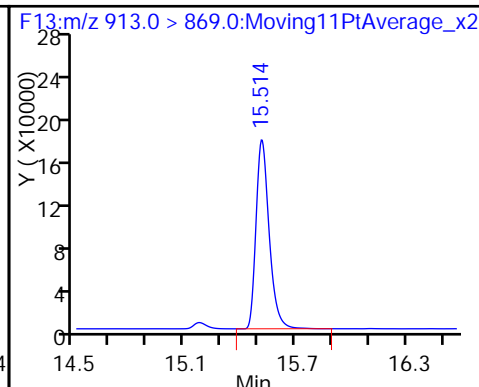
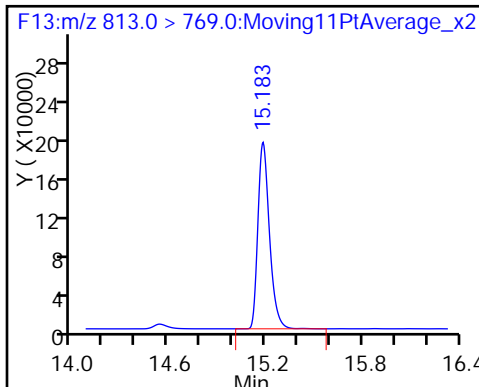
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 320-100147/2-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_062.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/10/2016 07:14  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 11:41  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0498		0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0476		0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0528		0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0412		0.0025	0.0020	0.00092
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0503		0.0025	0.0020	0.00087
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0529		0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	124		25-150
STL00990	13C4 PFOA	119		25-150
STL00995	13C5 PFNA	117		25-150
STL00994	18O2 PFHxS	108		25-150
STL00991	13C4 PFOS	99		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_062.d  
 Lims ID: LCS 320-100147/2-A  
 Client ID:  
 Sample Type: LCS  
 Inject. Date: 20-Feb-2016 11:41:26 ALS Bottle#: 15 Worklist Smp#: 60  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCS 320-100147/2-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:18:47

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.703	5.712	-0.009	865377	54.1		108	2525	
2 Perfluorobutyric acid	212.9 > 169.0	5.703	5.715	-0.012	588390	26.1		130	1526	
D 3 13C5-PFPeA	267.9 > 223.0	6.814	6.827	-0.013	1909548	62.1		124	5750	
4 Perfluoropentanoic acid	262.9 > 219.0	6.818	6.828	-0.010	950555	23.9		119	223	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.933	6.939	-0.006	329085	NC			329	
	298.9 > 99.0	6.929	6.939	-0.010	199360		1.65(0.00-0.00)		387	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.933	6.939	-0.006	329085	20.6		116		
7 Perfluorohexanoic acid	313.0 > 269.0	8.066	8.078	-0.012	916014	25.8		129	1757	
D 6 13C2 PFHxA	315.0 > 270.0	8.060	8.080	-0.020	1641733	58.0		116	6475	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.142	8.158	-0.016	291567	NC			989	
D 8 13C4-PFHpA	367.0 > 322.0	9.299	9.322	-0.023	1907376	61.9		124	5722	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.305	9.322	-0.017	989698	24.9		124	2636	
D 11 18O2 PFHxS	403.0 > 84.0	9.334	9.358	-0.024	761104	50.9		108	2298	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.340	9.364	-0.024	273522	NC			819	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.340	9.364	-0.024	273522	25.2		133		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.433	10.450	-0.017		2039439	59.7		119	6590	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.426	10.450	-0.024	1.000	974127	23.8		119	1129	
413.0 > 169.0	10.433	10.450	-0.017	1.001	312616		3.12(0.00-0.00)		711	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.433	10.457	-0.024	1.000	273837	NC			688	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.433	10.457	-0.024	1.000	273837	24.5		129		
D 16 13C4 PFOS										
503.0 > 80.0	11.392	11.414	-0.022		862175	47.4		99.2	2828	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.399	11.417	-0.018	1.000	463522	26.4		138	465	
499.0 > 99.0	11.392	11.417	-0.025	0.999	228306		2.03(0.00-0.00)		854	
D 17 13C5 PFNA										
468.0 > 423.0	11.415	11.434	-0.019		1649054	58.5		117	5820	
18 Perfluorononanoic acid										
463.0 > 419.0	11.415	11.439	-0.024	1.000	702858	26.4		132	1655	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.224	12.249	-0.025	1.000	212899	NC			941	
D 19 13C2 PFDA										
515.0 > 470.0	12.255	12.275	-0.020		1652251	57.7		115	2788	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.255	12.276	-0.021	1.000	769368	23.5		118	2247	
D 23 13C8 FOSA										
506.0 > 78.0	12.804	12.810	-0.006		1559046	37.6		75.1	2280	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.804	12.812	-0.008	1.000	710550	24.6		123	1893	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.929	12.949	-0.020	1.000	247870	NC			747	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.929	12.949	-0.020	1.000	247870	24.9		129		
D 26 13C2 PFUnA										
565.0 > 520.0	12.981	13.000	-0.019		2104757	56.3		113	6044	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.981	13.000	-0.019	1.000	876111	24.6		123	2372	
D 28 13C2 PFDaA										
615.0 > 570.0	13.596	13.609	-0.013		2271652	58.1		116	6437	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.596	13.611	-0.015	1.000	909345	27.4		137	1005	
31 PFDoS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.055	14.083	-0.028	1.000	290830	NC			998	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.108	14.130	-0.022	1.000	1123125	26.3		131	1020	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.552	14.571	-0.019	1.000	647031	23.4		117	223	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.552	14.571	-0.019		1962230	54.7		109	4945	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA	815.0 > 770.0	15.188	15.207	-0.019		2508425	58.2		116	5146
34 Perfluorohexadecanoic acid	813.0 > 769.0	15.188	15.208	-0.020	1.000	1372099	25.6		128	1589
36 Perfluorooctadecanoic acid	913.0 > 869.0	15.519	15.543	-0.024	1.000	1366360	27.1		135	1358

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_062.d

Injection Date: 20-Feb-2016 11:41:26

Instrument ID: A6

Lims ID: LCS 320-100147/2-A

Client ID:

Operator ID: JRB

ALS Bottle#: 15

Worklist Smp#: 60

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

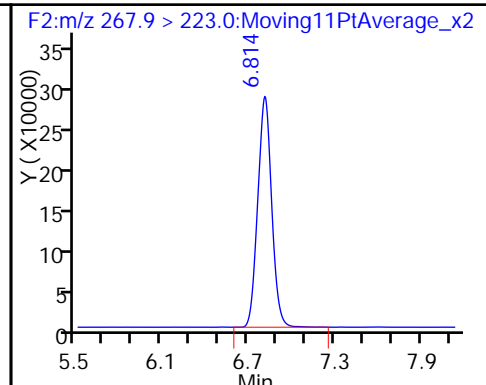
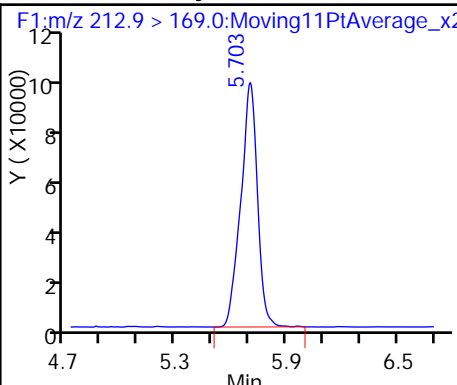
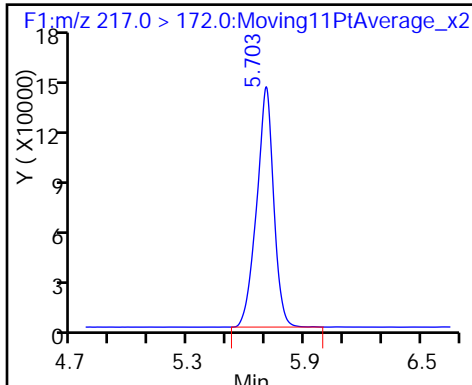
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

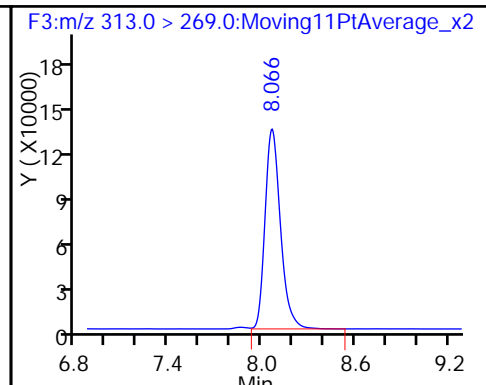
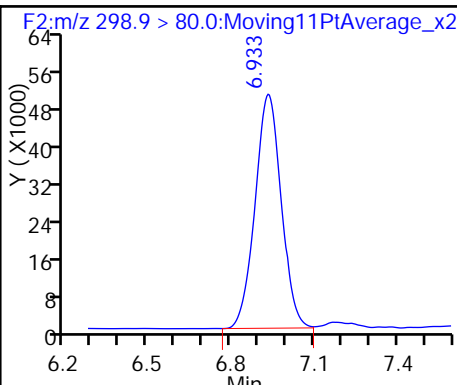
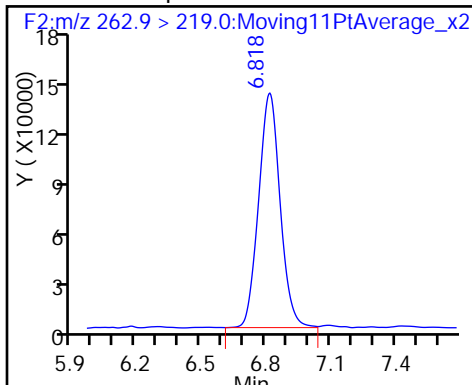
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

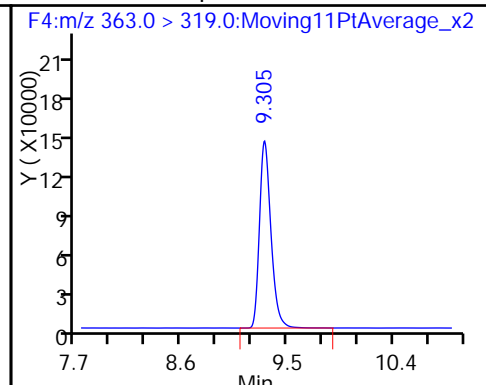
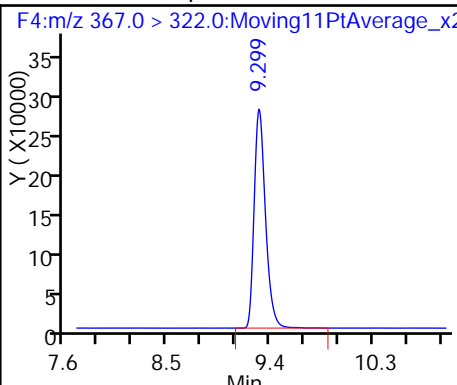
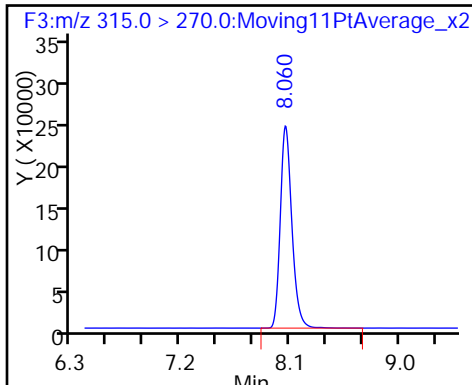
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

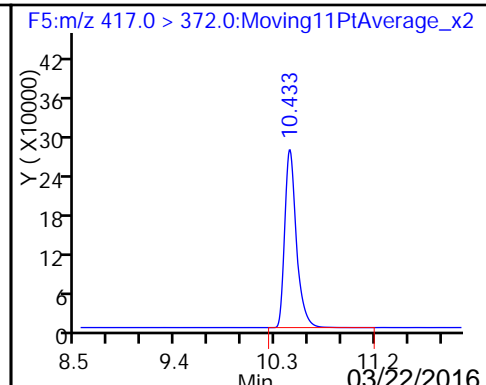
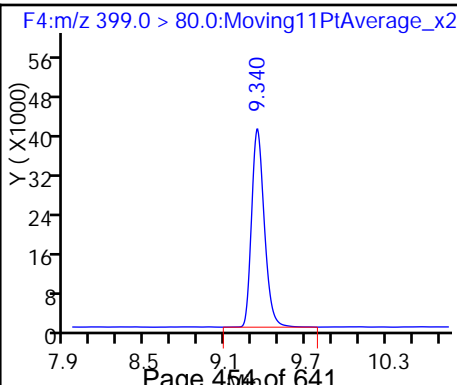
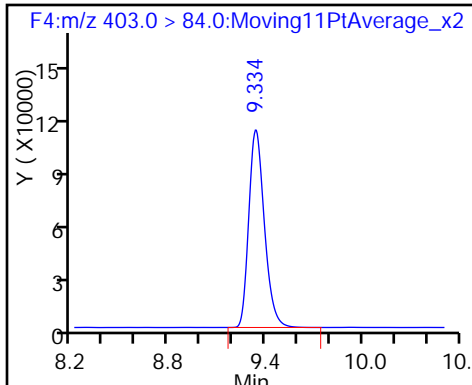
9 Perfluoroheptanoic acid

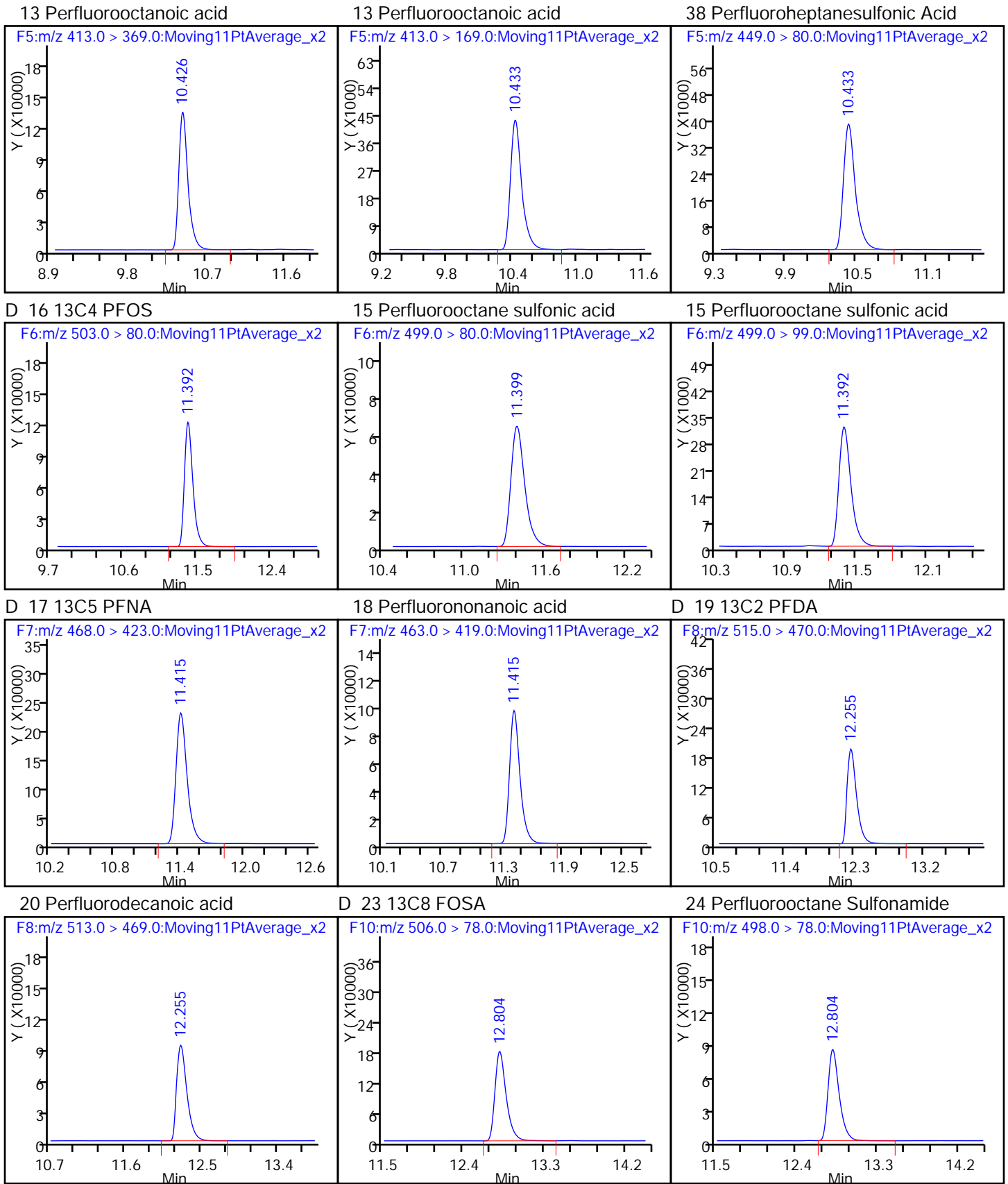


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

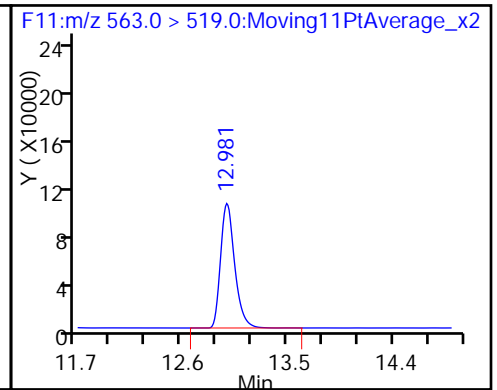
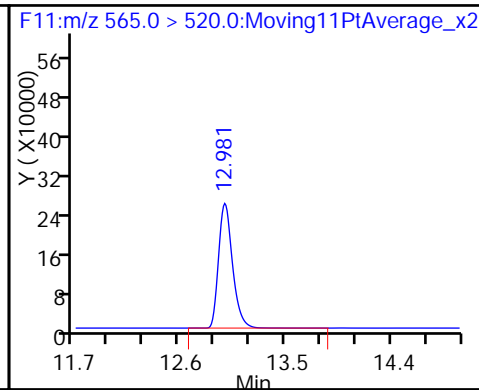
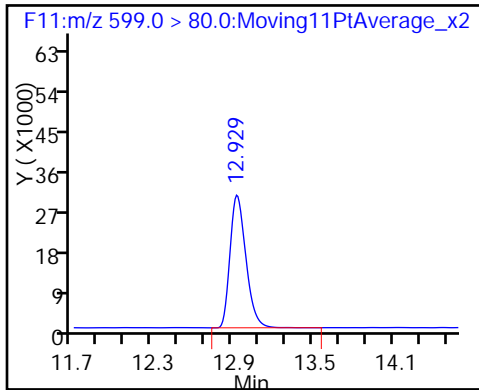




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

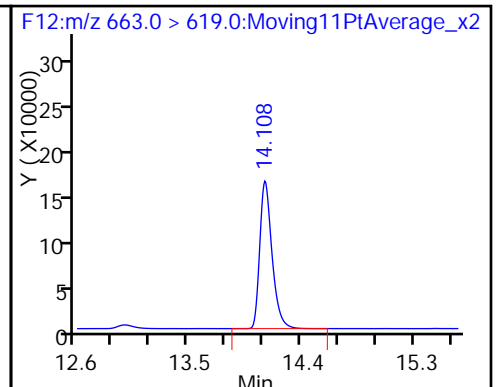
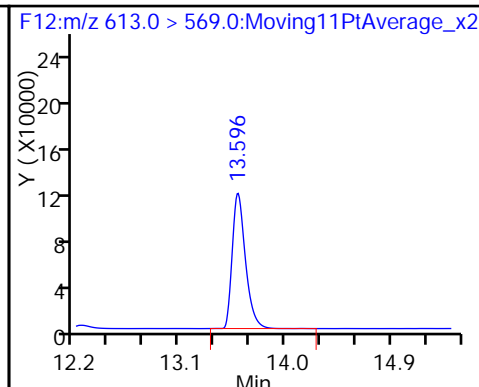
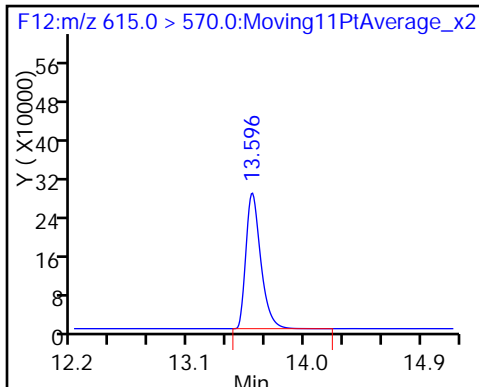
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

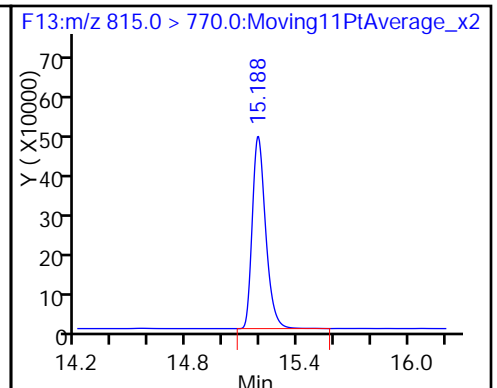
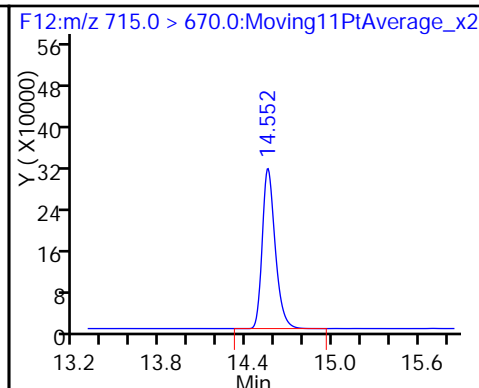
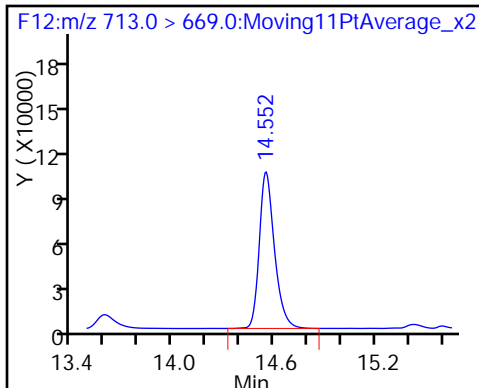
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

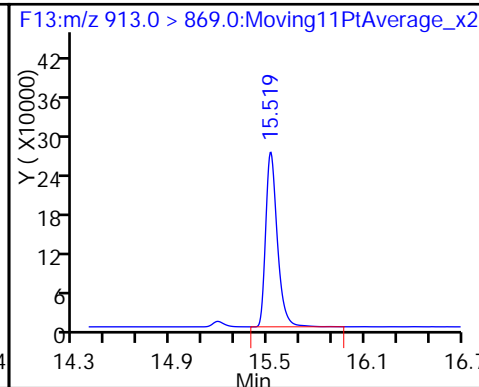
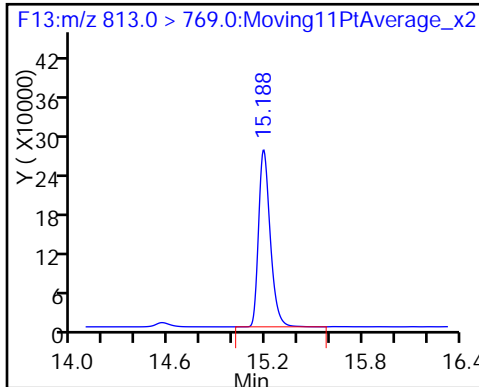
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid





FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCSD 320-100093/3-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_049.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/09/2016 10:23  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 07:05  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0506		0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0493		0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0517		0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0472		0.0025	0.0020	0.00092
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0486		0.0025	0.0020	0.00087
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0452		0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	95		25-150
STL00990	13C4 PFOA	88		25-150
STL00995	13C5 PFNA	91		25-150
STL00994	18O2 PFHxS	86		25-150
STL00991	13C4 PFOS	80		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_049.d  
 Lims ID: LCSD 320-100093/3-A  
 Client ID:  
 Sample Type: LCSD  
 Inject. Date: 20-Feb-2016 07:05:28 ALS Bottle#: 3 Worklist Smp#: 47  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCSD 320-100093/3-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 10:48:20 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 20-Feb-2016 13:11:09

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.699	5.712	-0.013	734641	45.9		91.8	2704	
2 Perfluorobutyric acid	212.9 > 169.0	5.696	5.715	-0.019	485146	25.3		127	1247	
D 3 13C5-PFPeA	267.9 > 223.0	6.803	6.827	-0.024	1540503	50.1		100	5533	
4 Perfluoropentanoic acid	262.9 > 219.0	6.803	6.828	-0.025	722067	22.5		112	178	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.918	6.939	-0.021	300270	NC			379	
	298.9 > 99.0	6.923	6.939	-0.016	149697		2.01(0.00-0.00)		240	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.918	6.939	-0.021	300270	23.6		133		
7 Perfluorohexanoic acid	313.0 > 269.0	8.050	8.078	-0.028	733184	25.9		130	1978	
D 6 13C2 PFHxA	315.0 > 270.0	8.050	8.080	-0.030	1305099	46.1		92.1	6340	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.126	8.158	-0.032	224529	NC			920	
D 8 13C4-PFHpA	367.0 > 322.0	9.288	9.322	-0.034	1471294	47.7		95.5	5622	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.288	9.322	-0.034	776183	25.3		126	2978	
D 11 18O2 PFHxS	403.0 > 84.0	9.318	9.358	-0.040	606231	40.5		85.6	2633	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.323	9.364	-0.041	210408	NC			687	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.323	9.364	-0.041	210408	24.3		128		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.413	10.450	-0.037		1501627	44.0		88.0	5693	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.413	10.450	-0.037	1.000	743776	24.7		123	1424	
413.0 > 169.0	10.420	10.450	-0.030	1.001	224572		3.31(0.00-0.00)		836	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.420	10.457	-0.037	1.000	220037	NC			800	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.420	10.457	-0.037	1.000	220037	24.6		129		
D 16 13C4 PFOS										
503.0 > 80.0	11.378	11.414	-0.036		691703	38.0		79.6	3976	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.385	11.417	-0.032	1.000	318089	22.6		118	1062	
499.0 > 99.0	11.385	11.417	-0.032	1.000	192858		1.65(0.00-0.00)		754	
D 17 13C5 PFNA										
468.0 > 423.0	11.401	11.434	-0.033		1281775	45.5		90.9	3447	
18 Perfluorononanoic acid										
463.0 > 419.0	11.401	11.439	-0.038	1.000	534851	25.8		129	1565	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.207	12.249	-0.042	1.000	199232	NC			591	
D 19 13C2 PFDA										
515.0 > 470.0	12.238	12.275	-0.037		1277810	44.6		89.3	4187	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.238	12.276	-0.038	1.000	617874	24.4		122	2237	
D 23 13C8 FOSA										
506.0 > 78.0	12.787	12.810	-0.023		187222	4.51		9.0	565	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.787	12.812	-0.025	1.000	75363	21.7		109	298	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.923	12.949	-0.026	1.000	208463	NC			652	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.923	12.949	-0.026	1.000	208463	26.1		136		
D 26 13C2 PFUnA										
565.0 > 520.0	12.964	13.000	-0.036		1543585	41.3		82.6	4747	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.964	13.000	-0.036	1.000	678501	26.0		130	2991	
D 28 13C2 PFDaA										
615.0 > 570.0	13.583	13.609	-0.026		1775672	45.4		90.8	5236	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.583	13.611	-0.028	1.000	659146	25.4		127	1290	
31 PFDoS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.044	14.083	-0.039	1.000	206389	NC			628	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.098	14.130	-0.032	1.000	827090	24.8		124	1247	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.537	14.571	-0.034	1.000	478423	22.1		111	277	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.543	14.571	-0.028		1463155	40.8		81.5	4057	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA										
815.0 > 770.0	15.179	15.207	-0.028		1834802	42.6		85.1	6199	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.179	15.208	-0.029	1.000	946964	22.3		112	1317	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.516	15.543	-0.027	1.000	838124	21.2		106	1517	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_049.d

Injection Date: 20-Feb-2016 07:05:28

Instrument ID: A6

Lims ID: LCSD 320-100093/3-A

Client ID:

Operator ID: JRB

ALS Bottle#: 3

Worklist Smp#: 47

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

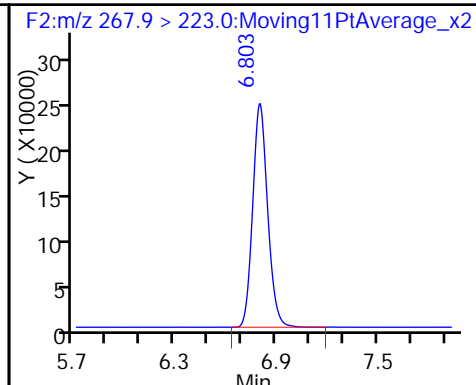
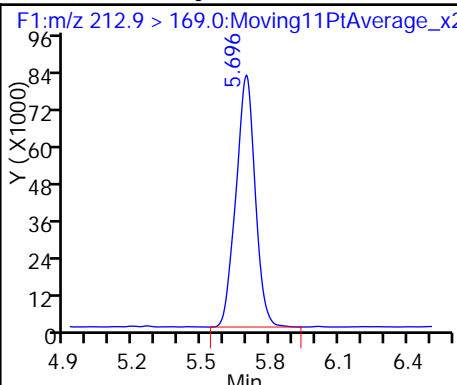
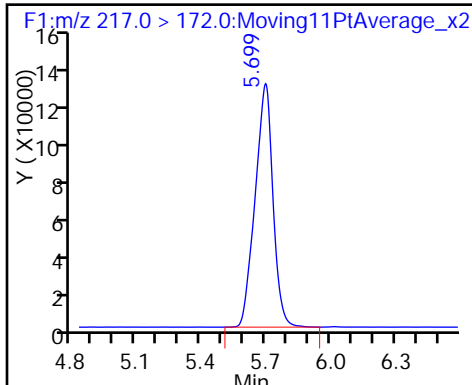
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

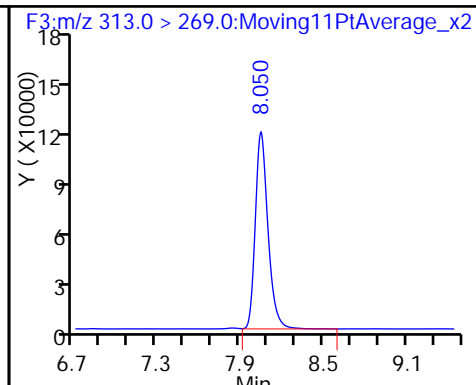
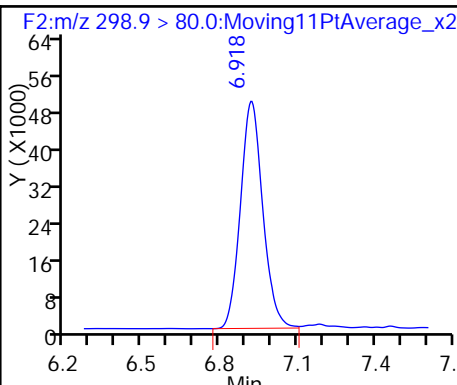
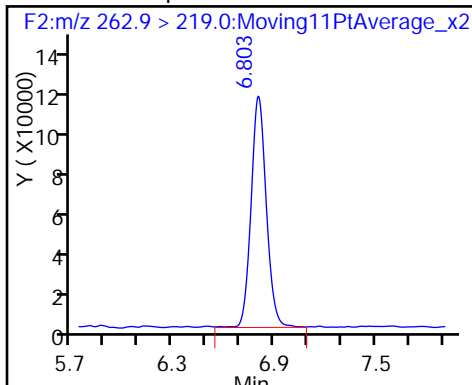
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

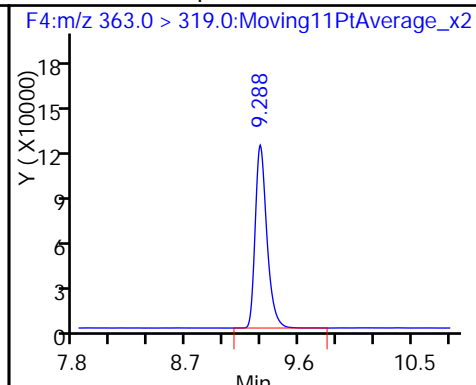
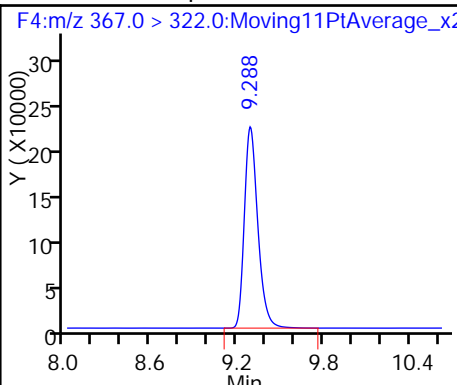
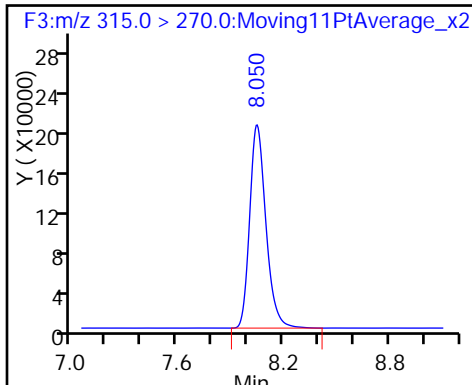
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

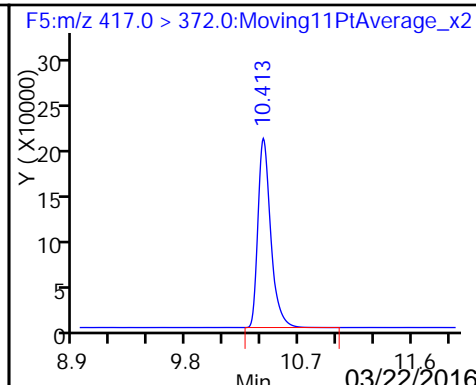
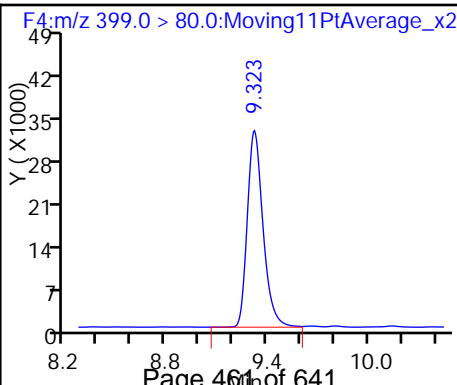
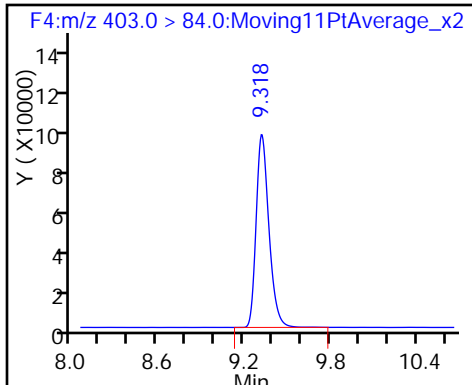
9 Perfluoroheptanoic acid

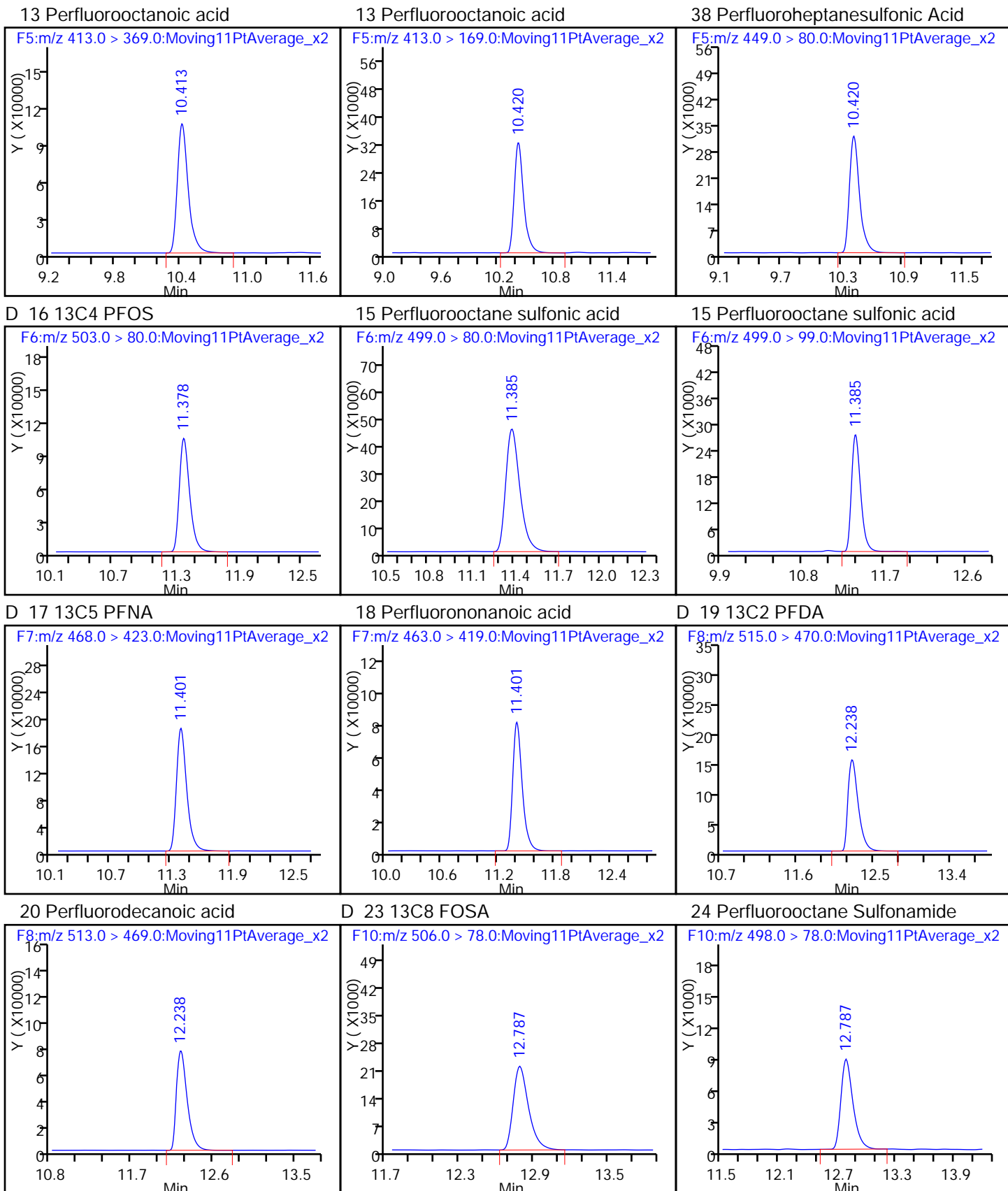


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

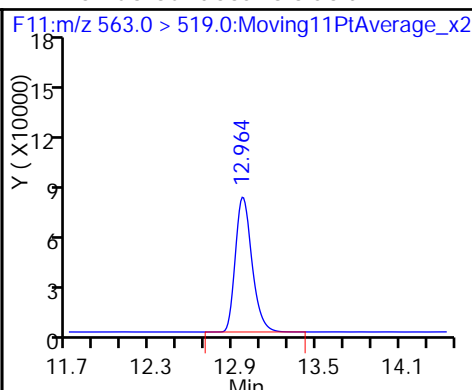
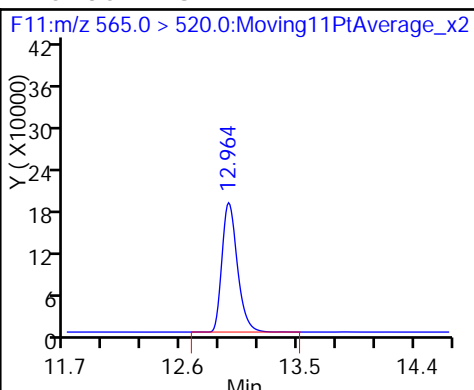
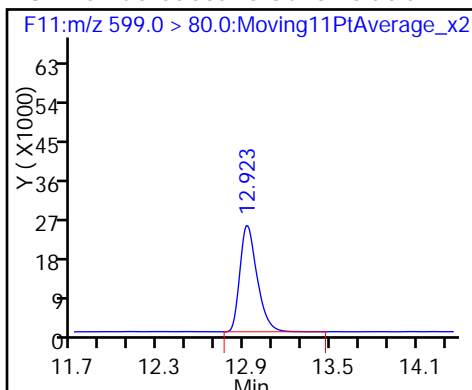




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUnA

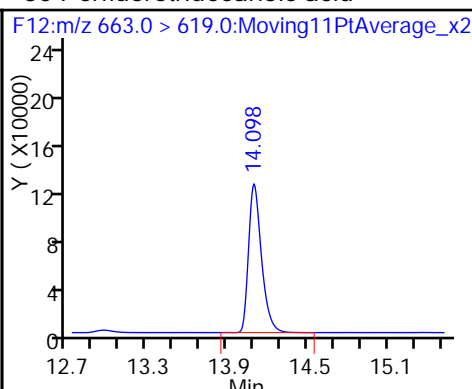
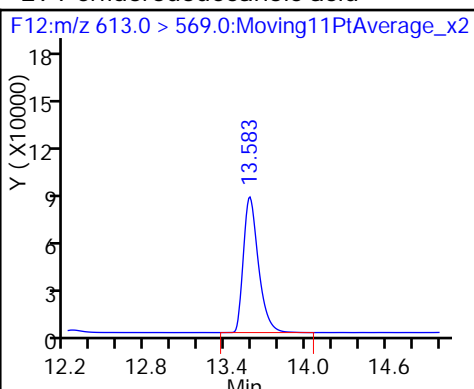
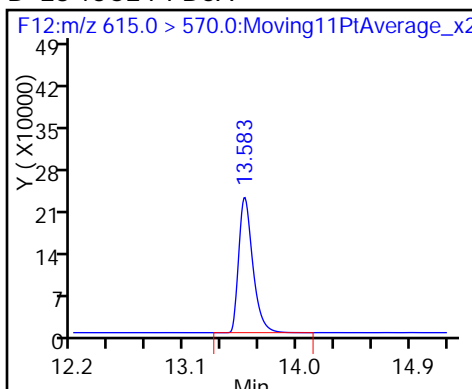
27 Perfluoroundecanoic acid



D 28 13C2 PFDaA

29 Perfluorododecanoic acid

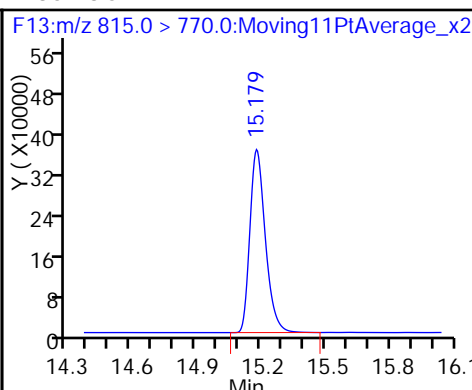
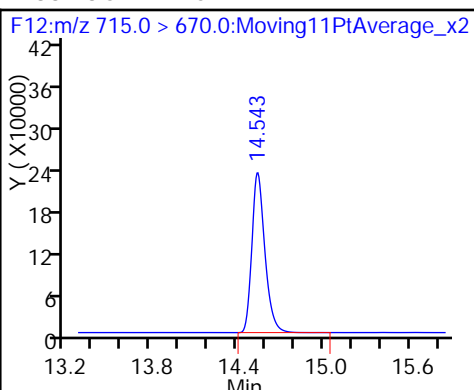
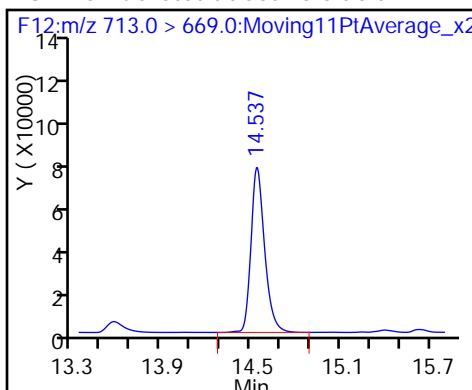
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

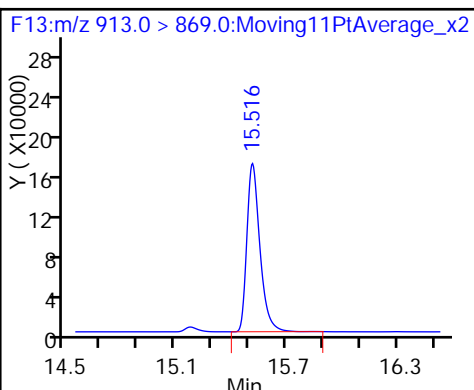
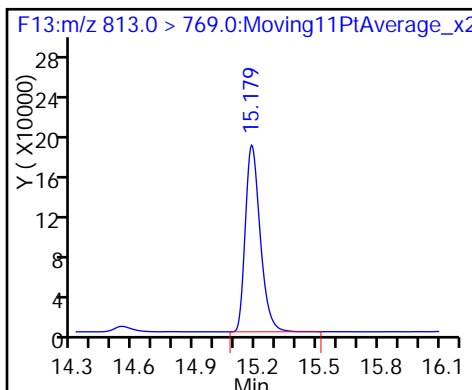
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1  
 SDG No.: CTO WE7G PFC Sampling  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCSD 320-100147/3-A  
 Matrix: Water Lab File ID: 19FEB2016A6A\_063.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 02/10/2016 07:14  
 Sample wt/vol: 500 (mL) Date Analyzed: 02/20/2016 12:02  
 Con. Extract Vol.: 1.00 (mL) Dilution Factor: 1  
 Injection Volume: 15 (uL) GC Column: Acquity ID: 2.1 (mm)  
 % Moisture: \_\_\_\_\_ GPC Cleanup: (Y/N) N  
 Analysis Batch No.: 100906 Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
375-85-9	Perfluoroheptanoic acid (PFHpA)	0.0461		0.0025	0.0020	0.00080
335-67-1	Perfluorooctanoic acid (PFOA)	0.0488		0.0025	0.0020	0.00075
375-95-1	Perfluorononanoic acid (PFNA)	0.0534		0.0025	0.0020	0.00065
375-73-5	Perfluorobutanesulfonic acid (PFBS)	0.0461		0.0025	0.0020	0.00092
355-46-4	Perfluorohexanesulfonic acid (PFHxS)	0.0473		0.0025	0.0020	0.00087
1763-23-1	Perfluorooctanesulfonic acid (PFOS)	0.0502		0.0040	0.0030	0.0013

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL01892	13C4-PFHpA	123		25-150
STL00990	13C4 PFOA	115		25-150
STL00995	13C5 PFNA	116		25-150
STL00994	18O2 PFHxS	107		25-150
STL00991	13C4 PFOS	100		25-150



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_063.d  
 Lims ID: LCSD 320-100147/3-A  
 Client ID:  
 Sample Type: LCSD  
 Inject. Date: 20-Feb-2016 12:02:40 ALS Bottle#: 16 Worklist Smp#: 61  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCSD 320-100147/3-A  
 Misc. Info.: Acquity BEH 1.7um, 3X150mm T=50\*C  
 Operator ID: JRB Instrument ID: A6  
 Method: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\PFAC\_A6.m  
 Limit Group: LC PFC\_DOD ICAL  
 Last Update: 22-Feb-2016 13:06:10 Calib Date: 19-Feb-2016 17:15:39  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_010.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK010

First Level Reviewer: barnettj Date: 22-Feb-2016 13:19:25

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 13C4 PFBA	217.0 > 172.0	5.703	5.712	-0.009	865315	54.1		108	3010	
2 Perfluorobutyric acid	212.9 > 169.0	5.703	5.715	-0.012	584537	25.9		130	1881	
D 3 13C5-PFPeA	267.9 > 223.0	6.814	6.827	-0.013	1896686	61.7		123	5073	
4 Perfluoropentanoic acid	262.9 > 219.0	6.814	6.828	-0.014	914284	23.1		115	180	
5 Perfluorobutane Sulfonate	298.9 > 80.0	6.933	6.939	-0.006	366254	NC			487	
	298.9 > 99.0	6.933	6.939	-0.006	179580		2.04(0.00-0.00)		418	
40 Perfluorobutanesulfonic acid	298.9 > 80.0	6.933	6.939	-0.006	366254	23.0		130		
7 Perfluorohexanoic acid	313.0 > 269.0	8.066	8.078	-0.012	915746	25.4		127	1635	
D 6 13C2 PFHxA	315.0 > 270.0	8.060	8.080	-0.020	1661145	58.6		117	6670	
22 PFPeS (Perflouro-1-pentanesulfonat	349.0 > 80.0	8.142	8.158	-0.016	272614	NC			1437	
D 8 13C4-PFHpA	367.0 > 322.0	9.299	9.322	-0.023	1899497	61.6		123	5479	
9 Perfluoroheptanoic acid	363.0 > 319.0	9.299	9.322	-0.023	913727	23.1		115	3225	
D 11 18O2 PFHxS	403.0 > 84.0	9.334	9.358	-0.024	757395	50.6		107	3894	
10 Perfluorohexane Sulfonate	399.0 > 80.0	9.334	9.364	-0.030	255878	NC			598	
41 Perfluorohexanesulfonic acid	399.0 > 80.0	9.334	9.364	-0.030	255878	23.7		125		03/22/2016

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 12 13C4 PFOA										
417.0 > 372.0	10.426	10.450	-0.024		1965668	57.6		115	8320	
13 Perfluorooctanoic acid										
413.0 > 369.0	10.426	10.450	-0.024	1.000	962302	24.4		122	1158	
413.0 > 169.0	10.426	10.450	-0.024	1.000	319130		3.02(0.00-0.00)		832	
14 Perfluoroheptane Sulfonate										
449.0 > 80.0	10.433	10.457	-0.024	1.000	279542	NC			913	
38 Perfluoroheptanesulfonic Acid										
449.0 > 80.0	10.433	10.457	-0.024	1.000	279542	24.9		131		
D 16 13C4 PFOS										
503.0 > 80.0	11.392	11.414	-0.022		866581	47.7		99.7	4049	
15 Perfluorooctane sulfonic acid										
499.0 > 80.0	11.392	11.417	-0.025	1.000	442810	25.1		131	418	
499.0 > 99.0	11.392	11.417	-0.025	1.000	235467		1.88(0.00-0.00)		777	
D 17 13C5 PFNA										
468.0 > 423.0	11.415	11.434	-0.019		1641246	58.2		116	7472	
18 Perfluorononanoic acid										
463.0 > 419.0	11.415	11.439	-0.024	1.000	707121	26.7		133	2987	
21 PFNS (Perflouro-1-nonanesulfonate)										
549.0 > 80.0	12.224	12.249	-0.025	1.000	240757	NC			732	
D 19 13C2 PFDA										
515.0 > 470.0	12.255	12.275	-0.020		1581281	55.2		110	3303	
20 Perfluorodecanoic acid										
513.0 > 469.0	12.255	12.276	-0.021	1.000	750409	24.0		120	2428	
D 23 13C8 FOSA										
506.0 > 78.0	12.804	12.810	-0.006		1783320	43.0		85.9	2421	
24 Perfluorooctane Sulfonamide										
498.0 > 78.0	12.804	12.812	-0.008	1.000	810653	24.5		123	1681	
25 Perfluorodecane Sulfonate										
599.0 > 80.0	12.929	12.949	-0.020	1.000	239484	NC			872	
39 Perfluorodecane Sulfonic acid										
599.0 > 80.0	12.929	12.949	-0.020	1.000	239484	24.0		124		
D 26 13C2 PFUnA										
565.0 > 520.0	12.981	13.000	-0.019		2055157	55.0		110	6847	
27 Perfluoroundecanoic acid										
563.0 > 519.0	12.981	13.000	-0.019	1.000	878074	25.3		126	1871	
D 28 13C2 PFDoA										
615.0 > 570.0	13.588	13.609	-0.021		2289189	58.5		117	6302	
29 Perfluorododecanoic acid										
613.0 > 569.0	13.588	13.611	-0.023	1.000	872443	26.1		130	953	
31 PFDoS (Perflouro-1-dodecanesulfona										
699.0 > 80.0	14.055	14.083	-0.028	1.000	264376	NC			887	
30 Perfluorotridecanoic acid										
663.0 > 619.0	14.109	14.130	-0.021	1.000	1087698	25.2		126	1007	
32 Perfluorotetradecanoic acid										
713.0 > 669.0	14.552	14.571	-0.019	1.000	639855	22.9		115	227	
D 33 13C2-PFTeDA										
715.0 > 670.0	14.552	14.571	-0.019		1928200	53.7		107	5038	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 35 13C2-PFHxDA										
815.0 > 770.0	15.188	15.207	-0.019		2519728	58.5		117	5345	
34 Perfluorohexadecanoic acid										
813.0 > 769.0	15.188	15.208	-0.020	1.000	1367316	25.3		126	1802	
36 Perfluorooctadecanoic acid										
913.0 > 869.0	15.519	15.543	-0.024	1.000	1259800	24.8		124	1158	

**QC Flag Legend**

Processing Flags

NC - Not Calibrated

TestAmerica Sacramento

Data File: \\ChromNA\Sacramento\ChromData\A6\20160219-28492.b\19FEB2016A6A\_063.d

Injection Date: 20-Feb-2016 12:02:40

Instrument ID: A6

Lims ID: LCSD 320-100147/3-A

Client ID:

Operator ID: JRB

ALS Bottle#: 16

Worklist Smp#: 61

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

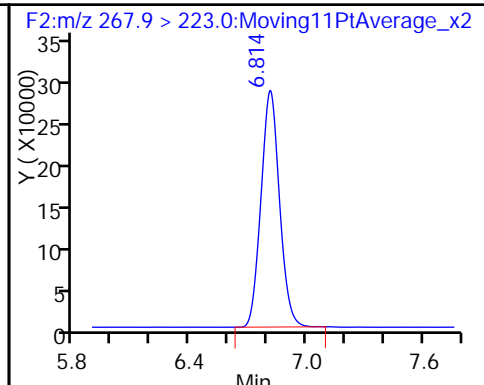
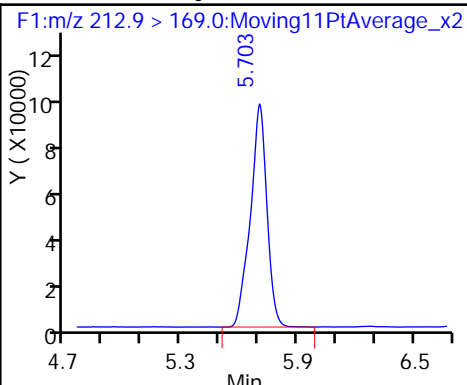
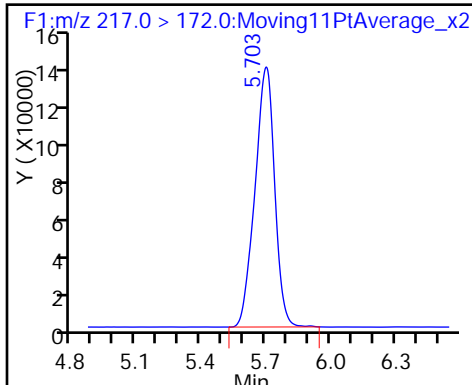
Method: PFAC\_A6

Limit Group: LC PFC\_DOD ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

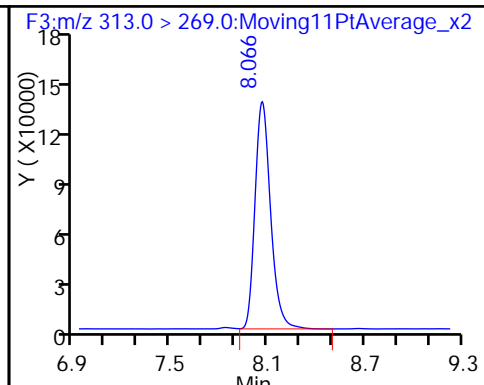
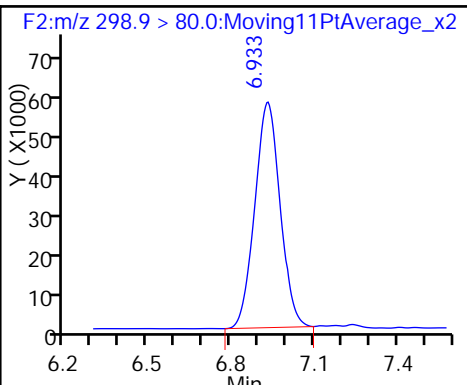
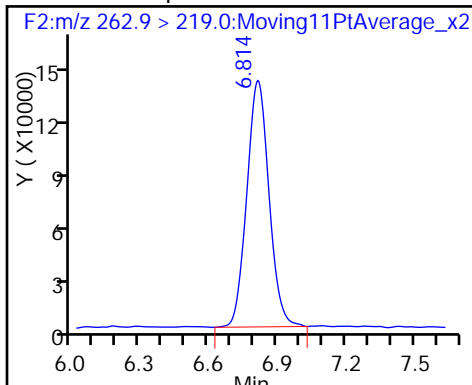
D 3 13C5-PFPeA



4 Perfluoropentanoic acid

40 Perfluorobutanesulfonic acid

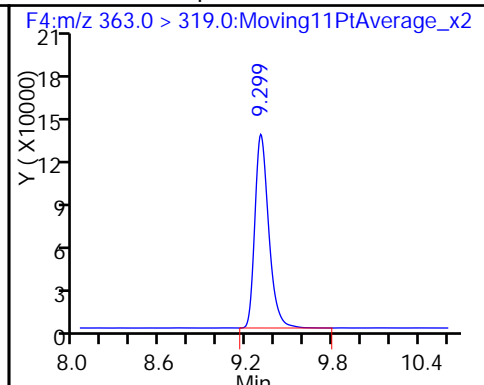
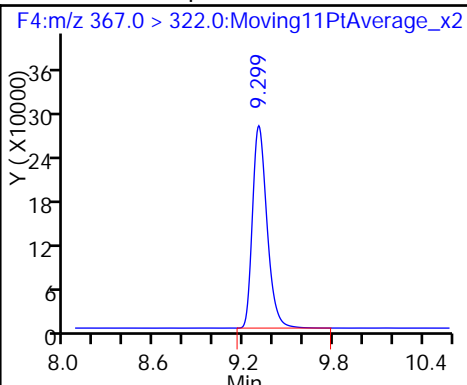
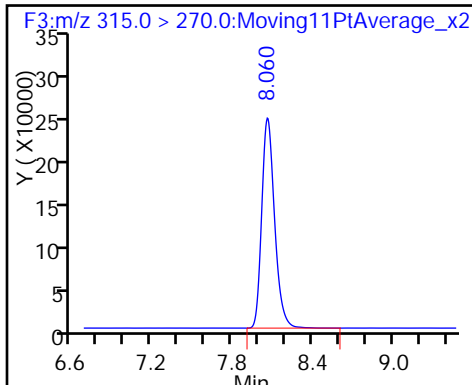
7 Perfluorohexanoic acid



D 6 13C2 PFHxA

D 8 13C4-PFHpA

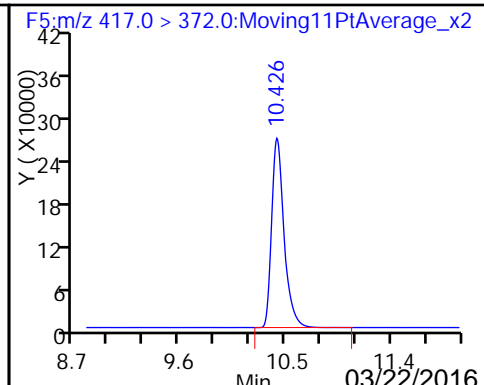
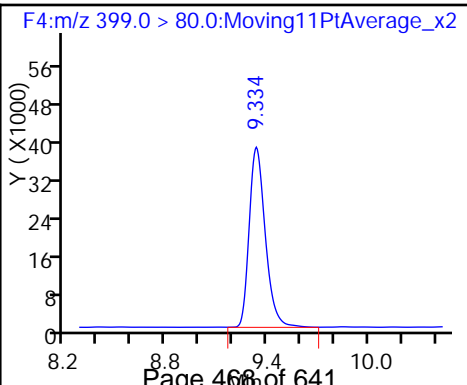
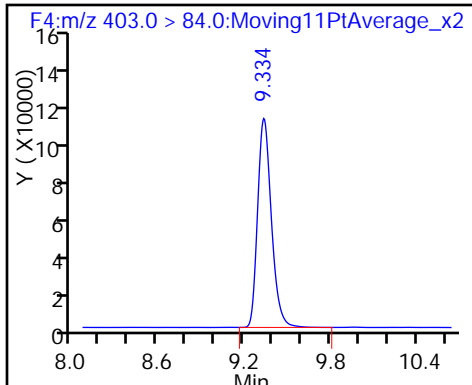
9 Perfluoroheptanoic acid

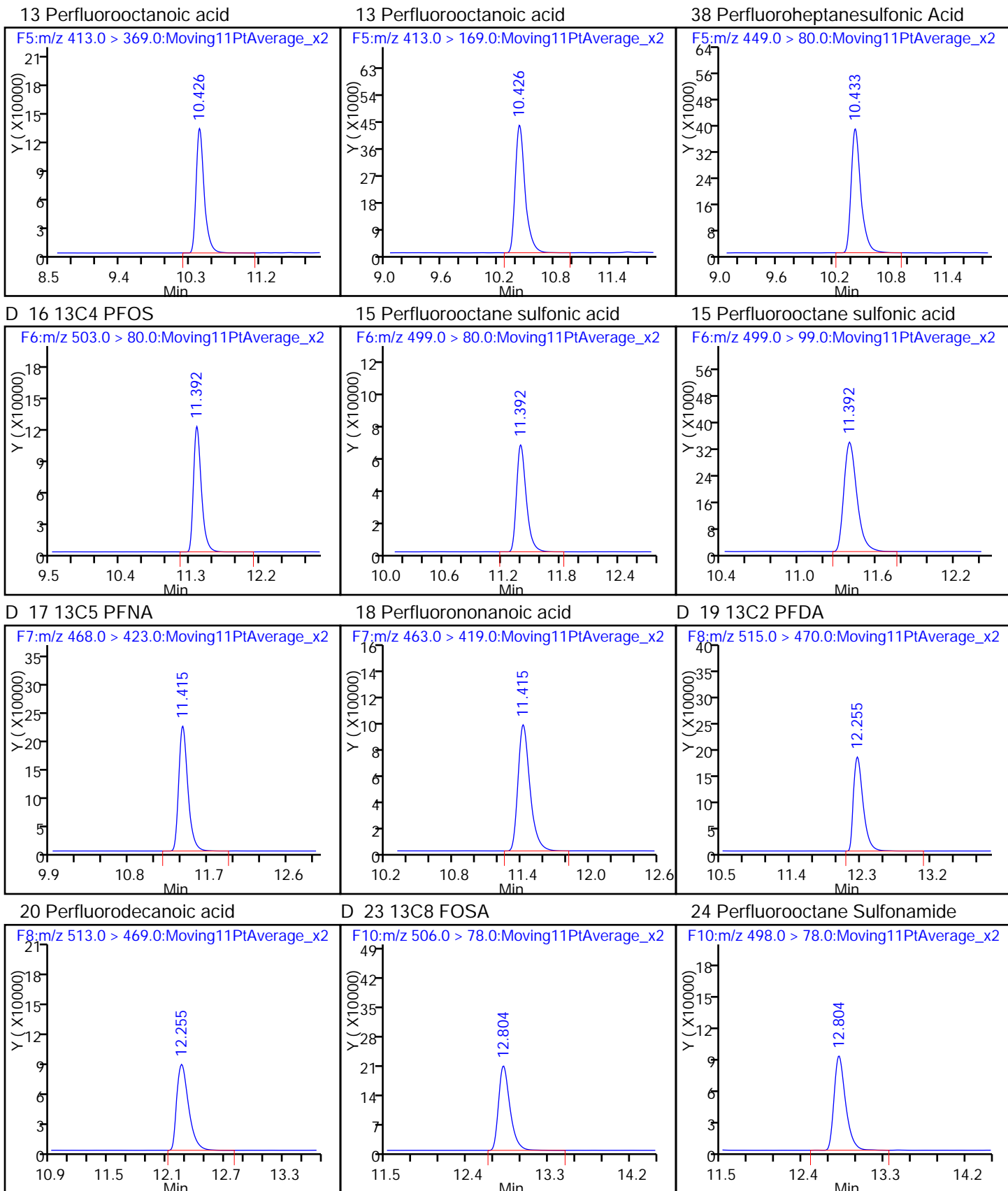


D 11 18O2 PFHxS

41 Perfluorohexanesulfonic acid

D 12 13C4 PFOA

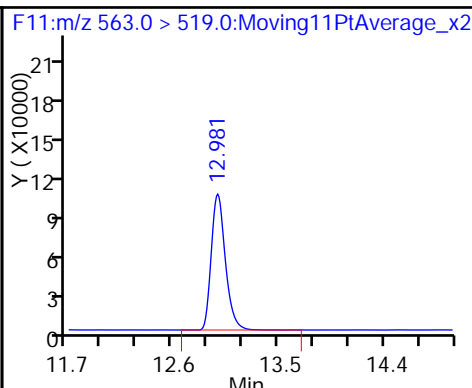
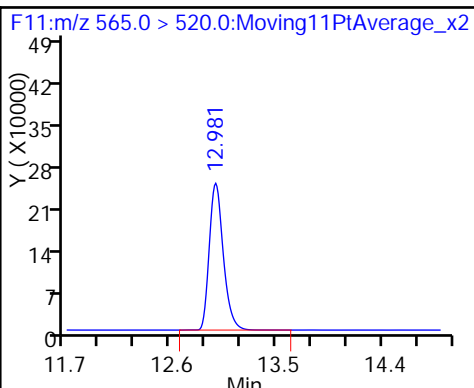
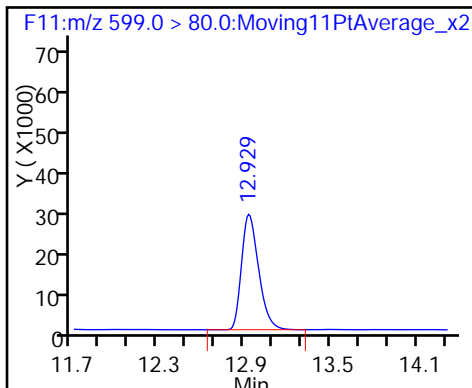




39 Perfluorodecane Sulfonic acid

D 26 13C2 PFUa

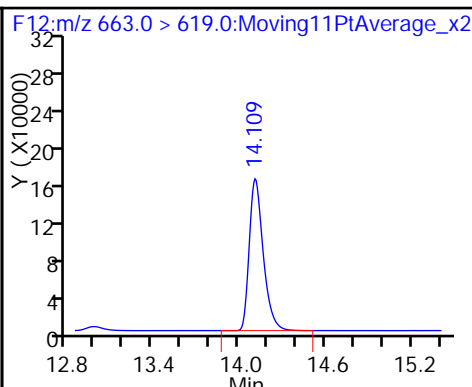
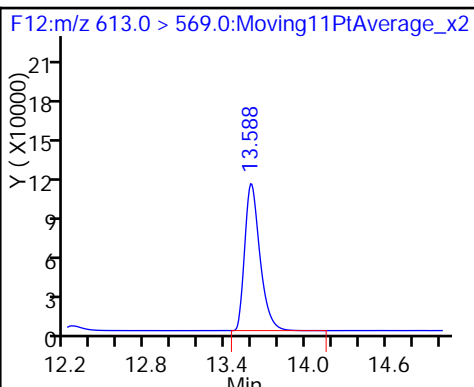
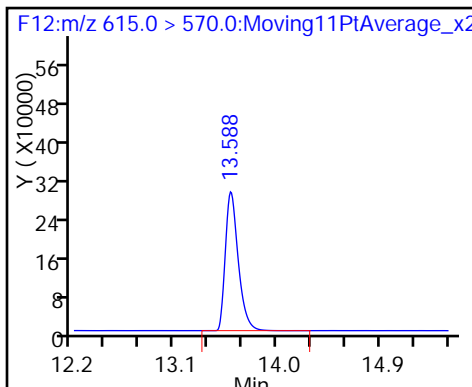
27 Perfluoroundecanoic acid



D 28 13C2 PFDa

29 Perfluorododecanoic acid

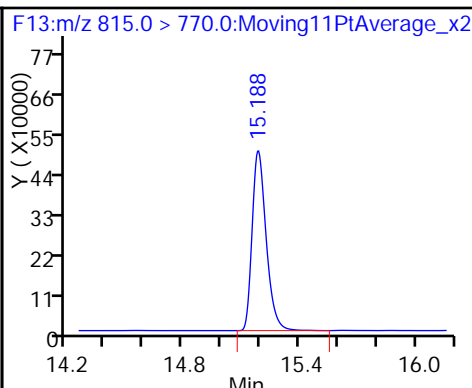
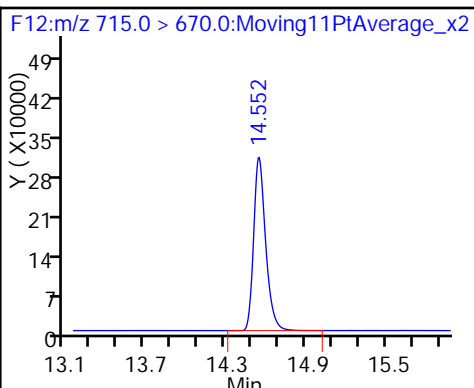
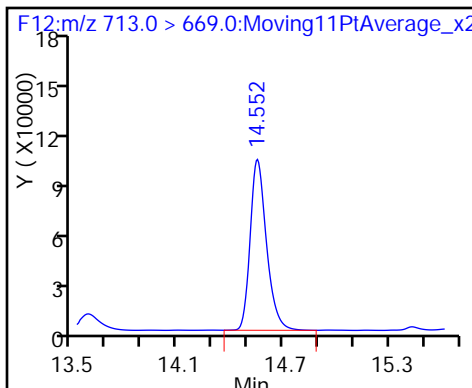
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid

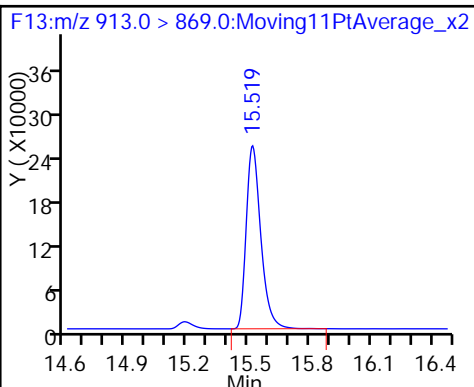
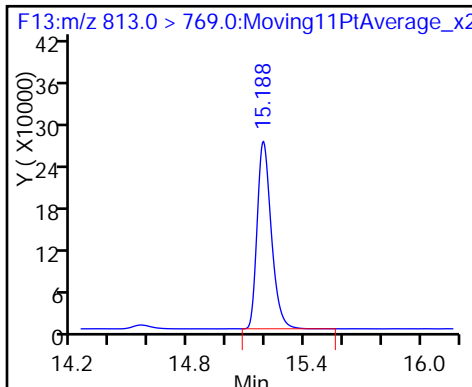
D 33 13C2-PFTeDA

D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid

36 Perfluorooctadecanoic acid



LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4 Start Date: 02/12/2016 10:35

Analysis Batch Number: 100305 End Date: 02/13/2016 15:33

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
STD 320-100305/2 IC		02/12/2016 10:35	1	12FEB2016A4A_00 3.d	Acquity 2.1(mm)
STD 320-100305/3 IC		02/12/2016 10:56	1	12FEB2016A4A_00 4.d	Acquity 2.1(mm)
STD 320-100305/4 IC		02/12/2016 11:17	1	12FEB2016A4A_00 5.d	Acquity 2.1(mm)
STD 320-100305/5 IC		02/12/2016 11:39	1	12FEB2016A4A_00 6.d	Acquity 2.1(mm)
STD 320-100305/6 IC		02/12/2016 12:00	1	12FEB2016A4A_00 7.d	Acquity 2.1(mm)
STD 320-100305/7 IC		02/12/2016 12:21	1	12FEB2016A4A_00 8.d	Acquity 2.1(mm)
STD 320-100305/8 IC		02/12/2016 12:42	1	12FEB2016A4A_00 9.d	Acquity 2.1(mm)
ZZZZZ		02/12/2016 13:03	1		Acquity 2.1(mm)
ICV 320-100305/10		02/12/2016 13:24	1	12FEB2016A4A_01 1.d	Acquity 2.1(mm)
CCV 320-100305/92		02/12/2016 16:57	1		Acquity 2.1(mm)
CCV 320-100305/40		02/13/2016 00:01	1	12FEB2016A4A_04 1.d	Acquity 2.1(mm)
MB 320-100093/1-A		02/13/2016 00:22	1	12FEB2016A4A_04 2.d	Acquity 2.1(mm)
ZZZZZ		02/13/2016 00:43	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 01:04	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 01:26	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 01:47	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 02:08	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 02:29	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 02:50	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 03:11	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 03:33	1		Acquity 2.1(mm)
CCV 320-100305/51		02/13/2016 03:54	1	12FEB2016A4A_05 2.d	Acquity 2.1(mm)
ZZZZZ		02/13/2016 04:15	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 04:36	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 04:57	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 05:19	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 05:40	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 06:01	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 06:22	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 06:43	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 07:04	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 07:26	1		Acquity 2.1(mm)
CCV 320-100305/62		02/13/2016 07:47	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 08:08	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 08:29	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 08:50	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 09:12	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 09:33	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 09:54	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 10:15	1		Acquity 2.1(mm)

LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A4 Start Date: 02/12/2016 10:35

Analysis Batch Number: 100305 End Date: 02/13/2016 15:33

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
ZZZZZ		02/13/2016 10:36	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 10:57	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 11:19	1		Acquity 2.1(mm)
CCV 320-100305/73		02/13/2016 11:40	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 12:01	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 12:22	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 12:43	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 13:04	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 13:26	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 13:47	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 14:08	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 14:29	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 14:50	1		Acquity 2.1(mm)
ZZZZZ		02/13/2016 15:12	1		Acquity 2.1(mm)
CCV 320-100305/84		02/13/2016 15:33	1		Acquity 2.1(mm)



LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Instrument ID: A6 Start Date: 02/19/2016 15:08

Analysis Batch Number: 100906 End Date: 02/20/2016 16:59

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
STD 320-100906/2 IC		02/19/2016 15:08	1	19FEB2016A6A_00 4.d	Acquity 2.1(mm)
STD 320-100906/3 IC		02/19/2016 15:29	1	19FEB2016A6A_00 5.d	Acquity 2.1(mm)
STD 320-100906/4 IC		02/19/2016 15:50	1	19FEB2016A6A_00 6.d	Acquity 2.1(mm)
STD 320-100906/5 IC		02/19/2016 16:11	1	19FEB2016A6A_00 7.d	Acquity 2.1(mm)
STD 320-100906/6 IC		02/19/2016 16:33	1	19FEB2016A6A_00 8.d	Acquity 2.1(mm)
STD 320-100906/7 IC		02/19/2016 16:54	1	19FEB2016A6A_00 9.d	Acquity 2.1(mm)
STD 320-100906/8 IC		02/19/2016 17:15	1	19FEB2016A6A_01 0.d	Acquity 2.1(mm)
ZZZZZ		02/19/2016 17:36	1		Acquity 2.1(mm)
ICV 320-100906/10		02/19/2016 17:58	1	19FEB2016A6A_01 2.d	Acquity 2.1(mm)
CCV 320-100906/43		02/20/2016 05:40	1	19FEB2016A6A_04 5.d	Acquity 2.1(mm)
MB 320-100093/1-A		02/20/2016 06:23	1	19FEB2016A6A_04 7.d	Acquity 2.1(mm)
LCS 320-100093/2-A		02/20/2016 06:44	1	19FEB2016A6A_04 8.d	Acquity 2.1(mm)
LCSD 320-100093/3-A		02/20/2016 07:05	1	19FEB2016A6A_04 9.d	Acquity 2.1(mm)
ZZZZZ		02/20/2016 07:26	1		Acquity 2.1(mm)
ZZZZZ		02/20/2016 07:47	1		Acquity 2.1(mm)
ZZZZZ		02/20/2016 08:09	1		Acquity 2.1(mm)
ZZZZZ		02/20/2016 08:30	1		Acquity 2.1(mm)
ZZZZZ		02/20/2016 08:51	1		Acquity 2.1(mm)
ZZZZZ		02/20/2016 09:12	1		Acquity 2.1(mm)
CCV 320-100906/54		02/20/2016 09:34	1	19FEB2016A6A_05 6.d	Acquity 2.1(mm)
ZZZZZ		02/20/2016 09:55	1		Acquity 2.1(mm)
320-17154-1		02/20/2016 10:16	1	19FEB2016A6A_05 8.d	Acquity 2.1(mm)
320-17154-3		02/20/2016 10:37	1	19FEB2016A6A_05 9.d	Acquity 2.1(mm)
320-17154-6		02/20/2016 10:58	1	19FEB2016A6A_06 0.d	Acquity 2.1(mm)
MB 320-100147/1-A		02/20/2016 11:20	1	19FEB2016A6A_06 1.d	Acquity 2.1(mm)
LCS 320-100147/2-A		02/20/2016 11:41	1	19FEB2016A6A_06 2.d	Acquity 2.1(mm)
LCSD 320-100147/3-A		02/20/2016 12:02	1	19FEB2016A6A_06 3.d	Acquity 2.1(mm)
320-17154-8		02/20/2016 12:23	1	19FEB2016A6A_06 4.d	Acquity 2.1(mm)
320-17154-10		02/20/2016 12:45	1	19FEB2016A6A_06 5.d	Acquity 2.1(mm)
CCV 320-100906/65		02/20/2016 13:27	1	19FEB2016A6A_06 7.d	Acquity 2.1(mm)
CCV 320-100906/75		02/20/2016 16:59	1		Acquity 2.1(mm)

LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento Job No.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Batch Number: 100093 Batch Start Date: 02/09/16 10:23 Batch Analyst: Arauz, Horacio J

Batch Method: 3535 Batch End Date: 02/09/16 20:35

Lab Sample ID	Client Sample ID	Method Chain	Basis	GrossWeight	TareWeight	InitialAmount	FinalAmount	LCMPFCSU 00025	LCPFCSP 00039
MB 320-100093/1		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	
LCS 320-100093/2		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	20 uL
LCSD 320-100093/3		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	20 uL
320-17154-A-1	OF-FB09-0216	3535, WS-LC-0025	T	549.36 g	44.49 g	504.9 mL	1.00 mL	50 uL	
320-17154-A-3	OF-FB37-0216	3535, WS-LC-0025	T	593.62 g	42.89 g	550.7 mL	1.00 mL	50 uL	
320-17154-A-6	OF-FB11-0216	3535, WS-LC-0025	T	596.20 g	44.78 g	551.4 mL	1.00 mL	50 uL	

Batch Notes	
Balance ID	QA-070
Batch Comment	0.1N Sodium Hydrox./H2O 565548; MeOH 572662; J.T.B Hexane 0000101243
H2O ID	02-09-16
Pipette ID	EC15219
Analyst ID - Reagent Drop	HJA
Analyst ID - SU Reagent Drop	HJA
Analyst ID - SU Reagent Drop Witness	SNE
Solvent Lot #	578365
Solvent Name	0.3% Ammonium hydroxide/MeOH
SOP Number	WS-LC-0025
SPE Cartridge Type	Wax 500mg
Solid Phase Extraction Disk ID	002635307A

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.: 320-17154-1

SDG No.: CTO WE7G PFC Sampling

Batch Number: 100147 Batch Start Date: 02/10/16 07:14 Batch Analyst: Arauz, Horacio J

Batch Method: 3535 Batch End Date: 02/10/16 20:25

Lab Sample ID	Client Sample ID	Method Chain	Basis	GrossWeight	TareWeight	InitialAmount	FinalAmount	LCMPFCSU 00025	LCPFCSP 00039
MB 320-100147/1		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	
LCS 320-100147/2		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	20 uL
LCSD 320-100147/3		3535, WS-LC-0025				500 mL	1.00 mL	50 uL	20 uL
320-17154-A-8	OF-FB28-0216	3535, WS-LC-0025	T	566.05 g	44.50 g	521.6 mL	1.00 mL	50 uL	
320-17154-A-10	OF-FB67-0216	3535, WS-LC-0025	T	578.68 g	42.81 g	535.9 mL	1.00 mL	50 uL	

Batch Notes	
Balance ID	QA-070
Batch Comment	0.1N Sodium Hydrox./H2O 565548; MeOH 572662; J.T.B Hexane 0000101243
H2O ID	02-09-16
Pipette ID	EC15219
Analyst ID - Reagent Drop	HJA
Analyst ID - SU Reagent Drop	HJA
Analyst ID - SU Reagent Drop Witness	CRF
Solvent Lot #	578365
Solvent Name	0.3% Ammonium hydroxide/MeOH
SOP Number	WS-LC-0025
SPE Cartridge Type	Wax 500mg
Solid Phase Extraction Disk ID	002635307A

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): 320-17150, 320-17154

Work List ID(s): 28353, 28492

Extraction Batch: 100093, 100147

Analysis Batch(es): 100305, 100906

Delivery Rank: 4

Due Date: 2/8/16

A. Calibration/Instrument Run QC	1 <sup>st</sup> Level	2 <sup>nd</sup> Level	N/A
1. ICAL locked in Chrom and TALS? ICAL Batch#	✓	✓	
2. ICAL, CCV Frequency & Criteria met.	✓	✓	
• RF <sub>average</sub> criteria appropriate for the method.	✓	✓	
• Linear Regression criteria appropriate if required ( $r > 0.995$ ).	✓	✓	
• Quadratic fit criteria appropriate if required ( $r^2 > 0.990$ ).			✓
• 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>B. QA/QC</b>			
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.	✓	✓	
<b>C. Sample Analysis</b>			
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)	✓	✓	
<b>D. Documentation</b>			
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.

1<sup>st</sup> Level (Analyst): JRB

Date: 2/22/16

2<sup>nd</sup> Level Reviewer: Mewall

Date: 2/22/2016

18

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100093

Analyst: Arauz, Horacio J

Batch Open: 2/9/2016 10:23:08AM

Method Code: 320-3535\_IVWT-320

Batch End: 2/09/16 20:35

①

## Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	InitAmnt FinAmnt	PHs Rcvd Adj1 Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
1 MB-320-100093/1 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		MB 320-100093/1-A
2 LCS-320-100093/2 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		LCS 320-100093/2-A
3 LCSD-320-100093/3 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		LCSD 320-100093/3-A
4 320-17150-A-1 (PFC_IDA_DOD5)	N/A (320-17150-1)	603.30 g 44.76 g	558.5 mL 1.00 mL		2/8/16	5_Days	4	cancel	320-17150-A-1-A
5 320-17150-A-2 (PFC_IDA_DOD5)	N/A (320-17150-1)	575.48 g 42.91 g	532.6 mL 1.00 mL		2/8/16	5_Days	4		320-17150-A-2-A
6 320-17150-A-3 (PFC_IDA_DOD5)	N/A (320-17150-1)	597.46 g 44.83 g	552.6 mL 1.00 mL		2/8/16	5_Days	4	cancel	320-17150-A-3-A
7 320-17150-A-4 (PFC_IDA_DOD5)	N/A (320-17150-1)	564.91 g 42.89 g	522 mL 1.00 mL		2/8/16	5_Days	4		320-17150-A-4-A
8 320-17150-A-5 (PFC_IDA_DOD5)	N/A (320-17150-1)	578.05 g 43.87 g	534.2 mL 1.00 mL		2/8/16	5_Days	4	cancel	320-17150-A-5-A
9 320-17150-A-6 (PFC_IDA_DOD5)	N/A (320-17150-1)	577.94 g 43.02 g	534.9 mL 1.00 mL		2/8/16	5_Days	4		320-17150-A-6-A
10 320-17150-A-7 (PFC_IDA_DOD5)	N/A (320-17150-1)	582.39 g 44.94 g	537.5 mL 1.00 mL		2/8/16	5_Days	4	cancel	320-17150-A-7-A

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)










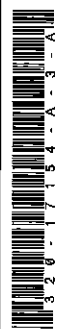
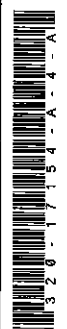

Batch Number: 320-100093

Analyst: Arauz, Horacio J

Batch Open: 2/9/2016 10:23:08AM

Method Code: 320-3535\_IVWT-320

Batch End:

Line	Sample ID	Weight	Volume	Concentration	Date	Days	Notes	Barcode
11	320-17150-A-8 (PFC_IDA_DOD5)	583.91 g 42.98 g	540.9 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4		
12	320-17150-A-9 (PFC_IDA_DOD5)	588.67 g 44.02 g	544.7 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4		
13	320-17150-A-10 (PFC_IDA_DOD5)	583.36 g 44.57 g	538.8 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4	Cancel	
14	320-17150-A-11 (PFC_IDA_DOD5)	597.25 g 42.99 g	554.3 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4		
15	320-17150-A-12 (PFC_IDA_DOD5)	596.09 g 44.90 g	551.2 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4	Cancel	
16	320-17150-A-13 (PFC_IDA_DOD5)	589.34 g 44.06 g	545.3 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4		
17	320-17150-A-14 (PFC_IDA_DOD5)	593.57 g 44.47 g	549.1 mL 1.00 mL	N/A (320-17150-1)	2/8/16	4	Cancel	
18	320-17154-A-1 (PFC_IDA_DOD5)	549.36 g 44.49 g	504.9 mL 1.00 mL	N/A (320-17154-1)	2/8/16	4		
19	320-17154-A-2 (PFC_IDA_DOD5)	577.08 g 45.84 g	531.2 mL 1.00 mL	N/A (320-17154-1)	2/8/16	4	Cancel	
20	320-17154-A-3 (PFC_IDA_DOD5)	593.62 g 42.89 g	550.7 mL 1.00 mL	N/A (320-17154-1)	2/8/16	4		
21	320-17154-A-4 (PFC_IDA_DOD5)	581.63 g 43.14 g	538.5 mL 1.00 mL	N/A (320-17154-1)	2/8/16	4	cancel	
22	320-17154-A-5 (PFC_IDA_DOD5)	586.86 g 45.91 g	541 mL 1.00 mL	N/A (320-17154-1)	2/8/16	4	Cancel	

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments),

Batch Number: 320-100093      Analyst: Arauz, Horacio J      Batch Open: 2/9/2016 10:23:08AM

Method Code: 320-3535\_IVWT-320      Batch End:

320-17154-A-6 (PFC_IDA_DOD5)	N/A (320-17154-1)	596.20 g	551.4 mL	2/8/16	5_Days	4	
		44.78 g	1.00 mL				

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100093

Analyst: Arauz, Horacio J

Batch Open: 2/9/2016 10:23:08AM

Method Code: 320-3535\_IVVWT-320

Batch End:

## Batch Notes

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type Wax 500mg

Solid Phase Extraction Disk Lot Number 002635307A

H2O Lot used 02-09-16

Pipette ID EC15219

Solvent Name 0.3% Ammonium hydroxide/MeOH

Solvent Lot # 578365

Analyst who added reagent HJA

SU Reagent Drop HJA

SU Reagent Drop Witness SNE

Acid Name NA

Acid Lot NA

Reagent ID NA

Reagent Lot Number NA

NaCl Lot # NA

SOP Number WS-LC-0025

Batch Comment 0.1N Sodium Hydrox./H2O 565548; MeOH 572662; J.T.B Hexane 0000101243



# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100093  
Method Code: 320-3535\_IWWT-320

Analyst: Arauz, Horacio J  
Batch Open: 2/9/2016 10:23:08AM  
Batch End:

Comments

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100093

Batch Open: 2/9/2016 10:23:08AM

Analyst: Arauz, Horacio J

Method Code: 320-3535\_IVWWT-320

Batch End:

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-100093/1	LCMPFCSU_00025	50 uL	1.00 mL	HSA 2-9-16	SAVE 2/9/16
LCS 320-100093/2	LCMPFCSU_00025	50 uL	1.00 mL		
LCS 320-100093/2	LCPFCSU_00039	20 uL	1.00 mL		
LCSD 320-100093/3	LCMPFCSU_00025	50 uL	1.00 mL		
LCSD 320-100093/3	LCPFCSU_00039	20 uL	1.00 mL		
320-17150-A-1	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-2	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-3	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-4	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-5	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-6	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-7	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-8	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-9	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-10	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-11	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-12	LCMPFCSU_00025	50 uL	1.00 mL		
320-17150-A-13	LCMPFCSU_00025	50 uL	1.00 mL		

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100093

Analyst: Arauz, Horacio J

Batch Open: 2/9/2016 10:23:08AM

Method Code: 320-3535\_IVWT-320

Batch End:

Sample ID	Reagent	Amount	Volume	Notes
320-17150-A-14	LCMPFCSU_00025	50 uL	1.00 mL	HZA 2-9-16 SUE 2/9/16
320-17154-A-1	LCMPFCSU_00025	50 uL	1.00 mL	
320-17154-A-2	LCMPFCSU_00025	50 uL	1.00 mL	
320-17154-A-3	LCMPFCSU_00025	50 uL	1.00 mL	
320-17154-A-4	LCMPFCSU_00025	50 uL	1.00 mL	
320-17154-A-5	LCMPFCSU_00025	50 uL	1.00 mL	
320-17154-A-6	LCMPFCSU_00025	50 uL	1.00 mL	

### Other Reagents:

Reagent

Amount/Units

Lot#:

Preparation Batch Number(s): 320-100093 Test: PFC-L

Earliest Holding Time: 2-10-16

Sample List Tab		1 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> Level Reviewer
Date and time accurate and entered into TALS correctly		/	✓
All necessary 'batch information' complete and entered into TALS correctly		/	✓

1<sup>st</sup> Level Reviewer: 

Date: 2/09/16

2<sup>nd</sup> Level Reviewer: SNE

Date: 2/10/16

Comments: \_\_\_\_\_

Box #18

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J








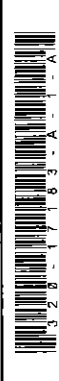


Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_IWWT-320

Batch End: 2/19/16 20:25

②

## Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG (Job #)	GrossWt TareWt	InitAmnt FinAmnt	PHs Adj1 Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
1 MB-320-100147/1 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		
2 LCS-320-100147/2 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		
3 LCSD-320-100147/3 N/A	N/A		500 mL 1.00 mL		N/A	N/A	N/A		
4 320-17154-A-7 (PFC_IDA_DOD5)	N/A (320-17154-1)	506.10 g 44.87 g	561.2 mL 1.00 mL		2/8/16	5_Days	4	Cancel	
5 320-17154-A-8 (PFC_IDA_DOD5)	N/A (320-17154-1)	566.05 g 44.50 g	521.6 mL 1.00 mL		2/8/16	5_Days	4		
6 320-17154-A-9 (PFC_IDA_DOD5)	N/A (320-17154-1)	572.25 g 45.31 g	526.9 mL 1.00 mL		2/8/16	5_Days	4	Cancel	
7 320-17154-A-10 (PFC_IDA_DOD5)	N/A (320-17154-1)	578.68 g 42.81 g	535.9 mL 1.00 mL		2/8/16	5_Days	4		
8 320-17183-A-1 (PFC_IDA_DOD5)	N/A (320-17183-1)	572.46 g 45.30 g	527.2 mL 1.00 mL		2/9/16	5_Days	4		
9 320-17183-A-2 (PFC_IDA_DOD5)	N/A (320-17183-1)	573.90 g 44.66 g	529.2 mL 1.00 mL		2/9/16	5_Days	4		
10 320-17183-A-3 (PFC_IDA_DOD5)	N/A (320-17183-1)	599.77 g 45.94 g	553.8 mL 1.00 mL		2/9/16	5_Days	4		

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)













Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_VVWT-320

Batch End:

11	320-17183-A-4 (PFC_IDA_DOD5)	N/A (320-17183-1)	591.89 g 44.22 g	547.7 mL 1.00 mL			2/9/16	5_Days	4	
12	320-17183-A-5 (PFC_IDA_DOD5)	N/A (320-17183-1)	595.34 g 45.96 g	549.4 mL 1.00 mL			2/9/16	5_Days	4	
13	320-17183-A-6 (PFC_IDA_DOD5)	N/A (320-17183-1)	582.18 g 44.82 g	537.4 mL 1.00 mL			2/9/16	5_Days	4	
14	320-17183-A-7 (PFC_IDA_DOD5)	N/A (320-17183-1)	504.2 g 43.25 g	561 mL 1.00 mL			2/9/16	5_Days	4	
15	320-17183-A-8 (PFC_IDA_DOD5)	N/A (320-17183-1)	593.22 g 44.26 g	549 mL 1.00 mL			2/9/16	5_Days	4	
16	320-17183-A-9 (PFC_IDA_DOD5)	N/A (320-17183-1)	598.43 g 44.88 g	553.6 mL 1.00 mL			2/9/16	5_Days	4	
17	320-17183-A-10 (PFC_IDA_DOD5)	N/A (320-17183-1)	574.41 g 44.85 g	529.6 mL 1.00 mL			2/9/16	5_Days	4	
18	320-17184-A-1 (PFC_IDA_DOD5)	N/A (320-17184-1)	573.03 g 44.43 g	528.6 mL 1.00 mL			2/10/16	8_Day_Rush	4	
19	320-17184-A-2 (PFC_IDA_DOD5)	N/A (320-17184-1)	582.88 g 44.14 g	538.7 mL 1.00 mL			2/10/16	8_Day_Rush	4	
20	320-17184-A-3 (PFC_IDA_DOD5)	N/A (320-17184-1)	574.41 g 44.44 g	530 mL 1.00 mL			2/10/16	8_Day_Rush	4	
21	320-17184-A-4 (PFC_IDA_DOD5)	N/A (320-17184-1)	591.29 g 44.50 g	546.8 mL 1.00 mL			2/10/16	8_Day_Rush	4	
22	320-17184-A-5 (PFC_IDA_DOD5)	N/A (320-17184-1)	582.37 g 45.04 g	537.3 mL 1.00 mL			2/10/16	8_Day_Rush	4	

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_IVWT-320

Batch End:

23	320-17184-A-6 (PFC_IDA_DOD5)	N/A (320-17184-1)	574.59 g	530 mL	2/10/16	8_Day_Rush	4	 <small>3 2 8 . 1 7 1 8 4 . A . 6 . A</small>
			44.63 g	1.00 mL				

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_IVWT-320

Batch End:

## Batch Notes

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type Wax 500mg

Solid Phase Extraction Disk Lot Number 002635307A

H2O Lot used 02-09-16

Pipette ID EC15219

Solvent Name 0.3% Ammonium hydroxide/MeOH

Solvent Lot # 578365

Analyst who added reagent HJA

SU Reagent Drop HJA

SU Reagent Drop Witness **CRF**

Acid Name NA

Acid Lot NA

Reagent ID NA

Reagent Lot Number NA

NaCl Lot # NA

SOP Number WS-LC-0025

Batch Comment 0.1N Sodium Hydrox./H2O 565548, MeOH 572662; J.T.B Hexane 0000101243



# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_IVWT-320

Batch End:

Comments

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_JVWT-320

Batch End:

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-100147/1	LCMPFCSU_00025	50 uL	1.00 mL	HSA 2-10-16	CFR z 16/16
LCS 320-100147/2	LCMPFCSU_00025	50 uL	1.00 mL		
LCS 320-100147/2	LCPFCSU_00039	20 uL	1.00 mL		
LCSD 320-100147/3	LCMPFCSU_00025	50 uL	1.00 mL		
LCSD 320-100147/3	LCPFCSU_00039	20 uL	1.00 mL		
320-17154-A-7	LCMPFCSU_00025	50 uCi	1.00 mL		
320-17154-A-8	LCMPFCSU_00025	50 uL	1.00 mL		
320-17154-A-9	LCMPFCSU_00025	50 uL	1.00 mL		
320-17154-A-10	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-1	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-2	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-3	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-4	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-5	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-6	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-7	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-8	LCMPFCSU_00025	50 uL	1.00 mL		
320-17183-A-9	LCMPFCSU_00025	50 uL	1.00 mL		

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-100147

Analyst: Arauz, Horacio J

Batch Open: 2/10/2016 7:14:29AM

Method Code: 320-3535\_IVWT-320

Batch End:

320-17183-A-10	LCMPFCSU_00025	50 uL	1.00 mL	HSA 2-10-16	CF 1/2	2 (10 116)
320-17184-A-1	LCMPFCSU_00025	50 uL	1.00 mL			
320-17184-A-2	LCMPFCSU_00025	50 uL	1.00 mL			
320-17184-A-3	LCMPFCSU_00025	50 uL	1.00 mL			
320-17184-A-4	LCMPFCSU_00025	50 uL	1.00 mL			
320-17184-A-5	LCMPFCSU_00025	50 uL	1.00 mL			
320-17184-A-6	LCMPFCSU_00025	50 uL	1.00 mL			

Reagent	Other Reagents:	Amount/Units	Lot#:

Preparation Batch Number(s): 320-100147 Test: PRC-L  
 Earliest Holding Time: 2-10-16 / 2-11-16

Sample List Tab		1 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> 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 <sup>st</sup> Level Reviewer	2 <sup>nd</sup> Level Reviewer
Date and time accurate and entered into TALS correctly		/	✓
All necessary 'batch information' complete and entered into TALS correctly		/	✓

1<sup>st</sup> Level Reviewer: 

Date: 2/10/16

2<sup>nd</sup> Level Reviewer: 

Date: 2/10/16

Comments: \_\_\_\_\_



# Subcontract Data



Prepared for: Test America

Project: 320-17154

# **Analytical Data Package (Level IV)**

Analysis: Low level PFOS and PFOA in water (Method 537)

Maxxam Job #: B630797

Maxxam Analytics International  
6740 Campobello Rd.  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)



## Table of Contents

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3. Analytical Results

3.1 Summary Report

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4. QA/QC Data

5. Initial Calibration

6. Continuing Calibration

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I hereby certify that to the best of my knowledge all analytical data presented in this report:

- Has been checked for completeness.
- Is accurate, legible and error free.
- Has been conducted in accordance with approved SOP's and that all deviations are clearly listed in the Case Narrative.
- This report has been generated in .pdf format.

Review Performed By:

Maxxam Analytics International  
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[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

# Glossary of Terms

- **Detection Limit (DL)** this can also be called **Method Detection Limit (MDL)**: The lowest concentration or amount of the target analyte that can be identified, measured, and reported with confidence that the analyte concentration is not a false positive value. (Clarification): The smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. At the DL, the false positive rate (Type I error) is 1%.
- **Limit of Detection (LOD)**: An estimate of the minimum amount of a substance that an analytical process can reliably detect. An LOD is analyte- and matrix-specific and may be laboratory-dependent. (Clarification): The smallest amount or concentration of a substance that must be present in a sample in order to be detected at a high level of confidence (99%). At the LOD, the false negative rate (Type II error) is 1%.
- **Limits of Quantitation (LOQ)** this can also be called **Reporting Detection Limit (RDL)**: The minimum levels, concentrations, or quantities of a target variable (e.g., target analyte) that can be reported with a specified degree of confidence. (Clarification): The lowest concentration that produces a quantitative result within specified limits of precision and bias. For DoD projects, the LOQ shall be set at or above the concentration of the lowest initial calibration standard.
- **Acceptance Criteria** are values used by the laboratory to determine that a process is in control.
- **Accuracy** is the degree of agreement of a measured value with the true or expected value.
- **Calibration Standards** are a set of solutions containing the analytes of interest at a specified concentration.
- **Calibration Verification Standard** consists of a calibration standard solution of intermediate concentration (mid-point initial calibration level) used to assess whether the initial calibration is still valid
- **Certified Reference Material** is a stable homogenous material that is certified by repetitive analysis from a supplier who is certified to generate said materials.

- **Internal Standard** a deuterated or  $^{13}\text{C}$ -labelled analyte that is added to a sample extract prior to instrumental analysis to compensate for injection variability.
- **Isomer** is a member of a group of compounds that differ from each other only in the locations of a specific number of common substituent atoms or groups of atoms on the parent compound.
- **Method Blank** is a laboratory control sample using reagents that are known to be free of contamination.
- **Precision** is the degree of agreement between the data generated from repetitive measurements under specific conditions.
- **Quality Assurance** is a system of activities whose purpose is to provide the producer or user of a product with the assurance that the product meets a defined standard of quality.
- **Quality Control** is the overall system of activities whose purpose is to control the quality of a product so that it meets the needs of the end user.
- **RSD** is the relative standard deviation.
- **Blank Spike** is a laboratory control sample that has been fortified with native analytes of interest.
- **Window Defining Mixture** is a solution containing only the earliest and latest eluting congeners within each homologous group of target analytes on a specified GC column.
- **RPD** or Relative Percent Difference. A measure used to compare duplicate sample analysis.
- **EMPC/NDR** – Peak detected does not meet ratio criteria and has resulted in a higher detection limit.



## 1.0 Project Narrative

Maxxam Analytics International  
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Sample Analysis

All samples were analyzed on QC batch 4386408 (2016/02/19). Recoveries of the Matrix Spike (MS) and/or Matrix Spike Duplicate (MSD) were above the upper control limit for Perfluorobutanesulfonate (PFBS), Perfluorohexanesulfonate (PFHxS) and Perfluoroheptanoic acid (PFHpA). Laboratory spiked water (LCS) resulted in satisfactory recovery of the compounds of interest. When considered together, these QC data suggest that matrix interferences may be biasing the data high. For results that were U-flagged, this potential bias has no impact.

Internal Standard Responses

Isotopically labeled  $^{13}\text{C}_4$ -Perfluoroheptanoic acid (MPFHpA) is used as an internal standard to quantify native Perfluoroheptanoic acid (PFHpA). The instrument response observed for this labeled compound was below the defined lower control limit (LCL) for the following sample:

BVX827 OF-RW37-0216

The instrument response for the injection standard ( $^{13}\text{C}_6$ -Perfluorohexanoic acid,  $^{13}\text{C}_6$ -PFHxA) fell within the required tolerance limits. This indicates that the response observed for this isotopically labeled internal standard was not a result of poor or inconsistent sample introduction into the liquid chromatograph/tandem mass spectrometer (LC/MS/MS).

Data Qualifiers

In the Results of Analyses, U-flags are applied to results that are less than the DL (MDL). J-flags are applied to results that are less than the RDL (LOQ) but greater than the DL (MDL). Due to limitations in LIMS, the results cannot be U-flagged to the LOD. The LODs for each analyte are presented in the table below.

Parameter	MDL (ng/L)	LOD (ng/L)	LOQ (ng/L)
Perfluorobutane sulfonate (PFBS)	0.27	1.0	2.0
Perfluorohexane sulfonate (PFHxS)	0.40	1.0	2.0
Perfluoroheptanoic acid (PFHpA)	0.39	1.0	2.0
Perfluorooctanoic acid (PFOA)	0.39	1.0	2.0
Perfluorooctane sulfonate (PFOS)	0.30	1.0	2.0
Perfluorononanoic acid (PFNA)	0.33	1.0	2.0

**Sin Chii Chia, B.Sc.**

schia@maxxam.ca

Office 905 817 5700

**PROJECT NARRATIVE**

**Maxxam Analytics**  
**Client Project #: 320-17154**



**Client: TestAmerica**  
**Client Project: 320-17154**

**I. SAMPLE RECEIPT/ANALYSIS**

a) Sample Listing

<b>Maxxam ID</b>	<b>Client Sample ID</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Date Prepped</b>	<b>Date Run</b>	<b>Initial Calibration</b>
<b>Low level PFOS and PFOA in water</b>						
BVX826	OF-RW09-0216	2016/02/03	2016/02/13	2016/02/18	2016/02/19	2016/02/19
BVX827	OF-RW37-0216	2016/02/03	2016/02/13	2016/02/18	2016/02/19	2016/02/19
BVX828	OF-RW11-0216	2016/02/03	2016/02/13	2016/02/18	2016/02/19	2016/02/19
BVX829	OF-RW28-0216	2016/02/03	2016/02/13	2016/02/18	2016/02/19	2016/02/19
BVX830	OF-RW67-0216	2016/02/03	2016/02/13	2016/02/18	2016/02/19	2016/02/19

Run Date is defined as the date of injection of the last calibration standard (12 hours or less) prior to the samples analyzed within that run sequence. Therefore the time of calibration injection that defines the run date is always within 12 hours of the time of sample injection.

b) Shipping Problems: none encountered

c) Documentation Problems: none encountered

**II. SAMPLE PREP:**

No problems encountered

**III. SAMPLE ANALYSIS:**

See also comments within the appropriate Certificate of Analysis

a) Hold Times: all within recommended hold times

b) Instrument Calibration: all within control limits

c) Quality Control: All applicable QC meets control criteria, except where otherwise noted.

d) All analytes requiring manual intergration(s) are noted on the sample chromatograms

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for other than the conditions detailed above.

In addition, I certify, that to the best of my knowledge and belief, the data as reported are true and accurate. Release of the data contained in this data package has been authorized by the cognizant laboratory official or his/her designee, as verified by this signature.

*M Di Grazia*

2016/03/08

Date



## 2. Sample Management Records

Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)





## 2.1 Sample Custody

Maxxam Analytics International  
6740 Campobello Rd  
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L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

# Chain of Custody Record

Temperature on Receipt \_\_\_\_\_

Drinking Water? Yes  No



320-17154 Chain of Custody

CTD WEF G

Due 2/22

Project Manager: [Redacted]  
 Telephone Number (Area Code)/Fax Number: [Redacted]  
 Site Contact: [Redacted]  
 Carrier/Waybill Number: **FEDEX**

Date: 02/03/16  
 Chain of Custody Number: 283602  
 Page 1 of 1

Analysis (Attach list if more space is needed):  
 Select PFGS

Special Instructions/Conditions of Receipt:

Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time	Matrix			Containers & Preservatives						Special Instructions/Conditions of Receipt		
			AP	Sed	Sol	Unpres.	HRSOA	HNO3	HCl	NH4OH	ZnAc		HORN	
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>16:10</del>												
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>16:45</del>												
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>17:20</del>												
<del>DF-RW137-0216</del>	<del>02/03/16</del>	<del>17:28</del>												
<del>DF-RW111-0216</del>	<del>02/03/16</del>	<del>16:04</del>												
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>16:10</del>												
<del>DF-RW28-0216</del>	<del>02/03/16</del>	<del>17:41</del>												
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>17:43</del>												
<del>DF-RW167-0216</del>	<del>02/03/16</del>	<del>18:19</del>												
<del>DF-RW109-0216</del>	<del>02/03/16</del>	<del>18:21</del>												

13-Feb-16 13:40  
 Hongmei Zhao (Grace)  
 B630797

RGN ENV-951

(A fee may be assessed if samples are retained longer than 1 month)

Possible Hazard Identification:  
 Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown  Other

Turn Around Time Required:  
 24 Hours  48 Hours  7 Days  14 Days  21 Days  Other

Sample Disposal:  
 Return To Client  Disposal By Lab  Archive For \_\_\_\_\_ Months

QC Requirements (Specify):

1. Released By: [Signature] Date: 02/03/16 Time: 19:30  
 2. Received By: [Signature] Date: 2/3/16 Time: 9:45  
 3. Received By: [Signature] Date: 2016/02/13 Time: 13:40

Comments: 0.2.0 REFER TO ACTE

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



### **3. Analytical Results**

Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)



## 3.1 Summary Report

Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)



Your Project #: 320-17154  
Your C.O.C. #: 283602

**Attention:PFC Reporting Group**

TestAmerica  
Sacramento  
880 Riverside Parkway  
West Sacramento, CA  
USA 95605

**Report Date: 2016/03/08**  
Report #: R3920386  
Version: 3 - Revision

**CERTIFICATE OF ANALYSIS – REVISED REPORT**

**MAXXAM JOB #: B630797**  
**Received: 2016/02/13, 13:40**

Sample Matrix: Water  
# Samples Received: 5

Analyses	Date		Laboratory Method	Reference
	Quantity Extracted	Date Analyzed		
Low level PFOS and PFOA in water	5	2016/02/18	2016/02/19 CAM SOP-00894	EPA 537 m

Reference Method suffix "m" indicates test methods incorporate validated modifications from specific reference methods to improve performance.  
\* RPDs calculated using raw data. The rounding of final results may result in the apparent difference.

**Encryption Key**

Please direct all questions regarding this Certificate of Analysis to your Project Manager.  
Melissa DiGrazia, Project Manager - ATUT  
Email: MDiGrazia@maxxam.ca  
Phone# (905) 817-5700

=====

Maxxam has procedures in place to guard against improper use of the electronic signature and have the required "signatories", as per section 5.10.2 of ISO/IEC 17025:2005(E), signing the reports. For Service Group specific validation please refer to the Validation Signature Page.  
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**RESULTS OF ANALYSES OF WATER**

Maxxam ID		BVX826	BVX827	BVX828	BVX829	BVX830			
Sampling Date		2016/02/03 16:45	2016/02/03 17:28	2016/02/03 16:04	2016/02/03 17:41	2016/02/03 18:19			
COC Number		283602	283602	283602	283602	283602			
	UNITS	OF-RW09-0216	OF-RW37-0216	OF-RW11-0216	OF-RW28-0216	OF-RW67-0216	RDL	MDL	QC Batch

Miscellaneous Parameters									
Perfluorobutane Sulfonate (PFBS)	ng/L	0.27 U	0.79 J	0.27 U	0.27 U	0.97 J	2.0	0.27	4386408
Perfluoroheptanoic Acid (PFHpA)	ng/L	0.39 U	0.46 J	0.39 U	0.39 U	0.39 U	2.0	0.39	4386408
Perfluorohexane Sulfonate (PFHxS)	ng/L	0.71 J	3.3	0.75 J	0.75 J	0.40 U	2.0	0.40	4386408
Perfluoro-n-Octanoic Acid (PFOA)	ng/L	0.41 J	7.2	0.39 U	0.39 U	0.39 U	2.0	0.39	4386408
Perfluorononanoic Acid (PFNA)	ng/L	0.33 U	0.33 U	0.33 U	0.37 J	0.33 U	2.0	0.33	4386408
Perfluorooctane Sulfonate (PFOS)	ng/L	0.32 J	0.81 J	0.87 J	0.30 U	0.30 U	2.0	0.30	4386408

Surrogate Recovery (%)									
13C4-Perfluoroheptanoic acid	%	54	52	55	54	48 (1)	N/A	N/A	4386408
13C4-Perfluorooctanesulfonate	%	59	64	54	59	51	N/A	N/A	4386408
13C4-Perfluorooctanoic acid	%	58	57	59	54	51	N/A	N/A	4386408
13C5-Perfluorononanoic acid	%	62	65	61	57	55	N/A	N/A	4386408
18O2-Perfluorohexanesulfonate	%	69	64	60	50	56	N/A	N/A	4386408

RDL = Reportable Detection Limit  
 QC Batch = Quality Control Batch  
 N/A = Not Applicable  
 (1) Surrogate recovery was below the defined lower control limit (LCL). Laboratory spiked water resulted in satisfactory recovery of the surrogate. When considered together, these QC data suggest that matrix interferences may be biasing the data low. Because quantitation is performed using isotope dilution techniques, any losses of the native compound that may occur during any of the sample preparation, extraction, cleanup or determinative steps will be mirrored by a similar loss of the labeled standard, and as such can be accounted for and corrected. Therefore, the quantification of these target compounds is not affected by the low surrogate recovery.

**TEST SUMMARY**

**Maxxam ID:** BVX826  
**Sample ID:** OF-RW09-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX827  
**Sample ID:** OF-RW37-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX828  
**Sample ID:** OF-RW11-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX829  
**Sample ID:** OF-RW28-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

**Maxxam ID:** BVX830  
**Sample ID:** OF-RW67-0216  
**Matrix:** Water

**Collected:** 2016/02/03  
**Shipped:**  
**Received:** 2016/02/13

Test Description	Instrumentation	Batch	Extracted	Date Analyzed	Analyst
Low level PFOS and PFOA in water	LCMS	4386408	2016/02/18	2016/02/19	Colm McNamara

### GENERAL COMMENTS

Report revised to remove qualifier legend

Perfluorinated Compounds (PFCs):

Perfluoroheptanoic acid (PFHpA) MDL = 0.39, LOD = 1.0, LOQ = 2.0  
Perfluorooctanoic acid (PFOA) MDL = 0.39, LOD = 1.0, LOQ = 2.0  
Perfluorononanoic acid (PFNA) MDL = 0.33, LOD = 1.0, LOQ = 2.0  
Perfluorobutane sulfonate (PFBS) MDL = 0.27, LOD = 1.0, LOQ = 2.0  
Perfluorohexane sulfonate (PFHxS) MDL = 0.40, LOD = 1.0, LOQ = 2.0  
Perfluorooctane sulfonate (PFOS) MDL = 0.30, LOD = 1.0, LOQ = 2.0  
All Units are in ng/L

**Results relate only to the items tested.**



**QUALITY ASSURANCE REPORT**

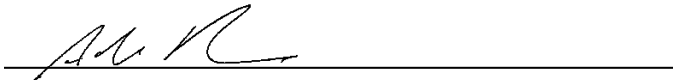
QA/QC Batch	Init	QC Type	Parameter	Date Analyzed	Value	Recovery	UNITS	QC Limits
4386408	CM5	Matrix Spike	13C4-Perfluoroheptanoic acid	2016/02/19		50	%	50 - 130
			13C4-Perfluorooctanesulfonate	2016/02/19		59	%	50 - 130
			13C4-Perfluorooctanoic acid	2016/02/19		58	%	50 - 130
			13C5-Perfluorononanoic acid	2016/02/19		61	%	50 - 130
			18O2-Perfluorohexanesulfonate	2016/02/19		53	%	50 - 130
			Perfluorobutane Sulfonate (PFBS)	2016/02/19		157 (1)	%	70 - 130
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19		133 (1)	%	70 - 130
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19		148 (1)	%	70 - 130
			Perfluorononanoic Acid (PFNA)	2016/02/19		120	%	70 - 130
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19		125	%	70 - 130
			Perfluorooctane Sulfonate (PFOS)	2016/02/19		116	%	70 - 130
4386408	CM5	RPD	Perfluorobutane Sulfonate (PFBS)	2016/02/19	4.2		%	30
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19	6.2		%	30
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19	11		%	30
			Perfluorononanoic Acid (PFNA)	2016/02/19	14		%	30
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19	8.5		%	30
			Perfluorooctane Sulfonate (PFOS)	2016/02/19	6.4		%	30
4386408	CM5	Spiked Blank	13C4-Perfluoroheptanoic acid	2016/02/19		67	%	50 - 130
			13C4-Perfluorooctanesulfonate	2016/02/19		77	%	50 - 130
			13C4-Perfluorooctanoic acid	2016/02/19		70	%	50 - 130
			13C5-Perfluorononanoic acid	2016/02/19		73	%	50 - 130
			18O2-Perfluorohexanesulfonate	2016/02/19		83	%	50 - 130
			Perfluorobutane Sulfonate (PFBS)	2016/02/19		107	%	70 - 130
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19		121	%	70 - 130
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19		115	%	70 - 130
			Perfluorononanoic Acid (PFNA)	2016/02/19		118	%	70 - 130
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19		124	%	70 - 130
			Perfluorooctane Sulfonate (PFOS)	2016/02/19		112	%	70 - 130
4386408	CM5	Method Blank	13C4-Perfluoroheptanoic acid	2016/02/19		74	%	50 - 130
			13C4-Perfluorooctanesulfonate	2016/02/19		82	%	50 - 130
			13C4-Perfluorooctanoic acid	2016/02/19		74	%	50 - 130
			13C5-Perfluorononanoic acid	2016/02/19		76	%	50 - 130
			18O2-Perfluorohexanesulfonate	2016/02/19		86	%	50 - 130
			Perfluorobutane Sulfonate (PFBS)	2016/02/19	0.27 U, MDL=0.27		ng/L	
			Perfluoroheptanoic Acid (PFHpA)	2016/02/19	0.39 U, MDL=0.39		ng/L	
			Perfluorohexane Sulfonate (PFHxS)	2016/02/19	0.40 U, MDL=0.40		ng/L	
			Perfluorononanoic Acid (PFNA)	2016/02/19	0.33 U, MDL=0.33		ng/L	
			Perfluoro-n-Octanoic Acid (PFOA)	2016/02/19	0.39 U, MDL=0.39		ng/L	

**QUALITY ASSURANCE REPORT(CONT'D)**

QA/QC Batch	Init	QC Type	Parameter	Date Analyzed	Value	Recovery	UNITS	QC Limits
			Perfluorooctane Sulfonate (PFOS)	2016/02/19	0.30 U, MDL=0.30		ng/L	
<p>Duplicate: Paired analysis of a separate portion of the same sample. Used to evaluate the variance in the measurement.</p> <p>Matrix Spike: A sample to which a known amount of the analyte of interest has been added. Used to evaluate sample matrix interference.</p> <p>Spiked Blank: A blank matrix sample to which a known amount of the analyte, usually from a second source, has been added. Used to evaluate method accuracy.</p> <p>Method Blank: A blank matrix containing all reagents used in the analytical procedure. Used to identify laboratory contamination.</p> <p>Surrogate: A pure or isotopically labeled compound whose behavior mirrors the analytes of interest. Used to evaluate extraction efficiency.</p> <p>(1) Recovery or RPD for this parameter is outside control limits. The overall quality control for this analysis meets acceptability criteria.</p>								

**VALIDATION SIGNATURE PAGE**

The analytical data and all QC contained in this report were reviewed and validated by the following individual(s).



Adam Robinson, Supervisor, LC/MS/MS

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Maxxam has procedures in place to guard against improper use of the electronic signature and have the required "signatories", as per section 5.10.2 of ISO/IEC 17025:2005(E), signing the reports. For Service Group specific validation please refer to the Validation Signature Page.



**METHOD 537**  
**DETERMINATION OF SELECTED PERFLUORINATED ALKYL**  
**ACIDS IN DRINKING WATER BY SOLID PHASE EXTRACTION**  
**AND LIQUID CHROMATOGRAPHY/TANDEM MASS SPECTROMETRY**  
**(LC/MS/MS)**

Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)



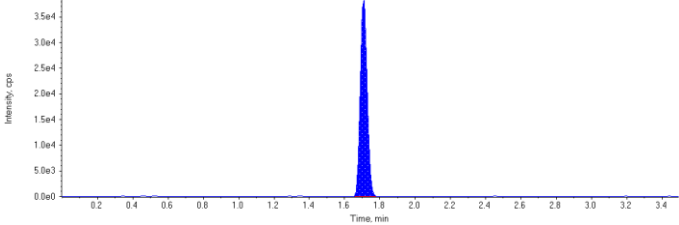
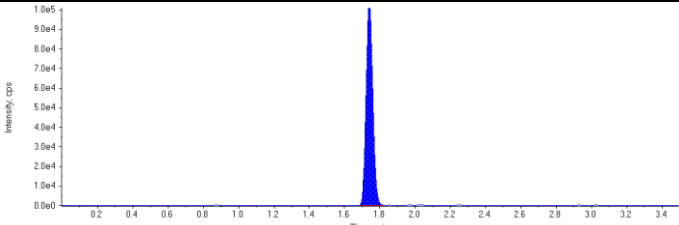
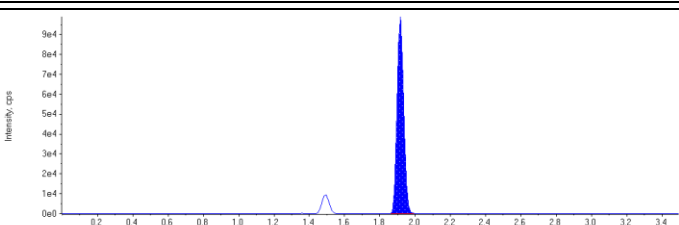
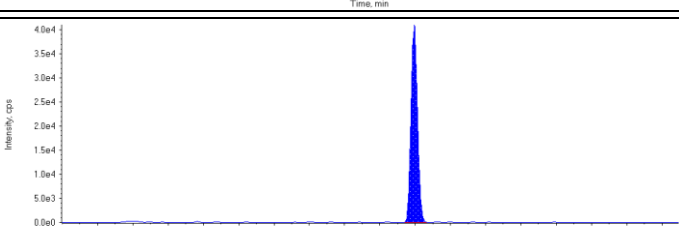
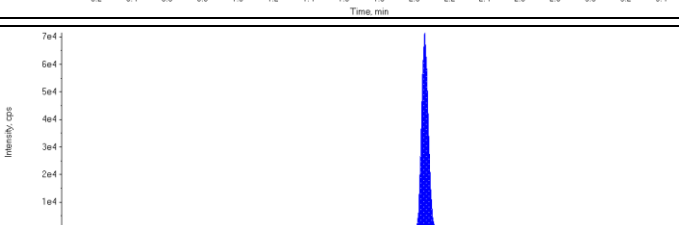
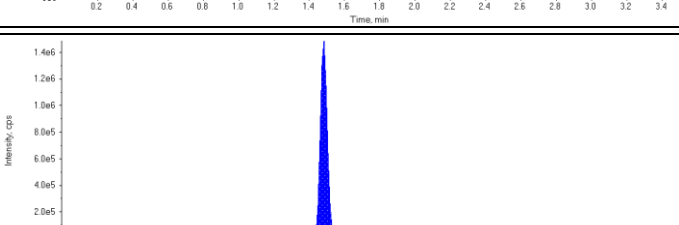
## 3.2 Sample Chromatograms

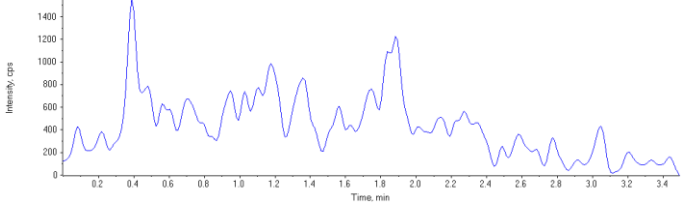
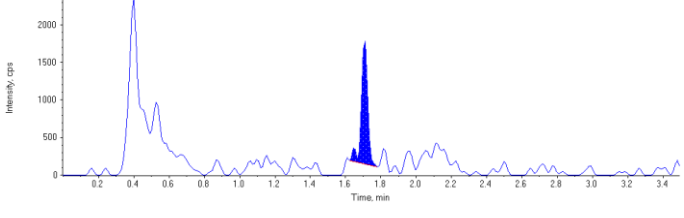
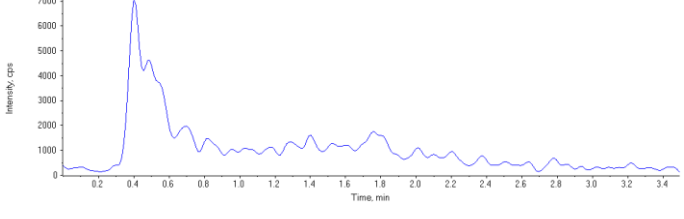
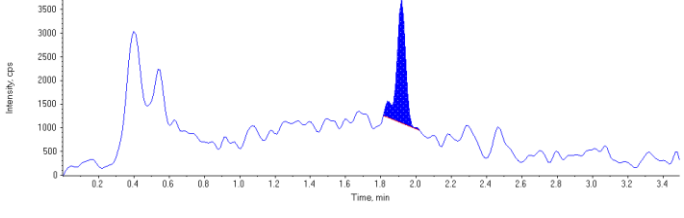
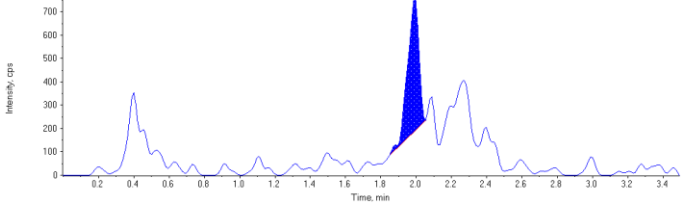
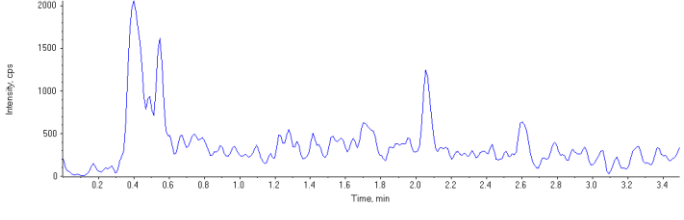
Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

<b>Sample Name</b>	<b>4386408~BVX826-01</b>	<b>Injection Vial</b>	28
<b>Sample ID</b>	4386408~BVX826-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:18:03 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
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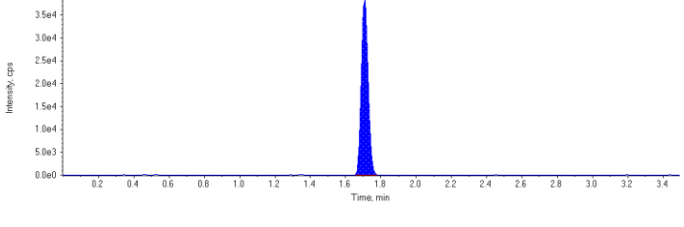
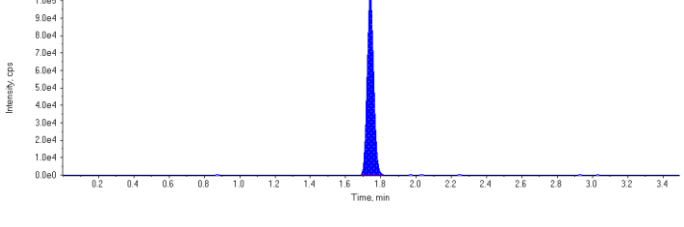
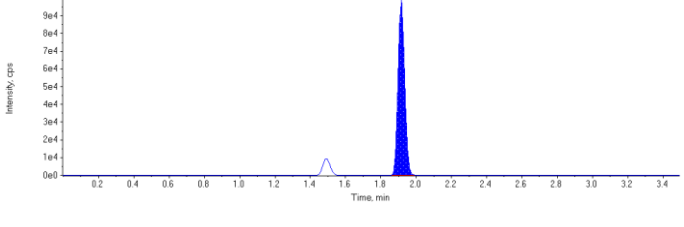
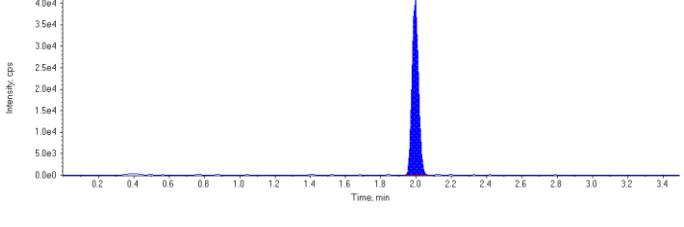
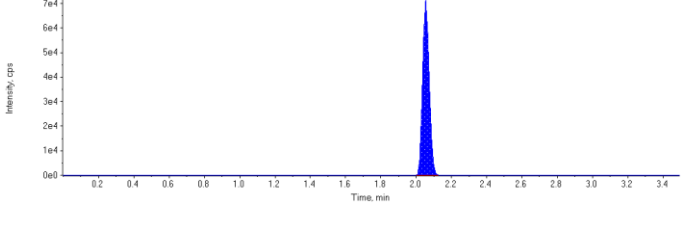
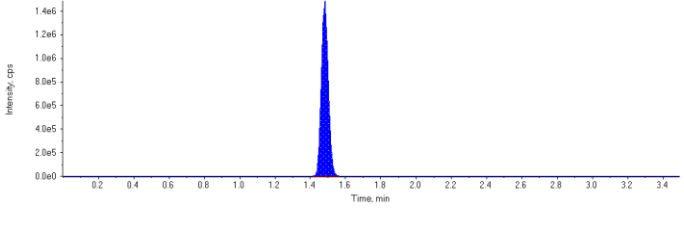
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	104000.	1.71	1.00	-
MPFHpA	272000.	1.74	1.00	-
MPFOA	256000.	1.92	1.00	-
MPFOS	105000.	1.99	1.00	-
MPFNA	186000.	2.05	1.00	-
13C6-PFHxA IS	4080000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	0	0.00	N/A	N/A	N/A
PFHxS 1	4150	1.71	N/A	0.713	N/A
PFHpA 1	0	0.00	N/A	N/A	N/A
PFOA 1	9130	1.92	N/A	0.412	N/A
PFOS 1	2530	1.99	N/A	0.320	N/A
PFNA 1	0	0.00	N/A	N/A	N/A
18O2-PFHxS	104000	1.71	N/A	69.2	N/A
13C4-PFHpA	272000	1.74	N/A	53.8	N/A
13C4-PFOA	256000	1.92	N/A	58.1	N/A
13C4-PFOS	105000	1.99	N/A	59.3	N/A
13C5-PFNA	186000	2.05	N/A	61.7	N/A
13C6-PFHxA	4080000	1.48	N/A	102.	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.15) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 0.713 µg/L</p> <p>Area Ratio: 0.0399</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.75) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 0.412 µg/L</p> <p>Area Ratio: 0.0356</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 0.320 µg/L</p> <p>Area Ratio: 0.0242</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

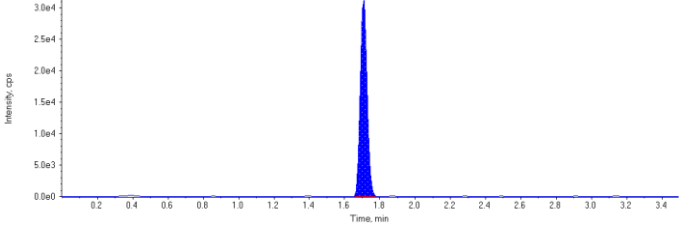
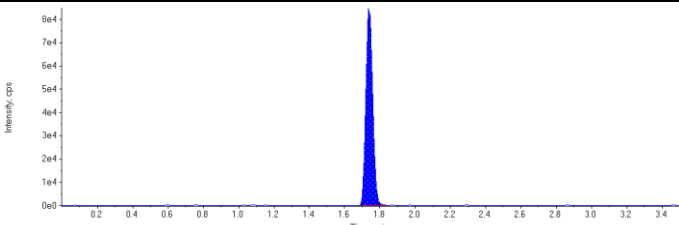
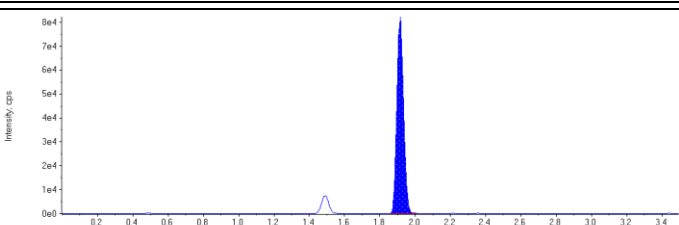
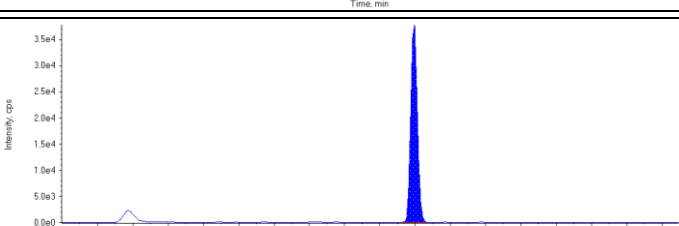
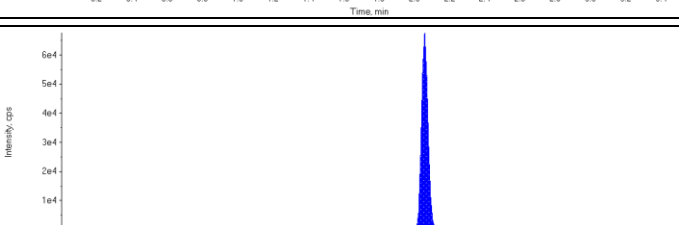
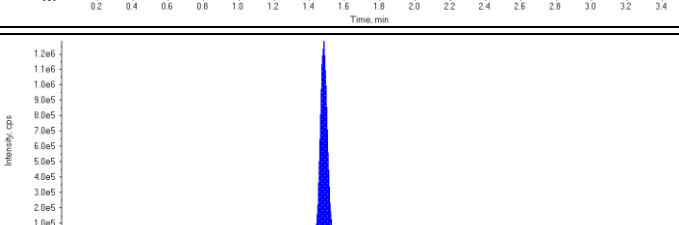


<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 69.2 µg/L</p> <p>Area Ratio: 0.0255</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 53.8 µg/L</p> <p>Area Ratio: 0.0665</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 58.1 µg/L</p> <p>Area Ratio: 0.0627</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 59.3 µg/L</p> <p>Area Ratio: 0.0256</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 61.7 µg/L</p> <p>Area Ratio: 0.0454</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 102. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

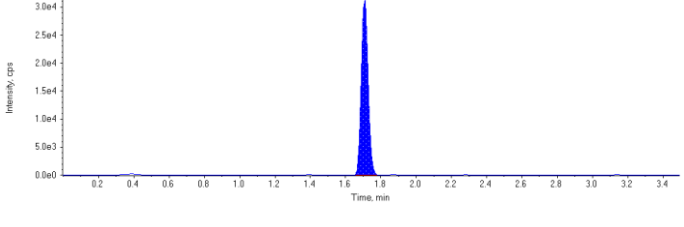
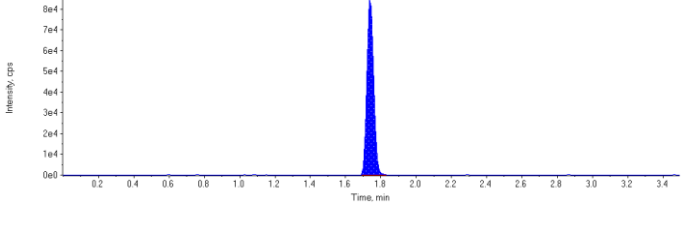
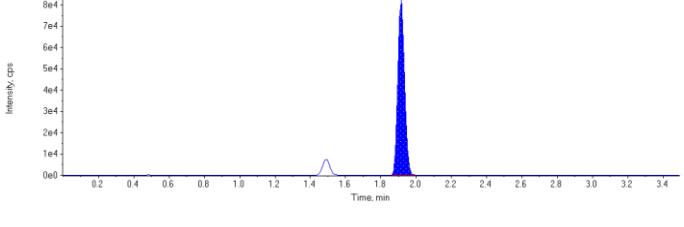
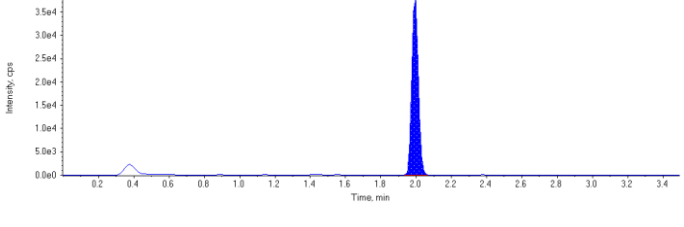
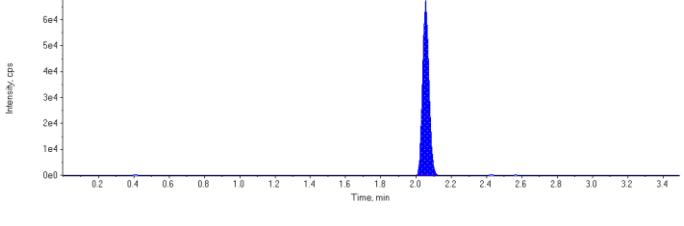
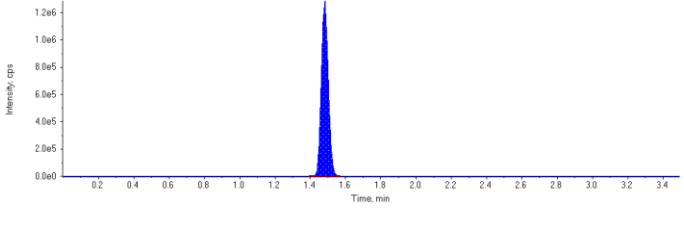
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<b>Sample ID</b>	4386408~BVX827-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:23:09 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	84400.	1.71	1.00	-
MPFHpA	229000.	1.74	1.00	-
MPFOA	217000.	1.91	1.00	-
MPFOS	98200.	1.99	1.00	-
MPFNA	170000.	2.05	1.00	-
13C6-PFHxA IS	3560000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	5820	1.16	N/A	0.792	N/A
PFHxS 1	74800	1.71	N/A	3.26	N/A
PFHpA 1	11800	1.72	N/A	0.455	N/A
PFOA 1	250000	1.92	N/A	7.15	N/A
PFOS 1	11200	1.95	N/A	0.812	N/A
PFNA 1	0	0.00	N/A	N/A	N/A
18O2-PFHxS	84400	1.71	N/A	64.4	N/A
13C4-PFHpA	229000	1.74	N/A	51.9	N/A
13C4-PFOA	217000	1.91	N/A	56.6	N/A
13C4-PFOS	98200	1.99	N/A	63.9	N/A
13C5-PFNA	170000	2.05	N/A	65.0	N/A
13C6-PFHxA	3560000	1.48	N/A	89.1	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

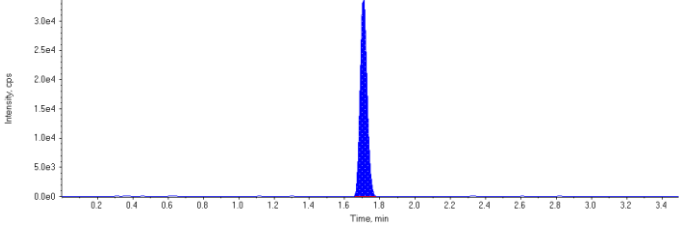
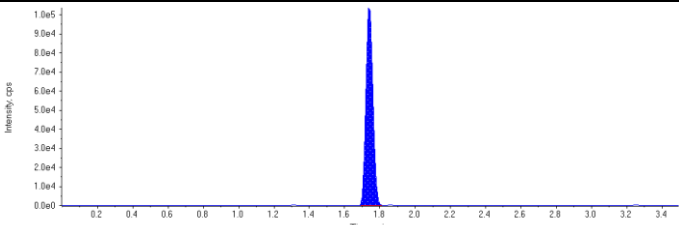
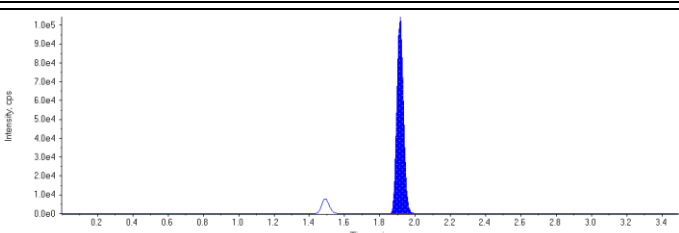
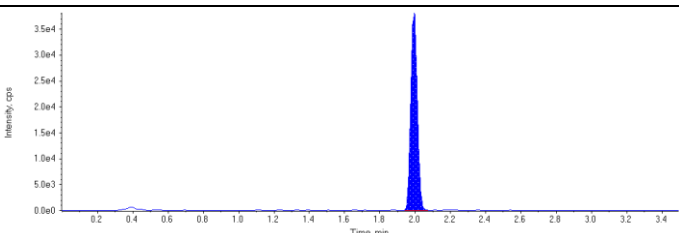
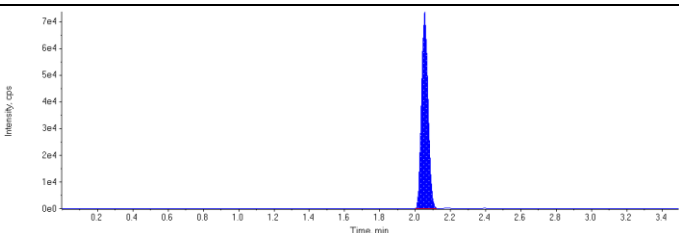
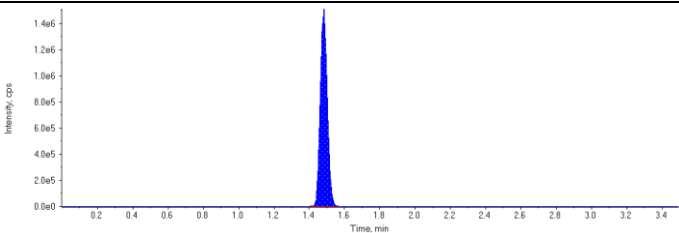
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 0.792 µg/L</p> <p>Area Ratio: 0.0689</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 3.26 µg/L</p> <p>Area Ratio: 0.886</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.72 (1.75) min</p> <p>Calculated Conc: 0.455 µg/L</p> <p>Area Ratio: 0.0515</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 7.15 µg/L</p> <p>Area Ratio: 1.15</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.95 (1.97) min</p> <p>Calculated Conc: 0.812 µg/L</p> <p>Area Ratio: 0.114</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

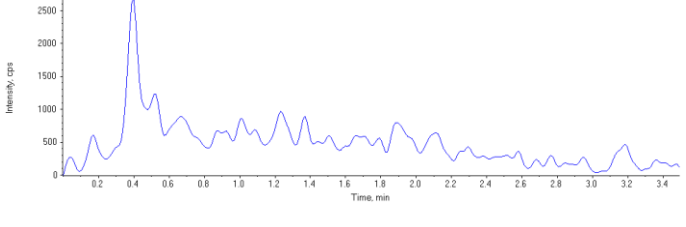
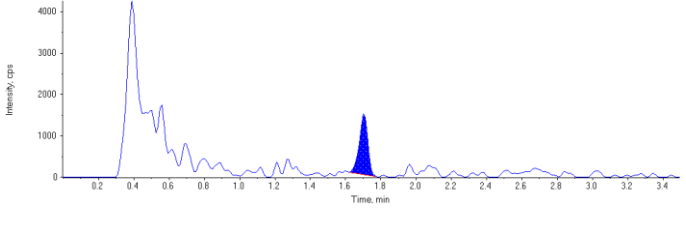
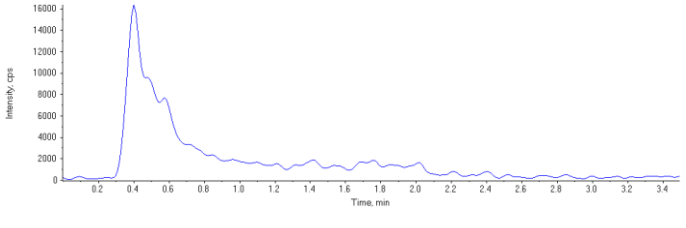
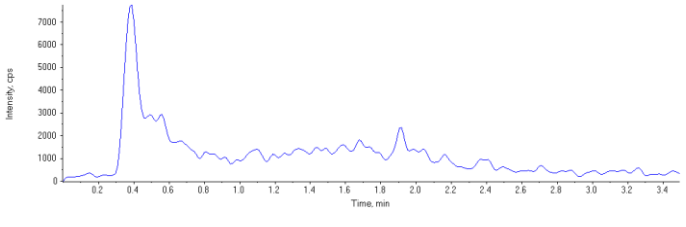
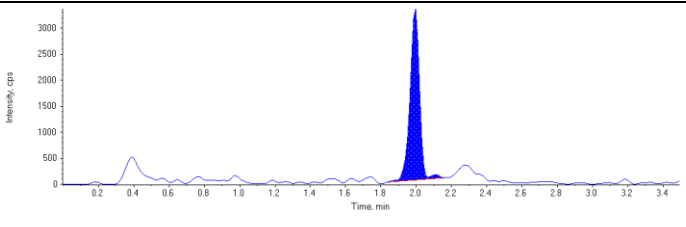
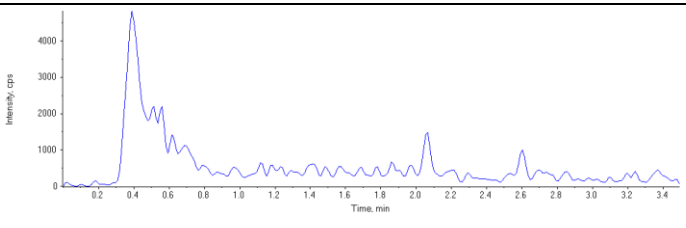
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 64.4 µg/L</p> <p>Area Ratio: 0.0237</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 51.9 µg/L</p> <p>Area Ratio: 0.0643</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 56.6 µg/L</p> <p>Area Ratio: 0.0611</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 63.9 µg/L</p> <p>Area Ratio: 0.0276</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 65.0 µg/L</p> <p>Area Ratio: 0.0479</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 89.1 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

<b>Sample Name</b>	<b>4386408~BVX828-01</b>	<b>Injection Vial</b>	30
<b>Sample ID</b>	4386408~BVX828-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:28:14 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
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<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

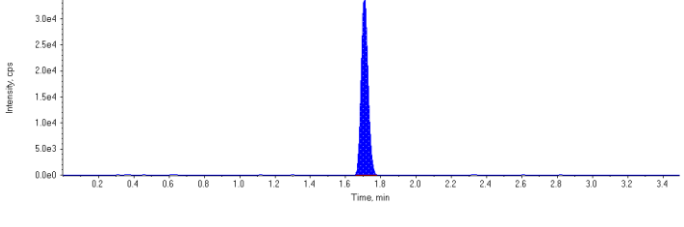
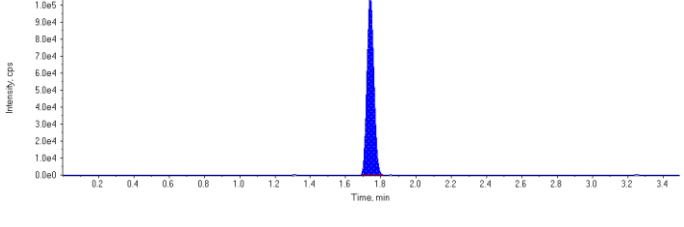
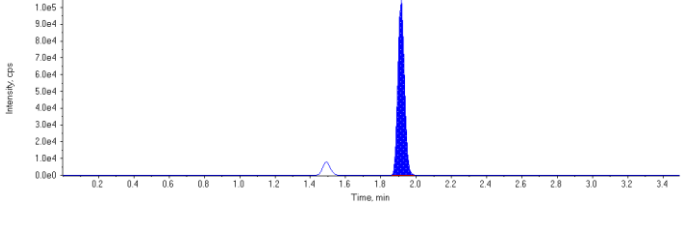
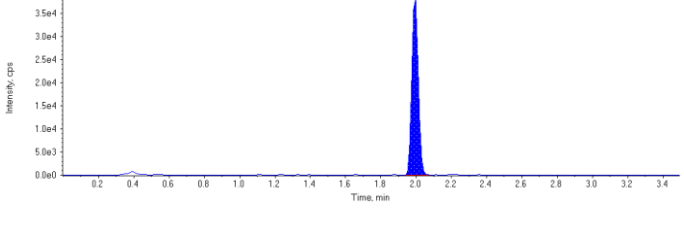
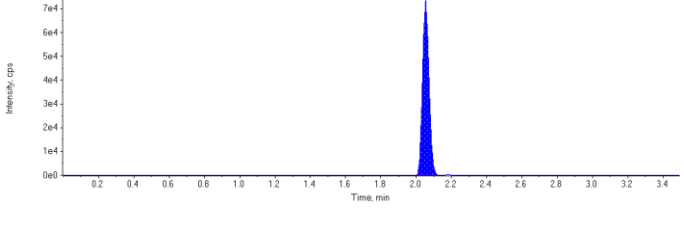
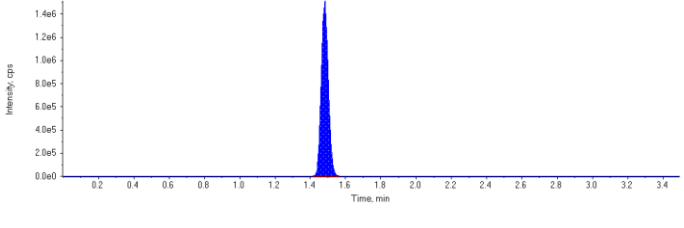
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	92800.	1.71	1.00	-
MPFHpA	289000.	1.74	1.00	-
MPFOA	269000.	1.91	1.00	-
MPFOS	98300.	1.99	1.00	-
MPFNA	188000.	2.05	1.00	-
13C6-PFHxA IS	4220000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	0	0.00	N/A	N/A	N/A
PFHxS 1	4910	1.70	N/A	0.752	N/A
PFHpA 1	0	0.00	N/A	N/A	N/A
PFOA 1	0	0.00	N/A	N/A	N/A
PFOS 1	12300	1.99	N/A	0.871	N/A
PFNA 1	0	0.00	N/A	N/A	N/A
18O2-PFHxS	92800	1.71	N/A	59.8	N/A
13C4-PFHpA	289000	1.74	N/A	55.3	N/A
13C4-PFOA	269000	1.91	N/A	59.2	N/A
13C4-PFOS	98300	1.99	N/A	53.9	N/A
13C5-PFNA	188000	2.05	N/A	60.5	N/A
13C6-PFHxA	4220000	1.48	N/A	106.	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.15) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.70 (1.68) min</p> <p>Calculated Conc: 0.752 µg/L</p> <p>Area Ratio: 0.0529</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.75) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.92) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 0.871 µg/L</p> <p>Area Ratio: 0.125</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

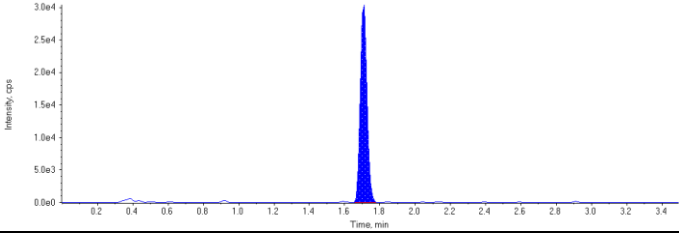
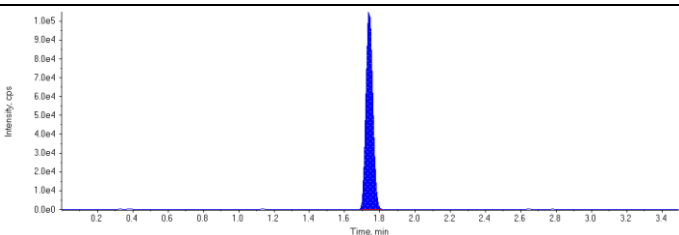
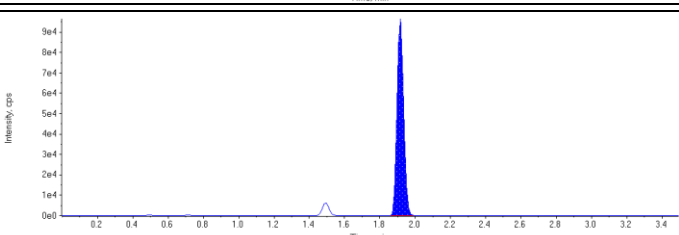
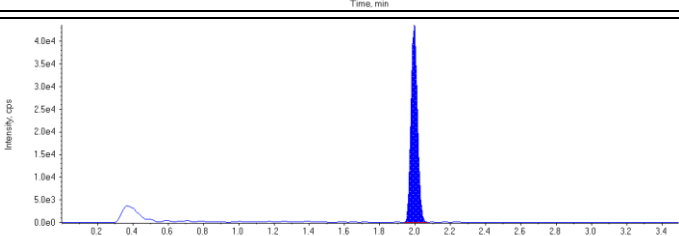
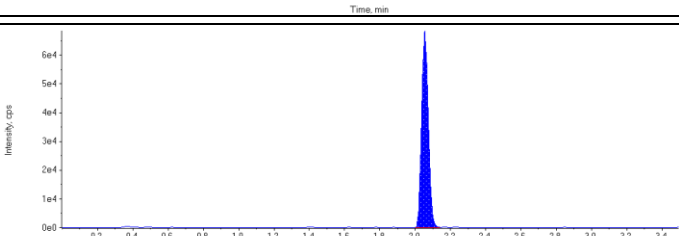
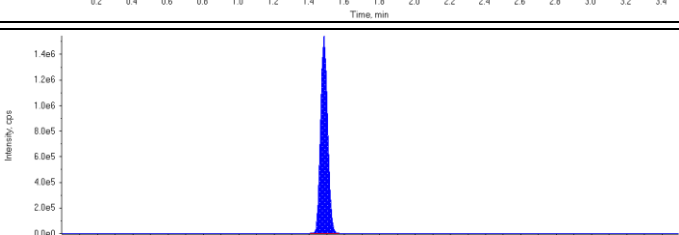


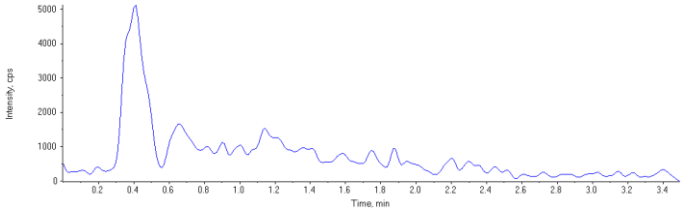
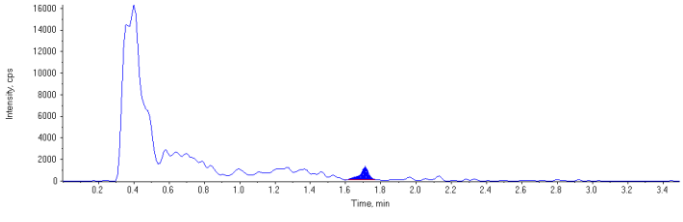
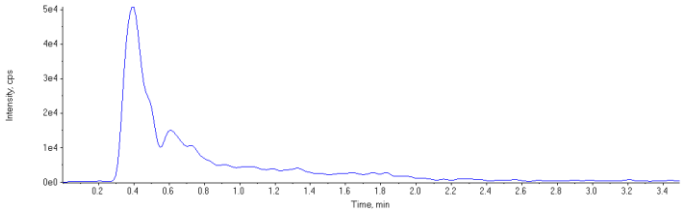
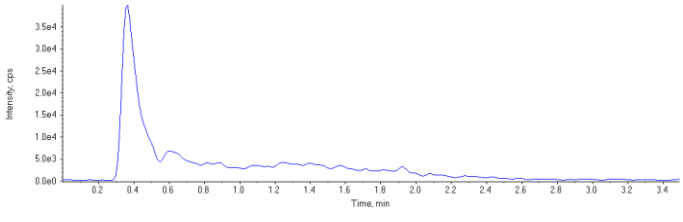
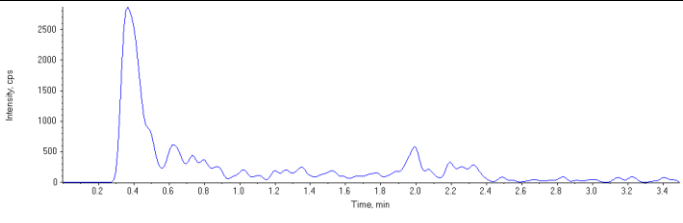
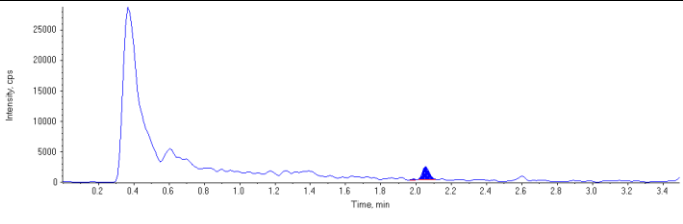
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 59.8 µg/L</p> <p>Area Ratio: 0.0220</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 55.3 µg/L</p> <p>Area Ratio: 0.0685</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 59.2 µg/L</p> <p>Area Ratio: 0.0639</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 53.9 µg/L</p> <p>Area Ratio: 0.0233</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 60.5 µg/L</p> <p>Area Ratio: 0.0445</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 106. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

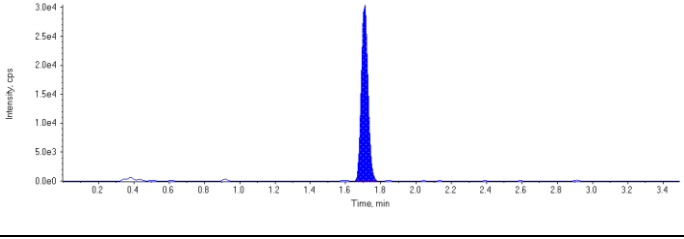
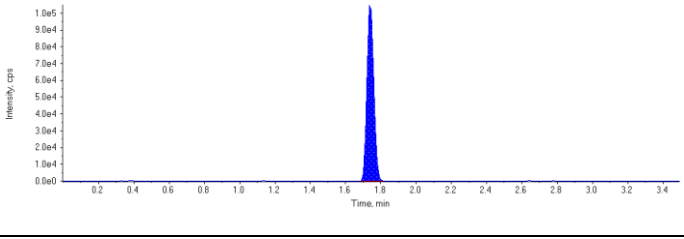
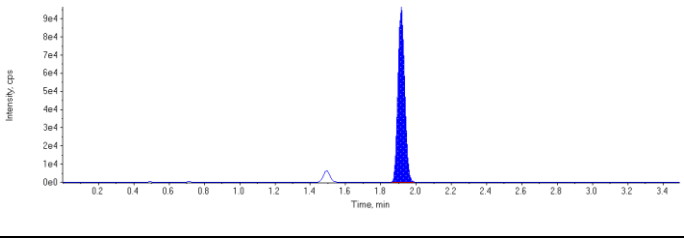
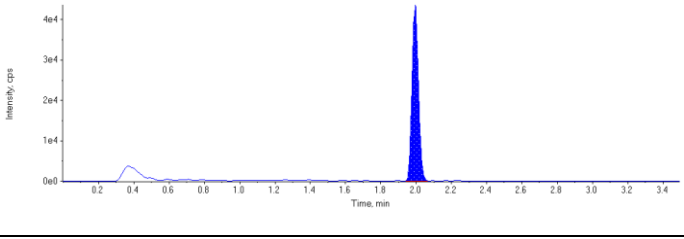
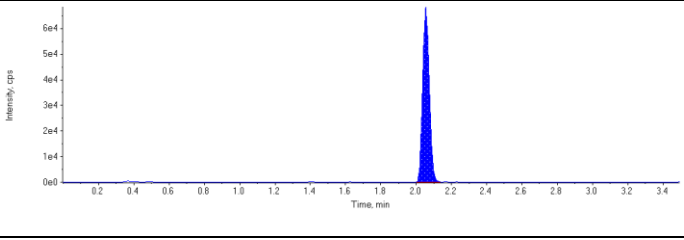
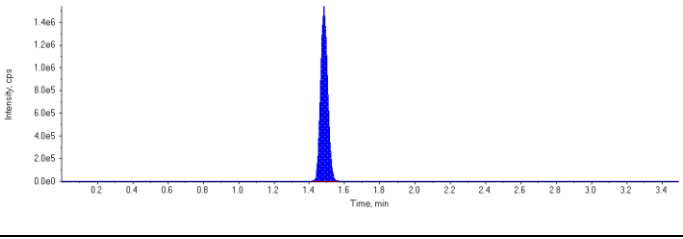
<b>Sample Name</b>	<b>4386408~BVX829-01</b>	<b>Injection Vial</b>	31
<b>Sample ID</b>	4386408~BVX829-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:33:19 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	79900.	1.71	1.00	-
MPFHpA	286000.	1.74	1.00	-
MPFOA	252000.	1.92	1.00	-
MPFOS	110000.	1.99	1.00	-
MPFNA	180000.	2.05	1.00	-
13C6-PFHxA IS	4310000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	0	0.00	N/A	N/A	N/A
PFHxS 1	4270	1.71	N/A	0.753	N/A
PFHpA 1	0	0.00	N/A	N/A	N/A
PFOA 1	0	0.00	N/A	N/A	N/A
PFOS 1	0	0.00	N/A	N/A	N/A
PFNA 1	6260	2.05	N/A	0.366	N/A
18O2-PFHxS	79900	1.71	N/A	50.3	N/A
13C4-PFHpA	286000	1.74	N/A	53.6	N/A
13C4-PFOA	252000	1.92	N/A	54.2	N/A
13C4-PFOS	110000	1.99	N/A	59.2	N/A
13C5-PFNA	180000	2.05	N/A	56.7	N/A
13C6-PFHxA	4310000	1.48	N/A	108.	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

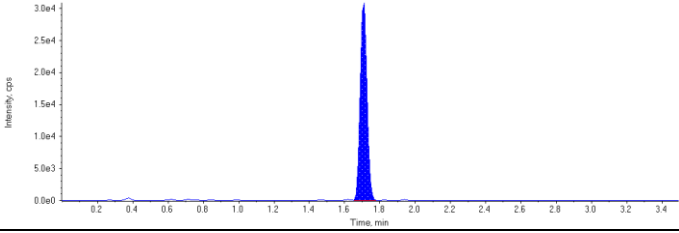
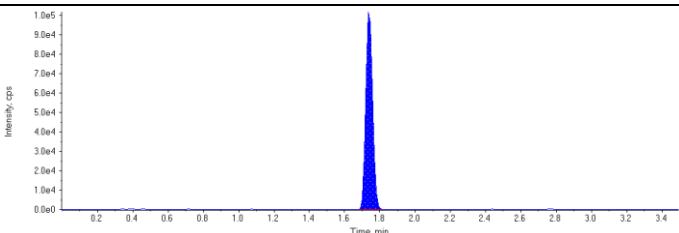
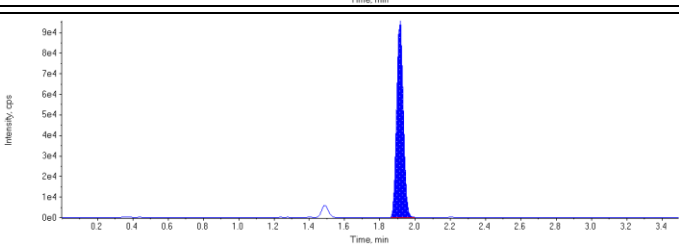
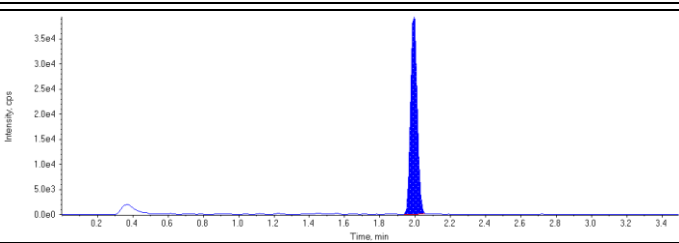
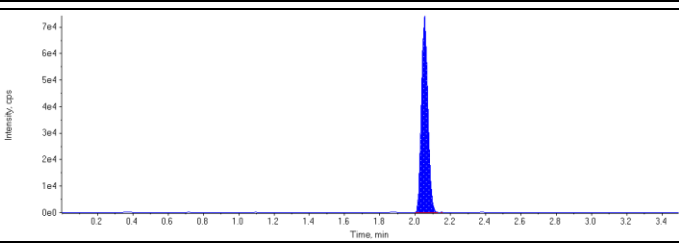
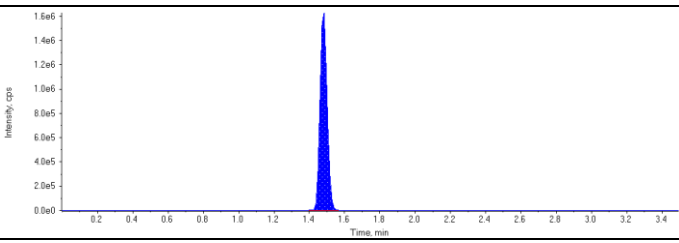
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.15) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 0.753 µg/L</p> <p>Area Ratio: 0.0534</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.75) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.92) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.97) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 0.366 µg/L</p> <p>Area Ratio: 0.0348</p> <p>Sample Type: (Unknown)</p>	

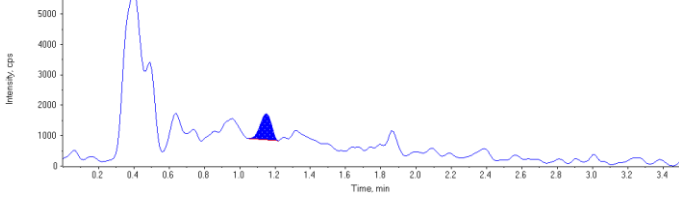
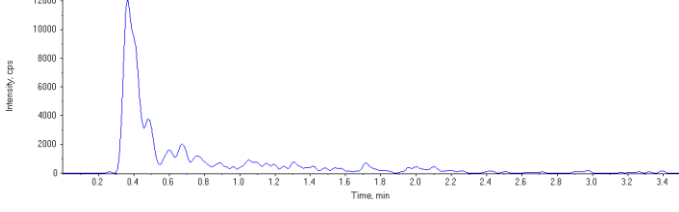
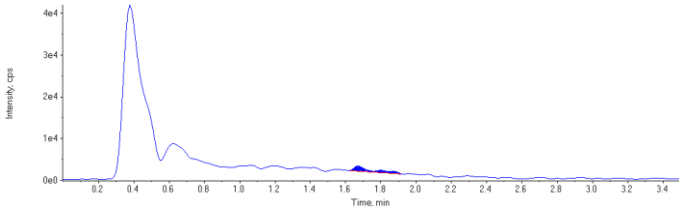
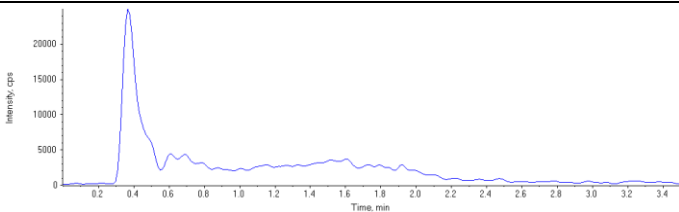
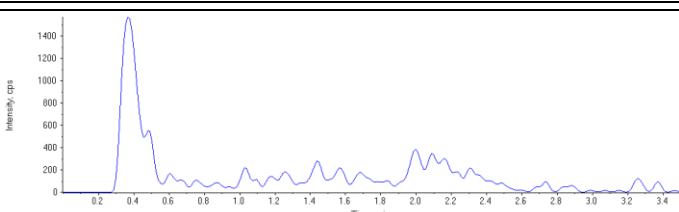
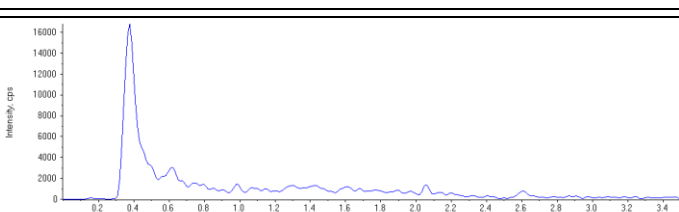
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 50.3 µg/L</p> <p>Area Ratio: 0.0185</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 53.6 µg/L</p> <p>Area Ratio: 0.0663</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 54.2 µg/L</p> <p>Area Ratio: 0.0584</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 59.2 µg/L</p> <p>Area Ratio: 0.0256</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 56.7 µg/L</p> <p>Area Ratio: 0.0417</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 108. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

<b>Sample Name</b>	<b>4386408~BVX830-01</b>	<b>Injection Vial</b>	32
<b>Sample ID</b>	4386408~BVX830-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:38:25 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	Re-injected
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	82200.	1.71	1.00	-
MPFHpA	275000.	1.74	1.00	-
MPFOA	249000.	1.91	1.00	-
MPFOS	102000.	1.99	1.00	-
MPFNA	189000.	2.05	1.00	-
13C6-PFHxA IS	4590000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	3770	1.15	N/A	0.719	N/A
PFHxS 1	0	0.00	N/A	N/A	N/A
PFHpA 1	9300	1.67	N/A	0.324	N/A
PFOA 1	0	0.00	N/A	N/A	N/A
PFOS 1	0	0.00	N/A	N/A	N/A
PFNA 1	0	0.00	N/A	N/A	N/A
18O2-PFHxS	82200	1.71	N/A	48.6	N/A
13C4-PFHpA	275000	1.74	N/A	48.4	N/A
13C4-PFOA	249000	1.91	N/A	50.3	N/A
13C4-PFOS	102000	1.99	N/A	51.6	N/A
13C5-PFNA	189000	2.05	N/A	55.9	N/A
13C6-PFHxA	4590000	1.48	N/A	115.	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.15 (1.15) min</p> <p>Calculated Conc: 0.719 µg/L</p> <p>Area Ratio: 0.0459</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.68) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.67 (1.75) min</p> <p>Calculated Conc: 0.324 µg/L</p> <p>Area Ratio: 0.0339</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.92) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.97) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

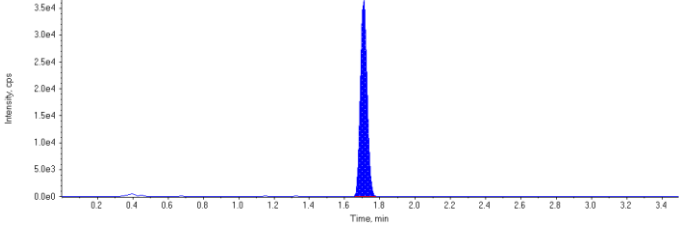
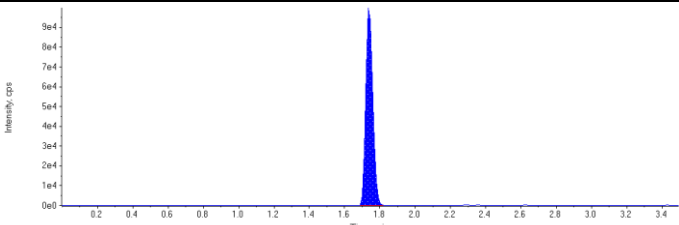
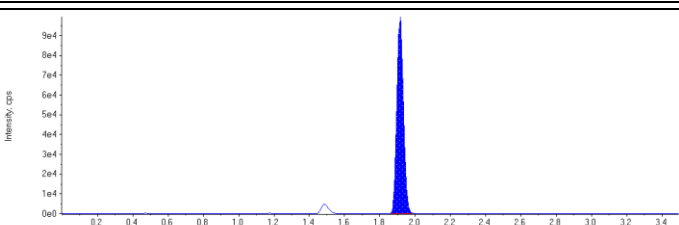
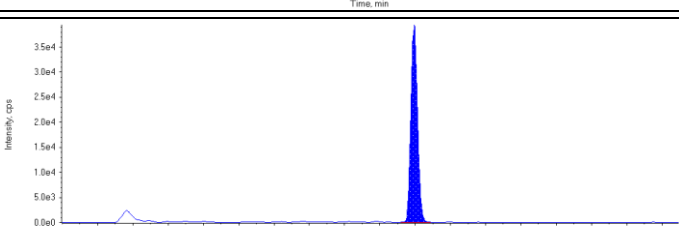
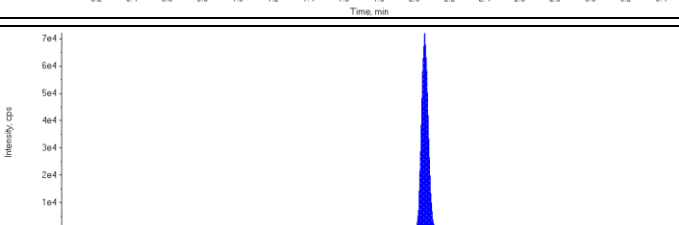
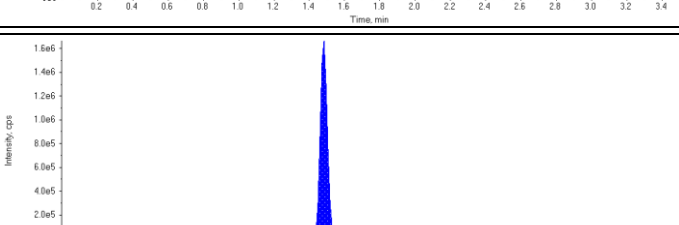


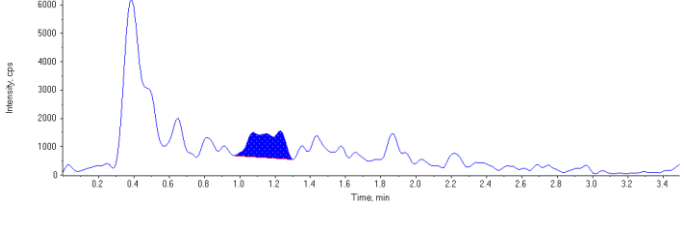
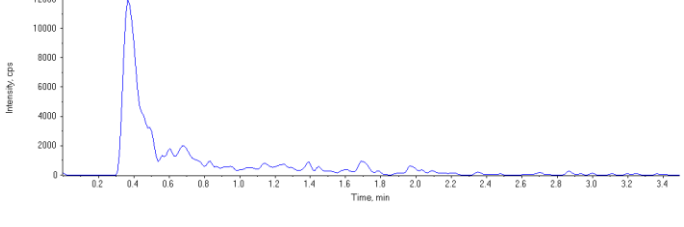
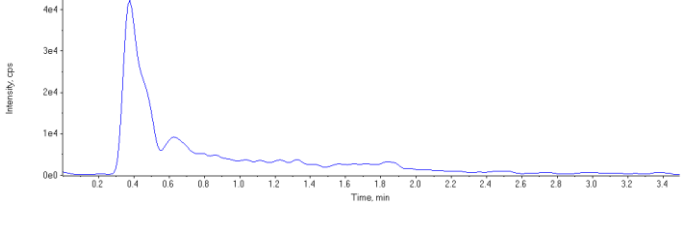
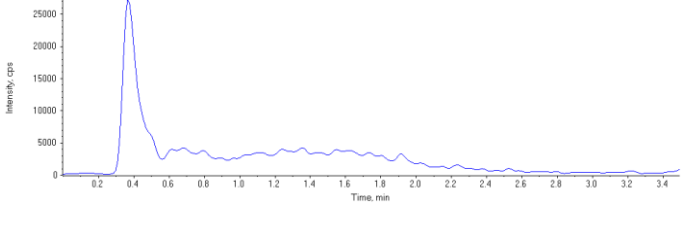
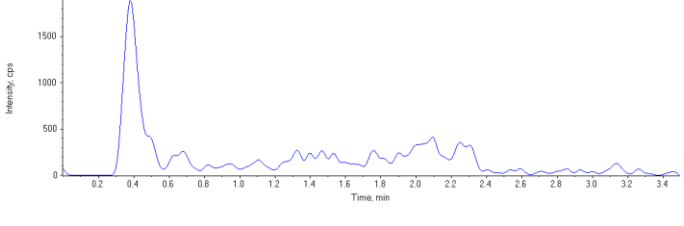
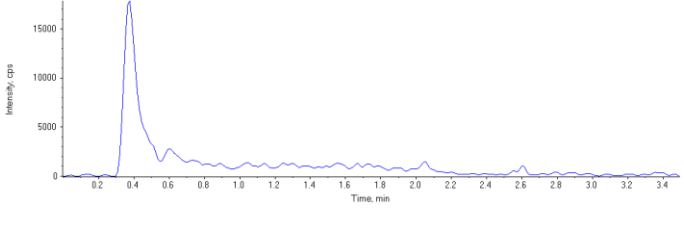
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 48.6 µg/L</p> <p>Area Ratio: 0.0179</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 48.4 µg/L</p> <p>Area Ratio: 0.0598</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 50.3 µg/L</p> <p>Area Ratio: 0.0543</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 51.6 µg/L</p> <p>Area Ratio: 0.0223</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 55.9 µg/L</p> <p>Area Ratio: 0.0411</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 115. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	

<b>Sample Name</b>	4386408~BVX830-01	<b>Injection Vial</b>	32
<b>Sample ID</b>	4386408~BVX830-01	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Unknown	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:58:48 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	Reported
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	96600.	1.71	1.00	-
MPFHpA	277000.	1.74	1.00	-
MPFOA	258000.	1.91	1.00	-
MPFOS	103000.	1.99	1.00	-
MPFNA	188000.	2.05	1.00	-
13C6-PFHxA IS	4700000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	11900	1.23	N/A	0.966	N/A
PFHxS 1	0	0.00	N/A	N/A	N/A
PFHpA 1	0	0.00	N/A	N/A	N/A
PFOA 1	0	0.00	N/A	N/A	N/A
PFOS 1	0	0.00	N/A	N/A	N/A
PFNA 1	0	0.00	N/A	N/A	N/A
18O2-PFHxS	96600	1.71	N/A	55.9	N/A
13C4-PFHpA	277000	1.74	N/A	47.6	N/A
13C4-PFOA	258000	1.91	N/A	50.8	N/A
13C4-PFOS	103000	1.99	N/A	50.8	N/A
13C5-PFNA	188000	2.05	N/A	54.5	N/A
13C6-PFHxA	4700000	1.48	N/A	118.	N/A

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Unknown)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Unknown)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.23 (1.15) min</p> <p>Calculated Conc: 0.966 µg/L</p> <p>Area Ratio: 0.123</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFBS 1. The y-axis is Intensity (cps) from 0 to 6000. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes, and a smaller peak is highlighted at 1.23 minutes.</p>
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.68) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFHxS 1. The y-axis is Intensity (cps) from 0 to 12000. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes.</p>
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.75) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFHpA 1. The y-axis is Intensity (cps) from 0 to 4e4. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes.</p>
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.92) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFOA 1. The y-axis is Intensity (cps) from 0 to 25000. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes.</p>
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.97) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFOS 1. The y-axis is Intensity (cps) from 0 to 1500. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes.</p>
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	 <p>Chromatogram for PFNA 1. The y-axis is Intensity (cps) from 0 to 15000. The x-axis is Time (min) from 0.2 to 3.4. A major peak is observed at approximately 0.4 minutes.</p>

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 55.9 µg/L</p> <p>Area Ratio: 0.0206</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 47.6 µg/L</p> <p>Area Ratio: 0.0589</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 50.8 µg/L</p> <p>Area Ratio: 0.0548</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 50.8 µg/L</p> <p>Area Ratio: 0.0220</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 54.5 µg/L</p> <p>Area Ratio: 0.0401</p> <p>Sample Type: (Unknown)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 118. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Unknown)</p>	



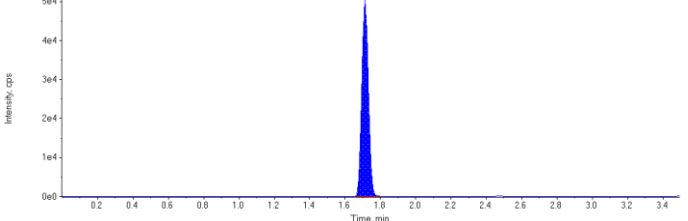
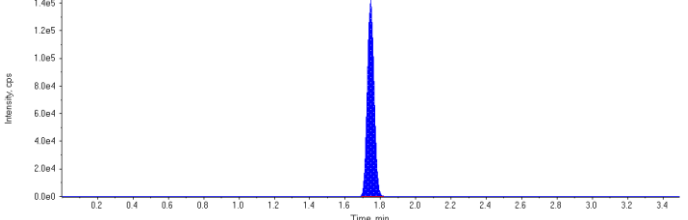
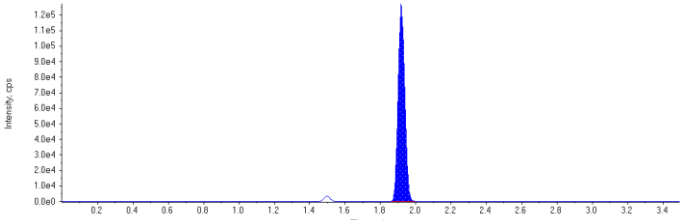
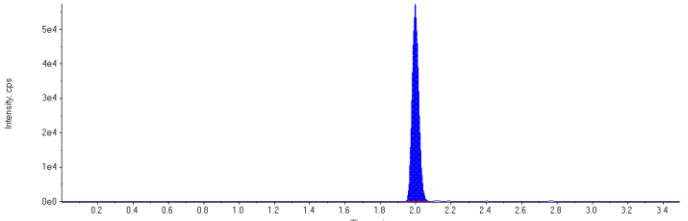
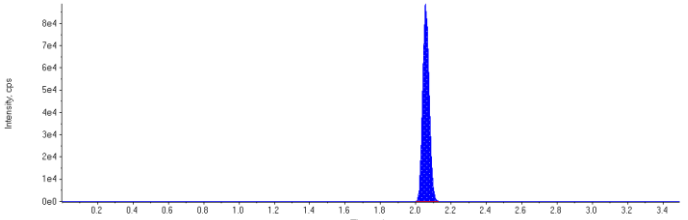
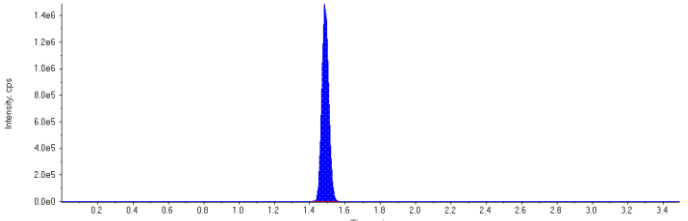
## 4. QA/QC Data

Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

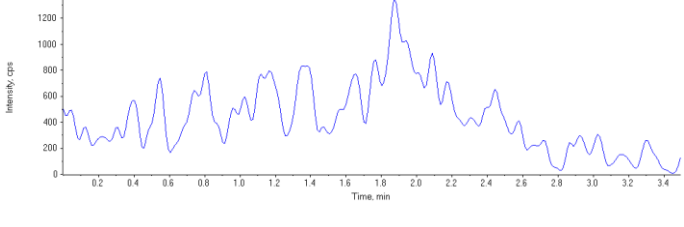
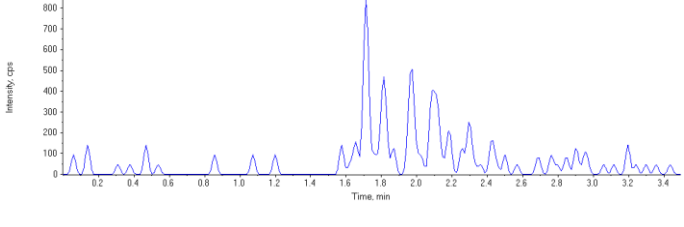
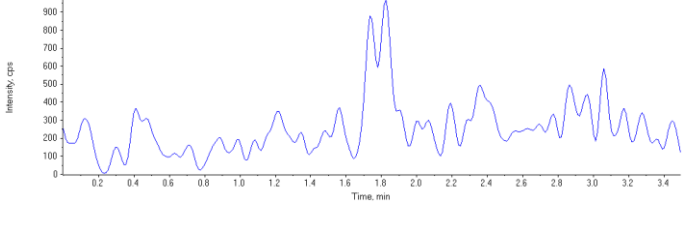
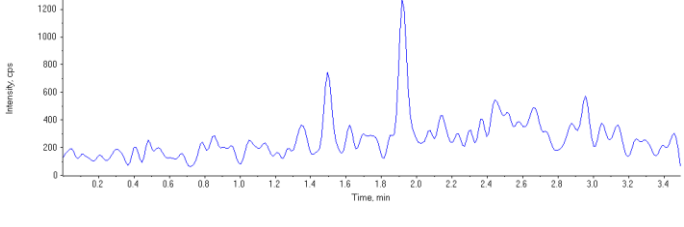
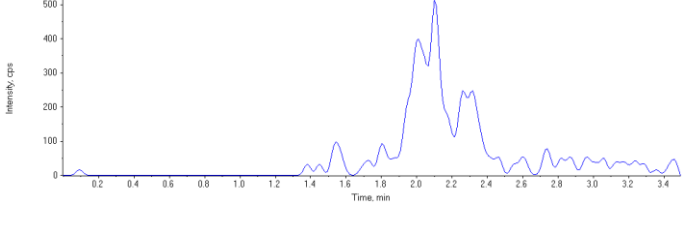
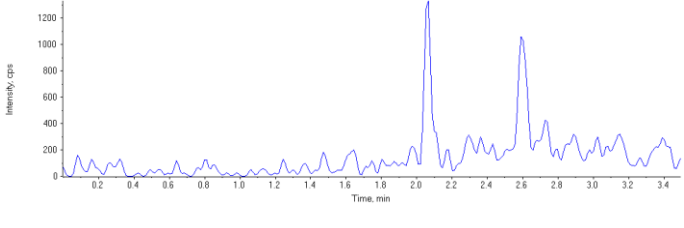
<b>Sample Name</b>	4386408~BLANK	<b>Injection Vial</b>	2
<b>Sample ID</b>	4386408~BLANK	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	130000.	1.71	1.00	-
MPFHpA	375000.	1.74	1.00	-
MPFOA	327000.	1.92	1.00	-
MPFOS	146000.	2.00	1.00	-
MPFNA	228000.	2.06	1.00	-
13C6-PFHxA IS	4100000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	0	0.00	0.00	N/A	0.0
PFHxS 1	0	0.00	0.00	N/A	0.0
PFHpA 1	0	0.00	0.00	N/A	0.0
PFOA 1	0	0.00	0.00	N/A	0.0
PFOS 1	0	0.00	0.00	N/A	0.0
PFNA 1	0	0.00	0.00	N/A	0.0
18O2-PFHxS	130000	1.71	100.	86.2	86.2
13C4-PFHpA	375000	1.74	100.	73.9	73.9
13C4-PFOA	327000	1.92	100.	73.9	73.9
13C4-PFOS	146000	2.00	100.	82.4	82.4
13C5-PFNA	228000	2.06	100.	75.7	75.7
13C6-PFHxA	4100000	1.49	100.	103.	103.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>



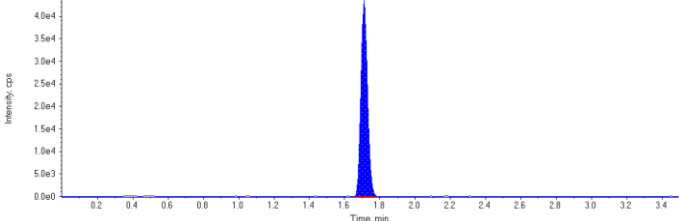
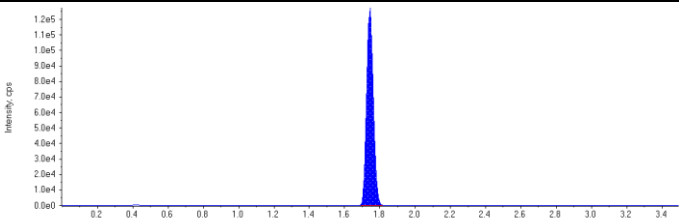
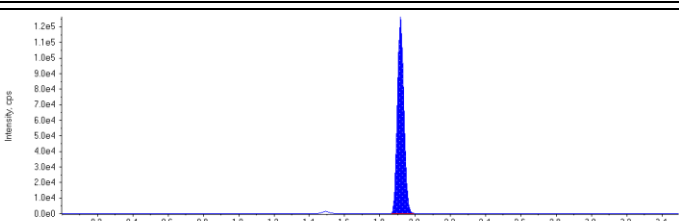
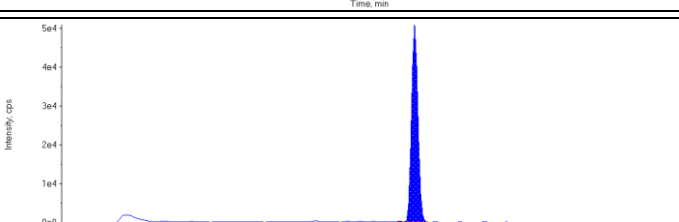
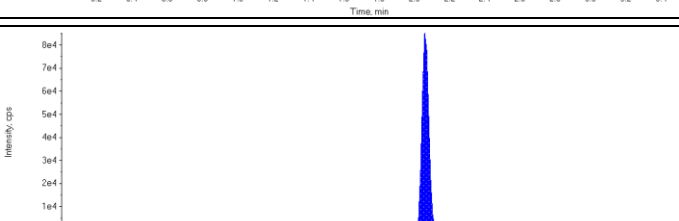
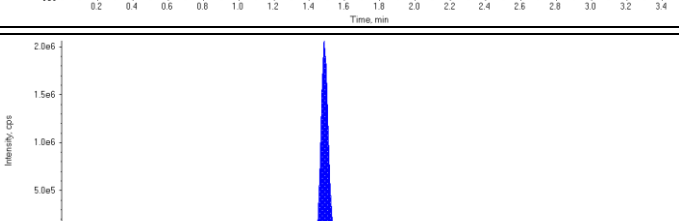
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.15) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.68) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.75) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 0.00 (1.92) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 0.00 (1.97) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 0.00 (2.02) min</p> <p>Calculated Conc: N/A µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

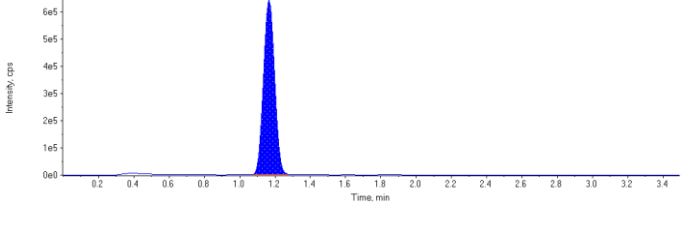
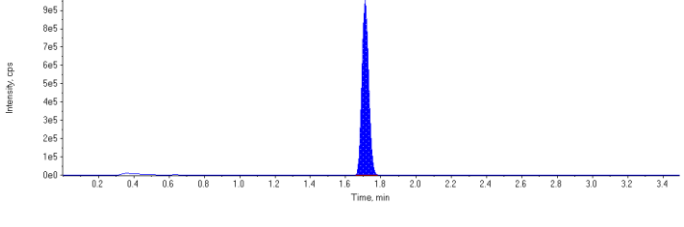
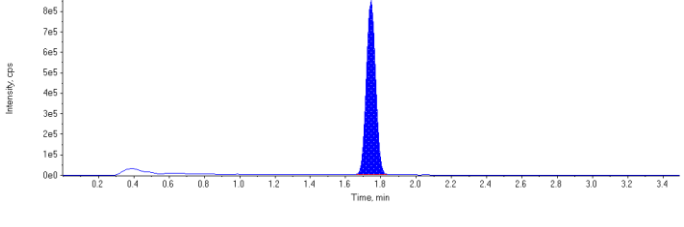
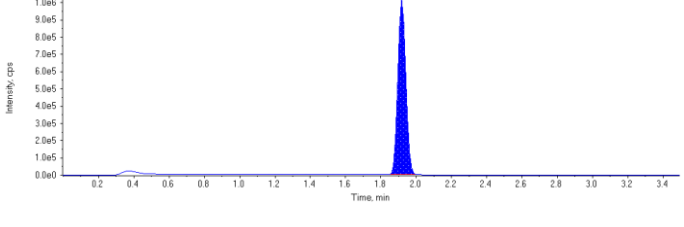
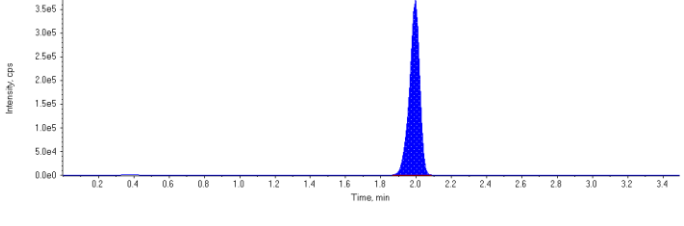
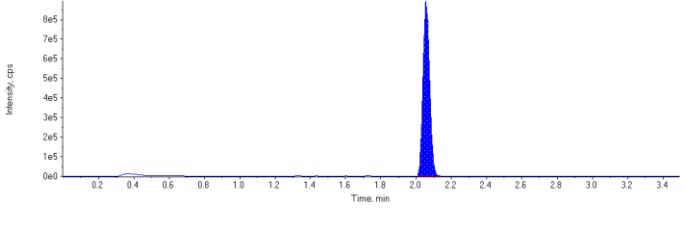
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 86.2 µg/L</p> <p>Area Ratio: 0.0317</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 73.9 µg/L</p> <p>Area Ratio: 0.0914</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 73.9 µg/L</p> <p>Area Ratio: 0.0798</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 82.4 µg/L</p> <p>Area Ratio: 0.0356</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 75.7 µg/L</p> <p>Area Ratio: 0.0557</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

<b>Sample Name</b>	<b>4386408~MTRX SPK</b>	<b>Injection Vial</b>	10
<b>Sample ID</b>	4386408~MTRX SPK	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:25:52 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	Re-injected
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	112000.	1.71	1.00	-
MPFHpA	335000.	1.74	1.00	-
MPFOA	321000.	1.92	1.00	-
MPFOS	127000.	2.00	1.00	-
MPFNA	227000.	2.06	1.00	-
13C6-PFHxA IS	5760000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2820000	1.17	50.0	81.2	162.0
PFHxS 1	2500000	1.71	50.0	68.1	136.0
PFHpA 1	3080000	1.74	50.0	68.3	137.0
PFOA 1	3200000	1.92	50.0	60.5	121.0
PFOS 1	1470000	1.99	50.0	64.1	128.0
PFNA 1	2340000	2.06	50.0	62.0	124.0
18O2-PFHxS	112000	1.71	100.	52.6	52.6
13C4-PFHpA	335000	1.74	100.	47.1	47.1
13C4-PFOA	321000	1.92	100.	51.7	51.7
13C4-PFOS	127000	2.00	100.	51.0	51.0
13C5-PFNA	227000	2.06	100.	53.5	53.5
13C6-PFHxA	5760000	1.49	100.	144.	144.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

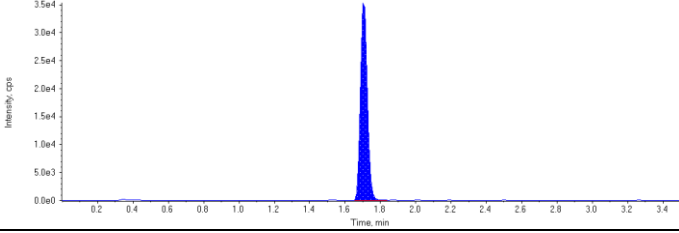
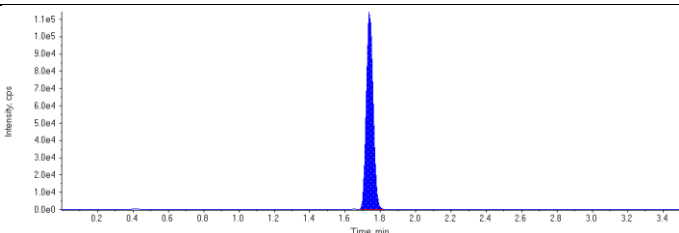
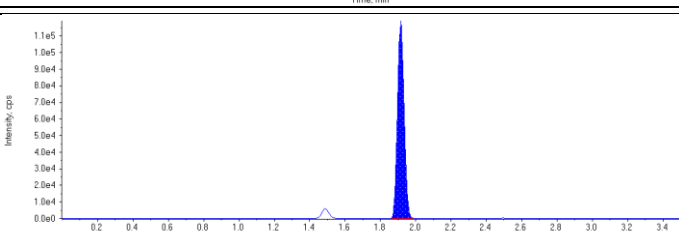
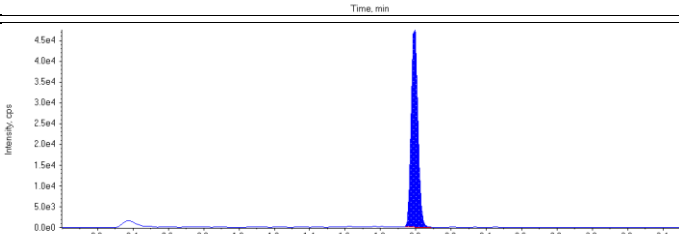
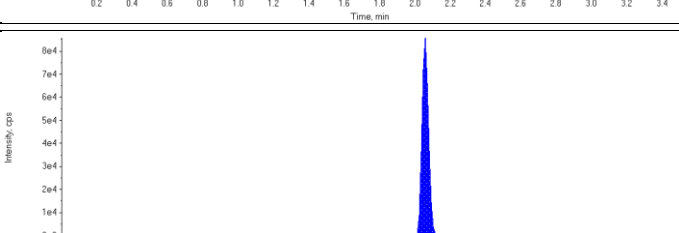
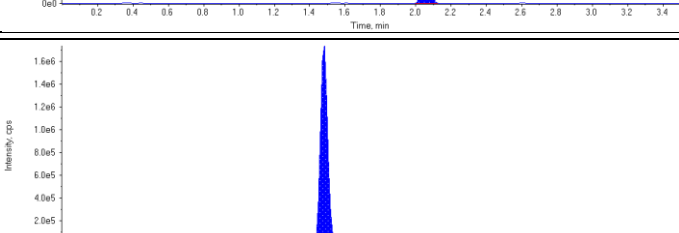
<p><b>PFBS 1 (298.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.17 (1.15) min</p> <p>Calculated Conc: 81.2 µg/L</p> <p>Area Ratio: 25.3</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1 (398.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 68.1 µg/L</p> <p>Area Ratio: 22.4</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1 (363.000/319.000 Da)</b></p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 68.3 µg/L</p> <p>Area Ratio: 9.19</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1 (413.100/369.000 Da)</b></p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 60.5 µg/L</p> <p>Area Ratio: 9.96</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1 (498.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 64.1 µg/L</p> <p>Area Ratio: 11.6</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1 (462.900/419.000 Da)</b></p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 62.0 µg/L</p> <p>Area Ratio: 10.3</p> <p>Sample Type: (Quality Control)</p>	

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 52.6 µg/L</p> <p>Area Ratio: 0.0194</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 47.1 µg/L</p> <p>Area Ratio: 0.0582</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 51.7 µg/L</p> <p>Area Ratio: 0.0557</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 51.0 µg/L</p> <p>Area Ratio: 0.0220</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 53.5 µg/L</p> <p>Area Ratio: 0.0394</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 144. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

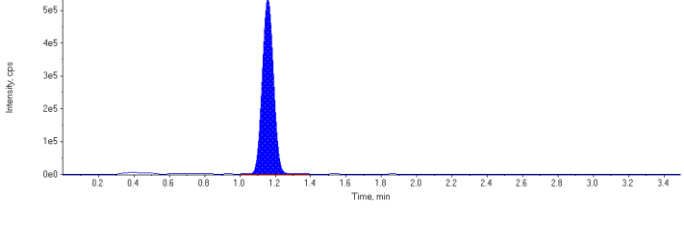
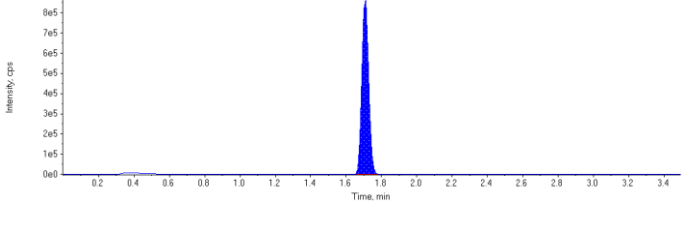
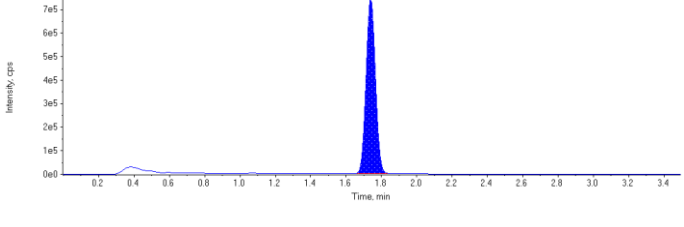
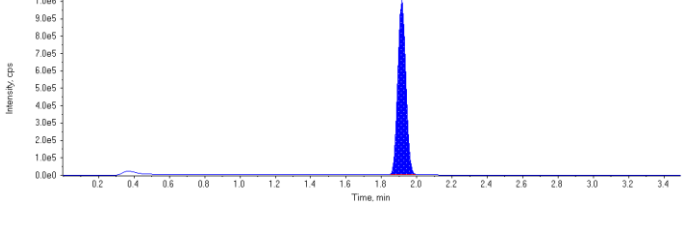
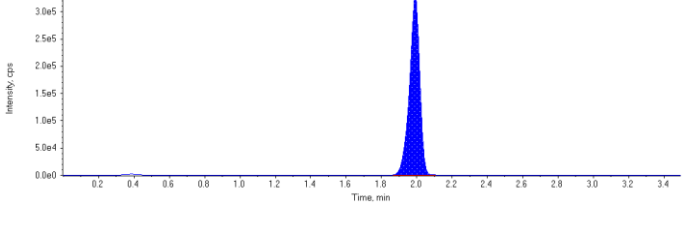
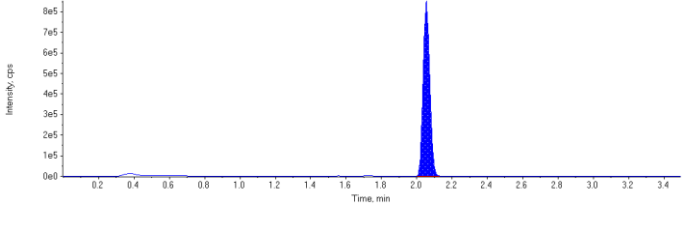
<b>Sample Name</b>	<b>4386408~MTRX SPK</b>	<b>Injection Vial</b>	10
<b>Sample ID</b>	4386408~MTRX SPK	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:48:37 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	Reported
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
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<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

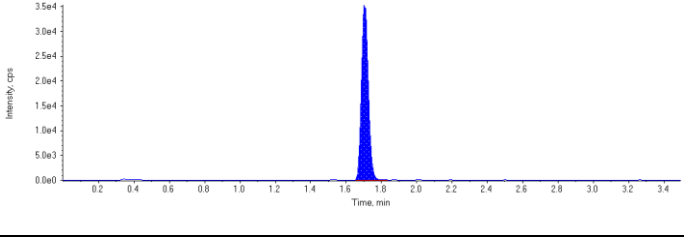
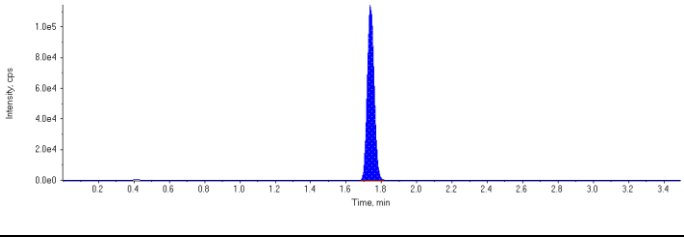
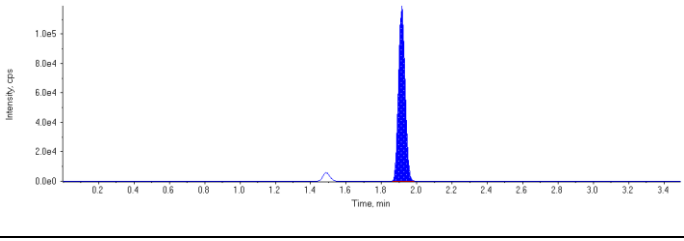
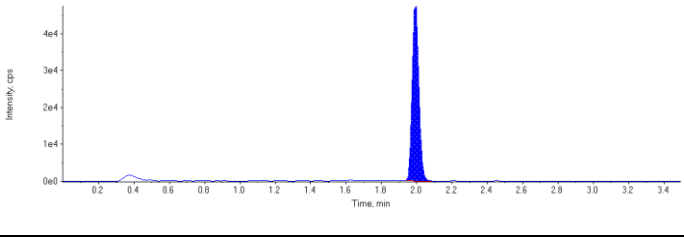
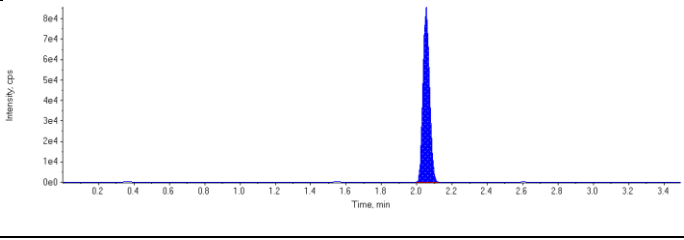
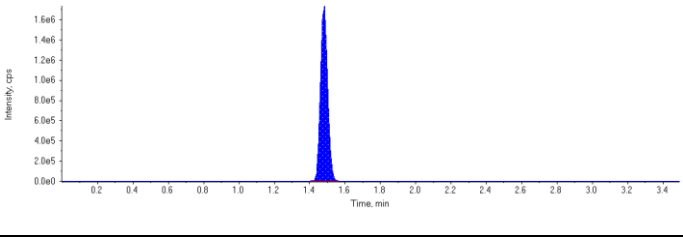
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	95600.	1.71	1.00	-
MPFHpA	303000.	1.74	1.00	-
MPFOA	307000.	1.91	1.00	-
MPFOS	126000.	1.99	1.00	-
MPFNA	222000.	2.05	1.00	-
13C6-PFHxA IS	4930000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2350000	1.16	50.0	79.2	158.0
PFHxS 1	2330000	1.71	50.0	73.8	148.0
PFHpA 1	2720000	1.74	50.0	66.7	133.0
PFOA 1	3150000	1.92	50.0	62.3	125.0
PFOS 1	1320000	1.99	50.0	58.0	116.0
PFNA 1	2220000	2.05	50.0	60.1	120.0
18O2-PFHxS	95600	1.71	100.	52.6	52.6
13C4-PFHpA	303000	1.74	100.	49.7	49.7
13C4-PFOA	307000	1.91	100.	57.7	57.7
13C4-PFOS	126000	1.99	100.	58.9	58.9
13C5-PFNA	222000	2.05	100.	61.2	61.2
13C6-PFHxA	4930000	1.48	100.	124.	124.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>



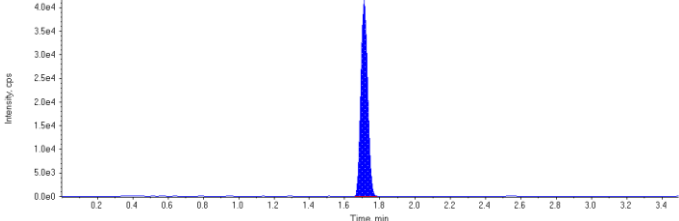
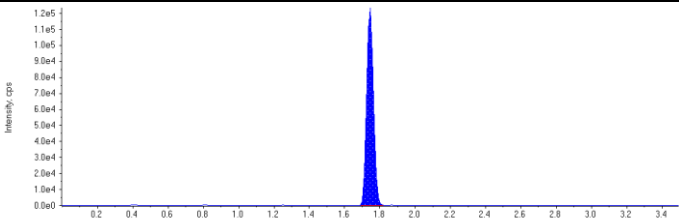
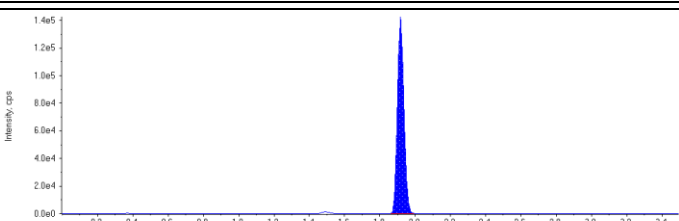
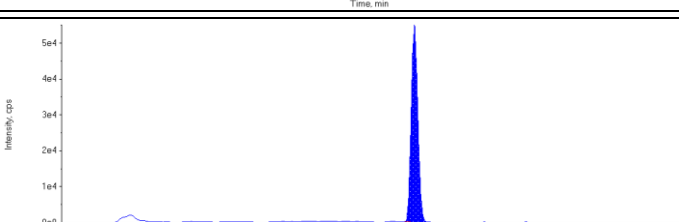
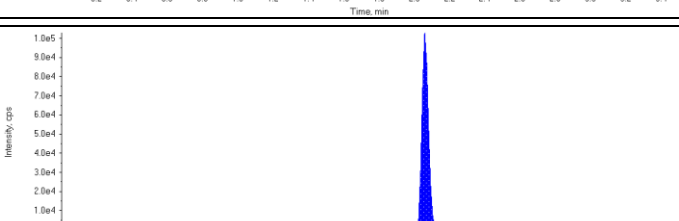
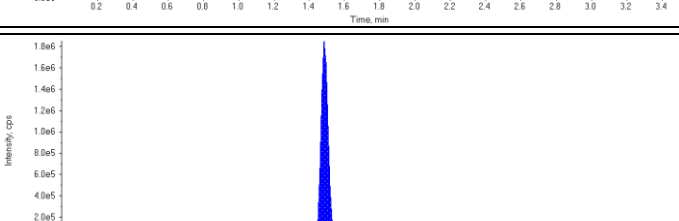
<p><b>PFBS 1 (298.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 79.2 µg/L</p> <p>Area Ratio: 24.6</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1 (398.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 73.8 µg/L</p> <p>Area Ratio: 24.3</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1 (363.000/319.000 Da)</b></p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 66.7 µg/L</p> <p>Area Ratio: 8.98</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1 (413.100/369.000 Da)</b></p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 62.3 µg/L</p> <p>Area Ratio: 10.3</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1 (498.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 58.0 µg/L</p> <p>Area Ratio: 10.5</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1 (462.900/419.000 Da)</b></p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 60.1 µg/L</p> <p>Area Ratio: 9.98</p> <p>Sample Type: (Quality Control)</p>	

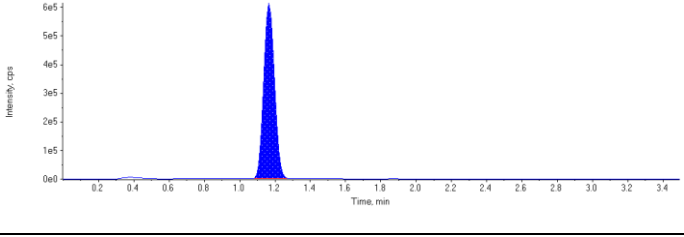
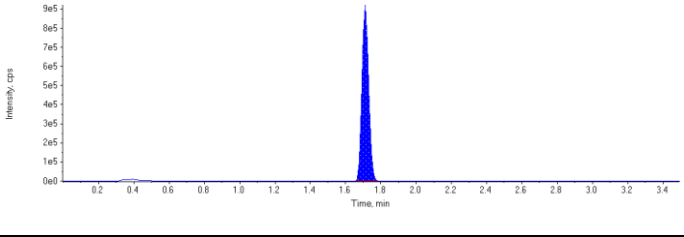
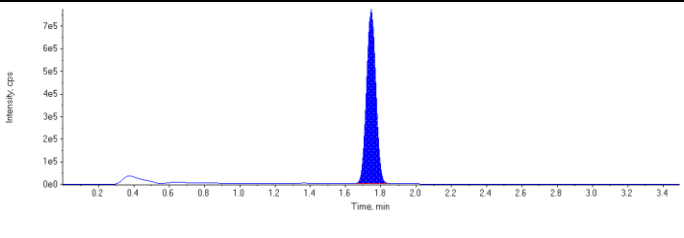
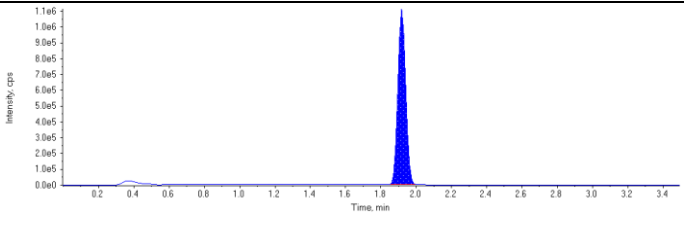
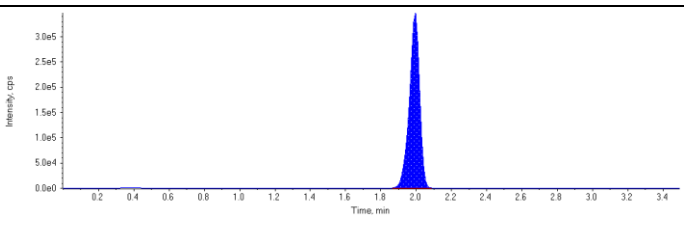
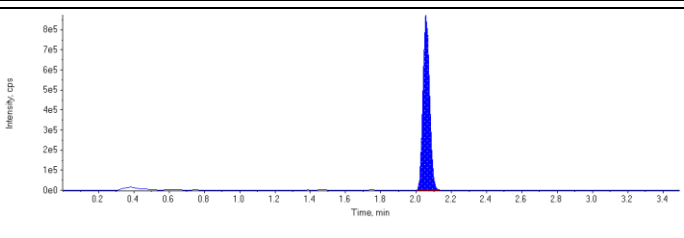
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 52.6 µg/L</p> <p>Area Ratio: 0.0194</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 49.7 µg/L</p> <p>Area Ratio: 0.0615</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 57.7 µg/L</p> <p>Area Ratio: 0.0623</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 58.9 µg/L</p> <p>Area Ratio: 0.0255</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 61.2 µg/L</p> <p>Area Ratio: 0.0450</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 124. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

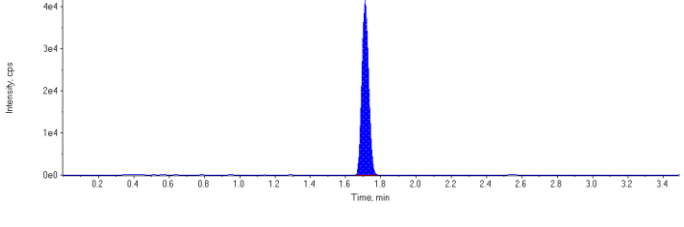
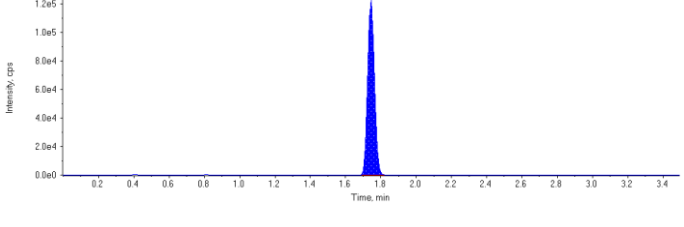
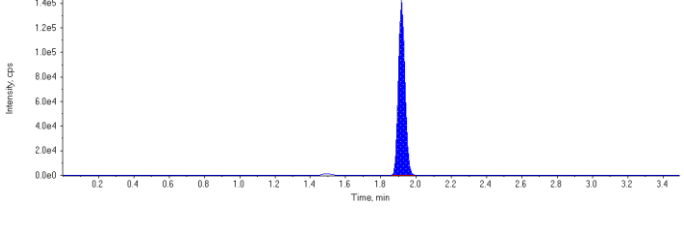
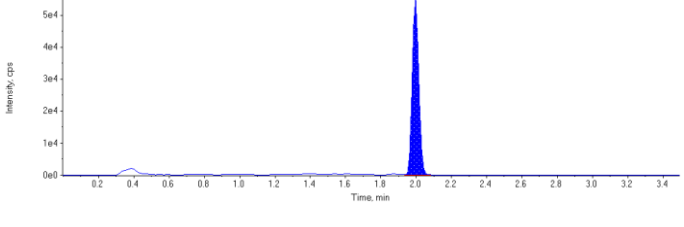
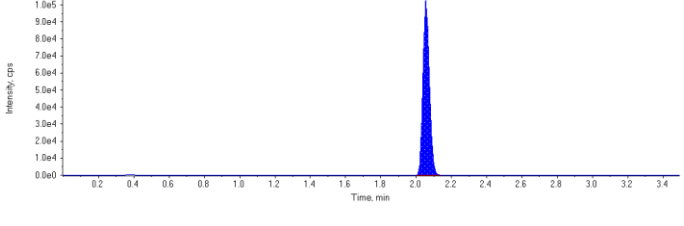
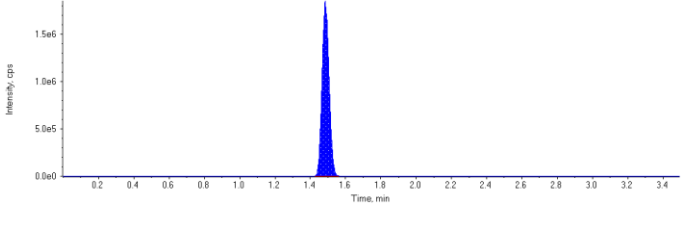
<b>Sample Name</b>	4386408~MTRX SPK:D1	<b>Injection Vial</b>	11
<b>Sample ID</b>	4386408~MTRX SPK:D1	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:30:58 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	110000.	1.71	1.00	-
MPFHpA	331000.	1.74	1.00	-
MPFOA	363000.	1.92	1.00	-
MPFOS	141000.	2.00	1.00	-
MPFNA	263000.	2.06	1.00	-
13C6-PFHxA IS	5140000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2590000	1.17	50.0	76.0	152.0
PFHxS 1	2400000	1.71	50.0	66.4	133.0
PFHpA 1	2790000	1.75	50.0	62.7	125.0
PFOA 1	3420000	1.92	50.0	57.2	114.0
PFOS 1	1400000	1.99	50.0	54.4	109.0
PFNA 1	2270000	2.06	50.0	52.0	104.0
18O2-PFHxS	110000	1.71	100.	58.1	58.1
13C4-PFHpA	331000	1.74	100.	52.1	52.1
13C4-PFOA	363000	1.92	100.	65.5	65.5
13C4-PFOS	141000	2.00	100.	63.7	63.7
13C5-PFNA	263000	2.06	100.	69.5	69.5
13C6-PFHxA	5140000	1.49	100.	129.	129.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

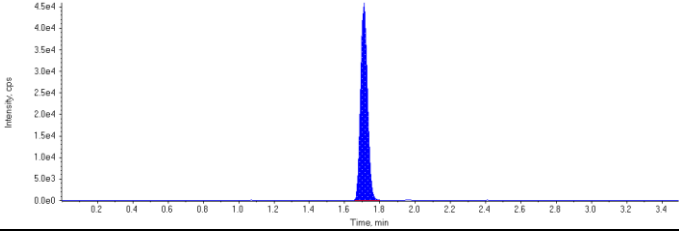
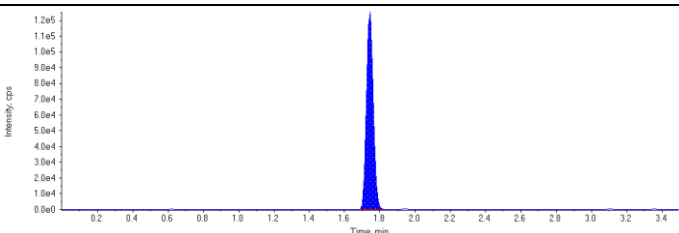
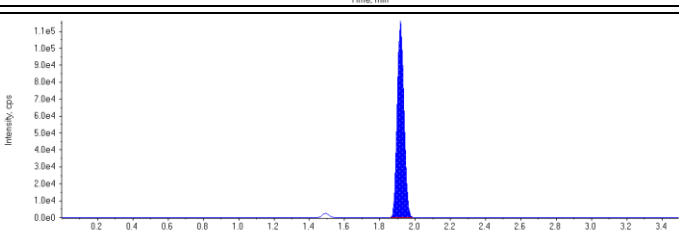
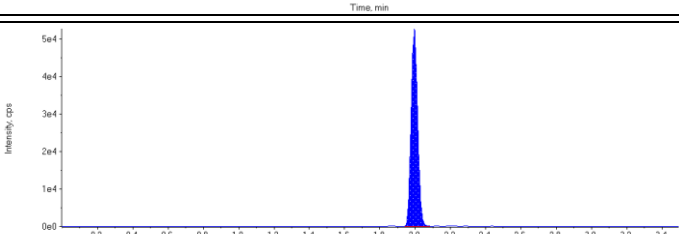
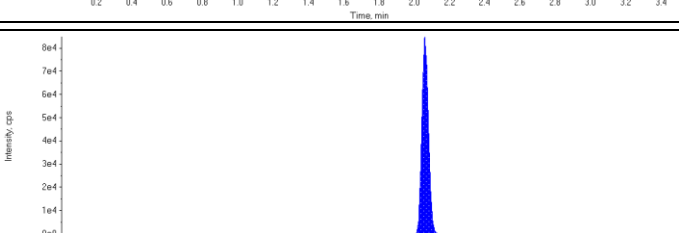
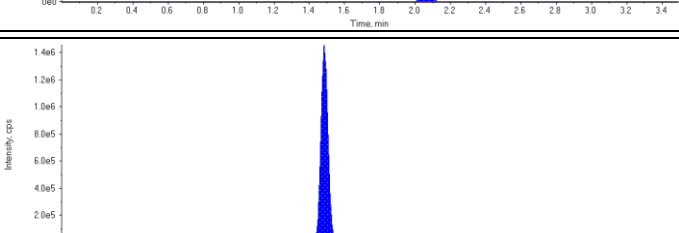
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.17 (1.15) min</p> <p>Calculated Conc: 76.0 µg/L</p> <p>Area Ratio: 23.6</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 66.4 µg/L</p> <p>Area Ratio: 21.9</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.75 (1.75) min</p> <p>Calculated Conc: 62.7 µg/L</p> <p>Area Ratio: 8.44</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 57.2 µg/L</p> <p>Area Ratio: 9.41</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 54.4 µg/L</p> <p>Area Ratio: 9.87</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 52.0 µg/L</p> <p>Area Ratio: 8.63</p> <p>Sample Type: (Quality Control)</p>	

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 58.1 µg/L</p> <p>Area Ratio: 0.0214</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 52.1 µg/L</p> <p>Area Ratio: 0.0644</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 65.5 µg/L</p> <p>Area Ratio: 0.0707</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 63.7 µg/L</p> <p>Area Ratio: 0.0275</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 69.5 µg/L</p> <p>Area Ratio: 0.0511</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 129. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

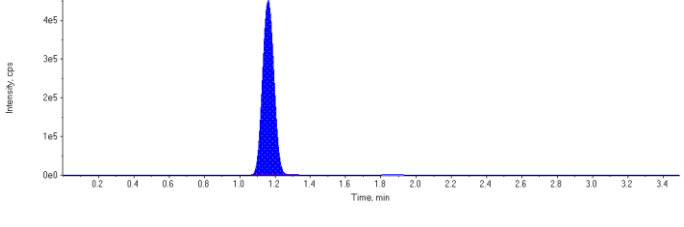
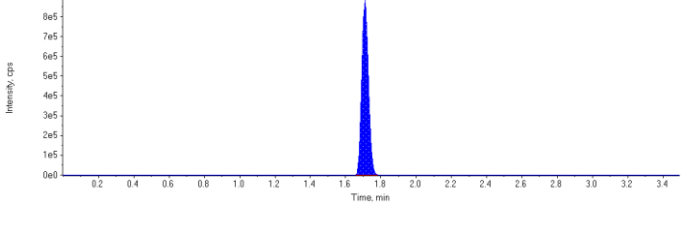
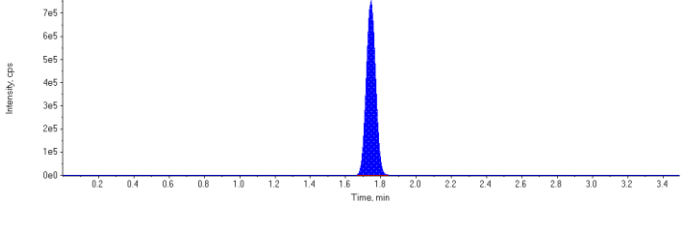
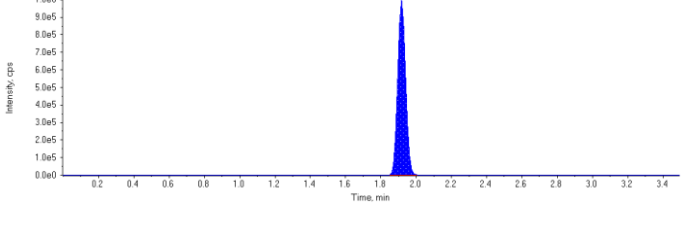
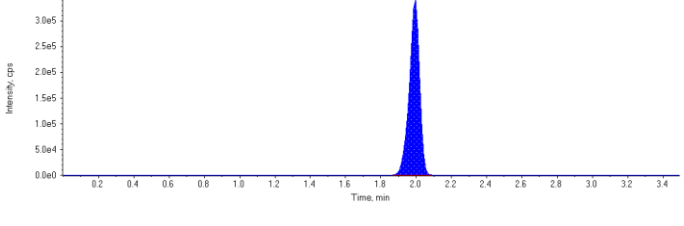
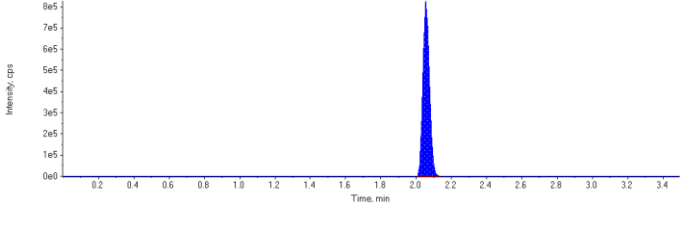
<b>Sample Name</b>	<b>4386408~SPIKE</b>	<b>Injection Vial</b>	12
<b>Sample ID</b>	4386408~SPIKE	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:36:03 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
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<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

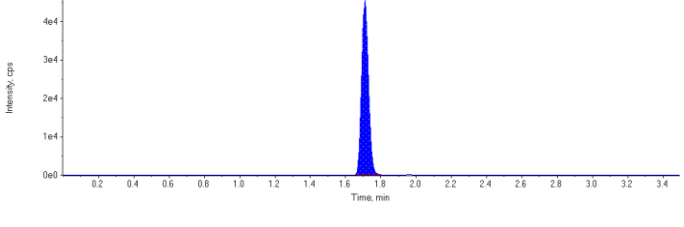
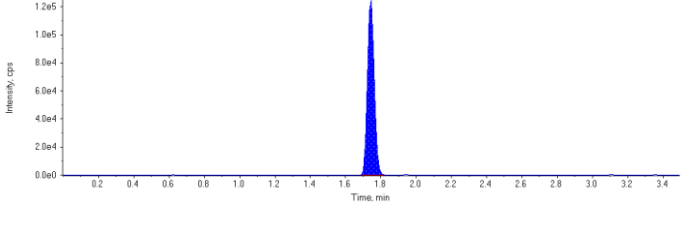
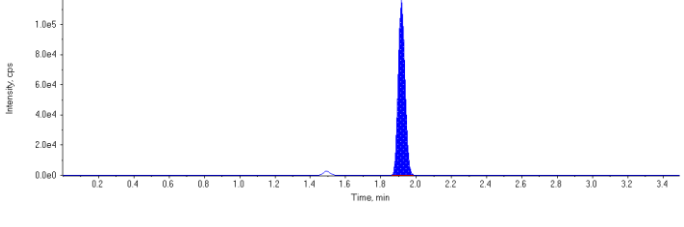
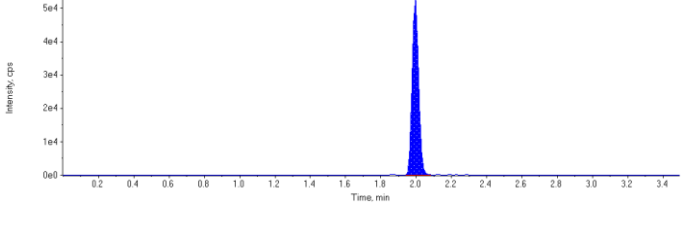
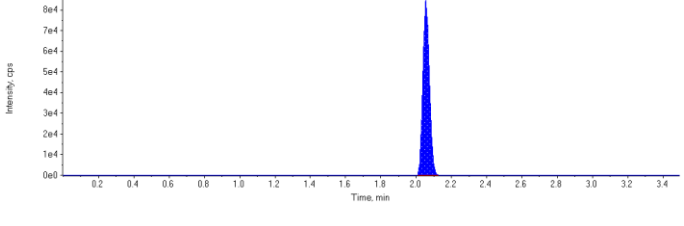
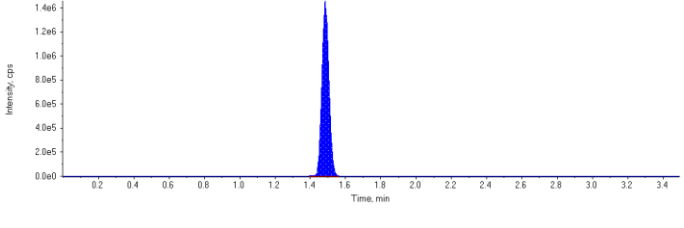
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	123000.	1.71	1.00	-
MPFHpA	336000.	1.74	1.00	-
MPFOA	305000.	1.92	1.00	-
MPFOS	135000.	2.00	1.00	-
MPFNA	218000.	2.06	1.00	-
13C6-PFHxA IS	4050000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2040000	1.16	50.0	53.6	107.0
PFHxS 1	2320000	1.71	50.0	57.3	115.0
PFHpA 1	2750000	1.74	50.0	60.7	121.0
PFOA 1	3120000	1.92	50.0	62.2	124.0
PFOS 1	1380000	1.99	50.0	56.2	112.0
PFNA 1	2130000	2.06	50.0	58.9	118.0
18O2-PFHxS	123000	1.71	100.	82.6	82.6
13C4-PFHpA	336000	1.74	100.	67.2	67.2
13C4-PFOA	305000	1.92	100.	69.9	69.9
13C4-PFOS	135000	2.00	100.	77.4	77.4
13C5-PFNA	218000	2.06	100.	73.2	73.2
13C6-PFHxA	4050000	1.49	100.	101.	101.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>



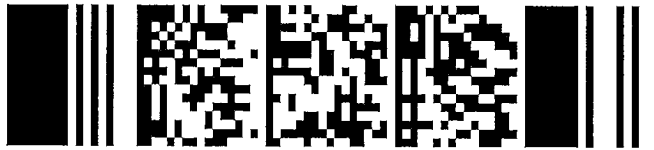
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 53.6 µg/L</p> <p>Area Ratio: 16.6</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 57.3 µg/L</p> <p>Area Ratio: 18.9</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 60.7 µg/L</p> <p>Area Ratio: 8.17</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 62.2 µg/L</p> <p>Area Ratio: 10.2</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 56.2 µg/L</p> <p>Area Ratio: 10.2</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 58.9 µg/L</p> <p>Area Ratio: 9.79</p> <p>Sample Type: (Quality Control)</p>	

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 82.6 µg/L</p> <p>Area Ratio: 0.0304</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 67.2 µg/L</p> <p>Area Ratio: 0.0831</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 69.9 µg/L</p> <p>Area Ratio: 0.0755</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 77.4 µg/L</p> <p>Area Ratio: 0.0335</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 73.2 µg/L</p> <p>Area Ratio: 0.0539</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 101. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

DoD Projects - Internal Data Validation Checklist					
Run date: 2016/02/19					
Worksheet # (s): 4386408					
Analysis: PFD5LOW-w			1st 100% review		*2nd 100% review
Primary review by the analyst - 1st 100 % analysis review			yes	no	n/a
1	Sample analyses meet hold time criteria			✓	✓
2	Analysis set-up meets method criteria	✓			✓
3	Tuning and correct calibration used - criteria meets method criteria	✓			✓
4	SQC/Control Charts updated, analysis in statistical/method control			✓	✓
5	Internal area counts checked (if applicable)	✓			✓
6	LCS, SRM are within acceptance criteria	✓			✓
7	Surrogate Recovery(s) is within acceptance criteria	✓			✓
8	Method Blank meets acceptance criteria	✓			✓
9	Matrix Spike recovery(s) meets acceptance criteria		✓		✓
10	Duplicate precision meets acceptance criteria	✓			✓
11	QC is documented on the run logs	✓			✓
12	Runs checked for carryover	✓			✓
13	Prep log / worksheet(s) are present, signed / dated by a prep / Instrument analysts	✓			✓
14	Initial weights, splits, imprinter volumes (where applicable) are documented	✓			✓
15	Standards and reagents traceable to Certificates of Analysis	✓			✓
16	Samples above calibration range diluted and reanalyzed	✓			✓
17	Dilution factors (where justified) have been checked for correctness and entered	✓			✓
18	Analytical observations/anomalies documented in LIMS	✓			✓
19	Random calculation checked and in correct units	✓			✓
20	If corrective actions were applied they are documented, initialed & dated			✓	✓
21	Manual integration – before & after data with a reason included, initialed & dated			✓	✓
22	Transferred data is validated in LIMS for correctness	✓			✓
23	Data package assembled (where required)	✓			✓
Reviewed by: <i>an</i>		Date: 2016/02/22			
Comments:					
Secondary Supervisor/Qualified Data Review Staff - 2nd 100% verification review					
1	Repeats documented and referenced	✓			
2	Method and sample deviations noted, anomalies described (if applicable)	✓			
3	Data and QC validated in LIMS	✓			
4	Random calculation checked	✓			
5	Benchsheet (s) signed and dated	✓			
6	Data Package (if required) checked for completeness	✓			
Reviewed by: <i>AN</i>		Date: 2016/02/22			
Comments: Samples BVX 770, BVX 771 and BVX 819 were re-worked due to high analyte concentration. The samples will be re-analyzed with an appropriate dilution. <i>AN 2016/02/23</i>					

\*Note: 2nd 100% verification review documented by secondary qualified data review  
 Primary and Secondary Internal Data Review Check must be performed by a different person

Worksheet Data Validation Checklist - Extractable Organics			
Worksheet # 4386408		Testcode: PFOs LOW TW	
<b>Sample Preparation</b>		yes	no
1	Samples extracted within hold time		/
2	Client sample ID verified against Lab ID (waters & oils)	/	
3	Parameter list and Client comments reviewed, (Spiking solutions matched to parameter list)	/	
4	Height of sediment or if sample was decanted, recorded on worksheet	/	
5	Method required QC processed with samples, maximum batch size = 20 client samples.	/	
6	Sample, duplicate, matrix spike appear similar, initial sample as well as final extract	/	
7	Sample weight or initial volume and extract final volume, aliquot factor clearly recorded.	/	
8	If performed any additional dilution clearly recorded		/
9	Matrix spike / Duplicate performed on IOL samples if present		/
10	Spiking solutions valid (haven't expired), ID and volume used clearly identified on worksheet	/	
11	Spiking process witnessed and signed off	/	
12	Extraction type recorded (N3A2B = neutral, 3 x acidic, 2 x basic)	/	GSZ
13	Sample prep deviations documented within CompliantPro as a Policy Deviation		/
14	Job Remarks reviewed on 2nd page of worksheet.	/	
15	Worksheet and reagent tracking record completed and authorized.	/	
Reviewed by: GSZ		Date: 2016/02/18	
Comments:			
<b>Worksheet Approval</b>		yes	no
1	Verified the position of the vials in autosampler against sequence list; signed off sequence list	/	
2	Calibration and CCV standards valid ( haven't expired)	/	
3	Initial calibration curve and DFTPP tune (if applicable) acceptable	/	
4	Continuing and Final CCV and DFTPP tune (if applicable) acceptable	/	
5	System performance check acceptable (if applicable)	/	
6	Internal standard responses acceptable	/	
7	Method blank meets acceptance criteria	/	
8	Lab Control Samples recoveries meets acceptance criteria	/	
9	Duplicate RPD meets acceptance criteria	/	
10	Matrix spike recoveries meets acceptance criteria	/	
11	Surrogate recoveries meets acceptance criteria	/	
12	Appropriate control charts updated		/
13	Samples above calibration range diluted and reanalyzed	/	
14	Dilutions clearly documented on tracking record, inst file and verified during data upload	/	
15	Samples following high level samples checked for carryover.	/	
16	Mass spectra ion ratios acceptable for positive results, hardcopy in file.		/
17	Analytical observations / anomalies documented	/	
18	DQW comments entered in LIMS, hardcopy in file		/
19	Sample Prep section (above) reviewed and verified.	/	
20	WS Approval performed in LIMS	/	
Reviewed by: <i>CR</i>		Date: 2016/02/22	
Comments:			
<b>Worksheet Validation</b>		yes	no
1	Calibration, QC and sample results reviewed and determined acceptable	/	
2	Manual integrations verified		/
3	Random calculation checked	/	
4	Data and QC validated in LIMS	/	
5	Comments reviewed for appropriateness	/	
6	Reworks / relogs documented in file	/	
7	Worksheet signed and dated,	/	
8	Worksheet approved and validated within LIMS	/	
Reviewed by: <i>APL</i>		Date: 2016/02/22	
Comments: Sample BUX 819 was sent for re-work due to high analyte concentration. The sample will be re-analyzed with an appropriate dilution.			
APL 2016/02/23			



**RUSH**

Report Name : Worksheet - (Liquids and Solids)

Assignment Date : Thursday, February 18, 2016

Assigned to : Geoffrey Sanchez

Test Code : PFOSLOW-W

Instrument Id:

Test Description : Low level PFOS and PFOA in water by LC-MS/MS

*Sediment*

Job Number	Sample Number	D	Sample ID	F	% Moisture	Wt or Vol	Final Vol	DF or AF	# Cont	Expiry Date	Test DeadLine	Criteria	Extract Date
	MTRX SPK	0	PFOSL BVX776-01		⊖	125	0.3	1X					2016/02/18
	MTRX SPK	1	PFOSL BVX776-01		⊖	125	0.3	1X					2016/02/18
	SPIKE		PFOSL		⊖	125	0.3	1X					2016/02/18
	BLANK				⊖	125	0.3	1X					2016/02/18
B630791*	*BVX769-01R		OF-FB08-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX770-01R		OF-RW08-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX771-01R		OF-RW08P-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX772-01R		OF-FB41-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX773-01R		OF-RW41-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX774-01R		OF-RW41P-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX775-01R		OF-FB56-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX776-01R		OF-RW56-0216		⊖	125	0.3	1X	3	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX779-01R		OF-FB51-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX780-01R		OF-RW51-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630791*	*BVX781-01R		OF-RW51P-0216		⊖	125	0.3	1X	1	2016/02/18	2016/02/23 17:00		2016/02/18
B630797*	*BVX826-01R		OF-RW09-0216		<0.1	125	0.3	1X	1	2016/02/17	2016/02/23 17:00		2016/02/18
B630797*	*BVX827-01R		OF-RW37-0216		⊖	125	0.3	1X	1	2016/02/17	2016/02/23 17:00		2016/02/18
B630797*	*BVX828-01R		OF-RW11-0216		<0.1	125	0.3	1X	1	2016/02/17	2016/02/23 17:00		2016/02/18
B630797*	*BVX829-01R		OF-RW28-0216		⊖	125	0.3	1X	1	2016/02/17	2016/02/23 17:00		2016/02/18
B630797*	*BVX830-01R		OF-RW67-0216		⊖	125	0.3	1X	1	2016/02/17	2016/02/23 17:00		2016/02/18
B630796	BVX814				<0.1	12.5	0.3	1X					
B630796	BVX817				<0.1	12.5	0.3	1X					
B630796	BVX818				<0.2	6.25	0.3	1X					
B630796	BVX819				<0.1	12.5	0.3	1X					
<i>GSZ 2016/02/18</i>													

Remarks: Job B630797 - extracted past the holding time (resulted from discrepancy between dates in TAT & summary; see attached email).

Job B630796 - rework; upfront dilution needed; lower volume used.

Samples extracted by: Geoffrey Sanchez, GSZ 2016/02/18

Instrumentation performed by: CS

Date: 2016/02/19

Calculations performed by: CS

Date: 2016/02/22

Validated by: AS

Date: 2016/02/23



Low level PFOS and PFOA in water - Water  
ng/L

Parameter Name	Units	MTRX SPK	DL	MTRX SPK Dup1	DL	SPIKE	DL
Perfluorobutanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorobutane Sulfonate (PFBS)	ng/L	11 156.786	2	11 150.386	2	107.20000	2
Perfluorodecane Sulfonate	ng/L	N/A*****		N/A*****		N/A*****	
Perfluoroheptanoic Acid (PFHpA)	ng/L	11 133.020	2	125.02000	2	121.40000	2
Perfluoroheptane sulfonate	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorohexanoic Acid (PFHxA)	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorohexane Sulfonate (PFHxS)	ng/L	11 147.600	2	11 132.800	2	114.60000	2
Perfluorononanoic Acid (PFNA)	ng/L	120.20000	2	104.00000	2	117.80000	2
Perfluoropentanoic Acid (PFPeA)	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorotetradecanoic Acid	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorotridecanoic Acid	ng/L	N/A*****		N/A*****		N/A*****	
Perfluoroundecanoic Acid (PFUnA)	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorodecanoic Acid (PFDA)	ng/L	N/A*****		N/A*****		N/A*****	
Perfluorododecanoic Acid (PFDoA)	ng/L	N/A*****		N/A*****		N/A*****	
Perfluoro-n-Octanoic Acid (PFOA)	ng/L	124.60000	2	114.40000	2	124.40000	2
Perfluorooctane Sulfonate (PFOS)	ng/L	116.00000	2	108.80000	2	112.40000	2
13C2-perfluorotetradecanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C4-Perfluorobutanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C5-Perfluorononanoic acid	ng/L	61.2		69.5		73.2	
13C2-Perfluorodecanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C2-Perfluorododecanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C2-Perfluorohexanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C2-Perfluoroundecanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
13C4-Perfluoroheptanoic acid	ng/L	49.7		52.1		67.2	
13C4-Perfluorooctanoic acid	ng/L	57.7		65.5		69.9	
13C4-Perfluorooctanesulfonate	ng/L	58.9		63.7		77.4	
13C5-Perfluoropentanoic acid	ng/L	N/A*****		N/A*****		N/A*****	
18O2-Perfluorohexanesulfonate	ng/L	52.6		58.1		82.6	

Parameter Name	BLANK	DL	B630791 BVX769	DL	B630791 BVX770	DL	B630791 BVX771
Perfluorobutanoic acid	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorobutane Sulfonate (PFBS)	0	2	0	2	5.03000	2	5.24000
Perfluorodecane Sulfonate	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluoroheptanoic Acid (PFHpA)	0	2	0	2	2.88000	2	2.95000
Perfluoroheptane sulfonate	N/A*****			2		2	
Perfluorohexanoic Acid (PFHxA)	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorohexane Sulfonate (PFHxS)	0	2	0	2	53.30000	2	59.70000
Perfluorononanoic Acid (PFNA)	0	2	0	2	0.42900	2	0.39200
Perfluoropentanoic Acid (PFPeA)	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorotetradecanoic Acid	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorotridecanoic Acid	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluoroundecanoic Acid (PFUnA)	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorodecanoic Acid (PFDA)	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluorododecanoic Acid (PFDoA)	N/A*****		N/A*****	2	N/A*****	2	N/A*****
Perfluoro-n-Octanoic Acid (PFOA)	0	2	0	2	13.20000	2	13.90000
Perfluorooctane Sulfonate (PFOS)	0	2	0	2	223.00000	2	225.00000
13C2-perfluorotetradecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluorobutanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C5-Perfluorononanoic acid	75.7		86.2		83.1		80.3
13C2-Perfluorodecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorododecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorohexanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluoroundecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluoroheptanoic acid	73.9		71.9		62.5		60.5
13C4-Perfluorooctanoic acid	73.9		79.1		69.6		67.8
13C4-Perfluorooctanesulfonate	82.4		89.8		87.8		82.8
13C5-Perfluoropentanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
18O2-Perfluorohexanesulfonate	86.2		79.4		87.0		78.1

Low level PFOS and PFOA in water - Water  
ng/L

Parameter Name	DL	B630791 BVX772	DL	B630791 BVX773	DL	B630791 BVX774	DL
Perfluorobutanoic acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorobutane Sulfonate (PFBS)	2	0	2	0	2	0.67400	2
Perfluorodecane Sulfonate	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoroheptanoic Acid (PFHpA)	2	0	2	0	2	0	2
Perfluoroheptane sulfonate	2		2		2		2
Perfluorohexanoic Acid (PFHxA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorohexane Sulfonate (PFHxS)	2	0	2	0.68700	2	0.68700	2
Perfluorononanoic Acid (PFNA)	2	0	2	0	2	0	2
Perfluoropentanoic Acid (PFPeA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorotetradecanoic Acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorotridecanoic Acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoroundecanoic Acid (PFUnA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorodecanoic Acid (PFDA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorododecanoic Acid (PFDoA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoro-n-Octanoic Acid (PFOA)	2	0	2	0.29800	2	0	2
Perfluorooctane Sulfonate (PFOS)	2	0	2	0	2	0	2
13C2-perfluorotetradecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluorobutanoic acid		N/A*****		N/A*****		N/A*****	
13C5-Perfluorononanoic acid		77.7		66.0		63.5	
13C2-Perfluorodecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorododecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorohexanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluoroundecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluoroheptanoic acid		64.8		59.4		54.9	
13C4-Perfluorooctanoic acid		67.5		64.8		60.0	
13C4-Perfluorooctanesulfonate		77.6		67.7		61.0	
13C5-Perfluoropentanoic acid		N/A*****		N/A*****		N/A*****	
18O2-Perfluorohexanesulfonate		72.2		62.2		62.2	

Parameter Name	B630791 BVX775	DL	B630791 BVX776	DL	B630791 BVX779	DL	B630791 BVX780
Perfluorobutanoic acid	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorobutane Sulfonate (PFBS)	0	2	0.80700	2	0.63000	2	0
Perfluorodecane Sulfonate	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluoroheptanoic Acid (PFHpA)	0	2	0.19000	2	0	2	6.61000
Perfluoroheptane sulfonate		2		2		2	
Perfluorohexanoic Acid (PFHxA)	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorohexane Sulfonate (PFHxS)	0	2	0	2	0	2	0
Perfluorononanoic Acid (PFNA)	0	2	0	2	0	2	0
Perfluoropentanoic Acid (PFPeA)	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorotetradecanoic Acid	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorotridecanoic Acid	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluoroundecanoic Acid (PFUnA)	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorodecanoic Acid (PFDA)	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluorododecanoic Acid (PFDoA)	N/A*****	2	N/A*****	2	N/A*****	2	N/A*****
Perfluoro-n-Octanoic Acid (PFOA)	0	2	0	2	0	2	0.30000
Perfluorooctane Sulfonate (PFOS)	0	2	0	2	0	2	0
13C2-perfluorotetradecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluorobutanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C5-Perfluorononanoic acid	62.9		74.4		88.7		62.6
13C2-Perfluorodecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorododecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorohexanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluoroundecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluoroheptanoic acid	53.0		59.7		75.7		50.5
13C4-Perfluorooctanoic acid	61.3		66.0		81.1		56.6
13C4-Perfluorooctanesulfonate	58.7		66.2		87.1		63.7
13C5-Perfluoropentanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
18O2-Perfluorohexanesulfonate	61.5		69.7		77.4		65.2



Low level PFOS and PFOA in water - Water  
ng/L

Parameter Name	DL	B630791 BVX781	DL	B630796 BVX814 ReWork	DL	B630796 BVX817 ReWork	DL
Perfluorobutanoic acid	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorobutane Sulfonate (PFBS)	2	0	2	9.58000	20	8.19000	20
Perfluorodecane Sulfonate	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluoroheptanoic Acid (PFHpA)	2	4.19000	2	17.70000	20	16.90000	20
Perfluoroheptane sulfonate	2		2	N/A*****	20	N/A*****	20
Perfluorohexanoic Acid (PFHxA)	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorohexane Sulfonate (PFHxS)	2	0	2	7.66000	20	7.48000	20
Perfluorononanoic Acid (PFNA)	2	0	2	3.74000	20	3.88000	20
Perfluoropentanoic Acid (PFPeA)	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorotetradecanoic Acid	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorotridecanoic Acid	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluoroundecanoic Acid (PFUnA)	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorodecanoic Acid (PFDA)	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluorododecanoic Acid (PFDoA)	2	N/A*****	2	N/A*****	20	N/A*****	20
Perfluoro-n-Octanoic Acid (PFOA)	2	0	2	493.00000	20	501.00000	20
Perfluorooctane Sulfonate (PFOS)	2	0	2	6.13000	20	5.90000	20
13C2-perfluorotetradecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluorobutanoic acid		N/A*****		N/A*****		N/A*****	
13C5-Perfluorononanoic acid		61.6		68.7		75.3	
13C2-Perfluorodecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorododecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorohexanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluoroundecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluoroheptanoic acid		53.4		58.9		69.6	
13C4-Perfluorooctanoic acid		57.3		64.0		72.1	
13C4-Perfluorooctanesulfonate		59.1		69.1		77.5	
13C5-Perfluoropentanoic acid		N/A*****		N/A*****		N/A*****	
18O2-Perfluorohexanesulfonate		61.7		66.0		84.1	

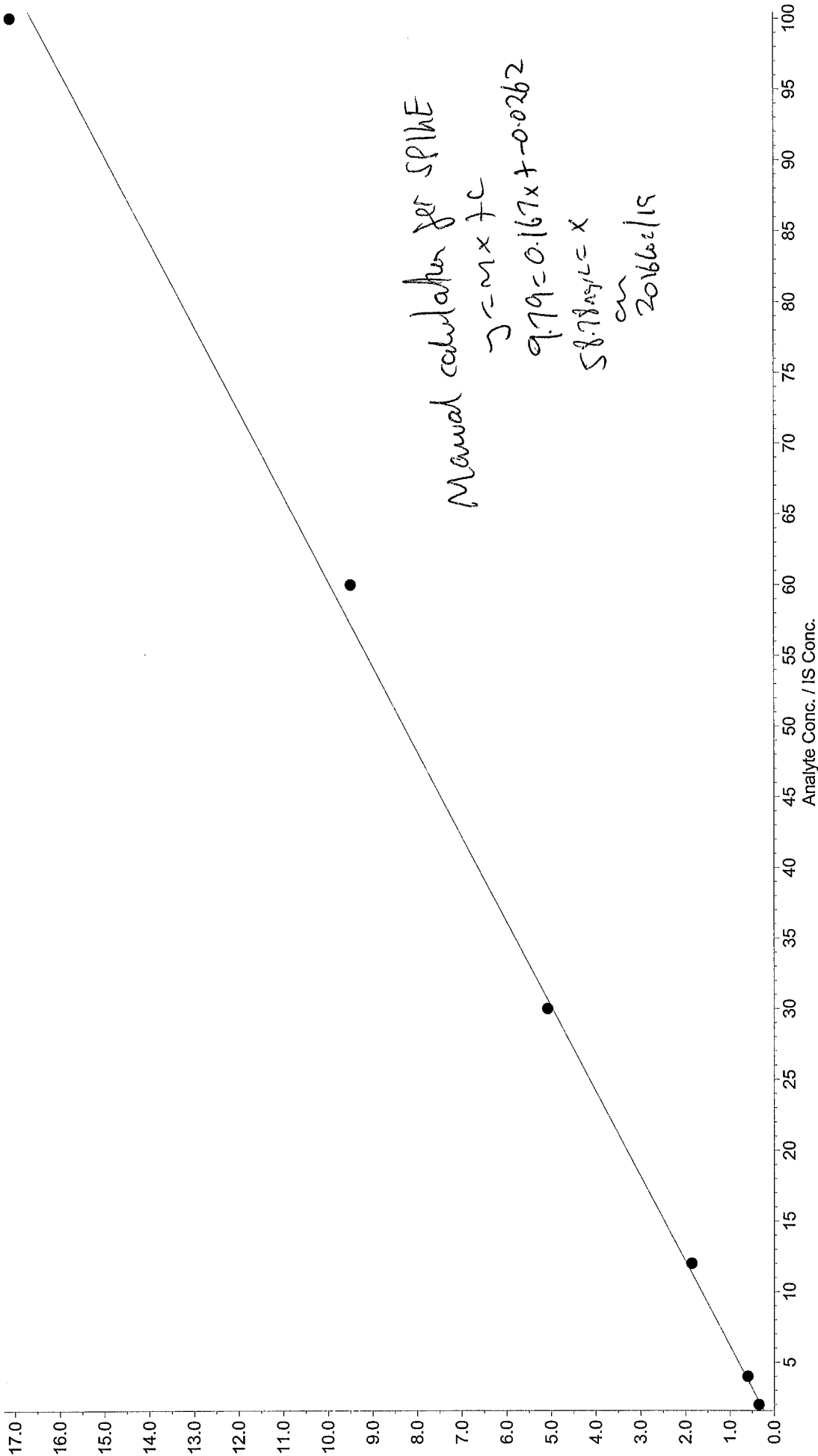
Parameter Name	B630796 BVX818 ReWork	DL	B630796 BVX819 ReWork	DL	B630797 BVX826	DL	B630797 BVX827
Perfluorobutanoic acid	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorobutane Sulfonate (PFBS)	N/A*****	40	8.62000	20	0	2	0.79200
Perfluorodecane Sulfonate	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluoroheptanoic Acid (PFHpA)	N/A*****	40	33.30000	20	0	2	0.45500
Perfluoroheptane sulfonate	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorohexanoic Acid (PFHxA)	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorohexane Sulfonate (PFHxS)	N/A*****	40	7.77000	20	0.71300	2	3.26000
Perfluorononanoic Acid (PFNA)	N/A*****	40	5.80000	20	0	2	0
Perfluoropentanoic Acid (PFPeA)	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorotetradecanoic Acid	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorotridecanoic Acid	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluoroundecanoic Acid (PFUnA)	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorodecanoic Acid (PFDA)	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluorododecanoic Acid (PFDoA)	N/A*****	40	N/A*****	20	N/A*****	2	N/A*****
Perfluoro-n-Octanoic Acid (PFOA)	1870	40	1100.00000	20	0.41200	2	7.15000
Perfluorooctane Sulfonate (PFOS)	N/A*****	40	8.57000	20	0.32000	2	0.81200
13C2-perfluorotetradecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluorobutanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C5-Perfluorononanoic acid	N/A*****		62.7		61.7		65.0
13C2-Perfluorodecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorododecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluorohexanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C2-Perfluoroundecanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
13C4-Perfluoroheptanoic acid	N/A*****		43.1		53.8		51.9
13C4-Perfluorooctanoic acid	77.6		53.3		58.1		56.6
13C4-Perfluorooctanesulfonate	N/A*****		79.5		59.3		63.9
13C5-Perfluoropentanoic acid	N/A*****		N/A*****		N/A*****		N/A*****
18O2-Perfluorohexanesulfonate	N/A*****		74.5		69.2		64.4

Low level PFOS and PFOA in water - Water  
ng/L

Parameter Name	DL	B630797 BVX828	DL	B630797 BVX829	DL	B630797 BVX830	DL
Perfluorobutanoic acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorobutane Sulfonate (PFBS)	2	0	2	0	2	0.96600	2
Perfluorodecane Sulfonate	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoroheptanoic Acid (PFHpA)	2	0	2	0	2	0	2
Perfluoroheptane sulfonate	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorohexanoic Acid (PFHxA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorohexane Sulfonate (PFHxS)	2	0.75200	2	0.75300	2	0	2
Perfluorononanoic Acid (PFNA)	2	0	2	0.36600	2	0	2
Perfluoropentanoic Acid (PFPeA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorotetradecanoic Acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorotridecanoic Acid	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoroundecanoic Acid (PFUnA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorodecanoic Acid (PFDA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluorododecanoic Acid (PFDoA)	2	N/A*****	2	N/A*****	2	N/A*****	2
Perfluoro-n-Octanoic Acid (PFOA)	2	0	2	0	2	0	2
Perfluorooctane Sulfonate (PFOS)	2	0.87100	2	0	2	0	2
13C2-perfluorotetradecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluorobutanoic acid		N/A*****		N/A*****		N/A*****	
13C5-Perfluorononanoic acid		60.5		56.7		54.5	
13C2-Perfluorodecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorododecanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluorohexanoic acid		N/A*****		N/A*****		N/A*****	
13C2-Perfluoroundecanoic acid		N/A*****		N/A*****		N/A*****	
13C4-Perfluoroheptanoic acid		55.3		53.6		47.6	
13C4-Perfluorooctanoic acid		59.2		54.2		50.8	
13C4-Perfluorooctanesulfonate		53.9		59.2		50.8	
13C5-Perfluoropentanoic acid		N/A*****		N/A*****		N/A*****	
18O2-Perfluorohexanesulfonate		59.8		50.3		55.9	

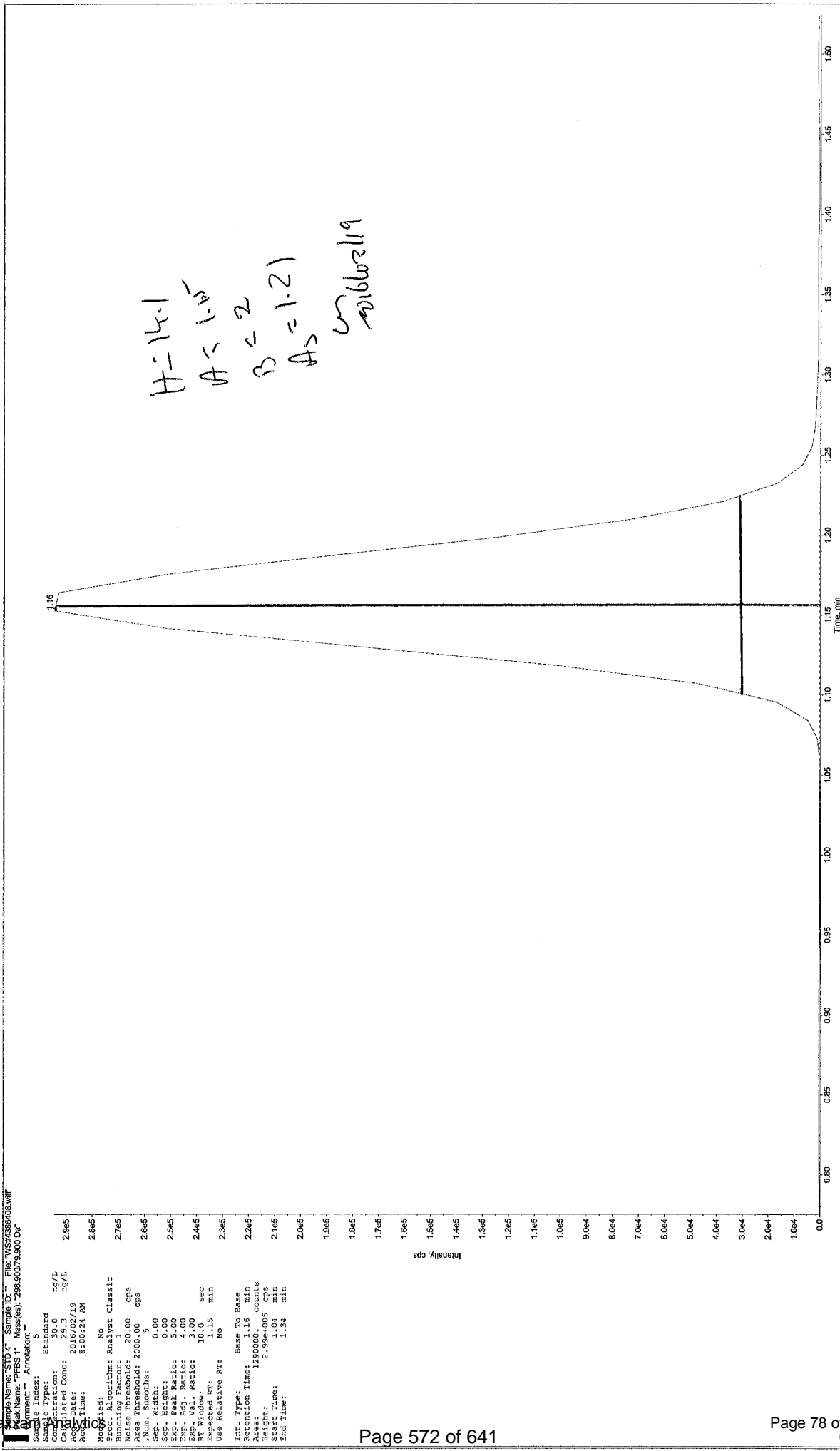
Parameter Name	RDL	MDL	IDL				
Perfluorobutanoic acid	2	0.41	0				
Perfluorobutane Sulfonate (PFBS)	2	0.27	0				
Perfluorodecane Sulfonate	2	0.38	0				
Perfluoroheptanoic Acid (PFHpA)	2	0.39	0				
Perfluoroheptane sulfonate	2	0.4	0				
Perfluorohexanoic Acid (PFHxA)	2	0.42	0				
Perfluorohexane Sulfonate (PFHxS)	2	0.4	0				
Perfluorononanoic Acid (PFNA)	2	0.33	0				
Perfluoropentanoic Acid (PFPeA)	2	0.46	0				
Perfluorotetradecanoic Acid	2	0.61	0				
Perfluorotridecanoic Acid	2	0.6	0				
Perfluoroundecanoic Acid (PFUnA)	2	0.5	0				
Perfluorodecanoic Acid (PFDA)	2	0.24	0				
Perfluorododecanoic Acid (PFDoA)	2	0.63	0				
Perfluoro-n-Octanoic Acid (PFOA)	2	0.39	0				
Perfluorooctane Sulfonate (PFOS)	2	0.3	0				
13C2-perfluorotetradecanoic acid							
13C4-Perfluorobutanoic acid							
13C5-Perfluorononanoic acid							
13C2-Perfluorodecanoic acid							
13C2-Perfluorododecanoic acid							
13C2-Perfluorohexanoic acid							
13C2-Perfluoroundecanoic acid							
13C4-Perfluoroheptanoic acid							
13C4-Perfluorooctanoic acid							
13C4-Perfluorooctanesulfonate							
13C5-Perfluoropentanoic acid							
18O2-Perfluorohexanesulfonate							

PFC\_Water\_160219\_4386408\_ULow.rdb (PFNA 1): "Linear" Regression ("1 / x" weighting): y = 0.167 x + -0.0262 (r = 0.9991)



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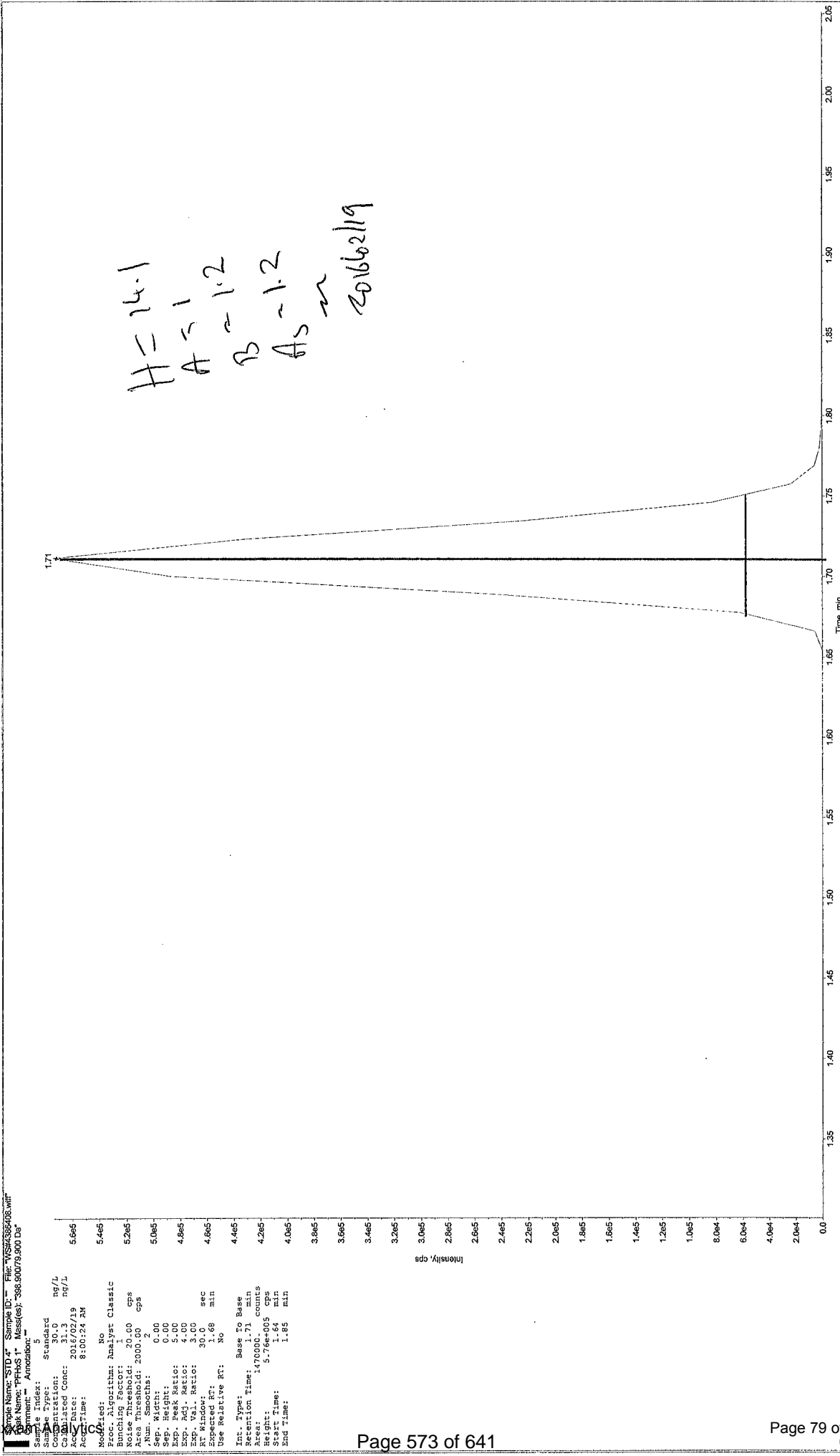
Printing Date: Friday, February 19, 2016



Sample Name: 20160219\_4386408\_ULow.rdb  
 Sample ID: 20160219\_4386408\_ULow.rdb  
 Sample Index: 5  
 Sample Type: Standard  
 Concentration: 20.0 mg/L  
 Acquisition Date: 2016/02/19  
 Acquisition Time: 8:00:24 AM  
 Method: No  
 Weights: Analyst Classic  
 Noise Threshold: 20.00 cps  
 Area Threshold: 2000.00 cps  
 Num. Scans: 5  
 Exp. Weight: 0.00  
 Exp. Peak Ratio: 5.00  
 Exp. Adj. Ratio: 4.00  
 Exp. Val. Ratio: 3.00  
 Expected RT: 1.15 min  
 Use Relative RT: No  
 Int. Type: Base To Base  
 Retention Time: 1.16 min  
 Area: 128000.0 cps  
 Height: 2.95e+05 cps  
 Start Time: 1.04 min  
 End Time: 1.34 min

Results Path: \\miss-netapp2\lcms\lcms3\Analyst  
Data\Projects\Enviro\PFOS\Results\PFC\_Water\_160219\_4386408\_ULow.rdb

Printing Date: Friday, February 19, 2016



Sample Name: STD 4\* Sample ID: File: WS#4386408.vmr  
 Comment: Annotation: PPHHS 1\* Mass(es): 388.90079.900 Da\*  
 Standard  
 Sample Type: Standard  
 Concentration: 30.0 ng/L  
 Calculated Conc: 31.3 ng/L  
 Acq Date: 2016/02/19  
 Acq Time: 8:00:24 AM  
 Modified: No  
 Proc. Algorithm: Analyst Classic  
 Bunching Factor: 1  
 Num. Peaks: 1  
 Area Threshold: 2000.00 cps  
 Num. Smoother: 2  
 Sep. Width: 0.00  
 Sep. Height: 0.00  
 Sep. Ratio: 4.00  
 EXP. Ret. Ratio: 4.00  
 EXP. Val. Ratio: 3.00  
 RT Window: 30.0 sec  
 Suspected RT: 1.68 min  
 Use Relative RT: No  
 Int. Type: Base To Base  
 Retention Time: 1.71 min  
 Area: 1470000 counts  
 Height: 5.665 cps  
 Start Time: 1.64 min  
 End Time: 1.85 min

Report Name: Worksheet - Parameter Lists

Report Date: 2016/02/18

Test Code: PFOSLOW-W

Worksheet Number: 4386408

<u>Sample Number</u>	<u>Parameter</u>
BVX769-01	Perfluorobutane Sulfonate (PFBS)
BVX770-01	Perfluoroheptanoic Acid (PFHpA)
BVX771-01	Perfluorohexane Sulfonate (PFHxS)
BVX772-01	Perfluorononanoic Acid (PFNA)
BVX773-01	Perfluoro-n-Octanoic Acid (PFOA)
BVX774-01	Perfluorooctane Sulfonate (PFOS)
BVX775-01	
BVX776-01	
BVX779-01	
BVX780-01	
BVX781-01	
BVX826-01	
BVX827-01	
BVX828-01	
BVX829-01	
BVX830-01	

WorkSheet 4386408 Instrument Sequences

1.	 4386408:MTRX SPK	MTRX SPK
2.	 4386408:MTRX SPK:D1	MTRX SPK :D1
3.	 4386408:SPIKE	SPIKE
4.	 4386408:BLANK	BLANK
5.	 4386408:BVX769-01	OF-FB08-0216
6.	 4386408:BVX770-01	OF-RW08-0216
7.	 4386408:BVX771-01	OF-RW08P-0216
8.	 4386408:BVX772-01	OF-FB41-0216
9.	 4386408:BVX773-01	OF-RW41-0216
10.	 4386408:BVX774-01	OF-RW41P-0216
11.	 4386408:BVX775-01	OF-FB56-0216
12.	 4386408:BVX776-01	OF-RW56-0216
13.	 4386408:BVX779-01	OF-FB51-0216
14.	 4386408:BVX780-01	OF-RW51-0216
15.	 4386408:BVX781-01	OF-RW51P-0216
16.	 4386408:BVX826-01	OF-RW09-0216
17.	 4386408:BVX827-01	OF-RW37-0216
18.	 4386408:BVX828-01	OF-RW11-0216
19.	 4386408:BVX829-01	OF-RW28-0216
20.	 4386408:BVX830-01	OF-RW67-0216

Worksheet Reagent Tracking Record

Worksheet # 4386408

Surrogate/Spike solutions	✓	Solution ID #		Conc.	Blk-Spk		MS		Samples	
					Solid	Liquid	Solid	Liquid	Solid	Liquid
DGT Spike				100 ug/mL	60	30	60	30	NA	NA
Diquat Dibromide				50 ug/mL	NA	350	NA	350	NA	NA
Explosives Spiking solution A				20 ug/mL	250	100	250	100	NA	NA
Explosives Spiking solution B				20/80 ug/mL	250	100	250	100	NA	NA
Formaldehyde Spike				100 ug/mL	25	25	25	25	NA	NA
Glyphosate Spike				25 ug/mL	500	20	500	20	NA	NA
Nonylphenol Ethoxylate Spike				100 ug/mL	100	100	100	100	NA	NA
Nonylphenol Spike				10 ug/mL	100	100	100	100	NA	NA
Paraquat Cl Tetrahydrate				20 ug/mL	NA	125	NA	125	NA	NA
Perchlorate Standard Spike				10 ng/mL	NA	100	NA	100	NA	NA
Perchlorate Standard Spike				500 ng/mL	40	NA	40	NA	NA	NA
Perchlorate O-18 Internal Standard				0.10 ng/uL	20	20	20	20	20	20
Morpholine Intermediate Std.				5 ug/mL	NA	50	NA	20	NA	NA
Morpholine-D8 Internal Standard				10 ug/mL	NA	100	NA	100	NA	100
Comp. PFC Spiking Solution A				1ug/mL	62.5	NA	62.5	NA	NA	NA
Comp. PFC Spiking Solution B				250 ng/mL	NA	40 .75	NA	40 .75	NA	NA
Comp. PFC Spiking Solution C	/	SK-6229		100 ng/mL	125	(62-5) NA	125	(62-5) NA	NA	NA
Internal Standard Solution A	/	SF-6010		50 ng/mL	50	(15) 100	50	(15) 100	50	(15) 100
Internal Standard Solution B				250ng/mL	50	NA	50	NA	50	NA
ICV/OCV PFC	/	I-4676		1ug/mL		(62-5)				
Solvent/Reagent	Suppller	✓	Lot No.	Date Opened	Solvent/Reagent	✓	Lot No.	Date Opened/ Prepared	*Spiked by: 65Z	
DCM	Fisher				50% NaOH				Spike Date 2016/02/18	
Hexane	Fisher				20mM TBAS				Spike Syringe ID# <i>Pijjarter</i>	
Acetone	Fisher				o-Phosphoric Acid				M23487B	
Ottawa Sand	Fisher				Borax				Int. Std Syringe ID# <i>Pijjarter</i>	
Methanol	<i>Wako</i> Fisher	/	SHB66076V	2016/02/18	Calcium Chloride				M23487B	
2-Propanol (IPA)	Fisher				EDTA				*Spiking Witnessed by: SSV	
Acetonitrile	Fisher				Phosphate Buffer				Final pH	
MTBE	Fisher				Sodium Thiosulphate				X	
Sodium Sulfate <i>60.40 H2O MeOH</i> Recon Solution	Fisher	/	87-190	2016/02/12	5M Acetate Buffer					
DCM:Ethyl Ether (75:25)					FMOc					
Hexane:IPA (98:2)					0.25M Na <sub>2</sub> CO <sub>3</sub>					
2% Formic Acid		/	87-131	2016/01/13	0.5M TBAS					
0.2% Formic Acid		/	87-114	2015/12/31	1% NH <sub>4</sub> OH 0.2%	/	87-198	2016/02/18		
0.05M KOH					Leachate Fluid					
0.05M HCl					Reagent Water	/	158399	2016/02/17		
Equipment	ID#	✓	Equipment	ID#	✓	Equipment	Lot #	Bottle Tracking		
Pipettor	M15944C P29746D K19609D	/	SPE Cartridge			10 mL Serological Pipet		Bottle#	14991	
		/	Filter			QC Balance ID		Cap#	14443	
Dispenser			Centrifuge			Thermometer ID &Temp	SN 140614017 45°C	Systems plus Lot#	16-01-06	
Syringe			Sonicator							

Comments: 125ml Ref Bottle (HPPB)  
Inj IS - SF-6017 - 2016/01/19  
HPLC Evaporator

\* - SPIKING OF REAGENTS MUST BE WITNESSED AT ALL TIMES.



Project: D:\Analyst Data\Projects\Enviro\FPOS Batch:FPC\_160219 Tab:Sample Set:SET1 AcqMethod:FPC\_Water\_Low.dam

Sample	Sample Name	Rack Code	Rack Position	Plate Code	Plate Position	Vial Position	Data File	Inj. Volume (µl)
1	Rinse	2 Well Plates	1	*54VialPlate*	1	1	PFC_160219WS#4386408	3.000
2	4386408-BLANK	2 Well Plates	1	*54VialPlate*	1	2	PFC_160219WS#4386408	3.000
3	STD 1	2 Well Plates	1	*54VialPlate*	1	3	PFC_160219WS#4386408	3.000
4	STD 2	2 Well Plates	1	*54VialPlate*	1	4	PFC_160219WS#4386408	3.000
5	STD 3	2 Well Plates	1	*54VialPlate*	1	5	PFC_160219WS#4386408	3.000
6	STD 4	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4386408	3.000
7	STD 5	2 Well Plates	1	*54VialPlate*	1	7	PFC_160219WS#4386408	3.000
8	STD 6	2 Well Plates	1	*54VialPlate*	1	8	PFC_160219WS#4386408	3.000
9	KCV	2 Well Plates	1	*54VialPlate*	1	9	PFC_160219WS#4386408	3.000
10	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4386408	3.000
11	4386408-MTRX SPK	2 Well Plates	1	*54VialPlate*	1	10	PFC_160219WS#4386408	3.000
12	4386408-MTRX SPK:D1	2 Well Plates	1	*54VialPlate*	1	11	PFC_160219WS#4386408	3.000
13	4386408-SPKK	2 Well Plates	1	*54VialPlate*	1	12	PFC_160219WS#4386408	3.000
14	4386408-BVX814-01 (10x)	2 Well Plates	1	*54VialPlate*	1	13	PFC_160219WS#4386408	3.000
15	4386408-BVX817-01 (10x)	2 Well Plates	1	*54VialPlate*	1	14	PFC_160219WS#4386408	3.000
16	4386408-BVX818-01 (20x)	2 Well Plates	1	*54VialPlate*	1	15	PFC_160219WS#4386408	3.000
17	4386408-BVX819-01 (10x)	2 Well Plates	1	*54VialPlate*	1	16	PFC_160219WS#4386408	3.000
18	4386408-BVX769-01	2 Well Plates	1	*54VialPlate*	1	17	PFC_160219WS#4386408	3.000
19	4386408-BVX770-01	2 Well Plates	1	*54VialPlate*	1	18	PFC_160219WS#4386408	3.000
20	4386408-BVX771-01	2 Well Plates	1	*54VialPlate*	1	19	PFC_160219WS#4386408	3.000
21	4386408-BVX772-01	2 Well Plates	1	*54VialPlate*	1	20	PFC_160219WS#4386408	3.000
22	4386408-BVX773-01	2 Well Plates	1	*54VialPlate*	1	21	PFC_160219WS#4386408	3.000
23	4386408-BVX774-01	2 Well Plates	1	*54VialPlate*	1	22	PFC_160219WS#4386408	3.000
24	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4386408	3.000
25	4386408-BVX775-01	2 Well Plates	1	*54VialPlate*	1	23	PFC_160219WS#4386408	3.000
26	4386408-BVX776-01	2 Well Plates	1	*54VialPlate*	1	24	PFC_160219WS#4386408	3.000
27	4386408-BVX779-01	2 Well Plates	1	*54VialPlate*	1	25	PFC_160219WS#4386408	3.000
28	4386408-BVX780-01	2 Well Plates	1	*54VialPlate*	1	26	PFC_160219WS#4386408	3.000
29	4386408-BVX781-01	2 Well Plates	1	*54VialPlate*	1	27	PFC_160219WS#4386408	3.000
30	4386408-BVX826-01	2 Well Plates	1	*54VialPlate*	1	28	PFC_160219WS#4386408	3.000
31	4386408-BVX827-01	2 Well Plates	1	*54VialPlate*	1	29	PFC_160219WS#4386408	3.000
32	4386408-BVX828-01	2 Well Plates	1	*54VialPlate*	1	30	PFC_160219WS#4386408	3.000
33	4386408-BVX829-01	2 Well Plates	1	*54VialPlate*	1	31	PFC_160219WS#4386408	3.000
34	4386408-BVX830-01	2 Well Plates	1	*54VialPlate*	1	32	PFC_160219WS#4386408	3.000
35	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4386408	3.000
36	4385924-BLANK	2 Well Plates	1	*54VialPlate*	1	33	PFC_160219WS#4385924	3.000
37	4385924-MTRX SPK	2 Well Plates	1	*54VialPlate*	1	34	PFC_160219WS#4385924	3.000
38	4385924-MTRX SPK:D1	2 Well Plates	1	*54VialPlate*	1	35	PFC_160219WS#4385924	3.000
39	4385924-SPKK	2 Well Plates	1	*54VialPlate*	1	36	PFC_160219WS#4385924	3.000
40	4385924-BVX947-01 (10x)	2 Well Plates	1	*54VialPlate*	1	37	PFC_160219WS#4385924	3.000
41	4385924-BVX792-01	2 Well Plates	1	*54VialPlate*	1	38	PFC_160219WS#4385924	3.000
42	4385924-BVX793-01	2 Well Plates	1	*54VialPlate*	1	39	PFC_160219WS#4385924	3.000
43	4385924-BVX794-01	2 Well Plates	1	*54VialPlate*	1	40	PFC_160219WS#4385924	3.000
44	4385924-BVX795-01	2 Well Plates	1	*54VialPlate*	1	41	PFC_160219WS#4385924	3.000
45	4385924-BVX796-01	2 Well Plates	1	*54VialPlate*	1	42	PFC_160219WS#4385924	3.000
46	4385924-BVX797-01	2 Well Plates	1	*54VialPlate*	1	43	PFC_160219WS#4385924	3.000
47	4385924-BVX798-01	2 Well Plates	1	*54VialPlate*	1	44	PFC_160219WS#4385924	3.000
48	4385924-BVX799-01	2 Well Plates	1	*54VialPlate*	1	45	PFC_160219WS#4385924	3.000
49	4385924-BVX800-01	2 Well Plates	1	*54VialPlate*	1	46	PFC_160219WS#4385924	3.000
50	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4385924	3.000
51	4385924-BVX801-01	2 Well Plates	1	*54VialPlate*	1	47	PFC_160219WS#4385924	3.000
52	4385924-BVX802-01	2 Well Plates	1	*54VialPlate*	1	48	PFC_160219WS#4385924	3.000
53	4385924-BVX803-01	2 Well Plates	1	*54VialPlate*	1	49	PFC_160219WS#4385924	3.000
54	4385924-BVX804-01	2 Well Plates	1	*54VialPlate*	1	50	PFC_160219WS#4385924	3.000
55	4385924-BVX805-01	2 Well Plates	1	*54VialPlate*	1	51	PFC_160219WS#4385924	3.000
56	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4385924	3.000
57	4385888-BLANK	2 Well Plates	1	*54VialPlate*	2	1	PFC_160219WS#4385888	3.000
58	4385888-MTRX SPK	2 Well Plates	1	*54VialPlate*	2	2	PFC_160219WS#4385888	3.000
59	4385888-MTRX SPK:D1	2 Well Plates	1	*54VialPlate*	2	3	PFC_160219WS#4385888	3.000
60	4385888-SPKK	2 Well Plates	1	*54VialPlate*	2	4	PFC_160219WS#4385888	3.000
61	4385888-BVX782-01	2 Well Plates	1	*54VialPlate*	2	5	PFC_160219WS#4385888	3.000
62	4385888-BVX783-01	2 Well Plates	1	*54VialPlate*	2	6	PFC_160219WS#4385888	3.000
63	4385888-BVX784-01	2 Well Plates	1	*54VialPlate*	2	7	PFC_160219WS#4385888	3.000
64	4385888-BVX785-01	2 Well Plates	1	*54VialPlate*	2	8	PFC_160219WS#4385888	3.000
65	4385888-BVX786-01	2 Well Plates	1	*54VialPlate*	2	9	PFC_160219WS#4385888	3.000
66	4385888-BVX787-01	2 Well Plates	1	*54VialPlate*	2	10	PFC_160219WS#4385888	3.000
67	4385888-BVX788-01	2 Well Plates	1	*54VialPlate*	2	11	PFC_160219WS#4385888	3.000
68	4385888-BVX789-01	2 Well Plates	1	*54VialPlate*	2	12	PFC_160219WS#4385888	3.000
69	4385888-BVX790-01	2 Well Plates	1	*54VialPlate*	2	13	PFC_160219WS#4385888	3.000
70	4385888-BVX791-01	2 Well Plates	1	*54VialPlate*	2	14	PFC_160219WS#4385888	3.000
71	CCV	2 Well Plates	1	*54VialPlate*	1	6	PFC_160219WS#4385888	3.000

Column #123  
 MPA = Soln # 506, FJRE17  
 MEB = MeOH, Fulo-L #1526B  
 on 201602/19

Results Name: PFC\_Water\_160219\_4386408\_ULow.rdb  
 Results Path: \\miss-netapp2\lcm\lcm3\Analyst  
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Sample ID	Acquisition Date	Acquisition Method	File Name	Rack Type	Rack Position	Vial Position	Plate Type	Plate Position
1	2016/02/19 7:39:53 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	2	*54VialPlate*	1
2	2016/02/19 7:45:04 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	3	*54VialPlate*	1
3	2016/02/19 7:50:14 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	4	*54VialPlate*	1
4	2016/02/19 7:55:19 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	5	*54VialPlate*	1
5	2016/02/19 8:00:24 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	6	*54VialPlate*	1
6	2016/02/19 8:05:30 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	7	*54VialPlate*	1
7	2016/02/19 8:10:36 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	8	*54VialPlate*	1
8	2016/02/19 8:15:41 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	9	*54VialPlate*	1
9	2016/02/19 8:20:47 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	6	*54VialPlate*	1
10	2016/02/19 8:25:52 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	10	*54VialPlate*	1
11	2016/02/19 8:30:58 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	11	*54VialPlate*	1
12	2016/02/19 8:36:03 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	12	*54VialPlate*	1
13	2016/02/19 8:51:19 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	13	*54VialPlate*	1
14	2016/02/19 8:56:24 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	14	*54VialPlate*	1
15	2016/02/19 9:01:31 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	15	*54VialPlate*	1
16	2016/02/19 9:06:37 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	16	*54VialPlate*	1
17	2016/02/19 9:11:43 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	17	*54VialPlate*	1
18	2016/02/19 9:16:48 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	18	*54VialPlate*	1
19	2016/02/19 9:21:58 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	19	*54VialPlate*	1
20	2016/02/19 9:27:04 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	20	*54VialPlate*	1
21	2016/02/19 9:32:10 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	21	*54VialPlate*	1
22	2016/02/19 9:37:16 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	22	*54VialPlate*	1
23	2016/02/19 9:42:22 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	6	*54VialPlate*	1
24	2016/02/19 9:52:33 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	23	*54VialPlate*	1
25	2016/02/19 9:57:39 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	24	*54VialPlate*	1
26	2016/02/19 10:02:45 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	25	*54VialPlate*	1
27	2016/02/19 10:07:51 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	26	*54VialPlate*	1
28	2016/02/19 10:12:57 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	27	*54VialPlate*	1
29	2016/02/19 10:18:03 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	28	*54VialPlate*	1
30	2016/02/19 10:23:09 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	29	*54VialPlate*	1
31	2016/02/19 10:28:14 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	30	*54VialPlate*	1
32	2016/02/19 10:33:19 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	31	*54VialPlate*	1
33	2016/02/19 10:38:25 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	32	*54VialPlate*	1
34	2016/02/19 10:43:32 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	6	*54VialPlate*	1
35	2016/02/19 10:48:37 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	10	*54VialPlate*	1
36	2016/02/19 10:58:48 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	32	*54VialPlate*	1
37	2016/02/19 11:03:54 AM	PFC_Water_Low.dam	PFC_160219\WS#\4386408.wiff	2 Well Plates	1	6	*54VialPlate*	1

Maxxim Analytics



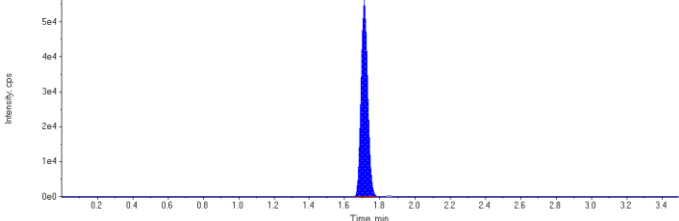
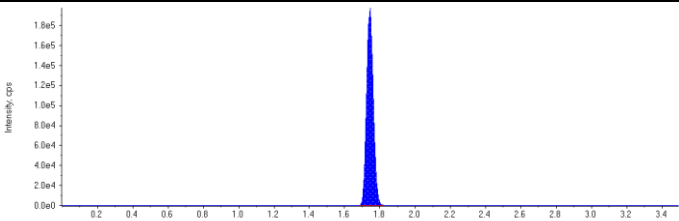
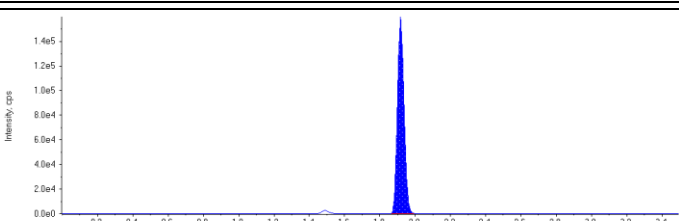
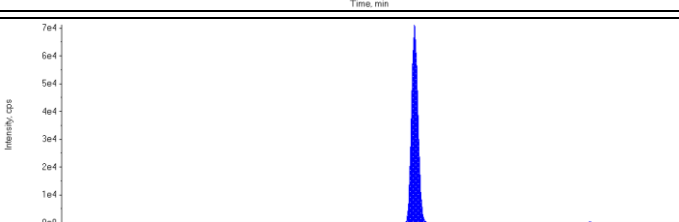
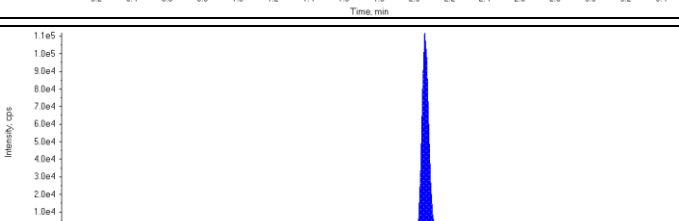
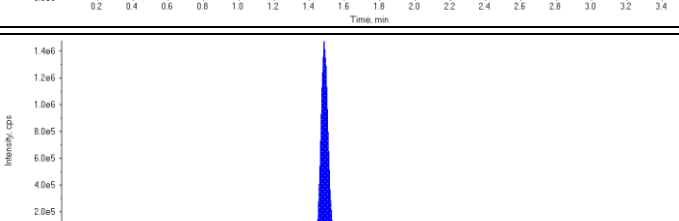
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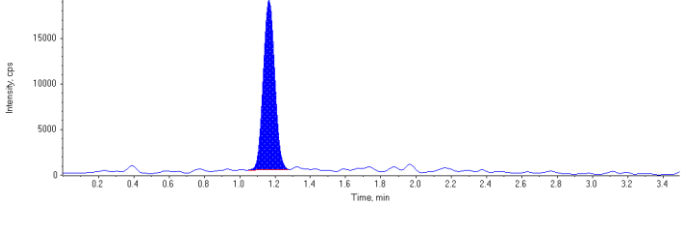
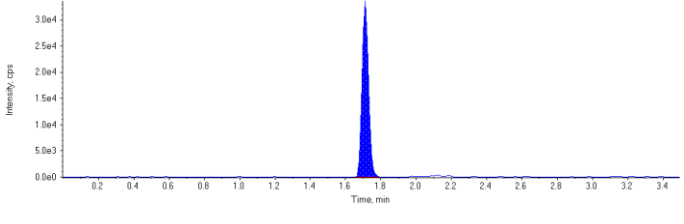
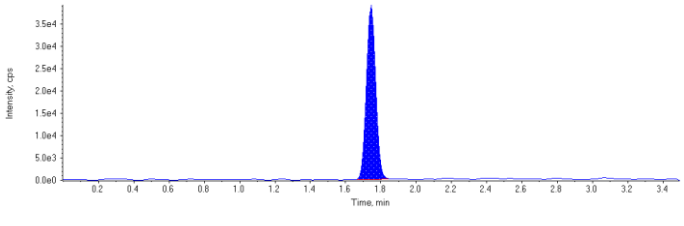
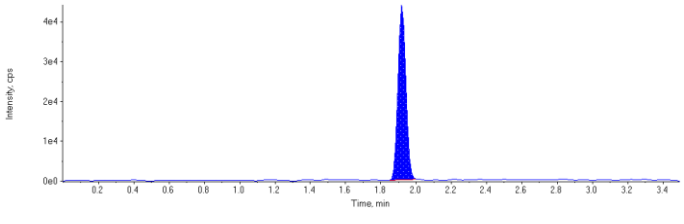
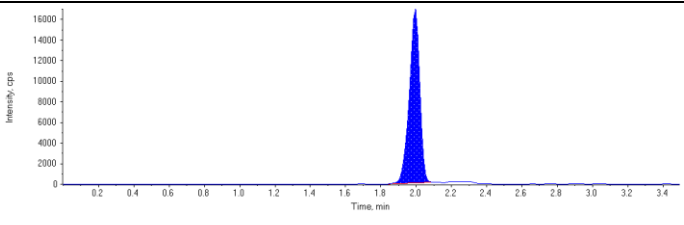
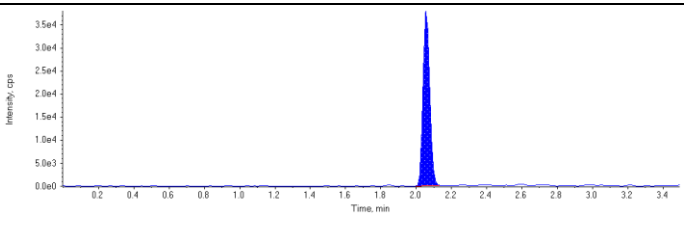
Maxxam Analytics International  
6740 Campobello Rd  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

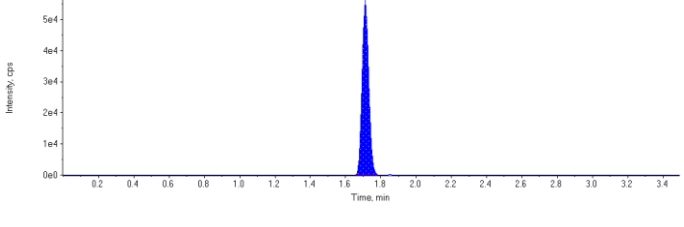
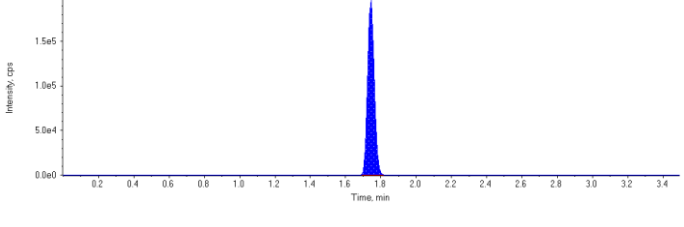
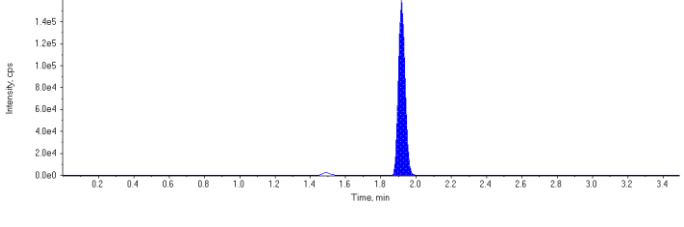
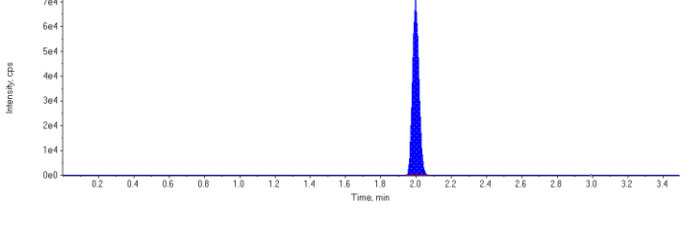
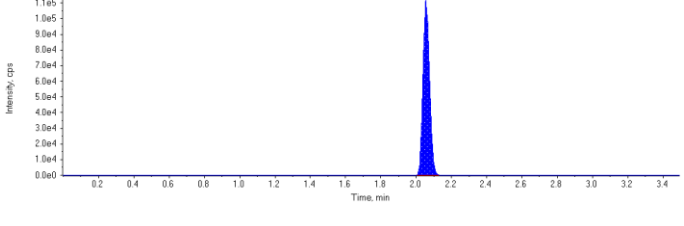
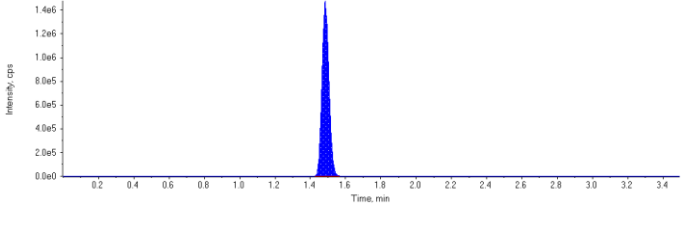
<b>Sample Name</b>	STD 1	<b>Injection Vial</b>	3
<b>Sample ID</b>	STD 1	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 7:45:04 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	146000.	1.71	1.00	-
MPFHpA	517000.	1.74	1.00	-
MPFOA	413000.	1.92	1.00	-
MPFOS	179000.	2.00	1.00	-
MPFNA	291000.	2.06	1.00	-
13C6-PFHxA IS	4050000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	83700	1.17	2.00	2.40	120.0
PFHxS 1	85100	1.71	2.00	2.35	117.0
PFHpA 1	142000	1.74	2.00	2.11	106.0
PFOA 1	136000	1.92	2.00	2.20	110.0
PFOS 1	69300	1.99	2.00	2.31	116.0
PFNA 1	101000	2.06	2.00	2.24	112.0
18O2-PFHxS	146000	1.71	100.	97.9	97.9
13C4-PFHpA	517000	1.74	100.	103.	103.0
13C4-PFOA	413000	1.92	100.	94.5	94.5
13C4-PFOS	179000	2.00	100.	102.	102.0
13C5-PFNA	291000	2.06	100.	97.6	97.6
13C6-PFHxA	4050000	1.49	100.	101.	101.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1 (298.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.17 (1.15) min</p> <p>Calculated Conc: 2.40 µg/L</p> <p>Area Ratio: 0.573</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1 (398.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 2.35 µg/L</p> <p>Area Ratio: 0.583</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1 (363.000/319.000 Da)</b></p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 2.11 µg/L</p> <p>Area Ratio: 0.275</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1 (413.100/369.000 Da)</b></p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 2.20 µg/L</p> <p>Area Ratio: 0.331</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1 (498.900/79.900 Da)</b></p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 2.31 µg/L</p> <p>Area Ratio: 0.387</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1 (462.900/419.000 Da)</b></p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 2.24 µg/L</p> <p>Area Ratio: 0.346</p> <p>Sample Type: (Standard)</p>	

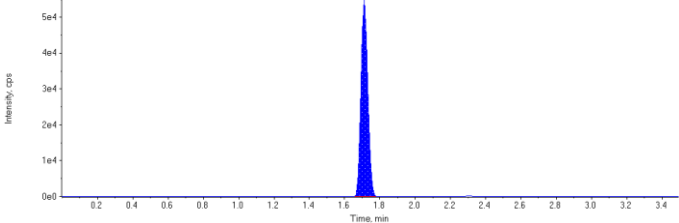
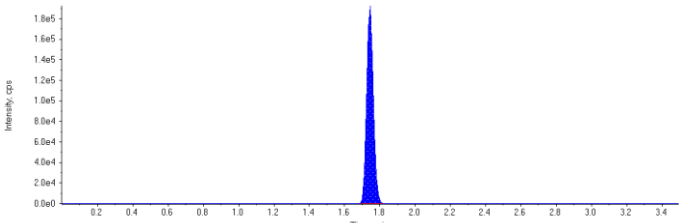
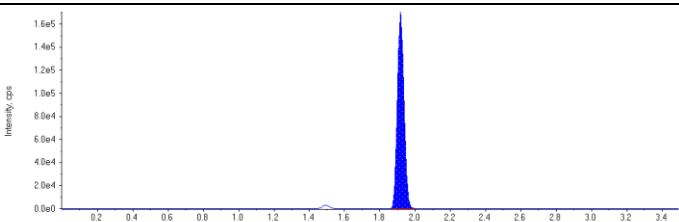
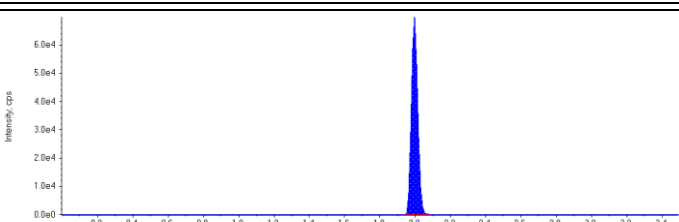
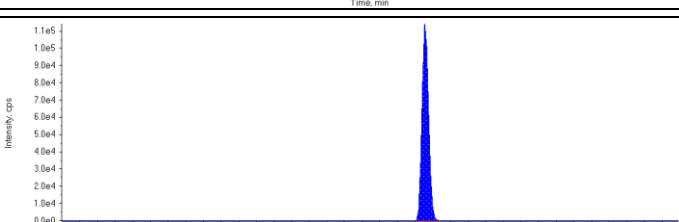
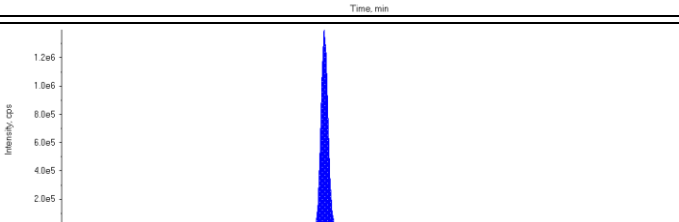
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 97.9 µg/L</p> <p>Area Ratio: 0.0361</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.128</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 94.5 µg/L</p> <p>Area Ratio: 0.102</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 102. µg/L</p> <p>Area Ratio: 0.0443</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 97.6 µg/L</p> <p>Area Ratio: 0.0719</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 101. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

<b>Sample Name</b>	STD 2	<b>Injection Vial</b>	4
<b>Sample ID</b>	STD 2	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 7:50:14 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
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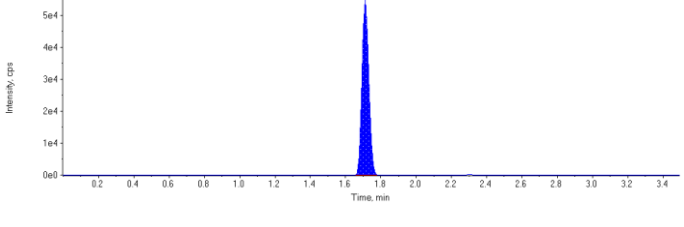
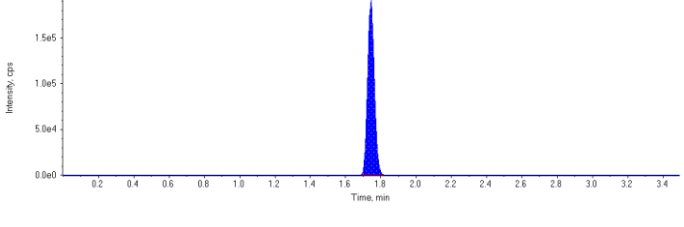
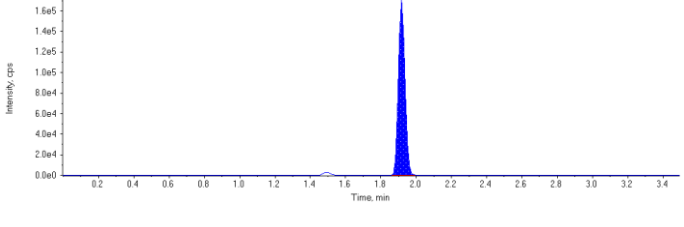
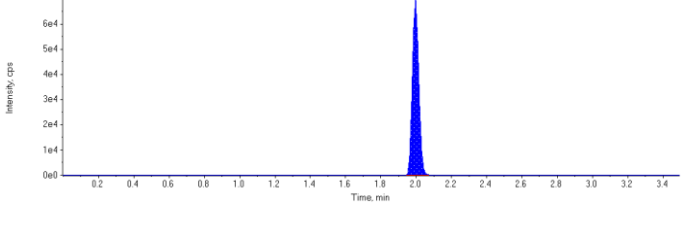
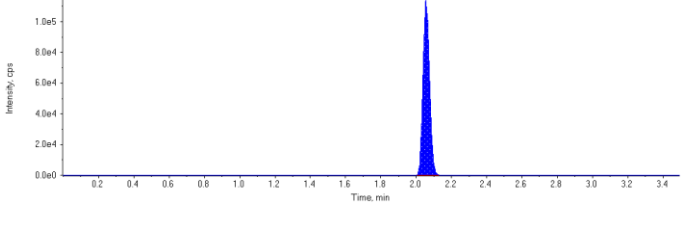
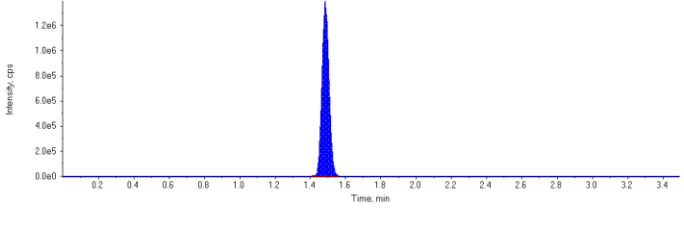
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	147000.	1.71	1.00	-
MPFHpA	503000.	1.74	1.00	-
MPFOA	442000.	1.92	1.00	-
MPFOS	177000.	2.00	1.00	-
MPFNA	296000.	2.06	1.00	-
13C6-PFHxA IS	3880000.	1.49	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	151000	1.16	4.00	3.86	96.4
PFHxS 1	151000	1.71	4.00	3.68	92.1
PFHpA 1	247000	1.74	4.00	3.72	92.9
PFOA 1	258000	1.92	4.00	3.73	93.3
PFOS 1	111000	1.99	4.00	3.63	90.7
PFNA 1	177000	2.06	4.00	3.74	93.6
18O2-PFHxS	147000	1.71	100.	103.	103.0
13C4-PFHpA	503000	1.74	100.	105.	105.0
13C4-PFOA	442000	1.92	100.	105.	105.0
13C4-PFOS	177000	2.00	100.	105.	105.0
13C5-PFNA	296000	2.06	100.	104.	104.0
13C6-PFHxA	3880000	1.49	100.	97.3	97.3



<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.49(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

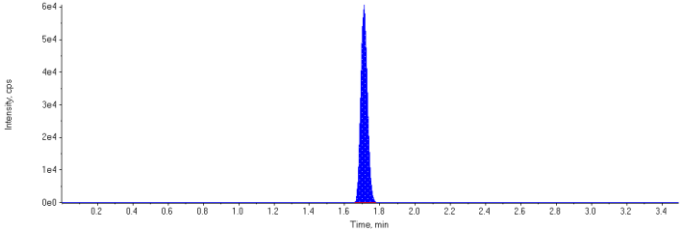
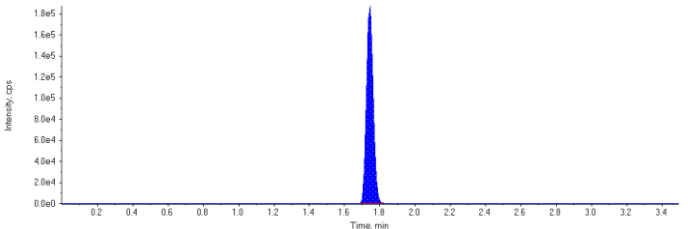
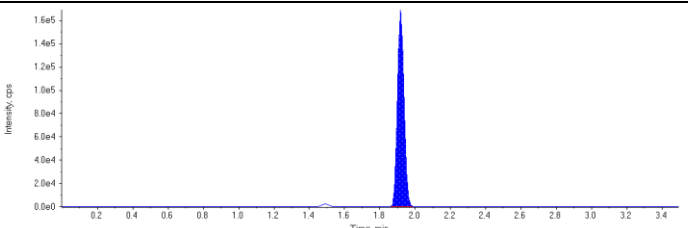
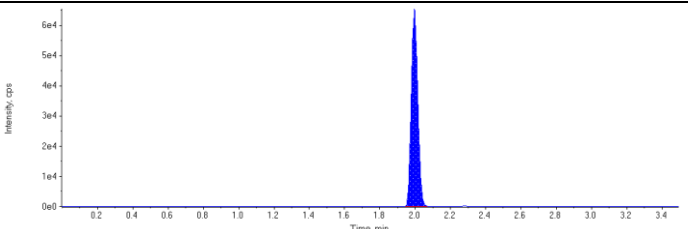
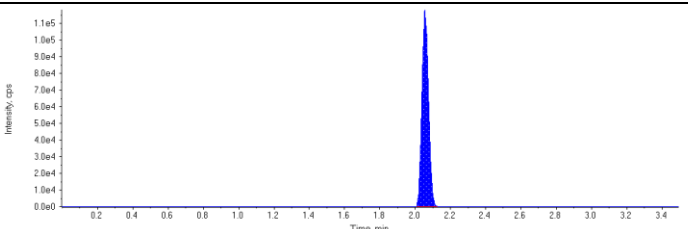
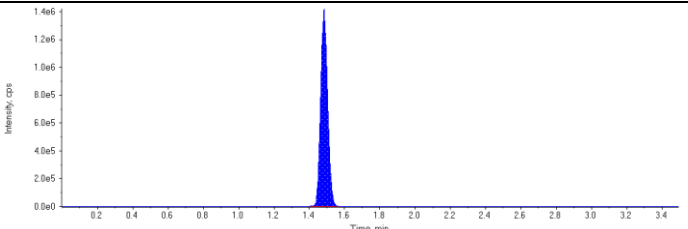
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 3.86 µg/L</p> <p>Area Ratio: 1.03</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 3.68 µg/L</p> <p>Area Ratio: 1.03</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 3.72 µg/L</p> <p>Area Ratio: 0.491</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 3.73 µg/L</p> <p>Area Ratio: 0.584</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 3.63 µg/L</p> <p>Area Ratio: 0.627</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 3.74 µg/L</p> <p>Area Ratio: 0.597</p> <p>Sample Type: (Standard)</p>	

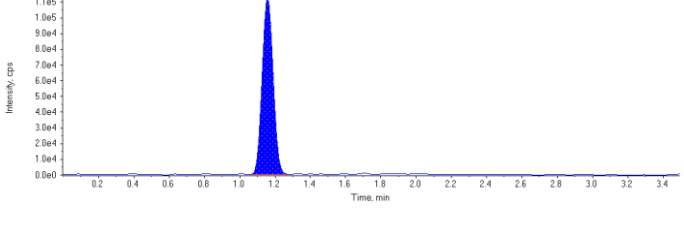
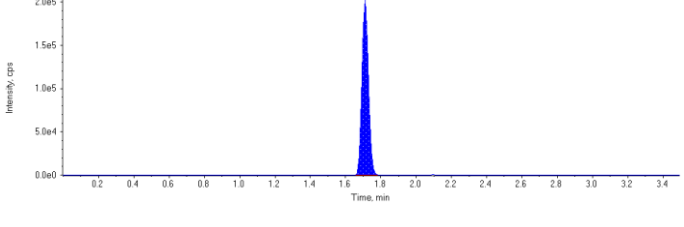
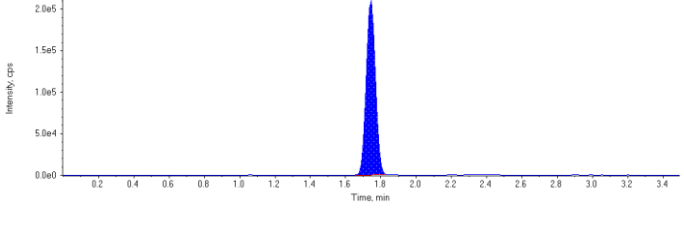
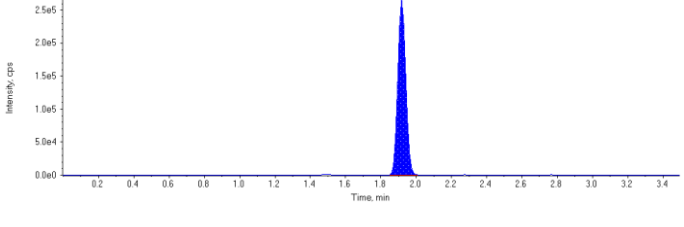
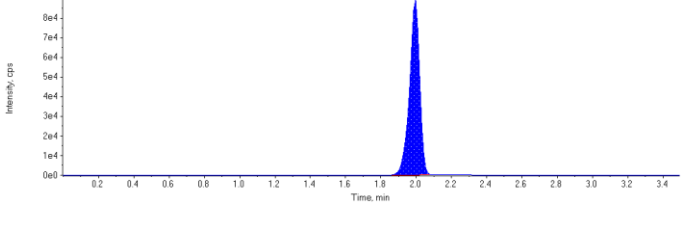
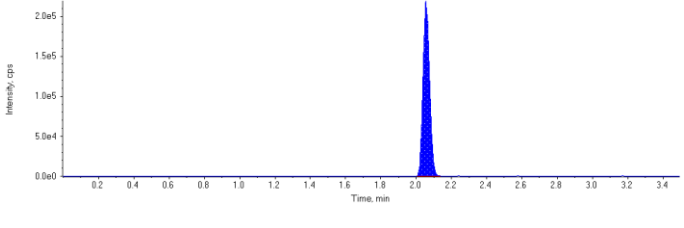
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.0378</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 105. µg/L</p> <p>Area Ratio: 0.129</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 105. µg/L</p> <p>Area Ratio: 0.114</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 105. µg/L</p> <p>Area Ratio: 0.0456</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.0763</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.49 (1.42) min</p> <p>Calculated Conc: 97.3 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

<b>Sample Name</b>	<b>STD 3</b>	<b>Injection Vial</b>	5
<b>Sample ID</b>	STD 3	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 7:55:19 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	156000.	1.71	1.00	-
MPFHpA	498000.	1.74	1.00	-
MPFOA	435000.	1.92	1.00	-
MPFOS	167000.	2.00	1.00	-
MPFNA	307000.	2.06	1.00	-
13C6-PFHxA IS	3890000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	482000	1.16	12.0	10.4	87.0
PFHxS 1	518000	1.71	12.0	10.6	88.1
PFHpA 1	764000	1.74	12.0	11.5	95.5
PFOA 1	835000	1.92	12.0	11.8	98.4
PFOS 1	356000	1.99	12.0	11.9	99.2
PFNA 1	570000	2.06	12.0	11.3	94.2
18O2-PFHxS	156000	1.71	100.	109.	109.0
13C4-PFHpA	498000	1.74	100.	104.	104.0
13C4-PFOA	435000	1.92	100.	104.	104.0
13C4-PFOS	167000	2.00	100.	99.4	99.4
13C5-PFNA	307000	2.06	100.	107.	107.0
13C6-PFHxA	3890000	1.48	100.	97.4	97.4

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 10.4 µg/L</p> <p>Area Ratio: 3.09</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 10.6 µg/L</p> <p>Area Ratio: 3.32</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 11.5 µg/L</p> <p>Area Ratio: 1.53</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 11.8 µg/L</p> <p>Area Ratio: 1.92</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 11.9 µg/L</p> <p>Area Ratio: 2.13</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 11.3 µg/L</p> <p>Area Ratio: 1.86</p> <p>Sample Type: (Standard)</p>	

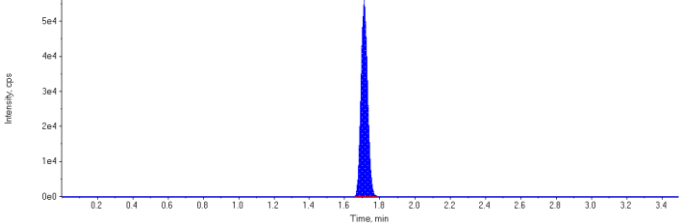
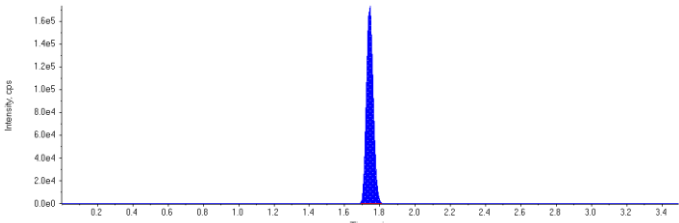
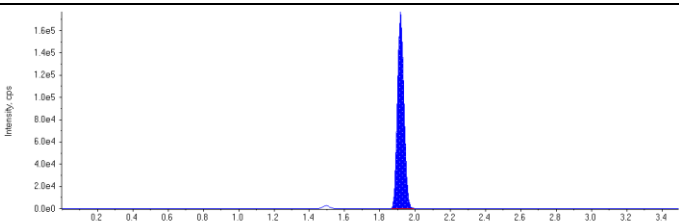
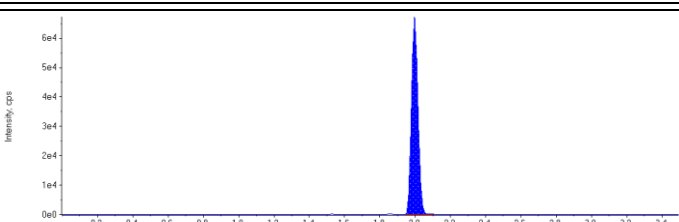
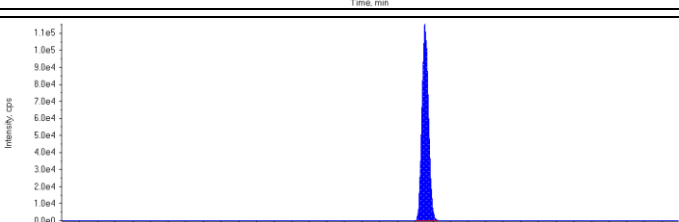
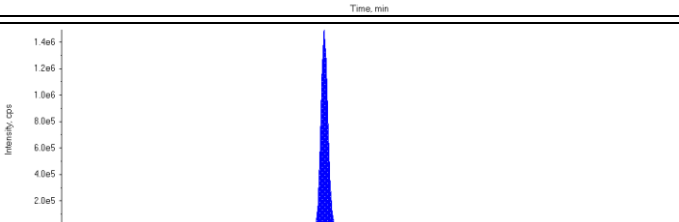
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 109. µg/L</p> <p>Area Ratio: 0.0401</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.128</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.112</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 99.4 µg/L</p> <p>Area Ratio: 0.0429</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 107. µg/L</p> <p>Area Ratio: 0.0791</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 97.4 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

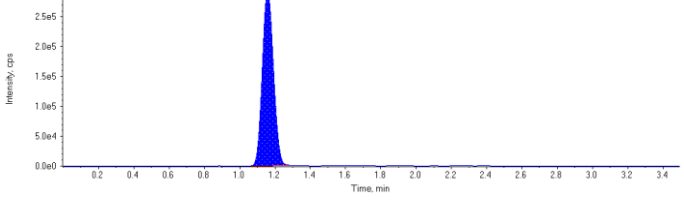
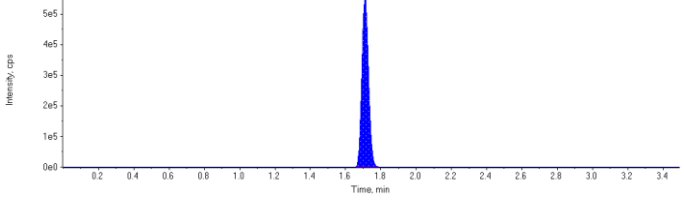
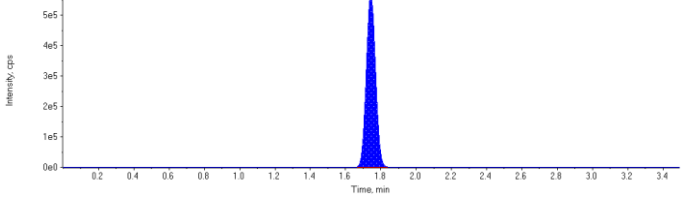
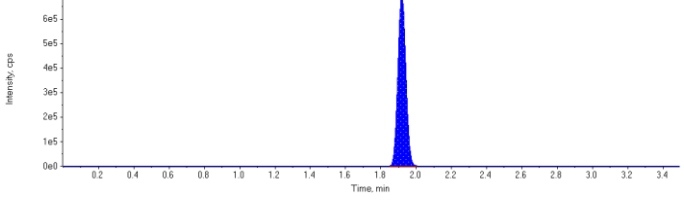
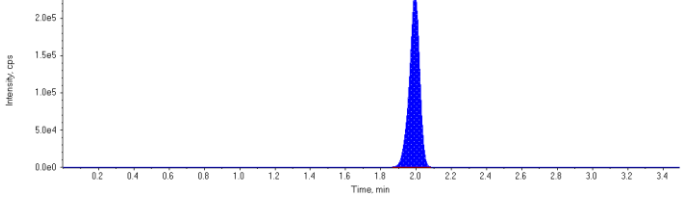
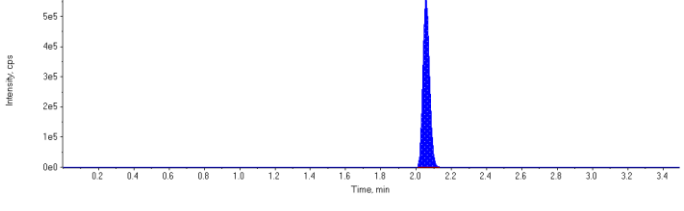
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<b>Sample ID</b>	STD 4	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:00:24 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

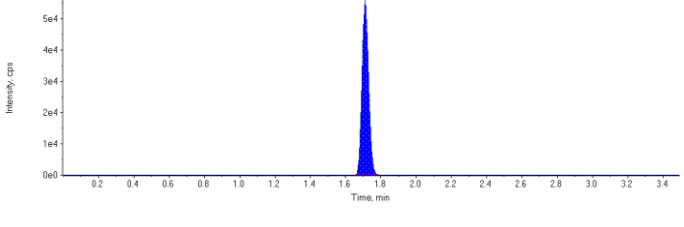
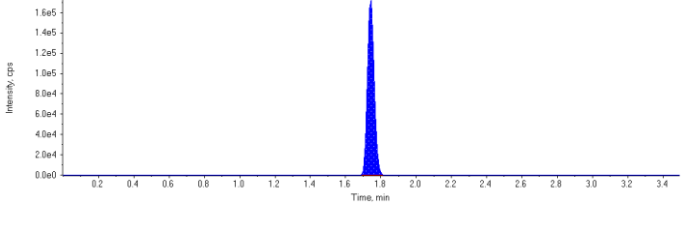
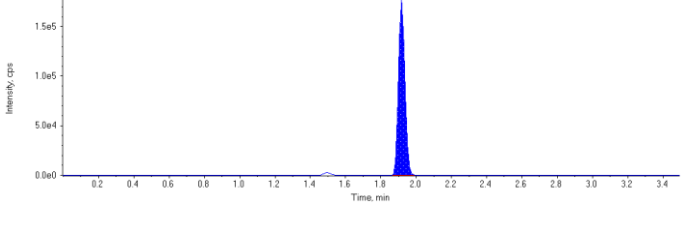
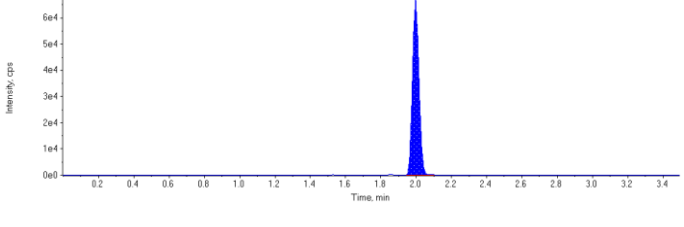
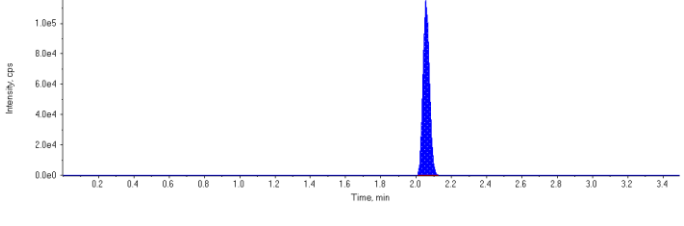
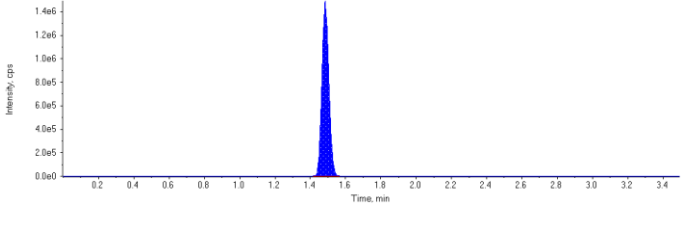
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	144000.	1.71	1.00	-
MPFHpA	467000.	1.74	1.00	-
MPFOA	447000.	1.92	1.00	-
MPFOS	172000.	2.00	1.00	-
MPFNA	296000.	2.06	1.00	-
13C6-PFHxA IS	4090000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	1290000	1.16	30.0	29.3	97.5
PFHxS 1	1470000	1.71	30.0	31.3	104.0
PFHpA 1	2080000	1.74	30.0	33.2	111.0
PFOA 1	2220000	1.92	30.0	30.3	101.0
PFOS 1	944000	1.99	30.0	30.3	101.0
PFNA 1	1500000	2.06	30.0	30.6	102.0
18O2-PFHxS	144000	1.71	100.	95.4	95.4
13C4-PFHpA	467000	1.74	100.	92.3	92.3
13C4-PFOA	447000	1.92	100.	101.	101.0
13C4-PFOS	172000	2.00	100.	97.3	97.3
13C5-PFNA	296000	2.06	100.	98.3	98.3
13C6-PFHxA	4090000	1.48	100.	102.	102.0



<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

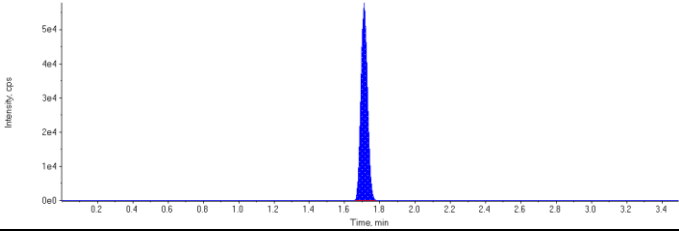
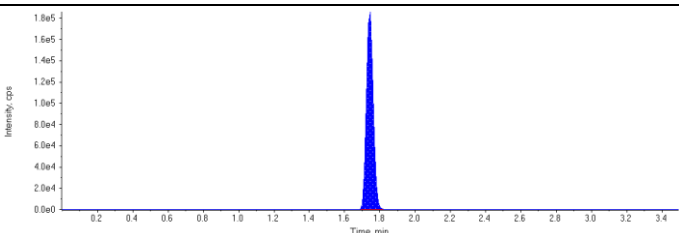
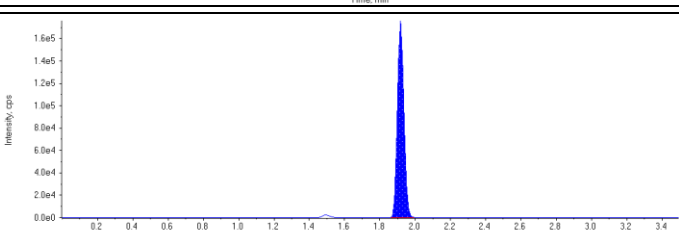
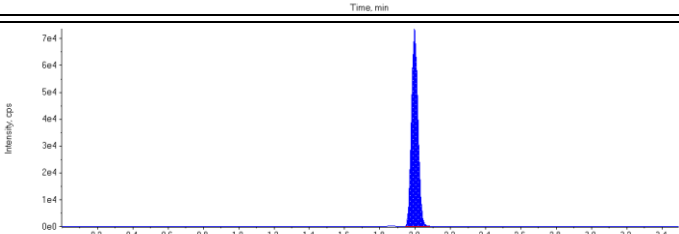
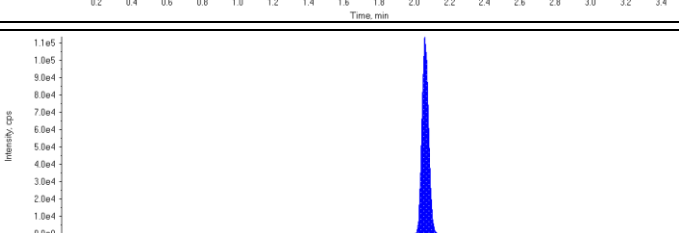
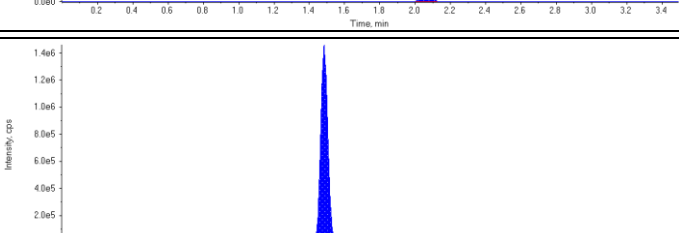
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 29.3 µg/L</p> <p>Area Ratio: 8.98</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 31.3 µg/L</p> <p>Area Ratio: 10.2</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 33.2 µg/L</p> <p>Area Ratio: 4.46</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 30.3 µg/L</p> <p>Area Ratio: 4.97</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 30.3 µg/L</p> <p>Area Ratio: 5.49</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 30.6 µg/L</p> <p>Area Ratio: 5.07</p> <p>Sample Type: (Standard)</p>	

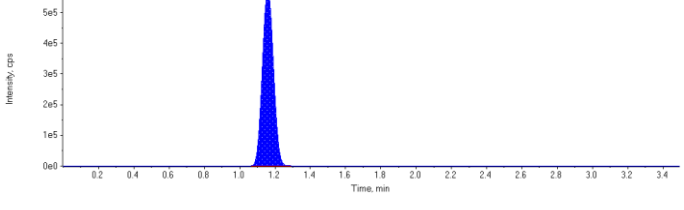
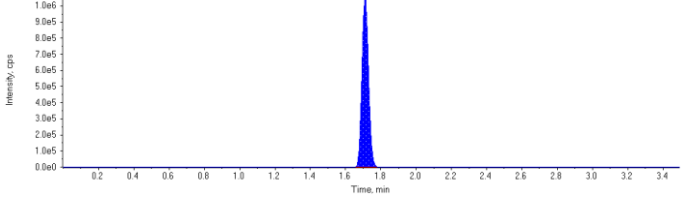
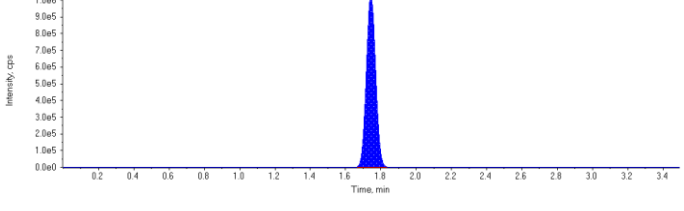
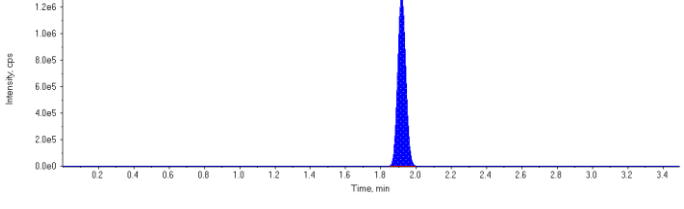
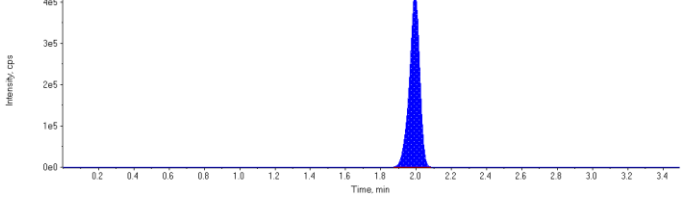
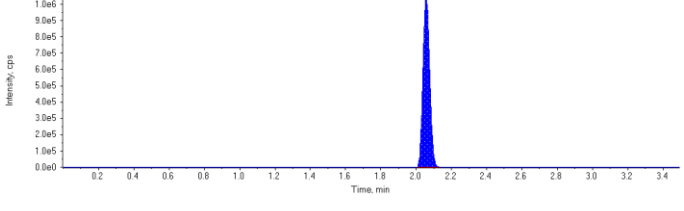
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 95.4 µg/L</p> <p>Area Ratio: 0.0351</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 92.3 µg/L</p> <p>Area Ratio: 0.114</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 101. µg/L</p> <p>Area Ratio: 0.109</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 97.3 µg/L</p> <p>Area Ratio: 0.0420</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 98.3 µg/L</p> <p>Area Ratio: 0.0723</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 102. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

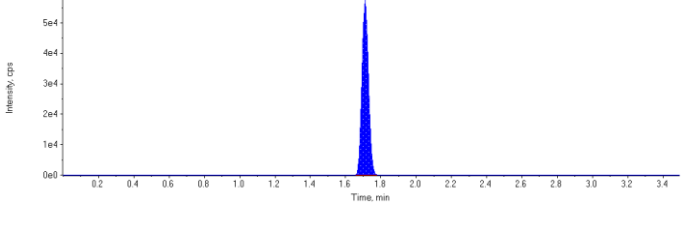
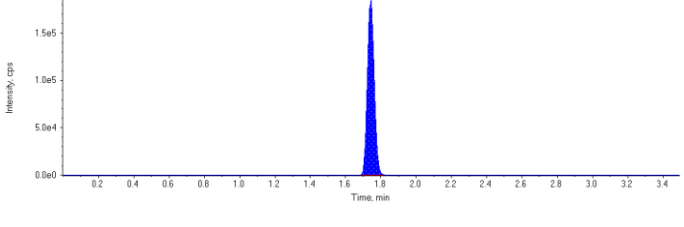
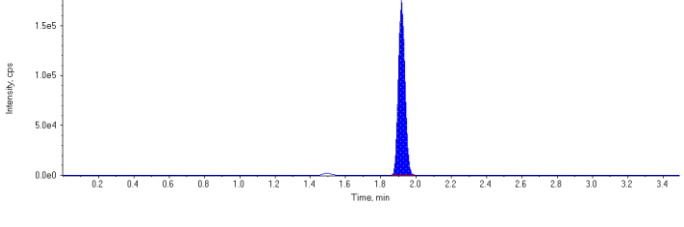
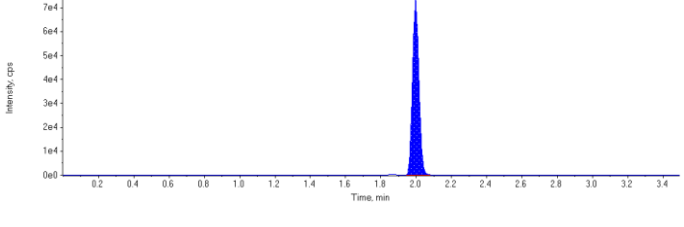
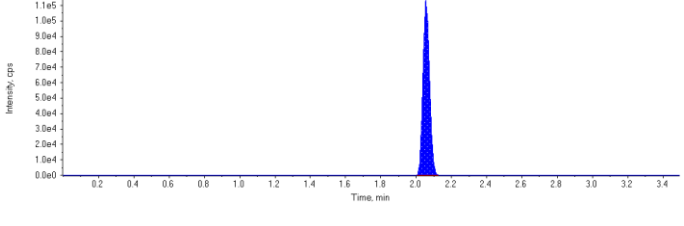
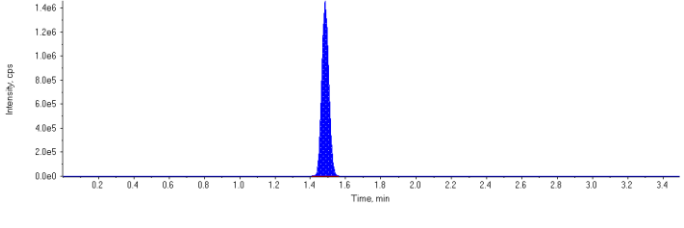
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<b>Sample ID</b>	STD 5	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:05:30 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	147000.	1.71	1.00	-
MPFHpA	499000.	1.74	1.00	-
MPFOA	448000.	1.92	1.00	-
MPFOS	186000.	2.00	1.00	-
MPFNA	296000.	2.06	1.00	-
13C6-PFHxA IS	4060000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2520000	1.16	60.0	55.2	92.0
PFHxS 1	2760000	1.71	60.0	57.0	95.1
PFHpA 1	3790000	1.74	60.0	56.5	94.2
PFOA 1	4130000	1.92	60.0	56.1	93.5
PFOS 1	1700000	1.99	60.0	50.5	84.1
PFNA 1	2810000	2.06	60.0	57.2	95.3
18O2-PFHxS	147000	1.71	100.	98.3	98.3
13C4-PFHpA	499000	1.74	100.	99.2	99.2
13C4-PFOA	448000	1.92	100.	102.	102.0
13C4-PFOS	186000	2.00	100.	106.	106.0
13C5-PFNA	296000	2.06	100.	99.1	99.1
13C6-PFHxA	4060000	1.48	100.	102.	102.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 55.2 µg/L</p> <p>Area Ratio: 17.1</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 57.0 µg/L</p> <p>Area Ratio: 18.8</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 56.5 µg/L</p> <p>Area Ratio: 7.60</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 56.1 µg/L</p> <p>Area Ratio: 9.23</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 50.5 µg/L</p> <p>Area Ratio: 9.16</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 57.2 µg/L</p> <p>Area Ratio: 9.50</p> <p>Sample Type: (Standard)</p>	

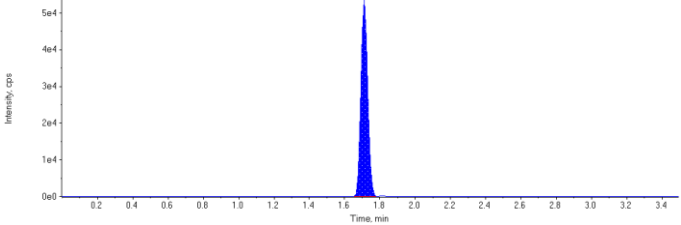
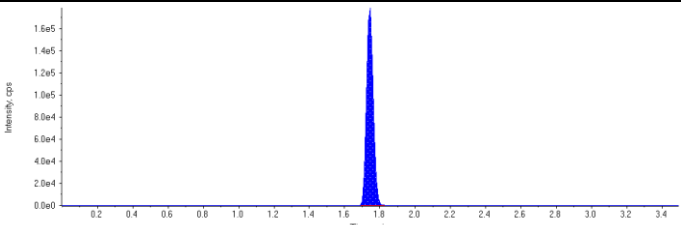
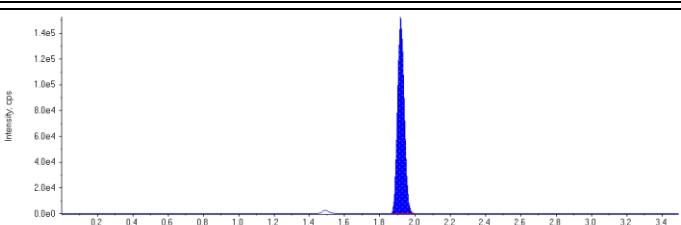
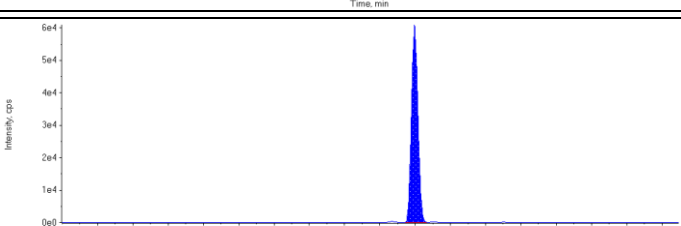
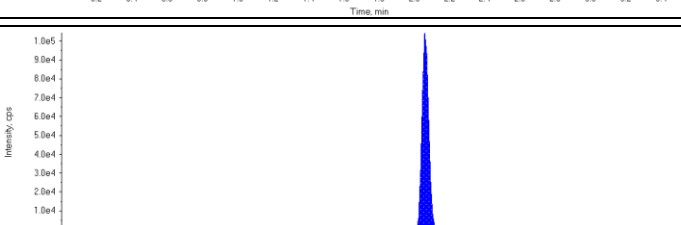
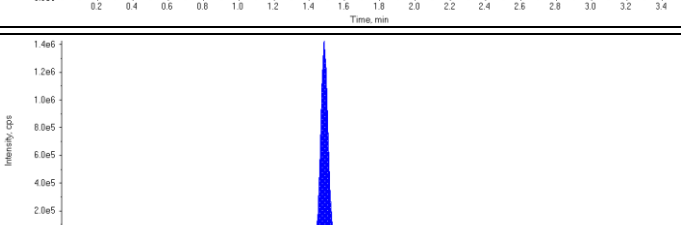
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 98.3 µg/L</p> <p>Area Ratio: 0.0362</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 99.2 µg/L</p> <p>Area Ratio: 0.123</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 102. µg/L</p> <p>Area Ratio: 0.110</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 106. µg/L</p> <p>Area Ratio: 0.0458</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 99.1 µg/L</p> <p>Area Ratio: 0.0729</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 102. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

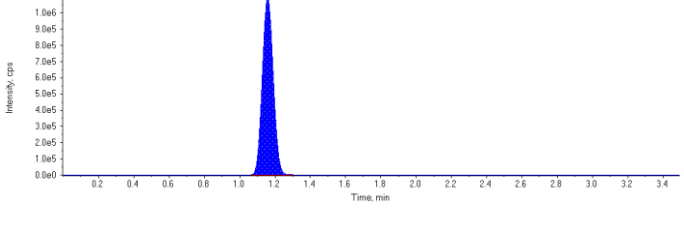
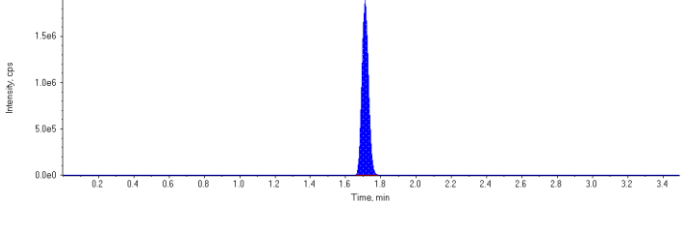
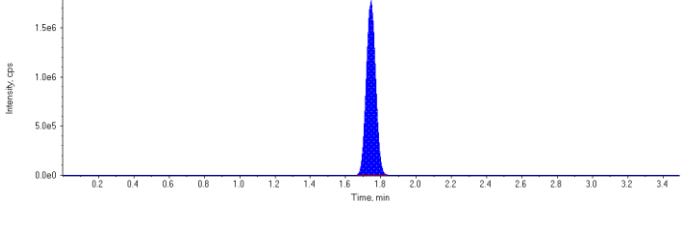
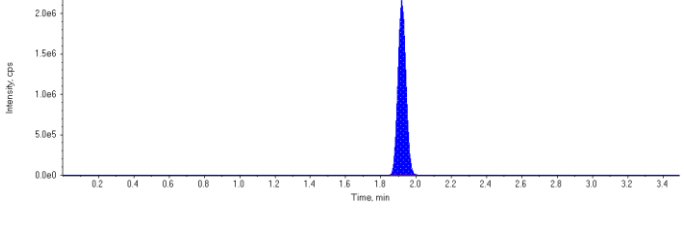
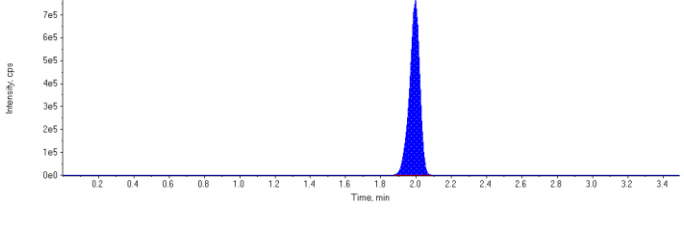
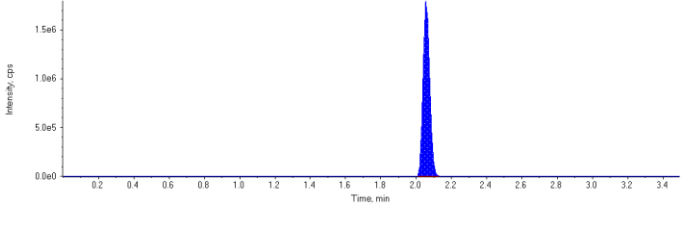
<b>Sample Name</b>	<b>STD 6</b>	<b>Injection Vial</b>	8
<b>Sample ID</b>	STD 6	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Standard	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:10:36 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

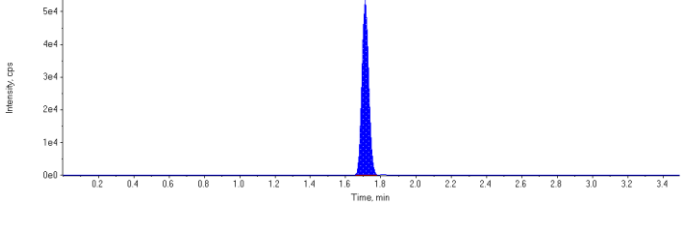
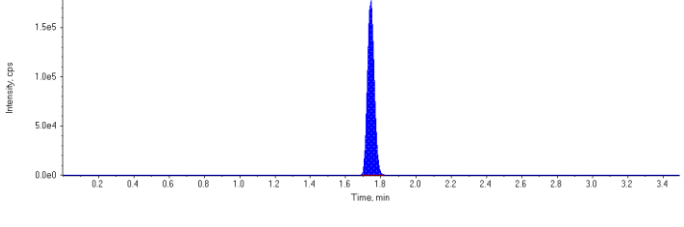
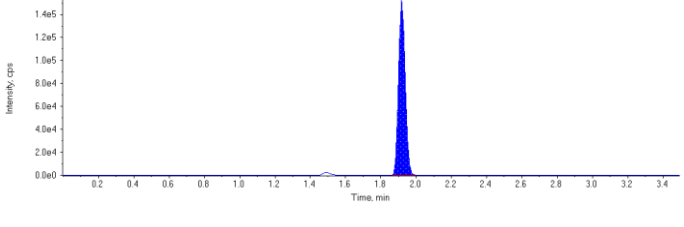
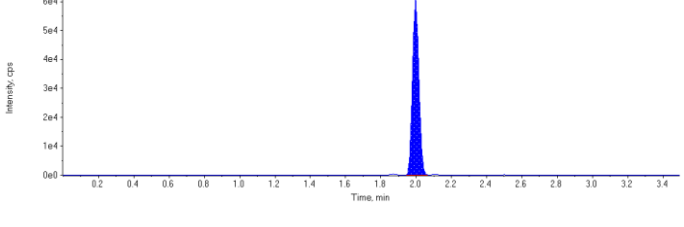
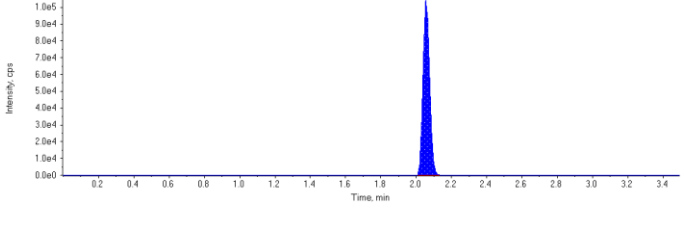
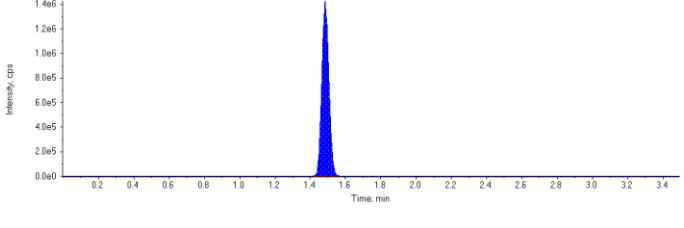
Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	142000.	1.71	1.00	-
MPFHpA	477000.	1.74	1.00	-
MPFOA	398000.	1.92	1.00	-
MPFOS	154000.	2.00	1.00	-
MPFNA	275000.	2.06	1.00	-
13C6-PFHxA IS	3980000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	4720000	1.16	100.	107.	107.0
PFHxS 1	4830000	1.71	100.	103.	103.0
PFHpA 1	6480000	1.74	100.	101.	101.0
PFOA 1	6810000	1.92	100.	104.	104.0
PFOS 1	3060000	2.00	100.	109.	109.0
PFNA 1	4710000	2.06	100.	103.	103.0
18O2-PFHxS	142000	1.71	100.	96.8	96.8
13C4-PFHpA	477000	1.74	100.	96.9	96.9
13C4-PFOA	398000	1.92	100.	92.7	92.7
13C4-PFOS	154000	2.00	100.	89.4	89.4
13C5-PFNA	275000	2.06	100.	94.0	94.0
13C6-PFHxA	3980000	1.48	100.	99.7	99.7



<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Standard)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Standard)</p>	<p style="text-align: center;">This image is not available</p>

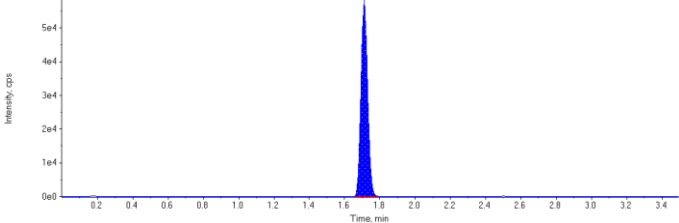
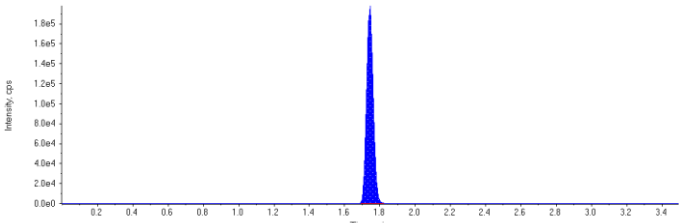
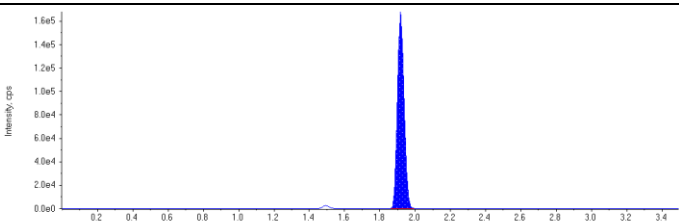
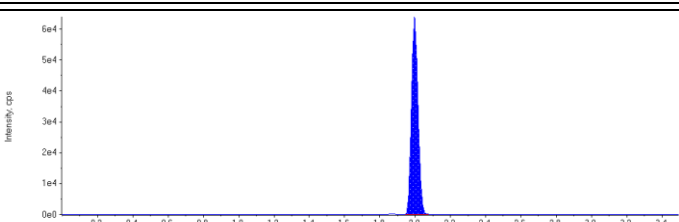
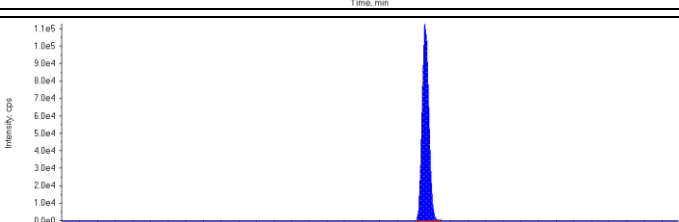
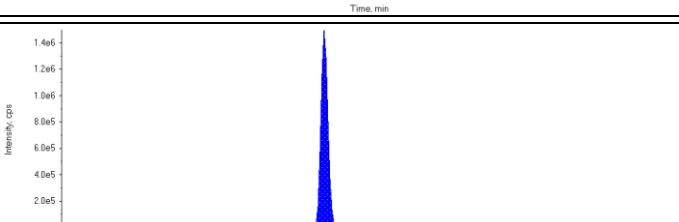
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 107. µg/L</p> <p>Area Ratio: 33.3</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 34.1</p> <p>Sample Type: (Standard)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 101. µg/L</p> <p>Area Ratio: 13.6</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 17.1</p> <p>Sample Type: (Standard)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 109. µg/L</p> <p>Area Ratio: 19.9</p> <p>Sample Type: (Standard)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 17.1</p> <p>Sample Type: (Standard)</p>	

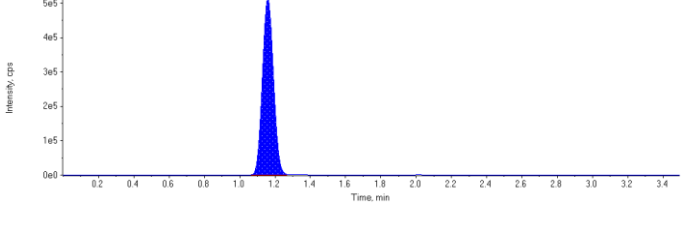
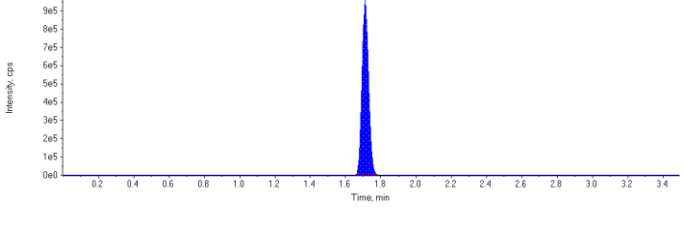
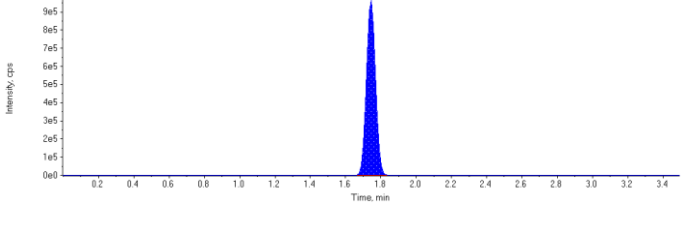
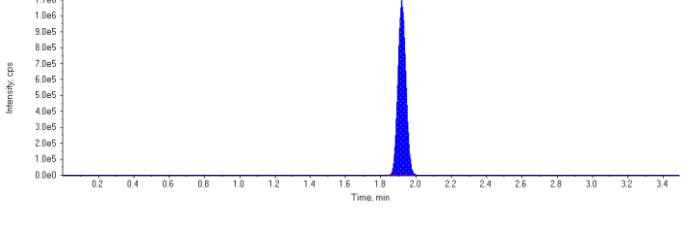
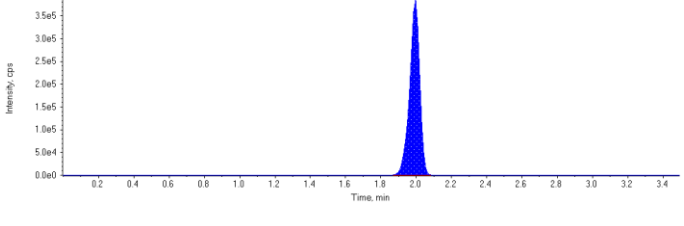
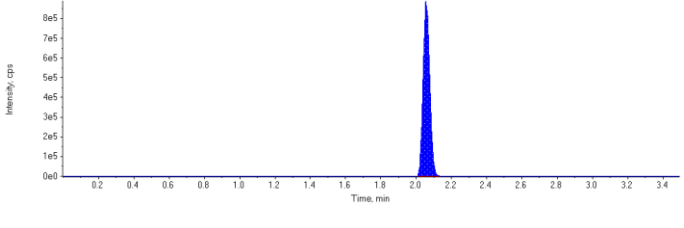
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 96.8 µg/L</p> <p>Area Ratio: 0.0356</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 96.9 µg/L</p> <p>Area Ratio: 0.120</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 92.7 µg/L</p> <p>Area Ratio: 0.100</p> <p>Sample Type: (Standard)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 89.4 µg/L</p> <p>Area Ratio: 0.0386</p> <p>Sample Type: (Standard)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 94.0 µg/L</p> <p>Area Ratio: 0.0692</p> <p>Sample Type: (Standard)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 99.7 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Standard)</p>	

<b>Sample Name</b>	<b>ICV</b>	<b>Injection Vial</b>	9
<b>Sample ID</b>	ICV	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:15:41 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	152000.	1.71	1.00	-
MPFHpA	517000.	1.74	1.00	-
MPFOA	437000.	1.92	1.00	-
MPFOS	163000.	2.00	1.00	-
MPFNA	297000.	2.06	1.00	-
13C6-PFHxA IS	4130000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	2240000	1.16	50.0	47.6	95.3
PFHxS 1	2450000	1.71	50.0	49.2	98.4
PFHpA 1	3460000	1.74	50.0	49.7	99.5
PFOA 1	3460000	1.92	50.0	48.2	96.3
PFOS 1	1530000	2.00	50.0	51.5	103.0
PFNA 1	2340000	2.06	50.0	47.6	95.1
18O2-PFHxS	152000	1.71	100.	99.7	99.7
13C4-PFHpA	517000	1.74	100.	101.	101.0
13C4-PFOA	437000	1.92	100.	98.0	98.0
13C4-PFOS	163000	2.00	100.	91.5	91.5
13C5-PFNA	297000	2.06	100.	97.6	97.6
13C6-PFHxA	4130000	1.48	100.	104.	104.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 47.6 µg/L</p> <p>Area Ratio: 14.7</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 49.2 µg/L</p> <p>Area Ratio: 16.2</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 49.7 µg/L</p> <p>Area Ratio: 6.69</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 48.2 µg/L</p> <p>Area Ratio: 7.92</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 51.5 µg/L</p> <p>Area Ratio: 9.34</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 47.6 µg/L</p> <p>Area Ratio: 7.89</p> <p>Sample Type: (Quality Control)</p>	

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 99.7 µg/L</p> <p>Area Ratio: 0.0367</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 101. µg/L</p> <p>Area Ratio: 0.125</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 98.0 µg/L</p> <p>Area Ratio: 0.106</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 91.5 µg/L</p> <p>Area Ratio: 0.0395</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 97.6 µg/L</p> <p>Area Ratio: 0.0719</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

Sample Name	Analyte	PFBS		PFHxS		PFHpA		PFOA		PFOS		PFNA	
	Mass labeled IS	MPFHxS		MPFHxS		MPFHpA		MPFOA		MPFOS		MPFNA	
	Average IS Area in ICAL	147000		147000		493500		430500		172500		293500	
		IS Peak Area	IS %R (ICAL)	IS Peak Area	IS %R (ICAL)	IS Peak Area	IS %R (ICAL)	IS Peak Area	IS %R (ICAL)	IS Peak Area	IS %R (ICAL)	IS Peak Area	IS %R (ICAL)
4386408~BLANK		130000	88	130000	88	375000	76	327000	76	146000	85	228000	78
STD 1		146000	99	146000	99	517000	105	413000	96	179000	104	291000	99
STD 2		147000	100	147000	100	503000	102	442000	103	177000	103	296000	101
STD 3		156000	106	156000	106	498000	101	435000	101	167000	97	307000	105
STD 4		144000	98	144000	98	467000	95	447000	104	172000	100	296000	101
STD 5		147000	100	147000	100	499000	101	448000	104	186000	108	296000	101
STD 6		142000	97	142000	97	477000	97	398000	92	154000	89	275000	94
ICV		152000	103	152000	103	517000	105	437000	102	163000	94	297000	101
CCV		156000	106	156000	106	500000	101	423000	98	177000	103	314000	107
4386408~MTRX SPK		112000	76	112000	76	335000	68	321000	75	127000	74	227000	77
4386408~MTRX SPK:D1		110000	75	110000	75	331000	67	363000	84	141000	82	263000	90
4386408~SPIKE		123000	84	123000	84	336000	68	305000	71	135000	78	218000	74
4386408~BVX814-01 (10x)		113000	77	113000	77	340000	69	322000	75	139000	81	236000	80
4386408~BVX817-01 (10x)		134000	91	134000	91	372000	75	336000	78	145000	84	240000	82
4386408~BVX818-01 (20x)		115000	78	115000	78	355000	72	335000	78	147000	85	235000	80
4386408~BVX819-01 (10x)		109000	74	109000	74	212000	43	228000	53	136000	79	183000	62
4386408~BVX769-01		103000	70	103000	70	315000	64	302000	70	137000	79	225000	77
4386408~BVX770-01		116000	79	116000	79	281000	57	273000	63	138000	80	222000	76
4386408~BVX771-01		101000	69	101000	69	263000	53	257000	60	126000	73	208000	71
4386408~BVX772-01		95400	65	95400	65	287000	58	261000	61	120000	70	205000	70
4386408~BVX773-01		97200	66	97200	66	312000	63	297000	69	124000	72	206000	70
4386408~BVX774-01		101000	69	101000	69	301000	61	286000	66	117000	68	207000	71
CCV		118000	80	118000	80	376000	76	350000	81	153000	89	275000	94
4386408~BVX775-01		83400	57	83400	57	241000	49	243000	56	93300	54	170000	58
4386408~BVX776-01		116000	79	116000	79	334000	68	322000	75	129000	75	248000	84
4386408~BVX779-01		97700	66	97700	66	321000	65	300000	70	129000	75	224000	76
4386408~BVX780-01		98900	67	98900	67	257000	52	252000	59	113000	66	190000	65
4386408~BVX781-01		89500	61	89500	61	261000	53	244000	57	101000	59	179000	61
4386408~BVX826-01		104000	71	104000	71	272000	55	256000	59	105000	61	186000	63
4386408~BVX827-01		84400	57	84400	57	229000	46	217000	50	98200	57	170000	58
4386408~BVX828-01		92800	63	92800	63	289000	59	269000	62	98300	57	188000	64
4386408~BVX829-01		79900	54	79900	54	286000	58	252000	59	110000	64	180000	61
4386408~BVX830-01		82200	56	82200	56	275000	56	249000	58	102000	59	189000	64
CCV		122000	83	122000	83	400000	81	381000	89	150000	87	266000	91
4386408~MTRX SPK		95600	65	95600	65	303000	61	307000	71	126000	73	222000	76
4386408~BVX830-01		96600	66	96600	66	277000	56	258000	60	103000	60	188000	64
CCV		123000	84	123000	84	413000	84	372000	86	165000	96	267000	91

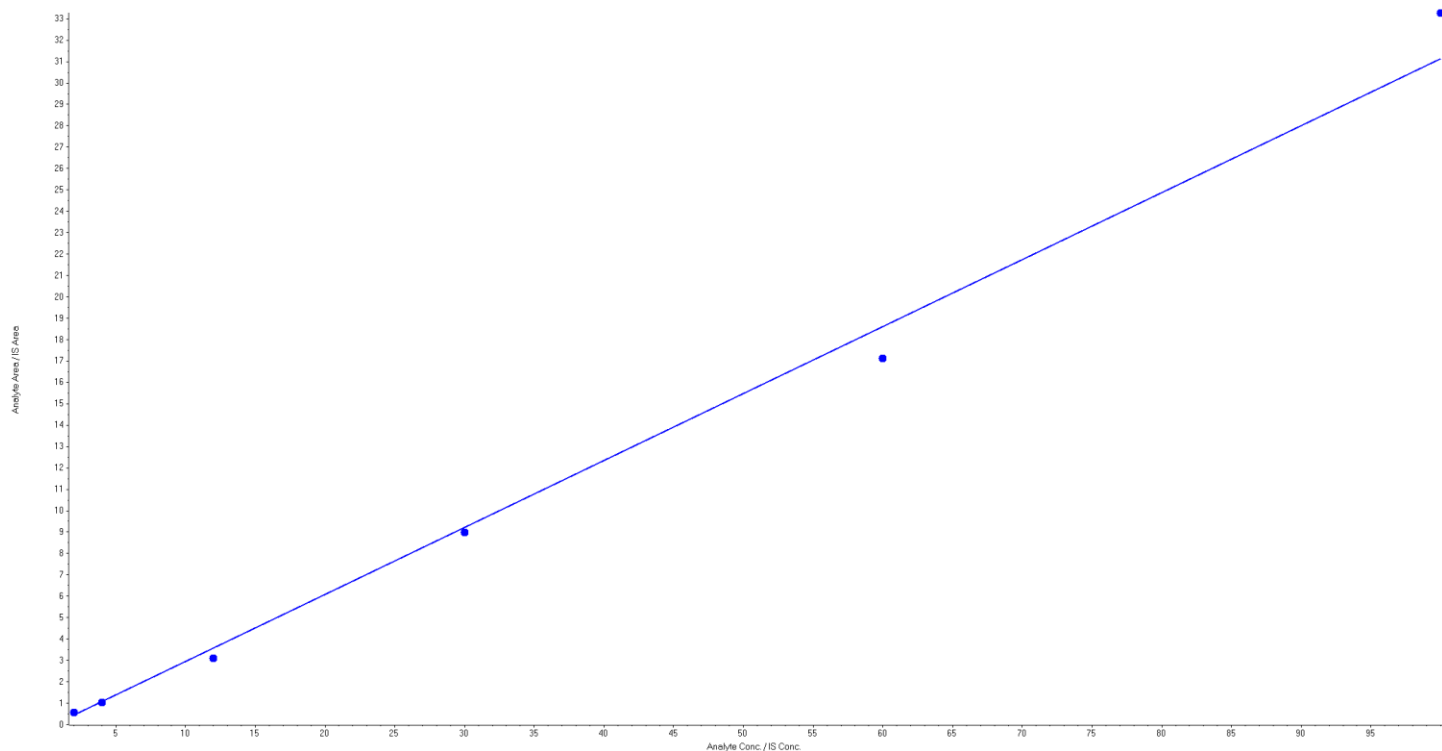


**Analyte Name:** PFBS 1  
**Internal Standard:** MPFHxS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.313 x + -0.179$  (r = 0.9966)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.40	120.2
4	1	3.86	96.4
12	1	10.44	87.0
30	1	29.26	97.5
60	1	55.22	92.0
100	1	106.82	106.8

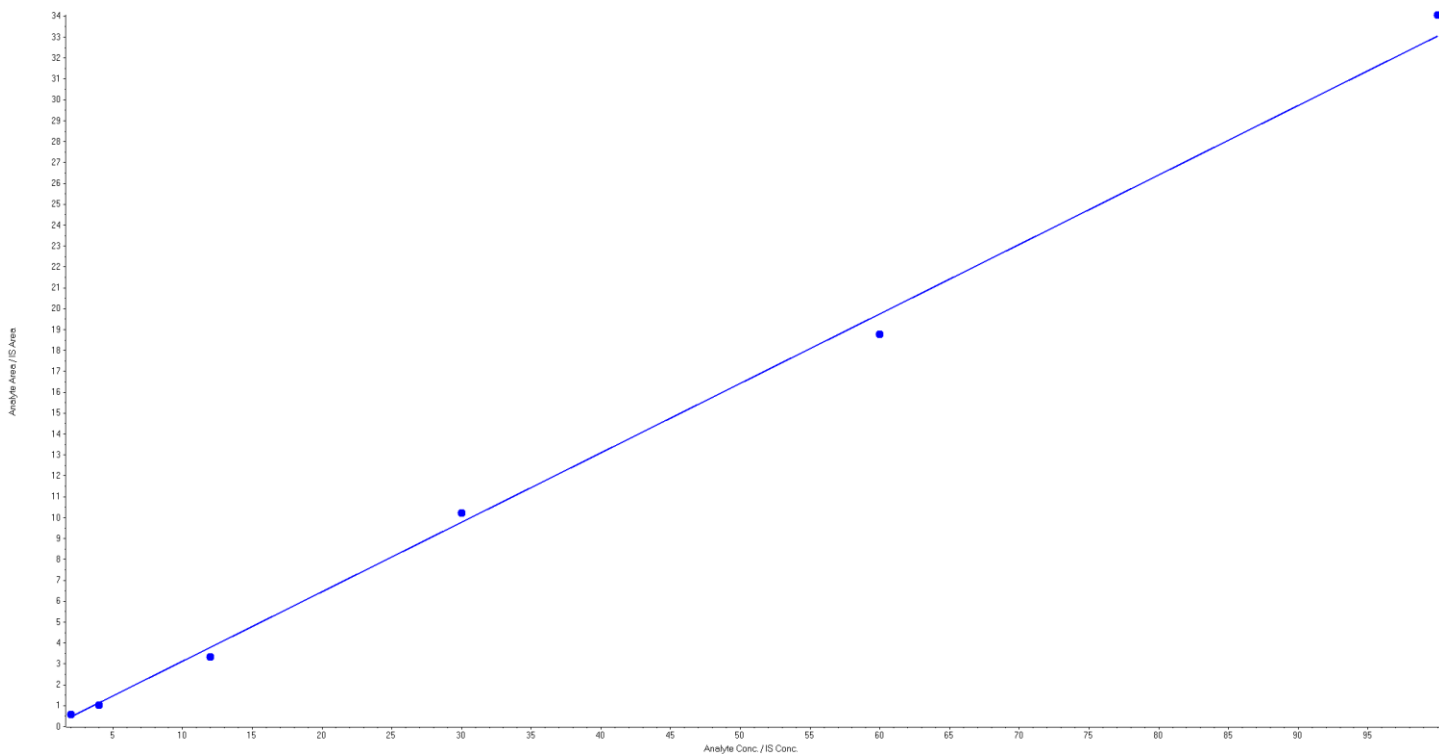


**Analyte Name:** PFHxS 1  
**Internal Standard:** MPFHxS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.333 x + -0.197$  (r = 0.9984)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.35	117.3
4	1	3.68	92.1
12	1	10.57	88.1
30	1	31.33	104.4
60	1	57.04	95.1
100	1	103.02	103.0

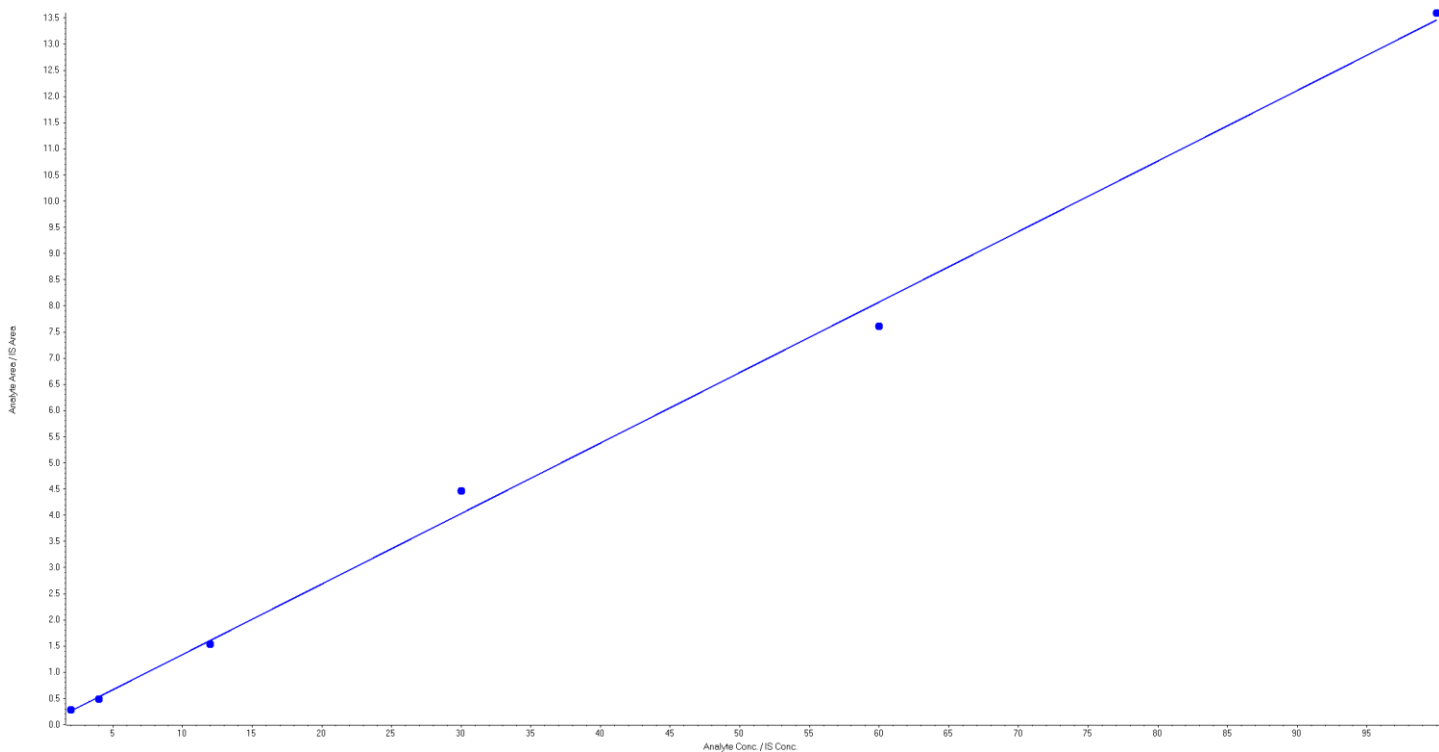


**Analyte Name:** PFHpA 1  
**Internal Standard:** MPFHpA

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.135 x + -0.0098$  (r = 0.9982)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.11	105.7
4	1	3.72	92.9
12	1	11.45	95.5
30	1	33.21	110.7
60	1	56.52	94.2
100	1	100.98	101.0

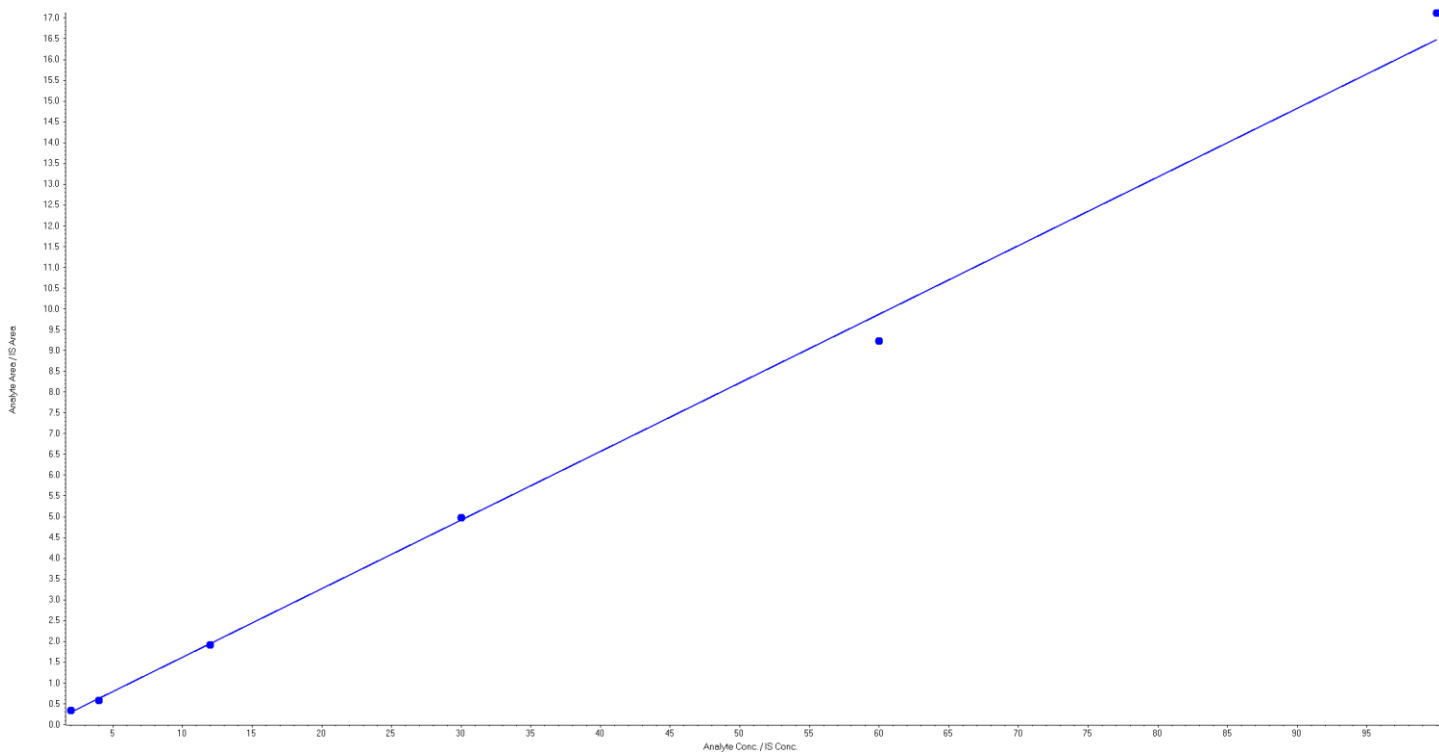


**Analyte Name:** PFOA 1  
**Internal Standard:** MPFOA

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.165 x + -0.0324$  (r = 0.9987)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.20	109.9
4	1	3.73	93.3
12	1	11.81	98.4
30	1	30.30	101.0
60	1	56.08	93.5
100	1	103.88	103.9

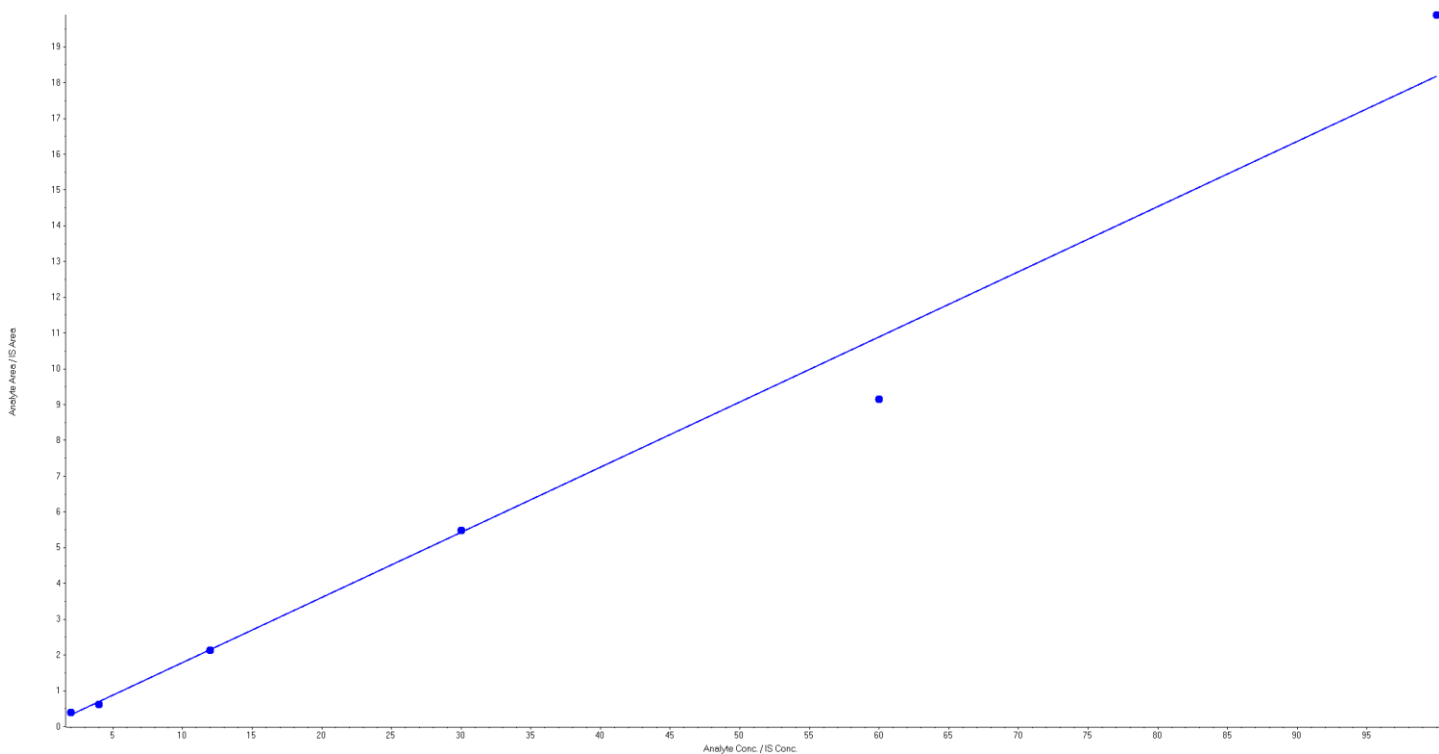


**Analyte Name:** PFOS 1  
**Internal Standard:** MPFOS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.182 x + -0.034$  (r = 0.9927)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.31	115.6
4	1	3.63	90.7
12	1	11.90	99.2
30	1	30.32	101.1
60	1	50.45	84.1
100	1	109.38	109.4

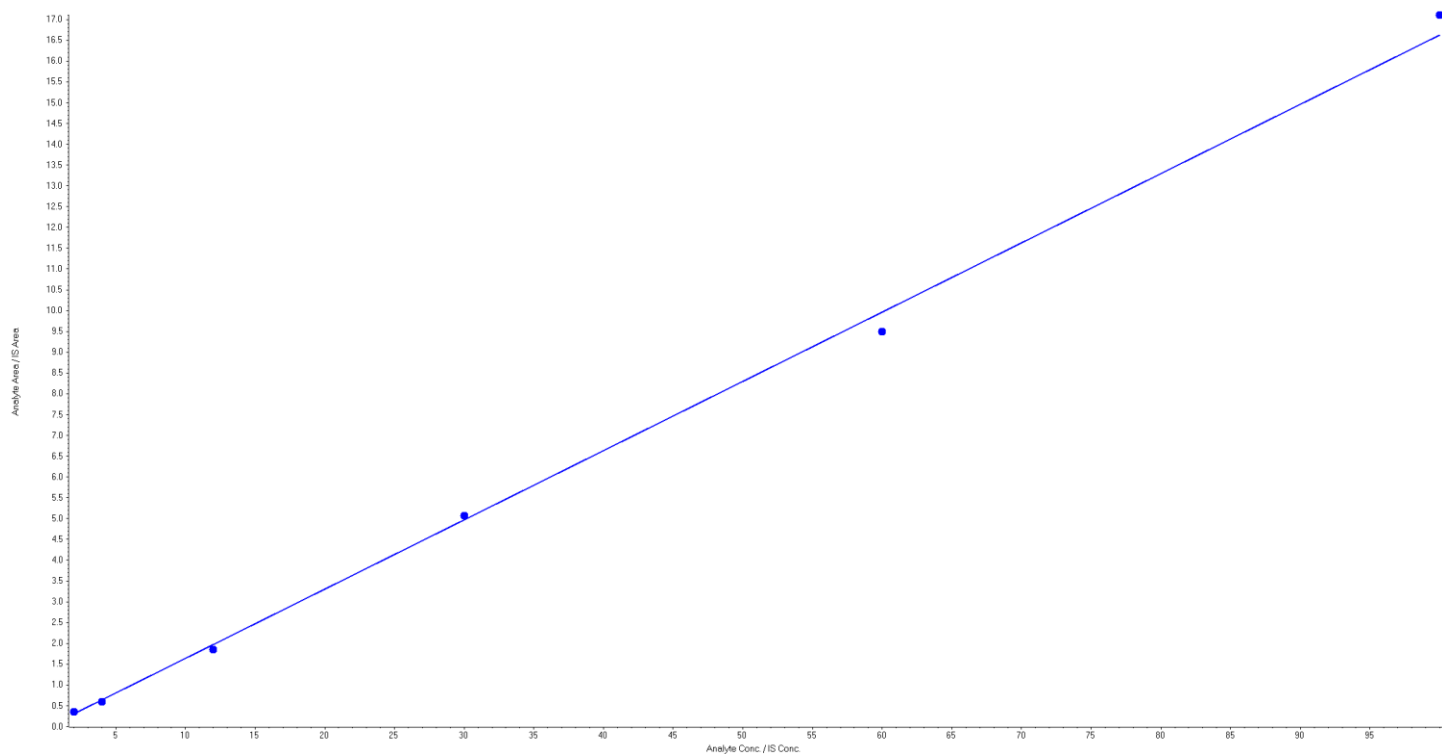


**Analyte Name:** PFNA 1  
**Internal Standard:** MPFNA

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.167 x + -0.0262$  (r = 0.9991)

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
2	1	2.24	111.9
4	1	3.74	93.6
12	1	11.30	94.2
30	1	30.63	102.1
60	1	57.19	95.3
100	1	102.89	102.9

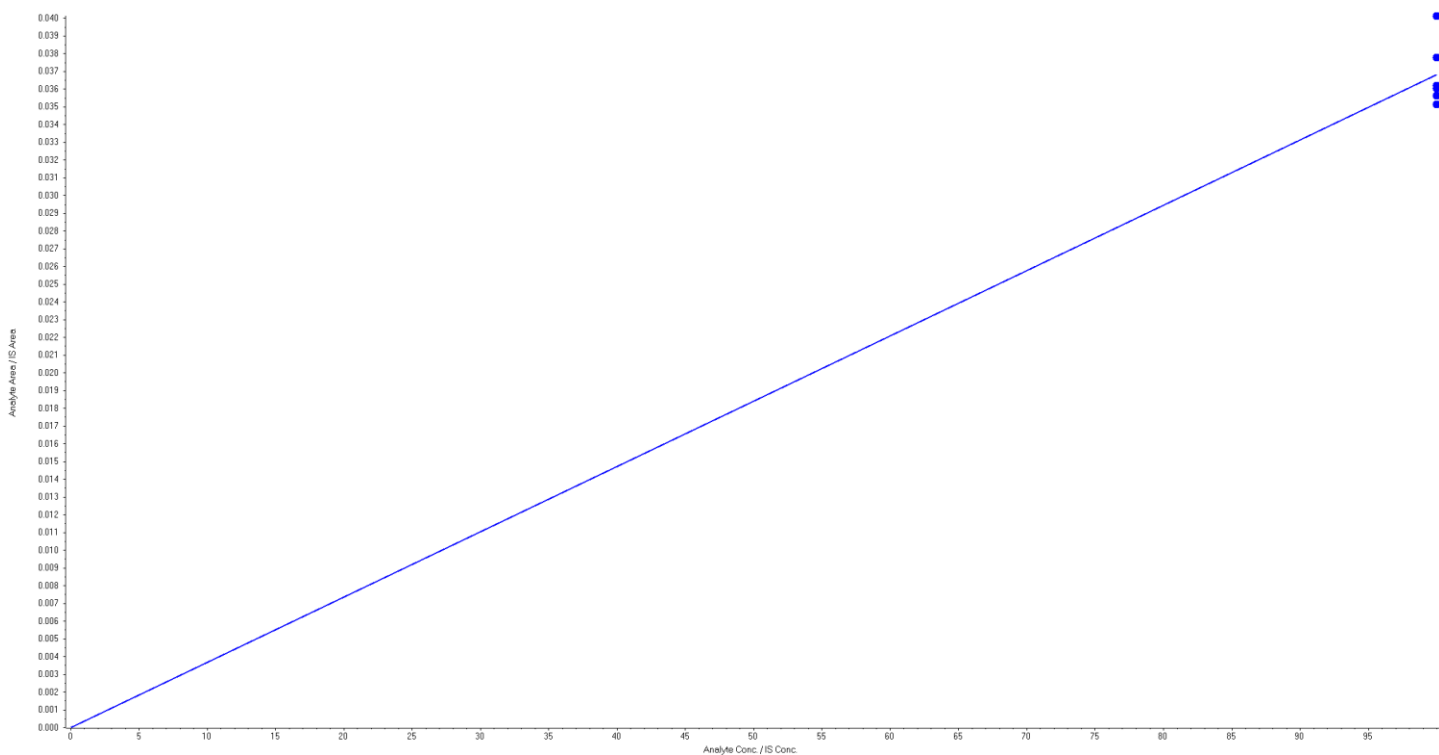


**Analyte Name:** 18O2-PFHxS  
**Internal Standard:** 13C6-PFHxA IS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.000368 x (r = 0.9990)$

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
100	6	100.00	100.0

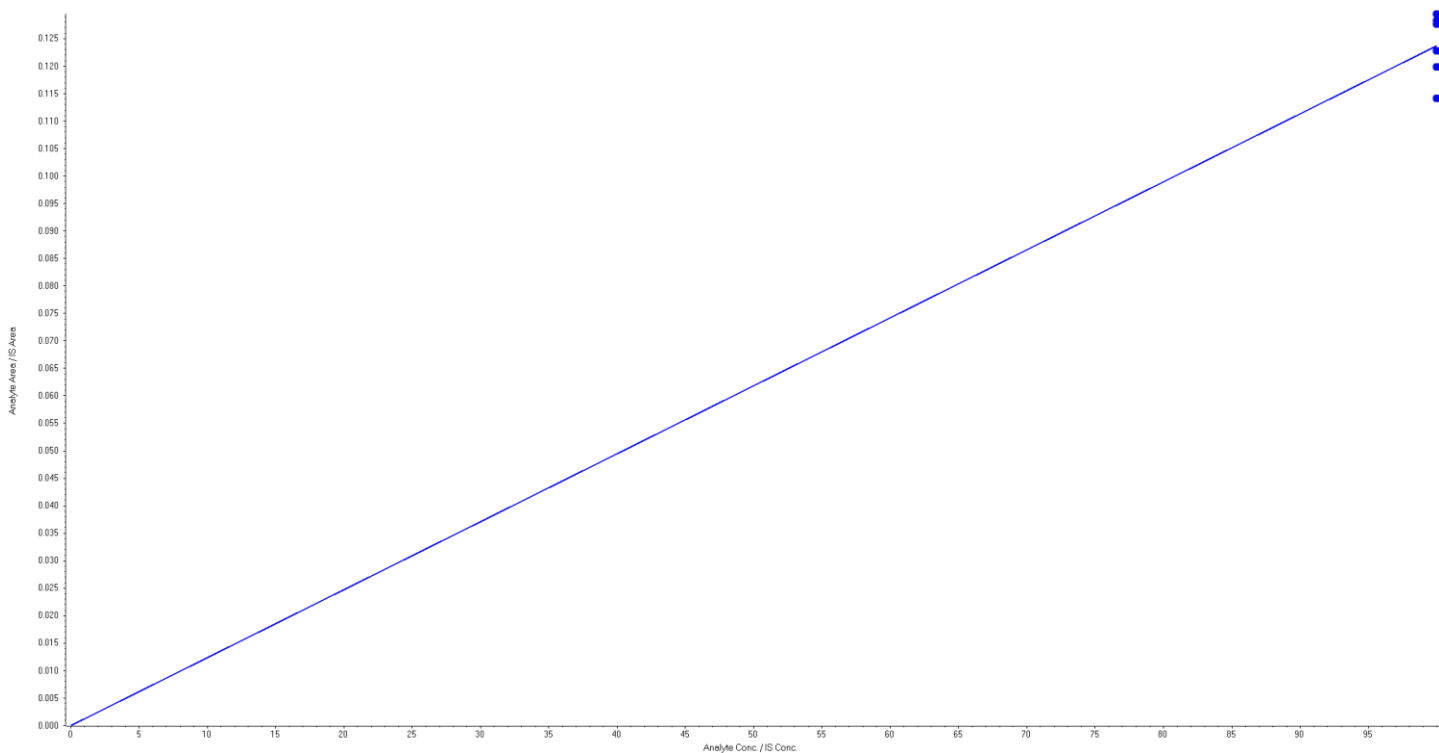


**Analyte Name:** 13C4-PFHpA  
**Internal Standard:** 13C6-PFHxA IS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.00124 x (r = 0.9990)$

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
100	6	100.00	100.0



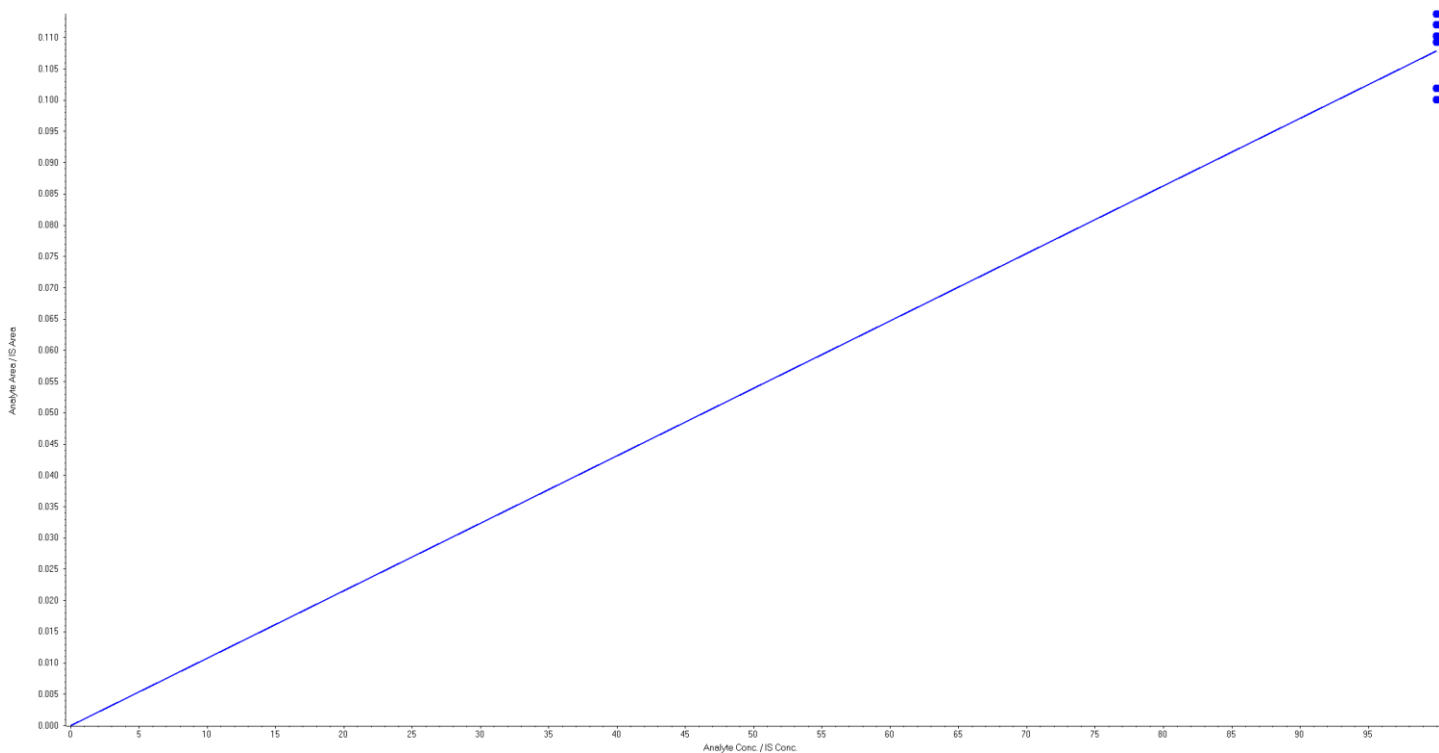


Analyte Name: 13C4-PFOA  
 Internal Standard: 13C6-PFHxA IS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

Regression Equation:  $y = 0.00108 x (r = 0.9989)$

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
100	6	100.00	100.0

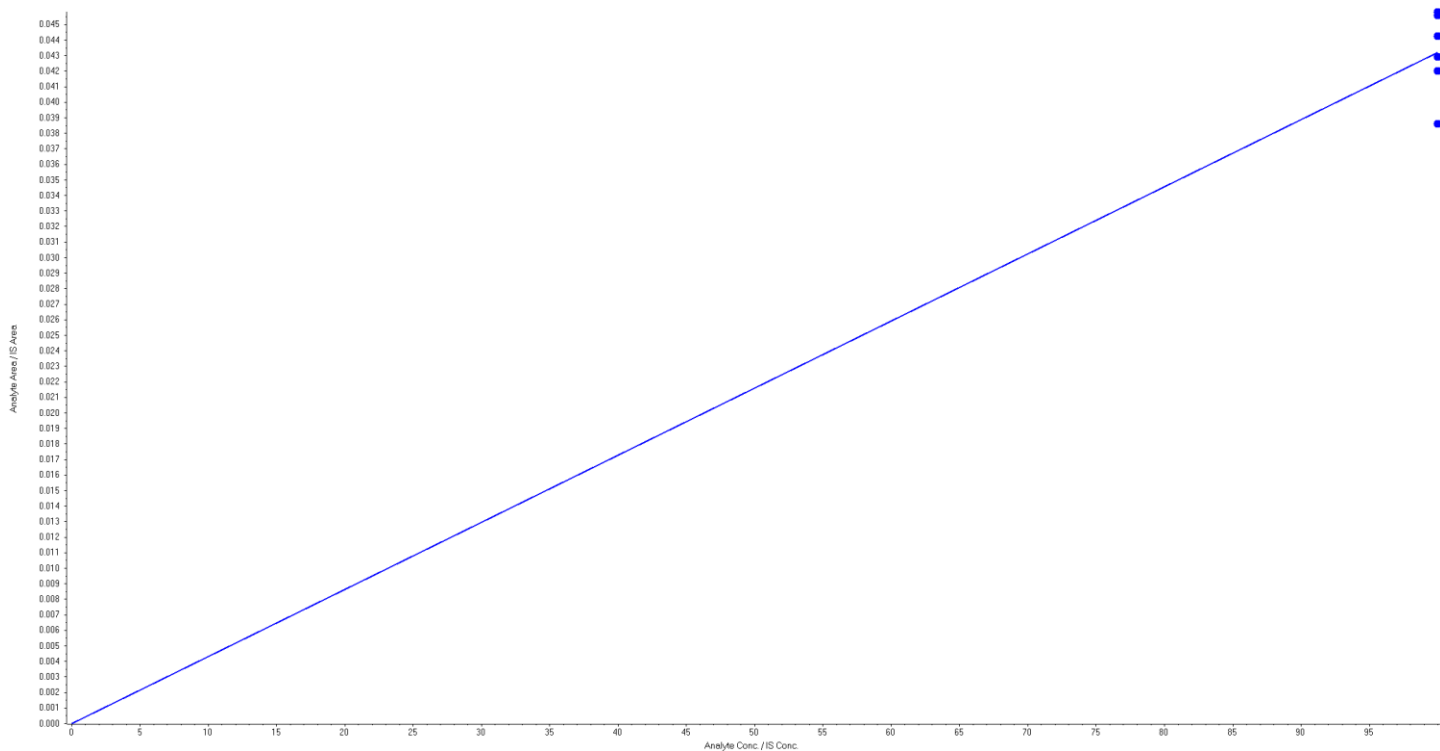


**Analyte Name:** 13C4-PFOS  
**Internal Standard:** 13C6-PFHxA IS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.000432 x (r = 0.9984)$

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
100	6	100.00	100.0

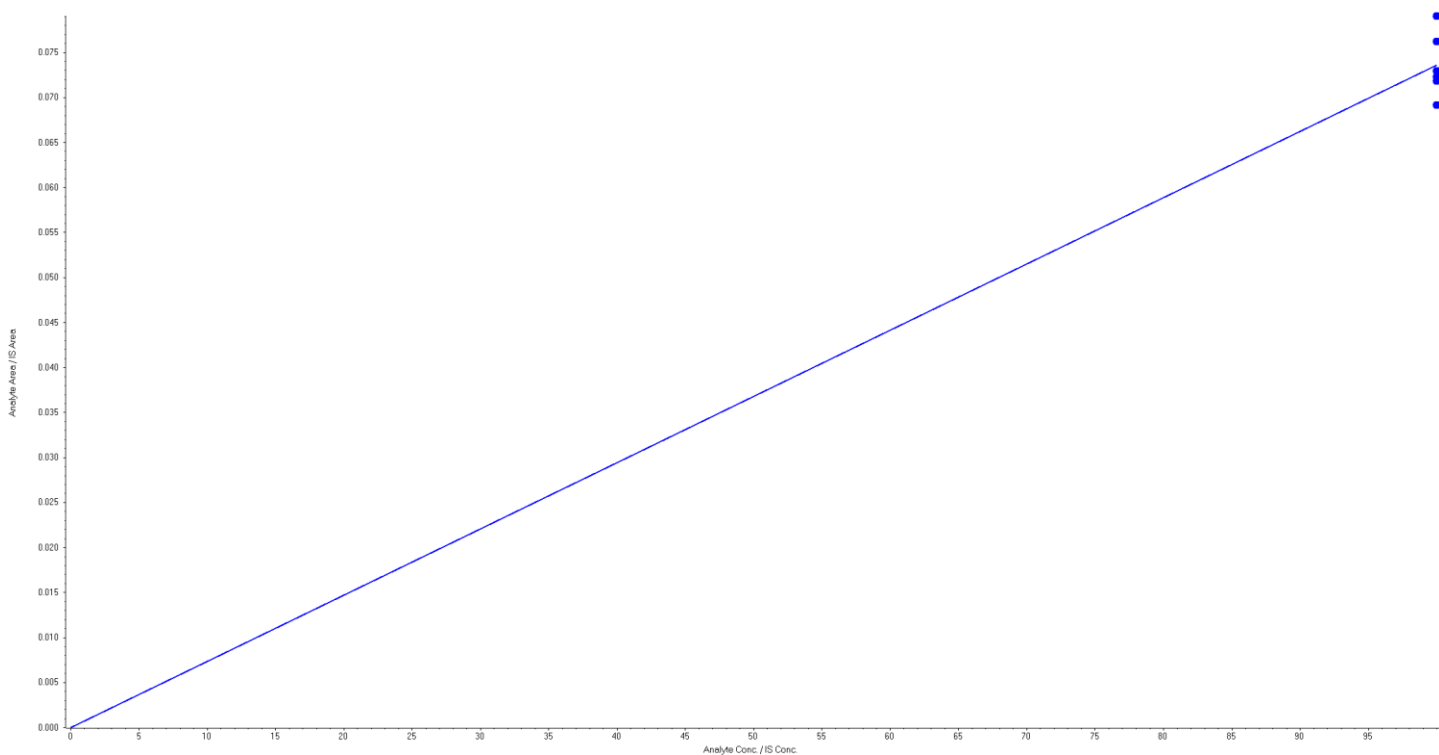


**Analyte Name:** 13C5-PFNA  
**Internal Standard:** 13C6-PFHxA IS

<b>Data File</b>	PFC_160219\WS#4386408.wiff	<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb
<b>Acquisition Date</b>	2016/02/19 7:39:53 AM	<b>Project</b>	Enviro\PFOS
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Instrument Name</b>	LCMS03

**Regression Equation:**  $y = 0.000736 x (r = 0.9991)$

Expected Concentration (µg/L)	Number of Values	Calculated Concentration (µg/L)	% Accuracy
100	6	100.00	100.0





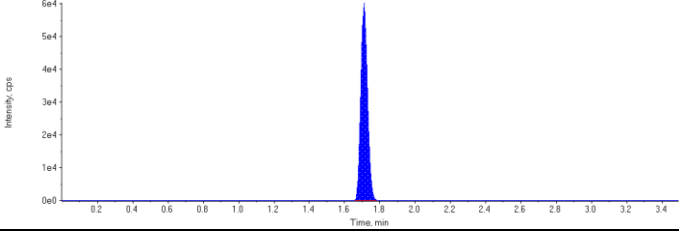
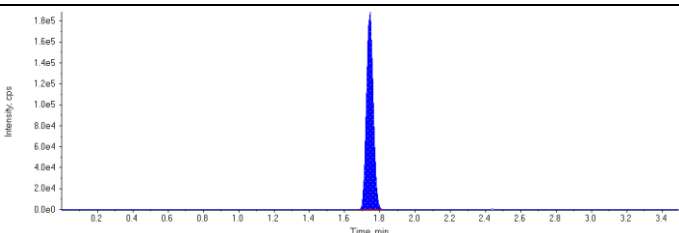
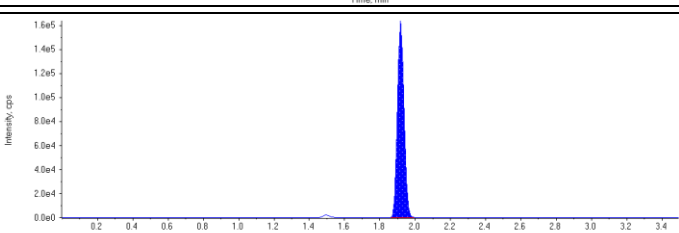
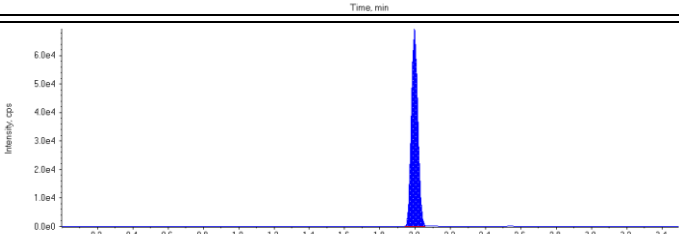
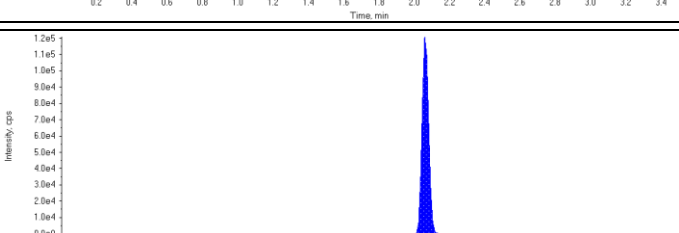
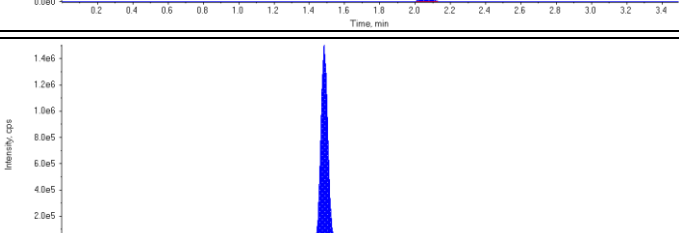
## 6. Continuing Calibration

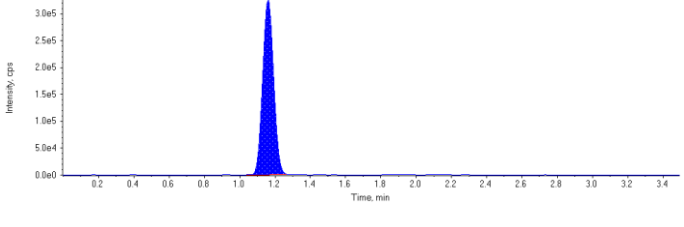
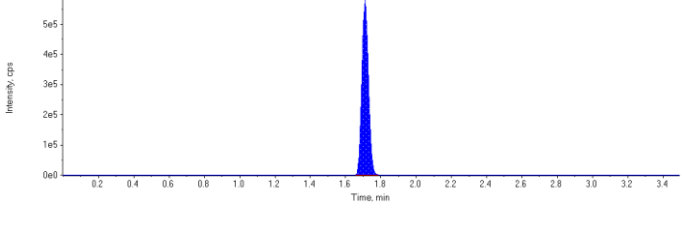
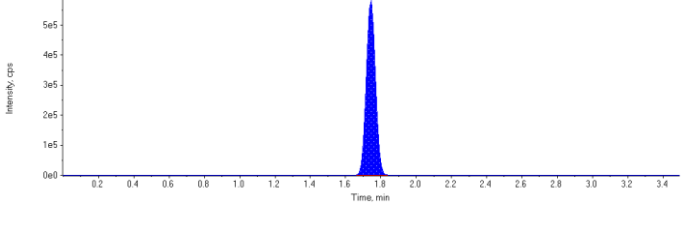
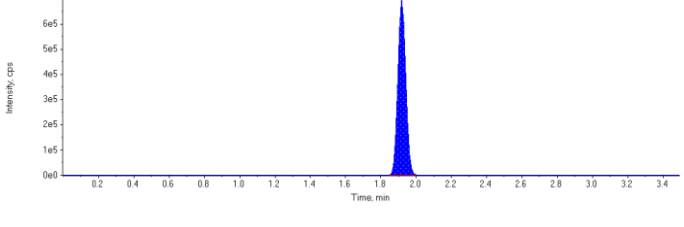
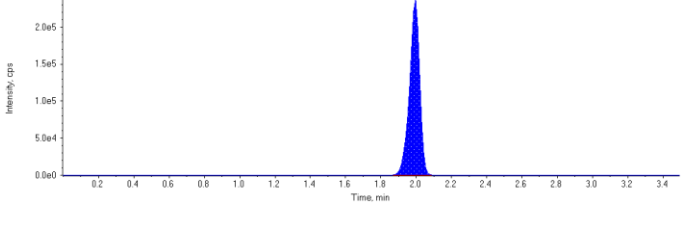
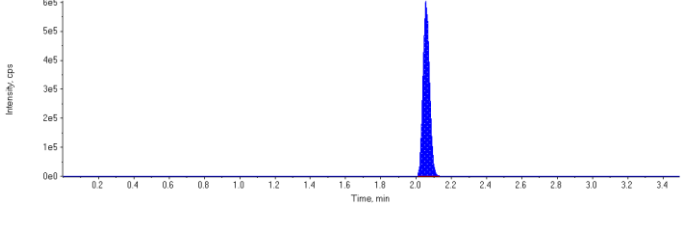
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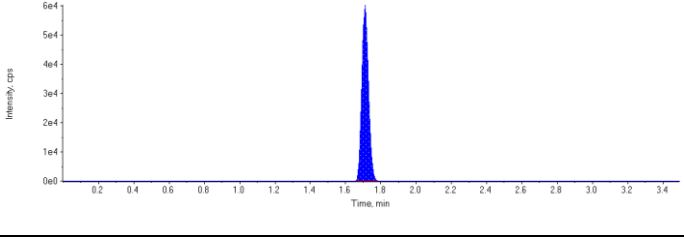
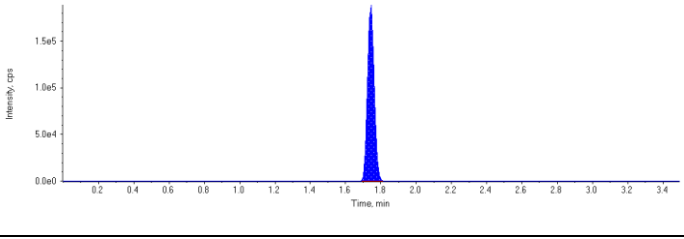
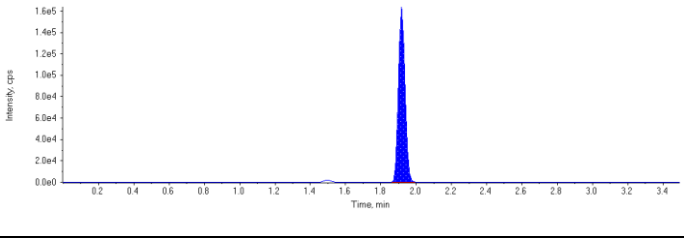
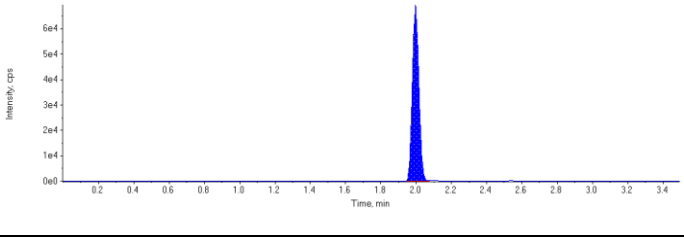
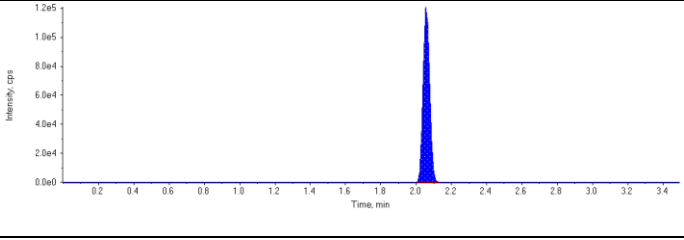
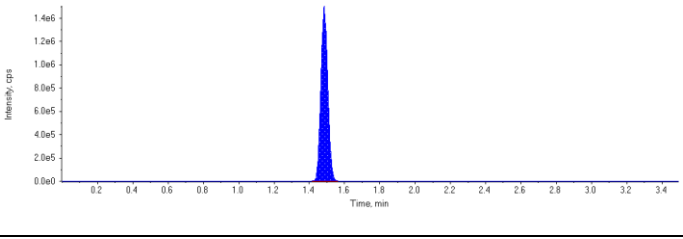
<b>Sample Name</b>	CCV	<b>Injection Vial</b>	6
<b>Sample ID</b>	CCV	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 8:20:47 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	156000.	1.71	1.00	-
MPFHpA	500000.	1.74	1.00	-
MPFOA	423000.	1.92	1.00	-
MPFOS	177000.	2.00	1.00	-
MPFNA	314000.	2.06	1.00	-
13C6-PFHxA IS	4100000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	1410000	1.16	30.0	29.3	97.8
PFHxS 1	1500000	1.71	30.0	29.3	97.8
PFHpA 1	2110000	1.74	30.0	31.4	105.0
PFOA 1	2180000	1.92	30.0	31.3	104.0
PFOS 1	950000	1.99	30.0	29.7	99.2
PFNA 1	1570000	2.06	30.0	30.1	100.0
18O2-PFHxS	156000	1.71	100.	104.	104.0
13C4-PFHpA	500000	1.74	100.	98.6	98.6
13C4-PFOA	423000	1.92	100.	95.7	95.7
13C4-PFOS	177000	2.00	100.	99.6	99.6
13C5-PFNA	314000	2.06	100.	104.	104.0
13C6-PFHxA	4100000	1.48	100.	103.	103.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.00(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 29.3 µg/L</p> <p>Area Ratio: 9.01</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 29.3 µg/L</p> <p>Area Ratio: 9.56</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 31.4 µg/L</p> <p>Area Ratio: 4.22</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 31.3 µg/L</p> <p>Area Ratio: 5.14</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 29.7 µg/L</p> <p>Area Ratio: 5.38</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 30.1 µg/L</p> <p>Area Ratio: 4.98</p> <p>Sample Type: (Quality Control)</p>	

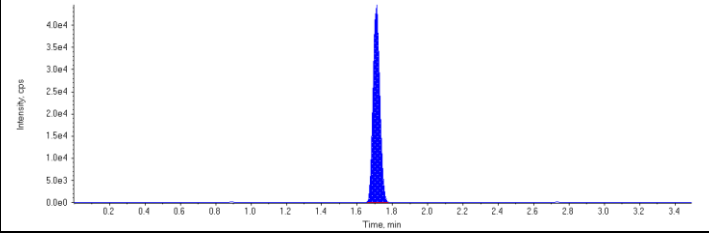
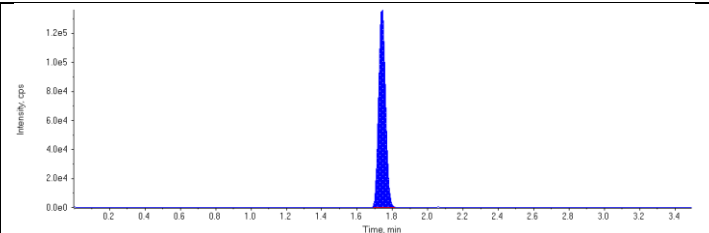
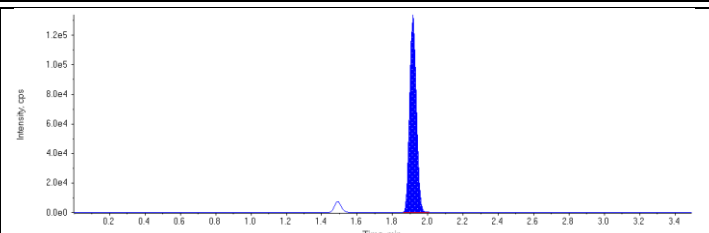
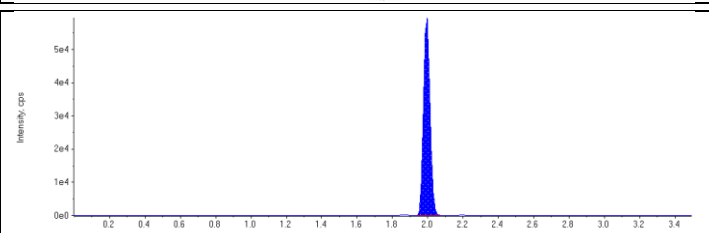
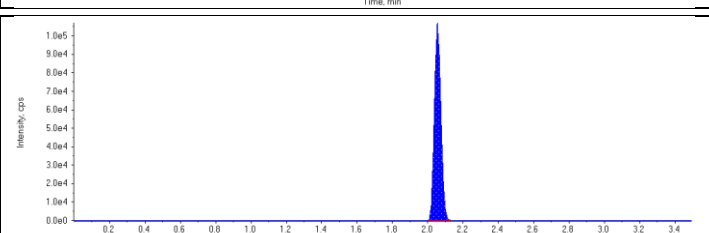
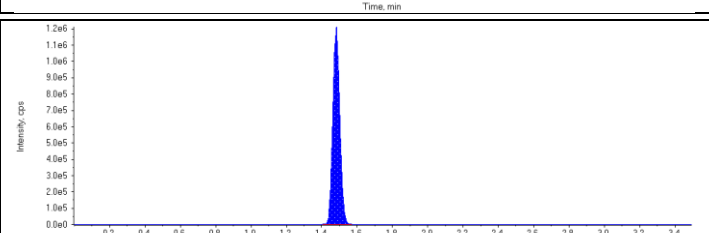
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.0382</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 98.6 µg/L</p> <p>Area Ratio: 0.122</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 95.7 µg/L</p> <p>Area Ratio: 0.103</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 2.00 (1.97) min</p> <p>Calculated Conc: 99.6 µg/L</p> <p>Area Ratio: 0.0431</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 104. µg/L</p> <p>Area Ratio: 0.0767</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

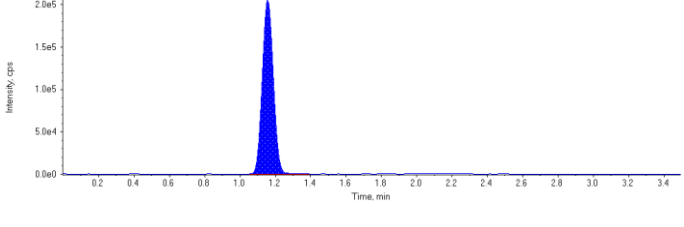
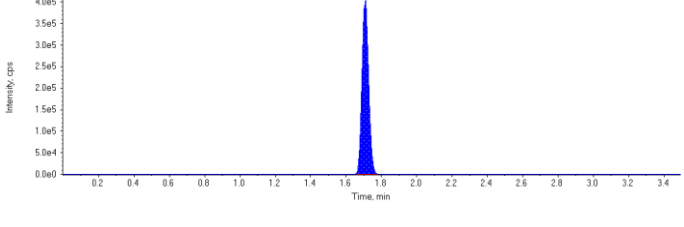
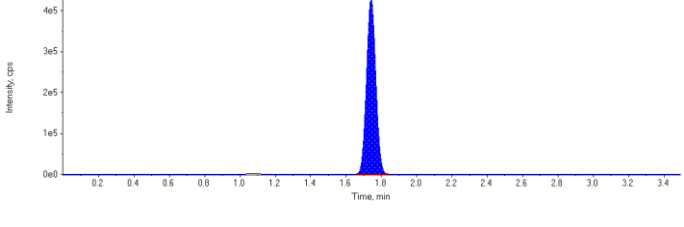
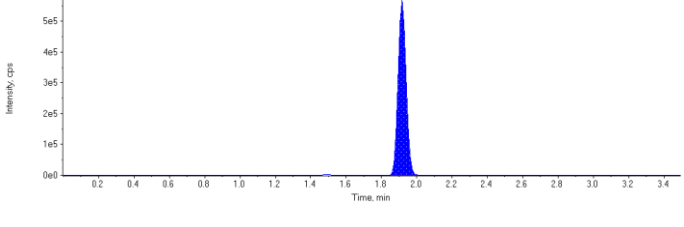
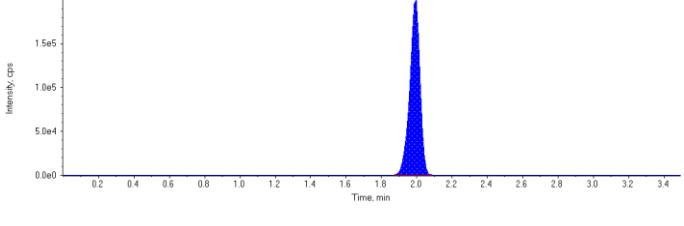
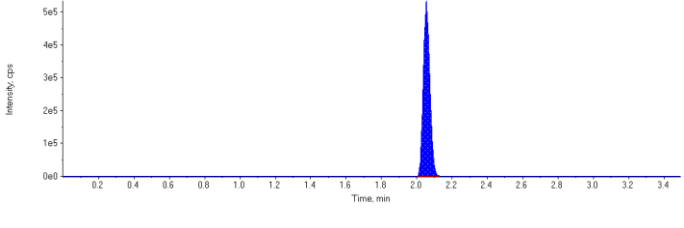


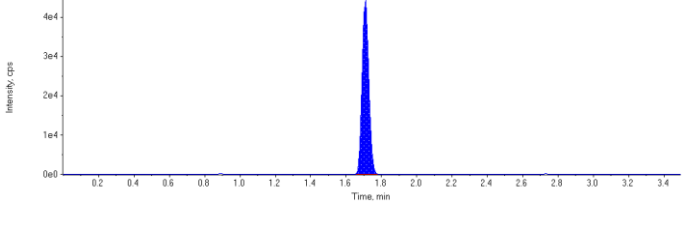
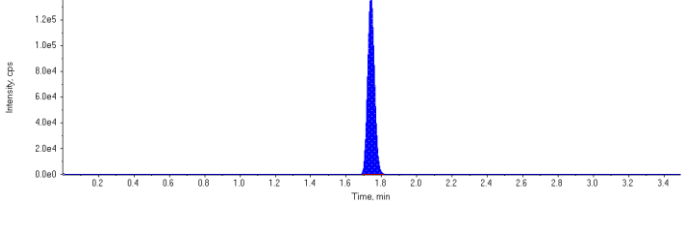
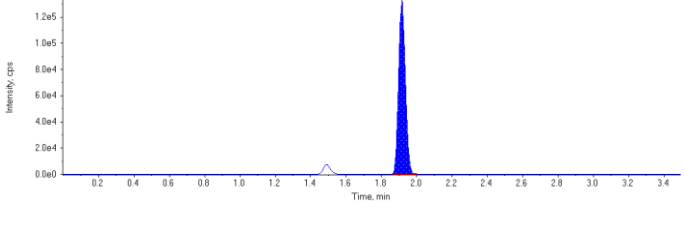
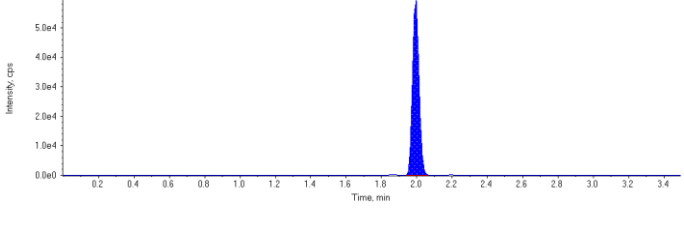
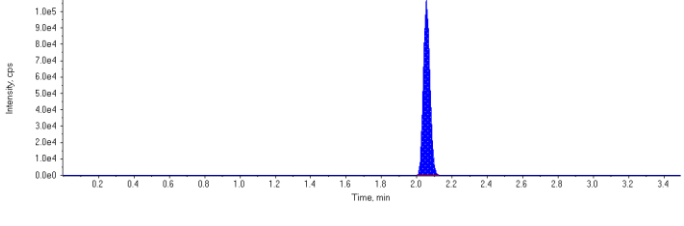
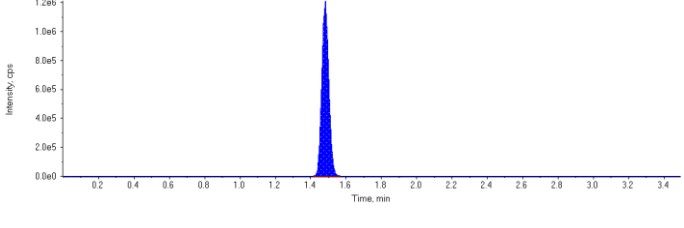
<b>Sample Name</b>	CCV	<b>Injection Vial</b>	6
<b>Sample ID</b>	CCV	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 9:42:22 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	118000.	1.71	1.00	-
MPFHpA	376000.	1.74	1.00	-
MPFOA	350000.	1.92	1.00	-
MPFOS	153000.	1.99	1.00	-
MPFNA	275000.	2.06	1.00	-
13C6-PFHxA IS	3350000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	894000	1.16	30.0	24.9	82.8
PFHxS 1	1060000	1.71	30.0	27.7	92.4
PFHpA 1	1570000	1.74	30.0	31.1	104.0
PFOA 1	1780000	1.92	30.0	31.0	103.0
PFOS 1	804000	1.99	30.0	29.0	96.6
PFNA 1	1370000	2.05	30.0	30.1	100.0
18O2-PFHxS	118000	1.71	100.	95.3	95.3
13C4-PFHpA	376000	1.74	100.	90.7	90.7
13C4-PFOA	350000	1.92	100.	96.7	96.7
13C4-PFOS	153000	1.99	100.	106.	106.0
13C5-PFNA	275000	2.06	100.	111.	111.0
13C6-PFHxA	3350000	1.48	100.	84.0	84.0

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.92(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.06(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

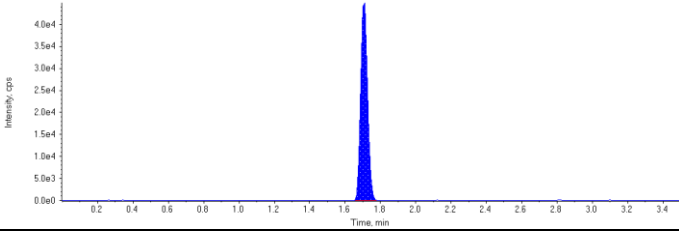
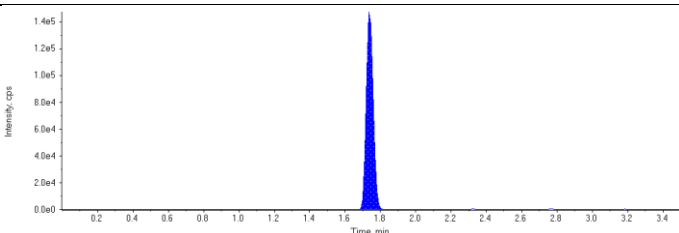
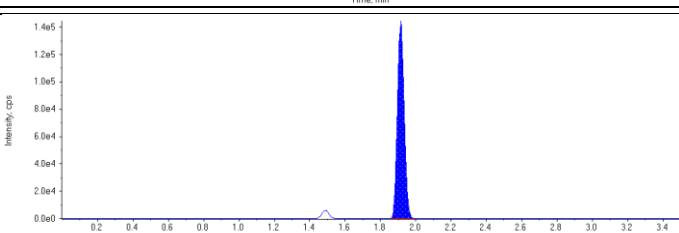
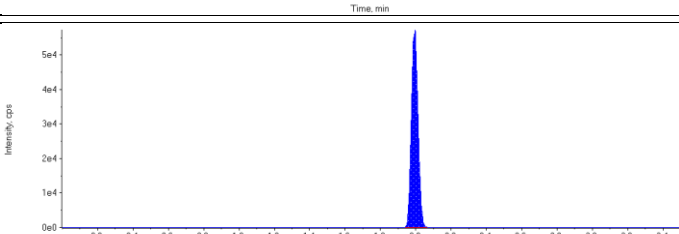
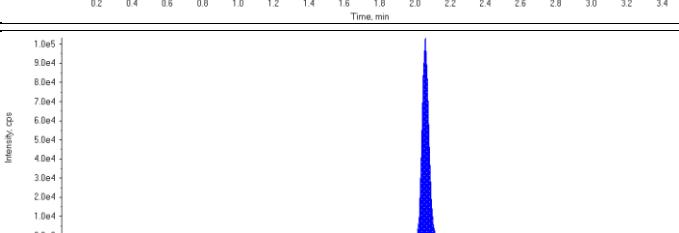
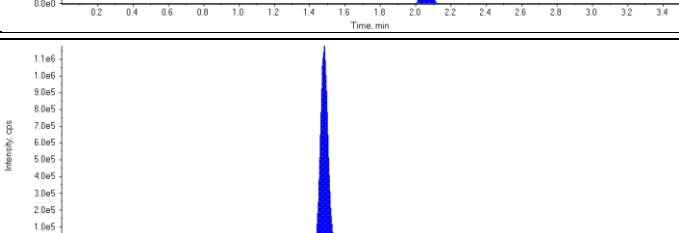
<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.16 (1.15) min</p> <p>Calculated Conc: 24.9 µg/L</p> <p>Area Ratio: 7.60</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 27.7 µg/L</p> <p>Area Ratio: 9.02</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 31.1 µg/L</p> <p>Area Ratio: 4.18</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 31.0 µg/L</p> <p>Area Ratio: 5.08</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 29.0 µg/L</p> <p>Area Ratio: 5.25</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 30.1 µg/L</p> <p>Area Ratio: 4.98</p> <p>Sample Type: (Quality Control)</p>	

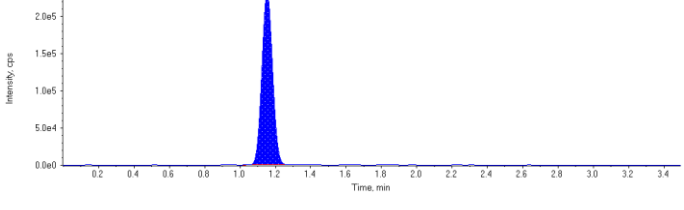
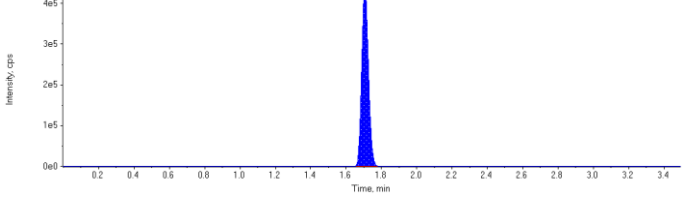
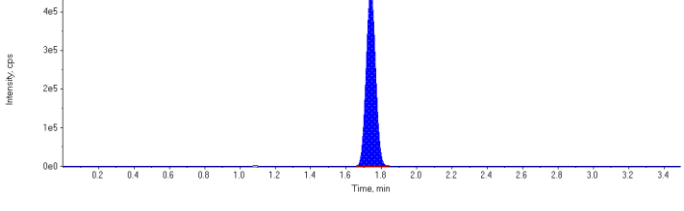
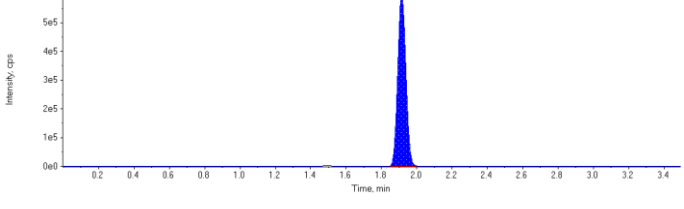
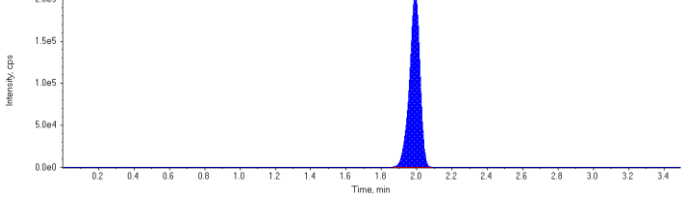
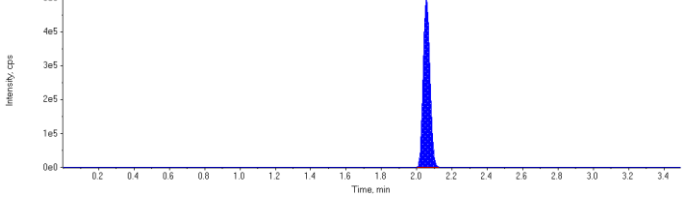
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 95.3 µg/L</p> <p>Area Ratio: 0.0351</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 90.7 µg/L</p> <p>Area Ratio: 0.112</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.88) min</p> <p>Calculated Conc: 96.7 µg/L</p> <p>Area Ratio: 0.104</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 106. µg/L</p> <p>Area Ratio: 0.0457</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.06 (2.02) min</p> <p>Calculated Conc: 111. µg/L</p> <p>Area Ratio: 0.0820</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 84.0 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

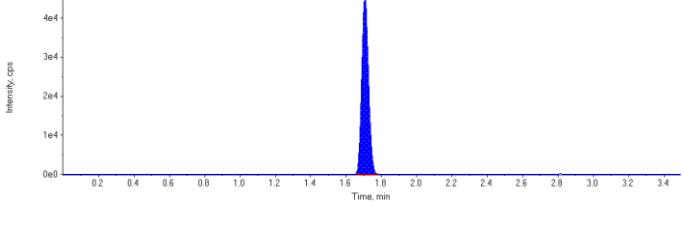
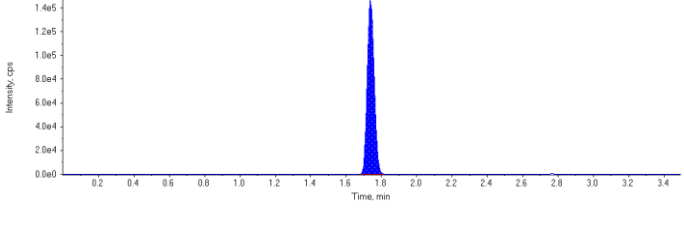
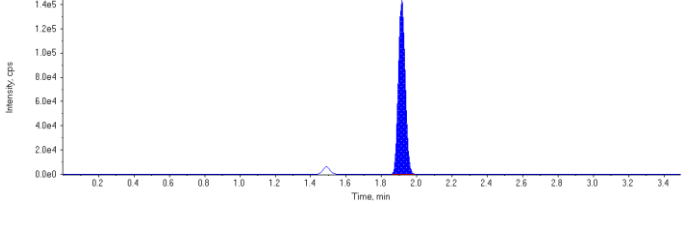
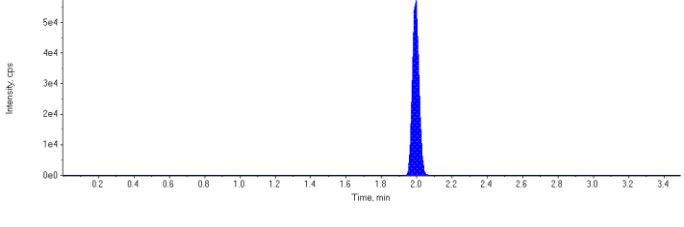
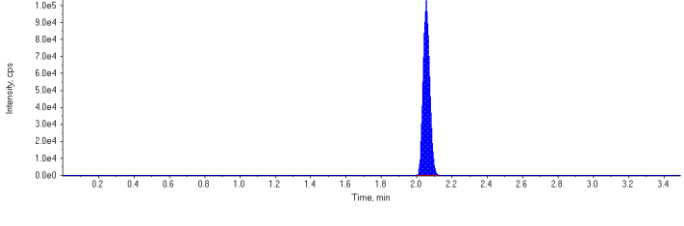
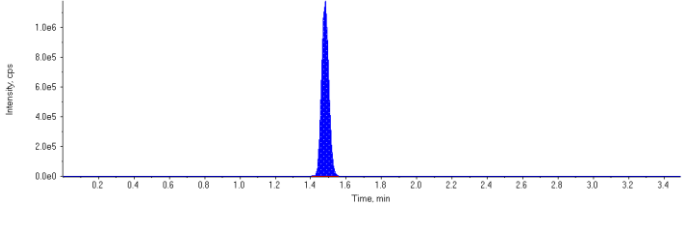
<b>Sample Name</b>	<b>CCV</b>	<b>Injection Vial</b>	6
<b>Sample ID</b>	CCV	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 10:43:32 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	122000.	1.71	1.00	-
MPFHpA	400000.	1.74	1.00	-
MPFOA	381000.	1.91	1.00	-
MPFOS	150000.	1.99	1.00	-
MPFNA	266000.	2.05	1.00	-
13C6-PFHxA IS	3360000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	975000	1.15	30.0	26.0	86.7
PFHxS 1	1160000	1.71	30.0	29.2	97.3
PFHpA 1	1680000	1.74	30.0	31.2	104.0
PFOA 1	1910000	1.92	30.0	30.5	102.0
PFOS 1	851000	1.99	30.0	31.3	104.0
PFNA 1	1330000	2.05	30.0	30.3	101.0
18O2-PFHxS	122000	1.71	100.	99.0	99.0
13C4-PFHpA	400000	1.74	100.	96.4	96.4
13C4-PFOA	381000	1.91	100.	105.	105.0
13C4-PFOS	150000	1.99	100.	103.	103.0
13C5-PFNA	266000	2.05	100.	108.	108.0
13C6-PFHxA	3360000	1.48	100.	84.1	84.1

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.15 (1.15) min</p> <p>Calculated Conc: 26.0 µg/L</p> <p>Area Ratio: 7.96</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 29.2 µg/L</p> <p>Area Ratio: 9.51</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 31.2 µg/L</p> <p>Area Ratio: 4.20</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 30.5 µg/L</p> <p>Area Ratio: 5.00</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 31.3 µg/L</p> <p>Area Ratio: 5.67</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 30.3 µg/L</p> <p>Area Ratio: 5.01</p> <p>Sample Type: (Quality Control)</p>	

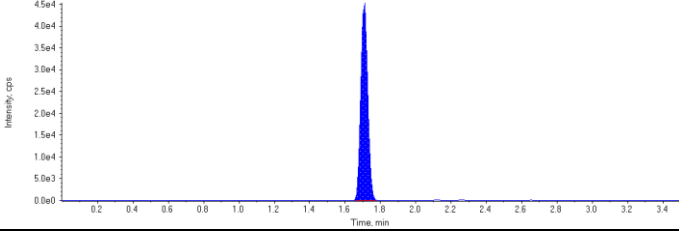
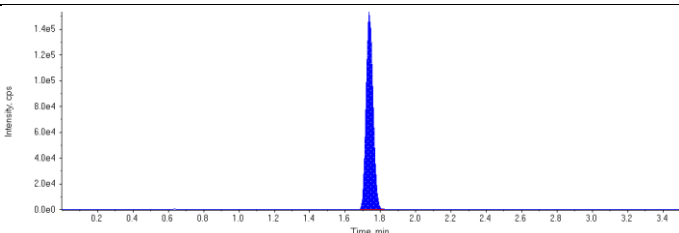
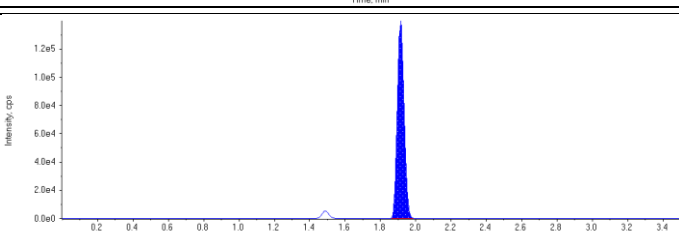
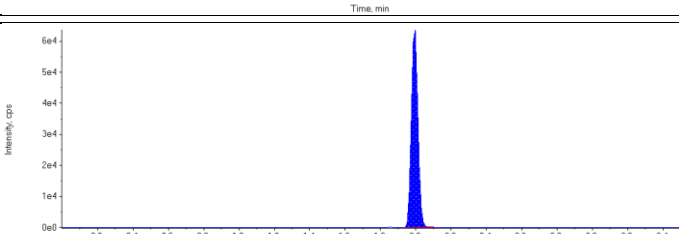
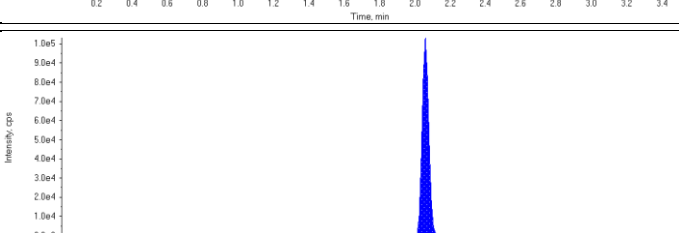
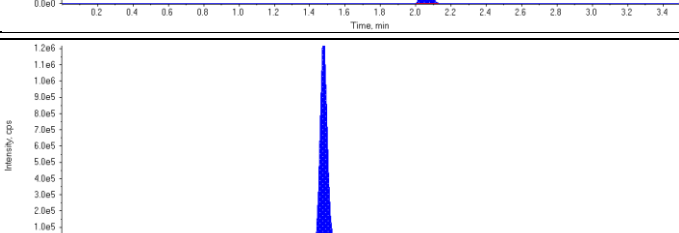
<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 99.0 µg/L</p> <p>Area Ratio: 0.0365</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 96.4 µg/L</p> <p>Area Ratio: 0.119</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 105. µg/L</p> <p>Area Ratio: 0.114</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.0447</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 108. µg/L</p> <p>Area Ratio: 0.0793</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 84.1 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	

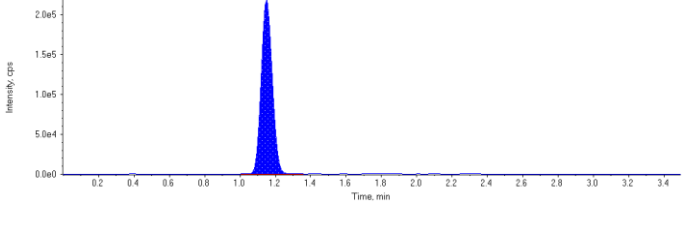
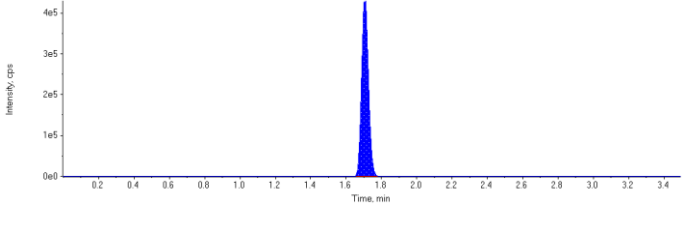
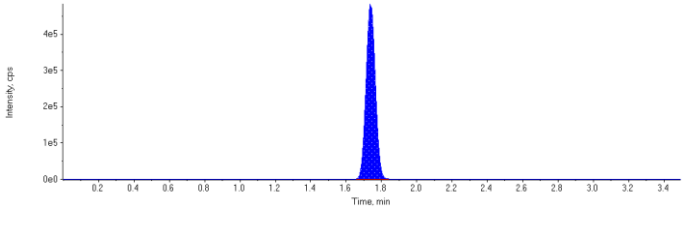
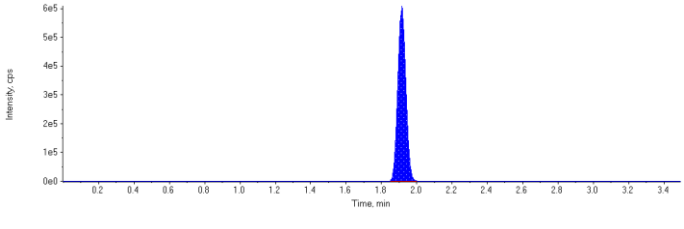
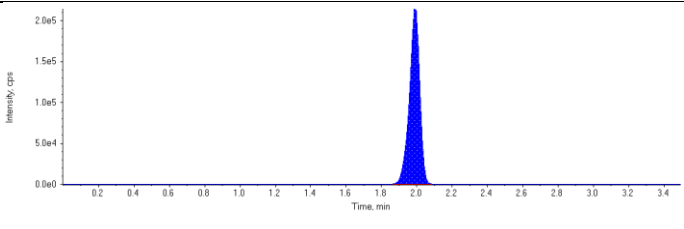
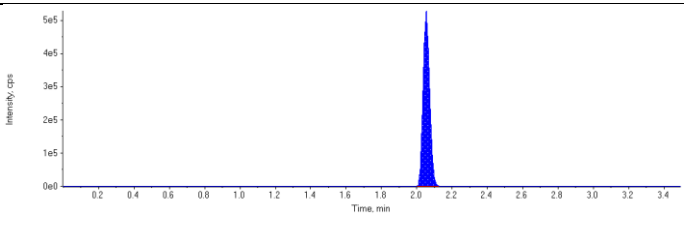


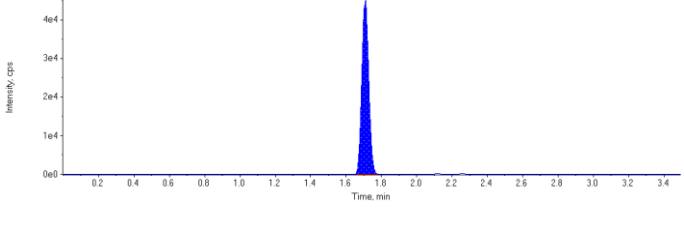
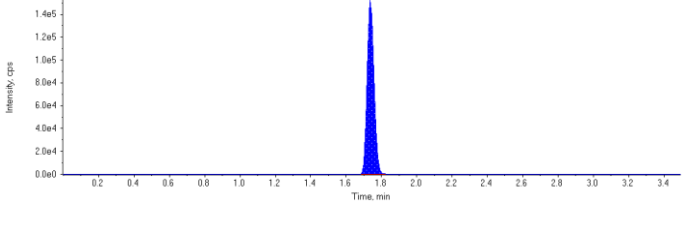
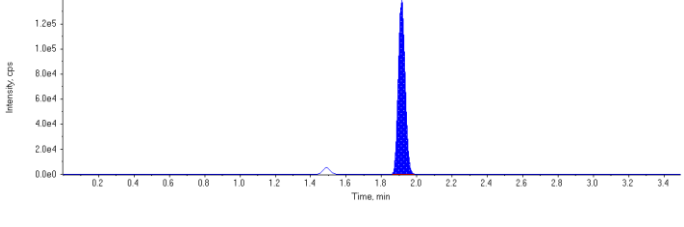
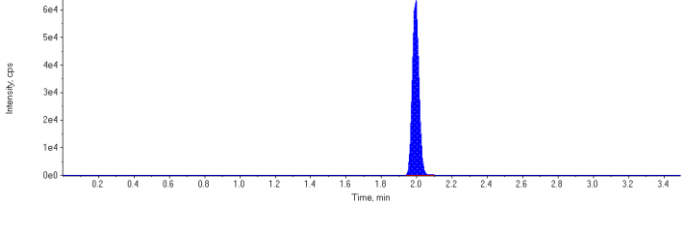
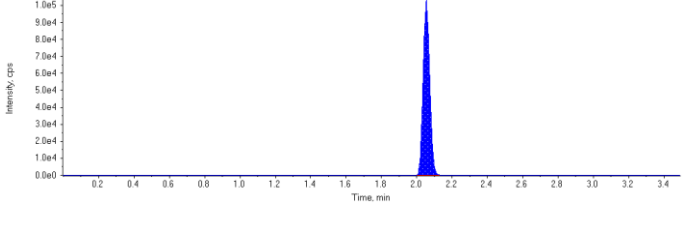
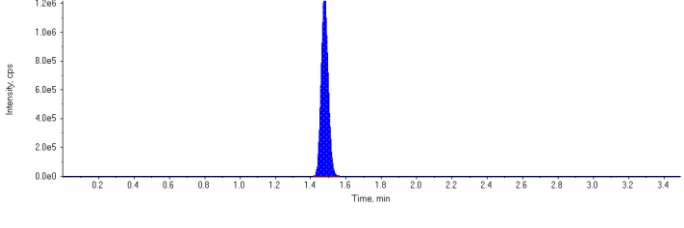
<b>Sample Name</b>	CCV	<b>Injection Vial</b>	6
<b>Sample ID</b>	CCV	<b>Injection Volume (µL)</b>	3
<b>Sample Type</b>	Quality Control	<b>Algorithm Used</b>	Analyst Classic
<b>Acquisition Date</b>	2016/02/19 11:03:54 AM	<b>Dilution Factor</b>	1.00
<b>Acquisition Method</b>	PFC_Water_Low.dam	<b>Sample Annotation</b>	-
<b>Project</b>	Enviro\PFOS	<b>Instrument Name</b>	LCMS03
<b>Data File</b>	PFC_160219\WS#4386408.wiff		
<b>Result Table</b>	PFC_Water_160219_4386408_ULow.rdb		

Internal Standard	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)
MPFHxS	123000.	1.71	1.00	-
MPFHpA	413000.	1.74	1.00	-
MPFOA	372000.	1.91	1.00	-
MPFOS	165000.	1.99	1.00	-
MPFNA	267000.	2.05	1.00	-
13C6-PFHxA IS	3510000.	1.48	1.00	-
N/A	N/A	N/A	N/A	-

Target Analyte	Area (cps)	RT (min)	Target Conc. (ng/L)	Calc. Conc. (ng/L)	Accuracy (%)
PFBS 1	950000	1.15	30.0	25.3	84.2
PFHxS 1	1180000	1.71	30.0	29.5	98.2
PFHpA 1	1770000	1.74	30.0	31.8	106.0
PFOA 1	1910000	1.92	30.0	31.4	105.0
PFOS 1	881000	1.99	30.0	29.4	98.1
PFNA 1	1370000	2.05	30.0	30.9	103.0
18O2-PFHxS	123000	1.71	100.	95.1	95.1
13C4-PFHpA	413000	1.74	100.	95.1	95.1
13C4-PFOA	372000	1.91	100.	98.3	98.3
13C4-PFOS	165000	1.99	100.	109.	109.0
13C5-PFNA	267000	2.05	100.	103.	103.0
13C6-PFHxA	3510000	1.48	100.	87.9	87.9

<p><b>MPFHxS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.71(1.68) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFHpA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.74(1.75) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOA</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.91(1.88) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFOS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.99(1.97) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>MPFNA</b> (Internal Standard)</p> <p>RT (Exp. RT): 2.05(2.02) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA IS</b> (Internal Standard)</p> <p>RT (Exp. RT): 1.48(1.42) min Concentration: 1.00 ng/L Sample Type: (Quality Control)</p>	
<p><b>N/A</b> (Internal Standard)</p> <p>RT (Exp. RT): N/A(N/A) min Concentration: N/A N/A Sample Type: (Quality Control)</p>	<p style="text-align: center;">This image is not available</p>

<p><b>PFBS 1</b> (298.900/79.900 Da)</p> <p>RT (Exp. RT): 1.15 (1.15) min</p> <p>Calculated Conc: 25.3 µg/L</p> <p>Area Ratio: 7.73</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHxS 1</b> (398.900/79.900 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 29.5 µg/L</p> <p>Area Ratio: 9.60</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFHpA 1</b> (363.000/319.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 31.8 µg/L</p> <p>Area Ratio: 4.28</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOA 1</b> (413.100/369.000 Da)</p> <p>RT (Exp. RT): 1.92 (1.92) min</p> <p>Calculated Conc: 31.4 µg/L</p> <p>Area Ratio: 5.14</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFOS 1</b> (498.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 29.4 µg/L</p> <p>Area Ratio: 5.33</p> <p>Sample Type: (Quality Control)</p>	
<p><b>PFNA 1</b> (462.900/419.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 30.9 µg/L</p> <p>Area Ratio: 5.12</p> <p>Sample Type: (Quality Control)</p>	

<p><b>18O2-PFHxS</b> (402.900/84.000 Da)</p> <p>RT (Exp. RT): 1.71 (1.68) min</p> <p>Calculated Conc: 95.1 µg/L</p> <p>Area Ratio: 0.0350</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFHpA</b> (366.900/322.000 Da)</p> <p>RT (Exp. RT): 1.74 (1.75) min</p> <p>Calculated Conc: 95.1 µg/L</p> <p>Area Ratio: 0.118</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOA</b> (416.900/372.000 Da)</p> <p>RT (Exp. RT): 1.91 (1.88) min</p> <p>Calculated Conc: 98.3 µg/L</p> <p>Area Ratio: 0.106</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C4-PFOS</b> (502.900/79.900 Da)</p> <p>RT (Exp. RT): 1.99 (1.97) min</p> <p>Calculated Conc: 109. µg/L</p> <p>Area Ratio: 0.0471</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C5-PFNA</b> (467.900/423.000 Da)</p> <p>RT (Exp. RT): 2.05 (2.02) min</p> <p>Calculated Conc: 103. µg/L</p> <p>Area Ratio: 0.0762</p> <p>Sample Type: (Quality Control)</p>	
<p><b>13C6-PFHxA</b> (318.900/274.000 Da)</p> <p>RT (Exp. RT): 1.48 (1.42) min</p> <p>Calculated Conc: 87.9 µg/L</p> <p>Area Ratio: 0.00</p> <p>Sample Type: (Quality Control)</p>	



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Maxxam Analytics International  
6740 Campobello Rd.  
Mississauga, Ontario, Canada  
L5N 2L8  
1-800-668-0639  
[www.maxxamanalytics.com](http://www.maxxamanalytics.com)

# Shipping and Receiving Documents

# Chain of Custody Record

TAL-4124 (1007)

Client: **CH2M Hill** Project Manager: **Bill Friedman** Chain of Custody Number: **283602**  
 Address: **5701 Cleveland St, Suite 200** Telephone Number (Area Code)/Fax Number: **757-671-6223** Date: **02/03/16** Page **1** of **1**  
 City: **Virginia Beach** State: **VA** Zip Code: **23462** Site Contact: \_\_\_\_\_ Lab Contact: \_\_\_\_\_  
 Project Name and Location (State): **CTD WFTG PFC Sampling** Carrier/Waybill Number: **FedEx**

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	Soil	Sed	Unpres.	H2SO4	HNO3	HCl	HOBr	ZnBr			HOBr
OF-FB09-0216	02/03/16	1646	X			X								
OF-RW09-0216		1645												
OF-FB37-0216		1720												
OF-RW37-0216		1728												
OF-RW11-0216		1604												
OF-FB11-0216		1610												
OF-RW28-0216		1741												
OF-FB28-0216		1743												
OF-RW67-0216		1819												
OF-FB67-0216		1821												

Possible Hazard Identification:  Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown  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	Date	Time
1. Relinquished By	02/03/16	19:30
2. Relinquished By		
3. Relinquished By		

Comments: **Yalman Smith** 1. Received By: **[Signature]** Date: **2/4/16** Time: **945**  
 2. Received By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 3. Received By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 Comments: **0.30**

**Chain of Custody Record**

Due 2/22

Temperature on Receipt \_\_\_\_\_



CTD WFTG

Drinking Water? Yes  No

320-17154 Chain of Custody

G

Project Manager: \_\_\_\_\_ Date: 02/03/16 Chain of Custody Number: 283602  
 Telephone Number (Area Code)/Fax Number: \_\_\_\_\_ Lab Number: \_\_\_\_\_  
 Site Contact: \_\_\_\_\_ Lab Contact: \_\_\_\_\_ Page 1 of 1

Carrier/Waybill Number: FedEx

Sample I.D. No. and Description  
 (Containers for each sample may be combined on one line)

Date	Time	Matrix						Containers & Preservatives						Select PEGs		
		Air	Aqueous	Solid	Soil	Urine	MSO4	HW03	HC1	NuOH	ZnAc2	NuOH				
<del>02/03/16</del>	<del>16:45</del>															
02/03/16	16:45														2	
<del>02/03/16</del>	<del>17:28</del>															
02/03/16	17:28														2	
<del>02/03/16</del>	<del>16:04</del>															
02/03/16	16:04														2	
<del>02/03/16</del>	<del>17:41</del>															
02/03/16	17:41														2	
<del>02/03/16</del>	<del>18:19</del>															
02/03/16	18:19														2	
<del>02/03/16</del>	<del>18:21</del>															

Special Instructions/Conditions of Receipt

13-Feb-16 13:40  
 Hongmei Zhao (Grace)  
  
 B630797  
 RGN ENV-951

Possible Hazard Identification:  Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown

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 \_\_\_\_\_

QC Requirements (Specify)

1. Relinquished By: <u>Yathram Smith</u>	Date: 02/03/16	Time: 19:30	1. Received By: <u>[Signature]</u>	Date: 2/4/16	Time: 9:45
2. Relinquished By: _____	Date: _____	Time: _____	2. Received By: <u>[Signature]</u>	Date: 2016/02/13	Time: 13:40
3. Relinquished By: _____	Date: _____	Time: _____	3. Received By: _____	Date: _____	Time: _____

Comments

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

0-3-16 REFER TO ACTR



# Login Sample Receipt Checklist

Client: CH2M Hill, Inc.

Job Number: 320-17154-1  
SDG Number: CTO WE7G PFC Sampling

**Login Number: 17154**  
**List Number: 1**  
**Creator: Nelson, Kym D**

**List Source: TestAmerica Sacramento**

Question	Answer	Comment
Radioactivity wasn't checked or is $\leq$ background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	
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.	True	
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.	N/A	
Sample Preservation Verified.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is $<6\text{mm}$ (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	



## Data Validation Summary

### Oceana CTO-WE44, NALF Fentress

TO: Juliana Dean/VBO  
Anita Dodson/VBO

FROM: Tiffany McGlynn/GNV

CC: Herb Kelly/GNV

DATE: March 18, 2016

#### Introduction

The following data validation report discusses the data validation process and findings for TestAmerica Laboratories and Maxxam Laboratories in the Sample Delivery Groups (SDGs) listed in the table below.

Samples were analyzed using the following analytical methods:

- WS-LC-0025 & 537 MOD Perfluorinated Hydrocarbons

The samples included in these SDGs are listed in the table below.

SDG	Sample Name	Matrix
320-17150	OF-RW42B-0216	Water
320-17150	OF-RW39-0216	Water
320-17150	OF-FB40-0216	Water
320-17150	OF-RW40-0216	Water
320-17150	OF-FB43-0216	Water
320-17150	OF-RW43-0216	Water
320-17150	OF-FB42B-0216	Water
320-17150	OF-RW42A-0216	Water
320-17150	OF-FB42A-0216	Water
320-17150	OF-RW35-0216	Water
320-17150	OF-FB35-0216	Water

<b>SDG</b>	<b>Sample Name</b>	<b>Matrix</b>
320-17150	OF-RW58-0216	Water
320-17150	OF-FB58-0216	Water
320-17150	OF-FB39-0216	Water
320-17154	OF-FB09-0216	Water
320-17154	OF-FB67-0216	Water
320-17154	OF-RW09-0216	Water
320-17154	OF-FB37-0216	Water
320-17154	OF-RW37-0216	Water
320-17154	OF-RW11-0216	Water
320-17154	OF-FB11-0216	Water
320-17154	OF-RW28-0216	Water
320-17154	OF-FB28-0216	Water
320-17154	OF-RW67-0216	Water
320-17183	OF-RW66-0216	Water
320-17183	OF-FB27-0216	Water
320-17183	OF-FB66-0216	Water
320-17183	OF-RW49-0216	Water
320-17183	OF-FB49-0216	Water
320-17183	OF-RW36A-0216	Water
320-17183	OF-FB36A-0216	Water
320-17183	OF-RW51A-0216	Water
320-17183	OF-FB51A-0216	Water
320-17183	OF-RW27-0216	Water
320-17184	OF-RW20-0216	Water
320-17184	OF-FB30-0216	Water
320-17184	OF-FB69-0216	Water
320-17184	OF-RW69-0216	Water
320-17184	OF-FB26-0216	Water
320-17184	OF-RW26-0216	Water
320-17184	OF-FB20-0216	Water
320-17184	OF-RW55-0216	Water
320-17184	OF-FB55-0216	Water
320-17184	OF-RW54-0216	Water
320-17184	OF-FB54-0216	Water
320-17184	OF-RW68-0216	Water
320-17184	OF-FB68-0216	Water
320-17184	OF-RW30-0216	Water
320-17185	OF-FB08-0216	Water
320-17185	OF-RW51-0216	Water
320-17185	OF-RW51P-0216	Water

<b>SDG</b>	<b>Sample Name</b>	<b>Matrix</b>
320-17185	OF-RW08-0216	Water
320-17185	OF-RW08P-0216	Water
320-17185	OF-FB41-0216	Water
320-17185	OF-RW41-0216	Water
320-17185	OF-RW41P-0216	Water
320-17185	OF-FB56-0216	Water
320-17185	OF-RW56-0216	Water
320-17185	OF-FB51-0216	Water
320-17190	OF-FB12-0216	Water
320-17190	OF-RW12-0216	Water
320-17190	OF-FB57-0216	Water
320-17190	OF-RW57-0216	Water
320-17190	OF-RW57P-0216	Water
320-17190	OF-FB25-0216	Water
320-17190	OF-RW25-0216	Water
320-17190	OF-FB16-0216	Water
320-17190	OF-RW16-0216	Water
320-17219	OF-FB47-0216	Water
320-17219	OF-RW47-0216	Water
320-17219	OF-FB47A-0216	Water
320-17219	OF-RW47A-0216	Water
320-17219	OF-FB48-0216	Water
320-17219	OF-RW48-0216	Water
320-17236	OF-FB70-0216	Water
320-17236	OF-RW70-0216	Water
320-17236	OF-FB44-0216	Water
320-17236	OF-RW44-0216	Water
320-17236	OF-RW44P-0216	Water
320-17236	OF-FB65-0216	Water
320-17236	OF-RW65-0216	Water
320-17236	OF-FB21-0216	Water
320-17236	OF-RW21-0216	Water
320-17241	OF-FB62-0216	Water
320-17241	OF-RW34-0216	Water
320-17241	OF-FB38-0216	Water
320-17241	OF-RW38-0216	Water
320-17241	OF-RW62-0216	Water
320-17241	OF-FB63-0216	Water
320-17241	OF-RW63-0216	Water
320-17241	OF-FB59-0216	Water

<b>SDG</b>	<b>Sample Name</b>	<b>Matrix</b>
320-17241	OF-RW59-0216	Water
320-17241	OF-FB50-0216	Water
320-17241	OF-RW50-0216	Water
320-17241	OF-FB34-0216	Water
320-17278	OF-FB24-0216	Water
320-17278	OF-RW24-0216	Water
320-17278	OF-FB31-0216	Water
320-17278	OF-RW31-0216	Water
320-17278	OF-FB60-0216	Water
320-17278	OF-RW60-0216	Water
320-17278	OF-RW60P-0216	Water
320-17278	OF-FB46-0216	Water
320-17278	OF-RW46-0216	Water
320-17321	OF-FB02-0216	Water
320-17321	OF-RW02-0216	Water
320-17321	OF-FB15-0216	Water
320-17321	OF-RW15-0216	Water
320-17321	OF-FB18-0216	Water
320-17321	OF-RW18-0216	Water
320-17859	OF-FB07-0316	Water
320-17859	OF-RW07-0316	Water
320-17859	OF-HPFB01-0316	Water
320-17859	OF-HP01-0316	Water

## **Data Evaluation**

Data was evaluated in accordance with the analytical methods and with the criteria found in the following guidance documents: Sampling and Analysis Plan Perfluorinated Compound Investigation, Naval Auxiliary Landing Field Fentress, Chesapeake, Virginia Contract Task Order WE44 (December 2015) and National Functional Guidelines for Organic Data Review (August 2014) with Region 3 Modification (Use of 'B' qualifier) as applicable. The samples were evaluated based on the following criteria:

- Data Completeness
- Technical Holding Times
- Tuning Instrument
- Initial/Continuing Calibrations
- Blanks

- Internal Standards
- Laboratory Control Samples
- Isotope Dilution Analyte
- Field Duplicates
- Identification/Quantitation
- Reporting Limits

## **Overall Evaluation of Data/Potential Usability Issues**

Specific details regarding qualification of the data are addressed in the sections below. If an issue is not addressed there were no actions required based on unmet quality criteria. When more than one qualifier is associated with a compound/analyte, the validator has chosen the qualifier that best indicates possible bias in the results and qualified these data accordingly.

### **Data Completeness**

The SDG was received complete and intact.

### **Technical Holding Times**

According to the chain of custody records, sampling was performed on 2/3/16 through 2/16/16. Samples were received at the laboratory 2/4/16 through 2/17/16. All sample preparation and analyses were performed within holding time requirements with the exception of the samples listed below. Affected data are summarized in **Attachment 1**.

<b>Sample Name</b>	<b>SDG</b>
OF-RW42B-0216	320-17150
OF-RW08-0216	320-17185
OF-RW08P-0216	320-17185
OF-FB62-0216	320-17241
OF-RW34-0216	320-17241
OF-FB38-0216	320-17241
OF-RW38-0216	320-17241
OF-RW62-0216	320-17241
OF-FB63-0216	320-17241
OF-RW63-0216	320-17241
OF-FB59-0216	320-17241

Sample Name	SDG
OF-RW59-0216	320-17241
OF-FB50-0216	320-17241
OF-RW50-0216	320-17241
OF-FB34-0216	320-17241

## Blanks

Several compounds were detected in the field blanks and method blanks as listed below. Affected data are summarized in **Attachment 1**.

SDG	Blank ID	Compound	Conc.	Units
320-17183	OF-FB49-0216	Perfluorohexanesulfonic acid (PFHxS)	0.00068	UG_L
320-17183	OF-FB36A-0216	Perfluorooctane Sulfonate (PFOS)	0.00042	UG_L
320-17185	OF-FB51-0216	Perfluorobutanesulfonic acid (PFBS)	0.00063	UG_L
320-17190	OF-FB12-0216	Perfluorohexanesulfonic acid (PFHxS)	0.00079	UG_L
320-17190	OF-FB57-0216	Perfluorohexanesulfonic acid (PFHxS)	0.00083	UG_L
320-17190	OF-FB25-0216	Perfluorobutanesulfonic acid (PFBS)	0.00092	UG_L
320-17190	OF-FB16-0216	Perfluorobutanesulfonic acid (PFBS)	0.0011	UG_L
320-17190	MB 320-100277/1-A	Perfluorobutanesulfonic acid (PFBS)	0.00103	UG_L
320-17190	MB 320-100277/1-A	Perfluorohexanesulfonic acid (PFHxS)	0.00102	UG_L
320-17190	MB 320-100277/1-A	Perfluorooctane Sulfonate (PFOS)	0.00144	UG_L
320-17859	MB 320-104553/1-A	Perfluorooctanoic acid (PFOA)	0.00217	UG_L

## Field Duplicate Precision

Perfluoroheptanoic acid (PFHpA) did not meet required precision criteria in native sample OF-RW51-0216 and field duplicate OF-RW51P-0216. Affected data are summarized in **Attachment 1**.

## Matrix Spike/Spike Duplicate

For spiked sample OF-RW56-0216 in SDG 320-17185, perfluorobutanesulfonic acid (PFBS) exhibited high recoveries in the MS/MSD. Affected data are summarized in **Attachment 1**.

## Surrogates

Surrogates for the samples listed below exhibited low recoveries. Affected data are summarized in **Attachment 1**.

Sample Name	SDG
OF-RW67-0216	320-17154
OF-RW47-0216	320-17219
OF-RW70-0216	320-17236



Sample Name	SDG
OF-RW24-0216	320-17278

### **Internal Standards**

Internal standards exhibited low recoveries for the samples listed below. Affected data are summarized in **Attachment 1**.

Sample Name	SDG
OF-RW37-0216	320-17154
OF-FB56-0216	320-17185

### **Conclusion**

These data can be used in the project decision-making process as qualified by the data quality evaluation process.

Please do not hesitate to contact us about this validation report.

Sincerely,



Tiffany McGlynn

## Qualification Flags

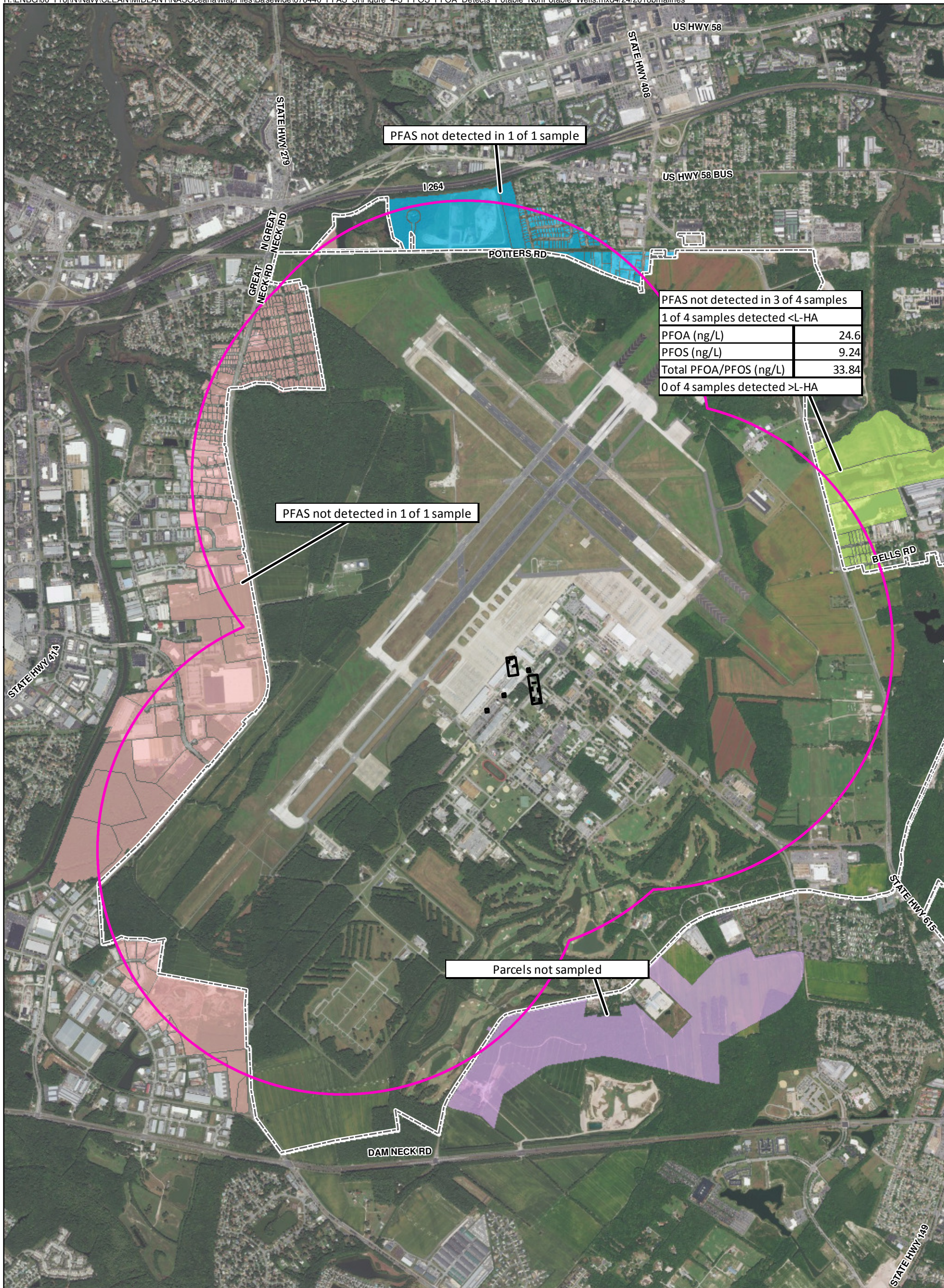
Exclude	More appropriate data exist for this analyte.
R	Data were rejected for use.
UL	Analyte not detected, quantitation limit is potentially biased low.
UJ	Analyte not detected, estimated quantitation limit.
U	Analyte not detected.
B	Not detected substantially above the level reported in laboratory or field blanks.
L	Analyte present, estimated value potentially biased low.
K	Analyte present, estimated value potentially biased high.
N	Analyte identification presumptive; no second column analysis performed or GC/MS tentative identification.
J	Analyte present, estimated value.
NJ	Analysis indicates the presence of an analyte that was "tentatively identified" and the associated value represents its approximate concentration.
None	Placeholder for calculating quality control issues that do not require flagging.
=	Analyte was detected at a concentration greater than the quantitation limit.

## Qualifier Code Reference

<b>Value</b>	<b>Description</b>
%SOL	High Moisture content
2C	Second Column – Poor Dual Column Reproducibility
2S	Second Source – Bad reproducibility between tandem detectors
BD	Blank Spike/Blank Spike Duplicate(LCS/LCSD) Precision
BRL	Below Reporting Limit
BSH	Blank Spike/LCS – High Recovery
BSL	Blank Spike/LCS – Low Recovery
CC	Continuing Calibration
CCBL	Continuing Calibration Blank Contamination
CCH	Continuing Calibration Verification – High Recovery
CCL	Continuing Calibration Verification – Low Recovery
DL	Redundant Result – due to Dilution
EBL	Equipment Blank Contamination
EMPC	Estimated Possible Maximum Concentration
ESH	Extraction Standard - High Recovery
ESL	Extraction Standard - Low Recovery
FBL	Field Blank Contamination
FD	Field Duplicate
HT	Holding Time
ICB	Initial Calibration – Bad Linearity or Curve Function
ICH	Initial Calibration – High Relative Response Factors
ICL	Initial Calibration – Low Relative Response Factors
IR15	Ion ratio exceeds +/- 15% difference
ISH	Internal Standard – High Recovery
ISL	Internal Standard – Low Recovery
LD	Lab Duplicate Reproducibility
LR	Concentration Exceeds Linear Range
MBL	Method Blank Contamination
MDP	Matrix Spike/Matrix Spike Duplicate Precision
MI	Matrix interference obscuring the raw data

<b>Value</b>	<b>Description</b>
MSH	Matrix Spike and/or Matrix Spike Duplicate – High Recovery
MSL	Matrix Spike and/or Matrix Spike Duplicate – Low Recovery
OT	Other
PD	Pesticide Degradation
RE	Redundant Result - due to Reanalysis or Re-extraction
SD	Serial Dilution Reproducibility
SSH	Spiked Surrogate – High Recovery
SSL	Spiked Surrogate – Low Recovery
TBL	Trip Blank Contamination
TN	Tune





- Legend**
- Non-Core Target Treatment Area (2004)
  - - Core Target Treatment Area (2004) (Core)
  - ▭ Sampling Area
  - ▭ Installation Boundary
  - Off-Base Parcels**
  - ▭ East
  - ▭ North
  - ▭ South
  - ▭ West

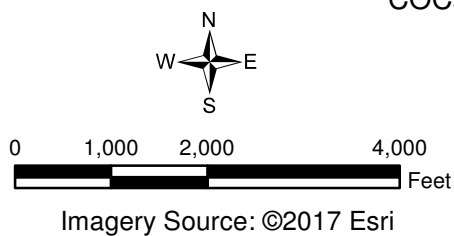


Figure 4-3  
COCs Detections in Potable Wells Sampled from Parcels Located Off-Base  
Basewide Per- and Polyfluoroalkyl Substances Site Inspection Report  
NAS Oceana, Virginia Beach, Virginia