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LABORATORY DATA PACKAGE, 320-10459-1, NAS BRUNSWICK ME  
12/12/2014  
TEST AMERICA

## ANALYTICAL REPORT

Job Number: 320-10459-1

Job Description: NAS Brunswick Maine 13-CTO WE09

For:

Tetra Tech, Inc.  
Foster Plaza VII  
661 Anderson Drive  
Pittsburgh, PA 15220

Attention: Jeff Orient



Approved for release.  
Karen M Sellers  
Project Manager II  
12/12/2014 4:08 PM

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12/12/2014

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# Definitions/Glossary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

## Qualifiers

### LCMS

Qualifier	Qualifier Description
H	Sample was prepped or analyzed beyond the specified holding time
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.

□	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: Tetra Tech, Inc.**

**Project: NAS Brunswick Maine 13-CTO WE09**

**Report Number: 320-10459-1**

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

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

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

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

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

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

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

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

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

### **RECEIPT**

The samples were received on 11/18/2014 9:50 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 0.5° C.

### **PFOA/PFOS**

#### **Method WS-LC-0025:**

The following samples were re-extracted past the 7 day extraction hold time due to contamination in the method blank (MB) and laboratory control sample (LCS) for the original prep batch: B7/10-GW02 (320-10459-1), B7/10-GW03S (320-10459-2), B7/10-GW06S (320-10459-4), EFA-FD01-1014 (320-10459-6), H123-GW02 (320-10459-5), QR4-GW02 (320-10459-3). There is no prescribed regulatory holding time for PFOA and PFOS. The 7 day extraction holding time is based on general EPA convention for the holding time of extractable organic compounds in water.

Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate/sample duplicate (MS/MSD/DUP) associated with batch 320-59712.

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

# Detection Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

## Client Sample ID: B7/10-GW02

## Lab Sample ID: 320-10459-1

Analyte	Result	Qualifier	LOQ	DL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	10	H	2.2	0.83	ng/L	1		WS-LC-0025	Total/NA
Perfluorooctane Sulfonate (PFOS)	6.6	H	2.2	1.4	ng/L	1		WS-LC-0025	Total/NA

## Client Sample ID: B7/10-GW03S

## Lab Sample ID: 320-10459-2

Analyte	Result	Qualifier	LOQ	DL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	2.0	J H	2.3	0.87	ng/L	1		WS-LC-0025	Total/NA

## Client Sample ID: QR4-GW02

## Lab Sample ID: 320-10459-3

No Detections.

## Client Sample ID: B7/10-GW06S

## Lab Sample ID: 320-10459-4

Analyte	Result	Qualifier	LOQ	DL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	20	H	2.2	0.84	ng/L	1		WS-LC-0025	Total/NA
Perfluorooctane Sulfonate (PFOS)	2.2	H	2.2	1.4	ng/L	1		WS-LC-0025	Total/NA

## Client Sample ID: H123-GW02

## Lab Sample ID: 320-10459-5

Analyte	Result	Qualifier	LOQ	DL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	17	H	2.2	0.82	ng/L	1		WS-LC-0025	Total/NA

## Client Sample ID: EFA-FD01-1014

## Lab Sample ID: 320-10459-6

Analyte	Result	Qualifier	LOQ	DL	Unit	Dil Fac	D	Method	Prep Type
Perfluorooctanoic acid (PFOA)	16	H	2.2	0.84	ng/L	1		WS-LC-0025	Total/NA

This Detection Summary does not include radiochemical test results.

TestAmerica Sacramento

# Client Sample Results

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

**Client Sample ID: B7/10-GW02**

**Lab Sample ID: 320-10459-1**

Date Collected: 11/15/14 12:05

Matrix: Water

Date Received: 11/18/14 09:50

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroctanoic acid (PFOA)	10	H	2.2	0.83	ng/L		12/03/14 13:38	12/04/14 20:45	1
Perfluorooctane Sulfonate (PFOS)	6.6	H	2.2	1.4	ng/L		12/03/14 13:38	12/04/14 20:45	1
<i>Isotope Dilution</i>	%Recovery	Qualifier	Limits						
13C4 PFOS	116		25 - 150				12/03/14 13:38	12/04/14 20:45	1
13C4 PFOA	85		25 - 150				12/03/14 13:38	12/04/14 20:45	1

**Client Sample ID: B7/10-GW03S**

**Lab Sample ID: 320-10459-2**

Date Collected: 11/15/14 15:20

Matrix: Water

Date Received: 11/18/14 09:50

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroctanoic acid (PFOA)	2.0	J H	2.3	0.87	ng/L		12/03/14 13:38	12/04/14 21:07	1
Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.5	ng/L		12/03/14 13:38	12/04/14 21:07	1
<i>Isotope Dilution</i>	%Recovery	Qualifier	Limits						
13C4 PFOS	108		25 - 150				12/03/14 13:38	12/04/14 21:07	1
13C4 PFOA	64		25 - 150				12/03/14 13:38	12/04/14 21:07	1

**Client Sample ID: QR4-GW02**

**Lab Sample ID: 320-10459-3**

Date Collected: 11/16/14 10:20

Matrix: Water

Date Received: 11/18/14 09:50

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctanoic acid (PFOA)	1.8	U H	2.3	0.88	ng/L		12/03/14 13:38	12/04/14 21:28	1
Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.5	ng/L		12/03/14 13:38	12/04/14 21:28	1
<i>Isotope Dilution</i>	%Recovery	Qualifier	Limits						
13C4 PFOS	109		25 - 150				12/03/14 13:38	12/04/14 21:28	1
13C4 PFOA	90		25 - 150				12/03/14 13:38	12/04/14 21:28	1

**Client Sample ID: B7/10-GW06S**

**Lab Sample ID: 320-10459-4**

Date Collected: 11/16/14 11:50

Matrix: Water

Date Received: 11/18/14 09:50

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroctanoic acid (PFOA)	20	H	2.2	0.84	ng/L		12/03/14 13:38	12/04/14 21:49	1
Perfluorooctane Sulfonate (PFOS)	2.2	H	2.2	1.4	ng/L		12/03/14 13:38	12/04/14 21:49	1
<i>Isotope Dilution</i>	%Recovery	Qualifier	Limits						
13C4 PFOS	108		25 - 150				12/03/14 13:38	12/04/14 21:49	1
13C4 PFOA	67		25 - 150				12/03/14 13:38	12/04/14 21:49	1

**Client Sample ID: H123-GW02**

**Lab Sample ID: 320-10459-5**

Date Collected: 11/16/14 15:00

Matrix: Water

Date Received: 11/18/14 09:50

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroctanoic acid (PFOA)	17	H	2.2	0.82	ng/L		12/03/14 13:38	12/04/14 22:10	1

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# Client Sample Results

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

## **Client Sample ID: H123-GW02**

Date Collected: 11/16/14 15:00

Date Received: 11/18/14 09:50

## **Lab Sample ID: 320-10459-5**

Matrix: Water

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluorooctane Sulfonate (PFOS)	1.6	U H	2.2	1.4	ng/L		12/03/14 13:38	12/04/14 22:10	1
<i>Isotope Dilution</i>									
	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOS	107		25 - 150				12/03/14 13:38	12/04/14 22:10	1
13C4 PFOA	89		25 - 150				12/03/14 13:38	12/04/14 22:10	1

## **Client Sample ID: EFA-FD01-1014**

Date Collected: 11/16/14 00:00

Date Received: 11/18/14 09:50

## **Lab Sample ID: 320-10459-6**

Matrix: Water

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

Analyte	Result	Qualifier	LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac
Perfluoroctanoic acid (PFOA)	16	H	2.2	0.84	ng/L		12/03/14 13:38	12/04/14 22:31	1
Perfluorooctane Sulfonate (PFOS)	1.7	U H	2.2	1.4	ng/L		12/03/14 13:38	12/04/14 22:31	1
<i>Isotope Dilution</i>									
	%Recovery	Qualifier	Limits				Prepared	Analyzed	Dil Fac
13C4 PFOS	109		25 - 150				12/03/14 13:38	12/04/14 22:31	1
13C4 PFOA	82		25 - 150				12/03/14 13:38	12/04/14 22:31	1

## Default Detection Limits

Client: Tetra Tech, Inc.

TestAmerica Job ID: 320-10459-1

Project/Site: NAS Brunswick Maine 13-CTO WE09

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

Analyte	LOQ	DL	Units	Method
Perfluorooctane Sulfonate (PFOS)	2.0	1.3	ng/L	WS-LC-0025
Perfluorooctanoic acid (PFOA)	2.0	0.75	ng/L	WS-LC-0025

## Isotope Dilution Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

Method: WS-LC-0025 - Perfluorinated Hydrocarbons

## Matrix: Water

## **Prep Type: Total/NA**

		Percent Isotope Dilution Recovery (Acceptance Limits)			
Lab Sample ID	Client Sample ID	13C4 PFOS		13C4 PFOA	
		(25-150)	(25-150)	(25-150)	(25-150)
320-10459-1	B7/10-GW02	116	85		
320-10459-2	B7/10-GW03S	108	64		
320-10459-3	QR4-GW02	109	90		
320-10459-4	B7/10-GW06S	108	67		
320-10459-5	H123-GW02	107	89		
320-10459-6	EFA-FD01-1014	109	82		
LCS 320-59712/2-A	Lab Control Sample	101	98		
MB 320-59712/1-A	Method Blank	111	109		

## **Surrogate Legend**

**13C4 PFOS = 13C4 PFOS**

13C4 PFOA = 13C4 PFOA

# QC Sample Results

Client: Tetra Tech, Inc.

TestAmerica Job ID: 320-10459-1

Project/Site: NAS Brunswick Maine 13-CTO WE09

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

**Lab Sample ID: MB 320-59712/1-A**

**Matrix: Water**

**Analysis Batch: 59779**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 59712**

Analyte	MB		LOQ	DL	Unit	D	Prepared	Analyzed	Dil Fac	
	Result	Qualifier								
Perfluorooctanoic acid (PFOA)	0.809	J	2.0	0.75	ng/L		12/03/14 13:38	12/04/14 20:03	1	
Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.3	ng/L		12/03/14 13:38	12/04/14 20:03	1	
<b>Isotope Dilution</b>										
Isotope Dilution	MB		MB		Limits		Prepared		Dil Fac	
	%Recovery	Qualifier								
13C4 PFOS	111				25 - 150					
13C4 PFOA	109				25 - 150					

**Lab Sample ID: LCS 320-59712/2-A**

**Matrix: Water**

**Analysis Batch: 59779**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 59712**

Analyte	Spike		LCS Result	LCS Qualifier	Unit	D	%Rec	%Rec.	Limits	
	Added									
Perfluorooctanoic acid (PFOA)	40.0		50.4		ng/L		126	60 - 140		
Perfluorooctane Sulfonate (PFOS)	38.2		39.7		ng/L		104	60 - 140		
<b>Isotope Dilution</b>										
Isotope Dilution	LCS		LCS		Limits					
	%Recovery	Qualifier								
13C4 PFOS	101				25 - 150					
13C4 PFOA	98				25 - 150					

# QC Association Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

## LCMS

### Prep Batch: 59712

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-10459-1	B7/10-GW02	Total/NA	Water	3535	
320-10459-2	B7/10-GW03S	Total/NA	Water	3535	
320-10459-3	QR4-GW02	Total/NA	Water	3535	
320-10459-4	B7/10-GW06S	Total/NA	Water	3535	
320-10459-5	H123-GW02	Total/NA	Water	3535	
320-10459-6	EFA-FD01-1014	Total/NA	Water	3535	
LCS 320-59712/2-A	Lab Control Sample	Total/NA	Water	3535	
MB 320-59712/1-A	Method Blank	Total/NA	Water	3535	

### Analysis Batch: 59779

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
320-10459-1	B7/10-GW02	Total/NA	Water	WS-LC-0025	59712
320-10459-2	B7/10-GW03S	Total/NA	Water	WS-LC-0025	59712
320-10459-3	QR4-GW02	Total/NA	Water	WS-LC-0025	59712
320-10459-4	B7/10-GW06S	Total/NA	Water	WS-LC-0025	59712
320-10459-5	H123-GW02	Total/NA	Water	WS-LC-0025	59712
320-10459-6	EFA-FD01-1014	Total/NA	Water	WS-LC-0025	59712
LCS 320-59712/2-A	Lab Control Sample	Total/NA	Water	WS-LC-0025	59712
MB 320-59712/1-A	Method Blank	Total/NA	Water	WS-LC-0025	59712

# Lab Chronicle

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

**Client Sample ID: B7/10-GW02**

**Lab Sample ID: 320-10459-1**

Date Collected: 11/15/14 12:05

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 20:45	JRB	TAL SAC

**Client Sample ID: B7/10-GW03S**

**Lab Sample ID: 320-10459-2**

Date Collected: 11/15/14 15:20

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 21:07	JRB	TAL SAC

**Client Sample ID: QR4-GW02**

**Lab Sample ID: 320-10459-3**

Date Collected: 11/16/14 10:20

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 21:28	JRB	TAL SAC

**Client Sample ID: B7/10-GW06S**

**Lab Sample ID: 320-10459-4**

Date Collected: 11/16/14 11:50

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 21:49	JRB	TAL SAC

**Client Sample ID: H123-GW02**

**Lab Sample ID: 320-10459-5**

Date Collected: 11/16/14 15:00

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 22:10	JRB	TAL SAC

**Client Sample ID: EFA-FD01-1014**

**Lab Sample ID: 320-10459-6**

Date Collected: 11/16/14 00:00

Matrix: Water

Date Received: 11/18/14 09:50

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared or Analyzed	Analyst	Lab
Total/NA	Prep	3535			59712	12/03/14 13:38	JER	TAL SAC
Total/NA	Analysis	WS-LC-0025		1	59779	12/04/14 22:31	JRB	TAL SAC

TestAmerica Sacramento

## Lab Chronicle

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

**Laboratory References:**

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

## Certification Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

### 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	01-31-16

## Method Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

Method	Method Description	Protocol	Laboratory
WS-LC-0025	Perfluorinated Hydrocarbons	TAL SOP	TAL SAC

**Protocol References:**

TAL SOP = TestAmerica Laboratories, Standard Operating Procedure

**Laboratory References:**

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

## Sample Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-10459-1	B7/10-GW02	Water	11/15/14 12:05	11/18/14 09:50
320-10459-2	B7/10-GW03S	Water	11/15/14 15:20	11/18/14 09:50
320-10459-3	QR4-GW02	Water	11/16/14 10:20	11/18/14 09:50
320-10459-4	B7/10-GW06S	Water	11/16/14 11:50	11/18/14 09:50
320-10459-5	H123-GW02	Water	11/16/14 15:00	11/18/14 09:50
320-10459-6	EFA-FD01-1014	Water	11/16/14 00:00	11/18/14 09:50

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
<b>LCMPFCSU_00010</b>	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
					LCM2PFTeDA_00001	0.2 mL	13C2-PFTeDA	0.4 ug/mL
					LCM4PFHPA_00001	0.2 mL	13C4-PFHpA	0.4 ug/mL
					LCM5PFPEA_00002	0.2 mL	13C5-PFPeA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C8 FOSA	1 ug/mL
					LCMPFBAA_00002	0.2 mL	13C4 PFBA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFDAA_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL
					LCMPFHxA_00004	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	18O2 PFHxS	0.3784 ug/mL
					LCMPFNA_00002	0.2 mL	13C5 PFNA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00006	0.5 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00003	0.2 mL	13C2 PFUna	0.4 ug/mL
.LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
.LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
.LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPA0112			(Purchased Reagent)		13C4-PFHpA	50 ug/mL
.LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
.LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
.LCMPFBAA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
.LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
.LCMPFDAA_00002	03/26/15	Wellington Laboratories, Lot MPFDAA0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
.LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
.LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
.LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL
.LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614			(Purchased Reagent)		13C4 PFOA	50 ug/mL
.LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
.LCMPFUdA_00003	09/29/16	Wellington Laboratories, Lot MPFUdA0911			(Purchased Reagent)		13C2 PFUna	50 ug/mL
<b>LCPFC-L1_00005</b>	01/15/15	11/06/14	MeOH/H2O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHpA	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	20 ng/mL
							13C2 PFDA	20 ng/mL
							13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							18O2 PFHxS	18.92 ng/mL
							13C5 PFNA	20 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUna	20 ng/mL
					LCPFCSP_00014	2.5 uL	Perfluorobutyric acid	0.5 ng/mL
							Perfluorobutane Sulfonate	0.442 ng/mL
							Perfluorodecanoic acid	0.5 ng/mL
							Perfluorododecanoic acid	0.5 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							PFDoS (Perfluoro-1-dodecanesulfonate)	0.484 ng/mL
							Perfluorodecane Sulfonate	0.482 ng/mL
							Perfluoroheptanoic acid	0.5 ng/mL
							Perfluoroheptane Sulfonate	0.476 ng/mL
							Perfluorohexanoic acid	0.5 ng/mL
							Perfluorohexadecanoic acid	0.5 ng/mL
							Perfluorohexane Sulfonate	0.473 ng/mL
							Perfluorononanoic acid	0.5 ng/mL
							PFNS (Perfluoro-1-nonanesulfonate)	0.48 ng/mL
							Perfluoroctanoic acid (PFOA)	0.5 ng/mL
							Perfluoroctandecanoic acid	0.5 ng/mL
							Perfluoroctane Sulfonate (PFOS)	0.478 ng/mL
							Perfluoroctane Sulfonamide	0.5 ng/mL
							Perfluoropentanoic acid	0.5 ng/mL
							PPPeS (Perfluoro-1-pentanesulfonate)	0.469 ng/mL
							Perfluorotetradecanoic acid	0.5 ng/mL
							Perfluorotridecanoic acid	0.5 ng/mL
							Perfluoroundecanoic acid	0.5 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001 LCM2PFTeDA_00001 LCM4PFHPA_00001 LCM5PFPEA_00002 LCM8FOSA_00004 LCMPFBa_00002 LCMPFDA_00003 LCMPFDaO_00002 LCMPFHxA_00004 LCMPFHxS_00002 LCMPFNA_00002 LCMPFOA_00005 LCMPFOS_00006 LCMPFUdA_00003	0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.5 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.5 mL 0.2 mL 0.2 mL 0.2 mL	13C2-PFHxDA 13C2-PFTeDA 13C4-PFHpA 13C5-PFPeA 13C8 FOSA 13C4 PFBA 13C2 PFDA 13C2 PFDoA 13C2 PFHxA 18O2 PFHxS 13C5 PFNA 13C4 PFOA 13C4 PFOS 13C2 PFUdA	0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 1 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.3784 ug/mL 0.4 ug/mL 1 ug/mL 0.956 ug/mL 0.4 ug/mL
.LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
.LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
.LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPa0112			(Purchased Reagent)		13C4-PFHpA	50 ug/mL
.LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
.LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
.LCMPFBa_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
.LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
.LCMPFDaO_00002	03/26/15	Wellington Laboratories, Lot MPFDaO0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
.LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
.LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
.LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCMPFOA_00005	07/04/19		Wellington Laboratories, Lot MPFOA0614		(Purchased Reagent)		13C4_PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19		Wellington Laboratories, Lot MPFOS0314		(Purchased Reagent)		13C4_PFOS	47.8 ug/mL
..LCMPFUdA_00003	09/29/16		Wellington Laboratories, Lot MPFUdA0911		(Purchased Reagent)		13C2_PFUuA	50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBA_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDa_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate )	0.968 ug/mL
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpA_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNAs_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nonanesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluoroctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluoroctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluoroctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOAs_00004	0.2 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PPPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBa_00002	02/22/15		Wellington Laboratories, Lot PFBA0212		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBs_00002	06/21/15		Wellington Laboratories, Lot LPFBS0612		(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
..LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDa_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDoS_00002	10/06/16		Wellington Laboratories, Lot LPFDoS1011		(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate )	48.4 ug/mL
..LCPFDS_00002	06/21/15		Wellington Laboratories, Lot LPFDS0612		(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFHpA_00002	04/18/15		Wellington Laboratories, Lot PFHpA0412		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15		Wellington Laboratories, Lot PFHxA1012		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00003	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15		Wellington Laboratories, Lot LPFHxS0312		(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFNAs_00002	06/14/15		Wellington Laboratories, Lot PFNA0612		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNS_00001	07/04/15		Wellington Laboratories, Lot LPFNS0712		(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluoroctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00003	07/13/15		Wellington Laboratories, Lot PFHODA0807		(Purchased Reagent)		Perfluoroctadecanoic acid	50 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFOS_00002	03/26/15		Wellington Laboratories, Lot LPFOS0312		(Purchased Reagent)		Perfluoroctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15		Wellington Laboratories, Lot FOSA0912M		(Purchased Reagent)		Perfluoroctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15		Wellington Laboratories, Lot LPFPeS0712		(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15		Wellington Laboratories, Lot PFTeDA0312		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15		Wellington Laboratories, Lot PFTrDA0312		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15		Wellington Laboratories, Lot PFUdA0312		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L2_00005	01/15/15	11/06/14	MeOH/H2O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHxA	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	20 ng/mL
							13C2 PFDA	20 ng/mL
							13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							18O2 PFHxS	18.92 ng/mL
							13C5 PFNA	20 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	20 ng/mL
					LCPFCSP_00014	5 uL	Perfluorobutyric acid	1 ng/mL
							Perfluorobutane Sulfonate	0.884 ng/mL
							Perfluorodecanoic acid	1 ng/mL
							Perfluorododecanoic acid	1 ng/mL
							PFDoS (Perflouro-1-dodecanesulfonate )	0.968 ng/mL
							Perfluorodecane Sulfonate	0.964 ng/mL
							Perfluoroheptanoic acid	1 ng/mL
							Perfluoroheptane Sulfonate	0.952 ng/mL
							Perfluorohehexanoic acid	1 ng/mL
							Perfluorohexadecanoic acid	1 ng/mL
							Perfluorohexane Sulfonate	0.946 ng/mL
							Perfluorononanoic acid	1 ng/mL
							PFNS (Perflouro-1-nonanesulfonate)	0.96 ng/mL
							Perfluoroctanoic acid (PFOA)	1 ng/mL
							Perfluoroctandecanoic acid	1 ng/mL
							Perfluoroctane Sulfonate	0.956 ng/mL
							Perfluoroctane Sulfonamide	1 ng/mL
							Perfluoropentanoic acid	1 ng/mL
							PFPeS (Perflouro-1-pentanesulfonate)	0.938 ng/mL
							Perfluorotetradecanoic acid	1 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	Perfluorotridecanoic acid	1 ng/mL
					LCM2PFTeDA_00001	0.2 mL	Perfluoroundecanoic acid	1 ng/mL
					LCM4PFHPA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
					LCM5PFPEA_00002	0.2 mL	13C2-PFTeDA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C4-PFHxPA	0.4 ug/mL
					LCMPFBA_00002	0.2 mL	13C5-PFPeA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C8 FOSA	1 ug/mL
					LCMPFDa_00002	0.2 mL	LCMPFBA	0.4 ug/mL
					LCMPFHxA_00004	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL
					LCMPFNA_00002	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	18O2 PFHXS	0.3784 ug/mL
					LCMPFOS_00006	0.5 mL	13C5 PFNA	0.4 ug/mL
					LCMPFUdA_00003	0.2 mL	13C4 PFOA	1 ug/mL
							13C4 PFOS	0.956 ug/mL
							13C2 PFUnA	0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPa0112			(Purchased Reagent)		13C4-PFHxPA	50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDa_00002	03/26/15	Wellington Laboratories, Lot MPFDa0A0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHXS	47.3 ug/mL
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00003	09/29/16	Wellington Laboratories, Lot MPFUdA0911			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBA_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDa_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDa_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDOS_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate )	0.968 ug/mL
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpa_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHps_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nananesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluoroctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluoroctandecanoic acid	1 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFOS_00002	0.2 mL	Perfluoroctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluoroctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PFPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
					(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBa_00002	02/22/15	Wellington Laboratories, Lot PFBA0212			(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
..LCPFBS_00002	06/21/15	Wellington Laboratories, Lot LPFBS0612			(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDA_00003	06/18/18	Wellington Laboratories, Lot PFDA0613			(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDaO_00003	01/03/18	Wellington Laboratories, Lot PFDoA0113			(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate )	48.4 ug/mL
..LCPFDoS_00002	10/06/16	Wellington Laboratories, Lot LPFDoS1011			(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFDS_00002	06/21/15	Wellington Laboratories, Lot LPFDS0612			(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpA_00002	04/18/15	Wellington Laboratories, Lot PFHpA0412			(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHpS_00004	11/21/17	Wellington Laboratories, Lot LPFHpS1112			(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxA_00002	10/29/15	Wellington Laboratories, Lot PFHxA1012			(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxDA_00003	11/28/17	Wellington Laboratories, Lot PFHxDA0707			(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFHxS_00002	03/27/15	Wellington Laboratories, Lot LPFHxS0312			(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	50 ug/mL
..LCPFNA_00002	06/14/15	Wellington Laboratories, Lot PFNA0612			(Purchased Reagent)		Perfluorononanoic acid	48 ug/mL
..LCPFNS_00001	07/04/15	Wellington Laboratories, Lot LPFNS0712			(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFOA_00004	10/11/18	Wellington Laboratories, Lot PFOA1013			(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFODA_00003	07/13/15	Wellington Laboratories, Lot PFHODA0807			(Purchased Reagent)		Perfluorooctadecanoic acid	47.8 ug/mL
..LCPFOS_00002	03/26/15	Wellington Laboratories, Lot LPFOS0312			(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	50 ug/mL
..LCPFOSA_00004	09/13/15	Wellington Laboratories, Lot FOSA0912M			(Purchased Reagent)		PFPeA (Perfluoropentanoic acid)	46.9 ug/mL
..LCPFPeA_00003	01/03/18	Wellington Laboratories, Lot PFPeA0113			(Purchased Reagent)		PFTeDA_00002	50 ug/mL
..LCPFPeS_00001	06/21/15	Wellington Laboratories, Lot LPFPeS0712			(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTeDA_00002	03/07/15	Wellington Laboratories, Lot PFTeDA0312			(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15	Wellington Laboratories, Lot PFTrDA0312			(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15	Wellington Laboratories, Lot PFUdA0312			(Purchased Reagent)		18O2 PFHxs	18.92 ng/mL
LCPFC-L3_00005	01/15/15	11/06/14	MeOH/H2O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHpA	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	20 ng/mL
							13C2 PFDA	20 ng/mL
							13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							13C5 PFNA	50 ng/mL
							13C4 PFOA	50 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCPFCSP_00014	25 uL	13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	20 ng/mL
							Perfluorobutyric acid	5 ng/mL
							Perfluorobutane Sulfonate	4.42 ng/mL
							Perfluorodecanoic acid	5 ng/mL
							Perfluorododecanoic acid	5 ng/mL
							PFDoS (Perflouro-1-dodecanesulfonate )	4.84 ng/mL
							Perfluorodecane Sulfonate	4.82 ng/mL
							Perfluoroheptanoic acid	5 ng/mL
							Perfluoroheptane Sulfonate	4.76 ng/mL
							Perfluorohexanoic acid	5 ng/mL
							Perfluorohexadecanoic acid	5 ng/mL
							Perfluorohexane Sulfonate	4.73 ng/mL
							Perfluorononanoic acid	5 ng/mL
							PFNS (Perflouro-1-nananesulfonate)	4.8 ng/mL
							Perfluoroctanoic acid (PFOA)	5 ng/mL
							Perfluoroctadecanoic acid	5 ng/mL
							Perfluoroctane Sulfonate (PFOS)	4.78 ng/mL
							Perfluoroctane Sulfonamide	5 ng/mL
							Perfluoropentanoic acid	5 ng/mL
							PPPeS (Perflouro-1-pentanesulfonate)	4.69 ng/mL
							Perfluorotetradecanoic acid	5 ng/mL
							Perfluorotridecanoic acid	5 ng/mL
							Perfluoroundecanoic acid	5 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
					LCM2PFTeDA_00001	0.2 mL	13C2-PFTeDA	0.4 ug/mL
					LCM4PFHPA_00001	0.2 mL	13C4-PFHxPA	0.4 ug/mL
					LCM5PFPEA_00002	0.2 mL	13C5-PFPeA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00002	0.2 mL	13C4 PFBA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFDa_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL
					LCMPFHxA_00004	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	18O2 PFHxS	0.3784 ug/mL
					LCMPFNA_00002	0.2 mL	13C5 PFNA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00006	0.5 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00003	0.2 mL	13C2 PFUnA	0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPa0112			(Purchased Reagent)		13C4-PFHxPA	50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCMPFBA_00002	08/13/15		Wellington Laboratories, Lot MPFBA0812		(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00003	09/13/17		Wellington Laboratories, Lot MPFDA0912		(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDaA_00002	03/26/15		Wellington Laboratories, Lot MPFDaA0312		(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00004	02/14/18		Wellington Laboratories, Lot MPFHxA0213		(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00002	03/08/15		Wellington Laboratories, Lot MPFHxS0312		(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00002	09/13/15		Wellington Laboratories, Lot MPFNA0912		(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00005	07/04/19		Wellington Laboratories, Lot MPFOA0614		(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19		Wellington Laboratories, Lot MPFOS0314		(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00003	09/29/16		Wellington Laboratories, Lot MPFUdA0911		(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBA_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDaA_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate)	0.968 ug/mL
					LCPFDs_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpA_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nonanesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluooctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluoroctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluoroctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBA_00002	02/22/15		Wellington Laboratories, Lot PFBA0212		(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBs_00002	06/21/15		Wellington Laboratories, Lot LPFBS0612		(Purchased Reagent)		Perfluorobutane Sulfonate	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 PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDoS_00002	10/06/16		Wellington Laboratories, Lot LPFDoS1011		(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate)	48.4 ug/mL
..LCPFDs_00002	06/21/15		Wellington Laboratories, Lot LPFDS0612		(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFHpA_00002	04/18/15		Wellington Laboratories, Lot PFHpA0412		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15		Wellington Laboratories, Lot PFHxA1012		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00003	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15		Wellington Laboratories, Lot LPFHxS0312		(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFNA_00002	06/14/15		Wellington Laboratories, Lot PFNA0612		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNS_00001	07/04/15		Wellington Laboratories, Lot LPFNS0712		(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluoroctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00003	07/13/15		Wellington Laboratories, Lot PFHODA0807		(Purchased Reagent)		Perfluoroctadecanoic acid	50 ug/mL
..LCPFOS_00002	03/26/15		Wellington Laboratories, Lot LPFOS0312		(Purchased Reagent)		Perfluoroctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15		Wellington Laboratories, Lot FOSA0912M		(Purchased Reagent)		Perfluoroctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15		Wellington Laboratories, Lot LPFPeS0712		(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15		Wellington Laboratories, Lot PFTeDA0312		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15		Wellington Laboratories, Lot PFTrDA0312		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15		Wellington Laboratories, Lot PFUdA0312		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L4_00005</b>	01/05/15	11/06/14	MeOH/H <sub>2</sub> O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHpA	20 ng/mL
							13C5-PFFeA	20 ng/mL
							13C8_FOSA	50 ng/mL
							13C4_PFBA	20 ng/mL
							13C2_PFDA	20 ng/mL
							13C2_PFDa	20 ng/mL
							13C2_PFHxA	20 ng/mL
							18O2_PFHxs	18.92 ng/mL
							13C5_PFNA	20 ng/mL
							13C4_PFOA	50 ng/mL
							13C4_PFOS	47.8 ng/mL
							13C2_PFUa	20 ng/mL
					LCPFCSP_00014	100 uL	Perfluorobutyric acid	20 ng/mL
							Perfluorobutane Sulfonate	17.68 ng/mL
							Perfluorodecanoic acid	20 ng/mL
							Perfluorododecanoic acid	20 ng/mL
							PFDoS (Perflouro-1-dodecanesulfonate )	19.36 ng/mL
							Perfluorodecane Sulfonate	19.28 ng/mL
							Perfluoroheptanoic acid	20 ng/mL
							Perfluoroheptane Sulfonate	19.04 ng/mL
							Perfluoroheptanoic acid	20 ng/mL
							Perfluoroheptadecanoic acid	20 ng/mL
							Perfluoroheptadecanoic acid	20 ng/mL
							Perfluorohexadecanoic acid	20 ng/mL
							Perfluorohexadecanoic acid	20 ng/mL
							Perfluorohexadecanoic acid	20 ng/mL
							Perfluorohexadecanoic acid	20 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Perfluoroctane Sulfonamide	20 ng/mL
							Perfluoropentanoic acid	20 ng/mL
							PFPeS (Perfluoro-1-pentanesulfonate)	18.76 ng/mL
							Perfluorotetradecanoic acid	20 ng/mL
							Perfluorotridecanoic acid	20 ng/mL
							Perfluoroundecanoic acid	20 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
					LCM2PFTeDA_00001	0.2 mL	13C2-PFTeDA	0.4 ug/mL
					LCM4PFHPA_00001	0.2 mL	13C4-PFHPA	0.4 ug/mL
					LCM5PFPEA_00002	0.2 mL	13C5-PFPeA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00002	0.2 mL	13C4 PFBA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFDa_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL
					LCMPFHxA_00004	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	18O2 PFHxS	0.3784 ug/mL
					LCMPFNA_00002	0.2 mL	13C5 PFNA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00006	0.5 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00003	0.2 mL	13C2 PFUnA	0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPA0112			(Purchased Reagent)		13C4-PFHPA	50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPEA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDa_00002	03/26/15	Wellington Laboratories, Lot MPFDa0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00003	09/29/16	Wellington Laboratories, Lot MPFUDa0911			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBA_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBS_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDa_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00002	0.2 mL	PFDoS (Perfluoro-1-dodecanesulfonate )	0.968 ug/mL
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpa_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHps_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nonanesulfonate)	0.96 ug/mL
					LCPFQA_00004	0.2 mL	Perfluoroctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluoroctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluoroctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluoroctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBa_00002	02/22/15	Wellington Laboratories, Lot PFBA0212			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBs_00002	06/21/15	Wellington Laboratories, Lot LPFBS0612			(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
..LCPFDA_00003	06/18/18	Wellington Laboratories, Lot PFDA0613			(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDa_00003	01/03/18	Wellington Laboratories, Lot PFDoA0113			(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDoS_00002	10/06/16	Wellington Laboratories, Lot LPFDoS1011			(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate)	48.4 ug/mL
..LCPFDS_00002	06/21/15	Wellington Laboratories, Lot LPFDS0612			(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFHpA_00002	04/18/15	Wellington Laboratories, Lot PFFhpA0412			(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17	Wellington Laboratories, Lot LFFhpS1112			(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15	Wellington Laboratories, Lot PFHxA1012			(Purchased Reagent)		Perfluorohehexanoic acid	50 ug/mL
..LCPFHxDa_00003	11/28/17	Wellington Laboratories, Lot PFHxDA0707			(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15	Wellington Laboratories, Lot LPFHxS0312			(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFNA_00002	06/14/15	Wellington Laboratories, Lot PFNA0612			(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNS_00001	07/04/15	Wellington Laboratories, Lot LPFNS0712			(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18	Wellington Laboratories, Lot PFOA1013			(Purchased Reagent)		Perfluoroctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00003	07/13/15	Wellington Laboratories, Lot PFHODA0807			(Purchased Reagent)		Perfluoroctadecanoic acid	50 ug/mL
..LCPFOS_00002	03/26/15	Wellington Laboratories, Lot LPFOS0312			(Purchased Reagent)		Perfluoroctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15	Wellington Laboratories, Lot FOSA0912M			(Purchased Reagent)		Perfluoroctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18	Wellington Laboratories, Lot PFPeA0113			(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15	Wellington Laboratories, Lot LPFPeS0712			(Purchased Reagent)		PPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15	Wellington Laboratories, Lot PFTeDA0312			(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15	Wellington Laboratories, Lot PFTrDA0312			(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15	Wellington Laboratories, Lot PFUdA0312			(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFC-L5_00006	01/15/15	11/06/14	MeOH/H <sub>2</sub> O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA 13C2-PFTeDA 13C4-PFHpA 13C5-PFPeA 13C8 FOSA 13C4 PFBA 13C2 PFDA	20 ng/mL 20 ng/mL 20 ng/mL 20 ng/mL 50 ng/mL 20 ng/mL 20 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFCSP_00014	250 uL	13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							18O2 PFHxS	18.92 ng/mL
							13C5 PFNA	20 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	20 ng/mL
							Perfluorobutyric acid	50 ng/mL
							Perfluorobutane Sulfonate	44.2 ng/mL
							Perfluorodecanoic acid	50 ng/mL
							Perfluorododecanoic acid	50 ng/mL
							PFDoS (Perfluoro-1-dodecanesulfonate)	48.4 ng/mL
							Perfluorodecane Sulfonate	48.2 ng/mL
							Perfluoroheptanoic acid	50 ng/mL
							Perfluoroheptane Sulfonate	47.6 ng/mL
							Perfluorohexanoic acid	50 ng/mL
							Perfluorohexadecanoic acid	50 ng/mL
							Perfluorohexane Sulfonate	47.3 ng/mL
							Perfluorononanoic acid	50 ng/mL
							PFNS (Perfluoro-1-nonanesulfonate)	48 ng/mL
							Perfluoroctanoic acid (PFOA)	50 ng/mL
							Perfluooctandecanoic acid	50 ng/mL
							Perfluooctane Sulfonate (PFOS)	47.8 ng/mL
							Perfluooctane Sulfonamide	50 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
							13C2-PFTeDA	0.4 ug/mL
							13C4-PFHxA	0.4 ug/mL
							13C5-PFPeA	0.4 ug/mL
							LCM8FOSA_00004	0.5 mL
							13C8 FOSA	1 ug/mL
							LCMPFBA_00002	0.2 mL
							13C4 PFBA	0.4 ug/mL
							LCMPFDA_00003	0.2 mL
							13C2 PFDA	0.4 ug/mL
							LCMPFDa_00002	0.2 mL
							13C2 PFDa	0.4 ug/mL
							LCMPFHxA_00004	0.2 mL
							13C2 PFHxA	0.4 ug/mL
							LCMPFHxS_00002	0.2 mL
							18O2 PFHxS	0.3784 ug/mL
							LCMPFNA_00002	0.2 mL
							13C5 PFNA	0.4 ug/mL
							LCMPFOA_00005	0.5 mL
							13C4 PFOA	1 ug/mL
							LCMPFOS_00006	0.5 mL
							13C4 PFOS	0.956 ug/mL
							LCMPFUdA_00003	0.2 mL
							13C2 PFUnA	0.4 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112		(Purchased Reagent)	13C2-PFHxDA		50 ug/mL	
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112		(Purchased Reagent)	13C2-PFTeDA		50 ug/mL	
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHpA0112		(Purchased Reagent)	13C4-PFHPA		50 ug/mL	
..LCM5PFPeA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312		(Purchased Reagent)	13C5-PFPeA		50 ug/mL	
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M		(Purchased Reagent)	13C8_FOSA		50 ug/mL	
..LCMPFBA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812		(Purchased Reagent)	13C4_PFBA		50 ug/mL	
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912		(Purchased Reagent)	13C2_PFDA		50 ug/mL	
..LCMPFDaA_00002	03/26/15	Wellington Laboratories, Lot MPFDaA0312		(Purchased Reagent)	13C2_PFDaA		50 ug/mL	
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213		(Purchased Reagent)	13C2_PFHxA		50 ug/mL	
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312		(Purchased Reagent)	18O2_PFHxS		47.3 ug/mL	
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912		(Purchased Reagent)	13C5_PFNA		50 ug/mL	
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614		(Purchased Reagent)	13C4_PFOA		50 ug/mL	
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314		(Purchased Reagent)	13C4_PFOS		47.8 ug/mL	
..LCMPFUdA_00003	09/29/16	Wellington Laboratories, Lot MPFUdA0911		(Purchased Reagent)	13C2_PFUdA		50 ug/mL	
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBa_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDaA_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate )	0.968 ug/mL
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpA_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNAs_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nananesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluoroctanoic acid (PFOA)	1 ug/mL
					LCPFDaA_00003	0.2 mL	Perfluoroctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluoroctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluoroctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PFPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBa_00002	02/22/15	Wellington Laboratories, Lot PFBA0212		(Purchased Reagent)	Perfluorobutyric acid		50 ug/mL	
..LCPFBs_00002	06/21/15	Wellington Laboratories, Lot LPFBS0612		(Purchased Reagent)	Perfluorobutane Sulfonate		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 PFDoA0113		(Purchased Reagent)	Perfluorododecanoic acid		50 ug/mL	
..LCPFDoS_00002	10/06/16	Wellington Laboratories, Lot LPFDoS1011		(Purchased Reagent)	PFDoS (Perflouro-1-dodecanesulfonate )		48.4 ug/mL	
..LCPFDS_00002	06/21/15	Wellington Laboratories, Lot LPFDS0612		(Purchased Reagent)	Perfluorodecane Sulfonate		48.2 ug/mL	

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFHpA_00002	04/18/15		Wellington Laboratories, Lot PFHpA0412		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15		Wellington Laboratories, Lot PFHxA1012		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00003	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15		Wellington Laboratories, Lot LPFHxS0312		(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFNAs_00002	06/14/15		Wellington Laboratories, Lot PFNA0612		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNS_00001	07/04/15		Wellington Laboratories, Lot LPFNS0712		(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluoroctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00003	07/13/15		Wellington Laboratories, Lot PFHODA0807		(Purchased Reagent)		Perfluoroctandecanoic acid	50 ug/mL
..LCPFOS_00002	03/26/15		Wellington Laboratories, Lot LPFOS0312		(Purchased Reagent)		Perfluoroctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15		Wellington Laboratories, Lot FOSA0912M		(Purchased Reagent)		Perfluoroctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15		Wellington Laboratories, Lot LPFPeS0712		(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15		Wellington Laboratories, Lot PFTeDA0312		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15		Wellington Laboratories, Lot PFTrDA0312		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15		Wellington Laboratories, Lot PFUdA0312		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L6_00005</b>	01/15/15	11/06/14	MeOH/H2O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHxA	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8_FOSA	50 ng/mL
							13C4_PFFBA	20 ng/mL
							13C2_PFDA	20 ng/mL
							13C2_PFDa	20 ng/mL
							13C2_PFFhxA	20 ng/mL
							18O2_PFHxS	18.92 ng/mL
							13C5_PFNAs	20 ng/mL
							13C4_PFOA	50 ng/mL
							13C4_PFOS	47.8 ng/mL
							13C2_PFUuNa	20 ng/mL
					LCPFCSP_00014	1000 uL	Perfluorobutyric acid	200 ng/mL
							Perfluorobutane Sulfonate	176.8 ng/mL
							Perfluorodecanoic acid	200 ng/mL
							Perfluorododecanoic acid	200 ng/mL
							PFDoS (Perflouro-1-dodecanesulfonate )	193.6 ng/mL
							Perfluorodecane Sulfonate	192.8 ng/mL
							Perfluoroheptanoic acid	200 ng/mL
							Perfluoroheptane Sulfonate	190.4 ng/mL
							Perfluorohexanoic acid	200 ng/mL
							Perfluorohexadecanoic acid	200 ng/mL
							Perfluorohexane Sulfonate	189.2 ng/mL
							Perfluorononanoic acid	200 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							PFNS (Perfluoro-1-nonanesulfonate)	192 ng/mL
							Perfluorooctanoic acid (PFOA)	200 ng/mL
							Perfluorooctadecanoic acid	200 ng/mL
							Perfluorooctane Sulfonate (PFOS)	191.2 ng/mL
							Perfluorooctane Sulfonamide	200 ng/mL
							Perfluoropentanoic acid	200 ng/mL
							PFPeS (Perfluoro-1-pentanesulfonate)	187.6 ng/mL
							Perfluorotetradecanoic acid	200 ng/mL
							Perfluorotridecanoic acid	200 ng/mL
							Perfluoroundecanoic acid	200 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001 LCM2PFTeDA_00001 LCM4PFHPA_00001 LCM5PFPEA_00002 LCM8FOSA_00004 LCMPFBa_00002 LCMPFDA_00003 LCMPFDaO_00002 LCMPFHxA_00004 LCMPFHxS_00002 LCMPFNA_00002 LCMPFOA_00005 LCMPFOS_00006 LCMPFUdA_00003	0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.5 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.5 mL 0.5 mL 0.2 mL	13C2-PFHxDA 13C2-PFTeDA 13C4-PFHpA 13C5-PFPeA 13C8 FOSA 13C4 PFBA 13C2 PFDA 13C2 PFDoA 13C2 PFHxA 18O2 PFHxS 13C5 PFNA 13C4 PFOA 13C4 PFOS 13C2 PFUna	0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 1 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 0.4 ug/mL 1 ug/mL 0.956 ug/mL 0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112	(Purchased Reagent)		13C2-PFHxDA			50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112	(Purchased Reagent)		13C2-PFTeDA			50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPA0112	(Purchased Reagent)		13C4-PFHpA			50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312	(Purchased Reagent)		13C5-PFPeA			50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M	(Purchased Reagent)		13C8 FOSA			50 ug/mL
..LCMPFBa_00002	08/13/15	Wellington Laboratories, Lot MPFBa0812	(Purchased Reagent)		13C4 PFBA			50 ug/mL
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912	(Purchased Reagent)		13C2 PFDA			50 ug/mL
..LCMPFDaO_00002	03/26/15	Wellington Laboratories, Lot MPFDoA0312	(Purchased Reagent)		13C2 PFDoA			50 ug/mL
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213	(Purchased Reagent)		13C2 PFHxA			50 ug/mL
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312	(Purchased Reagent)		18O2 PFHxS			47.3 ug/mL
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912	(Purchased Reagent)		13C5 PFNA			50 ug/mL
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614	(Purchased Reagent)		13C4 PFOA			50 ug/mL
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314	(Purchased Reagent)		13C4 PFOS			47.8 ug/mL
..LCMPFUdA_00003	09/29/16	Wellington Laboratories, Lot MPFUdA0911	(Purchased Reagent)		13C2 PFUna			50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBa_00002 LCPFBs_00002 LCPFDA_00003 LCPFDaO_00003 LCPFDs_00002	0.2 mL 0.2 mL 0.2 mL 0.2 mL 0.2 mL	Perfluorobutyric acid Perfluorobutane Sulfonate Perfluorodecanoic acid Perfluorododecanoic acid PFDoS (Perfluoro-1-dodecanesulfonate )	1 ug/mL 0.884 ug/mL 1 ug/mL 1 ug/mL 0.968 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpA_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNNS_00001	0.2 mL	PFNS (Perflouro-1-nonanesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluorooctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PFPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBa_00002	02/22/15	Wellington Laboratories, Lot PFBA0212			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL
..LCPFBs_00002	06/21/15	Wellington Laboratories, Lot LPFBS0612			(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
..LCPFDA_00003	06/18/18	Wellington Laboratories, Lot PFDA0613			(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDaO_00003	01/03/18	Wellington Laboratories, Lot PFDaO0113			(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDoS_00002	10/06/16	Wellington Laboratories, Lot LPFDoS1011			(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate)	48.4 ug/mL
..LCPFDS_00002	06/21/15	Wellington Laboratories, Lot LPFDS0612			(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFHpA_00002	04/18/15	Wellington Laboratories, Lot PFFHpA0412			(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17	Wellington Laboratories, Lot LPFHpS1112			(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15	Wellington Laboratories, Lot PFFHxA1012			(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDA_00003	11/28/17	Wellington Laboratories, Lot PFFHxDA0707			(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15	Wellington Laboratories, Lot LPFHxS0312			(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFNNA_00002	06/14/15	Wellington Laboratories, Lot FFNA0612			(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNNS_00001	07/04/15	Wellington Laboratories, Lot LPFNS0712			(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18	Wellington Laboratories, Lot PFOA1013			(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFODA_00003	07/13/15	Wellington Laboratories, Lot PFHODA0807			(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS_00002	03/26/15	Wellington Laboratories, Lot LPFOS0312			(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15	Wellington Laboratories, Lot FOSA0912M			(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18	Wellington Laboratories, Lot PFPeA0113			(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15	Wellington Laboratories, Lot LFFPeS0712			(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15	Wellington Laboratories, Lot PFTeDA0312			(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15	Wellington Laboratories, Lot PFTTrDA0312			(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15	Wellington Laboratories, Lot PFUdA0312			(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
<b>LCPFC-L7_00005</b>	01/15/15	11/06/14 MeOH/H <sub>2</sub> O, Lot 042259		5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDa	20 ng/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							13C2-PFTeDA	20 ng/mL
							13C4-PFH <sub>2</sub> A	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	20 ng/mL
							13C2 PFDA	20 ng/mL
							13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							18O2 PFHxS	18.92 ng/mL
							13C5 PFNA	20 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFU <sub>n</sub> A	20 ng/mL
					LCPFCSP_00014	2500 uL	Perfluorobutyric acid	500 ng/mL
							Perfluorobutane Sulfonate	442 ng/mL
							Perfluorodecanoic acid	500 ng/mL
							Perfluorododecanoic acid	500 ng/mL
							PFDoS (Perflouro-1-dodecanesulfonate )	484 ng/mL
							Perfluorodecane Sulfonate	482 ng/mL
							Perfluoroheptanoic acid	500 ng/mL
							Perfluoroheptane Sulfonate	476 ng/mL
							Perfluorohexanoic acid	500 ng/mL
							Perfluorohexadecanoic acid	500 ng/mL
							Perfluorohexane Sulfonate	473 ng/mL
							Perfluorononanoic acid	500 ng/mL
							PFNS (Perflouro-1-nonanesulfonate)	480 ng/mL
							Perfluoroctanoic acid (PFOA)	500 ng/mL
							Perfluoroctadecanoic acid	500 ng/mL
							Perfluoroctane Sulfonate (PFOS)	478 ng/mL
							Perfluoroctane Sulfonamide	500 ng/mL
							Perfluoropentanoic acid	500 ng/mL
							PFPeS (Perflouro-1-pentanesulfonate)	469 ng/mL
							Perfluorotetradecanoic acid	500 ng/mL
							Perfluorotridecanoic acid	500 ng/mL
							Perfluoroundecanoic acid	500 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	13C2-PFHxDA	0.4 ug/mL
					LCM2PFTeDA_00001	0.2 mL	13C2-PFTeDA	0.4 ug/mL
					LCM4PFHPA_00001	0.2 mL	13C4-PFH <sub>2</sub> A	0.4 ug/mL
					LCM5PFPEA_00002	0.2 mL	13C5-PFPeA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C8 FOSA	1 ug/mL
					LCMPFBA_00002	0.2 mL	13C4 PFBA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFDoA_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCMPFHxA_00004	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	18O2 PFHxS	0.3784 ug/mL
					LCMPFNA_00002	0.2 mL	13C5 PFNA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00006	0.5 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUdA_00003	0.2 mL	13C2 PFUnA	0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHpA0112			(Purchased Reagent)		13C4-PFHPA	50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDaA_00002	03/26/15	Wellington Laboratories, Lot MPFDaA0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUdA_00003	09/29/16	Wellington Laboratories, Lot MPFUdA0911			(Purchased Reagent)		13C2 PFUnA	50 ug/mL
.LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBa_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDaA_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDoS_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate )	0.968 ug/mL
					LCPFDs_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpa_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHps_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nonanesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluorooctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL
					LCPFPeS_00001	0.2 mL	PFPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL
..LCPFBa_00002	02/22/15	Wellington Laboratories, Lot PFBA0212			(Purchased Reagent)		Perfluorobutyric acid	50 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..LCPFBS_00002	06/21/15		Wellington Laboratories, Lot LPFBS0612		(Purchased Reagent)		Perfluorobutane Sulfonate	44.2 ug/mL
..LCPFDA_00003	06/18/18		Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid	50 ug/mL
..LCPFDaO_00003	01/03/18		Wellington Laboratories, Lot PFDoA0113		(Purchased Reagent)		Perfluorododecanoic acid	50 ug/mL
..LCPFDoS_00002	10/06/16		Wellington Laboratories, Lot LPFDoS1011		(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate )	48.4 ug/mL
..LCPFDS_00002	06/21/15		Wellington Laboratories, Lot LPFDS0612		(Purchased Reagent)		Perfluorodecane Sulfonate	48.2 ug/mL
..LCPFHpA_00002	04/18/15		Wellington Laboratories, Lot PFHpA0412		(Purchased Reagent)		Perfluoroheptanoic acid	50 ug/mL
..LCPFHpS_00004	11/21/17		Wellington Laboratories, Lot LPFHpS1112		(Purchased Reagent)		Perfluoroheptane Sulfonate	47.6 ug/mL
..LCPFHxA_00002	10/29/15		Wellington Laboratories, Lot PFHxA1012		(Purchased Reagent)		Perfluorohexanoic acid	50 ug/mL
..LCPFHxDa_00003	11/28/17		Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid	50 ug/mL
..LCPFHxS_00002	03/27/15		Wellington Laboratories, Lot LPFHxS0312		(Purchased Reagent)		Perfluorohexane Sulfonate	47.3 ug/mL
..LCPFNA_00002	06/14/15		Wellington Laboratories, Lot PFNA0612		(Purchased Reagent)		Perfluorononanoic acid	50 ug/mL
..LCPFNS_00001	07/04/15		Wellington Laboratories, Lot LPFNS0712		(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)	48 ug/mL
..LCPFOA_00004	10/11/18		Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)	50 ug/mL
..LCPFDaO_00003	07/13/15		Wellington Laboratories, Lot PFHODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid	50 ug/mL
..LCPFOS_00002	03/26/15		Wellington Laboratories, Lot LPFOS0312		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	47.8 ug/mL
..LCPFOSA_00004	09/13/15		Wellington Laboratories, Lot FOSA0912M		(Purchased Reagent)		Perfluorooctane Sulfonamide	50 ug/mL
..LCPFPeA_00003	01/03/18		Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid	50 ug/mL
..LCPFPeS_00001	06/21/15		Wellington Laboratories, Lot LPFPeS0712		(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)	46.9 ug/mL
..LCPFTeDA_00002	03/07/15		Wellington Laboratories, Lot PFTeDA0312		(Purchased Reagent)		Perfluorotetradecanoic acid	50 ug/mL
..LCPFTrDA_00002	03/26/15		Wellington Laboratories, Lot PFTrDA0312		(Purchased Reagent)		Perfluorotridecanoic acid	50 ug/mL
..LCPFUdA_00002	03/19/15		Wellington Laboratories, Lot PFUdA0312		(Purchased Reagent)		Perfluoroundecanoic acid	50 ug/mL
LCPFCIC_00005	01/20/15	11/06/14	MeOH/H2O, Lot 042259	5 mL	LCMPFCSU_00010	250 uL	13C2-PFHxDA	20 ng/mL
							13C2-PFTeDA	20 ng/mL
							13C4-PFHpA	20 ng/mL
							13C5-PFPeA	20 ng/mL
							13C8 FOSA	50 ng/mL
							13C4 PFBA	20 ng/mL
							13C2 PFDA	20 ng/mL
							13C2 PFDoA	20 ng/mL
							13C2 PFHxA	20 ng/mL
							18O2 PFHxS	18.92 ng/mL
					LCPFACMXB_00002	125 uL	13C5 PFNA	20 ng/mL
							13C4 PFOA	50 ng/mL
							13C4 PFOS	47.8 ng/mL
							13C2 PFUnA	20 ng/mL
							Perfluorooctane Sulfonate (PFOS)	47.75 ng/mL
.LCMPFCSU_00010	01/20/15	09/09/14	Methanol, Lot 067374	25 mL	LCM2PFHxDA_00001	0.2 mL	Perfluorooctanoic acid (PFOA)	50 ng/mL
							13C2-PFHxDA	0.4 ug/mL
							13C2-PFTeDA	0.4 ug/mL
							13C4-PFHpA	0.4 ug/mL
							13C5-PFPeA	0.4 ug/mL
					LCM8FOSA_00004	0.5 mL	13C8 FOSA	1 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					LCMPFBA_00002	0.2 mL	13C4 PFBA	0.4 ug/mL
					LCMPFDA_00003	0.2 mL	13C2 PFDA	0.4 ug/mL
					LCMPFDa_00002	0.2 mL	13C2 PFDoA	0.4 ug/mL
					LCMPFHxA_00004	0.2 mL	13C2 PFHxA	0.4 ug/mL
					LCMPFHxS_00002	0.2 mL	18O2 PFHxS	0.3784 ug/mL
					LCMPFNA_00002	0.2 mL	13C5 PFNA	0.4 ug/mL
					LCMPFOA_00005	0.5 mL	13C4 PFOA	1 ug/mL
					LCMPFOS_00006	0.5 mL	13C4 PFOS	0.956 ug/mL
					LCMPFUDa_00003	0.2 mL	13C2 PFUna	0.4 ug/mL
..LCM2PFHxDA_00001	11/29/15	Wellington Laboratories, Lot M2PFHxDA1112			(Purchased Reagent)		13C2-PFHxDA	50 ug/mL
..LCM2PFTeDA_00001	11/29/15	Wellington Laboratories, Lot M2PFTeDA1112			(Purchased Reagent)		13C2-PFTeDA	50 ug/mL
..LCM4PFHPA_00001	01/20/15	Wellington Laboratories, Lot M4PFHPA0112			(Purchased Reagent)		13C4-PFHpA	50 ug/mL
..LCM5PFPEA_00002	03/07/15	Wellington Laboratories, Lot M5PFPeA0312			(Purchased Reagent)		13C5-PFPeA	50 ug/mL
..LCM8FOSA_00004	10/11/18	Wellington Laboratories, Lot M8FOSA1013M			(Purchased Reagent)		13C8 FOSA	50 ug/mL
..LCMPFBA_00002	08/13/15	Wellington Laboratories, Lot MPFBA0812			(Purchased Reagent)		13C4 PFBA	50 ug/mL
..LCMPFDA_00003	09/13/17	Wellington Laboratories, Lot MPFDA0912			(Purchased Reagent)		13C2 PFDA	50 ug/mL
..LCMPFDa_00002	03/26/15	Wellington Laboratories, Lot MPFDa0A0312			(Purchased Reagent)		13C2 PFDoA	50 ug/mL
..LCMPFHxA_00004	02/14/18	Wellington Laboratories, Lot MPFHxA0213			(Purchased Reagent)		13C2 PFHxA	50 ug/mL
..LCMPFHxS_00002	03/08/15	Wellington Laboratories, Lot MPFHxS0312			(Purchased Reagent)		18O2 PFHxS	47.3 ug/mL
..LCMPFNA_00002	09/13/15	Wellington Laboratories, Lot MPFNA0912			(Purchased Reagent)		13C5 PFNA	50 ug/mL
..LCMPFOA_00005	07/04/19	Wellington Laboratories, Lot MPFOA0614			(Purchased Reagent)		13C4 PFOA	50 ug/mL
..LCMPFOS_00006	03/10/19	Wellington Laboratories, Lot MPFOS0314			(Purchased Reagent)		13C4 PFOS	47.8 ug/mL
..LCMPFUDa_00003	09/29/16	Wellington Laboratories, Lot MPFUDa0911			(Purchased Reagent)		13C2 PFUna	50 ug/mL
.LCPFACMXB_00002	03/08/15	Wellington Laboratories, Lot PFACMXB0312			(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)	1.91 ug/mL
							Perfluorooctanoic acid (PFOA)	2 ug/mL
LCPFCSP_00014	01/15/15	07/15/14	Methanol, Lot 042259	10 mL	LCPFBa_00002	0.2 mL	Perfluorobutyric acid	1 ug/mL
					LCPFBs_00002	0.2 mL	Perfluorobutane Sulfonate	0.884 ug/mL
					LCPFDA_00003	0.2 mL	Perfluorodecanoic acid	1 ug/mL
					LCPFDa_00003	0.2 mL	Perfluorododecanoic acid	1 ug/mL
					LCPFDos_00002	0.2 mL	PFDoS (Perflouro-1-dodecanesulfonate)	0.968 ug/mL
					LCPFDS_00002	0.2 mL	Perfluorodecane Sulfonate	0.964 ug/mL
					LCPFHpA_00002	0.2 mL	Perfluoroheptanoic acid	1 ug/mL
					LCPFHpS_00004	0.2 mL	Perfluoroheptane Sulfonate	0.952 ug/mL
					LCPFHxA_00002	0.2 mL	Perfluorohexanoic acid	1 ug/mL
					LCPFHxDA_00003	0.2 mL	Perfluorohexadecanoic acid	1 ug/mL
					LCPFHxS_00002	0.2 mL	Perfluorohexane Sulfonate	0.946 ug/mL
					LCPFNA_00002	0.2 mL	Perfluorononanoic acid	1 ug/mL
					LCPFNS_00001	0.2 mL	PFNS (Perflouro-1-nananesulfonate)	0.96 ug/mL
					LCPFOA_00004	0.2 mL	Perfluorooctanoic acid (PFOA)	1 ug/mL
					LCPFODA_00003	0.2 mL	Perfluorooctadecanoic acid	1 ug/mL
					LCPFOS_00002	0.2 mL	Perfluorooctane Sulfonate (PFOS)	0.956 ug/mL
					LCPFOSA_00004	0.2 mL	Perfluorooctane Sulfonamide	1 ug/mL
					LCPFPeA_00003	0.2 mL	Perfluoropentanoic acid	1 ug/mL

## REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration	
					Reagent ID	Volume Added			
					LCPFPeS_00001	0.2 mL	PFPeS (Perflouro-1-pentanesulfonate)	0.938 ug/mL	
					LCPFTeDA_00002	0.2 mL	Perfluorotetradecanoic acid	1 ug/mL	
					LCPFTrDA_00002	0.2 mL	Perfluorotridecanoic acid	1 ug/mL	
					LCPFUdA_00002	0.2 mL	Perfluoroundecanoic acid	1 ug/mL	
.LCPFBA_00002	02/22/15	Wellington Laboratories, Lot PFBA0212		(Purchased Reagent)		Perfluorobutyric acid		50 ug/mL	
.LCPFBS_00002	06/21/15	Wellington Laboratories, Lot LPFBS0612		(Purchased Reagent)		Perfluorobutane Sulfonate		44.2 ug/mL	
.LCPFDA_00003	06/18/18	Wellington Laboratories, Lot PFDA0613		(Purchased Reagent)		Perfluorodecanoic acid		50 ug/mL	
.LCPFDaO_00003	01/03/18	Wellington Laboratories, Lot PFDaO0113		(Purchased Reagent)		Perfluorododecanoic acid		50 ug/mL	
.LCPFDoS_00002	10/06/16	Wellington Laboratories, Lot LPFDoS1011		(Purchased Reagent)		PFDoS (Perflouro-1-dodecanesulfonate )		48.4 ug/mL	
.LCPFDS_00002	06/21/15	Wellington Laboratories, Lot LPFDS0612		(Purchased Reagent)		Perfluorodecane Sulfonate		48.2 ug/mL	
.LCPFHpA_00002	04/18/15	Wellington Laboratories, Lot PFHpA0412		(Purchased Reagent)		Perfluoroheptanoic acid		50 ug/mL	
.LCPFHpS_00004	11/21/17	Wellington Laboratories, Lot LFFHpS1112		(Purchased Reagent)		Perfluoroheptane Sulfonate		47.6 ug/mL	
.LCPFHxA_00002	10/29/15	Wellington Laboratories, Lot PFHxA1012		(Purchased Reagent)		Perfluorohexanoic acid		50 ug/mL	
.LCPFHxDA_00003	11/28/17	Wellington Laboratories, Lot PFHxDA0707		(Purchased Reagent)		Perfluorohexadecanoic acid		50 ug/mL	
.LCPFHxS_00002	03/27/15	Wellington Laboratories, Lot LPFHxS0312		(Purchased Reagent)		Perfluorohexane Sulfonate		47.3 ug/mL	
.LCPFNAs_00002	06/14/15	Wellington Laboratories, Lot PFNA0612		(Purchased Reagent)		Perfluorononanoic acid		50 ug/mL	
.LCPFNs_00001	07/04/15	Wellington Laboratories, Lot LPFNS0712		(Purchased Reagent)		PFNS (Perflouro-1-nonanesulfonate)		48 ug/mL	
.LCPFOA_00004	10/11/18	Wellington Laboratories, Lot PFOA1013		(Purchased Reagent)		Perfluorooctanoic acid (PFOA)		50 ug/mL	
.LCPFODA_00003	07/13/15	Wellington Laboratories, Lot PFHODA0807		(Purchased Reagent)		Perfluorooctadecanoic acid		50 ug/mL	
.LCPFOS_00002	03/26/15	Wellington Laboratories, Lot LPFOS0312		(Purchased Reagent)		Perfluorooctane Sulfonate (PFOS)		47.8 ug/mL	
.LCPFOSA_00004	09/13/15	Wellington Laboratories, Lot FOSA0912M		(Purchased Reagent)		Perfluorooctane Sulfonamide		50 ug/mL	
.LCPFPeA_00003	01/03/18	Wellington Laboratories, Lot PFPeA0113		(Purchased Reagent)		Perfluoropentanoic acid		50 ug/mL	
.LCPFPeS_00001	06/21/15	Wellington Laboratories, Lot LPFPeS0712		(Purchased Reagent)		PFPeS (Perflouro-1-pentanesulfonate)		46.9 ug/mL	
.LCPFTeDA_00002	03/07/15	Wellington Laboratories, Lot PFTeDA0312		(Purchased Reagent)		Perfluorotetradecanoic acid		50 ug/mL	
.LCPFTrDA_00002	03/26/15	Wellington Laboratories, Lot PFTrDA0312		(Purchased Reagent)		Perfluorotridecanoic acid		50 ug/mL	
.LCPFUdA_00002	03/19/15	Wellington Laboratories, Lot PFUdA0312		(Purchased Reagent)		Perfluoroundecanoic acid		50 ug/mL	

Reagent

---

**LCM2PFHxDA\_00001**



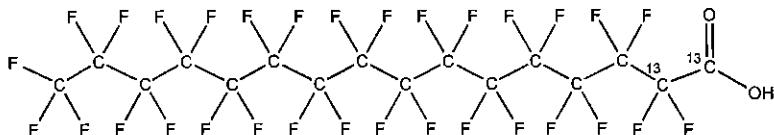
# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

12-19-12 New lot.

PRODUCT CODE: M2PFHxDA      LOT NUMBER: M2PFHxDA1112  
COMPOUND: Perfluoro-n-[1,2-<sup>13</sup>C]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  
LAST TESTED: (mm/dd/yyyy) 11/29/2012      (1,2-<sup>13</sup>C<sub>2</sub>)  
EXPIRY DATE: (mm/dd/yyyy) 11/29/2015  
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:

B.G. Chittim

Date: 12/05/2012

(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.

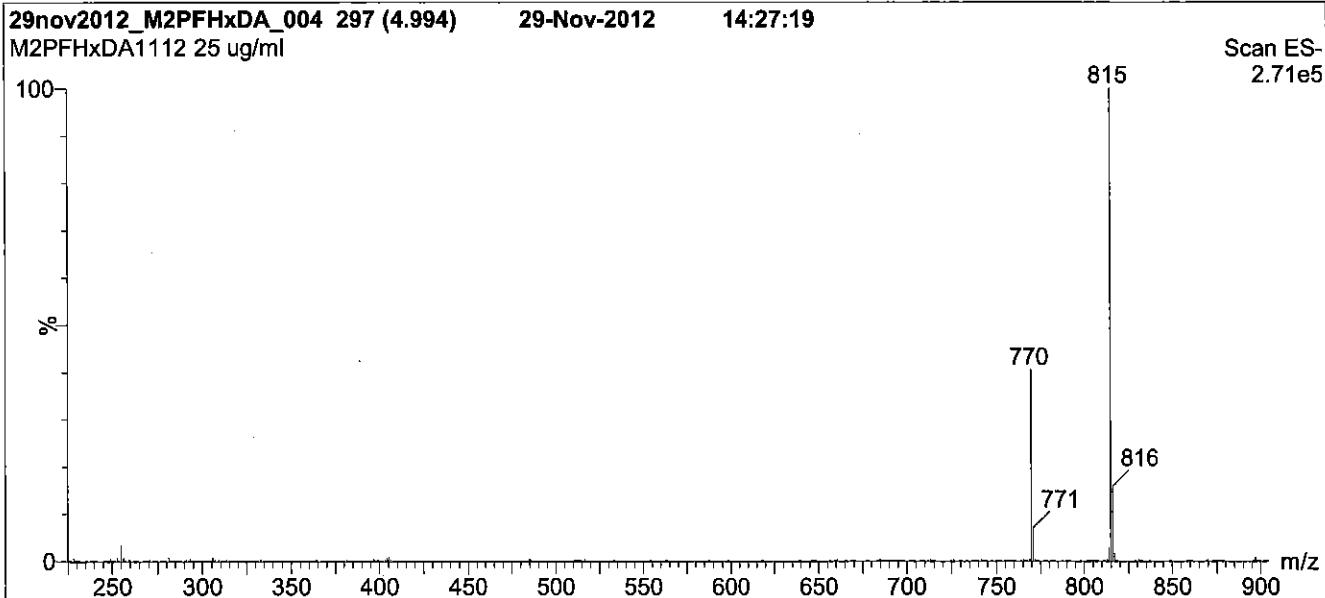
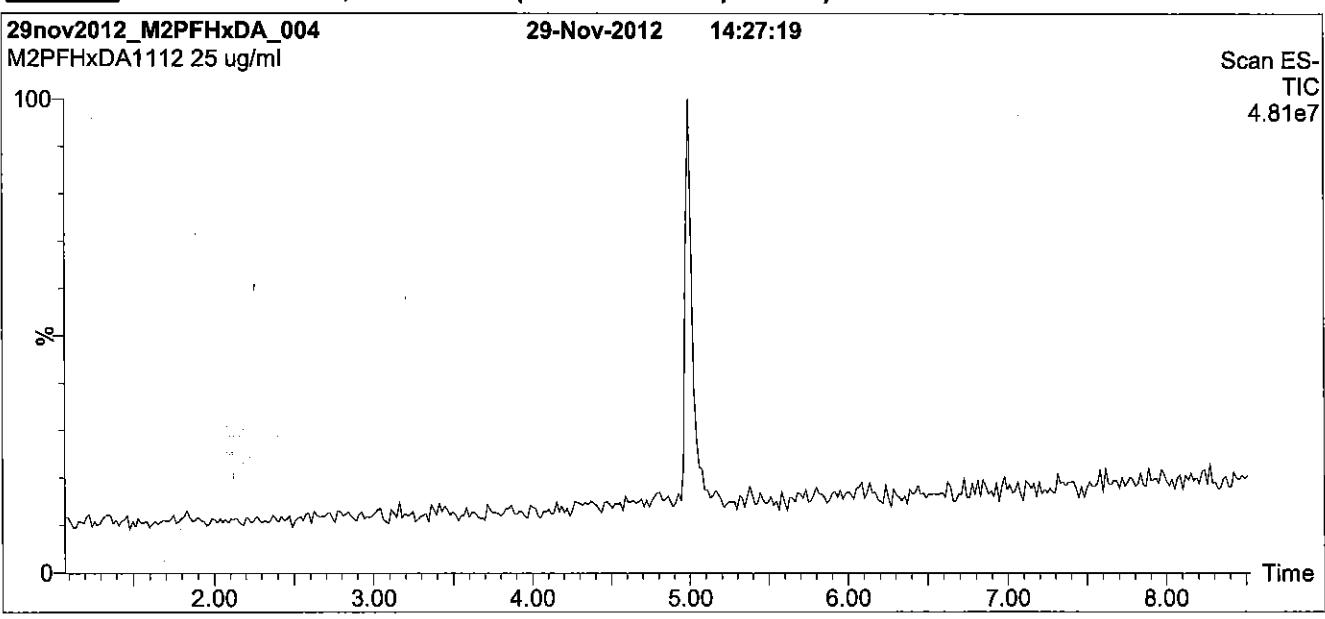
#### **QUALITY MANAGEMENT:**

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 ACCLASS (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 Acuity Ultra Performance LC  
**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity 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

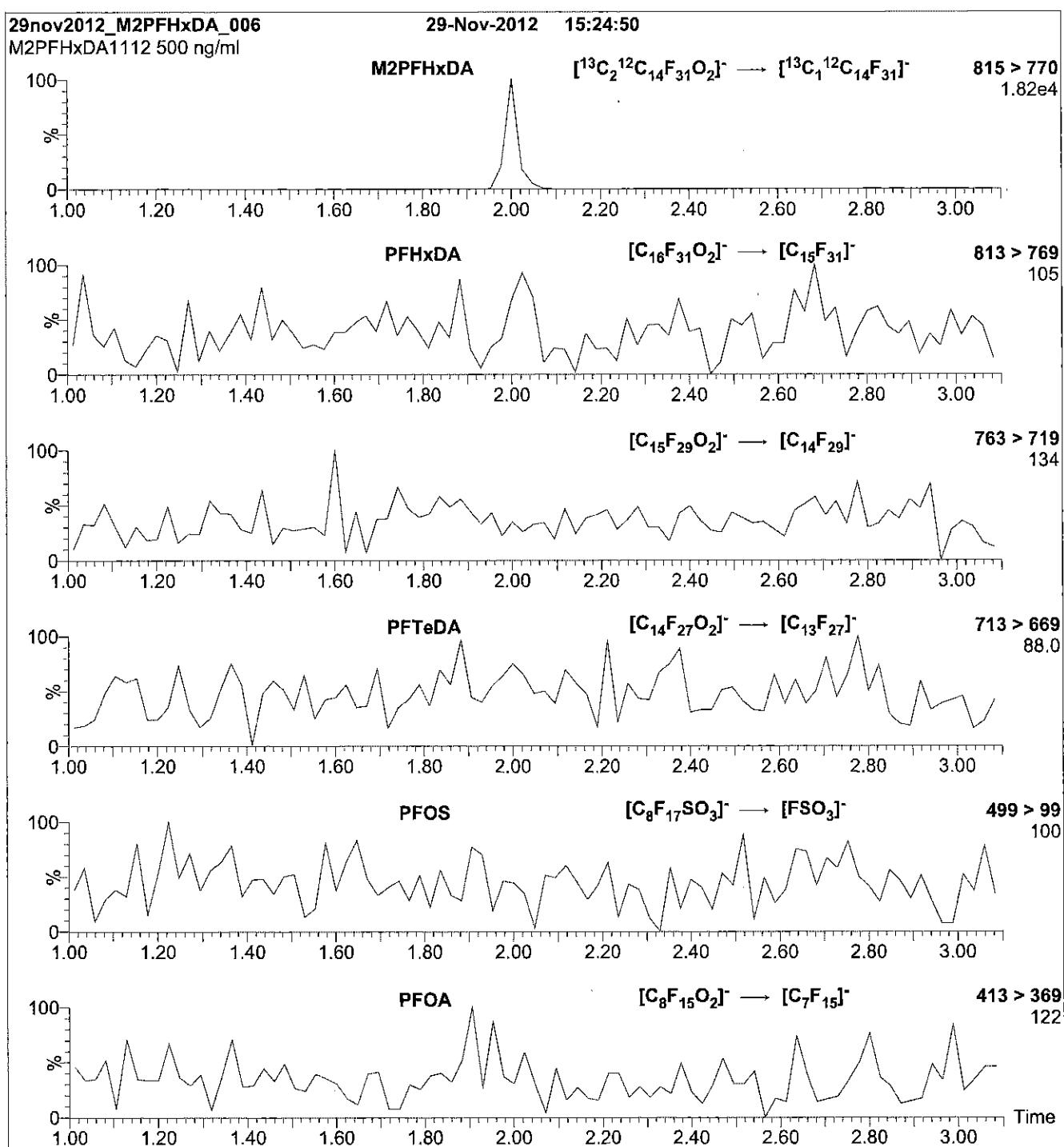
Flow: 300 µl/min

**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)

**MS Parameters**

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

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  $\mu\text{l}/\text{min}$

Reagent

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**LCM2PFTeDA\_00001**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

12-29-12 New pdf.

PRODUCT CODE:

M2PFTeDA

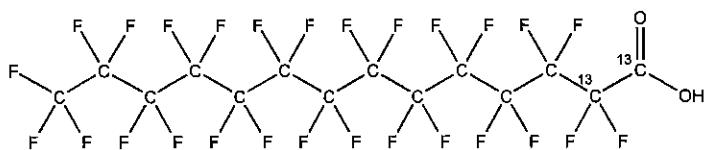
LOT NUMBER: M2PFTeDA1112

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

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: ≥99% <sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

11/29/2012

(1,2-<sup>13</sup>C<sub>2</sub>)

EXPIRY DATE: (mm/dd/yyyy)

11/29/2015

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

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

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

ADDITIONAL INFORMATION:

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

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

Certified By:

B.G. Chittim

Date: 12/04/2012

(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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**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:

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

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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.

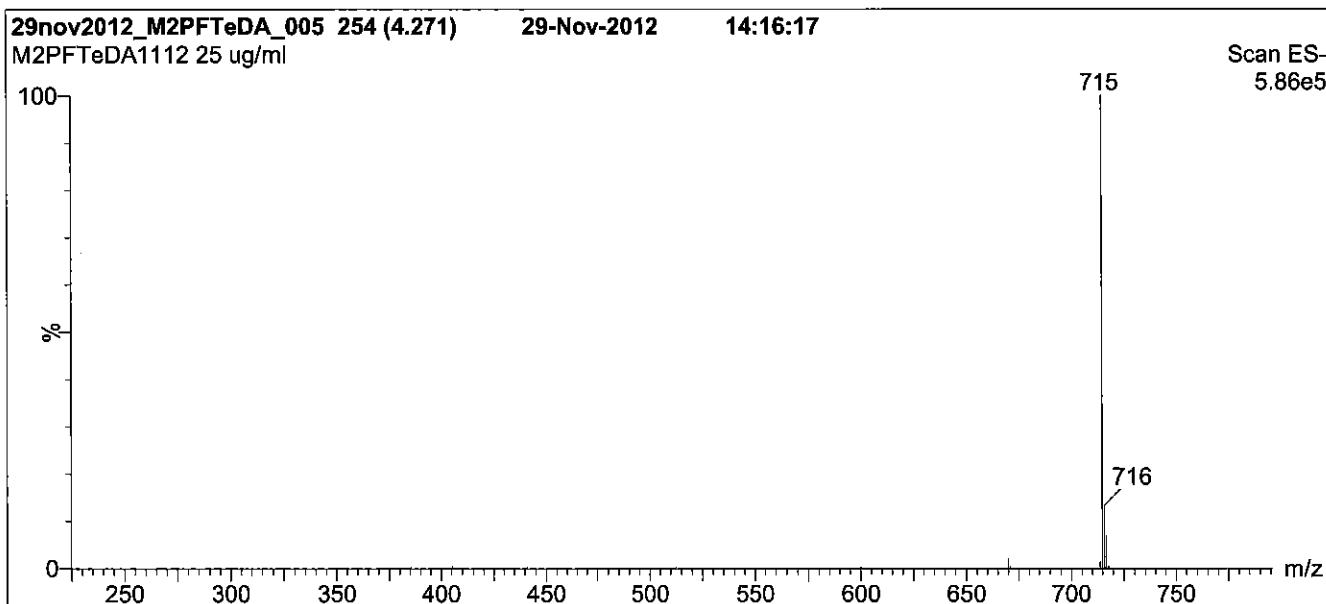
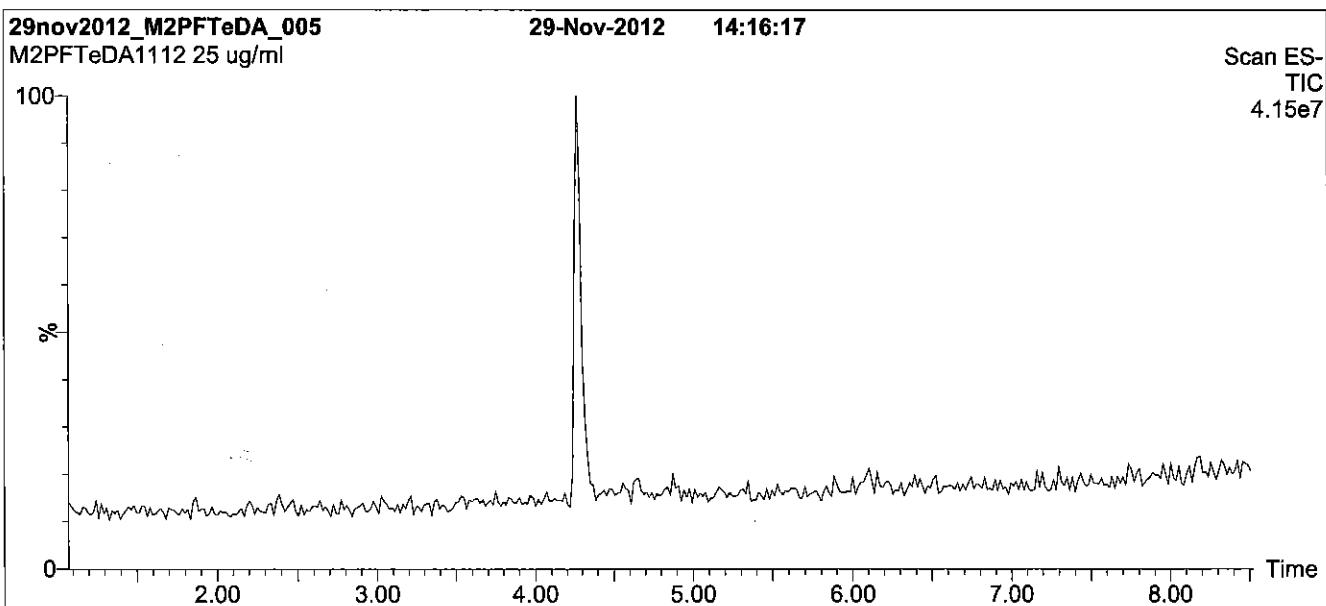
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**Figure 1:** M2PFTeDA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity 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

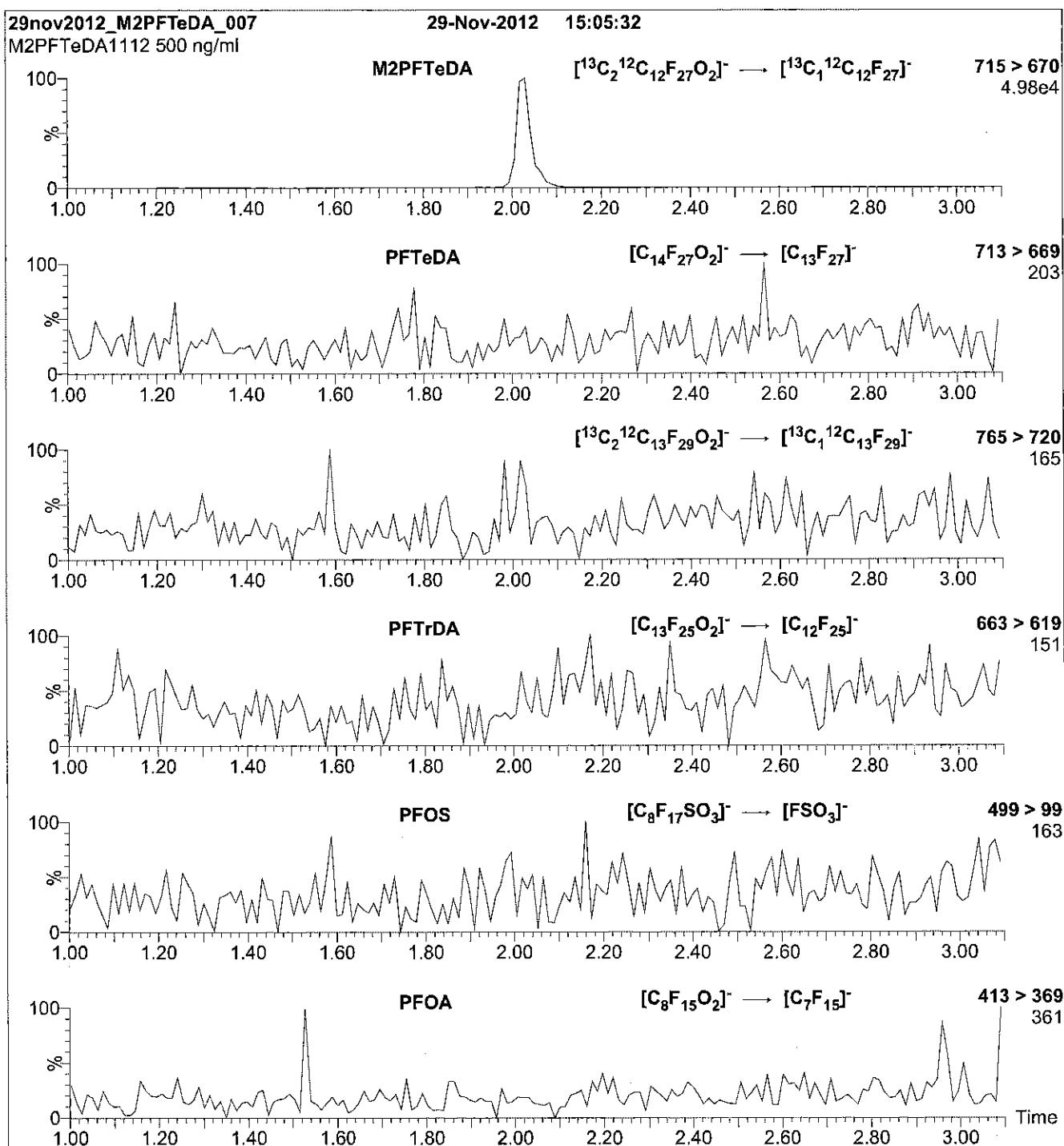
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (225 - 1200 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 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)

**MS Parameters**

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

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  $\mu\text{l}/\text{min}$

Reagent

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**LCM4PFHPA\_00001**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R 12-14-12

LCM4PFHPA-00001

PRODUCT CODE:

M4PFHpA

LOT NUMBER: M4PFHpA0112

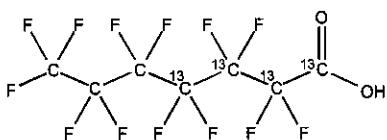
COMPOUND:

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

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: ≥99%<sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

01/20/2012

EXPIRY DATE: (mm/dd/yyyy)

01/20/2015

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: 01/23/2012

(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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$x_1, x_2, \dots, x_n$  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.

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**EXPIRY DATE / PERIOD OF VALIDITY:**

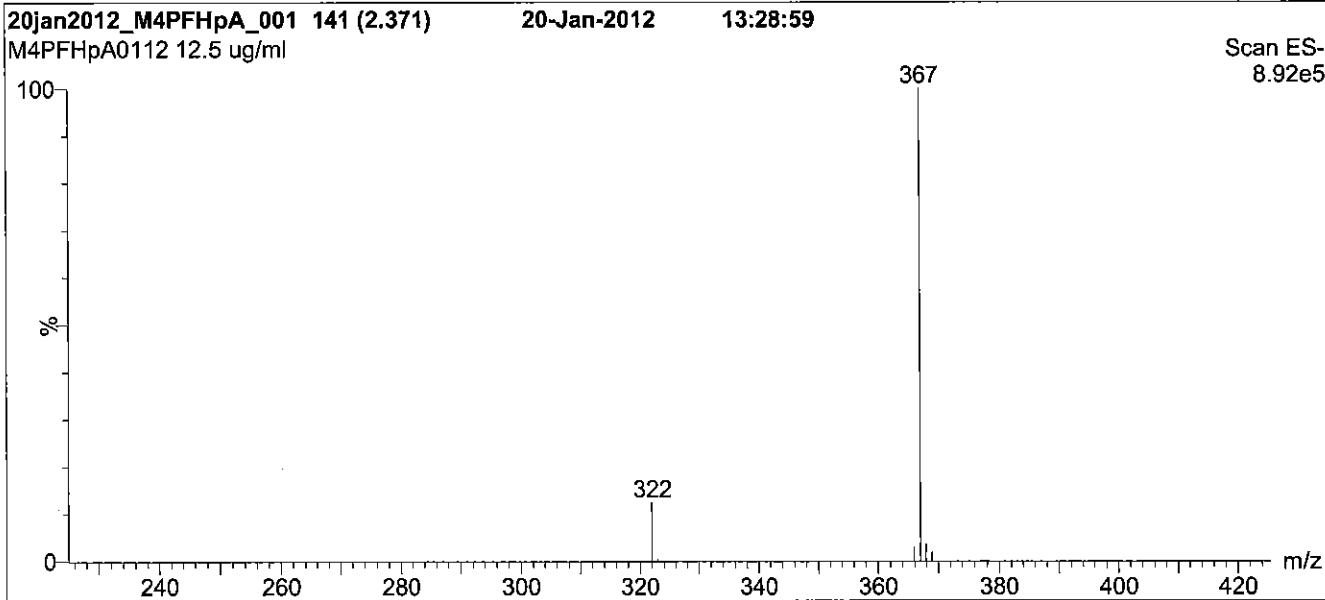
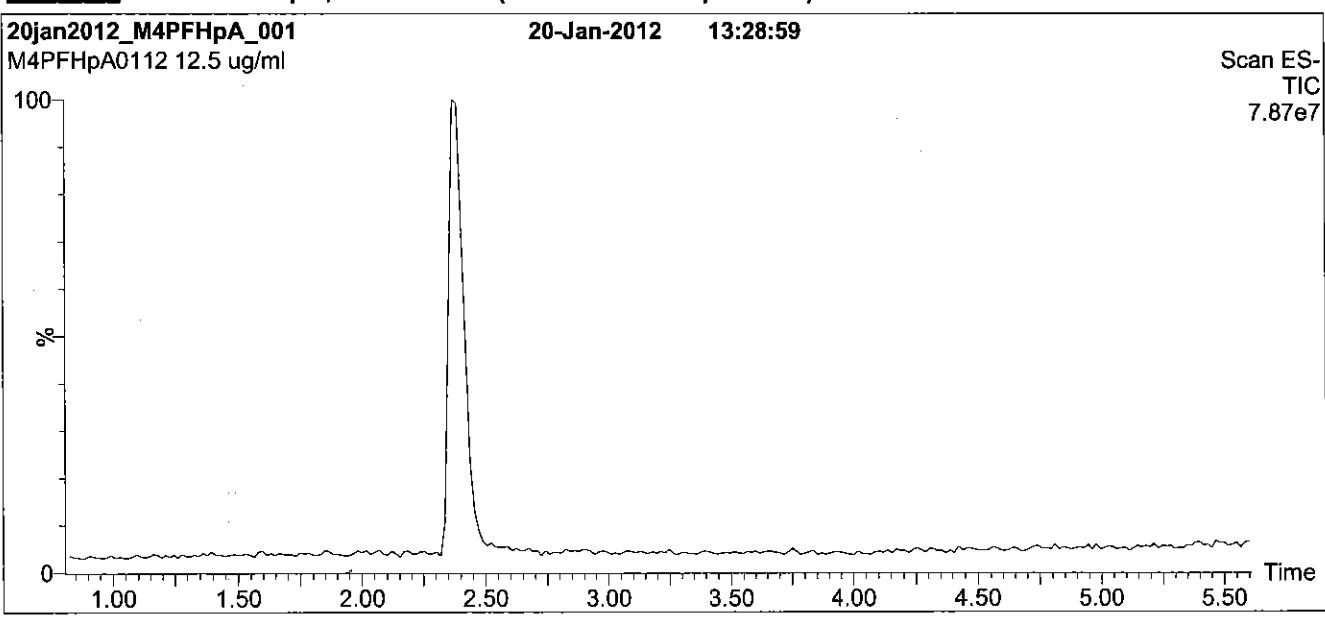
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**Figure 1:** M4PFHpA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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

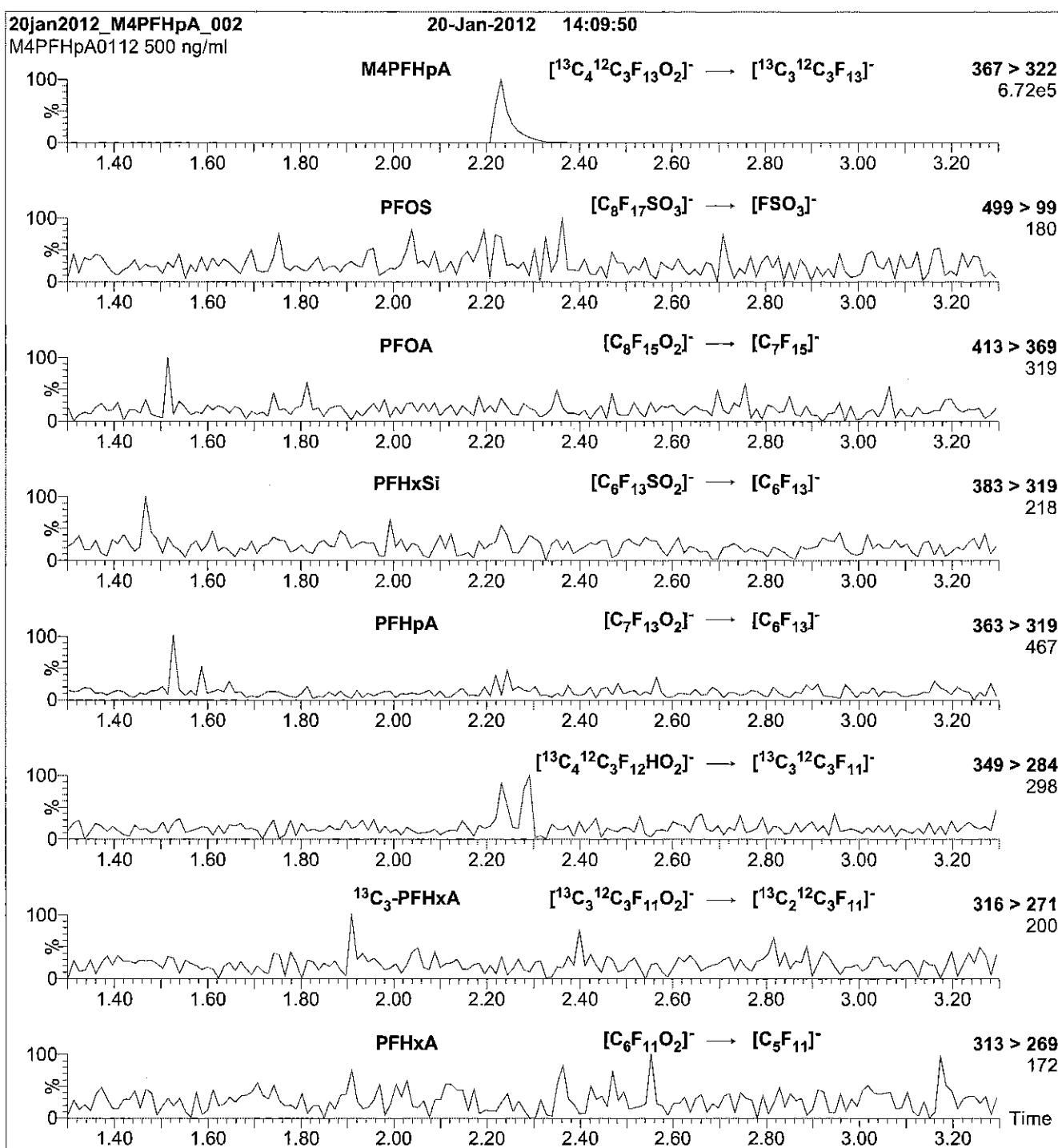
Flow: 300 µl/min

**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  $\mu$ l (500 ng/ml M4PFHpA)

**MS Parameters**

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

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

Flow: 300  $\mu$ l/min

Reagent

---

**LCM5PFPEA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LCPF LCM5PFPeA-0000 Z  
DEL 12-19-12

PRODUCT CODE:

M5PFPeA

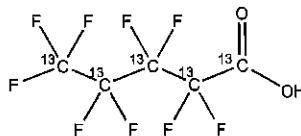
LOT NUMBER: M5PFPeA0312

COMPOUND:

Perfluoro-n-[<sup>13</sup>C<sub>5</sub>]pentanoic acid

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA:

<sup>13</sup>C<sub>5</sub>HF<sub>9</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 269.01

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: >99% <sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

03/07/2012

EXPIRY DATE: (mm/dd/yyyy)

03/07/2015

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

DOCUMENTATION/ DATA ATTACHED:

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

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

ADDITIONAL INFORMATION:

- See page 2 for further details.
- Contains 4 mole eq. of NaOH to prevent conversion of the carboxylic acid to the methyl ester.
- Contains < 0.1% of perfluoro-n-pentanoic acid.

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

Certified By:

B.G. Chittim

Date: 03/19/2012

(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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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:**

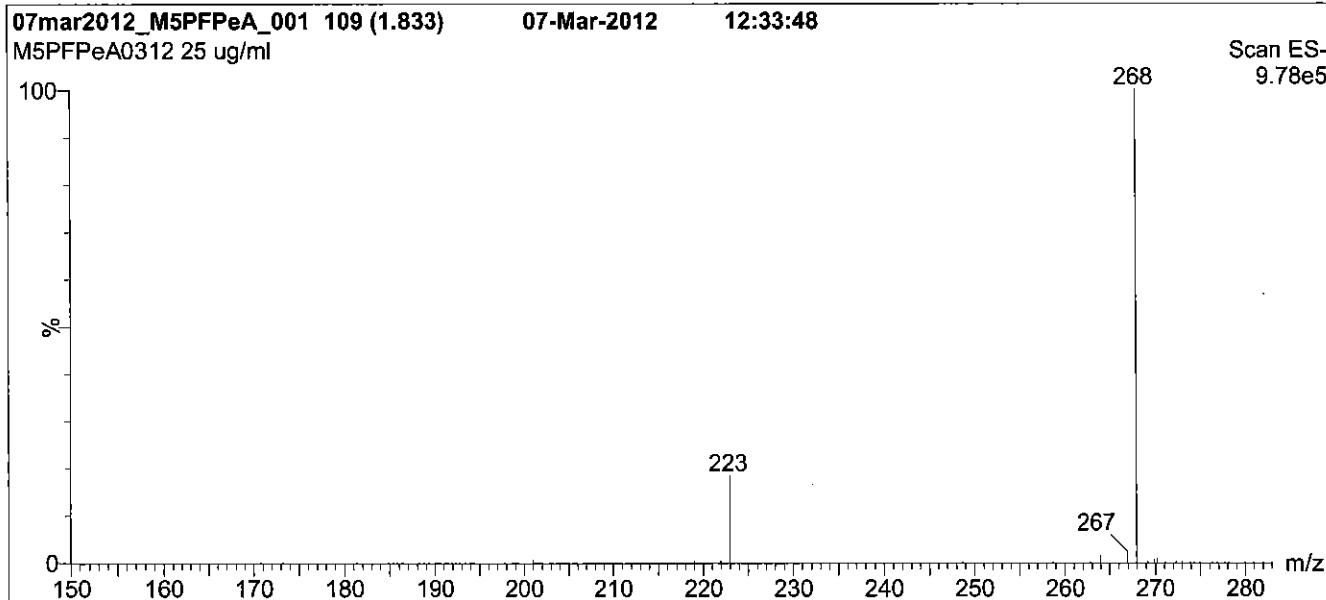
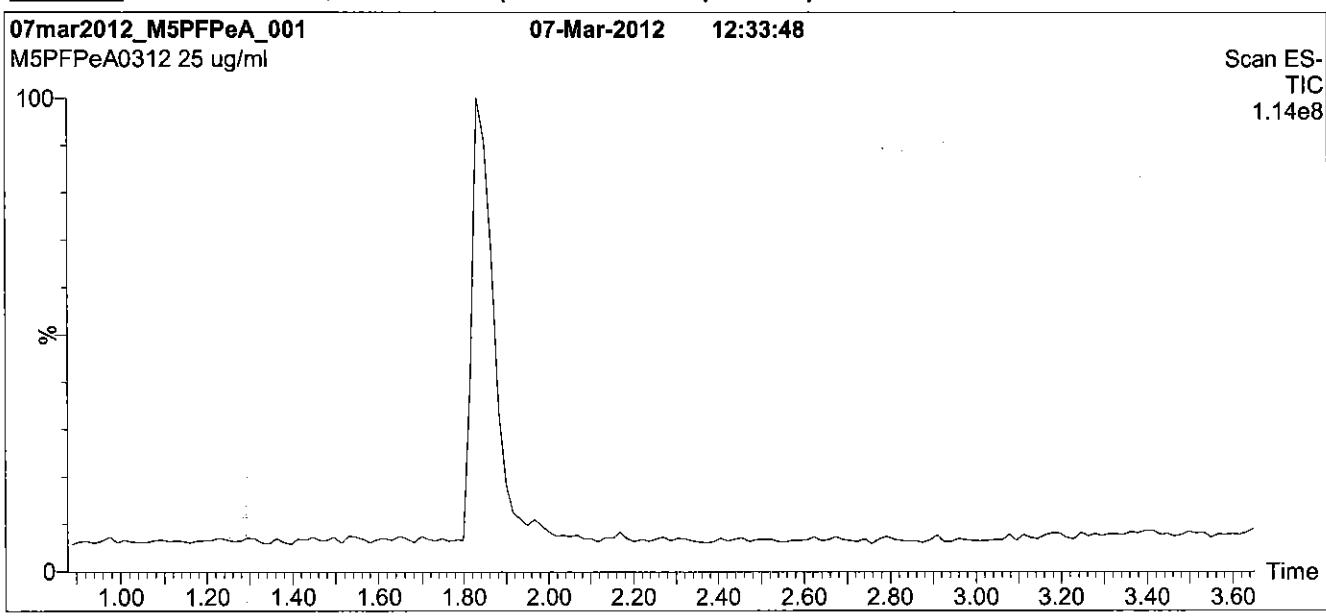
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**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.

\*\*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 Acuity Ultra Performance LC  
**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 35% (80:20 MeOH:ACN) / 65% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc)  
Ramp to 90% organic over 6.5 min and hold for  
2 min before returning to initial conditions in 0.5 min.  
Time: 10 min

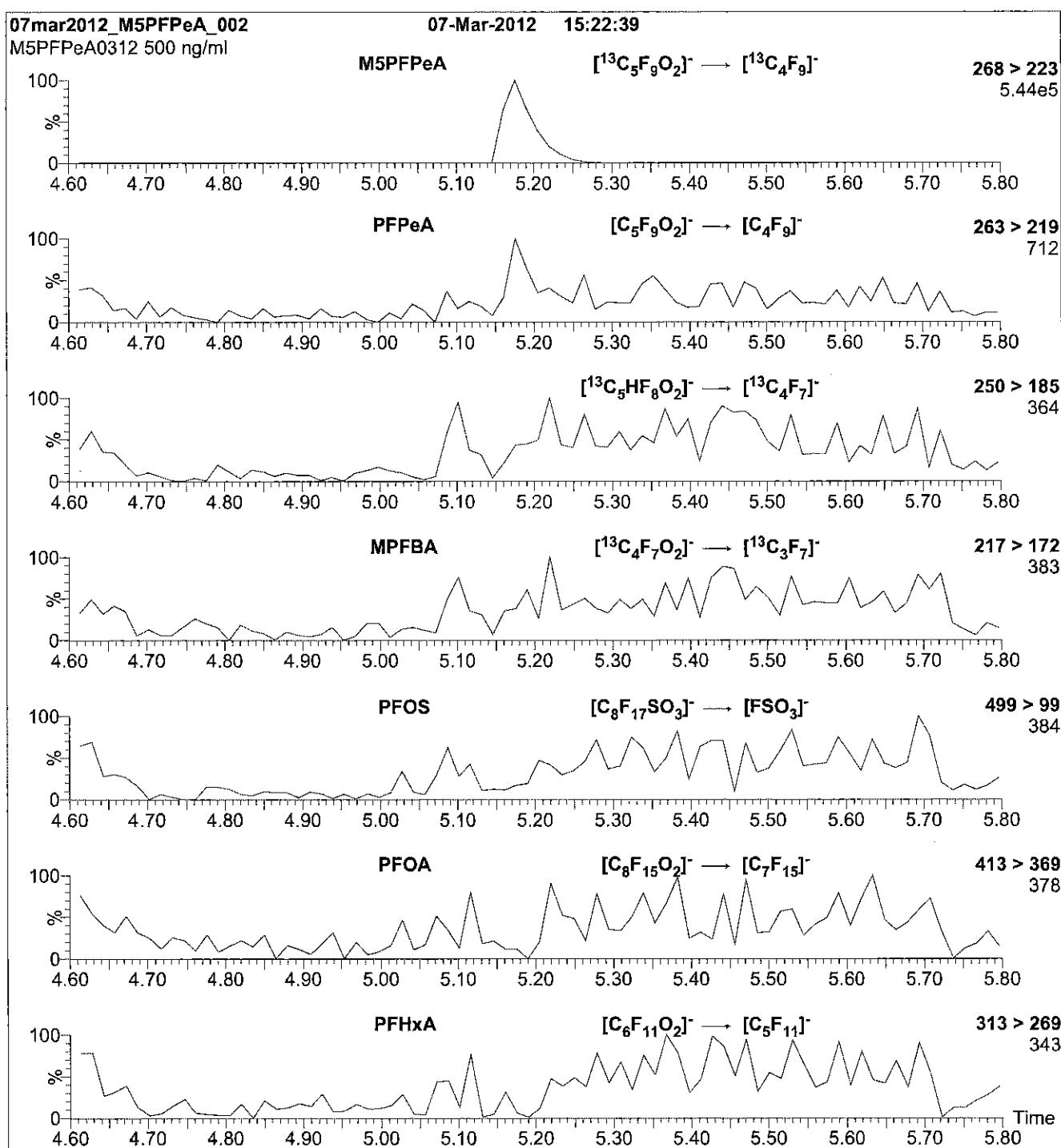
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (150 - 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)

**MS Parameters**

Collision Gas (mbar) = 3.39e-3  
Collision Energy (eV) = 9

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  $\mu\text{l}/\text{min}$

Reagent

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**LCM8FOSA\_00004**

Rec 8/14/14 SIV

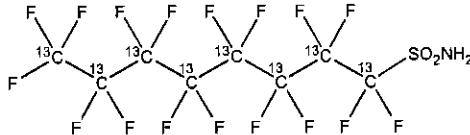


**WELLINGTON  
LABORATORIES**

**CERTIFICATE OF ANALYSIS  
DOCUMENTATION**

**PRODUCT CODE:** M8FOSA-M      **LOT NUMBER:** M8FOSA1013M  
**COMPOUND:** Perfluoro-1-[<sup>13</sup>C<sub>8</sub>]octanesulfonamide

**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>      **MOLECULAR WEIGHT:** 507.09  
**CONCENTRATION:** 50 ± 2.5 µg/ml      **SOLVENT(S):** Methanol  
**CHEMICAL PURITY:** >98%      **ISOTOPIC PURITY:** ≥99% <sup>13</sup>C  
**LAST TESTED:** (mm/dd/yyyy) 10/11/2013      **(<sup>13</sup>C<sub>8</sub>)**  
**EXPIRY DATE:** (mm/dd/yyyy) 10/11/2018  
**RECOMMENDED STORAGE:** Refrigerate ampoule

**DOCUMENTATION/ DATA ATTACHED:**

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

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

**ADDITIONAL INFORMATION:**

- See page 2 for further details.

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

Certified By:

A handwritten signature in black ink, appearing to read "B.G. Chittim".

Date: 10/25/2013

(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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#### **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.

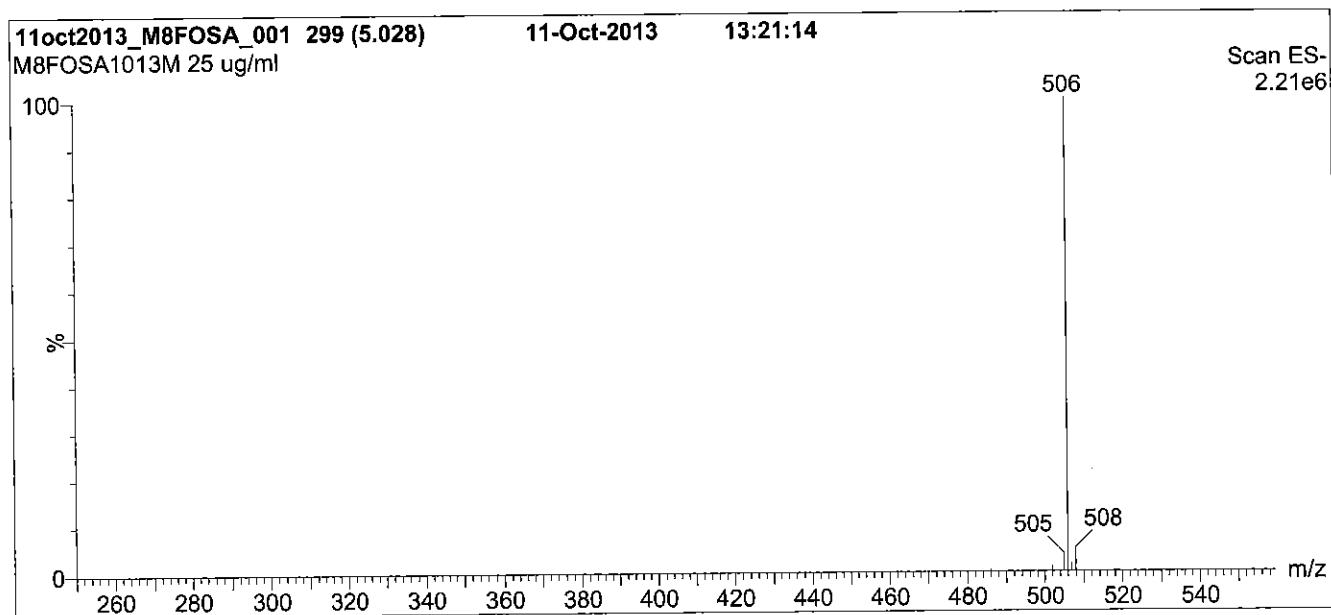
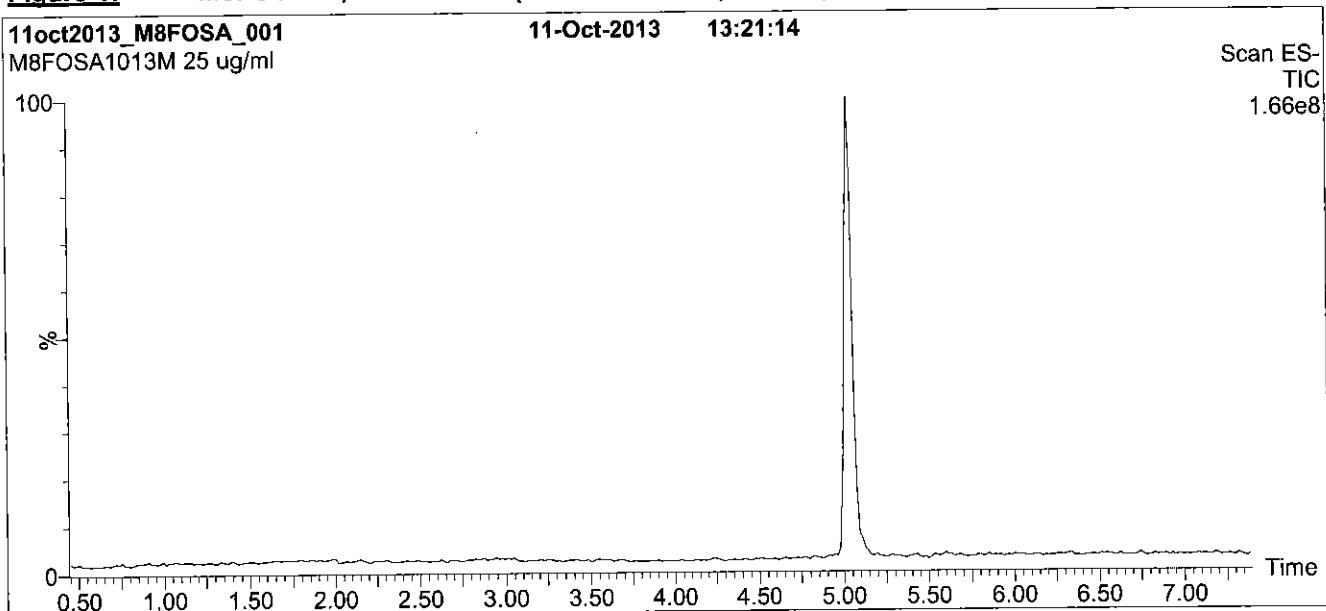
#### **QUALITY MANAGEMENT:**

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 ACCLASS (certificate number AR-1523).



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**Figure 1:** M8FOSA-M; LC/MS Data (TIC and Mass Spectrum)



#### **Conditions for Figure 1:**

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

### **Chromatographic Conditions**

Column: Acquity UPLC BEH Shield RP<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 7.5 min and hold for 1 min  
 before returning to initial conditions in 0.5 min.  
 Time: 10 min

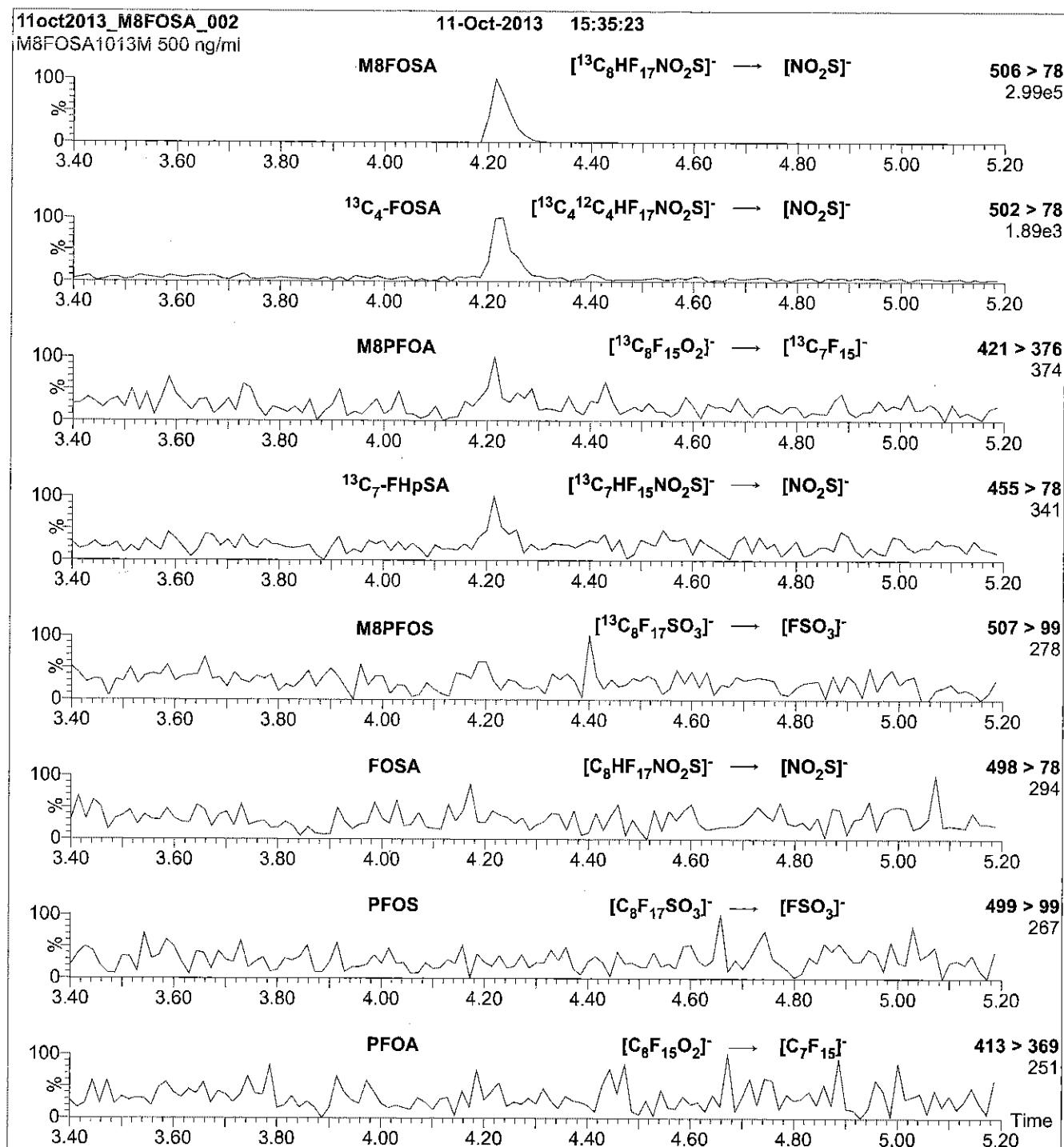
Flow: 300  $\mu$ l/min

## MS Parameters

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

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

**Figure 2:** M8FOSA-M; LC/MS/MS Data (Selected MRM Transitions)



**Conditions for Figure 2:**

Injection: Direct loop injection  
10  $\mu$ l (500 ng/ml M8FOSA-M)

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 30

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

Reagent

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**LCMPFBA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R 12-14-12

LCMPFBA-00002

PRODUCT CODE:

MPFBA

LOT NUMBER: MPFBA0812

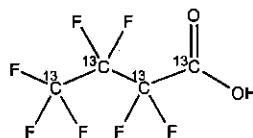
COMPOUND:

Perfluoro-n-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]butanoic acid

STRUCTURE:

CAS #:

Not available



MOLECULAR FORMULA:

<sup>13</sup>C<sub>4</sub>HF<sub>7</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 218.01

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

08/13/2012

EXPIRY DATE: (mm/dd/yyyy)

08/13/2015

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 ~ 0.7% of native PFBA.

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

Certified By:

B.G. Chittim

Date: 08/23/2012

(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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**SYNTHESIS / CHARACTERIZATION:**

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**HOMOGENEITY:**

Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS 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:**

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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:**

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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.

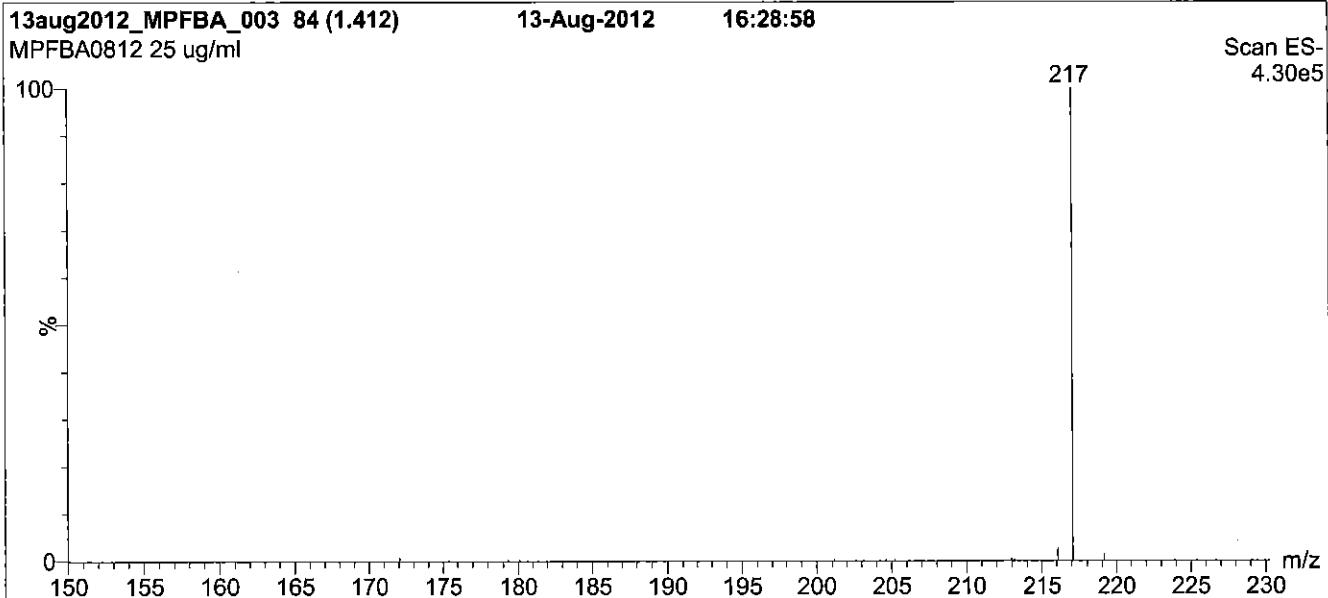
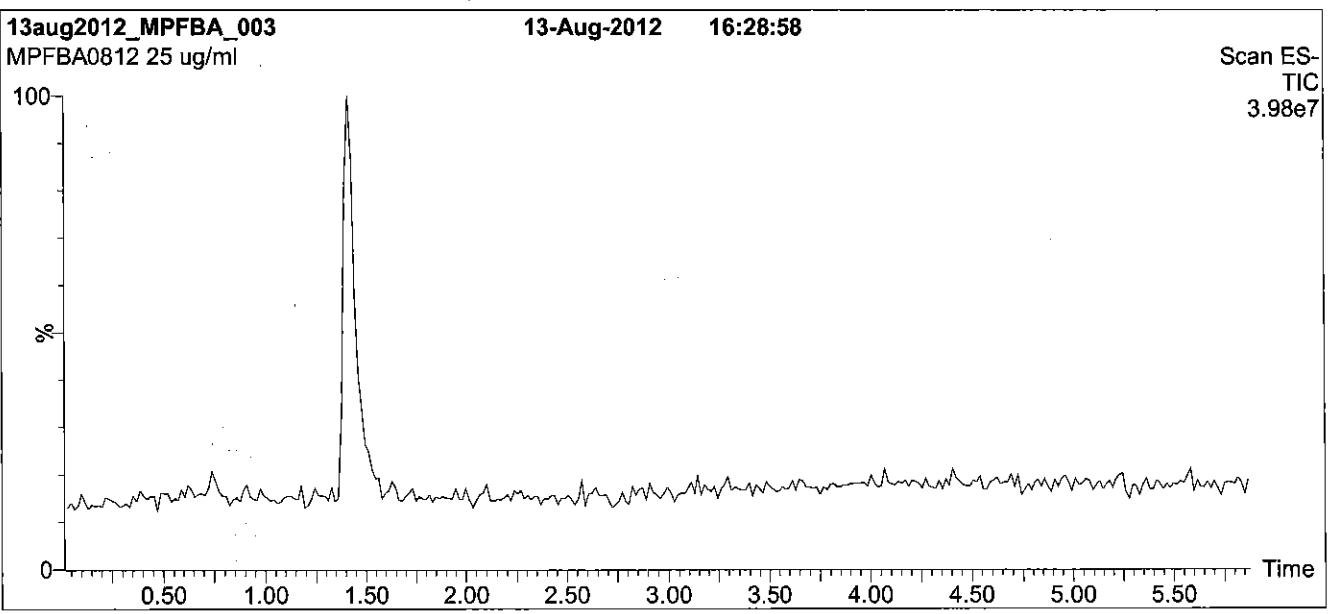
**QUALITY MANAGEMENT:**

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 ACCLASS (certificate number AR-1523).



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



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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 7.5 min and hold for 1 min  
before returning to initial conditions in 0.5 min.

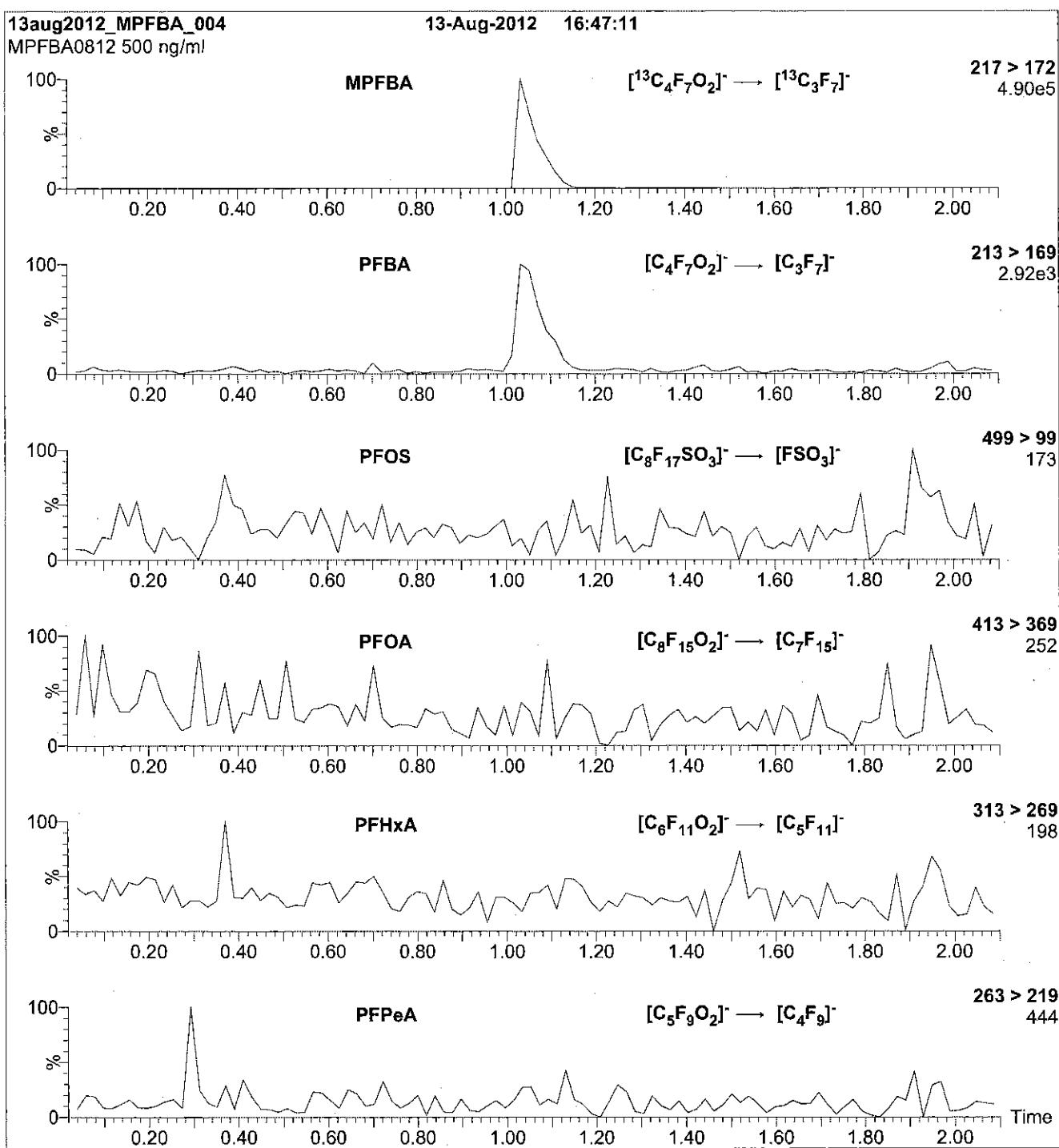
Time: 10 min  
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 10.00  
Cone Gas Flow (l/hr) = 60  
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)

**MS Parameters**

Collision Gas (mbar) = 3.62e-3  
Collision Energy (eV) = 10

Mobile phase: Isocratic 60% (80:20 MeOH:ACN) / 40%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

Reagent

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**LCMPFDA\_00003**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE:

MPFDA

LOT NUMBER: MPFDA0912

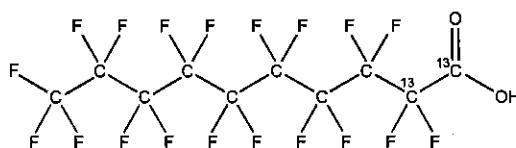
COMPOUND:

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

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: >99% <sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

09/13/2012

EXPIRY DATE: (mm/dd/yyyy)

09/13/2017

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 ~ 0.1% 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: 01/09/2013

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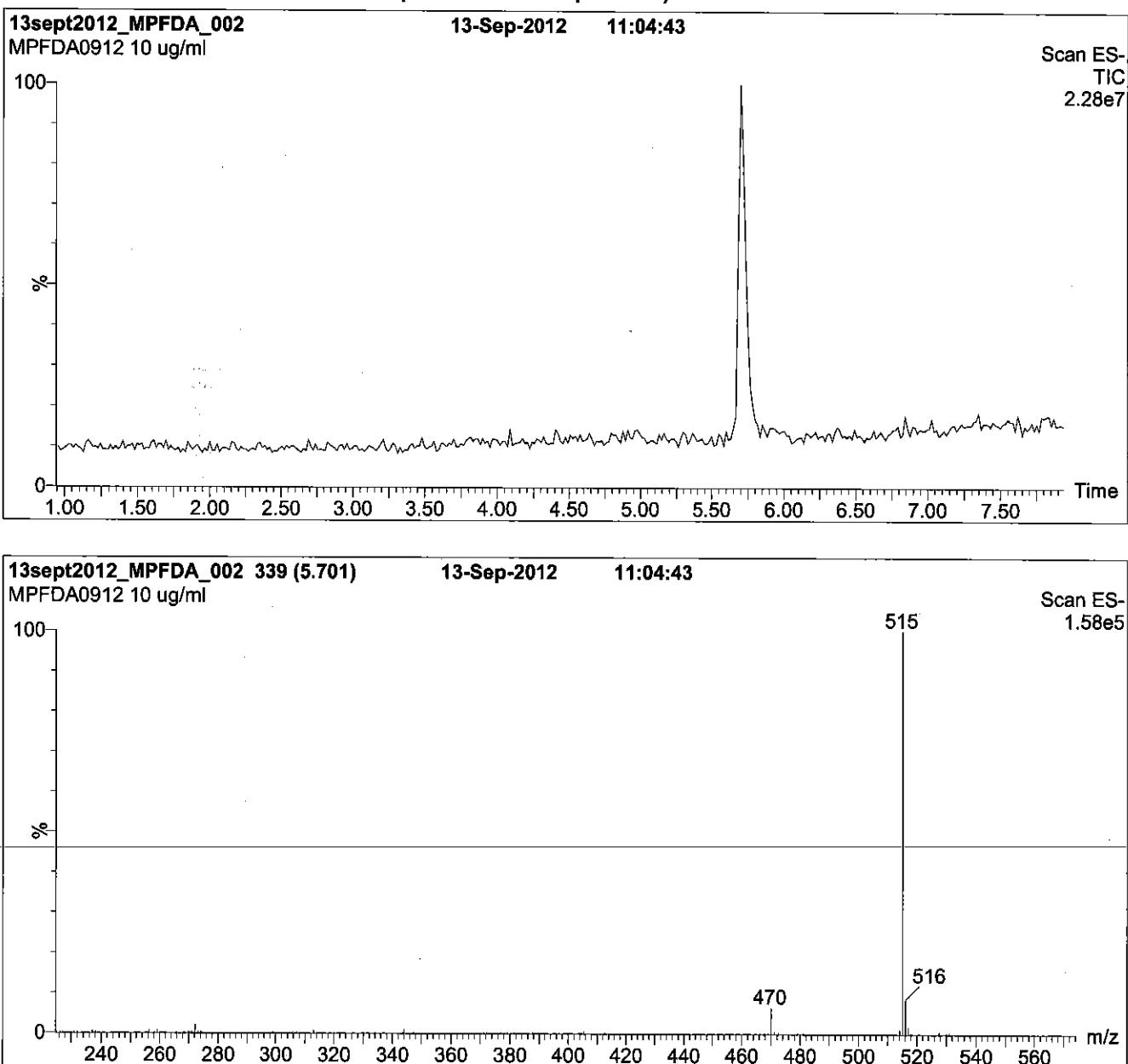
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**Figure 1:** MPFDA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>,  
1.7 µm, 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

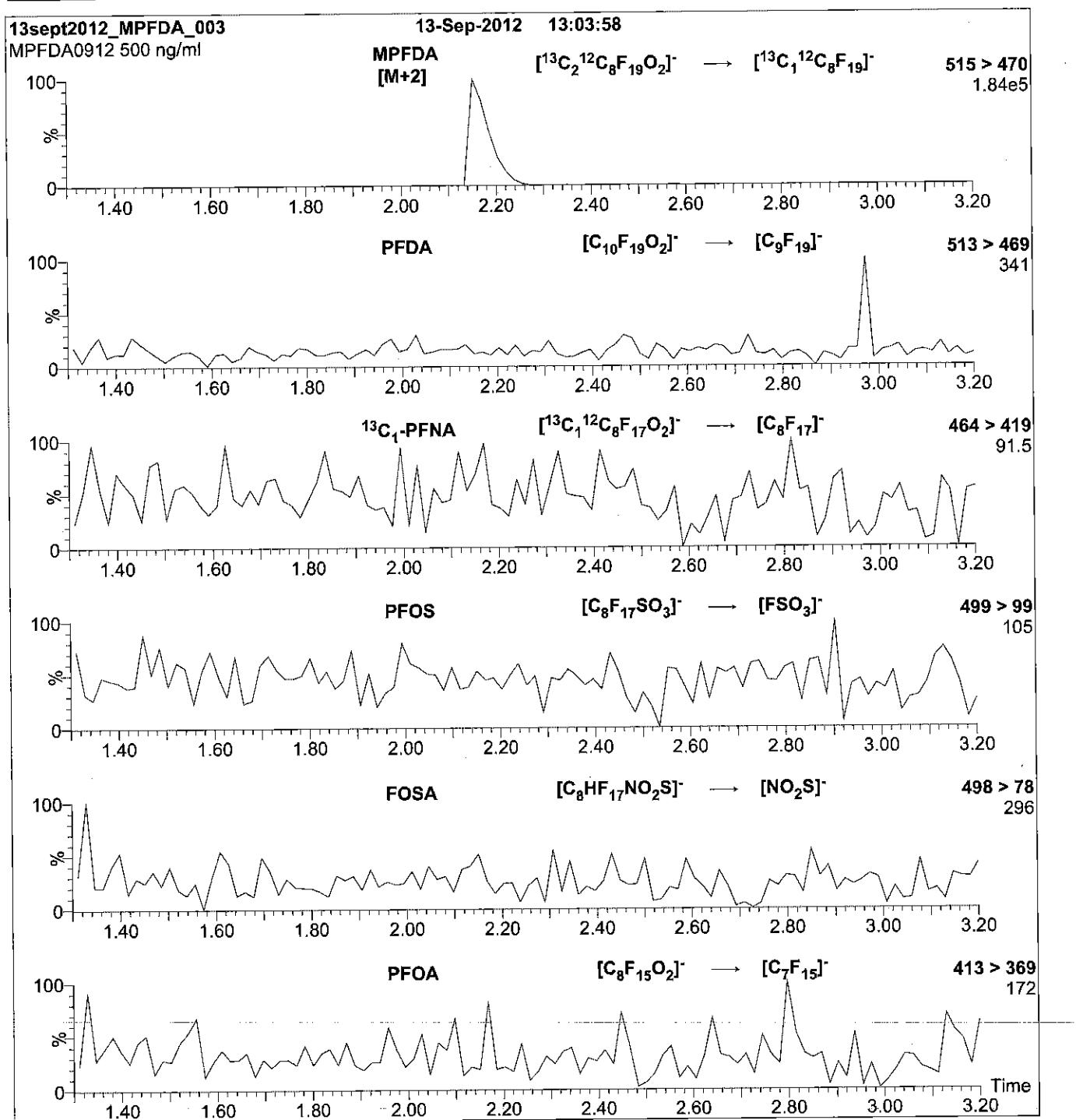
Flow: 300 µl/min

**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)

**MS Parameters**

Collision Gas (mbar) = 3.62e-3  
Collision Energy (eV) = 13

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  $\mu\text{l}/\text{min}$

Reagent

---

**LCMPFDoA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LCMPFDoA\_00002

PRODUCT CODE:

MPFDoA

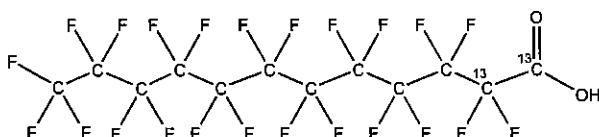
LOT NUMBER: MPFDoA0312

COMPOUND:

Perfluoro-n-[1,2-<sup>13</sup>C]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

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: >99% <sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

03/26/2012

EXPIRY DATE: (mm/dd/yyyy)

03/26/2015

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: 04/04/2012

(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:**

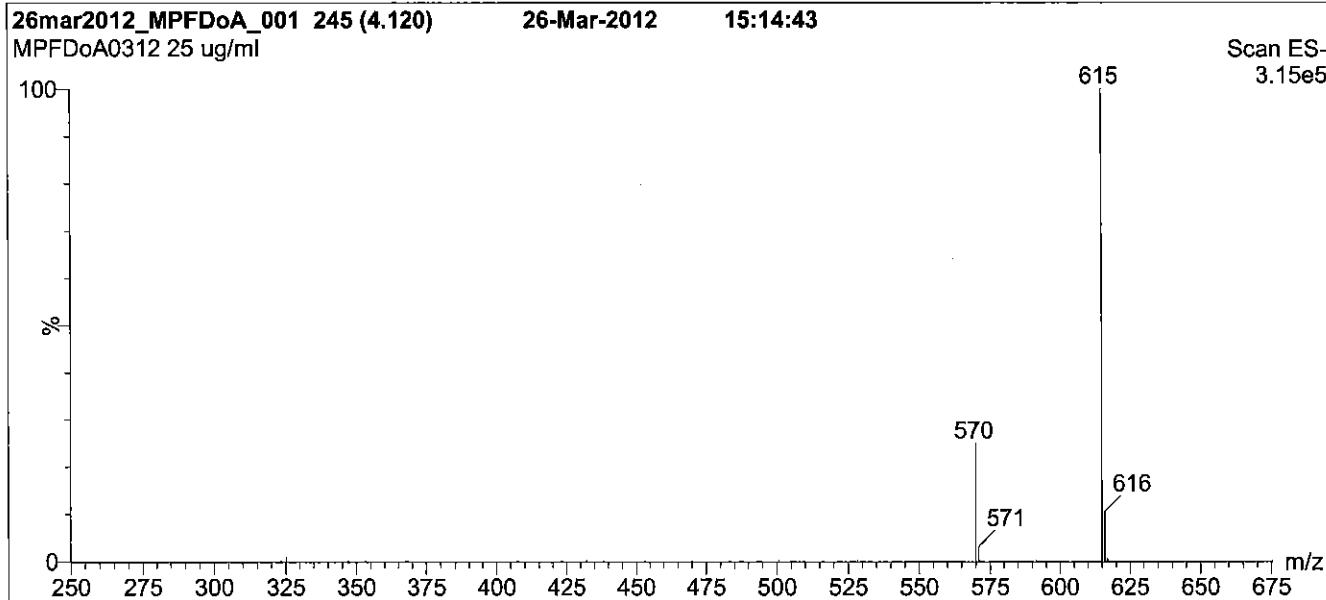
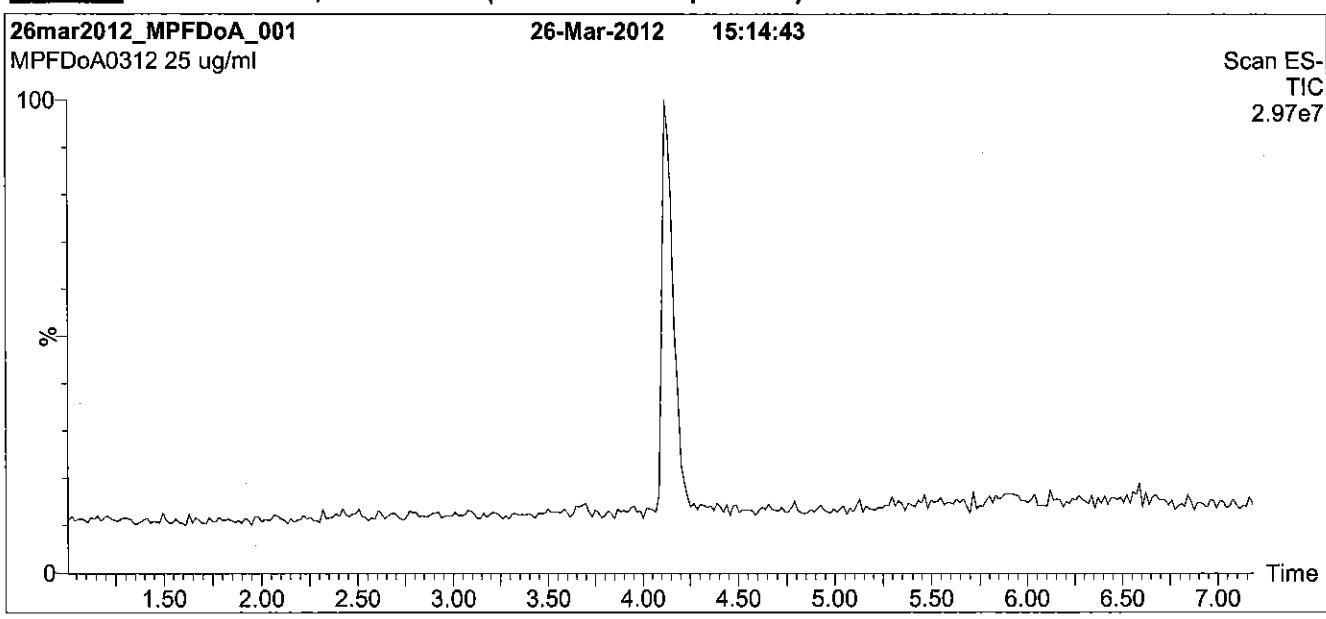
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**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.

\*\*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 Acuity Ultra Performance LC  
**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity 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 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 µl/min

**MS Parameters**

Experiment: Full Scan (250 - 850 amu)

Source: Electrospray (negative)

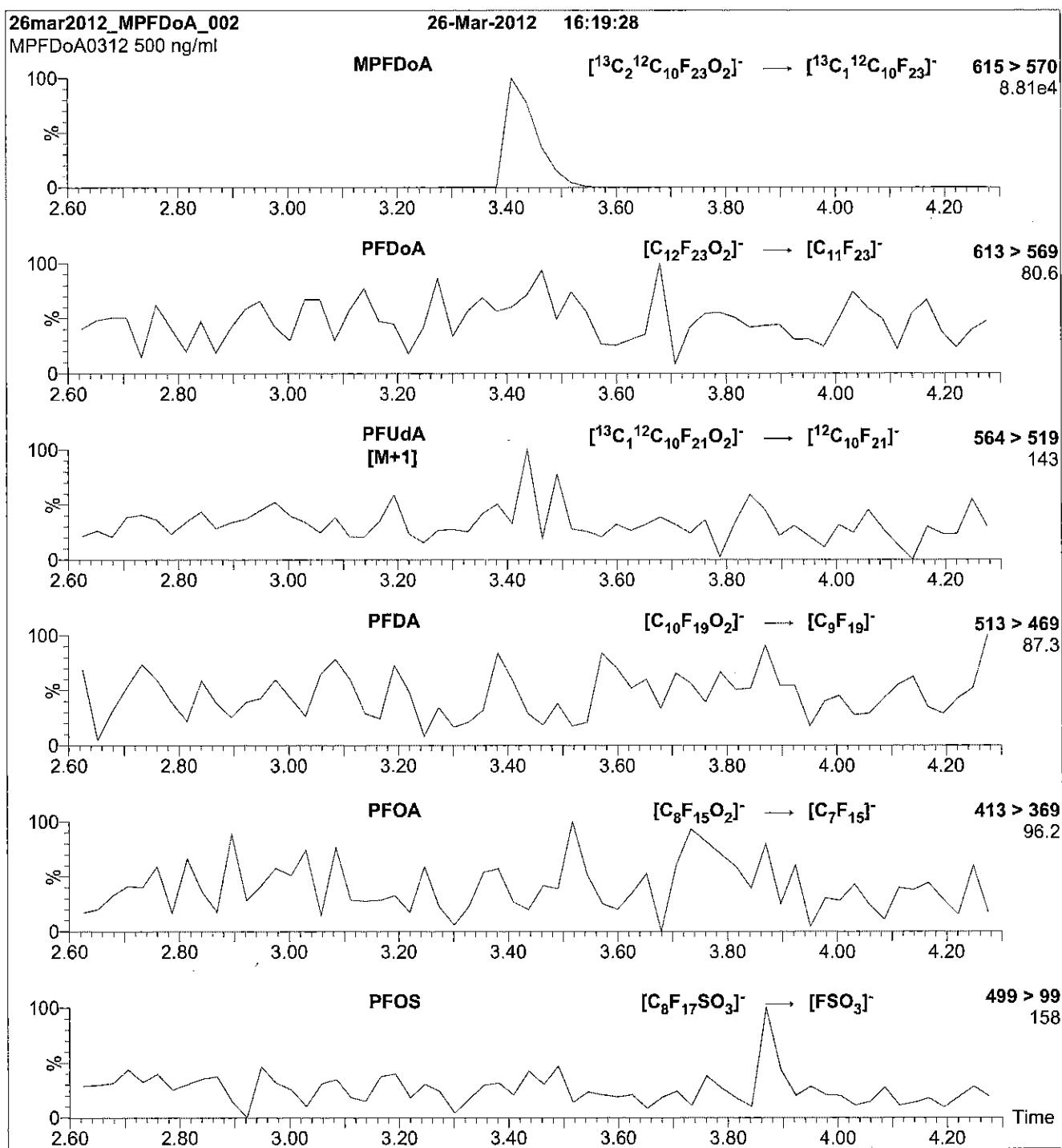
Capillary Voltage (kV) = 2.00

Cone Voltage (V) = 20.00

Cone Gas Flow (l/hr) = 100

Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
 Collision Energy (eV) = 13

Mobile phase: Isocratic 60% (80:20 MeOH:ACN) / 40%  $\text{H}_2\text{O}$   
 (both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

Reagent

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**LCMPFHxA  00004**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE:

MPFHxA

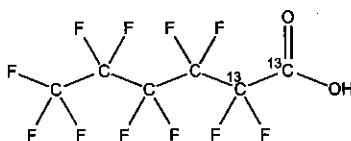
LOT NUMBER: MPFHxA0213

COMPOUND:

Perfluoro-n-[1,2-<sup>13</sup>C<sub>2</sub>]hexanoic acid

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>

MOLECULAR WEIGHT: 316.04

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol  
Water (<1%)

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: >99%<sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

02/14/2013

(1,2-<sup>13</sup>C<sub>2</sub>)

EXPIRY DATE: (mm/dd/yyyy)

02/14/2018

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 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: 02/20/2013

(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.

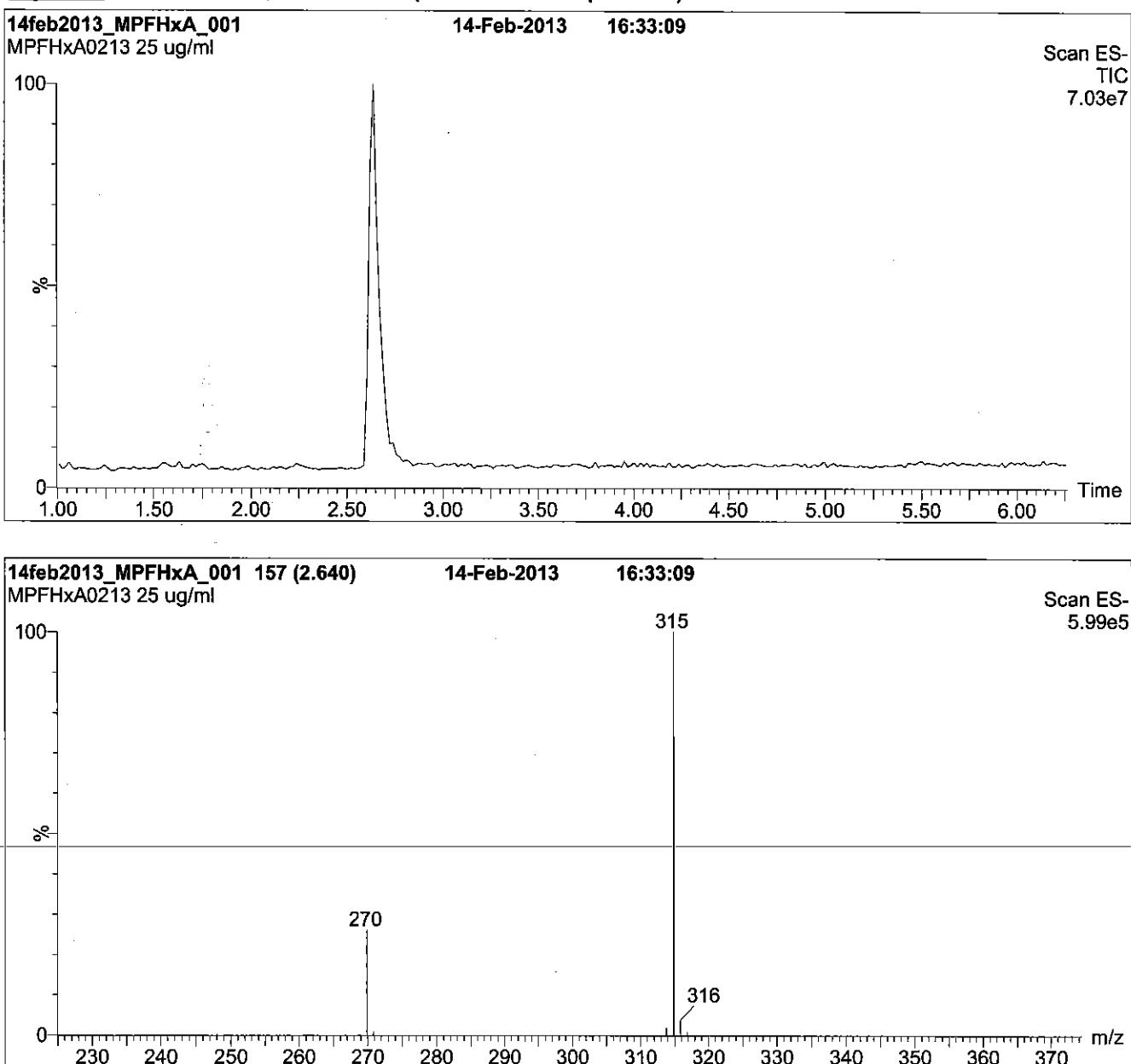
#### **QUALITY MANAGEMENT:**

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 ACCLASS (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 Acuity Ultra Performance LC  
**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP,  
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 over 0.5 min.  
Time: 10 min

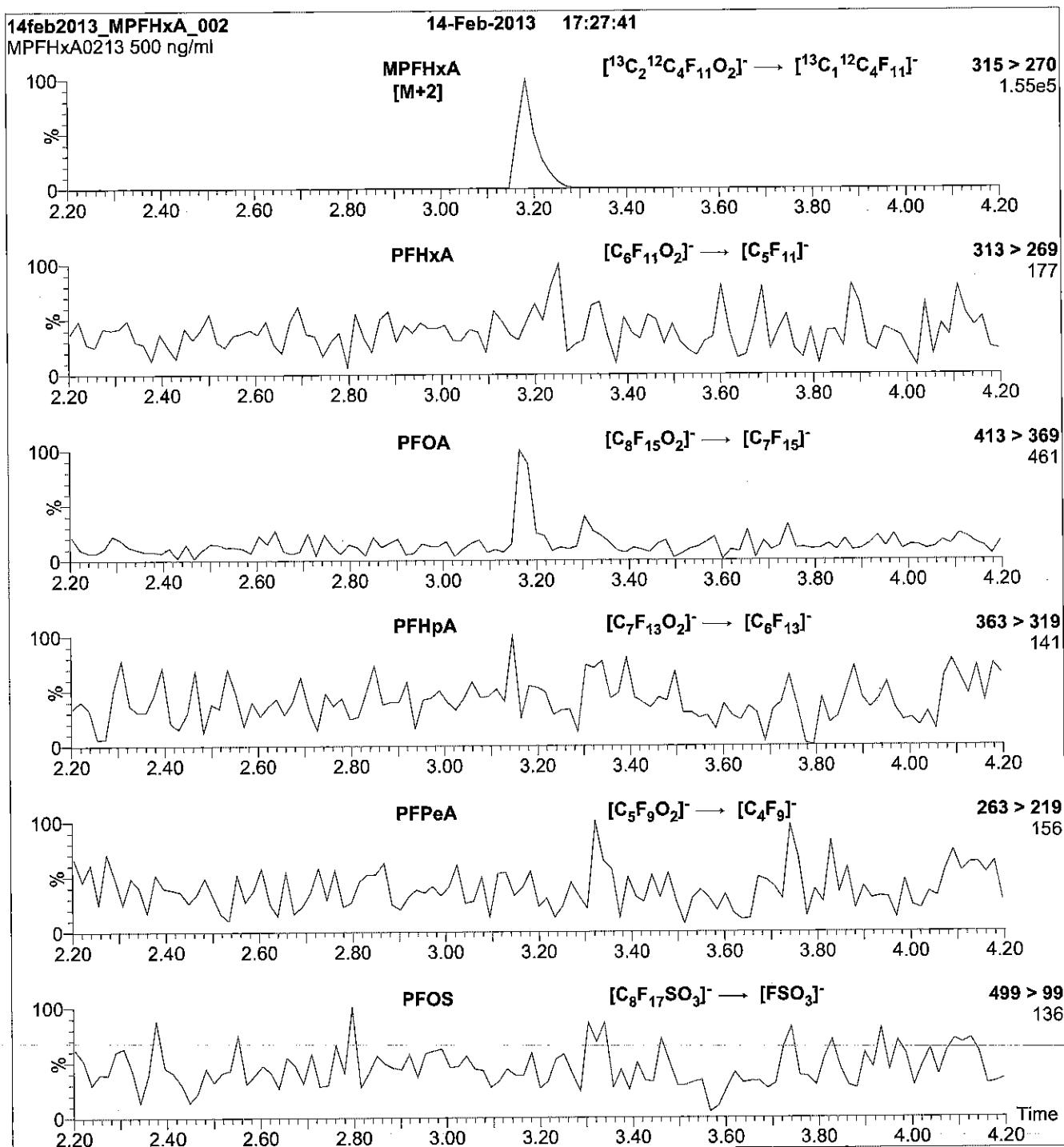
Flow: 300  $\mu$ l/min

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



**Conditions for Figure 2:**

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

**MS Parameters**

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

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  $\mu\text{l}/\text{min}$

Reagent

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**LCMPFHxS  00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LCMPFHxS-00002

PRODUCT CODE:

MPFHxS

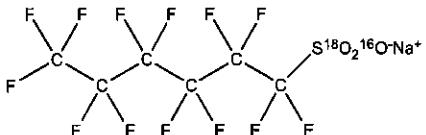
LOT NUMBER: MPFHxS0312

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<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)

03/08/2012

EXPIRY DATE: (mm/dd/yyyy)

03/08/2015

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.

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

Certified By:

B.G. Chittim

Date: 03/19/2012

(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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**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:**

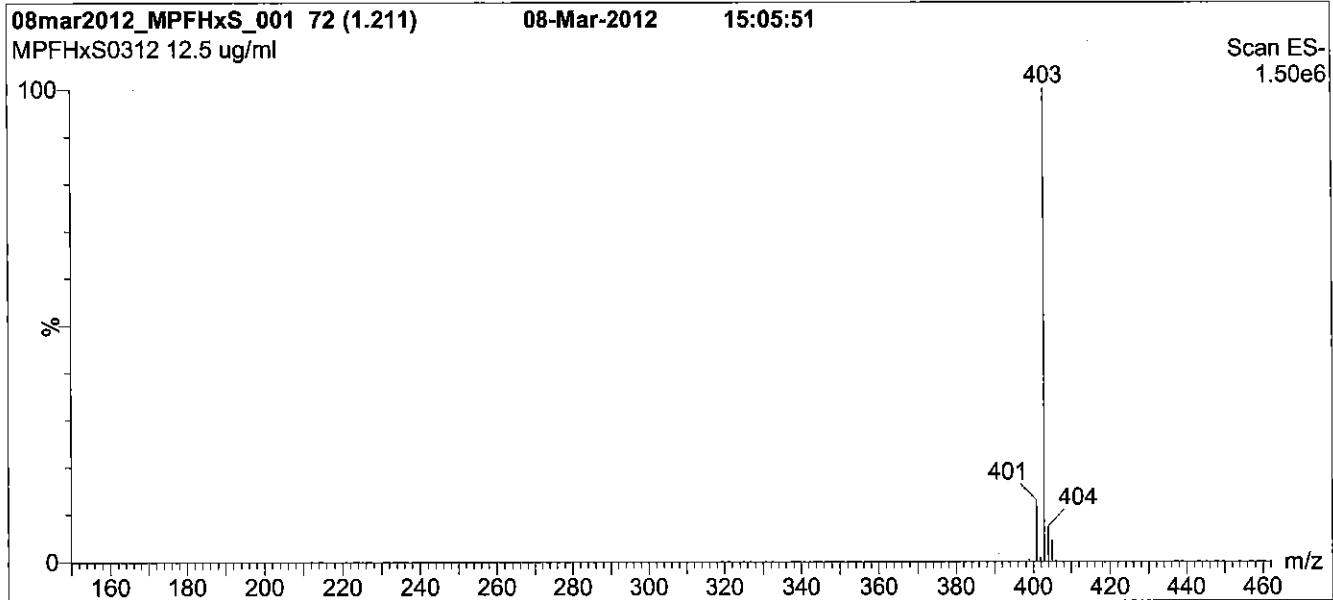
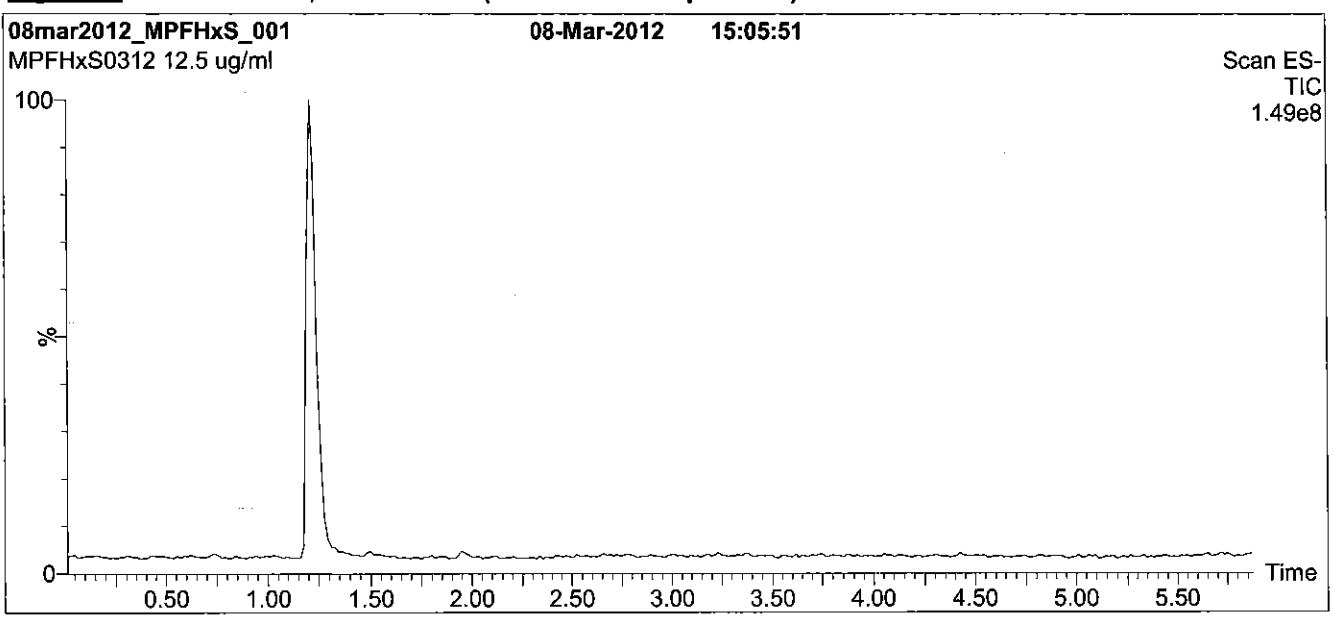
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**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.

\*\*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:** MPFHxS; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 2.1 x 100 mm

Mobile phase: Isocratic with flush  
Start: 65% (80:20 MeOH:ACN) / 35% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Hold for 6.5 min. Ramp to 90% organic over 0.5 min. Hold  
for 1.5 min before returning to initial conditions in 0.5 min.  
Time: 10 min

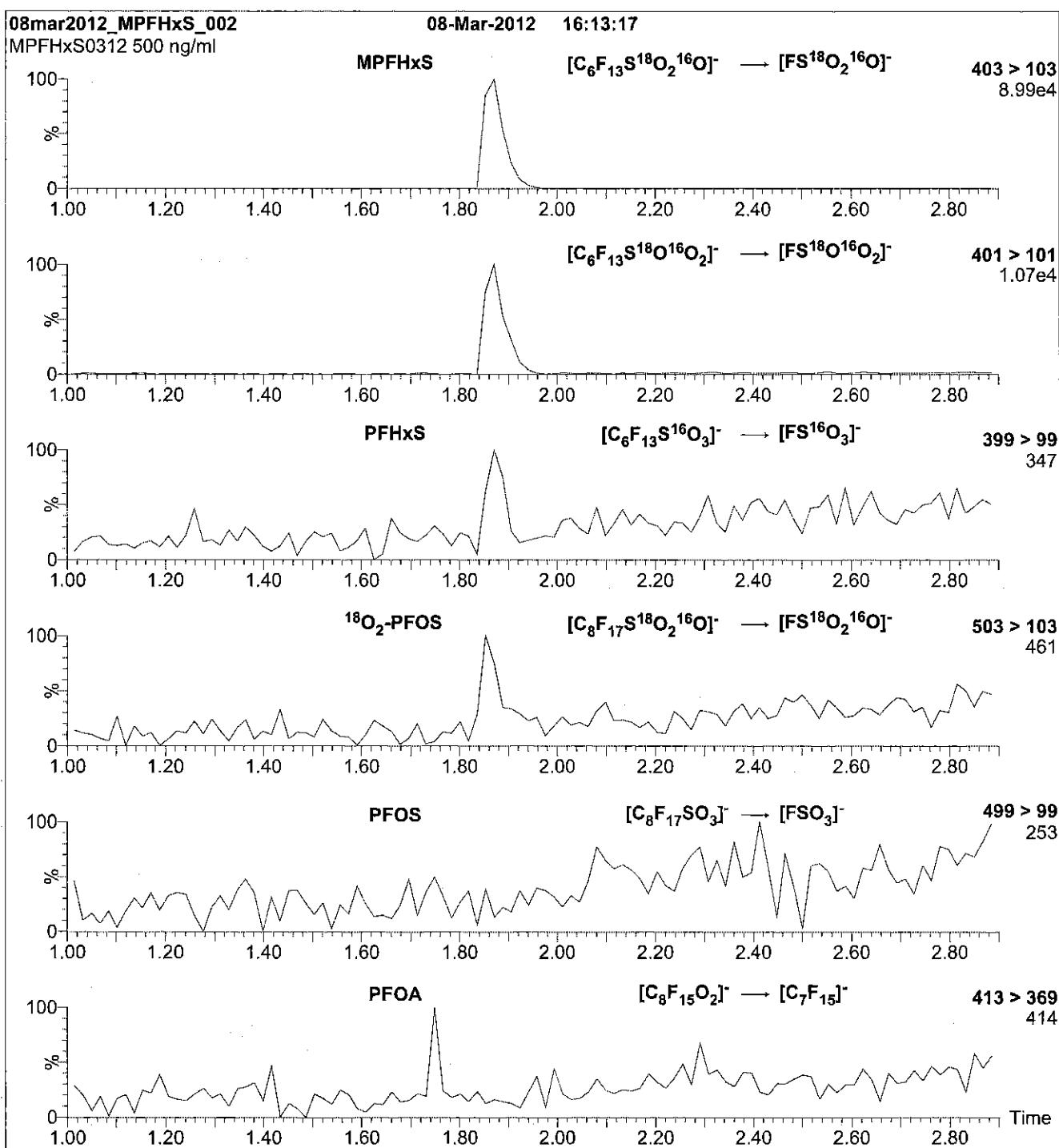
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (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)

**MS Parameters**

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

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

Reagent

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**LCMPFNA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

12-14-12

LCMPFNA\_00002

PRODUCT CODE:

MPFNA

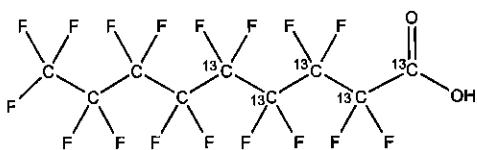
LOT NUMBER: MPFNA0912

COMPOUND:

Perfluoro-n-[1,2,3,4,5-<sup>13</sup>C<sub>5</sub>]nonanoic acid

STRUCTURE:

CAS #: Not available



MOLECULAR FORMULA:

<sup>13</sup>C<sub>5</sub><sup>12</sup>C<sub>4</sub>HF<sub>17</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 469.04

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: >99%<sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

09/13/2012

(1,2,3,4,5-<sup>13</sup>C<sub>5</sub>)

EXPIRY DATE: (mm/dd/yyyy)

09/13/2015

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: 09/26/2012

(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.

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**LIMITED WARRANTY:**

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

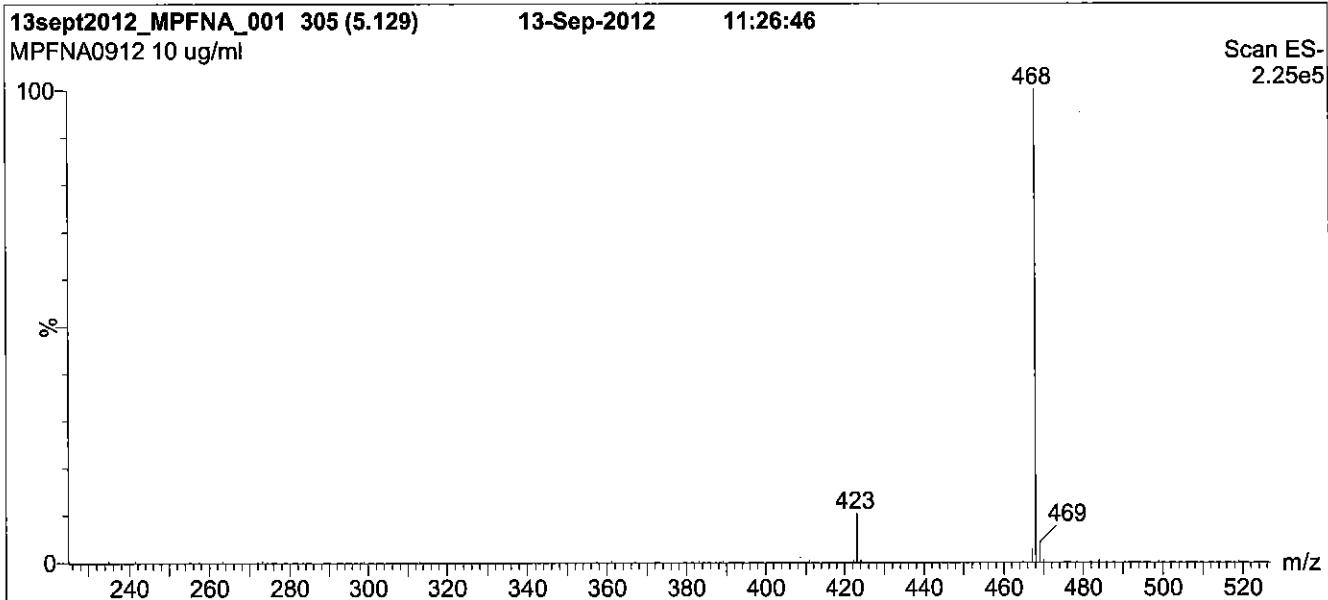
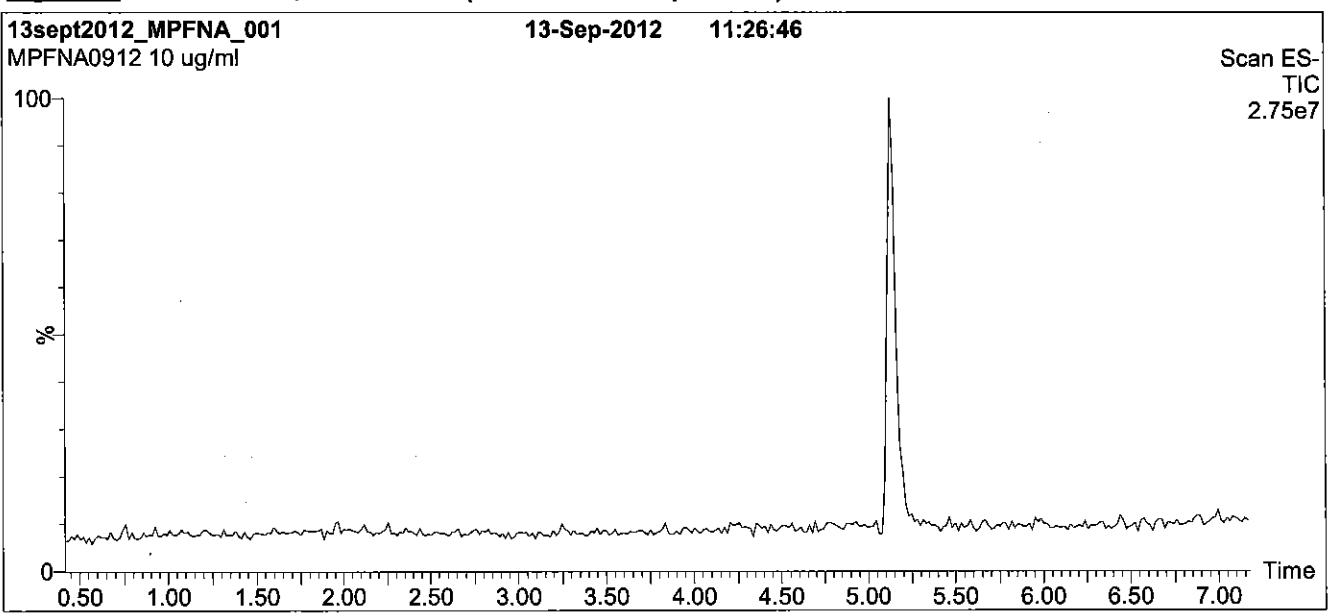
**QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to 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 ACCLASS (certificate number AR-1523).



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



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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

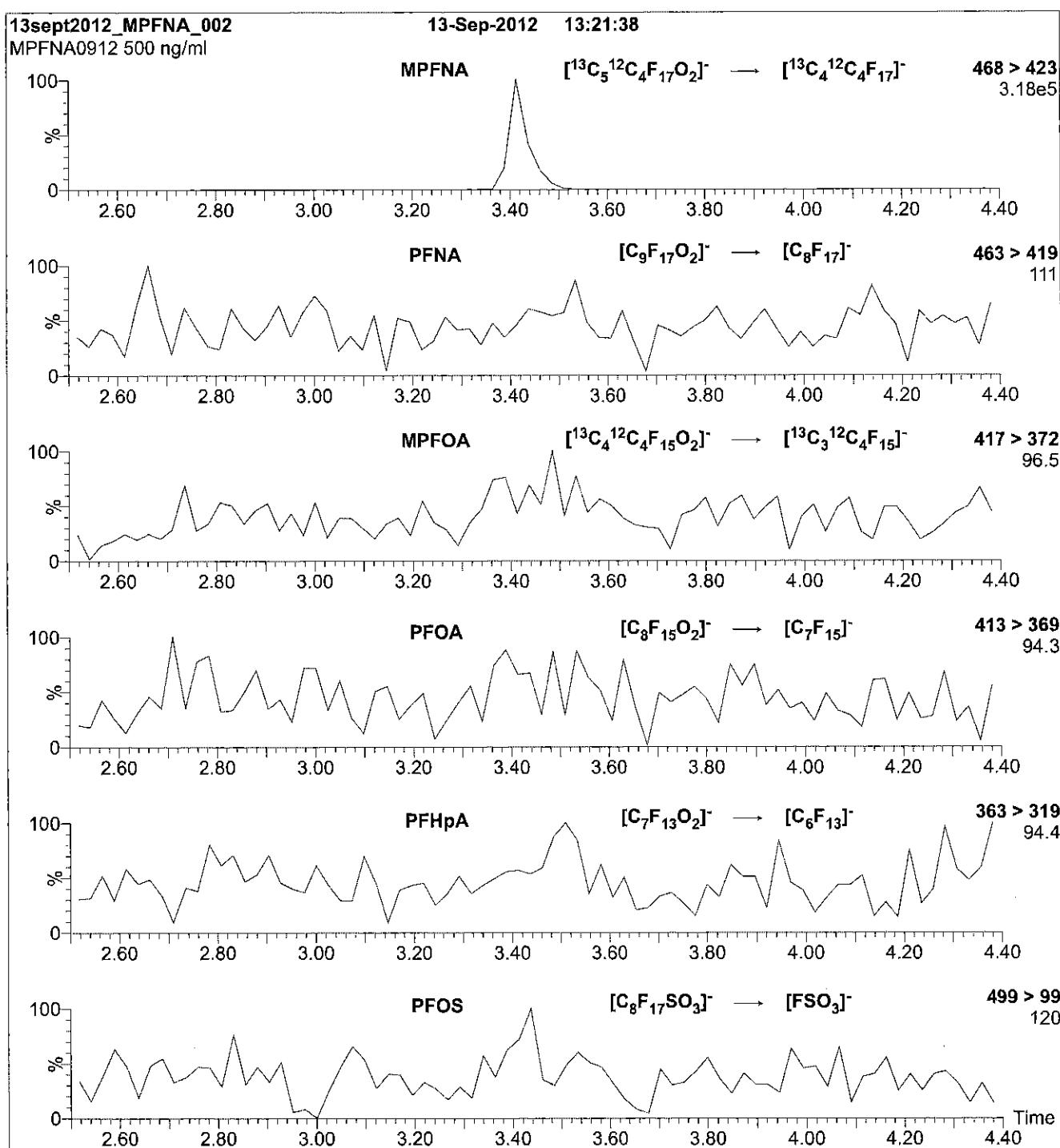
Flow: 300 µl/min

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



**Conditions for Figure 2:**

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

**MS Parameters**

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

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  $\mu\text{l}/\text{min}$

Reagent

---

**LCMPFOA\_00005**

Rec 8/14/14 SK

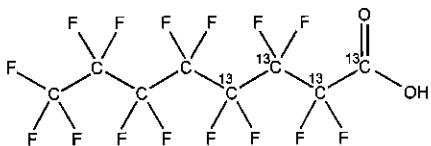


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: MPFOA      LOT NUMBER: MPFOA0614  
COMPOUND: Perfluoro-n-[1,2,3,4-<sup>13</sup>C<sub>4</sub>]octanoic acid

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>      MOLECULAR WEIGHT: 418.04  
CONCENTRATION: 50 ± 2.5 µg/ml      SOLVENT(S): Methanol  
Water (<1%)  
CHEMICAL PURITY: >98%      ISOTOPIC PURITY: ≥99% <sup>13</sup>C  
LAST TESTED: (mm/dd/yyyy) 07/04/2014      (1,2,3,4-<sup>13</sup>C<sub>4</sub>)  
EXPIRY DATE: (mm/dd/yyyy) 07/04/2019  
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 native perfluoro-n-octanoic acid (PFOA).

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

Certified By:

B.G. Chittim

Date: 07/08/2014

(mm/dd/yyyy)

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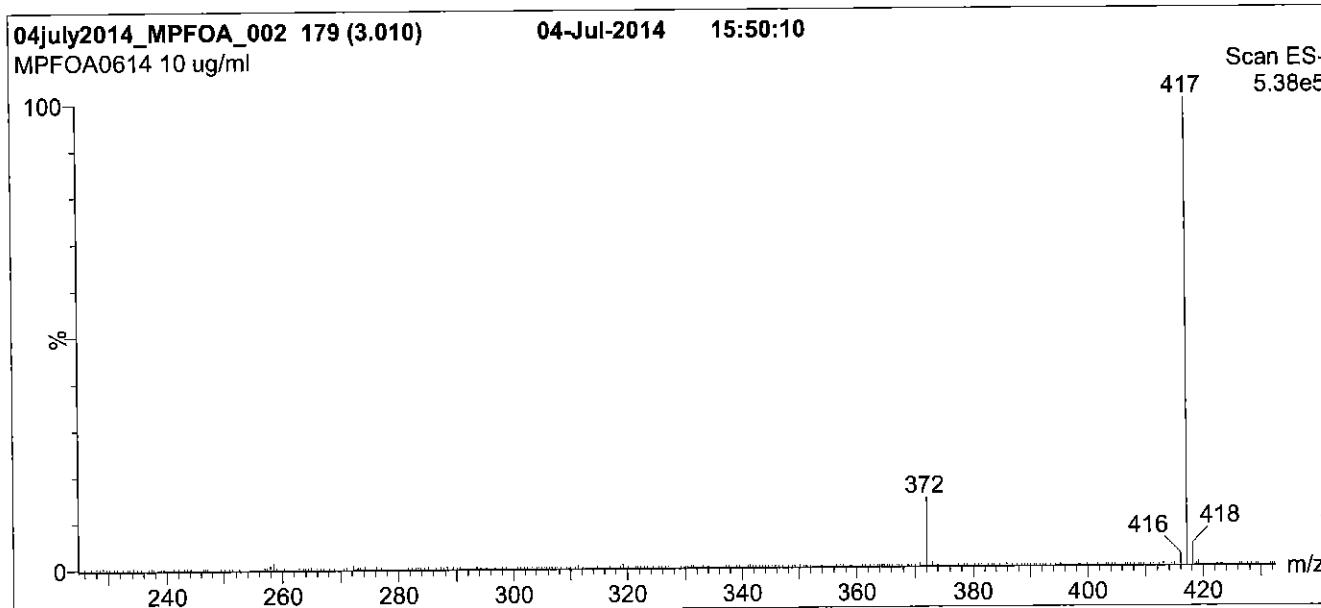
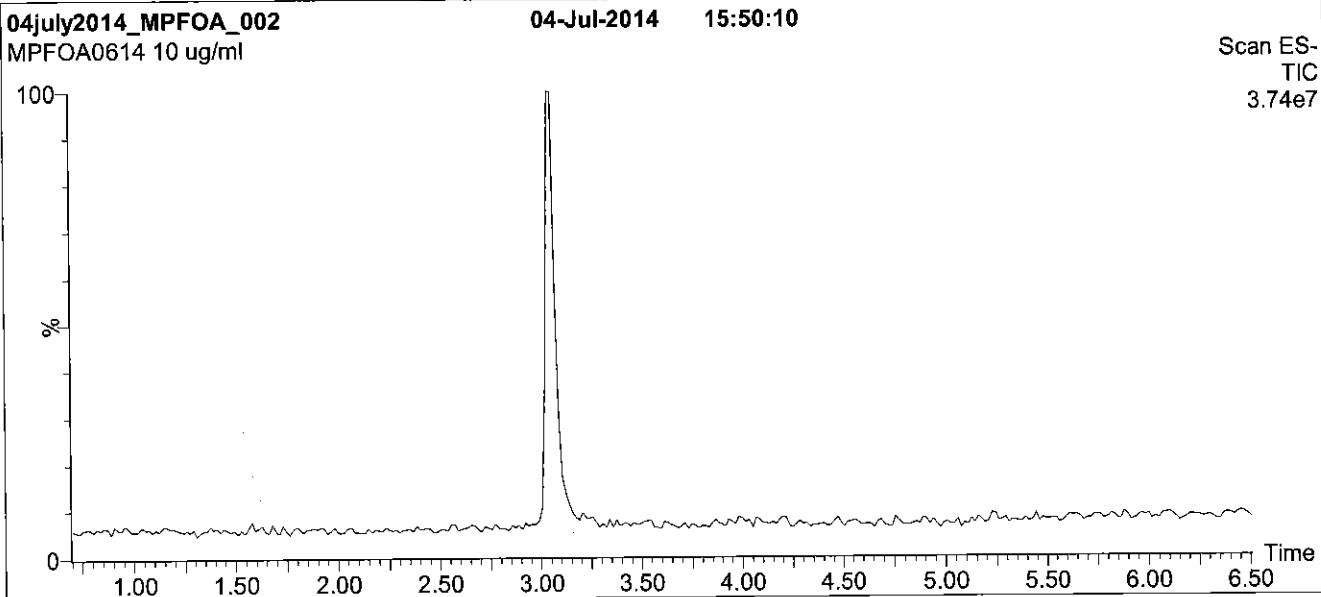
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**Figure 1:** MPFOA; LC/MS Data (TIC and Mass Spectrum)



**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 µm, 2.1 x 100 mm

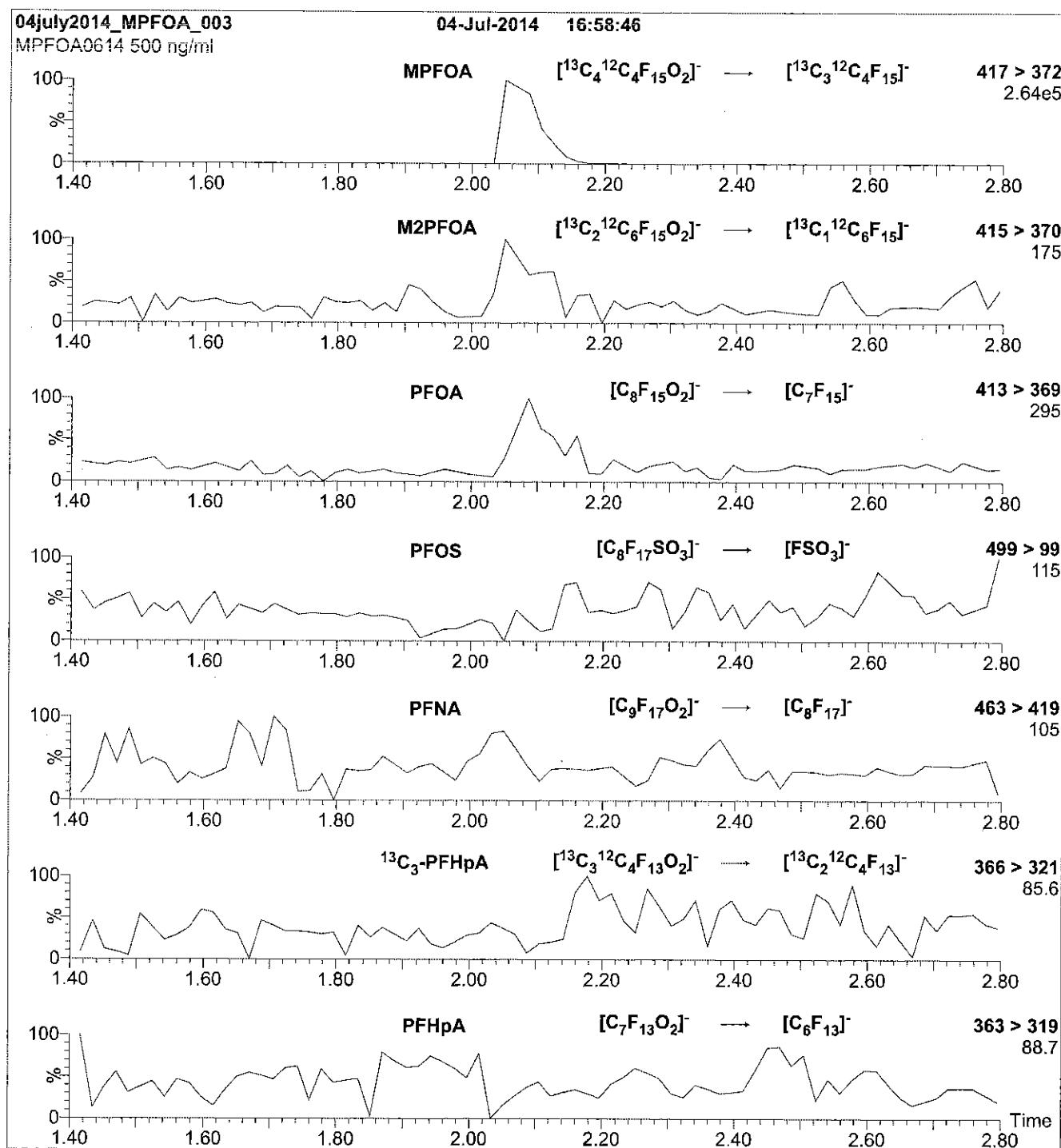
Mobile phase: Gradient  
Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7.5 min and hold for 1 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min

Flow: 300 µl/min

**MS Parameters**

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.62e-3  
Collision Energy (eV) = 11

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  $\mu\text{l}/\text{min}$

Reagent

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**LCMPFOS\_00006**

Rec'd 8/14/14 SV

Scanned: 8/18/14 ✓

318140

ID: LCMPPFOS\_00006  
 Exp: 03/10/19 Prpd: SKV  
 13C4-Perfluorooctanesulfonate

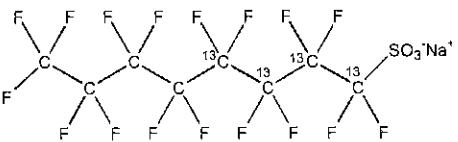


**WELLINGTON**  
LABORATORIES

**CERTIFICATE OF ANALYSIS**  
**DOCUMENTATION**

**PRODUCT CODE:**

MPFOS

**LOT NUMBER:** MPFOS0314**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)

**ISOTOPIC PURITY:** ≥99% <sup>13</sup>C**CHEMICAL PURITY:**

&gt;98%

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

03/10/2014

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

03/10/2019

**RECOMMENDED STORAGE:** Store ampoule in a cool, dark place**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 ~ 0.5% 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: 03/19/2014

(mm/dd/yyyy)

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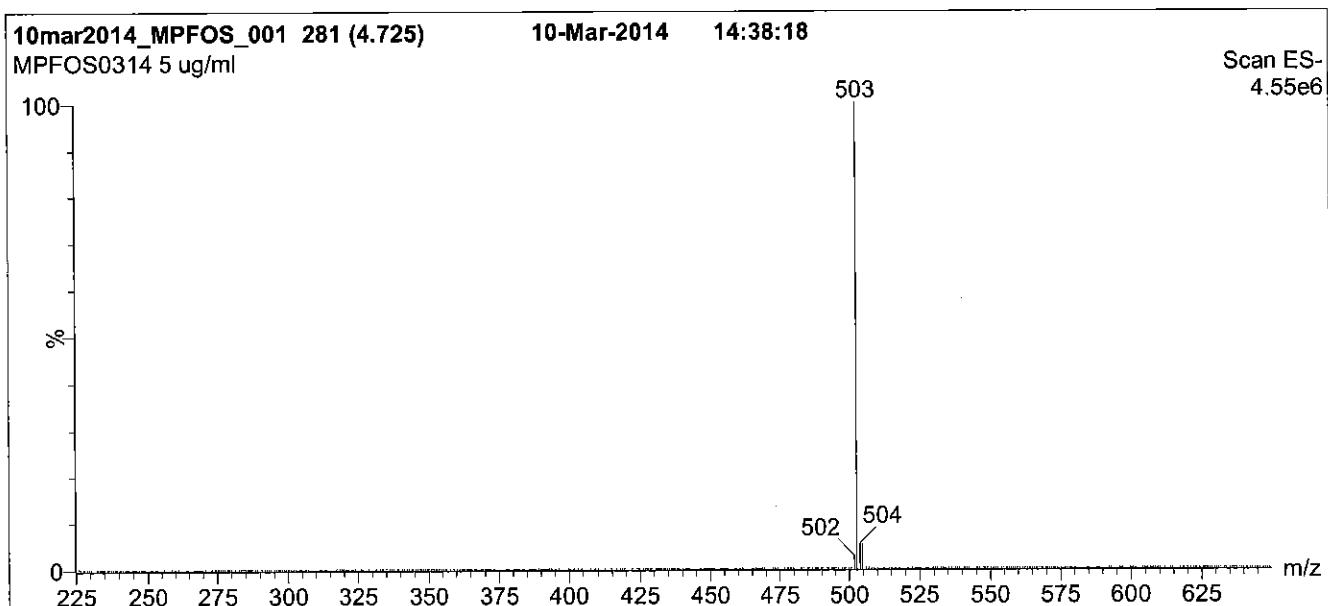
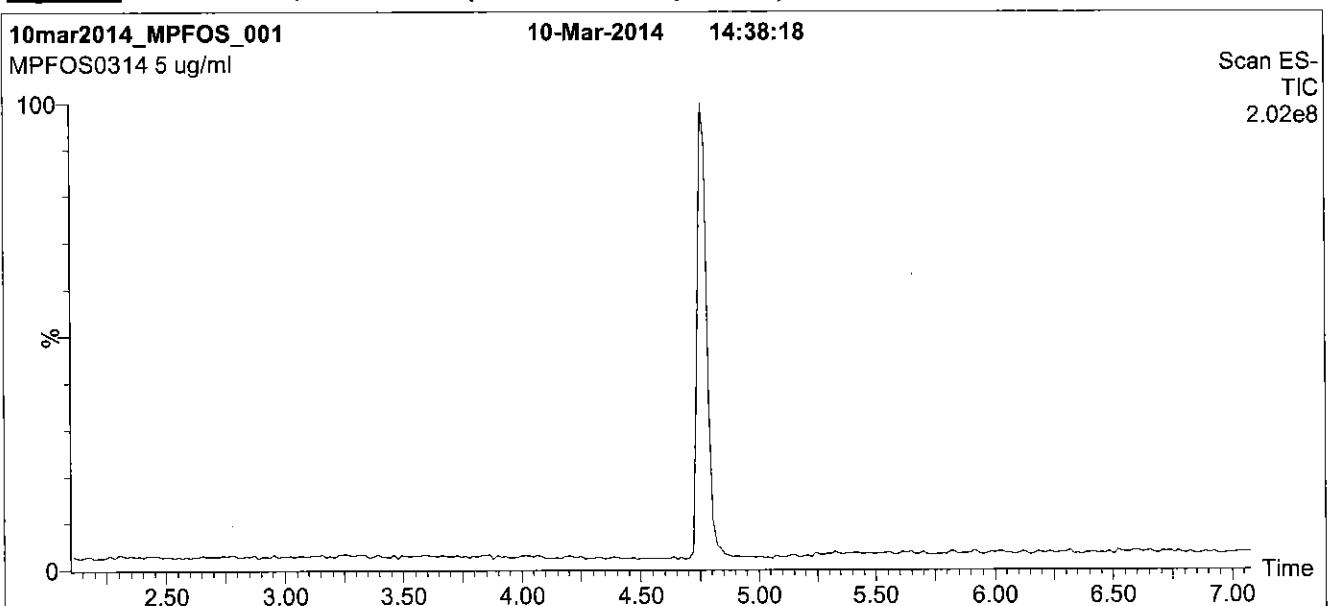
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**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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

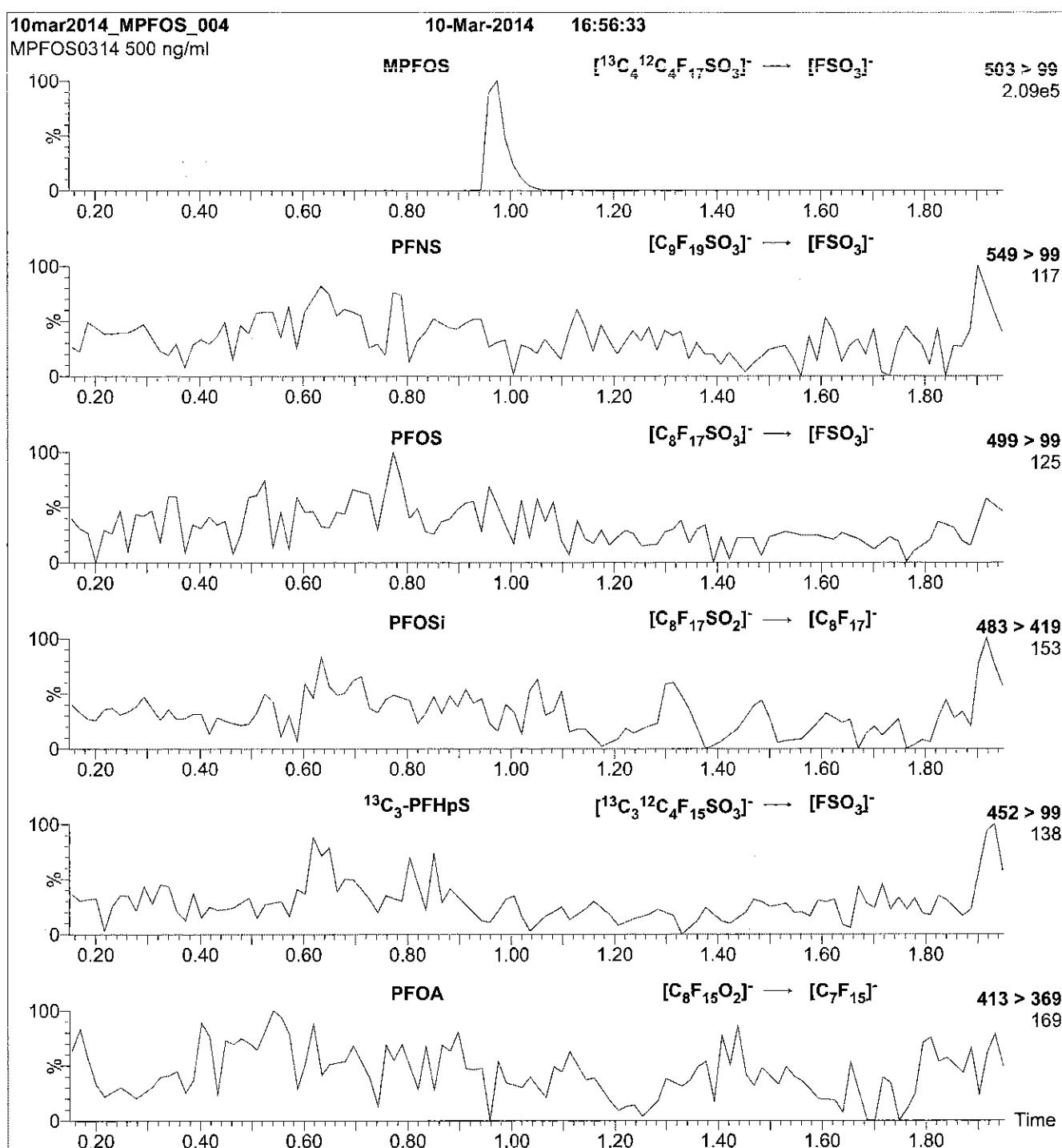
Flow: 300 µl/min

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.46e-3  
Collision Energy (eV) = 40

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  $\mu\text{l}/\text{min}$

Reagent

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**LCMPFUdA  00003**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE:

MPFUdA

LOT NUMBER: MPFUdA0911

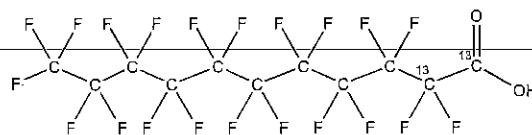
COMPOUND:

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

CHEMICAL PURITY:

>98%

ISOTOPIC PURITY: ≥99% <sup>13</sup>C

LAST TESTED: (mm/dd/yyyy)

09/29/2011

Water (<1%)

EXPIRY DATE: (mm/dd/yyyy)

09/29/2016

(1,2-<sup>13</sup>C<sub>2</sub>)

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-PFUdA (~1%; see Figure 2), 2-<sup>13</sup>C-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: 01/09/2013

(mm/dd/yyyy)

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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.

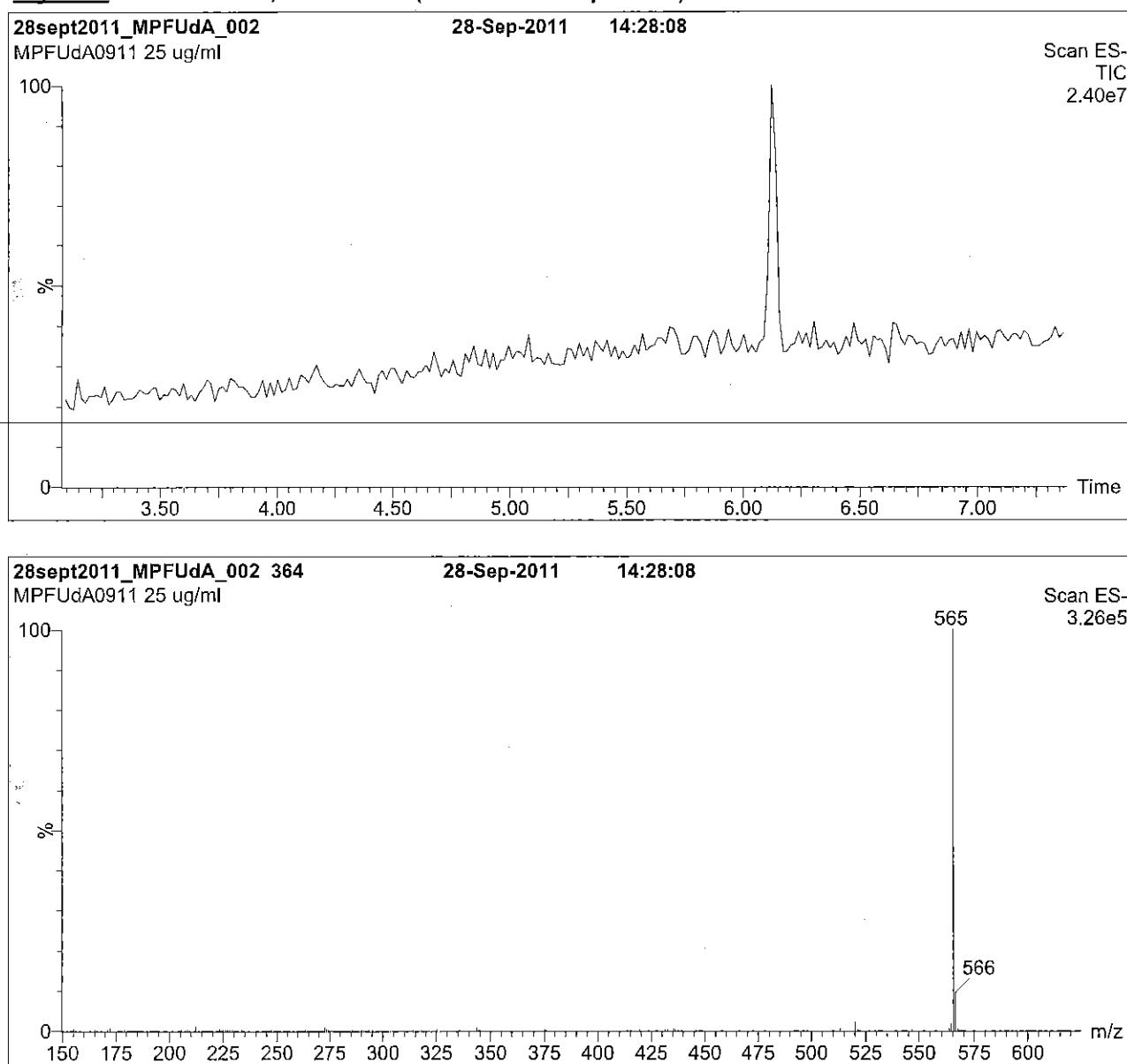
**QUALITY MANAGEMENT:**

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 ACCLASS (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: Kinetex PFP  
2.6 µm, 4.6 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)  
Hold for 2 min. 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

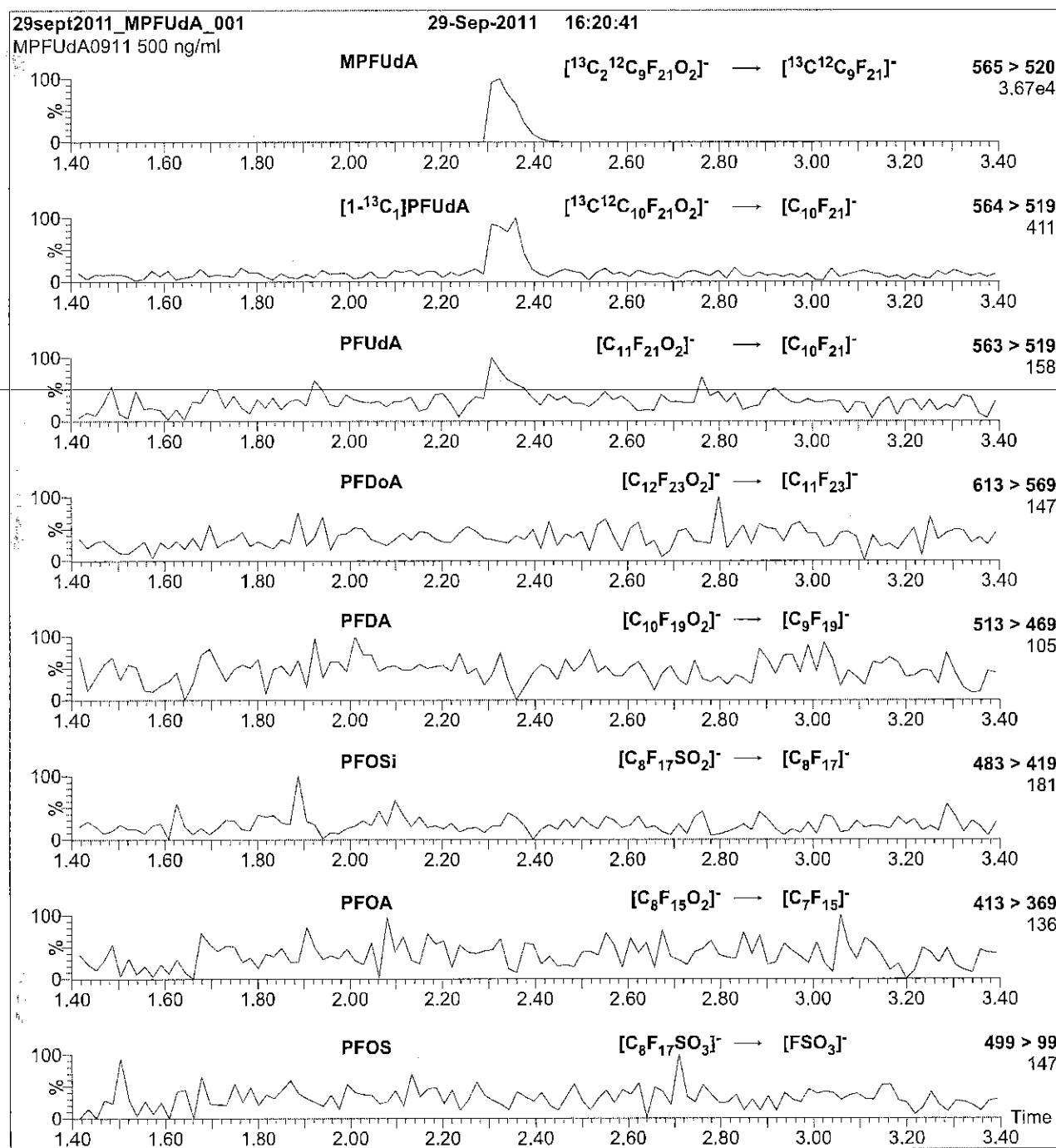
Flow: 1.0 ml/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 3.00  
Cone Voltage (V) = 15.00  
Cone Gas Flow (l/hr) = 65  
Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

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

**MS Parameters**

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

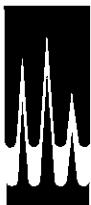
Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20%  $\text{H}_2\text{O}$   
(both with 10 mM  $\text{NH}_4\text{OAc}$  buffer)

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

Reagent

---

**LCPFACMXB\_00002**



**WELLINGTON  
LABORATORIES**

**CERTIFICATE OF ANALYSIS  
DOCUMENTATION**

R: 12-14-12

LCPFC MXB\_00002

**PFAC-MXB**

**Solution/Mixture of Native  
Perfluoroalkylcarboxylic Acids and  
Native Perfluoroalkylsulfonates**

**PRODUCT CODE:**

PFAC-MXB

**LOT NUMBER:**

PFACMXB0312

**SOLVENT(S):**

Methanol / Water (<1%)

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

03/07/2012

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

03/08/2012

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

03/08/2015

**RECOMMENDED STORAGE:**

Store ampoule in a cool, dark place

**DESCRIPTION:**

PFAC-MXB is a solution/mixture of thirteen native perfluoroalkylcarboxylic acids ( $C_4$ - $C_{14}$ ,  $C_{16}$ , and  $C_{18}$ ) and four native perfluoroalkylsulfonates ( $C_4$ ,  $C_6$ ,  $C_8$  and  $C_{10}$ ). The full name, abbreviation and concentration for each of the components are given in Table A.

The individual perfluoroalkylcarboxylic acids and perfluoroalkylsulfonates all have chemical purities of >98%.

**DOCUMENTATION/ DATA ATTACHED:**

Table A: Components and Concentrations of the Solution/Mixture

Figure 1: LC/MS Data (SIR)

Figure 2: LC/MS/MS data (Selected MRM Transitions)

Figure 3: LC/MS/MS data (Selected MRM Transitions)

**ADDITIONAL INFORMATION:**

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

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

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

**INTENDED USE:**

The products prepared by Wellington Laboratories Inc. are for laboratory use only. 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.

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**Table A: PFAC-MXB; Components and Concentrations (ng/ml, ± 5% in Methanol / Water (<1%))**

Name	Abbreviation	Concentration (ng/ml)		Peak Assignment in Figure 1
Perfluoro-n-butanoic acid	PFBA	2000		A
Perfluoro-n-pentanoic acid	PPPeA	2000		B
Perfluoro-n-hexanoic acid	PFHxA	2000		C
Perfluoro-n-heptanoic acid	PFHpA	2000		D
Perfluoro-n-octanoic acid	PFOA	2000		E
Perfluoro-n-nonanoic acid	PFNA	2000		F
Perfluoro-n-decanoic acid	PFDA	2000		G
Perfluoro-n-undecanoic acid	PFUdA	2000		H
Perfluoro-n-dodecanoic acid	PFDoA	2000		I
Perfluoro-n-tridecanoic acid	PFTrDA	2000		J
Perfluoro-n-tetradecanoic acid	PFTeDA	2000		K
Perfluoro-n-hexadecanoic acid	PFHxDA	2000		L
Perfluoro-n-octadecanoic acid	PFODA	2000		M
Name	Abbreviation	Concentration (ng/ml)		Peak Assignment in Figure 1
		as the salt	as the anion	
Potassium perfluoro-1-butanesulfonate	L-PFBS	2000	1770	N
Sodium perfluoro-1-hexamersulfonate	L-PFHxS	2000	1890	O
Sodium perfluoro-1-octanesulfonate	L-PFOS	2000	1910	P
Sodium perfluoro-1-decanesulfonate	L-PFDS	2000	1930	Q

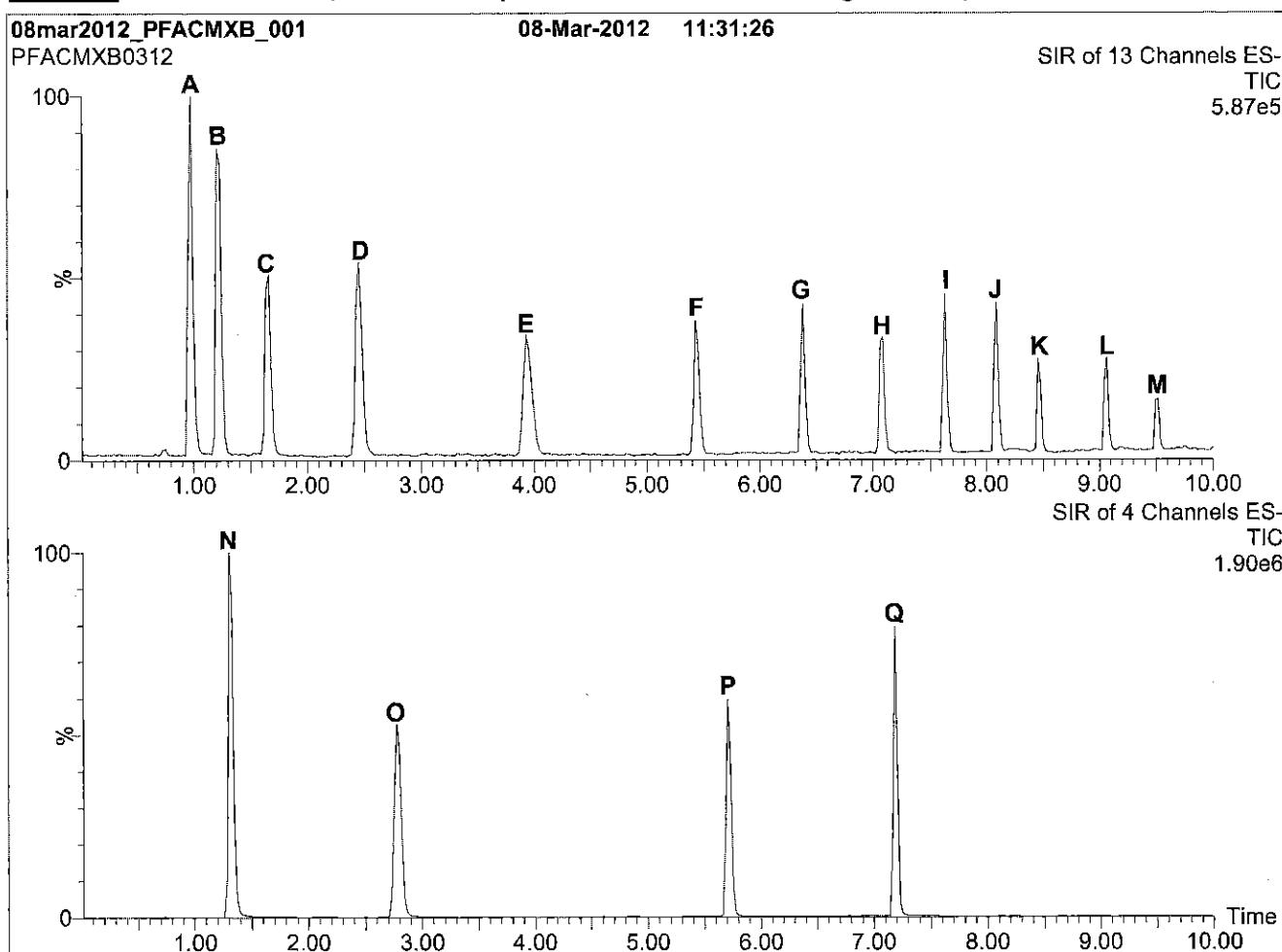
Certified By:

  
B.G. Chittim

Date: 03/19/2012

(mm/dd/yyyy)

**Figure 1:** PFAC-MXB; LC/MS Data (Total Ion Current Chromatogram; SIR)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
 1.7 µm, 2.1 x 100 mm

Mobile phase: Gradient  
 Start: 50% H<sub>2</sub>O / 50% (80:20 MeOH:ACN)  
 (both with 10 mM NH<sub>4</sub>OAc buffer)  
 Hold for 3 min; ramp to 100% organic over 7 min,  
 hold for 2 min, then return to initial conditions in 0.5 min.

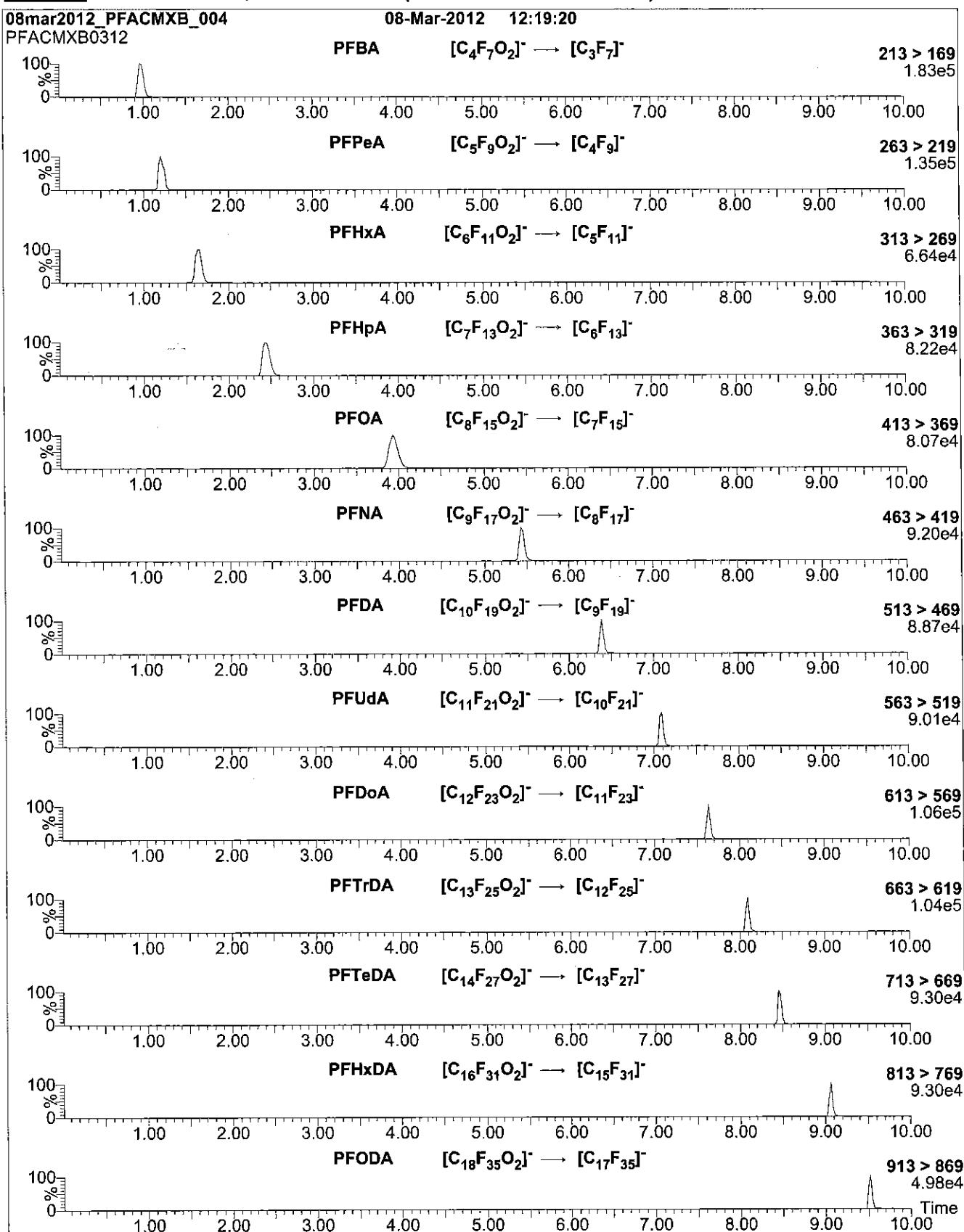
Time: 15 min

Flow: 300 µl/min

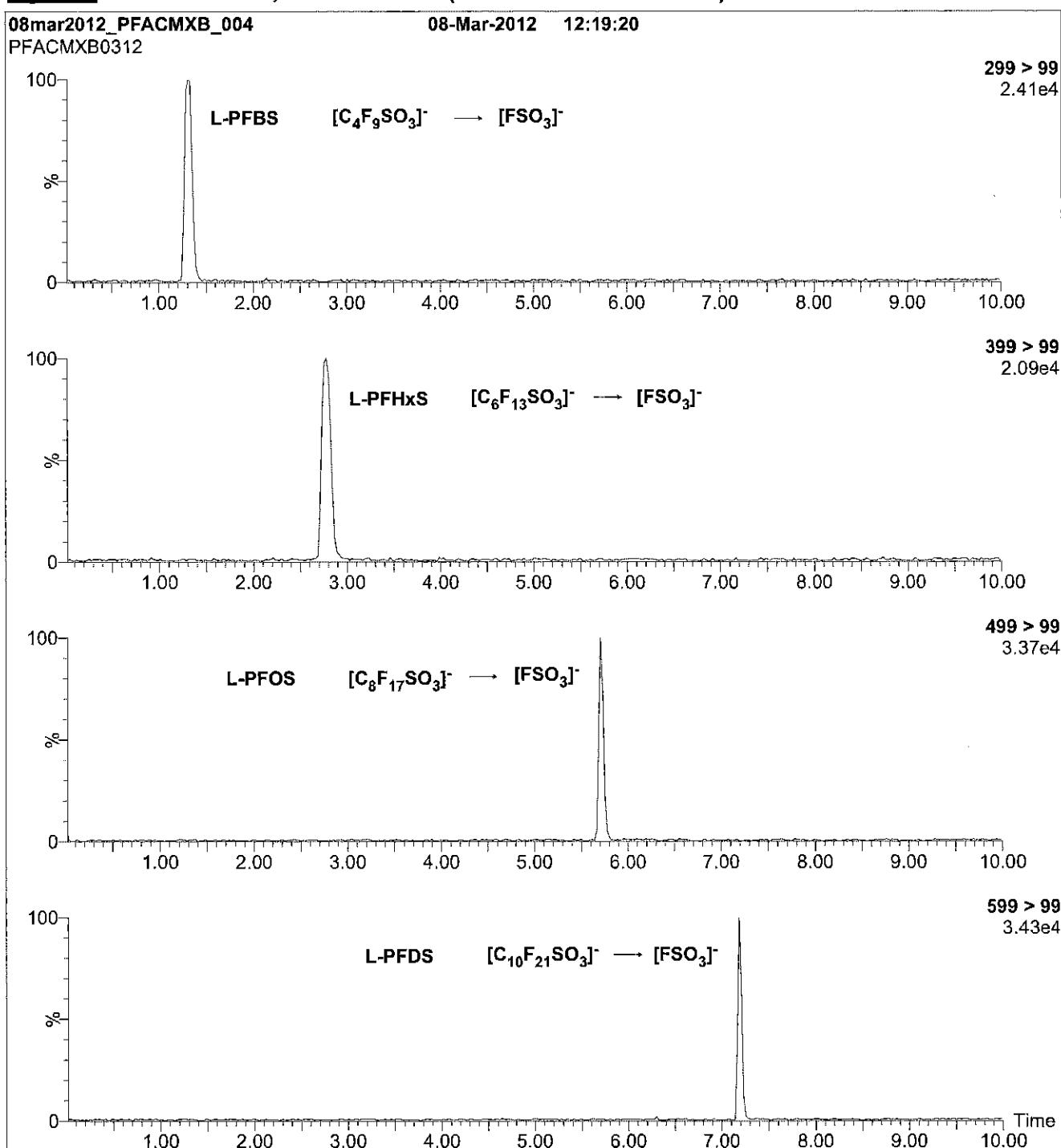
**MS Parameters**

Experiment:  
 Function 1: SIR of 13 Channels  
 Function 2: SIR of 4 Channels  
 Source: Electrospray (negative)  
 Capillary Voltage (kV) = 2.50  
 Cone Voltage (V) = variable (10-70)  
 Cone Gas Flow (l/hr) = 100  
 Desolvation Gas Flow (l/hr) = 750

**Figure 2: PFAC-MXB; LC/MS/MS Data (Selected MRM Transitions)**



**Figure 3: PFAC-MXB; LC/MS/MS Data (Selected MRM Transitions)**



**Conditions for Figures 2 and 3:**

Injection: on-column (PFAC-MXB)

**MS Parameters**

Collision Gas (mbar) = 3.17e-3

Mobile phase: Same as Figure 1

Collision Energy (eV) = 9-50 (variable)

Flow: 300  $\mu$ l/min

Reagent

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**LCPFBA\_00002**

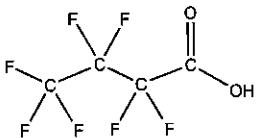


# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE: PFBA      LOT NUMBER: PFBA0212  
COMPOUND: Perfluoro-n-butanoic acid

STRUCTURE:      CAS #: 375-22-4



<u>MOLECULAR FORMULA:</u>	C <sub>4</sub> H <sub>7</sub> F <sub>7</sub> O <sub>2</sub>	<u>MOLECULAR WEIGHT:</u>	214.04
<u>CONCENTRATION:</u>	50 ± 2.5 µg/ml	<u>SOLVENT(S):</u>	Methanol Water (<1%)
<u>CHEMICAL PURITY:</u>	>98%		
<u>LAST TESTED:</u> (mm/dd/yyyy)	02/22/2012		
<u>EXPIRY DATE:</u> (mm/dd/yyyy)	02/22/2015		
<u>RECOMMENDED STORAGE:</u>	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: 02/23/2012

(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:**

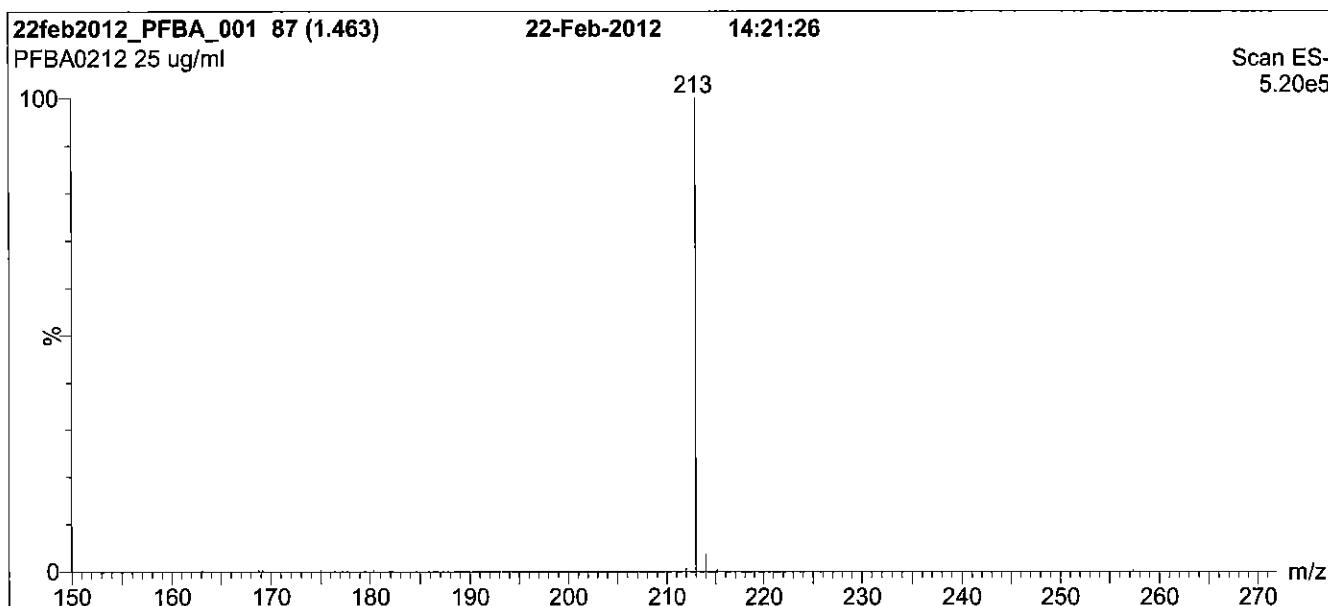
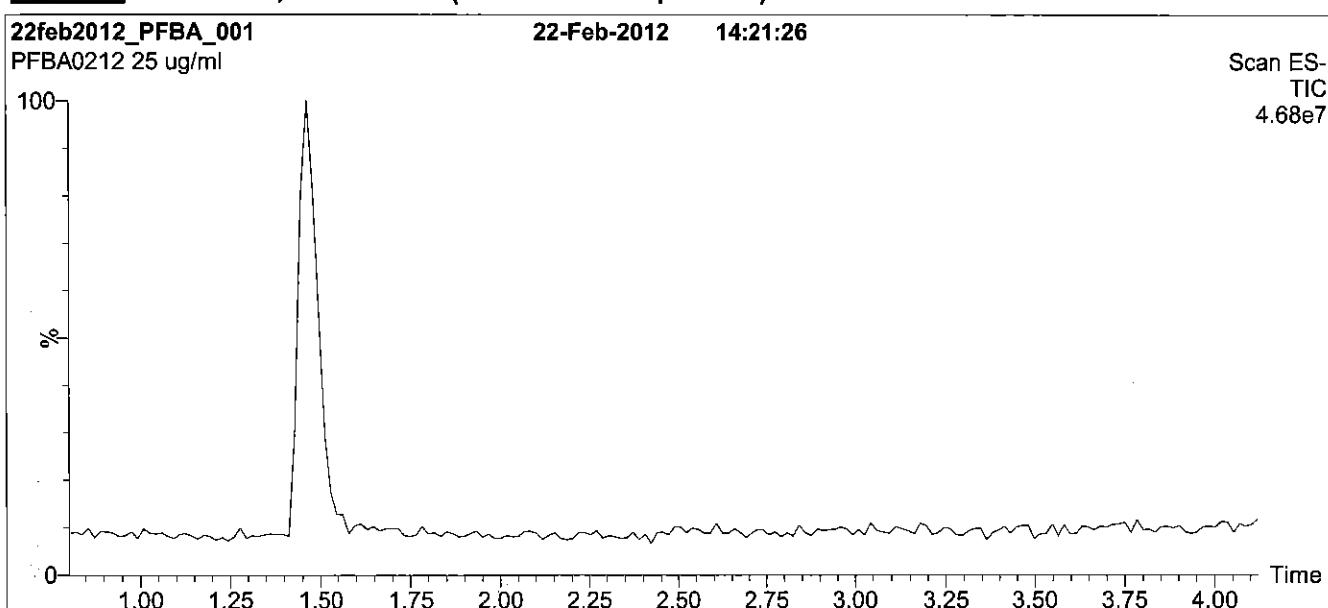
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

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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**

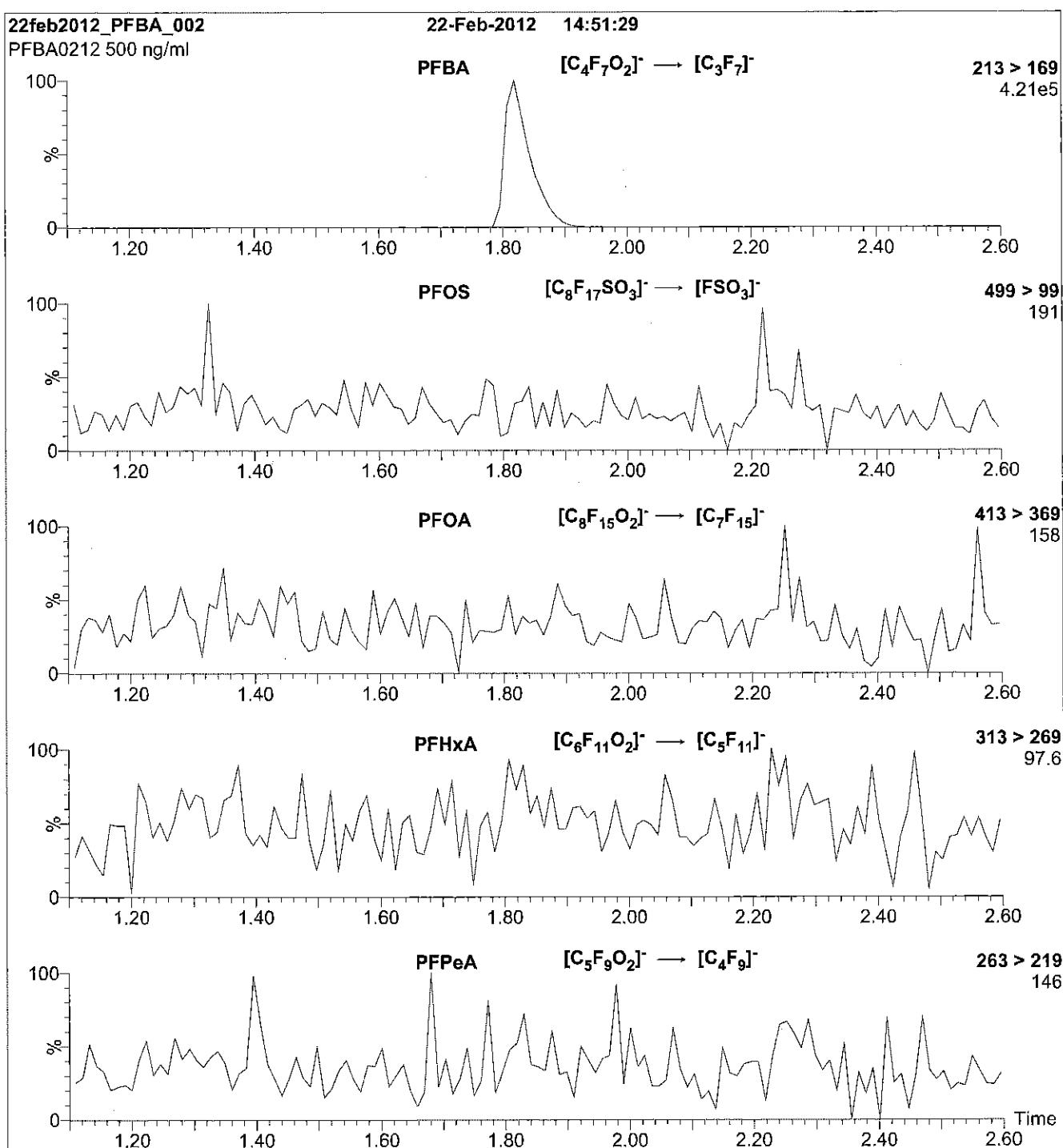
Column: Acquity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 2.1 x 100 mm

Mobile phase: Gradient  
Start: 30% (80:20 MeOH:ACN) / 70% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc)  
Ramp to 90% organic over 7.5 min and hold for 1.5 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min  
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)  
Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 10.00  
Cone Gas Flow (l/hr) = 60  
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)

**MS Parameters**

Collision Gas (mbar) = 3.54e-3  
Collision Energy (eV) = 10

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

Flow: 300  $\mu$ l/min

Reagent

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**LCPFBS\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LC PFBS - 00002

PRODUCT CODE:

L-PFBS

LOT NUMBER: LPFBS0612

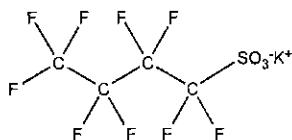
COMPOUND:

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)

06/21/2012

EXPIRY DATE: (mm/dd/yyyy)

06/21/2015

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: 07/09/2012

(mm/dd/yyyy)

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

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

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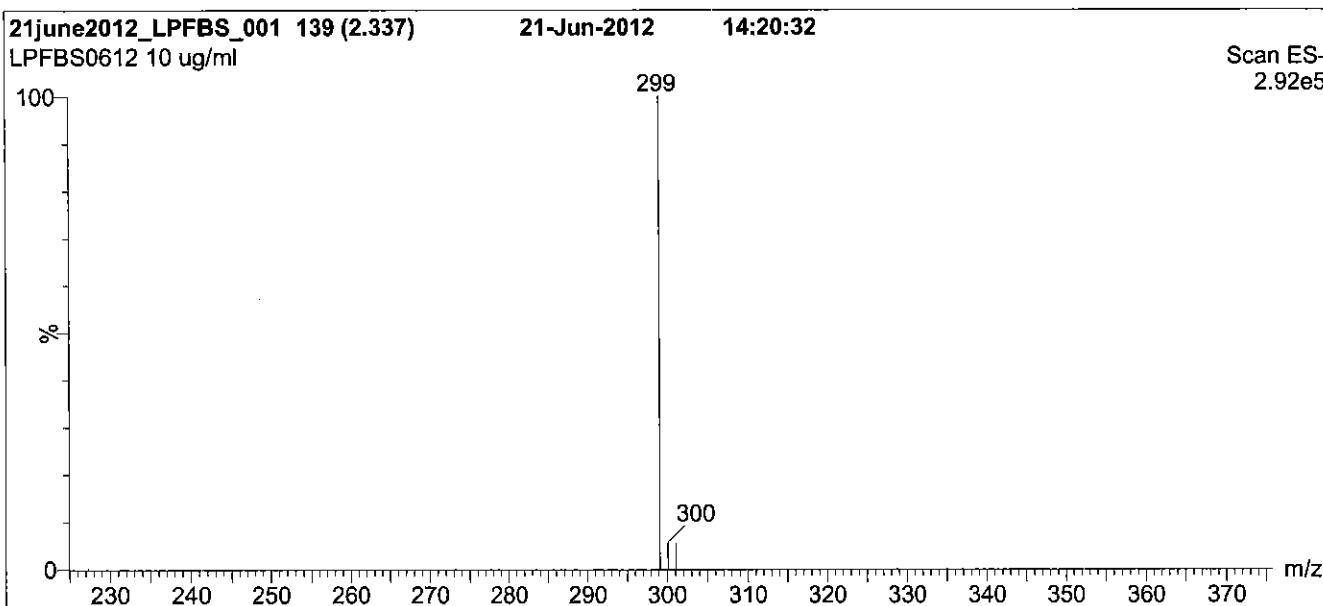
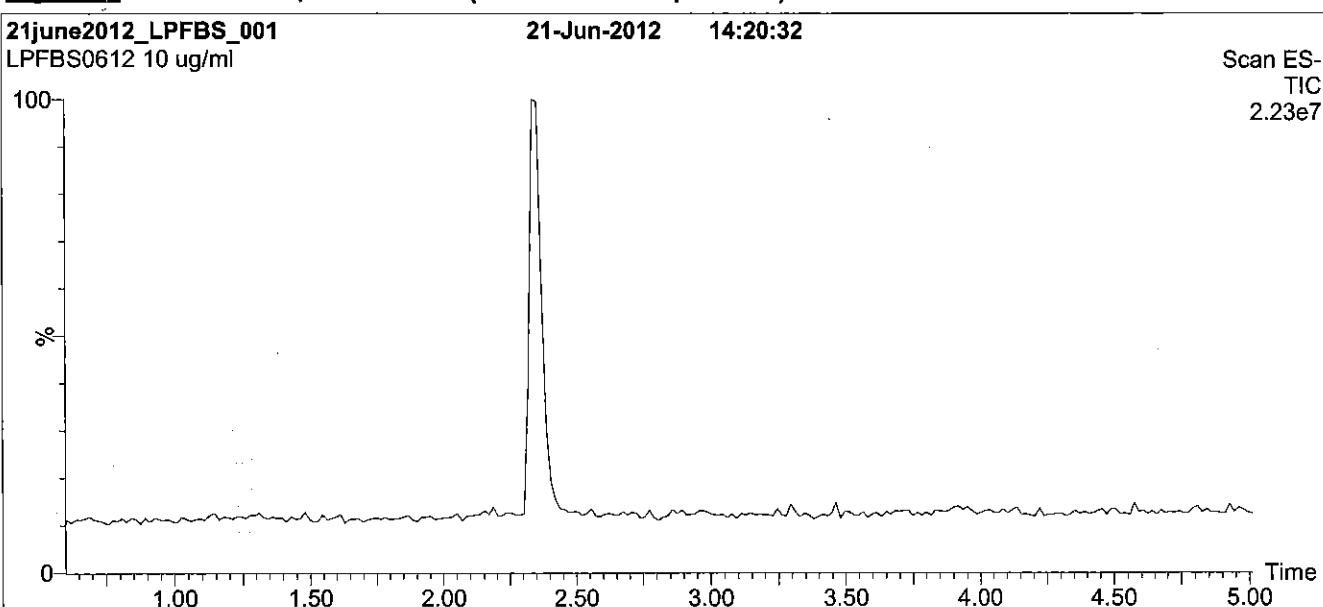
**QUALITY MANAGEMENT:**

This product was produced using a Quality Management System registered to 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 ACCLASS (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-PFBS; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity 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.

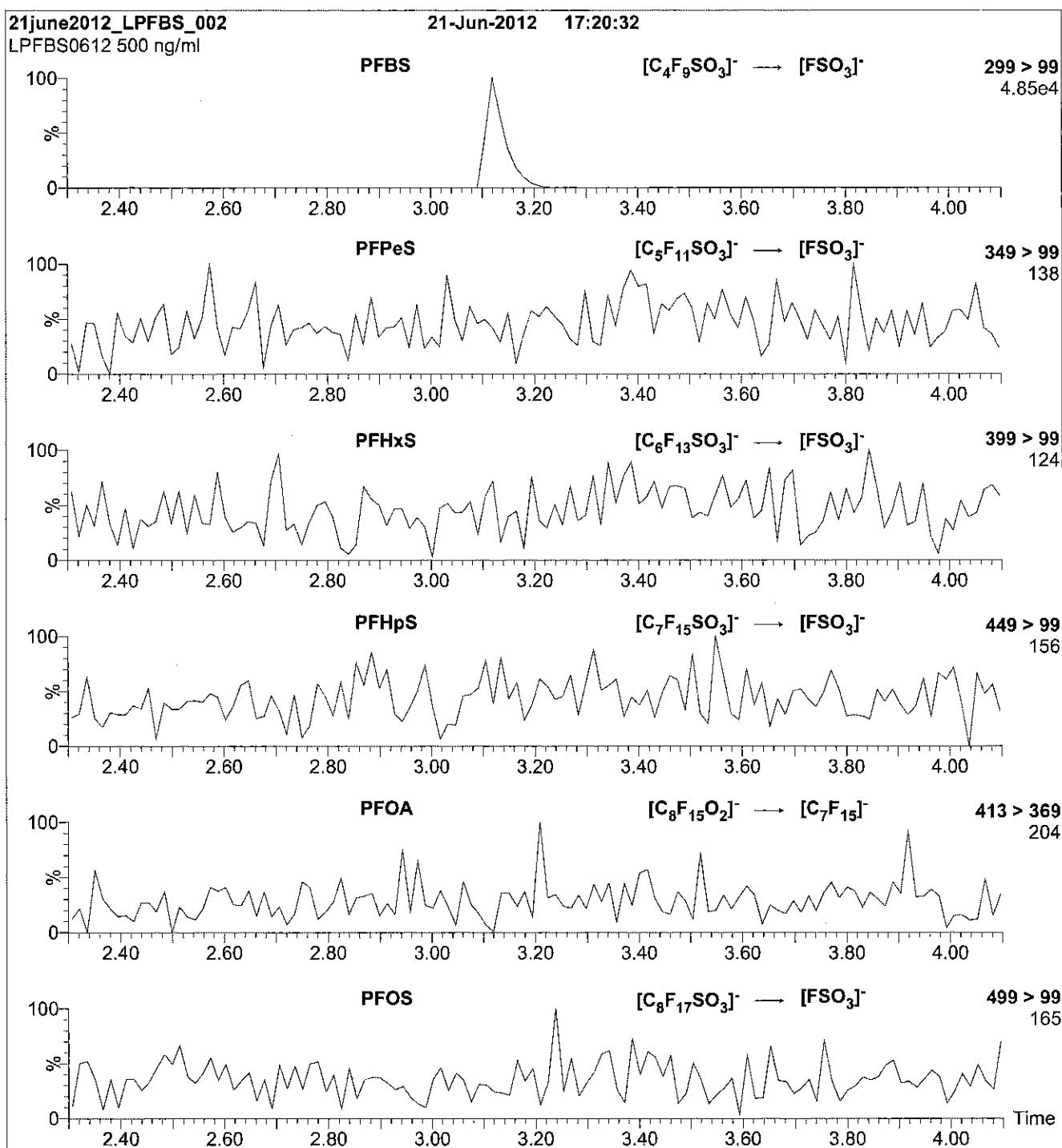
Time: 10 min  
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (225 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 40.00  
Cone Gas Flow (l/hr) = 50  
Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

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

**MS Parameters**

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

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

Reagent

---

**LCPFDA\_00003**

rec 7/16/14

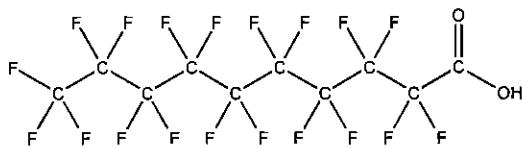


# WELLINGTON LABORATORIES

## 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>HF<sub>19</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:

A handwritten signature in black ink, appearing to read "B.G. Chittim".

Date: 07/03/2013

(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. 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.

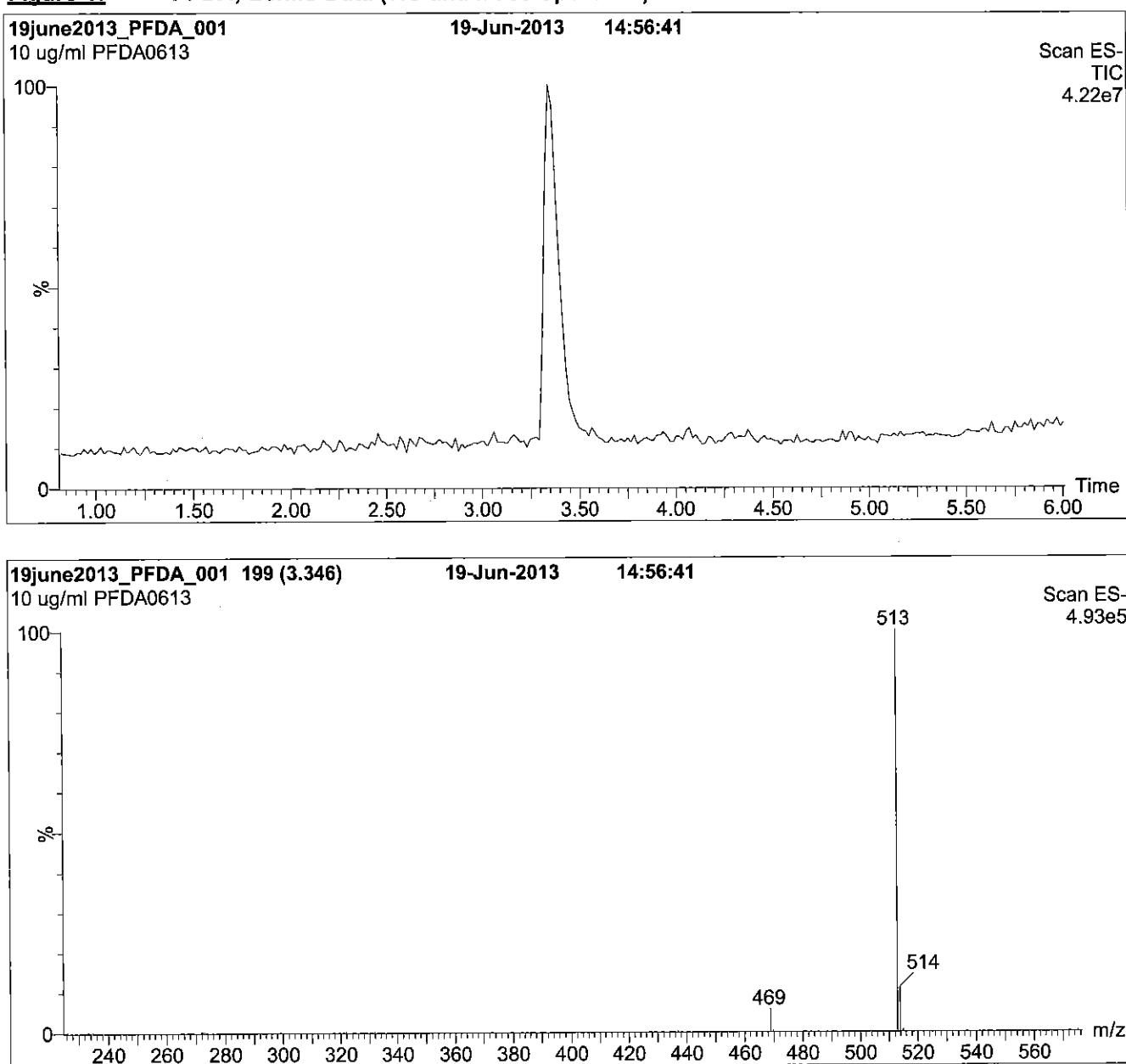
### **QUALITY MANAGEMENT:**

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 ACCLASS (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 Acuity Ultra Performance LC  
**MS:** Micromass Quattro *micro* API MS

**Chromatographic Conditions**

Column: Acuity 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 90% organic over 7 min and hold for  
1.5 min before returning to initial conditions in 0.5 min.  
Time: 10 min

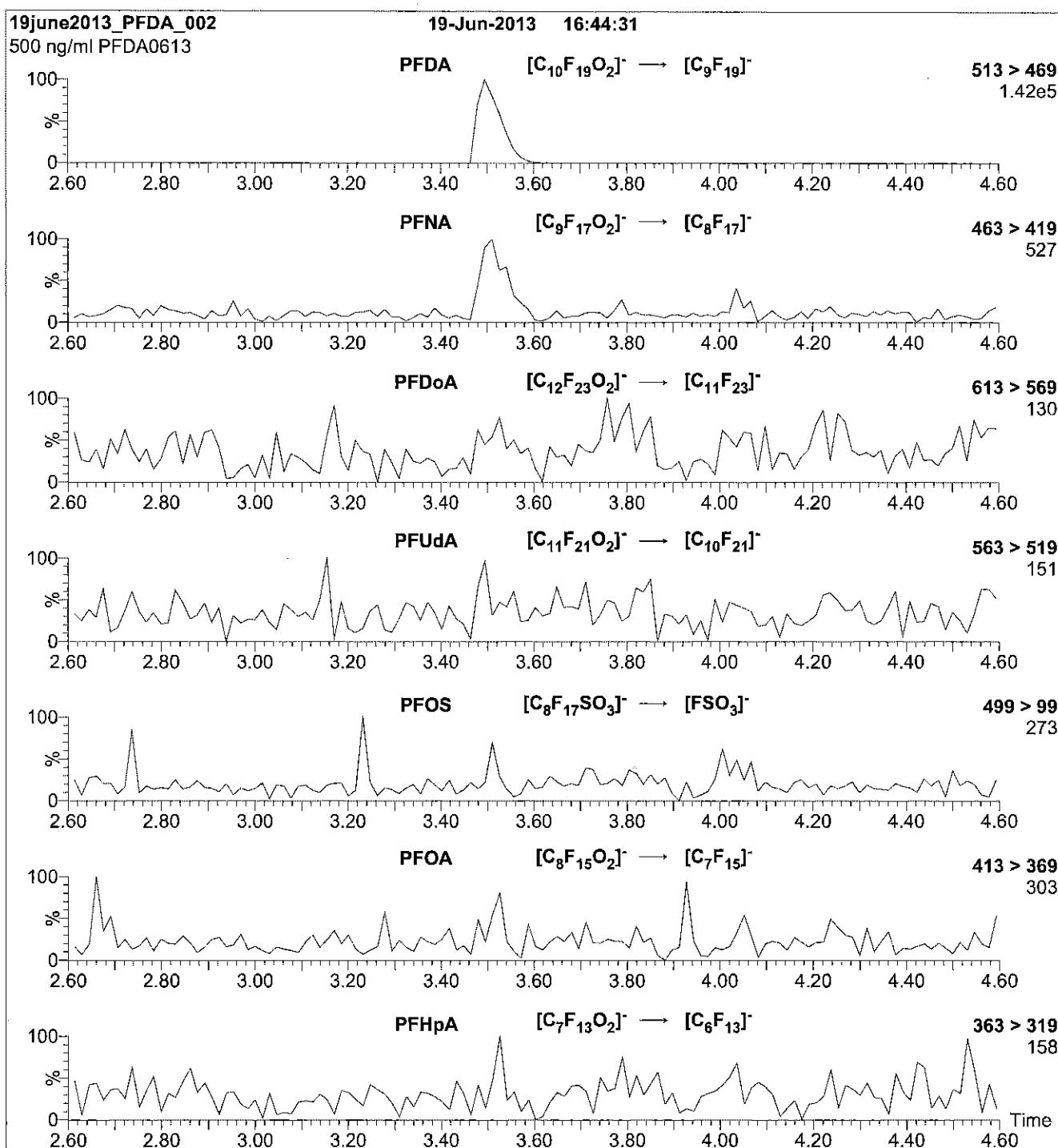
Flow: 300 µl/min

**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)

**MS Parameters**

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

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

Reagent

---

**LCPFD**DoA**\_00003**



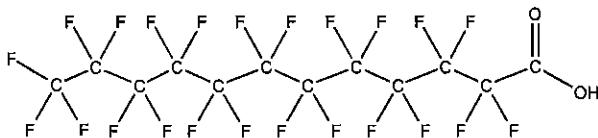
# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Rec Thulis

PRODUCT CODE: PFDoA      LOT NUMBER: PFDoA0113  
COMPOUND: Perfluoro-n-dodecanoic acid

STRUCTURE:      CAS #: 307-55-1



MOLECULAR FORMULA: C<sub>12</sub>HF<sub>23</sub>O<sub>2</sub>      MOLECULAR WEIGHT: 614.10  
CONCENTRATION: 50 ± 2.5 µg/ml      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:

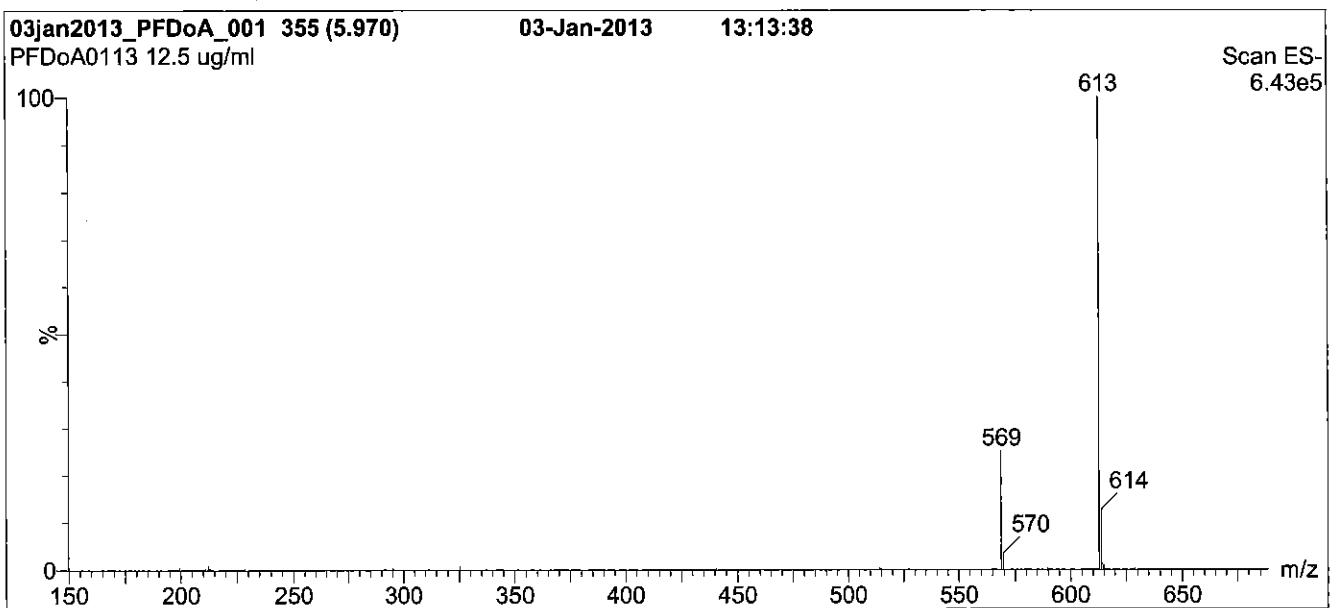
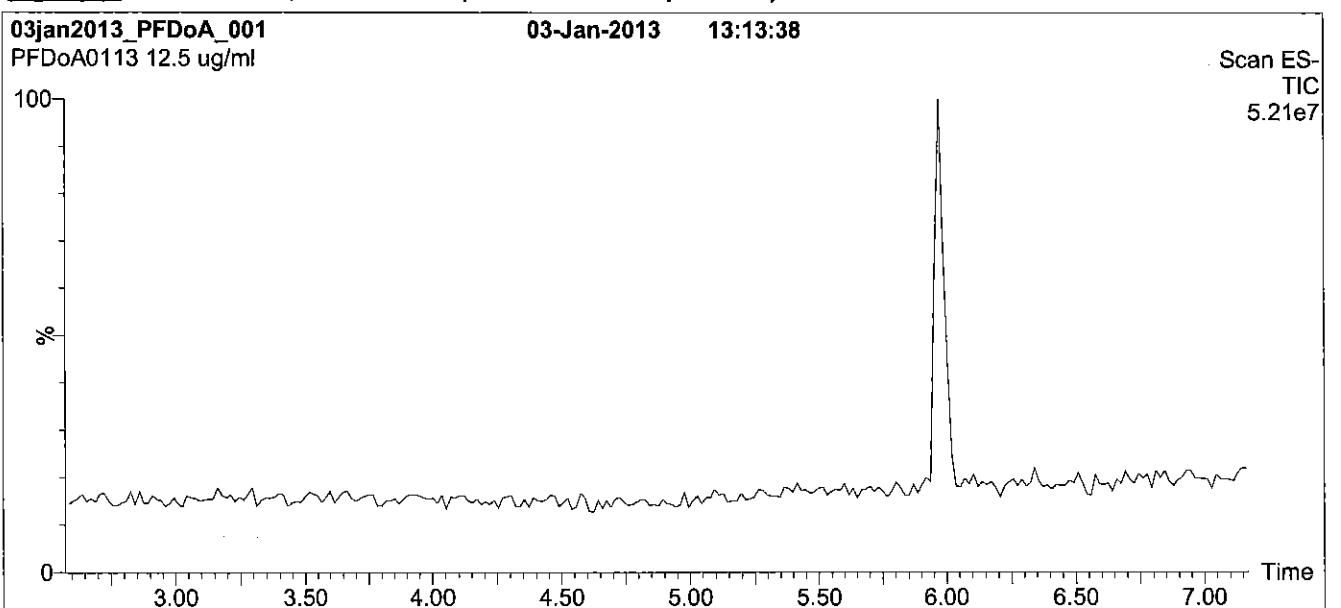
B.G. Chittim

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: PFDoA; 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\_00002**

Rec His/14



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

**PRODUCT CODE:**

L-PFDoS

**LOT NUMBER:** LPFDoS1011

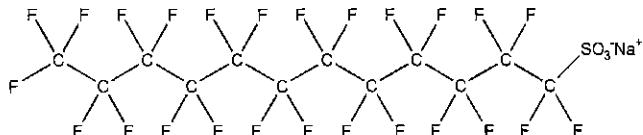
**COMPOUND:**

Sodium perfluoro-1-dodecanesulfonate

**STRUCTURE:**

**CAS #:**

Not available



**MOLECULAR FORMULA:**

$\text{C}_{12}\text{F}_{25}\text{SO}_3\text{Na}$

**MOLECULAR WEIGHT:** 722.14

**CONCENTRATION:**

$50.0 \pm 2.5 \mu\text{g/ml}$  (Na salt)

**SOLVENT(S):** Methanol

$48.4 \pm 2.4 \mu\text{g/ml}$  (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:

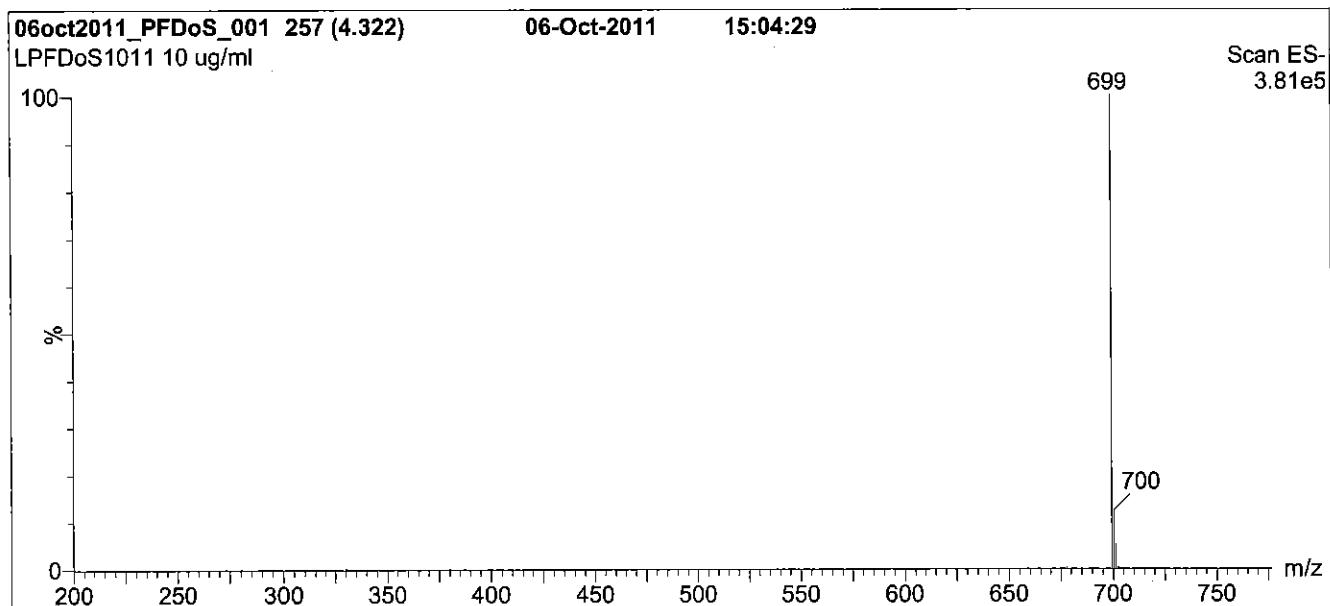
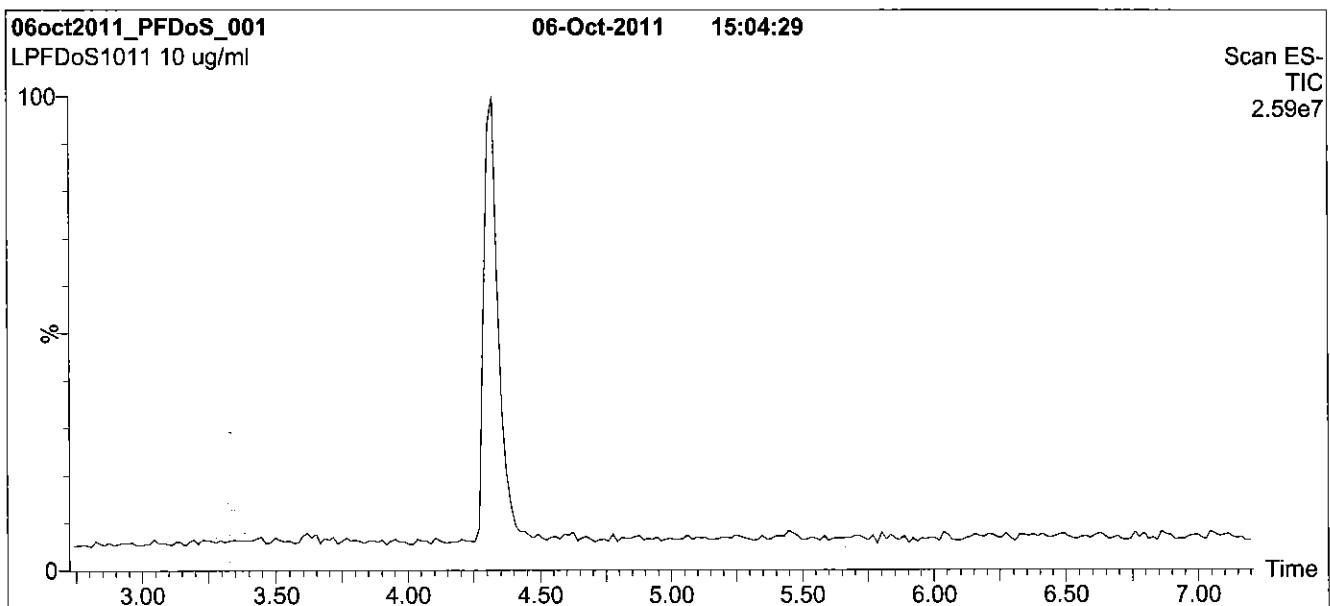
B.G. Chittim

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

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



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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 µl/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

Reagent

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**LCPFDS\_00002**



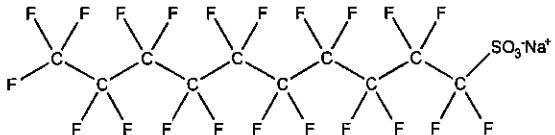
# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R 12-14-12  
LPFDS\_00002

PRODUCT CODE: L-PFDS      LOT NUMBER: LPFDS0612  
COMPOUND: Sodium perfluoro-1-decanesulfonate

STRUCTURE:      CAS #: Not available



MOLECULAR FORMULA: C<sub>10</sub>F<sub>21</sub>SO<sub>3</sub>Na      MOLECULAR WEIGHT: 622.13  
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt)      SOLVENT(S): Methanol  
48.2 ± 2.4 µg/ml (PFDS anion)  
CHEMICAL PURITY: >98%  
LAST TESTED: (mm/dd/yyyy) 06/21/2012  
EXPIRY DATE: (mm/dd/yyyy) 06/21/2015  
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: 07/12/2012

(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.

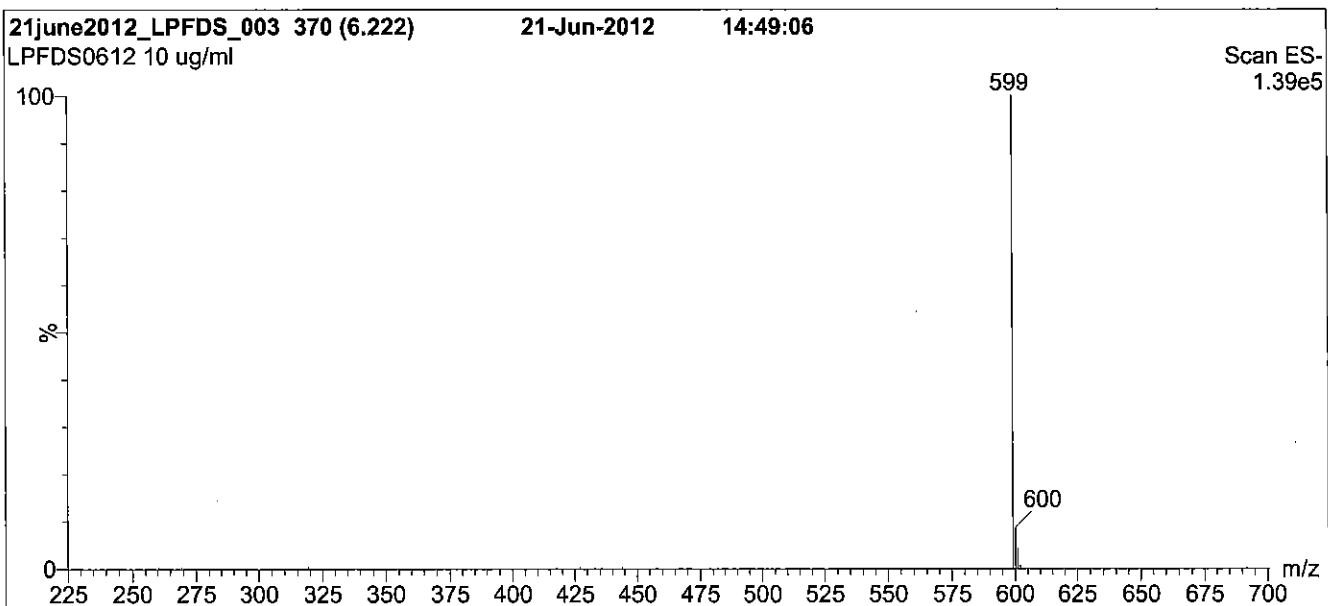
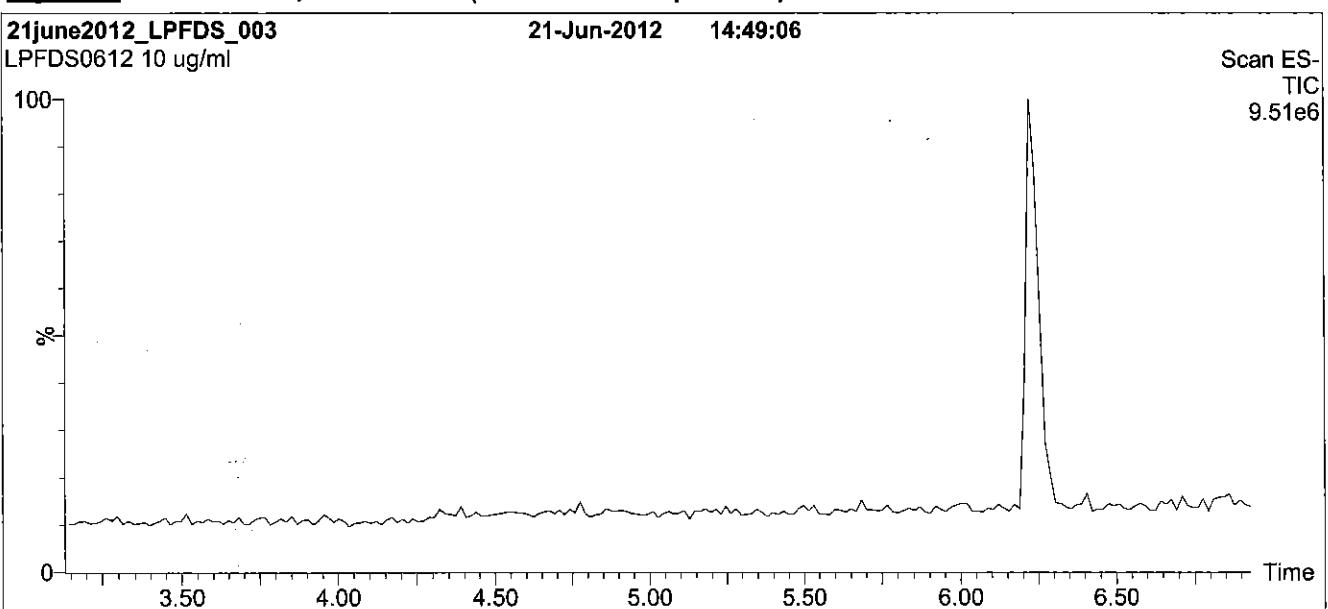
#### **QUALITY MANAGEMENT:**

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 ACCLASS (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-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 μm, 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

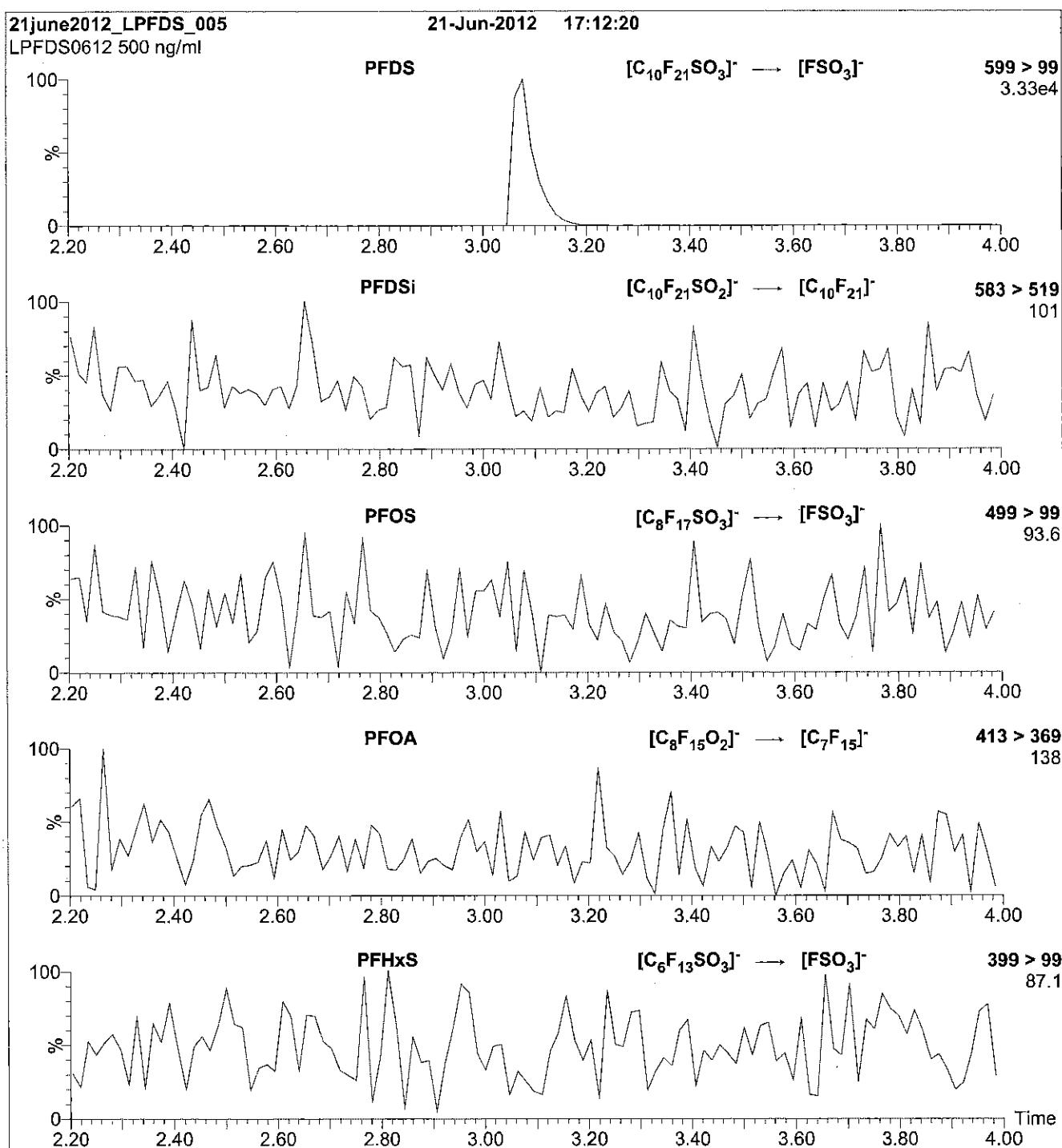
Flow: 300 μl/min

**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)

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 50

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

Reagent

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**LCPFHpA\_00002**

R 9-17-12 DE4



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Dup LC PFHpA-00001

new i.+ LC PFHpA -00002

PRODUCT CODE:

PFHpA

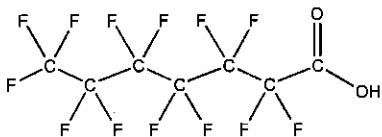
LOT NUMBER: PFHpA0412

COMPOUND:

Perfluoro-n-heptanoic acid

STRUCTURE:

CAS #: 375-85-9



MOLECULAR FORMULA:

C<sub>7</sub>HF<sub>13</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 364.06

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

04/18/2012

EXPIRY DATE: (mm/dd/yyyy)

04/18/2015

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: 05/03/2012

(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:**

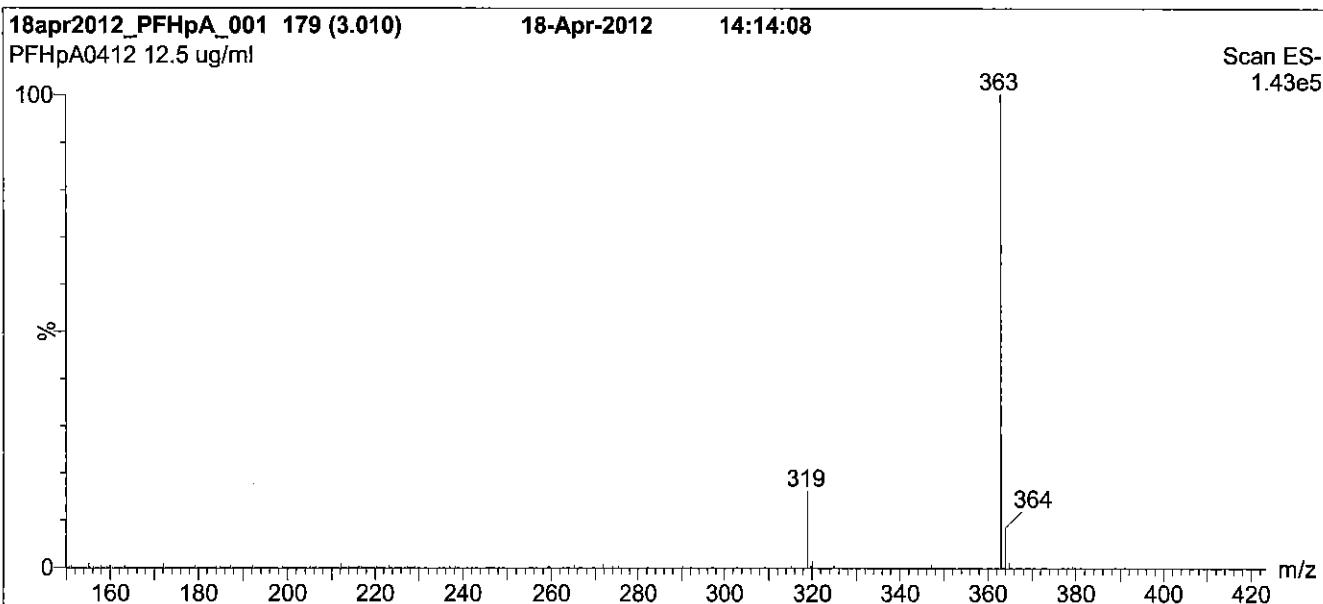
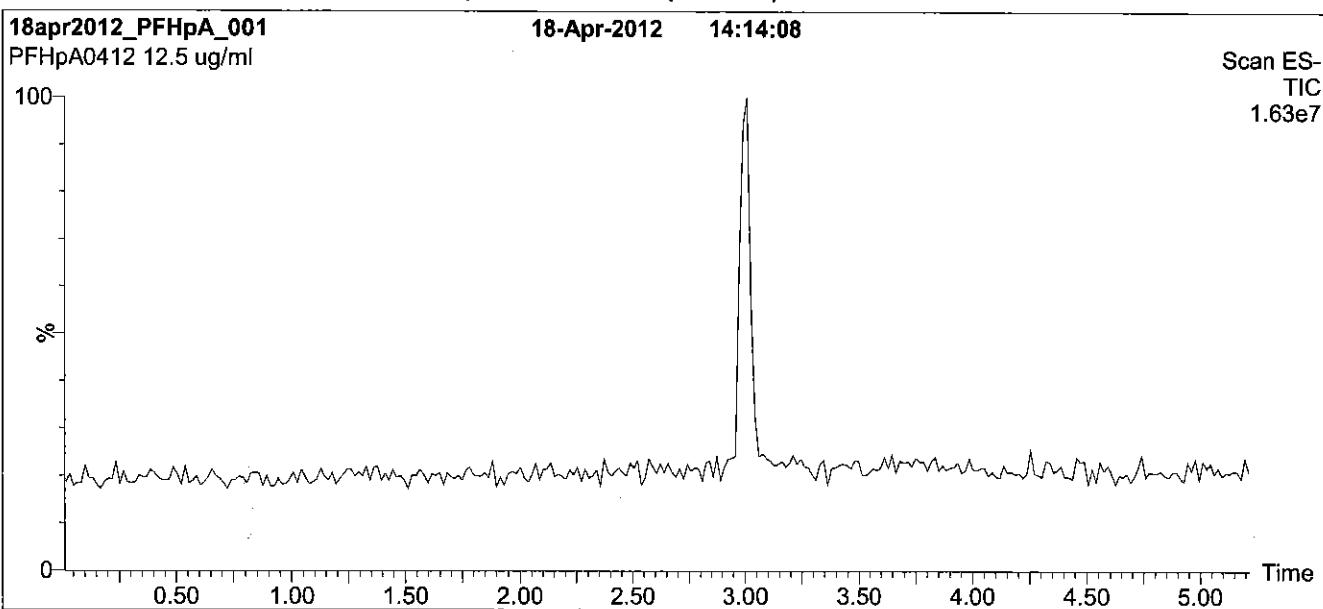
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**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.

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

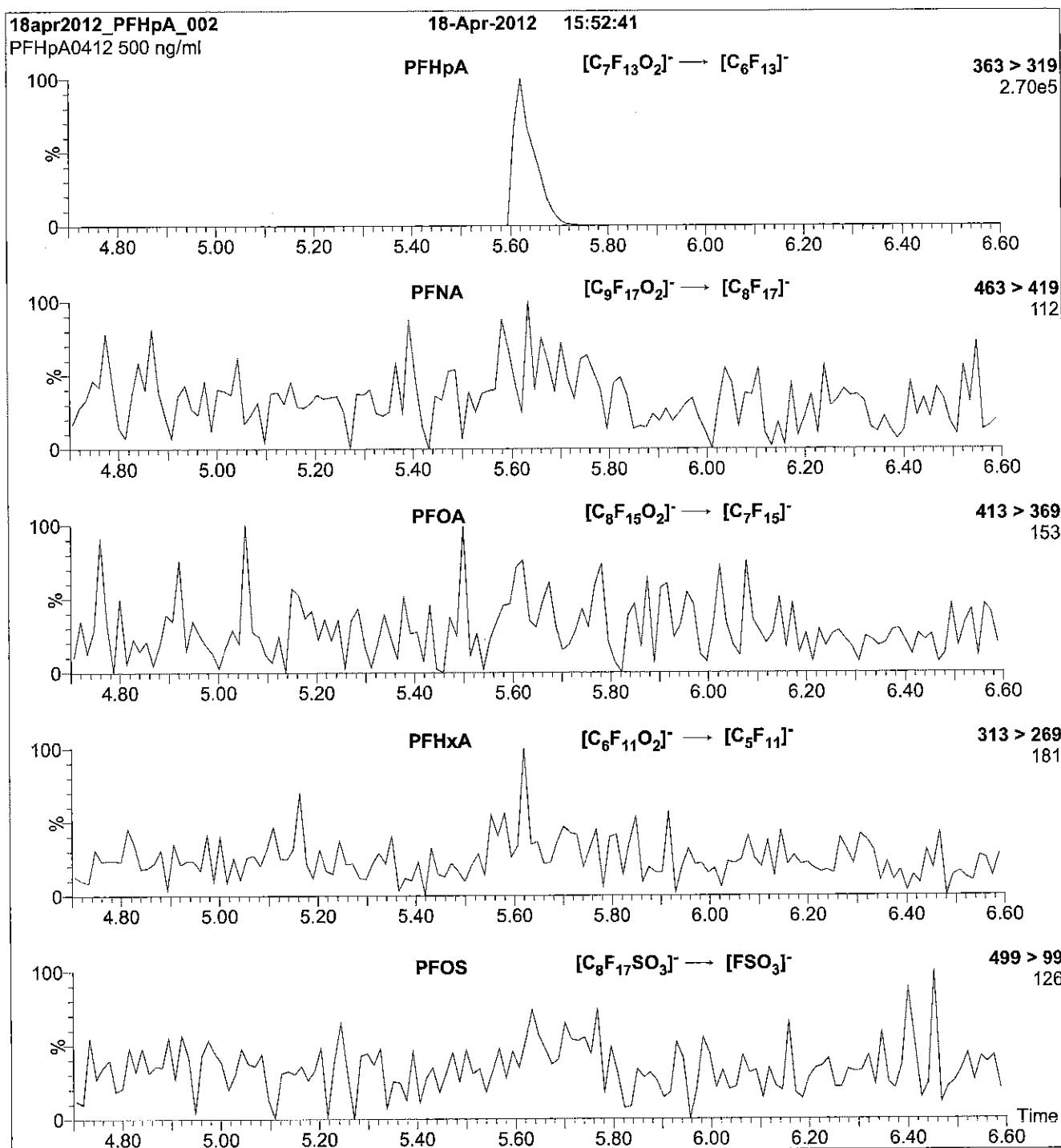
Mobile phase: Gradient  
Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min and hold for  
2 min before returning to initial conditions in 0.5 min.  
Time: 10 min

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) = 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)

**MS Parameters**

Collision Gas (mbar) = 3.54e-3  
Collision Energy (eV) = 11

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

Reagent

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**LCPFHxA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LC PFHxA - 00002

PRODUCT CODE:

PFHxA

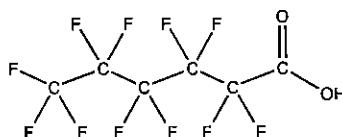
LOT NUMBER: PFHxA1012

COMPOUND:

Perfluoro-n-hexanoic acid

STRUCTURE:

CAS #: 307-24-4



MOLECULAR FORMULA:

C<sub>6</sub>HF<sub>11</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 314.05

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

10/29/2012

EXPIRY DATE: (mm/dd/yyyy)

10/29/2015

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

### DOCUMENTATION/ DATA ATTACHED:

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

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

### ADDITIONAL INFORMATION:

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

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

Certified By:

B.G. Chittim

Date: 11/01/2012

(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:**

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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.

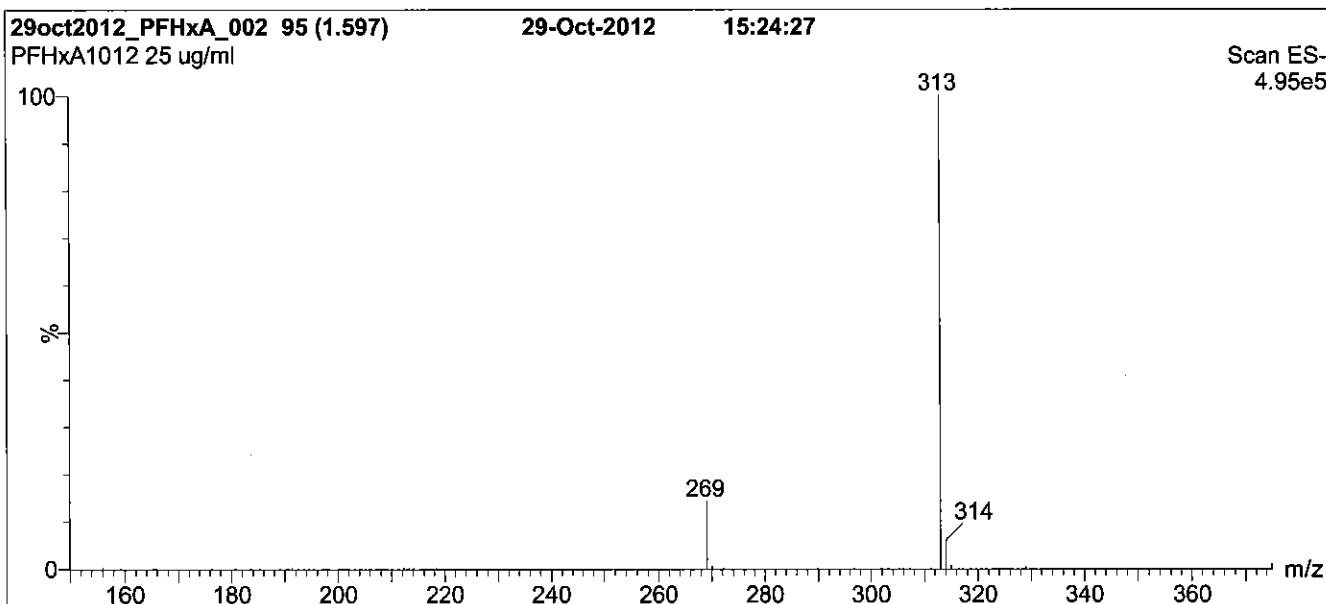
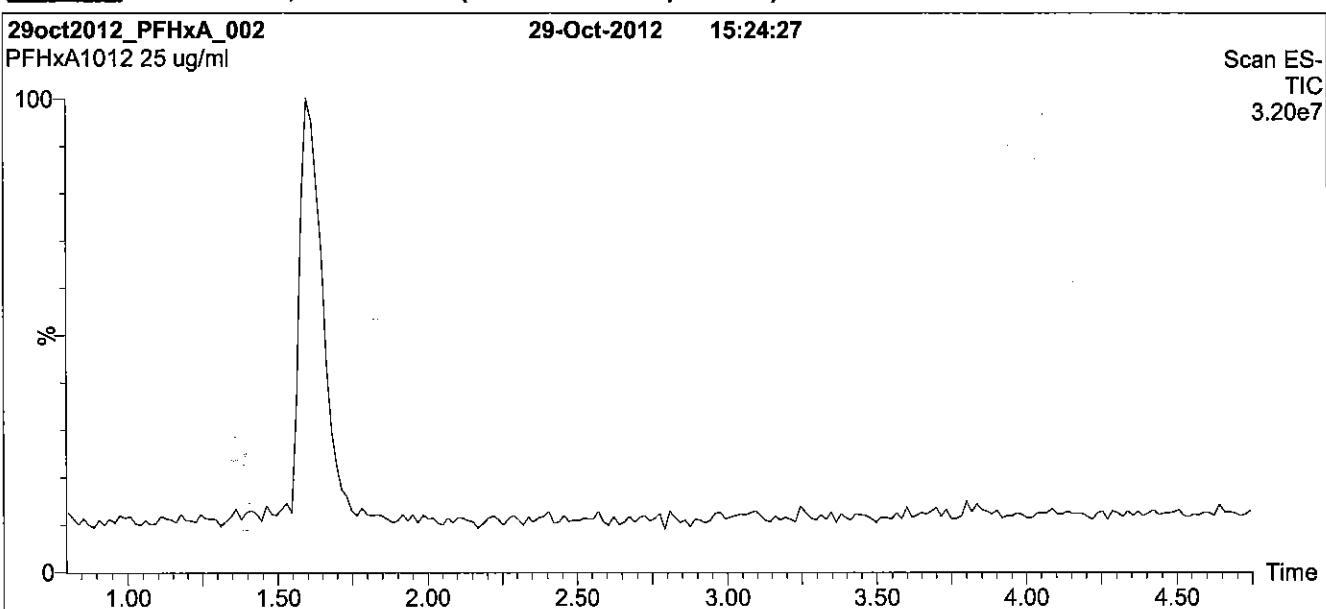
#### **QUALITY MANAGEMENT:**

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 ACCLASS (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:** PFHxA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 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.

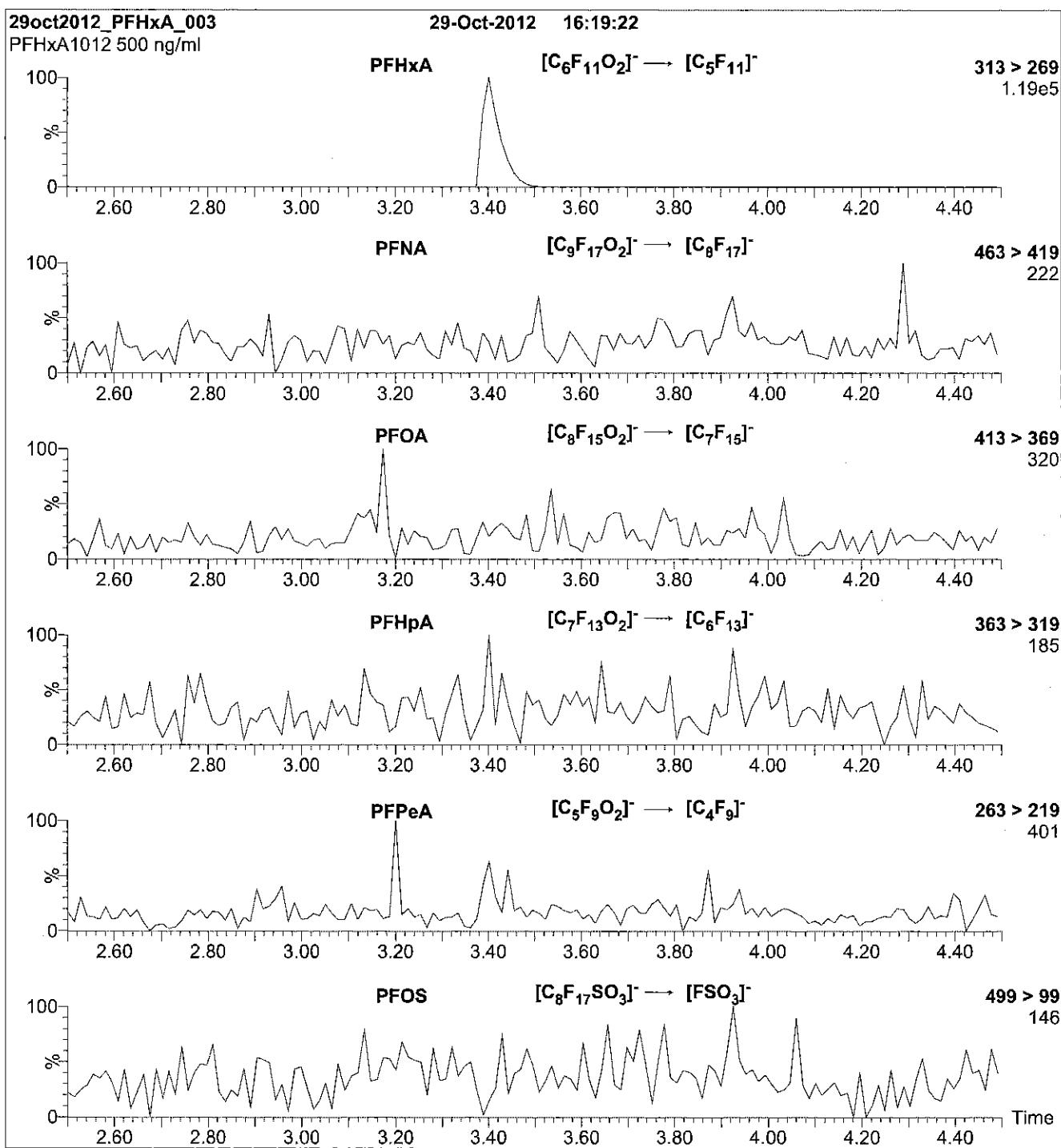
Time: 10 min  
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 15.00  
Cone Gas Flow (l/hr) = 100  
Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.13e-3  
 Collision Energy (eV) = 10

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

Reagent

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**LCPFHxS\_00002**

R 9-17-12 DE4



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Dup LCPFHxS-00001

new lot LCPFHxS-00002

PRODUCT CODE:

L-PFHxS

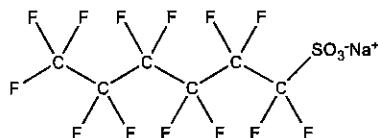
LOT NUMBER: LPFHxS0312

COMPOUND:

Sodium perfluoro-1-hexanesulfonate

STRUCTURE:

CAS #: 82382-12-5



MOLECULAR FORMULA:

C<sub>6</sub>F<sub>13</sub>SO<sub>3</sub>Na

MOLECULAR WEIGHT: 422.10

CONCENTRATION:

50.0 ± 2.5 µg/ml (Na salt)

SOLVENT(S): Methanol

47.3 ± 2.4 µg/ml (PFHxS anion) ✓

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

03/27/2012

EXPIRY DATE: (mm/dd/yyyy)

03/27/2015

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

### DOCUMENTATION/ DATA ATTACHED:

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

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

### ADDITIONAL INFORMATION:

- See page 2 for further details.

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

Certified By:

B.G. Chittim

Date: 04/18/2012

(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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**SYNTHESIS / CHARACTERIZATION:**

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**HOMOGENEITY:**

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**UNCERTAINTY:**

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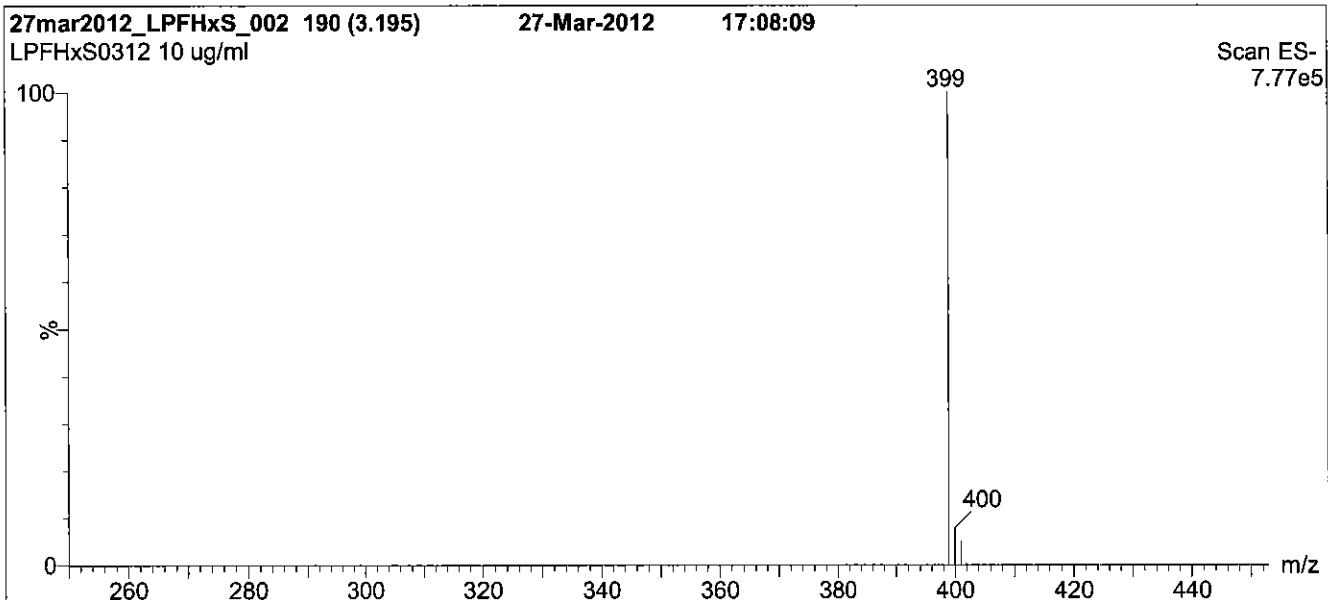
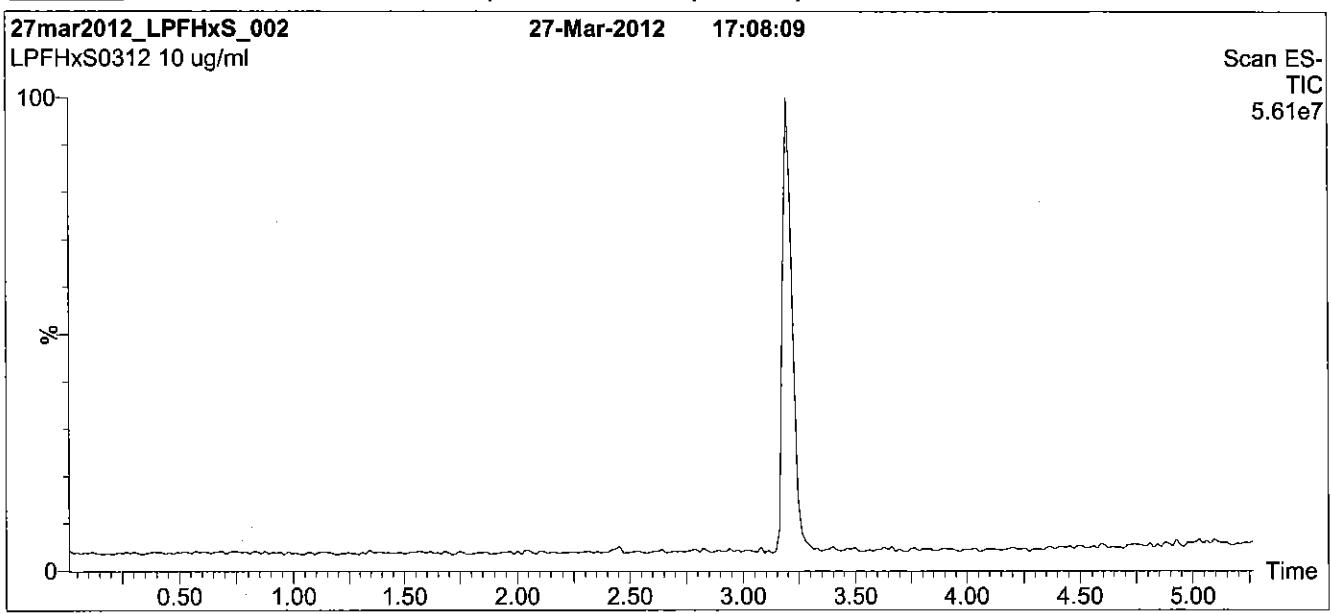
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**LIMITED WARRANTY:**

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**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 Shield RP<sub>18</sub>  
1.7 µm, 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

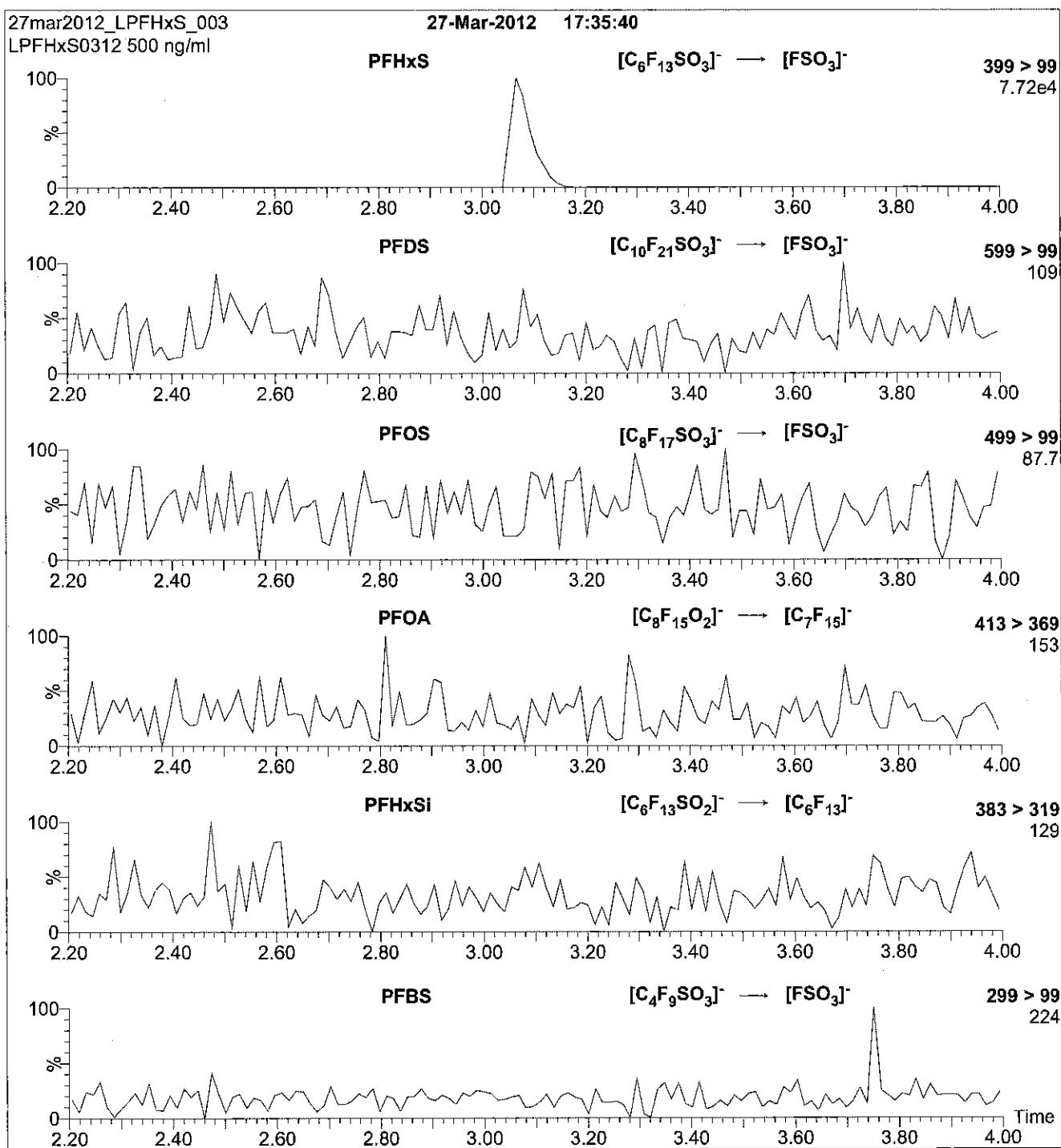
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (250 - 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:** L-PFHzS; LC/MS/MS Data (Selected MRM Transitions)



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.35e-3  
Collision Energy (eV) = 30

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

Flow: 300  $\mu$ l/min

Reagent

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**LCPFNA\_00002**

R 9-17-12 DE4



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Dup LCPFDA-00001

New lot LCPFOA-00002

PRODUCT CODE:

PFNA

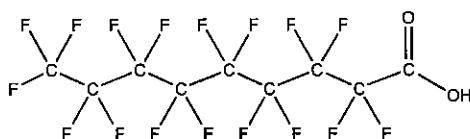
LOT NUMBER: PFNA0612

COMPOUND:

Perfluoro-n-nonanoic acid

STRUCTURE:

CAS #: 375-95-1



MOLECULAR FORMULA:

C<sub>9</sub>HF<sub>17</sub>O<sub>2</sub>

MOLECULAR WEIGHT: 464.08

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

06/14/2012

EXPIRY DATE: (mm/dd/yyyy)

06/14/2015

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: 06/15/2012

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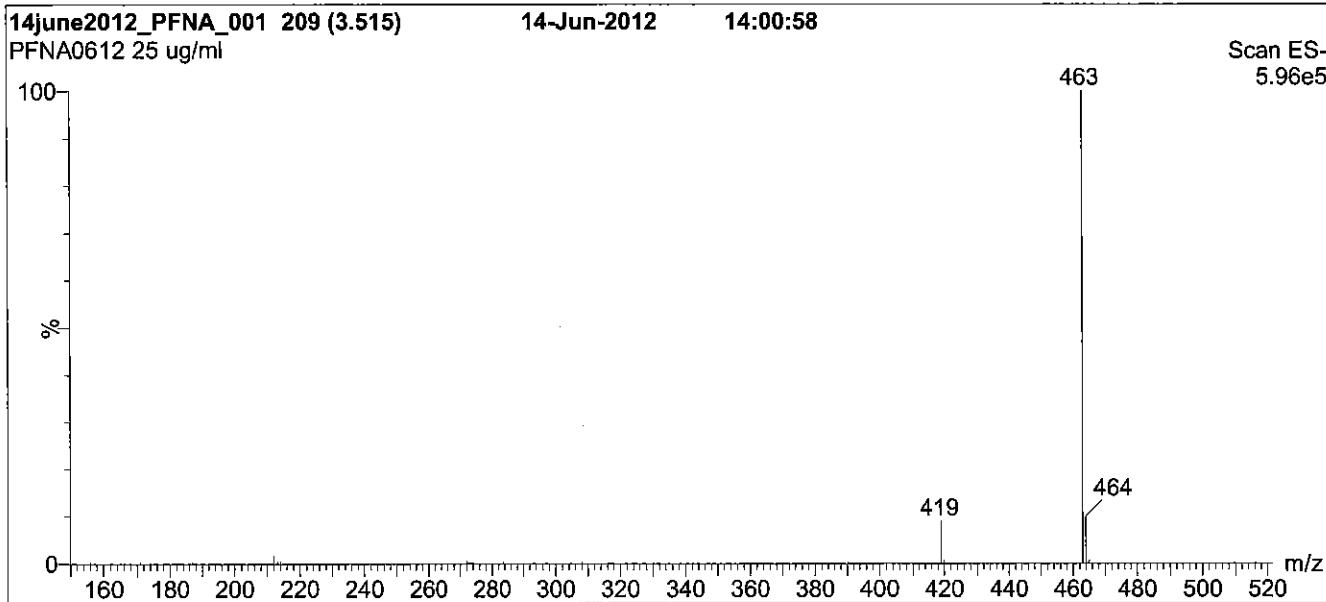
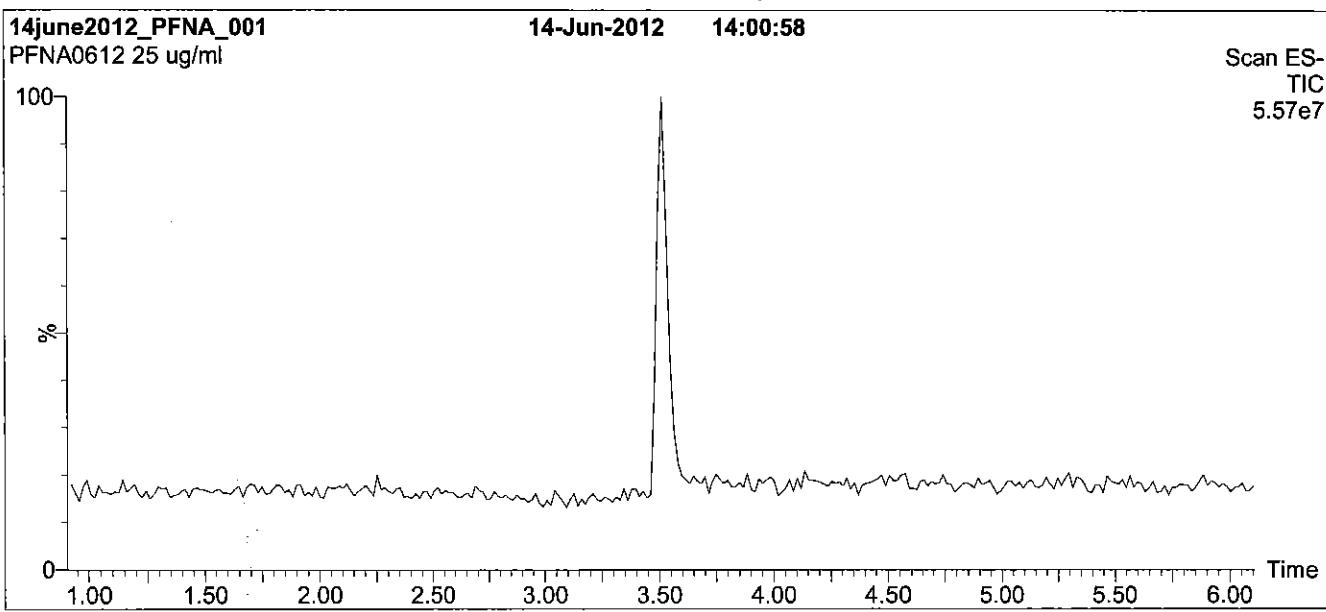
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**Figure 1:** PFNA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<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 1.5 min  
before returning to initial conditions in 0.5 min.

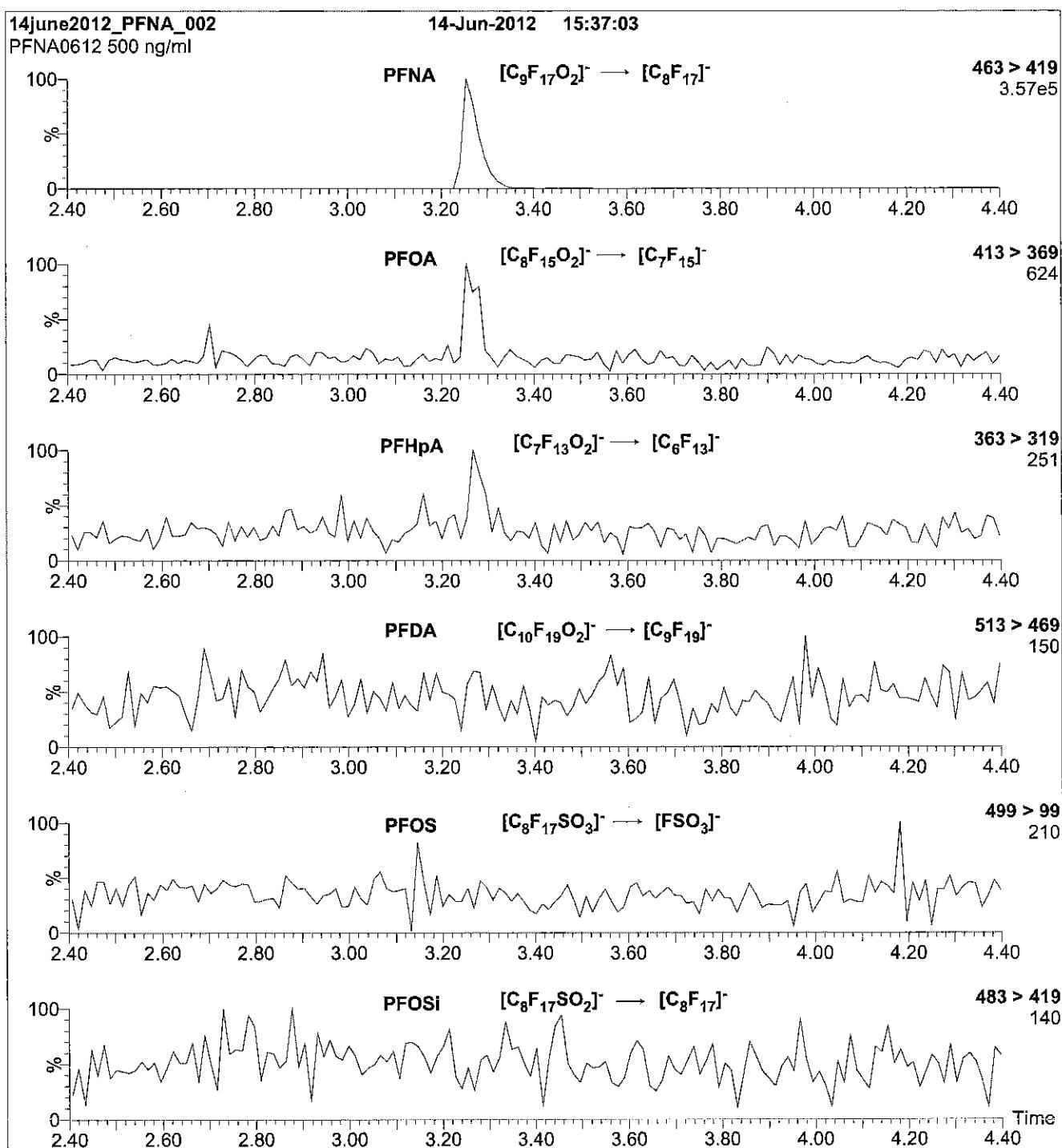
Time: 10 min  
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (150 - 850 amu)

Source: Electrospray (negative)  
Capillary Voltage (kV) = 2.00  
Cone Voltage (V) = 15.00  
Cone Gas Flow (l/hr) = 50  
Desolvation Gas Flow (l/hr) = 750

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



**Conditions for Figure 2:**

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

**MS Parameters**

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

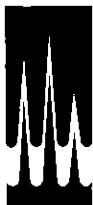
Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

Reagent

---

**LCPFNS\_00001**



**WELLINGTON  
LABORATORIES**

**CERTIFICATE OF ANALYSIS  
DOCUMENTATION**

R = (2-14-12)

**PRODUCT CODE:**

L-PFNS

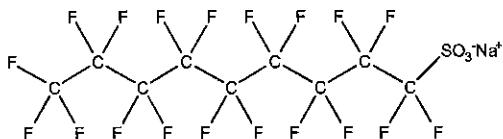
**LOT NUMBER:** LPFNS0712

**COMPOUND:**

Sodium perfluoro-1-nananesulfonate

**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)

07/04/2015

**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:** 07/12/2012

(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.

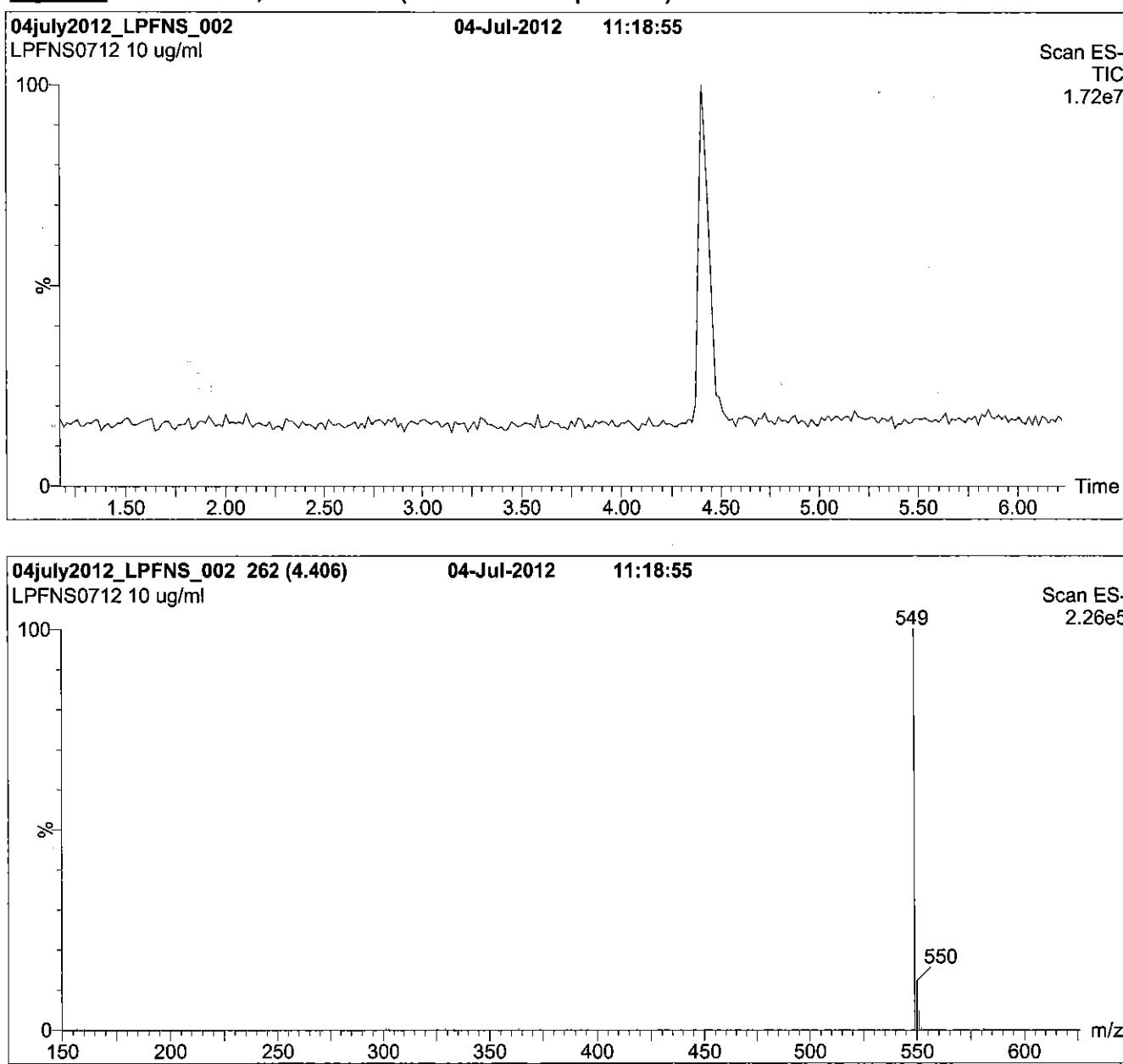
### **QUALITY MANAGEMENT:**

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 ACCLASS (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-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 μm, 2.1 x 100 mm

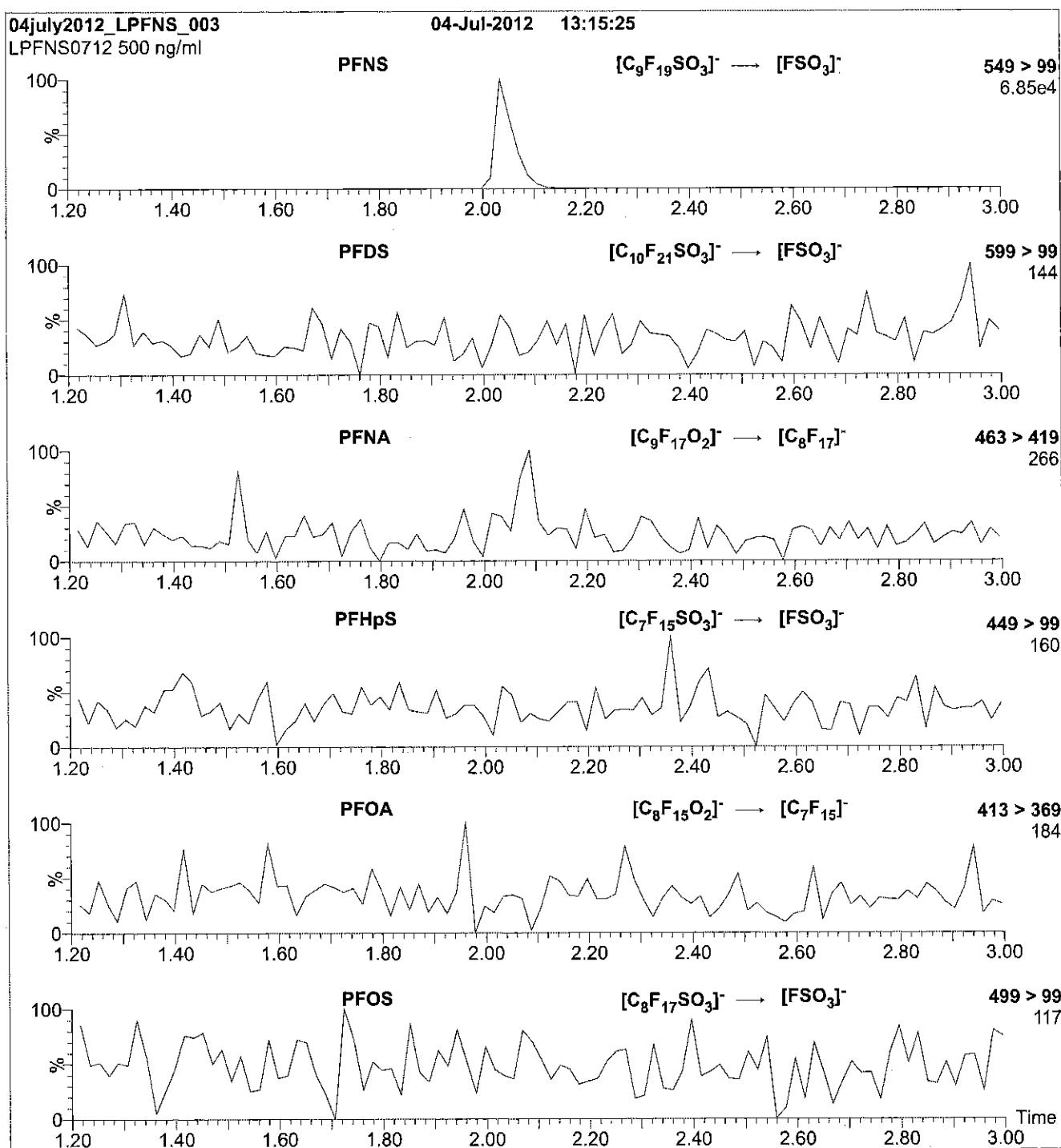
Mobile phase: Gradient  
Start: 55% (80:20 MeOH:ACN) / 45% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min  
and hold for 1.5 min before returning  
to initial conditions in 0.5 min.  
Time: 10 min  
Flow: 300 μl/min

**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)

**MS Parameters**

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

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

Reagent

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**LCPFOA\_00004**

Rec 7/15/14



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE:

PFOA

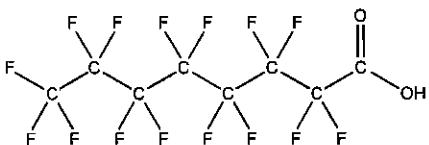
LOT NUMBER: PFOA1013

COMPOUND:

Perfluoro-n-octanoic acid

STRUCTURE:

CAS #: 335-67-1



MOLECULAR FORMULA:

$C_8HF_{15}O_2$

MOLECULAR WEIGHT: 414.07

CONCENTRATION:

50 ± 2.5 µg/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:

A handwritten signature in black ink.

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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### **EXPIRY DATE / PERIOD OF VALIDITY:**

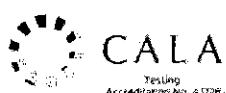
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

### **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.

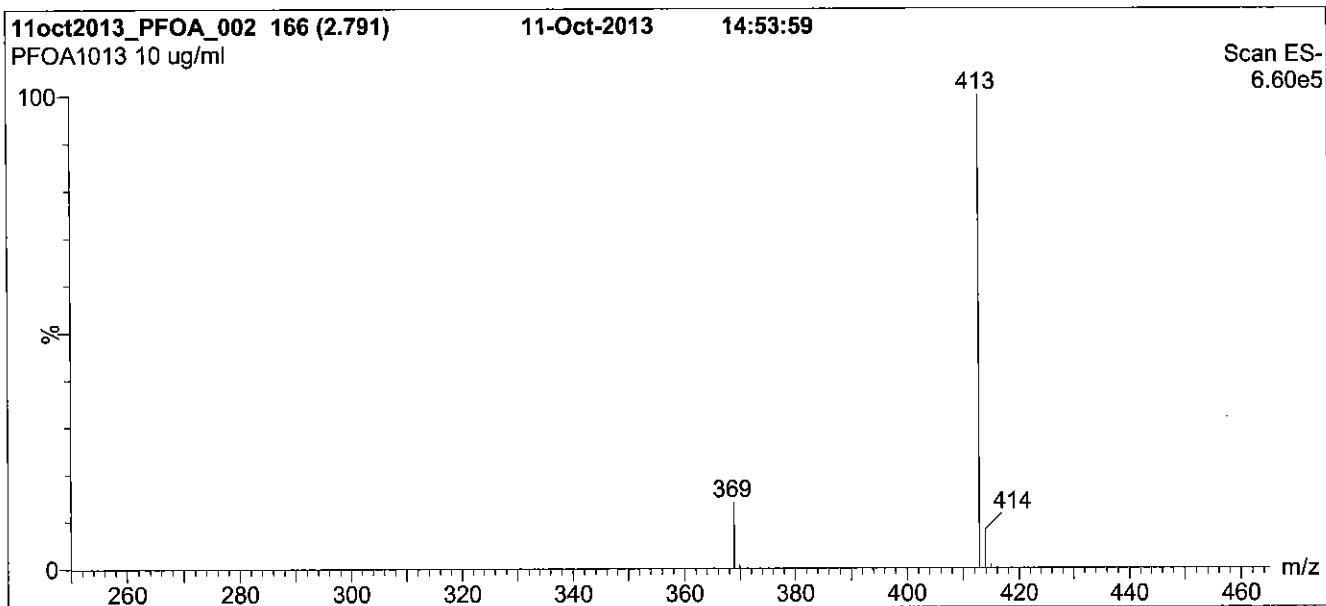
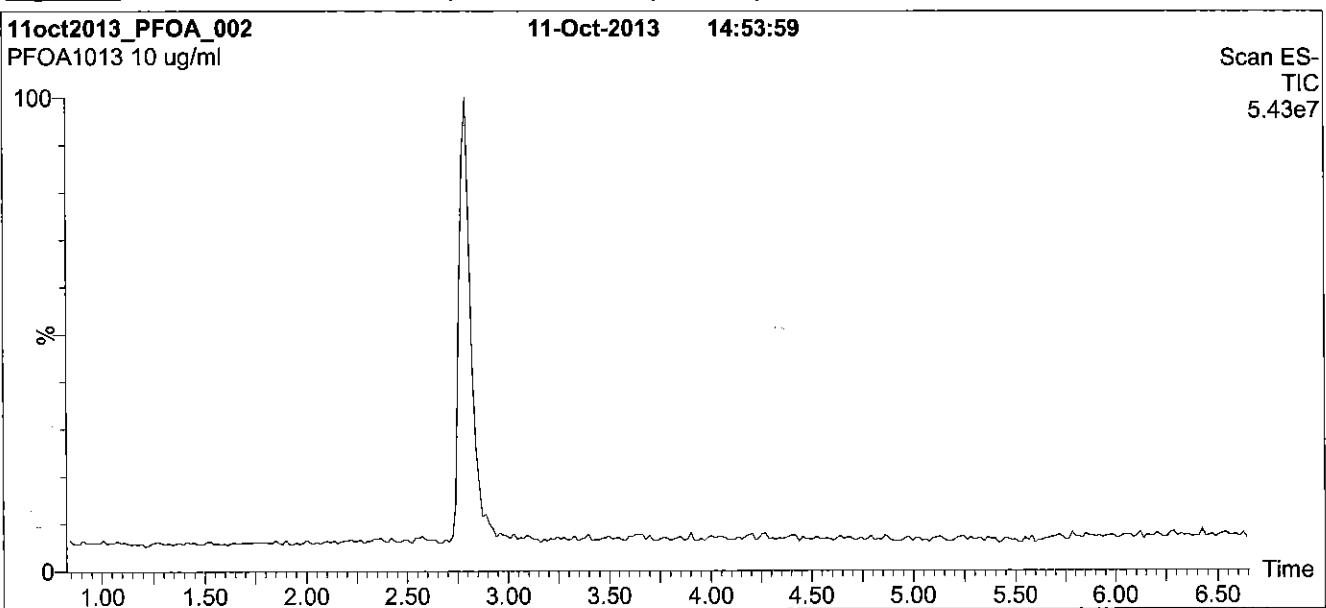
### **QUALITY MANAGEMENT:**

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



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<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 7.5 min and hold for  
1 min before returning to initial conditions in 0.5 min.  
Time: 10 min

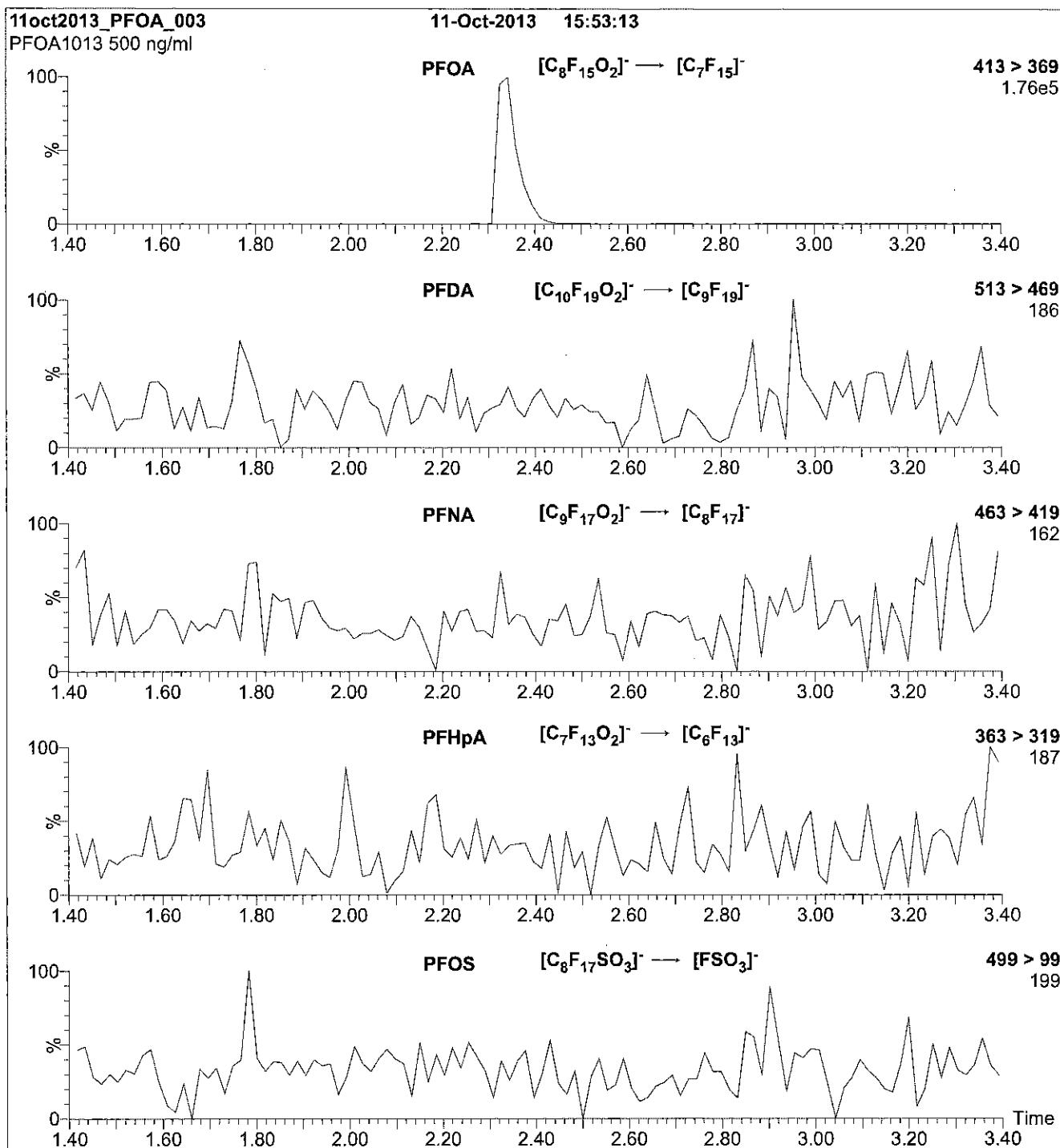
Flow: 300 µl/min

**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)

**MS Parameters**

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

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

Reagent

---

**LCPFOS\_00002**

R: 9-17-12 DE6



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

Dup LPFOS-00001  
new lot LPFOS-00002

PRODUCT CODE:

L-PFOS

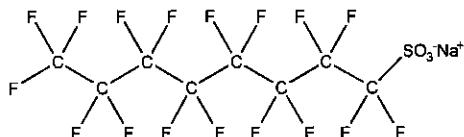
LOT NUMBER: LPFOS0312

COMPOUND:

Sodium perfluoro-1-octanesulfonate

STRUCTURE:

CAS #: 4021-47-0



MOLECULAR FORMULA:

$C_8F_{17}SO_3Na$

MOLECULAR WEIGHT: 522.11

CONCENTRATION:

$50.0 \pm 2.5 \mu\text{g/ml}$  (Na salt)

SOLVENT(S): Methanol

$47.8 \pm 2.4 \mu\text{g/ml}$  (PFOS anion) ✓

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

03/26/2012

EXPIRY DATE: (mm/dd/yyyy)

03/26/2015

RECOMMENDED STORAGE: Store ampoule in a cool, dark place

### DOCUMENTATION/ DATA ATTACHED:

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

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

### ADDITIONAL INFORMATION:

- See page 2 for further details.

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

Certified By:

B.G. Chittim

Date: 04/04/2012

(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:**

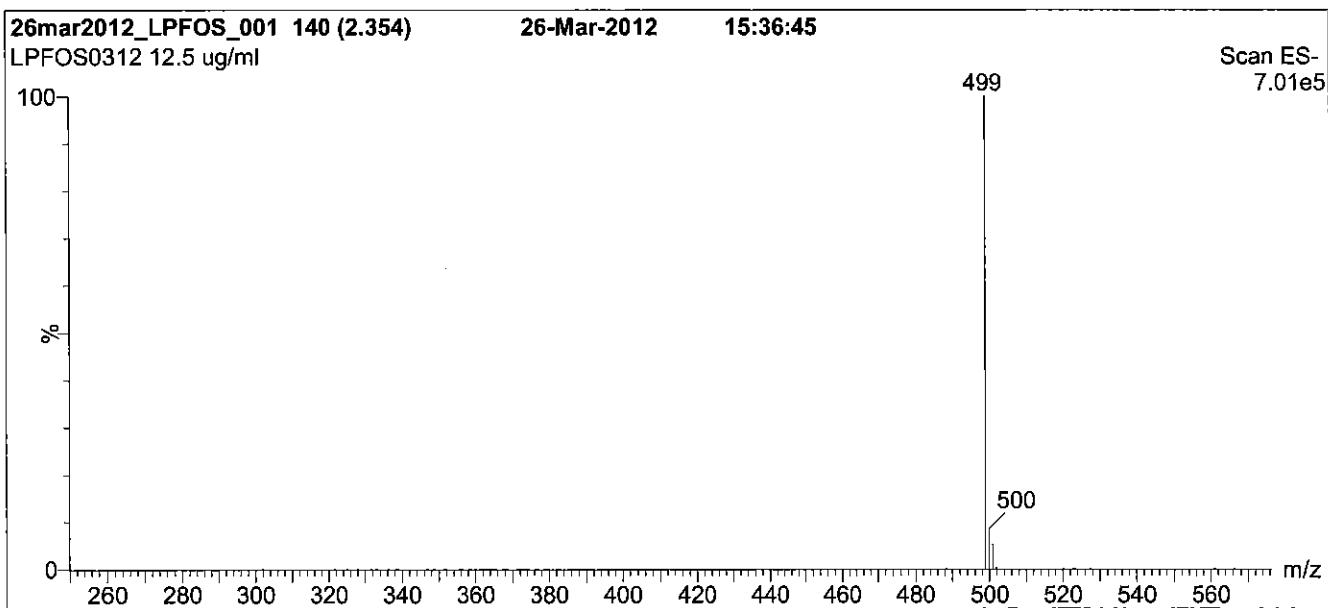
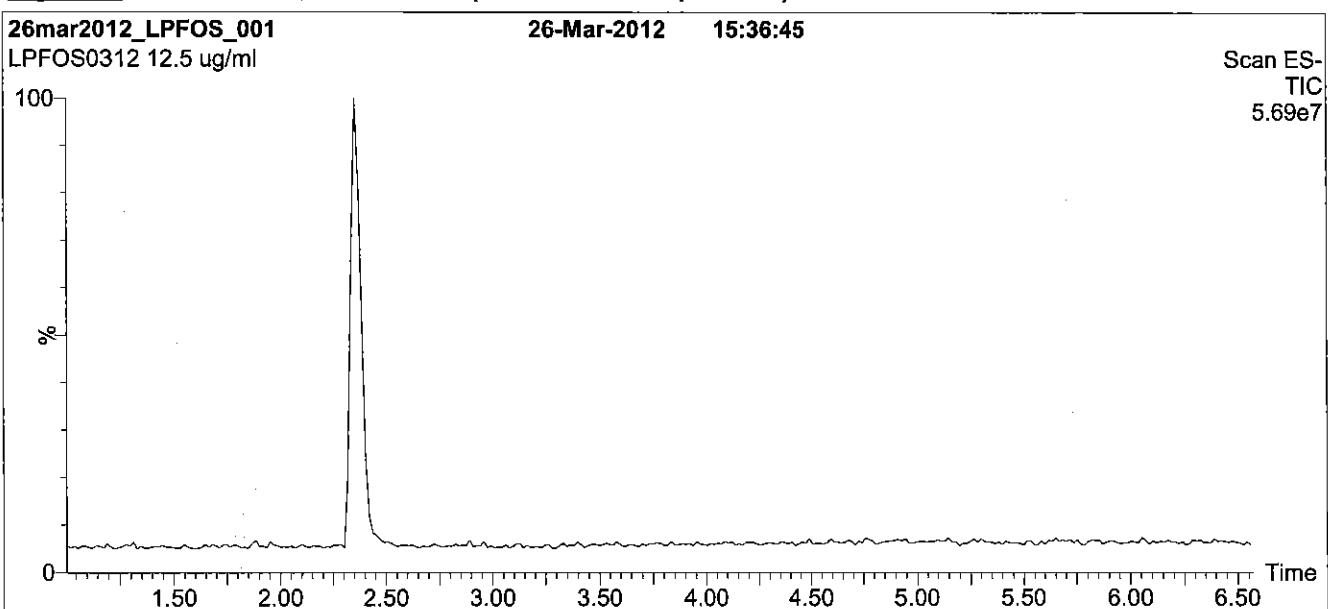
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**Figure 1:** L-PFOS; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity 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)  
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

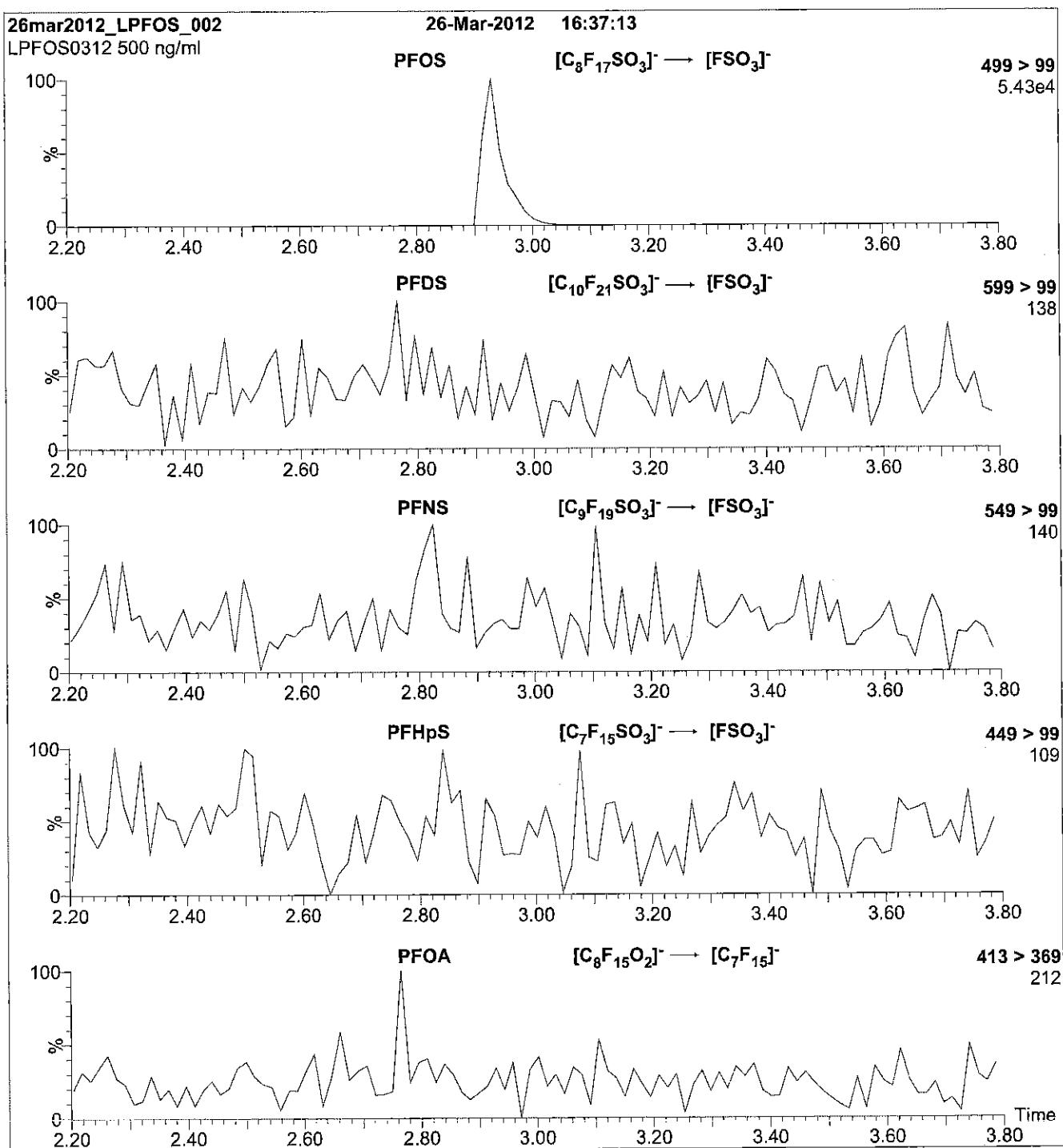
Flow: 300  $\mu$ l/min

**MS Parameters**

Experiment: Full Scan (250 - 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) = 700

**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)

**MS Parameters**

Collision Gas (mbar) = 3.31e-3  
Collision Energy (eV) = 40

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

Flow: 300  $\mu$ l/min

Reagent

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**LCPFOSA\_00004**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R : 12-14-12  
LCPFOSA-00004

PRODUCT CODE:

FOSA-M

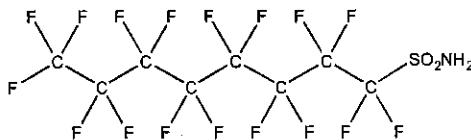
LOT NUMBER: FOSA0912M

COMPOUND:

Perfluoro-1-octanesulfonamide

STRUCTURE:

CAS #: 754-91-6



MOLECULAR FORMULA:

C<sub>8</sub>H<sub>17</sub>NO<sub>2</sub>S

MOLECULAR WEIGHT: 499.14

CONCENTRATION:

50 ± 2.5 µg/ml

SOLVENT(S): Methanol

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

09/13/2012

EXPIRY DATE: (mm/dd/yyyy)

09/13/2015

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: 10/03/2012

(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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#### **LIMITED WARRANTY:**

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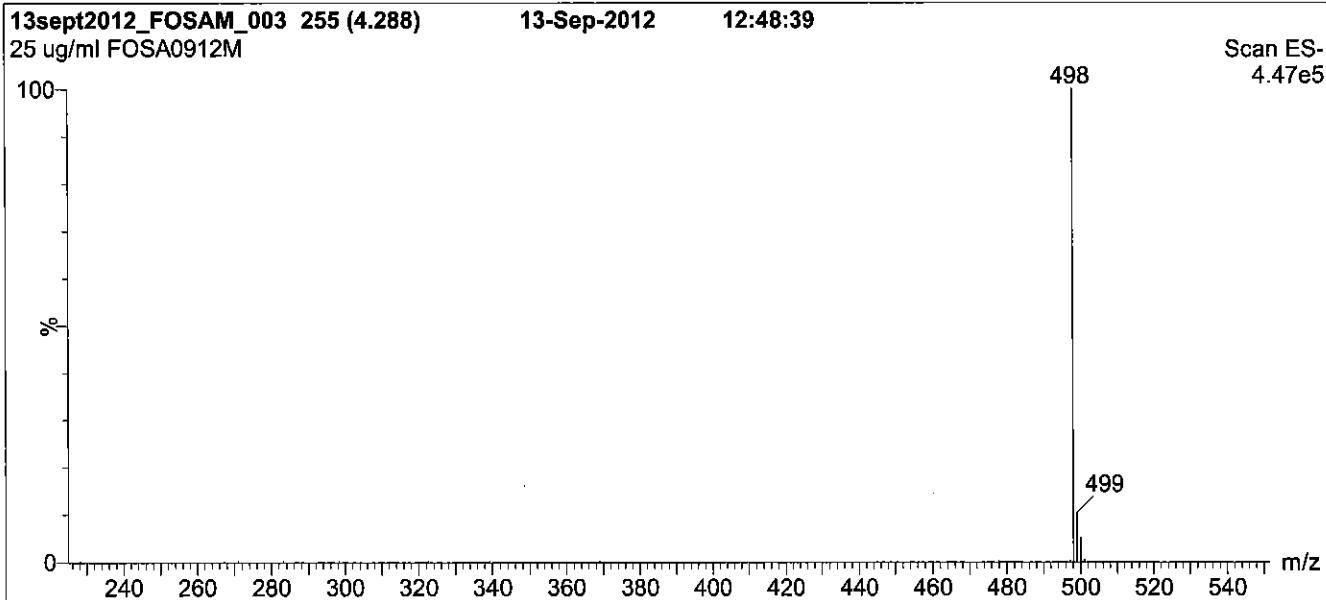
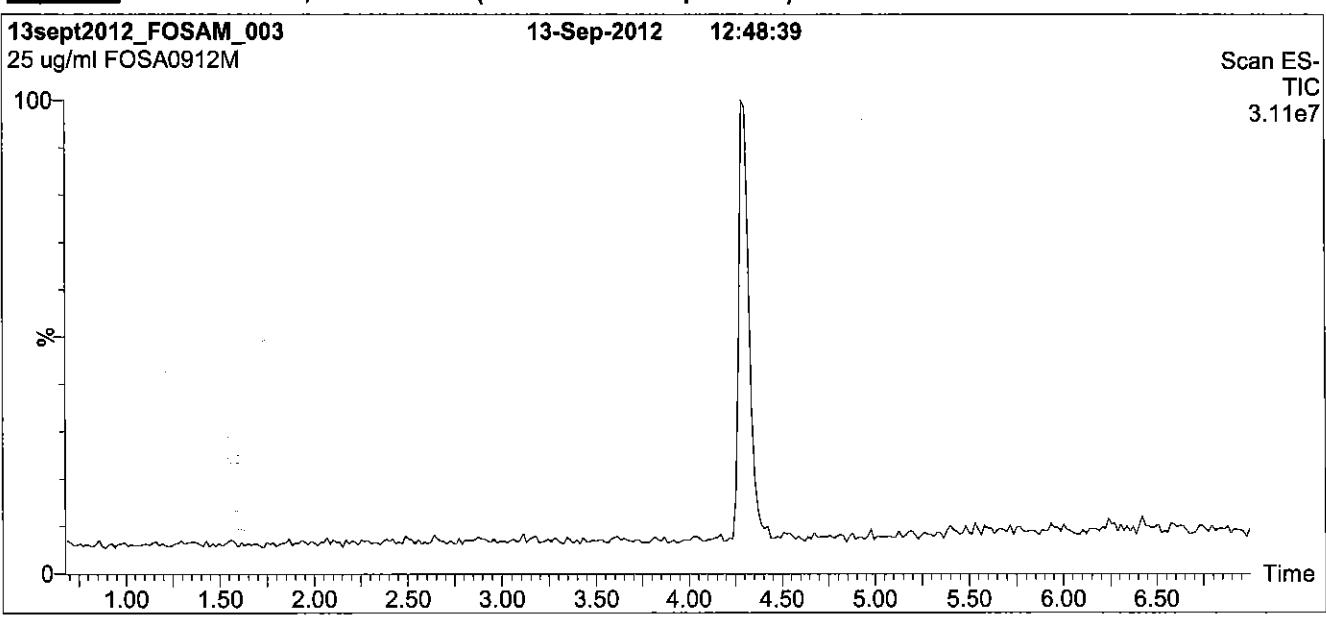
#### **QUALITY MANAGEMENT:**

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 ACCLASS (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-M; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity 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 90% organic over 7 min and hold for 2 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min

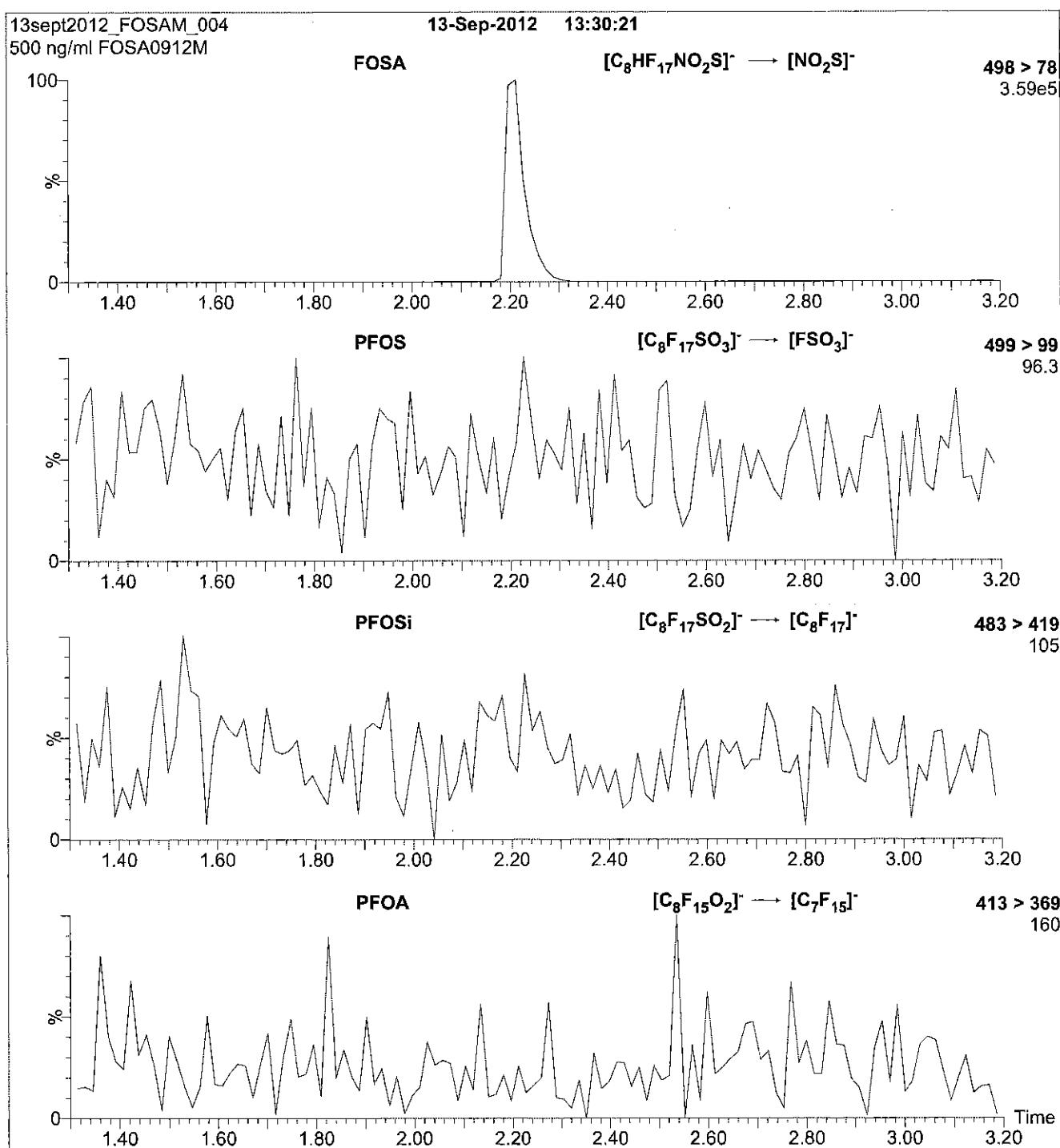
Flow: 300 μl/min

**MS Parameters**

Experiment: Full Scan (225 - 850 amu)

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

**Figure 2:** FOSA-M; LC/MS/MS Data (Selected MRM Transitions)



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.43e-3  
Collision Energy (eV) = 30

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

Reagent

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**LCPFPeA\_00003**

Rec 7/15/14



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

PRODUCT CODE:

PFPeA

LOT NUMBER: PFPeA0113

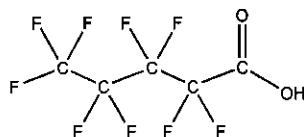
COMPOUND:

Perfluoro-n-pentanoic acid

STRUCTURE:

CAS #:

2706-90-3



MOLECULAR FORMULA:

$C_5HF_9O_2$

MOLECULAR WEIGHT: 264.05

CONCENTRATION:

$50 \pm 2.5 \mu\text{g/ml}$

SOLVENT(S): Methanol  
Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

01/03/2013

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  $C_5H_2F_8O_2$  (hydrido - derivative) as measured by  $^{19}F$  NMR.

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

Certified By:

A handwritten signature in black ink.

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.

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Prior to solution preparation, crystalline material is tested for homogeneity using a variety of techniques (as stated above) and its solubility in a given diluent is taken into consideration. Duplicate solutions of a new product are prepared from the same crystalline lot and, after the addition of an appropriate internal standard, they are compared by GC/MS 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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$$u_c(y(x_1, x_2, \dots, x_n)) = \sqrt{\sum_{i=1}^n u(y, x_i)^2}$$

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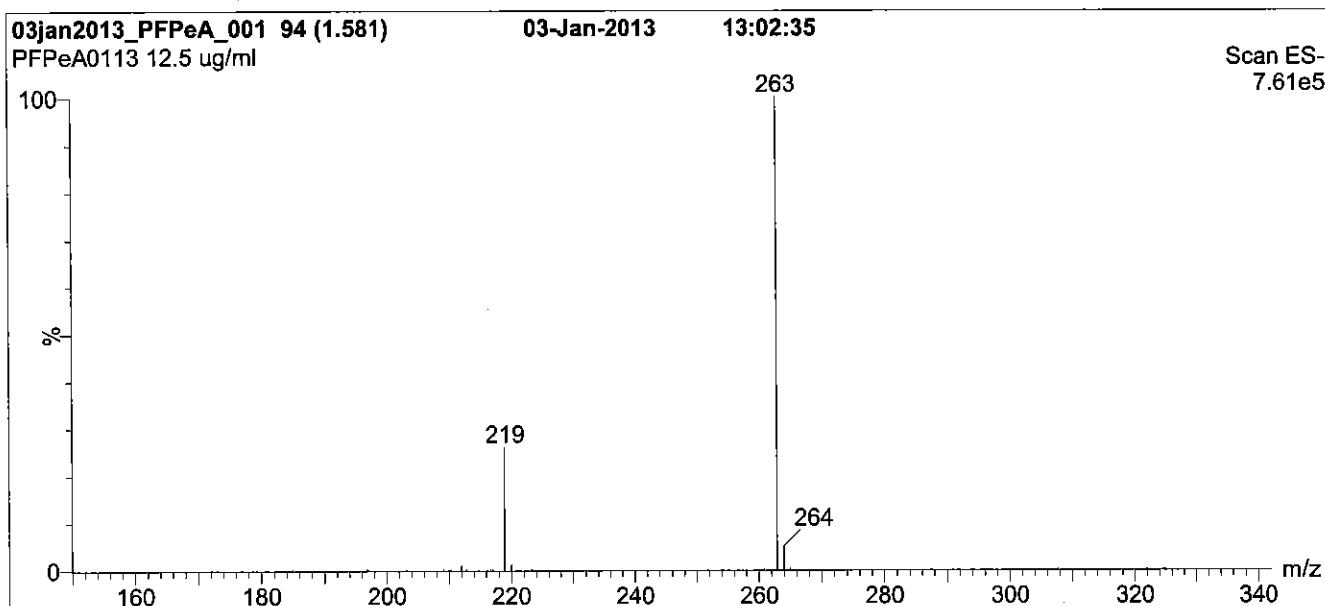
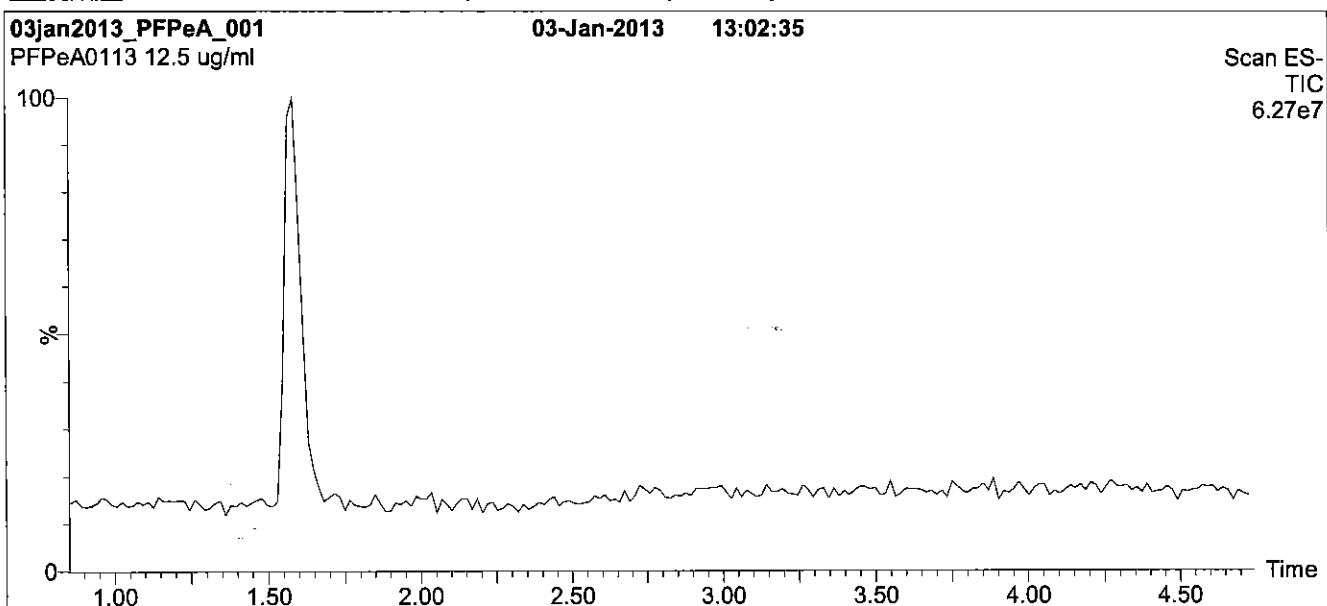
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**Figure 1: PFPeA; LC/MS Data (TIC and Mass Spectrum)**



**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 µm, 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

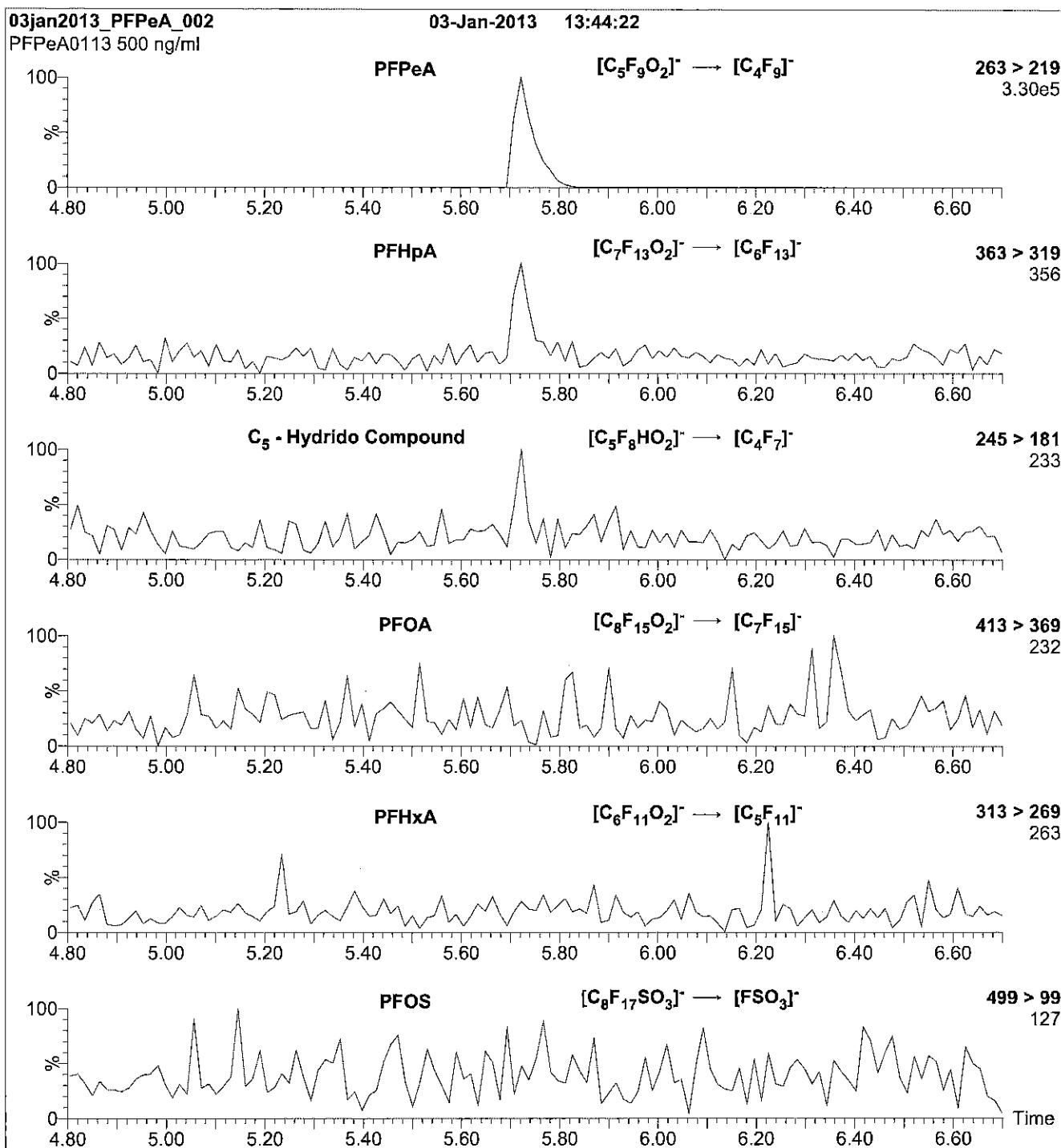
Flow: 300 µl/min

**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)

**MS Parameters**

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

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

Reagent

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**LCPFPeS\_00001**



# WELLINGTON LABORATORIES

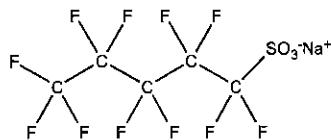
## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12 New part  
LC PFPeS-00001

PRODUCT CODE: L-PFPeS LOT NUMBER: LPFPeS0712

COMPOUND: Sodium perfluoro-1-pentanesulfonate

STRUCTURE: CAS #: Not available



MOLECULAR FORMULA: C<sub>5</sub>F<sub>11</sub>SO<sub>3</sub>Na MOLECULAR WEIGHT: 372.09  
CONCENTRATION: 50.0 ± 2.5 µg/ml (Na salt) SOLVENT(S): Methanol  
46.9 ± 2.3 µg/ml (PFPeS anion)  
CHEMICAL PURITY: >98%  
LAST TESTED: (mm/dd/yyyy) 07/04/2012  
EXPIRY DATE: (mm/dd/yyyy) 07/04/2015  
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: 07/09/2012

(mm/dd/yyyy)

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

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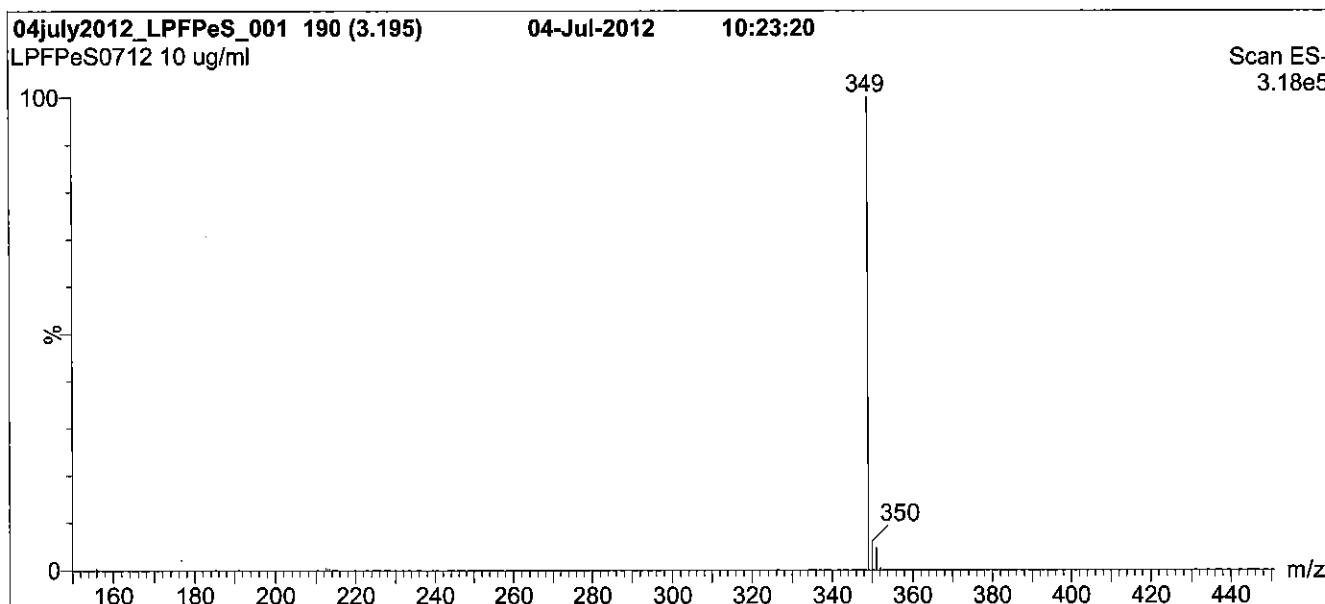
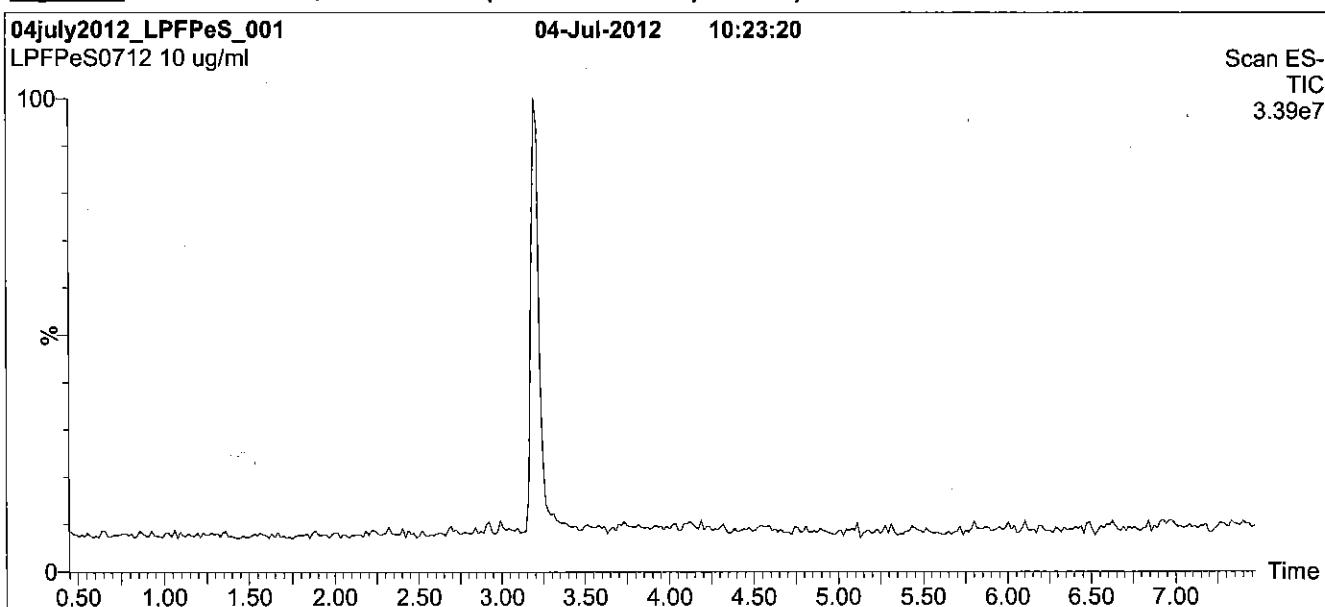
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**Figure 1:** L-PFPeS; LC/MS Data (TIC and Mass Spectrum)



**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 μm, 2.1 x 100 mm

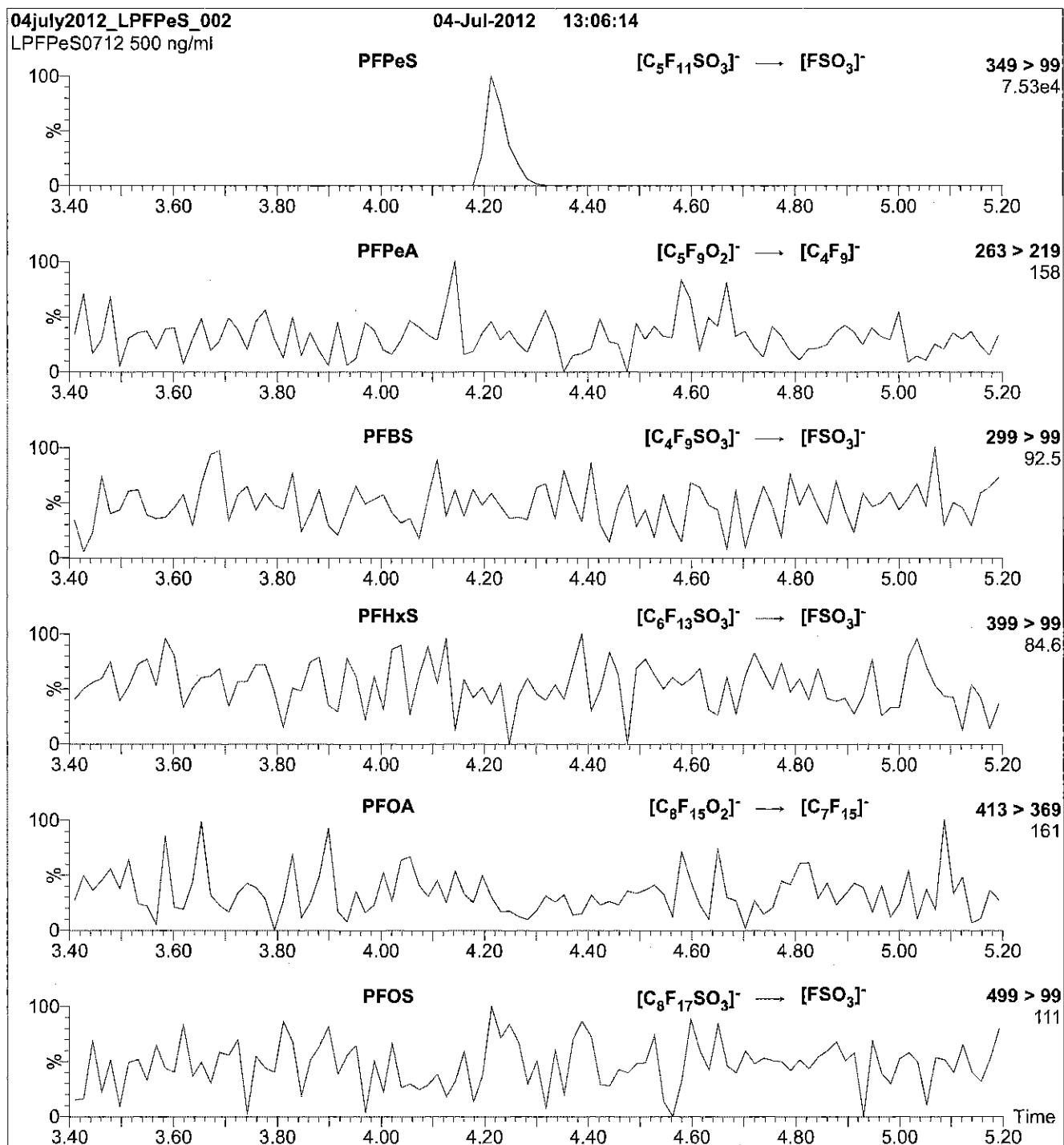
Mobile phase: Gradient  
Start: 40% (80:20 MeOH:ACN) / 60% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 90% organic over 7 min  
and hold for 1.5 min before returning  
to initial conditions in 0.5 min.  
Time: 10 min  
Flow: 300 μl/min

**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)

**MS Parameters**

Collision Gas (mbar) = 3.66e-3  
Collision Energy (eV) = 30

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

Reagent

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**LCPFTeDA\_00002**



# WELLINGTON LABORATORIES

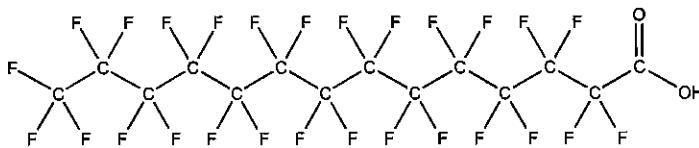
## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LC PFTe DA - 00002

PRODUCT CODE: PFTeDA      LOT NUMBER: PFTeDA0312  
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) 03/07/2012  
EXPIRY DATE: (mm/dd/yyyy) 03/07/2015  
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: 03/09/2012

(mm/dd/yyyy)

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

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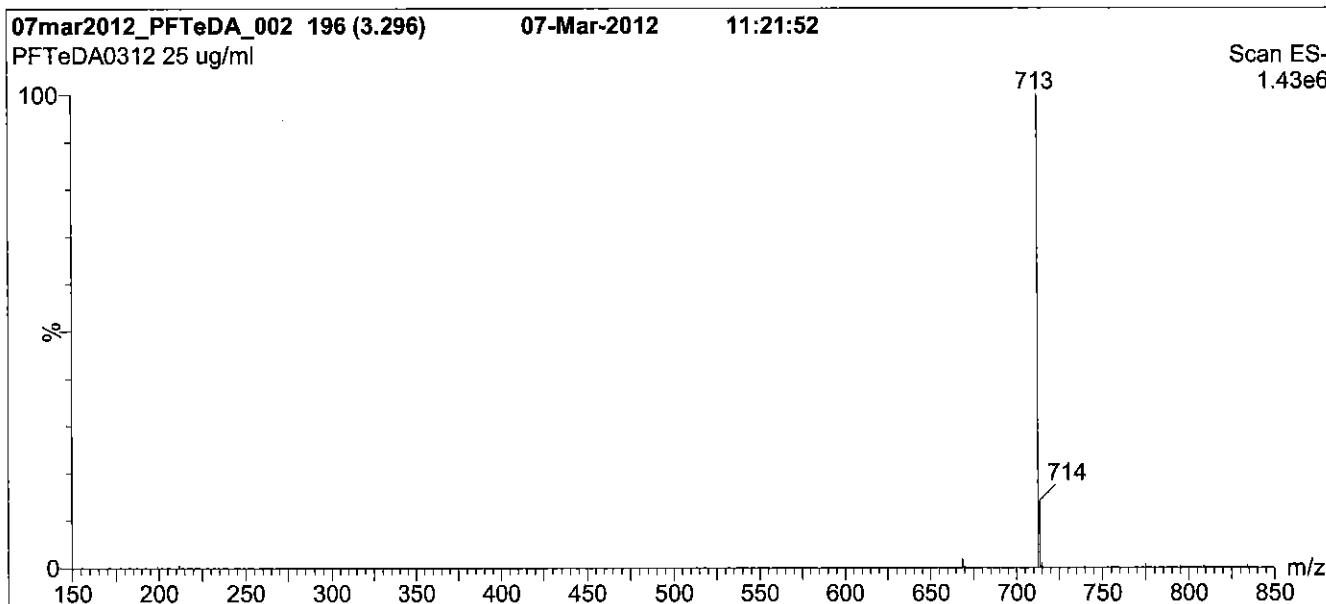
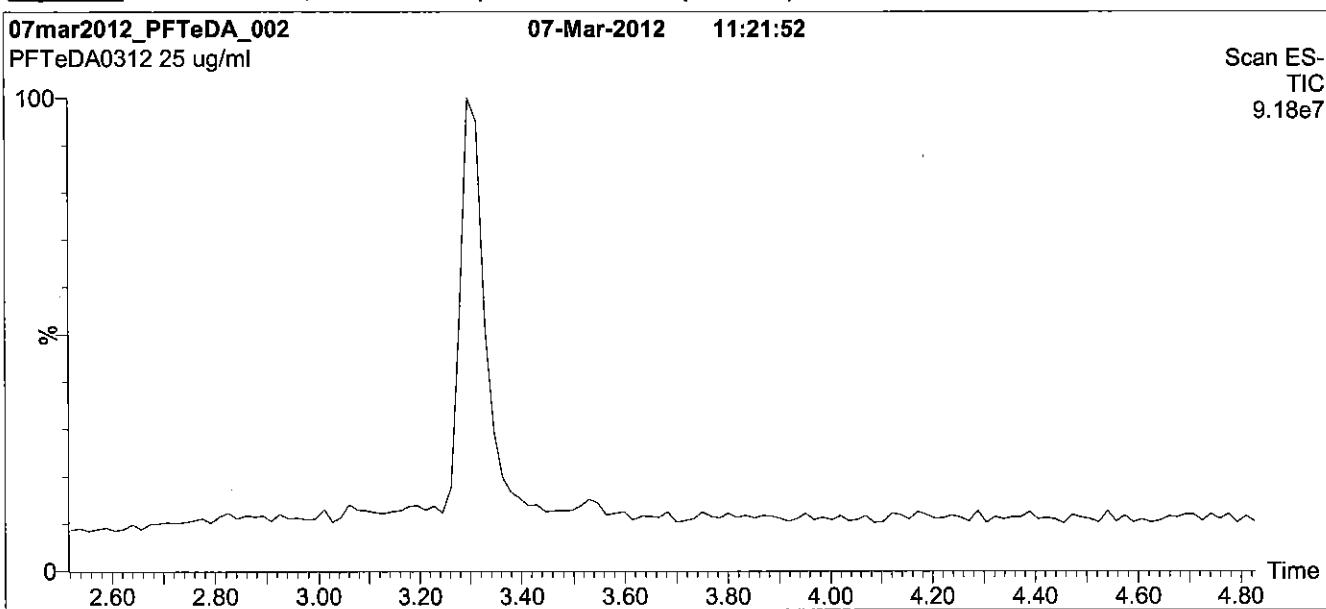
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**Figure 1:** PFTeDA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity UPLC BEH Shield RP<sub>18</sub>  
1.7 µm, 2.1 x 100 mm

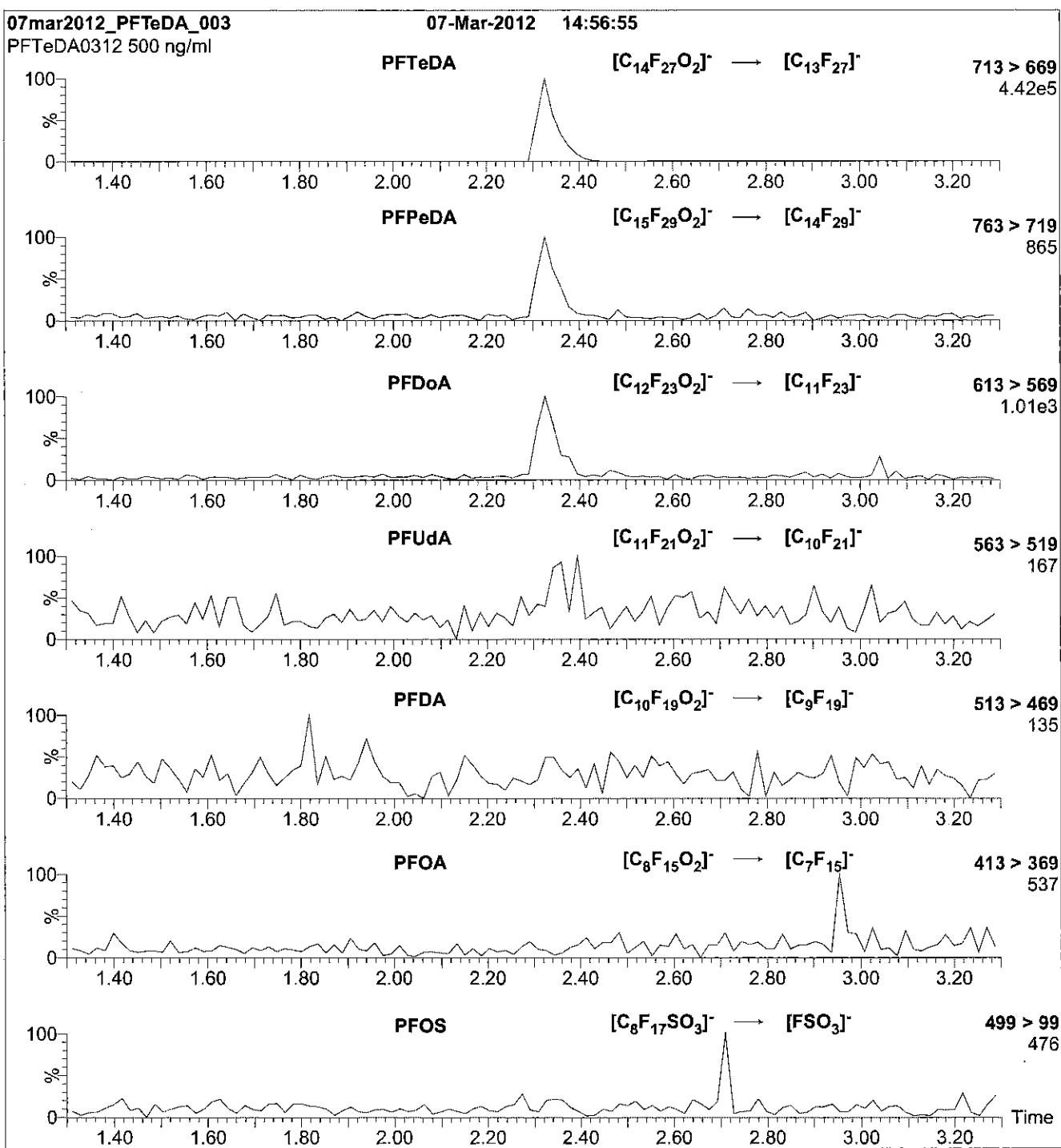
Mobile phase: Gradient  
Start: 70% (80:20 MeOH:ACN) / 30% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)  
Ramp to 100% organic over 7 min and hold 1.5 min  
before returning to initial conditions in 0.50 min.  
Time: 10 min

Flow: 300 µl/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) = 650

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.31e-3  
Collision Energy (eV) = 14

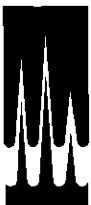
Mobile phase: Isocratic 80% (80:20 MeOH:ACN) / 20% H<sub>2</sub>O  
(both with 10 mM NH<sub>4</sub>OAc buffer)

Flow: 300  $\mu$ l/min

Reagent

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**LCPFTrDA  00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LCPFTrDA-00002

PRODUCT CODE:

PFTrDA

LOT NUMBER: PFTrDA0312

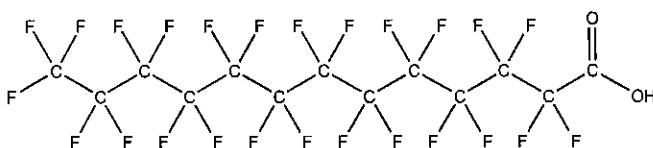
COMPOUND:

Perfluoro-n-tridecanoic acid

STRUCTURE:

CAS #:

72629-94-8



MOLECULAR FORMULA:

$C_{13}HF_{25}O_2$

MOLECULAR WEIGHT: 664.11

CONCENTRATION:

$50 \pm 2.5 \mu\text{g/ml}$

SOLVENT(S): Methanol

Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

03/26/2012

EXPIRY DATE: (mm/dd/yyyy)

03/26/2015

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 PFDa ( $C_{12}HF_{23}O_2$ ), 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: 04/04/2012

(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:**

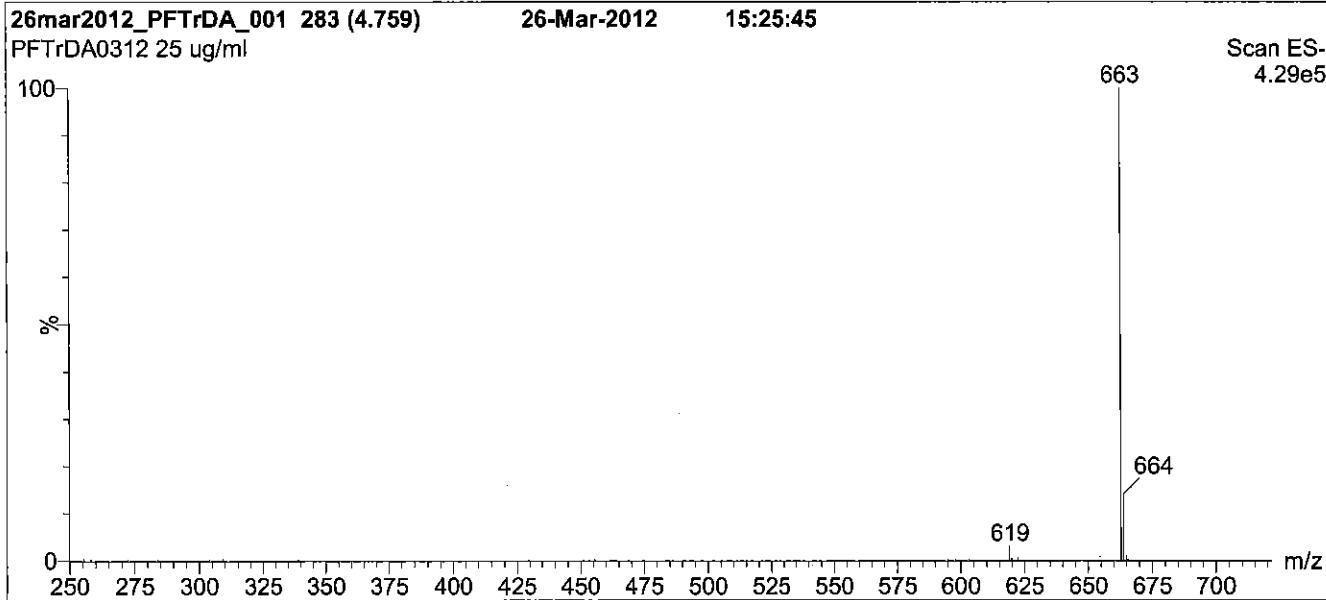
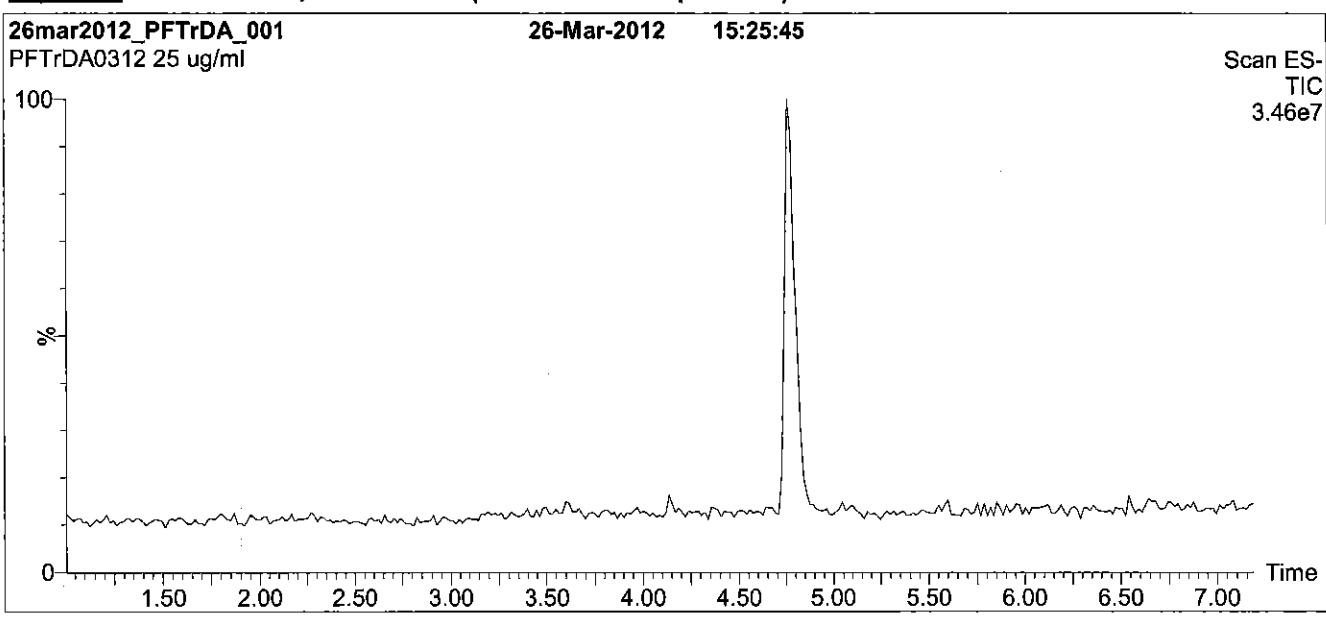
Ongoing stability studies of this product have demonstrated stability in its composition and concentration for the period of time specified by the expiry date in the unopened ampoule. Monitoring for any degradation or change in concentration of the listed analyte(s) is performed on a routine basis.

**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.

\*\*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:** PFTrDA; LC/MS Data (TIC and Mass Spectrum)



**Conditions for Figure 1:**

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

**Chromatographic Conditions**

Column: Acuity 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 90% organic over 7 min and hold for 1.5 min  
before returning to initial conditions in 0.5 min.

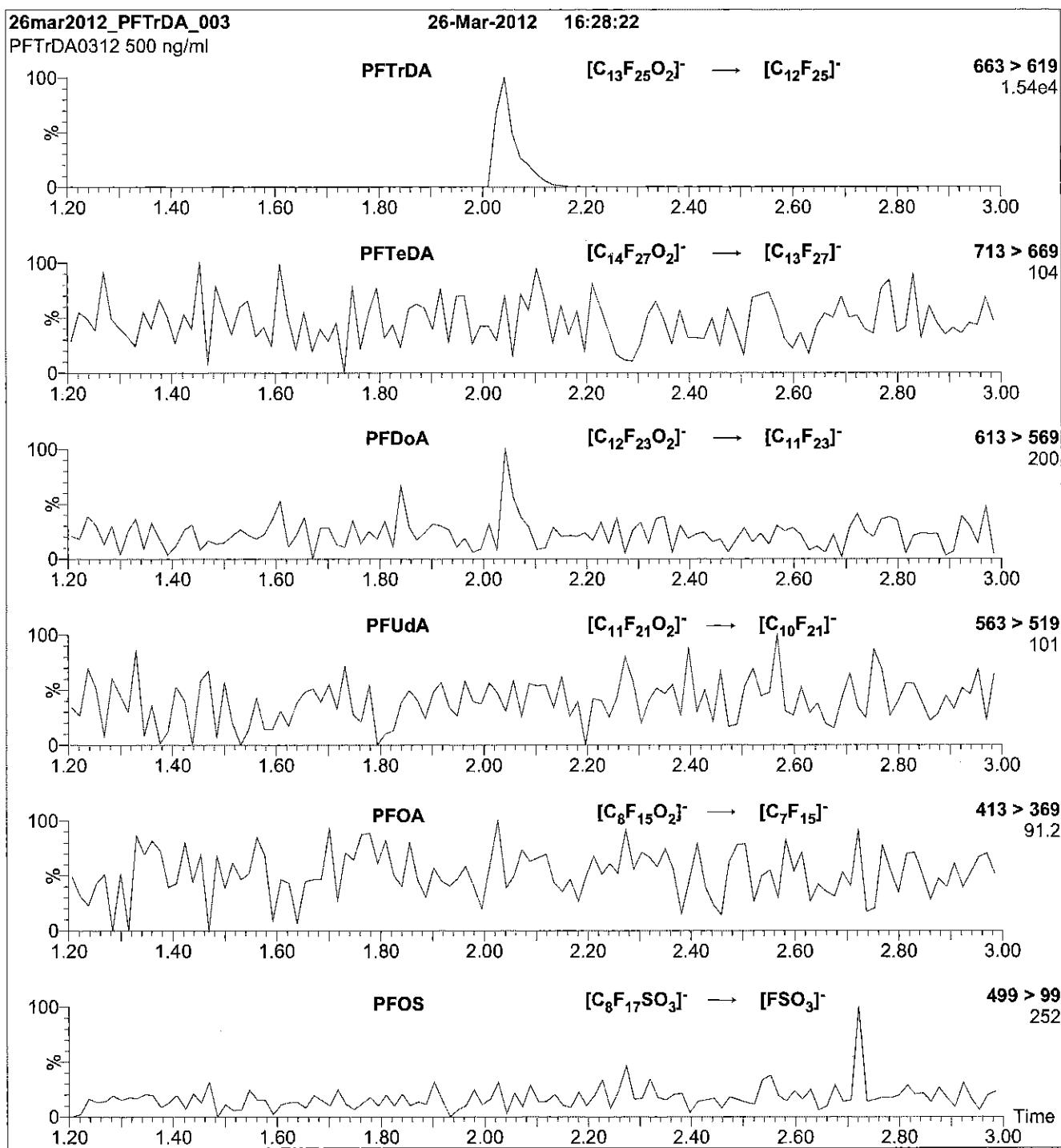
Time: 10 min  
Flow: 300 µl/min

**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) = 60  
Desolvation Gas Flow (l/hr) = 650

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



**Conditions for Figure 2:**

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

**MS Parameters**

Collision Gas (mbar) = 3.20e-3  
Collision Energy (eV) = 15

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

Flow: 300  $\mu$ l/min

Reagent

---

**LCPFUdA\_00002**



# WELLINGTON LABORATORIES

## CERTIFICATE OF ANALYSIS DOCUMENTATION

R: 12-14-12

LC PFUdA-00002

PRODUCT CODE:

PFUdA

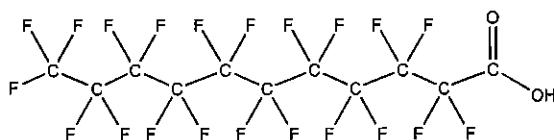
LOT NUMBER: PFUdA0312

COMPOUND:

Perfluoro-n-undecanoic acid

STRUCTURE:

CAS #: 2058-94-8



MOLECULAR FORMULA:

$C_{11}HF_{21}O_2$

MOLECULAR WEIGHT: 564.09

CONCENTRATION:

$50 \pm 2.5 \mu\text{g/ml}$

SOLVENT(S): Methanol

Water (<1%)

CHEMICAL PURITY:

>98%

LAST TESTED: (mm/dd/yyyy)

03/19/2012

EXPIRY DATE: (mm/dd/yyyy)

03/19/2015

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.

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Date: 03/27/2012

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

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**EXPIRY DATE / PERIOD OF VALIDITY:**

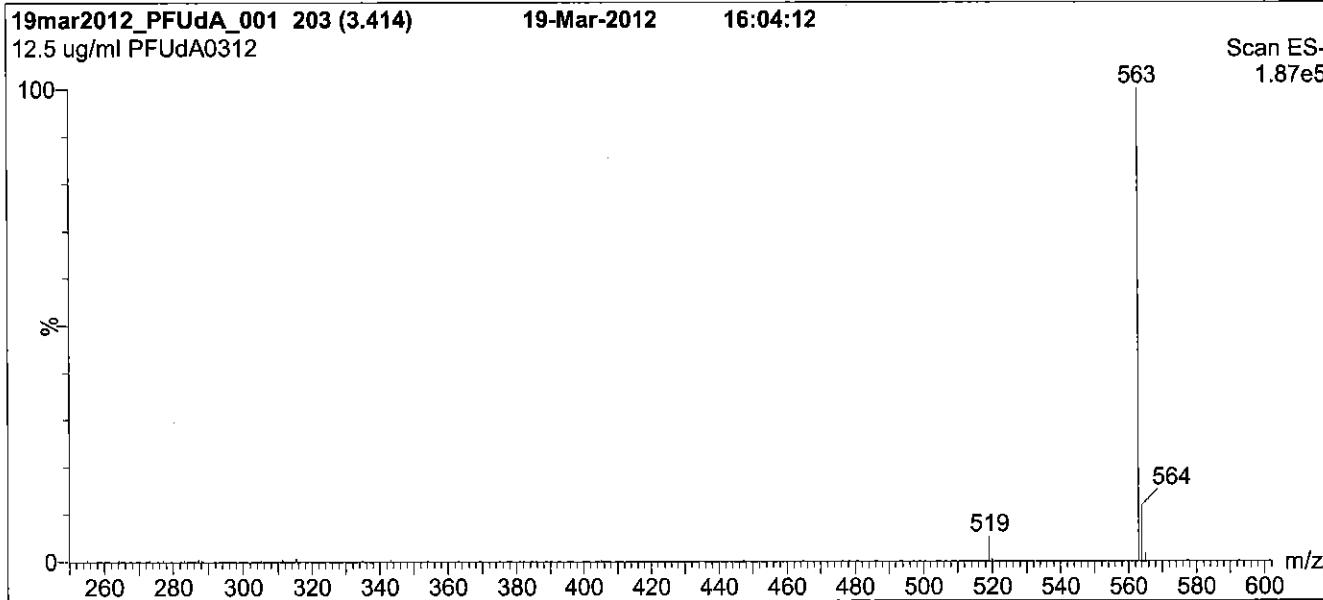
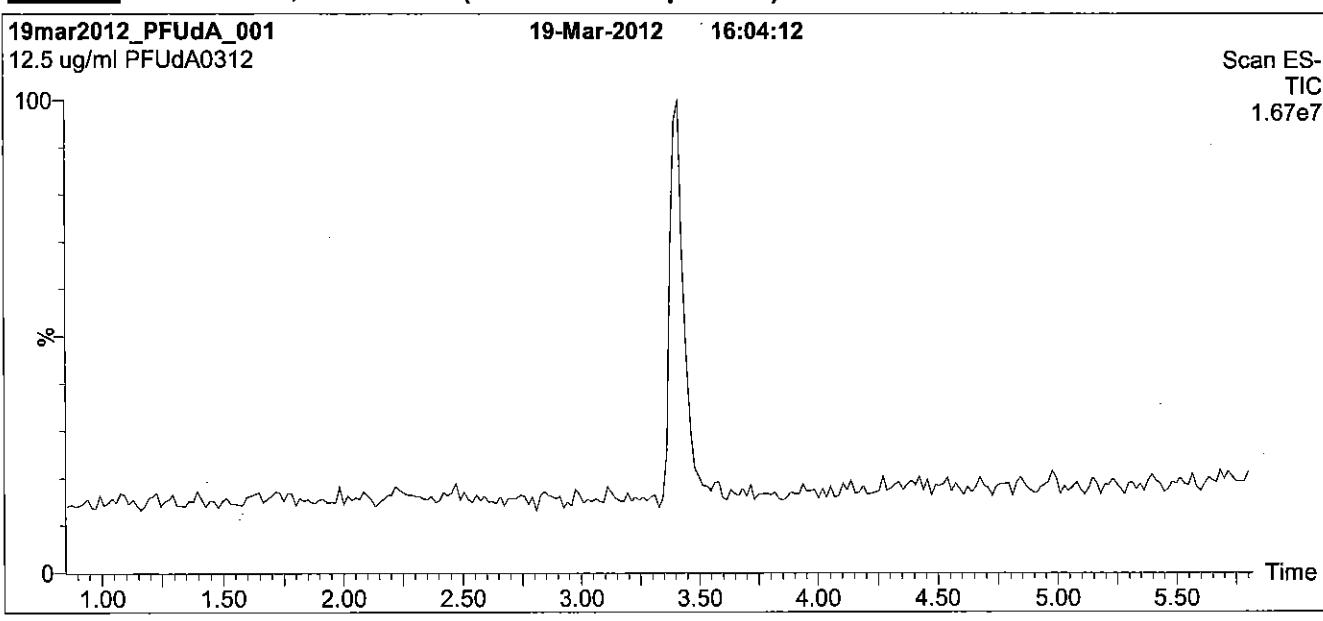
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**Figure 1:** PFUdA; LC/MS Data (TIC and Mass Spectrum)



**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 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 90% organic over 6.5 min and hold for 2 min  
before returning to initial conditions in 0.5 min.  
Time: 10 min

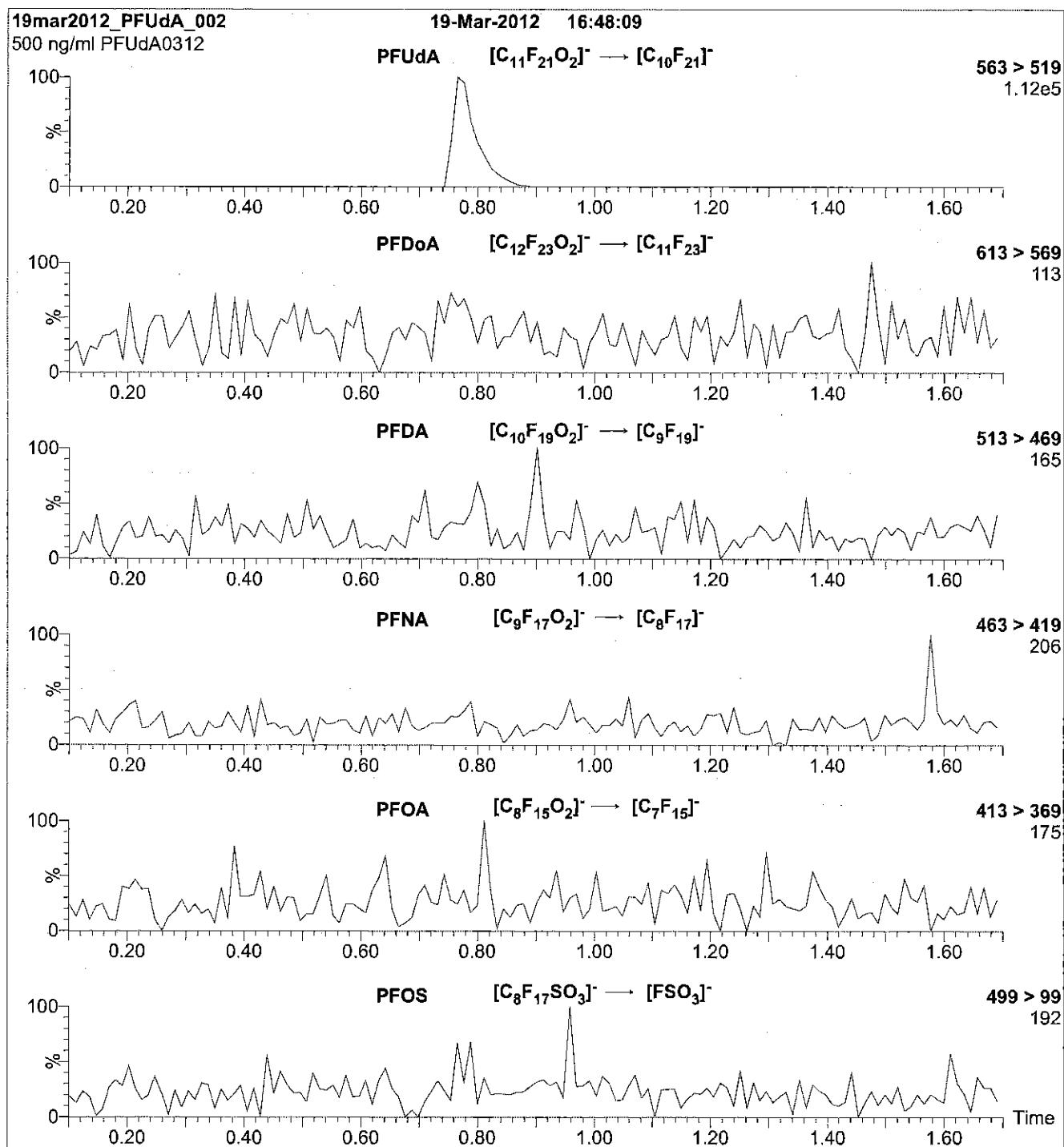
Flow: 300 µl/min

**MS Parameters**

Experiment: Full Scan (250 - 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)

**MS Parameters**

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

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

# **PFC\_IDA**

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**Perfluorinated Hydrocarbons**

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFOA #	PFOS #
B7/10-GW02	320-10459-1	85	116
B7/10-GW03S	320-10459-2	64	108
QR4-GW02	320-10459-3	90	109
B7/10-GW06S	320-10459-4	67	108
H123-GW02	320-10459-5	89	107
EFA-FD01-1014	320-10459-6	82	109
	MB 320-59712/1-A	109	111
	LCS 320-59712/2-A	98	101

PFOA = 13C4 PFOA  
PFOS = 13C4 PFOS

QC LIMITS  
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.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Matrix: Water Level: Low Lab File ID: 03DEC14A4A\_083.d  
Lab ID: LCS 320-59712/2-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ng/L)	LCS CONCENTRATION (ng/L)	LCS % REC	QC LIMITS REC	#
Perfluorooctanoic acid (PFOA)	40.0	50.4	126	60-140	
Perfluorooctane Sulfonate (PFOS)	38.2	39.7	104	60-140	
13C4 PFOS	95.6	96.3	101	25-150	
13C4 PFOA	100	98.0	98	25-150	

# Column to be used to flag recovery and RPD values

FORM III WS-LC-0025

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Lab File ID: 03DEC14A4A\_082.d Lab Sample ID: MB 320-59712/1-A  
Matrix: Water Date Extracted: 12/03/2014 13:38  
Instrument ID: A4 Date Analyzed: 12/04/2014 20:03  
Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-59712/2-A	03DEC14A4A_083.d	12/04/2014 20:24
B7/10-GW02	320-10459-1	03DEC14A4A_084.d	12/04/2014 20:45
B7/10-GW03S	320-10459-2	03DEC14A4A_085.d	12/04/2014 21:07
QR4-GW02	320-10459-3	03DEC14A4A_086.d	12/04/2014 21:28
B7/10-GW06S	320-10459-4	03DEC14A4A_087.d	12/04/2014 21:49
H123-GW02	320-10459-5	03DEC14A4A_088.d	12/04/2014 22:10
EFA-FD01-1014	320-10459-6	03DEC14A4A_089.d	12/04/2014 22:31

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>B7/10-GW02</u>	Lab Sample ID: <u>320-10459-1</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_084.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/15/2014 12:05</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>448.35 (mL)</u>	Date Analyzed: <u>12/04/2014 20:45</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	10	H	2.2	1.7	0.83
1763-23-1	Perfluorooctane Sulfonate (PFOS)	6.6	H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	116		25-150
STL00990	13C4 PFOA	85		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_084.d  
 Lims ID: 320-10459-A-1-B Lab Sample ID: 320-10459-1  
 Client ID: B7/10-GW02  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 20:45:53 ALS Bottle#: 24 Worklist Smp#: 48  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-1-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICAL File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

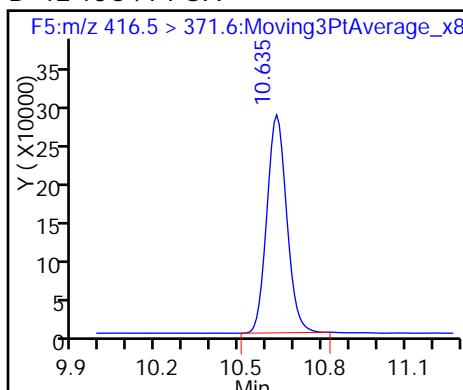
416.5 > 371.6	10.635	10.628	0.007		1440812	42.6		85.2	1974
13 Perfluorooctanoic acid									
412.8 > 368.8	10.635	10.628	0.007	1.000	226107	4.70			101
412.8 > 168.7	10.635	10.628	0.007	1.000	70963		3.19(0.00-0.00)		69.2

## D 16 13C4 PFOS

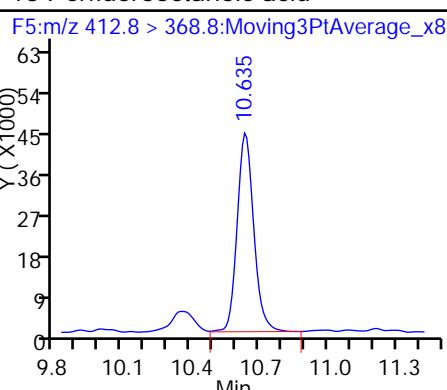
502.4 > 79.7	11.602	11.595	0.007		2500324	55.4		116	4350
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	279742	2.96			25.9
498.9 > 98.7	11.602	11.595	0.007	1.000	170535		1.64(0.00-0.00)		56.4

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_084.d  
 Injection Date: 04-Dec-2014 20:45:53 Instrument ID: A4  
 Lims ID: 320-10459-A-1-B Lab Sample ID: 320-10459-1  
 Client ID: B7/10-GW02  
 Operator ID: JRB ALS Bottle#: 24 Worklist Smp#: 48  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Method: PFAC\_A4 Limit Group: LC PFC ICAL

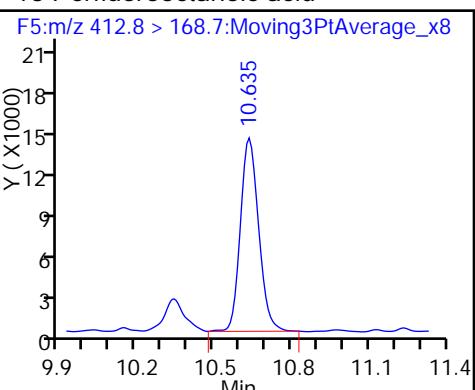
## D 12 13C4 PFOA



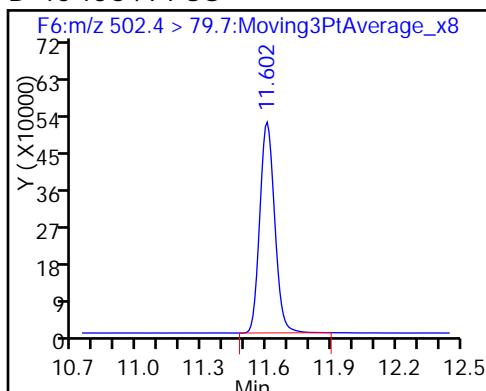
## 13 Perfluorooctanoic acid



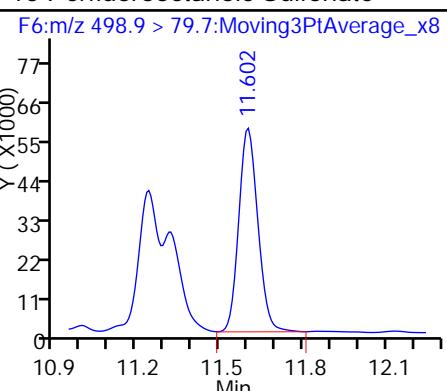
## 13 Perfluorooctanoic acid



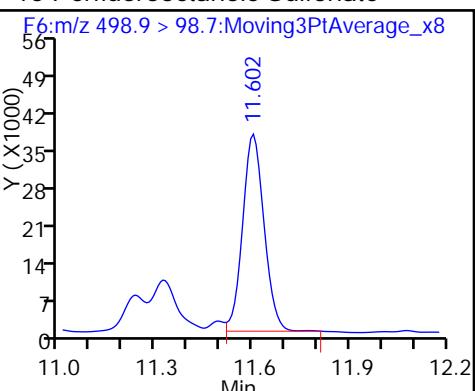
## D 16 13C4 PFOS



## 15 Perfluorooctanoic Sulfonate



## 15 Perfluorooctanoic Sulfonate



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>B7/10-GW03S</u>	Lab Sample ID: <u>320-10459-2</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_085.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/15/2014 15:20</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>427.75 (mL)</u>	Date Analyzed: <u>12/04/2014 21:07</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	2.0	J H	2.3	1.8	0.87
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.8	1.5

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	108		25-150
STL00990	13C4 PFOA	64		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_085.d  
 Lims ID: 320-10459-A-2-B Lab Sample ID: 320-10459-2  
 Client ID: B7/10-GW03S  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 21:07:07 ALS Bottle#: 25 Worklist Smp#: 49  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-2-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

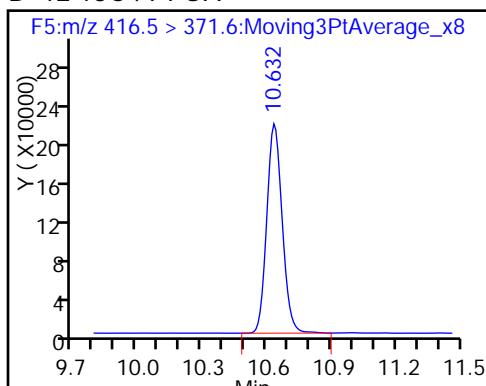
416.5 > 371.6	10.632	10.628	0.004		1083559	32.0		64.1	2381
13 Perfluorooctanoic acid									
412.8 > 368.8	10.641	10.628	0.013	1.000	30849	0.8525		19.0	
412.8 > 168.7	10.641	10.628	0.013	1.000	10104		3.05(0.00-0.00)		10.9

## D 16 13C4 PFOS

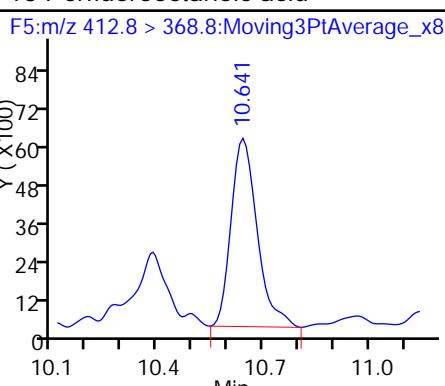
502.4 > 79.7	11.602	11.595	0.007		2323120	51.5		108	3478
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	35729	0.4074		23.1	
498.9 > 98.7	11.602	11.595	0.007	1.000	15556		2.30(0.00-0.00)		18.4

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_085.d  
 Injection Date: 04-Dec-2014 21:07:07 Instrument ID: A4  
 Lims ID: 320-10459-A-2-B Lab Sample ID: 320-10459-2  
 Client ID: B7/10-GW03S  
 Operator ID: JRB ALS Bottle#: 25 Worklist Smp#: 49  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Method: PFAC\_A4 Limit Group: LC PFC ICAL

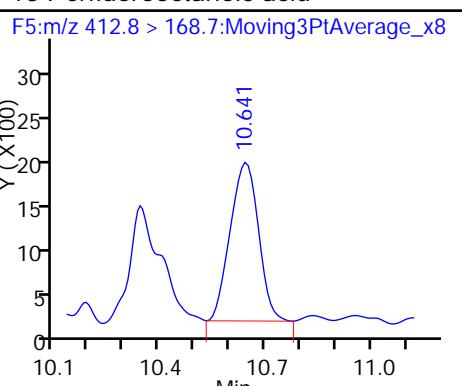
## D 12 13C4 PFOA



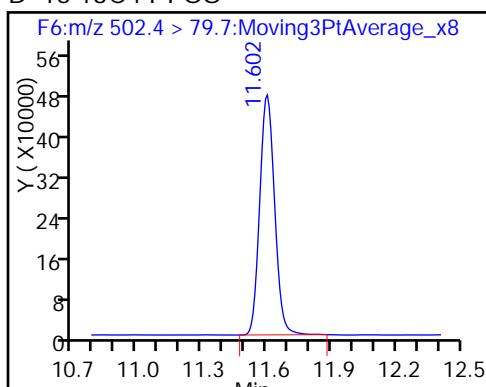
## 13 Perfluorooctanoic acid



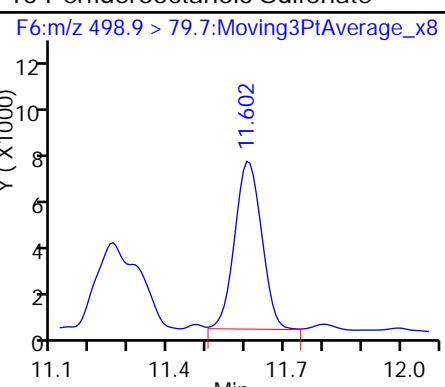
## 13 Perfluorooctanoic acid



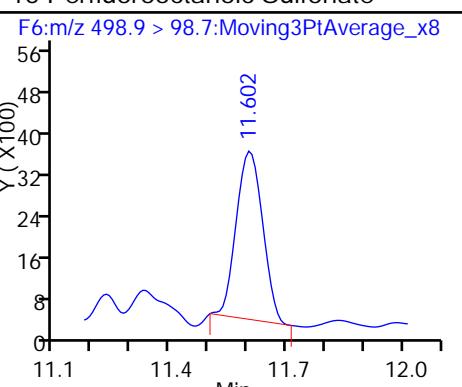
## D 16 13C4 PFOS



## 15 Perfluorooctanoic Sulfonate



## 15 Perfluorooctanoic Sulfonate



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>QR4-GW02</u>	Lab Sample ID: <u>320-10459-3</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_086.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 10:20</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>427.31 (mL)</u>	Date Analyzed: <u>12/04/2014 21:28</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	1.8	U H	2.3	1.8	0.88
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.8	1.5

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	109		25-150
STL00990	13C4 PFOA	90		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_086.d  
 Lims ID: 320-10459-A-3-B Lab Sample ID: 320-10459-3  
 Client ID: QR4-GW02  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 21:28:19 ALS Bottle#: 26 Worklist Smp#: 50  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-3-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICAL File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

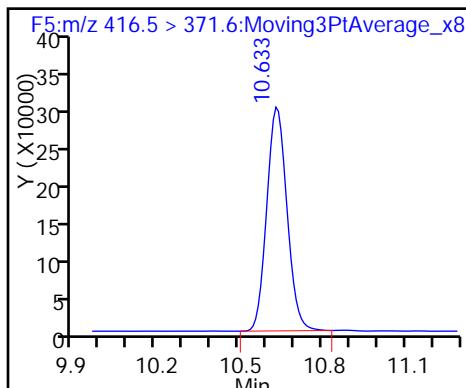
416.5 > 371.6	10.633	10.628	0.005		1518512	44.9		89.8	2377
13 Perfluorooctanoic acid									
412.8 > 368.8	10.633	10.628	0.005	1.000	17704	0.3491			9.9
412.8 > 168.7	10.642	10.628	0.014	1.001	7303		2.42(0.00-0.00)		14.9

## D 16 13C4 PFOS

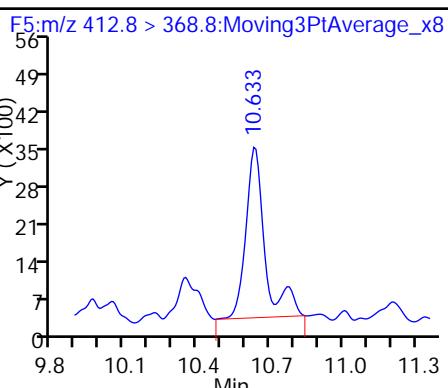
502.4 > 79.7	11.602	11.595	0.007		2356329	52.2		109	4306
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	22351	0.2512			29.9
498.9 > 98.7	11.602	11.595	0.007	1.000	11635		1.92(0.00-0.00)		10.4

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_086.d  
 Injection Date: 04-Dec-2014 21:28:19      Instrument ID: A4  
 Lims ID: 320-10459-A-3-B      Lab Sample ID: 320-10459-3  
 Client ID: QR4-GW02  
 Operator ID: JRB      ALS Bottle#: 26      Worklist Smp#: 50  
 Injection Vol: 15.0 ul      Dil. Factor: 1.0000  
 Method: PFAC\_A4      Limit Group: LC PFC ICAL

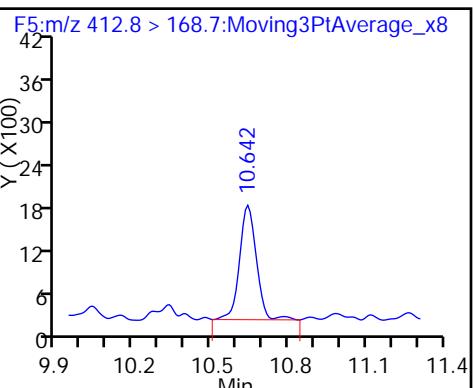
## D 12 13C4 PFOA



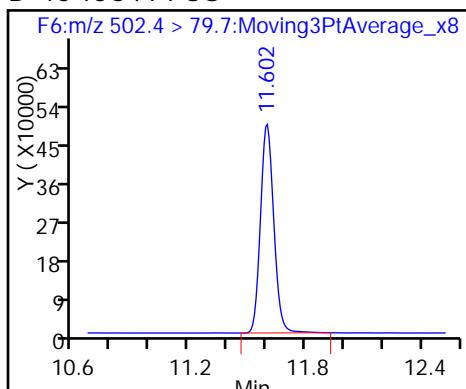
## 13 Perfluorooctanoic acid



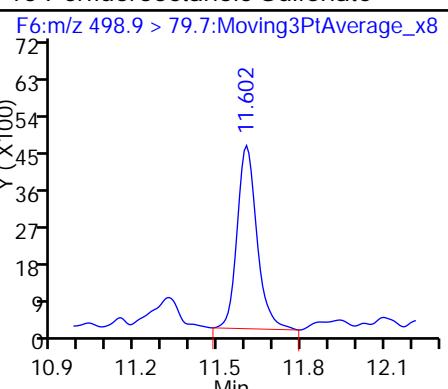
## 13 Perfluorooctanoic acid



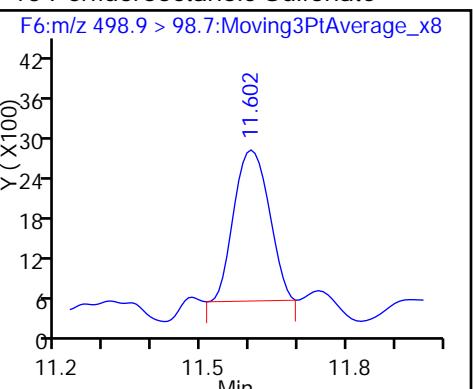
## D 16 13C4 PFOS



## 15 Perfluorooctanoic Sulfonate



## 15 Perfluorooctanoic Sulfonate



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>B7/10-GW06S</u>	Lab Sample ID: <u>320-10459-4</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_087.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 11:50</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>446.26 (mL)</u>	Date Analyzed: <u>12/04/2014 21:49</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	20	H	2.2	1.7	0.84
1763-23-1	Perfluorooctane Sulfonate (PFOS)	2.2	H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	108		25-150
STL00990	13C4 PFOA	67		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_087.d  
 Lims ID: 320-10459-A-4-B Lab Sample ID: 320-10459-4  
 Client ID: B7/10-GW06S  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 21:49:30 ALS Bottle#: 27 Worklist Smp#: 51  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-4-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

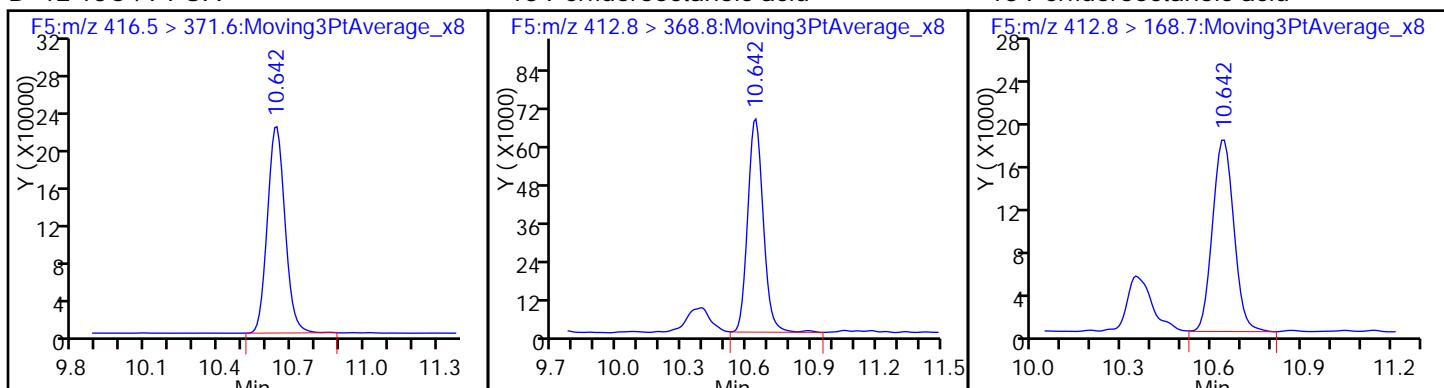
416.5 > 371.6	10.642	10.628	0.014		1138168	33.6		67.3	1719
13 Perfluorooctanoic acid									
412.8 > 368.8	10.642	10.628	0.014	1.000	340395	8.96			107
412.8 > 168.7	10.642	10.628	0.014	1.000	89169		3.82(0.00-0.00)		61.5

## D 16 13C4 PFOS

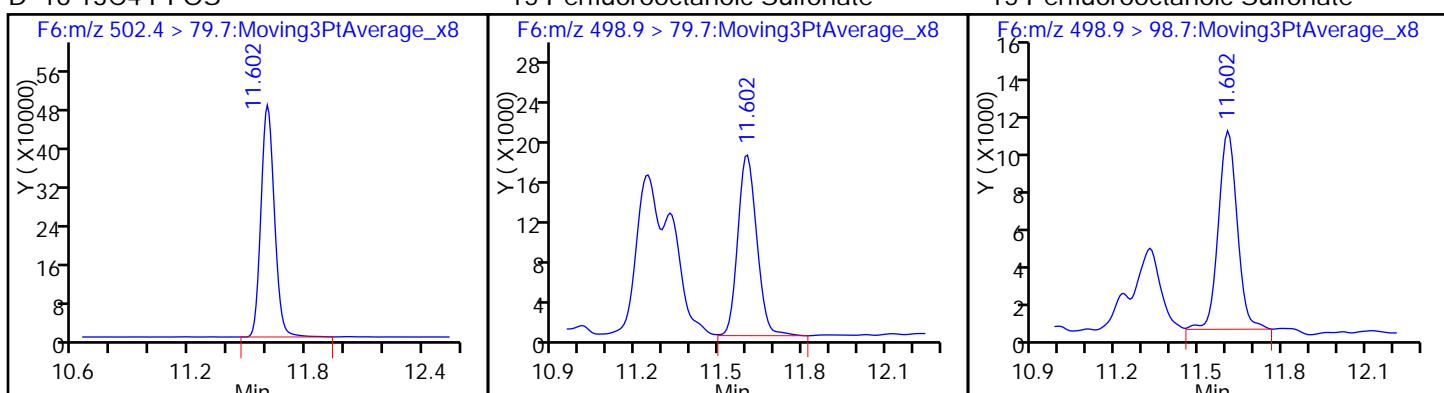
502.4 > 79.7	11.602	11.595	0.007		2338394	51.9		108	3471
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	87815	0.99			17.4
498.9 > 98.7	11.602	11.595	0.007	1.000	51215		1.71(0.00-0.00)		31.1

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_087.d  
 Injection Date: 04-Dec-2014 21:49:30 Instrument ID: A4  
 Lims ID: 320-10459-A-4-B Lab Sample ID: 320-10459-4  
 Client ID: B7/10-GW06S  
 Operator ID: JRB ALS Bottle#: 27 Worklist Smp#: 51  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Method: PFAC\_A4 Limit Group: LC PFC ICAL

## D 12 13C4 PFOA



## D 16 13C4 PFOS



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>H123-GW02</u>	Lab Sample ID: <u>320-10459-5</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_088.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 15:00</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>455.47 (mL)</u>	Date Analyzed: <u>12/04/2014 22:10</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	17	H	2.2	1.6	0.82
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.6	U H	2.2	1.6	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	107		25-150
STL00990	13C4 PFOA	89		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_088.d  
 Lims ID: 320-10459-A-5-B Lab Sample ID: 320-10459-5  
 Client ID: H123-GW02  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 22:10:42 ALS Bottle#: 28 Worklist Smp#: 52  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-5-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

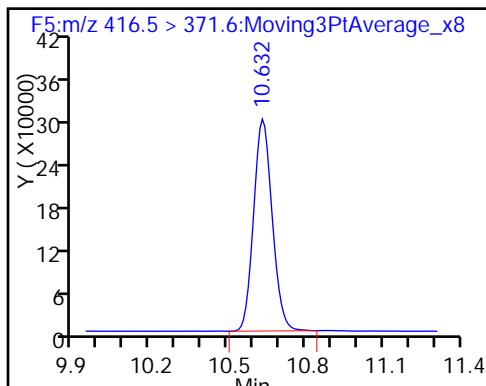
416.5 > 371.6	10.632	10.628	0.004		1513765	44.7		89.5	2011
13 Perfluorooctanoic acid									
412.8 > 368.8	10.632	10.628	0.004	1.000	391049	7.74			109
412.8 > 168.7	10.632	10.628	0.004	1.000	117288		3.33(0.00-0.00)		61.6

## D 16 13C4 PFOS

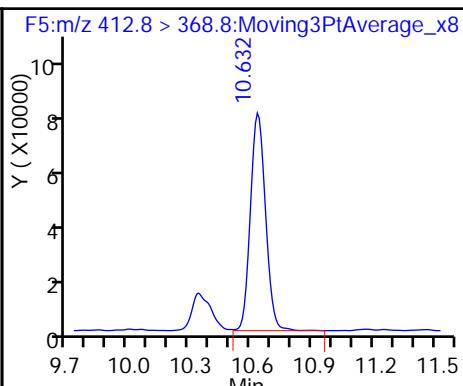
502.4 > 79.7	11.594	11.595	-0.001		2301909	51.0		107	5627
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	46613	0.5364			15.9
498.9 > 98.7	11.594	11.595	-0.001	0.999	22984		2.03(0.00-0.00)		22.7

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_088.d  
 Injection Date: 04-Dec-2014 22:10:42      Instrument ID: A4  
 Lims ID: 320-10459-A-5-B      Lab Sample ID: 320-10459-5  
 Client ID: H123-GW02  
 Operator ID: JRB      ALS Bottle#: 28      Worklist Smp#: 52  
 Injection Vol: 15.0 ul      Dil. Factor: 1.0000  
 Method: PFAC\_A4      Limit Group: LC PFC ICAL

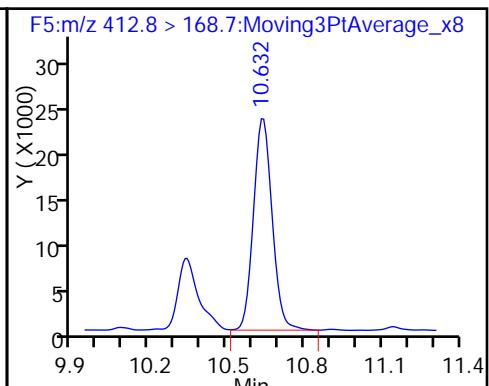
## D 12 13C4 PFOA



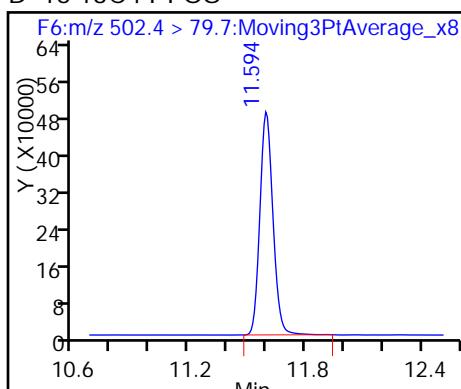
## 13 Perfluorooctanoic acid



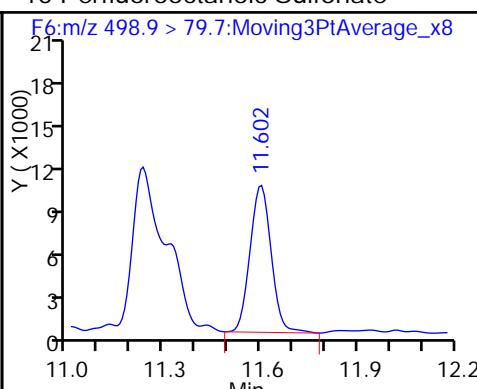
## 13 Perfluorooctanoic acid



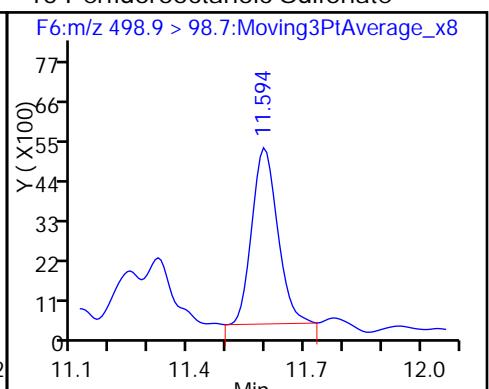
## D 16 13C4 PFOS



## 15 Perfluorooctanoic Sulfonate



## 15 Perfluorooctanoic Sulfonate



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>EFA-FD01-1014</u>	Lab Sample ID: <u>320-10459-6</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_089.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 00:00</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>447.90 (mL)</u>	Date Analyzed: <u>12/04/2014 22:31</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	16	H	2.2	1.7	0.84
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.7	U H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	109		25-150
STL00990	13C4 PFOA	82		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_089.d  
 Lims ID: 320-10459-A-6-B Lab Sample ID: 320-10459-6  
 Client ID: EFA-FD01-1014  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 22:31:53 ALS Bottle#: 29 Worklist Smp#: 53  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-6-b 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

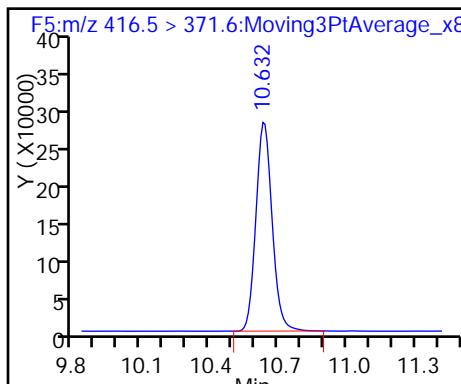
416.5 > 371.6	10.632	10.628	0.004		1381817	40.8		81.7	1956
13 Perfluorooctanoic acid									
412.8 > 368.8	10.632	10.628	0.004	1.000	332759	7.21			105
412.8 > 168.7	10.642	10.628	0.014	1.001	99306		3.35(0.00-0.00)		57.8

## D 16 13C4 PFOS

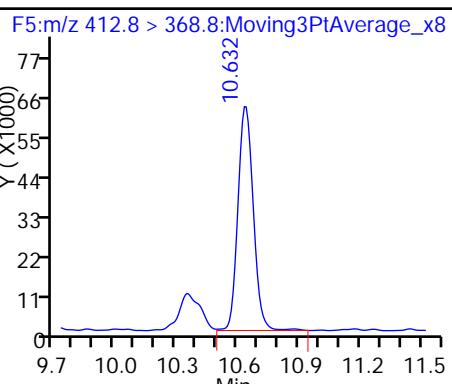
502.4 > 79.7	11.602	11.595	0.007		2348037	52.1		109	4014
15 Perfluorooctanoic Sulfonate									
498.9 > 79.7	11.602	11.595	0.007	1.000	48141	0.5431			14.7
498.9 > 98.7	11.610	11.595	0.015	1.001	30527		1.58(0.00-0.00)		28.2

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_089.d  
 Injection Date: 04-Dec-2014 22:31:53 Instrument ID: A4  
 Lims ID: 320-10459-A-6-B Lab Sample ID: 320-10459-6  
 Client ID: EFA-FD01-1014  
 Operator ID: JRB ALS Bottle#: 29 Worklist Smp#: 53  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Method: PFAC\_A4 Limit Group: LC PFC ICAL

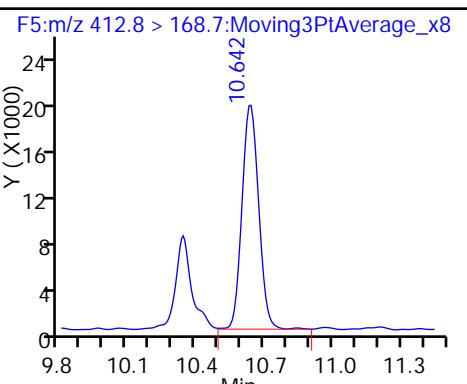
## D 12 13C4 PFOA



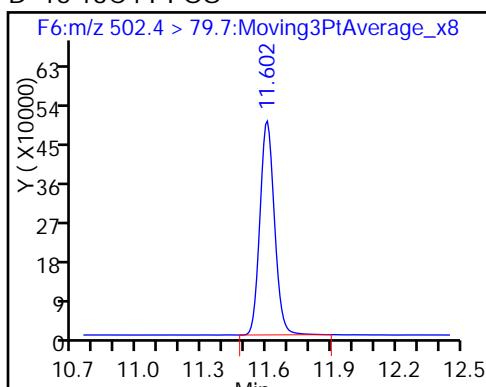
## 13 Perfluorooctanoic acid



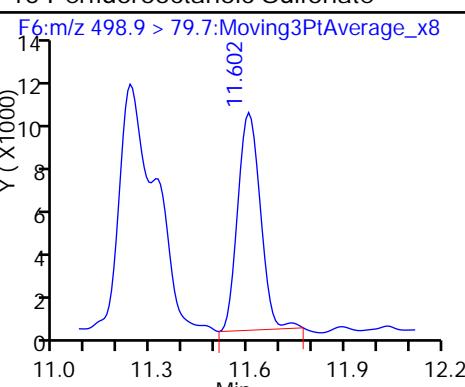
## 13 Perfluorooctanoic acid



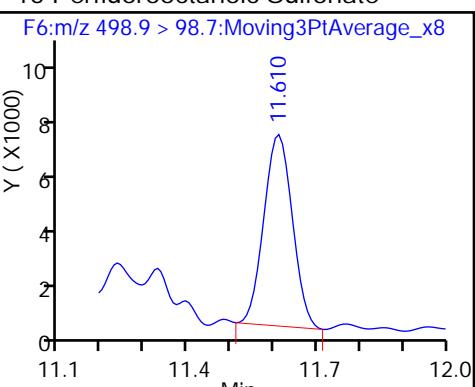
## D 16 13C4 PFOS



## 15 Perfluorooctanoic Sulfonate



## 15 Perfluorooctanoic Sulfonate



FORM VI  
LCMS INITIAL CALIBRATION DATA  
EXTERNAL STANDARD RETENTION TIME SUMMARY

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1 Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.d

ANALYTE	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6	LVL 7				RT WINDOW	AVG RT
13C4 PFBA	5.934	5.937	5.928	5.934	5.934	5.934	5.934				5.684 - 6.184	5.934
13C5 PFFeA	7.035	7.035	7.029	7.043	7.046	7.036	7.040				6.788 - 7.288	7.038
13C2 PFHxA	8.273	8.284	8.273	8.290	8.284	8.284	8.286				8.032 - 8.532	8.282
13C4-PFHxA	9.505	9.510	9.499	9.505	9.510	9.505	9.510				9.256 - 9.756	9.506
18O2 PFHxS	9.538	9.545	9.538	9.545	9.545	9.545	9.552				9.294 - 9.794	9.544
13C4 PFOA	10.623	10.632	10.623	10.632	10.633	10.632	10.623				10.378 - 10.878	10.628
13C4 PFOS	11.585	11.602	11.594	11.594	11.594	11.602	11.594				11.345 - 11.845	11.595
13C5 PFNA	11.603	11.620	11.611	11.612	11.620	11.620	11.612				11.364 - 11.864	11.614
13C2 PFDA	12.455	12.466	12.455	12.466	12.466	12.466	12.466				12.213 - 12.713	12.463
13C8 FOSA	12.994	13.003	12.994	12.994	12.995	13.003	13.004				12.748 - 13.248	12.998
13C2 PFUnA	13.189	13.198	13.189	13.198	13.198	13.198	13.206				12.948 - 13.448	13.198
13C2 PFDmA	13.802	13.811	13.802	13.811	13.811	13.811	13.813				13.559 - 14.059	13.809
13C2-PFTeDA	14.765	14.765	14.765	14.765	14.765	14.772	14.766				14.516 - 15.016	14.766
13C2-PFHxDA	15.414	15.414	15.415	15.415	15.415	15.415	15.415				15.165 - 15.665	15.415

FORM VI  
LCMS INITIAL CALIBRATION DATA  
EXTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 5	LVL 2 LVL 6	LVL 3 LVL 7	LVL 4		B	M1	M2								
13C4 PFBA	5268.6 3806.7	5812.9 3203.9	5254.9 2254.3	4338.2	Ave		42777.06429				30.0		50.0			
13C5 PFPeA	9419.7 7332.0	9580.9 5721.7	8177.5 4869.4	8903.4	Ave		7714.90000				24.0		50.0			
13C2 PFHxA	21470 18141	22784 16317	24835 11422	18449	Ave		19059.8143				23.0		50.0			
13C4-PFHxA	21643 17882	21390 11698	21748 10393	19442	Ave		17742.3500				27.0		50.0			
18O2 PFHxS	53128 47816	52233 33678	50966 25618	51213	Ave		44950.3322				24.0		50.0			
13C4 PFOA	41322 31379	41400 24644	40475 19047	38546	Ave		33830.5286				27.0		50.0			
13C4 PFOS	54398 44130	54385 33330	56166 23197	50088	Ave		45099.0735				28.0		50.0			
13C5 PFNA	44879 37219	45410 30155	48879 22012	41831	Ave		38626.2857				25.0		50.0			
13C2 PFDA	63235 48806	61430 37175	59297 32265	58211	Ave		51488.2929				24.0		50.0			
13C8 FOSA	150505 147550	155647 126255	158874 106829	155831	Ave		143069.894				14.0		50.0			
13C2 PFUnA	66511 59868	66532 58833	66894 69853	64772	Ave		64751.6643				6.2		50.0			
13C2 PFDoA	76991 59932	66196 56050	76655 52807	68862	Ave		65356.2357				15.0		50.0			
13C2-PFTeDA	77915 69707	76286 61208	80240 58365	76834	Ave		71507.9143				12.0		50.0			
13C2-PFHxDA	96546 111182	105842 105984	112387 93820	103841	Ave		104228.936				6.6		50.0			

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

Analy Batch No.: 59779

SDG No.:

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorobutanoic acid (PFBA)	10678 16623	22829 12169	21990	19527	19477	AveID		2.3514				25.0		35.0			
Perfluoropentanoic acid (PFPeA)	8732.0 11185	14944 9458.0	14198	14832	13130	AveID		1.6538				21.0		35.0			
Perfluorobutane Sulfonate (PFBS)	14683 39320	46758 32165	49055	48250	67174	AveID		0.9863				37.0		50.0			
Perfluorohexanoic acid (PFHxA)	9952.0 16870	21004 12521	26818	19955	20968	AveID		0.9761				24.0		35.0			
PFPeS (Perflouro-1-pentanesulfonate)	12755 32980	40229 24854	44086	42428	40598	AveID		0.7860				32.0		50.0			
Perfluoroheptanoic acid (PFHpA)	14724 21344	28535 18892	35883	33256	29606	AveID		1.5247				27.0		35.0			
Perfluorohexane Sulfonate (PFHxS)	24632 33051	52072 26648	53418	49529	45799	AveID		0.9222				22.0		35.0			
Perfluorooctanoic acid (PFOA)	54920 43078	74889 33366	68131	62497	54775	AveID		1.6697				9.7		35.0			
Perfluoro-1-heptanesulfonate (PFHps)	12752 33076	45349 24051	51122	49900	46278	AveID		0.8647				33.0		50.0			
Perfluorooctane Sulfonate (PFOS)	143245 49883	124703 36721	92010	73648	66987	AveID		1.8046				26.0		35.0			
Perfluorononanoic acid (PFNA)	22096 39342	60944 27244	61254	52411	44283	AveID		1.1532				26.0		35.0			
PFNS (Perflouro-1-nonanesulfonate)	16554 25936	40576 18531	47018	41535	37537	AveID		0.7349				26.0		50.0			
Perfluorodecanoic acid (PFDA)	16398 41090	58072 33184	67070	63478	54711	AveID		0.9544				33.0		35.0			
Perfluorooctane Sulfonamide (FOSA)	60576 132015	157092 116050	155459	165246	157618	AveID		0.9501				26.0		35.0			
Perfluorododecane sulfonate (PFDS)	10237 23088	37158 16598	42628	36833	33233	AveID		0.6467				32.0		50.0			
Perfluoroundecanoic acid (PFUnA)	25356 49828	75233 40491	72252	72984	62389	AveID		0.8839				34.0		35.0			
Perfluorododecanoic acid (PFDoA)	25100 62748	61887 50239	92219	78804	73401	AveID		0.9863				32.0		35.0			
PFDoS (Perflouro-1-dodecanesulfonate)	7778.9 23770	32535 17735	43717	40638	35319	AveID		0.6584				36.0		50.0			
Perfluorotridecanoic Acid (PFTriA)	20248 45661	65903 40208	67545	66220	62650	AveID		0.8175				32.0		50.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1 Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorotetradecanoic acid (PFTeA)	15018 30898	36572 27633	43161	42497	38214	AveID		0.5200				29.0		50.0			
Perfluoro-n-hexadecanoic acid (PFHxDA)	44032 92953	98406 78170	100981	86858	96066	AveID		1.3398				27.0		50.0			
Perfluoro-n-octadecanoic acid (PFODA)	22762 104163	94742 95748	99443	104957	102973	AveID		1.4197				38.0		50.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
EXTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1 Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.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	105372 64078	116258 45086	105097	86764	76134	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C5 PFPeA	Ave	188393 114434	191617 97387	163549	178067	146639	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C2 PFHxA	Ave	429395 326332	455689 228440	496708	368981	362829	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C4-PFHxA	Ave	432854 233963	427809 207868	434956	388840	357639	20.0 20.0	20.0 20.0	20.0	20.0	20.0
18O2 PFHxS	Ave	1005188 637190	988251 484686	964272	968949	904686	18.9 18.9	18.9 18.9	18.9	18.9	18.9
13C4 PFOA	Ave	2066105 1232205	2070020 952358	2023751	1927321	1568925	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C4 PFOS	Ave	2600203 1593174	2599599 1108797	2684739	2394219	2109419	47.8 47.8	47.8 47.8	47.8	47.8	47.8
13C5 PFNA	Ave	897580 603092	908195 440238	977581	836616	744378	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C2 PFDA	Ave	1264690 743492	1228601 645298	1185934	1164228	976118	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C8 FOSA	Ave	7525230 6312732	7782345 5341425	7943692	7791526	7377513	50.0 50.0	50.0 50.0	50.0	50.0	50.0
13C2 PFUnA	Ave	1330212 1176651	1330638 1397054	1337877	1295448	1197353	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C2 PFDoA	Ave	1539828 1120992	1323923 1056139	1533105	1377246	1198640	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C2-PFTeDA	Ave	1558303 1224157	1525719 1167304	1604800	1536682	1394143	20.0 20.0	20.0 20.0	20.0	20.0	20.0
13C2-PFHxDA	Ave	1930922 2119674	2116849 1876393	2247748	2076818	2223647	20.0 20.0	20.0 20.0	20.0	20.0	20.0

Curve Type Legend:

Ave = Average

FORM VI  
LCMS INITIAL CALIBRATION DATA  
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.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	5339 3324570	22829 6084630	109952	390544	973858	0.500 200	1.00 500	5.00	20.0	50.0
Perfluoropentanoic acid (PFPeA)		AveID	4366 2237092	14944 4729024	70990	296633	656489	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorobutane Sulfonate (PFBS)		AveID	6490 6951713	41334 14216894	216823	853055	2969104	0.442 177	0.884 442	4.42	17.7	44.2
Perfluorohexanoic acid (PFHxA)		AveID	4976 3374043	21004 6260320	134092	399091	1048415	0.500 200	1.00 500	5.00	20.0	50.0
PFPes (Perflouro-1-pentanesulfonate)		AveID	5982 6187089	37735 11656614	206763	795956	1904029	0.469 188	0.938 469	4.69	18.8	46.9
Perfluoroheptanoic acid (PFHpA)		AveID	7362 4268821	28535 9446001	179416	665119	1480278	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorohexane Sulfonate (PFHxS)		AveID	11651 6253308	49260 12604320	252665	937096	2166299	0.473 189	0.946 473	4.73	18.9	47.3
Perfluoroctanoic acid (PFOA)		AveID	27460 8615547	74889 16683058	340656	1249944	2738772	0.500 200	1.00 500	5.00	20.0	50.0
Perfluoro-1-heptanesulfonate (PFHpS)		AveID	6070 6297711	43172 11448389	243339	950094	2202840	0.476 190	0.952 476	4.76	19.0	47.6
Perfluorooctane Sulfonate (PFOS)		AveID	68471 9537705	119216 17552751	439807	1408150	3201983	0.478 191	0.956 478	4.78	19.1	47.8
Perfluorononanoic acid (PFNA)		AveID	11048 7868399	60944 13621836	306271	1048226	2214146	0.500 200	1.00 500	5.00	20.0	50.0
PFNS (Perflouro-1-nonanesulfonate)		AveID	7946 4979770	38953 8895061	225687	797477	1801756	0.480 192	0.960 480	4.80	19.2	48.0
Perfluorodecanoic acid (PFDA)		AveID	8199 8217900	58072 16591989	335348	1269551	2735529	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorooctane Sulfonamide (FOSA)		AveID	30288 26403028	157092 58025028	777293	3304921	7880884	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorodecane sulfonate (PFDS)		AveID	4934 4451333	35820 8000344	205469	710141	1601818	0.482 193	0.964 482	4.82	19.3	48.2

FORM VI  
LCMS INITIAL CALIBRATION DATA  
RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1 Analy Batch No.: 59779

SDG No.: \_\_\_\_\_

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

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
Perfluoroundecanoic acid (PFUnA)		AveID	12678 9965686	75233 20245258	361261	1459686	3119455	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorododecanoic acid (PFDa)		AveID	12550 12549584	61887 25119659	461093	1576082	3670027	0.500 200	1.00 500	5.00	20.0	50.0
PFDoS (Perflouro-1-dodecanesulfonate)		AveID	3765 4601893	31494 8583795	211588	786746	1709417	0.484 194	0.968 484	4.84	19.4	48.4
Perfluorotridecanoic Acid (PFTriA)		AveID	10124 9132281	65903 20104094	337727	1324401	3132510	0.500 200	1.00 500	5.00	20.0	50.0
Perfluorotetradecanoic acid (PFTeA)		AveID	7509 6179569	36572 13816653	215805	849938	1910691	0.500 200	1.00 500	5.00	20.0	50.0
Perfluoro-n-hexadecanoic acid (PFHxDA)		AveID	22016 18590504	98406 39084771	504906	1737164	4803324	0.500 200	1.00 500	5.00	20.0	50.0
Perfluoro-n-octadecanoic acid (PFODA)		AveID	11381 20832627	94742 47874246	497216	2099144	5148634	0.500 200	1.00 500	5.00	20.0	50.0

Curve Type Legend:

AveID = Average isotope dilution

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_040.d  
 Lims ID: Std L1  
 Client ID:  
 Sample Type: IC Calib Level: 1  
 Inject. Date: 04-Dec-2014 03:44:16 ALS Bottle#: 4 Worklist Smp#: 4  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L1 PFC 0.50/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:51 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 1 113C4 PFBA

216.7 > 171.5	5.934	5.934	0.0		105372	24.6		123	409
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## 2 Perfluorobutyric acid

212.7 > 168.6	5.934	5.935	-0.001	1.000	5339	0.2410		48.2	19.7
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## 4 Perfluoropentanoic acid

262.9 > 218.7	7.039	7.037	0.002	1.000	4366	0.2803		56.1	7.2
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## D 3 113C5-PFPeA

267.6 > 222.7	7.035	7.038	-0.003		188393	24.4		122	572
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## 5 Perfluorobutane Sulfonate

298.8 > 79.6	7.154	7.157	-0.003	1.000	6490	0.1239		28.0	28.9
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298.8 > 98.6	7.147	7.157	-0.010	0.999	5511	1.18(0.00-0.00)		28.0	19.8
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## D 6 113C2 PFHxA

314.6 > 269.7	8.273	8.282	-0.009		429395	22.5		113	1739
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## 7 Perfluorohexanoic acid

312.9 > 268.7	8.279	8.284	-0.005	1.000	4976	0.2374		47.5	15.1
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## 22 PFPeS (Perflouro-1-pentanesulfonat

348.7 > 79.5	8.360	8.360	0.0	0.877	5982	0.1432		30.5	22.5
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## D 8 113C4-PFHxA

366.6 > 321.6	9.505	9.506	-0.001		432854	24.4		122	1684
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## 9 Perfluoroheptanoic acid

362.8 > 318.7	9.510	9.508	0.002	1.000	7362	0.2231		44.6	21.3
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## 10 Perfluorohexane Sulfonate

398.8 > 79.7	9.531	9.543	-0.012	1.000	11651	0.2378		50.3	36.2
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## D 11 118O2 PFHxA

402.5 > 83.6	9.538	9.544	-0.006		1005188	22.4		118	2708
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## D 12 113C4 PFOA

416.5 > 371.6	10.623	10.628	-0.005		2066105	61.1		122	3807
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Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.623	10.628	-0.005	1.000	27460	0.3980		79.6	41.1	
412.8 > 168.7	10.632	10.628	0.004	1.001	11257		2.44(0.00-0.00)	79.6	18.2	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.641	10.635	0.006	1.000	6070	0.1291		27.1	12.0	
D 16 13C4 PFOS										
502.4 > 79.7	11.585	11.595	-0.010		2600203	57.7		121	6261	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.585	11.595	-0.010	1.000	68471	0.6975		146	82.9	
498.9 > 98.7	11.585	11.595	-0.010	1.000	37166		1.84(0.00-0.00)	146	84.2	
D 17 13C5 PFNA										
467.5 > 422.6	11.603	11.614	-0.011		897580	23.2		116	2143	
18 Perfluorononanoic acid										
462.5 > 418.6	11.611	11.616	-0.005	1.000	11048	0.2135		42.7	21.7	
21 PFNS (Perflouro-1-nonanesulfonate)										
548.6 > 79.6	12.414	12.429	-0.015	1.000	7946	0.1988		41.4	18.1	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.445	12.461	-0.016	1.000	8199	0.1359		27.2	21.7	
D 19 13C2 PFDA										
514.4 > 469.5	12.455	12.463	-0.008		1264690	24.6		123	1953	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	12.984	12.998	-0.014	1.000	30288	0.2118		42.4	45.9	
D 23 13C8 FOSA										
505.4 > 77.6	12.994	12.998	-0.004		7525230	52.6		105	4617	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	4934	0.1402		29.1	8.8	
D 26 13C2 PFUnA										
564.3 > 519.5	13.189	13.198	-0.009		1330212	20.5		103	2524	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.189	13.198	-0.009	1.000	12678	0.2156		43.1	25.7	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.793	13.808	-0.015	1.000	12550	0.1653		33.1	7.8	
D 28 13C2 PFDoA										
614.4 > 569.4	13.802	13.809	-0.007		1539828	23.6		118	2135	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.265	14.272	-0.007	1.000	3765	0.1051		21.7	6.9	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.319	14.327	-0.008	1.000	10124	0.1608		32.2	7.6	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	7509	0.1876		37.5	7.0	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1558303	21.8		109	2502	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.414	15.415	-0.001		1930922	18.5		92.6	2349	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.414	15.415	-0.001	1.000	22016	0.2134		42.7	26.9	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.756	15.752	0.004	1.000	Page 11381 of 350	0.1041		20.8	151	12/12/2014

Report Date: 05-Dec-2014 10:01:51

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L1\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:01:51

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_040.d

Injection Date: 04-Dec-2014 03:44:16

Instrument ID: A4

Lims ID: Std L1

Client ID:

Operator ID: JRB

ALS Bottle#: 4 Worklist Smp#: 4

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

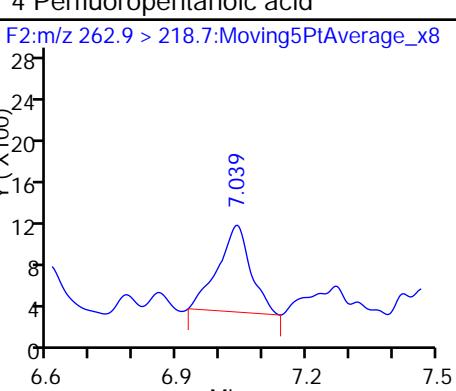
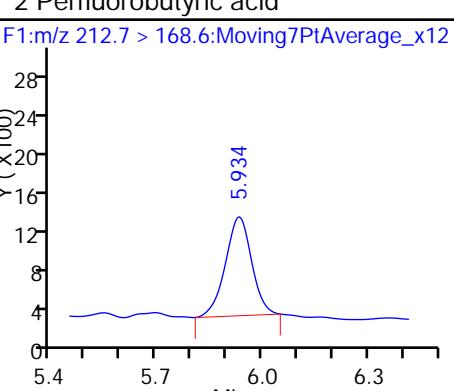
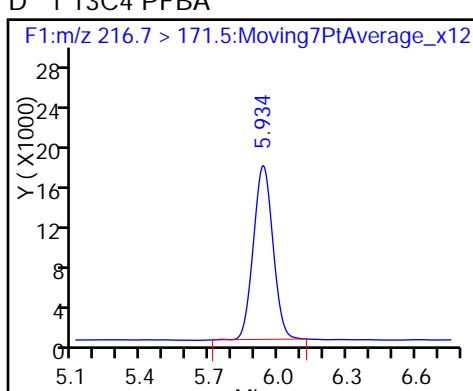
Method: PFAC\_A4

Limit Group: LC PFC ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

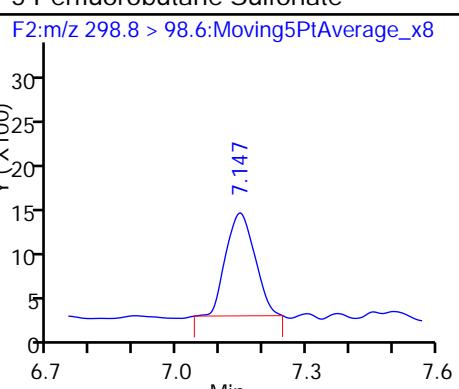
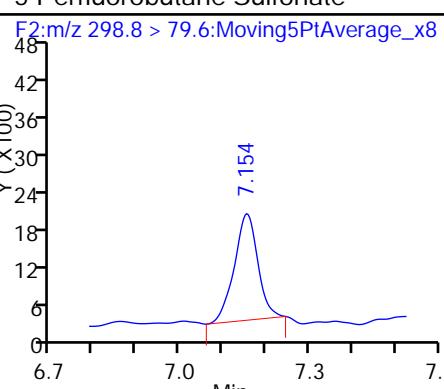
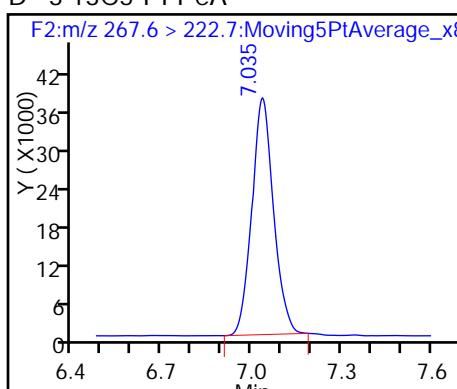
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

5 Perfluorobutane Sulfonate

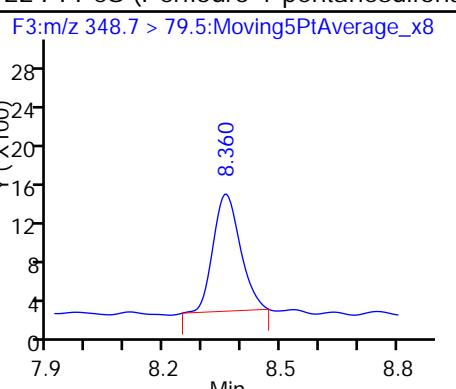
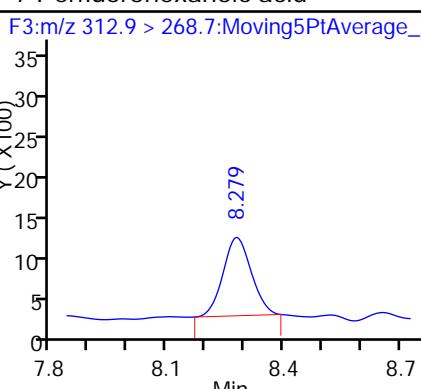
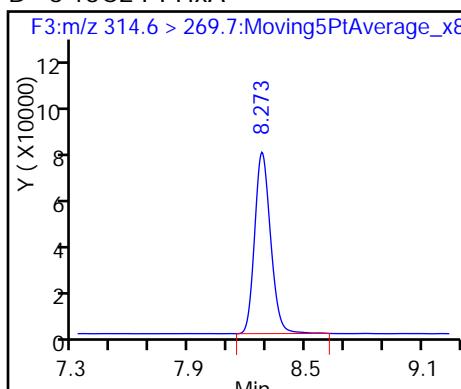
5 Perfluorobutane Sulfonate



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

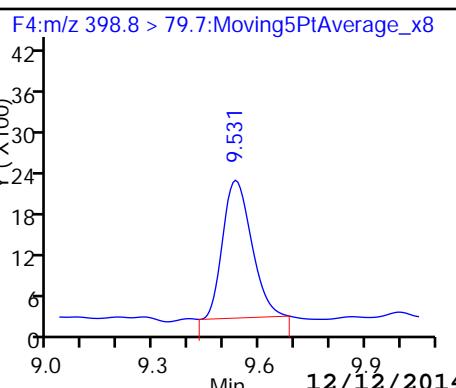
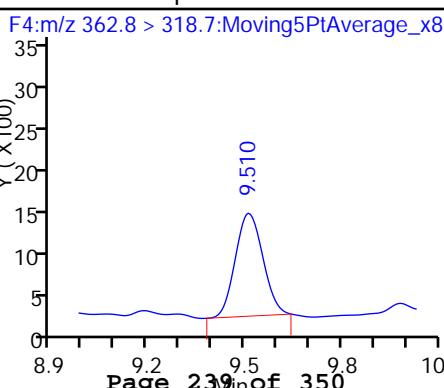
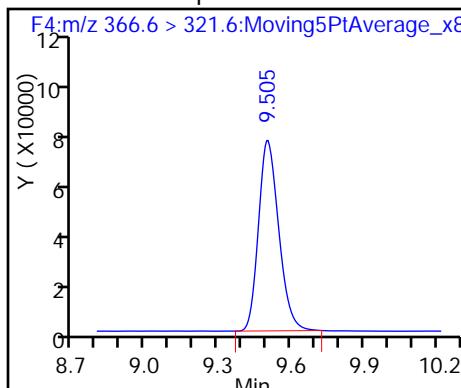
22 PFPeS (Perflouro-1-pentanesulfonat)



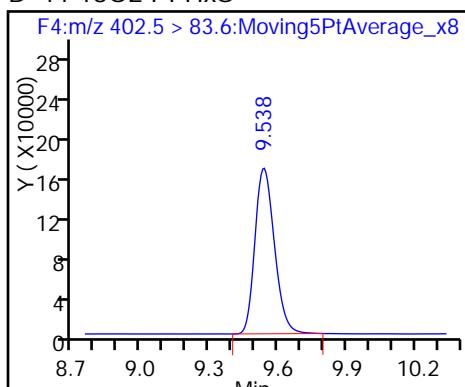
D 8 13C4-PFHxA

9 Perfluoroheptanoic acid

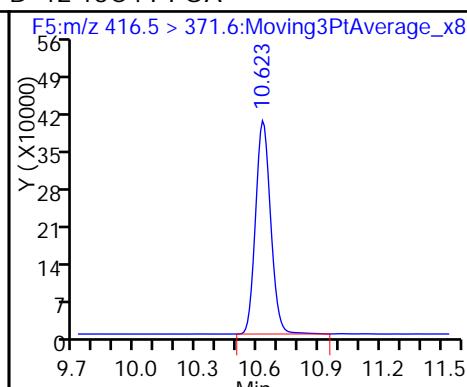
10 Perfluorohexane Sulfonate



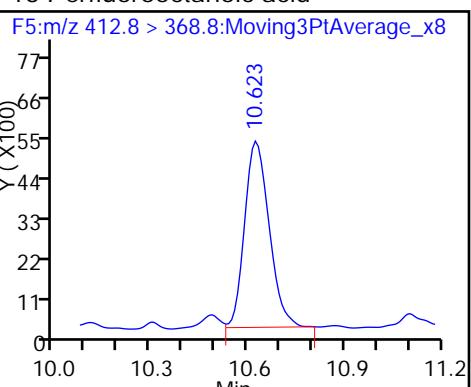
D 11 18O2 PFHxS



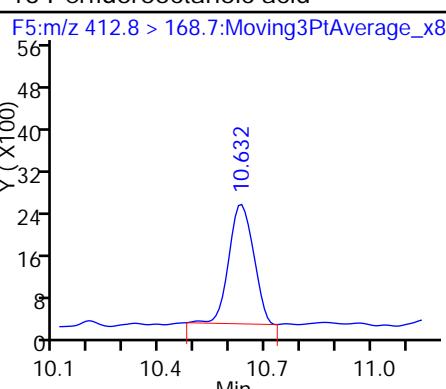
D 12 13C4 PFOA



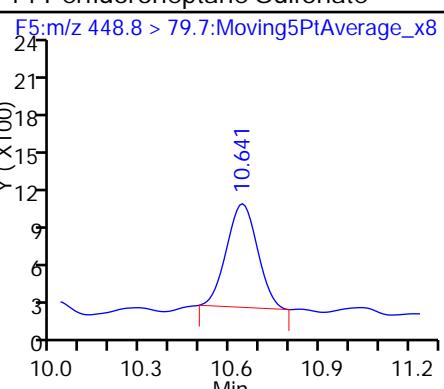
13 Perfluorooctanoic acid



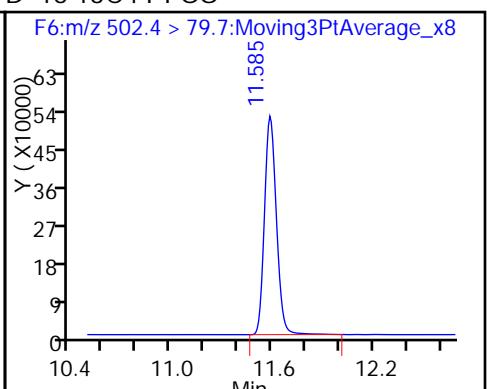
13 Perfluorooctanoic acid



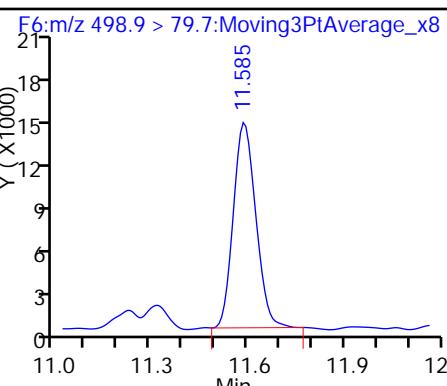
14 Perfluoroheptane Sulfonate



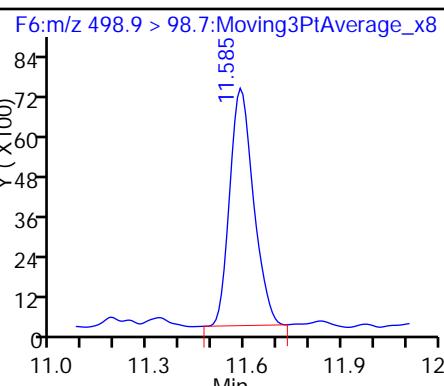
D 16 13C4 PFOS



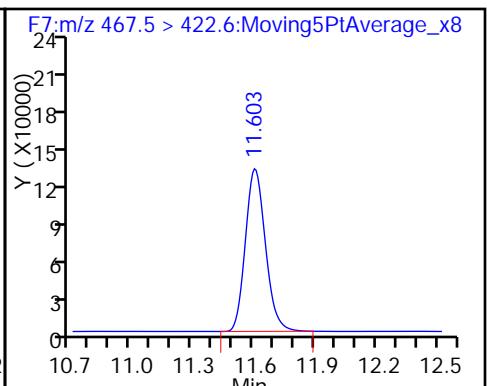
15 Perfluorooctanoic Sulfonate



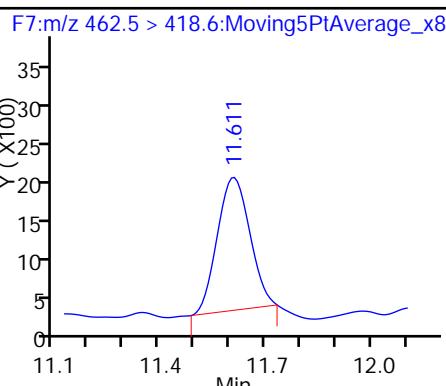
15 Perfluorooctanoic Sulfonate



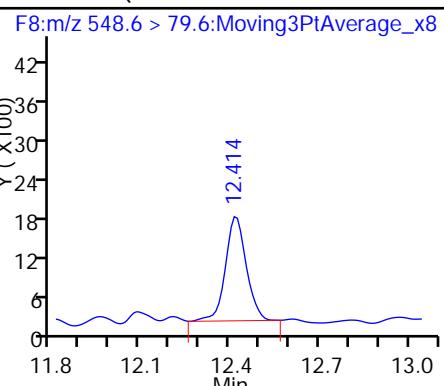
D 17 13C5 PFNA



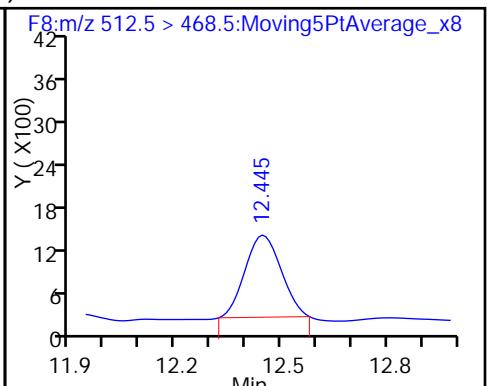
18 Perfluorononanoic acid



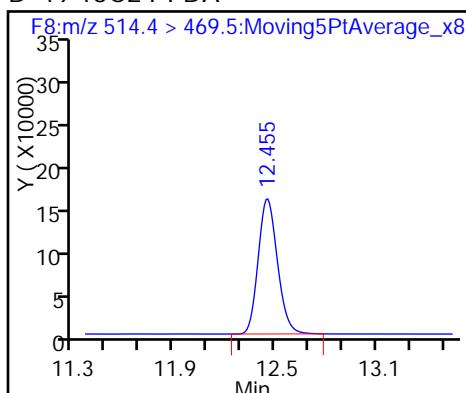
21 PFNS (Perflouro-1-nananesulfonate)



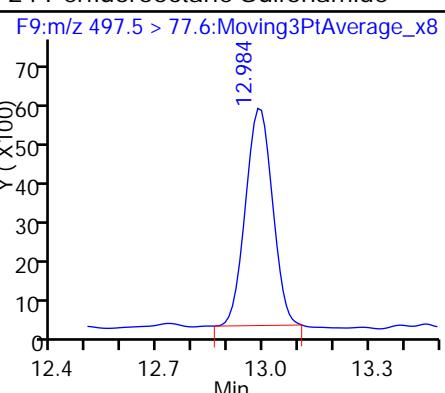
20 Perfluorodecanoic acid



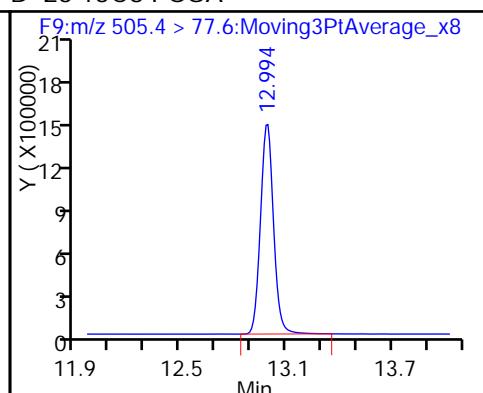
## D 19 13C2 PFDA



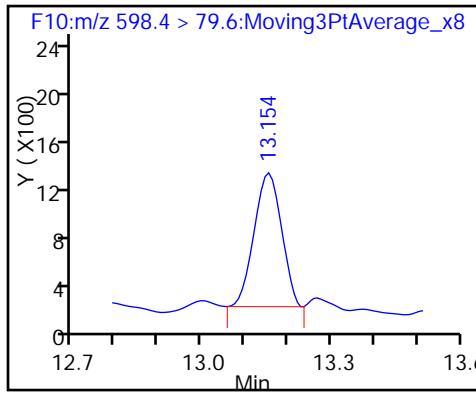
## 24 Perfluorooctane Sulfonamide



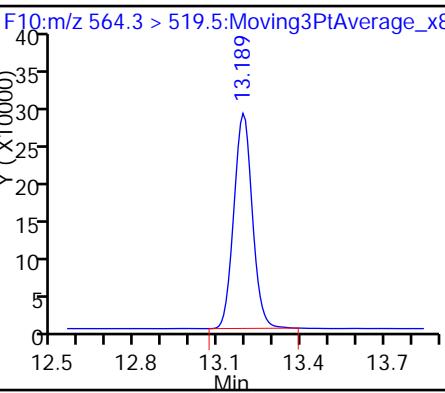
## D 23 13C8 FOSA



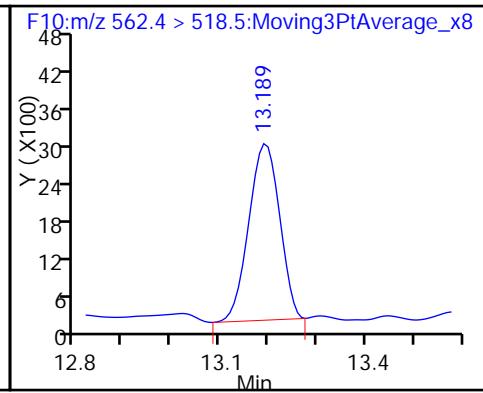
## 25 Perfluorodecane Sulfonate



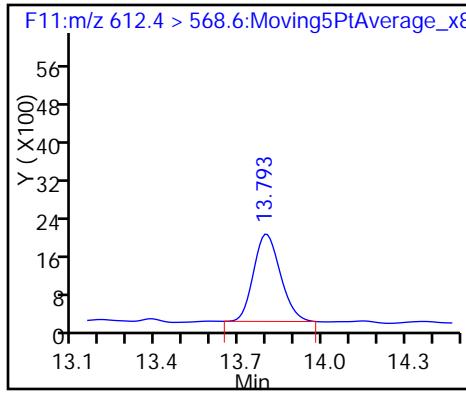
## D 26 13C2 PFUna



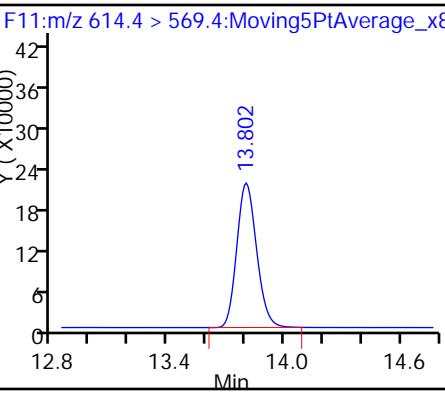
## 27 Perfluoroundecanoic acid



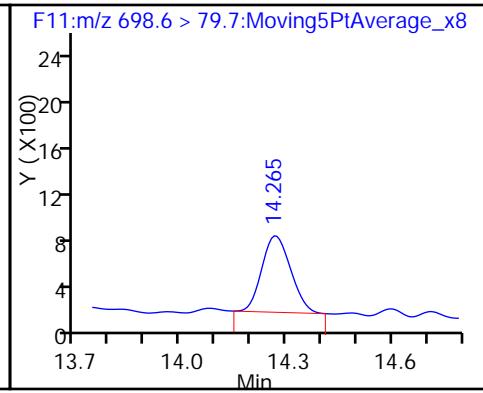
## 29 Perfluorododecanoic acid



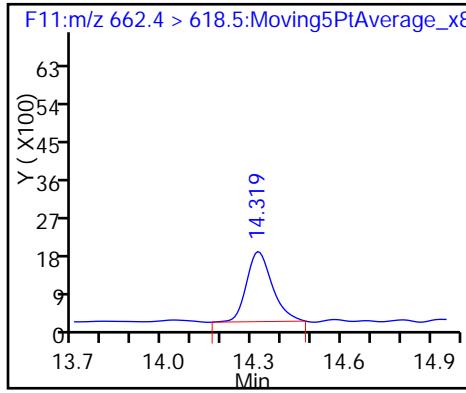
## D 28 13C2 PFDoA



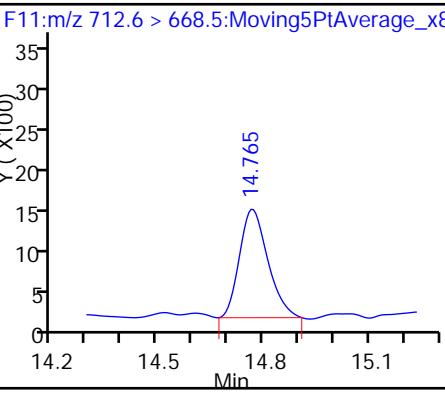
## 31 PFDoS (Perflouro-1-dodecanesulfona



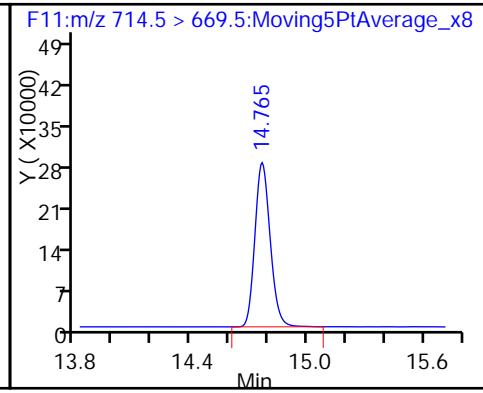
## 30 Perfluorotridecanoic acid



## 32 Perfluorotetradecanoic acid



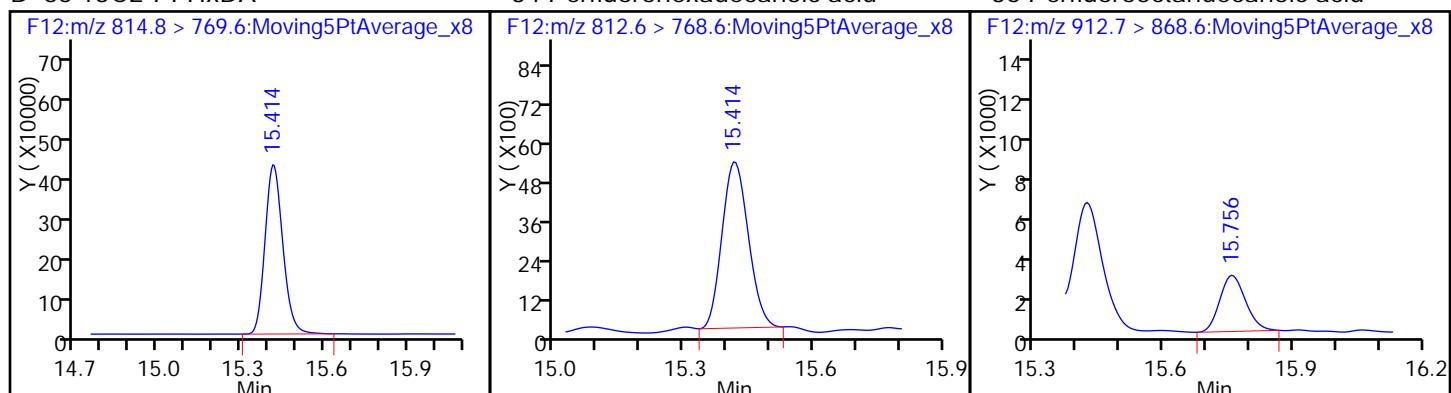
## D 33 13C2-PFTeDA



D 35 13C2-PFHxD

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_041.d  
 Lims ID: Std L2  
 Client ID:  
 Sample Type: IC Calib Level: 2  
 Inject. Date: 04-Dec-2014 04:05:29 ALS Bottle#: 5 Worklist Smp#: 5  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L2 PFC 1.00/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:52 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 113C4 PFBA										
216.7 > 171.5	5.937	5.934	0.003		116258	27.2		136	401	
2 Perfluorobutyric acid										
212.7 > 168.6	5.937	5.935	0.002	1.000	22829	1.01		101	80.4	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.035	7.037	-0.002	1.000	14944	0.9431		94.3	25.5	
D 3 113C5-PFPeA										
267.6 > 222.7	7.035	7.038	-0.003		191617	24.8		124	619	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.157	7.157	0.0	1.000	41334	0.8023		90.8	163	
298.8 > 98.6	7.161	7.157	0.004	1.001	23851	1.73(0.00-0.00)		90.8	77.3	
D 6 113C2 PFHxA										
314.6 > 269.7	8.284	8.282	0.002		455689	23.9		120	1334	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.284	8.284	0.0	1.000	21004	0.9444		94.4	87.1	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.366	8.360	0.006	0.876	37735	0.9191		98.0	147	
D 8 113C4-PFHxA										
366.6 > 321.6	9.510	9.506	0.004		427809	24.1		121	930	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.517	9.508	0.009	1.000	28535	0.8749		87.5	83.2	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	49260	1.02		108	173	
D 11 118O2 PFHxA										
402.5 > 83.6	9.545	9.544	0.001		988251	22.0		116	3013	
D 12 113C4 PFOA										
416.5 > 371.6	10.632	10.628	0.004		2070020	61.2		122	3758	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.632	10.628	0.004	1.000	74889	1.08		108	111	
412.8 > 168.7	10.632	10.628	0.004	1.000	23992		3.12(0.00-0.00)	108	29.9	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.632	10.635	-0.003	1.000	43172	0.9181		96.4	74.5	
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.595	0.007		2599599	57.6		121	4817	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.602	11.595	0.007	1.000	119216	1.21		127	146	
498.9 > 98.7	11.602	11.595	0.007	1.000	73803		1.62(0.00-0.00)	127	131	
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.614	0.006		908195	23.5		118	2426	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	60944	1.16		116	96.0	
21 PFNS (Perflouro-1-nonenesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	38953	0.9746		102	90.8	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	58072	0.99		99.0	88.3	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		1228601	23.9		119	1888	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.003	12.998	0.005	1.000	157092	1.06		106	308	
D 23 13C8 FOSA										
505.4 > 77.6	13.003	12.998	0.005		7782345	54.4		109	4481	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	35820	1.02		106	83.6	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1330638	20.5		103	3038	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.206	13.198	0.008	1.000	75233	1.28		128	150	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	61887	0.9479		94.8	41.7	
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1323923	20.3		101	1630	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	31494	0.8795		90.9	68.8	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.326	14.327	-0.001	1.000	65903	1.22		122	56.4	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	36572	1.06		106	47.2	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1525719	21.3		107	2687	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.414	15.415	-0.001		2116849	20.3		102	2541	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.414	15.415	-0.001	1.000	98406	1.11		111	135	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.751	15.752	-0.001	1.000	Page 94742 of 350	1.01		101	110	12/12/2014

Report Date: 05-Dec-2014 10:01:53

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L2\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:01:53

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_041.d

Injection Date: 04-Dec-2014 04:05:29

Instrument ID: A4

Lims ID: Std L2

Client ID:

Operator ID: JRB

ALS Bottle#: 5 Worklist Smp#: 5

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

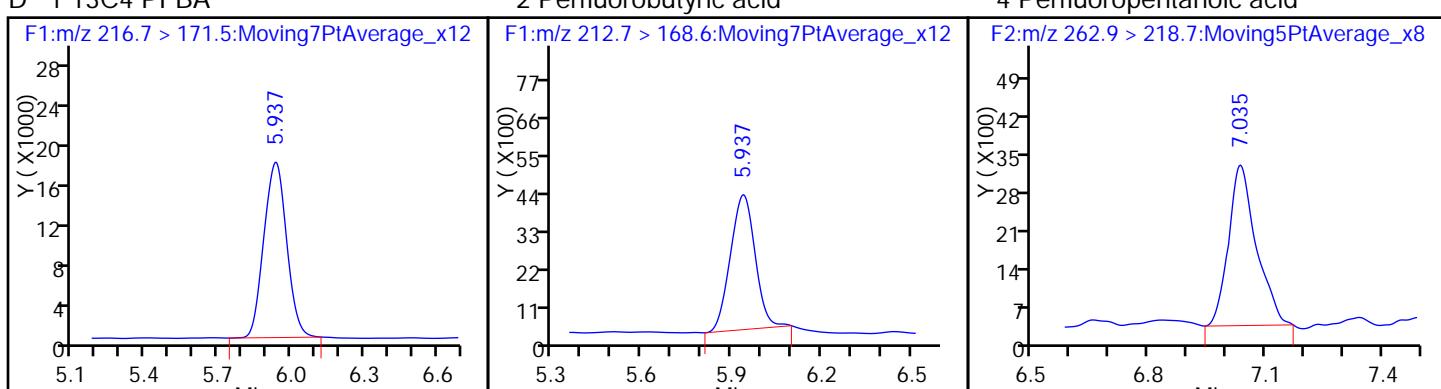
Method: PFAC\_A4

Limit Group: LC PFC ICAL

## D 1 13C4 PFBA

## 2 Perfluorobutyric acid

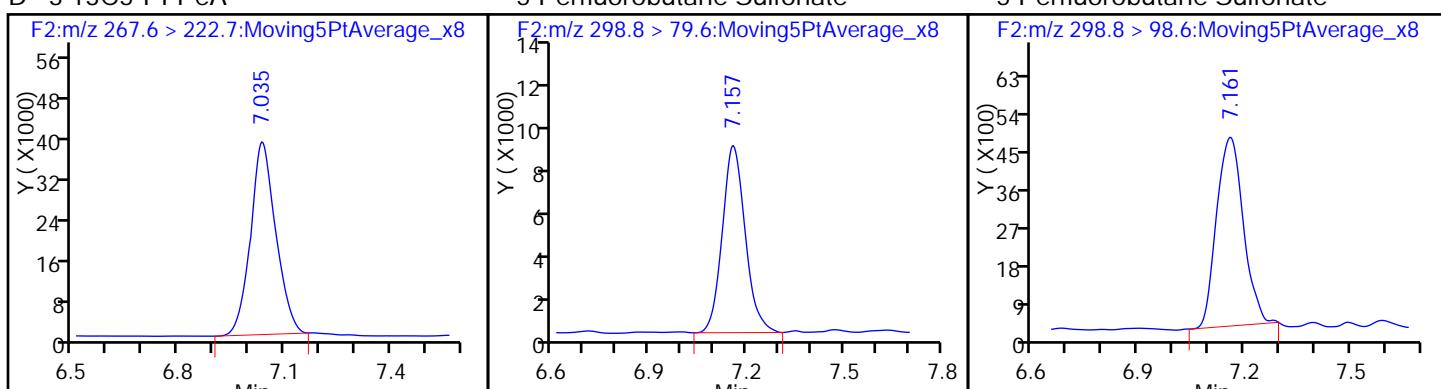
## 4 Perfluoropentanoic acid



## D 3 13C5-PFPeA

## 5 Perfluorobutane Sulfonate

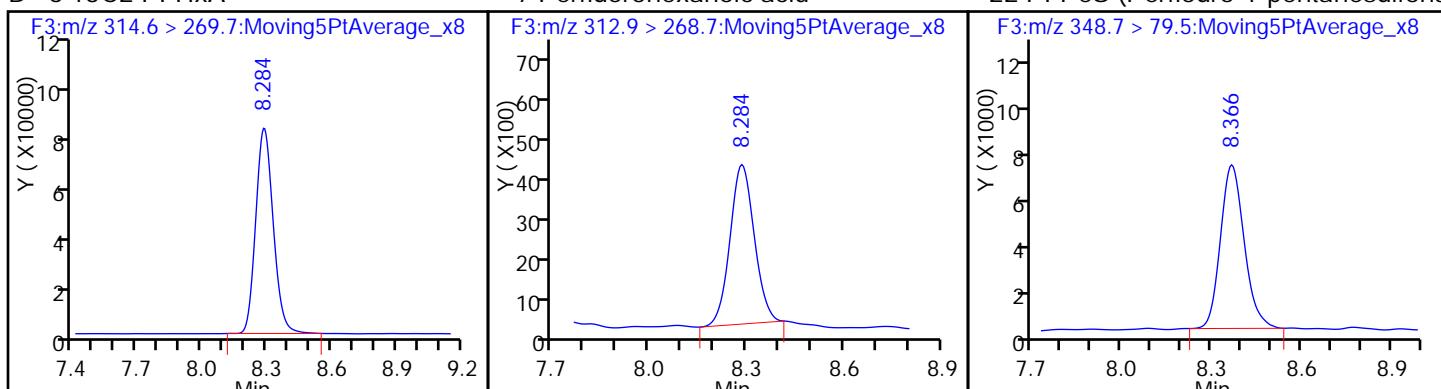
## 5 Perfluorobutane Sulfonate



## D 6 13C2 PFHxA

## 7 Perfluorohexanoic acid

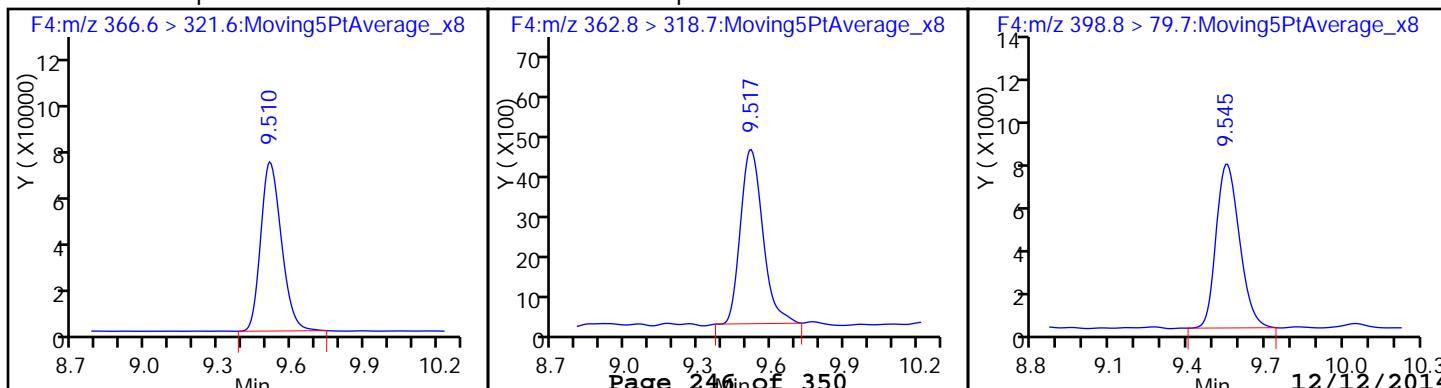
## 22 PFPeS (Perflouro-1-pentanesulfonat)



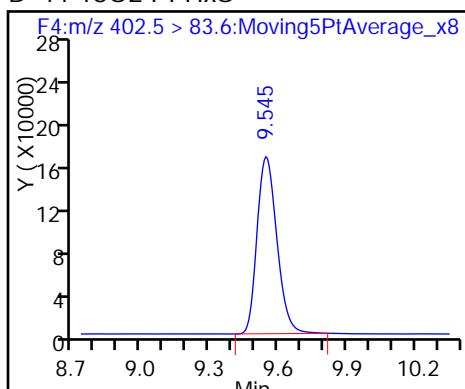
## D 8 13C4-PFHxA

## 9 Perfluoroheptanoic acid

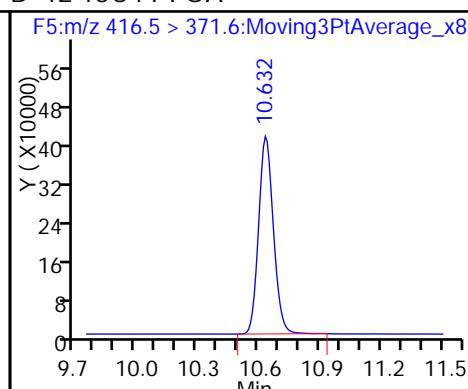
## 10 Perfluorohexane Sulfonate



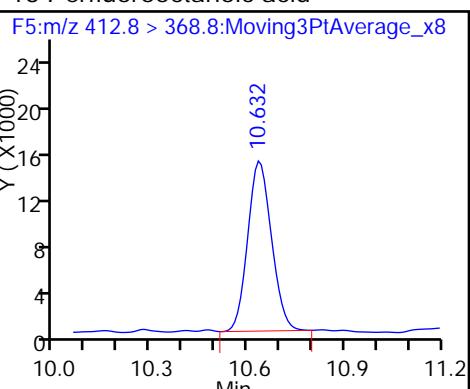
D 11 18O2 PFHxS



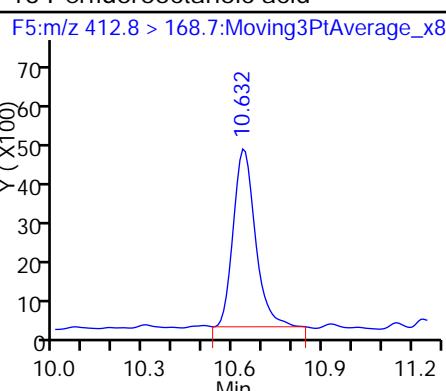
D 12 13C4 PFOA



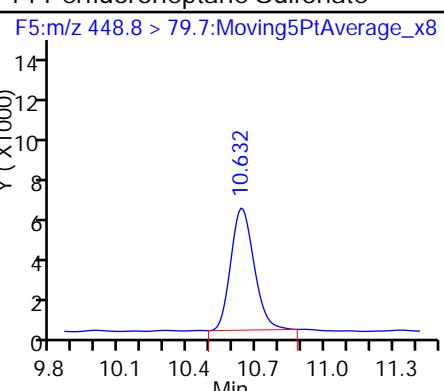
13 Perfluorooctanoic acid



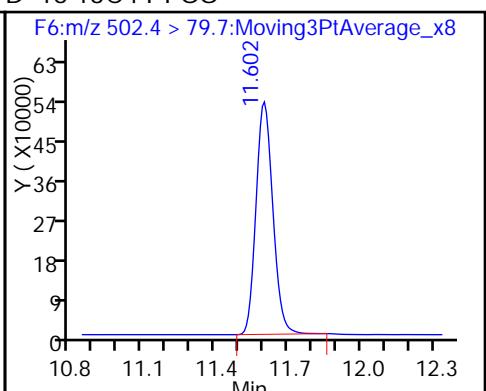
13 Perfluorooctanoic acid



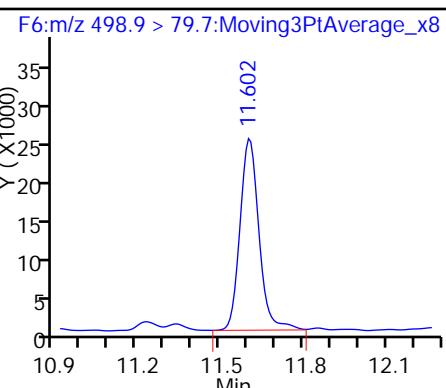
14 Perfluoroheptane Sulfonate



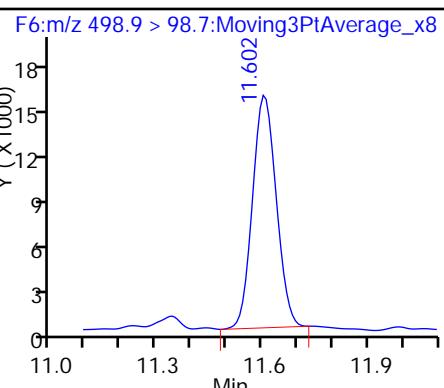
D 16 13C4 PFOS



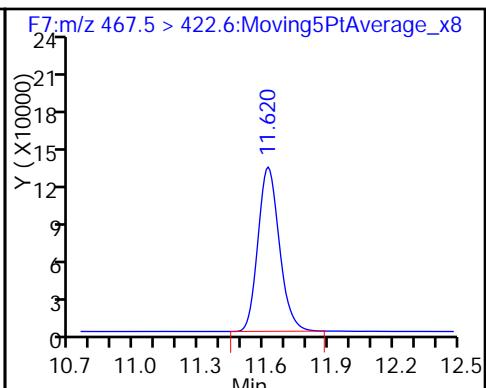
15 Perfluorooctanoic Sulfonate



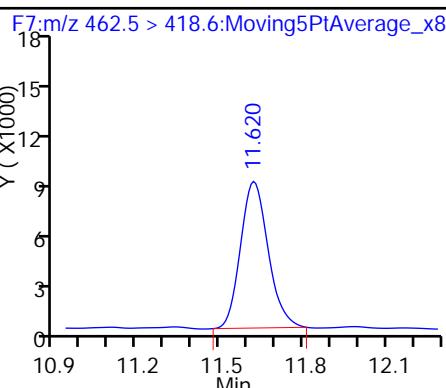
15 Perfluorooctanoic Sulfonate



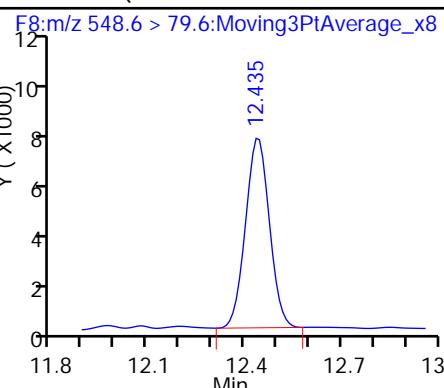
D 17 13C5 PFNA



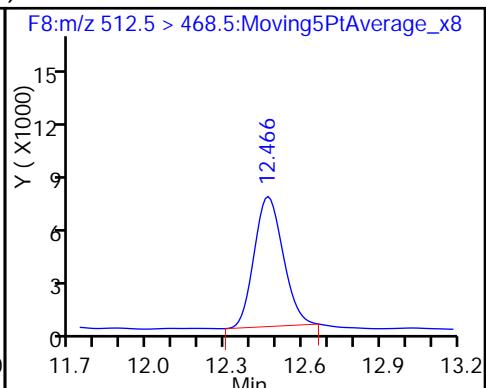
18 Perfluorononanoic acid



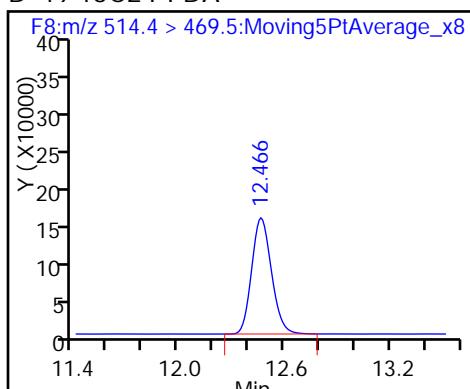
21 PFNS (Perflouro-1-nananesulfonate)



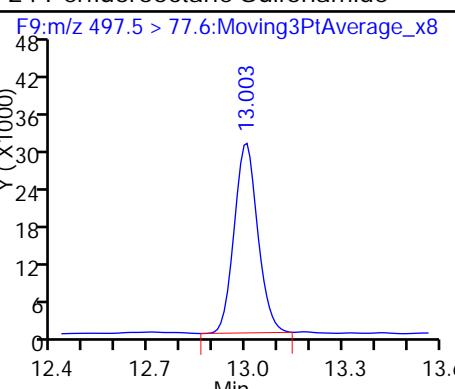
20 Perfluorodecanoic acid



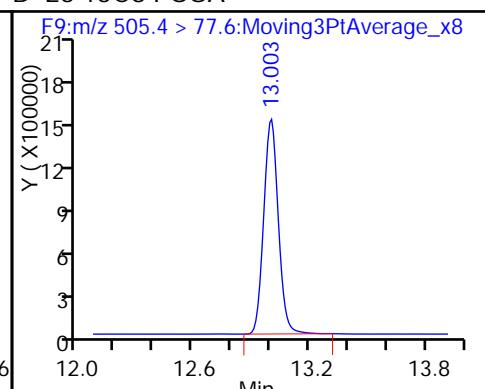
D 19 13C2 PFDA



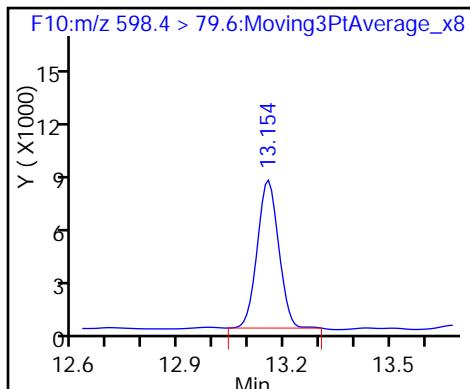
24 Perfluorooctane Sulfonamide



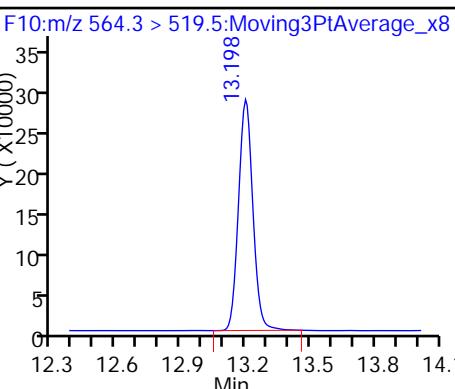
D 23 13C8 FOSA



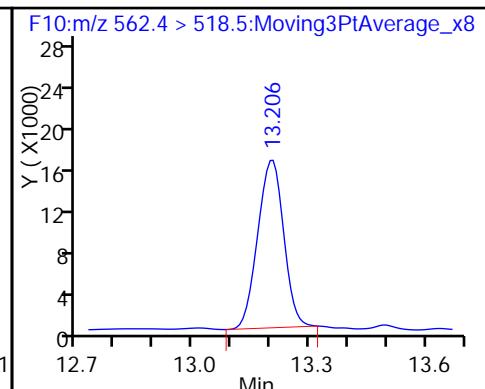
25 Perfluorodecane Sulfonate



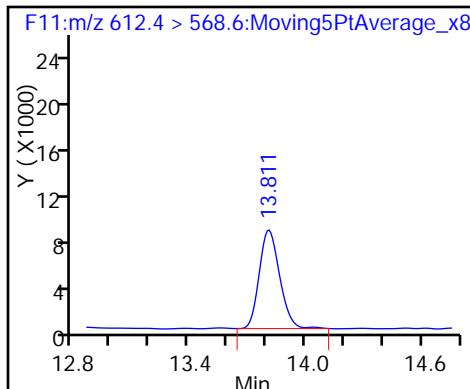
D 26 13C2 PFUna



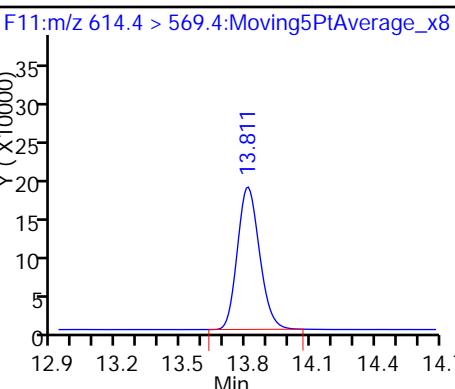
27 Perfluoroundecanoic acid



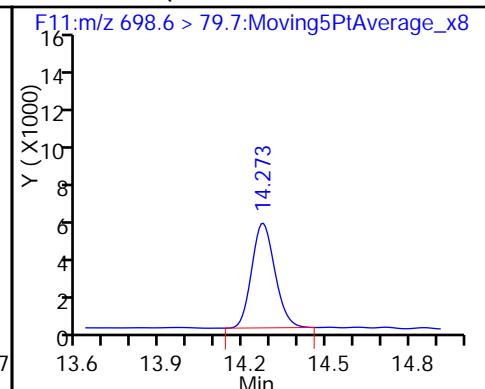
29 Perfluorododecanoic acid



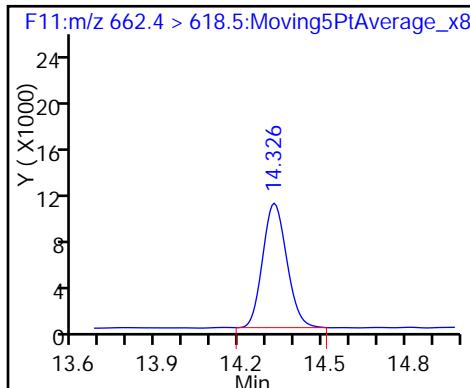
D 28 13C2 PFDoA



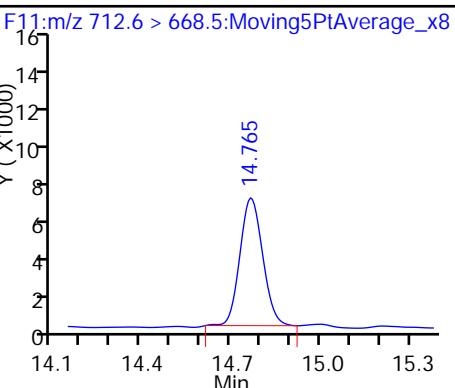
31 PFDoS (Perflouro-1-dodecanesulfona



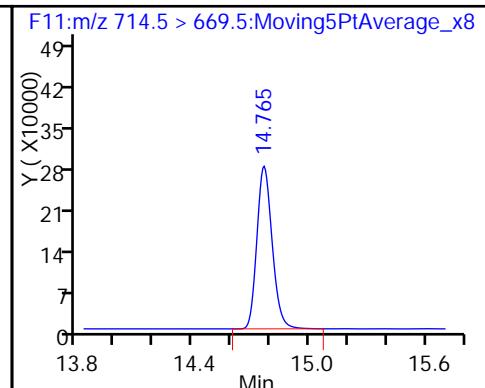
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



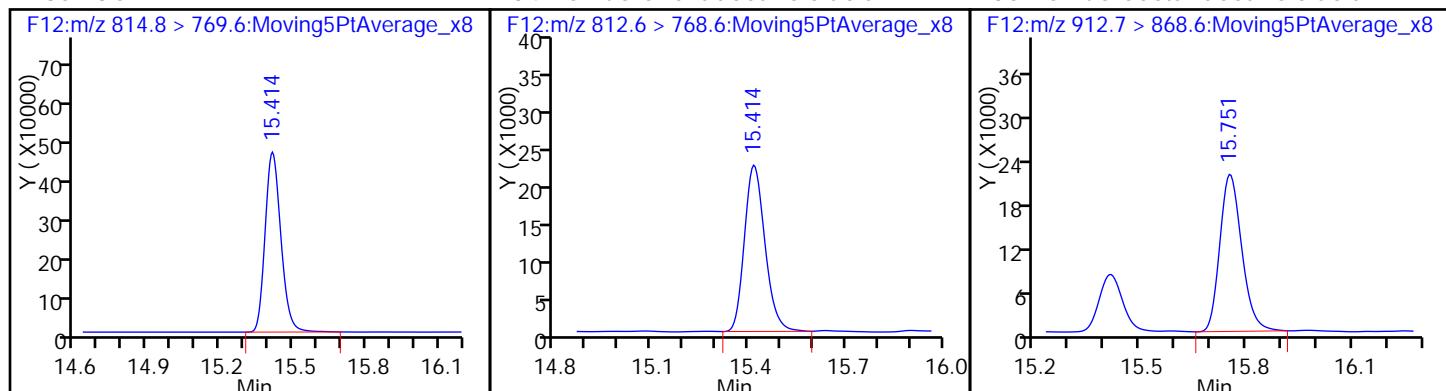
D 33 13C2-PFTeDA



D 35 13C2-PFHxD

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_042.d  
 Lims ID: Std L3  
 Client ID:  
 Sample Type: IC Calib Level: 3  
 Inject. Date: 04-Dec-2014 04:26:41 ALS Bottle#: 6 Worklist Smp#: 6  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L3 PFC 5.0/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:53 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
<b>D 1 113C4 PFBA</b>										
216.7 > 171.5	5.928	5.934	-0.006		105097	24.6		123	350	
2 Perfluorobutyric acid										
212.7 > 168.6	5.928	5.935	-0.007	1.000	109952	5.72		114	320	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.029	7.037	-0.008	1.000	70990	5.25		105	61.6	
<b>D 3 113C5-PFPeA</b>										
267.6 > 222.7	7.029	7.038	-0.009		163549	21.2		106	563	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.144	7.157	-0.013	1.000	216823	4.31		97.6	758	
298.8 > 98.6	7.147	7.157	-0.010	1.000	138038		1.57(0.00-0.00)	97.6	610	
<b>D 6 113C2 PFHxA</b>										
314.6 > 269.7	8.273	8.282	-0.009		496708	26.1		130	1850	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.273	8.284	-0.011	1.000	134092	5.53		111	318	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.349	8.360	-0.011	0.875	206763	5.16		110	739	
<b>D 8 113C4-PFHxA</b>										
366.6 > 321.6	9.499	9.506	-0.007		434956	24.5		123	1400	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.499	9.508	-0.009	1.000	179416	5.41		108	509	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.538	9.543	-0.005	1.000	252665	5.38		114	1071	
<b>D 11 118O2 PFHxS</b>										
402.5 > 83.6	9.538	9.544	-0.006		964272	21.5		113	2820	
<b>D 12 113C4 PFOA</b>										
416.5 > 371.6	10.623	10.628	-0.005		2023751	59.8		120	2824	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.623	10.628	-0.005	1.000	340656	5.04		101	470	
412.8 > 168.7	10.632	10.628	0.004	1.001	113441		3.00(0.00-0.00)	101	146	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.632	10.635	-0.003	1.000	243339	5.01		105	509	
D 16 13C4 PFOS										
502.4 > 79.7	11.594	11.595	-0.001		2684739	59.5		125	5305	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.594	11.595	-0.001	1.000	439807	4.34		90.8	593	
498.9 > 98.7	11.594	11.595	-0.001	1.000	243413		1.81(0.00-0.00)	90.8	455	
D 17 13C5 PFNA										
467.5 > 422.6	11.611	11.614	-0.003		977581	25.3		127	2917	
18 Perfluorononanoic acid										
462.5 > 418.6	11.611	11.616	-0.005	1.000	306271	5.43		109	581	
21 PFNS (Perflouro-1-nonanesulfonate)										
548.6 > 79.6	12.425	12.429	-0.004	1.000	225687	5.47		114	394	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.455	12.461	-0.006	1.000	335348	5.93		119	613	
D 19 13C2 PFDA										
514.4 > 469.5	12.455	12.463	-0.008		1185934	23.0		115	2273	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	12.994	12.998	-0.004	1.000	777293	5.15		103	922	
D 23 13C8 FOSA										
505.4 > 77.6	12.994	12.998	-0.004		7943692	55.5		111	5034	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.146	13.152	-0.006	1.000	205469	5.66		117	431	
D 26 13C2 PFUnA										
564.3 > 519.5	13.189	13.198	-0.009		1337877	20.7		103	2258	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.189	13.198	-0.009	1.000	361261	6.11		122	720	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.802	13.808	-0.006	1.000	461093	6.10		122	358	
D 28 13C2 PFDoA										
614.4 > 569.4	13.802	13.809	-0.007		1533105	23.5		117	2111	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	211588	5.72		118	424	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.326	14.327	-0.001	1.000	337727	5.39		108	250	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	215805	5.41		108	218	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1604800	22.4		112	2550	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.415	15.415	0.0		2247748	21.6		108	2844	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.415	15.415	0.0	1.000	504906	4.92		98.3	523	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.751	15.752	-0.001	1.000	Page 497 of 457	4.57		91.4	462	12/12/2014

Report Date: 05-Dec-2014 10:01:54

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L3\_00005

Amount Added: 1.00

Units: mL

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_042.d

Injection Date: 04-Dec-2014 04:26:41

Instrument ID: A4

Lims ID: Std L3

Client ID:

Operator ID: JRB

ALS Bottle#: 6 Worklist Smp#: 6

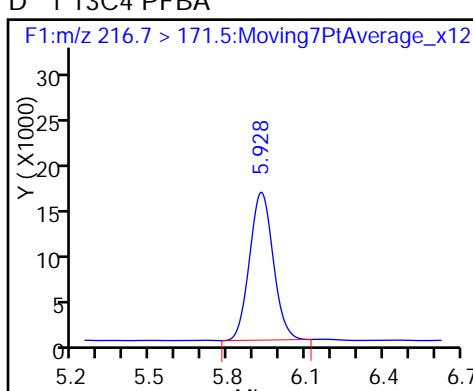
Injection Vol: 15.0 ul

Dil. Factor: 1.0000

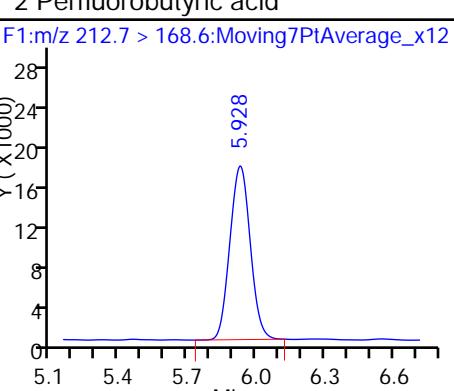
Method: PFAC\_A4

Limit Group: LC PFC ICAL

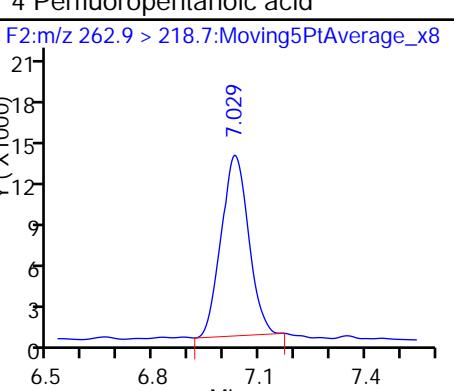
## D 1 13C4 PFBA



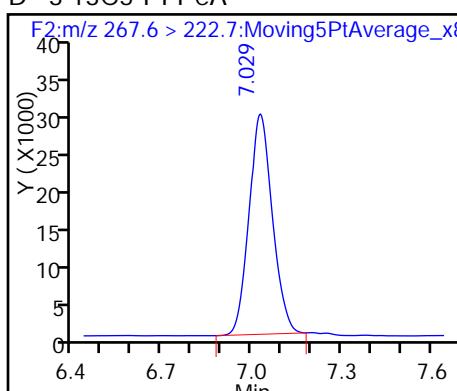
## 2 Perfluorobutyric acid



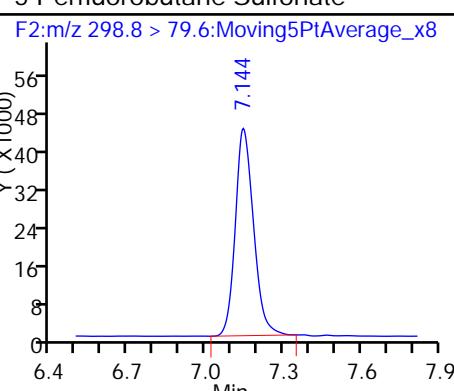
## 4 Perfluoropentanoic acid



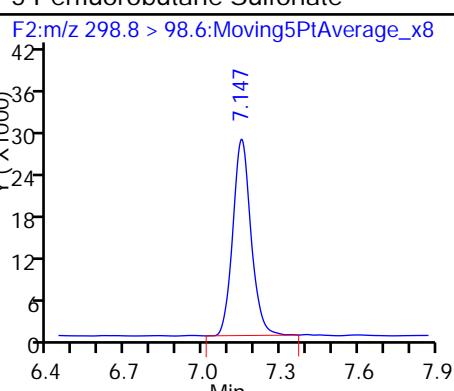
## D 3 13C5-PFPeA



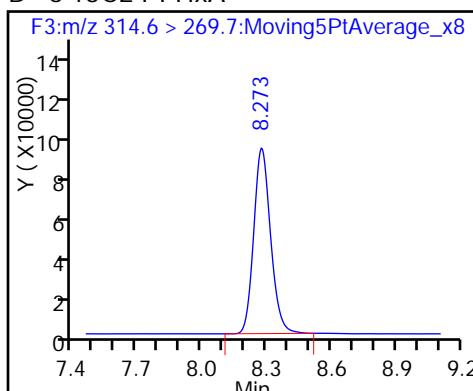
## 5 Perfluorobutane Sulfonate



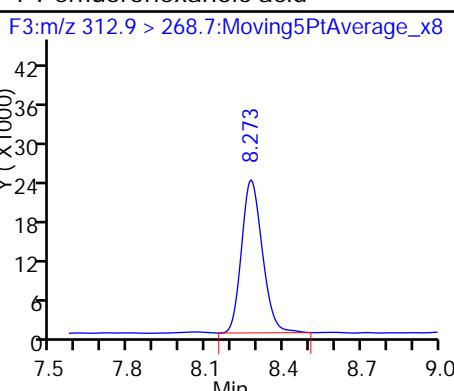
## 5 Perfluorobutane Sulfonate



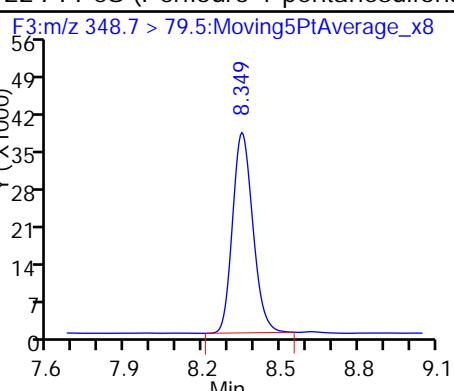
## D 6 13C2 PFHxA



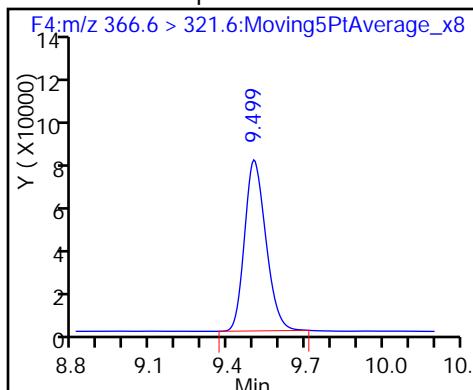
## 7 Perfluorohexanoic acid



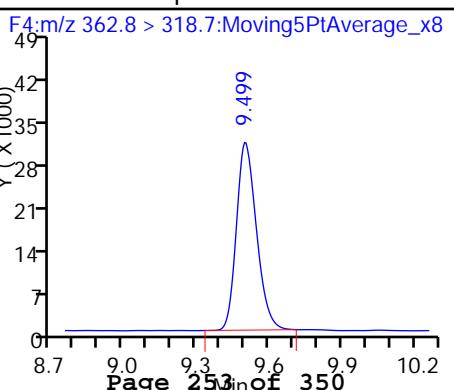
## 22 PFPeS (Perflouro-1-pentanesulfonat)



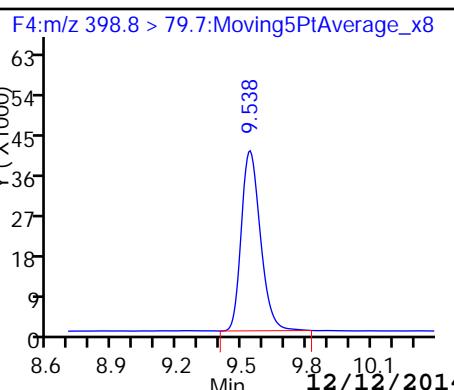
## D 8 13C4-PFHxA



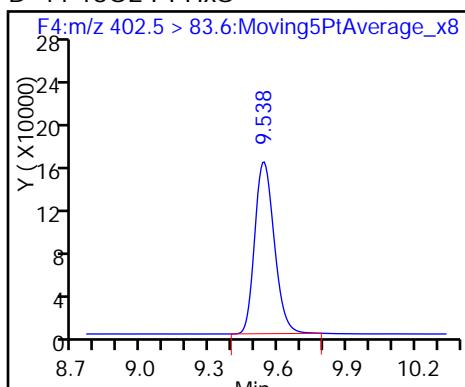
## 9 Perfluoroheptanoic acid



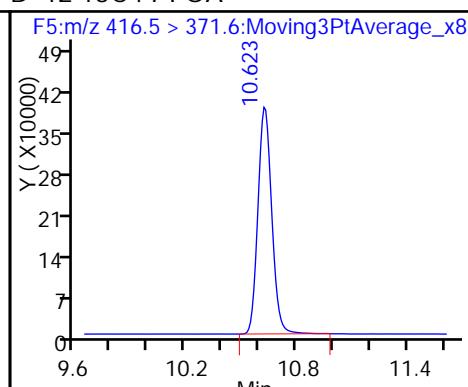
## 10 Perfluorohexane Sulfonate



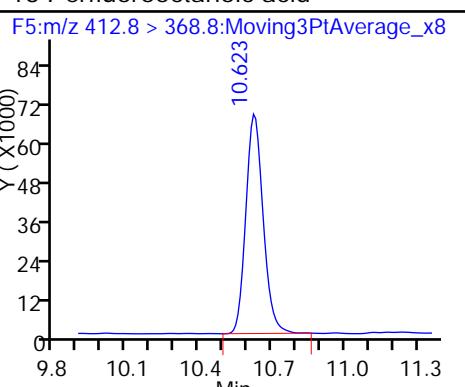
D 11 18O2 PFHxS



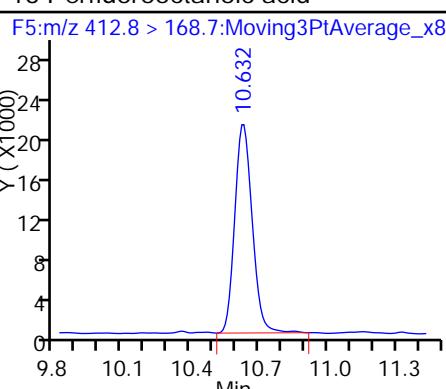
D 12 13C4 PFOA



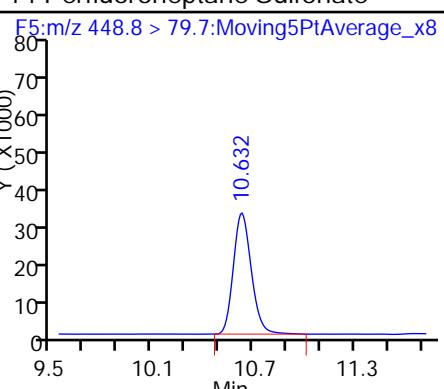
13 Perfluorooctanoic acid



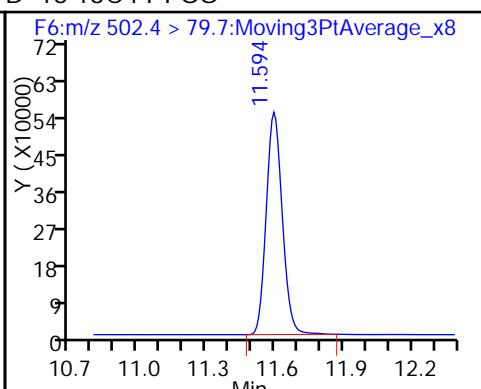
13 Perfluorooctanoic acid



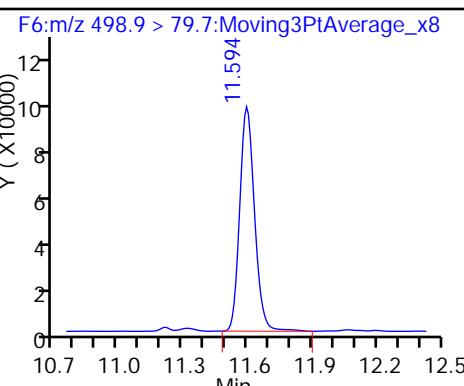
14 Perfluoroheptane Sulfonate



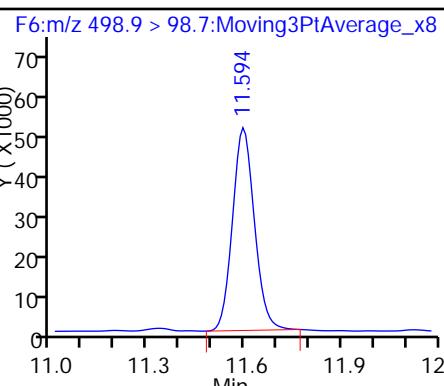
D 16 13C4 PFOS



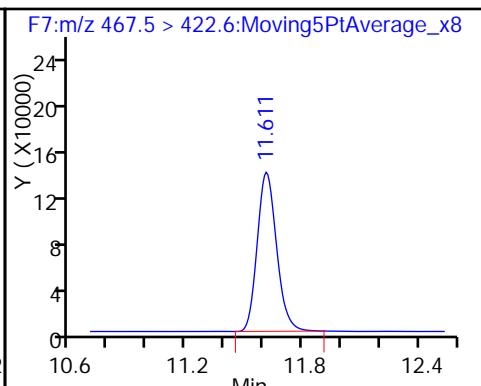
15 Perfluorooctanoic Sulfonate



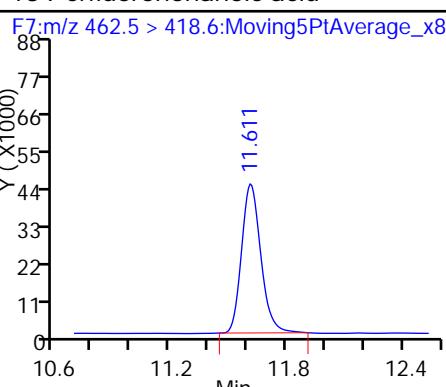
15 Perfluorooctanoic Sulfonate



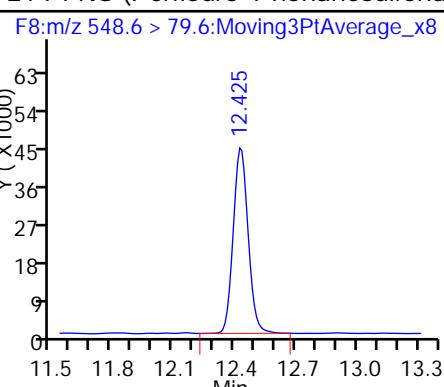
D 17 13C5 PFNA



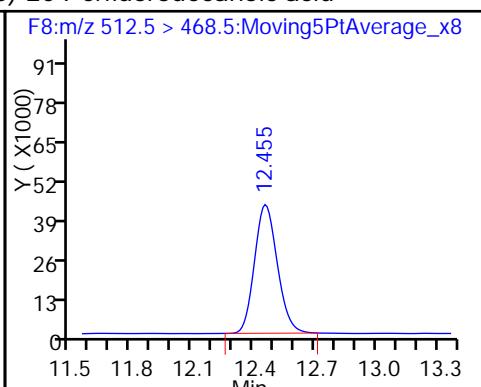
18 Perfluorononanoic acid



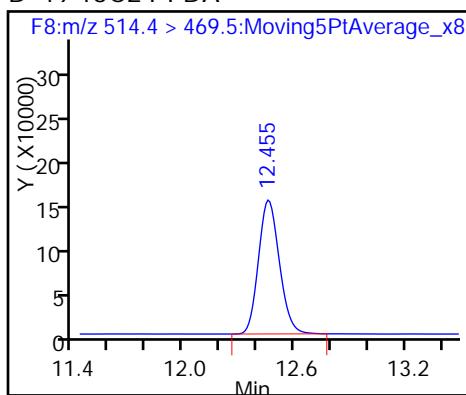
21 PFNS (Perflouro-1-nananesulfonate)



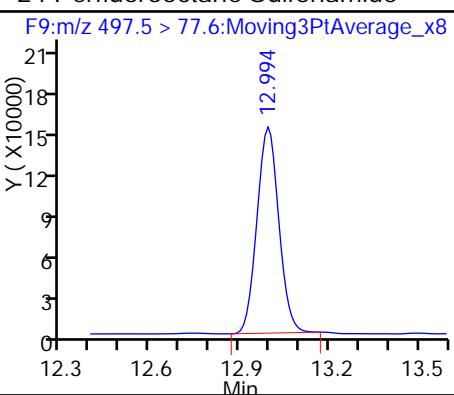
20 Perfluorodecanoic acid



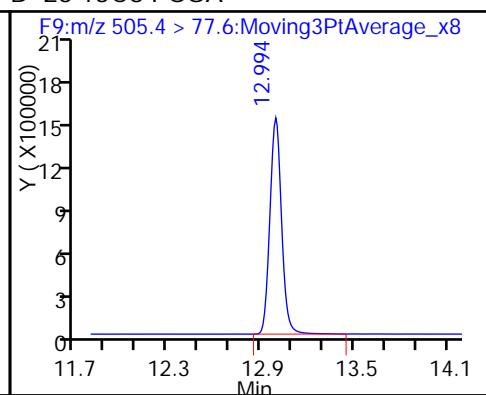
## D 19 13C2 PFDA



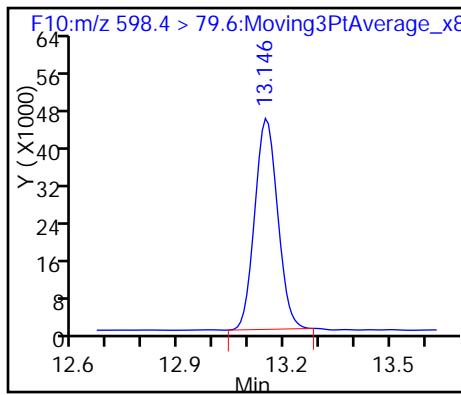
## 24 Perfluorooctane Sulfonamide



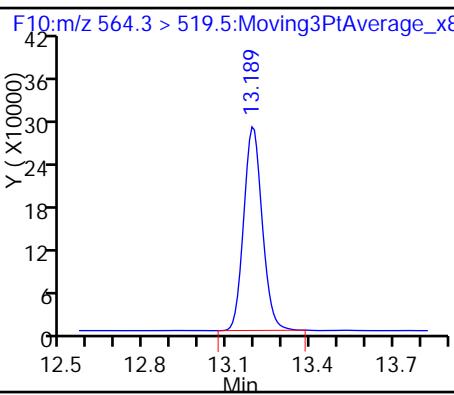
## D 23 13C8 FOSA



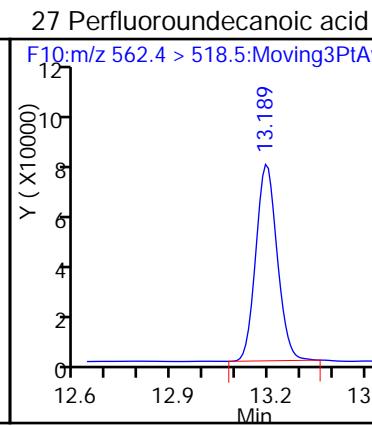
## 25 Perfluorodecane Sulfonate



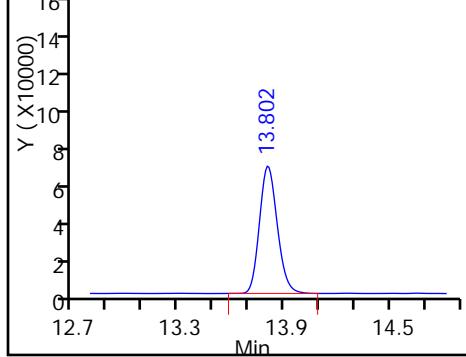
## D 26 13C2 PFUna



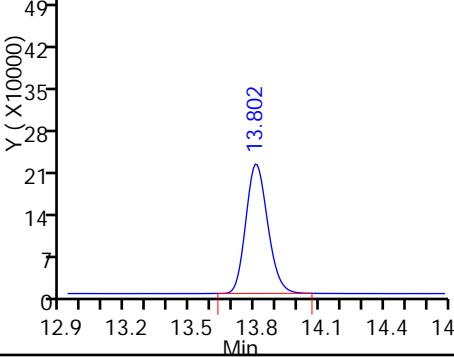
## 27 Perfluoroundecanoic acid



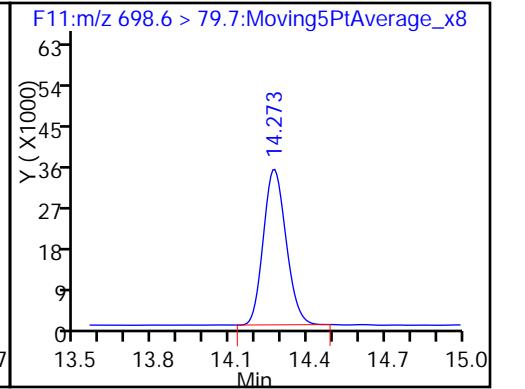
## 29 Perfluorododecanoic acid



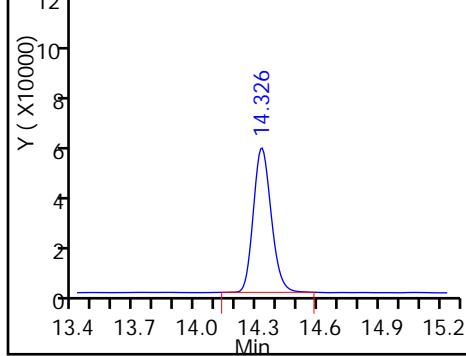
## D 28 13C2 PFDoA



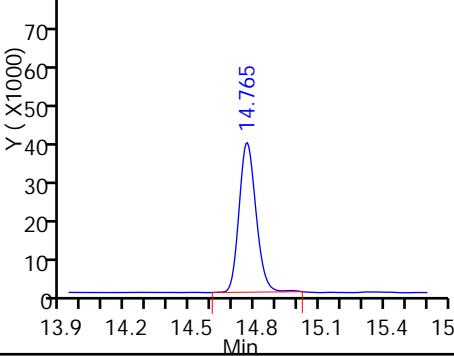
## 31 PFDoS (Perflouro-1-dodecanesulfona



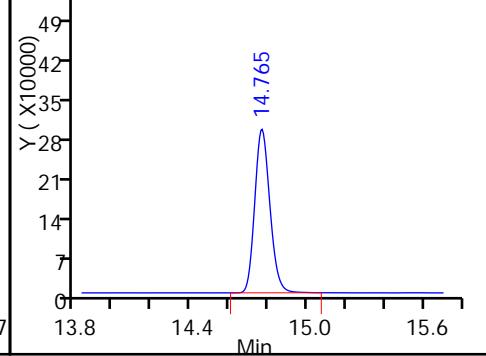
## 30 Perfluorotridecanoic acid



## 32 Perfluorotetradecanoic acid



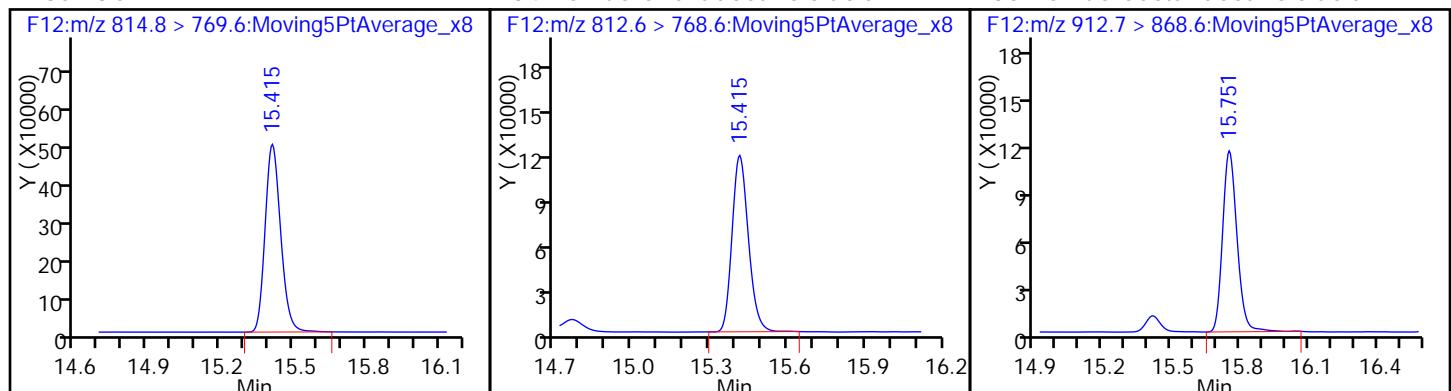
## D 33 13C2-PFTeDA



D 35 13C2-PFHxD

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_043.d  
 Lims ID: Std L4  
 Client ID:  
 Sample Type: IC Calib Level: 4  
 Inject. Date: 04-Dec-2014 04:47:55 ALS Bottle#: 7 Worklist Smp#: 7  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L4 PFC 20/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:54 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

First Level Reviewer: barnettj Date: 04-Dec-2014 09:44:55

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 113C4 PFBA										
216.7 > 171.5	5.934	5.934	0.0		86764	20.3		101	423	
2 Perfluorobutyric acid										
212.7 > 168.6	5.934	5.935	-0.001	1.000	390544	18.7		93.3	1329	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.039	7.037	0.002	1.000	296633	20.1		101	416	
D 3 113C5-PFPeA										
267.6 > 222.7	7.043	7.038	0.005		178067	23.1		115	848	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.161	7.157	0.004	1.000	853055	16.9		95.5	1804	
298.8 > 98.6	7.161	7.157	0.004	1.000	535360	1.59(0.00-0.00)		95.5	1767	
D 6 113C2 PFHxA										
314.6 > 269.7	8.290	8.282	0.008		368981	19.4		96.8	1065	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.290	8.284	0.006	1.000	399091	22.2		111	811	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.360	8.360	0.0	0.876	795956	19.8		105	3427	
D 8 113C4-PFHpA										
366.6 > 321.6	9.505	9.506	-0.001		388840	21.9		110	934	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.511	9.508	0.003	1.000	665119	22.4		112	1608	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	937096	19.8		105	2463	
D 11 18O2 PFHxS										
402.5 > 83.6	9.545	9.544	0.001		968949	21.6		114	3044	
D 12 113C4 PFOA										
416.5 > 371.6	10.632	10.628	0.004		1927321	57.0		114	3772	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.632	10.628	0.004	1.000	1249944	19.4		97.1	1444	
412.8 > 168.7	10.632	10.628	0.004	1.000	400743		3.12(0.00-0.00)	97.1	637	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.632	10.635	-0.003	1.000	950094	21.9		115	1551	
D 16 13C4 PFOS										
502.4 > 79.7	11.594	11.595	-0.001		2394219	53.1		111	3990	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.594	11.595	-0.001	1.000	1408150	15.6		81.5	1300	
498.9 > 98.7	11.594	11.595	-0.001	1.000	854068		1.65(0.00-0.00)	81.5	1324	
D 17 13C5 PFNA										
467.5 > 422.6	11.612	11.614	-0.002		836616	21.7		108	2209	
18 Perfluorononanoic acid										
462.5 > 418.6	11.612	11.616	-0.004	1.000	1048226	21.7		109	1672	
21 PFNS (Perflouro-1-nananesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	797477	21.7		113	1391	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	1269551	22.9		114	2193	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		1164228	22.6		113	1500	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	12.994	12.998	-0.004	1.000	3304921	22.3		112	2521	
D 23 13C8 FOSA										
505.4 > 77.6	12.994	12.998	-0.004		7791526	54.5		109	3157	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	710141	21.9		114	1659	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1295448	20.0		100	2289	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	1459686	25.5		127	1907	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	1576082	23.2		116	912	
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1377246	21.1		105	1763	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	786746	23.9		123	1478	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.327	14.327	0.0	1.000	1324401	23.5		118	1066	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	849938	23.7		119	889	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1536682	21.5		107	3721	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.415	15.415	0.0		2076818	19.9		99.6	2593	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.415	15.415	0.0	1.000	1737164	18.8		94.1	1317	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.751	15.752	-0.001	1.000	2099144	21.5		107	1834	

Report Date: 05-Dec-2014 10:01:55

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L4\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:01:55

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_043.d

Injection Date: 04-Dec-2014 04:47:55

Instrument ID: A4

Lims ID: Std L4

Client ID:

Operator ID: JRB

ALS Bottle#: 7 Worklist Smp#: 7

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

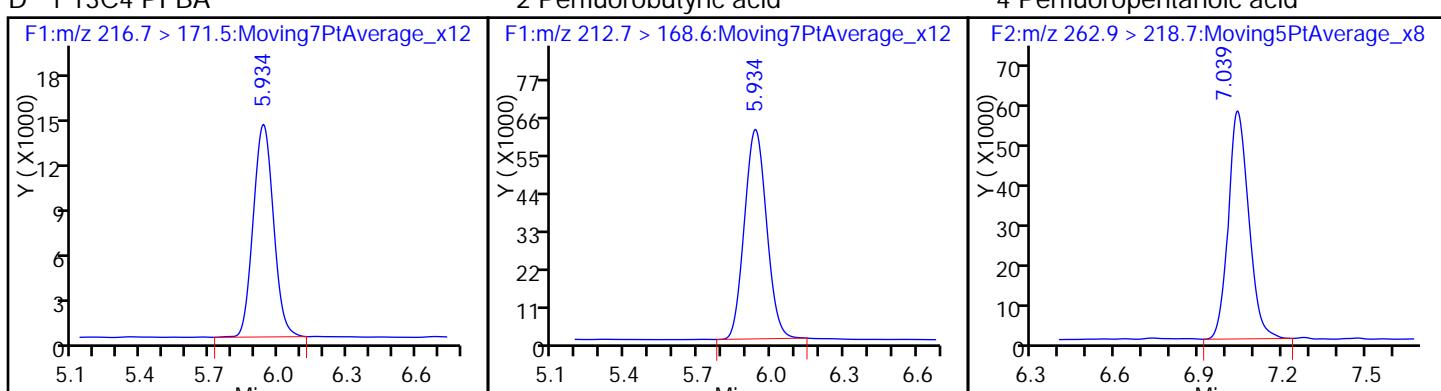
Method: PFAC\_A4

Limit Group: LC PFC ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

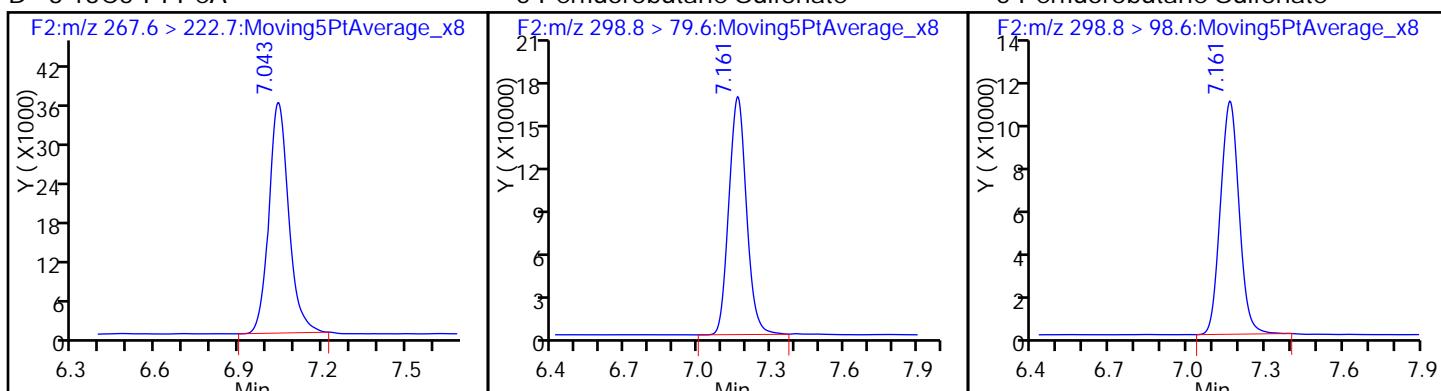
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

5 Perfluorobutane Sulfonate

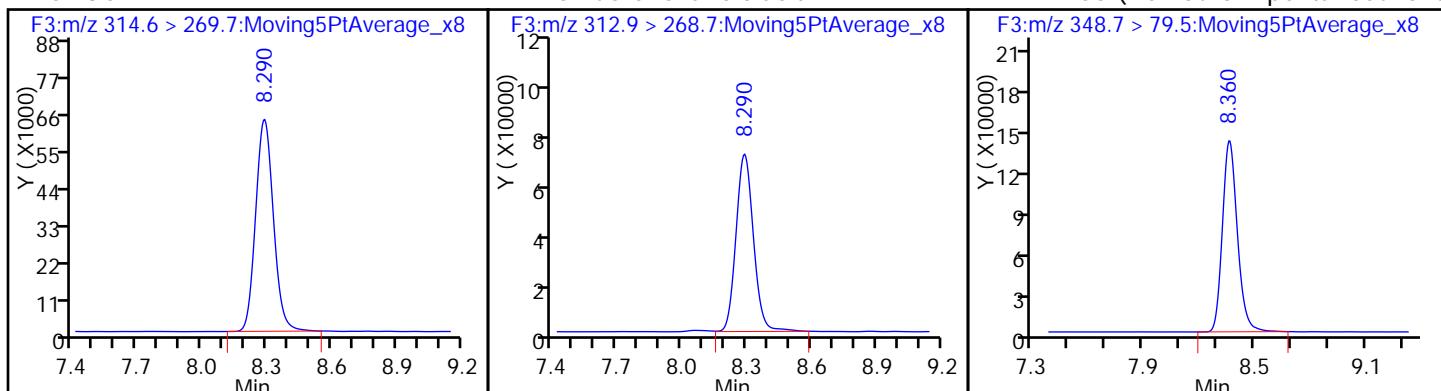
5 Perfluorobutane Sulfonate



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

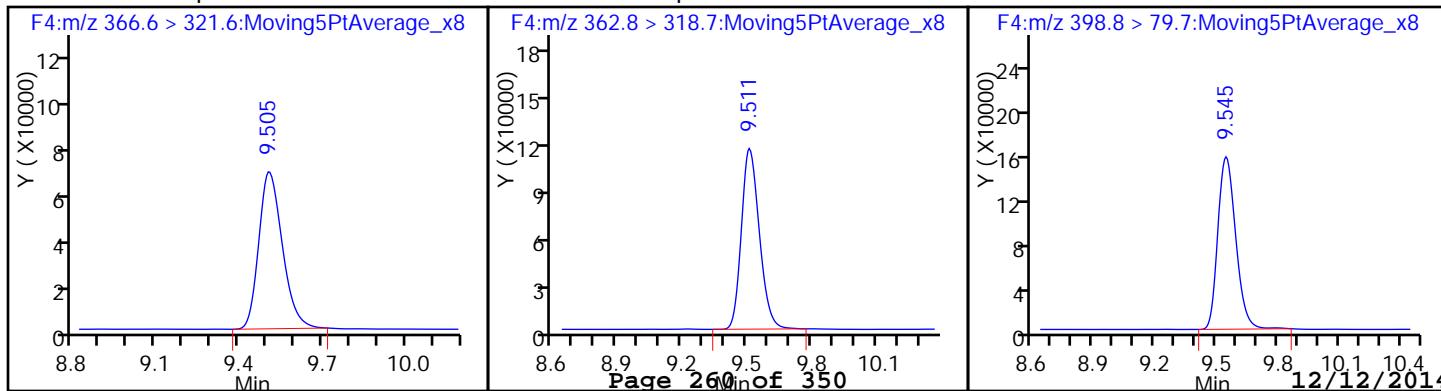
22 PFPeS (Perflouro-1-pentanesulfonat)



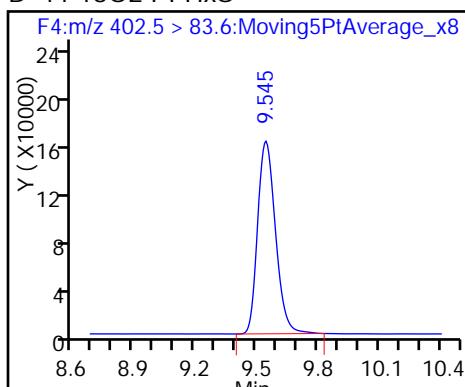
D 8 13C4-PFHxA

9 Perfluoroheptanoic acid

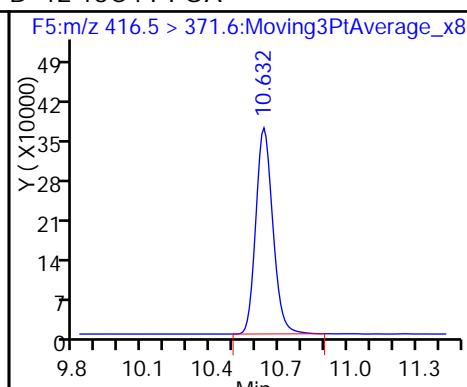
10 Perfluorohexane Sulfonate



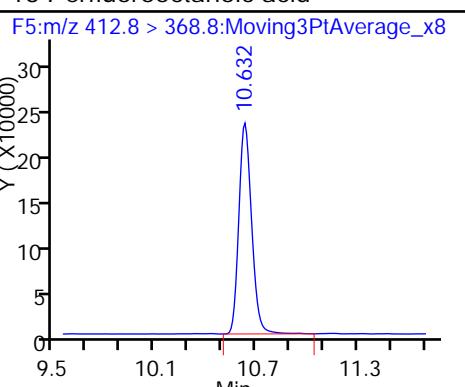
D 11 18O2 PFHxS



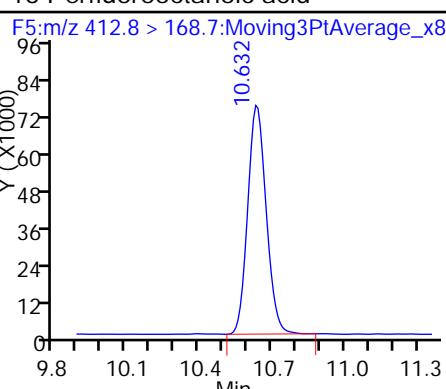
D 12 13C4 PFOA



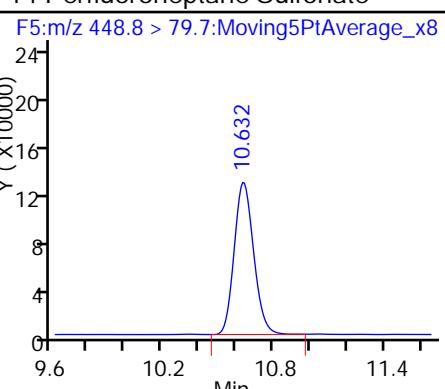
13 Perfluorooctanoic acid



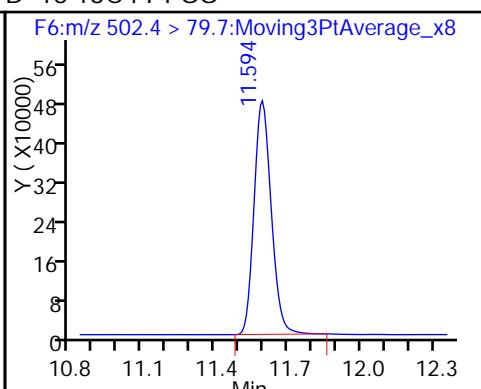
13 Perfluorooctanoic acid



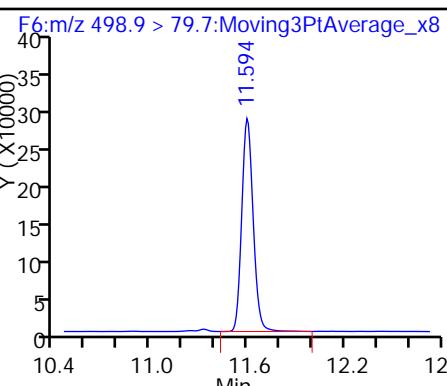
14 Perfluoroheptane Sulfonate



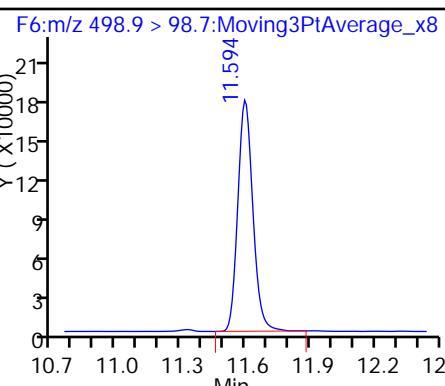
D 16 13C4 PFOS



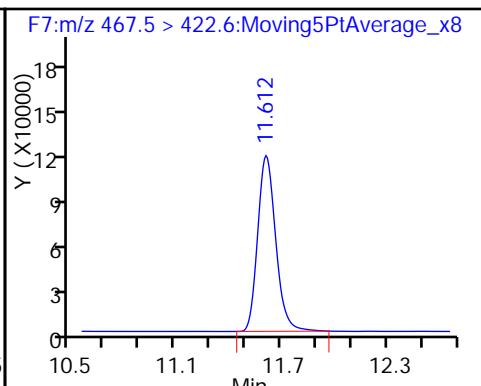
15 Perfluorooctanoic Sulfonate



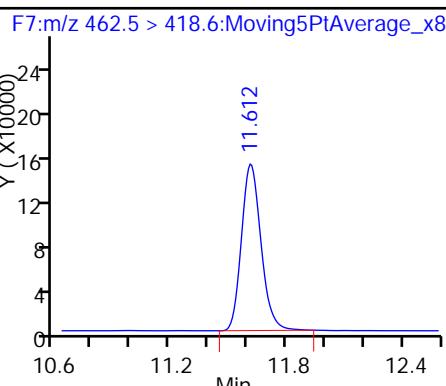
15 Perfluorooctanoic Sulfonate



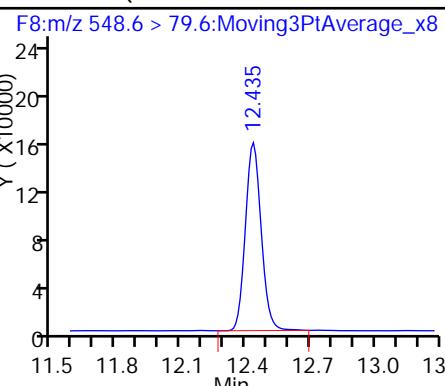
D 17 13C5 PFNA



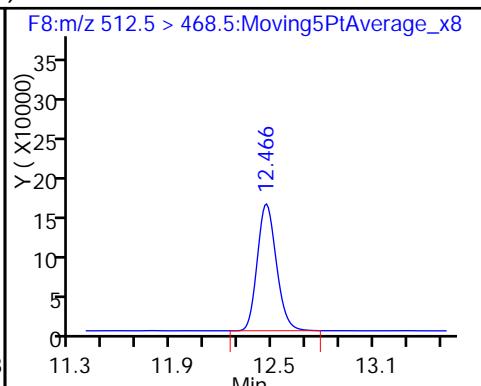
18 Perfluorononanoic acid



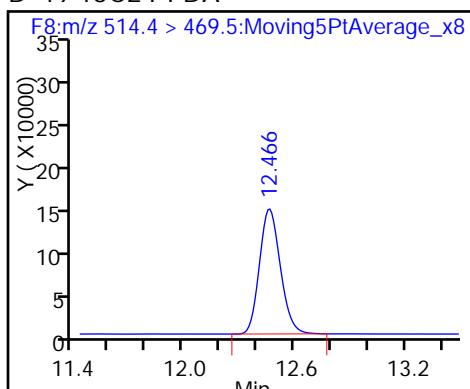
21 PFNS (Perflouro-1-nananesulfonate)



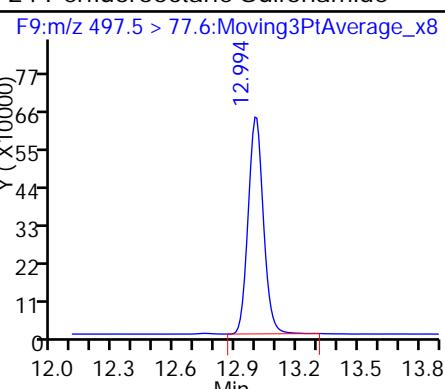
20 Perfluorodecanoic acid



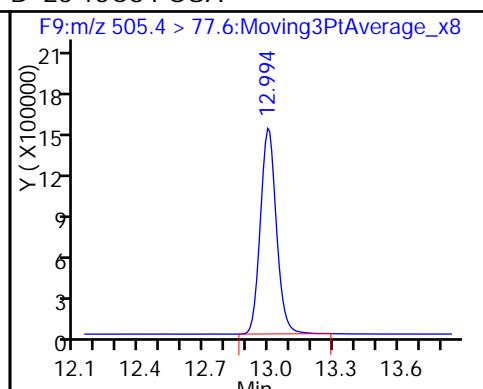
## D 19 13C2 PFDA



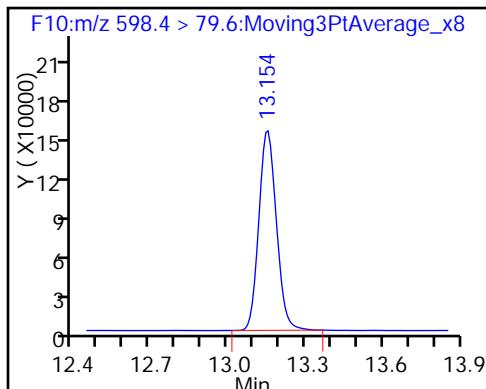
## 24 Perfluorooctane Sulfonamide



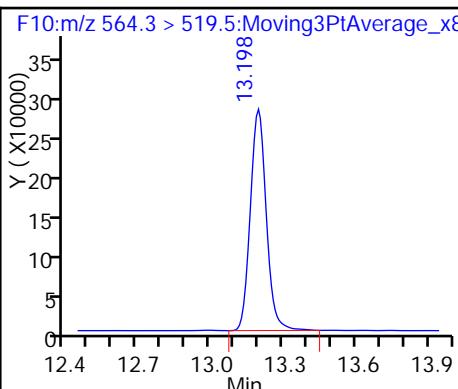
## D 23 13C8 FOSA



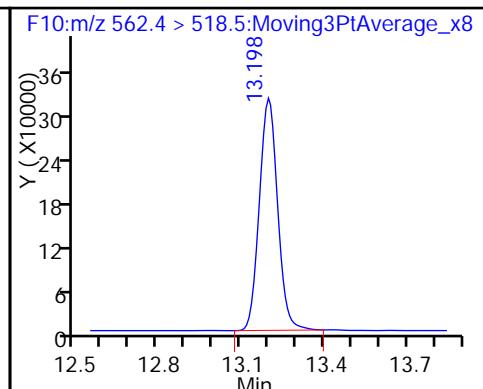
## 25 Perfluorodecane Sulfonate



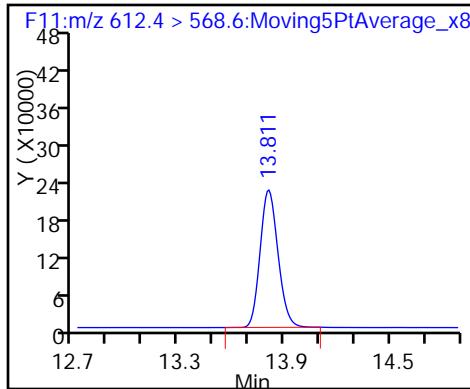
## D 26 13C2 PFUna



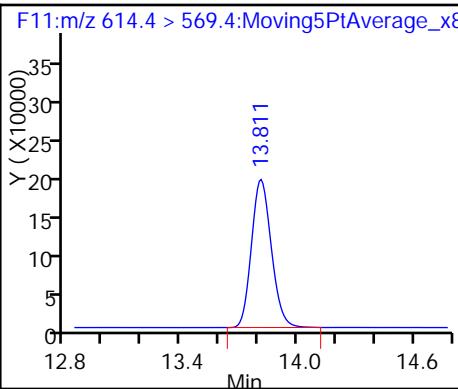
## 27 Perfluoroundecanoic acid



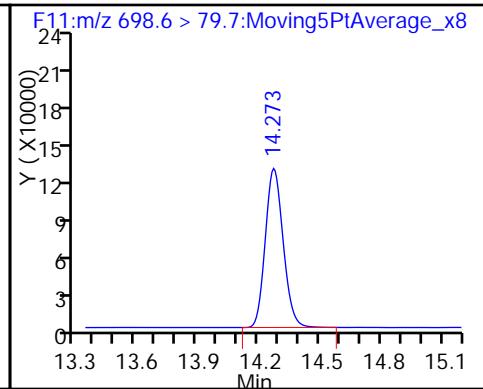
## 29 Perfluorododecanoic acid



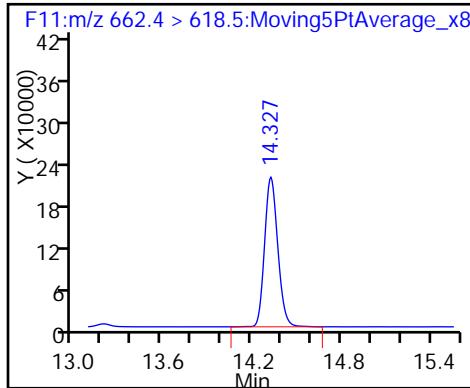
## D 28 13C2 PFDoA



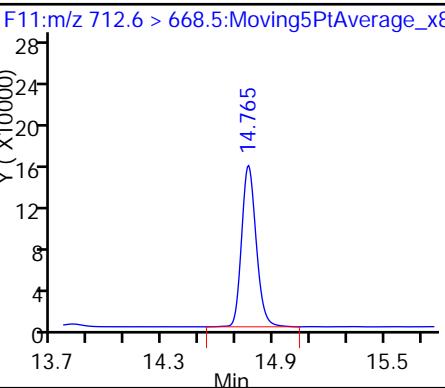
## 31 PFDoS (Perflouro-1-dodecanesulfona



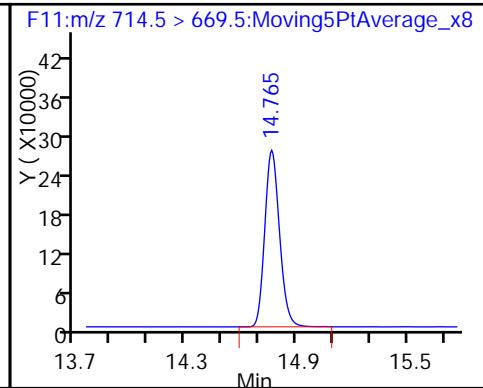
## 30 Perfluorotridecanoic acid



## 32 Perfluorotetradecanoic acid



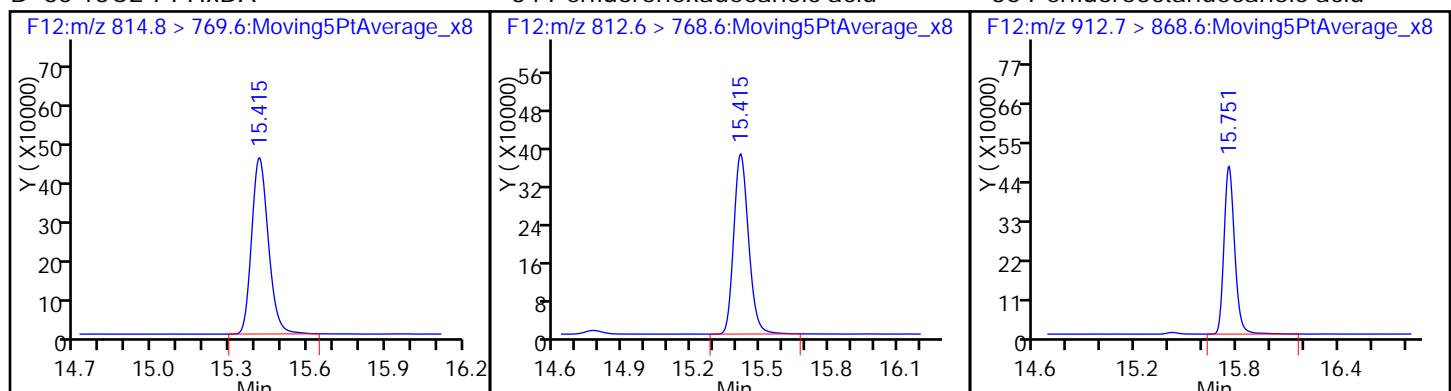
## D 33 13C2-PFTeDA



D 35 13C2-PFHxDa

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_044.d  
 Lims ID: Std L5  
 Client ID:  
 Sample Type: IC Calib Level: 5  
 Inject. Date: 04-Dec-2014 05:09:08 ALS Bottle#: 8 Worklist Smp#: 8  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L5 PFC 50/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:55 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

First Level Reviewer: barnettj Date: 04-Dec-2014 14:16:16

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 113C4 PFBA										
216.7 > 171.5	5.934	5.934	0.0		76134	17.8		89.0	270	
2 Perfluorobutyric acid										
212.7 > 168.6	5.934	5.935	-0.001	1.000	973858	56.5		113	2549	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.039	7.037	0.002	1.000	656489	54.1		108	922	
D 3 113C5-PFPeA										
267.6 > 222.7	7.046	7.038	0.008		146639	19.0		95.0	503	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.161	7.157	0.004	1.000	2969104	63.0		142	6743	
298.8 > 98.6	7.158	7.157	0.001	0.999	1828732	1.62(0.00-0.00)		142	4401	
D 6 113C2 PFHxA										
314.6 > 269.7	8.284	8.282	0.002		362829	19.0		95.2	950	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.284	8.284	0.0	1.000	1048415	59.2		118	1383	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.360	8.360	0.0	0.876	1904029	50.7		108	3852	
D 8 113C4-PFHxA										
366.6 > 321.6	9.510	9.506	0.004		357639	20.2		101	1026	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.510	9.508	0.002	1.000	1480278	54.3		109	2660	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	2166299	49.1		104	3954	
D 11 118O2 PFHxA										
402.5 > 83.6	9.545	9.544	0.001		904686	20.1		106	2013	
D 12 113C4 PFOA										
416.5 > 371.6	10.633	10.628	0.005		1568925	46.4		92.8	2473	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.633	10.628	0.005	1.000	2738772	52.3		105	2366	
412.8 > 168.7	10.633	10.628	0.005	1.000	911992		3.00(0.00-0.00)	105	1636	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.633	10.635	-0.002	1.000	2202840	57.7		121	3728	
D 16 13C4 PFOS										
502.4 > 79.7	11.594	11.595	-0.001		2109419	46.8		97.9	3107	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.594	11.595	-0.001	1.000	3201983	40.2		84.1	2416	
498.9 > 98.7	11.602	11.595	0.007	1.001	1798361		1.78(0.00-0.00)	84.1	1906	
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.614	0.006		744378	19.3		96.4	2568	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	2214146	51.6		103	3103	
21 PFNS (Perflouro-1-nananesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	1801756	55.6		116	2778	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	2735529	58.7		117	2845	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		976118	19.0		94.8	1886	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.004	12.998	0.006	1.000	7880884	56.2		112	3240	
D 23 13C8 FOSA										
505.4 > 77.6	12.995	12.998	-0.003		7377513	51.6		103	3100	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	1601818	56.1		116	2268	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1197353	18.5		92.5	2787	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	3119455	58.9		118	3940	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	3670027	62.1		124	2134	
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1198640	18.3		91.7	1577	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	1709417	58.8		122	2751	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.327	14.327	0.0	1.000	3132510	63.9		128	2299	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	1910691	61.3		123	1772	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1394143	19.5		97.5	2441	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.415	15.415	0.0		2223647	21.3		107	2662	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.415	15.415	0.0	1.000	4803324	59.8		120	2343	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.751	15.752	-0.001	1.000	Page 5148634 of 350	60.5		121	1859	12/12/2014

Report Date: 05-Dec-2014 10:01:56

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L5\_00006

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:01:56

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_044.d

Injection Date: 04-Dec-2014 05:09:08

Instrument ID: A4

Lims ID: Std L5

Client ID:

Operator ID: JRB

ALS Bottle#: 8 Worklist Smp#: 8

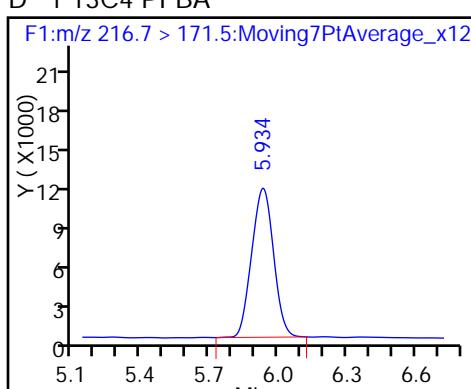
Injection Vol: 15.0 ul

Dil. Factor: 1.0000

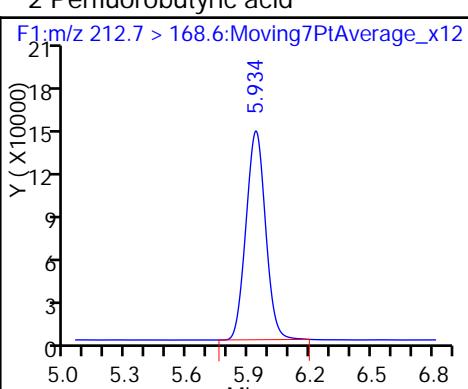
Method: PFAC\_A4

Limit Group: LC PFC ICAL

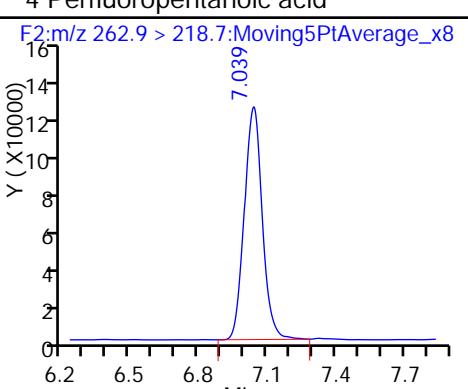
## D 1 13C4 PFBA



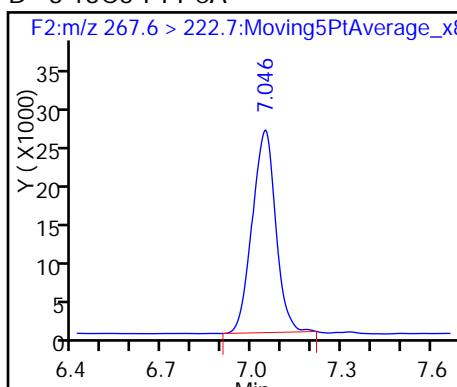
## 2 Perfluorobutyric acid



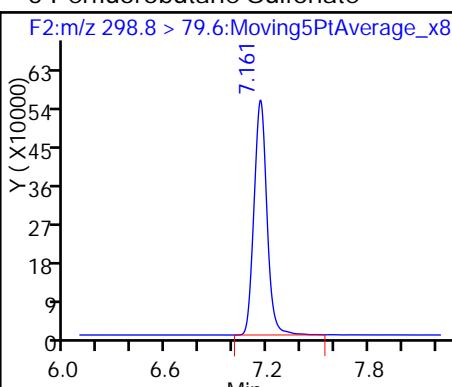
## 4 Perfluoropentanoic acid



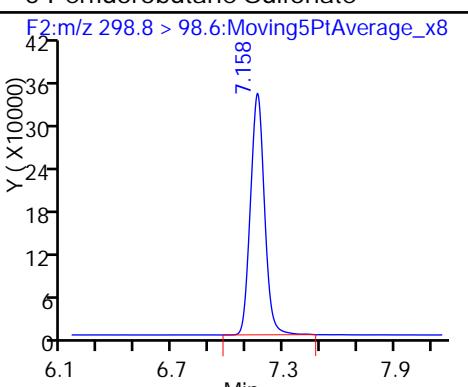
## D 3 13C5-PFPeA



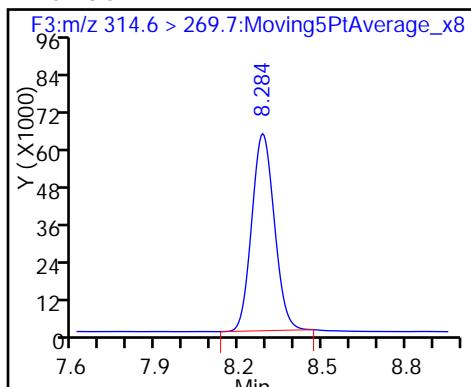
## 5 Perfluorobutane Sulfonate



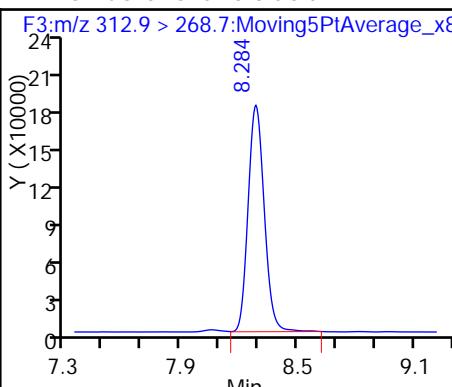
## 5 Perfluorobutane Sulfonate



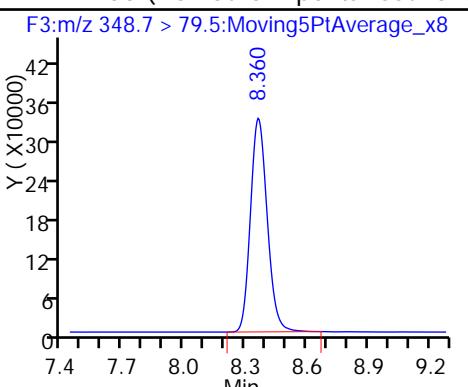
## D 6 13C2 PFHxA



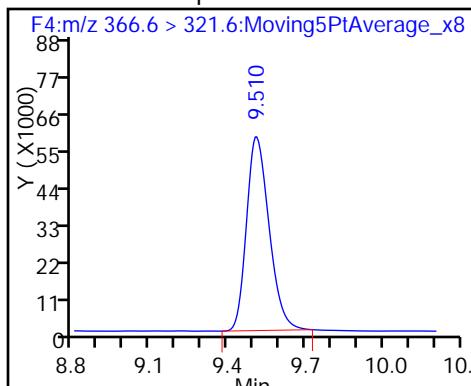
## 7 Perfluorohexanoic acid



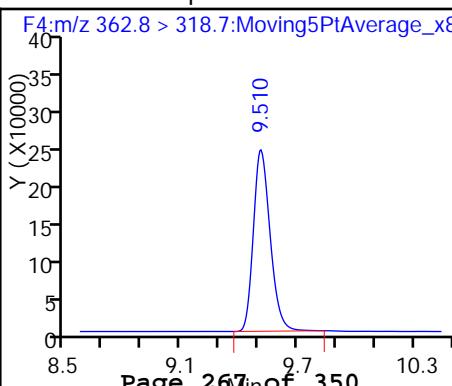
## 22 PFPeS (Perflouro-1-pentanesulfonat)



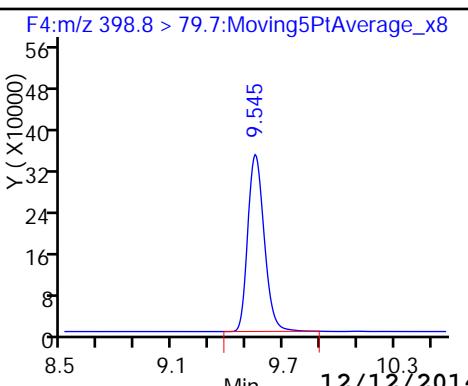
## D 8 13C4-PFHxA



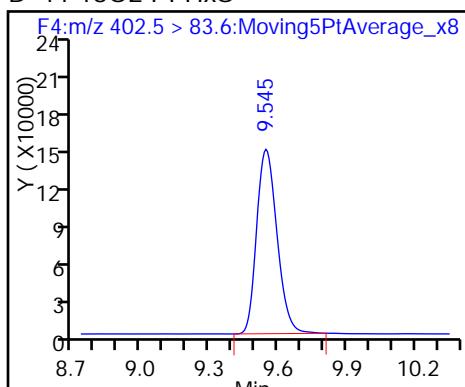
## 9 Perfluoroheptanoic acid



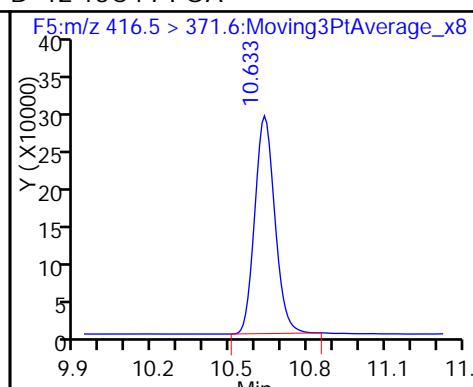
## 10 Perfluorohexane Sulfonate



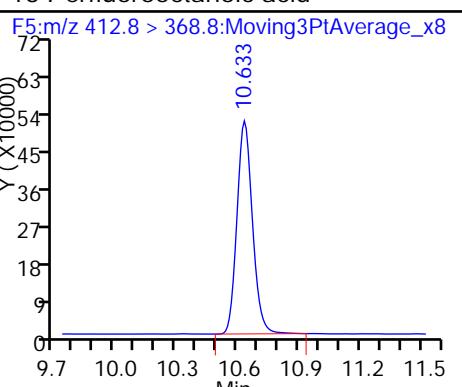
D 11 18O2 PFHxS



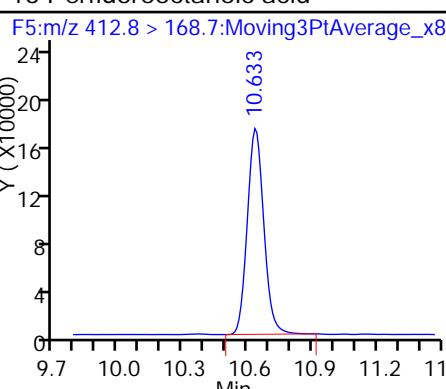
D 12 13C4 PFOA



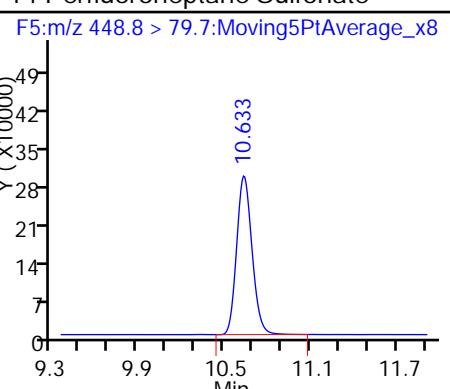
13 Perfluorooctanoic acid



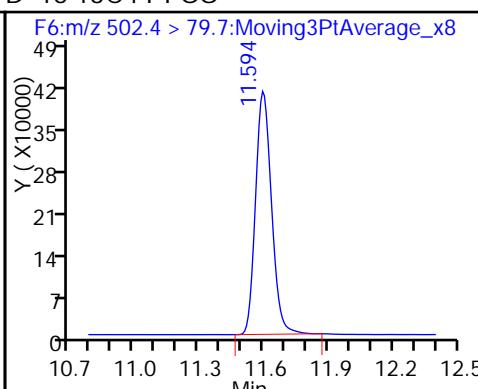
13 Perfluorooctanoic acid



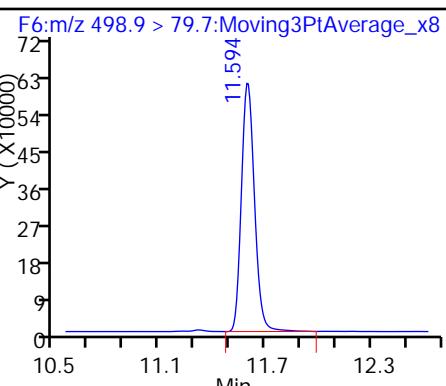
14 Perfluoroheptane Sulfonate



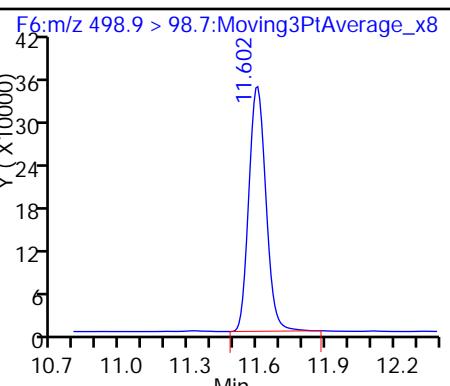
D 16 13C4 PFOS



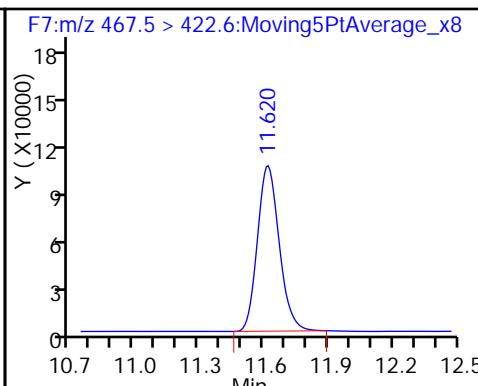
15 Perfluorooctanoic Sulfonate



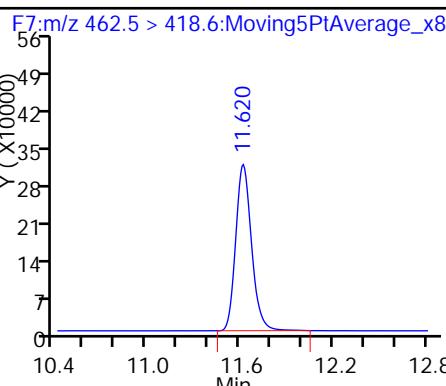
15 Perfluorooctanoic Sulfonate



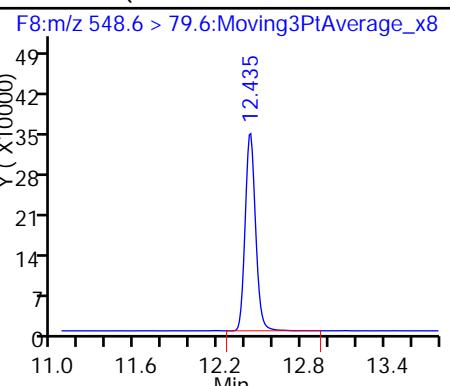
D 17 13C5 PFNA



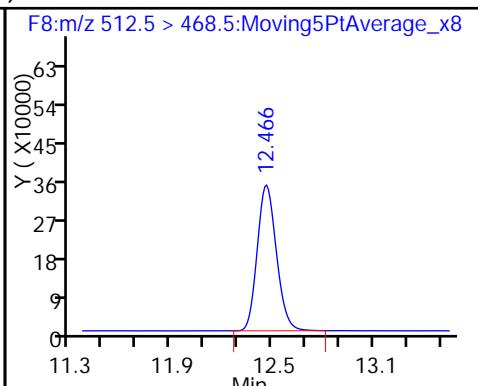
18 Perfluorononanoic acid



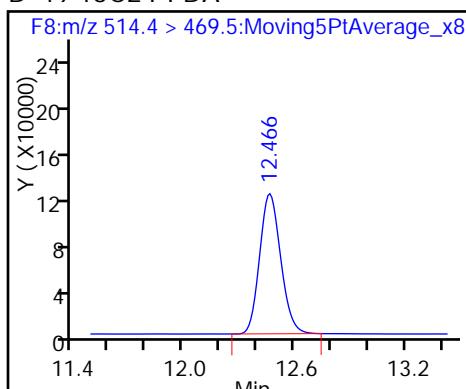
21 PFNS (Perflouro-1-nananesulfonate)



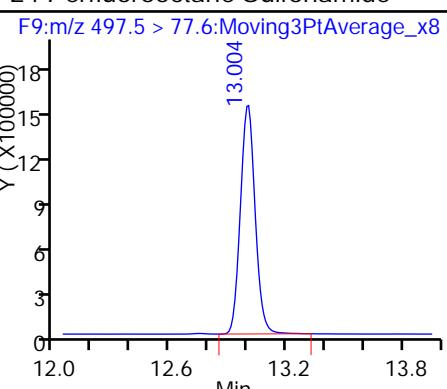
20 Perfluorodecanoic acid



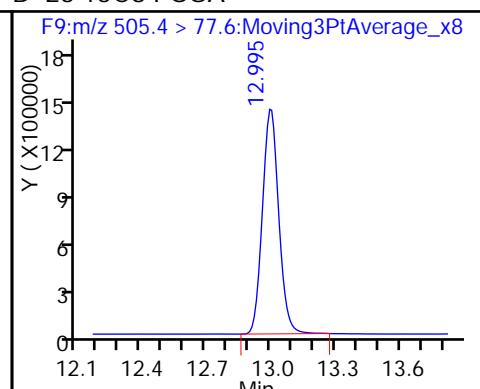
## D 19 13C2 PFDA



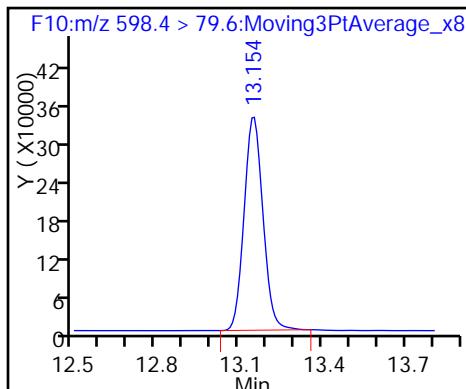
## 24 Perfluorooctane Sulfonamide



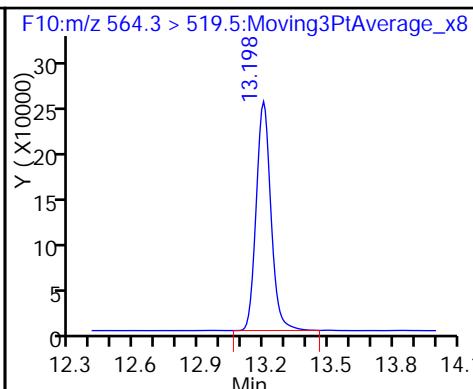
## D 23 13C8 FOSA



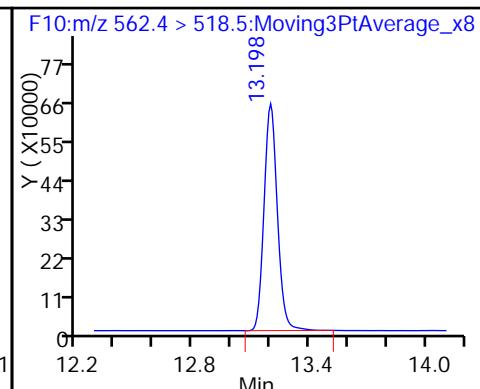
## 25 Perfluorodecane Sulfonate



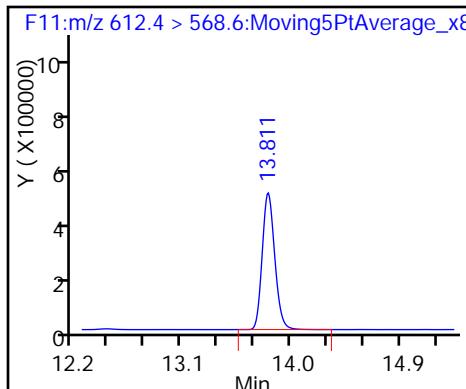
## D 26 13C2 PFUna



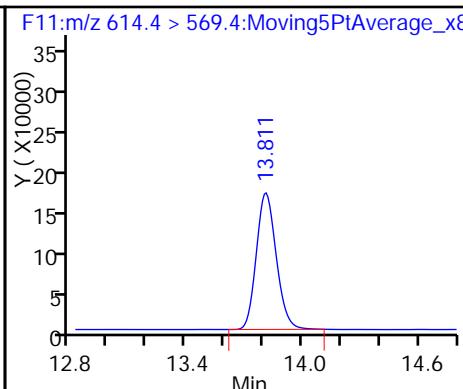
## 27 Perfluoroundecanoic acid



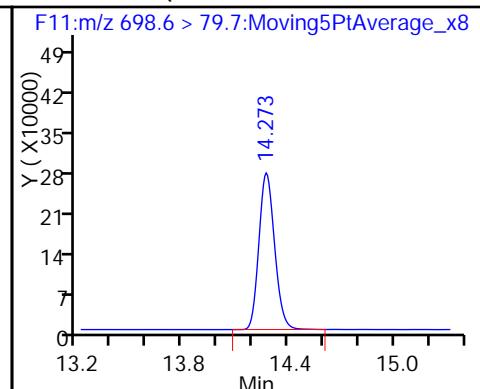
## 29 Perfluorododecanoic acid



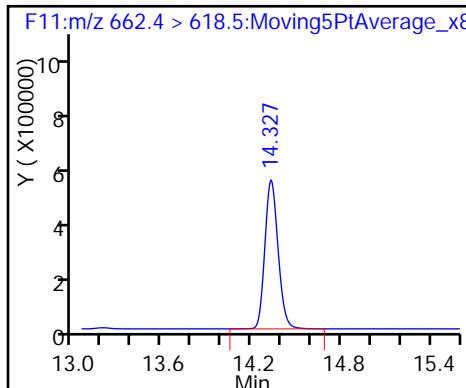
## D 28 13C2 PFDoA



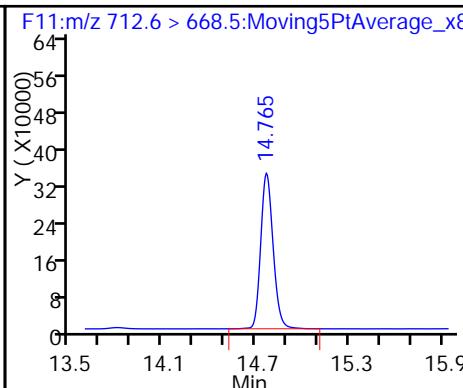
## 31 PFDoS (Perflouro-1-dodecanesulfona



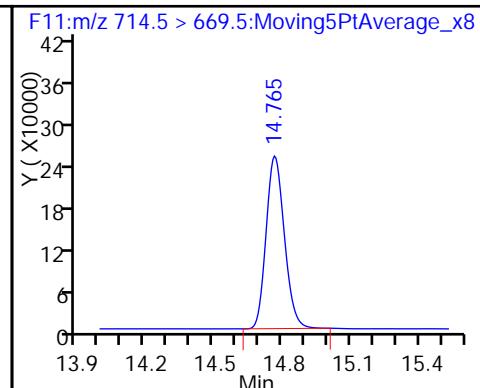
## 30 Perfluorotridecanoic acid



## 32 Perfluorotetradecanoic acid



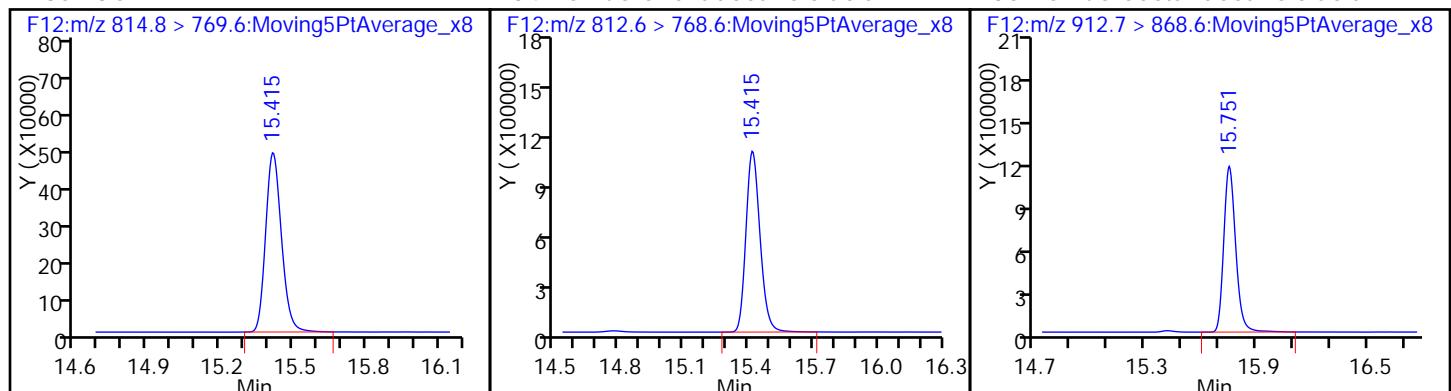
## D 33 13C2-PFTeDA



D 35 13C2-PFHxD

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_045.d  
 Lims ID: Std L6  
 Client ID:  
 Sample Type: IC Calib Level: 6  
 Inject. Date: 04-Dec-2014 05:30:22 ALS Bottle#: 9 Worklist Smp#: 9  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L6 PFC 200/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:57 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

First Level Reviewer: barnettj Date: 04-Dec-2014 10:17:35

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 113C4 PFBA										
216.7 > 171.5	5.934	5.934	0.0		64078	15.0		74.9	172	
2 Perfluorobutyric acid										
212.7 > 168.6	5.934	5.935	-0.001	1.000	3324570	247.1		124	11134	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.036	7.037	-0.001	1.000	2237092	236.4		118	2477	
D 3 113C5-PFPeA										
267.6 > 222.7	7.036	7.038	-0.002		114434	14.8		74.2	390	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.158	7.157	0.001	1.000	6951713	209.3		118	10653	
298.8 > 98.6	7.158	7.157	0.001	1.000	4325906	1.61(0.00-0.00)		118	11278	
D 6 113C2 PFHxA										
314.6 > 269.7	8.284	8.282	0.002		326332	17.1		85.6	1067	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.284	8.284	0.0	1.000	3374043	211.8		106	2254	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.360	8.360	0.0	0.876	6187089	233.7		125	7653	
D 8 113C4-PFHpA										
366.6 > 321.6	9.505	9.506	-0.001		233963	13.2		65.9	638	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.505	9.508	-0.003	1.000	4268821	239.3		120	5537	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	6253308	201.3		106	10496	
D 11 118O2 PFHxS										
402.5 > 83.6	9.545	9.544	0.001		637190	14.2		74.9	1671	
D 12 113C4 PFOA										
416.5 > 371.6	10.632	10.628	0.004		1232205	36.4		72.8	1691	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.632	10.628	0.004	1.000	8615547	209.4		105	5932	
412.8 > 168.7	10.632	10.628	0.004	1.000	2789255		3.09(0.00-0.00)	105	3340	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.642	10.635	0.007	1.000	6297711	218.5		115	5936	
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.595	0.007		1593174	35.3		73.9	2573	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.602	11.595	0.007	1.000	9537705	158.6		82.9	3559	
498.9 > 98.7	11.602	11.595	0.007	1.000	5595082		1.70(0.00-0.00)	82.9	3373	
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.614	0.006		603092	15.6		78.1	1329	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	7868399	226.3		113	7645	
21 PFNS (Perflouro-1-nonenesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	4979770	203.3		106	5351	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	8217900	231.6		116	8527	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		743492	14.4		72.2	1688	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.003	12.998	0.005	1.000	26403028	220.1		110	3107	
D 23 13C8 FOSA										
505.4 > 77.6	13.003	12.998	0.005		6312732	44.1		88.2	3932	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	4451333	206.5		107	4554	
D 26 13C2 PFUnA										
564.3 > 519.5	13.206	13.198	0.008		1176651	18.2		90.9	1960	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.206	13.198	0.008	1.000	9965686	191.6		95.8	6752	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	12549584	227.0		114	3636	
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1120992	17.2		85.8	1041	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	4601893	209.7		108	4870	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.334	14.327	0.007	1.000	9132281	199.3		99.6	3101	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.772	14.766	0.006	1.000	6179569	212.0		106	2829	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.772	14.766	0.006		1224157	17.1		85.6	2164	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.415	15.415	0.0		2119674	20.3		102	2572	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.415	15.415	0.0	1.000	18590504	247.6		124	2880	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.751	15.752	-0.001	1.000	20832627	261.8		131	2455	12/12/2014

Report Date: 05-Dec-2014 10:01:57

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L6\_00005

Amount Added: 1.00

Units: mL

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_045.d

Injection Date: 04-Dec-2014 05:30:22

Instrument ID: A4

Lims ID: Std L6

Client ID:

Operator ID: JRB

ALS Bottle#: 9 Worklist Smp#: 9

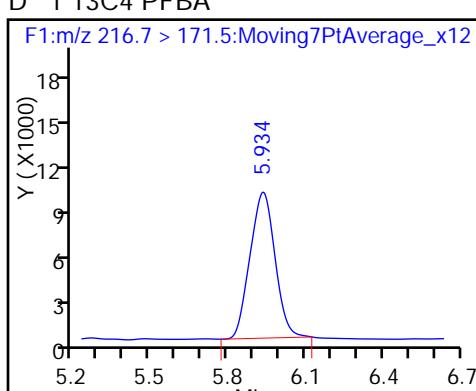
Injection Vol: 15.0 ul

Dil. Factor: 1.0000

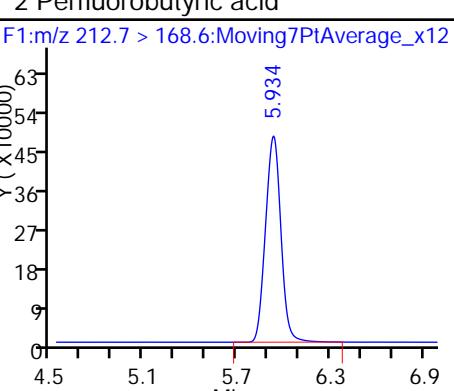
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Limit Group: LC PFC ICAL

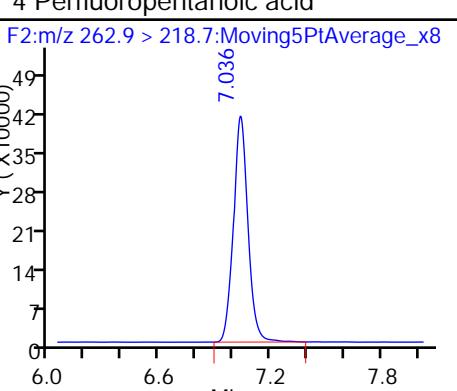
## D 1 13C4 PFBA



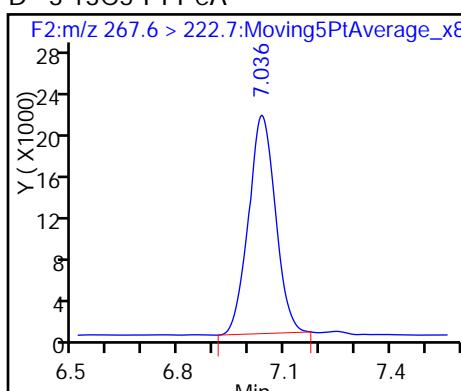
## 2 Perfluorobutyric acid



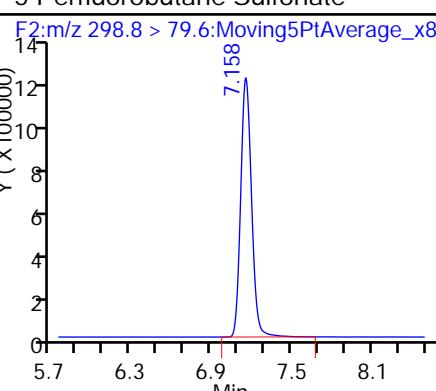
## 4 Perfluoropentanoic acid



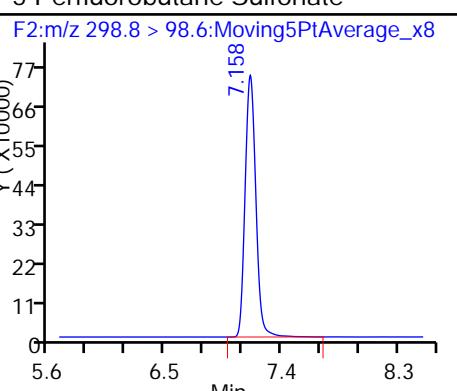
## D 3 13C5-PFPeA



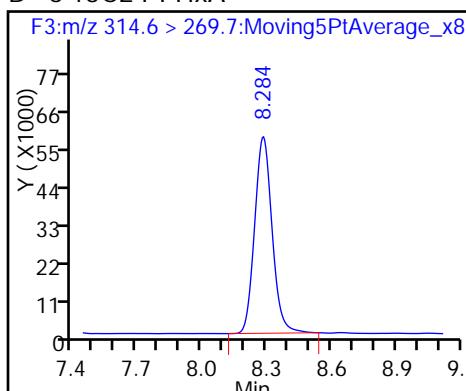
## 5 Perfluorobutane Sulfonate



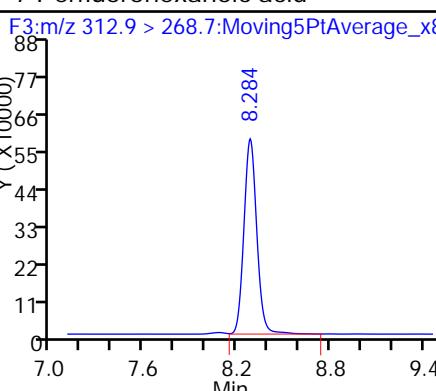
## 5 Perfluorobutane Sulfonate



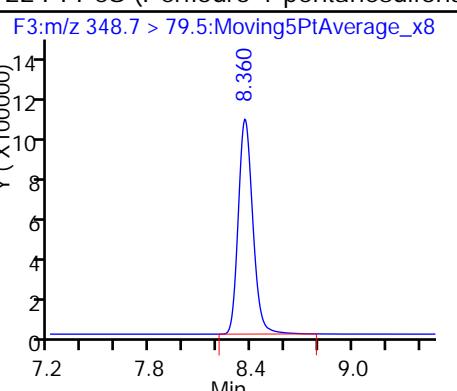
## D 6 13C2 PFHxA



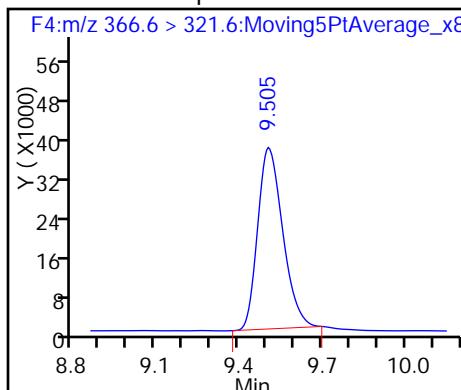
## 7 Perfluorohexanoic acid



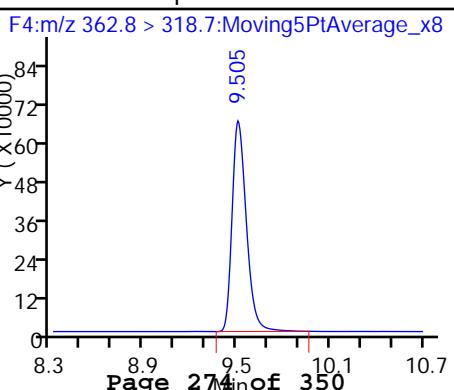
## 22 PFPeS (Perflouro-1-pentanesulfonat)



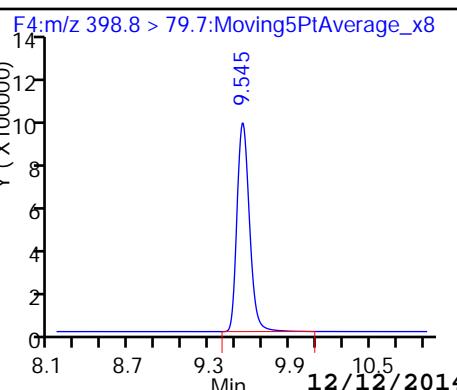
## D 8 13C4-PFHxA



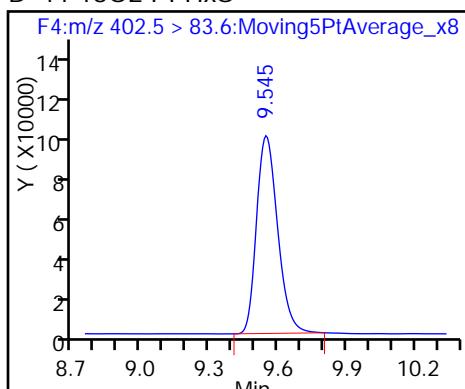
## 9 Perfluoroheptanoic acid



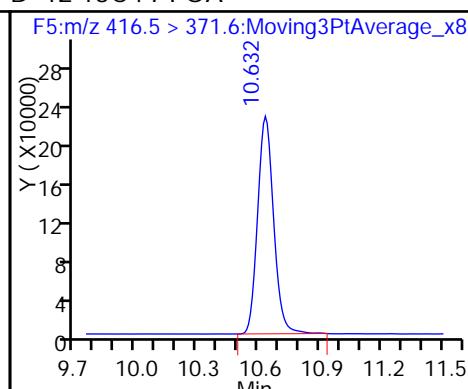
## 10 Perfluorohexane Sulfonate



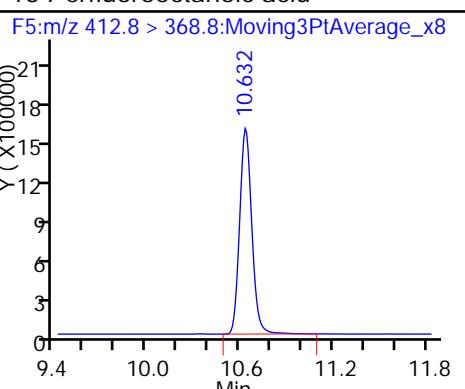
D 11 18O2 PFHxS



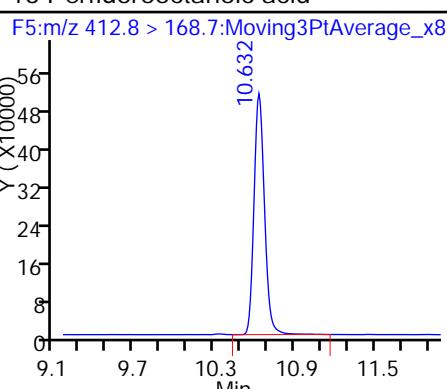
D 12 13C4 PFOA



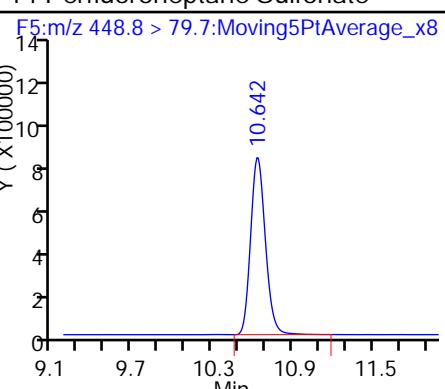
13 Perfluorooctanoic acid



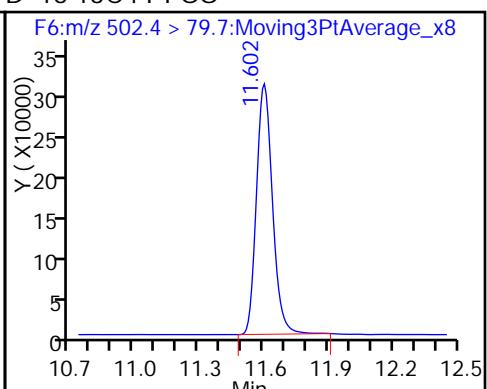
13 Perfluorooctanoic acid



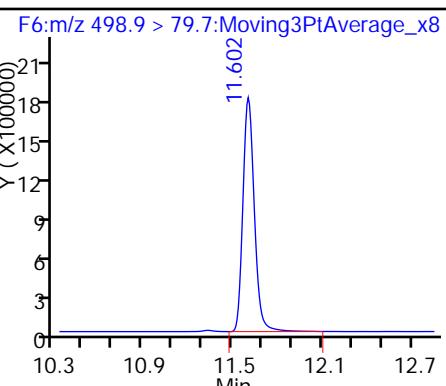
14 Perfluoroheptane Sulfonate



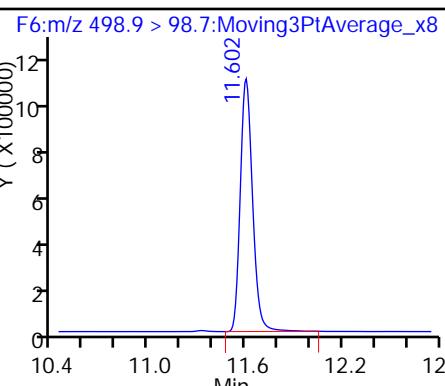
D 16 13C4 PFOS



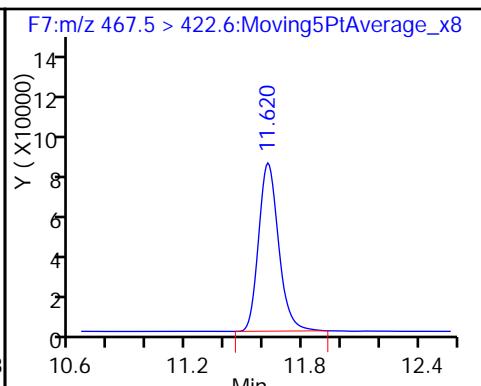
15 Perfluorooctanoic Sulfonate



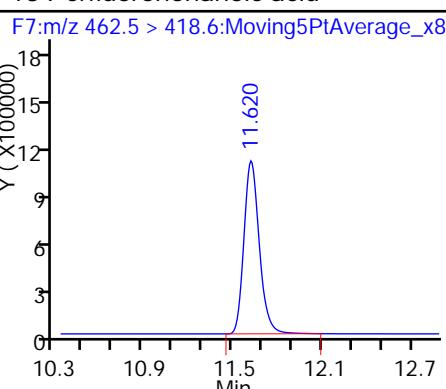
15 Perfluorooctanoic Sulfonate



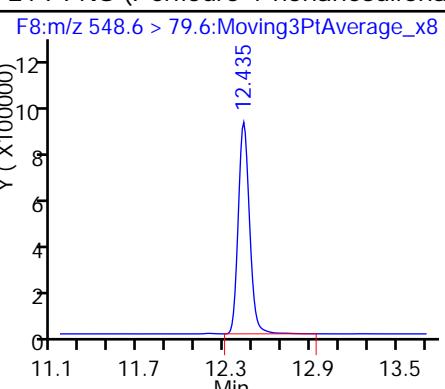
D 17 13C5 PFNA



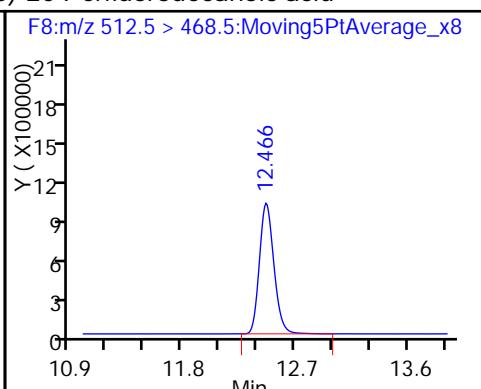
18 Perfluorononanoic acid



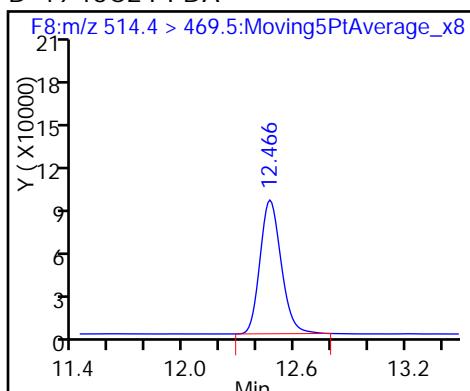
21 PFNS (Perflouro-1-nananesulfonate)



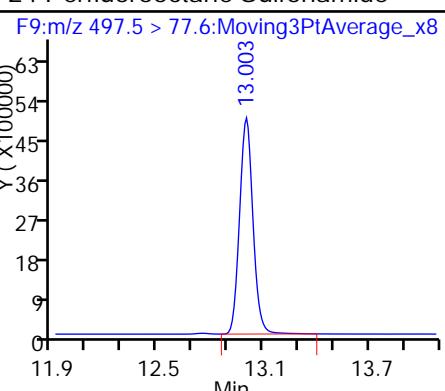
20 Perfluorodecanoic acid



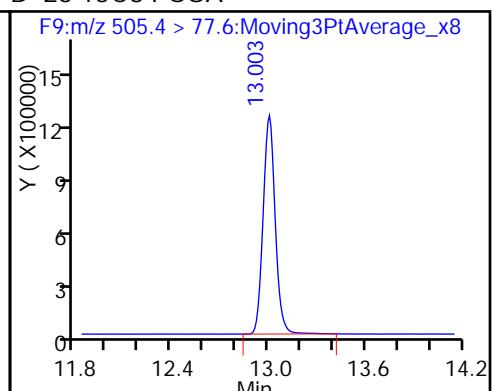
D 19 13C2 PFDA



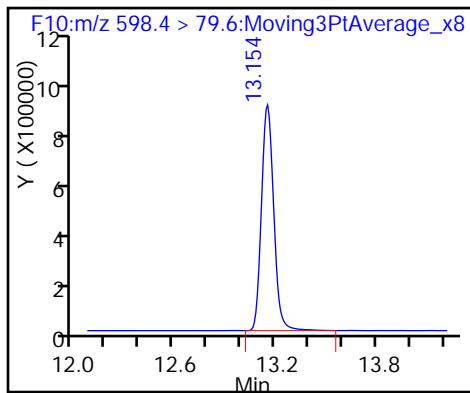
24 Perfluorooctane Sulfonamide



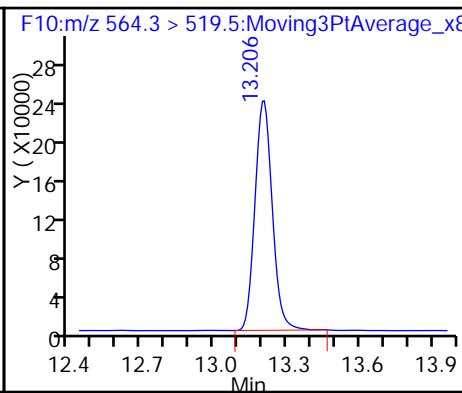
D 23 13C8 FOSA



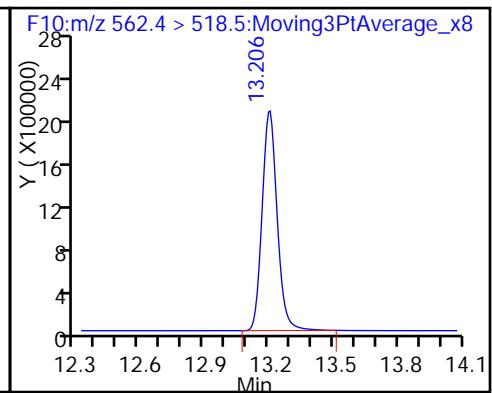
25 Perfluorodecane Sulfonate



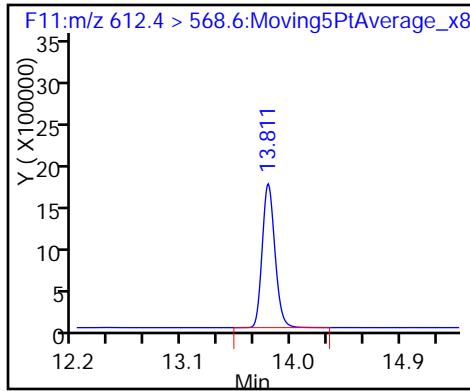
D 26 13C2 PFUna



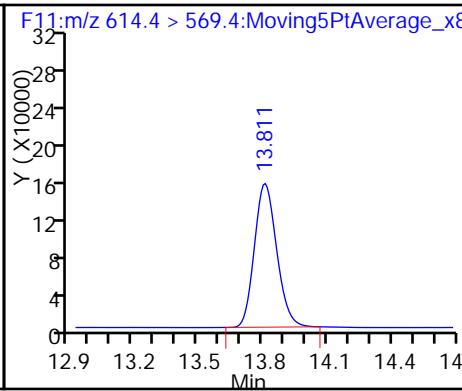
27 Perfluoroundecanoic acid



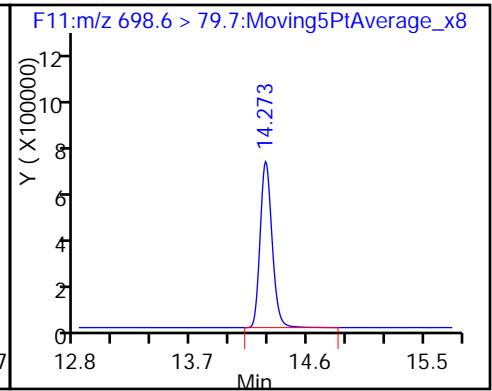
29 Perfluorododecanoic acid



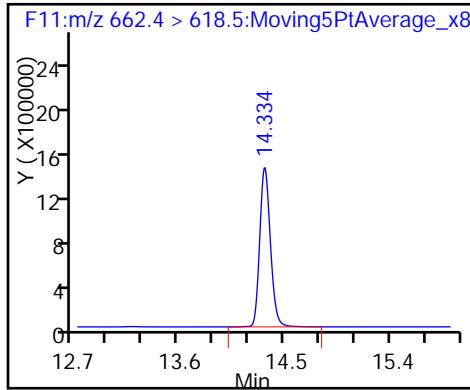
D 28 13C2 PFDoA



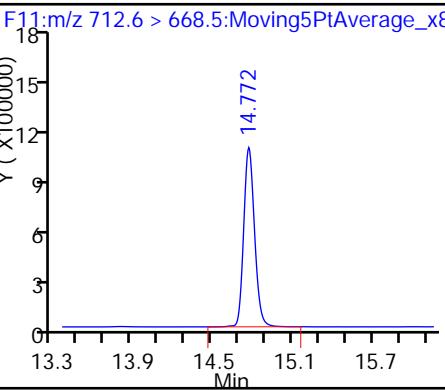
31 PFDoS (Perflouro-1-dodecanesulfona



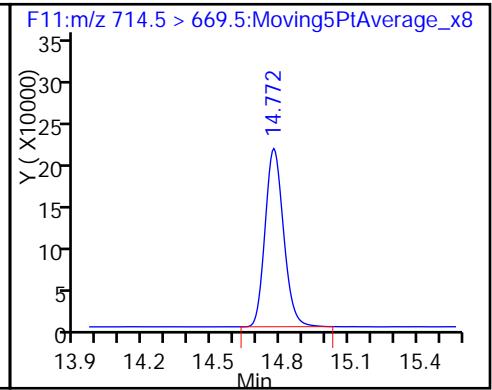
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



D 33 13C2-PFTeDA



Report Date: 05-Dec-2014 10:01:57

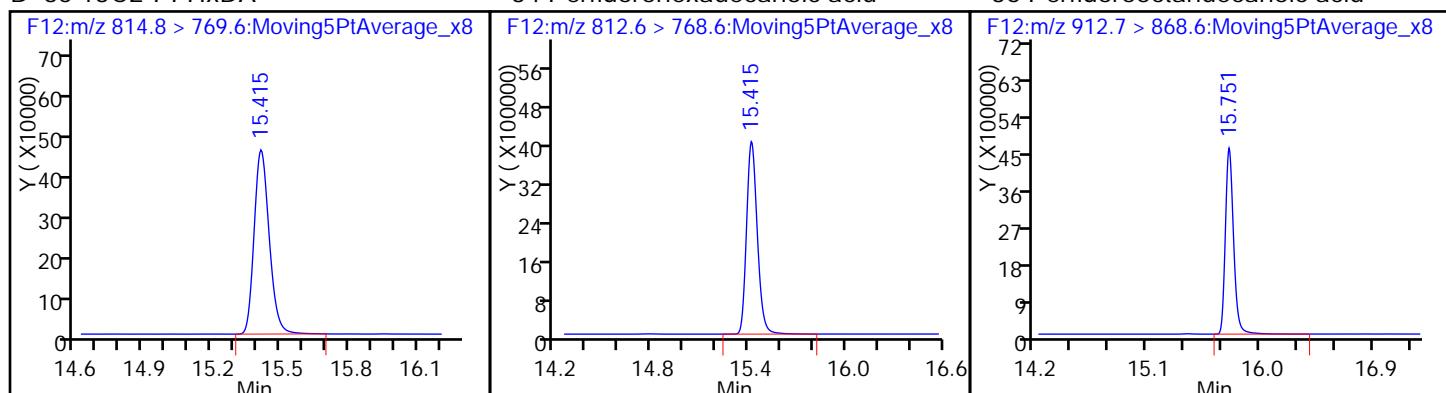
Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_045.d

D 35 13C2-PFHxDA

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Lims ID: Std L7  
 Client ID:  
 Sample Type: IC Calib Level: 7  
 Inject. Date: 04-Dec-2014 05:51:34 ALS Bottle#: 10 Worklist Smp#: 10  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: LCPFC-L7 PFC 500/50ng/mL  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:01:58 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 113C4 PFBA										
216.7 > 171.5	5.934	5.934	0.0		45086	10.5		52.7	168	
2 Perfluorobutyric acid										
212.7 > 168.6	5.941	5.935	0.006	1.000	6084630	531.4		106	12919	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.040	7.037	0.003	1.000	4729024	587.2		117	5934	
D 3 113C5-PFPeA										
267.6 > 222.7	7.040	7.038	0.002		97387	12.6		63.1	247	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.162	7.157	0.005	1.000	14216894	562.7		127	13614	
298.8 > 98.6	7.162	7.157	0.005	1.000	8687902	1.64(0.00-0.00)		127	15376	
D 6 113C2 PFHxA										
314.6 > 269.7	8.286	8.282	0.004		228440	12.0		59.9	527	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.291	8.284	0.007	1.000	6260320	561.5		112	1620	
22 PPoS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.362	8.360	0.002	0.875	11656614	578.9		123	13351	
D 8 113C4-PFHxA										
366.6 > 321.6	9.510	9.506	0.004		207868	11.7		58.6	427	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.505	9.508	-0.003	1.000	9446001	596.1		119	9459	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.552	9.543	0.009	1.000	12604320	533.5		113	11248	
D 11 118O2 PFHxA										
402.5 > 83.6	9.552	9.544	0.008		484686	10.8		57.0	923	
D 12 113C4 PFOA										
416.5 > 371.6	10.623	10.628	-0.005		952358	28.2		56.3	1312	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.623	10.628	-0.005	1.000	16683058	524.6		105	6845	
412.8 > 168.7	10.632	10.628	0.004	1.001	5344475		3.12(0.00-0.00)	105	3569	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.632	10.635	-0.003	1.000	11448389	570.8		120	5166	
D 16 13C4 PFOS										
502.4 > 79.7	11.594	11.595	-0.001		1108797	24.6		51.4	1556	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.594	11.595	-0.001	1.000	17552751	419.3		87.7	2797	
498.9 > 98.7	11.594	11.595	-0.001	1.000	10235840		1.71(0.00-0.00)	87.7	3533	
D 17 13C5 PFNA										
467.5 > 422.6	11.612	11.614	-0.002		440238	11.4		57.0	1058	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	13621836	536.6		107	8608	
21 PFNS (Perflouro-1-nonenesulfonate)										
548.6 > 79.6	12.425	12.429	-0.004	1.000	8895061	521.8		109	5811	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	16591989	538.8		108	7125	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		645298	12.5		62.7	912	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.004	12.998	0.006	1.000	58025028	571.7		114	3532	
D 23 13C8 FOSA										
505.4 > 77.6	13.004	12.998	0.006		5341425	37.3		74.7	2000	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.146	13.152	-0.006	1.000	8000344	533.3		111	5011	
D 26 13C2 PFUnA										
564.3 > 519.5	13.206	13.198	0.008		1397054	21.6		108	2197	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	20245258	327.9		65.6	9226	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.813	13.808	0.005	1.000	25119659	482.3		96.5	3670	
D 28 13C2 PFDoA										
614.4 > 569.4	13.813	13.809	0.004		1056139	16.2		80.8	956	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.274	14.272	0.002	1.000	8583795	562.0		116	3466	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.328	14.327	0.001	1.000	20104094	465.7		93.1	3264	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.766	14.766	0.0	1.000	13816653	503.2		101	3017	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.766	14.766	0.0		1167304	16.3		81.6	1880	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.415	15.415	0.0		1876393	18.0		90.0	1772	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.415	15.415	0.0	1.000	39084771	552.4		110	2331	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.752	15.752	0.0	1.000	47874246	638.6	Page 279 of 350	128	2236	12/12/2014

Report Date: 05-Dec-2014 10:01:58

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L7\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:01:58

Chrom Revision: 2.2 06-Nov-2014 14:50:32

## TestAmerica Sacramento

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_046.d

Injection Date: 04-Dec-2014 05:51:34

Instrument ID: A4

Lims ID: Std L7

Client ID:

Operator ID: JRB

ALS Bottle#: 10 Worklist Smp#: 10

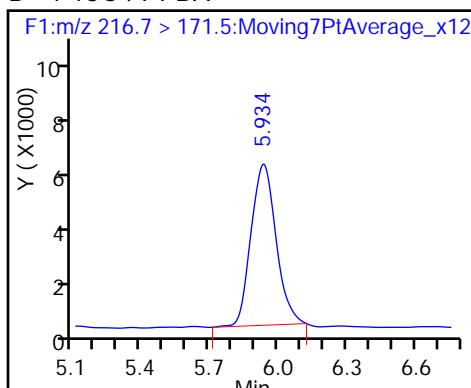
Injection Vol: 15.0 ul

Dil. Factor: 1.0000

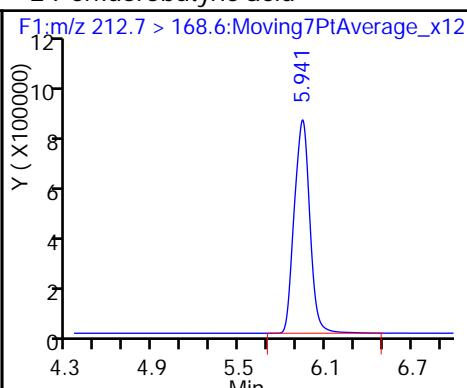
Method: PFAC\_A4

Limit Group: LC PFC ICAL

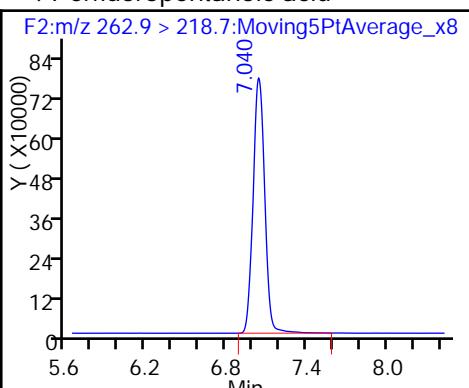
## D 1 13C4 PFBA



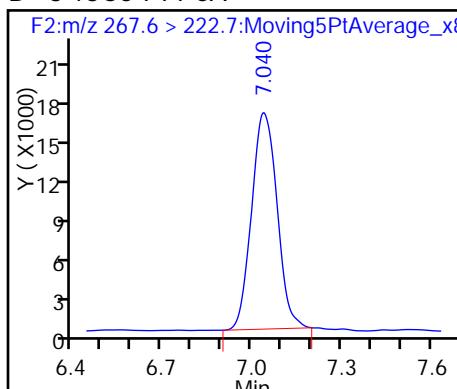
## 2 Perfluorobutyric acid



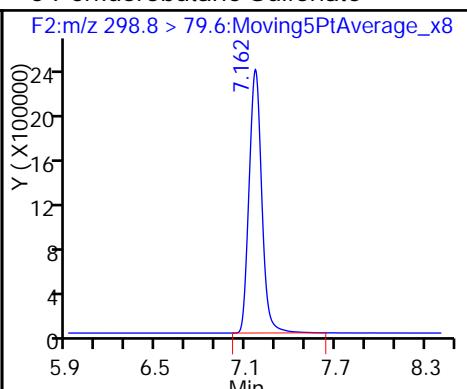
## 4 Perfluoropentanoic acid



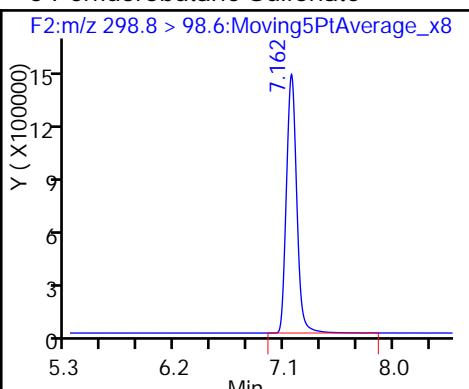
## D 3 13C5-PFPeA



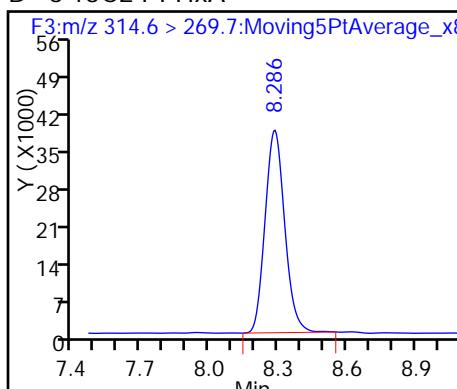
## 5 Perfluorobutane Sulfonate



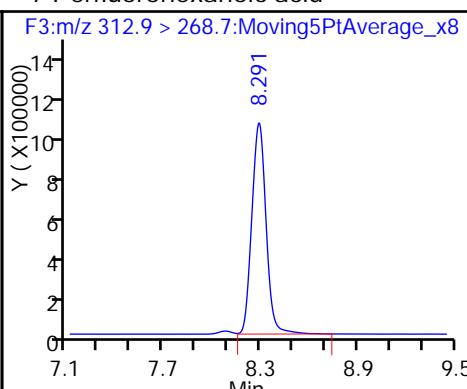
## 5 Perfluorobutane Sulfonate



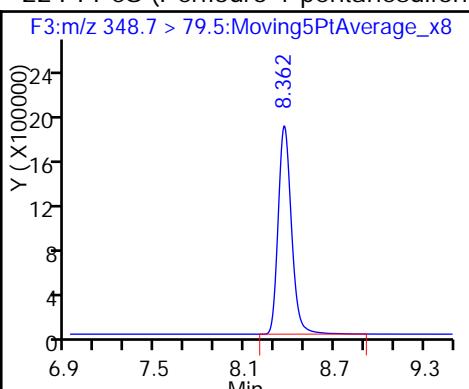
## D 6 13C2 PFHxA



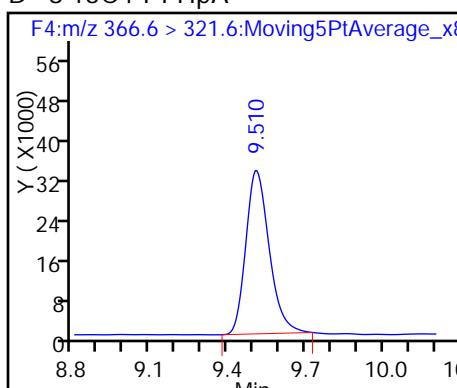
## 7 Perfluorohexanoic acid



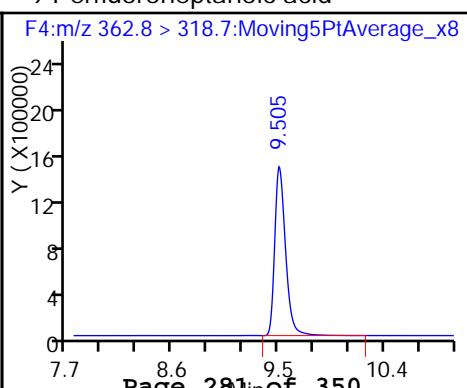
## 22 PFPeS (Perflouro-1-pentanesulfonat)



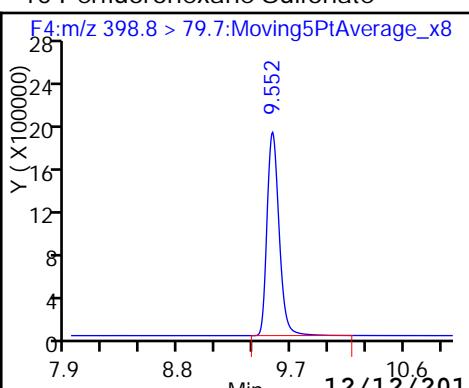
## D 8 13C4-PFHxA



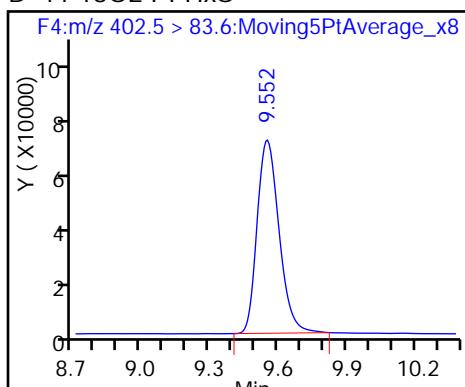
## 9 Perfluoroheptanoic acid



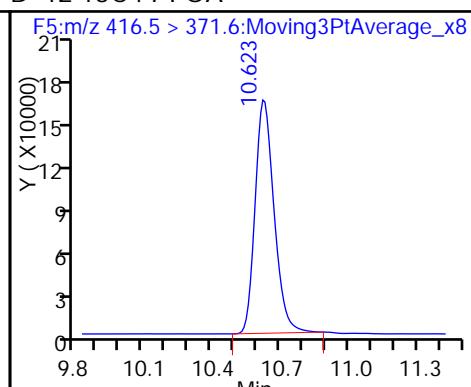
## 10 Perfluorohexane Sulfonate



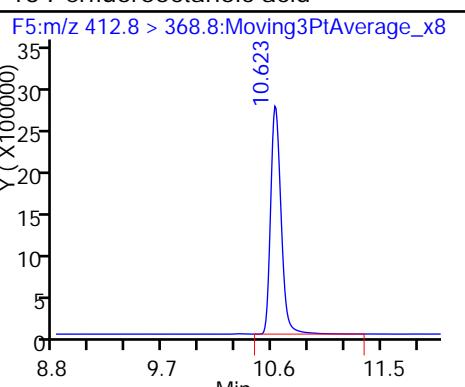
D 11 18O2 PFHxS



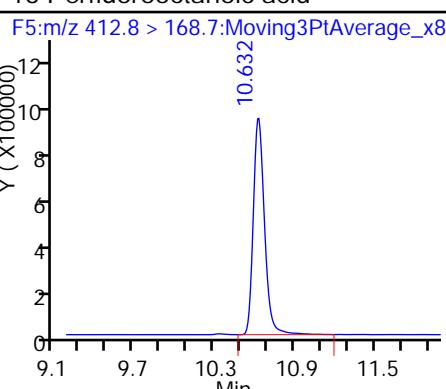
D 12 13C4 PFOA



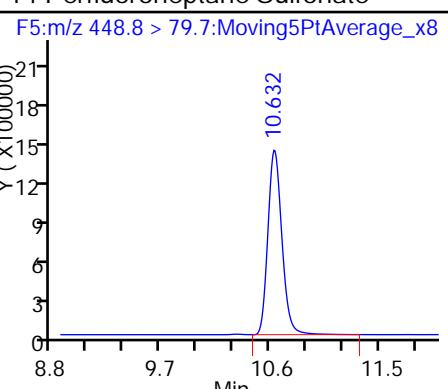
13 Perfluorooctanoic acid



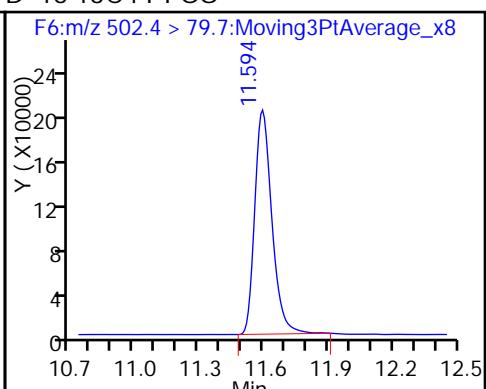
13 Perfluorooctanoic acid



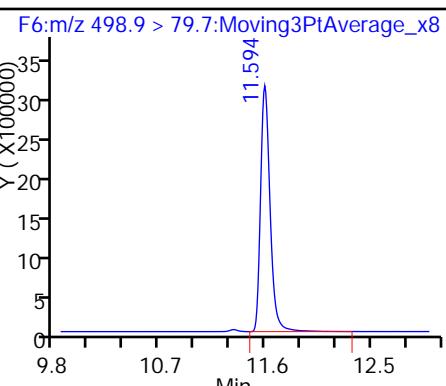
14 Perfluoroheptane Sulfonate



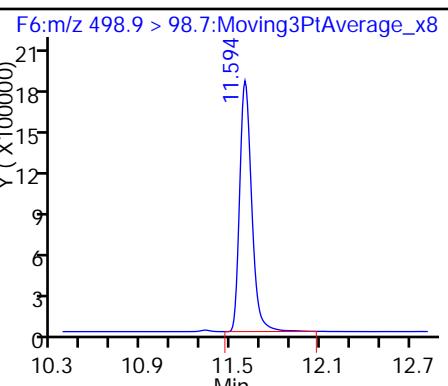
D 16 13C4 PFOS



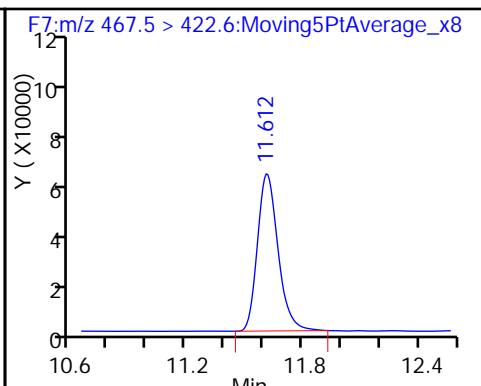
15 Perfluorooctanoic Sulfonate



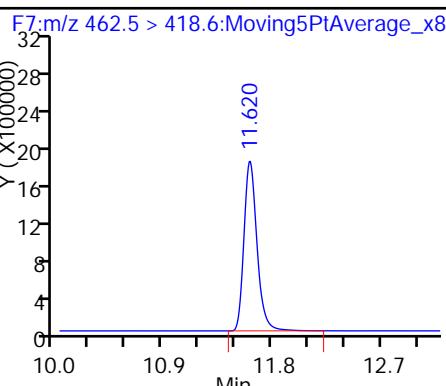
15 Perfluorooctanoic Sulfonate



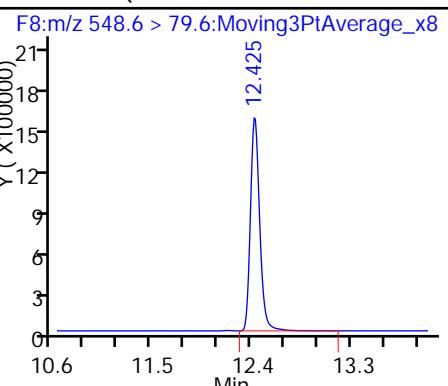
D 17 13C5 PFNA



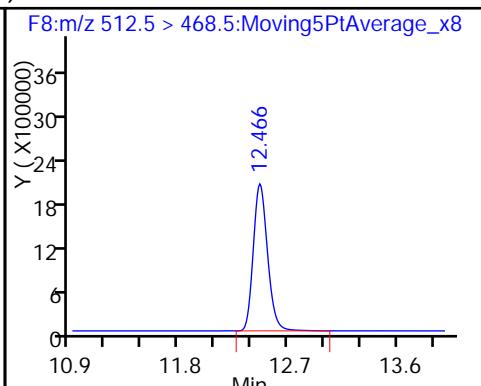
18 Perfluorononanoic acid



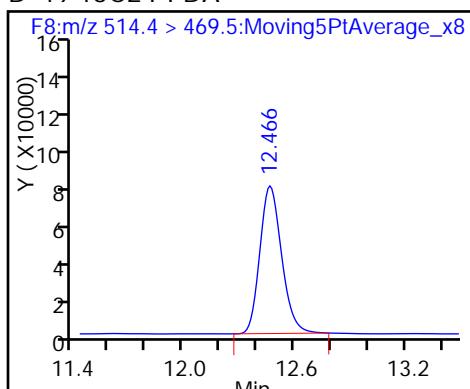
21 PFNS (Perflouro-1-nananesulfonate)



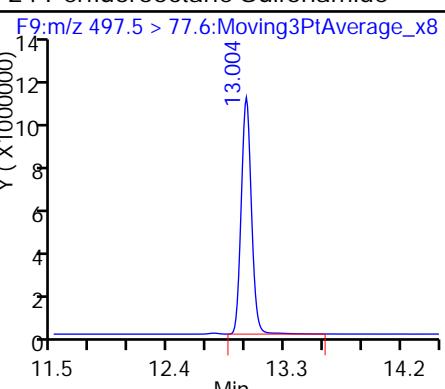
20 Perfluorodecanoic acid



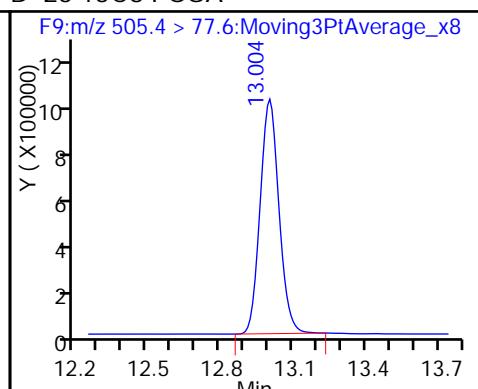
D 19 13C2 PFDA



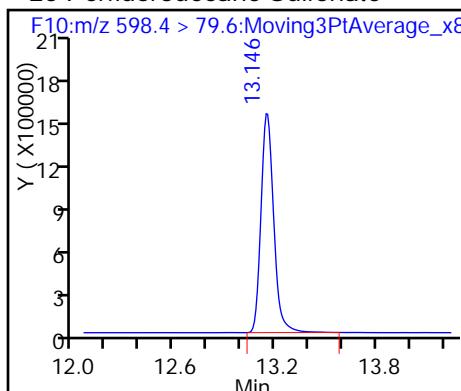
24 Perfluorooctane Sulfonamide



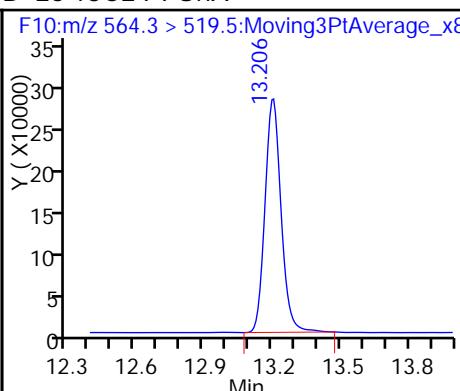
D 23 13C8 FOSA



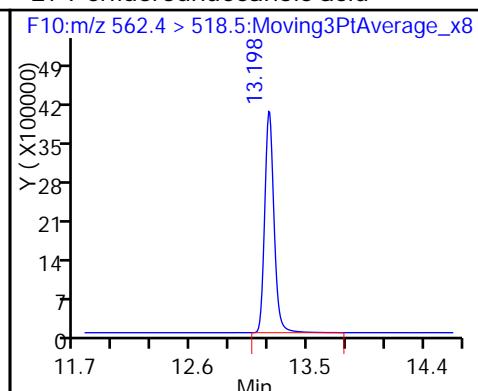
25 Perfluorodecane Sulfonate



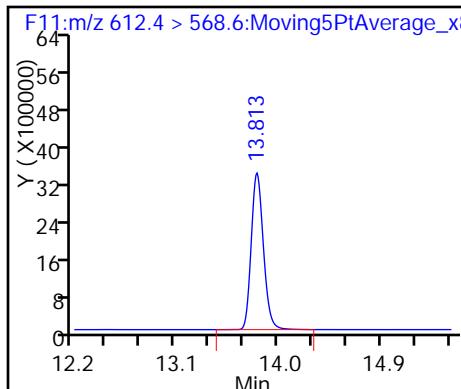
D 26 13C2 PFUna



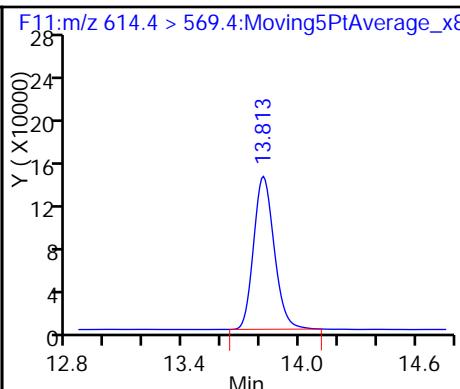
27 Perfluoroundecanoic acid



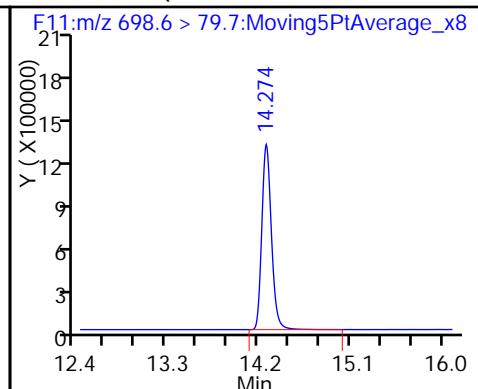
29 Perfluorododecanoic acid



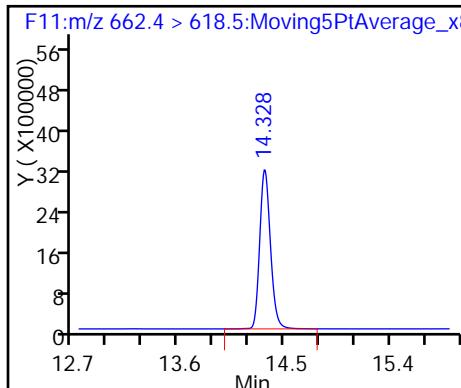
D 28 13C2 PFDoA



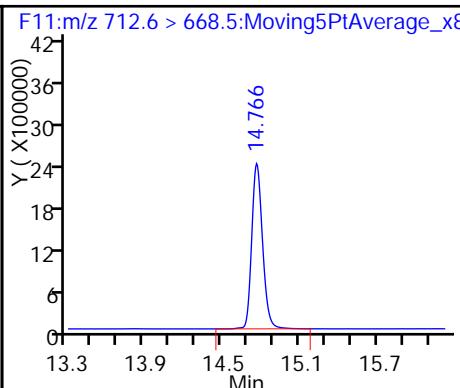
31 PFDoS (Perflouro-1-dodecanesulfona



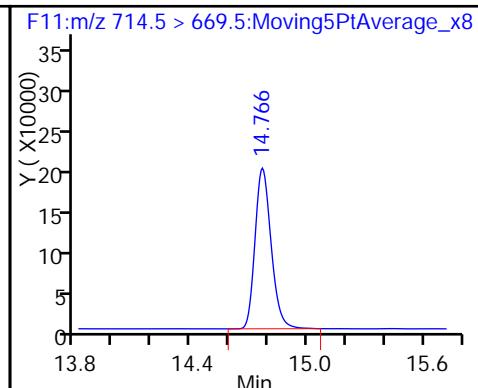
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



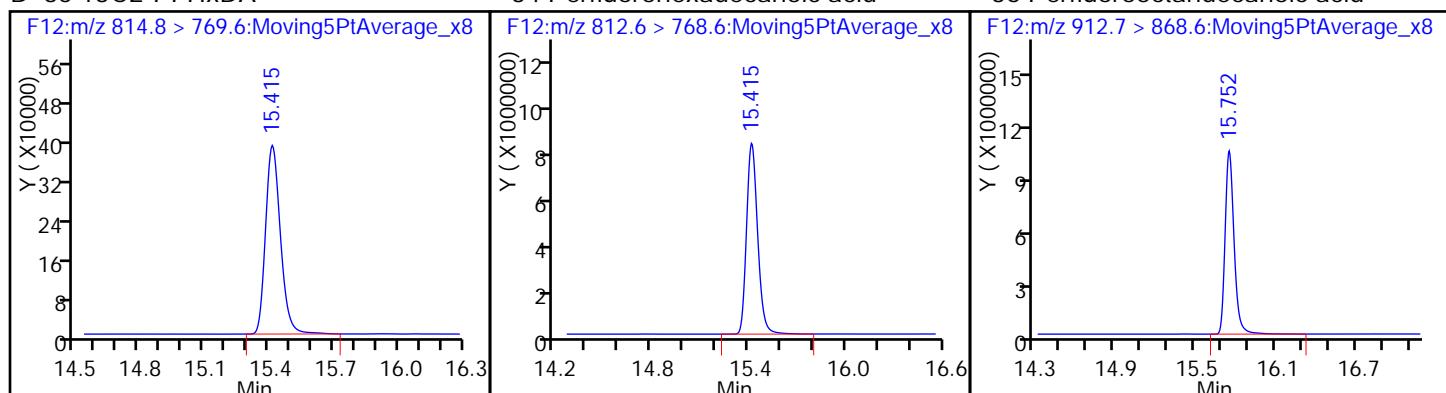
D 33 13C2-PFTeDA



D 35 13C2-PFHxDa

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Lab Sample ID: ICV 320-59779/22 Calibration Date: 12/04/2014 10:23  
Instrument ID: A4 Calib Start Date: 12/04/2014 03:44  
GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51  
Lab File ID: 03DEC14A4A\_058.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	2.637		56.1	50.0	12.2	50.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	1.886		57.0	50.0	14.0	50.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.236		55.5	44.3	25.3	50.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.190		61.0	50.0	21.9	50.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.160		70.8	50.0	41.7	50.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	1.119		57.4	47.3	21.4	50.0
Perfluoro-1-heptanesulfonate (PFHps)	AveID	0.8647	0.9109		50.1	47.6	5.3	50.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.812		54.2	50.0	8.5	50.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.655		43.8	47.8	-8.3	50.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.206		52.3	50.0	4.6	50.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.211		63.4	50.0	26.8	50.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	0.9856		51.9	50.0	3.7	50.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7475		55.8	48.3	15.6	50.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.141		64.5	50.0	29.1	50.0
Perfluorododecanoic acid (PFDa)	AveID	0.9863	1.176		59.6	50.0	19.2	50.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9516		58.2	50.0	16.4	50.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.6011		57.8	50.0	15.6	50.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.644		61.3	50.0	22.7	50.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.763		62.1	50.0	24.2	50.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_058.d  
 Lims ID: ICV  
 Client ID:  
 Sample Type: ICV  
 Inject. Date: 04-Dec-2014 10:23:34 ALS Bottle#: 11 Worklist Smp#: 22  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: ICV  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist:  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:02:31 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
--------	----	--------	--------	--------	----------	--------------	---------------	------	-----	-------

D 1 113C4 PFBA										
216.7 > 171.5	5.950	5.934	0.016		78611	18.4		91.9	276	
2 Perfluorobutyric acid										
212.7 > 168.6	5.950	5.935	0.015	1.000	998130	56.1			3772	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.070	7.037	0.033	1.000	713887	57.0			819	
D 3 113C5-PFPeA										
267.6 > 222.7	7.067	7.038	0.029		151391	19.6		98.1	382	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.186	7.157	0.029	1.000	2103191	55.5			4835	
298.8 > 98.6	7.186	7.157	0.029	1.000	1253987	1.68(0.00-0.00)			2670	
D 6 113C2 PFHxA										
314.6 > 269.7	8.317	8.282	0.035		380517	20.0		99.8	1366	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.317	8.284	0.033	1.000	1132047	61.0			1862	
D 8 113C4-PFHxA										
366.6 > 321.6	9.538	9.506	0.032		290495	16.4		81.9	1032	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.538	9.508	0.030	1.000	1568563	70.8			3709	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.573	9.543	0.030	1.000	2033871	57.4			4859	
D 11 118O2 PFHxS										
402.5 > 83.6	9.573	9.544	0.029		727555	16.2		85.5	2121	
D 12 113C4 PFOA										
416.5 > 371.6	10.660	10.628	0.032		1537054	45.4		90.9	2980	
13 Perfluorooctanoic acid										
412.8 > 368.8	10.660	10.628	0.032	1.000	2784461	54.2			2059	
412.8 > 168.7	10.660	10.628	0.032	1.000	901089	3.09(0.00-0.00)			1834	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.660	10.635	0.025	1.000	1849156	50.1				2439
D 16 13C4 PFOS										
502.4 > 79.7	11.610	11.595	0.015		2038537	45.2		94.6	3811	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.619	11.595	0.024	1.000	3369395	43.8				2694
498.9 > 98.7	11.619	11.595	0.024	1.000	1930634		1.75(0.00-0.00)			2401
D 17 13C5 PFNA										
467.5 > 422.6	11.637	11.614	0.023		804258	20.8		104	2517	
18 Perfluorononanoic acid										
462.5 > 418.6	11.637	11.616	0.021	1.000	2424305	52.3				2893
20 Perfluorodecanoic acid										
512.5 > 468.5	12.476	12.461	0.015	1.000	3129689	63.4				3766
D 19 13C2 PFDA										
514.4 > 469.5	12.476	12.463	0.013		1034113	20.1		100	2116	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.012	12.998	0.014	1.000	7036514	51.9				2509
D 23 13C8 FOSA										
505.4 > 77.6	13.012	12.998	0.014		7139099	49.9		99.8	5003	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	1538194	55.8				2286
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1158043	17.9		89.4	2092	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	3303471	64.5				3542
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	3801675	59.6				2789
D 28 13C2 PFDaA										
614.4 > 569.4	13.811	13.809	0.002		1293615	19.8		99.0	1668	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.327	14.327	0.0	1.000	3077366	58.2				1847
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	1943943	57.8				1926
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1424288	19.9		99.6	2616	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.404	15.415	-0.011		2266653	21.7		109	2388	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.404	15.415	-0.011	1.000	5315127	61.3				2456
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.736	15.752	-0.016	1.000	5702352	62.1				2021

**Reagents:**

LCPFCIC\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:02:31

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_058.d

Injection Date: 04-Dec-2014 10:23:34

Instrument ID: A4

Lims ID: ICV

Client ID:

Operator ID: JRB

ALS Bottle#: 11 Worklist Smp#: 22

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

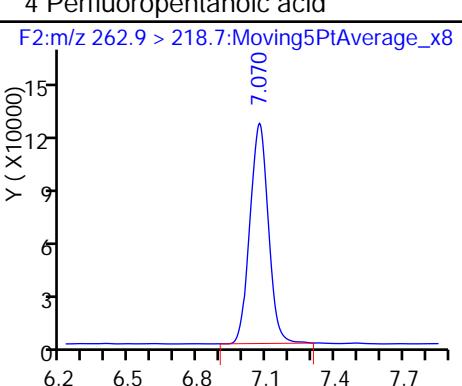
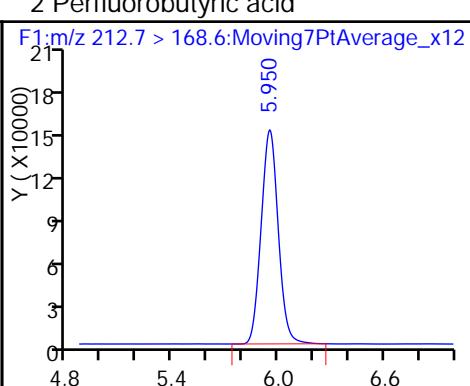
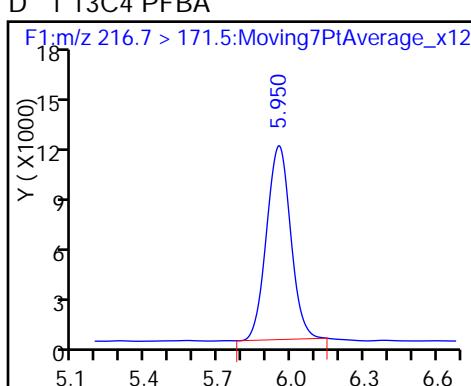
Method: PFAC\_A4

Limit Group: LC PFC ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

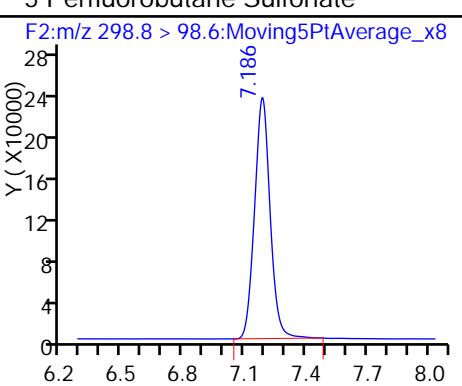
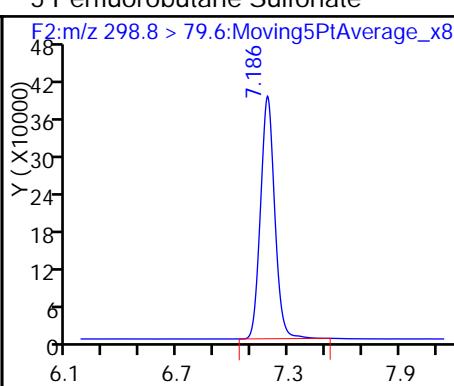
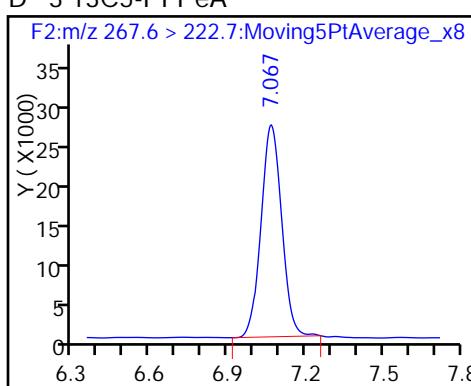
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

5 Perfluorobutane Sulfonate

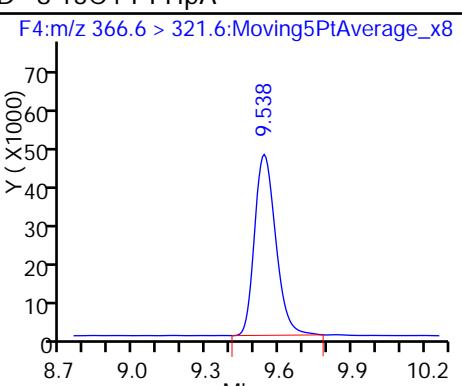
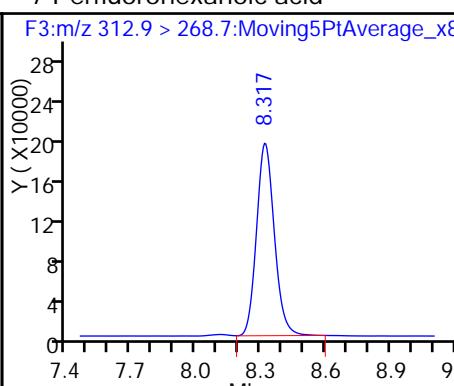
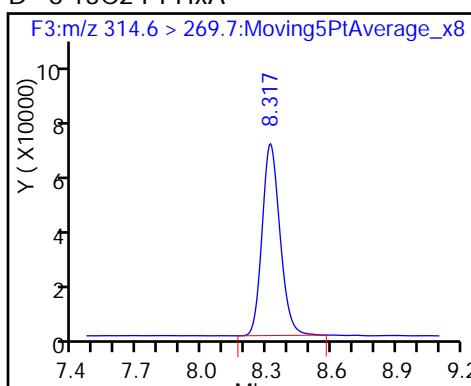
5 Perfluorobutane Sulfonate



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

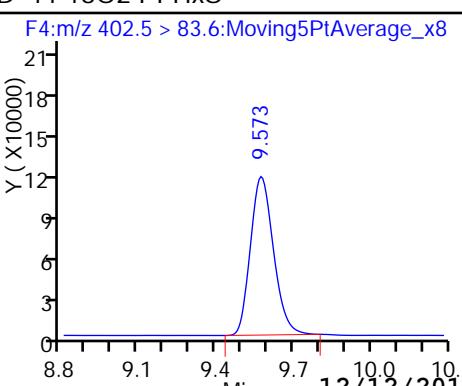
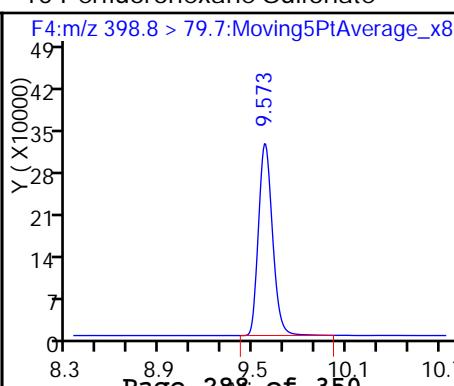
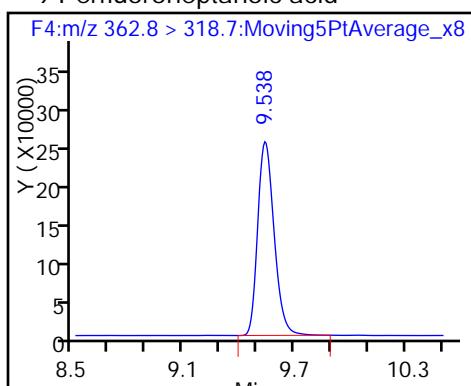
D 8 13C4-PFHxA



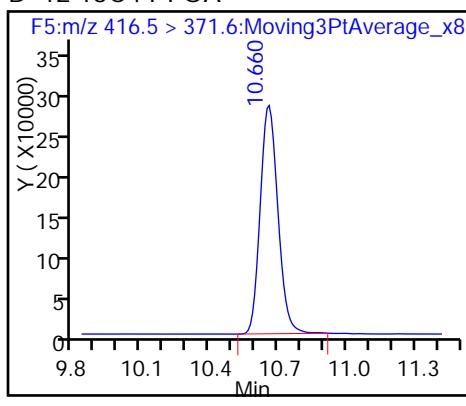
9 Perfluoroheptanoic acid

10 Perfluorohexane Sulfonate

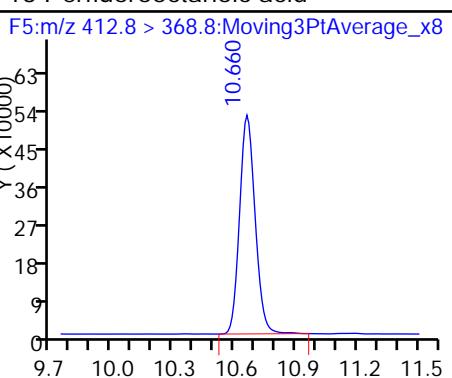
D 11 18O2 PFHxS



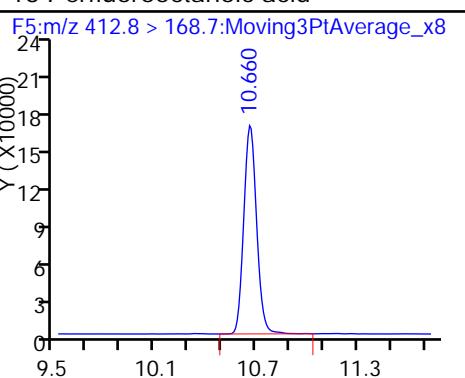
D 12 13C4 PFOA



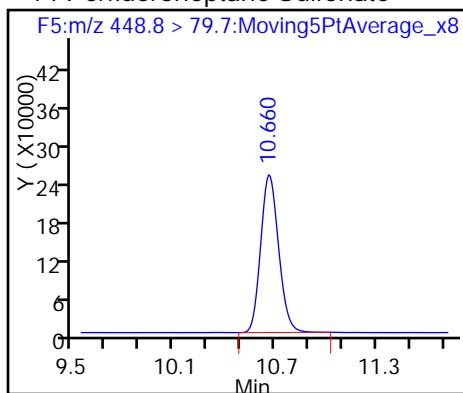
13 Perfluorooctanoic acid



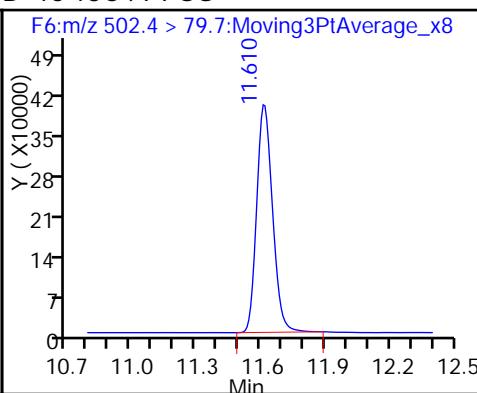
13 Perfluorooctanoic acid



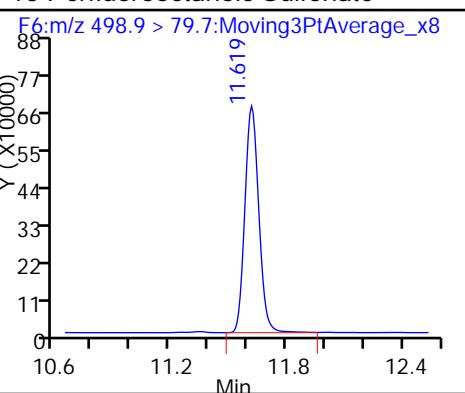
14 Perfluoroheptane Sulfonate



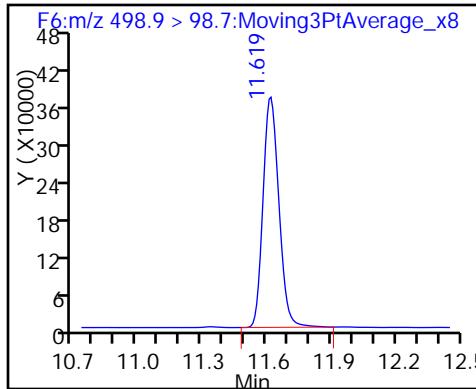
D 16 13C4 PFOS



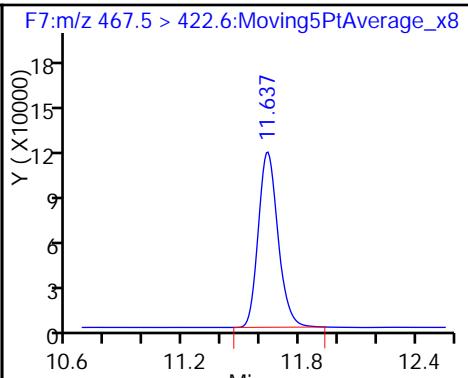
15 Perfluorooctanoic Sulfonate



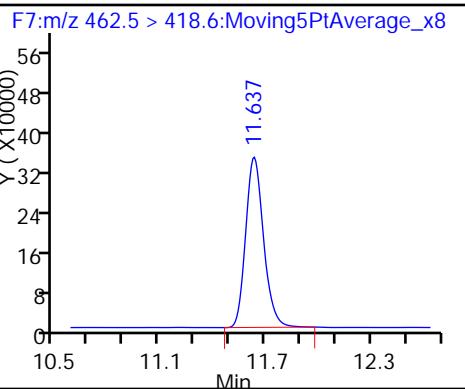
15 Perfluorooctanoic Sulfonate



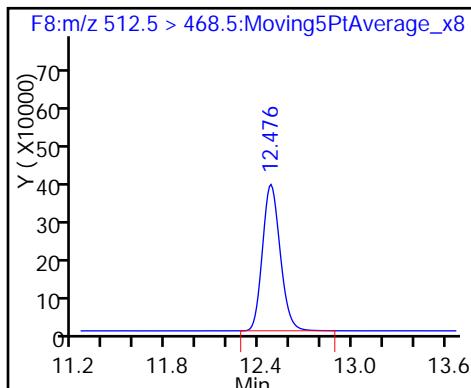
D 17 13C5 PFNA



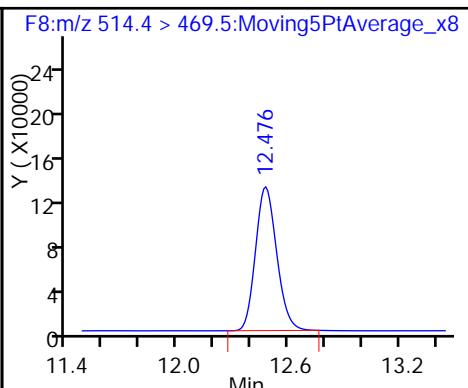
18 Perfluorononanoic acid



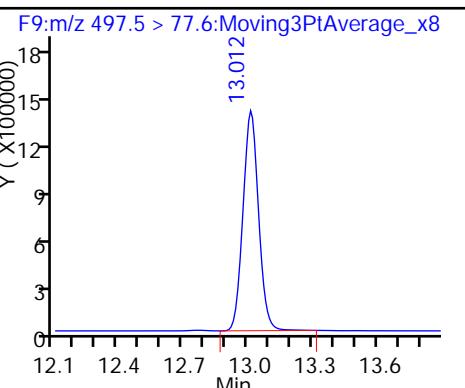
20 Perfluorodecanoic acid



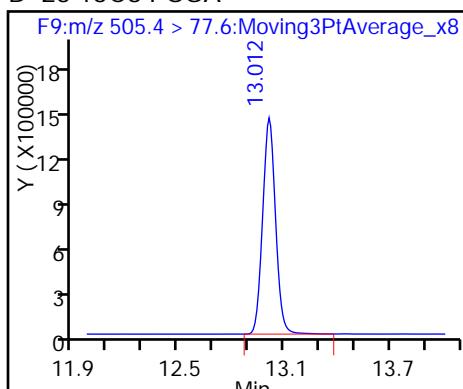
D 19 13C2 PFDA



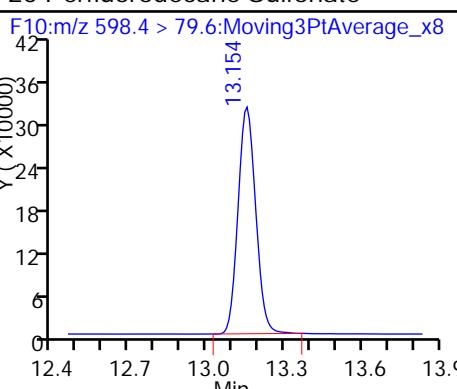
24 Perfluorooctane Sulfonamide



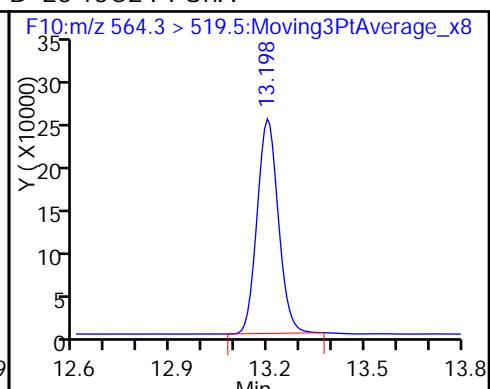
D 23 13C8 FOSA



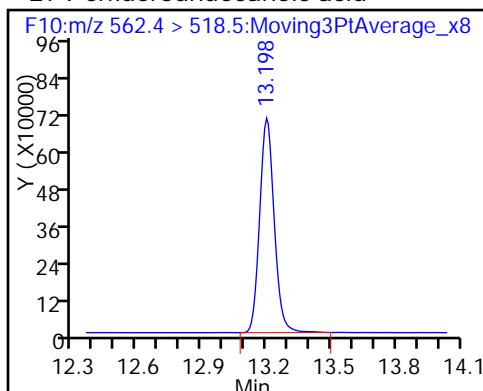
25 Perfluorodecane Sulfonate



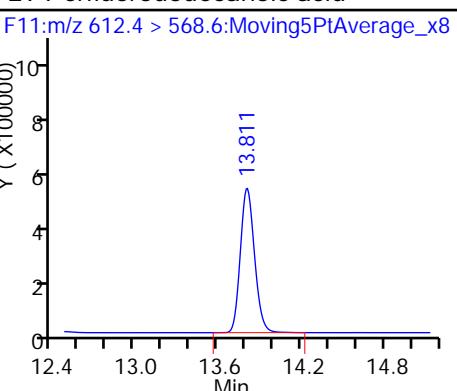
D 26 13C2 PFUnA



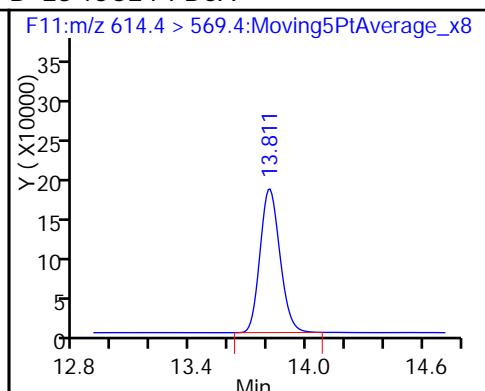
27 Perfluoroundecanoic acid



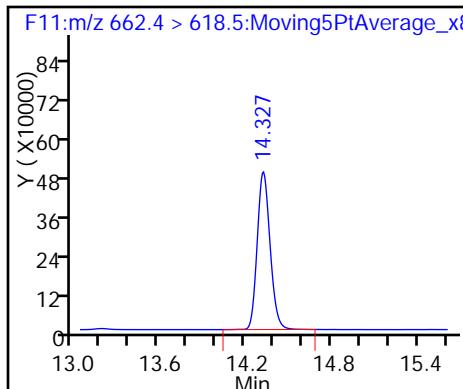
29 Perfluorododecanoic acid



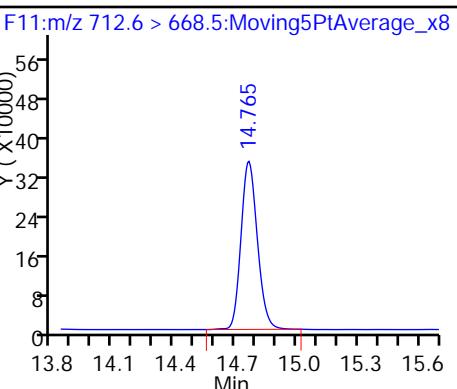
D 28 13C2 PFDaA



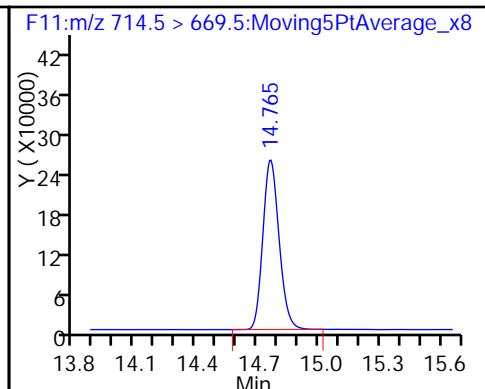
30 Perfluorotridecanoic acid



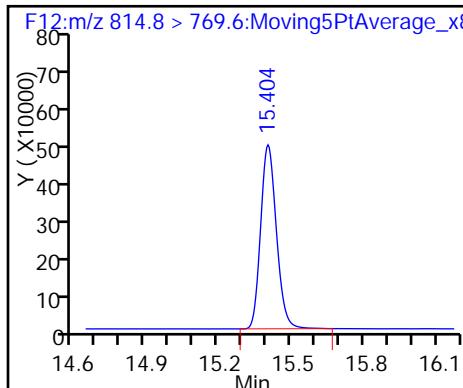
32 Perfluorotetradecanoic acid



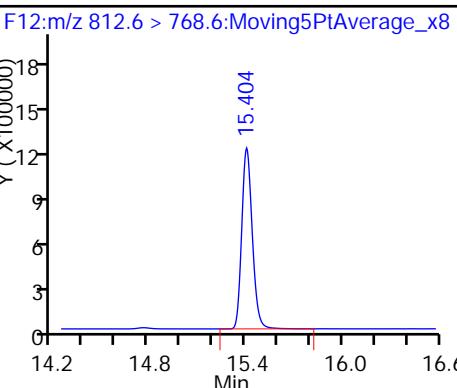
D 33 13C2-PFTeDA



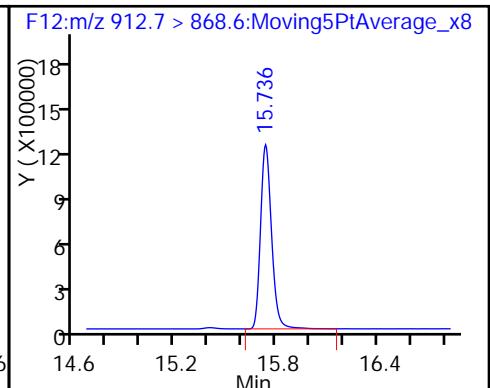
D 35 13C2-PFHxDA



34 Perfluorohexadecanoic acid



36 Perfluoroctadecanoic acid



Report Date: 05-Dec-2014 10:02:31

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_058.d

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Lab Sample ID: CCV 320-59779/45 Calibration Date: 12/04/2014 19:42  
Instrument ID: A4 Calib Start Date: 12/04/2014 03:44  
GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51  
Lab File ID: 03DEC14A4A\_081.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	3.794		32.3	20.0	61.4*	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	2.282		27.6	20.0	38.0	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.023		18.3	17.7	3.7	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.252		25.6	20.0	28.2	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.068		27.1	20.0	35.7	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	0.998		20.5	18.9	8.2	40.0
Perfluoro-1-heptanesulfonate (PFHps)	AveID	0.8647	0.9528		21.0	19.0	10.2	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.850		22.2	20.0	10.8	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.488		15.8	19.1	-17.6	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.142		19.8	20.0	-1.0	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.018		21.3	20.0	6.6	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.009		21.2	20.0	6.2	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7121		21.2	19.3	10.1	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.135		25.7	20.0	28.4	40.0
Perfluorododecanoic acid (PFDa)	AveID	0.9863	1.182		24.0	20.0	19.9	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9328		22.8	20.0	14.1	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5912		22.7	20.0	13.7	40.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.503		22.4	20.0	12.2	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.589		22.4	20.0	11.9	40.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_081.d  
 Lims ID: CCV L4  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 04-Dec-2014 19:42:11 ALS Bottle#: 13 Worklist Smp#: 45  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L4  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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D 1 113C4 PFBA										
216.7 > 171.5	5.928	5.934	-0.006		37702	8.81		44.1	164	
2 Perfluorobutyric acid										
212.7 > 168.6	5.928	5.935	-0.007	1.000	368051	32.3		161	1250	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.036	7.037	-0.001	1.000	221395	27.6		138	299	
D 3 113C5-PFPeA										
267.6 > 222.7	7.036	7.038	-0.002		97006	12.6		62.9	407	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.154	7.157	-0.003	1.000	809307	18.3		104	2339	
298.8 > 98.6	7.154	7.157	-0.003	1.000	515857		1.57(0.00-0.00)		1627	
D 6 113C2 PFHxA										
314.6 > 269.7	8.284	8.282	0.002		284568	14.9		74.7	990	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.284	8.284	0.0	1.000	356142	25.6		128	727	
22 PPoS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.360	8.360	0.0	0.876	678416	19.3		103	1522	
D 8 113C4-PFHxA										
366.6 > 321.6	9.510	9.506	0.004		292541	16.5		82.4	597	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.510	9.508	0.002	1.000	605064	27.1		136	1746	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	844262	20.5		108	2060	
D 11 118O2 PFHxA										
402.5 > 83.6	9.545	9.544	0.001		846358	18.8		99.5	2006	
D 12 113C4 PFOA										
416.5 > 371.6	10.632	10.628	0.004		1785879	52.8		106	2174	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.642	10.628	0.014	1.000	1321432	22.2		111	1760	
412.8 > 168.7	10.642	10.628	0.014	1.000	399862		3.30(0.00-0.00)		592	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.642	10.635	0.007	1.000	862776	21.0		110	1352	
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.595	0.007		2273365	50.4		105	5010	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.602	11.595	0.007	1.000	1352981	15.8		82.4	1845	
498.9 > 98.7	11.602	11.595	0.007	1.000	777457		1.74(0.00-0.00)		1773	
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.614	0.006		882241	22.8		114	2428	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	1007484	19.8		99.0	1627	
21 PFNS (Perflouro-1-nananesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	709832	20.3		106	1096	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	1152984	21.3		107	1270	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		1132812	22.0		110	2279	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.004	12.998	0.006	1.000	2996256	21.2		106	2169	
D 23 13C8 FOSA										
505.4 > 77.6	13.004	12.998	0.006		7425537	51.9		104	4287	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	652922	21.2		110	985	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1234387	19.1		95.3	2050	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	1400526	25.7		128	1903	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	1627337	24.0		120	1143	
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1376207	21.1		105	1822	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.273	14.272	0.001	1.000	732714	23.4		121	1367	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.327	14.327	0.0	1.000	1283705	22.8		114	847	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	813561	22.7		114	728	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		1526321	21.3		107	2478	
D 35 13C2-PFHxDa										
814.8 > 769.6	15.404	15.415	-0.011		2277015	21.8		109	2792	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.404	15.415	-0.011	1.000	2068581	22.4		112	1315	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.736	15.752	-0.016	1.000	2186469	22.4		112	1210	12/12/2014

Report Date: 05-Dec-2014 10:03:22

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

LCPFC-L4\_00005

Amount Added: 1.00

Units: mL

Report Date: 05-Dec-2014 10:03:22

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_081.d

Injection Date: 04-Dec-2014 19:42:11

Instrument ID: A4

Lims ID: CCV L4

Client ID:

Operator ID: JRB

ALS Bottle#: 13 Worklist Smp#: 45

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

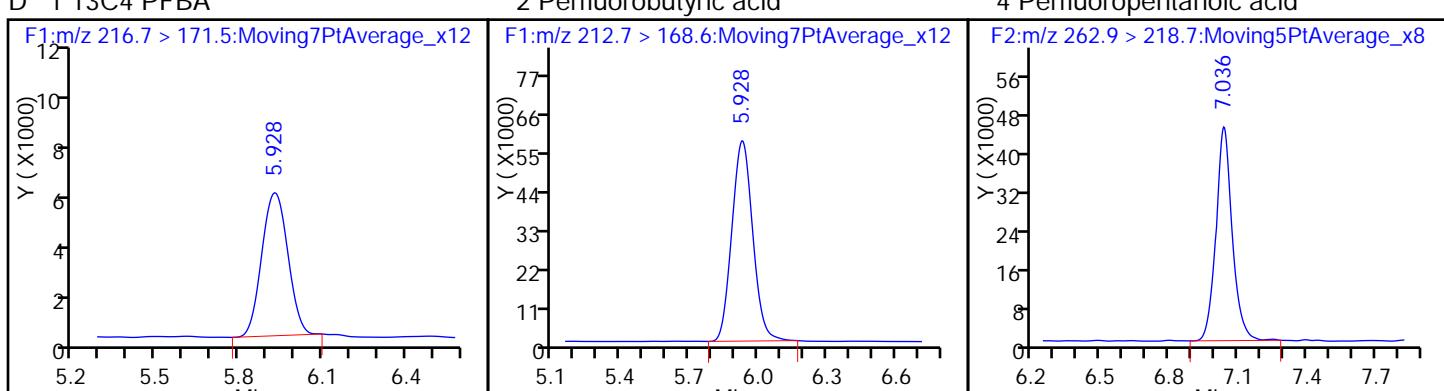
Method: PFAC\_A4

Limit Group: LC PFC ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid

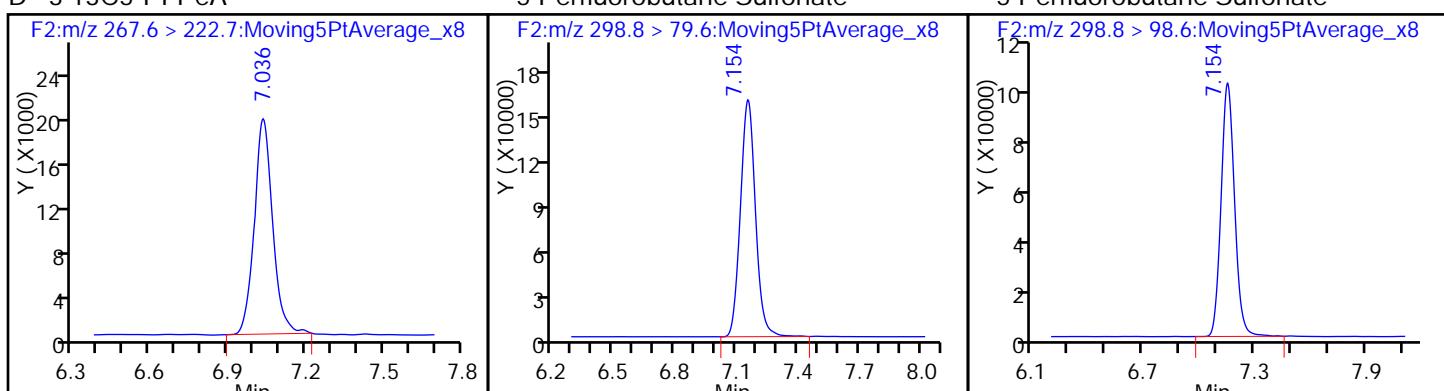
4 Perfluoropentanoic acid



D 3 13C5-PFPeA

5 Perfluorobutane Sulfonate

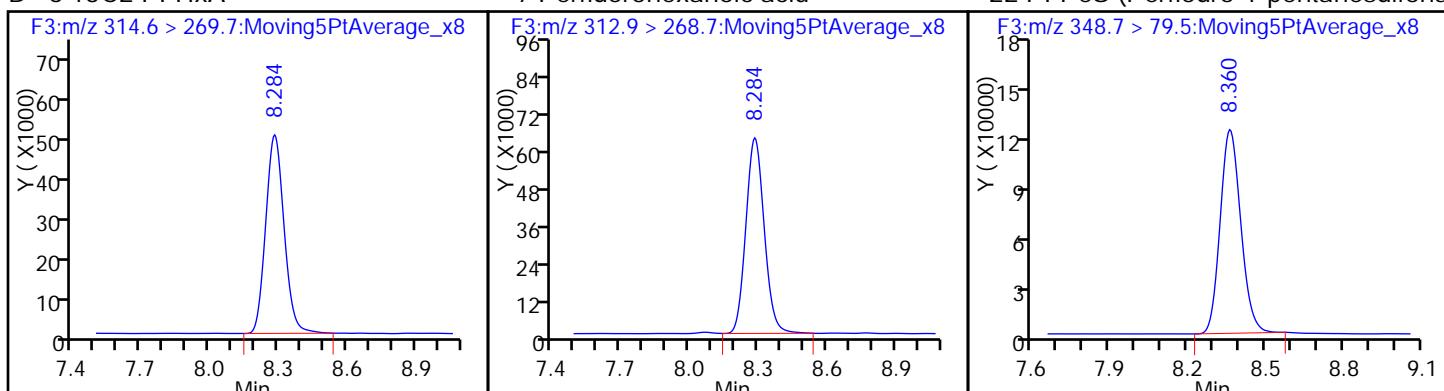
5 Perfluorobutane Sulfonate



D 6 13C2 PFHxA

7 Perfluorohexanoic acid

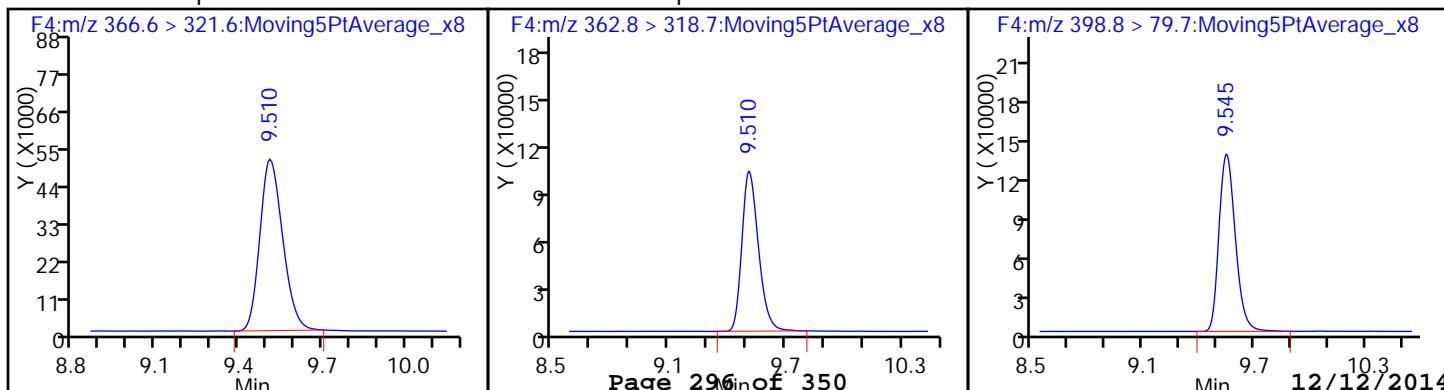
22 PFPeS (Perflouro-1-pentanesulfonat)



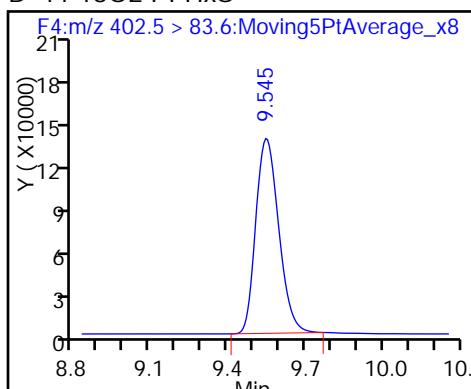
D 8 13C4-PFHxA

9 Perfluoroheptanoic acid

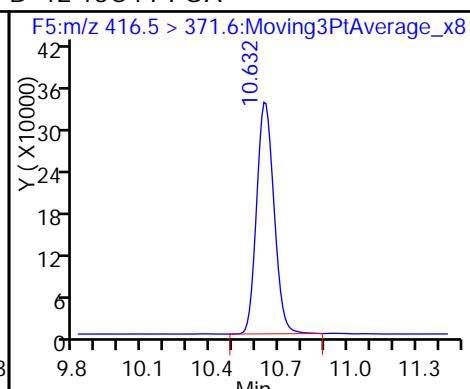
10 Perfluorohexane Sulfonate



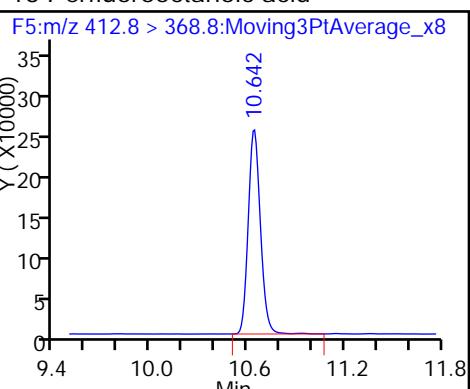
D 11 18O2 PFHxS



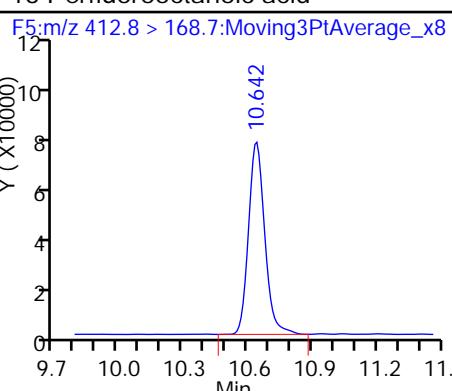
D 12 13C4 PFOA



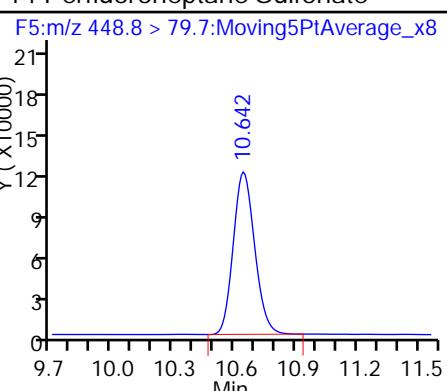
13 Perfluorooctanoic acid



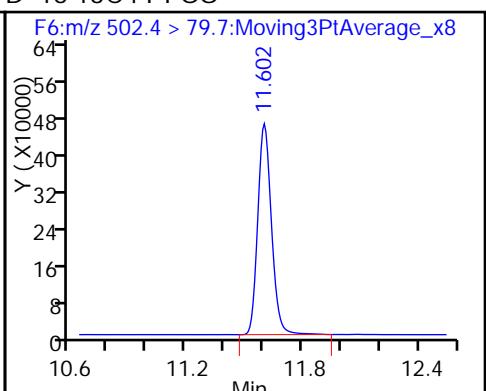
13 Perfluorooctanoic acid



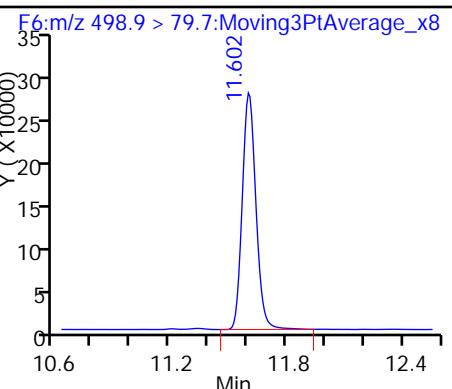
14 Perfluoroheptane Sulfonate



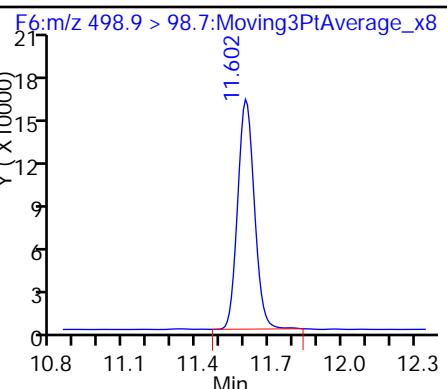
D 16 13C4 PFOS



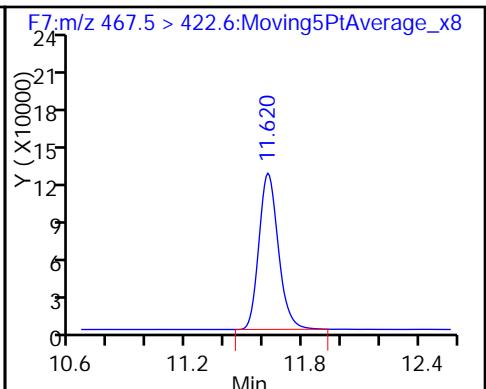
15 Perfluorooctanoic Sulfonate



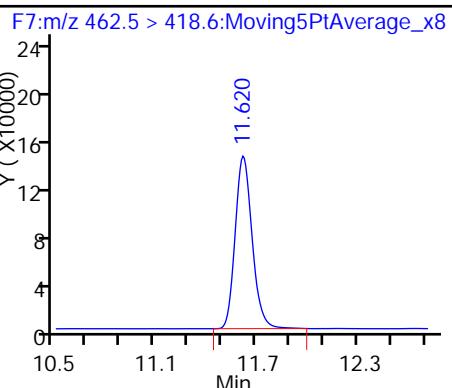
15 Perfluorooctanoic Sulfonate



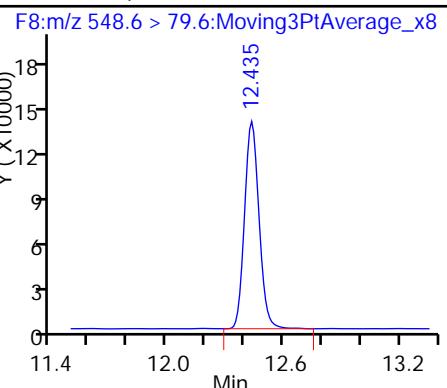
D 17 13C5 PFNA



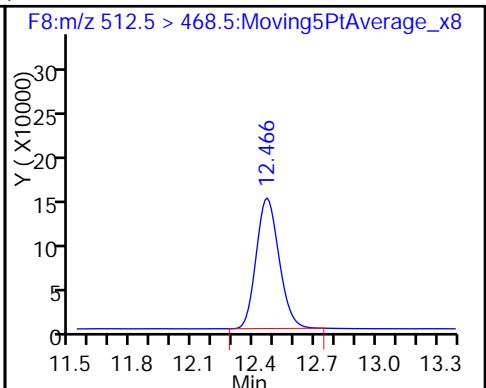
18 Perfluorononanoic acid



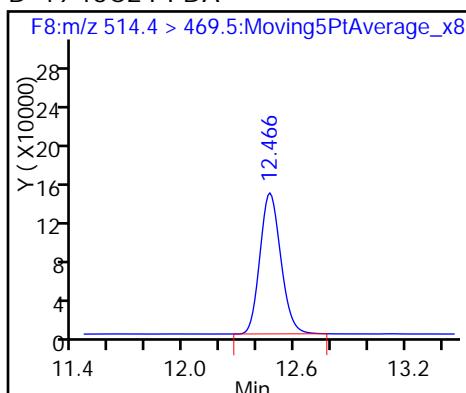
21 PFNS (Perflouro-1-nananesulfonate)



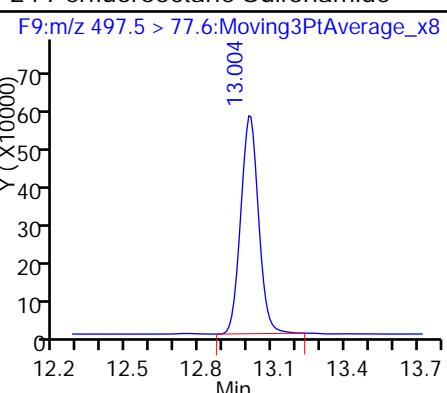
20 Perfluorodecanoic acid



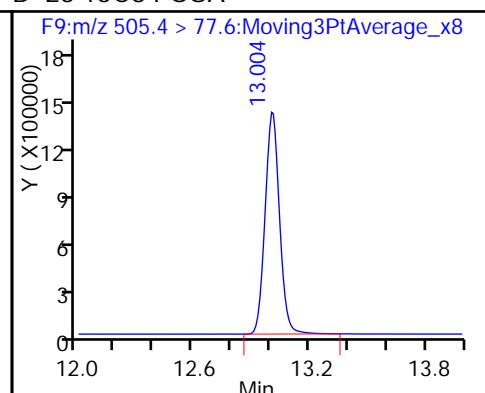
D 19 13C2 PFDA



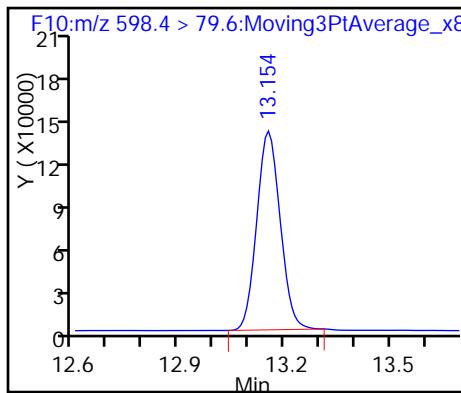
24 Perfluorooctane Sulfonamide



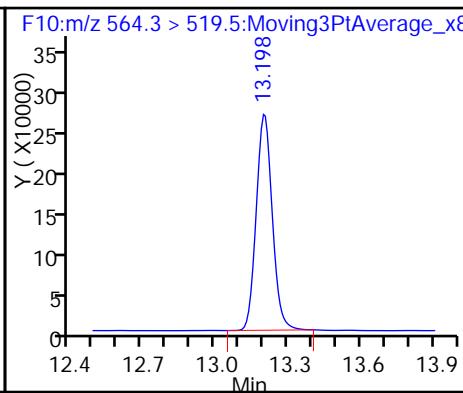
D 23 13C8 FOSA



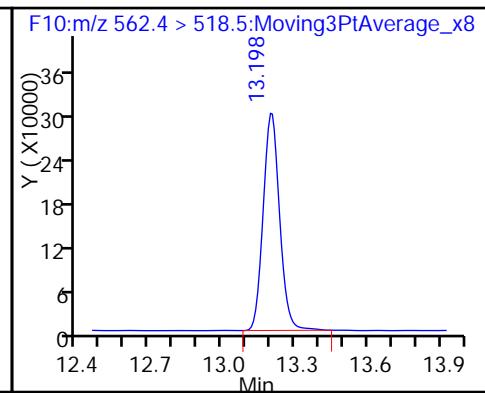
25 Perfluorodecane Sulfonate



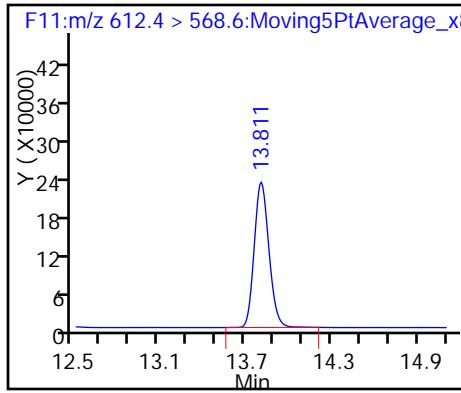
D 26 13C2 PFUna



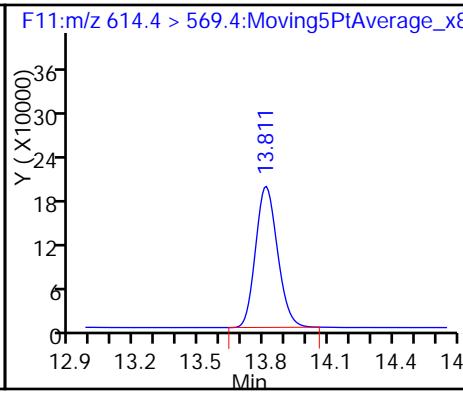
27 Perfluoroundecanoic acid



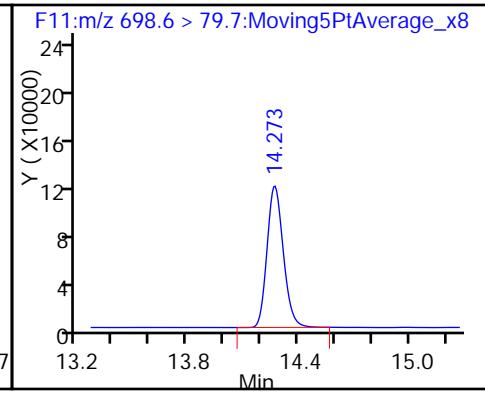
29 Perfluorododecanoic acid



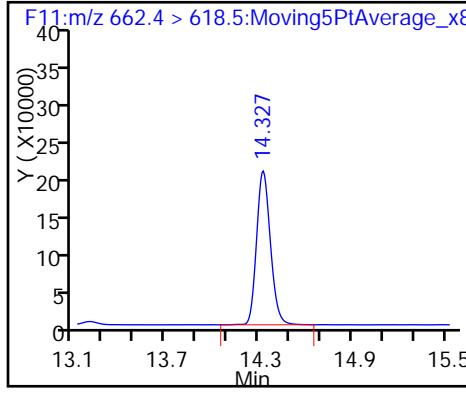
D 28 13C2 PFDoA



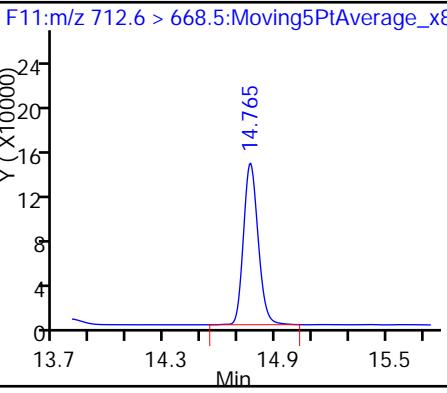
31 PFDoS (Perflouro-1-dodecanesulfona



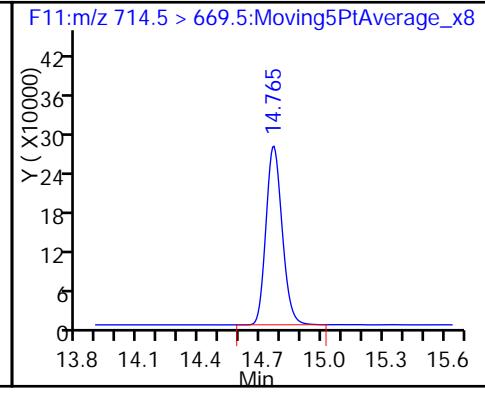
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



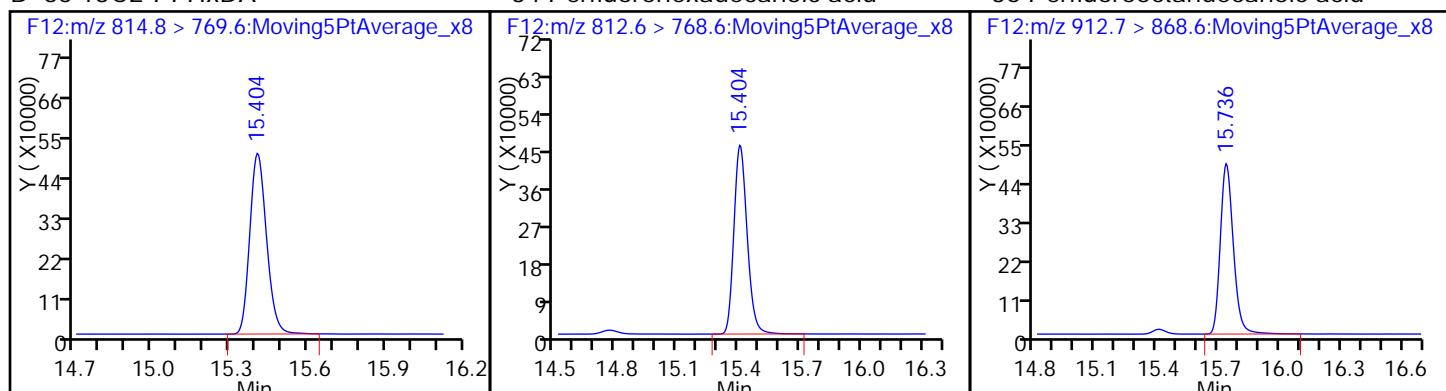
D 33 13C2-PFTeDA



D 35 13C2-PFHxDa

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Lab Sample ID: CCV 320-59779/57 Calibration Date: 12/04/2014 23:56  
Instrument ID: A4 Calib Start Date: 12/04/2014 03:44  
GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51  
Lab File ID: 03DEC14A4A\_093.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	3.249		69.1	50.0	38.2	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	2.193		66.3	50.0	32.6	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.639		73.5	44.2	66.2*	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.185		60.7	50.0	21.4	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	1.987		65.1	50.0	30.3	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	1.045		53.6	47.3	13.3	40.0
Perfluoro-1-heptanesulfonate (PFHps)	AveID	0.8647	0.9556		52.6	47.6	10.5	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.657		49.6	50.0	-0.7	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.512		40.0	47.8	-16.2	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.219		52.8	50.0	5.7	40.0
Perfluorodecanoic acid (PFDa)	AveID	0.9544	1.069		56.0	50.0	12.0	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.050		55.3	50.0	10.5	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.6689		49.8	48.2	3.4	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.054		59.6	50.0	19.3	40.0
Perfluorododecanoic acid (PFDa)	AveID	0.9863	1.179		59.8	50.0	19.5	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9404		57.5	50.0	15.0	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5983		57.5	50.0	15.1	40.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.598		59.6	50.0	19.3	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.784		62.8	50.0	25.7	40.0

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_093.d  
 Lims ID: CCV L5  
 Client ID:  
 Sample Type: CCV  
 Inject. Date: 04-Dec-2014 23:56:45 ALS Bottle#: 14 Worklist Smp#: 57  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: CCV L5  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Sublist: chrom-PFAC\_A4\*sub5  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 11:43:09 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

First Level Reviewer: barnettj Date: 05-Dec-2014 11:43:09

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 113C4 PFBA										
216.7 > 171.5	5.934	5.934	0.0		43898	10.3		51.3	214	
2 Perfluorobutyric acid										
212.7 > 168.6	5.934	5.935	-0.001	1.000	918655	69.1		138	2962	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.040	7.037	0.003	1.000	619975	66.3		133	784	
D 3 113C5-PFPeA										
267.6 > 222.7	7.040	7.038	0.002		113092	14.7		73.3	417	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.155	7.157	-0.002	1.000	2883515	73.5		166	7024	
298.8 > 98.6	7.155	7.157	-0.002	1.000	1831083	1.57(0.00-0.00)		5544		
D 6 113C2 PFHxA										
314.6 > 269.7	8.286	8.283	0.003		308162	16.2		80.8	907	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.291	8.285	0.006	1.000	912713	60.7		121	1752	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.367	8.361	0.006	0.876	1689870	54.0		115	4635	
D 8 113C4-PFHpA										
366.6 > 321.6	9.511	9.507	0.004		274996	15.5		77.5	698	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.518	9.509	0.009	1.000	1365733	65.1		130	3237	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.552	9.544	0.008	1.000	1966484	53.6		113	3383	
D 11 18O2 PFHxS										
402.5 > 83.6	9.552	9.545	0.007		752958	16.8		88.5	2077	
D 12 113C4 PFOA										
416.5 > 371.6	10.642	10.630	0.012		1624638	48.0		96.0	2184	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
13 Perfluorooctanoic acid										
412.8 > 368.8	10.642	10.630	0.012	1.000	2692481	49.6		99.3	3143	
412.8 > 168.7	10.642	10.630	0.012	1.000	885834		3.04(0.00-0.00)		1573	
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.642	10.636	0.006	1.000	1989674	52.6		111	4178	
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.596	0.006		2090959	46.4		97.0	3565	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.602	11.596	0.006	1.000	3160526	40.0		83.8	2974	
498.9 > 98.7	11.602	11.596	0.006	1.000	1809472		1.75(0.00-0.00)		2399	
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.615	0.005		760335	19.7		98.4	1997	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.617	0.003	1.000	2316836	52.8		106	3784	
21 PFNS (Perflouro-1-nananesulfonate)										
548.6 > 79.6	12.435	12.430	0.005	1.000	1676333	52.1		109	3571	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.462	0.004	1.000	2455672	56.0		112	3689	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		918867	17.8		89.2	1282	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.003	12.999	0.004	1.000	7228791	55.3		111	2906	
D 23 13C8 FOSA										
505.4 > 77.6	13.003	12.999	0.004		6884880	48.1		96.2	2814	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.145	13.151	-0.006	1.000	1410239	49.8		103	2332	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1048630	16.2		81.0	1912	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	2764337	59.6		119	3977	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.802	13.807	-0.005	1.000	3664061	59.8		120	2495	
D 28 13C2 PFDoA										
614.4 > 569.4	13.802	13.808	-0.006		1243065	19.0		95.1	1371	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.265	14.271	-0.006	1.000	1614752	56.1		116	2339	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.319	14.326	-0.007	1.000	2922334	57.5		115	2022	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.758	14.765	-0.007	1.000	1859210	57.5		115	1394	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.758	14.765	-0.007		1401500	19.6		98.0	2419	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.404	15.413	-0.009		2197476	21.1		105	2391	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.404	15.413	-0.009	1.000	4965341	59.6		119	2268	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.731	15.749	-0.018	1.000	Page 5545303 of 350	62.8		126	1916	12/12/2014

Report Date: 05-Dec-2014 11:43:09

Chrom Revision: 2.2 06-Nov-2014 14:50:32

**Reagents:**

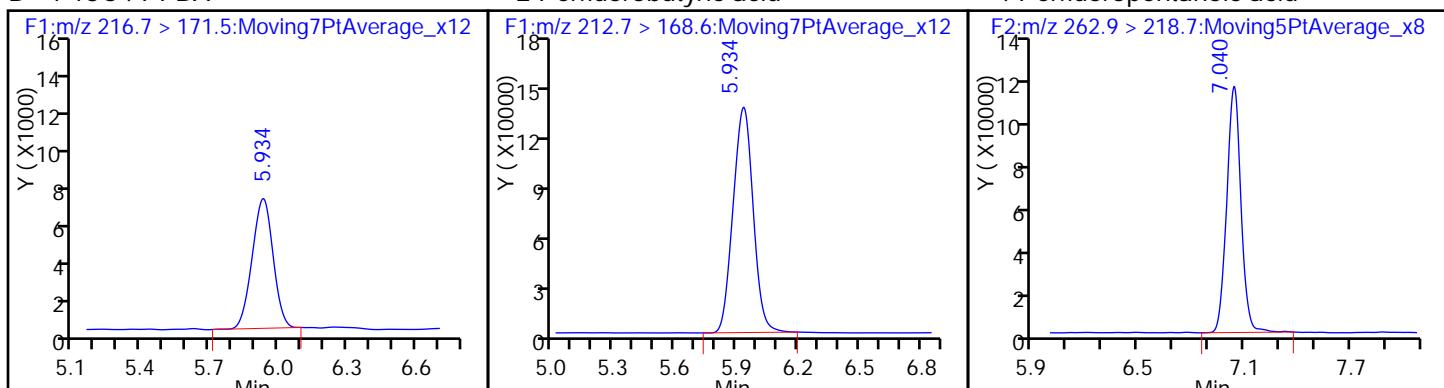
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Amount Added: 1.00

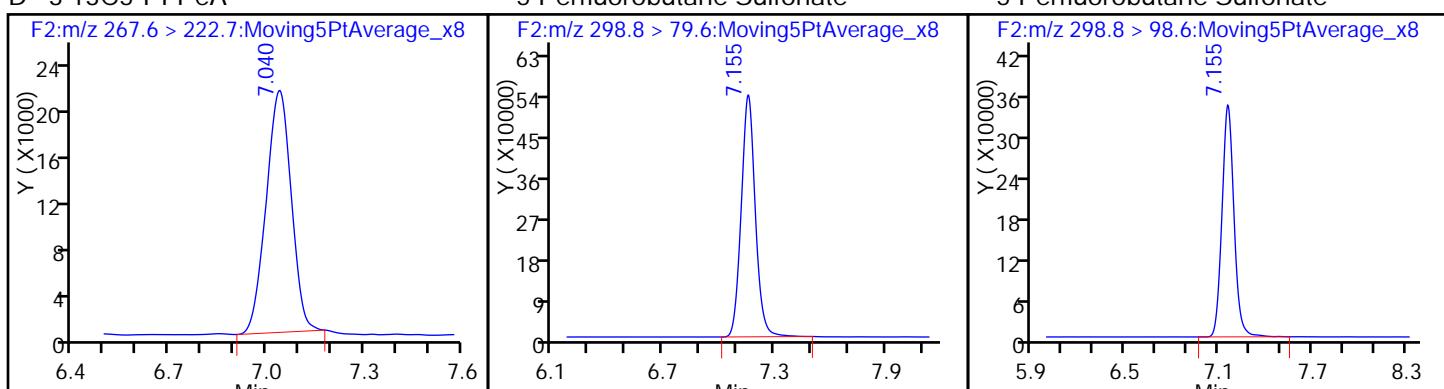
Units: mL

TestAmerica Sacramento  
 Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_093.d  
 Injection Date: 04-Dec-2014 23:56:45 Instrument ID: A4  
 Lims ID: CCV L5  
 Client ID:  
 Operator ID: JRB ALS Bottle#: 14 Worklist Smp#: 57  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Method: PFAC\_A4 Limit Group: LC PFC ICAL

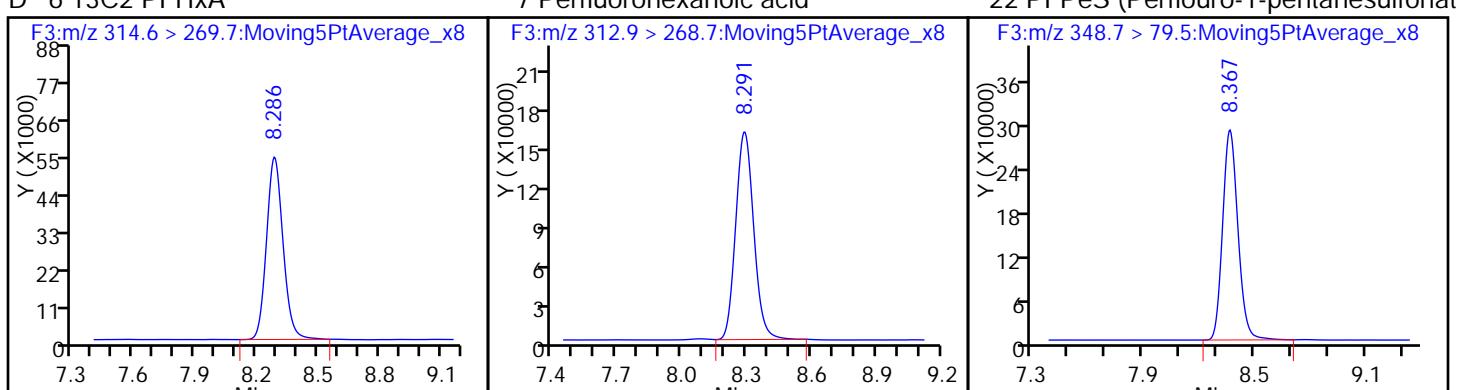
## D 1 13C4 PFBA



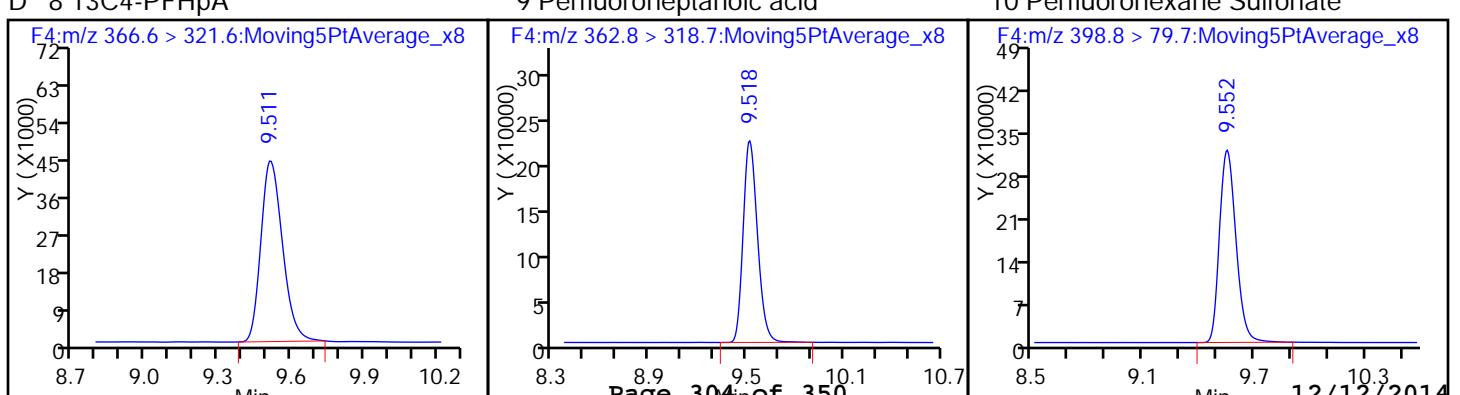
## D 3 13C5-PFPeA



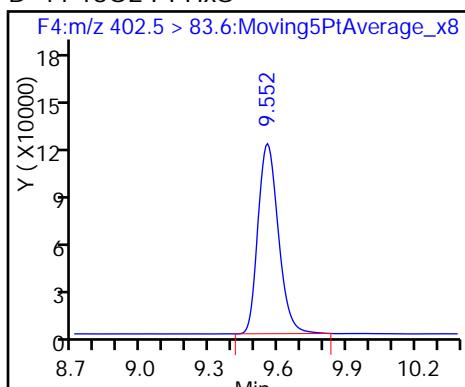
## D 6 13C2 PFHxA



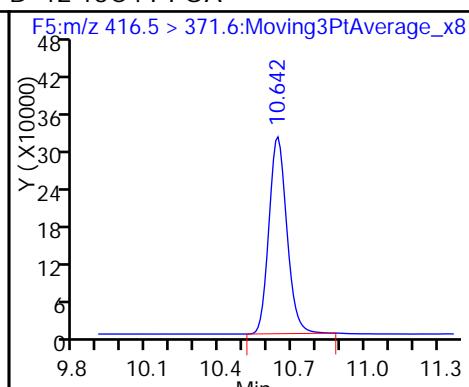
## D 8 13C4-PFHxA



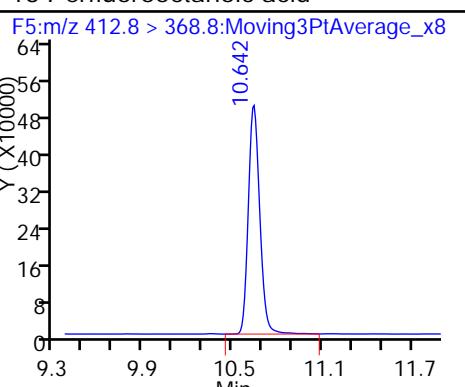
D 11 18O2 PFHxS



D 12 13C4 PFOA



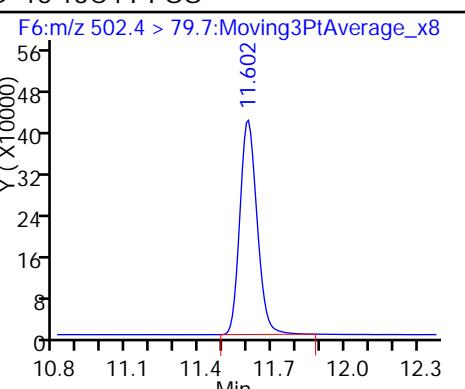
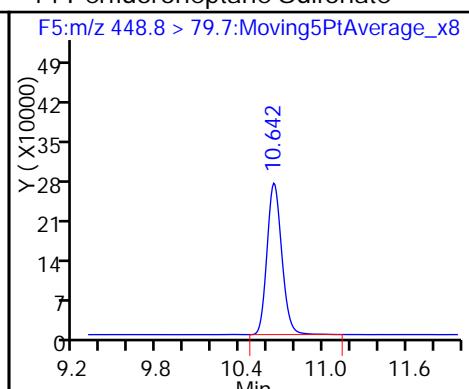
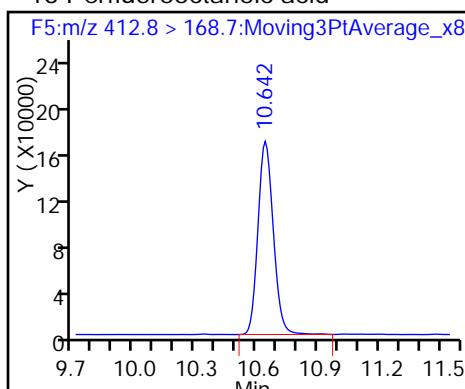
13 Perfluorooctanoic acid



13 Perfluorooctanoic acid

14 Perfluoroheptane Sulfonate

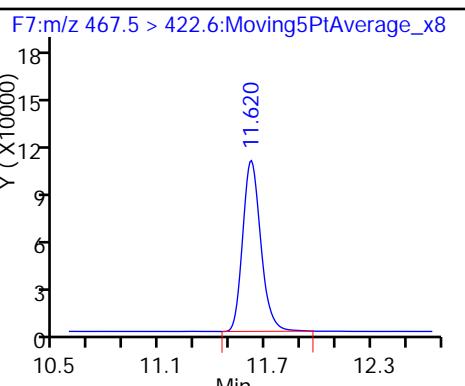
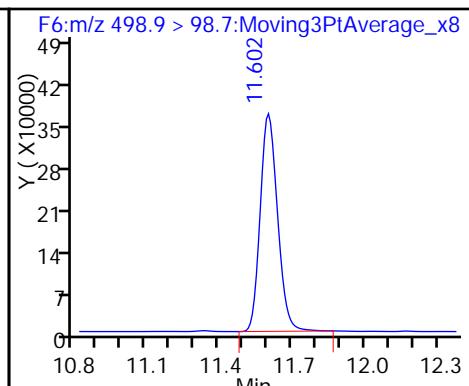
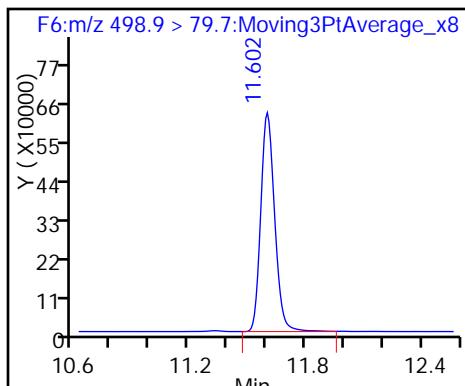
D 16 13C4 PFOS



15 Perfluorooctanoic Sulfonate

15 Perfluorooctanoic Sulfonate

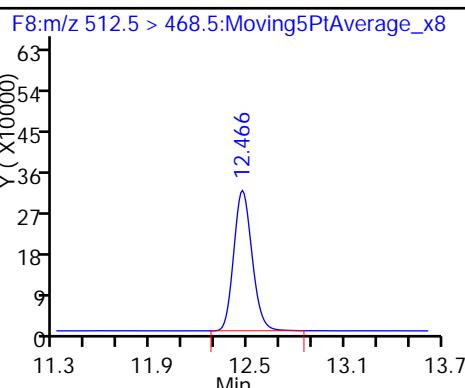
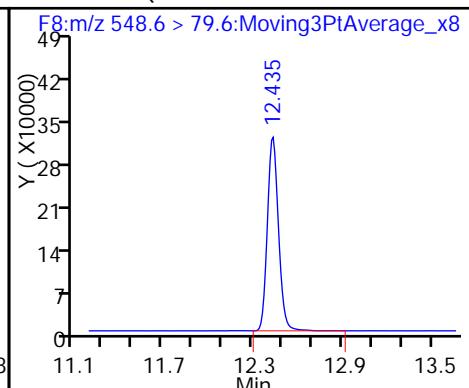
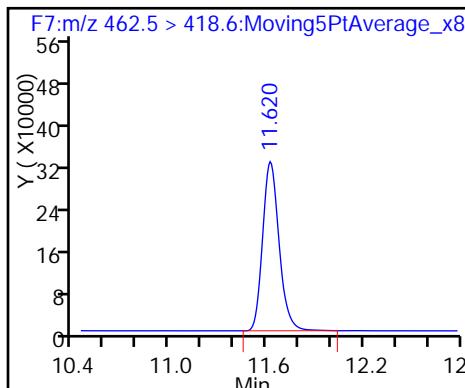
D 17 13C5 PFNA



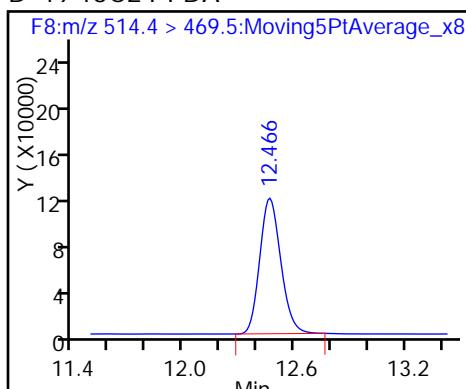
18 Perfluorononanoic acid

21 PFNS (Perflouro-1-nananesulfonate)

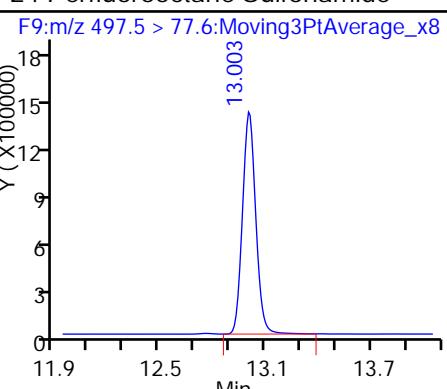
20 Perfluorodecanoic acid



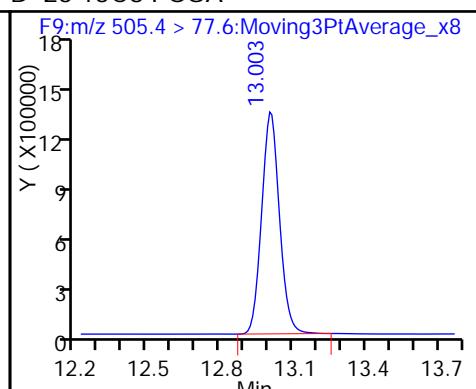
D 19 13C2 PFDA



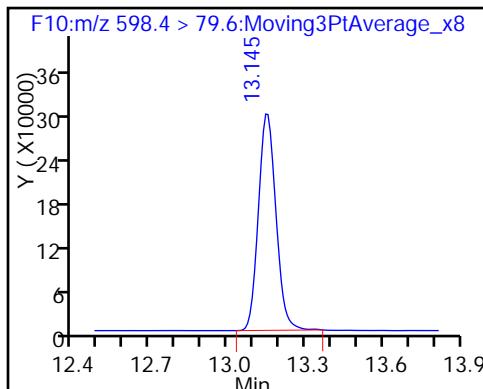
24 Perfluorooctane Sulfonamide



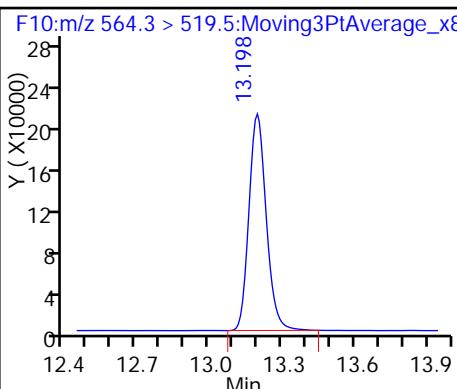
D 23 13C8 FOSA



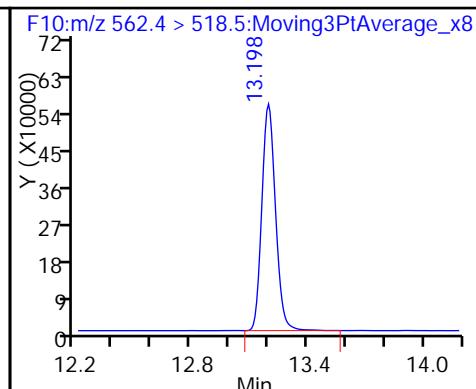
25 Perfluorodecane Sulfonate



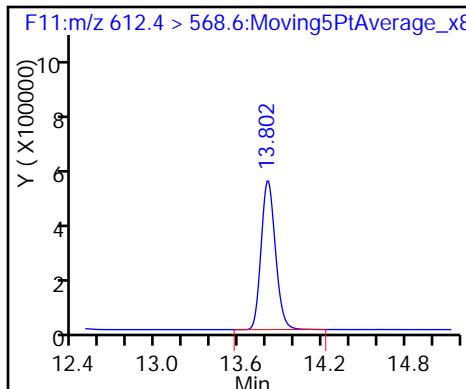
D 26 13C2 PFUna



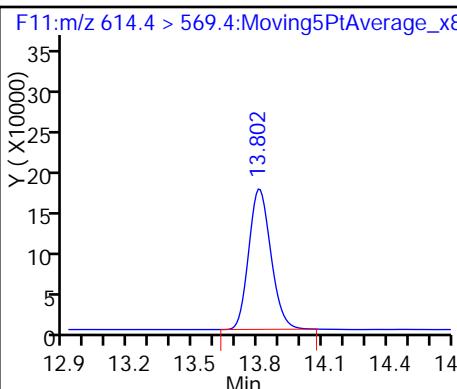
27 Perfluoroundecanoic acid



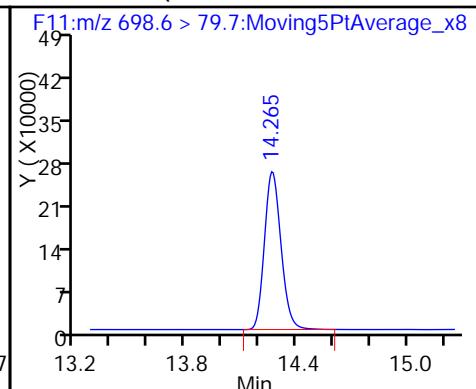
29 Perfluorododecanoic acid



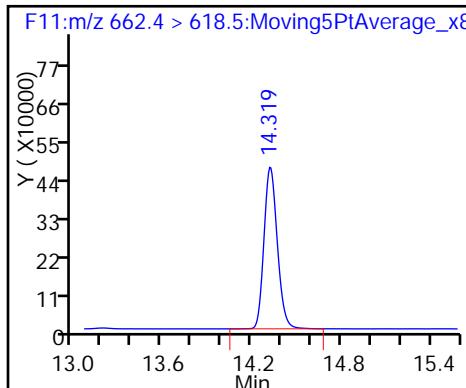
D 28 13C2 PFDoA



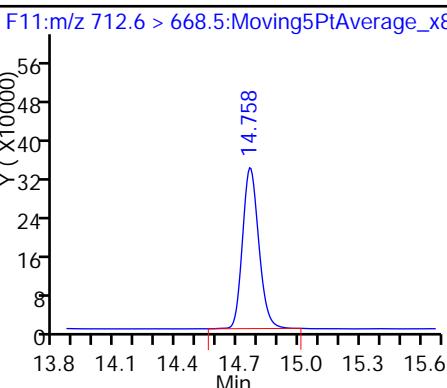
31 PFDoS (Perflouro-1-dodecanesulfona



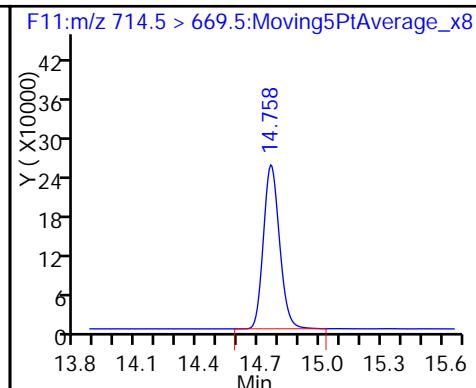
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



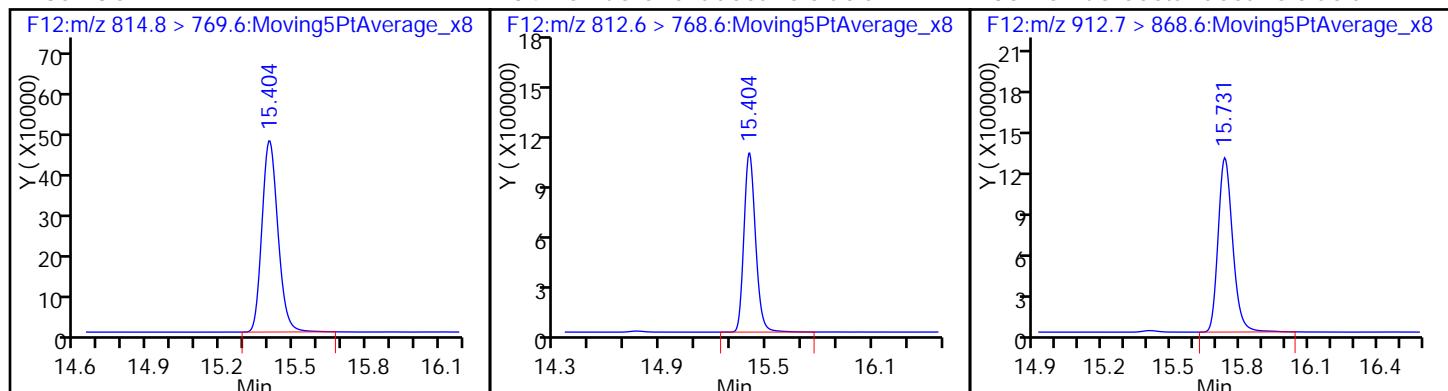
D 33 13C2-PFTeDA



D 35 13C2-PFHxDa

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 320-59712/1-A  
Matrix: Water Lab File ID: 03DEC14A4A\_082.d  
Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
Extraction Method: 3535 Date Extracted: 12/03/2014 13:38  
Sample wt/vol: 500.00 (mL) Date Analyzed: 12/04/2014 20:03  
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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	0.809	J	2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	111		25-150
STL00990	13C4 PFOA	109		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_082.d  
 Lims ID: MB 320-59712/1-A  
 Client ID:  
 Sample Type: MB  
 Inject. Date: 04-Dec-2014 20:03:26 ALS Bottle#: 22 Worklist Smp#: 46  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: mb 320-59712/1-a 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICAL File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
D 1 113C4 PFBA										
216.7 > 171.5	5.928	5.934	-0.006		27219	6.36		31.8	83.8	
D 3 13C5-PFPeA										
267.6 > 222.7	7.039	7.038	0.001		89401	11.6		57.9	377	
D 6 13C2 PFHxA										
314.6 > 269.7	8.284	8.282	0.002		261844	13.7		68.7	874	
D 8 13C4-PFHxA										
366.6 > 321.6	9.510	9.506	0.004		290618	16.4		81.9	882	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	2915	0.0737				10.4
D 11 18O2 PFHxS										
402.5 > 83.6	9.545	9.544	0.001		810985	18.0		95.4	2247	
D 12 13C4 PFOA										
416.5 > 371.6	10.642	10.628	0.014		1843821	54.5		109	3092	
13 Perfluorooctanoic acid										
412.8 > 368.8	10.642	10.628	0.014	1.000	24900	0.4044				26.5
412.8 > 168.7	10.632	10.628	0.004	0.999	3853		6.46(0.00-0.00)			6.6
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.595	0.007		2393416	53.1		111	4947	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.610	11.595	0.015	1.000	20748	0.2296				39.7
498.9 > 98.7	11.602	11.595	0.007	0.999	10394		2.00(0.00-0.00)			19.6
D 17 13C5 PFNA										
467.5 > 422.6	11.620	11.614	0.006		905590	23.4		117	2613	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		1121599	21.8		109	2647	
D 23 13C8 FOSA										
505.4 > 77.6	13.004	12.998	0.006		4931589	34.5		68.9	3402	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1117477	17.3		86.3	2721	

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.189	13.198	-0.009	1.000	2362	0.0478				6.5
D 28 13C2 PFDoA										
614.4 > 569.4	13.811	13.809	0.002		1008439	15.4		77.1	1856	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.765	14.766	-0.001		749399	10.5		52.4	1763	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.404	15.415	-0.011		1255809	12.0		60.2	1924	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.410	15.415	-0.005	1.000	6386	0.0945				9.2
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.741	15.752	-0.011	1.000	6696	0.0935				9.5

Report Date: 05-Dec-2014 10:03:23

Chrom Revision: 2.2 06-Nov-2014 14:50:32

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_082.d

Injection Date: 04-Dec-2014 20:03:26

Instrument ID: A4

Lims ID: MB 320-59712/1-A

Client ID:

Operator ID: JRB

ALS Bottle#: 22 Worklist Smp#: 46

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

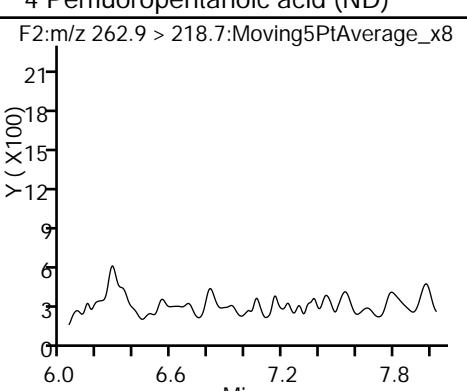
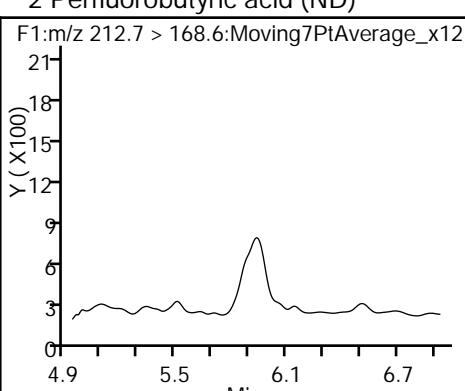
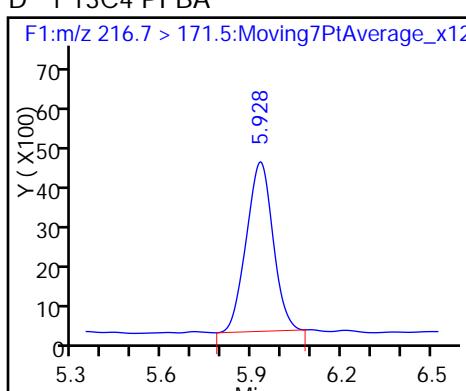
Method: PFAC\_A4

Limit Group: LC PFC ICAL

D 1 13C4 PFBA

2 Perfluorobutyric acid (ND)

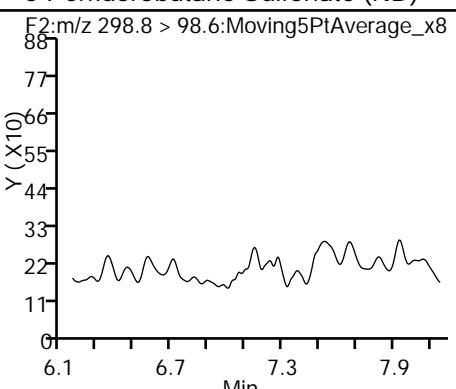
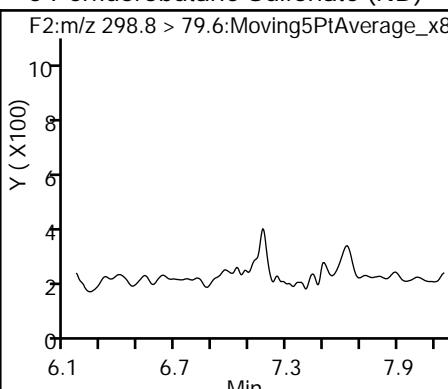
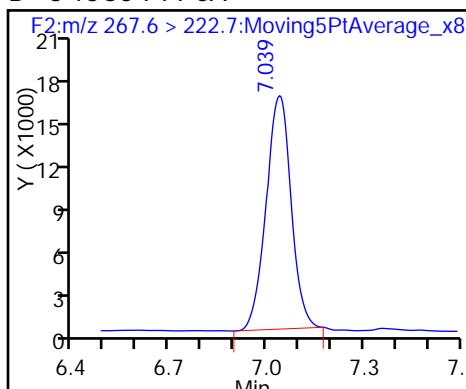
4 Perfluoropentanoic acid (ND)



D 3 13C5-PFPeA

5 Perfluorobutane Sulfonate (ND)

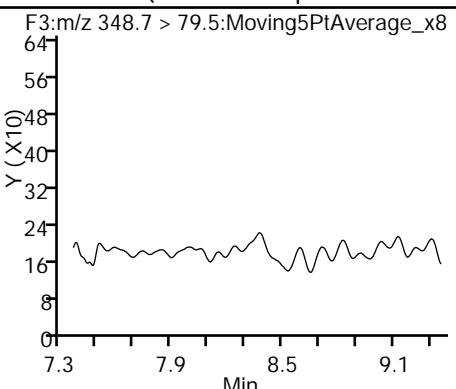
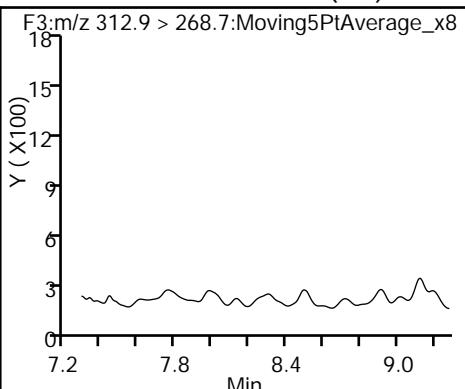
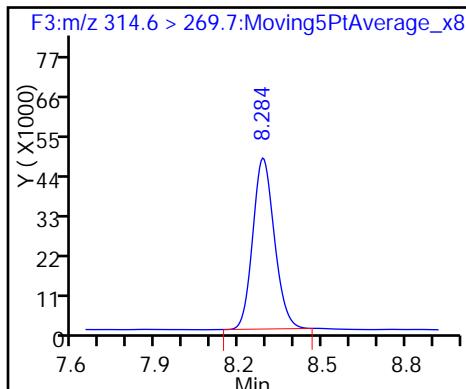
5 Perfluorobutane Sulfonate (ND)



D 6 13C2 PFHxA

7 Perfluorohexanoic acid (ND)

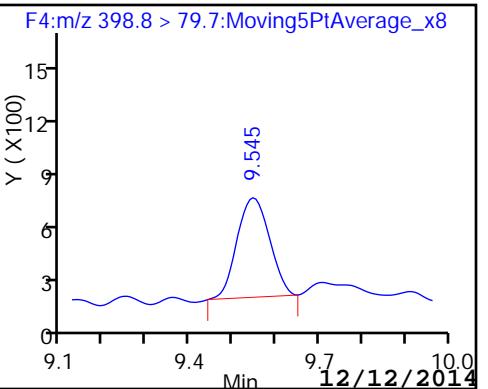
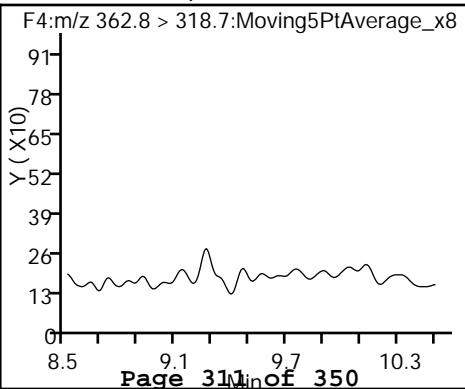
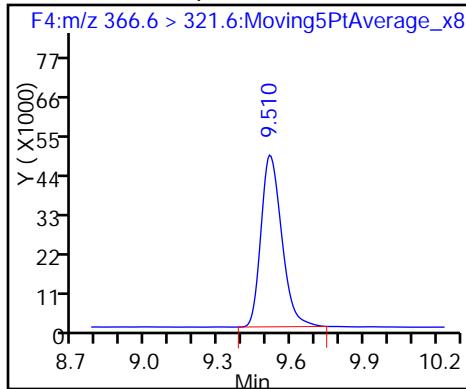
22 PFPeS (Perflouro-1-pentanesulfonat (ND))



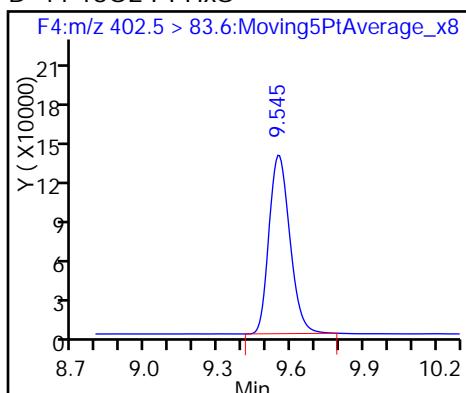
D 8 13C4-PFHxA

9 Perfluoroheptanoic acid (ND)

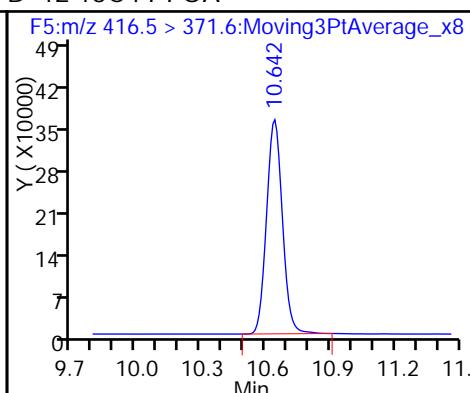
10 Perfluorohexane Sulfonate



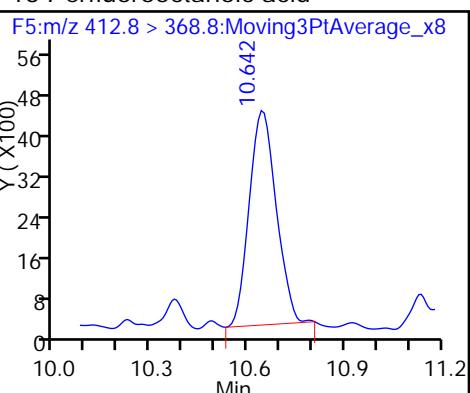
D 11 18O2 PFHxS



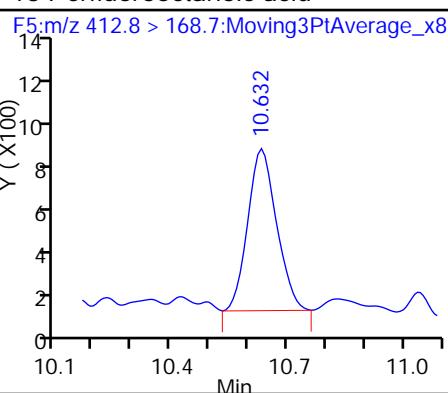
D 12 13C4 PFOA



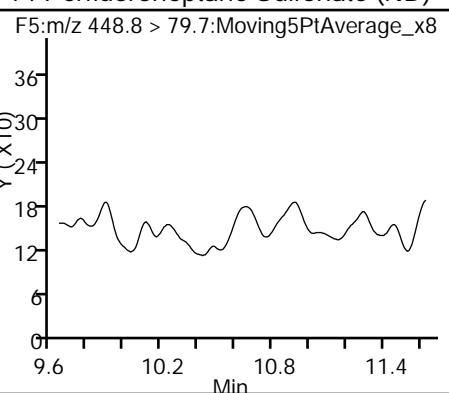
13 Perfluorooctanoic acid



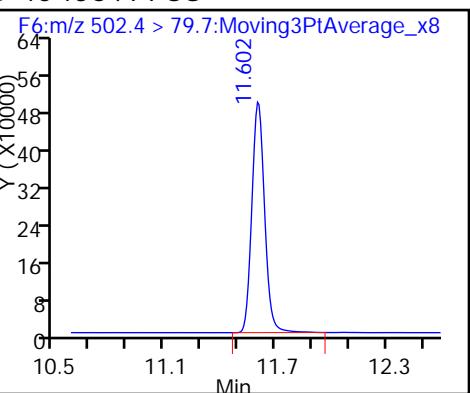
13 Perfluorooctanoic acid



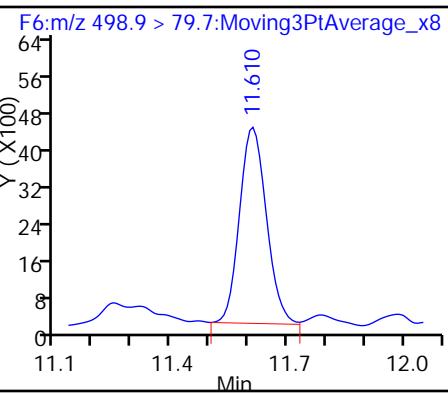
14 Perfluoroheptane Sulfonate (ND)



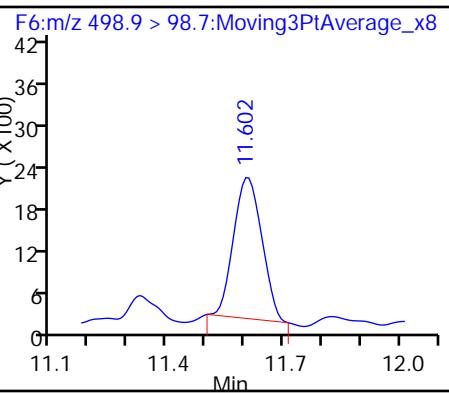
D 16 13C4 PFOS



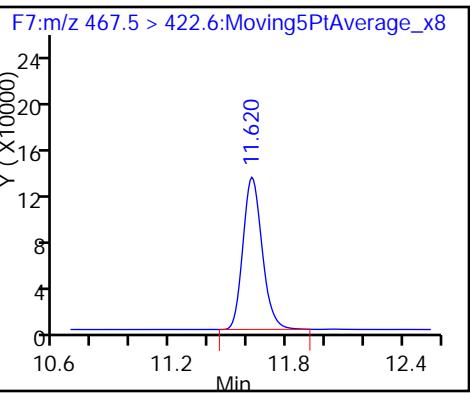
15 Perfluorooctanoic Sulfonate



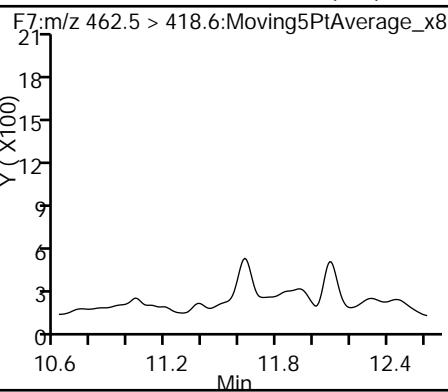
15 Perfluorooctanoic Sulfonate



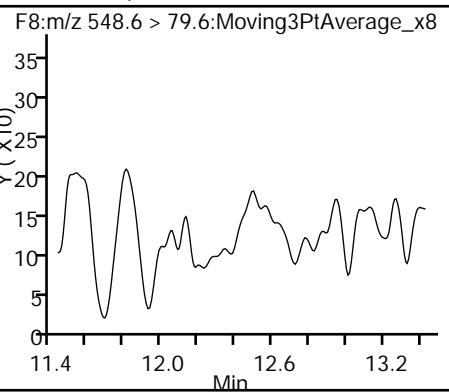
D 17 13C5 PFNA



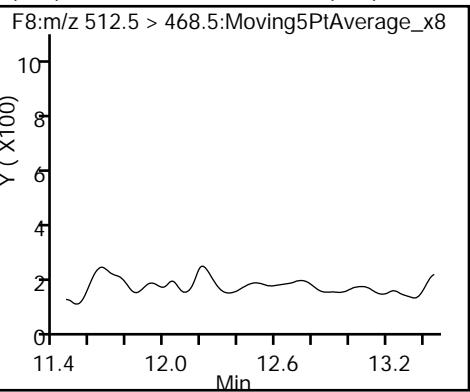
18 Perfluorononanoic acid (ND)



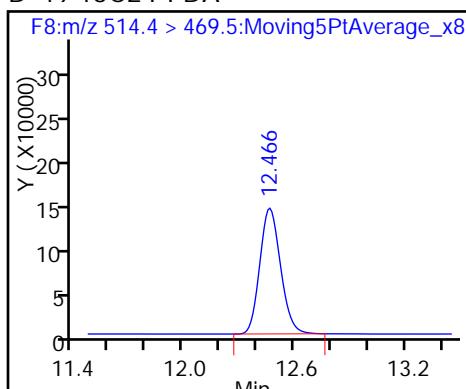
21 PFNS (Perflouro-1-nananesulfonate) (ND)



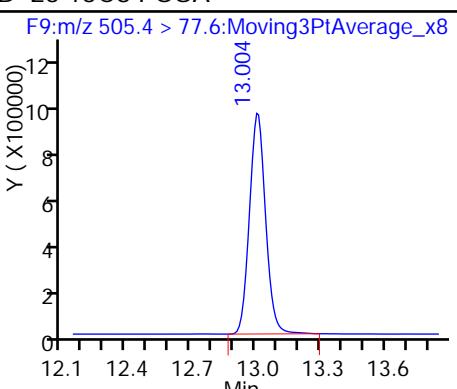
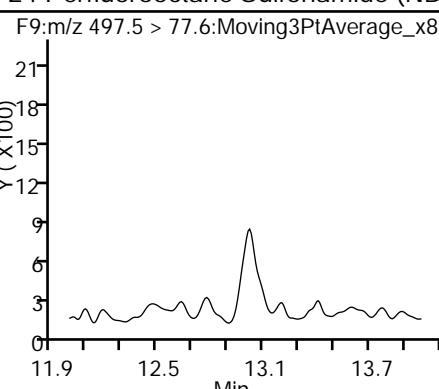
D 18 Perfluorodecanoic acid (ND)



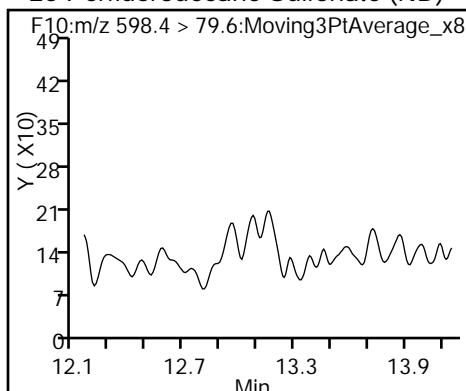
D 19 13C2 PFDA



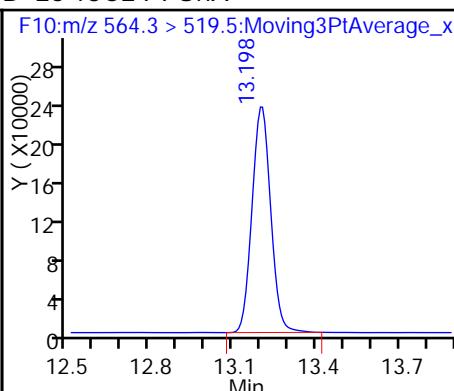
24 Perfluorooctane Sulfonamide (ND) D 23 13C8 FOSA



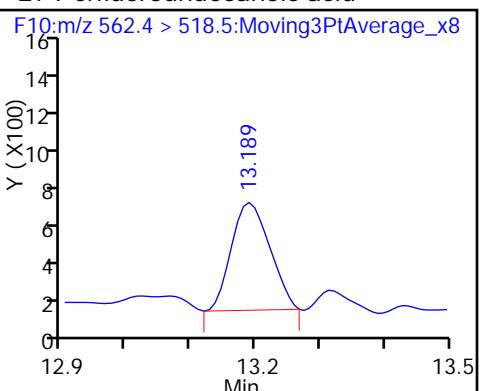
25 Perfluorodecane Sulfonate (ND)



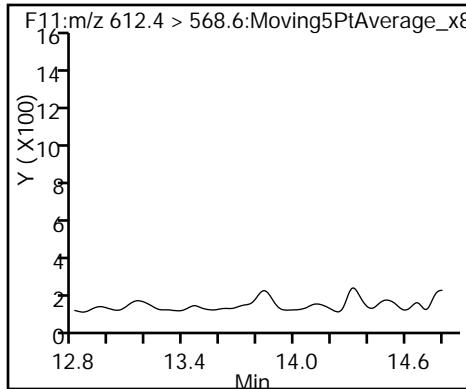
D 26 13C2 PFUna



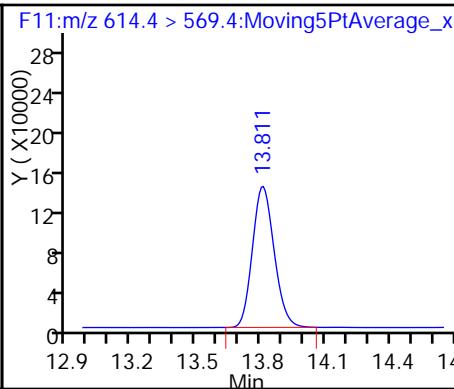
27 Perfluoroundecanoic acid



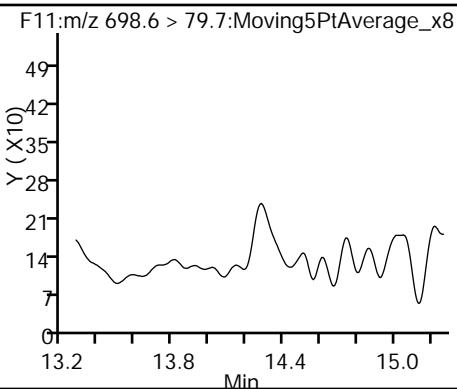
29 Perfluorododecanoic acid (ND)



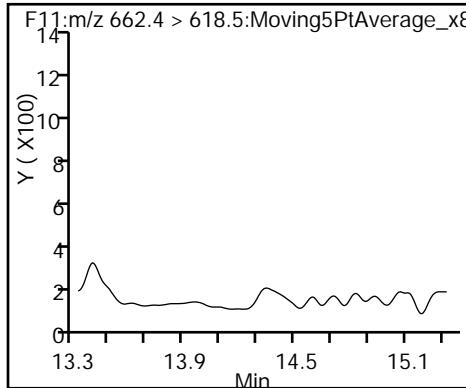
D 28 13C2 PFDoA



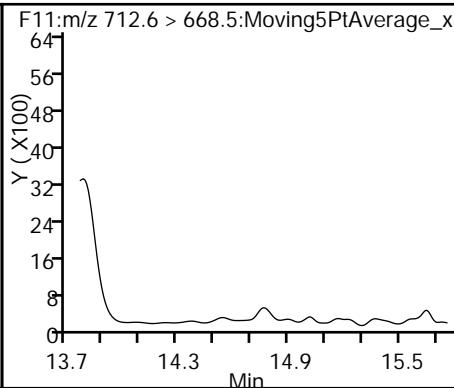
31 PFDoS (Perflouro-1-dodecanesulfona (ND))



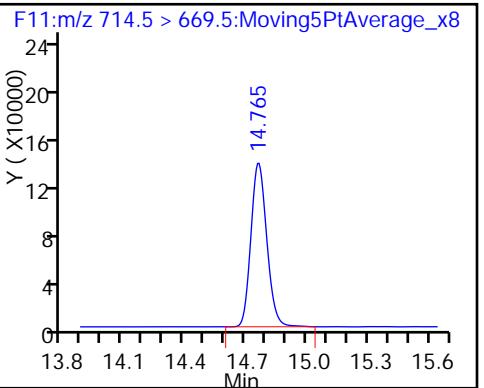
30 Perfluorotridecanoic acid (ND)



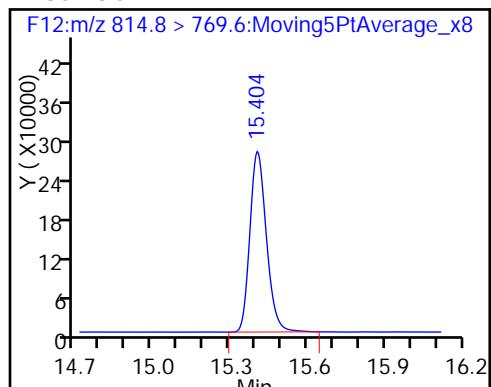
32 Perfluorotetradecanoic acid (ND)



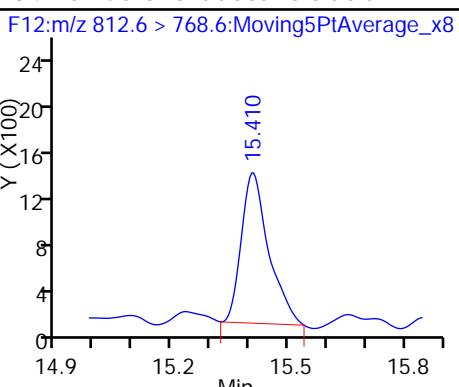
D 33 13C2-PFTeDA



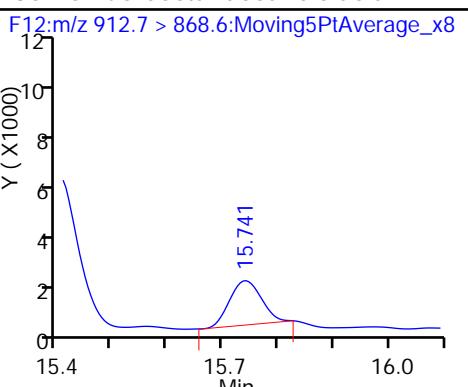
## D 35 13C2-PFHxDa



## 34 Perfluorohexadecanoic acid



## 36 Perfluoroctadecanoic acid



FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: \_\_\_\_\_ Lab Sample ID: LCS 320-59712/2-A  
Matrix: Water Lab File ID: 03DEC14A4A\_083.d  
Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
Extraction Method: 3535 Date Extracted: 12/03/2014 13:38  
Sample wt/vol: 500.00 (mL) Date Analyzed: 12/04/2014 20:24  
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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	50.4		2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	39.7		2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	101		25-150
STL00990	13C4 PFOA	98		25-150

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_083.d  
 Lims ID: LCS 320-59712/2-A  
 Client ID:  
 Sample Type: LCS  
 Inject. Date: 04-Dec-2014 20:24:40 ALS Bottle#: 23 Worklist Smp#: 47  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: lcs 320-59712/2-a 59712  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1 : Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
<b>D 1 113C4 PFBA</b>										
216.7 > 171.5	5.922	5.934	-0.012		25563	5.98		29.9	94.1	
2 Perfluorobutyric acid										
212.7 > 168.6	5.928	5.935	-0.007	1.000	404010	48.1		241	1217	
4 Perfluoropentanoic acid										
262.9 > 218.7	7.037	7.037	0.0	1.000	220687	37.4		187	351	
<b>D 3 13C5-PFPeA</b>										
267.6 > 222.7	7.040	7.038	0.002		71436	9.26		46.3	285	
5 Perfluorobutane Sulfonate										
298.8 > 79.6	7.152	7.157	-0.005	1.000	902925	21.2		120	1704	
298.8 > 98.6	7.152	7.157	-0.005	1.000	587476	1.54(0.00-0.00)		1433		
<b>D 6 13C2 PFHxA</b>										
314.6 > 269.7	8.286	8.282	0.004		240697	12.6		63.1	723	
7 Perfluorohexanoic acid										
312.9 > 268.7	8.286	8.284	0.002	1.000	317763	27.0		135	699	
22 PFPeS (Perflouro-1-pentanesulfonat										
348.7 > 79.5	8.362	8.360	0.002	0.876	816453	24.0		128	2181	
<b>D 8 13C4-PFHxA</b>										
366.6 > 321.6	9.511	9.506	0.005		273562	15.4		77.1	1033	
9 Perfluoroheptanoic acid										
362.8 > 318.7	9.511	9.508	0.003	1.000	721283	34.6		173	1713	
10 Perfluorohexane Sulfonate										
398.8 > 79.7	9.545	9.543	0.002	1.000	939802	23.6		125	2436	
<b>D 11 18O2 PFHxA</b>										
402.5 > 83.6	9.545	9.544	0.001		817188	18.2		96.1	1996	
<b>D 12 13C4 PFOA</b>										
416.5 > 371.6	10.632	10.628	0.004		1657511	49.0		98.0	2628	
13 Perfluorooctanoic acid										
412.8 > 368.8	10.632	10.628	0.004	1.000	1395384	25.2		126	1416	
412.8 > 168.7	10.632	10.628	0.004	1.000	432499	3.23(0.00-0.00)		806		

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
14 Perfluoroheptane Sulfonate										
448.8 > 79.7	10.632	10.635	-0.003	1.000	1068143	27.2		143	1398	
D 16 13C4 PFOS										
502.4 > 79.7	11.602	11.595	0.007		2170616	48.1		101	4250	
15 Perfluorooctanoic Sulfonate										
498.9 > 79.7	11.602	11.595	0.007	1.000	1628323	19.9		104	2281	
498.9 > 98.7	11.602	11.595	0.007	1.000	894977		1.82(0.00-0.00)		1434	
D 17 13C5 PFNA										
467.5 > 422.6	11.612	11.614	-0.002		828770	21.5		107	2153	
18 Perfluorononanoic acid										
462.5 > 418.6	11.620	11.616	0.004	1.000	1102178	23.1		115	1939	
21 PFNS (Perflouro-1-nananesulfonate)										
548.6 > 79.6	12.435	12.429	0.006	1.000	792745	23.8		124	1225	
20 Perfluorodecanoic acid										
512.5 > 468.5	12.466	12.461	0.005	1.000	1288911	26.2		131	1742	
D 19 13C2 PFDA										
514.4 > 469.5	12.466	12.463	0.003		1029276	20.0		100.0	1444	
24 Perfluorooctane Sulfonamide										
497.5 > 77.6	13.003	12.998	0.005	1.000	2119886	27.0		135	1726	
D 23 13C8 FOSA										
505.4 > 77.6	13.003	12.998	0.005		4137743	28.9		57.8	3273	
25 Perfluorodecane Sulfonate										
598.4 > 79.6	13.154	13.152	0.002	1.000	712044	24.2		126	1243	
D 26 13C2 PFUnA										
564.3 > 519.5	13.198	13.198	0.0		1129572	17.4		87.2	2189	
27 Perfluoroundecanoic acid										
562.4 > 518.5	13.198	13.198	0.0	1.000	1575629	31.6		158	1970	
29 Perfluorododecanoic acid										
612.4 > 568.6	13.811	13.808	0.003	1.000	1585953	29.6		148	1178	
D 28 13C2 PFDa										
614.4 > 569.4	13.811	13.809	0.002		1087143	16.6		83.2	1451	
31 PFDoS (Perflouro-1-dodecanesulfona										
698.6 > 79.7	14.265	14.272	-0.007	1.000	474162	15.9		81.9	812	
30 Perfluorotridecanoic acid										
662.4 > 618.5	14.326	14.327	-0.001	1.000	887921	20.0		99.9	609	
32 Perfluorotetradecanoic acid										
712.6 > 668.5	14.765	14.766	-0.001	1.000	470152	16.6		83.2	488	
D 33 13C2-PFTeDA										
714.5 > 669.5	14.758	14.766	-0.008		730624	10.2		51.1	1256	
D 35 13C2-PFHxDA										
814.8 > 769.6	15.404	15.415	-0.011		1315788	12.6		63.1	2100	
34 Perfluorohexadecanoic acid										
812.6 > 768.6	15.404	15.415	-0.011	1.000	1400292	19.2		96.1	1675	
36 Perfluorooctadecanoic acid										
912.7 > 868.6	15.736	15.752	-0.016	1.000	1742284	22.6		113	1367	

Data File: \\Sacchrom\\ChromData\\A4\\20141204-17820.b\\03DEC14A4A\_083.d

Injection Date: 04-Dec-2014 20:24:40

Instrument ID: A4

Lims ID: LCS 320-59712/2-A

Client ID:

Operator ID: JRB

ALS Bottle#: 23 Worklist Smp#: 47

Injection Vol: 15.0 ul

Dil. Factor: 1.0000

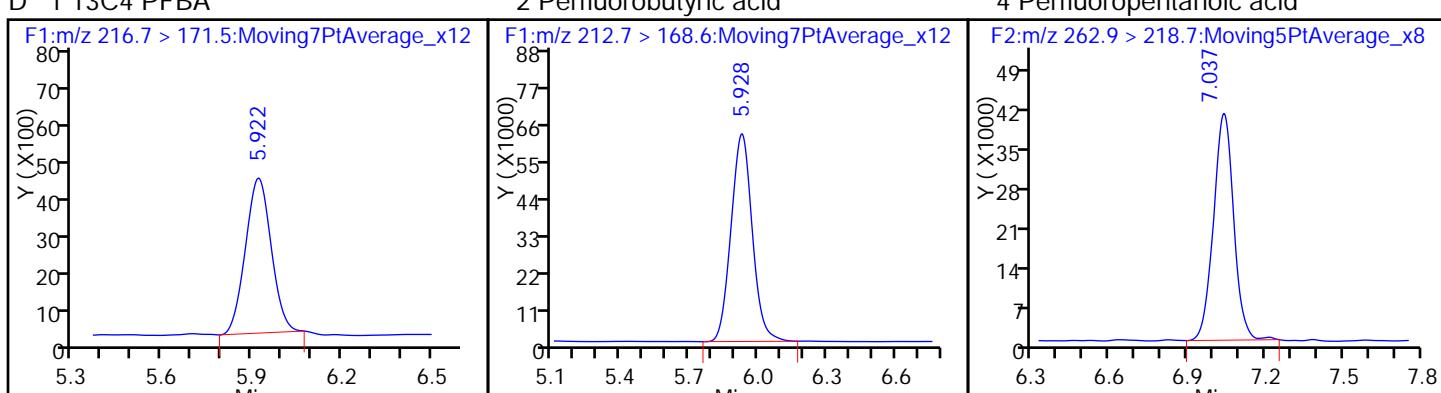
Method: PFAC\_A4

Limit Group: LC PFC ICAL

## D 1 13C4 PFBA

## 2 Perfluorobutyric acid

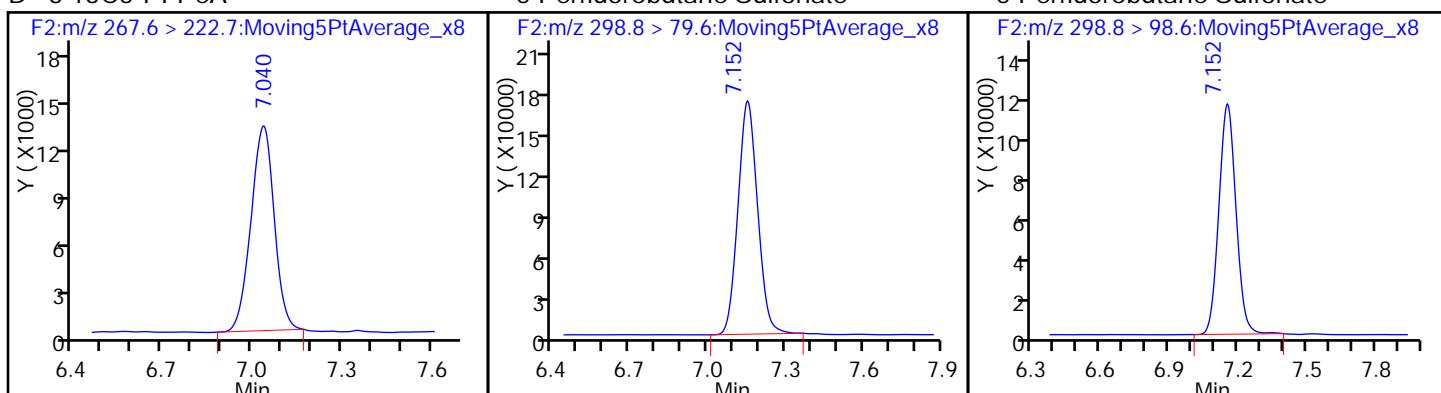
## 4 Perfluoropentanoic acid



## D 3 13C5-PFPeA

## 5 Perfluorobutane Sulfonate

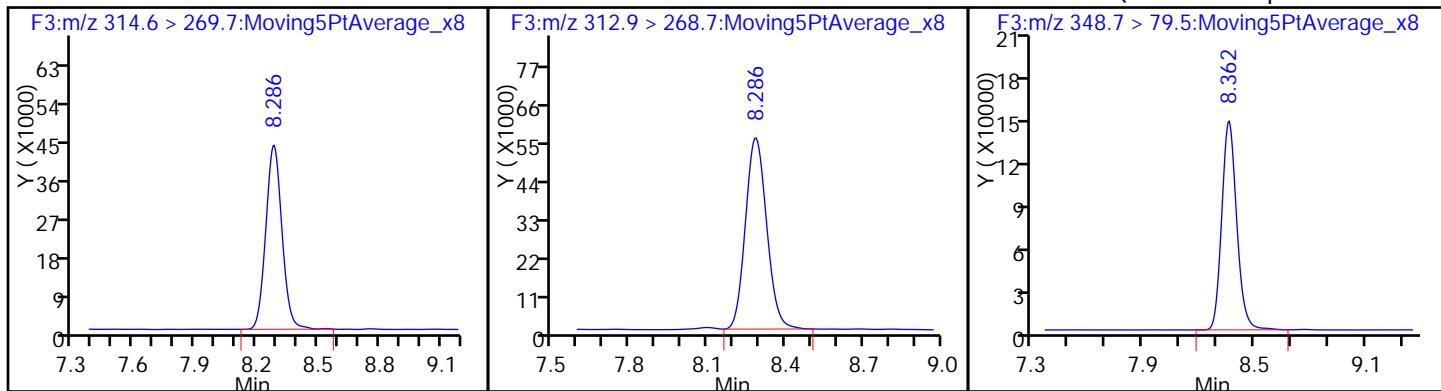
## 5 Perfluorobutane Sulfonate



## D 6 13C2 PFHxA

## 7 Perfluorohexanoic acid

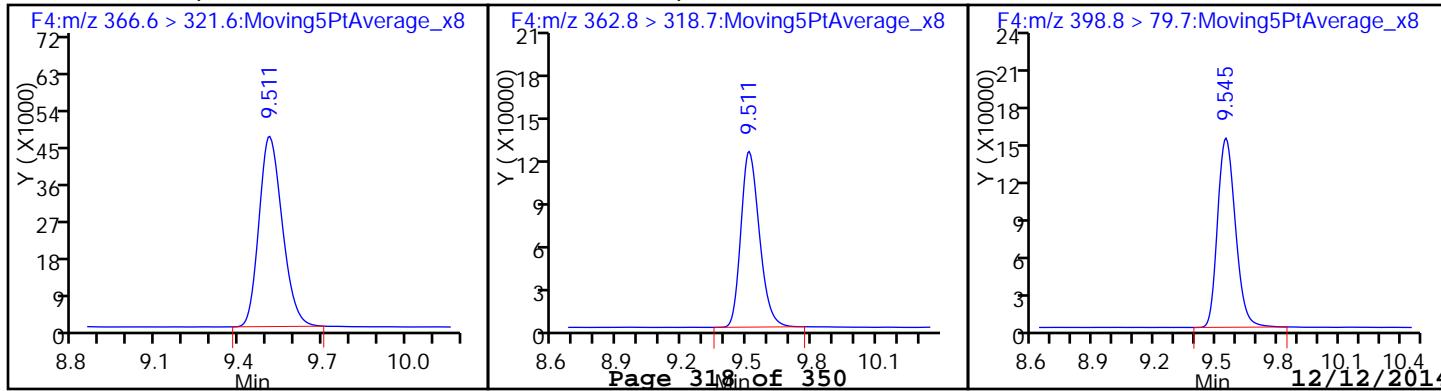
## 22 PFPeS (Perflouro-1-pentanesulfonat)



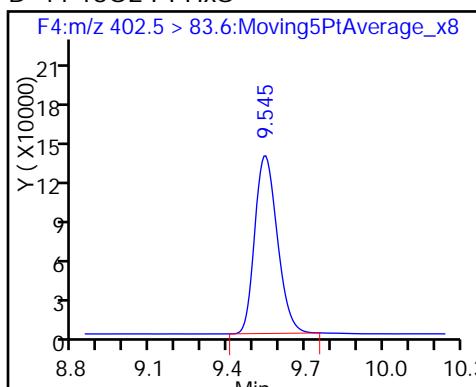
## D 8 13C4-PFHxA

## 9 Perfluoroheptanoic acid

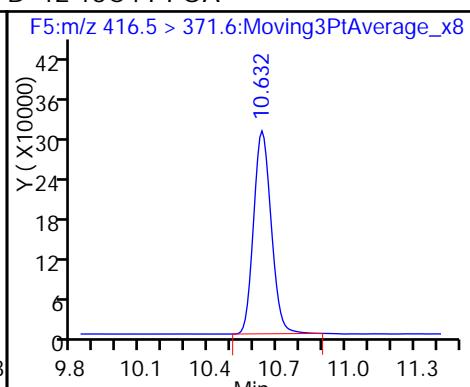
## 10 Perfluorohexane Sulfonate



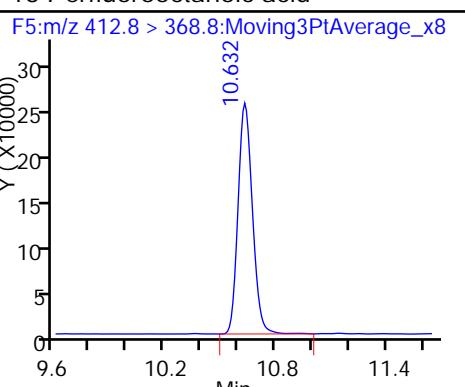
D 11 18O2 PFHxS



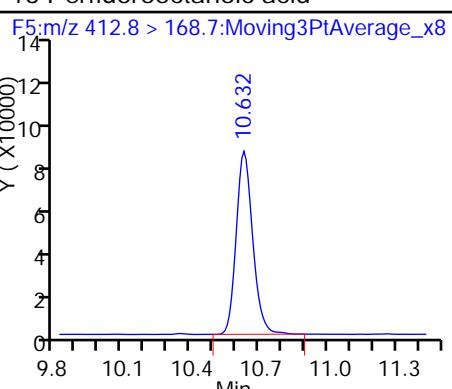
D 12 13C4 PFOA



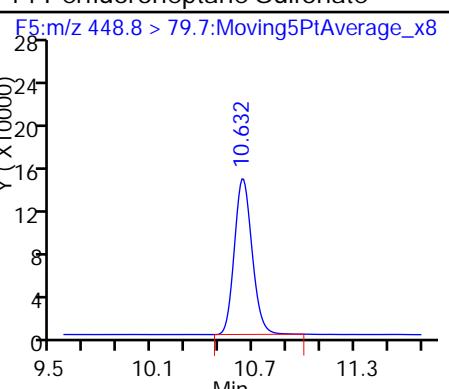
13 Perfluorooctanoic acid



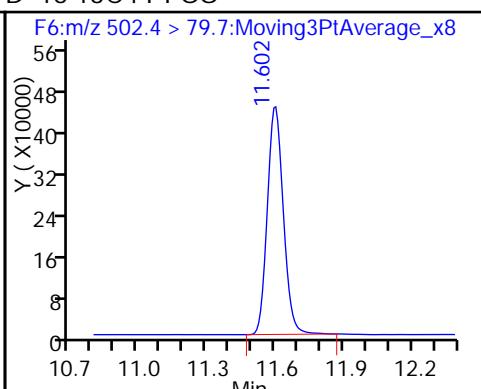
13 Perfluorooctanoic acid



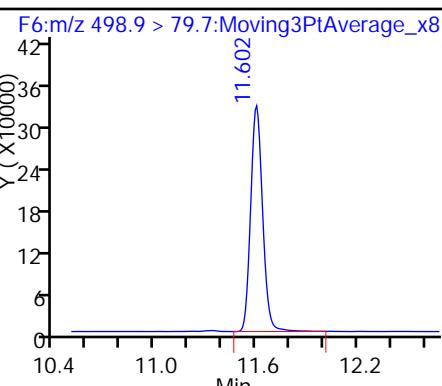
14 Perfluoroheptane Sulfonate



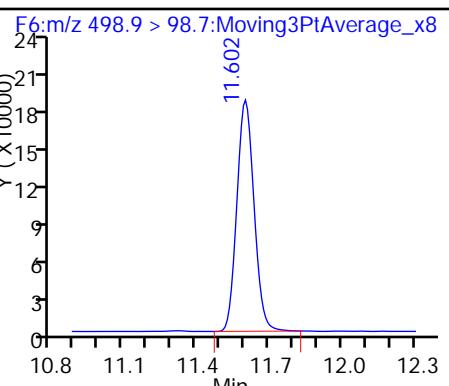
D 16 13C4 PFOS



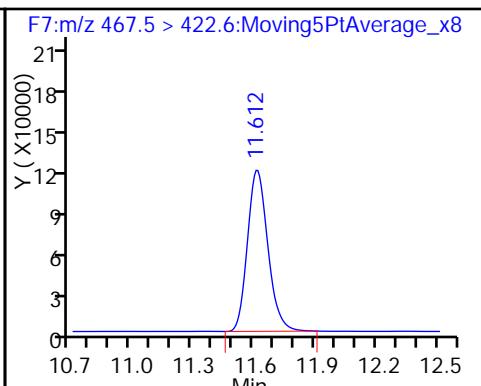
15 Perfluorooctanoic Sulfonate



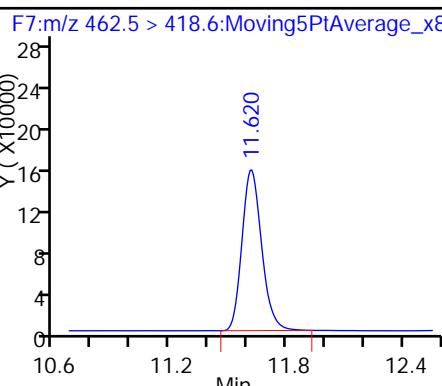
15 Perfluorooctanoic Sulfonate



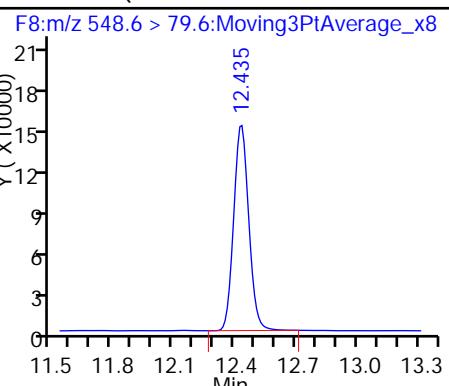
D 17 13C5 PFNA



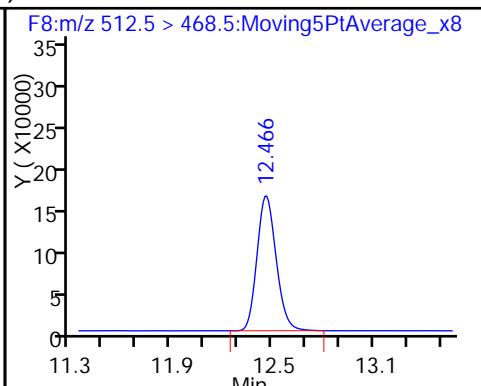
18 Perfluorononanoic acid



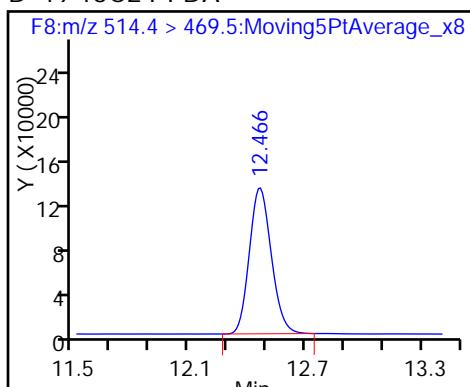
21 PFNS (Perflouro-1-nananesulfonate)



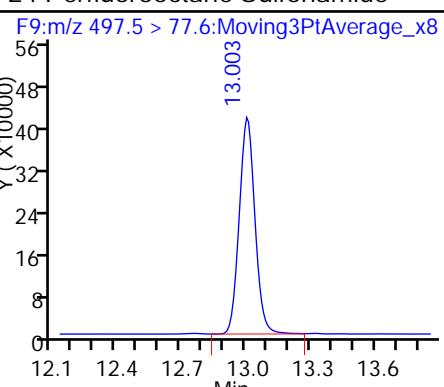
20 Perfluorodecanoic acid



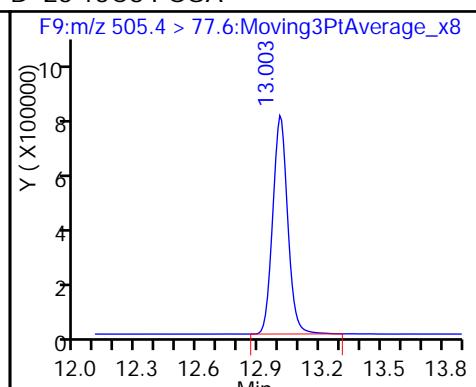
D 19 13C2 PFDA



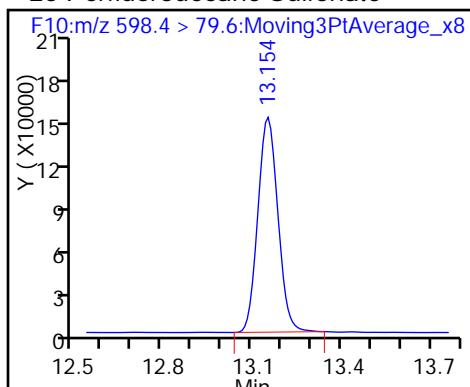
24 Perfluorooctane Sulfonamide



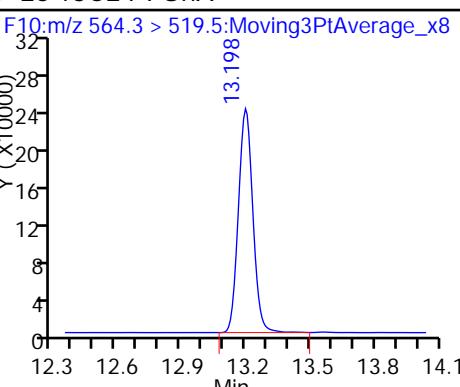
D 23 13C8 FOSA



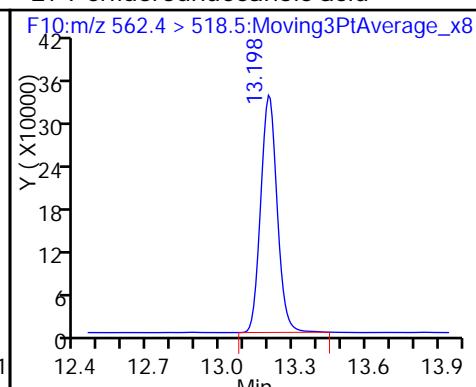
25 Perfluorodecane Sulfonate



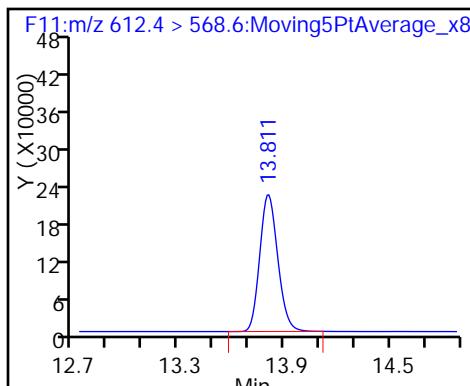
D 26 13C2 PFUna



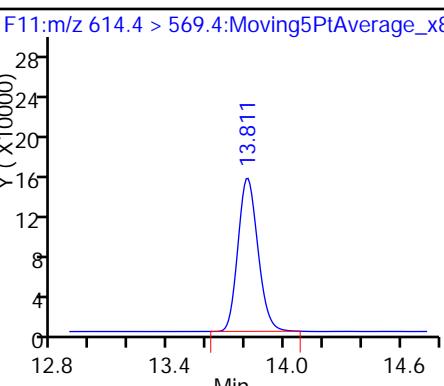
27 Perfluoroundecanoic acid



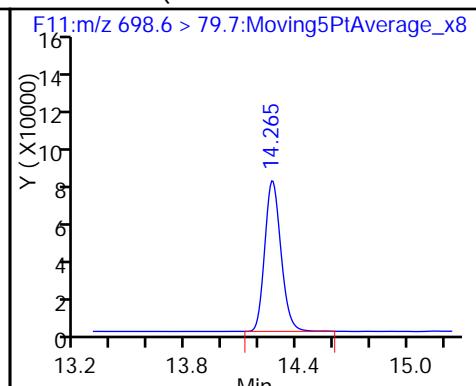
29 Perfluorododecanoic acid



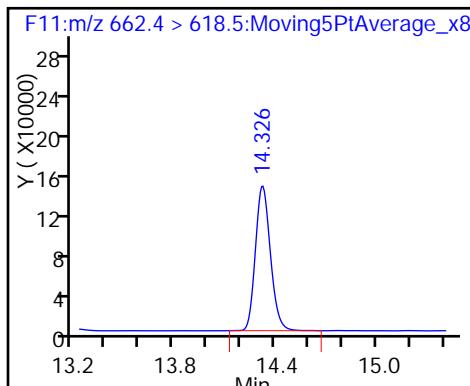
D 28 13C2 PFDoA



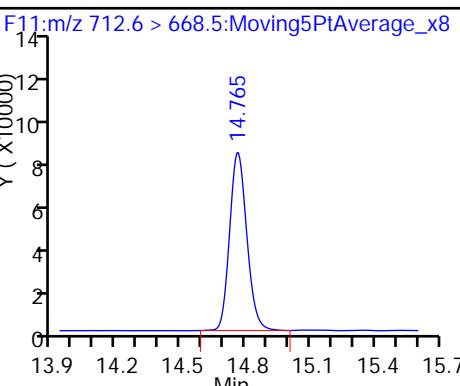
31 PFDoS (Perflouro-1-dodecanesulfona



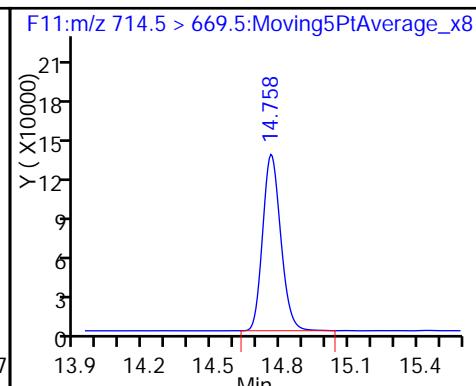
30 Perfluorotridecanoic acid



32 Perfluorotetradecanoic acid



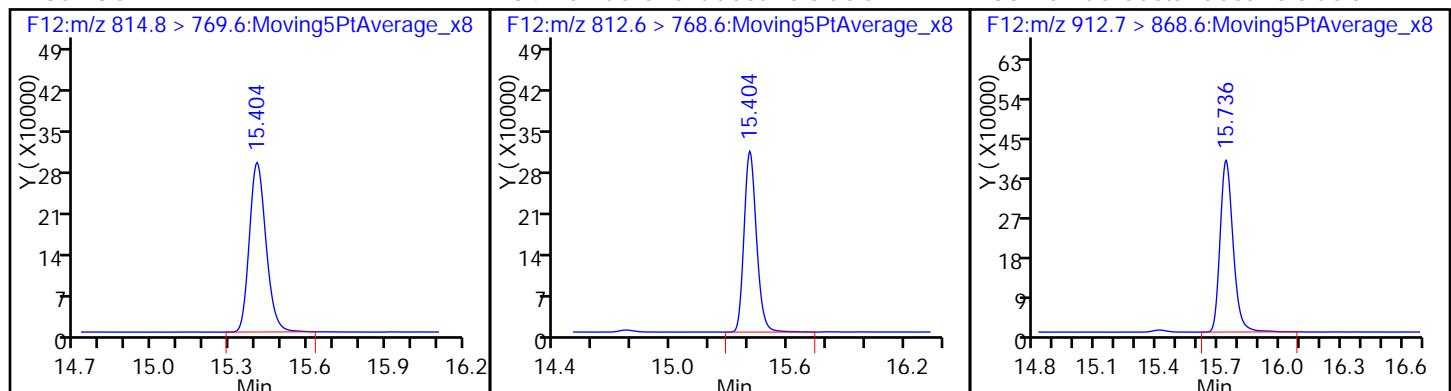
D 33 13C2-PFTeDA



D 35 13C2-PFHxD

34 Perfluorohexadecanoic acid

36 Perfluoroctadecanoic acid



## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Instrument ID: A4

Start Date: 12/04/2014 03:44

Analysis Batch Number: 59779

End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
STD 320-59779/4 IC		12/04/2014 03:44	1	03DEC14A4A_040.d	Acquity 2.1(mm)
STD 320-59779/5 IC		12/04/2014 04:05	1	03DEC14A4A_041.d	Acquity 2.1(mm)
STD 320-59779/6 IC		12/04/2014 04:26	1	03DEC14A4A_042.d	Acquity 2.1(mm)
STD 320-59779/7 IC		12/04/2014 04:47	1	03DEC14A4A_043.d	Acquity 2.1(mm)
STD 320-59779/8 IC		12/04/2014 05:09	1	03DEC14A4A_044.d	Acquity 2.1(mm)
STD 320-59779/9 IC		12/04/2014 05:30	1	03DEC14A4A_045.d	Acquity 2.1(mm)
STD 320-59779/10 IC		12/04/2014 05:51	1	03DEC14A4A_046.d	Acquity 2.1(mm)
CCV 320-59779/20		12/04/2014 09:23	1		Acquity 2.1(mm)
ICV 320-59779/22		12/04/2014 10:23	1	03DEC14A4A_058.d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 11:50	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:11	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:37	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:59	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:20	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:41	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:02	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:24	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:44	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:06	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:27	1		Acquity 2.1(mm)
CCV 320-59779/34		12/04/2014 15:48	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:10	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:31	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:52	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:13	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:34	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:56	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:17	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:38	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:59	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 19:20	1		Acquity 2.1(mm)
CCV 320-59779/45		12/04/2014 19:42	1	03DEC14A4A_081.d	Acquity 2.1(mm)
MB 320-59712/1-A		12/04/2014 20:03	1	03DEC14A4A_082.d	Acquity 2.1(mm)
LCS 320-59712/2-A		12/04/2014 20:24	1	03DEC14A4A_083.d	Acquity 2.1(mm)
320-10459-1	B7/10-GW02	12/04/2014 20:45	1	03DEC14A4A_084.d	Acquity 2.1(mm)
320-10459-2	B7/10-GW03S	12/04/2014 21:07	1	03DEC14A4A_085.d	Acquity 2.1(mm)
320-10459-3	QR4-GW02	12/04/2014 21:28	1	03DEC14A4A_086.d	Acquity 2.1(mm)
320-10459-4	B7/10-GW06S	12/04/2014 21:49	1	03DEC14A4A_087.d	Acquity 2.1(mm)

## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica SacramentoJob No.: 320-10459-1

SDG No.:

Instrument ID: A4Start Date: 12/04/2014 03:44Analysis Batch Number: 59779End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
320-10459-5	H123-GW02	12/04/2014 22:10	1	03DEC14A4A_088.d	Acquity 2.1 (mm)
320-10459-6	EFA-FD01-1014	12/04/2014 22:31	1	03DEC14A4A_089.d	Acquity 2.1 (mm)
ZZZZZ		12/04/2014 22:53	20		Acquity 2.1 (mm)
ZZZZZ		12/04/2014 23:14	20		Acquity 2.1 (mm)
ZZZZZ		12/04/2014 23:35	20		Acquity 2.1 (mm)
CCV 320-59779/57		12/04/2014 23:56	1	03DEC14A4A_093.d	Acquity 2.1 (mm)

## LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Batch Number: 59712

Batch Start Date: 12/03/14 13:38

Batch Analyst: Reed, Jonathan E

Batch Method: 3535

Batch End Date: 12/03/14 19:10

Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount	LCMPFCSU 00010	LCPPFCSP 00014		
MB 320-59712/1		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL			
LCS 320-59712/2		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL	20 uL		
320-10459-A-1	B7/10-GW02	3535, WS-LC-0025	T	448.35 mL	1.00 mL	50 uL			
320-10459-A-2	B7/10-GW03S	3535, WS-LC-0025	T	427.75 mL	1.00 mL	50 uL			
320-10459-A-3	QR4-GW02	3535, WS-LC-0025	T	427.31 mL	1.00 mL	50 uL			
320-10459-A-4	B7/10-GW06S	3535, WS-LC-0025	T	446.26 mL	1.00 mL	50 uL			
320-10459-A-5	H123-GW02	3535, WS-LC-0025	T	455.47 mL	1.00 mL	50 uL			
320-10459-A-6	EFA-FD01-1014	3535, WS-LC-0025	T	447.90 mL	1.00 mL	50 uL			

## Batch Notes

Balance ID	QA-070
H2O Lot used	11/24/14
Pipette ID	EC15219
Analyst who added reagent	JER
SU Reagent Drop	JER
SU Reagent Drop Witness	SNE
Solvent Lot #	0.3%NH4OH-Me_00006
Solvent Name	0.3%NH4OH/MeOH
SOP Number	WS-LC-0005
SPE Cartridge Type	WAX 500mg
Solid Phase Extraction Disk Lot Number	002334148A

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.

WS-LC-0025

Page 1 of 1

## HPLC/LCMS Data Review Checklist

Job Number(s): 8957, 9021, 10572, 10114, Work List ID(s): 17820  
10591, 10459

Extraction Batch: 51087, 58953, 59090, 59712 Analysis Batch(es): 59779

Delivery Rank 2, 4 Due Date: Varies

	1 <sup>st</sup> Level	2 <sup>nd</sup> Level	N/A
A. Calibration/Instrument Run QC			
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 ( $r \geq 0.995$ ).	✓	/	
• Quadratic fit criteria appropriate if required ( $r^2 \geq 0.990$ ).			✓
• For Linear Regression and Quadratic fit – Does the y-intercept support $\frac{1}{2}$ the reporting limit as described in CA-Q-S-005?			✓
• All curve points show calculated concentrations.			
3. Peaks correctly ID'd by data system.	✓	/	
5. Tune check frequency & criteria met and Tune check report attached.		/	
B. QA/QC			
1. Are all QC samples properly linked in TALS?	✓	/	
2. Method blank, LCS/LCSD and MS/SD frequencies met.	✓	/	
3. LCS/LCSD and MB data are within control limits. If not, NCM is present.	✓	/	
4. Are MS/MSD recoveries and RPD within control limits?			
5. Holding Times were met for prep and analytical. <i>In House samples + NCM</i>	✓	/	
6. IS/Surrogate recoveries meet criteria or properly noted.	✓	/	
C. Sample Analysis			
1. Was correct analysis performed and were project instructions followed?			
2. If required, are compounds within RT windows?	✓	/	
3. If required, are positive hits confirmed and >40% RPD flagged?	✓	/	
4. Manual Integrations reviewed and appropriate.			
5. All analytes correctly reported. (Primary, secondary, acceptable status)	✓		
6. Correct reporting limits used. (based on client request, prep factors, and dilutions)	✓	/	
D. Documentation			
1. Are all non-conformances documented/attached? NCM#	✓	/	
2. Do results make sense (e.g. dilutions, etc.)?	✓	/	
3. Have all flags been reviewed for appropriateness?	✓	/	
4. For level 3 and 4 reports, have forms and raw data been reviewed?			
5. Was QC Checker run for this job?	✓	/	

\*Upon completion of this checklist, the reviewer must scan and attach the checklist to the TALS job.

1<sup>st</sup> Level (Analyst): JRB

Date: 12/5/14

2<sup>nd</sup> Level Reviewer: JW

Date: 12/12/14

# Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59087

Method Code: 320-Shake\_Bath\_14D-320

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:06:49PM

Batch End: 11/26/2014 8:40:00PM

## Shake Extraction with Ultrasonic Bath Extraction

Input Sample Lab ID (Analytical Method)	SDG	Initial Amount	Final Amount	Due Date	Analytical TAT	DIV Rank	Comments	Output Sample Lab ID
MB-320-59087/1 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		 M B 3 2 0 - 5 9 0 8 7 / 1 - A
LCS-320-59087/2 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		 L C S 3 2 0 - 5 9 0 8 7 / 2 - A
320-10515-B-1 (PFC_IDA)	10502757.5501080 1	5.09 g	1.00 mL	12/9/14	12_Days	4		 3 2 0 - 1 0 5 1 5 - B - 1 - A
320-10515-B-1~MS (PFC_IDA)	10502757.5501080 1	5.13 g	1.00 mL	12/9/14	12_Days	4		 3 2 0 - 1 0 5 1 5 - B - 1 - B - M S
320-10515-B-1~MSD (PFC_IDA)	10502757.5501080 1	5.09 g	1.00 mL	12/9/14	12_Days	4		 3 2 0 - 1 0 5 1 5 - B - 1 - C - M S D
320-8957-G-1 (PFC_IDA)	N/A	5.00 g	1.00 mL	9/4/14	12_Days	2		 3 2 0 - 8 9 5 7 - G - 1 - A
320-9021-W-1 (PFC_IDA)	N/A	5.00 g	1.00 mL	9/9/14	12_Days	2		 3 2 0 - 9 0 2 1 - W - 1 - A

## Batch Notes

Blank Sand Lot # 142675

Filter Lot # NA

Millipore Water Dispense Date 11/24/14

Person's name who witnessed reagent drop SNE

SPE Cartridge Lot # 014933351A

SPE Cartridge Type WAX 150mg

Batch Comment SOLVENT: 0.3%NH4OH-Me\_00006

12/2/2014

Printed : 11/26/2014

# Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59087

Method Code: 320-Shake\_Bath\_14D-320

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:06:49PM

Batch End: 11/26/2014 8:40:00PM

## Comments

320-10515-B-1

Method Comments: See QAS, decant excess water & take sample from solids if possible

320-10515-B-1-MS

Method Comments: See QAS, decant excess water & take sample from solids if possible

320-10515-B-1~MSD

Method Comments: See QAS, decant excess water & take sample from solids if possible

320-8957-G-1

Method Comments: Please add surrogate to the container and then rinse with 50mL of solvent. See Jeremy with questions

320-9021-W-1

Method Comments: Please add surrogate to the container and then rinse with 50mL of solvent. See Jeremy with questions

# Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59087

Method Code: 320-Shake\_Bath\_14D-320

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:06:49PM

Batch End: 11/26/14 8:40pm

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-59087/1	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>Reed 11/25/14</i>	SNE 11/25/14
LCS 320-59087/2	LCMPFCSU_00010	50.00 uL	1.00 mL		
LCS 320-59087/2	LCPFCSP_00014	20.00 uL	1.00 mL		
320-10515-B-1	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10515-B-1 MS	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10515-B-1 MS	LCPFCSP_00014	20.00 uL	1.00 mL		
320-10515-B-1 MSD	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10515-B-1 MSD	LCPFCSP_00014	20.00 uL	1.00 mL		
320-8957-G-1	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-9021-W-1	LCMPFCSU_00010	50.00 uL	1.00 mL		

# **Solid SW-846-3500 Analysis Sheet**

(To Accompany Samples to Instruments)

Batch Number: 320-59087

Method Code: 320-Shake\_Bath\_14D-320

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:06:49PM

Batch End: 11/26/2014 8:40:00PM

## **Other Reagents:**

<b>Reagent</b>	<b>Amount/Units</b>	<b>Lot#:</b>



THE LEADER IN ENVIRONMENTAL TESTING

**Sacramento  
Preparation Data Review Checklist**

Preparation Batch Number(s): 59087 Test: PFC-S  
 Earliest Holding Time: 12/02/14

<b>Sample List Tab</b>		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)		NA	NA
Method/sample/login/QAS checked and correct		✓	
<b>Worksheet Tab</b>		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		✓	✓
<b>Reagents Tab</b>		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		✓	✓
<b>Batch Information</b>		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: JLWDate: 11/26/142<sup>nd</sup> Level Reviewer: HJADate: 12/1/14

Comments:

2016  
X

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-59712

Method Code: 320-3535\_IVWT-320

Batch Open: 12/3/2014 1:38:55PM

Batch End: 12/3/2014 7:10:00PM

## Solid-Phase Extraction (SPE)

Input Sample Lab ID (Analytical Method)	SDG	GrossWt TareWt	InitAmnt FinAmnt	Rcvd	PHs Adj1	Adj2	Due Date	Analytical TAT	Dlv Rank	Comments	Output Sample Lab ID
1 MB~320-59712/1 N/A	N/A		500.00 mL				N/A	N/A	N/A		 M B 3 2 0 - 5 9 7 1 2 / 1 - A
			1.00 mL								
2 LCS~320-59712/2 N/A	N/A		500.00 mL				N/A	N/A	N/A		 L C S 3 2 0 - 5 9 7 1 2 / 2 - A
			1.00 mL								
3 320-10459-A-1 (PFC_IDA)	N/A		448.35 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 1 - B
			1.00 mL								
4 320-10459-A-2 (PFC_IDA)	N/A		427.75 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 2 - B
			1.00 mL								
5 320-10459-A-3 (PFC_IDA)	N/A		427.31 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 3 - B
			1.00 mL								
6 320-10459-A-4 (PFC_IDA)	N/A		446.26 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 4 - B
			1.00 mL								
7 320-10459-A-5 (PFC_IDA)	N/A		455.47 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 5 - B
			1.00 mL								
8 320-10459-A-6 (PFC_IDA)	N/A		447.90 mL				12/6/14	18_Days	4		 3 2 0 - 1 0 4 5 9 - A - 6 - B
			1.00 mL								

12/12/2014

Printed : 12/3/2014

Page 1 of 4

TestAmerica Sacramento

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59712

Method Code: 320-3535\_IVWT-320

Analyst: Reed, Jonathan E

Batch Open: 12/3/2014 1:38:55PM

Batch End: 12/3/2014 7:10:00PM

## Batch Notes

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type WAX 500mg

Solid Phase Extraction Disk Lot 002334148A

Number

H<sub>2</sub>O Lot used 11/24/14

Pipette ID EC15219

Solvent Name 0.3%NH4OH/MeOH

Solvent Lot # 0.3%NH4OH-Me\_00006

Analyst who added reagent JER

SU Reagent Drop JER

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-0005

Batch Comment NA

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59712

Method Code: 320-3535\_IVWT-320

Analyst: Reed, Jonathan E

Batch Open: 12/3/2014 1:38:55PM

Batch End: 12/3/2014 7:10:00PM

## Comments

320-10459-A-1

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

320-10459-A-2

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

320-10459-A-3

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

320-10459-A-4

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

320-10459-A-5

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

320-10459-A-6

Method Comments: Q5  
Rework Comments: MB and LCS contaminated.

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-59712

Method Code: 320-3535\_IVWT-320

Batch Open: 12/3/2014 1:38:55PM

Batch End: 12/03/14 7:10 PM

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-59712/1	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>Reed 12/03/14</i>	SNE 12/3/14
LCS 320-59712/2	LCMPFCSU_00010	50.00 uL	1.00 mL		
LCS 320-59712/2	LCPFCSP_00014	20.00 uL	1.00 mL		
320-10459-A-1	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10459-A-2	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10459-A-3	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10459-A-4	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10459-A-5	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10459-A-6	LCMPFCSU_00010	50.00 uL	1.00 mL		

## Other Reagents:

Reagent

Amount/Units

Lot#:

Preparation Batch Number(s): 59712

Test: PFC-L

Earliest Holding Time: 4/22/14

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: \_\_\_\_\_

2<sup>nd</sup> Level Reviewer: \_\_\_\_\_

Comments: \_\_\_\_\_

Date: 12/03/14

Date: 12-4-14

*Box 11*

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-59090

Batch Open: 11/25/2014 2:15:36PM

Method Code: 320-3535\_IVWT-320

Batch End: 11/26/2014 2:38:00PM

## Solid-Phase Extraction (SPE)

		Input Sample Lab ID (Analytical Method)	SDG	GrossWt TareWt	InitAmnt FinAmnt	Rcvd	PHs Adj1	Adj2	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
1	N/A	MB~320-59090/1 N/A	N/A		500.00 mL				N/A	N/A	N/A		 M B 3 2 0 - 5 9 0 9 0 / 1 - A
					1.00 mL								 L C S 3 2 0 - 5 9 0 9 0 / 2 - A
2	N/A	LCS~320-59090/2 N/A	N/A		500.00 mL				N/A	N/A	N/A		 3 2 0 - 1 0 5 9 1 - B - 1 - A
					1.00 mL								 3 2 0 - 1 0 5 9 1 - B - 1 - A
3	N/A	320-10591-B-1 (PFC_IDA)	N/A	552.44 g	512.3 mL				12/5/14	8_Days	4		 3 2 0 - 1 0 5 9 1 - B - 1 - A
				40.19 g	1.00 mL								 3 2 0 - 1 0 5 9 1 - B - 1 - A

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59090

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:15:36PM

Method Code: 320-3535\_IVWT-320

Batch End: 11/26/2014 2:38:00PM

## Batch Notes

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type WAX 500mg

Solid Phase Extraction Disk Lot 002334148A

Number

H<sub>2</sub>O Lot used 11/24/14

Pipette ID EC15219

Solvent Name 0.3%NH4OH/MeOH

Solvent Lot # 0.3%NH4OH-Me\_00006

Analyst who added reagent JER

SU Reagent Drop JER

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 NA

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-59090

Method Code: 320-3535\_IVWT-320

Batch Open: 11/25/2014 2:15:36PM

Batch End: 11/26/2014 2:38:00PM

## Comments

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-59090

Batch Open: 11/25/2014 2:15:36PM

Method Code: 320-3535\_IVWT-320

Batch End: 11/26/14 2:38 PM

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-59090/1	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>J Reed</i> 11/25/14	SNE 11/25/14
LCS 320-59090/2	LCMPFCSU_00010	50.00 uL	1.00 mL		
LCS 320-59090/2	LCPFCSP_00014	20.00 uL	1.00 mL		
320-10591-B-1	LCMPFCSU_00010	50.00 uL	1.00 mL		

## Other Reagents:

Reagent	Amount/Units	Lot#:

Preparation Batch Number(s): 59090

Test: PFC-L

Earliest Holding Time: 11/21/14

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: J. Keay

Date: 11/26/14

2<sup>nd</sup> Level Reviewer: HJA

Date: 11-26-14

Comments: \_\_\_\_\_

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/2014 5:08:00PM

## Solid-Phase Extraction (SPE)

Page 341	Input Sample Lab ID (Analytical Method)	SDG	GrossWt TareWt	InitAmnt FinAmnt	Rcvd	PHs Adj1	Adj2	Due Date	Analytical TAT	Dlv Rank	Comments	Output Sample Lab ID
1	MB~320-58953/1 N/A	N/A		500.00 mL				N/A	N/A	N/A		 M B 3 2 0 - 5 8 9 5 3 / 1 - A
2	LCS~320-58953/2 N/A			1.00 mL								 L C S 3 2 0 - 5 8 9 5 3 / 2 - A
3	320-10572-A-1 (PFC_IDA)	N/A		500.25 mL				12/9/14	18_Days	4		 3 2 0 - 1 0 5 7 2 - A - 1 - A
4	N/A			1.00 mL								
5	N/A	N/A						N/A	N/A	N/A		
6	N/A			1.00 mL								
7	320-10572-A-2 (PFC_IDA)	N/A		499.85 mL				12/9/14	18_Days	4		 3 2 0 - 1 0 5 7 2 - A - 2 - A
8	320-10572-A-3 (PFC_IDA)			1.00 mL								
9	320-10572-A-4 (PFC_IDA)	N/A		502.05 mL				12/9/14	18_Days	4		 3 2 0 - 1 0 5 7 2 - A - 3 - A
10	320-10114-A-16 (PFC_IDA)		LCMS PFC Aq A4	500.00 mL				1/7/15	47_Days	4	LOD-1	 3 2 0 - 1 0 1 1 4 - A - 1 6 - A
11	320-10114-A-17 (PFC_IDA)	LCMS PFC Aq A4	500.00 mL	1.00 mL								
12	12/12/2014			1.00 mL				1/7/15	47_Days	4	LOD-2	 3 2 0 - 1 0 1 1 4 - A - 1 7 - A

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Analyst: Reed, Jonathan E

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/2014 5:08:00PM

11	320-10114-A-18 (PFC_IDA)	LCMS PFC Aq A4		500.00 mL			1/7/15	47_Days	4	LOD-3	 3 2 0 - 1 0 1 1 4 - A - 1 8 - A
12	320-10114-A-19 (PFC_IDA)	LCMS PFC Aq A4		500.00 mL			1/7/15	47_Days	4	LOQ-1	 3 2 0 - 1 0 1 1 4 - A - 1 9 - A
13	320-10114-A-20 (PFC_IDA)	LCMS PFC Aq A4		500.00 mL			1/7/15	47_Days	4	LOQ-2	 3 2 0 - 1 0 1 1 4 - A - 2 0 - A
				1.00 mL							
				1.00 mL							

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Analyst: Reed, Jonathan E

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/2014 5:08:00PM

## Batch Notes

First Start time NA

First End time NA

Balance ID QA-070

SPE Cartridge Type WAX 500mg

Solid Phase Extraction Disk Lot 0023341484

Number

H<sub>2</sub>O Lot used 11/14/14

Pipette ID EC15219

Solvent Name 0.3%NH4OH/MeOH

Solvent Lot # 0.3%NH4OH-Me\_00006

Analyst who added reagent JER

SU Reagent Drop JER

SU Reagent Drop Witness AVM

Acid Name NA

Acid Lot NA

Reagent ID NA

Reagent Lot Number NA

NaCl Lot # NA

SOP Number WS-LC-0025

Batch Comment NA

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Analyst: Reed, Jonathan E

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/2014 5:08:00PM

## Comments

320-10572-A-1

Method Comments: Q5

320-10572-A-2

Method Comments: Q5

320-10572-A-3

Method Comments: Q5

320-10572-A-4

Method Comments: Q5

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/14 5:08 PM

## Reagent Additions Worksheet

Lab ID	Reagent Code	Amount Added	Final Amount	By	Witness
MB 320-58953/1	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>J Reed 11/24/14</i>	A/M 11/24/14
LCS 320-58953/2	LCMPFCSU_00010	50.00 uL	1.00 mL		
LCS 320-58953/2	LCPFCSP_00014	20.00 uL	1.00 mL		
320-10572-A-1	LCMPFCSU_00010	50.00 uL	1.00 mL		
	LCMPFCSU_00010	50.00 uL	1.00 mL		
	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>J Reed 11/24/14</i>	
320-10572-A-2	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10572-A-3	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10572-A-4	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10114-A-16	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10114-A-16	LCPFCSP_00016	50.00 uL	1.00 mL		
320-10114-A-17	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10114-A-17	LCPFCSP_00016	100.00 uL	1.00 mL		
320-10114-A-18	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10114-A-18	LCPFCSP_00016	150.00 uL	1.00 mL		
320-10114-A-19	LCMPFCSU_00010	50.00 uL	1.00 mL		
320-10114-A-19	LCPFCSP_00016	125.00 uL	1.00 mL		
320-10114-A-20	LCMPFCSU_00010	50.00 uL	1.00 mL	<i>✓</i>	

# Aqueous Extraction Analysis Sheet

(To Accompany Samples to Instruments)

Analyst: Reed, Jonathan E

Batch Number: 320-58953

Method Code: 320-3535\_IVWT-320

Batch Open: 11/24/2014 4:22:13PM

Batch End: 11/25/14 5:08 PM

320-10114-A-20	LCPFCSP_00016	200.00 $\mu$ L	1.00 mL	<i>JReed 11/24/14</i>	<i>ADM 11/24/14</i>
----------------	---------------	----------------	---------	-----------------------	---------------------

## Other Reagents:

Reagent	Amount/Units	Lot#:

Preparation Batch Number(s): 58953

Test: PFC-L

Earliest Holding Time: 11/25/14

<b>Sample List Tab</b>	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	✓	
<b>Worksheet Tab</b>	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	✓	✓
<b>Reagents Tab</b>	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	✓	
<b>Batch Information</b>	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: 11/25/14

2<sup>nd</sup> Level Reviewer: \_\_\_\_\_

Date: 11/25/14

Comments: \_\_\_\_\_

# **Shipping and Receiving Documents**



**Tetra Tech, Inc.**

**CHAIN OF CUSTODY**

**NUMBER** No. 0533

PAGE 1 OF 1

**DISTRIBUTION:**

DISTRIBUTION: WHITE (ACCOMPANIES SAMPLE)      YELLOW (FIELD COPY)      PINK (FILE COPY)      4/02R  
FORM NO. TINUS-001

4/02R

FORM NO. TtNUS-001

## Login Sample Receipt Checklist

Client: Tetra Tech, Inc.

Job Number: 320-10459-1

**Login Number: 10459**

**List Source: TestAmerica Sacramento**

**List Number: 1**

**Creator: Sadler, Jeremy**

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.	N/A	
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?	True	
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	N/A	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	

"B7/10-GW02","WS-LC-0025","RES","320-10459-1","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)","6.6","ng/L","H","1.4","DL","","TRG","","","2.2","LOQ","YES",-99,"","448.35","1.00","1.7","","" "B7/10-GW02","WS-LC-0025","RES","320-10459-1","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)","10","ng/L","H","0.83","DL","","TRG","","","2.2","LOQ","YES",-99,"","448.35","1.00","1.7","","" "B7/10-GW02","WS-LC-0025","RES","320-10459-1","TALSAC","STL00990","13C4  
PFOA","95","ng/L","","-99","DL","","TRG","85","","-99","LOQ","YES","112","","448.35","1.00","0","","" "B7/10-GW02","WS-LC-0025","RES","320-10459-1","TALSAC","STL00991","13C4  
PFOS","120","ng/L","","-99","DL","","TRG","116","","-99","LOQ","YES","107","","448.35","1.00","0","","" "B7/10-GW03S","WS-LC-0025","RES","320-10459-2","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)","1.8","ng/L","U H","1.5","DL","","TRG","","","2.3","LOQ","YES",-99,"","427.75","1.00","1.8","","" "B7/10-GW03S","WS-LC-0025","RES","320-10459-2","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)","2.0","ng/L","J H","0.87","DL","","TRG","","","2.3","LOQ","YES",-99,"","427.75","1.00","1.8","","" "B7/10-GW03S","WS-LC-0025","RES","320-10459-2","TALSAC","STL00990","13C4  
PFOA","75","ng/L","","-99","DL","","TRG","64","","-99","LOQ","YES","117","","427.75","1.00","0","","" "B7/10-GW03S","WS-LC-0025","RES","320-10459-2","TALSAC","STL00991","13C4  
PFOS","120","ng/L","","-99","DL","","TRG","108","","-99","LOQ","YES","112","","427.75","1.00","0","","" "QR4-GW02","WS-LC-0025","RES","320-10459-3","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)","1.8","ng/L","U H","1.5","DL","","TRG","","","2.3","LOQ","YES",-99,"","427.31","1.00","1.8","","" "QR4-GW02","WS-LC-0025","RES","320-10459-3","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)","1.8","ng/L","U H",0.88,"DL","","TRG","","","2.3","LOQ","YES",-99,"","427.31","1.00","1.8","","" "QR4-GW02","WS-LC-0025","RES","320-10459-3","TALSAC","STL00990","13C4  
PFOA","110","ng/L","","-99","DL","","TRG","90","","-99","LOQ","YES","117","","427.31","1.00","0","","" "QR4-GW02","WS-LC-0025","RES","320-10459-3","TALSAC","STL00991","13C4  
PFOS","120","ng/L","","-99","DL","","TRG","109","","-99","LOQ","YES","112","","427.31","1.00","0","","" "B7/10-GW06S","WS-LC-0025","RES","320-10459-4","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)","2.2","ng/L","H","1.4","DL","","TRG","","","2.2","LOQ","YES",-99,"","446.26","1.00","1.7","","" "B7/10-GW06S","WS-LC-0025","RES","320-10459-4","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)","20","ng/L","H",0.84,"DL","","TRG","","","2.2","LOQ","YES",-99,"","446.26","1.00","1.7","","" "B7/10-GW06S","WS-LC-0025","RES","320-10459-4","TALSAC","STL00990","13C4  
PFOA","75","ng/L","","-99","DL","","TRG","67","","-99","LOQ","YES","112","","446.26","1.00","0","","" "B7/10-GW06S","WS-LC-0025","RES","320-10459-4","TALSAC","STL00991","13C4  
PFOS","120","ng/L","","-99","DL","","TRG","108","","-99","LOQ","YES","107","","446.26","1.00","0","","" "H123-GW02","WS-LC-0025","RES","320-10459-5","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)","1.6","ng/L","U H",1.4,"DL","","TRG","","","2.2","LOQ","YES",-99,"","455.47","1.00","1.6","","" "H123-GW02","WS-LC-0025","RES","320-10459-5","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)","17","ng/L","H",0.82,"DL","","TRG","","","2.2","LOQ","YES",-99,"","455.47","1.00","1.6","","" "H123-GW02","WS-LC-0025","RES","320-10459-5","TALSAC","STL00990","13C4  
PFOA","98","ng/L","","-99","DL","","TRG","89","","-99","LOQ","YES","110","","455.47","1.00","0","","" "H123-GW02","WS-LC-0025","RES","320-10459-5","TALSAC","STL00991","13C4  
PFOS","110","ng/L","","-99","DL","","TRG","107","","-99","LOQ","YES","105","","455.47","1.00","0","","" "EFA-FD01-1014","WS-LC-0025","RES","320-10459-6","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)",1.7,ng/L,U H,1.4,DL,,"TRG","","","2.2","LOQ","YES",-99,"",447.90,"1.00",1.7,"" "EFA-FD01-1014","WS-LC-0025","RES","320-10459-6","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)",16,ng/L,H,0.84,DL,,"TRG","","","2.2","LOQ","YES",-99,"",447.90,"1.00",1.7,"" "EFA-FD01-1014","WS-LC-0025","RES","320-10459-6","TALSAC","STL00990","13C4  
PFOA","91","ng/L","","-99","DL","","TRG","82","","-99","LOQ","YES","112","","447.90","1.00","0","","" "EFA-FD01-1014","WS-LC-0025","RES","320-10459-6","TALSAC","STL00991","13C4  
PFOS","120","ng/L","","-99","DL","","TRG","109","","-99","LOQ","YES","107","","447.90","1.00","0","","" "LCS 320-59712/2-A","WS-LC-0025","RES","LCS 320-59712/2-A","TALSAC","1763-23-1","Perfluorooctane Sulfonate (PFOS)",39.7,ng/L,"",1.3,DL,,"SPK",104,"",2.0,LOQ,YES,38.2,"",500.00,"1.00",1.5,"" "LCS 320-59712/2-A","WS-LC-0025","RES","LCS 320-59712/2-A","TALSAC","335-67-1","Perfluorooctanoic acid (PFOA)",50.4,ng/L,"",0.75,DL,,"SPK",126,"",2.0,LOQ,YES,40.0,"",500.00,"1.00",1.5,""

"LCS 320-59712/2-A","WS-LC-0025","RES","LCS 320-59712/2-A","TALSAC","STL00990","13C4  
PFOA","98.0","ng/L","","-99","DL","","SPK","98","","-99","LOQ","YES","100","","500.00","1.00","0","","  
"LCS 320-59712/2-A","WS-LC-0025","RES","LCS 320-59712/2-A","TALSAC","STL00991","13C4  
PFOS","96.3","ng/L","","-99","DL","","SPK","101","","-99","LOQ","YES","95.6","","500.00","1.00","0","","  
"MB 320-59712/1-A","WS-LC-0025","RES","MB 320-59712/1-A","TALSAC","1763-23-1","Perfluorooctane  
Sulfonate (PFOS)","1.5","ng/L","U","1.3","DL","","TRG","","2.0","LOQ","YES",-99,"500.00","1.00","1.5","","  
"MB 320-59712/1-A","WS-LC-0025","RES","MB 320-59712/1-A","TALSAC","335-67-1","Perfluorooctanoic acid  
(PFOA)","0.809","ng/L","J","0.75","DL","","TRG","","2.0","LOQ","YES",-99,"500.00","1.00","1.5","","  
"MB 320-59712/1-A","WS-LC-0025","RES","MB 320-59712/1-A","TALSAC","STL00990","13C4  
PFOA","109","ng/L","","-99","DL","","TRG","109","","-99","LOQ","YES","100","","500.00","1.00","0","","  
"MB 320-59712/1-A","WS-LC-0025","RES","MB 320-59712/1-A","TALSAC","STL00991","13C4  
PFOS","106","ng/L","","-99","DL","","TRG","111","","-99","LOQ","YES","95.6","","500.00","1.00","0","","  
"13-CTO WE09","NAS, Brunswick","B7/10-GW02","11/15/2014 12:05","AQ","320-10459-1","NM","","0.00","WS-  
LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
20:45","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","B7/10-GW03S","11/15/2014 15:20","AQ","320-10459-2","NM","","0.00","WS-  
LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
21:07","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","QR4-GW02","11/16/2014 10:20","AQ","320-10459-3","NM","","0.00","WS-  
LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
21:28","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","B7/10-GW06S","11/16/2014 11:50","AQ","320-10459-4","NM","","0.00","WS-  
LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
21:49","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","H123-GW02","11/16/2014 15:00","AQ","320-10459-5","NM","","0.00","WS-  
LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
22:10","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","EFA-FD01-1014","11/16/2014 00:00","AQ","320-10459-  
6","NM","","0.00","WS-LC-0025","3535","RES","12/03/2014 13:38","12/04/2014  
22:31","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","11/18/2014 09:50","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","LCS 320-59712/2-A","","AQ","LCS 320-59712/2-A","LCS","","-99","WS-LC-  
0025","3535","RES","12/03/2014 13:38","12/04/2014  
20:24","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","12/03/2014 13:38","11/18/2014 15:12","","  
"13-CTO WE09","NAS, Brunswick","MB 320-59712/1-A","","AQ","MB 320-59712/1-A","MB","","-99","WS-LC-  
0025","3535","RES","12/03/2014 13:38","12/04/2014  
20:03","TALSAC","COA","WET","NA","1","NA","NA","","100","320-59712","320-59712","NA","320-59779","320-  
10459-1","12/03/2014 13:38","11/18/2014 15:12","","

**TETRA TECH****INTERNAL CORRESPONDENCE**

TO: J. ORIENT DATE: FEBRUARY 17, 2015

FROM: MICHELLE L. ALLEN COPIES: DV FILE

**SUBJECT:** ORGANIC DATA VALIDATION – PFOA/PFOS  
FORMER NAVAL AIR STATION (NAS) BRUNSWICK, BRUNSWICK, ME  
CTO WE09  
SAMPLE DELIVERY GROUPS (SDGs) 320-10459-1 & 320-10572-1**SAMPLES:** 10/Aqueous/PFOA/PFOSSDG 320-10459-1B7/10-GW02  
EFA-FD01-1014B7/10-GW03S  
H123-GW02B7/10-GW06S  
QR4-GW02SDG 320-10572-1EFA-GW01  
EFA-GW09

EFA-GW02

EFA-GW03

Overview

The sample set for former NAS Brunswick, SDGs 320-10459-1 & 320-10572-1 consisted of ten (10) aqueous environmental samples. All ten (1) aqueous samples were analyzed for perfluorooctanoic acid (PFOA) and perfluorooctane sulfonate (PFOS). One field duplicate pair was included in these Sample Delivery Groups (SDGs): EFA-FD01-1014/H123-GW02.

The samples were collected by Tetra Tech, Inc. on November 15, 16, 18, 19, and 20, 2014 and analyzed by TestAmerica. The PFOA & PFOS analyses were conducted using Test America Laboratories (TAL) Standard Operating Procedure (SOP) WS-LC-0025 analytical and reporting protocols. The data was evaluated based on the following parameters:

- \*     • Data Completeness
- \*     • Holding Times/Sample Preservation
- \*     • LC/MS Tuning
- \*     • Initial and Continuing Calibration Results
- \*     • Laboratory Method Blank Results
- \*     • Surrogate Spike Recoveries
- \*     • Internal Standard Recoveries
- \*     • Laboratory Control Sample Results
- \*     • Field Duplicate Precision
- \*     • Detection Limits
- \*     • Compound Identification and Quantification

The asterisk (\*) indicates that all quality control criteria were met for this parameter. Qualified (if applicable) analytical results are summarized in Appendix A. Results as reported by the laboratory are presented in Appendix B, and Appendix C contains the documentation to support the findings as discussed in this data validation report. An EPA Region 1 tier II validation was performed on the data in these SDGs. The text of this report has been formulated to address only those areas affecting data quality.

TO: J. ORIENT  
SDGs: 320-10459-1 & 320-10572-1

PAGE 2

### **LABORATORY METHOD AND TRIP BLANK RESULTS**

The following compound was detected in the laboratory method blank at the following maximum concentration:

Compound	Maximum Concentration (ng/L)	Action Level (ng/L)
PFOA <sup>(1)</sup>	0.809	4.05

- <sup>(1)</sup> Maximum concentration detected in the laboratory method blank, MB 320-59712/1-A, from preparation batch # 320-59712 affecting the all samples from SDG 320-10459-1.

An action level of 5X the maximum contaminant concentration was established to evaluate blank contamination. Dilution factor and sample aliquot were taken into consideration during the application of the action level. The detected result reported for PFOA below the established action level in sample B7/10-GW03S was qualified as non-detected, (U).

### **NOTES**

EPA Region 1 worksheets were not applicable to this methodology and were not included in this tier II validation.

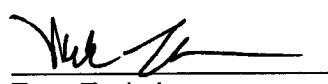
Detected results reported below the Limit of Quantitation (LOQ) but above the Detection Limit (DL) were qualified as estimated, (J). Non-detected results are reported to the Limit of Detection (LOD).

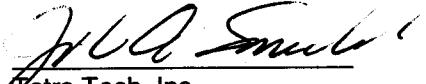
### **EXECUTIVE SUMMARY**

**Laboratory Performance:** PFOA was detected in the laboratory method blank.

**Other Factors Affecting Data Quality:** Detected results below the LOQ were estimated.

The data for these analyses were reviewed with reference to the EPA New England Environmental Data Review Supplement for Regional Data Review Elements Superfund Guidance/Procedures (April 2013), National Functional Guidelines for Organic Data Validation (January 2008), and the Department of Defense (DoD) document entitled, "Quality Systems Manual (QSM) for Environmental Laboratories" (July 2013). The text of this report has been formulated to address only those areas affecting data quality.

  
Tetra Tech, Inc.  
Michelle L. Allen  
Environmental Chemist

  
Tetra Tech, Inc.  
Joseph A. Samchuck  
Data Validation Manager

**Attachments:**

1. Appendix A - Qualified Analytical Results
2. Appendix B - Results as reported by the Laboratory
3. Appendix C - Support Documentation

**APPENDIX A**  
**QUALIFIED LABORATORY RESULTS**

**Qualifier Codes:**

- A = Lab Blank Contamination
- B = Field Blank Contamination
- C = Calibration Noncompliance (i.e., % RSDs, %Ds, ICVs, CCVs, RRFs, etc.)
- C01 = GC/MS Tuning Noncompliance
- D = MS/MSD Recovery Noncompliance
- E = LCS/LCSD Recovery Noncompliance
- F = Lab Duplicate Imprecision
- G = Field Duplicate Imprecision
- H = Holding Time Exceedance
- I = ICP Serial Dilution Noncompliance
- J = ICP PDS Recovery Noncompliance; MSA's r < 0.995
- K = ICP Interference - includes ICS % R Noncompliance
- L = Instrument Calibration Range Exceedance
- M = Sample Preservation Noncompliance
- N = Internal Standard Noncompliance
- N01 = Internal Standard Recovery Noncompliance Dioxins
- N02 = Recovery Standard Noncompliance Dioxins
- N03 = Clean-up Standard Noncompliance Dioxins
- O = Poor Instrument Performance (i.e., base-time drifting)
- P = Uncertainty near detection limit (< 2 x IDL for inorganics and <CRQL for organics)
- Q = Other problems (can encompass a number of issues; i.e. chromatography, interferences, etc.)
- R = Surrogates Recovery Noncompliance
- S = Pesticide/PCB Resolution
- T = % Breakdown Noncompliance for DDT and Endrin
- U = RPD between columns/detectors >40% for positive results determined via GC/HPLC
- V = Non-linear calibrations; correlation coefficient r < 0.995
- W = EMPC result
- X = Signal to noise response drop
- Y = Percent solids <30%
- Z = Uncertainty at 2 standard deviations is greater than sample activity
- Z1 = Tentatively Identified Compound considered presumptively present
- Z2 = Tentatively Identified Compound column bleed
- Z3 = Tentatively Identified Compound aldol condensate

PROJ_NO: 02063	NSAMPLE	B7/10-GW02		B7/10-GW03S		B7/10-GW06S		EFA-FD01-1014				
SDG: 320-10459-1	LAB_ID	320-10459-1		320-10459-2		320-10459-4		320-10459-6				
FRACTION: OS	SAMP_DATE	11/15/2014		11/15/2014		11/16/2014		11/16/2014				
MEDIA: WATER	QC_TYPE	NM		NM		NM		NM				
	UNITS	NG/L		NG/L		NG/L		NG/L				
	PCT_SOLIDS	0.0		0.0		0.0		0.0				
	DUP_OF							H123-GW02				
PARAMETER	RESULT	VQL	QLCD	RESULT	VQL	QLCD	RESULT	VQL	QLCD	RESULT	VQL	QLCD
PENTADECAFLUOROOCTANOIC ACID	10			2	U	A	20			16		
PERFLUOROOCTANE SULFONIC ACID	6.6			1.8	U		2.2			1.7	U	

PROJ_NO: 02063	NSAMPLE	H123-GW02	QRY-GW02			
SDG: 320-10459-1	LAB_ID	320-10459-5	320-10459-3			
FRACTION: OS	SAMP_DATE	11/16/2014	11/16/2014			
MEDIA: WATER	QC_TYPE	NM	NM			
	UNITS	NG/L	NG/L			
	PCT_SOLIDS	0.0	0.0			
	DUP_OF					
PARAMETER	RESULT	VQL	QLCD	RESULT	VQL	QLCD
PENTADECAFLUOROOCTANOIC ACID	17			1.8	U	
PERFLUOROOCTANE SULFONIC ACID	1.6	U		1.8	U	

<b>PROJ_NO:</b> 02063 <b>SDG:</b> 320-10572-1 <b>FRACTION:</b> OS <b>MEDIA:</b> WATER	<b>NSAMPLE</b>	EFA-GW01			EFA-GW02			EFA-GW03			EFA-GW09		
	<b>LAB_ID</b>	320-10572-1			320-10572-2			320-10572-3			320-10572-4		
	<b>SAMP_DATE</b>	11/18/2014			11/18/2014			11/19/2014			11/20/2014		
	<b>QC_TYPE</b>	NM			NM			NM			NM		
	<b>UNITS</b>	NG/L			NG/L			NG/L			NG/L		
	<b>PCT_SOLIDS</b>	0.0			0.0			0.0			0.0		
	<b>DUP_OF</b>												
PARAMETER	RESULT	VQL	QLCD	RESULT	VQL	QLCD	RESULT	VQL	QLCD	RESULT	VQL	QLCD	
PENTADECAFLUOROOCTANOIC ACID	1.8	J	P	1.5	U		11			31			
PERFLUOROOCTANE SULFONIC ACID	1.5	U		1.5	U		14			7.6			

**APPENDIX B**  
**RESULTS AS REPORTED BY THE LABORATORY**

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.: _____	
Client Sample ID: <u>B7/10-GW02</u>	Lab Sample ID: <u>320-10459-1</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_084.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/15/2014 12:05</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>448.35 (mL)</u>	Date Analyzed: <u>12/04/2014 20:45</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	10	H	2.2	1.7	0.83
1763-23-1	Perfluorooctane Sulfonate (PFOS)	6.6	H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	116		25-150
STL00990	13C4 PFOA	85		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.: _____	
Client Sample ID: <u>B7/10-GW03S</u>	Lab Sample ID: <u>320-10459-2</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_085.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/15/2014 15:20</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>427.75 (mL)</u>	Date Analyzed: <u>12/04/2014 21:07</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	2.0	J H	2.3	1.8	0.87
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.8	1.5

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	108		25-150
STL00990	13C4 PFOA	64		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.: _____	
Client Sample ID: <u>B7/10-GW06S</u>	Lab Sample ID: <u>320-10459-4</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_087.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 11:50</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>446.26 (mL)</u>	Date Analyzed: <u>12/04/2014 21:49</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	20	H	2.2	1.7	0.84
1763-23-1	Perfluorooctane Sulfonate (PFOS)	2.2	H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	108		25-150
STL00990	13C4 PFOA	67		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10459-1</u>
SDG No.:	
Client Sample ID: <u>EFA-FD01-1014</u>	Lab Sample ID: <u>320-10459-6</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_089.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/16/2014 00:00</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>12/03/2014 13:38</u>
Sample wt/vol: <u>447.90 (mL)</u>	Date Analyzed: <u>12/04/2014 22:31</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	16	H	2.2	1.7	0.84
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.7	U H	2.2	1.7	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	109		25-150
STL00990	13C4 PFOA	82		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Client Sample ID: H123-GW02 Lab Sample ID: 320-10459-5

Matrix: Water Lab File ID: 03DEC14A4A\_088.d

Analysis Method: WS-LC-0025 Date Collected: 11/16/2014 15:00

Extraction Method: 3535 Date Extracted: 12/03/2014 13:38

Sample wt/vol: 455.47 (mL) Date Analyzed: 12/04/2014 22:10

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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	17	H	2.2	1.6	0.82
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.6	U H	2.2	1.6	1.4

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	107		25-150
STL00990	13C4 PFOA	89		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento	Job No.: 320-10459-1
SDG No.:	
Client Sample ID: QR4-GW02	Lab Sample ID: 320-10459-3
Matrix: Water	Lab File ID: 03DEC14A4A_086.d
Analysis Method: WS-LC-0025	Date Collected: 11/16/2014 10:20
Extraction Method: 3535	Date Extracted: 12/03/2014 13:38
Sample wt/vol: 427.31(mL)	Date Analyzed: 12/04/2014 21:28
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.: 59779	Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	1.8	U H	2.3	1.8	0.88
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.8	U H	2.3	1.8	1.5

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	109		25-150
STL00990	13C4 PFOA	90		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10572-1</u>
SDG No.: _____	
Client Sample ID: <u>EFA-GW01</u>	Lab Sample ID: <u>320-10572-1</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_074.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/18/2014 11:00</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>11/24/2014 16:22</u>
Sample wt/vol: <u>500.25 (mL)</u>	Date Analyzed: <u>12/04/2014 17:13</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	1.8	J	2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	114		25-150
STL00990	13C4 PFOA	113		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: EFA-GW02 Lab Sample ID: 320-10572-2  
Matrix: Water Lab File ID: 03DEC14A4A\_075.d  
Analysis Method: WS-LC-0025 Date Collected: 11/18/2014 16:25  
Extraction Method: 3535 Date Extracted: 11/24/2014 16:22  
Sample wt/vol: 499.85 (mL) Date Analyzed: 12/04/2014 17:34  
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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	1.5	U	2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	105		25-150
STL00990	13C4 PFOA	97		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1  
SDG No.: \_\_\_\_\_  
Client Sample ID: EFA-GW03 Lab Sample ID: 320-10572-3  
Matrix: Water Lab File ID: 03DEC14A4A\_076.d  
Analysis Method: WS-LC-0025 Date Collected: 11/19/2014 11:10  
Extraction Method: 3535 Date Extracted: 11/24/2014 16:22  
Sample wt/vol: 502.05 (mL) Date Analyzed: 12/04/2014 17:56  
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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	11		2.0	1.5	0.74
1763-23-1	Perfluorooctane Sulfonate (PFOS)	14		2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	109		25-150
STL00990	13C4 PFOA	93		25-150

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10572-1</u>
SDG No.: _____	
Client Sample ID: <u>EFA-GW09</u>	Lab Sample ID: <u>320-10572-4</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_077.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected: <u>11/20/2014 10:40</u>
Extraction Method: <u>3535</u>	Date Extracted: <u>11/24/2014 16:22</u>
Sample wt/vol: <u>501.98 (mL)</u>	Date Analyzed: <u>12/04/2014 18:17</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	31		2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	7.6		2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	113		25-150
STL00990	13C4 PFOA	72		25-150

**APPENDIX C**  
**SUPPORT DOCUMENTATION**

TestAmerica Sacramento  
Target Compound Quantitation Report

Data File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_087.d  
 Lims ID: 320-10459-A-4-B Lab Sample ID: 320-10459-4 RRF = 1.6697  
 Client ID: B7/10-GW06S  
 Sample Type: Client  
 Inject. Date: 04-Dec-2014 21:49:30 ALS Bottle#: 27 Worklist Smp#: 51  
 Injection Vol: 15.0 ul Dil. Factor: 1.0000  
 Sample Info: 320-10459-a-4-b 59712 Extraction: 446.26 ml  
 Misc. Info.: AcquityBEH 1.7u C18, 150x3.0mm, T=50C,A(4967-113F)  
 Operator ID: JRB Instrument ID: A4 1 ml  
 Method: \\Sacchrom\ChromData\A4\20141204-17820.b\PFAC\_A4.m  
 Limit Group: LC PFC ICAL df = 1  
 Last Update: 05-Dec-2014 10:03:21 Calib Date: 04-Dec-2014 05:51:34  
 Integrator: Picker  
 Quant Method: Isotopic Dilution Quant By: Initial Calibration  
 Last ICal File: \\Sacchrom\ChromData\A4\20141204-17820.b\03DEC14A4A\_046.d  
 Column 1: Acquity BEH C18 ( 2.10 mm) Det: F1:MRM  
 Process Host: XAWRK022

Signal	RT	EXP RT	DLT RT	REL RT	Response	Amount ng/ml	Ratio(Limits)	%Rec	S/N	Flags
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## D 12 13C4 PFOA

416.5 > 371.6 10.642 10.628 0.014 1138168 33.6 67.3 1719

13 Perfluorooctanoic acid

→ 412.8 > 368.8 10.642 10.628 0.014 1.000 340395 8.96 3.82(0.00-0.00) 107  
 412.8 > 168.7 10.642 10.628 0.014 1.000 89169 61.5

## D 16 13C4 PFOS

502.4 > 79.7 11.602 11.595 0.007 2338394 51.9 108 3471

15 Perfluorooctanoic Sulfonate

498.9 > 79.7 11.602 11.595 0.007 1.000 87815 17.4  
 498.9 > 98.7 11.602 11.595 0.007 1.000 51215 31.1

$$\begin{aligned}
 & \frac{340395 \times 30 \text{ ng/ml}}{1138168 \times 1.6697} \times \frac{1 \text{ ml}}{446.26} \times \frac{1000 \text{ ml}}{1 \text{ L}} \\
 & = 20.07 \text{ ng/L}
 \end{aligned}$$

**NAS BRUNSWICK**  
**WATER DATA**  
**320-10459-1**

FRACTION	CHEMICAL	H123-GW02	UNITS	EPA-FD01-1014	RPD	D
OS	PENTADECAFLUOROOCTANOIC ACID	17 J	NG/L	16 J	6.06	1.00

**Current RPD Quality Control Limit: 30 %.**

**Shaded cells indicate RPDs that exceed the applicable quality control limit.**



Tetra Tech, Inc.

## CHAIN OF CUSTODY

NUMBER No. 0533

PAGE 1 OF 1

PROJECT NO: 112502063		FACILITY: NAS Brunswick	PROJECT MANAGER Jeff Orient	PHONE NUMBER 412 921 7090	LABORATORY NAME AND CONTACT: Test America						
SAMPLERS (SIGNATURE) <i>T. Smith</i> <i>R. Evans</i>		FIELD OPERATIONS LEADER Tim Evans	PHONE NUMBER 412 921 7090	CARRIER/WAYBILL NUMBER 8066 1897 7115	ADDRESS 880 Riverside Pkwy						
STANDARD TAT <input checked="" type="checkbox"/> RUSH TAT <input type="checkbox"/> <input type="checkbox"/> 24 hr. <input type="checkbox"/> 48 hr. <input type="checkbox"/> 72 hr. <input type="checkbox"/> 7 day <input type="checkbox"/> 14 day		CONTAINER TYPE PLASTIC (P) or GLASS (G)		CITY, STATE West Sacramento CA							
DATE YEAR 2014	TIME	SAMPLE ID	LOCATION ID	TOP DEPTH (FT)	BOTTOM DEPTH (FT)	MATRIX (GW, SO, SW, SD, QC, ETC.)	COLLECTION METHOD GRAB (G) COMP (C)	NO. OF CONTAINERS	PRESERVATIVE USED	TYPE OF ANALYSIS PFDA/PFOS	COMMENTS
11/15	1205	B7/10 - Gw02	B7/10- mw02	-	-	GW	G	1	1		
↓	1520	B7/10 - Gw03S	B7/10- mw03S	-	-	GW	G	1	1		
11/16	1020	QRY - Gw02	QRY- mw02	-	-	GW	G	1	1		
↓	1150	B7/10 - Gw06S	B7/10- mw06S	-	-	GW	G	1	1		
↓	1500	H123 - Gw02	H123- mw02	-	-	GW	G	1	1		
↓	0000	PFDA - FD01-1014	Field Duplicate	-	-	GW	G	1	1		
1. RELINQUISHED BY <i>T. Smith</i>		DATE 11/17/14	TIME 17:30	1. RECEIVED BY FedEx		DATE	TIME				
2. RELINQUISHED BY		DATE	TIME	2. RECEIVED BY <i>Mr. R. Evans</i>		DATE 11/17/14	TIME 19:30				
3. RELINQUISHED BY		DATE	TIME	3. RECEIVED BY		DATE	TIME				
COMMENTS											

320-10459 Chain of Custody

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11/17/2014

DISTRIBUTION:

WHITE (ACCOMPANIES SAMPLE)

YELLOW (FIELD COPY)

PINK (FILE COPY)

4/02R

FORM NO. TINUS-001

## Login Sample Receipt Checklist

Client: Tetra Tech, Inc.

Job Number: 320-10459-1

**Login Number: 10459**

**List Source: TestAmerica Sacramento**

**List Number: 1**

**Creator: Sadler, Jeremy**

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.	N/A	
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?	True	
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	N/A	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	

## CASE NARRATIVE

**Client: Tetra Tech, Inc.**

**Project: NAS Brunswick Maine 13-CTO WE09**

**Report Number: 320-10459-1**

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

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

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

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

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

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

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

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

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

### **RECEIPT**

The samples were received on 11/18/2014 9:50 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 0.5° C.

### **PFOA/PFOS**

#### **Method WS-LC-0025:**

The following samples were re-extracted past the 7 day extraction hold time due to contamination in the method blank (MB) and laboratory control sample (LCS) for the original prep batch: B7/10-GW02 (320-10459-1), B7/10-GW03S (320-10459-2), B7/10-GW06S (320-10459-4), EFA-FD01-1014 (320-10459-6), H123-GW02 (320-10459-5), QR4-GW02 (320-10459-3). There is no prescribed regulatory holding time for PFOA and PFOS. The 7 day extraction holding time is based on general EPA convention for the holding time of extractable organic compounds in water.

Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate/sample duplicate (MS/MSD/DUP) associated with batch 320-59712.

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

## Method Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

Method	Method Description	Protocol	Laboratory
WS-LC-0025	Perfluorinated Hydrocarbons	TAL SOP	TAL SAC

**Protocol References:**

TAL SOP = TestAmerica Laboratories, Standard Operating Procedure

**Laboratory References:**

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

## Sample Summary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-10459-1	B7/10-GW02	Water	11/15/14 12:05	11/18/14 09:50
320-10459-2	B7/10-GW03S	Water	11/15/14 15:20	11/18/14 09:50
320-10459-3	QR4-GW02	Water	11/16/14 10:20	11/18/14 09:50
320-10459-4	B7/10-GW06S	Water	11/16/14 11:50	11/18/14 09:50
320-10459-5	H123-GW02	Water	11/16/14 15:00	11/18/14 09:50
320-10459-6	EFA-FD01-1014	Water	11/16/14 00:00	11/18/14 09:50

## Definitions/Glossary

Client: Tetra Tech, Inc.

Project/Site: NAS Brunswick Maine 13-CTO WE09

TestAmerica Job ID: 320-10459-1

### Qualifiers

#### LCMS

Qualifier	Qualifier Description
H	Sample was prepped or analyzed beyond the specified holding time
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.
D	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)

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFOA #	PFOS #
B7/10-GW02	320-10459-1	85	116
B7/10-GW03S	320-10459-2	64	108
QR4-GW02	320-10459-3	90	109
B7/10-GW06S	320-10459-4	67	108
H123-GW02	320-10459-5	89	107
EFA-FD01-1014	320-10459-6	82	109
	MB 320-59712/1-A	109	111
	LCS 320-59712/2-A	98	101

PFOA = 13C4 PFOA

PFOS = 13C4 PFOS

QC LIMITS

25-150

25-150

# Column to be used to flag recovery values

FORM II WS-LC-0025

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Lab File ID: 03DEC14A4A\_082.d Lab Sample ID: MB 320-59712/1-A

Matrix: Water Date Extracted: 12/03/2014 13:38

Instrument ID: A4 Date Analyzed: 12/04/2014 20:03

Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-59712/2-A	03DEC14A4A_083.d	12/04/2014 20:24
B7/10-GW02	320-10459-1	03DEC14A4A_084.d	12/04/2014 20:45
B7/10-GW03S	320-10459-2	03DEC14A4A_085.d	12/04/2014 21:07
QR4-GW02	320-10459-3	03DEC14A4A_086.d	12/04/2014 21:28
B7/10-GW06S	320-10459-4	03DEC14A4A_087.d	12/04/2014 21:49
H123-GW02	320-10459-5	03DEC14A4A_088.d	12/04/2014 22:10
EFA-FD01-1014	320-10459-6	03DEC14A4A_089.d	12/04/2014 22:31

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
 SDG No.: \_\_\_\_\_  
 Client Sample ID: \_\_\_\_\_ Lab Sample ID: MB 320-59712/1-A  
 Matrix: Water Lab File ID: 03DEC14A4A\_082.d  
 Analysis Method: WS-LC-0025 Date Collected: \_\_\_\_\_  
 Extraction Method: 3535 Date Extracted: 12/03/2014 13:38  
 Sample wt/vol: 500.00 (mL) Date Analyzed: 12/04/2014 20:03  
 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.: 59779 Units: ng/L

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	0.809	J	2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	111		25-150
STL00990	13C4 PFOA	109		25-150

FORM III  
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low Lab File ID: 03DEC14A4A\_083.d

Lab ID: LCS 320-59712/2-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ng/L)	LCS CONCENTRATION (ng/L)	LCS % REC	QC LIMITS REC	#
Perfluorooctanoic acid (PFOA)	40.0	50.4	126	60-140	
Perfluorooctane Sulfonate (PFOS)	38.2	39.7	104	60-140	
13C4 PFOS	95.6	96.3	101	25-150	
13C4 PFOA	100	98.0	98	25-150	

# Column to be used to flag recovery and RPD values

FORM III WS-LC-0025

## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Instrument ID: A4

Start Date: 12/04/2014 03:44

Analysis Batch Number: 59779

End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
STD 320-59779/4 IC		12/04/2014 03:44	1	03DEC14A4A_040. d	Acquity 2.1(mm)
STD 320-59779/5 IC		12/04/2014 04:05	1	03DEC14A4A_041. d	Acquity 2.1(mm)
STD 320-59779/6 IC		12/04/2014 04:26	1	03DEC14A4A_042. d	Acquity 2.1(mm)
STD 320-59779/7 IC		12/04/2014 04:47	1	03DEC14A4A_043. d	Acquity 2.1(mm)
STD 320-59779/8 IC		12/04/2014 05:09	1	03DEC14A4A_044. d	Acquity 2.1(mm)
STD 320-59779/9 IC		12/04/2014 05:30	1	03DEC14A4A_045. d	Acquity 2.1(mm)
STD 320-59779/10 IC		12/04/2014 05:51	1	03DEC14A4A_046. d	Acquity 2.1(mm)
CCV 320-59779/20		12/04/2014 09:23	1		Acquity 2.1(mm)
ICV 320-59779/22		12/04/2014 10:23	1	03DEC14A4A_058. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 11:50	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:11	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:37	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:59	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:20	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:41	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:02	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:24	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:44	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:06	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:27	1		Acquity 2.1(mm)
CCV 320-59779/34		12/04/2014 15:48	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:10	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:31	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:52	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:13	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:34	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 17:56	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:17	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:38	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:59	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 19:20	1		Acquity 2.1(mm)
CCV 320-59779/45		12/04/2014 19:42	1	03DEC14A4A_081. d	Acquity 2.1(mm)
MB 320-59712/1-A		12/04/2014 20:03	1	03DEC14A4A_082. d	Acquity 2.1(mm)
LCS 320-59712/2-A		12/04/2014 20:24	1	03DEC14A4A_083. d	Acquity 2.1(mm)
320-10459-1	B7/10-GW02	12/04/2014 20:45	1	03DEC14A4A_084. d	Acquity 2.1(mm)
320-10459-2	B7/10-GW03S	12/04/2014 21:07	1	03DEC14A4A_085. d	Acquity 2.1(mm)
320-10459-3	QR4-GW02	12/04/2014 21:28	1	03DEC14A4A_086. d	Acquity 2.1(mm)
320-10459-4	B7/10-GW06S	12/04/2014 21:49	1	03DEC14A4A_087. d	Acquity 2.1(mm)

## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1  
SDG No.:  
Instrument ID: A4 Start Date: 12/04/2014 03:44  
Analysis Batch Number: 59779 End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
320-10459-5	H123-GW02	12/04/2014 22:10	1	03DEC14A4A_088. d	Acquity 2.1(mm)
320-10459-6	EFA-FD01-1014	12/04/2014 22:31	1	03DEC14A4A_089. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 22:53	20		Acquity 2.1(mm)
ZZZZZ		12/04/2014 23:14	20		Acquity 2.1(mm)
ZZZZZ		12/04/2014 23:35	20		Acquity 2.1(mm)
CCV 320-59779/57		12/04/2014 23:56	1	03DEC14A4A_093. d	Acquity 2.1(mm)

FORM VI  
LCMS INITIAL CALIBRATION DATA  
EXTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-10459-1 Analy Batch No.: 59779  
SDG No.: \_\_\_\_\_  
Instrument ID: A4 GC Column: Acuity ID: 2.1(mm) Heated Purge: (Y/N) N  
Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4		B	M1	M2								
13C4 PFBA	5268.6 3806.7	5812.9 3203.9	5254.9 2254.3	4338.2	Ave		4277.06429				30.0		50.0			
13C5 PFPeA	9419.7 7332.0	9580.9 5721.7	8177.5 4869.4	8903.4	Ave		7714.90000				24.0		50.0			
13C2 PFHxA	21470 18141	22784 16317	24835 11422	18449	Ave		19059.8143				23.0		50.0			
13C4-PFHxA	21643 17882	21390 11698	21748 10393	19442	Ave		17742.3500				27.0		50.0			
18O2 PFHxS	53128 47816	52233 33678	50966 25618	51213	Ave		44950.3322				24.0		50.0			
13C4 PFOA	41322 31379	41400 24644	40475 19047	38546	Ave		33830.5286				27.0		50.0			
13C4 PFOS	54398 44130	54385 33330	56166 23197	50088	Ave		45099.0735				28.0		50.0			
13C5 PFNA	44879 37219	45410 30155	48879 22012	41831	Ave		38626.2857				25.0		50.0			
13C2 PFDA	63235 48806	61430 37175	59297 32265	58211	Ave		51488.2929				24.0		50.0			
13C8 FOSA	150505 147550	155647 126255	158874 106829	155831	Ave		143069.894				14.0		50.0			
13C2 PFUnA	66511 59868	66532 58833	66894 69853	64772	Ave		64751.6643				6.2		50.0			
13C2 PFDoA	76991 59932	66196 56050	76655 52807	68862	Ave		65356.2357				15.0		50.0			
13C2-PFTeDA	77915 69707	76286 61208	80240 58365	76834	Ave		71507.9143				12.0		50.0			
13C2-PFHxDA	96546 111182	105842 105984	112387 93820	103841	Ave		104228.936				6.6		50.0			

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento	Job No.: 320-10459-1	Analy Batch No.: 59779
SDG No.:		
Instrument ID: A4	GC Column: Acquity ID: 2.1 (mm)	Heated Purge: (Y/N) N
Calibration Start Date: 12/04/2014 03:44	Calibration End Date: 12/04/2014 05:51	Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorobutanoic acid (PFBA)	10678 16623	22829 12169	21990	19527	19477	AveID		2.3514				25.0		35.0			
Perfluoropentanoic acid (PFPeA)	8732.0 11185	14944 9458.0	14198	14832	13130	AveID		1.6538				21.0		35.0			
Perfluorobutane Sulfonate (PFBS)	14683 39320	46758 32165	49055	48250	67174	AveID		0.9863				37.0		50.0			
Perfluorohexanoic acid (PFHxA)	9952.0 16870	21004 12521	26818	19955	20968	AveID		0.9761				24.0		35.0			
PFPoS (Perflouro-1-pentanesulfonate)	12755 32980	40229 24854	44086	42428	40598	AveID		0.7860				32.0		50.0			
Perfluoroheptanoic acid (PFHpA)	14724 21344	28535 18892	35883	33256	29606	AveID		1.5247				27.0		35.0			
Perfluorohexane Sulfonate (PFHxS)	24632 33051	52072 26648	53418	49529	45799	AveID		0.9222				22.0		35.0			
Perfluorooctanoic acid (PFOA)	54920 43078	74889 33366	68131	62497	54775	AveID		1.6697				9.7		35.0			
Perfluoro-1-heptanesulfonate (PFHpS)	12752 33076	45349 24051	51122	49900	46278	AveID		0.8647				33.0		50.0			
Perfluorooctane Sulfonate (PFOS)	143245 49883	124703 36721	92010	73648	66987	AveID		1.8046				26.0		35.0			
Perfluorononanoic acid (PFNA)	22096 39342	60944 27244	61254	52411	44283	AveID		1.1532				26.0		35.0			
PFNS (Perflouro-1-nananesulfonate)	16554 25936	40576 18531	47018	41535	37537	AveID		0.7349				26.0		50.0			
Perfluorodecanoic acid (PFDA)	16398 41090	58072 33184	67070	63478	54711	AveID		0.9544				33.0		35.0			
Perfluorooctane Sulfonamide (FOSA)	60576 132015	157092 116050	155459	165246	157618	AveID		0.9501				26.0		35.0			
Perfluorodecane sulfonate (PFDS)	10237 23088	37158 16598	42628	36833	33233	AveID		0.6467				32.0		50.0			
Perfluoroundecanoic acid (PFUnA)	25356 49828	75233 40491	72252	72984	62389	AveID		0.8839				34.0		35.0			
Perfluorododecanoic acid (PFDoA)	25100 62748	61887 50239	92219	78804	73401	AveID		0.9863				32.0		35.0			
PFDoS (Perflouro-1-dodecanesulfonate)	7778.9 23770	32535 17735	43717	40638	35319	AveID		0.6584				36.0		50.0			
Perfluorotridecanoic Acid (PFTriA)	20248 45661	65903 40208	67545	66220	62650	AveID		0.8175				32.0		50.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

Analy Batch No.: 59779

SDG No.:

Instrument ID: A4

GC Column: Acquity ID: 2.1(mm)

Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44

Calibration End Date: 12/04/2014 05:51

Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorotetradecanoic acid (PFTeA)	15018 30898	36572 27633	43161	42497	38214	AveID		0.5200			29.0		50.0				
Perfluoro-n-hexadecanoic acid (PFHxDA)	44032 92953	98406 78170	100981	86858	96066	AveID		1.3398			27.0		50.0				
Perfluoro-n-octadecanoic acid (PFODA)	22762 104163	94742 95748	99443	104957	102973	AveID		1.4197			38.0		50.0				

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Lab Sample ID: ICV 320-59779/22

Calibration Date: 12/04/2014 10:23

Instrument ID: A4

Calib Start Date: 12/04/2014 03:44

GC Column: Acquity ID: 2.10 (mm)

Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_058.d

Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	2.637		56.1	50.0	12.2	50.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	1.886		57.0	50.0	14.0	50.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.236		55.5	44.3	25.3	50.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.190		61.0	50.0	21.9	50.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.160		70.8	50.0	41.7	50.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	1.119		57.4	47.3	21.4	50.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9109		50.1	47.6	5.3	50.0
Perfluoroctanoic acid (PFOA)	AveID	1.670	1.812		54.2	50.0	8.5	50.0
Perfluoroctane Sulfonate (PFOS)	AveID	1.805	1.655		43.8	47.8	-8.3	50.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.206		52.3	50.0	4.6	50.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.211		63.4	50.0	26.8	50.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	0.9856		51.9	50.0	3.7	50.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7475		55.8	48.3	15.6	50.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.141		64.5	50.0	29.1	50.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.176		59.6	50.0	19.2	50.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9516		58.2	50.0	16.4	50.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.6011		57.8	50.0	15.6	50.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.644		61.3	50.0	22.7	50.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.763		62.1	50.0	24.2	50.0

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCV 320-59779/45 Calibration Date: 12/04/2014 19:42

Instrument ID: A4 Calib Start Date: 12/04/2014 03:44

GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_081.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	3.794		32.3	20.0	61.4*	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	2.282		27.6	20.0	38.0	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.023		18.3	17.7	3.7	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.252		25.6	20.0	28.2	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.068		27.1	20.0	35.7	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	0.998		20.5	18.9	8.2	40.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9528		21.0	19.0	10.2	40.0
Perfluoroctanoic acid (PFOA)	AveID	1.670	1.850		22.2	20.0	10.8	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.488		15.8	19.1	-17.6	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.142		19.8	20.0	-1.0	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.018		21.3	20.0	6.6	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.009		21.2	20.0	6.2	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7121		21.2	19.3	10.1	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.135		25.7	20.0	28.4	40.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.182		24.0	20.0	19.9	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9328		22.8	20.0	14.1	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5912		22.7	20.0	13.7	40.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.503		22.4	20.0	12.2	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.589		22.4	20.0	11.9	40.0

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCV 320-59779/57 Calibration Date: 12/04/2014 23:56

Instrument ID: A4 Calib Start Date: 12/04/2014 03:44

GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_093.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	3.249		69.1	50.0	38.2	40.0
Perfluoropentanoic acid (FFPeA)	AveID	1.654	2.193		66.3	50.0	32.6	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.639		73.5	44.2	66.2*	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.185		60.7	50.0	21.4	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	1.987		65.1	50.0	30.3	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	1.045		53.6	47.3	13.3	40.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9556		52.6	47.6	10.5	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.657		49.6	50.0	-0.7	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.512		40.0	47.8	-16.2	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.219		52.8	50.0	5.7	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.069		56.0	50.0	12.0	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.050		55.3	50.0	10.5	40.0
Perfluorododecane sulfonate (PFDS)	AveID	0.6467	0.6689		49.8	48.2	3.4	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.054		59.6	50.0	19.3	40.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.179		59.8	50.0	19.5	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9404		57.5	50.0	15.0	40.0
Perfluorotetradecanoic acid (FFTTeA)	AveID	0.5200	0.5983		57.5	50.0	15.1	40.0
Perfluoro-n-hexadecanoic acid (FFHxDA)	AveID	1.340	1.598		59.6	50.0	19.3	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.784		62.8	50.0	25.7	40.0

## LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento

Job No.: 320-10459-1

SDG No.:

Batch Number: 59712

Batch Start Date: 12/03/14 13:38

Batch Analyst: Reed, Jonathan E

Batch Method: 3535

Batch End Date: 12/03/14 19:10

Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount	LCMPFCSU 00010	LCPFCSP 00014		
MB 320-59712/1		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL			
LCS 320-59712/2		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL	20 uL		
320-10459-A-1	B7/10-GW02	3535, WS-LC-0025	T	448.35 mL	1.00 mL	50 uL			
320-10459-A-2	B7/10-GW03S	3535, WS-LC-0025	T	427.75 mL	1.00 mL	50 uL			
320-10459-A-3	QR4-GW02	3535, WS-LC-0025	T	427.31 mL	1.00 mL	50 uL			
320-10459-A-4	B7/10-GW06S	3535, WS-LC-0025	T	446.26 mL	1.00 mL	50 uL			
320-10459-A-5	H123-GW02	3535, WS-LC-0025	T	455.47 mL	1.00 mL	50 uL			
320-10459-A-6	EFA-FD01-1014	3535, WS-LC-0025	T	447.90 mL	1.00 mL	50 uL			

Batch Notes	
Balance ID	QA-070
H2O Lot used	11/24/14
Pipette ID	EC15219
Analyst who added reagent	JER
SU Reagent Drop	JER
SU Reagent Drop Witness	SNE
Solvent Lot #	0.3%NH4OH-Me_00006
Solvent Name	0.3%NH4OH/MeOH
SOP Number	WS-LC-0005
SPE Cartridge Type	WAX 500mg
Solid Phase Extraction Disk Lot Number	002334148A

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.

WS-LC-0025

Page 1 of 1

## HPLC/LCMS Data Review Checklist

Job Number(s): 8957, 9021, 10572, 10114,  
10541, 10459, Work List ID(s): 17820

Extraction Batch: 51087, 58153, 58090, 59718, Analysis Batch(es): 59779

Delivery Rank 2, 4 Due Date: Varies

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 $\frac{1}{2}$ the reporting limit as described in CA-Q-S-005?			/
• All curve points show calculated concentrations.	/	/	
3. Peaks correctly ID'd by data system.	/	/	
5. Tune check frequency & criteria met and Tune check report attached.			/
B. QA/QC			
1. Are all QC samples properly linked in TALS?	/	/	
2. Method blank, LCS/LCSD and MS/SD frequencies met.	/	/	
3. LCS/LCSD and MB data are within control limits. If not, NCM is present.	/	/	
4. Are MS/MSD recoveries and RPD within control limits?			/
5. Holding Times were met for prep and analytical. <i>In House samples + NCM</i>	/	/	
6. IS/Surrogate recoveries meet criteria or properly noted.	/	/	
C. Sample Analysis			
1. Was correct analysis performed and were project instructions followed?	/	/	
2. If required, are compounds within RT windows?	/	/	
3. If required, are positive hits confirmed and >40% RPD flagged?			/
4. Manual Integrations reviewed and appropriate.			
5. All analytes correctly reported. (Primary, secondary, acceptable status)	/	/	
6. Correct reporting limits used. (based on client request, prep factors, and dilutions)	/	/	
D. Documentation			
1. Are all non-conformances documented/attached? NCM#	/	/	
2. Do results make sense (e.g. dilutions, etc.)?	/	/	
3. Have all flags been reviewed for appropriateness?	/	/	
4. For level 3 and 4 reports, have forms and raw data been reviewed?	/	/	
5. Was QC Checker run for this job?	/	/	

\*Upon completion of this checklist, the reviewer must scan and attach the checklist to the TALS job.

1<sup>st</sup> Level (Analyst): JRB Date: 12/5/14

2<sup>nd</sup> Level Reviewer: JWJ Date: 12/12/14

*Act 12*

# Solid SW-846-3500 Analysis Sheet

(To Accompany Samples to Instruments)

Batch Number: 320-59087

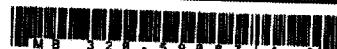
Method Code: 320-Shake\_Bath\_14D-320

Analyst: Reed, Jonathan E

Batch Open: 11/25/2014 2:06:49PM

Batch End: 11/26/2014 8:40:00PM

## Shake Extraction with Ultrasonic Bath Extraction

	Input Sample Lab ID (Analytical Method)	SDG	Initial Amount	Final Amount	Due Date	Analytical TAT	Div Rank	Comments	Output Sample Lab ID
1	MB-320-59087/1 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		
2	LCS-320-59087/2 N/A	N/A	5.00 g	1.00 mL	N/A	N/A	N/A		
3	320-10515-B-1 (PFC_IDA)	10502757.5501080 1	5.09 g	1.00 mL	12/9/14	12_Days	4		
4	320-10515-B-1-MS (PFC_IDA)	10502757.5501080 1	5.13 g	1.00 mL	12/9/14	12_Days	4		
5	320-10515-B-1-MSD (PFC_IDA)	10502757.5501080 1	5.09 g	1.00 mL	12/9/14	12_Days	4		
6	320-8957-G-1 (PFC_IDA)	N/A	5.00 g	1.00 mL	9/4/14	12_Days	2		
7	320-9021-W-1 (PFC_IDA)	N/A	5.00 g	1.00 mL	9/9/14	12_Days	2		

### Batch Notes

Blank Sand Lot # 142675

Filter Lot # NA

Millipore Water Dispense Date 11/24/14

Person's name who witnessed reagent drop SNE

SPE Cartridge Lot # 014933351A

SPE Cartridge Type WAX 150mg

Batch Comment SOLVENT: 0.3%NH4OH-Me\_00006

12/12/2014

Printed: 11/26/2014



Tetra Tech, Inc.

## CHAIN OF CUSTODY

NUMBER No. 0537

PAGE 1 OF 1

PROJECT NO: 112G-020C 3	FACILITY: EFA/ NAS Brunswick	PROJECT MANAGER Jeff Orient	PHONE NUMBER 412 921-7090	LABORATORY NAME AND CONTACT: Test America							
SAMPLERS (SIGNATURE) 	FIELD OPERATIONS LEADER Tim Evans	PHONE NUMBER 412 921 7090	ADDRESS 880 Riverside Pkwy								
	CARRIER/WAYBILL NUMBER FedEx 8065 18977159		CITY, STATE West Sacramento CA								
STANDARD TAT <input checked="" type="checkbox"/> RUSH TAT <input type="checkbox"/> <input type="checkbox"/> 24 hr. <input type="checkbox"/> 48 hr. <input type="checkbox"/> 72 hr. <input type="checkbox"/> 7 day <input type="checkbox"/> 14 day		CONTAINER TYPE PLASTIC (P) or GLASS (G)	P								
DATE YEAR 2014	TIME	SAMPLE ID	LOCATION ID	TOP DEPTH (FT)	BOTTOM DEPTH (FT)	MATRIX (GW, SO, SW, SD, QC, ETC.)	COLLECTION METHOD GRAB (G) COMP (C)	No. OF CONTAINERS	PRESERVATIVE USED	TYPE OF ANALYSIS PFOA/PFOS	COMMENTS
11/18	1100	EFA-Gw01	EFA-01	-	-	GW	G	1	1		
11/18	1625	EFA-Gw02	EFA-02	-	-	GW	G	1	1		
11/19	1110	EFA-Gw03	EFA-03	-	-	GW	G	1	1		
11/20	1040	EFA-Gw09	EFA-09	-	-	GW	G	1	1		
Page 349 of 350											
 320-10572 Chain of Custody											
1. RELINQUISHED BY 	DATE 11/20/14	TIME 19:30	1. RECEIVED BY FedEx	DATE 11/20/14	TIME 19:30						
2. RELINQUISHED BY	DATE	TIME	2. RECEIVED BY Km Nelson	DATE 11/20/14	TIME 19:35						
3. RELINQUISHED BY	DATE	TIME	3. RECEIVED BY	DATE	TIME 0:75						
COMMENTS											

DISTRIBUTION:

WHITE (ACCOMPANIES SAMPLE)

YELLOW (FIELD COPY)

PINK (FILE COPY)

4/02R

FORM NO. TiNUS-001

## Login Sample Receipt Checklist

Client: Tetra Tech, Inc.

Job Number: 320-10572-1

**Login Number: 10572**

**List Source: TestAmerica Sacramento**

**List Number: 1**

**Creator: Nelson, Kym D**

Question	Answer	Comment
Radioactivity wasn't checked or is </= background as measured by a survey meter.	True	
The cooler's custody seal, if present, is intact.	True	
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.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	N/A	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	

## CASE NARRATIVE

**Client: Tetra Tech, Inc.**

**Project: Naval Air Station Brunswick Maine**

**Report Number: 320-10572-1**

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

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

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

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

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

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

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

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

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

### **RECEIPT**

The samples were received on 11/21/2014 9:15 AM; the samples arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 0.7° C.

### **PFOA/PFOS**

Insufficient sample volume was available to perform a matrix spike/matrix spike duplicate/sample duplicate (MS/MSD/DUP) associated with batch 320-58953.

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

## Method Summary

Client: Tetra Tech, Inc.

Project/Site: Naval Air Station Brunswick Maine

TestAmerica Job ID: 320-10572-1

Method	Method Description	Protocol	Laboratory
WS-LC-0025	Perfluorinated Hydrocarbons	TAL SOP	TAL SAC

**Protocol References:**

TAL SOP = TestAmerica Laboratories, Standard Operating Procedure

**Laboratory References:**

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

## Sample Summary

Client: Tetra Tech, Inc.

Project/Site: Naval Air Station Brunswick Maine

TestAmerica Job ID: 320-10572-1

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
320-10572-1	EFA-GW01	Water	11/18/14 11:00	11/21/14 09:15
320-10572-2	EFA-GW02	Water	11/18/14 16:25	11/21/14 09:15
320-10572-3	EFA-GW03	Water	11/19/14 11:10	11/21/14 09:15
320-10572-4	EFA-GW09	Water	11/20/14 10:40	11/21/14 09:15

## Definitions/Glossary

Client: Tetra Tech, Inc.

Project/Site: Naval Air Station Brunswick Maine

TestAmerica Job ID: 320-10572-1

### 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.
D	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)

FORM II  
LCMS SURROGATE RECOVERY

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low

GC Column (1): Acquity ID: 2.1 (mm)

Client Sample ID	Lab Sample ID	PFOA #	PFOS #
EFA-GW01	320-10572-1	113	114
EFA-GW02	320-10572-2	97	105
EFA-GW03	320-10572-3	93	109
EFA-GW09	320-10572-4	72	113
	MB 320-58953/1-A	105	106
	LCS 320-58953/2-A	97	99

PFOA = 13C4 PFOA

PFOS = 13C4 PFOS

QC LIMITS

25-150

25-150

# Column to be used to flag recovery values

FORM II WS-LC-0025

FORM IV  
LCMS METHOD BLANK SUMMARY

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.: \_\_\_\_\_

Lab File ID: 03DEC14A4A\_066.d

Lab Sample ID: MB 320-58953/1-A

Matrix: Water

Date Extracted: 11/24/2014 16:22

Instrument ID: A4

Date Analyzed: 12/04/2014 14:24

Level: (Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 320-58953/2-A	03DEC14A4A_067.d	12/04/2014 14:44
EFA-GW01	320-10572-1	03DEC14A4A_074.d	12/04/2014 17:13
EFA-GW02	320-10572-2	03DEC14A4A_075.d	12/04/2014 17:34
EFA-GW03	320-10572-3	03DEC14A4A_076.d	12/04/2014 17:56
EFA-GW09	320-10572-4	03DEC14A4A_077.d	12/04/2014 18:17

FORM I  
LCMS ORGANICS ANALYSIS DATA SHEET

Lab Name: <u>TestAmerica Sacramento</u>	Job No.: <u>320-10572-1</u>
SDG No.:	
Client Sample ID:	Lab Sample ID: <u>MB 320-58953/1-A</u>
Matrix: <u>Water</u>	Lab File ID: <u>03DEC14A4A_066.d</u>
Analysis Method: <u>WS-LC-0025</u>	Date Collected:
Extraction Method: <u>3535</u>	Date Extracted: <u>11/24/2014 16:22</u>
Sample wt/vol: <u>500.00 (mL)</u>	Date Analyzed: <u>12/04/2014 14:24</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>59779</u>	Units: <u>ng/L</u>

CAS NO.	COMPOUND NAME	RESULT	Q	LOQ	LOD	DL
335-67-1	Perfluorooctanoic acid (PFOA)	1.5	U	2.0	1.5	0.75
1763-23-1	Perfluorooctane Sulfonate (PFOS)	1.5	U	2.0	1.5	1.3

CAS NO.	ISOTOPE DILUTION	%REC	Q	LIMITS
STL00991	13C4 PFOS	106		25-150
STL00990	13C4 PFOA	105		25-150

FORM III  
LCMS LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1

SDG No.: \_\_\_\_\_

Matrix: Water Level: Low Lab File ID: 03DEC14A4A\_067.d

Lab ID: LCS 320-58953/2-A Client ID: \_\_\_\_\_

COMPOUND	SPIKE ADDED (ng/L)	LCS CONCENTRATION (ng/L)	LCS % REC	QC LIMITS REC	#
Perfluorooctanoic acid (PFOA)	40.0	46.3	116	60-140	
Perfluorooctane Sulfonate (PFOS)	38.2	35.8	94	60-140	
13C4 PFOS	95.6	95.0	99	25-150	
13C4 PFOA	100	96.6	97	25-150	

# Column to be used to flag recovery and RPD values

FORM III WS-LC-0025

## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.:

Instrument ID: A4

Start Date: 12/04/2014 03:44

Analysis Batch Number: 59779

End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
STD 320-59779/4 IC		12/04/2014 03:44	1	03DEC14A4A_040. d	Acquity 2.1(mm)
STD 320-59779/5 IC		12/04/2014 04:05	1	03DEC14A4A_041. d	Acquity 2.1(mm)
STD 320-59779/6 IC		12/04/2014 04:26	1	03DEC14A4A_042. d	Acquity 2.1(mm)
STD 320-59779/7 IC		12/04/2014 04:47	1	03DEC14A4A_043. d	Acquity 2.1(mm)
STD 320-59779/8 IC		12/04/2014 05:09	1	03DEC14A4A_044. d	Acquity 2.1(mm)
STD 320-59779/9 IC		12/04/2014 05:30	1	03DEC14A4A_045. d	Acquity 2.1(mm)
STD 320-59779/10 IC		12/04/2014 05:51	1	03DEC14A4A_046. d	Acquity 2.1(mm)
CCV 320-59779/20		12/04/2014 09:23	1	03DEC14A4A_056. d	Acquity 2.1(mm)
ICV 320-59779/22		12/04/2014 10:23	1	03DEC14A4A_058. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 11:50	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:11	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:37	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 12:59	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:20	100		Acquity 2.1(mm)
ZZZZZ		12/04/2014 13:41	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 14:02	1		Acquity 2.1(mm)
MB 320-58953/1-A		12/04/2014 14:24	1	03DEC14A4A_066. d	Acquity 2.1(mm)
LCS 320-58953/2-A		12/04/2014 14:44	1	03DEC14A4A_067. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:06	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 15:27	1		Acquity 2.1(mm)
CCV 320-59779/34		12/04/2014 15:48	1	03DEC14A4A_070. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:10	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:31	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 16:52	1		Acquity 2.1(mm)
320-10572-1	EFA-GW01	12/04/2014 17:13	1	03DEC14A4A_074. d	Acquity 2.1(mm)
320-10572-2	EFA-GW02	12/04/2014 17:34	1	03DEC14A4A_075. d	Acquity 2.1(mm)
320-10572-3	EFA-GW03	12/04/2014 17:56	1	03DEC14A4A_076. d	Acquity 2.1(mm)
320-10572-4	EFA-GW09	12/04/2014 18:17	1	03DEC14A4A_077. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:38	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 18:59	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 19:20	1		Acquity 2.1(mm)
CCV 320-59779/45		12/04/2014 19:42	1	03DEC14A4A_081. d	Acquity 2.1(mm)
ZZZZZ		12/04/2014 20:03	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 20:24	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 20:45	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 21:07	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 21:28	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 21:49	1		Acquity 2.1(mm)

## LCMS ANALYSIS RUN LOG

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1

SDG No.: \_\_\_\_\_

Instrument ID: A4 Start Date: 12/04/2014 03:44

Analysis Batch Number: 59779 End Date: 12/04/2014 23:56

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
ZZZZZ		12/04/2014 22:10	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 22:31	1		Acquity 2.1(mm)
ZZZZZ		12/04/2014 22:53	20		Acquity 2.1(mm)
ZZZZZ		12/04/2014 23:14	20		Acquity 2.1(mm)
ZZZZZ		12/04/2014 23:35	20		Acquity 2.1(mm)
CCV 320-59779/57		12/04/2014 23:56	1		Acquity 2.1(mm)

FORM VI  
LCMS INITIAL CALIBRATION DATA  
EXTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1 Analy Batch No.: 59779

SDG No.:

Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD 320-59779/4	03DEC14A4A_040.d
Level 2	STD 320-59779/5	03DEC14A4A_041.d
Level 3	STD 320-59779/6	03DEC14A4A_042.d
Level 4	STD 320-59779/7	03DEC14A4A_043.d
Level 5	STD 320-59779/8	03DEC14A4A_044.d
Level 6	STD 320-59779/9	03DEC14A4A_045.d
Level 7	STD 320-59779/10	03DEC14A4A_046.d

ANALYTE	CF				CURVE TYPE	COEFFICIENT			#	MIN CF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 5	LVL 2 LVL 6	LVL 3 LVL 7	LVL 4		B	M1	M2								
13C4 PFBA	5268.6 3806.7	5812.9 3203.9	5254.9 2254.3	4338.2	Ave		4277.06429				30.0		50.0			
13C5 PFPeA	9419.7 7332.0	9580.9 5721.7	8177.5 4869.4	8903.4	Ave		7714.90000				24.0		50.0			
13C2 PFHxA	21470 18141	22784 16317	24835 11422	18449	Ave		19059.8143				23.0		50.0			
13C4-PFHxA	21643 17882	21390 11698	21748 10393	19442	Ave		17742.3500				27.0		50.0			
18O2 PFHxS	53128 47816	52233 33678	50966 25618	51213	Ave		44950.3322				24.0		50.0			
13C4 PFOA	41322 31379	41400 24644	40475 19047	38546	Ave		33830.5286				27.0		50.0			
13C4 PFOS	54398 44130	54385 33330	56166 23197	50088	Ave		45099.0735				28.0		50.0			
13C5 PFNA	44879 37219	45410 30155	48879 22012	41831	Ave		38626.2857				25.0		50.0			
13C2 PFDA	63235 48806	61430 37175	59297 32265	58211	Ave		51488.2929				24.0		50.0			
13C8 FOSA	150505 147550	155647 126255	158874 106829	155831	Ave		143069.894				14.0		50.0			
13C2 PFUnA	66511 59868	66532 58833	66894 69853	64772	Ave		64751.6643				6.2		50.0			
13C2 PFDoA	76991 59932	66196 56050	76655 52807	68862	Ave		65356.2357				15.0		50.0			
13C2-PFTeDA	77915 69707	76286 61208	80240 58365	76834	Ave		71507.9143				12.0		50.0			
13C2-PFHxDA	96546 111182	105842 105984	112387 93820	103841	Ave		104228.936				6.6		50.0			

Note: The m1 coefficient is the same as Ave CF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento Job No.: 320-10572-1 Analy Batch No.: 59779  
SDG No.: \_\_\_\_\_  
Instrument ID: A4 GC Column: Acquity ID: 2.1 (mm) Heated Purge: (Y/N) N \_\_\_\_\_  
Calibration Start Date: 12/04/2014 03:44 Calibration End Date: 12/04/2014 05:51 Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorobutanoic acid (PFBA)	10678 16623	22829 12169	21990	19527	19477	AveID		2.3514				25.0		35.0			
Perfluoropentanoic acid (PFPeA)	8732.0 11185	14944 9458.0	14198	14832	13130	AveID		1.6538				21.0		35.0			
Perfluorobutane Sulfonate (PFBS)	14683 39320	46758 32165	49055	48250	67174	AveID		0.9863				37.0		50.0			
Perfluorohexanoic acid (PFHxA)	9952.0 16870	21004 12521	26818	19955	20968	AveID		0.9761				24.0		35.0			
PFPeS (Perflouro-1-pentanesulfonate)	12755 32980	40229 24854	44086	42428	40598	AveID		0.7860				32.0		50.0			
Perfluoroheptanoic acid (PFHpA)	14724 21344	28535 18892	35883	33256	29606	AveID		1.5247				27.0		35.0			
Perfluorohexane Sulfonate (PFHxS)	24632 33051	52072 26648	53418	49529	45799	AveID		0.9222				22.0		35.0			
Perfluoroctanoic acid (PFOA)	54920 43078	74889 33366	68131	62497	54775	AveID		1.6697				9.7		35.0			
Perfluoro-1-heptanesulfonate (PFHpS)	12752 33076	45349 24051	51122	49900	46278	AveID		0.8647				33.0		50.0			
Perfluorooctane Sulfonate (PFOS)	143245 49883	124703 36721	92010	73648	66987	AveID		1.8046				26.0		35.0			
Perfluorononanoic acid (PFNA)	22096 39342	60944 27244	61254	52411	44283	AveID		1.1532				26.0		35.0			
PFNS (Perflouro-1-nonanesulfonate)	16554 25936	40576 18531	47018	41535	37537	AveID		0.7349				26.0		50.0			
Perfluorodecanoic acid (PFDA)	16398 41090	58072 33184	67070	63478	54711	AveID		0.9544				33.0		35.0			
Perfluorooctane Sulfonamide (FOSA)	60576 132015	157092 116050	155459	165246	157618	AveID		0.9501				26.0		35.0			
Perfluorodecane sulfonate (PFDS)	10237 23088	37158 16598	42628	36833	33233	AveID		0.6467				32.0		50.0			
Perfluoroundecanoic acid (PFUnA)	25356 49828	75233 40491	72252	72984	62389	AveID		0.8839				34.0		35.0			
Perfluorododecanoic acid (PFDoA)	25100 62748	61887 50239	92219	78804	73401	AveID		0.9863				32.0		35.0			
PFDoS (Perflouro-1-dodecanesulfonate)	7778.9 23770	32535 17735	43717	40638	35319	AveID		0.6584				36.0		50.0			
Perfluorotridecanoic Acid (PFTriA)	20248 45661	65903 40208	67545	66220	62650	AveID		0.8175				32.0		50.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI  
LCMS INITIAL CALIBRATION DATA  
CURVE EVALUATION

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

Analy Batch No.: 59779

SDG No.:

Instrument ID: A4

GC Column: Acuity

ID: 2.1 (mm)

Heated Purge: (Y/N) N

Calibration Start Date: 12/04/2014 03:44

Calibration End Date: 12/04/2014 05:51

Calibration ID: 10479

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3	LVL 4	LVL 5		B	M1	M2								
Perfluorotetradecanoic acid (PFTeA)	15018 30898	36572 27633	43161	42497	38214	AveID		0.5200				29.0		50.0			
Perfluoro-n-hexadecanoic acid (PFHxDA)	44032 92953	98406 78170	100981	86858	96066	AveID		1.3398				27.0		50.0			
Perfluoro-n-octadecanoic acid (PFODA)	22762 104163	94742 95748	99443	104957	102973	AveID		1.4197				38.0		50.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.:

Lab Sample ID: CCV 320-59779/20

Calibration Date: 12/04/2014 09:23

Instrument ID: A4

Calib Start Date: 12/04/2014 03:44

GC Column: Acquity

ID: 2.10 (mm)

Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_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	2.351	2.886		24.5	20.0	22.7	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	1.983		24.0	20.0	19.9	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.029		18.5	17.7	4.4	40.0
Perfluoroheptanoic acid (PFHxA)	AveID	0.9761	1.064		21.8	20.0	9.0	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.016		26.4	20.0	32.2	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	0.9830		20.2	18.9	6.6	40.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	1.083		23.9	19.0	25.3	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.896		22.7	20.0	13.5	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.558		16.5	19.1	-13.6	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.188		20.6	20.0	3.0	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.105		23.1	20.0	15.7	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.027		21.6	20.0	8.1	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7177		21.4	19.3	11.0	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.156		26.1	20.0	30.7	40.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.170		23.7	20.0	18.6	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9363		22.9	20.0	14.5	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5680		21.8	20.0	9.2	40.0
Perfluoro-n-hexadecanoic acid (FFHxDA)	AveID	1.340	1.393		20.8	20.0	3.9	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.440		20.3	20.0	1.4	40.0

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.:

Lab Sample ID: ICV 320-59779/22

Calibration Date: 12/04/2014 10:23

Instrument ID: A4

Calib Start Date: 12/04/2014 03:44

GC Column: Acquity ID: 2.10 (mm)

Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_058.d

Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	2.637		56.1	50.0	12.2	50.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	1.886		57.0	50.0	14.0	50.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.236		55.5	44.3	25.3	50.0
Perfluoroheptanoic acid (PFHxA)	AveID	0.9761	1.190		61.0	50.0	21.9	50.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	2.160		70.8	50.0	41.7	50.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	1.119		57.4	47.3	21.4	50.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9109		50.1	47.6	5.3	50.0
Perfluoroctanoic acid (PFOA)	AveID	1.670	1.812		54.2	50.0	8.5	50.0
Perfluoroctane Sulfonate (PFOS)	AveID	1.805	1.655		43.8	47.8	-8.3	50.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.206		52.3	50.0	4.6	50.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.211		63.4	50.0	26.8	50.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	0.9856		51.9	50.0	3.7	50.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7475		55.8	48.3	15.6	50.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.141		64.5	50.0	29.1	50.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.176		59.6	50.0	19.2	50.0
Perfluorotridecanoic Acid (PFTrIA)	AveID	0.8175	0.9516		58.2	50.0	16.4	50.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.6011		57.8	50.0	15.6	50.0
Perfluoro-n-hexadecanoic acid (FFHxDA)	AveID	1.340	1.644		61.3	50.0	22.7	50.0
Perfluoro-n-octadecanoic acid (FFODA)	AveID	1.420	1.763		62.1	50.0	24.2	50.0

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.: \_\_\_\_\_

Lab Sample ID: CCV 320-59779/34 Calibration Date: 12/04/2014 15:48

Instrument ID: A4 Calib Start Date: 12/04/2014 03:44

GC Column: Acquity ID: 2.10 (mm) Calib End Date: 12/04/2014 05:51

Lab File ID: 03DEC14A4A\_070.d Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	2.527		21.5	20.0	7.5	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	1.839		22.2	20.0	11.2	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	0.9250		16.6	17.7	-6.2	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.166		23.9	20.0	19.5	40.0
Perfluoroheptanoic acid (PFHpA)	AveID	1.525	1.731		22.7	20.0	13.6	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	0.8886		18.2	18.9	-3.6	40.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9356		20.6	19.0	8.2	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.742		20.9	20.0	4.3	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.442		15.3	19.1	-20.1	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.235		21.4	20.0	7.1	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.137		23.8	20.0	19.1	40.0
Perfluoroctane Sulfonamide (FOSA)	AveID	0.9501	1.031		21.7	20.0	8.5	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.6365		19.0	19.3	-1.6	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.159		26.2	20.0	31.1	40.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.149		23.3	20.0	16.5	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9686		23.7	20.0	18.5	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5747		22.1	20.0	10.5	40.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.470		21.9	20.0	9.7	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.601		22.5	20.0	12.7	40.0

FORM VII  
LCMS CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Sacramento	Job No.: 320-10572-1
SDG No.:	
Lab Sample ID: CCV 320-59779/45	Calibration Date: 12/04/2014 19:42
Instrument ID: A4	Calib Start Date: 12/04/2014 03:44
GC Column: Acquity ID: 2.10 (mm)	Calib End Date: 12/04/2014 05:51
Lab File ID: 03DEC14A4A_081.d	Conc. Units: ng/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Perfluorobutanoic acid (PFBA)	AveID	2.351	3.794		32.3	20.0	61.4*	40.0
Perfluoropentanoic acid (PFPeA)	AveID	1.654	2.282		27.6	20.0	38.0	40.0
Perfluorobutane Sulfonate (PFBS)	AveID	0.9863	1.023		18.3	17.7	3.7	40.0
Perfluorohexanoic acid (PFHxA)	AveID	0.9761	1.252		25.6	20.0	28.2	40.0
Perfluorohethanoic acid (PFHpA)	AveID	1.525	2.068		27.1	20.0	35.7	40.0
Perfluorohexane Sulfonate (PFHxS)	AveID	0.9222	0.998		20.5	18.9	8.2	40.0
Perfluoro-1-heptanesulfonate (PFHpS)	AveID	0.8647	0.9528		21.0	19.0	10.2	40.0
Perfluorooctanoic acid (PFOA)	AveID	1.670	1.850		22.2	20.0	10.8	40.0
Perfluorooctane Sulfonate (PFOS)	AveID	1.805	1.488		15.8	19.1	-17.6	40.0
Perfluorononanoic acid (PFNA)	AveID	1.153	1.142		19.8	20.0	-1.0	40.0
Perfluorodecanoic acid (PFDA)	AveID	0.9544	1.018		21.3	20.0	6.6	40.0
Perfluorooctane Sulfonamide (FOSA)	AveID	0.9501	1.009		21.2	20.0	6.2	40.0
Perfluorodecane sulfonate (PFDS)	AveID	0.6467	0.7121		21.2	19.3	10.1	40.0
Perfluoroundecanoic acid (PFUnA)	AveID	0.8839	1.135		25.7	20.0	28.4	40.0
Perfluorododecanoic acid (PFDoA)	AveID	0.9863	1.182		24.0	20.0	19.9	40.0
Perfluorotridecanoic Acid (PFTriA)	AveID	0.8175	0.9328		22.8	20.0	14.1	40.0
Perfluorotetradecanoic acid (PFTeA)	AveID	0.5200	0.5912		22.7	20.0	13.7	40.0
Perfluoro-n-hexadecanoic acid (PFHxDA)	AveID	1.340	1.503		22.4	20.0	12.2	40.0
Perfluoro-n-octadecanoic acid (PFODA)	AveID	1.420	1.589		22.4	20.0	11.9	40.0

## LCMS BATCH WORKSHEET

Lab Name: TestAmerica Sacramento

Job No.: 320-10572-1

SDG No.:

Batch Number: 58953

Batch Start Date: 11/24/14 16:22

Batch Analyst: Reed, Jonathan E

Batch Method: 3535

Batch End Date: 11/25/14 17:08

Lab Sample ID	Client Sample ID	Method Chain	Basis	InitialAmount	FinalAmount	LCMPFCSU 00010	LCPFCSP 00014		
MB 320-58953/1		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL			
LCS 320-58953/2		3535, WS-LC-0025		500.00 mL	1.00 mL	50 uL	20 uL		
320-10572-A-1	EFA-GW01	3535, WS-LC-0025	T	500.25 mL	1.00 mL	50 uL			
320-10572-A-2	EFA-GW02	3535, WS-LC-0025	T	499.85 mL	1.00 mL	50 uL			
320-10572-A-3	EFA-GW03	3535, WS-LC-0025	T	502.05 mL	1.00 mL	50 uL			
320-10572-A-4	EFA-GW09	3535, WS-LC-0025	T	501.98 mL	1.00 mL	50 uL			

## Batch Notes

Balance ID	QA-070
H2O Lot used	11/14/14
Pipette ID	EC15219
Analyst who added reagent	JER
SU Reagent Drop	JER
SU Reagent Drop Witness	AVM
Solvent Lot #	0.3%NH4OH-Me_00006
Solvent Name	0.3%NH4OH/MeOH
SOP Number	WS-LC-0025
SPE Cartridge Type	WAX 500mg
Solid Phase Extraction Disk Lot Number	0023341484

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.

WS-LC-0025

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DODCMD_ID	INSTALLATION_ID	SDG	SITE_NAME	NORM_SITE_NAME	LOCATION_NAME	LOCATION_TYPE_DESC	COORD_X	COORD_Y	CONTRACT_ID	DO_CTO_NUMBER	CONTR_NAME	SAMPLE_NAME	SAMPLE_MATRIX_DESC	SAMPLE_TYPE_DESC	COLLECT_DATE	ANALYTICAL_METHOD	ANALYTICAL_METHOD_GRP_DESC
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	RCRA CLOSURE	SITE 00011	NASB-H123-MW02	Monitoring well	3014401.09	387341.48	N6247008D1001	WE09	TETRA TECH NUS, INC.	H123-GW02-20141116-D	Ground water	Field duplicate	16-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	RCRA CLOSURE	SITE 00011	NASB-H123-MW02	Monitoring well	3014401.09	387341.48	N6247008D1001	WE09	TETRA TECH NUS, INC.	H123-GW02-20141116	Ground water	Normal (Regular)	16-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	BLDG 7/10	B7/10-MW065	Monitoring well	3014954.96	387415.51	N6247008D1001	WE09	TETRA TECH NUS, INC.	B7/10-GW065-20141116	Ground water	Normal (Regular)	16-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds	
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	BLDG 7/10	B7/10-MW03S	Monitoring well	3014935.31	387512.11	N6247008D1001	WE09	TETRA TECH NUS, INC.	B7/10-GW03S-20141115	Ground water	Normal (Regular)	15-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds	
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	BLDG 7/10	B7/10-MW02	Monitoring well	3014857.41	387526.42	N6247008D1001	WE09	TETRA TECH NUS, INC.	B7/10-GW02-20141115	Ground water	Normal (Regular)	15-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds	
MID_ATLANTIC	BRUNSWICK_NAS	320-10459-1	MRP QS	MW-QRY-02	Monitoring well	3010220.926	382943.47	N6247008D1001	WE09	TETRA TECH NUS, INC.	QRY-GW02-20141116	Ground water	Normal (Regular)	16-Nov-14	TA_WS-LC-0025	Perfluoroalkyl Compounds	