

Optimizing analytical strategies for the quantification and identification of PFASs in water

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'Fit for Purpose' Analysis



Factors in deciding analytical method selection

Considerations when selecting analytical technique to employ

- I. Types of analytes
- II. Concentrations required (ppm, ppb, ppt etc.)
- III. Type of matrix (drinking water, surface water, wastewater, sludge, food etc.)
- IV. Type of sample clean-up required
- V. Regulatory considerations (EPA methods, EU & ASTM methods etc.)



Local Contamination

Concern Grows Over Tainted Drinking Water

Vermont, New Hampshire and New York expand efforts to find out how much of a potentially toxic chemical is in drinking water



Residents look for their homes on a map showing PFOA-contaminated areas at a public information meeting in Litchfield, N.H., earlier this month. PHOTO: CHERYL SENTER FOR THE WALL STREET JOURNAL

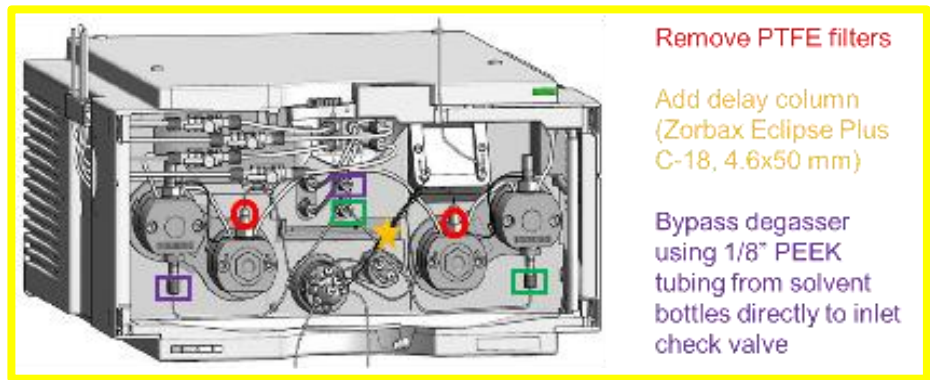
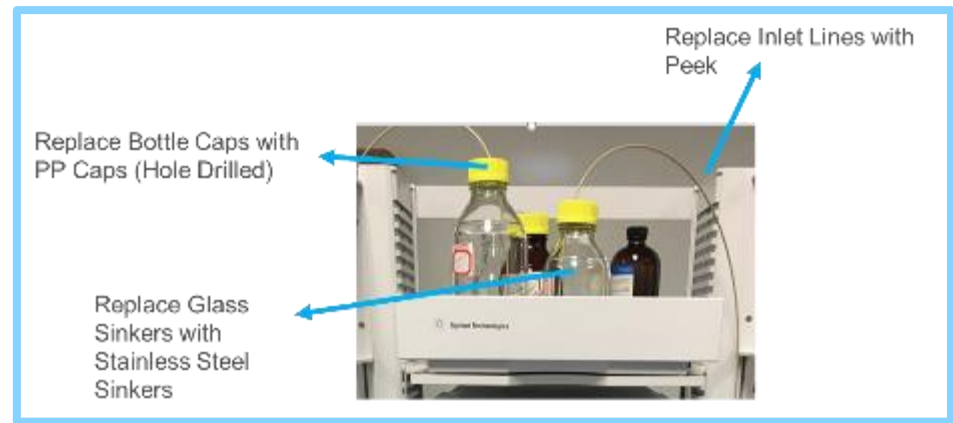
Public concern over PFOA has spread through upstate New York and New England since August 2014, when a resident of Hoosick Falls, N.Y., near the Vermont border, tested his drinking water and found high levels of the acid. The man was concerned because his father, a former employee of the town’s plastics plant that used PFOA, died of cancer.

of PFOA contamination. The state in March sampled PFOA levels up to 620 parts per trillion in private Litchfield wells, well above the 100-parts-per-trillion level at which New Hampshire officials start to consider the amount unsafe. Tests in Merrimack measured as high as 1,600 parts per trillion.

State	Number of systems with water samples with PFOA	Number of water samples with PFOA	Percent of samples with PFOA detected	Population served by drinking water systems with PFOA ^a	Range of average level of PFOA in samples (parts per billion) ^b	Average level of PFOA compared to safe level in new research
Total	94	287	23%	6,560,087		

PFAS Analysis – LC Instrument Setup

Eliminate Background Contamination

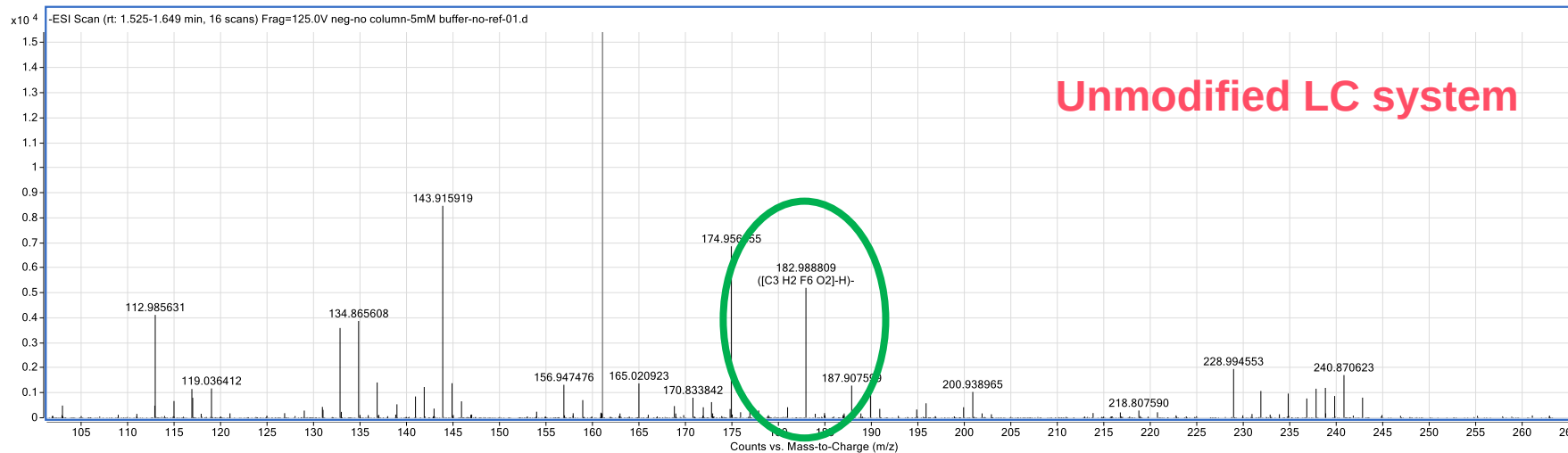


Application Note: Recommended Plumbing Configurations for Reduction in Per/Polyfluoroalkyl Substance Background with Agilent 1260/1290 Infinity (II) LC's (5991-7863EN)

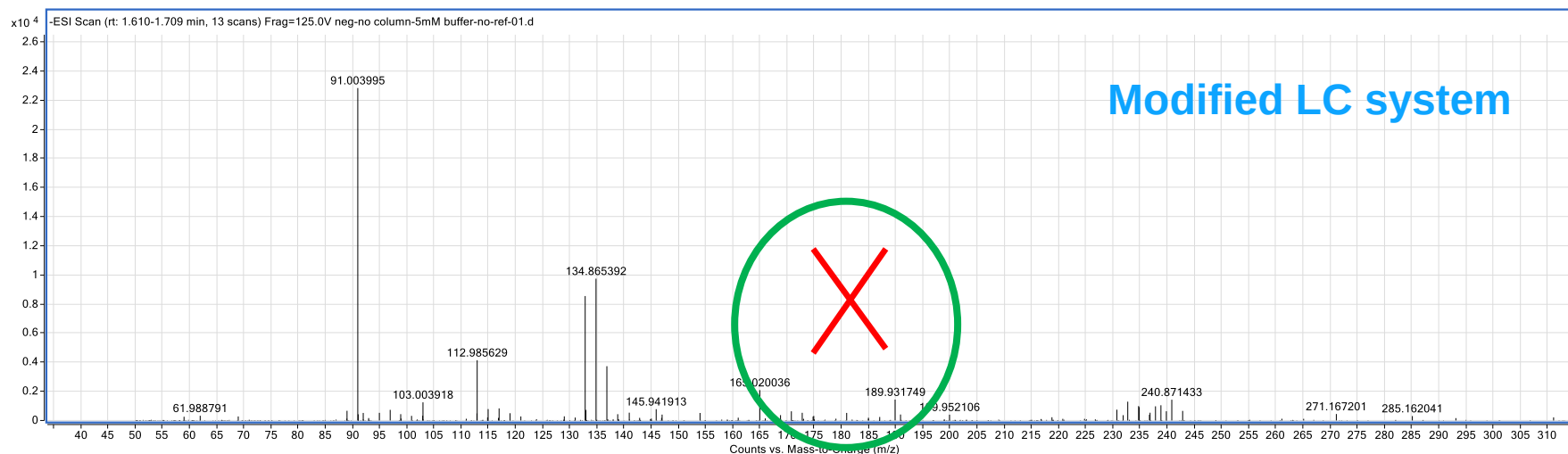


PFAS Analysis – LC Setup

Eliminate Background Contamination



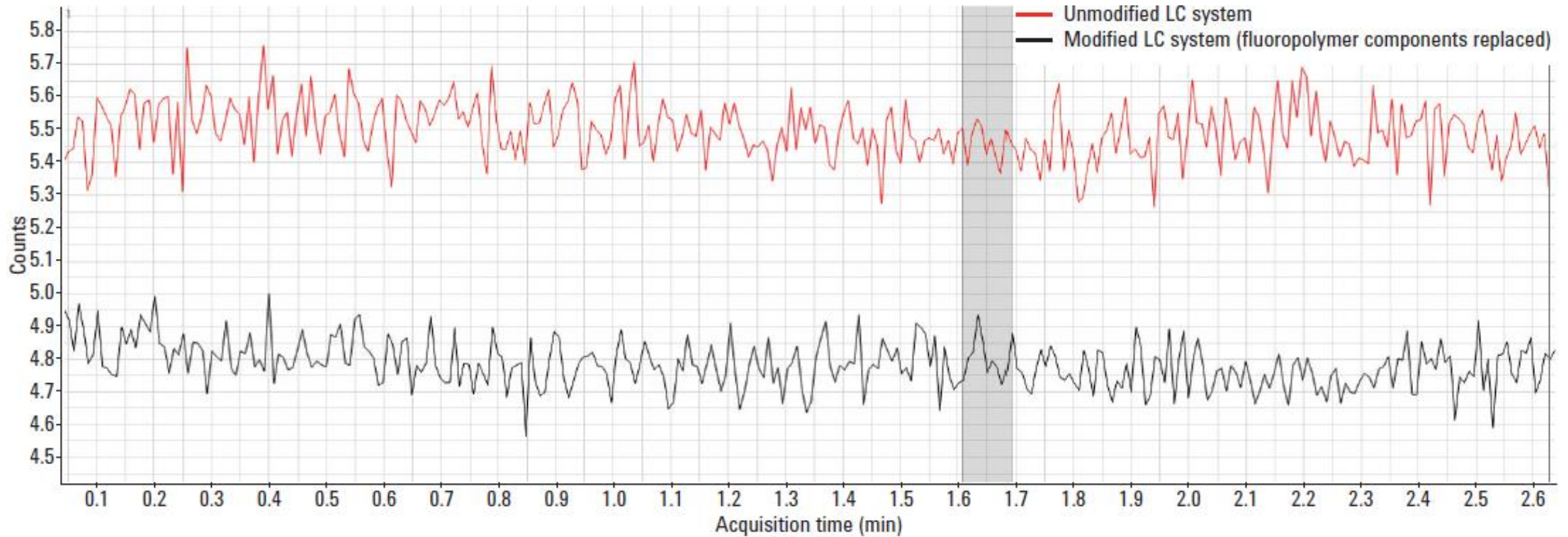
Fluoropolymer additive
(Hexafluoropropanediol)
absent in the modified
LC system



Data collected on a
6530 LC-Q/TOF

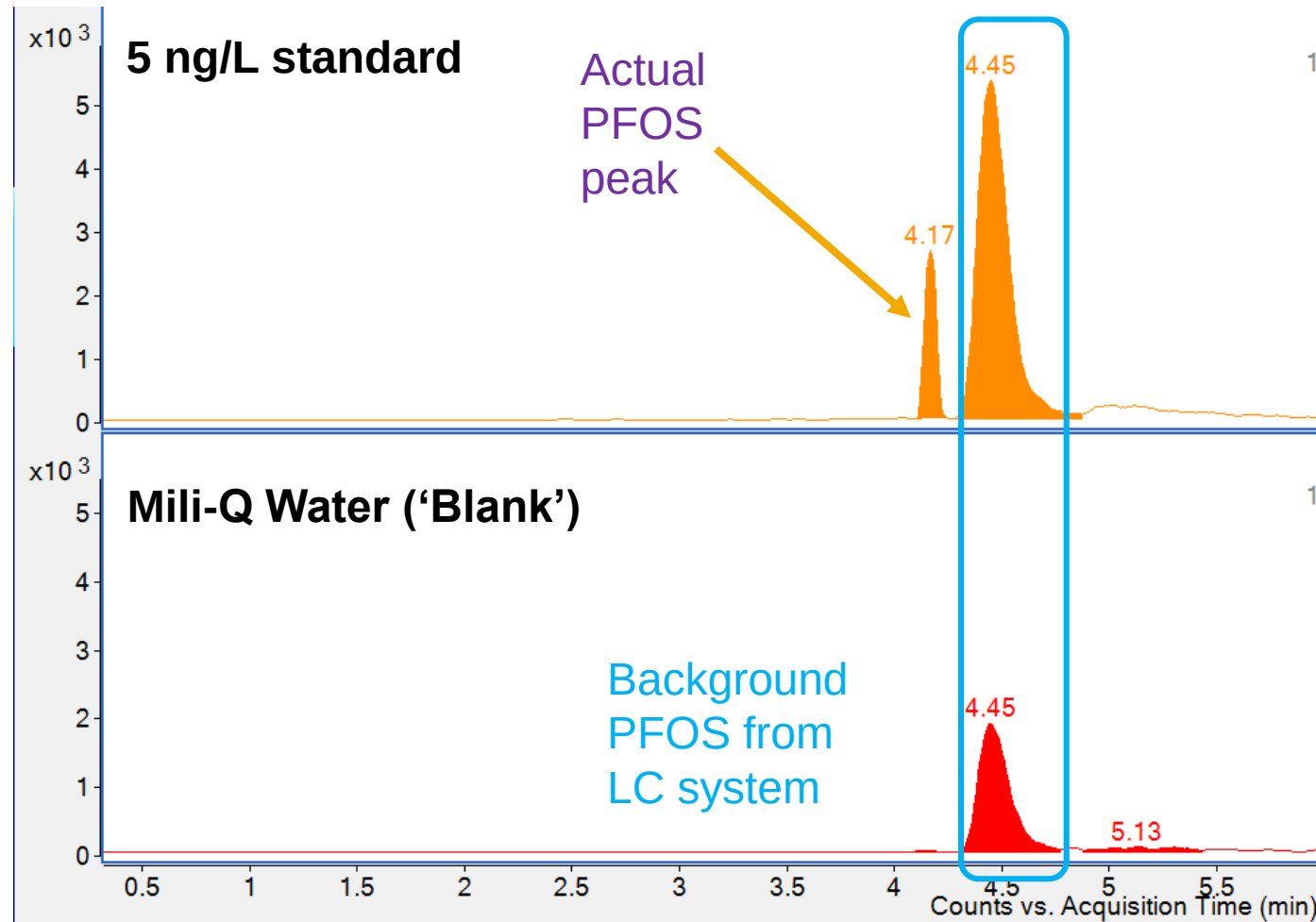
PFAS Analysis – LC Setup

Eliminate Background Contamination



PFAS Instrument Setup

Background Contamination



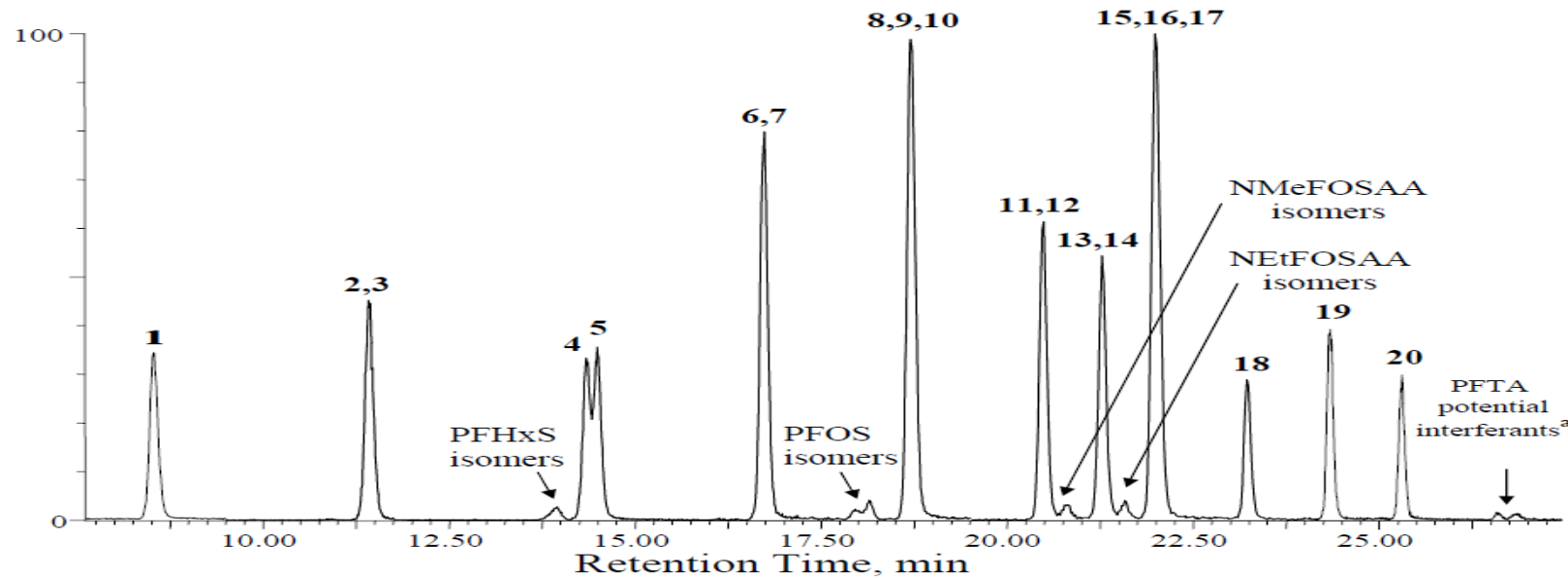
- The use of a Delay column is convenient when running multiple methods on the same instrument.
- “Delay” column and not a “Trapping” column. I.e. the background peak will be retention time separated but will be present.
- The delay column does not account for any contamination after the pump (the autosampler)

EPA 537: Perfluorinated Alkyl Acids in Drinking Water by SPE-LC/MS/MS

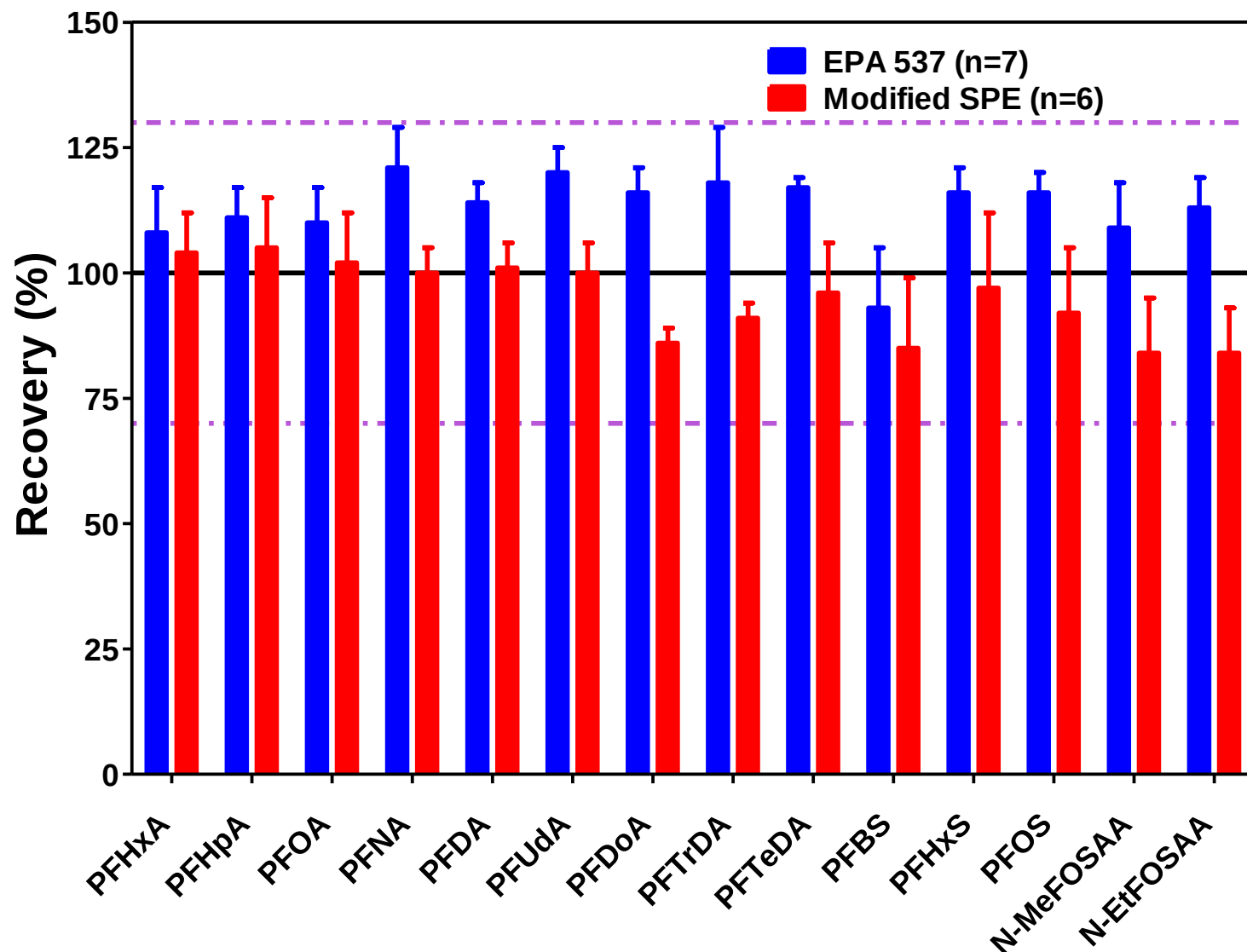
250 mL sample collected in **Polypropylene Bottles** with 5 g/L Trizma (to quench residual chlorine)

Solid Phase Extraction and evaporation to final **1 mL extract** in ~96/4 MeOH/Water

Analysis by **LC-MS/MS** in **Negative Electrospray** mode for **14 analytes** (9 PFCAs, 3 PFSAAs, 2 FOSAAs) – **37 min analysis (10 µL injection)**



Solid Phase Extraction USEPA 537 vs Modified



- Recoveries calculated in Reagent Water
- Modified SPE cartridge: SampliQ WAX cartridge (6 cc, 150 mg)
- All recoveries between 70-130 %

Spiking concentrations:

Modified SPE: 4 ng/L in 250 mL
EPA 537: 5-21 ng/L in 250 mL

Final extract: 1 mL ~95% MeOH

Typical LC/TQ Lab Layout



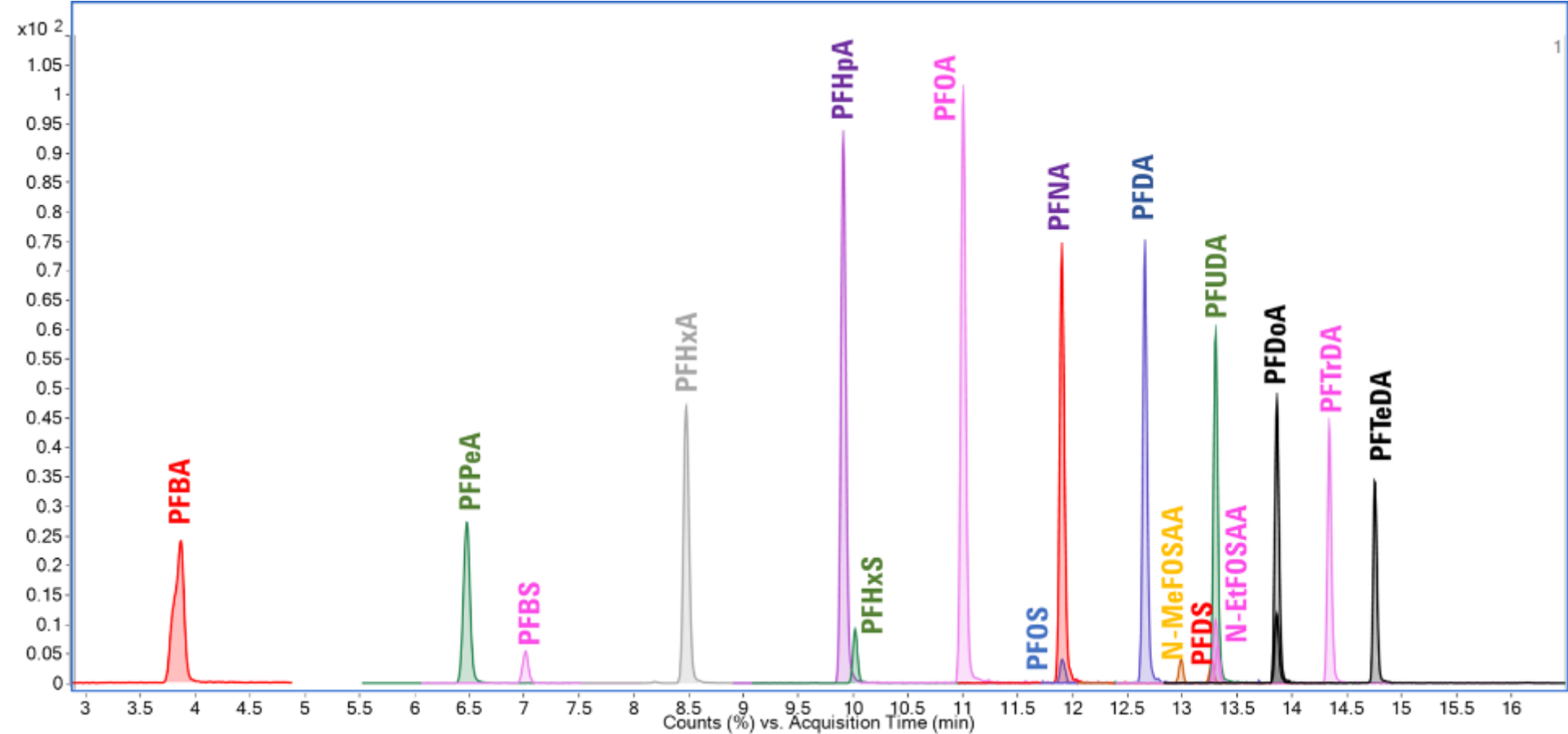
6 feet
1.83 meters



Analysis of PFAS in Drinking Water with Ultivo

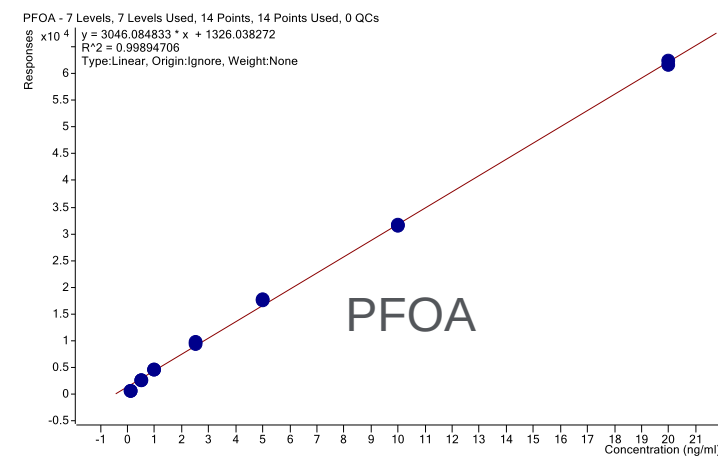
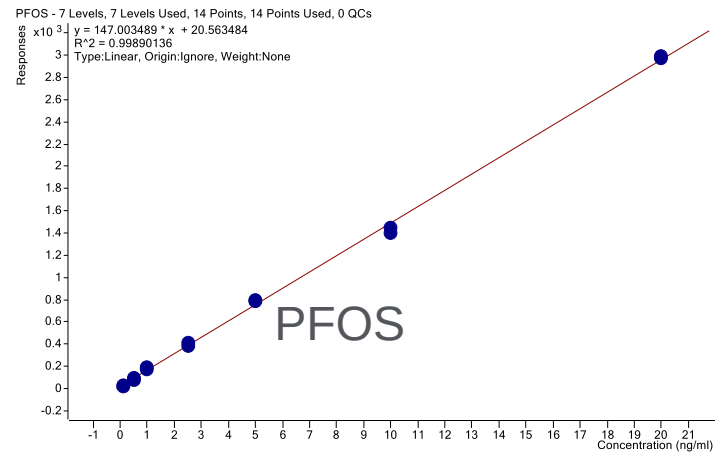
EPA Method 537

Analysis of 17 PFASs in drinking water with <0.1 ng/mL DLs in extract (includes all comps in EPA 537)

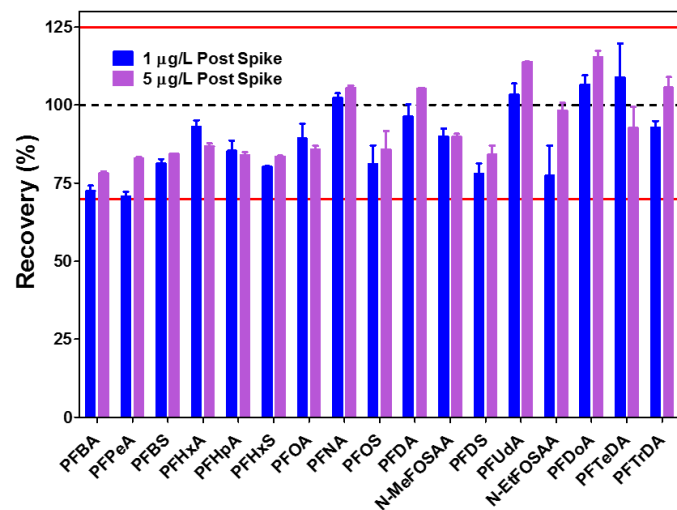


Analysis of PFAS in Drinking Water with Ultivo EPA Method 537

- All 17 PFASs had linear calibration curves, with $R^2 > 0.99$
- Overall recoveries were between 70–125% for both spiking levels
- %RSD was 0.3–10.8 % for all compounds at both spiking levels (1.0 µg/L and 5.0 µg/L)



Calibration Curves from 0.1-20 ng/mL for PFOA and PFOS



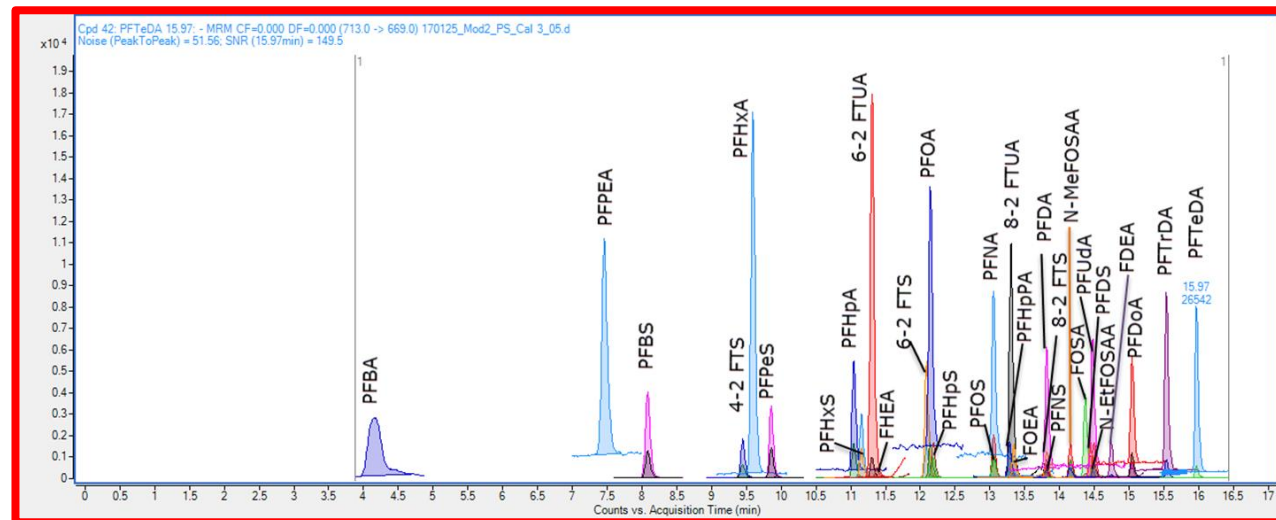
Post-spike recoveries of 17 PFASs in a water sample extracted according to EPA 537 with Agilent SampliQ WAX cartridges



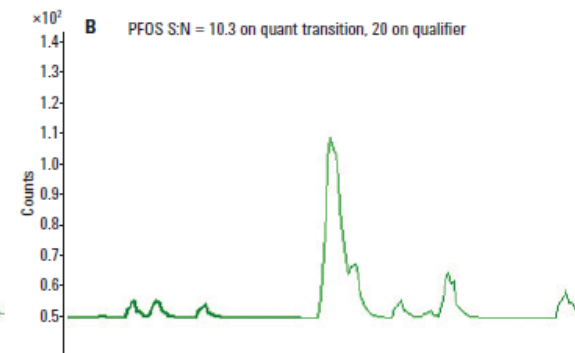
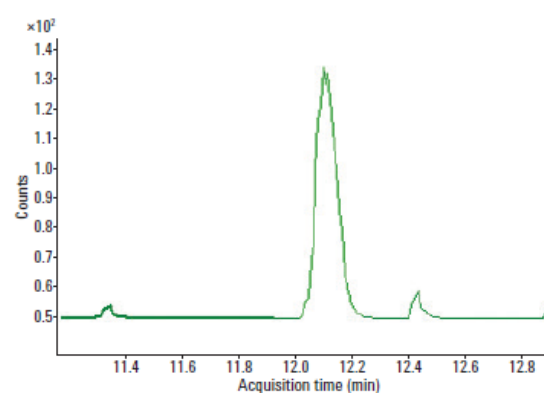
Ultivo Triple Quad LC/MS

Analysis of 30 PFASs in Drinking Water

Compound	Acronym	Class	Fluorinated C chain	EPA 537 or Additional
Perfluorobutanoate	PFBA	Acid	C4	Additional
Perfluoropentanoate	PFPeA	Acid	C5	Additional
Perfluorohexanoate	PFHxA	Acid	C6	537
Perfluoroheptanoate	PFHpA	Acid	C7	537
Perfluorooctanoate	PFOA	Acid	C8	537
Perfluorononanoate	PFNA	Acid	C9	537
Perfluorodecanoate	PFDA	Acid	C10	537
Perfluoroundecanoate	PFUdA	Acid	C11	537
Perfluorododecanoate	PFDoA	Acid	C12	537
Perfluorotridecanoate	PFTriDA	Acid	C13	537
Perfluorotetradecanoate	PFTeDA	Acid	C14	537
Perfluorooctanesulfonamide	FOSA	FOSA	C8	Additional
N-Ethyl-N-((heptadecafluorooctyl)sulfonyl)glycine	N-EtFOSAA	FOSAA	C8	537
N-((Heptadecafluorooctyl)sulfonyl)-N-methylglycine	N-MeFOSAA	FOSAA	C8	537
2-Perfluorohexyl ethanoic acid	FHEA	FTA-e	C6	Additional
2-Perfluorooctyl ethanoic acid	FOEA	FTA-e	C8	Additional
2-Perfluorodecyl ethanoic acid	FDEA	FTA-e	C10	Additional
3-Perfluoroheptyl propanoic acid (FHpPA)	PFHpPA	FTA-p	C7	Additional
4:2 Fluorotelomer sulfonate	4-2 FTS	FTS	C4	Additional
6:2 Fluorotelomer sulfonate	6-2 FTS	FTS	C6	Additional
8:2 Fluorotelomer sulfonate	8-2 FTS	FTS	C8	Additional
2H-Perfluoro-2-octanoic acid (FHUEA)	6-2 FTUA	FTUA	C6	Additional
2H-Perfluoro-2-decanoic acid (FOUEA)	8-2 FTUA	FTUA	C8	Additional
Perfluorobutylsulfonate	PFBS	Sulfonate	C4	537
Perfluoropentylsulfonate	PFPeS	Sulfonate	C5	Additional
Perfluorohexylsulfonate	PFHxS	Sulfonate	C6	537
Perfluoroheptylsulfonate	PFHpS	Sulfonate	C7	Additional
Perfluorooctylsulfonate	PFOS	Sulfonate	C8	537
Perfluorononylsulfonate	PFNS	Sulfonate	C9	Additional
Perfluorodecylsulfonate	PFDS	Sulfonate	C10	Additional



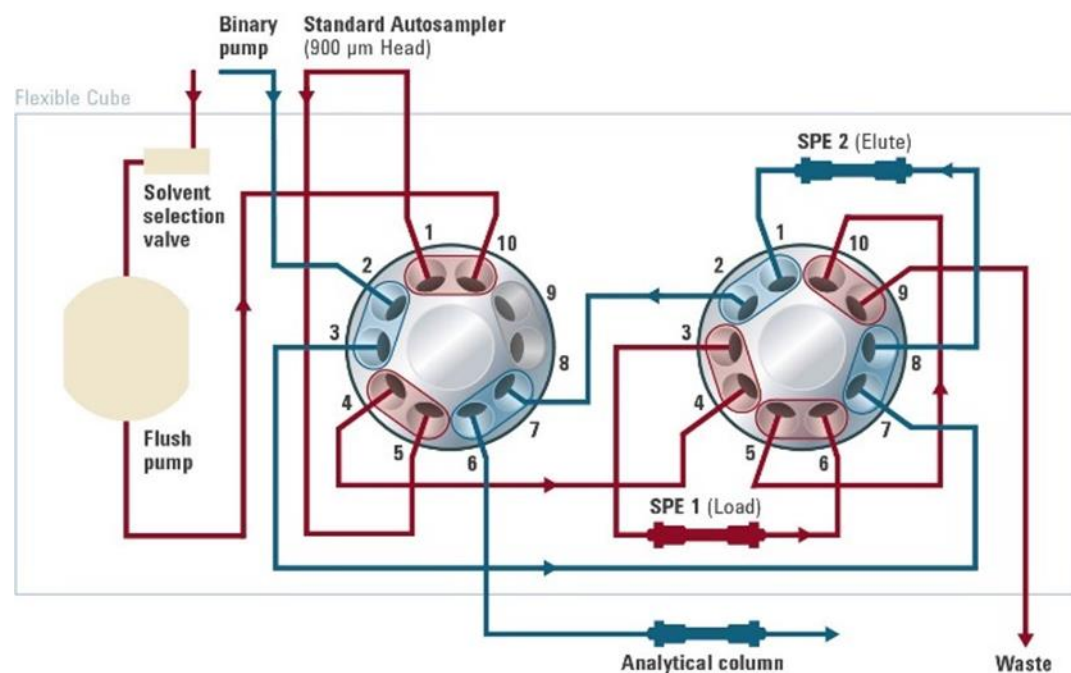
Sensitivity (0.08 ng/L)



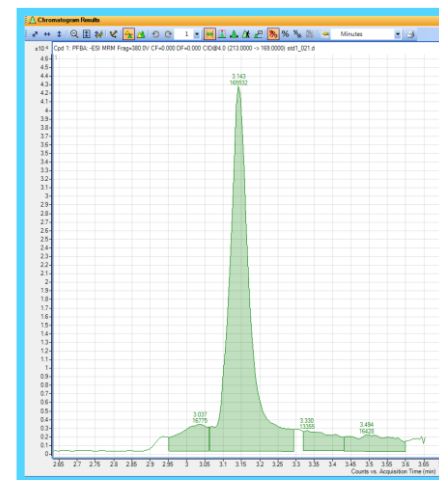
Automated Online SPE for PFASs in Potable Water



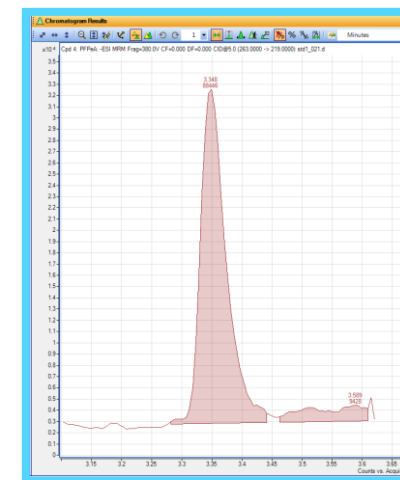
Lower human
intervention steps:
↑ Accuracy
↑ Reproducibility



1 ng/L spike into tap water



PFBA



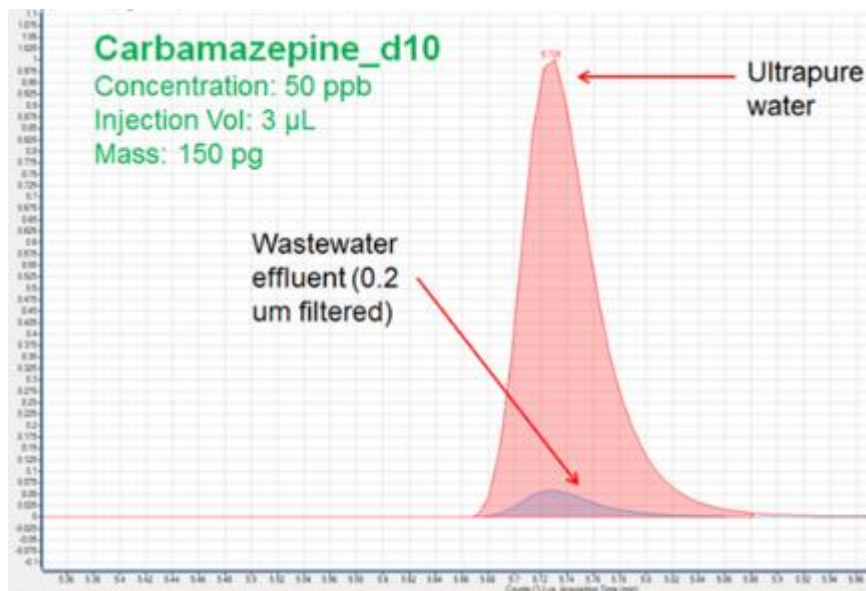
PFPeA

Online SPE method highlights:

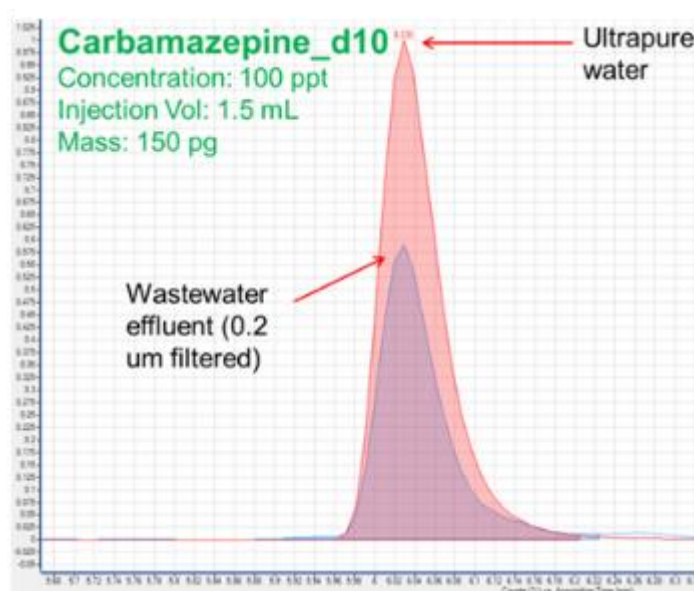
- 1290 Infinity LC + 6470 LC/TQ
- 900 µL injection
- <20 min cycle time
- DLs: 0.4-3 ng/L

Analysis of Environmental Contaminants by LC-MS/MS

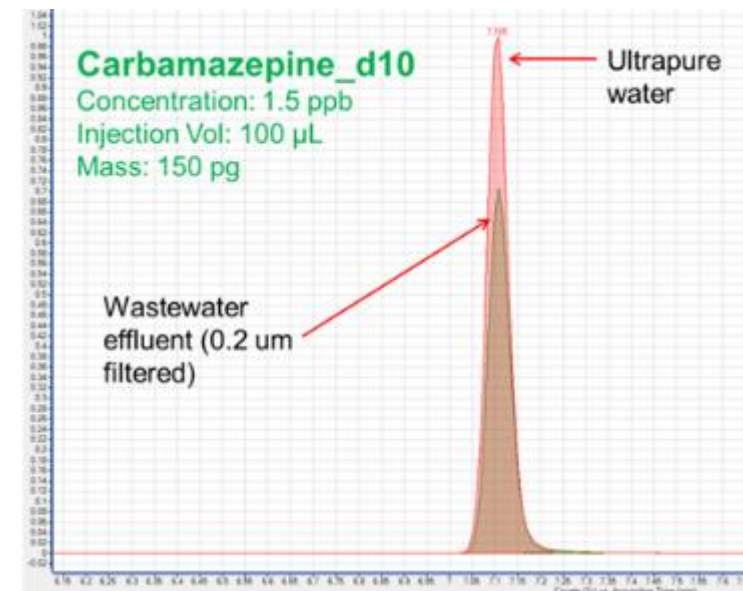
Ion Suppression Comparison by Sample Prep Method



Conventional SPE
Mass on Column :150 pg



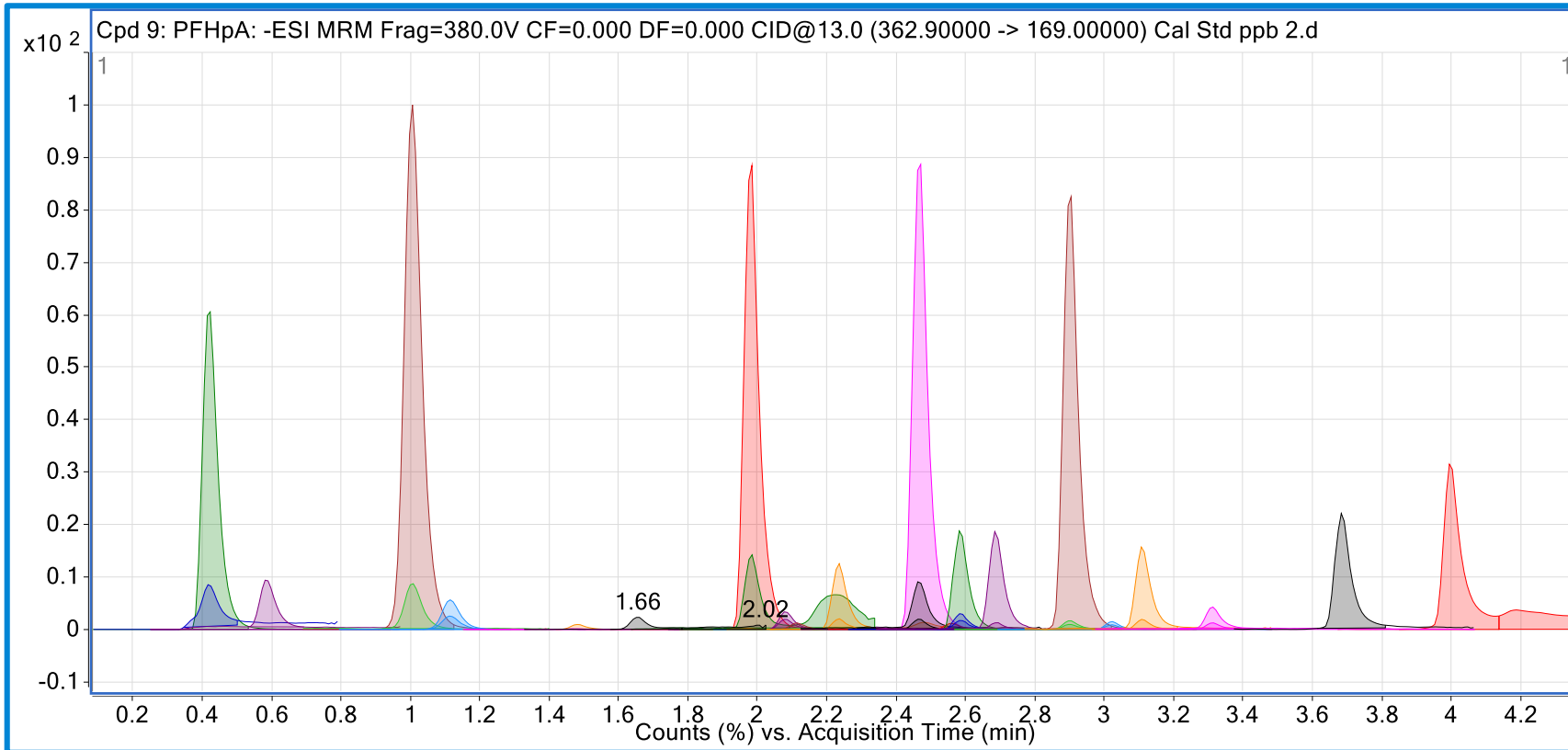
Online SPE
Mass on Column :150 pg



Direct Injection
Mass on Column :150 pg

Perfluorinated Alkyl Acids in Drinking Water by LC/TQ

Collaborative work with Suffolk County Water Authority



System

- 1290 Infinity LC coupled to a 6490 LC-MS/MS
- 14 PFAAs in 6 min

MDLs:

- PFOA, PFOS: <1 ng/L
- Others: 1-10 ng/L
- 80 μ L aqueous injection

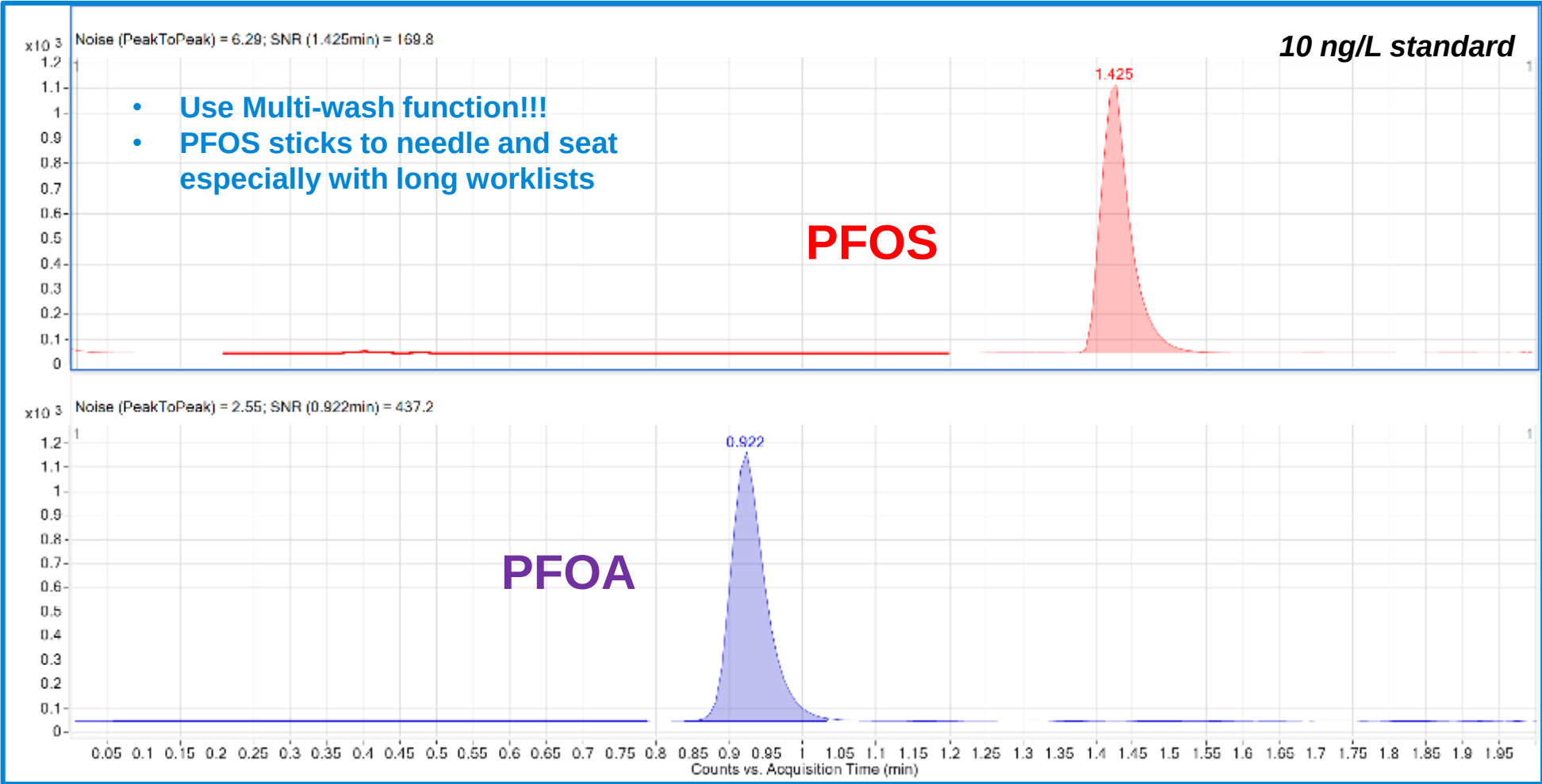
Matrices:

- Drinking Water
- Surface Water
- Ground Water

Analysis of PFOA/PFOS

Direct Injection; Rapid Analysis (1290 II UHPLC + 6495 LC-MS/MS)

Liquid Chromatograph	Agilent 1290 Infinity II Binary Pump
Mass Spectrometer	Agilent 6495 triple quadrupole LC-MS
Analytical Column	Agilent Poroshell 120 EC-C18; 3.0 x 50mm; 2.7um
Delay Column	Agilent Eclipse Plus C18, 4.6 x 50mm; 3.5um
Mobile Phase	A: Water+5mM Amm. Acetate B: Acetonitrile
Run Time	2 min
Injection Volume	80 uL (Water)



Are we measuring the right PFASs?

8-2

FTAL PFHpA

FOSA

PFT_rD PF

THE TRUTH IS OUT THERE

THE X FILES

THE TRUTH IS OUT THERE

SAA

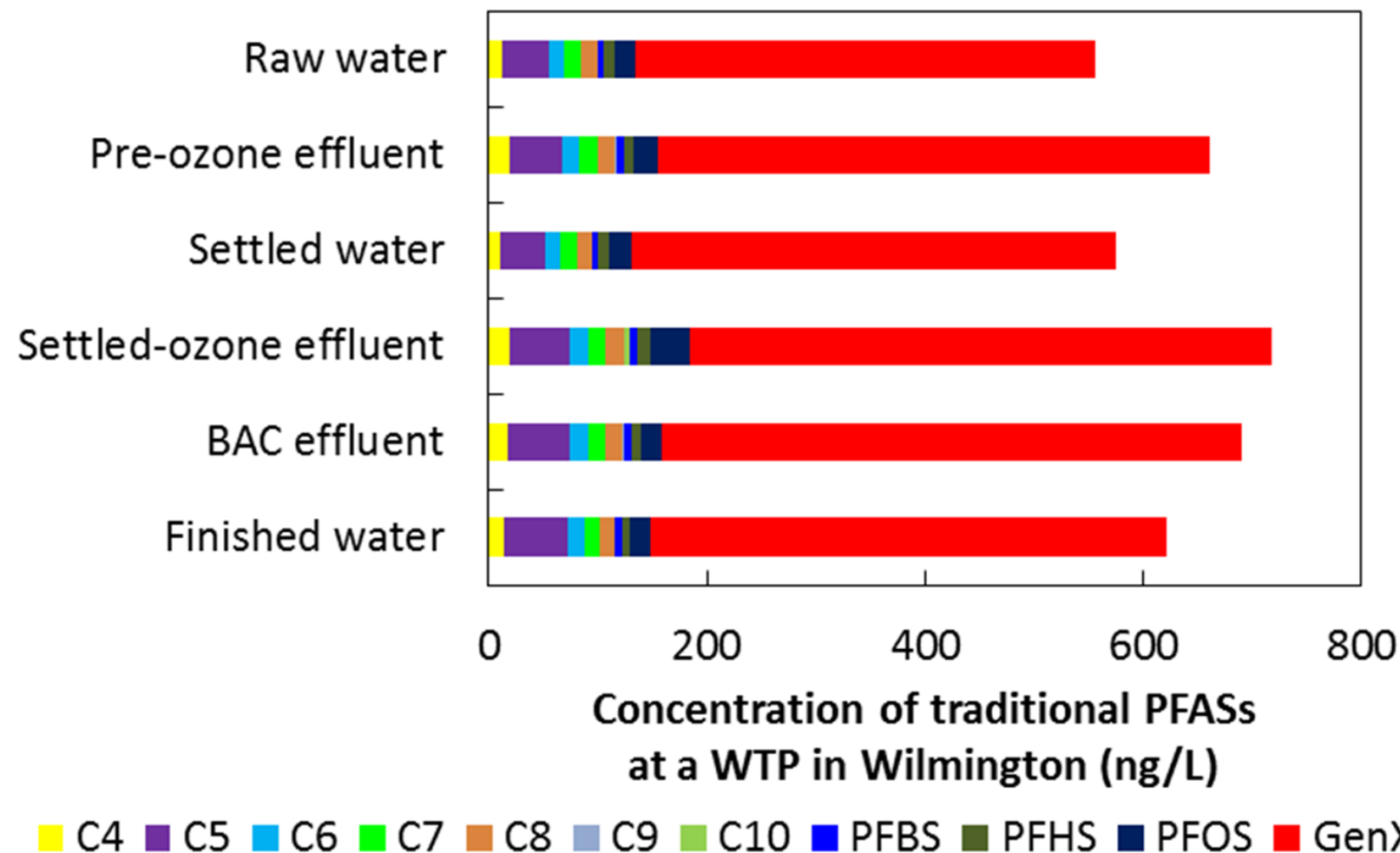
What else

???

???

Are we looking for the right PFASs?

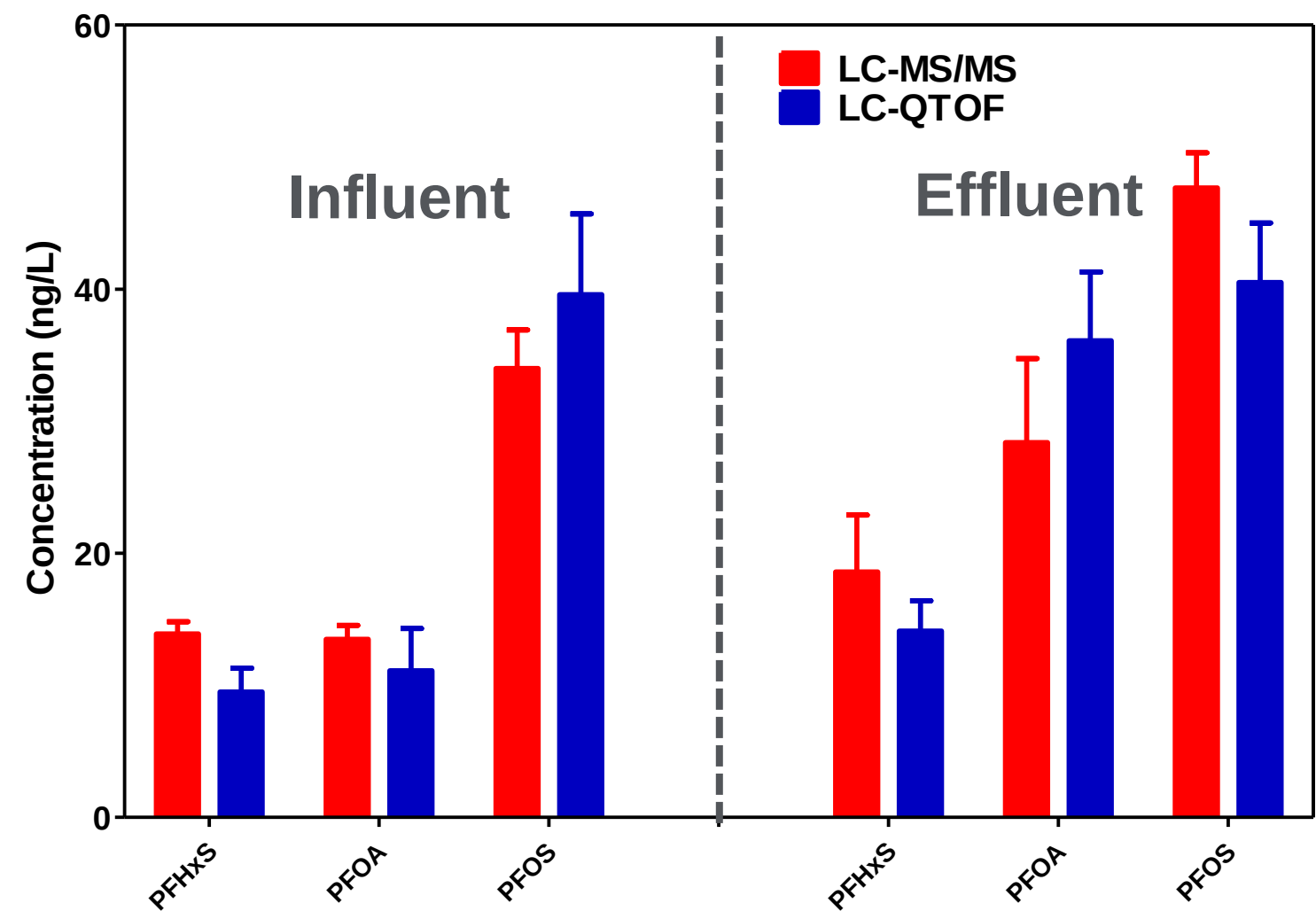
New PFASs being produced and released into the Environment



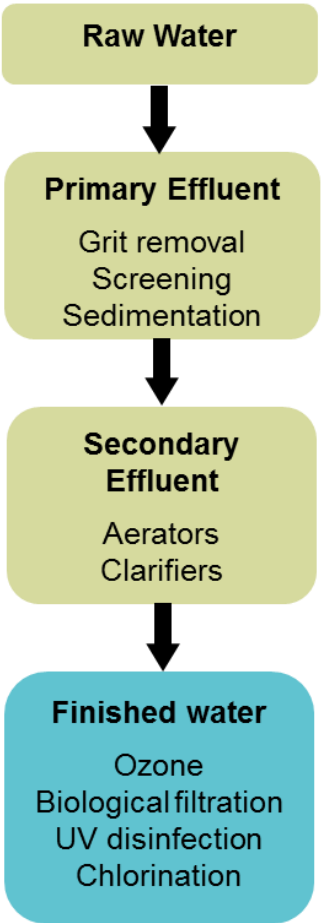
Slide Credit: Dr. Detlef Knappe, North Carolina State University

Targeted Analysis

LC-MS/MS vs LC-Q/TOF



WWTP 1 Activated sludge



Collaborative project
with RMIT University,
Melbourne (Timothy
Coggan ; Prof. Bradley
Clarke)

Suspect Screening Personal Compounds Database and Library with Spectra

- Custom PFAS database with 106 compounds
- MS/MS spectra and Retention time data available for a subset of compounds

MassHunter PCDL Manager for Forensics and Toxicology - C:\MassHunter\PCDL\PFAS with some RT_03282017.cdb

File Edit View PCDL Links Help

Find Spectra

Single Search Batch Search Batch Summary Edit Compounds Spectral Search Browse Spectra Edit Spectra

Mass

Precursor ion: Ion polarity: (Any)

Tolerance: 200 ppm mDa Ionization mode: (Any)

Collision energy

Tolerance: 2.0 eV

Additional Filters

Added Filters

Add Remove

Spectra for compound: PFUnDA / Perfluoroundecanoic acid (PFUnA)

Compound Name	Ion Species	Precursor Ion	CE (V)	Polarity	Ionization	Instrument
PFUnDA / Perfluoroundecanoic acid (PFUnA)	(M-H)-	562.95685	10	Negative	ESI	QTOF
PFUnDA / Perfluoroundecanoic acid (PFUnA)	(M-H)-	562.95685	20	Negative	ESI	QTOF
PFUnDA / Perfluoroundecanoic acid (PFUnA)	(M-H)-	562.95685	40	Negative	ESI	QTOF

Graphic Mass List

Library spectrum

Abundance

m/z

518.96704 100.00

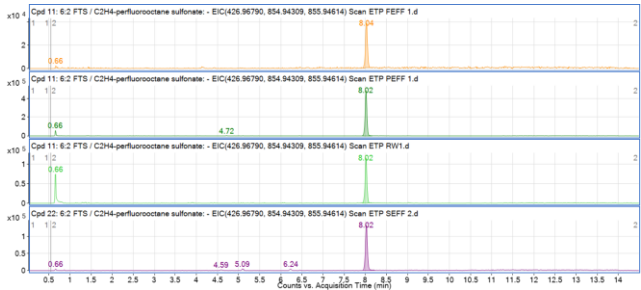
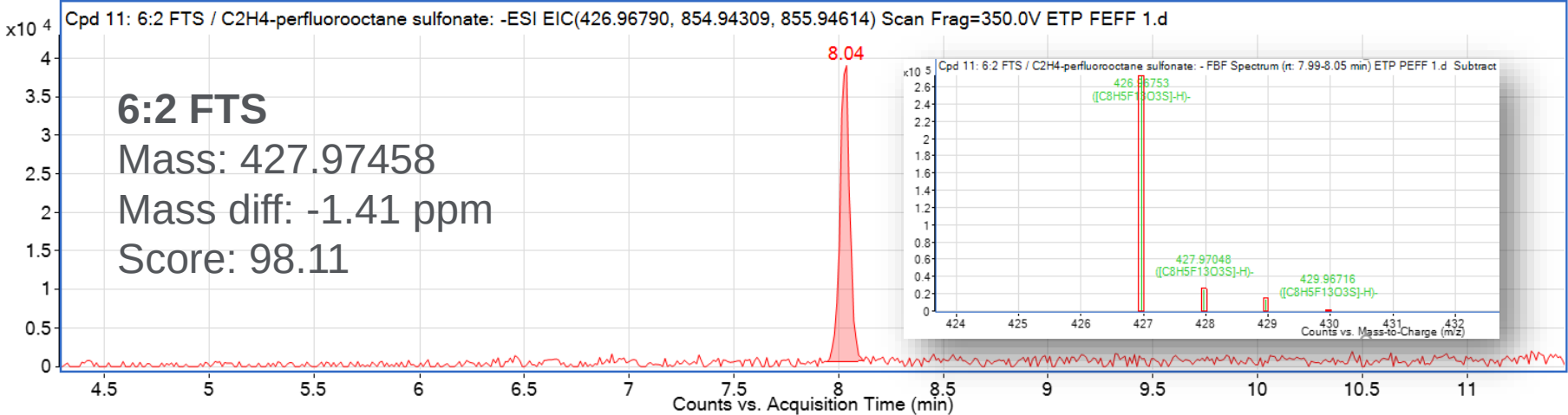
268.98428 10.39

368.97659 1.05

Compound Name	Formula	Mass	Anion	Cation	RT (min)	CAS	ChemSpider	IUPAC Name	Spectra	ChapmanHallID	CH
PFBS / Perfluorobutanesulfonic acid (PFBS)	C4HF9O3S	299.95027	☐	☐	5.660	375-73-5	61132	1,1,2,2,3,3,4,4,4-Nonafluoro-1-butananesulfonic acid	3		
PFHxA / Perfluorohexanoic acid	C6HF11O2	313.98009	☐	☐		307-24-4	60864	Undecafluorohexanoic acid	3		
PFHpA / Perfluoroheptanoic acid	C7HF13O2	363.97690	☐	☐	7.300	375-85-9	61135	Tridecafluoroheptanoic acid	3		
PFHxS / Perfluorohexanesulfonic acid	C6HF13O3S	399.94388	☐	☐	7.350	355-46-4	61053	1,1,2,2,3,3,4,4,4,5,5,6,6,6-Tridecafluoro-1-hexanes...	3		
PFOA / Perfluorooctanoic acid	C8HF15O2	413.97370	☐	☐	8.070	335-67-1	9180	Pentadecafluorooctanoic acid	3		
PFOS / Perfluorooctanesulfonic acid	C8HF17O3S	499.93749	☐	☐	8.730	1763-23-1	67068	1,1,2,2,3,3,4,4,5,5,6,6,7,7,8,8,8-Heptafluoro...	3		
PFDA / Perfluorodecanoic acid	C10HF19O2	513.96732	☐	☐	9.330	335-76-2	9181	Nonadecafluorodecanoic acid	3		
PFUnDA / Perfluoroundecanoic acid (PFUnA)	C11HF21O2	563.96412	☐	☐	9.830	2058-94-8	69649	Henicosfluoroundecanoic acid	3		
PFBA / Perfluorobutanoic acid (Heptafluorobutyri...	C4HF7O2	213.98648	☐	☐	3.370	375-22-4	9394	Heptafluorobutanoic acid	2		

Suspect Screening Personal Compounds Database and Library with Spectra

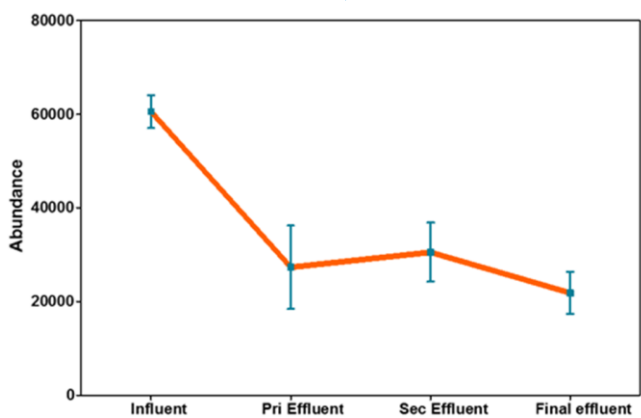
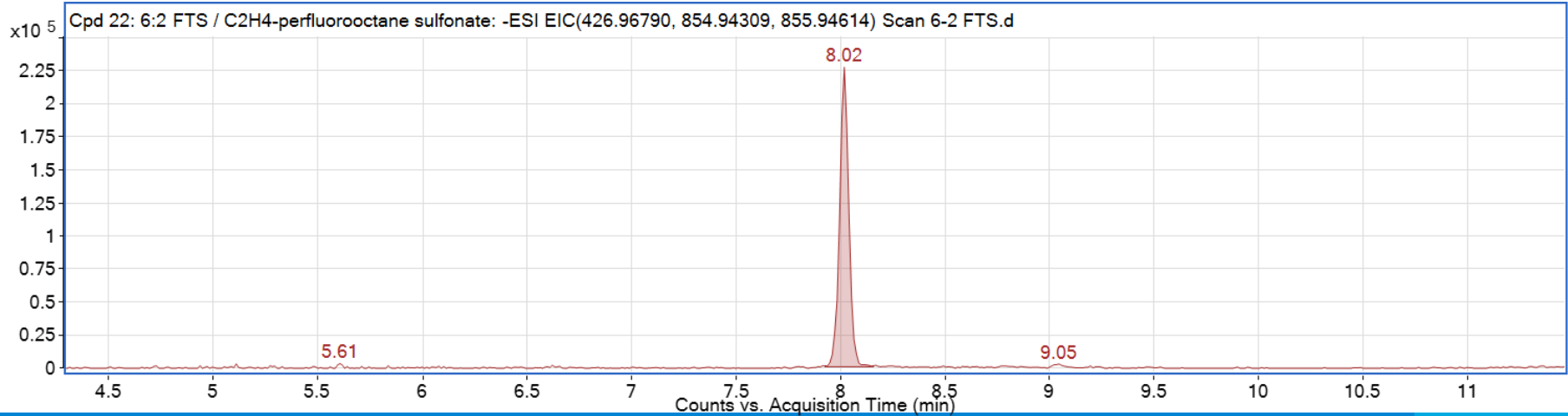
WWTP1 Final Effluent



6:2 FTS ID'd in all samples
of WWTP



Standard



Untargeted Analysis

Find by Auto MS/MS

Find by Targeted MS/MS

Find by Molecular Feature

Find by MRM

Find by Integration



Feature Identification

Method Editor: Find Compounds by Molecular Feature

Find Compounds by Molecular Feature

Extraction Ion Species Charge State Compound Filters Mass Filters **Mass Defect** Peak Filters (MS/MS) Results

Mass defect filtering

☒ Filter results on mass defects

Expected mass defect

Variable

-0.00160 Da + (0.00000 per 19.00 Da)

F

Calculate from formula

Mass defect tolerance

Constant (asymmetric)

+ 0 - 0.09000 Da

Mass Defect for Fluorinated substances



Method Editor: Generate Formulas

Generate Formulas from Compound

Allowed Species Limits Charge State Fragment Formulas

Charge carrier to be assumed if not known

Positive ions:

☒ -electron

☒ +H

☐ +Na

☐ +K

☐ +NH4

☐ +C2H5

☐ +C3H5

Negative ions:

☒ +electron

☒ -H

☐ -Cl

☐ -Br

☐ -COOH

☐ -CF3COO

☐ -CH3COOH

MS ion electron state: allow both even and odd

☒ Group hits with same formula (but different charge carriers)

Elements and limits

Element	Minimum	Maximum
C	3	40
H	0	70
O	0	15
N	0	15
F	1	30



Formula Generation

Chromatogram Results

TCC Scan ETP RW3.d

Scan	Label	Mass	Height	RT	Score (MFE)	Std Dev	Base Peak	Polarity	Max Z	Max L	Start	End	Width	Vol	Vol %	Z	
1	Cap 1 1.25	115.0919	114.1315	4	152.9992	327362	1.25	87	112.9992	Negative	-1	-1	1.25	1.48	0.11	12581	0.74
2	Cap 2 1.36	244.97035	245.05479	2	243.97027	196204	1.36	100	243.97027	Negative	-1	-1	1.25	1.48	0.11	12581	0.74
3	Cap 3 1.41	246.99056	247.28664	4	245.98328	172221	1.41	100	245.98328	Negative	-1	-1	1.27	1.59	0.11	12931	0.76
4	Cap 4 1.42	277.93436	278.47779	5	276.92709	93295	1.42	99.7	276.92709	Negative	-1	-1	1.29	1.52	0.11	877347	0.52
5	Cap 5 1.44	182.99316	183.18791	4	181.99588	274996	1.44	100	181.99588	Negative	-1	-1	1.32	1.6	0.11	20050	11.79
6	Cap 6 1.68	182.99806	183.22036	3	181.99078	181064	1.68	100	181.99078	Negative	-1	-1	1.6	1.76	0.09	11644	0.68
7	Cap 7 1.72	260.97148	261.08226	2	259.96421	119703	1.72	100	259.96421	Negative	-1	-1	1.63	1.8	0.08	894412	0.41
8	Cap 8 1.74	268.98914	268.98973	5	264.92786	1388984	1.74	97.2	264.92786	Negative	-1	-1	1.68	1.78	0.04	42331	2.49
9	Cap 9 1.74	262.98316	263.23173	3	261.97589	1297365	1.74	100	261.97589	Negative	-1	-1	1.69	1.78	0.04	34248	2.01
10	Cap 10 1.75	217.96543	218.29556	4	216.95816	198815	1.75	100	216.95816	Negative	-1	-1	1.67	1.79	0.04	626270	0.37
11	Cap 11 1.82	262.94134	263.23937	3	261.93426	142203	1.82	100	261.93426	Negative	-1	-1	1.77	1.99	0.08	696489	0.41
12	Cap 12 1.87	217.96689	218.24775	3	216.96162	96022	1.87	99.8	216.96162	Negative	-1	-1	1.8	2.01	0.11	681832	0.4
13	Cap 13 2.05	127.96439	128.14546	3	126.95711	293071	2.05	80	126.95711	Negative	-1	-1	1.97	2.21	0.08	14255	0.84
14	Cap 14 2.07	80.96997	81.12375	3	79.9587	702419	2.07	97.6	79.9587	Negative	-1	-1	1.93	2.28	0.13	91455	5.38
15	Cap 15 2.12	374.95666	375.44275	5	373.95079	200937	2.12	100	373.95079	Negative	-1	-1	1.95	2.31	0.13	18120	1.1
16	Cap 16 2.25	260.97685	261.24121	3	259.96958	65494	2.25	82.3	259.96958	Negative	-1	-1	2.16	2.38	0.13	591352	0.35
17	Cap 17 2.43	235.92983	236.31845	5	234.92296	369536	2.43	100	234.92296	Negative	-1	-1	2.29	2.65	0.13	33156	1.95
18	Cap 18 2.51	291.9959	292.42082	5	290.94862	170236	2.51	100	290.94862	Negative	-1	-1	2.37	2.68	0.06	789327	0.48
19	Cap 19 2.63	247.98378	248.07183	2	246.97851	145289	2.63	100	246.97851	Negative	-1	-1	2.47	2.8	0.11	971873	0.57

Feature list with mass defect filters

Untargeted Analysis

Identification of Novel PFAS

Show/Hide	Cpd	Label	Name	Formula	Score	Mass	Avg Mass	Ions	m/z	Height	RT
☒	73	Cpd 73: C9...		C9 H5 F13...	97.81	565.96073	566.05939	3	564.95342	870578	9.82
☒	64	Cpd 64: C8...		C8 H4 F11...	92.3	372.99366	373.03874	3	371.98639	1991326	8.05
☒	69	Cpd 69: C9...		C9 H F15...	90.92	505.94803	506.00974	3	504.94081	201819	8.72
☒	78	Cpd 78: C1...		C11 H2 F2...	85.79	615.95987	616.05175	2	614.95267	126557	10.25
☒	77	Cpd 77: C9...		C9 H2 F17...	84.95	570.96636	571.06182	2	569.95916	137880	10.25
☒	19	Cpd 19: C7...		C7 H3 F N...	84.42	247.98374	248.07183	2	246.97651	145289	2.63
☒	61	Cpd 61: C6...		C6 H2 F8...	83.91	401.94769	402.01514	2	400.94045	574068	7.33
☒	41	Cpd 41: C3...		C3 H F O3	82.62	103.9908	104.14269	3	102.98394	84059	3.95
☒	59	Cpd 59: C7...		C7 H2 F6...	81.12	405.94909	406.00834	2	404.94184	212333	7.33
☒	1	Cpd 1: C6...		C6 H3 F O5	80.14	173.99683	174.17875	4	172.99002	327062	1.25
☒	2	Cpd 2: C5...		C5 H2 F3...	79.17	244.97817	245.05479	2	243.97107	196284	1.36
☒	68	Cpd 68: C6...		C6 H7 F9...	79.14	423.99377	424.03414	2	422.9865	1705215	8.72
☒	5	Cpd 5: C4...		C4 H4 F O7	77.73	182.99288	183.18751	4	181.98588	2744996	1.44
☒	7	Cpd 7: C8...		C8 H F2 N...	76.3	260.97134	261.08326	2	259.96421	118703	1.72
☒	32	Cpd 32: C1...		C12 H4 F2...	75.8	329.98252	330.32489	4	328.97589	98662	3.43
☒	42	Cpd 42: C5...		C5 H F3 N...	74.04	257.97381	258.25562	4	256.9671	184002	3.96
☒	36	Cpd 36: C4...		C4 H4 F O6	74.03	166.99785	167.16441	3	165.99097	135303	3.72
☒	43	Cpd 43: C7...		C7 H F3 N...	74	219.98596	220.24417	4	218.97938	136151	4.03
☒	40	Cpd 40: C1...		C10 H F3...	73.9	329.98497	330.32165	4	328.97832	112221	3.93
☒	39	Cpd 39: C1...		C13 H F2...	73.49	302.98512	303.26718	3	301.9786	114865	3.91
☒	9	Cpd 9: C6...		C6 H2 F N...	71.71	262.98287	263.23173	3	261.97589	1097365	1.74
☒	3	Cpd 3: C8...		C8 H4 F O8	69.64	246.9896	247.28664	4	245.98328	172221	1.41
☒	48	Cpd 48: C8...		C8 H2 F N...	68.78	290.96606	291.26063	3	289.95962	89006	4.39
☒	16	Cpd 16: C5...		C5 H F4 N...	68.15	260.97674	261.24121	3	259.96958	83456	2.23
☒	12	Cpd 12: C4...		C4 H F3 O7	65.37	217.96735	218.24775	3	216.96162	96022	1.87
☒	21	Cpd 21: C7...		C7 H2 F N...	64.49	294.9623	295.31824	5	293.95612	353418	2.67
☒	79	Cpd 79: C3...		C3 H F5 O3	64.01	179.9837	180.27145	4	178.97717	161022	10.6
☒	52	Cpd 52: C7...		C7 H5 F2...	63.42	332.97801	333.3129	3	331.97118	101205	5.44
☒	15	Cpd 15: C8...		C8 H2 F3...	62.58	374.95645	375.44275	5	373.95079	200983	2.12
☒	87	Cpd 87: C4...		C4 H3 F N...	60.74	179.99326	180.24644	3	178.98673	85689	12.55
☒	6	Cpd 6: C7...		C7 H2 F N...	59.54	182.99709	183.22036	3	181.99078	181064	1.68
☒	18	Cpd 18: C5...		C5 H F3 N...	59.02	291.95427	292.42082	5	290.94862	170036	2.51
☒	75	Cpd 75: C3...		C3 H F N O7	52.07	181.97636	182.0593	2	180.96904	133909	9.86
☒	49	Cpd 49: C4...		C4 H F3 N...	51.82	263.95848	264.38963	5	262.95261	104212	4.52
☒	63	Cpd 63: PF...	PFOA 13...	C4 [13C]4...	49.76	417.98671	418.04076	4	416.97945	886238	8.05
☒	67	Cpd 67: PF...	PFNA 13...	C4 [13C]5...	49.56	468.98666	469.03418	4	467.9794	978955	8.72

Inclusion list for targeted MS/MS and structural elucidation using fragment information using software tools (Molecular structure correlator)

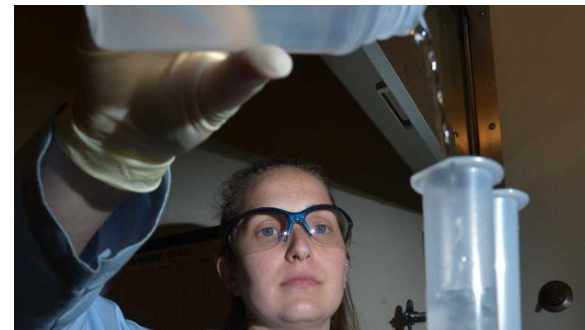
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