

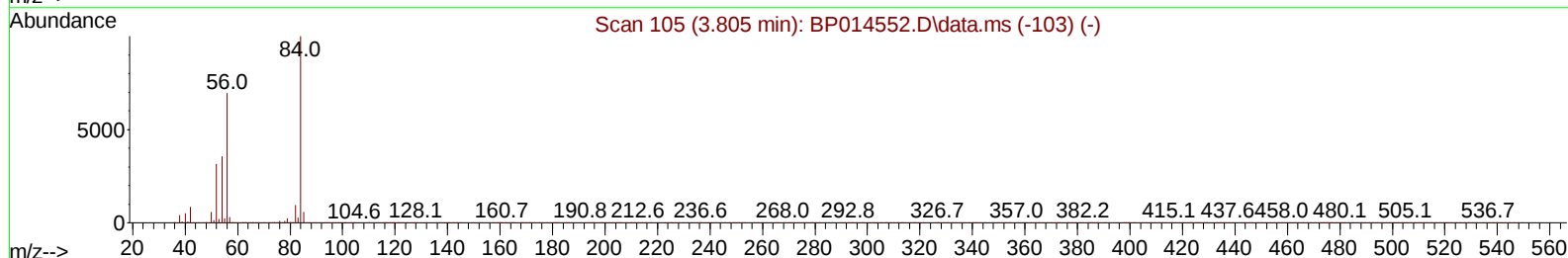
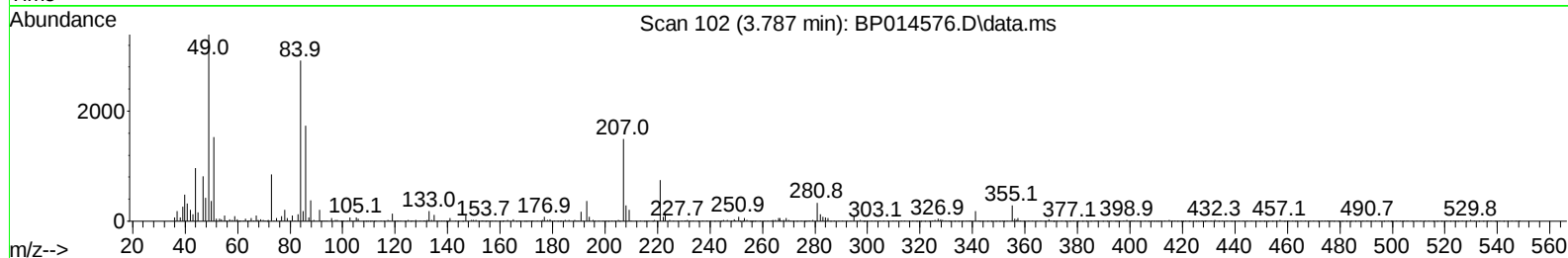
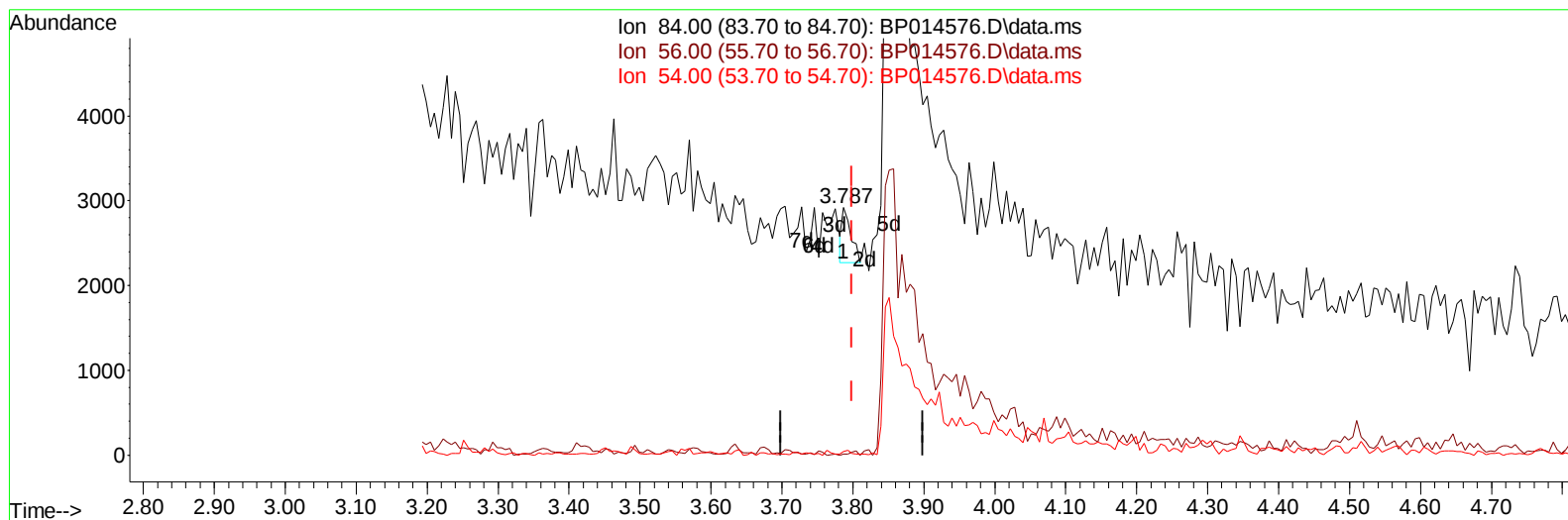
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014576.D
 Acq On : 06 Apr 2023 02:42
 Operator : CG/JU
 Sample : 02194-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHDT9

Manual IntegrationsAPPROVED

Quant Time: Apr 06 03:30:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 06 00:55:13 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/06/2023
 Supervised By :Jagrut Upadhyay 04/06/2023



TIC: BP014576.D\data.ms

(4) Pyridine-d5 (S)

3.787min (-0.012) 0.08 ng/u1

response 571

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.45#
54.00	32.60	1.37#
0.00	0.00	0.00

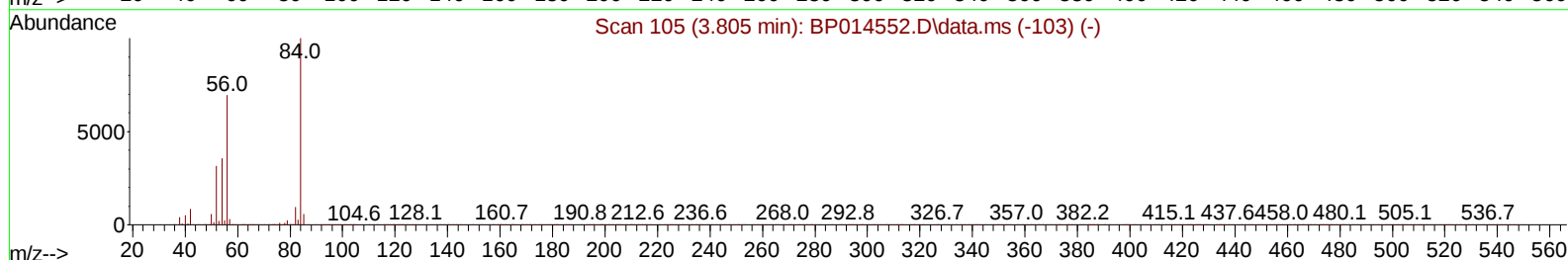
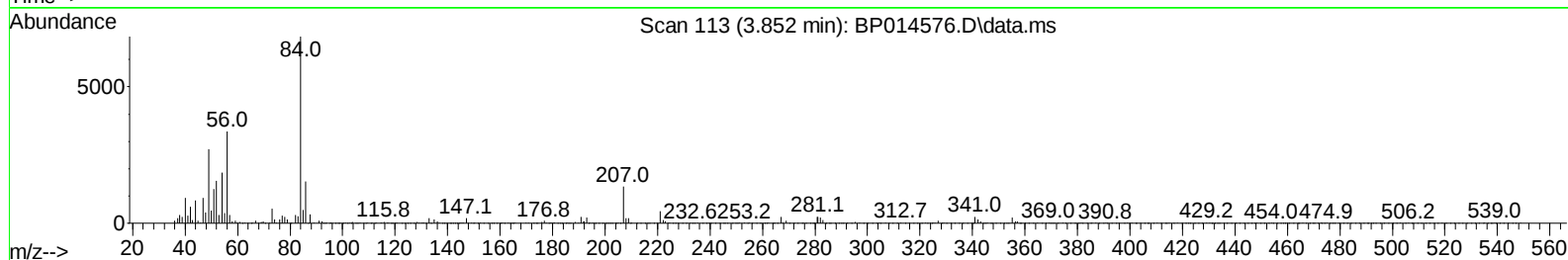
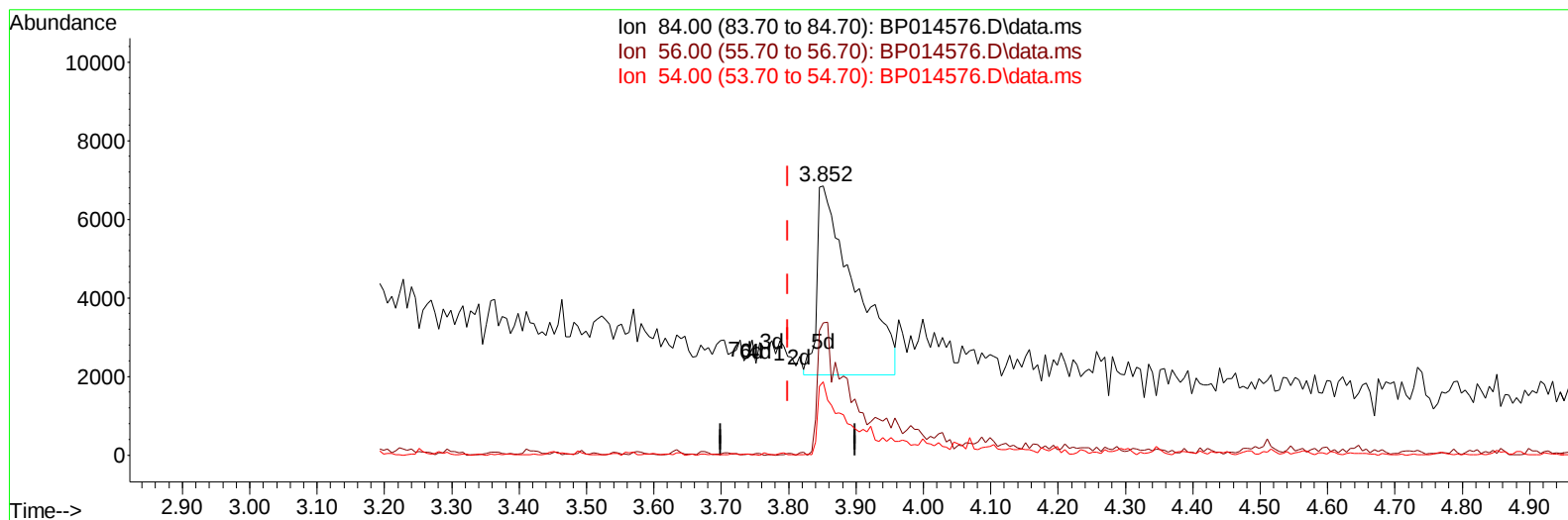
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(4) Pyridine-d5 (S)

3.852min (+ 0.053) 2.53 ng/u1 m

response 18372

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	49.17#
54.00	32.60	27.14
0.00	0.00	0.00

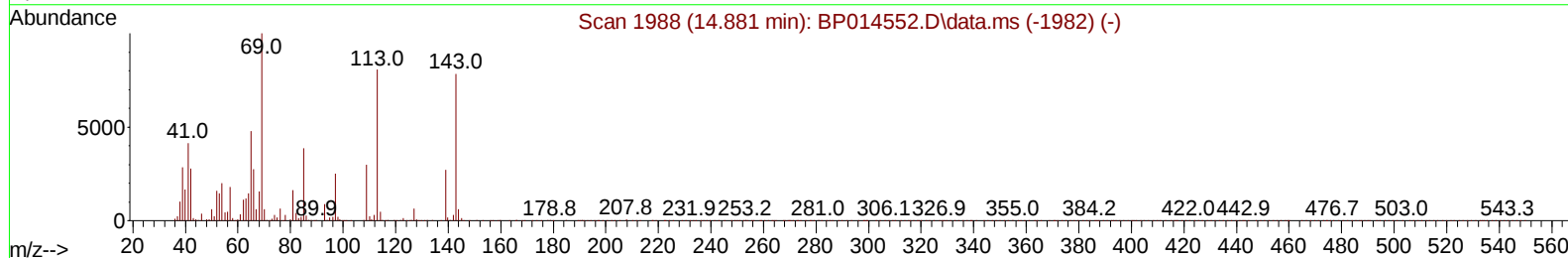
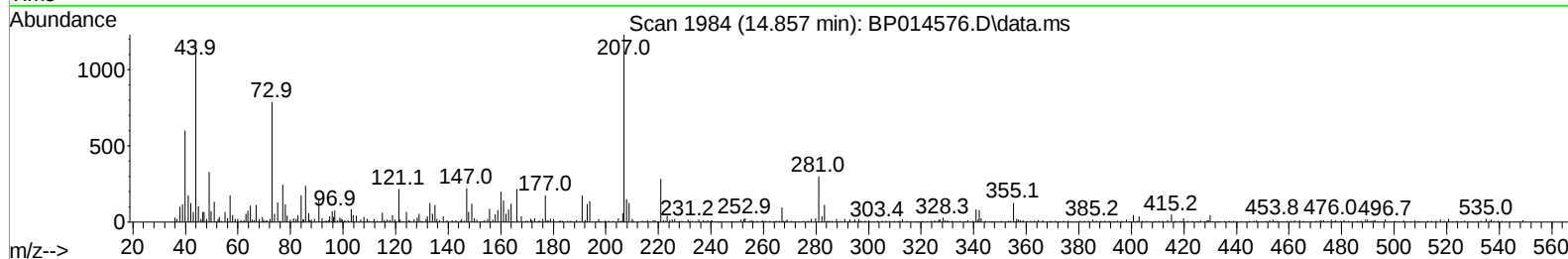
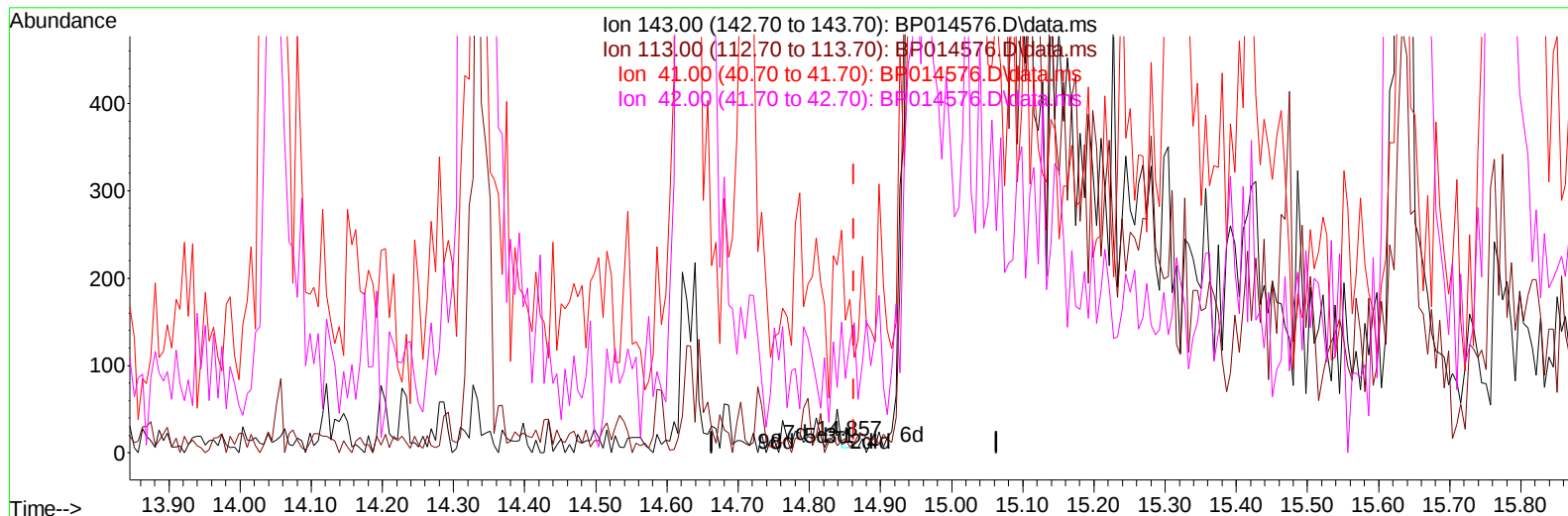
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Sample : 02194-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014576.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.006) 0.00 ng/u1

response 5

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	50.00#
41.00	44.20	1257.14#
42.00	31.50	907.14#

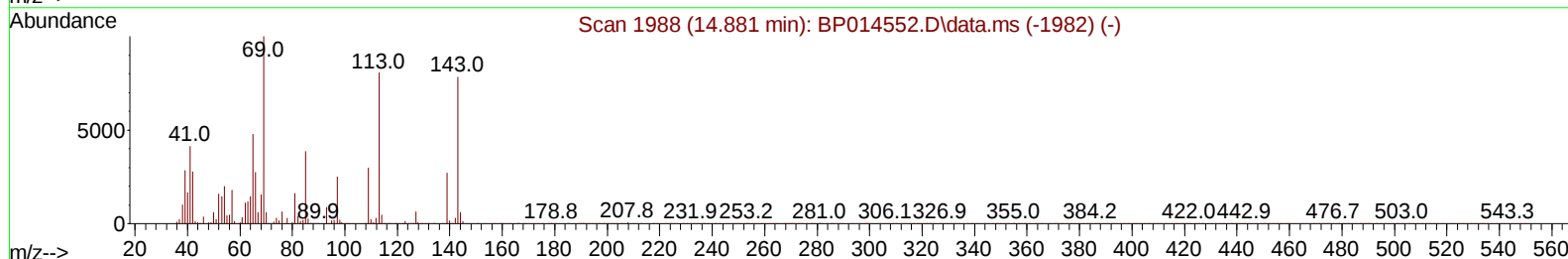
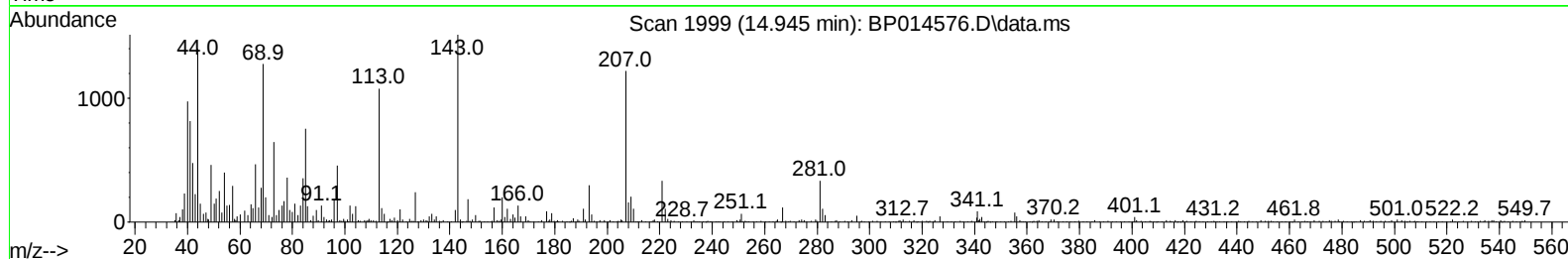
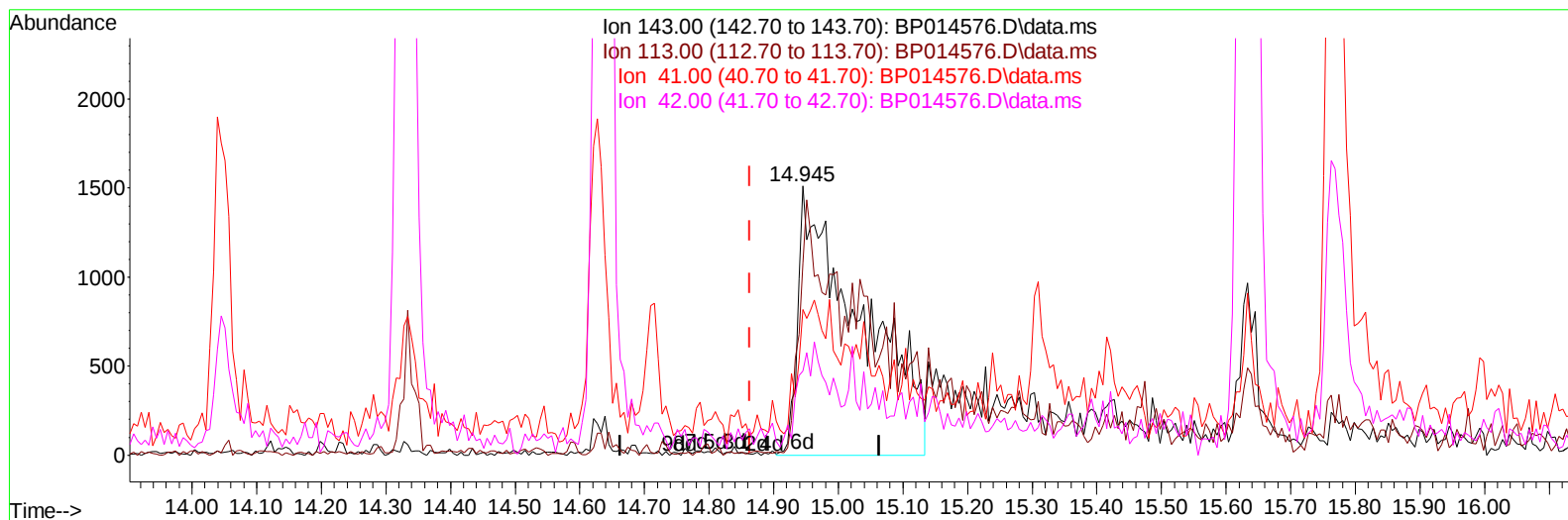
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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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TIC: BP014576.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.082) 2.33 ng/ul m

response 10034

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	71.34#
41.00	44.20	54.07#
42.00	31.50	31.37

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	109897	20.000	ng/ul	0.00
20) Naphthalene-d8	10.804	136	518093	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	352028	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	763957	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	626697	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	665332	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	12221	4.312	ng/uL	0.00
4) Pyridine-d5	3.852	84	18372m	2.530	ng/ul	0.05
7) Phenol-d5	7.163	99	41052	4.050	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.310	67	184243	28.333	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	139615	19.059	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	86810	10.618	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	100146	23.944	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	102126	21.324	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	171591	19.673	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	147581	12.846	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	789267	27.652	ng/ul	0.00
49) Acenaphthylene-d8	14.333	160	848525	26.791	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	10034m	2.327	ng/ul	0.08
60) Fluorene-d10	15.633	176	661412	27.787	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	101513	18.059	ng/ul	0.00
73) Anthracene-d10	17.498	188	1163371	31.466	ng/ul	0.00
81) Pyrene-d10	19.727	212	1280365	30.416	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1122580	31.757	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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