

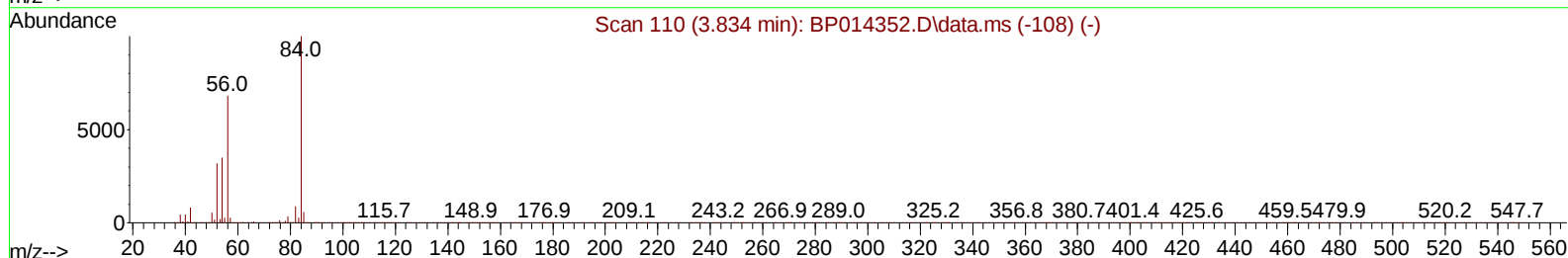
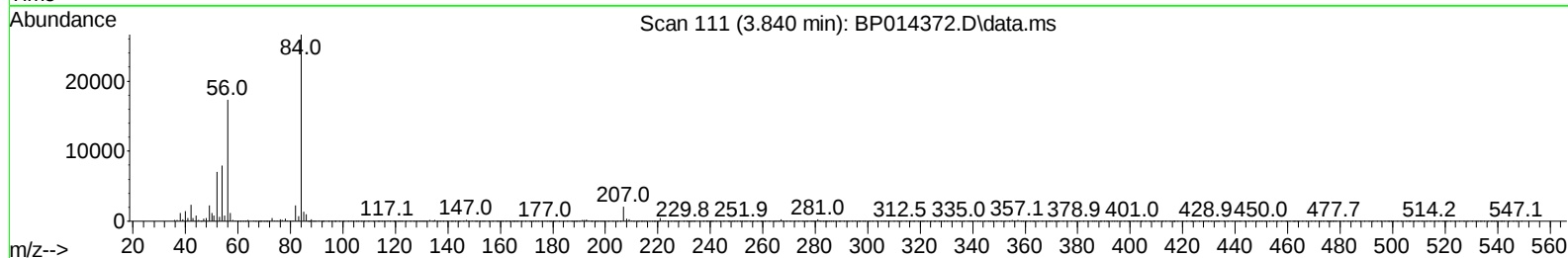
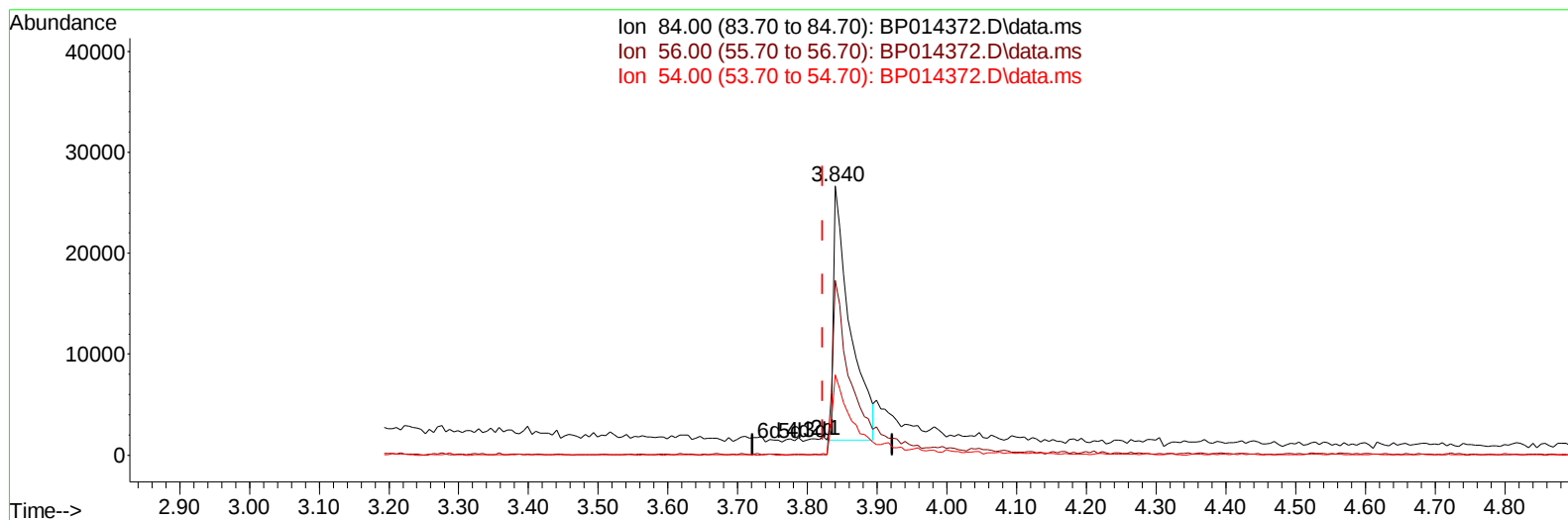
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032923\
 Data File : BP014372.D
 Acq On : 29 Mar 2023 20:26
 Operator : CG/JU
 Sample : 02067-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFZ79

Manual IntegrationsAPPROVED

Quant Time: Mar 30 01:05:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 30 00:56:20 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/30/2023
 Supervised By :Jagrut Upadhyay 03/30/2023



TIC: BP014372.D\data.ms

(4) Pyridine-d5 (S)

3.840min (+ 0.018) 3.72 ng/u1

response 40760

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	64.94
54.00	32.60	29.78
0.00	0.00	0.00

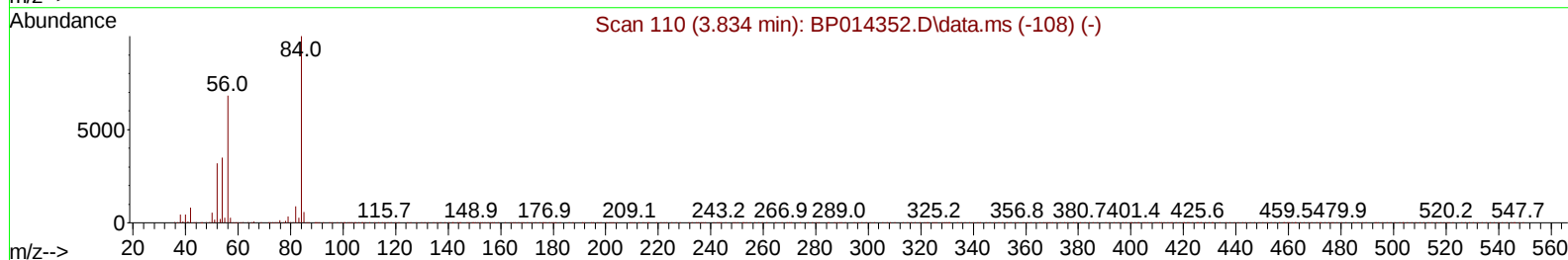
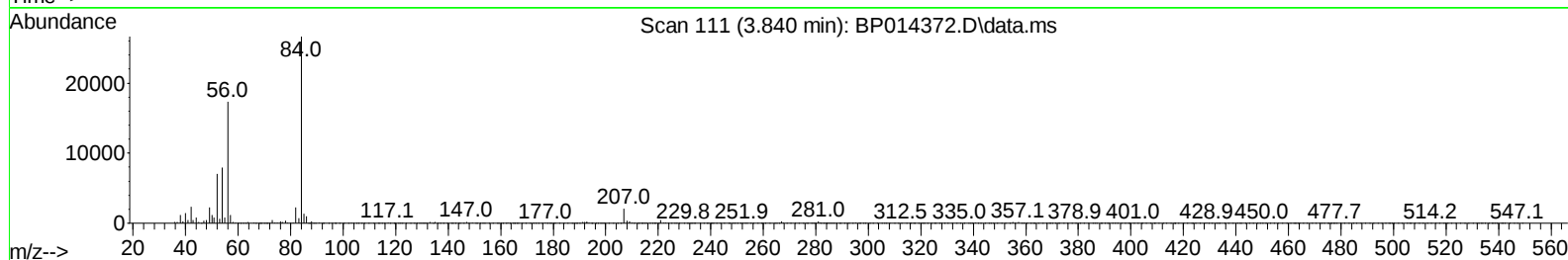
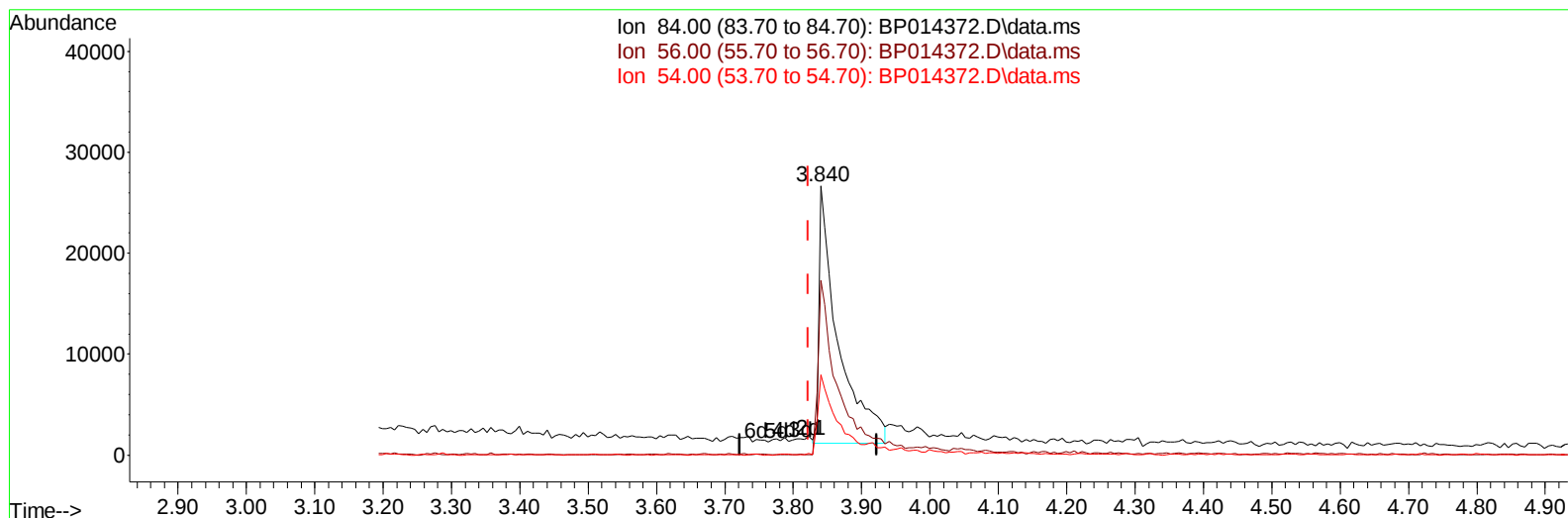
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(4) Pyridine-d5 (S)

3.840min (+ 0.018) 4.49 ng/ul m

response 49216

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	64.94
54.00	32.60	29.78
0.00	0.00	0.00

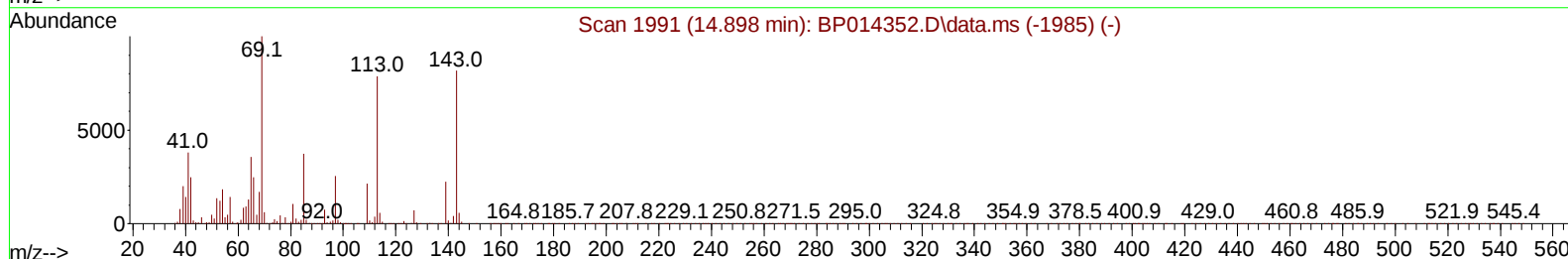
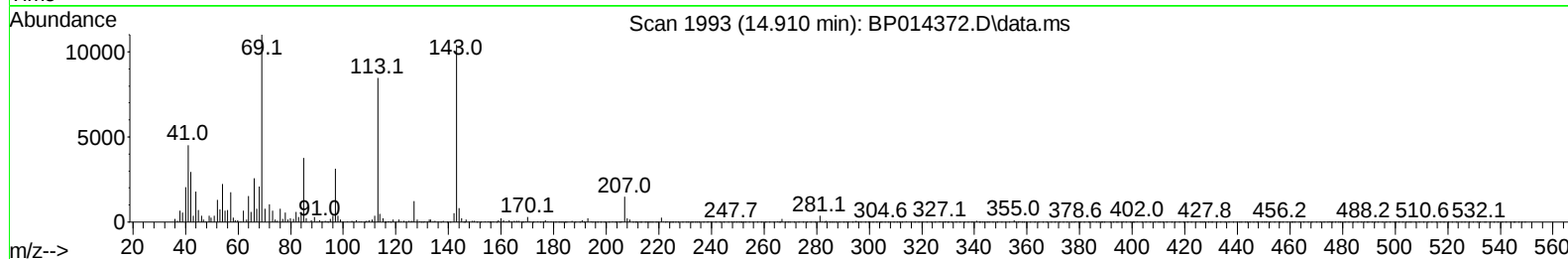
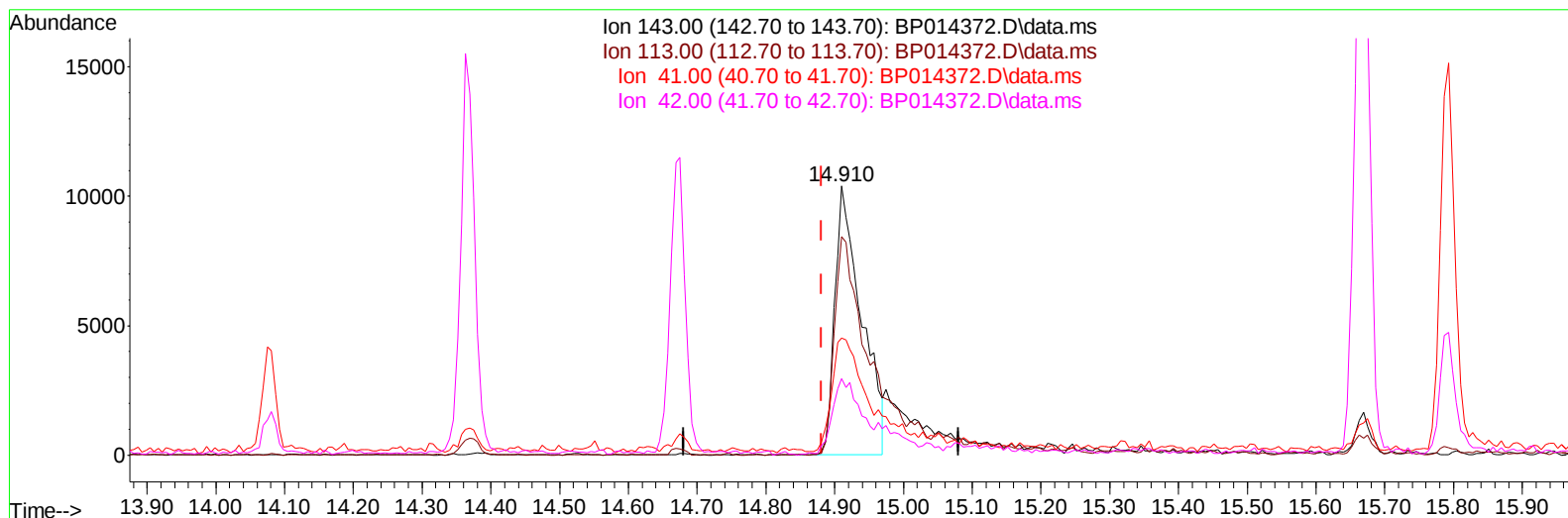
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TIC: BP014372.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.910min (+ 0.030) 4.65 ng/u1****response 27892**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	81.32
41.00	44.20	43.49
42.00	31.50	28.44

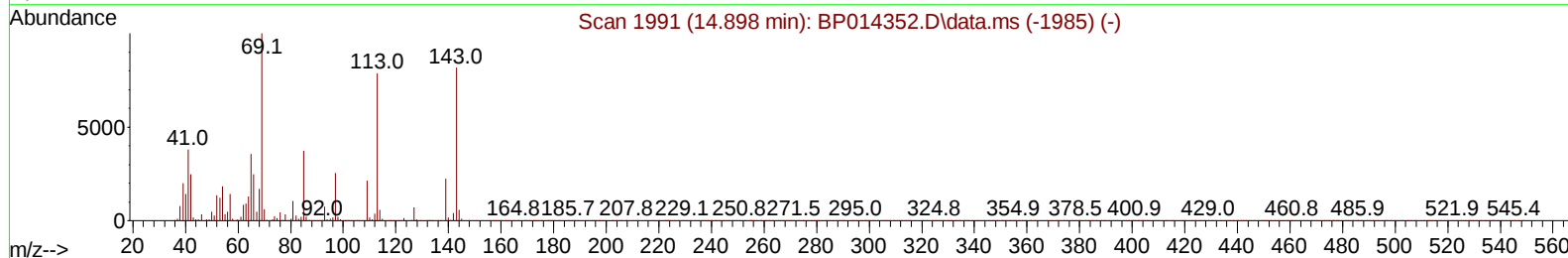
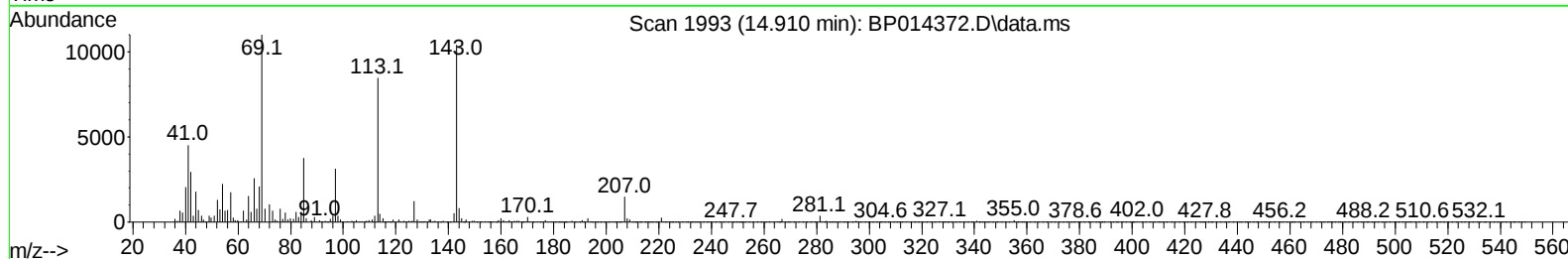
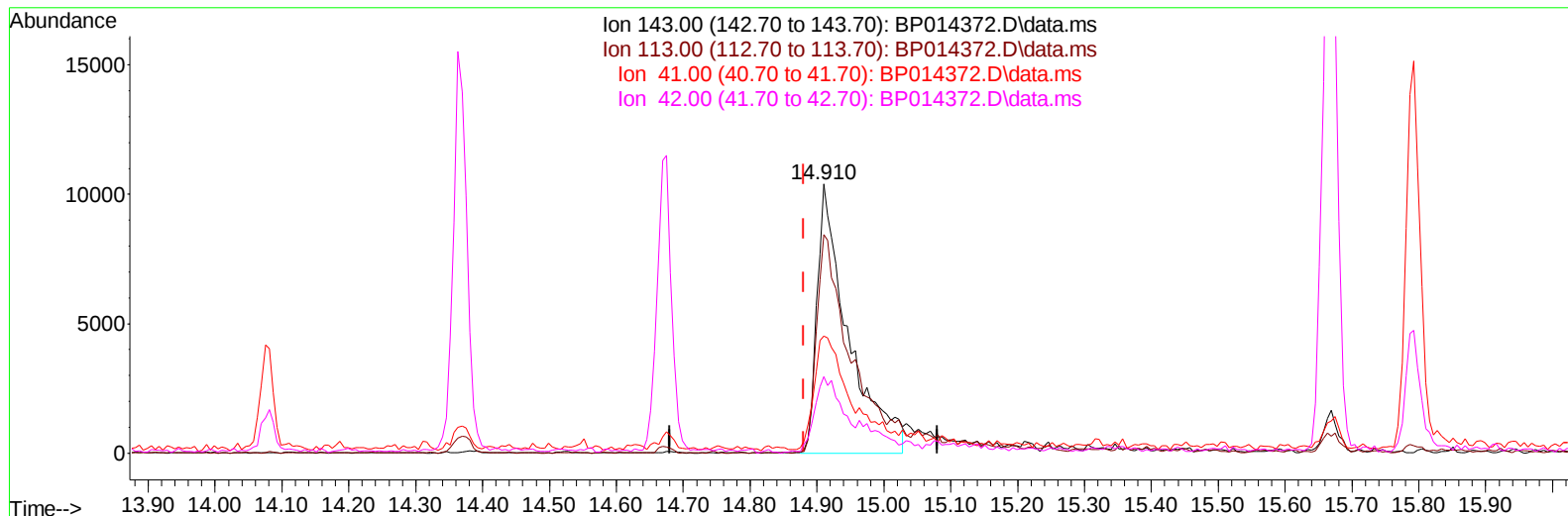
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Sample : 02067-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP014372.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.030) 5.63 ng/ul m

response 33783

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	81.32
41.00	44.20	43.49
42.00	31.50	28.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	165725	20.000	ng/ul	0.00
20) Naphthalene-d8	10.846	136	729562	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	490009	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.434	188	1051887	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	870636	20.000	ng/ul	0.00
88) Perylene-d12	24.045	264	767277	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	20894	4.889	ng/uL	0.00
4) Pyridine-d5	3.840	84	49216m	4.494	ng/ul	0.02
7) Phenol-d5	7.181	99	115044	7.527	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	360934	36.806	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	309633	28.029	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	227122	18.422	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	206857	35.122	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	235471	34.915	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	386842	31.496	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	230230	14.231	ng/ul	0.00
46) Dimethylphthalate-d6	14.081	166	1538113	38.713	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	1577228	35.777	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	33783m	5.628	ng/ul	0.03
60) Fluorene-d10	15.669	176	1275342	38.492	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.793	200	263000	33.980	ng/ul	0.00
73) Anthracene-d10	17.534	188	2023465	39.748	ng/ul	0.00
81) Pyrene-d10	19.763	212	2330940	39.858	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	1693255	41.537	ng/ul	0.00

Target Compounds Qvalue

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

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