

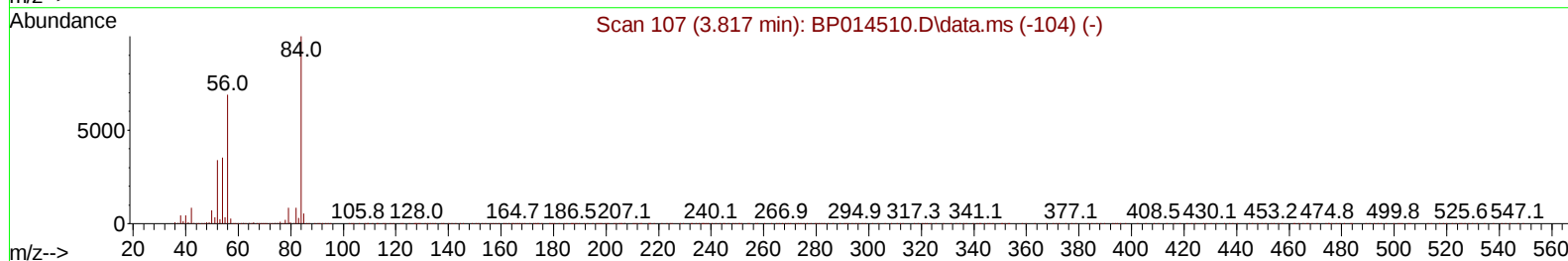
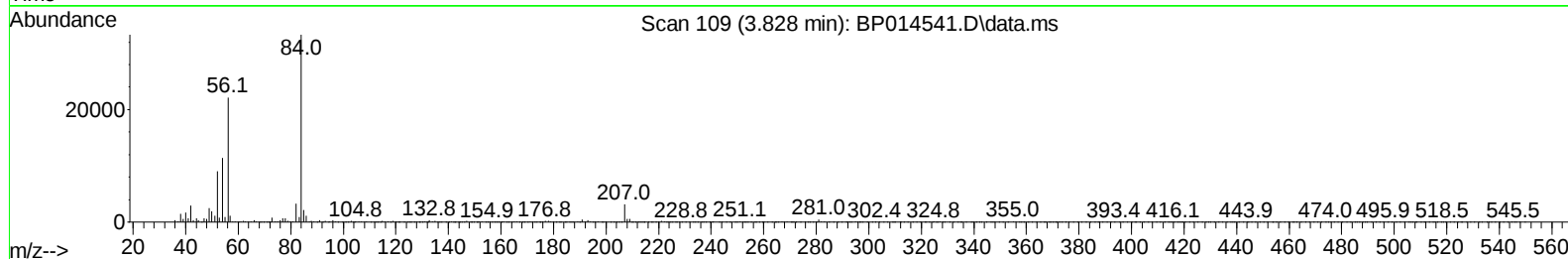
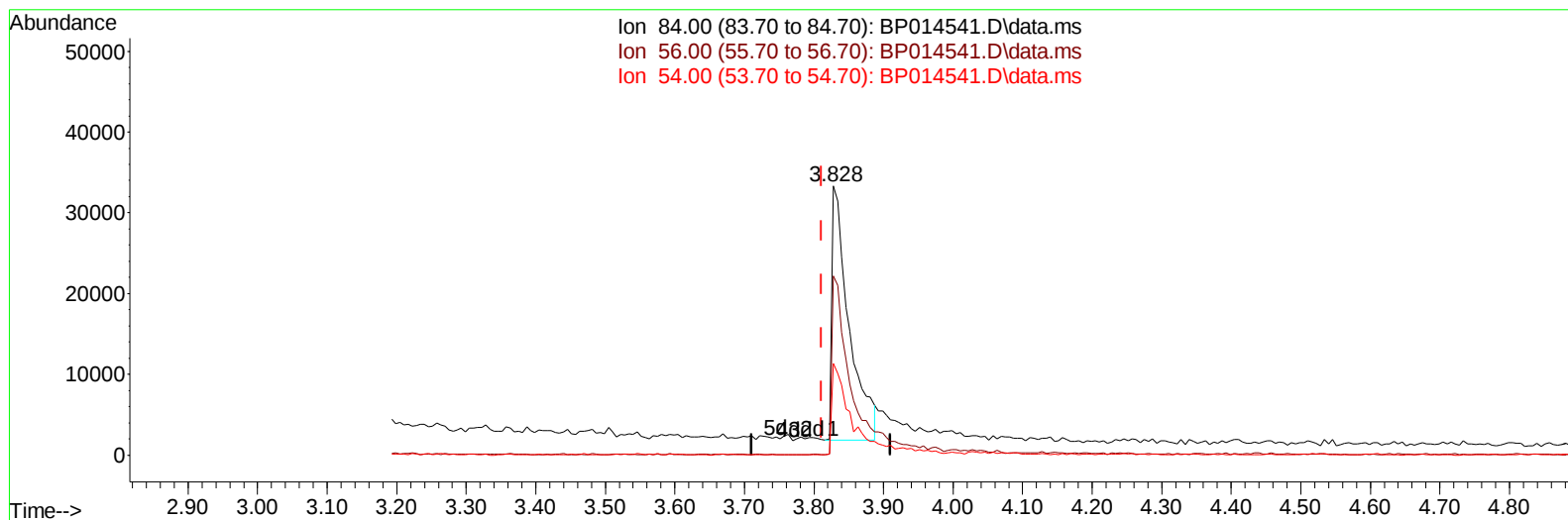
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014541.D
 Acq On : 04 Apr 2023 22:34
 Operator : CG/JU
 Sample : 02158-24
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-1(20230331)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:48:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 04 23:43:56 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/05/2023
 Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014541.D\data.ms

(4) Pyridine-d5 (S)

3.828min (+ 0.017) 6.77 ng/u1

response 53988

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	66.51
54.00	32.60	34.10
0.00	0.00	0.00

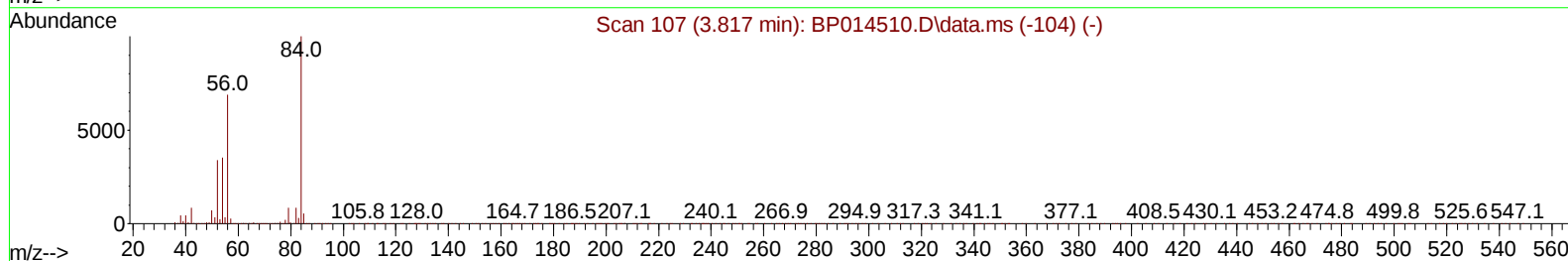
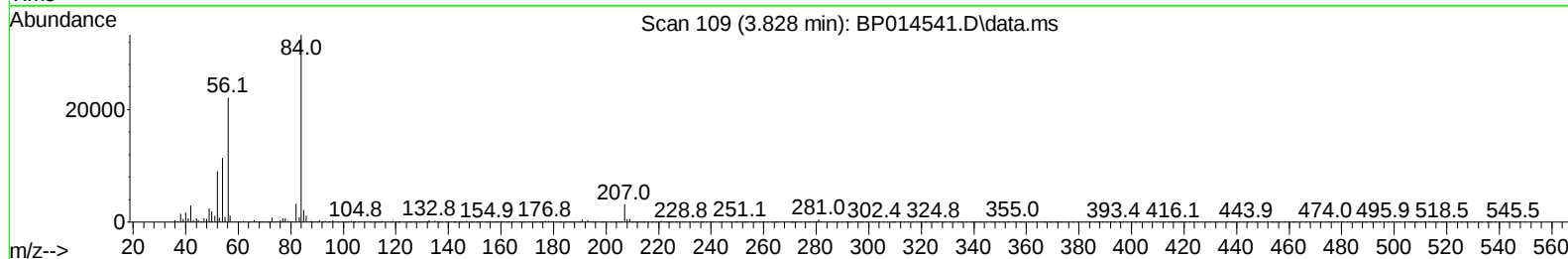
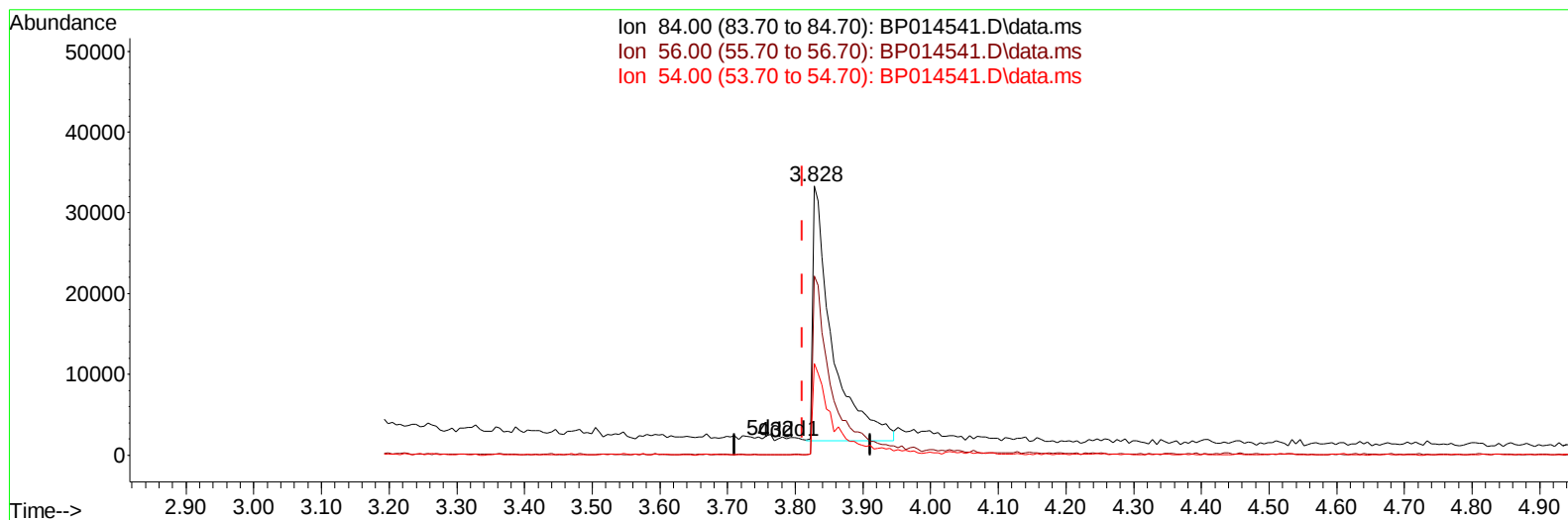
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TIC: BP014541.D\data.ms

(4) Pyridine-d5 (S)

3.828min (+ 0.017) 7.92 ng/u1 m

response 63151

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	66.51
54.00	32.60	34.10
0.00	0.00	0.00

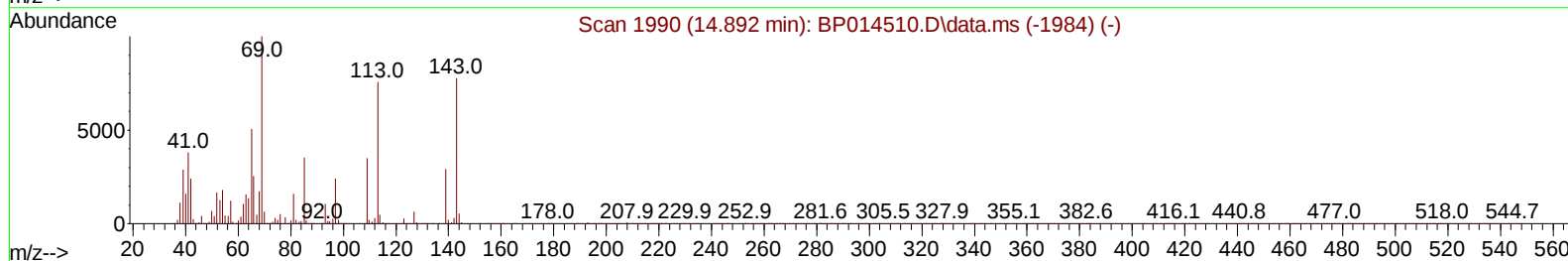
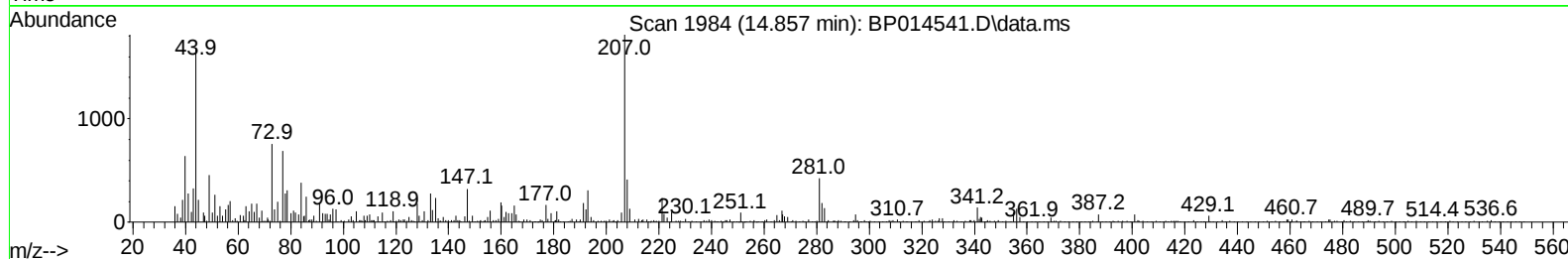
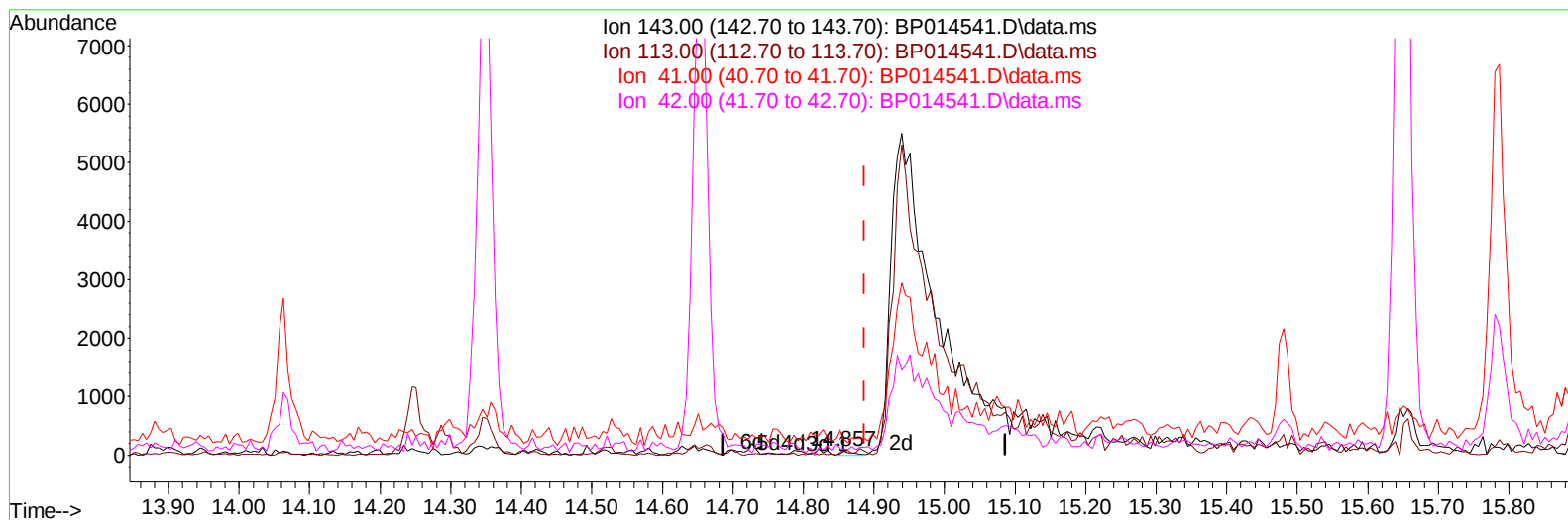
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Instrument :
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TIC: BP014541.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.030) 0.02 ng/u1

response 92

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.33#
41.00	44.20	370.67#
42.00	31.50	133.33#

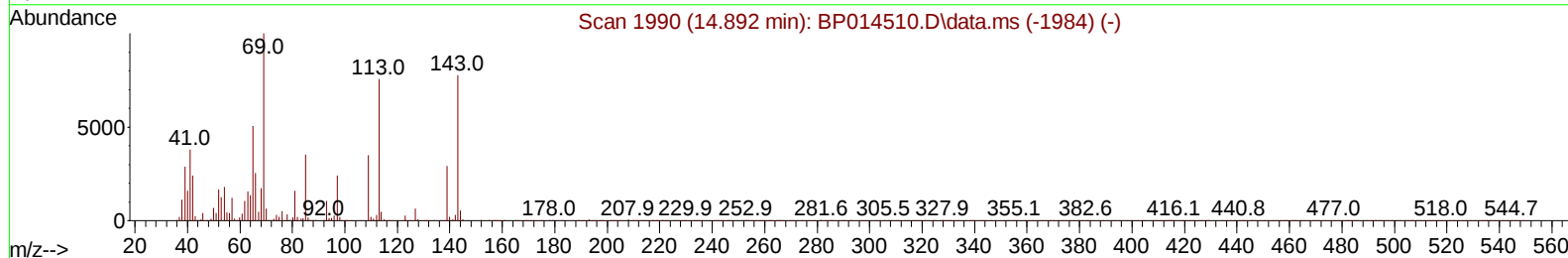
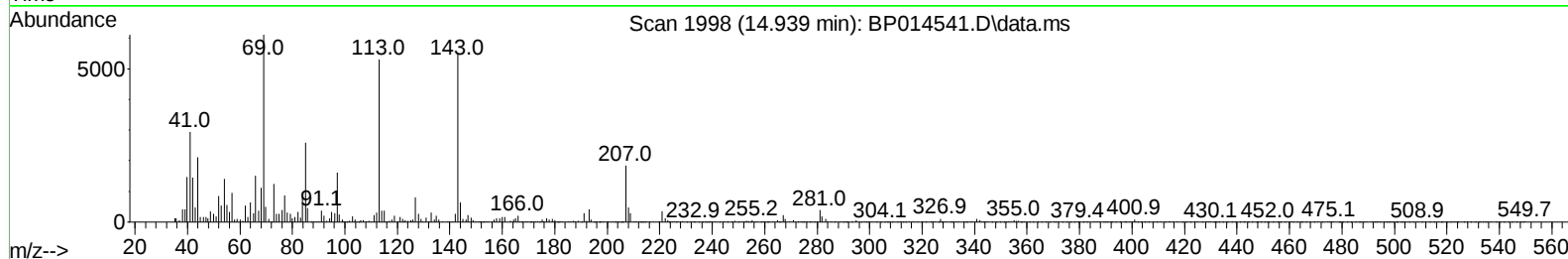
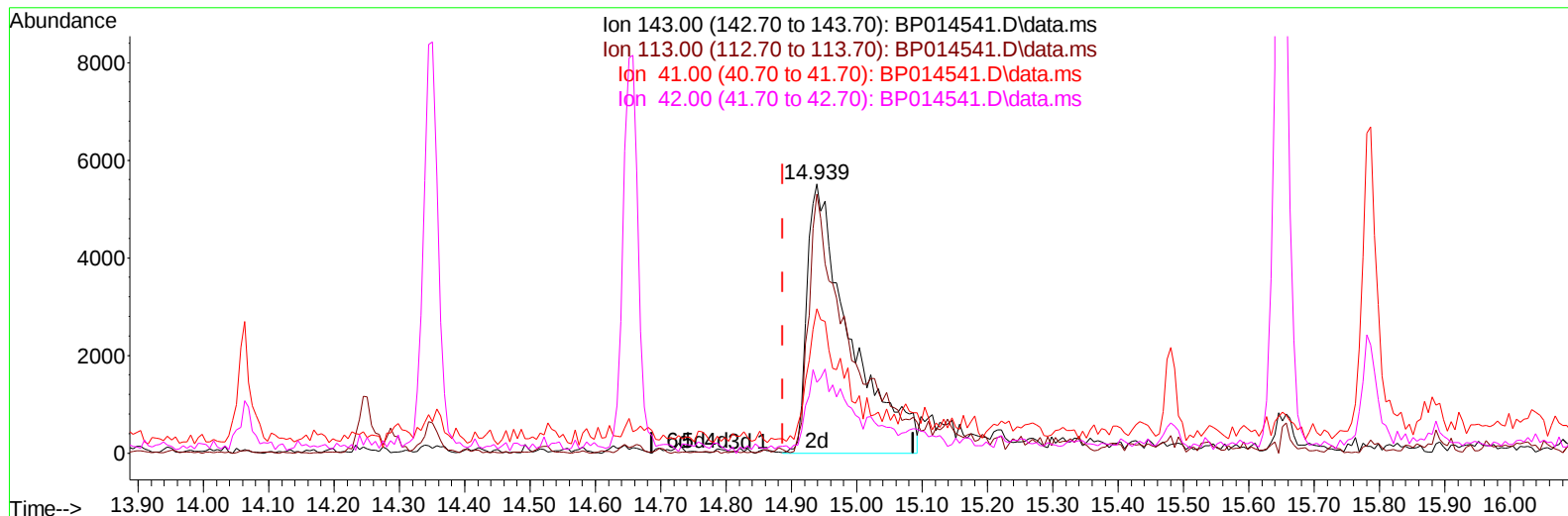
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Instrument :
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TIC: BP014541.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (+ 0.053) 6.06 ng/ul m

response 24986

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	96.41
41.00	44.20	53.55#
42.00	31.50	26.21

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	120594	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	510997	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	336376	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	797436	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	834690	20.000	ng/ul	0.00
88) Perylene-d12	24.009	264	956327	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13954	4.487	ng/uL	0.00
4) Pyridine-d5	3.828	84	63151m	7.924	ng/ul	0.02
7) Phenol-d5	7.181	99	72719	6.538	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.328	67	203259	28.484	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	184842	22.995	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	131266	14.632	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	116671	28.283	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	128016	27.101	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	202666	23.559	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	250392	22.097	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	786258	28.828	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	868544	28.700	ng/ul	0.00
54) 4-Nitrophenol-d4	14.939	143	24986m	6.064	ng/ul	0.05
60) Fluorene-d10	15.651	176	688290	30.262	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	113620	19.364	ng/ul	0.00
73) Anthracene-d10	17.516	188	1248007	32.338	ng/ul	0.00
81) Pyrene-d10	19.745	212	1551766	27.678	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.850	264	1643448	32.346	ng/ul	0.00

Target Compounds Qvalue

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

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