

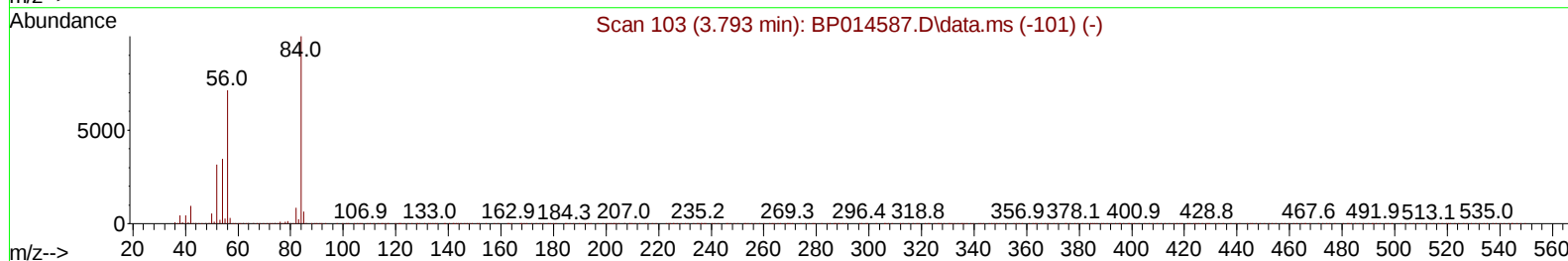
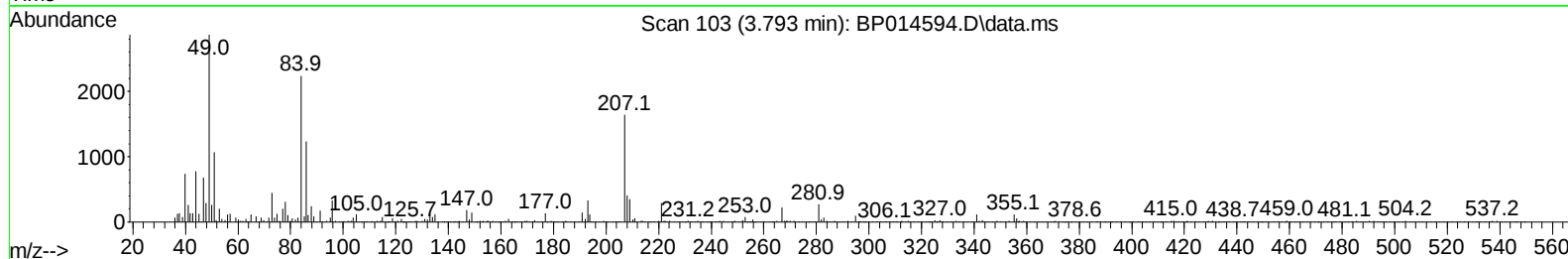
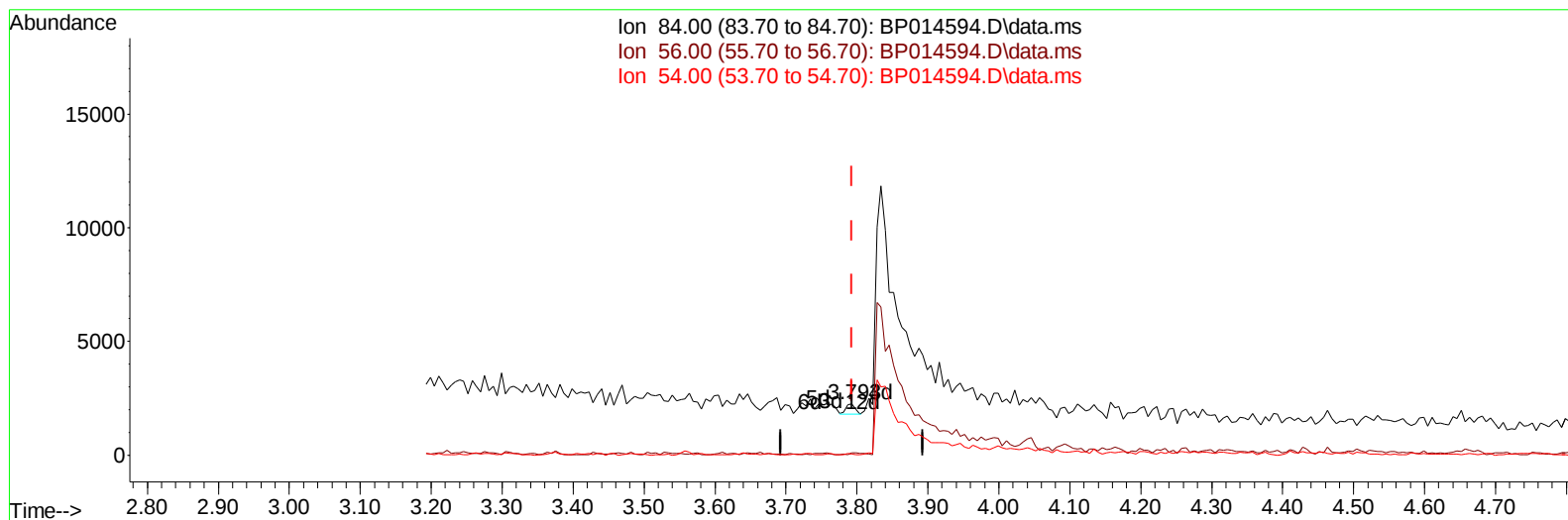
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014594.D
 Acq On : 06 Apr 2023 18:57
 Operator : CG/JU
 Sample : 02194-06
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHDW4

Manual IntegrationsAPPROVED

Quant Time: Apr 07 00:47:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 07 00:45:34 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/07/2023
 Supervised By :Jagrut Upadhyay 04/07/2023



TIC: BP014594.D\data.ms

(4) Pyridine-d5 (S)

3.793min (-0.000) 0.04 ng/u1

response 383

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	5.05#
54.00	32.60	2.14#
0.00	0.00	0.00

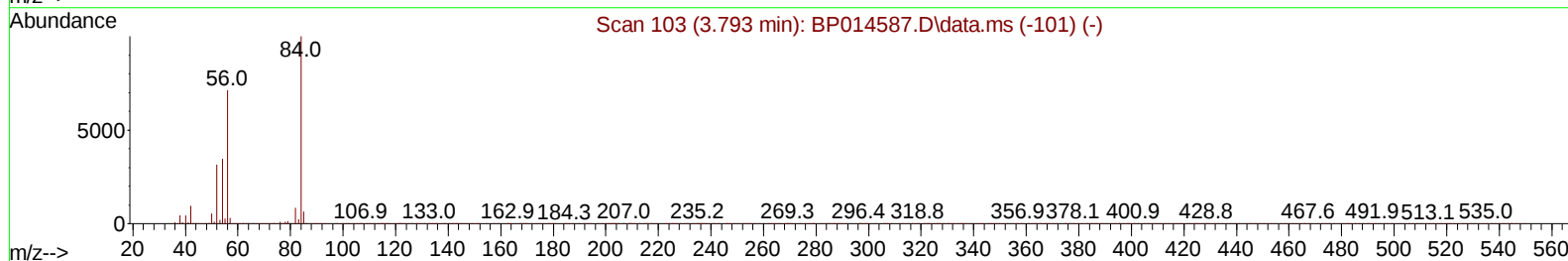
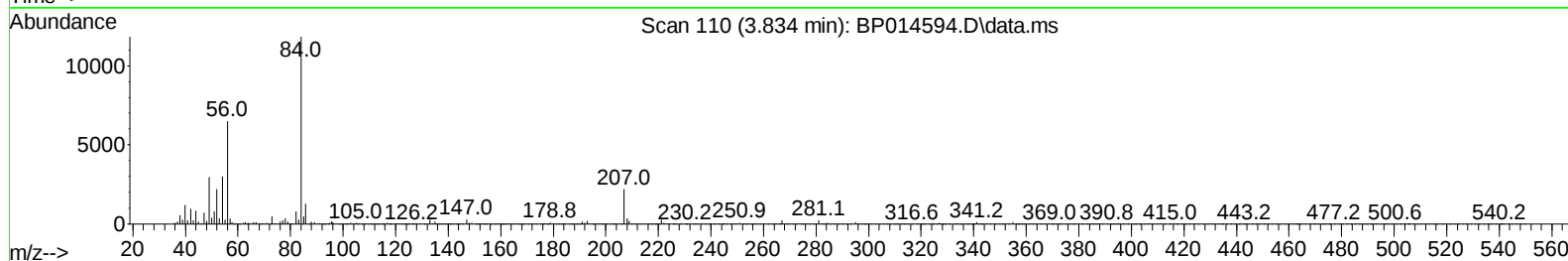
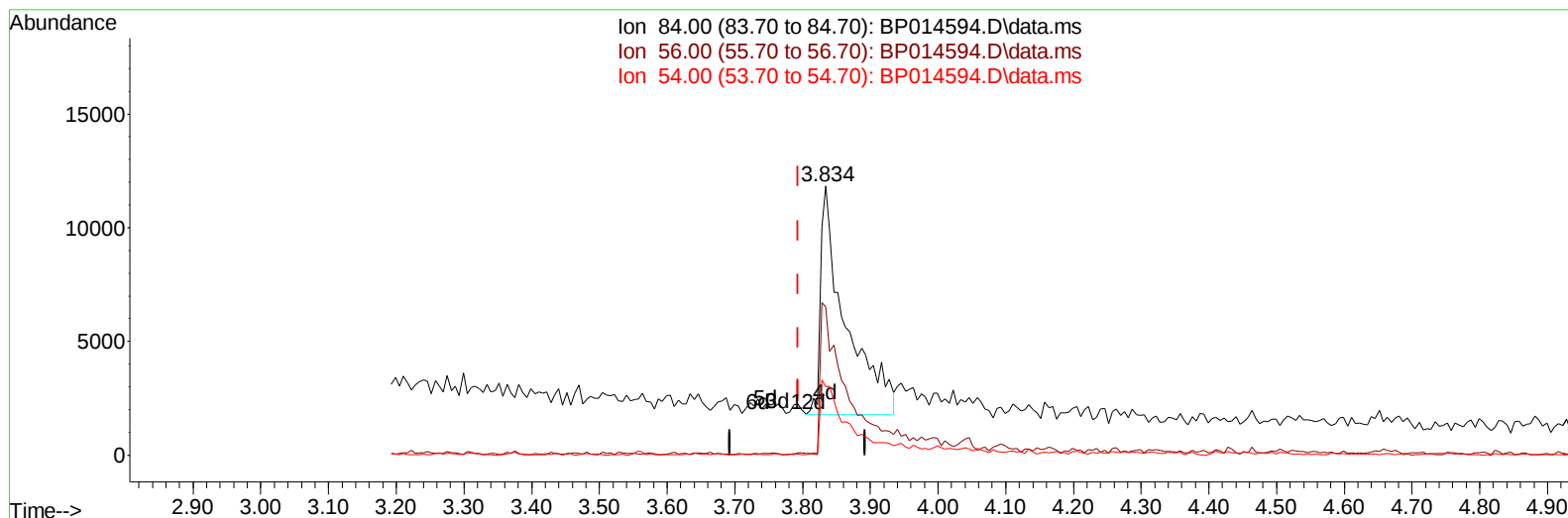
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(4) Pyridine-d5 (S)

3.834min (+ 0.041) 2.97 ng/ul m

response 25811

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	55.01
54.00	32.60	25.39#
0.00	0.00	0.00

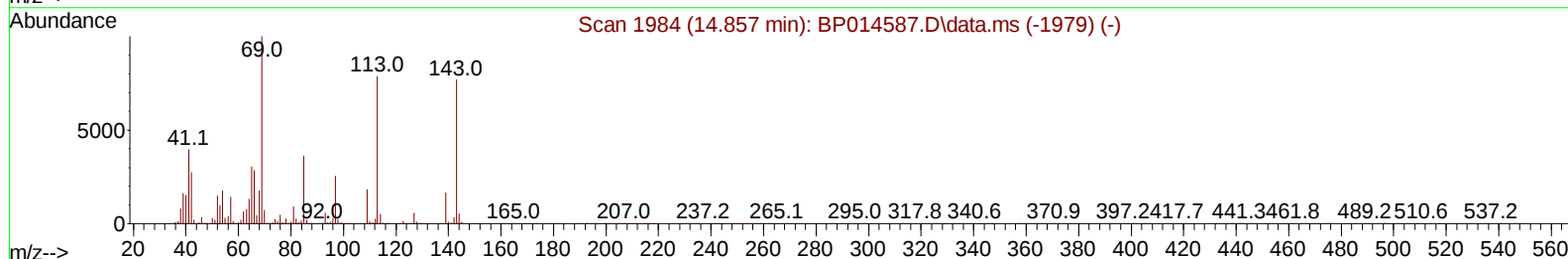
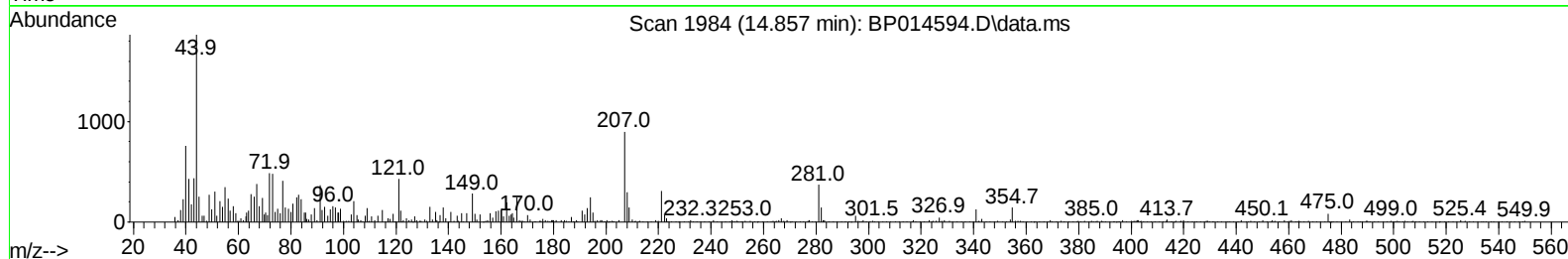
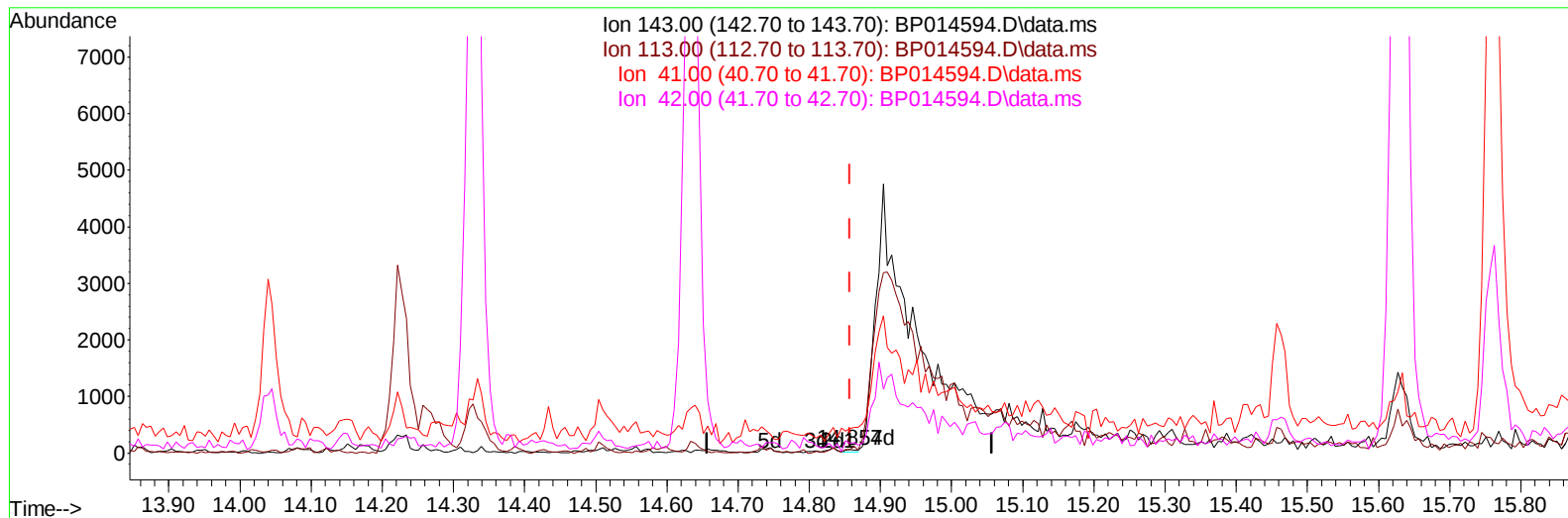
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ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014594.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.000) 0.01 ng/u1

response 60

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	103.28
41.00	44.20	701.64#
42.00	31.50	285.25#

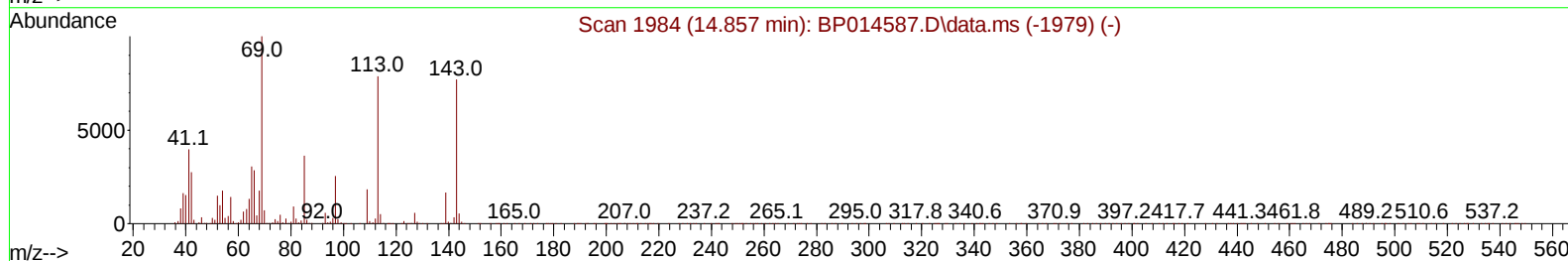
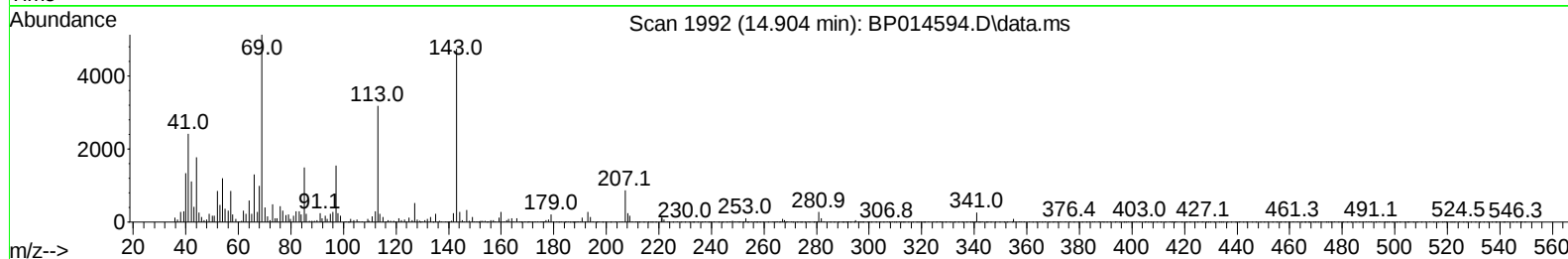
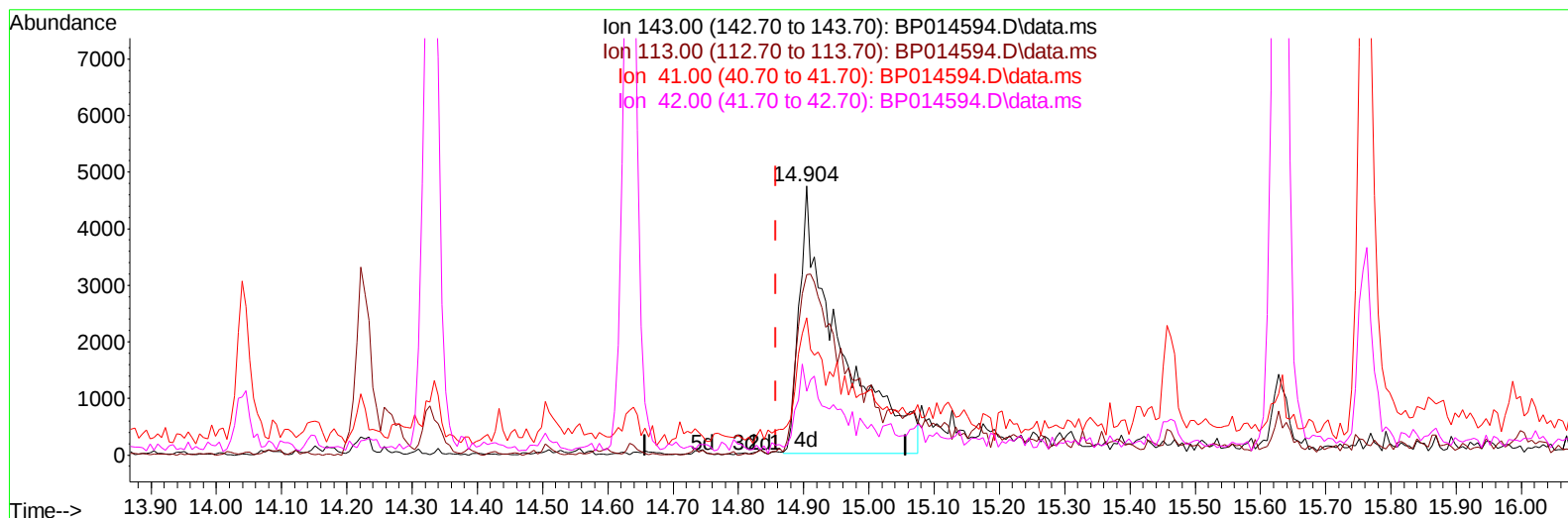
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Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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TIC: BP014594.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.047) 3.90 ng/ul m

response 19827

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	67.17#
41.00	44.20	50.99
42.00	31.50	23.63#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	131628	20.000	ng/ul	0.00
20) Naphthalene-d8	10.804	136	596845	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	414702	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	945527	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	857655	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	725591	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	15180	4.472	ng/uL	0.00
4) Pyridine-d5	3.834	84	25811m	2.967	ng/ul	0.04
7) Phenol-d5	7.163	99	61930	5.102	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.310	67	230390	29.580	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	184820	21.065	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	129495	13.225	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	127869	26.539	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	134826	24.437	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	231761	23.066	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	160357	12.116	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	1067125	31.736	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1105530	29.631	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	19827m	3.903	ng/ul	0.05
60) Fluorene-d10	15.633	176	902242	32.176	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	175860	25.277	ng/ul	0.00
73) Anthracene-d10	17.498	188	1561606	34.126	ng/ul	0.00
81) Pyrene-d10	19.727	212	1899783	32.978	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1368237	35.492	ng/ul	0.00

Target Compounds Qvalue

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

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