

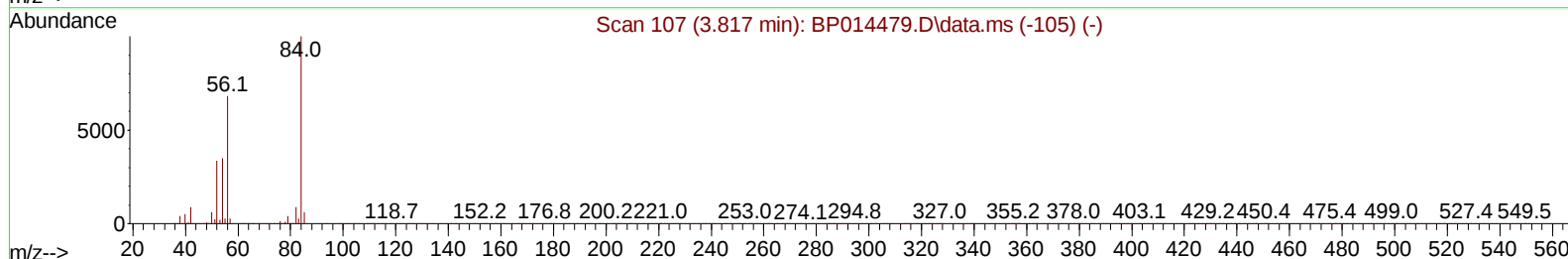
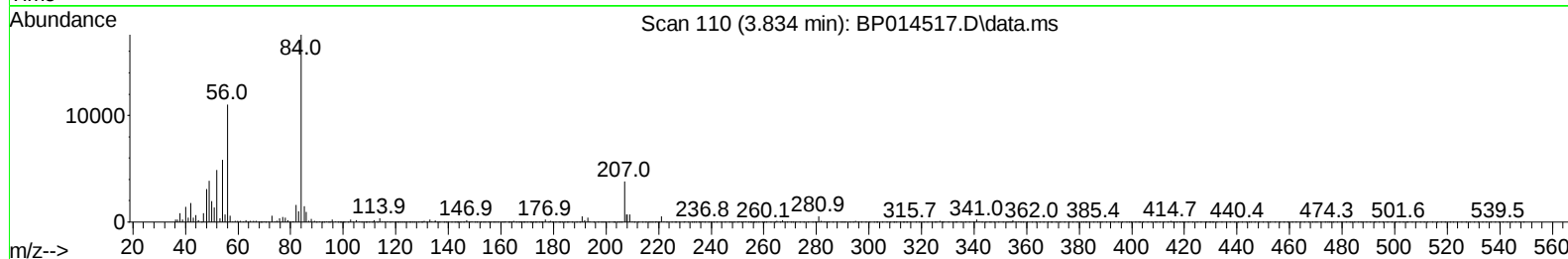
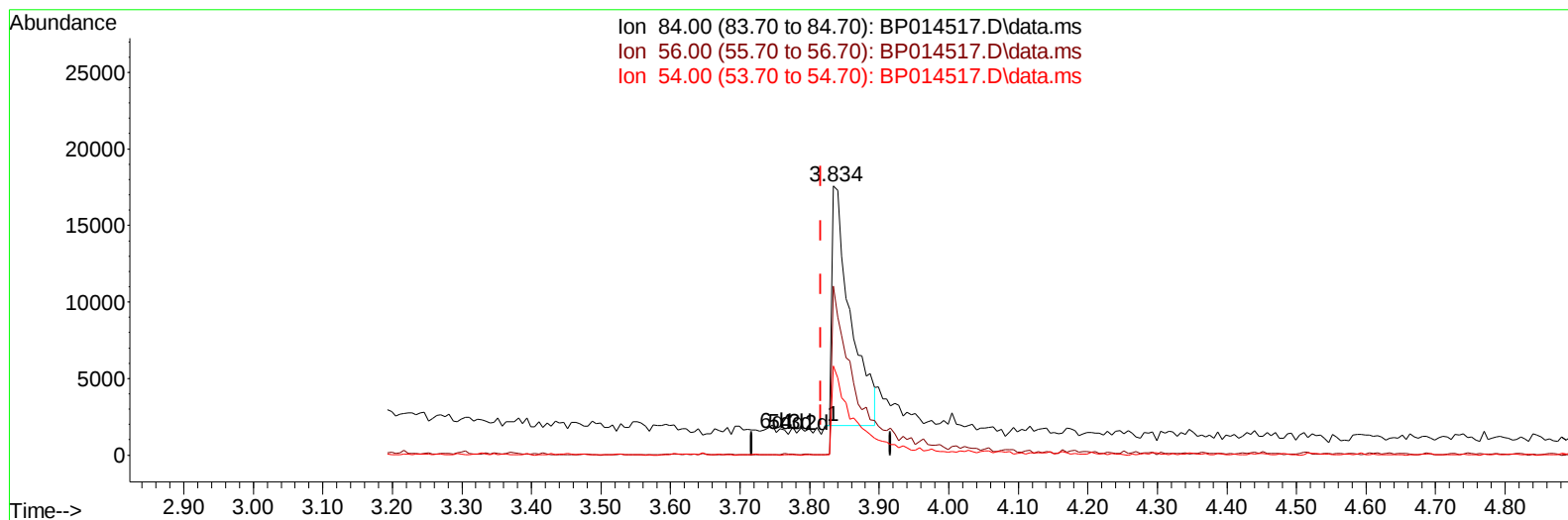
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014517.D
 Acq On : 04 Apr 2023 08:42
 Operator : CG/JU
 Sample : 02158-16
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-2(20230330)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 09:06:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 04 01:34:36 2023
 Response via : Initial Calibration

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



TIC: BP014517.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 3.36 ng/u1

response 28942

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	62.74
54.00	32.60	33.16
0.00	0.00	0.00

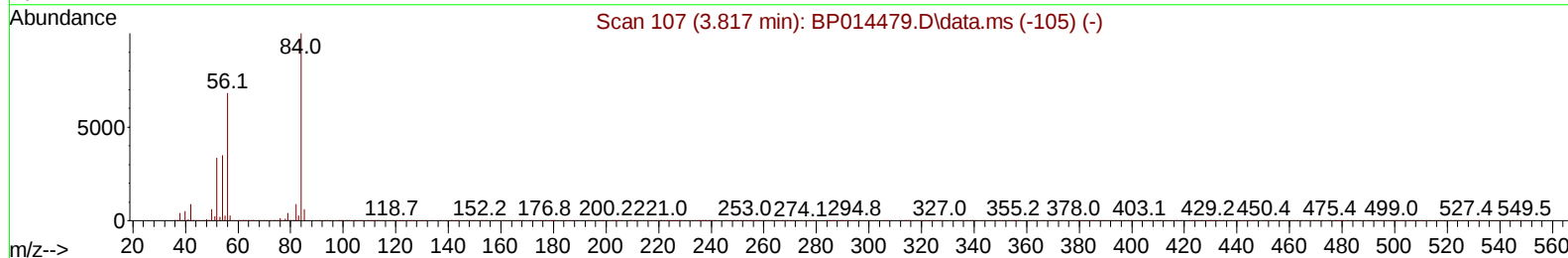
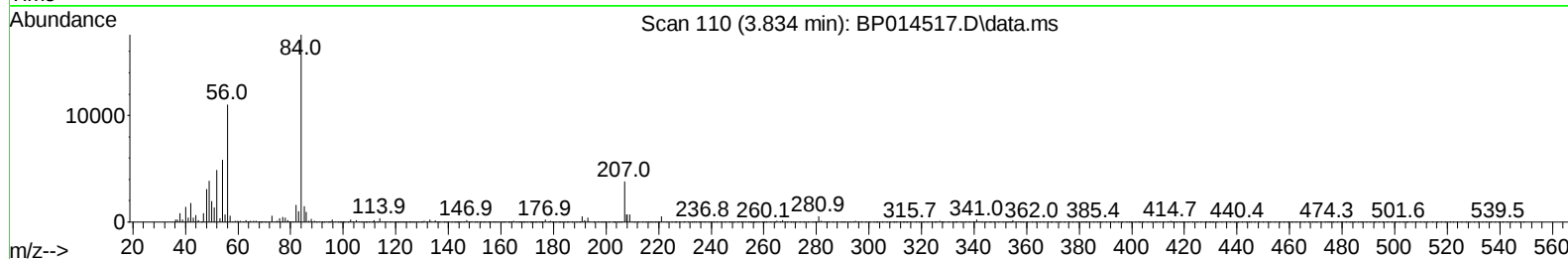
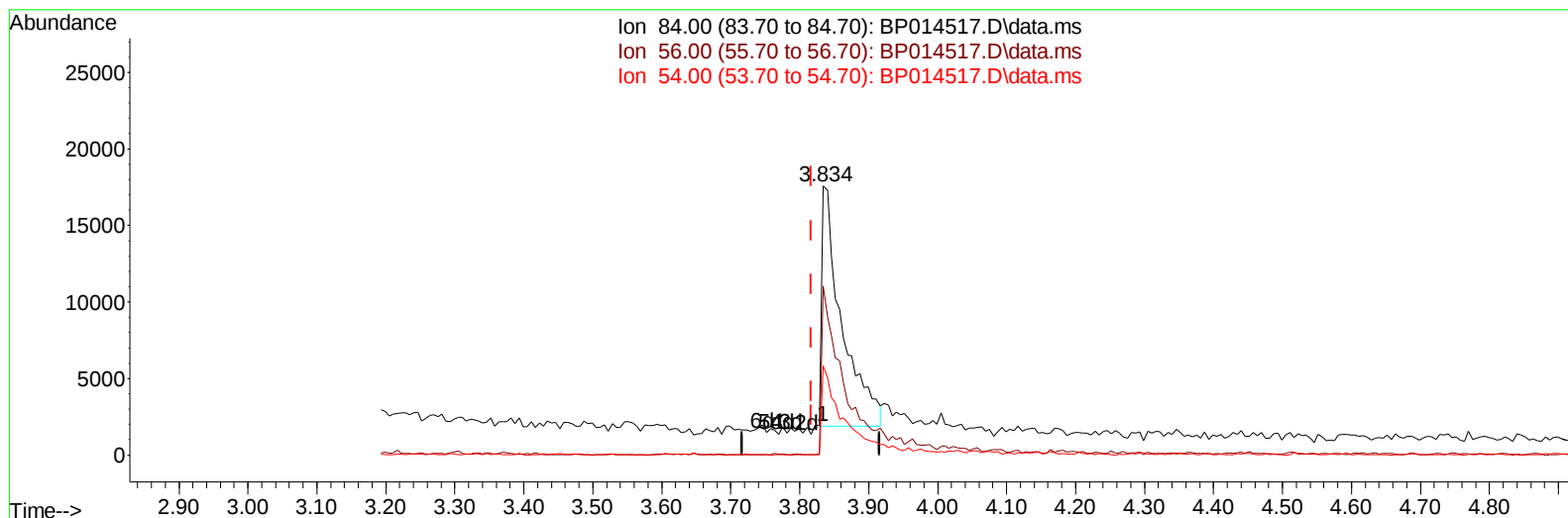
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TIC: BP014517.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 3.68 ng/ul m

response 31675

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	62.74
54.00	32.60	33.16
0.00	0.00	0.00

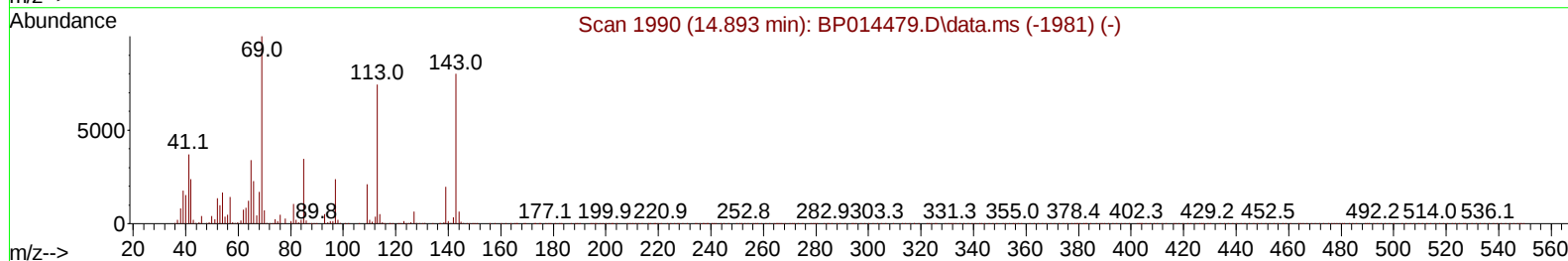
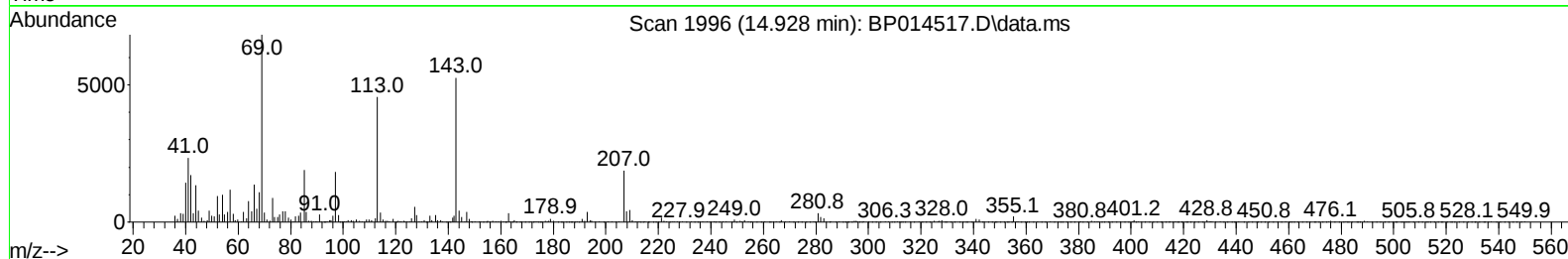
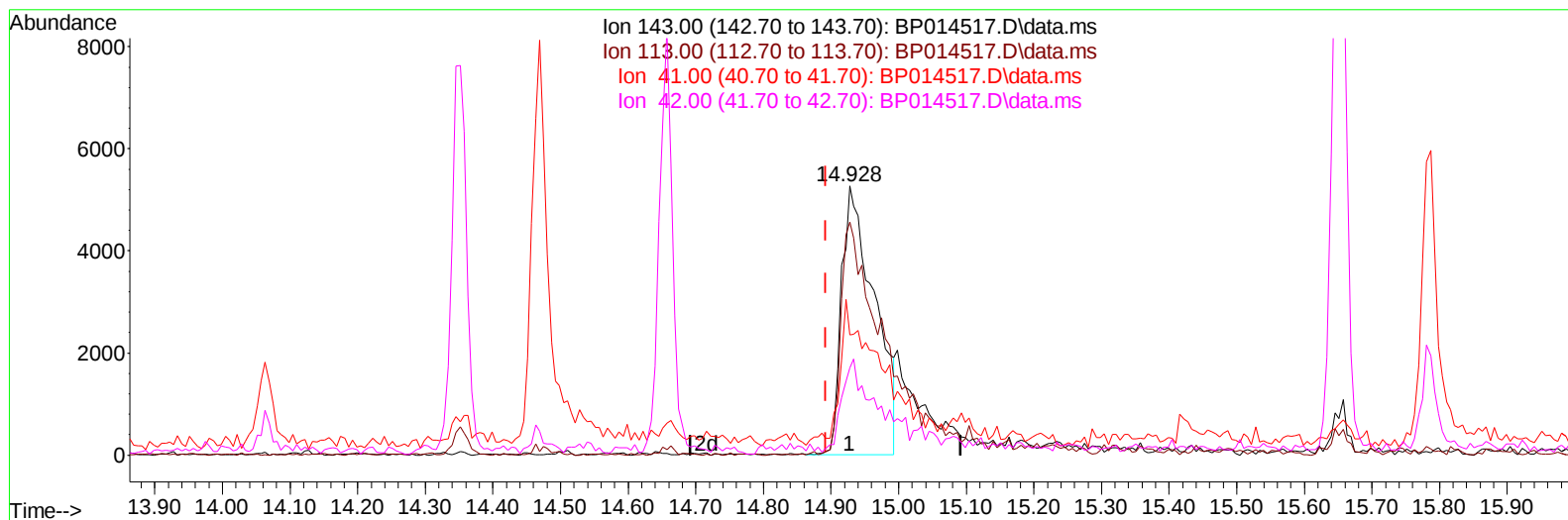
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Instrument :
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TIC: BP014517.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.928min (+ 0.035) 5.15 ng/u1****response 17783**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.71
41.00	44.20	44.61
42.00	31.50	32.33

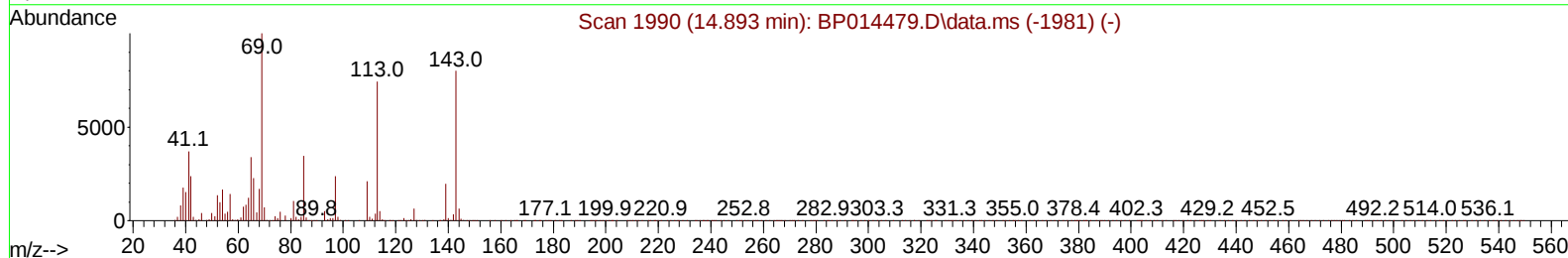
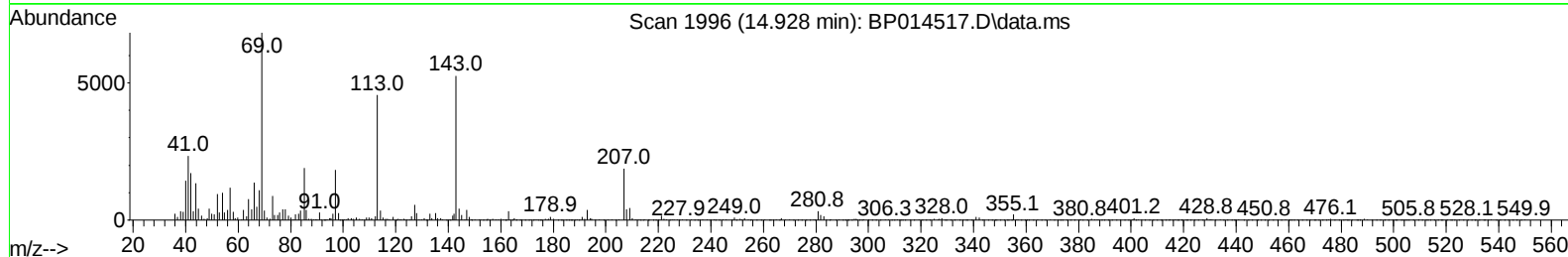
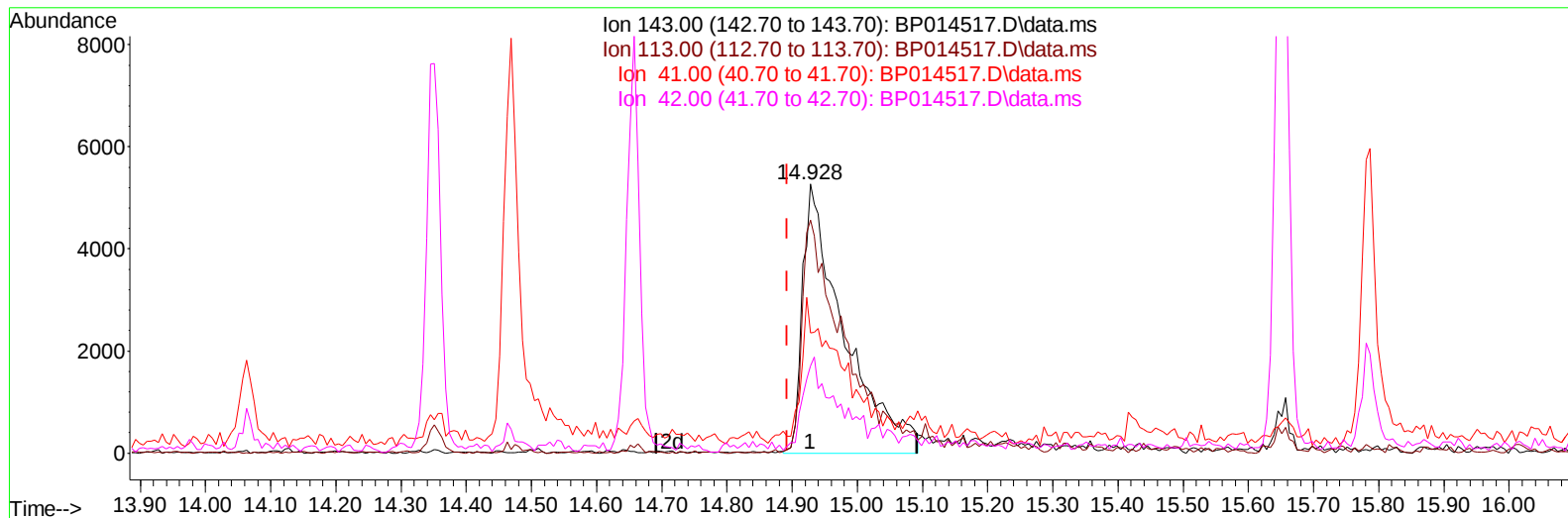
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Sample : 02158-16
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP014517.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.035) 6.71 ng/ul m

response 23183

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.71
41.00	44.20	44.61
42.00	31.50	32.33

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	130321	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	498288	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	281979	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	580297	20.000	ng/ul	-0.01
79) Chrysene-d12	21.504	240	621911	20.000	ng/ul	-0.01
88) Perylene-d12	24.015	264	734985	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	7687	2.287	ng/uL	0.00
4) Pyridine-d5	3.834	84	31675m	3.678	ng/ul	0.02
7) Phenol-d5	7.181	99	72119	6.001	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	206985	26.842	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	185989	21.410	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	125250	12.919	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	111435	27.702	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	121564	26.392	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	192674	22.968	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	145421	13.161	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	668083	29.221	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	731949	28.852	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	23183m	6.712	ng/ul	0.04
60) Fluorene-d10	15.651	176	581905	30.520	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	98208	23.000	ng/ul	0.00
73) Anthracene-d10	17.516	188	900208	32.054	ng/ul	0.00
81) Pyrene-d10	19.745	212	1118837	26.783	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.857	264	1285304	32.915	ng/ul	0.00

Target Compounds Qvalue

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

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