

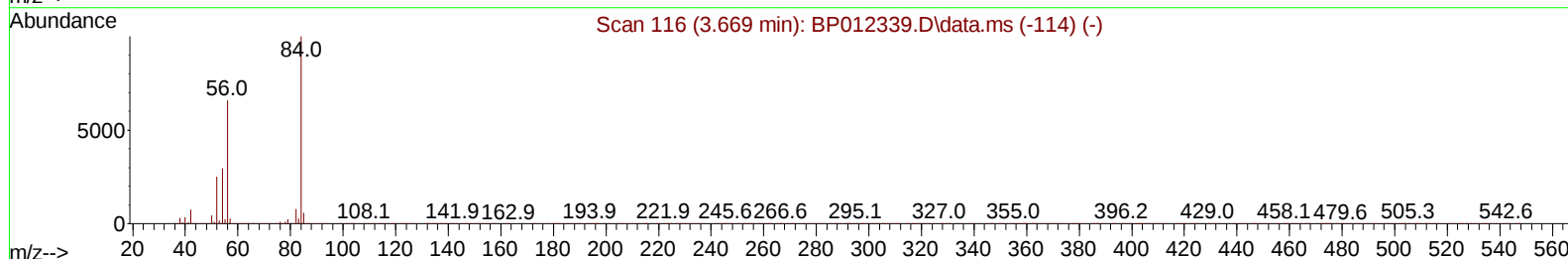
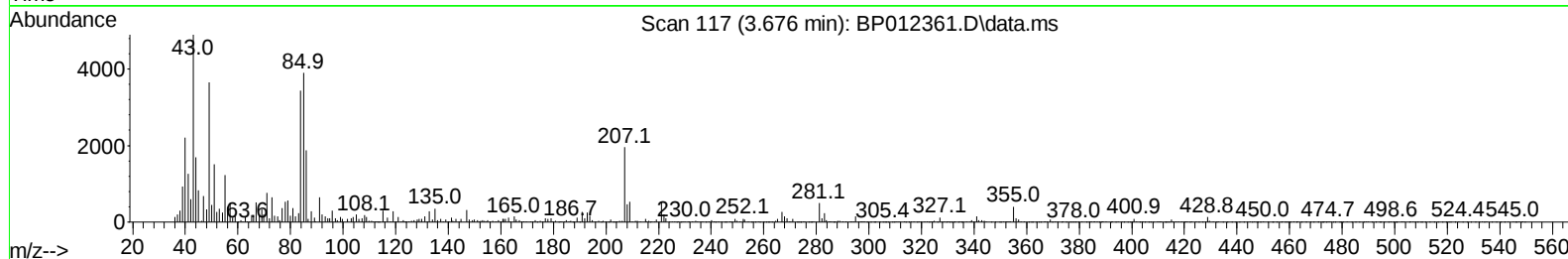
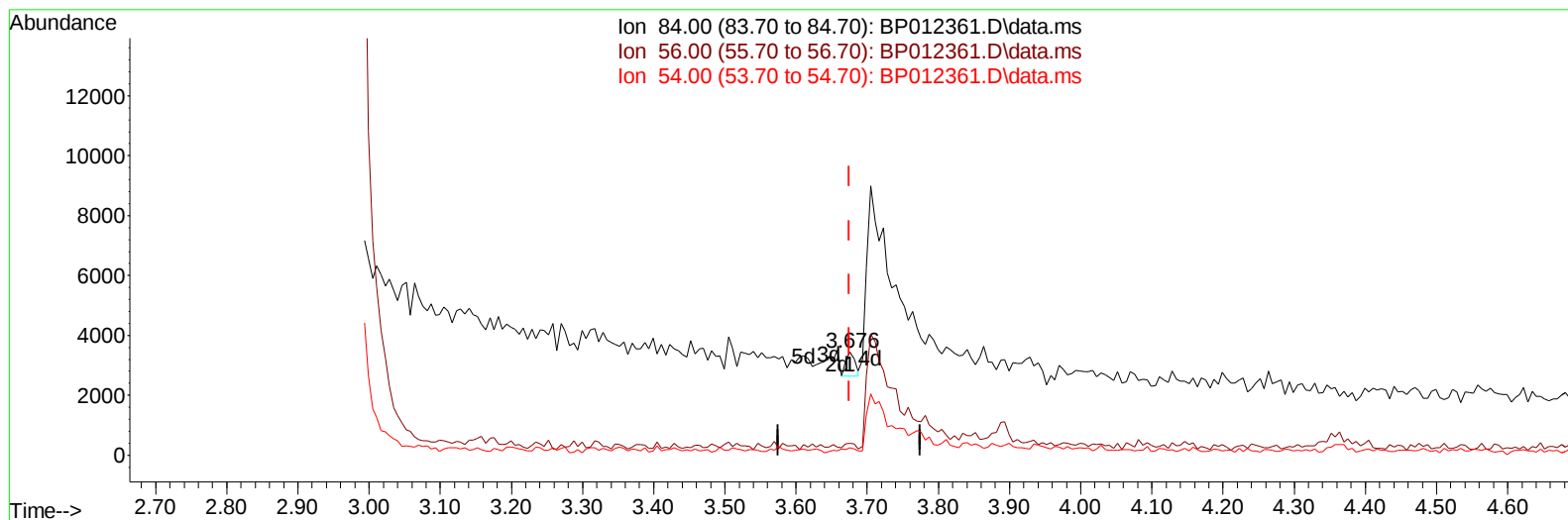
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012361.D
 Acq On : 28 Oct 2022 01:47
 Operator : CG/JU
 Sample : N5144-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BR0

Manual IntegrationsAPPROVED

Quant Time: Oct 28 02:19:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
 Supervised By :mohammad ahmed 10/31/2022



TIC: BP012361.D\data.ms

(4) Pyridine-d5 (S)

3.676min (+ 0.001) 0.03 ng/u1

response 725

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	11.43#
54.00	29.80	7.30#
0.00	0.00	0.00

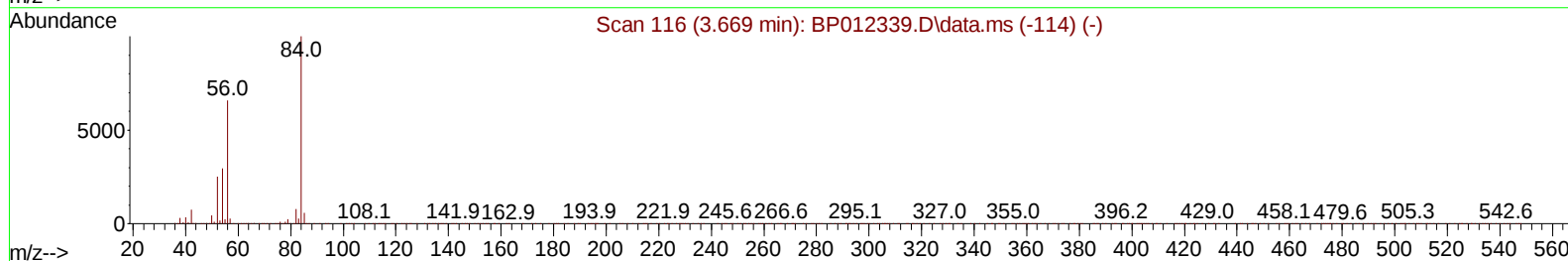
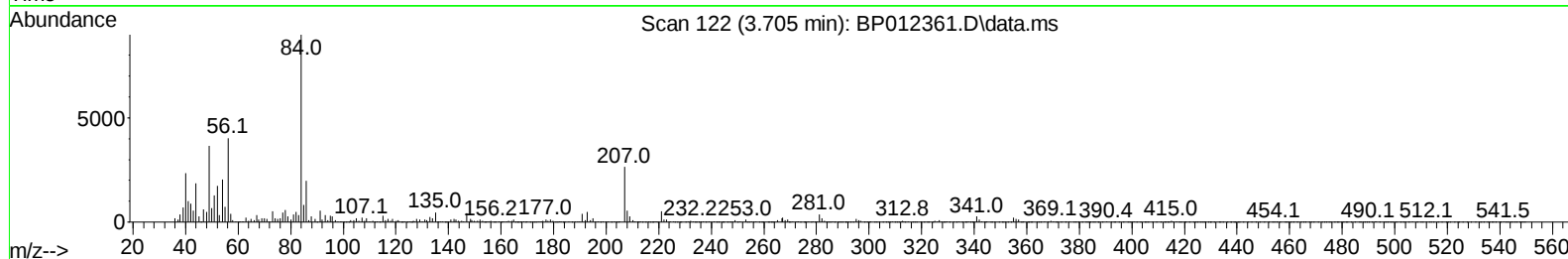
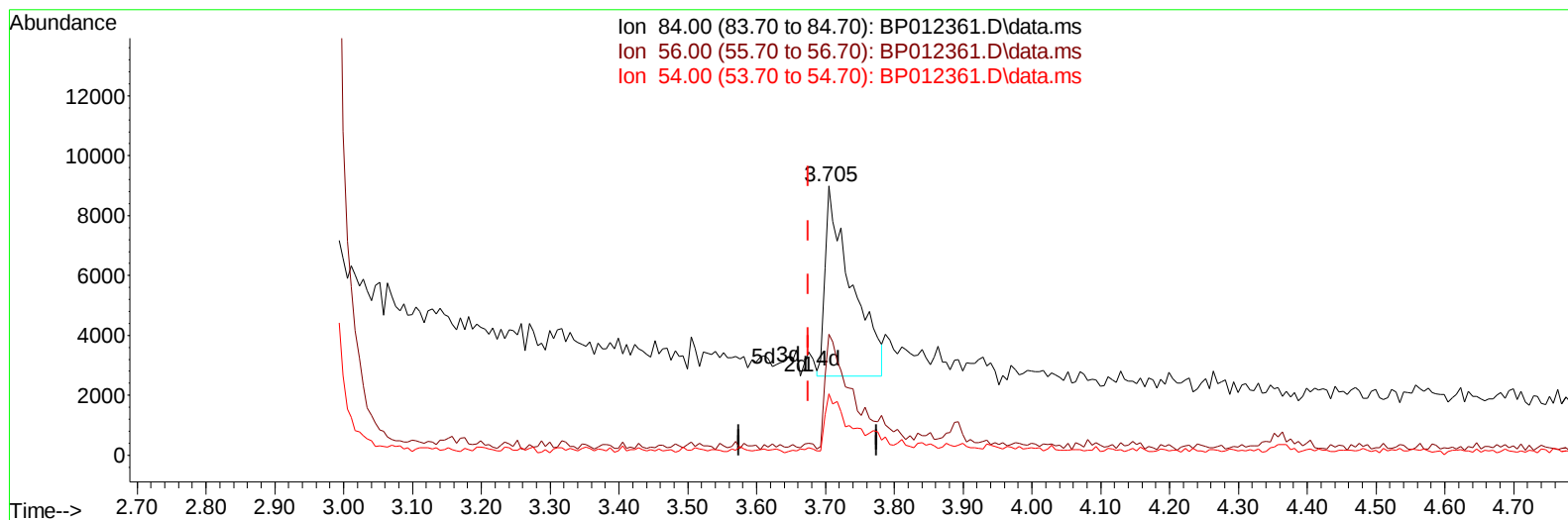
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(4) Pyridine-d5 (S)

3.705min (+ 0.030) 0.78 ng/ul m

response 16746

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	44.89#
54.00	29.80	22.86#
0.00	0.00	0.00

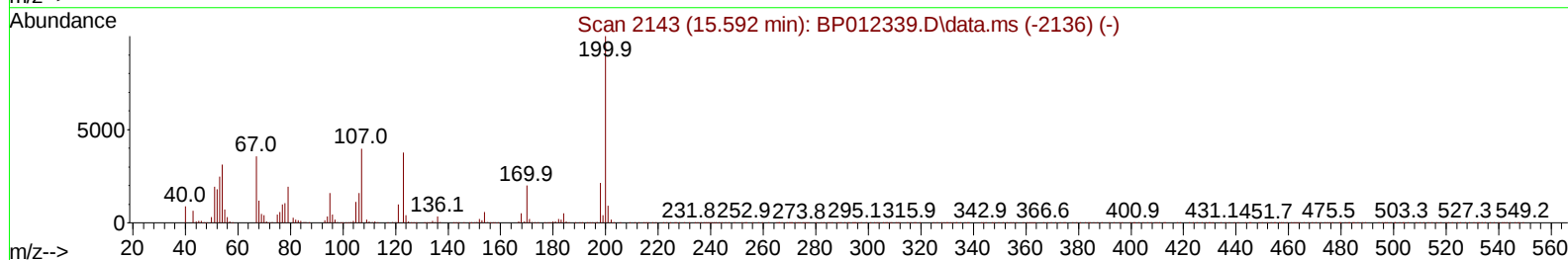
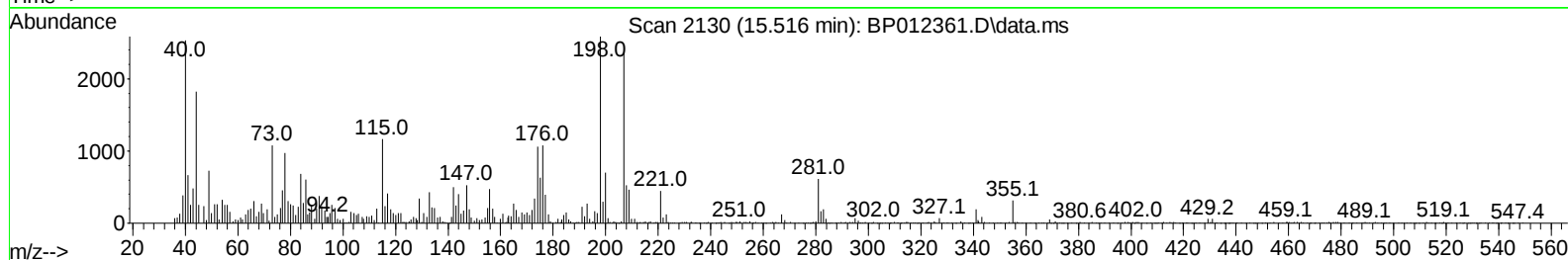
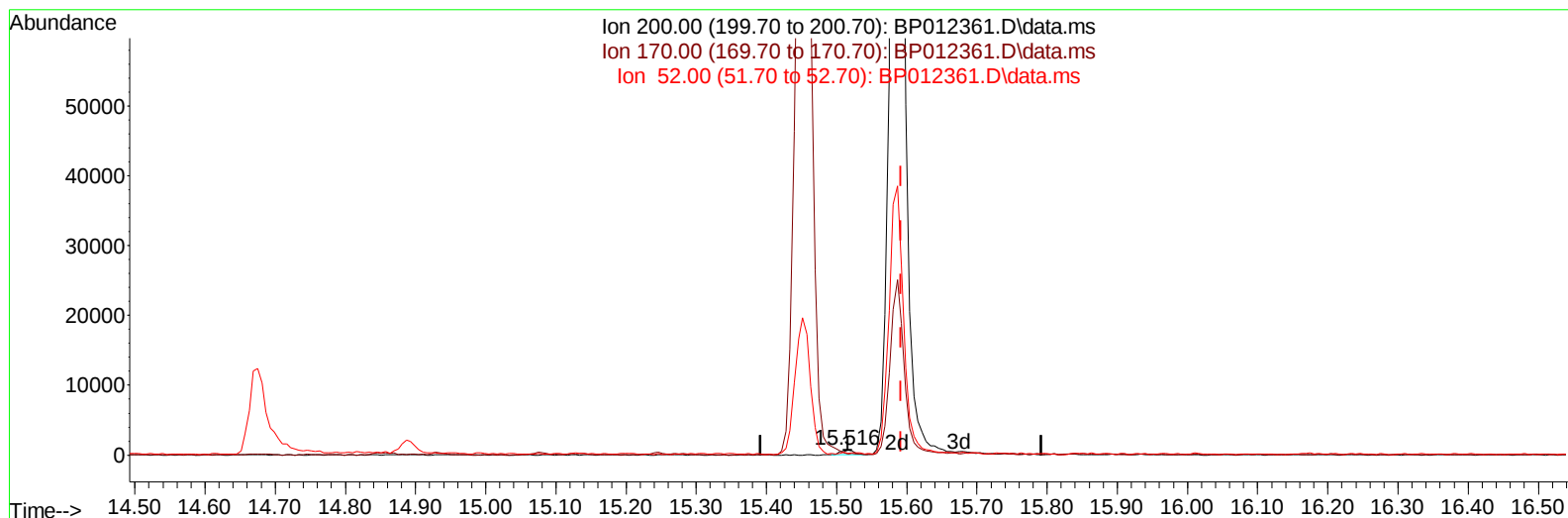
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TIC: BP012361.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.516min (-0.076) 0.09 ng/u1

response 931

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.90	21.54
52.00	37.70	37.23
0.00	0.00	0.00

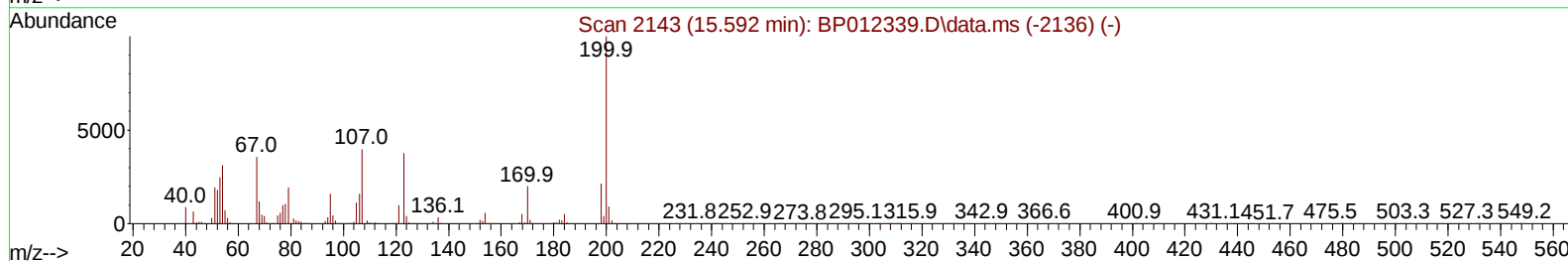
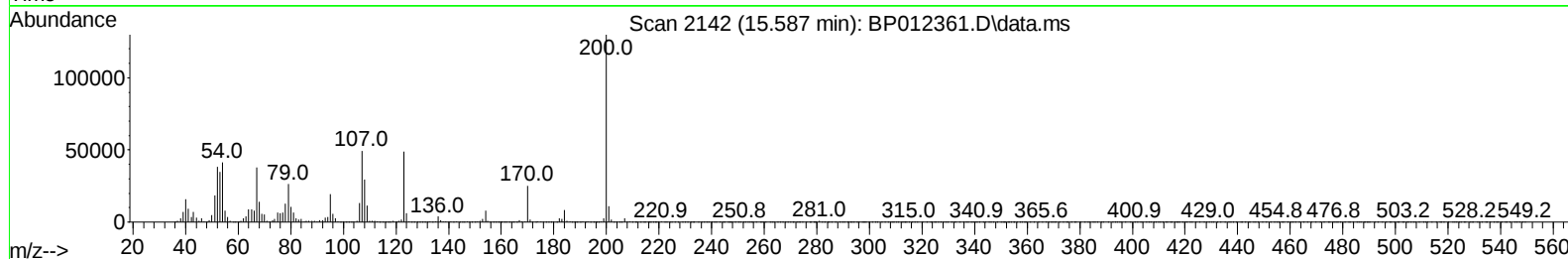
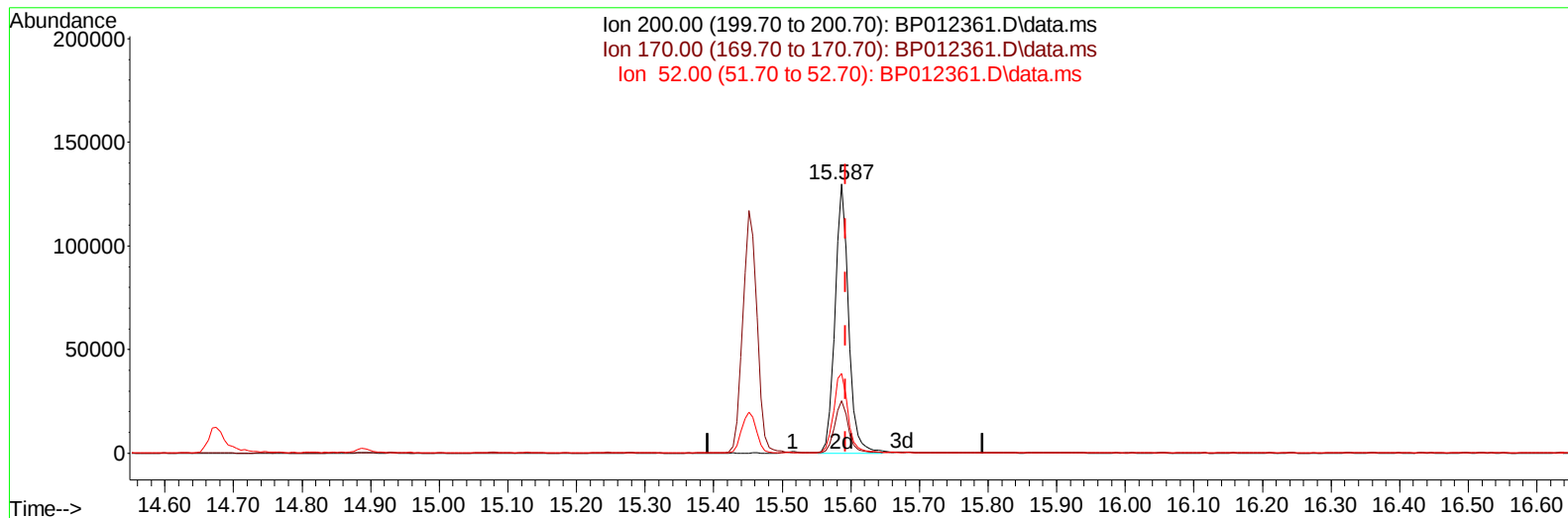
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TIC: BP012361.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.587min (-0.005) 16.89 ng/ul m

response 178056

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.90	19.38
52.00	37.70	29.67#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	301152	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.605	136	1370679	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.457	164	916026	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	2025053	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	2193767	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	2066324	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	27423	3.667	ng/uL	0.00
4) Pyridine-d5	3.705	84	16746m	0.779	ng/ul	0.03
7) Phenol-d5	6.975	99	569863	21.097	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	357063	22.144	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	456060	23.121	ng/ul	-0.01
15) 4-Methylphenol-d8	8.522	113	318544	14.911	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	241911	27.216	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.693	143	254011	27.533	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.234	165	452443	22.629	ng/ul	0.00
31) 4-Chloroaniline-d4	10.758	131	437220	14.687	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1535699	24.490	ng/ul	-0.01
49) Acenaphthylene-d8	14.146	160	1764890	24.629	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.675	143	214737	18.094	ng/ul	0.00
60) Fluorene-d10	15.451	176	1398993	26.058	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.587	200	178056m	16.888	ng/ul	0.00
73) Anthracene-d10	17.316	188	2326657	26.665	ng/ul	0.00
81) Pyrene-d10	19.557	212	3021768	27.435	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	2792057	27.654	ng/ul	0.00

Target Compounds Qvalue

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

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