

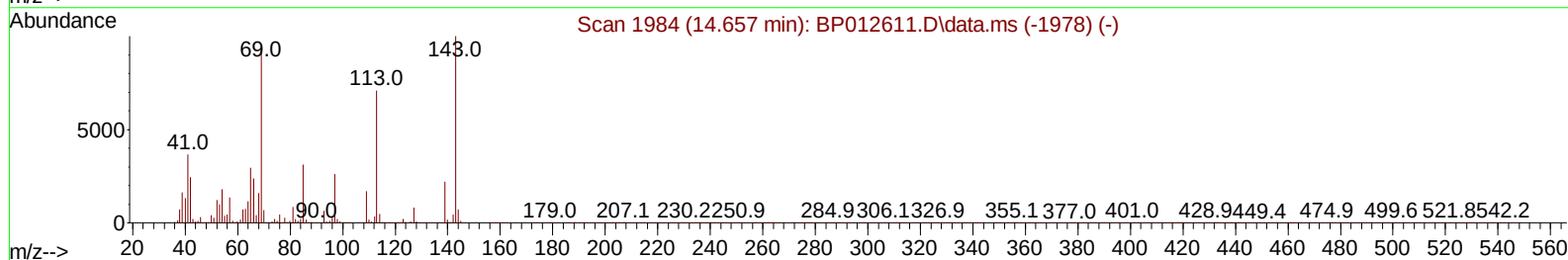
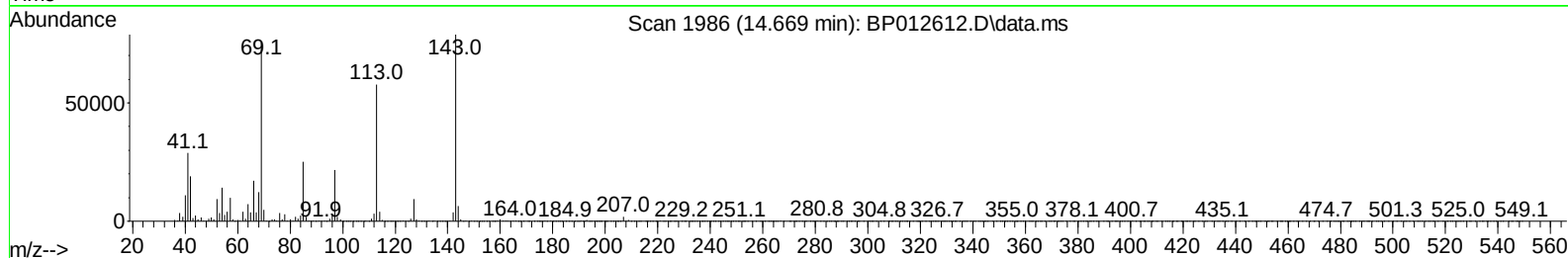
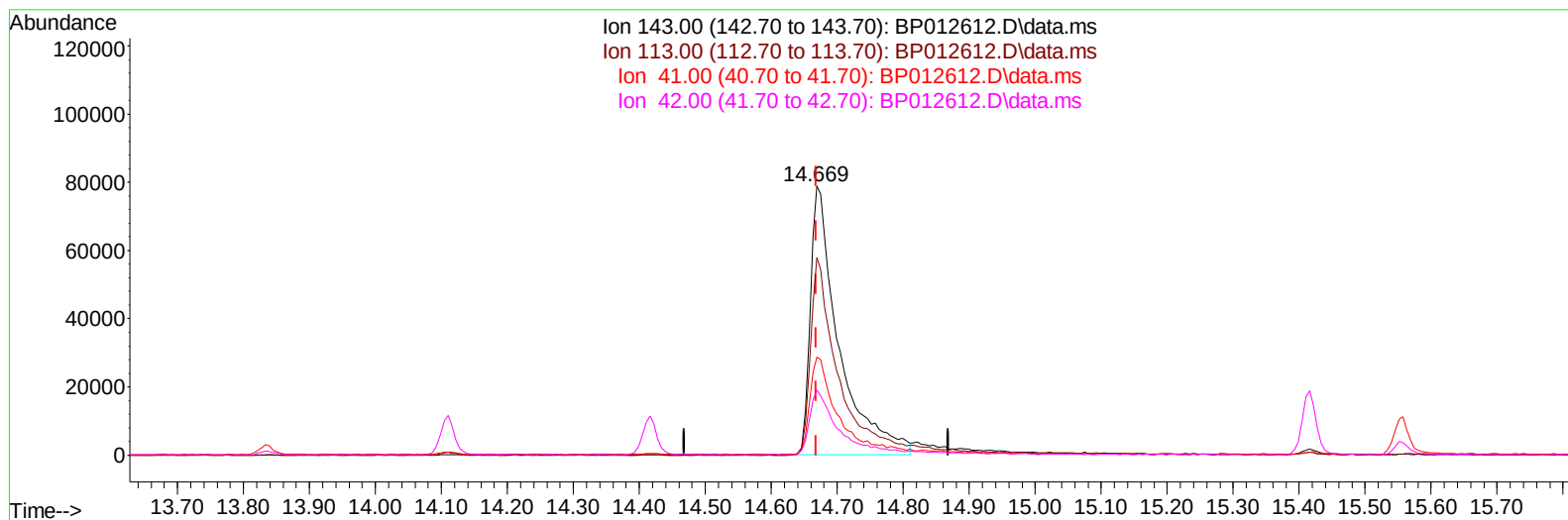
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111122\
 Data File : BP012612.D
 Acq On : 11 Nov 2022 11:45
 Operator : CG/JU
 Sample : PB148940BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK940

Manual IntegrationsAPPROVED

Quant Time: Nov 11 21:40:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 09 23:28:28 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/12/2022
 Supervised By :mohammad ahmed 11/15/2022



TIC: BP012612.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 32.46 ng/ul

response 237307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.39
41.00	35.20	36.51
42.00	24.00	24.24

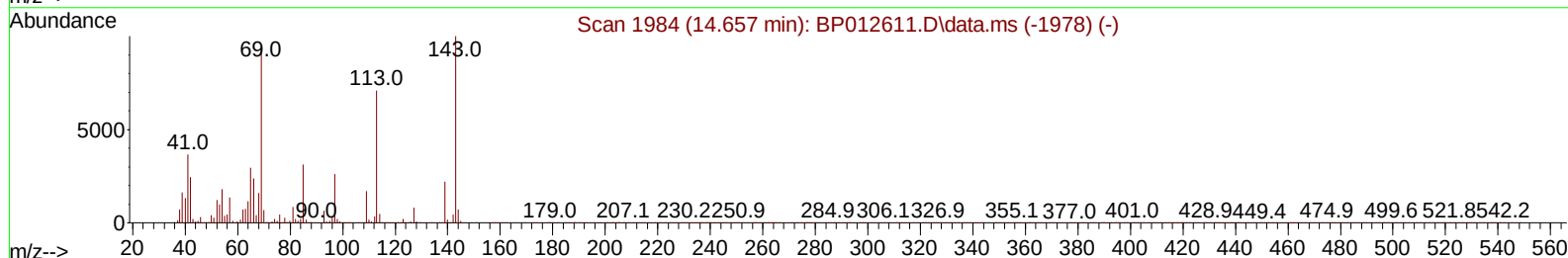
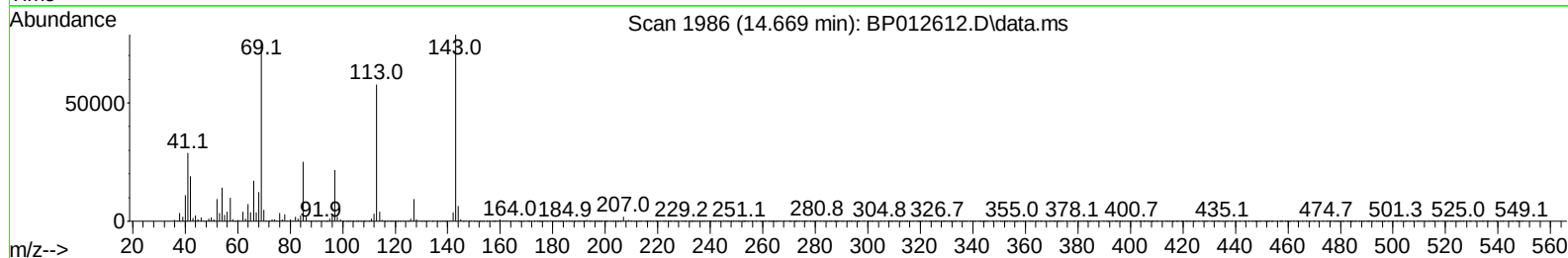
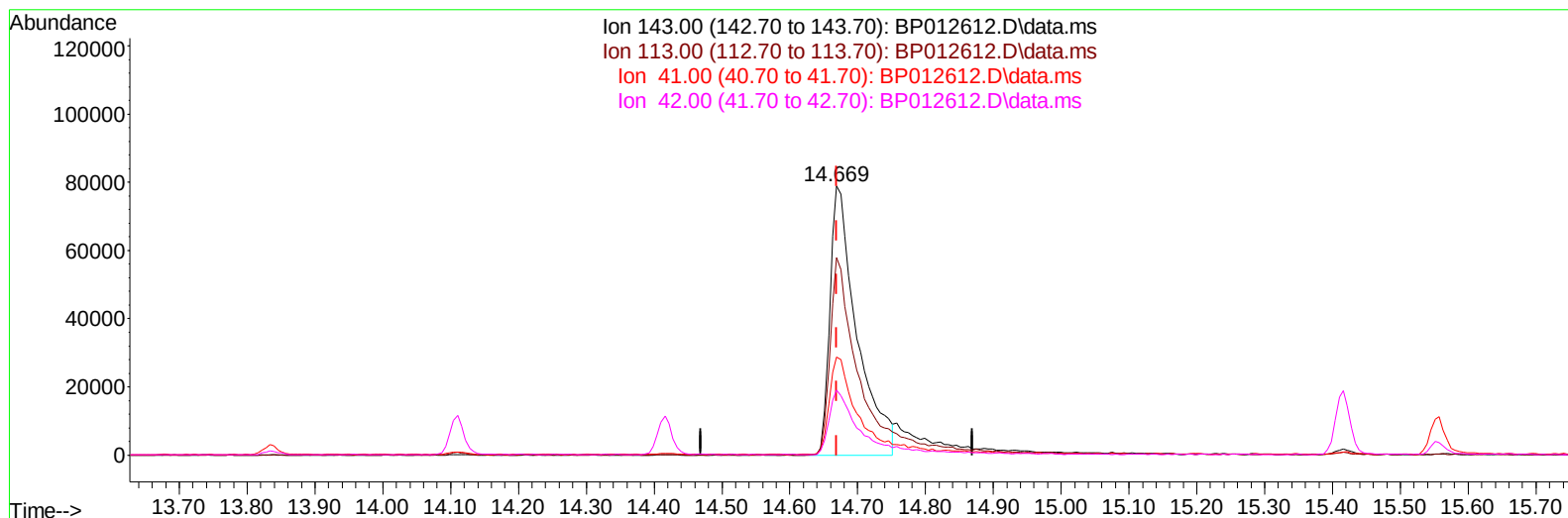
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TIC: BP012612.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.000) 29.69 ng/ul m

response 217034

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.39
41.00	35.20	36.51
42.00	24.00	24.24

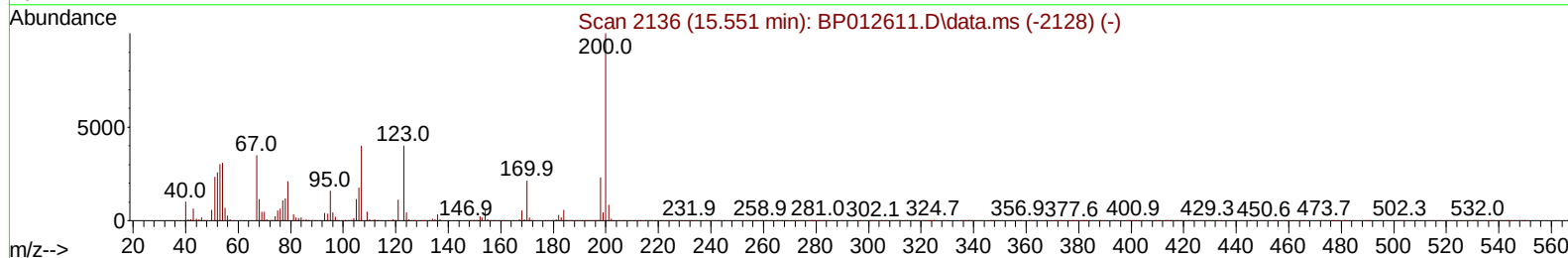
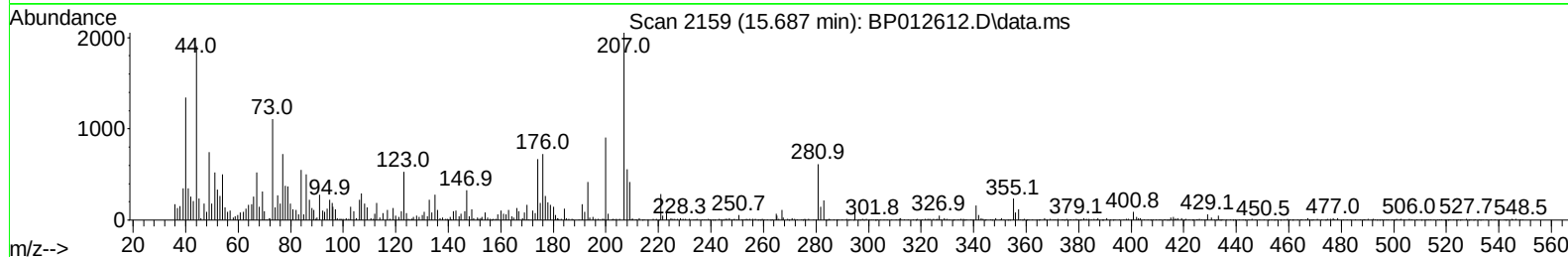
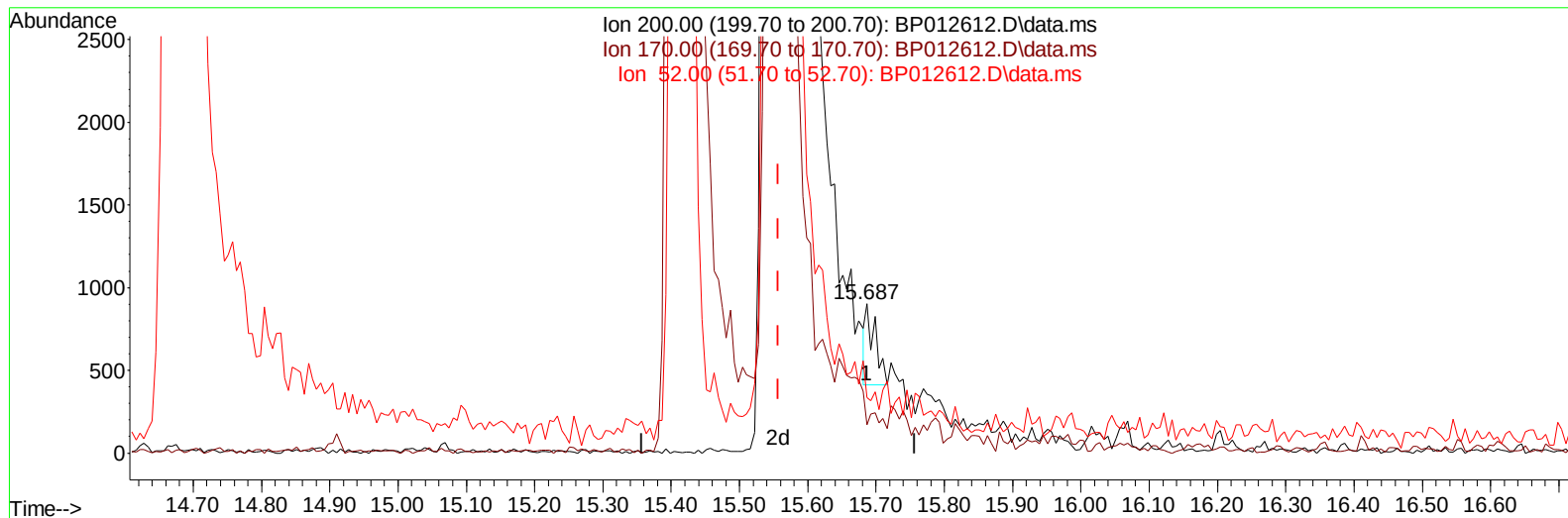
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TIC: BP012612.D\data.ms

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

15.687min (+ 0.129) 0.06 ng/ul

response 486

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	19.27
52.00	41.00	36.99
0.00	0.00	0.00

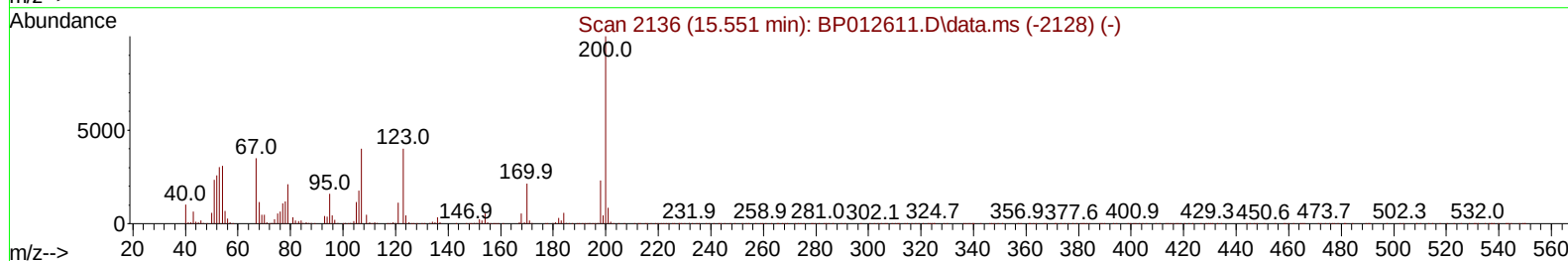
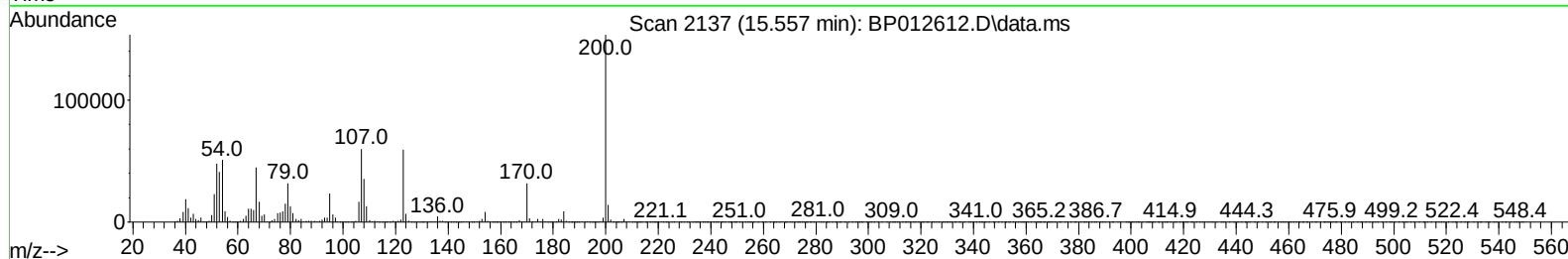
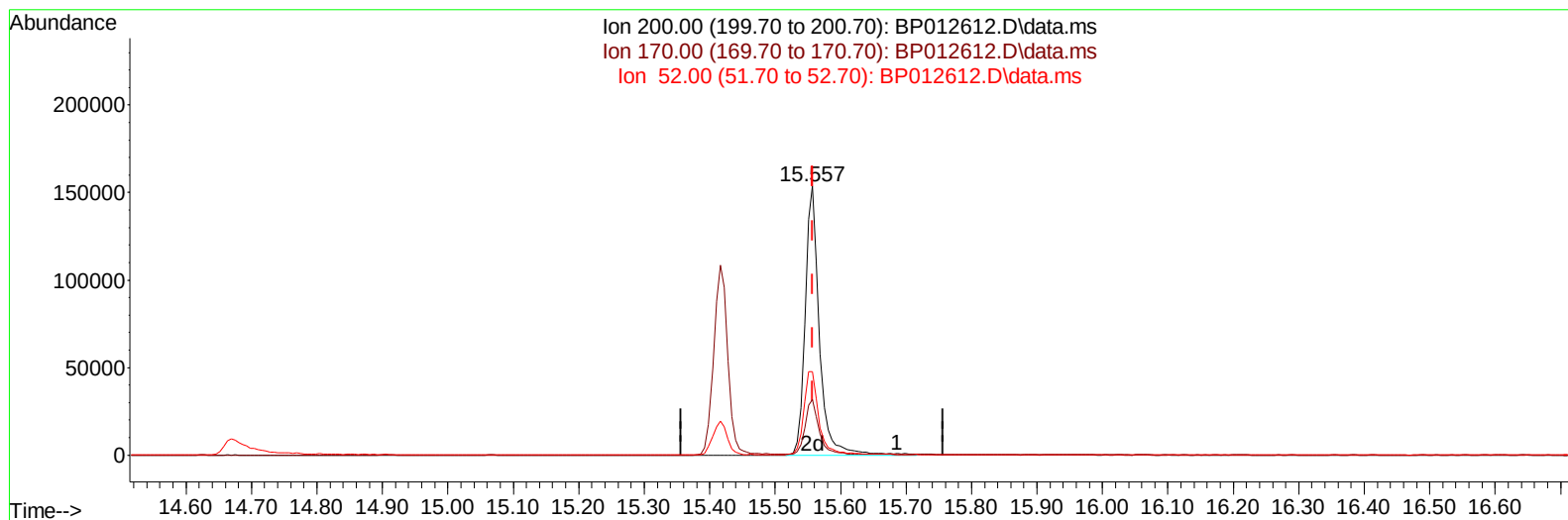
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Instrument :
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TIC: BP012612.D\data.ms

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

15.557min (-0.000) 28.11 ng/ul m

response 231287

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.80	20.73
52.00	41.00	31.13#
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	7.763	152	219489	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1011795	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	673584	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1487840	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1693727	20.000	ng/ul	0.00
88) Perylene-d12	23.657	264	1752425	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	34305	6.370	ng/uL	0.00
4) Pyridine-d5	3.646	84	192585	12.655	ng/ul	0.00
7) Phenol-d5	6.952	99	586338	31.689	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	373212	33.712	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	464701	32.636	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	473448	32.347	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	239333	31.969	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	262043	31.869	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	433750	29.281	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	456658	21.885	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1522697	32.797	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1691732	32.209	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	217034m	29.689	ng/ul	0.00
60) Fluorene-d10	15.416	176	1339653	34.030	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	231287m	28.106	ng/ul	0.00
73) Anthracene-d10	17.281	188	2270561	35.142	ng/ul	0.00
81) Pyrene-d10	19.528	212	2685907	28.298	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	3035101	35.133	ng/ul	0.00

Target Compounds Qvalue

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

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