

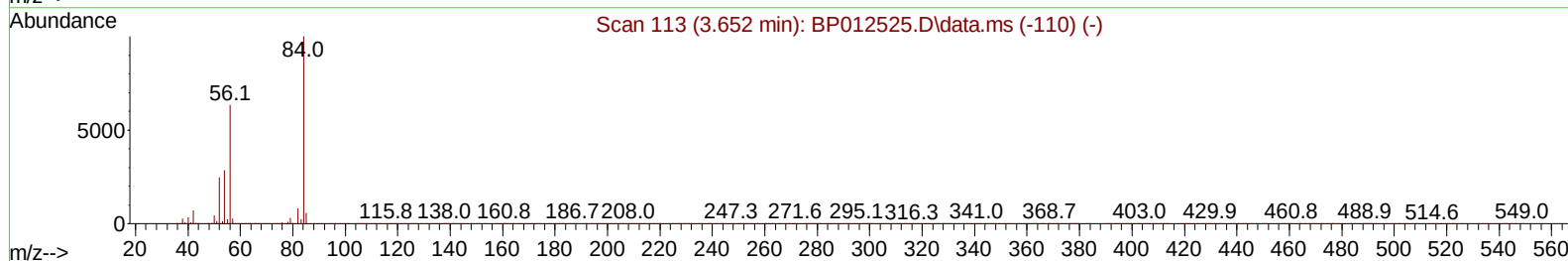
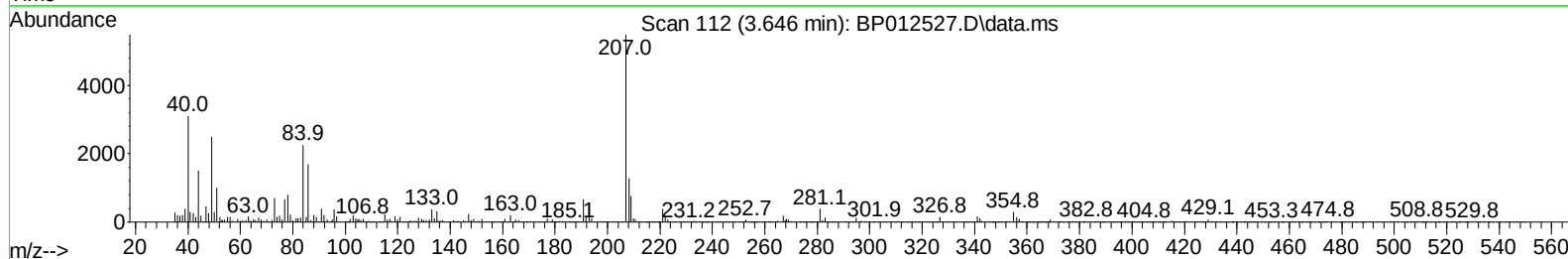
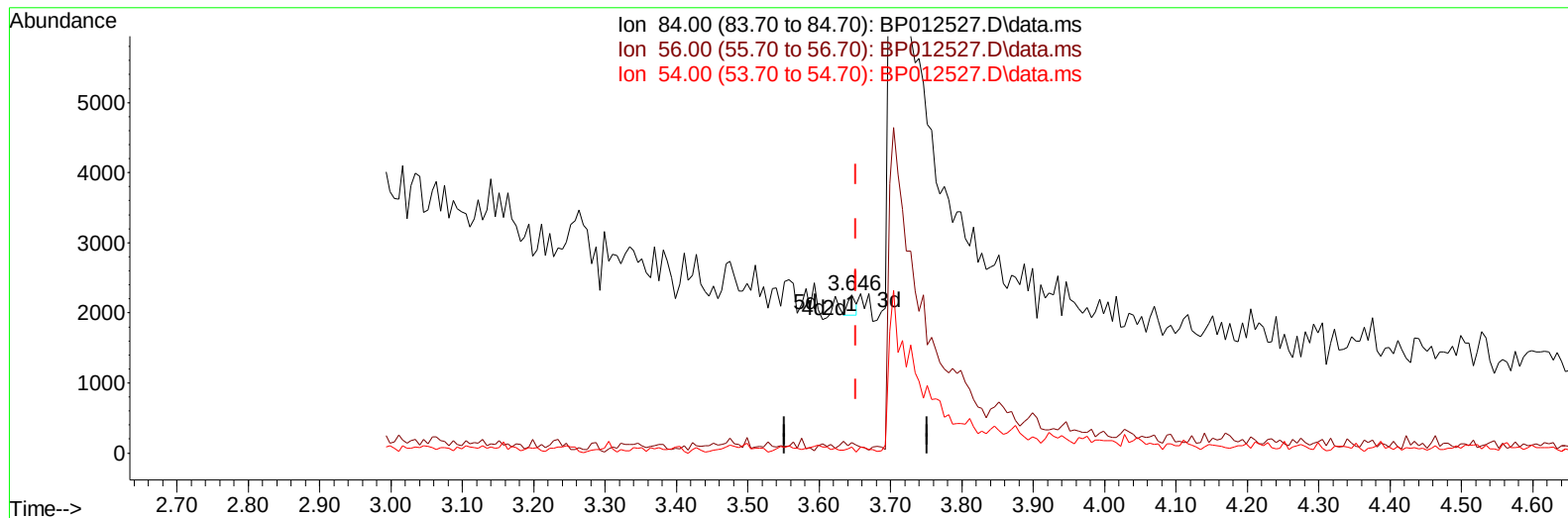
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
 Data File : BP012527.D
 Acq On : 07 Nov 2022 12:39
 Operator : CG/JU
 Sample : PB148750BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK750

Manual IntegrationsAPPROVED

Quant Time: Nov 07 21:39:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:53:36 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/08/2022
 Supervised By :mohammad ahmed 11/14/2022



TIC: BP012527.D\data.ms

(4) Pyridine-d5 (S)

3.646min (-0.006) 0.01 ng/u1

response 229

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	6.47#
54.00	29.80	3.68#
0.00	0.00	0.00

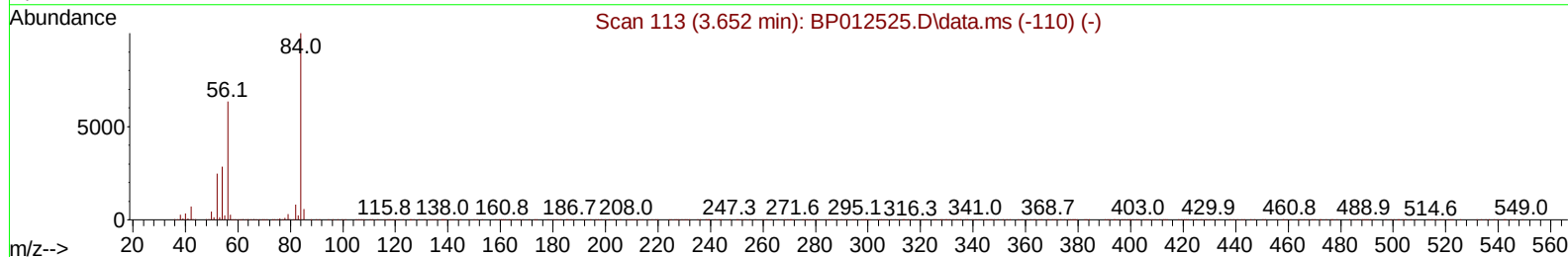
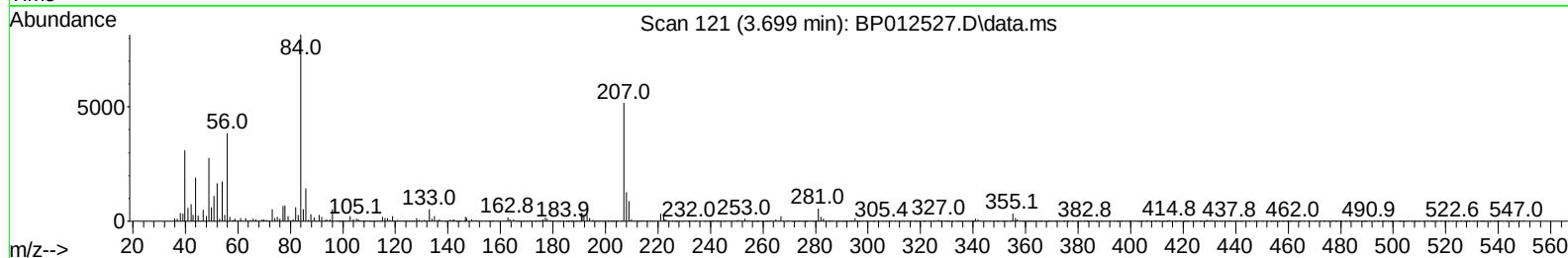
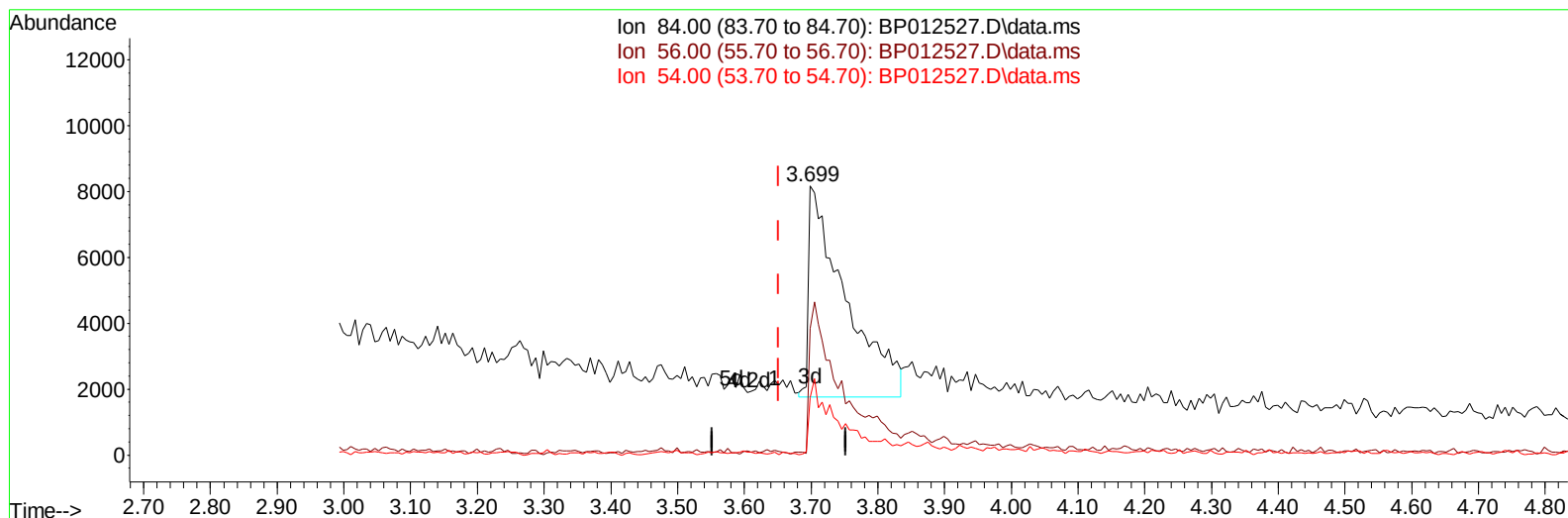
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(4) Pyridine-d5 (S)

3.699min (+ 0.047) 1.23 ng/ul m

response 24453

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	47.00#
54.00	29.80	21.37#
0.00	0.00	0.00

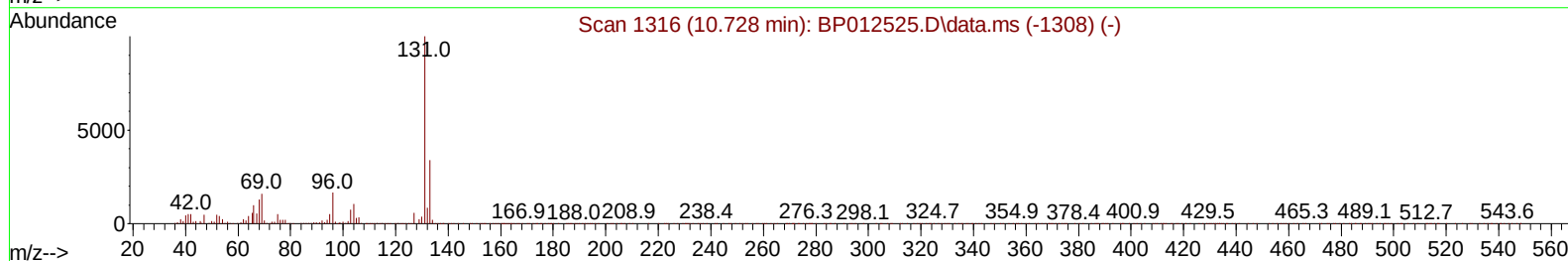
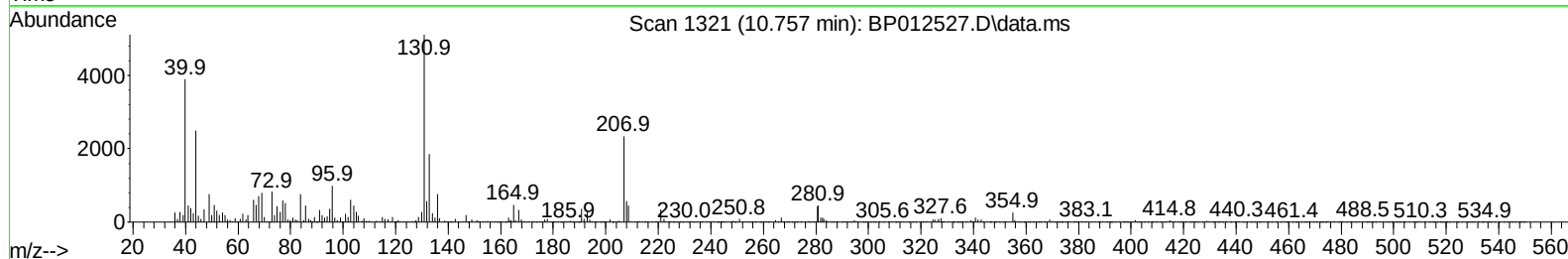
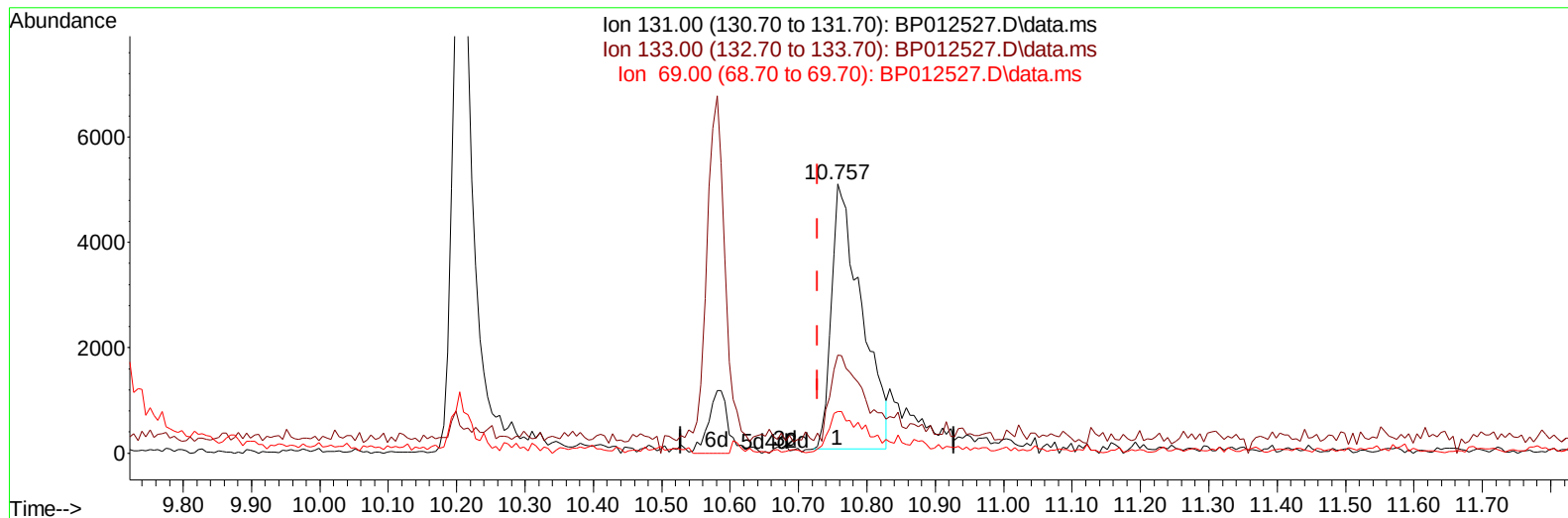
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TIC: BP012527.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.757min (+ 0.029) 0.47 ng/u1

response 15350

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	36.43
69.00	16.50	15.55
0.00	0.00	0.00

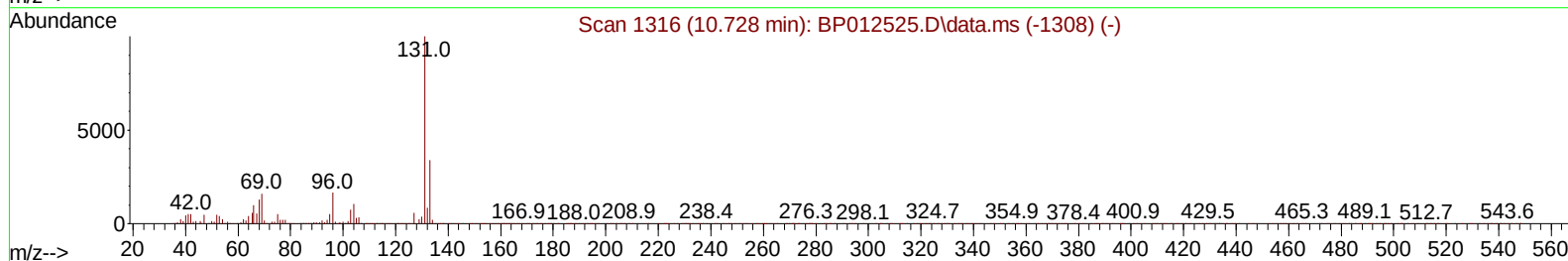
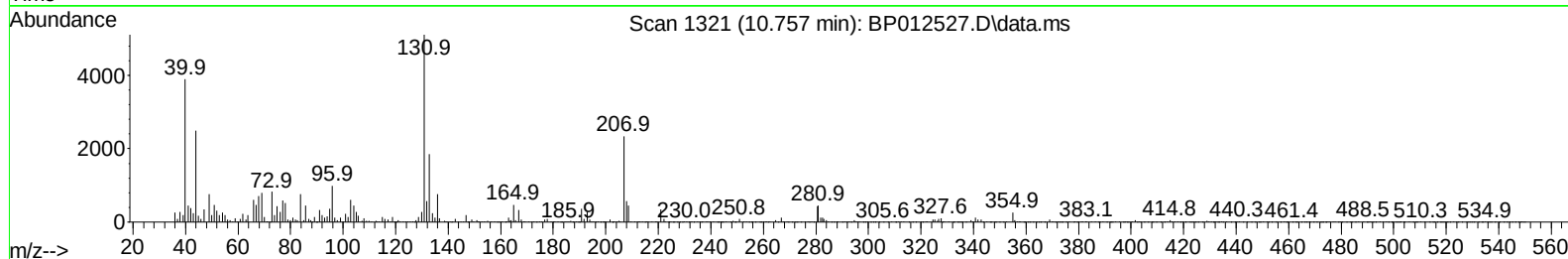
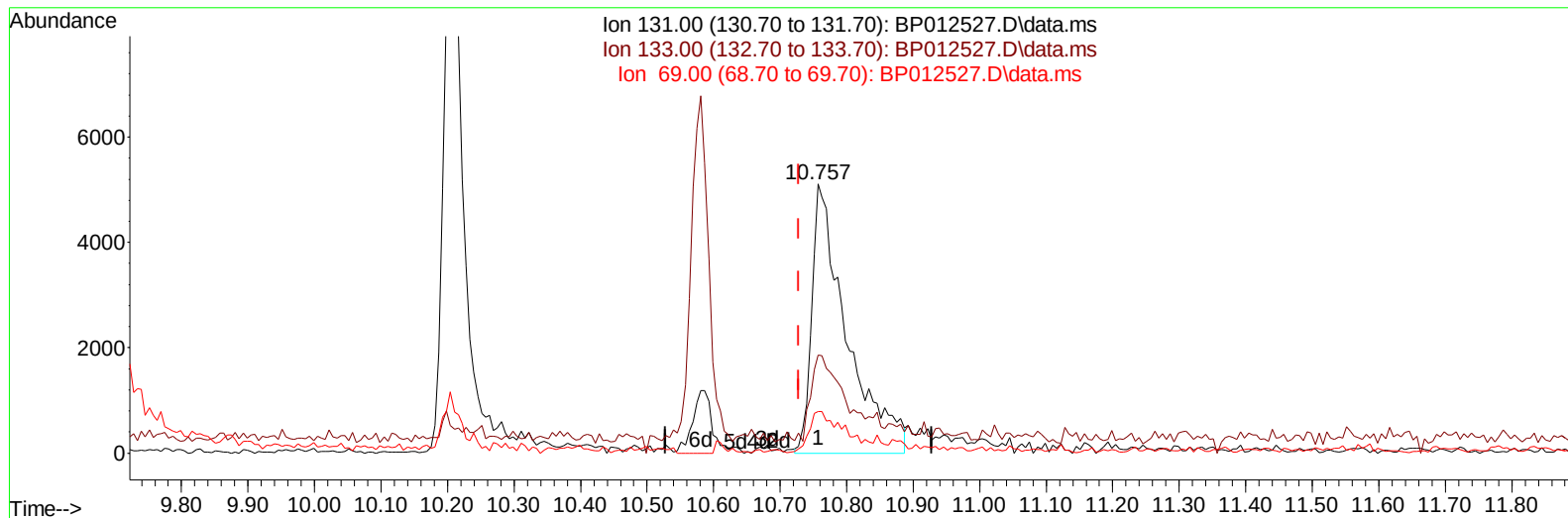
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Misc :
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Instrument :
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TIC: BP012527.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.757min (+ 0.029) 0.57 ng/ul m

response 18589

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	36.43
69.00	16.50	15.55
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	7.775	152	277860	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	1497969	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	1068196	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2479743	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	2207403	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	2015603	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	38955	5.645	ng/uL	0.00
4) Pyridine-d5	3.699	84	24453m	1.234	ng/ul	0.05
7) Phenol-d5	6.957	99	724247	29.061	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	447516	30.080	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	590588	32.451	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	655538	33.258	ng/ul	0.00
21) Nitrobenzene-d5	8.951	128	339543	34.954	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	395776	39.254	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	730986	33.454	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	18589m	0.571	ng/ul	0.03
46) Dimethylphthalate-d6	13.845	166	2534230	34.657	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	1708867	20.450	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	324190	23.425	ng/ul	0.00
60) Fluorene-d10	15.428	176	2143903	34.244	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	211218	16.360	ng/ul	0.00
73) Anthracene-d10	17.292	188	3269925	30.604	ng/ul	0.00
81) Pyrene-d10	19.533	212	3904425	35.230	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	2780400	28.232	ng/ul	0.00

Target Compounds Qvalue

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

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