

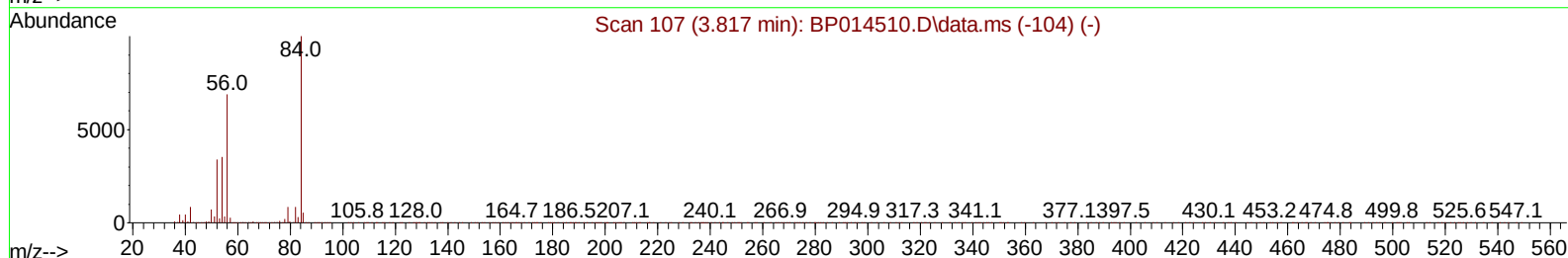
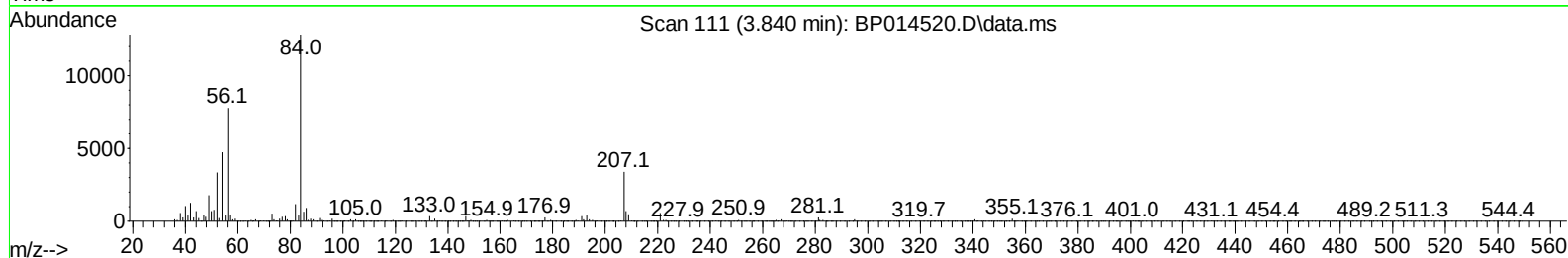
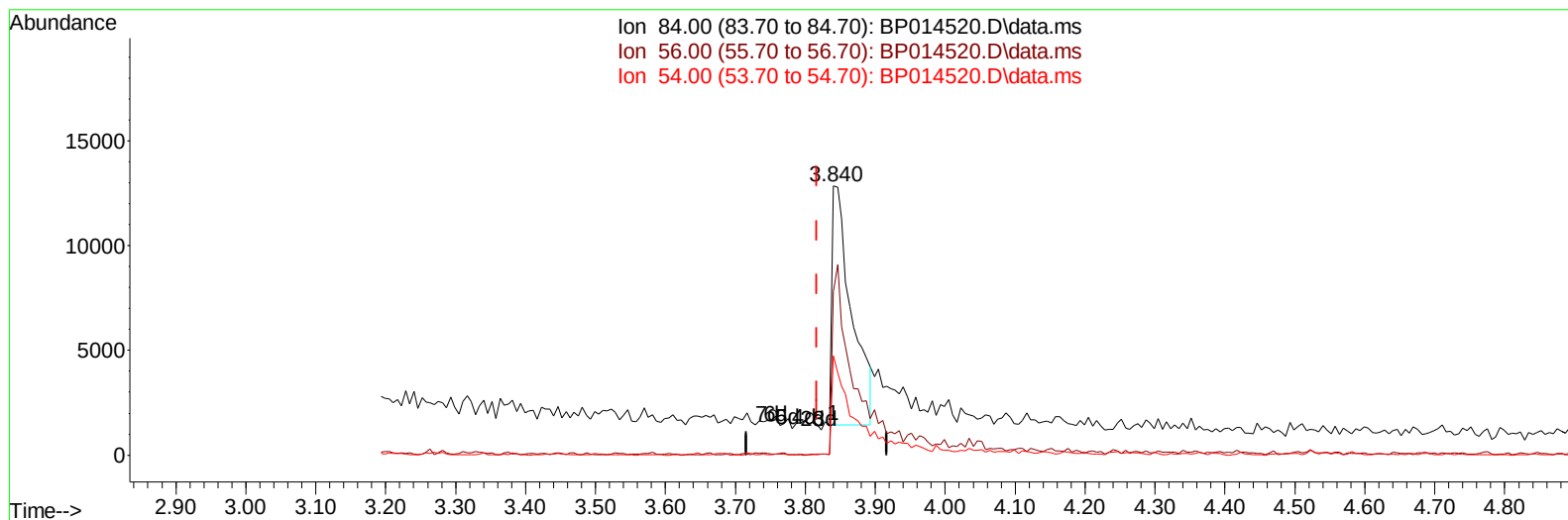
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014520.D
 Acq On : 04 Apr 2023 10:23
 Operator : CG/JU
 Sample : 02158-20
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB(20230330)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:08:48 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 04/05/2023
 Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014520.D\data.ms

(4) Pyridine-d5 (S)

3.840min (+ 0.023) 2.34 ng/ul

response 22305

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	60.75
54.00	32.60	36.82
0.00	0.00	0.00

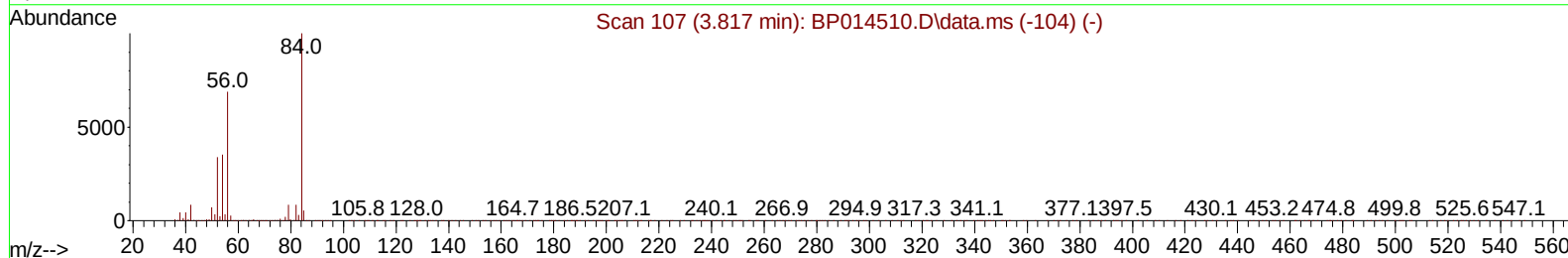
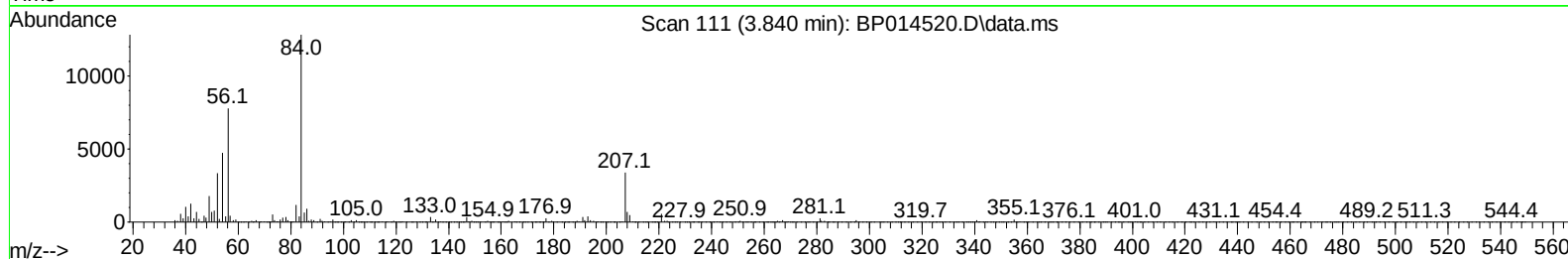
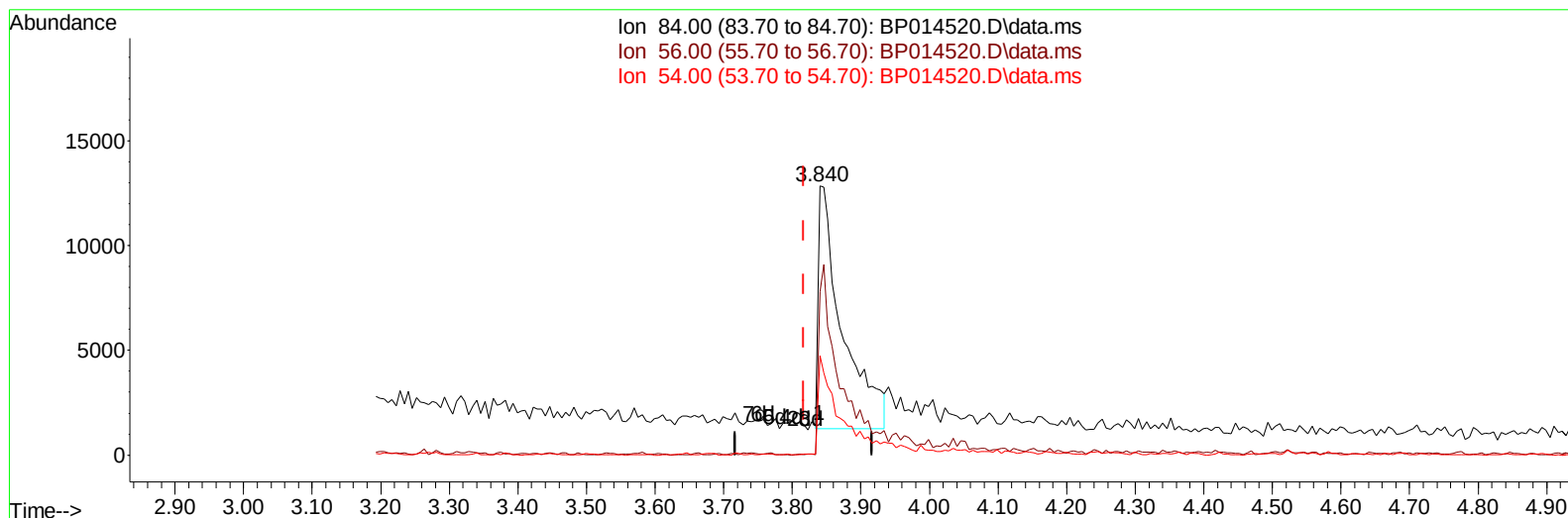
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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(4) Pyridine-d5 (S)

3.840min (+ 0.023) 2.95 ng/ul m

response	28119	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	60.75
54.00	32.60	36.82
0.00	0.00	0.00

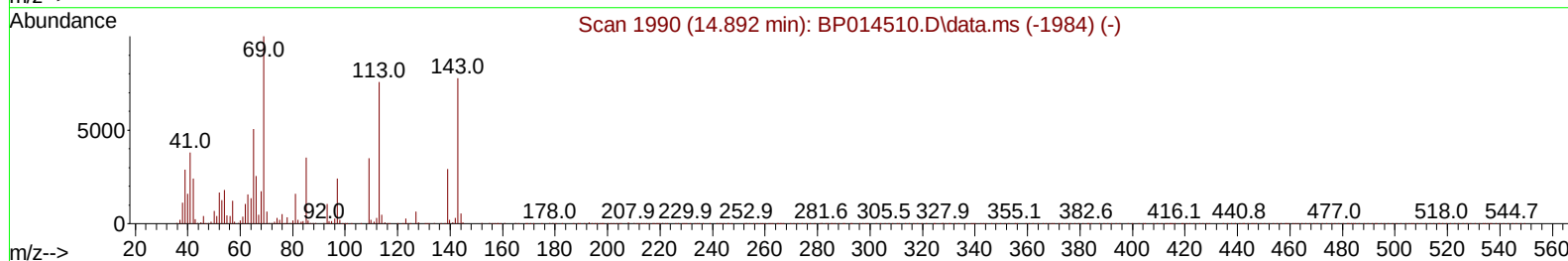
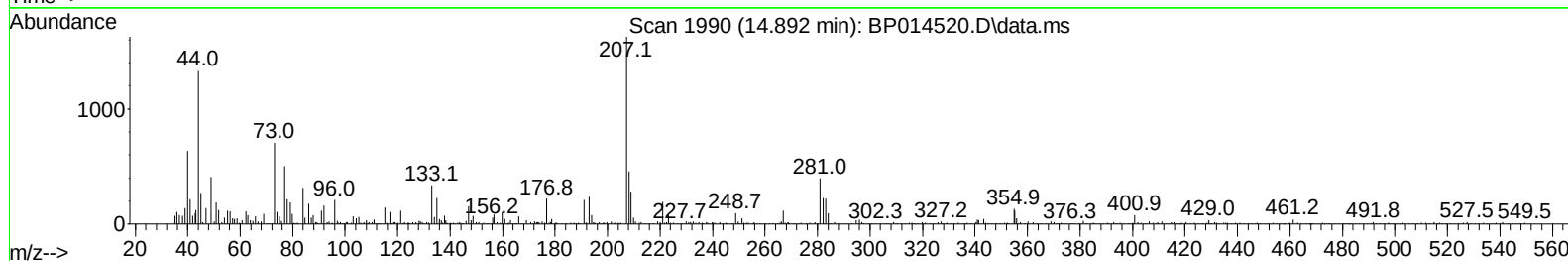
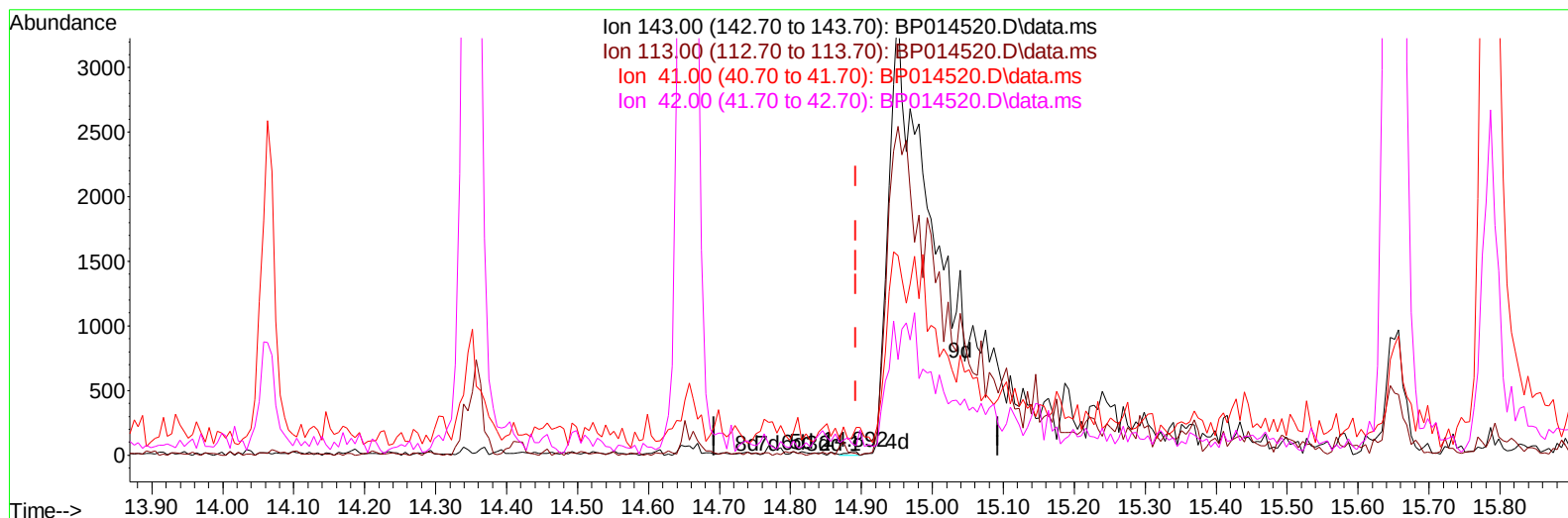
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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TIC: BP014520.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.892min (-0.000) 0.01 ng/u1

response 28

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	44.44#
41.00	44.20	788.89#
42.00	31.50	274.07#

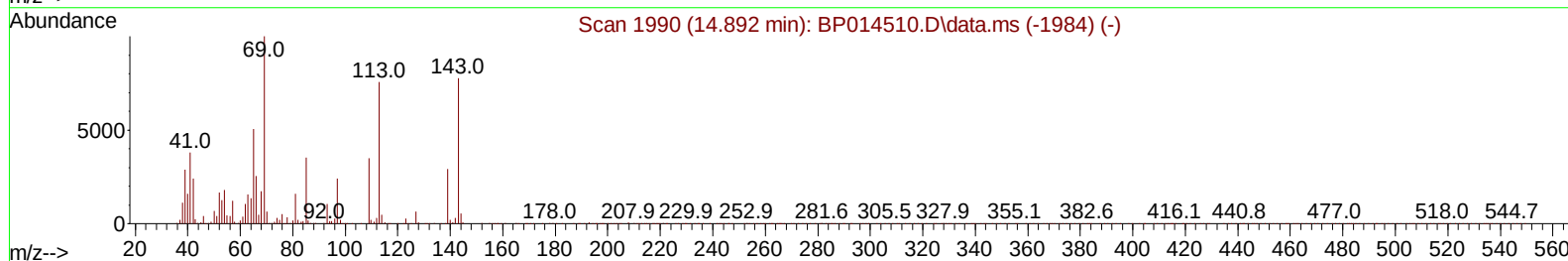
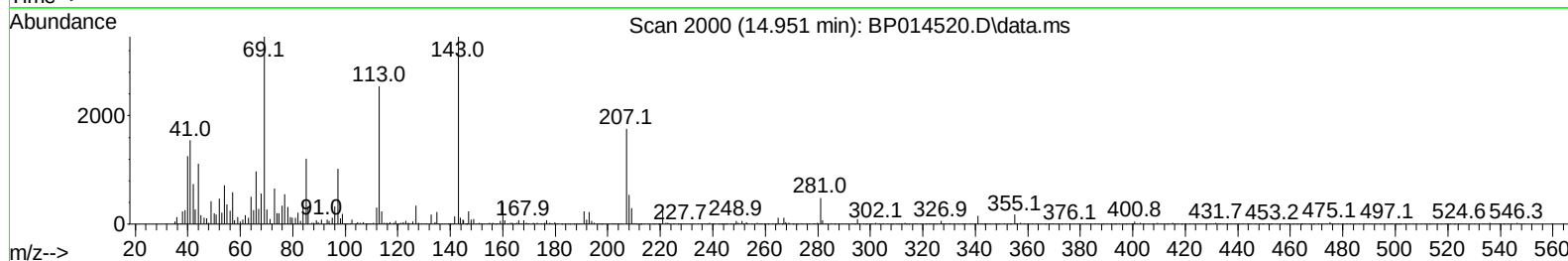
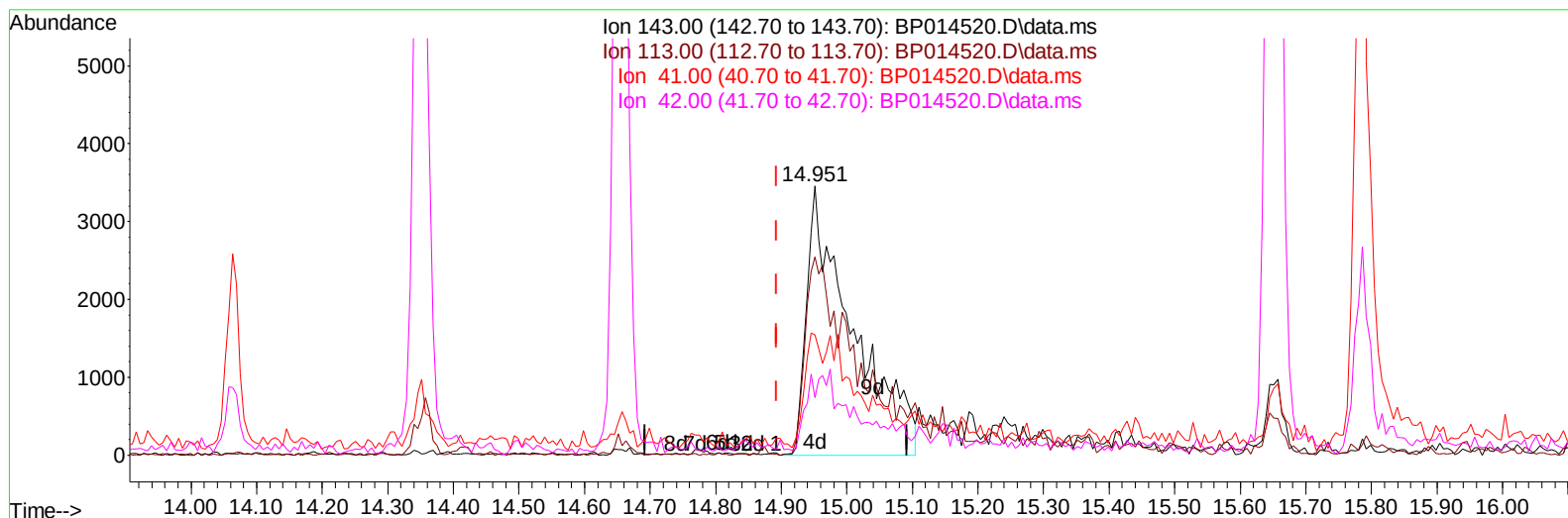
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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TIC: BP014520.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.059) 3.29 ng/ul m

response 16726

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	73.65#
41.00	44.20	44.67
42.00	31.50	21.40#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	144478	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	627743	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	415408	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	849612	20.000	ng/ul	-0.01
79) Chrysene-d12	21.504	240	608577	20.000	ng/ul	-0.01
88) Perylene-d12	24.015	264	638810	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	15763	4.231	ng/uL	0.00
4) Pyridine-d5	3.840	84	28119m	2.945	ng/ul	0.02
7) Phenol-d5	7.181	99	66488	4.990	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	236917	27.713	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	199675	20.734	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	130803	12.170	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	135234	26.686	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	148510	25.593	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	229201	21.688	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	224502	16.128	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	915500	27.181	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	984717	26.348	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	16726m	3.287	ng/ul	0.06
60) Fluorene-d10	15.651	176	768868	27.373	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	125391	20.058	ng/ul	0.00
73) Anthracene-d10	17.522	188	1282367	31.188	ng/ul	0.00
81) Pyrene-d10	19.745	212	1326371	32.447	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	1054893	31.081	ng/ul	-0.01

Target Compounds Qvalue

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

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