

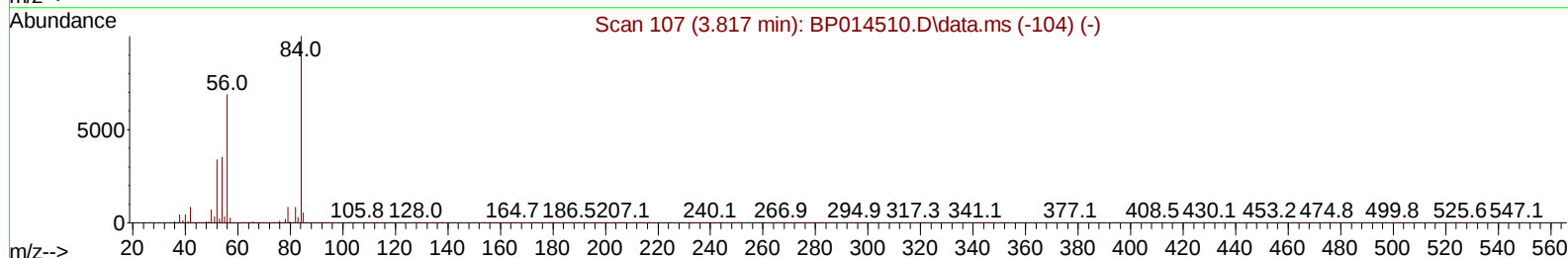
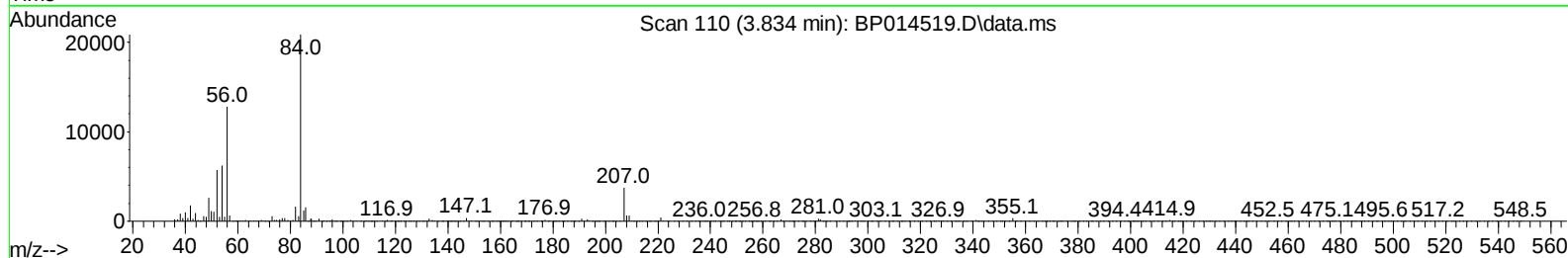
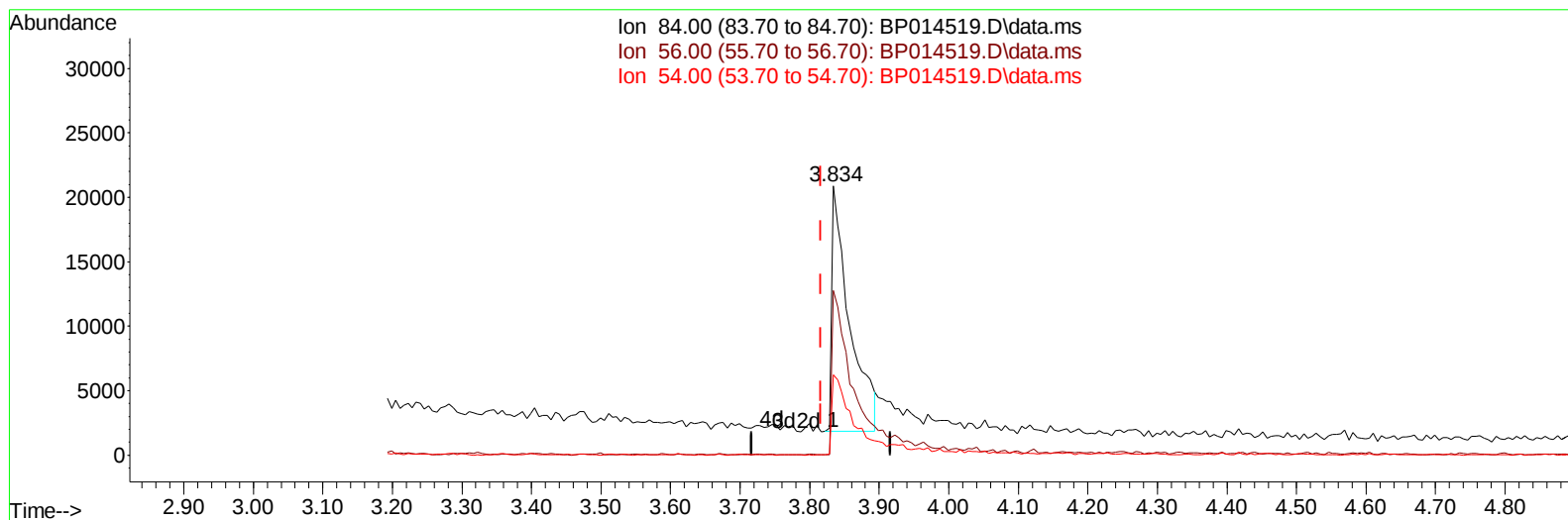
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014519.D
 Acq On : 04 Apr 2023 09:49
 Operator : CG/JU
 Sample : 02158-19
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB(20230329)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:08:36 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: BP014519.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 4.20 ng/u1

response 33436

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	61.21
54.00	32.60	29.91
0.00	0.00	0.00

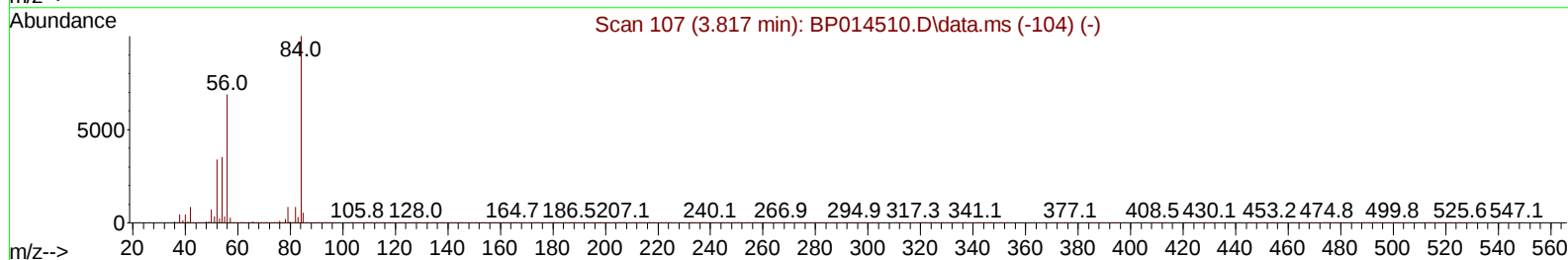
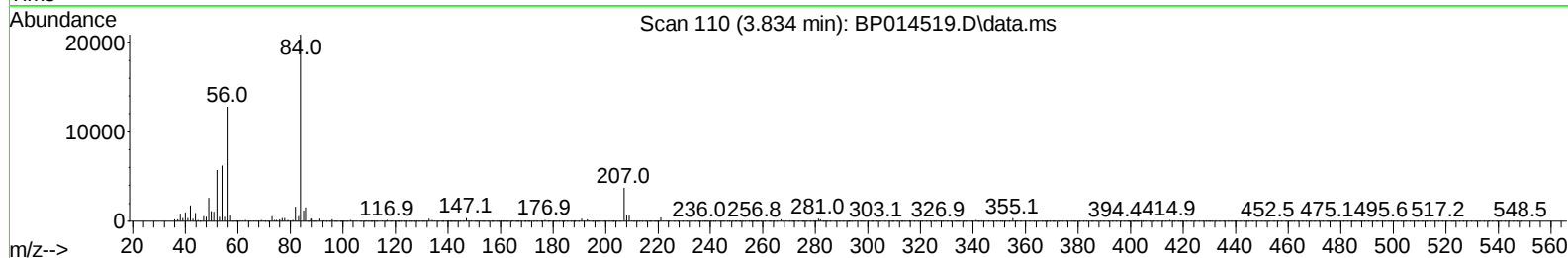
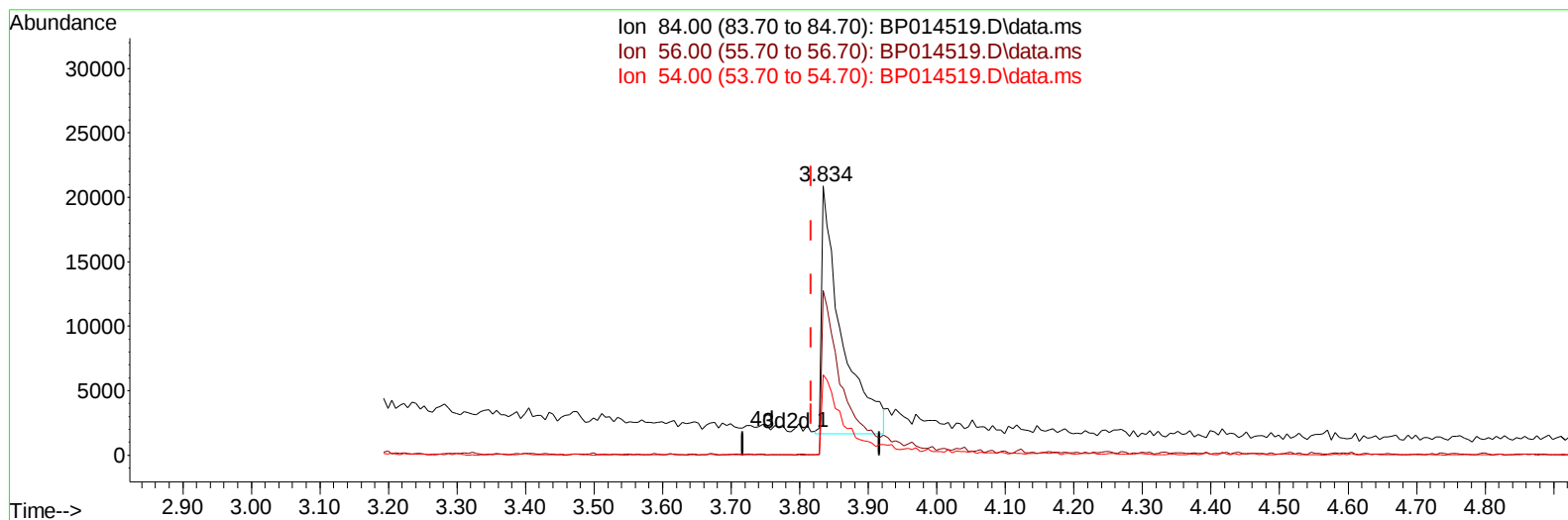
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TIC: BP014519.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 4.86 ng/ul m

response 38677

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	61.21
54.00	32.60	29.91
0.00	0.00	0.00

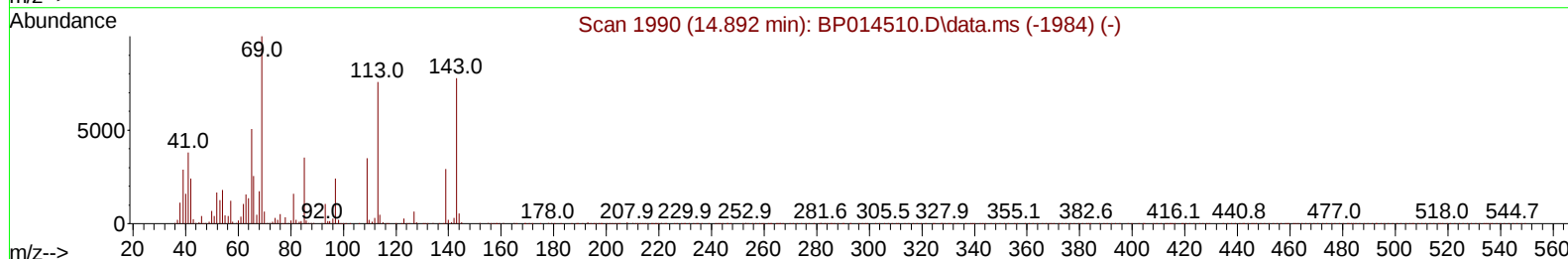
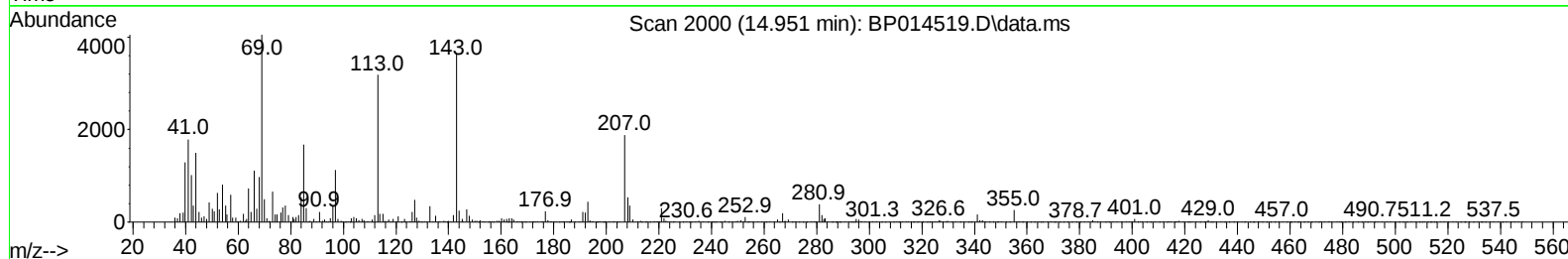
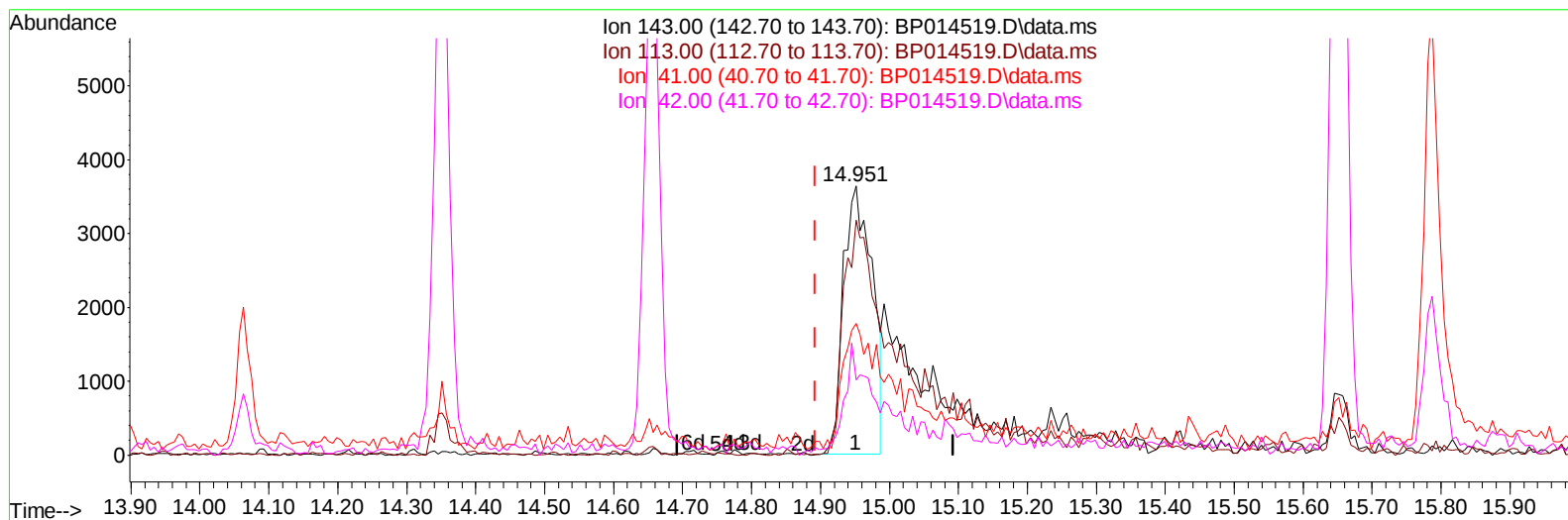
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ALS Vial : 42 Sample Multiplier: 1

Instrument :
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TIC: BP014519.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.059) 2.86 ng/u1

response 10514

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.32
41.00	44.20	49.05
42.00	31.50	27.94

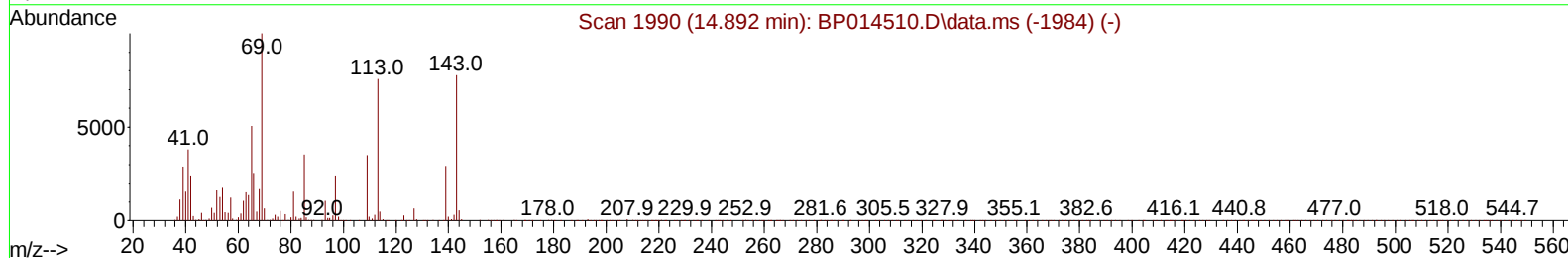
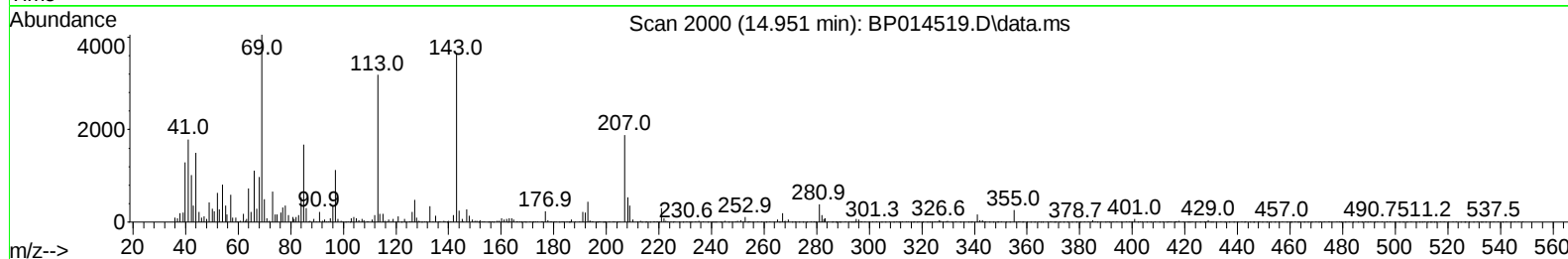
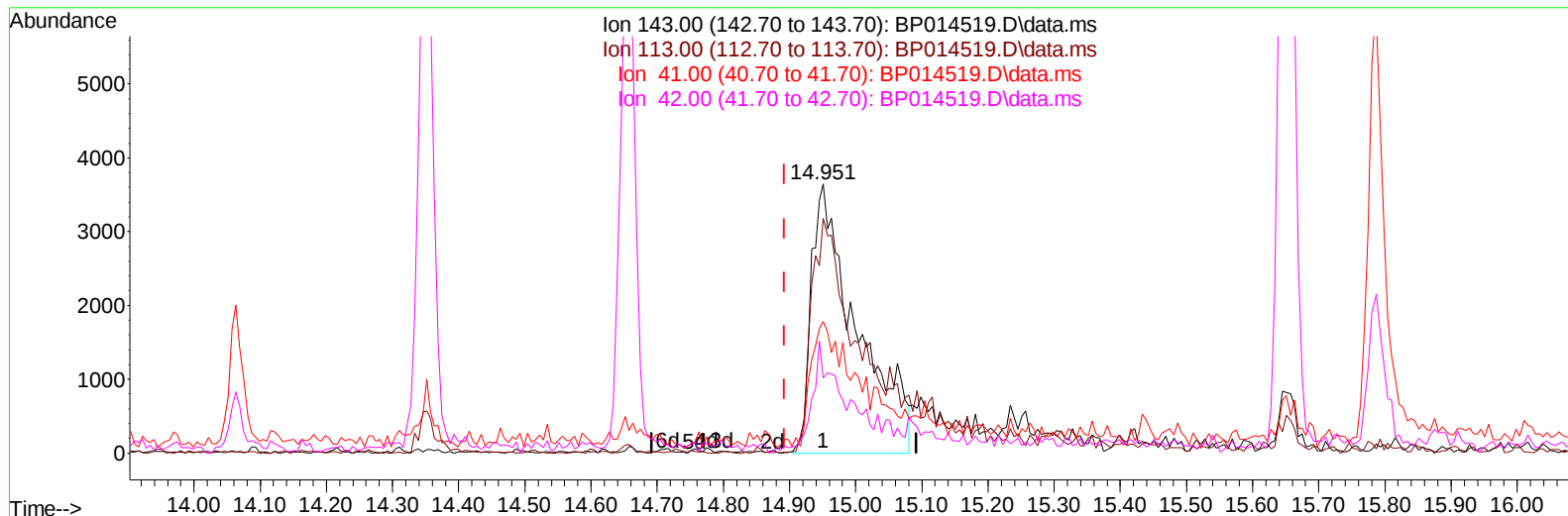
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Sample : 02158-19
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP014519.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.059) 4.72 ng/ul m

response 17374

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.32
41.00	44.20	49.05
42.00	31.50	27.94

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	120413	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	482979	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	300226	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	708172	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	765468	20.000	ng/ul	0.00
88) Perylene-d12	24.016	264	873509	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	14358	4.624	ng/uL	0.00
4) Pyridine-d5	3.834	84	38677m	4.861	ng/ul	0.02
7) Phenol-d5	7.181	99	64932	5.847	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	205633	28.860	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	180511	22.490	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	116163	12.968	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	110803	28.418	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	119243	26.708	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	191680	23.574	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	205380	19.176	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	734410	30.169	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	803939	29.763	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	17374m	4.724	ng/ul	0.06
60) Fluorene-d10	15.651	176	640856	31.569	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	108810	20.882	ng/ul	0.00
73) Anthracene-d10	17.522	188	1169616	34.127	ng/ul	0.00
81) Pyrene-d10	19.751	212	1483297	28.849	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.857	264	1600964	34.497	ng/ul	0.00

Target Compounds Qvalue

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

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