

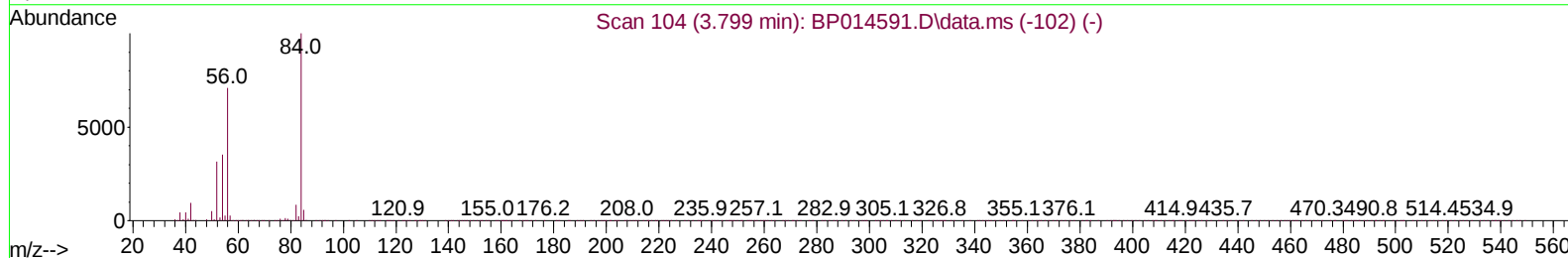
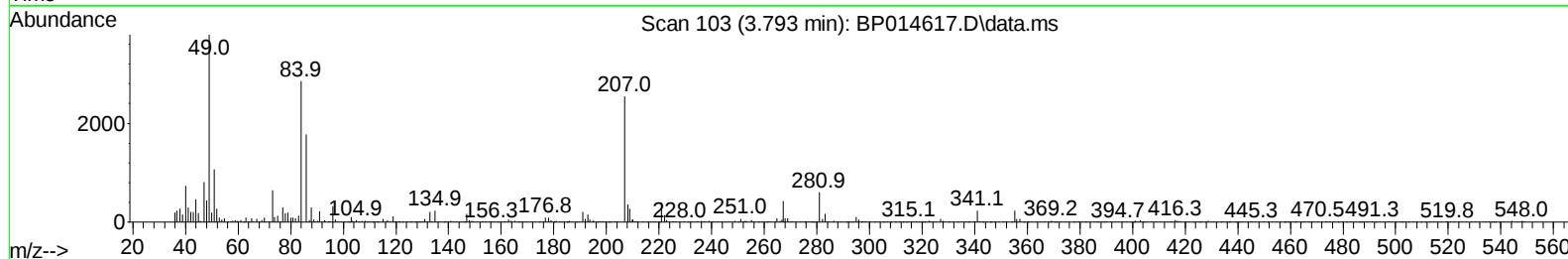
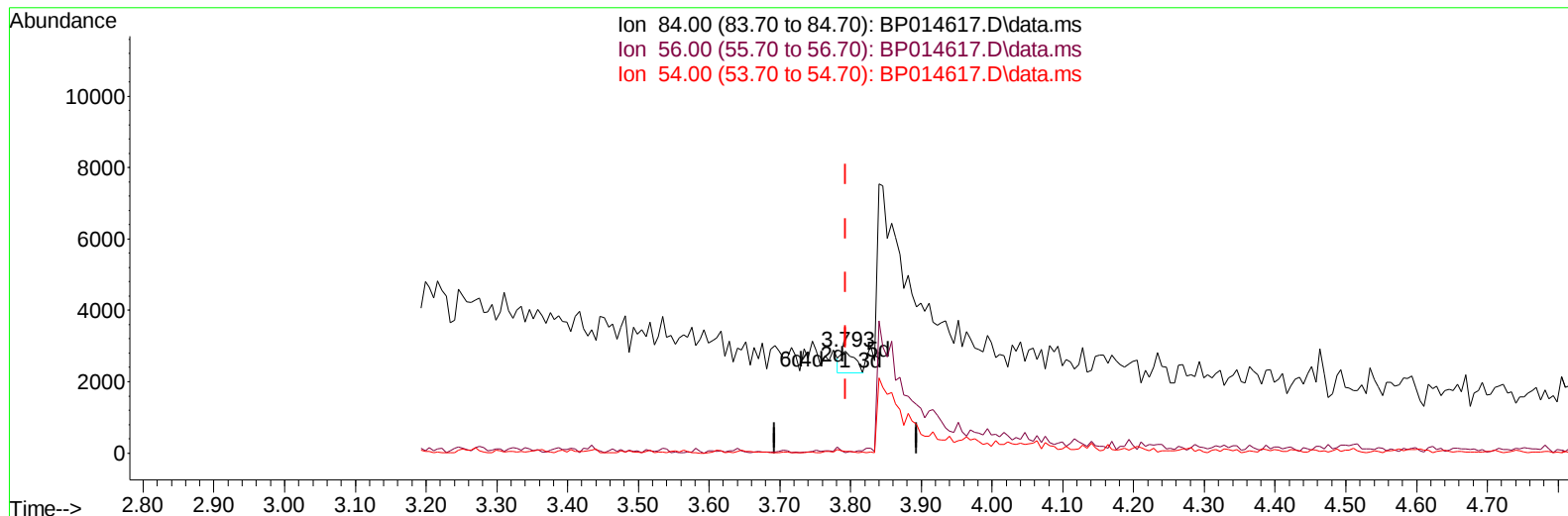
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
Data File : BP014617.D
Acq On : 07 Apr 2023 22:21
Operator : CG/JU
Sample : 02225-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD4

Manual IntegrationsAPPROVED

Quant Time: Apr 08 01:51:28 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 07 00:45:34 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/10/2023
Supervised By :Jagrut Upadhyay 04/10/2023



TIC: BP014617.D\data.ms

(4) Pyridine-d5 (S)

3.793min (-0.000) 0.09 ng/u1

response 826

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.59#
54.00	32.60	1.68#
0.00	0.00	0.00

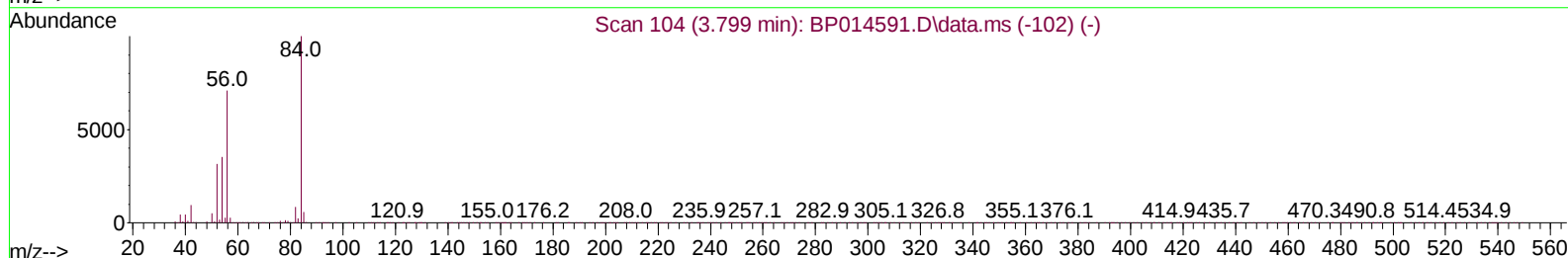
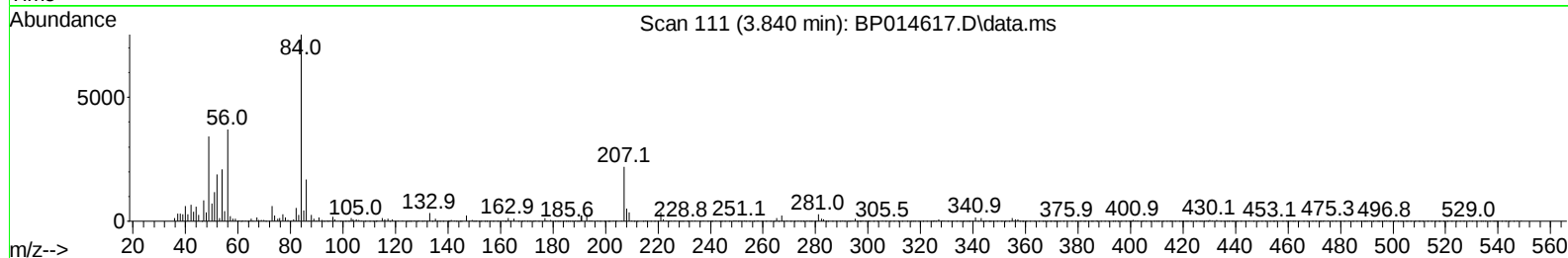
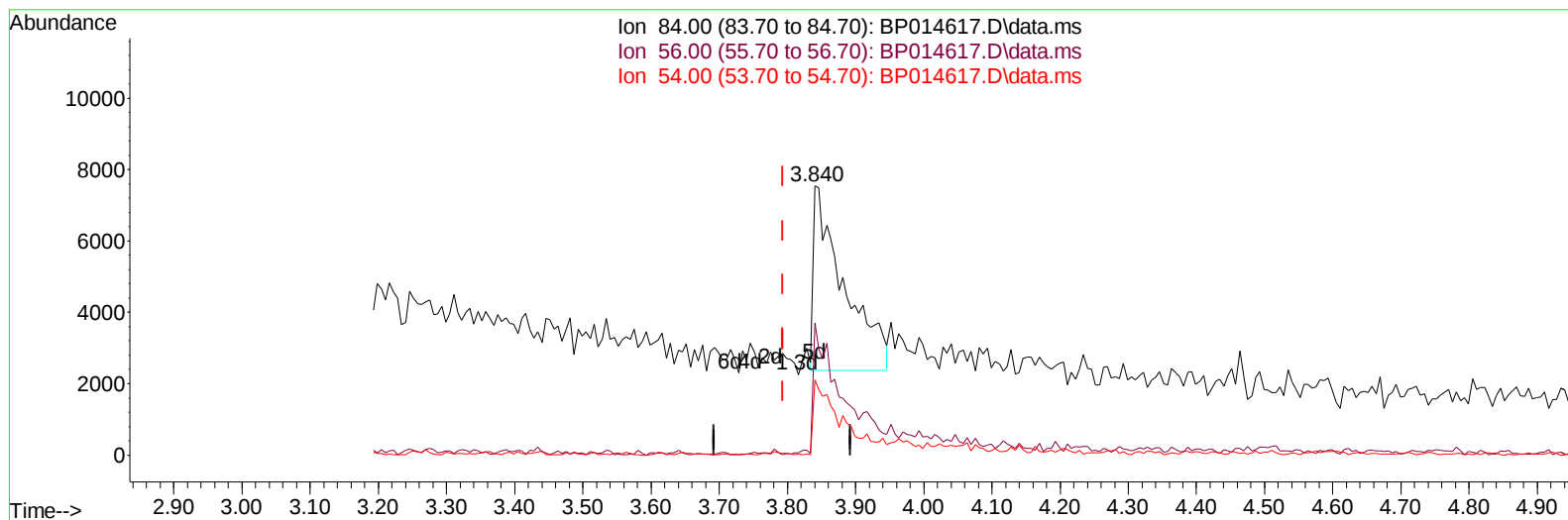
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(4) Pyridine-d5 (S)

3.840min (+ 0.047) 1.79 ng/ul m

response 16104

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	49.18#
54.00	32.60	27.94
0.00	0.00	0.00

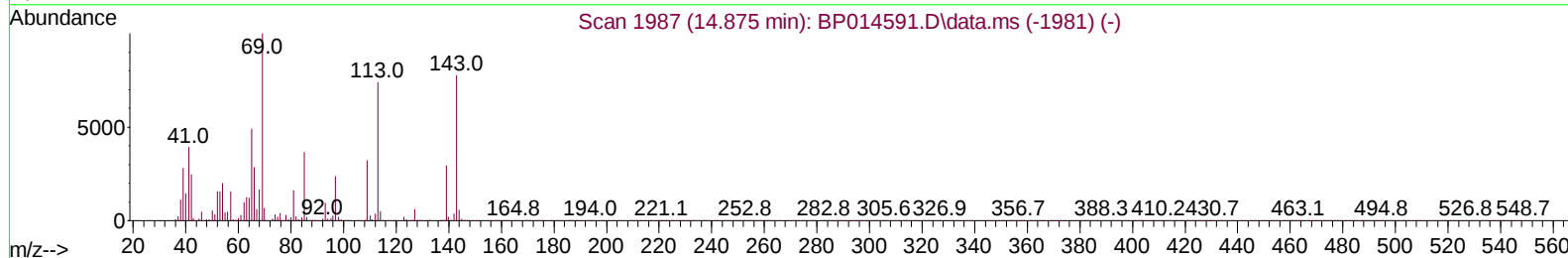
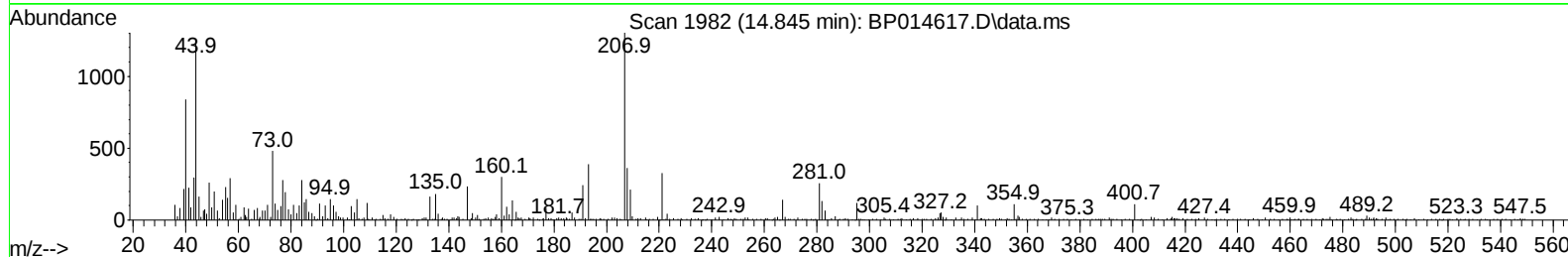
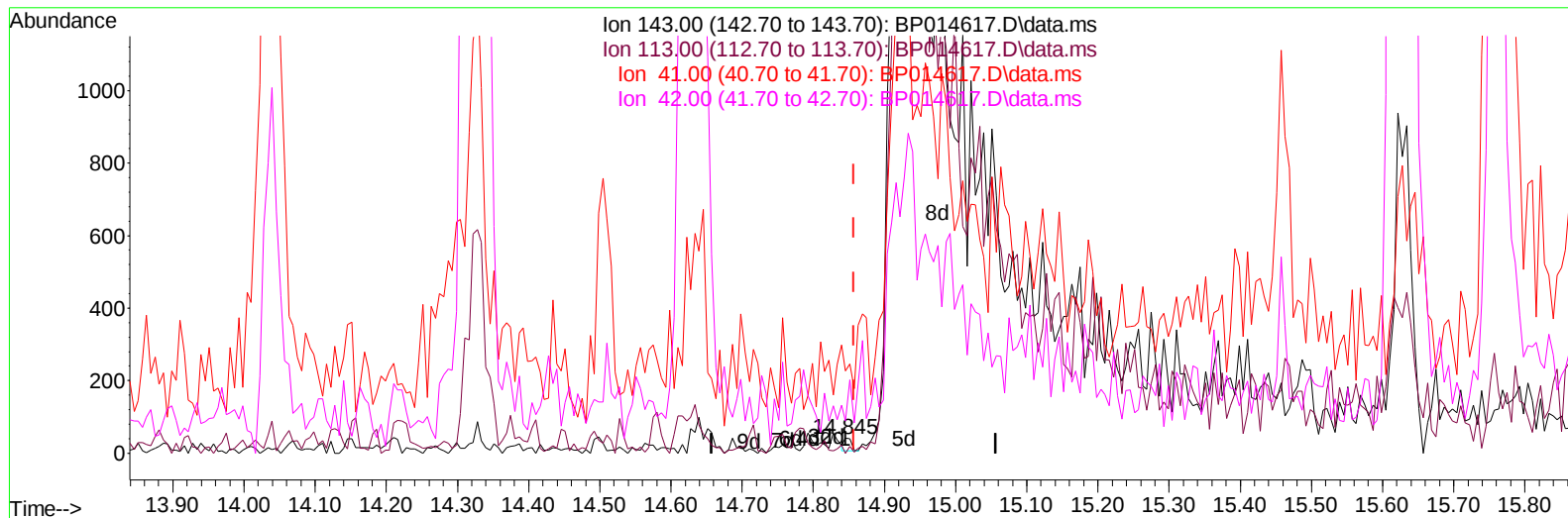
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TIC: BP014617.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.845min (-0.012) 0.01 ng/u1

response 24

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	0.00#
41.00	44.20	553.66#
42.00	31.50	217.07#

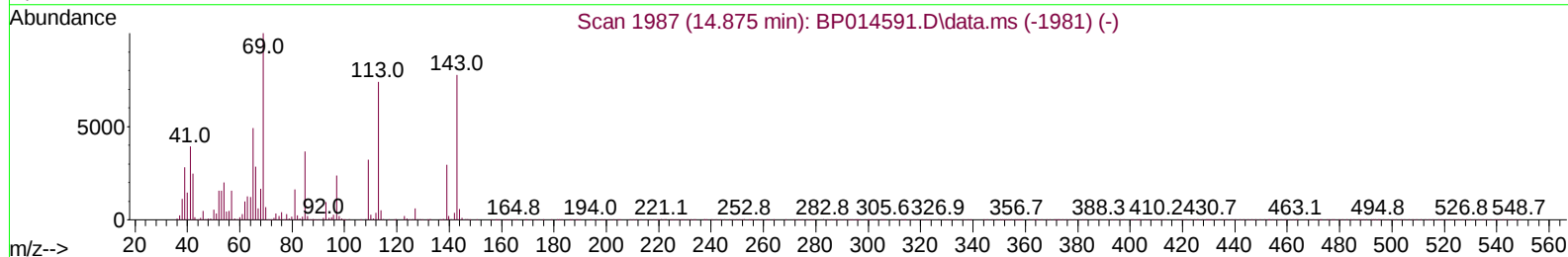
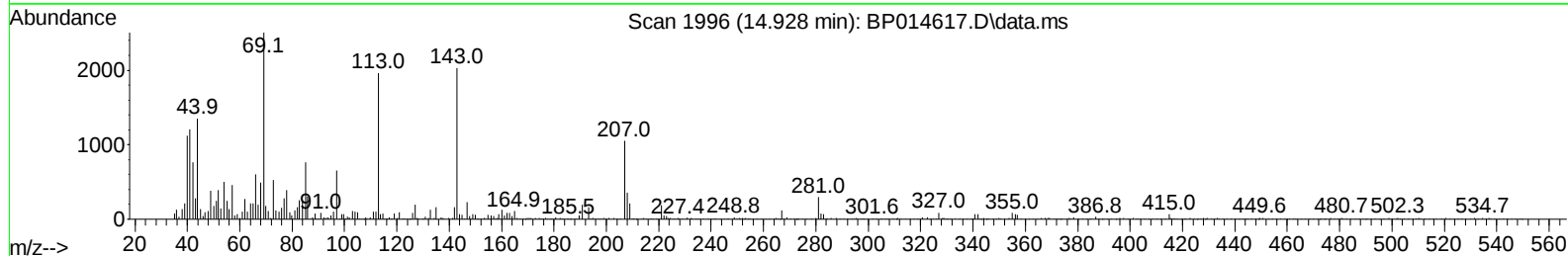
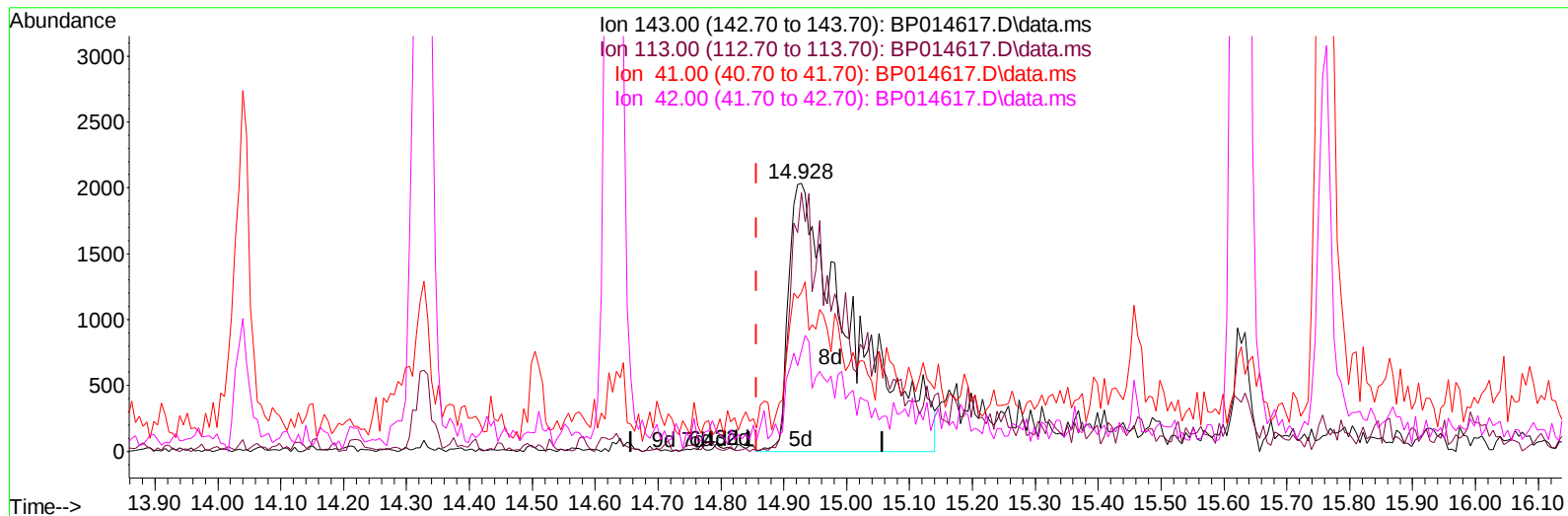
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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TIC: BP014617.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.071) 2.95 ng/ul m

response 13938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	96.56
41.00	44.20	59.37#
42.00	31.50	37.53

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	136102	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	607246	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	385281	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	840256	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	628454	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	676465	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	14314	4.078	ng/uL	0.00
4) Pyridine-d5	3.840	84	16104m	1.791	ng/ul	0.05
7) Phenol-d5	7.152	99	55199	4.398	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	216906	26.933	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	172356	18.998	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	111290	10.992	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	114988	23.456	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	127674	22.745	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	198926	19.459	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	120852	8.975	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	890993	28.522	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	930351	26.840	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	13938m	2.953	ng/ul	0.07
60) Fluorene-d10	15.628	176	724947	27.828	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	129533	20.951	ng/ul	0.00
73) Anthracene-d10	17.492	188	1154761	28.397	ng/ul	0.00
81) Pyrene-d10	19.722	212	1287262	30.494	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1080120	30.053	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.399	88	8602	2.230	ng/uL	93

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

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