

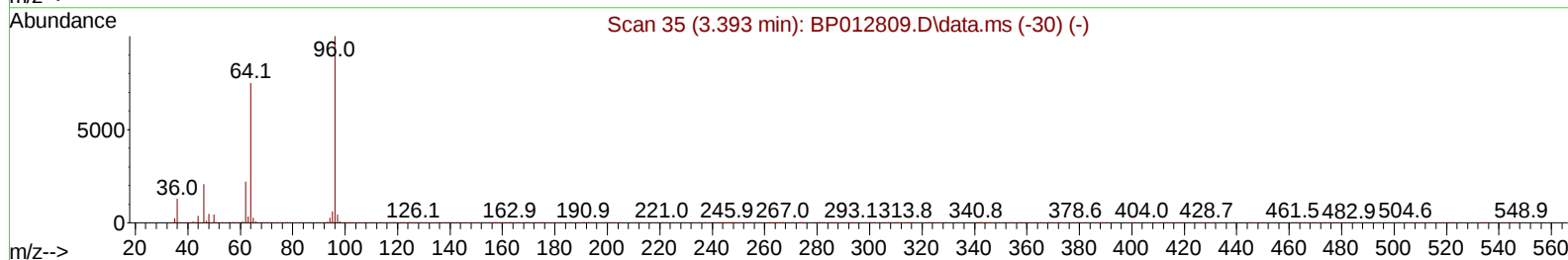
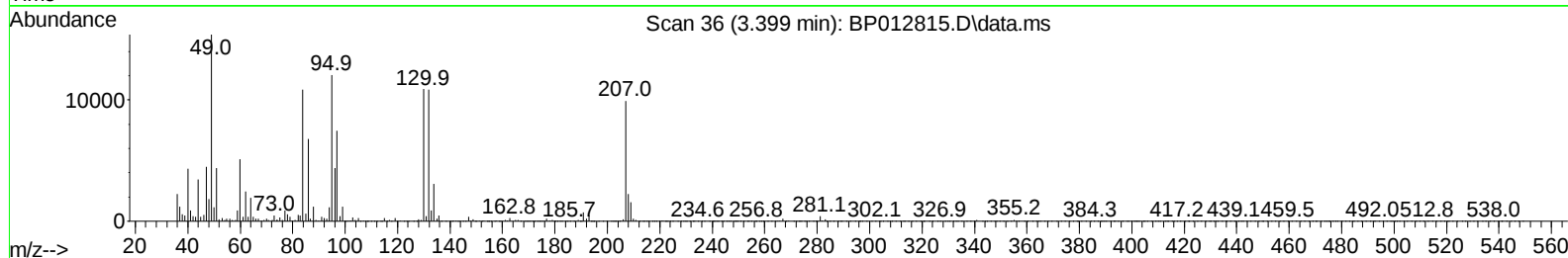
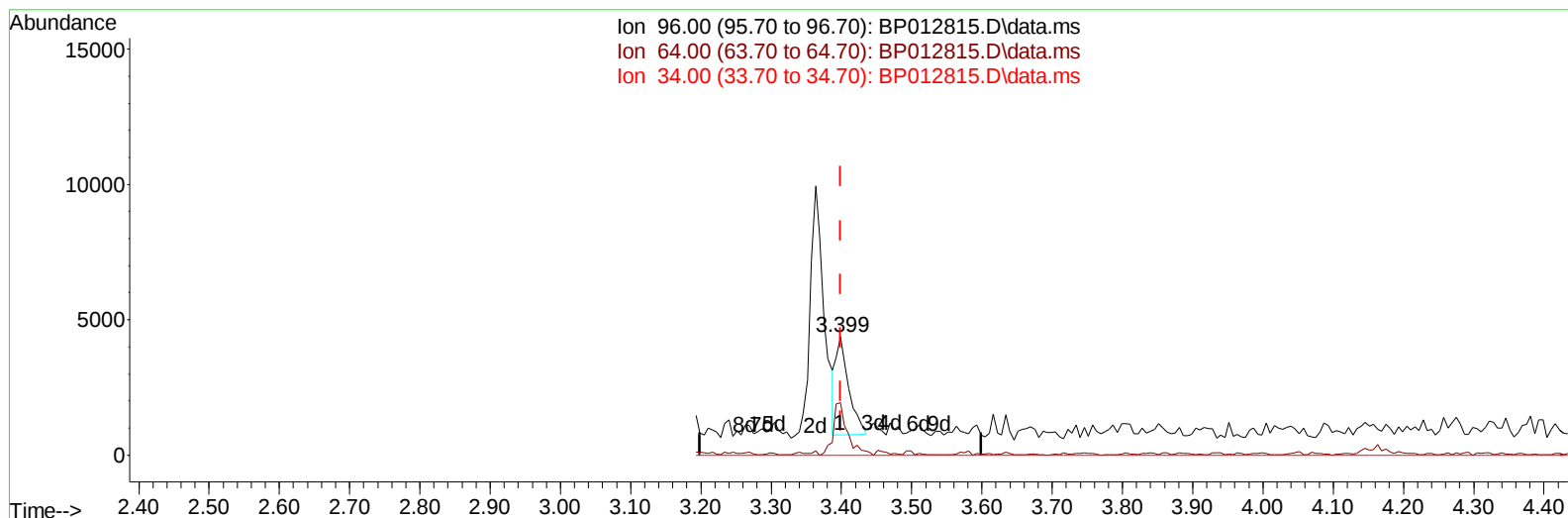
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012815.D
 Acq On : 28 Nov 2022 14:13
 Operator : CG/JU
 Sample : N5710-11DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KZ0DL

Manual IntegrationsAPPROVED

Quant Time: Nov 28 22:39:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/29/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012815.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.399min (-0.000) 0.37 ng/uL

response 4652

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	44.45#
34.00	0.00	0.00
0.00	0.00	0.00

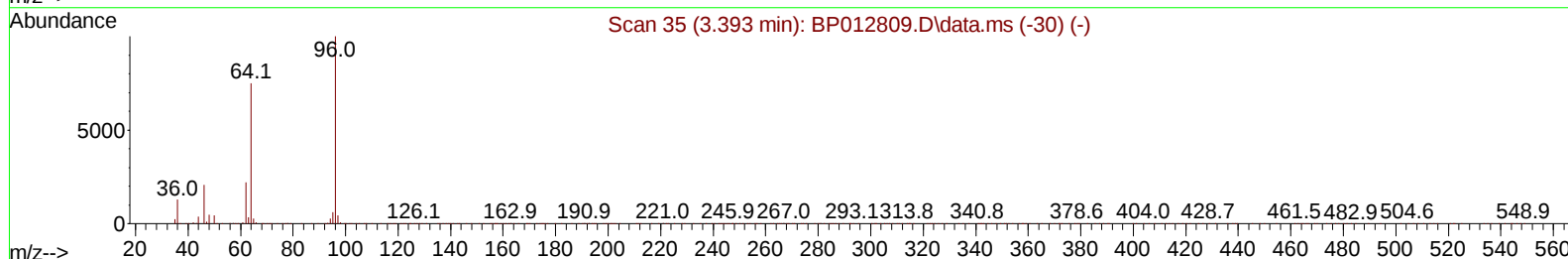
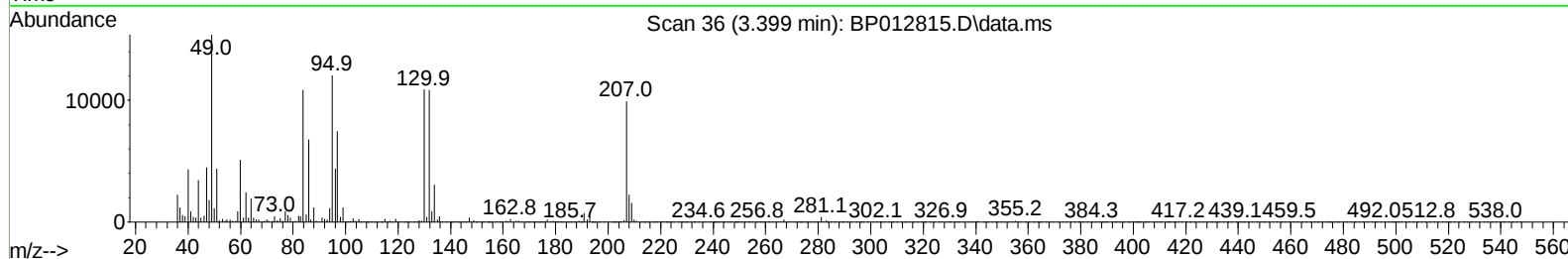
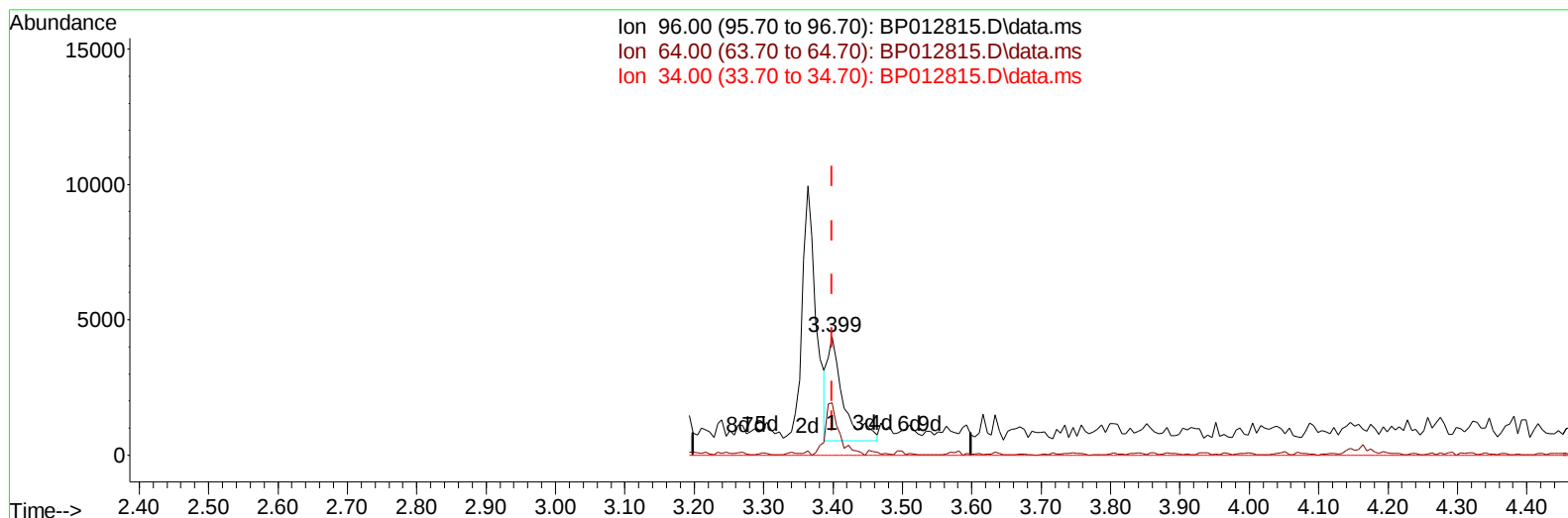
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TIC: BP012815.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.399min (-0.000) 0.48 ng/uL m

response 6097

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	44.45#
34.00	0.00	0.00
0.00	0.00	0.00

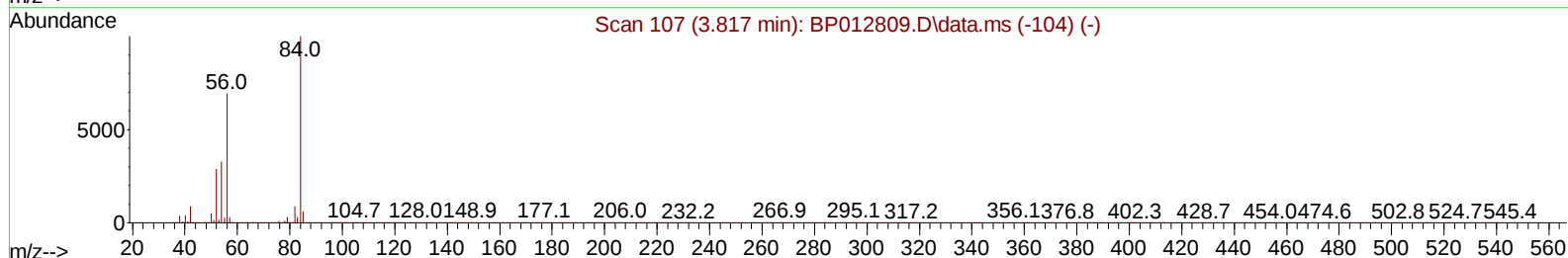
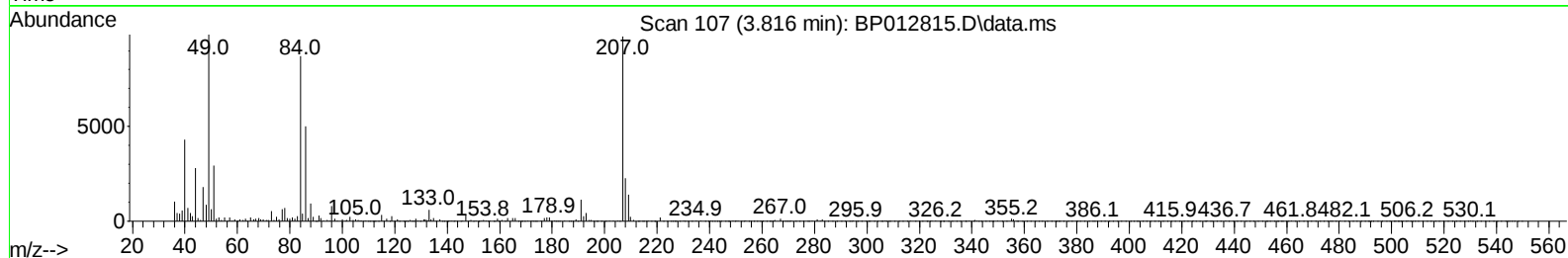
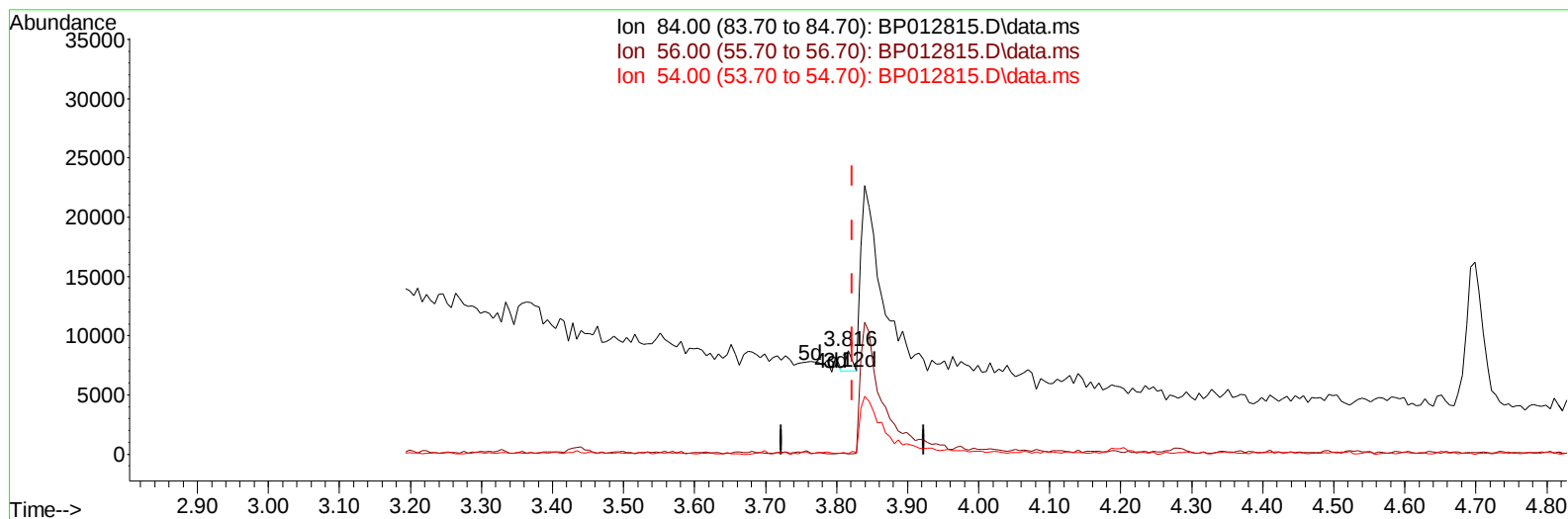
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TIC: BP012815.D\data.ms

(4) Pyridine-d5 (S)

3.816min (-0.006) 0.03 ng/u1

response 1035

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	0.45#
54.00	31.30	0.64#
0.00	0.00	0.00

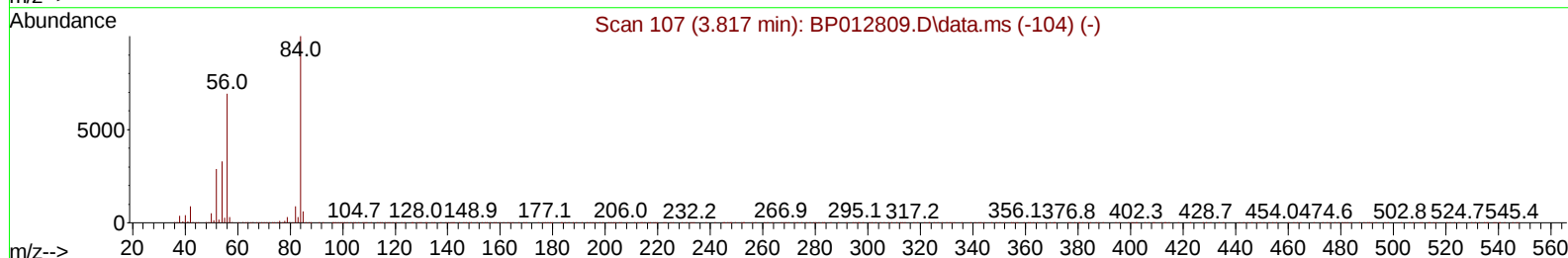
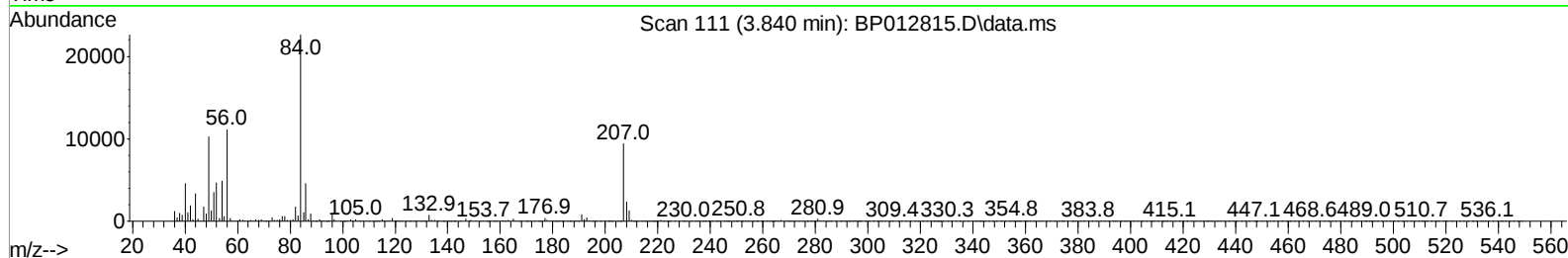
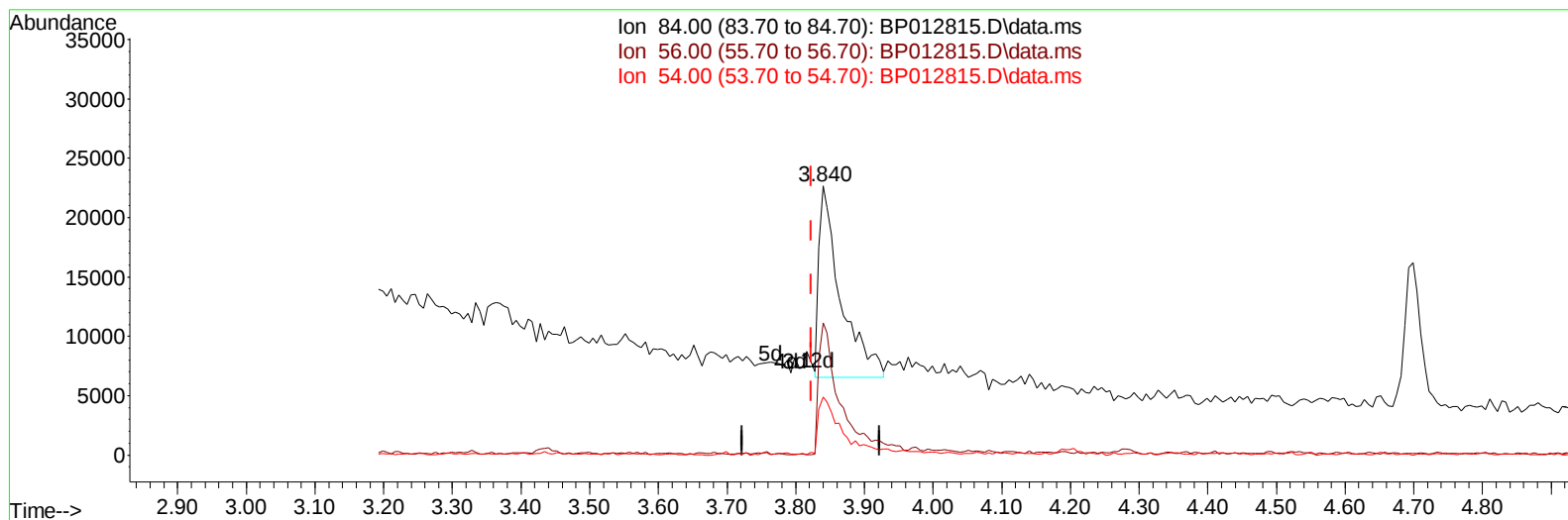
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TIC: BP012815.D\data.ms

(4) Pyridine-d5 (S)

3.840min (+ 0.017) 0.93 ng/u1 m

response 35294

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	49.25#
54.00	31.30	21.56#
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	8.016	152	545085	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	2380192	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	1579181	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	3526674	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	3612526	20.000	ng/ul	0.00
88) Perylene-d12	24.009	264	3540103	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	6097m	0.479	ng/uL	0.00
4) Pyridine-d5	3.840	84	35294m	0.927	ng/ul	0.02
7) Phenol-d5	7.163	99	23692	0.529	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	90260	3.270	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	70548	2.067	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	39059	1.082	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	35304	2.423	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	34296	2.217	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	80812	2.239	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	98707	1.997	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	347836	3.200	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	358053	2.823	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.645	176	321485	3.377	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.510	188	510284	3.287	ng/ul	0.00
81) Pyrene-d10	19.733	212	614937	2.930	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	490299	2.820	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.434	88	339329	24.839	ng/uL	98

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

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