

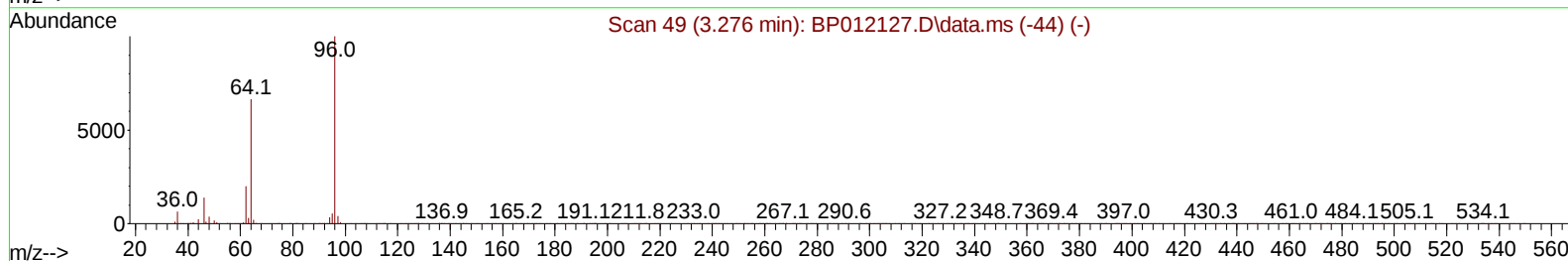
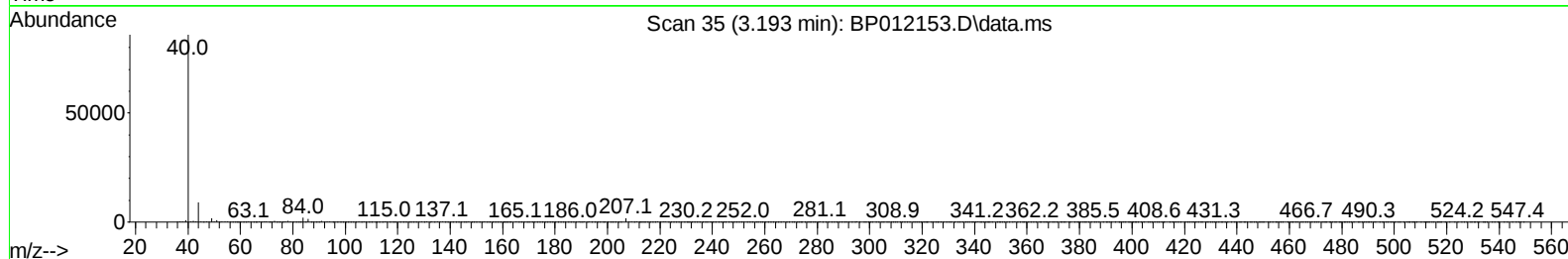
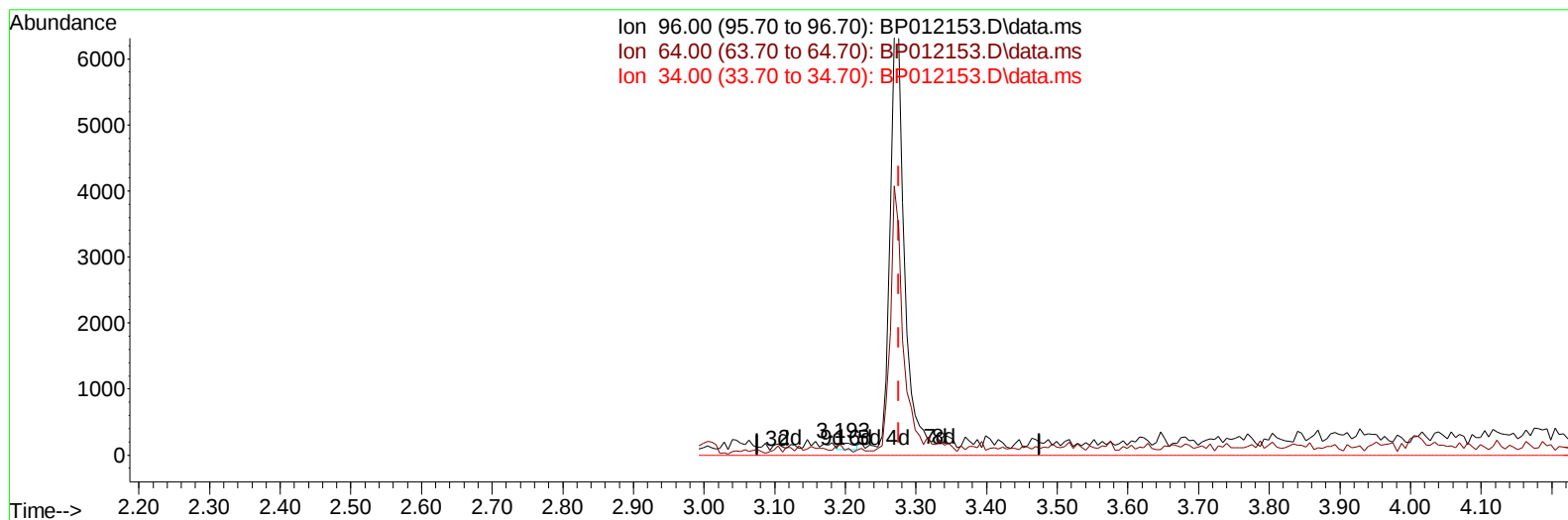
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012153.D
 Acq On : 15 Oct 2022 00:27
 Operator : CG/JU
 Sample : N4984-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGNB5

Manual IntegrationsAPPROVED

Quant Time: Oct 15 02:22:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/15/2022
 Supervised By :mohammad ahmed 10/16/2022



TIC: BP012153.D\data.ms

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

3.193min (-0.082) 0.02 ng/uL

response	155	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	84.95
34.00	0.00	0.00
0.00	0.00	0.00

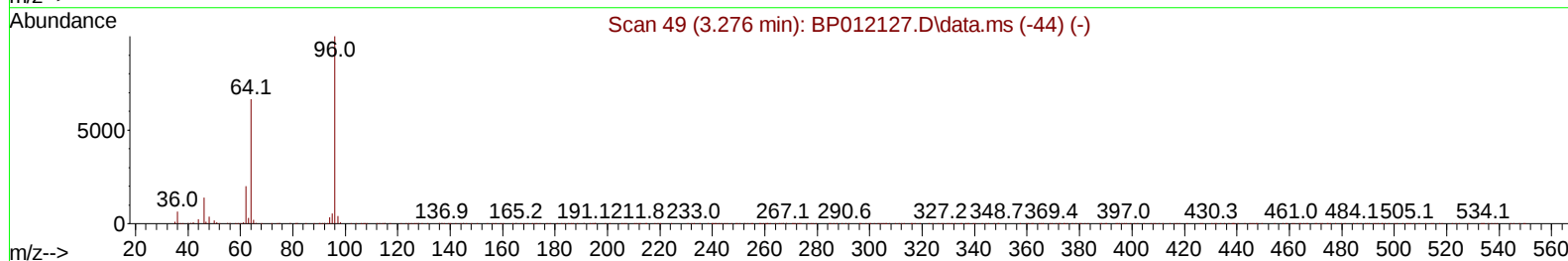
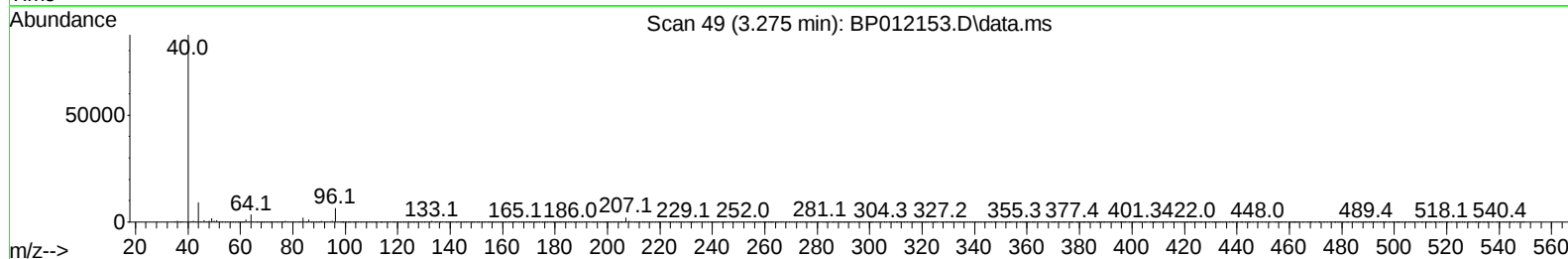
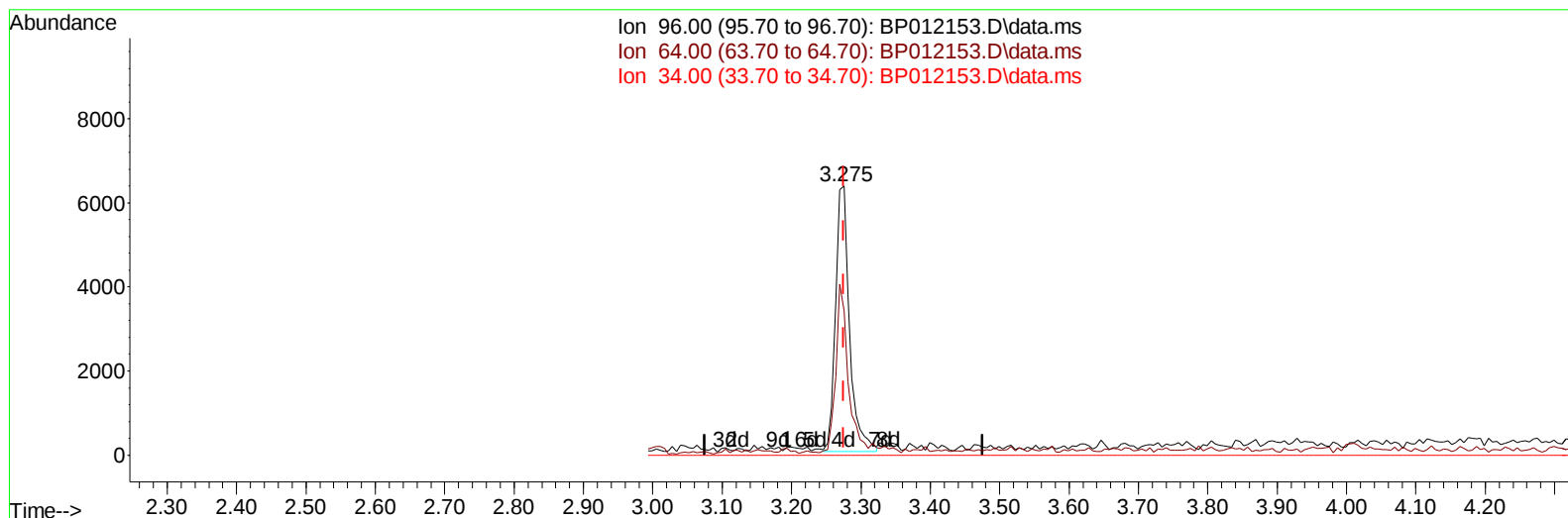
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(3) 1,4-Dioxane-d8 (S)

3.275min (-0.000) 1.34 ng/uL m

response 8852

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	53.72#
34.00	0.00	0.00
0.00	0.00	0.00

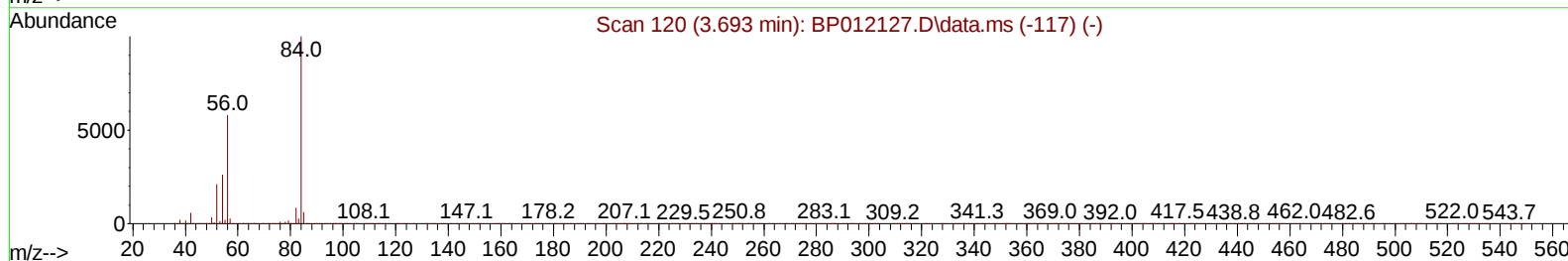
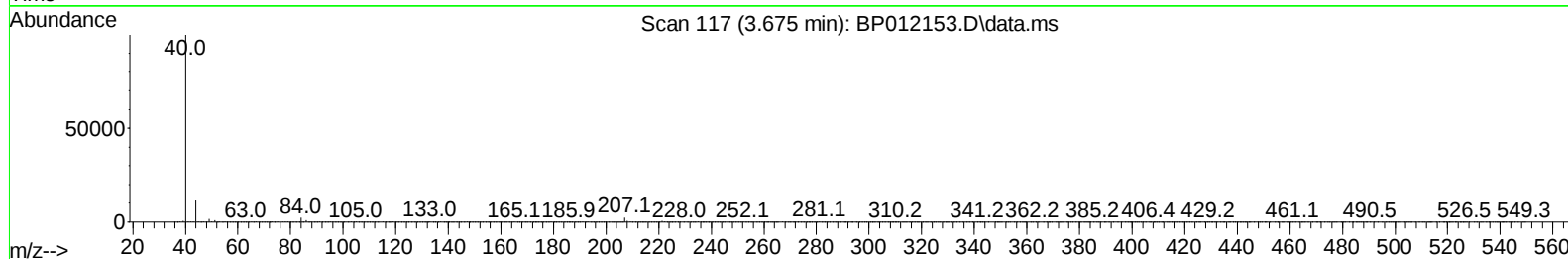
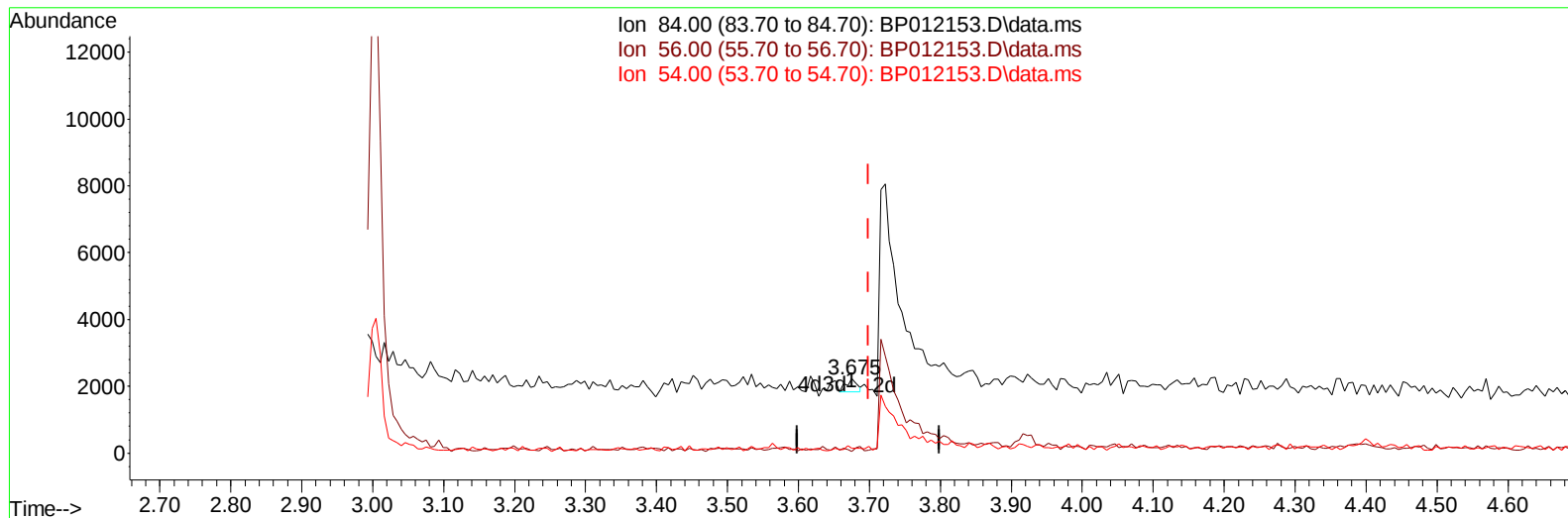
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TIC: BP012153.D\data.ms

(4) Pyridine-d5 (S)

3.675min (-0.024) 0.02 ng/u1

response 408

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	1.93#
54.00	26.00	6.41#
0.00	0.00	0.00

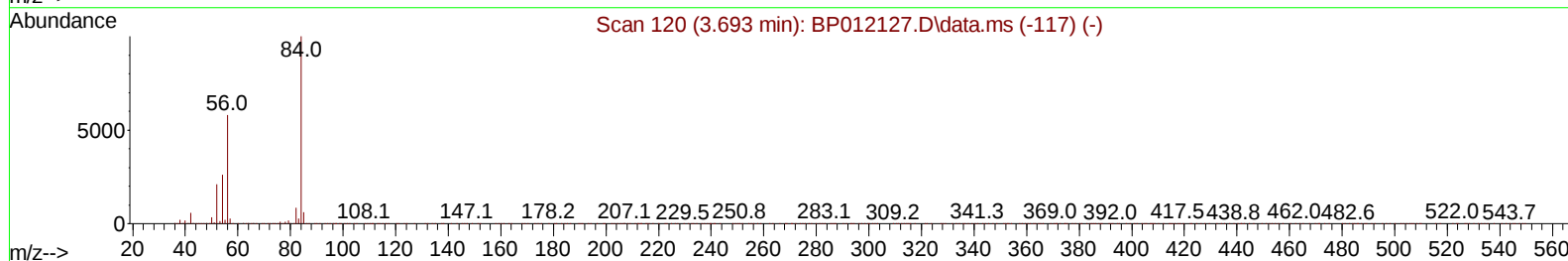
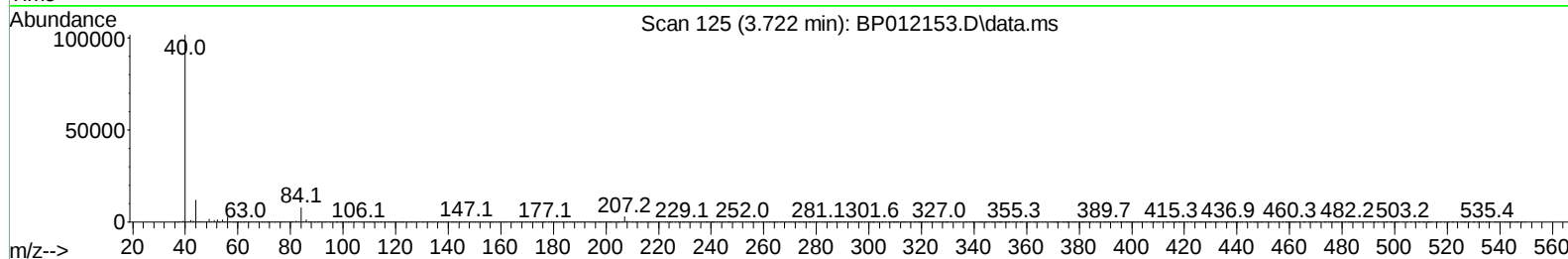
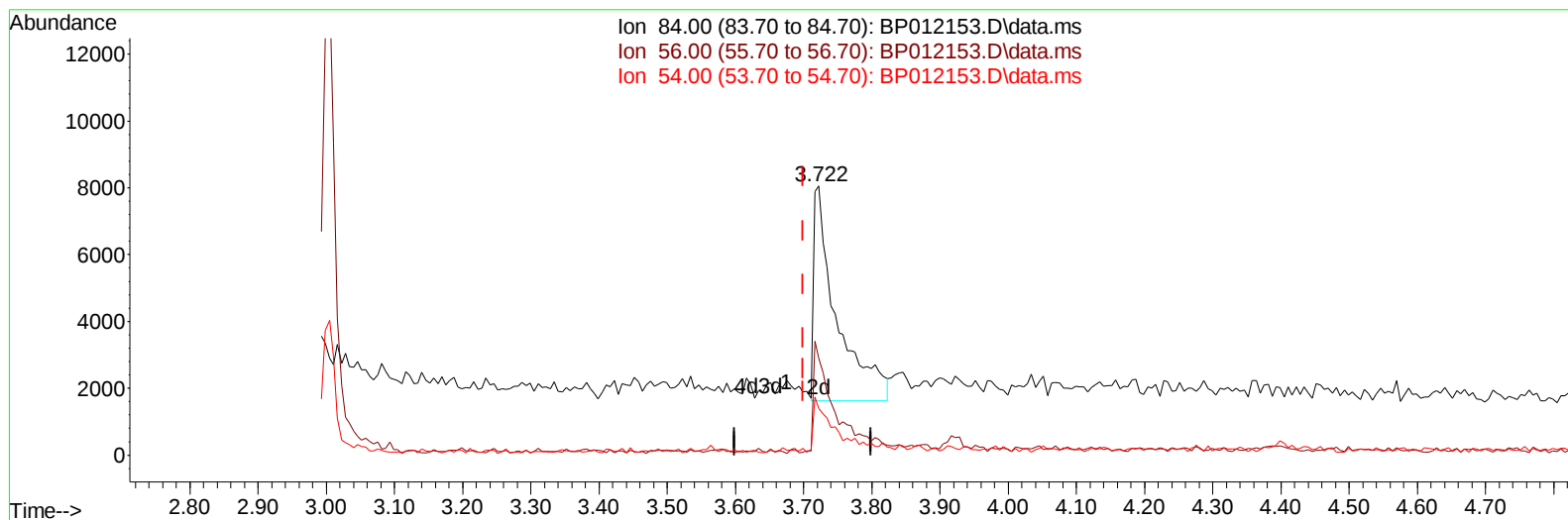
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TIC: BP012153.D\data.ms

(4) Pyridine-d5 (S)

3.722min (+ 0.023) 0.86 ng/u1 m

response 15042

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	36.00#
54.00	26.00	17.37#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	259626	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.646	136	1087324	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.492	164	636010	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1359526	20.000	ng/ul	0.00
79) Chrysene-d12	21.339	240	1371512	20.000	ng/ul	# 0.00
88) Perylene-d12	23.757	264	1441671	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	8852m	1.344	ng/uL	0.00
4) Pyridine-d5	3.722	84	15042m	0.859	ng/ul	0.02
7) Phenol-d5	7.010	99	394034	18.871	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	232455	19.613	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.369	132	317950	19.183	ng/ul	-0.01
15) 4-Methylphenol-d8	8.557	113	303184	18.904	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	147943	19.265	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	147583	18.953	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.269	165	271722	18.006	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.793	131	334699	15.051	ng/ul	-0.01
46) Dimethylphthalate-d6	13.904	166	809368	20.094	ng/ul	-0.01
49) Acenaphthylene-d8	14.181	160	977048	19.616	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.704	143	81600	10.319	ng/ul	0.00
60) Fluorene-d10	15.486	176	764299	20.972	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.610	200	43230	6.022	ng/ul	-0.01
73) Anthracene-d10	17.351	188	1243229	21.059	ng/ul	0.00
81) Pyrene-d10	19.586	212	1475563	21.000	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.598	264	1467400	20.938	ng/ul	-0.01

Target Compounds Qvalue

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

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