Quantitation Report (Qedit)

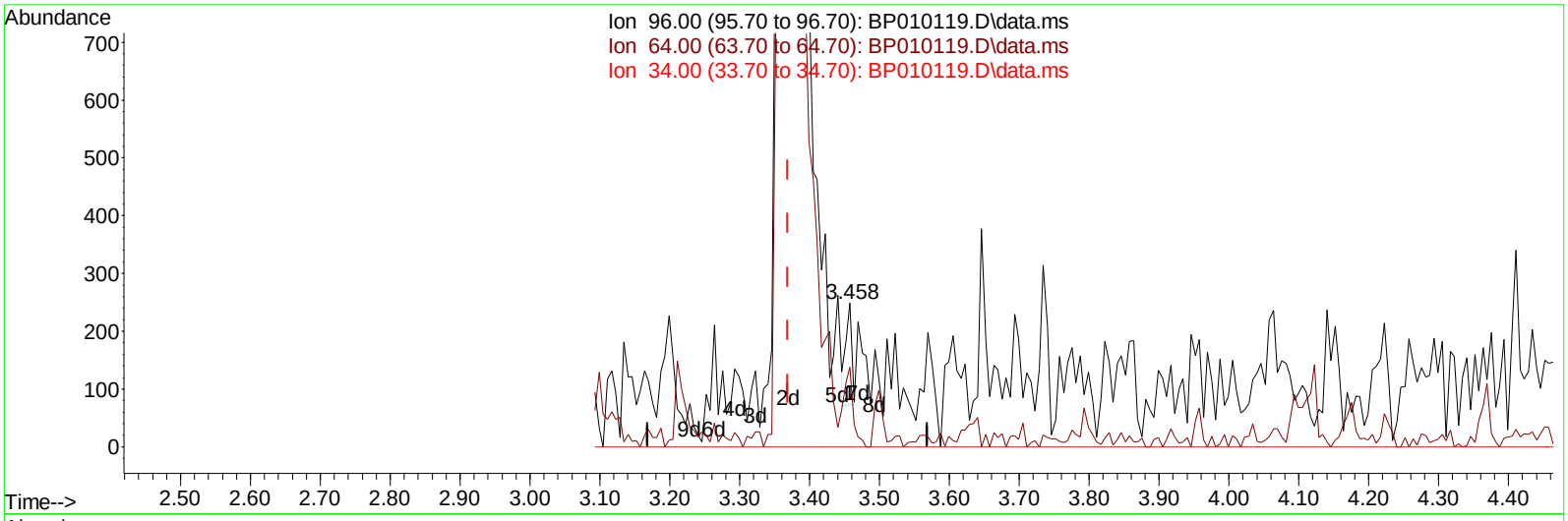
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042722\
Data File : BP010119.D
Acq On : 27 Apr 2022 12:44
Operator : CG/JU
Sample : N2470-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

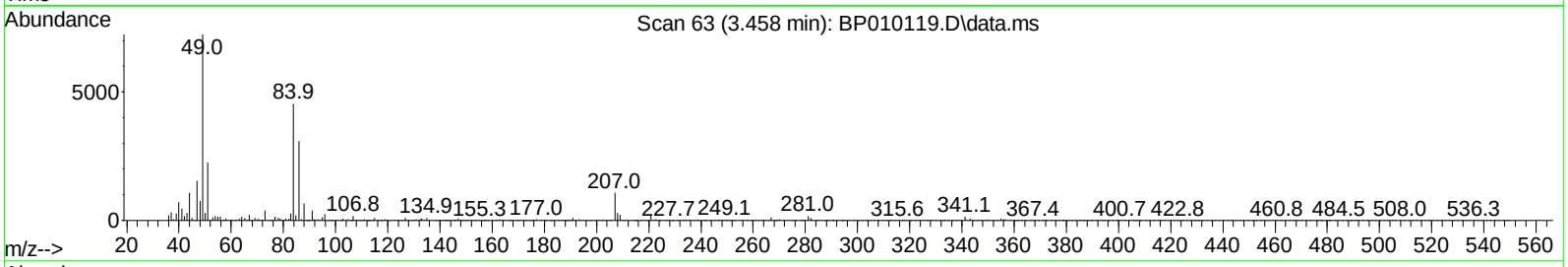
Quant Time: Apr 28 03:24:38 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

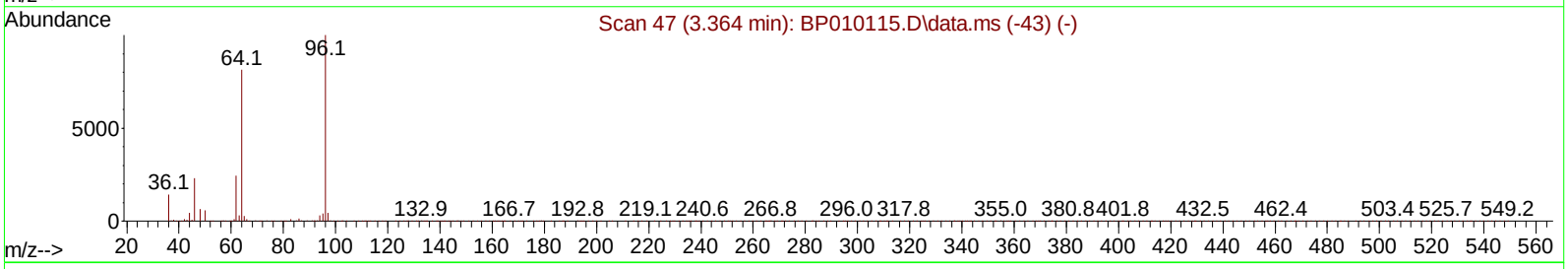
Instrument :
BNA_P
ClientSampleId :
BFFG9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/28/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010119.D\data.ms

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

3.458min (+ 0.089) 0.03 ng/uL

response 103

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	55.42
34.00	0.00	0.00
0.00	0.00	0.00

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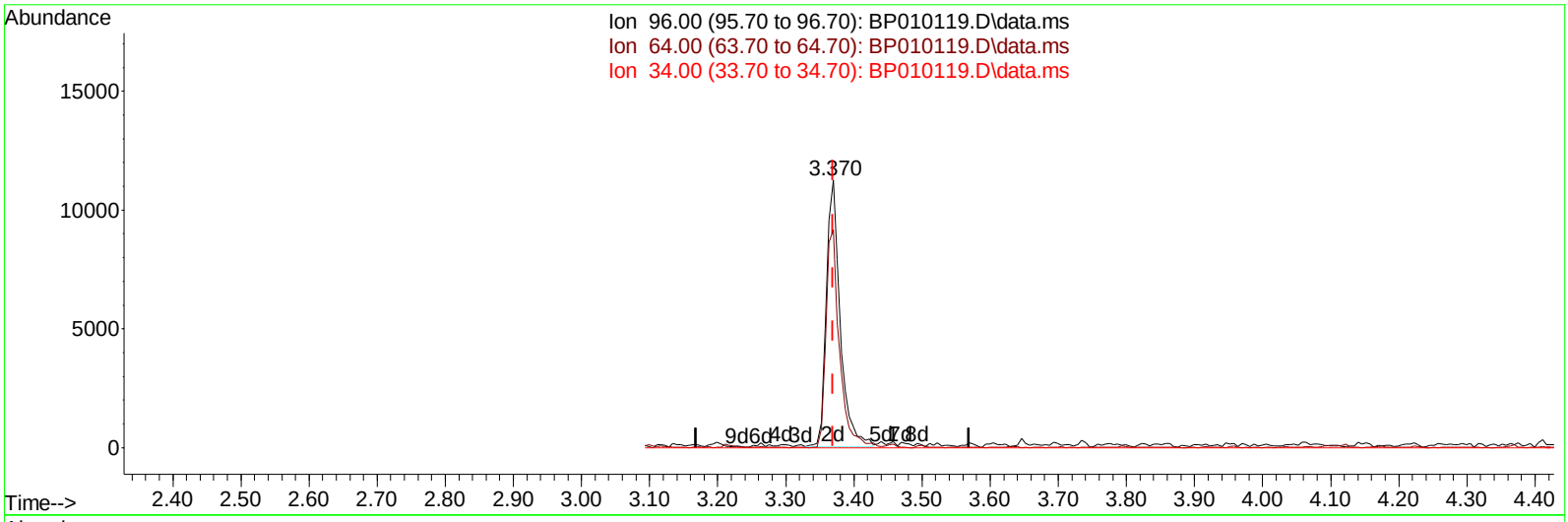

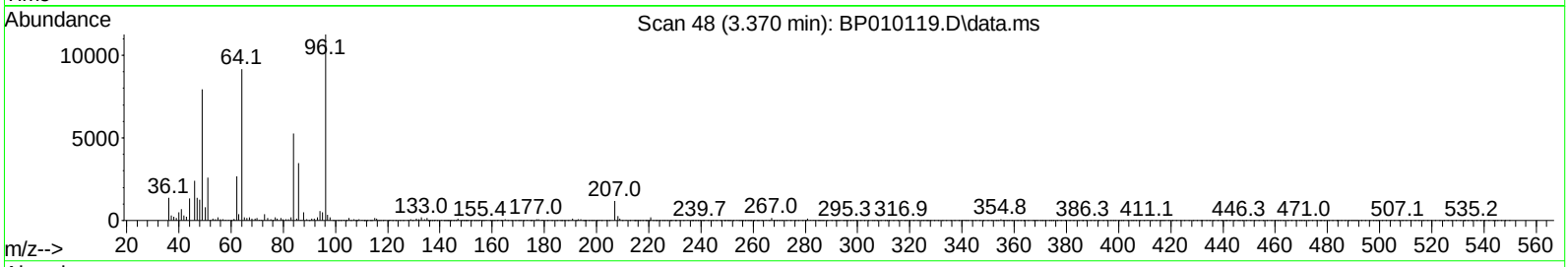

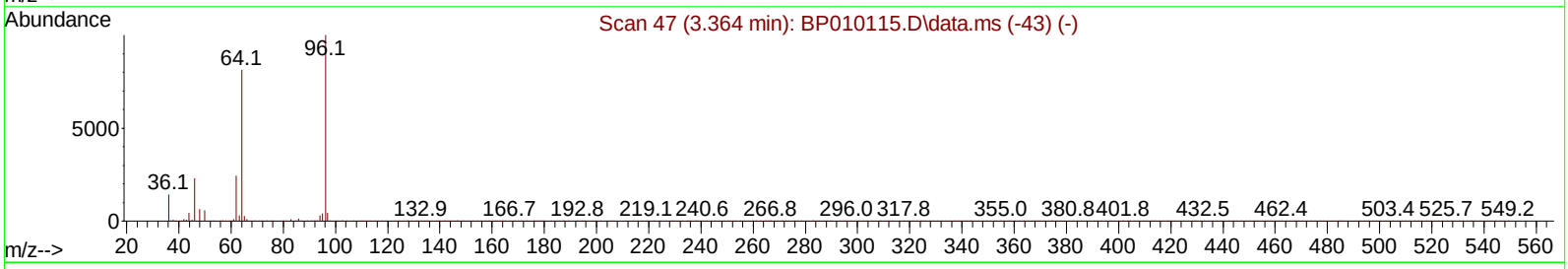

TIC: BP010119.D\data.ms

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

3.370min (+ 0.001) 3.98 ng/uL m

response 15860

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.53#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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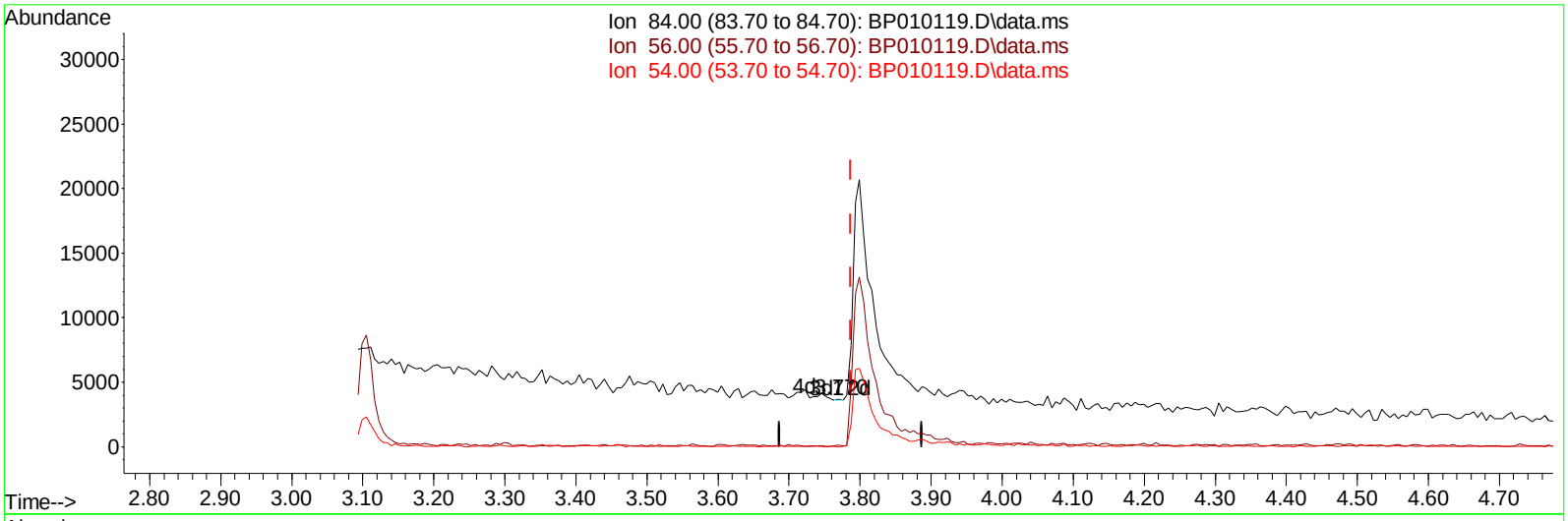

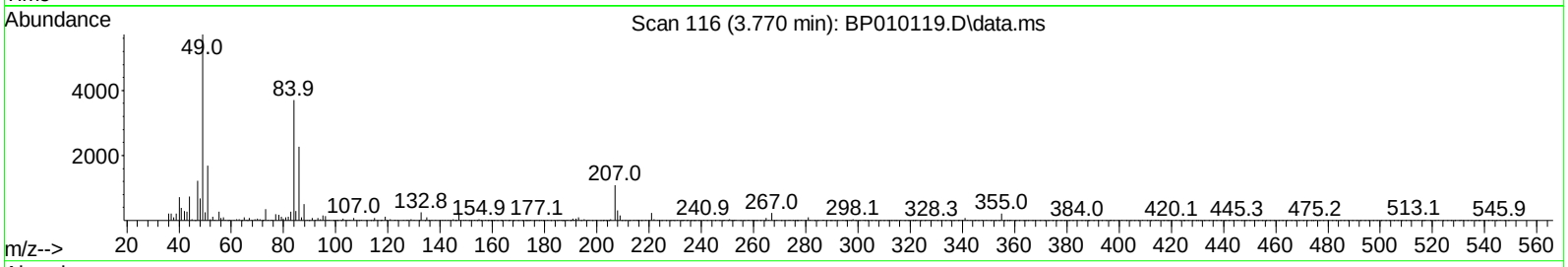

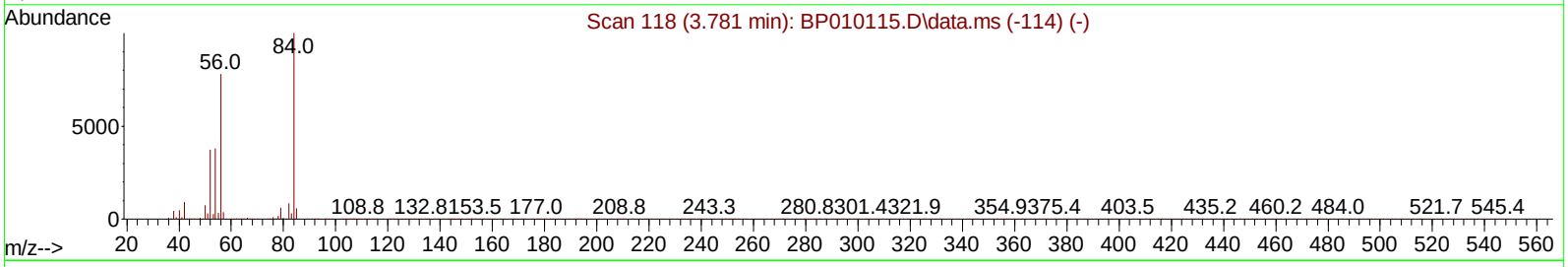

TIC: BP010119.D\data.ms

(4) Pyridine-d5 (S)

3.770min (-0.017) 0.00 ng/ul

response 40

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	2.89#
54.00	24.80	0.00#
0.00	0.00	0.00

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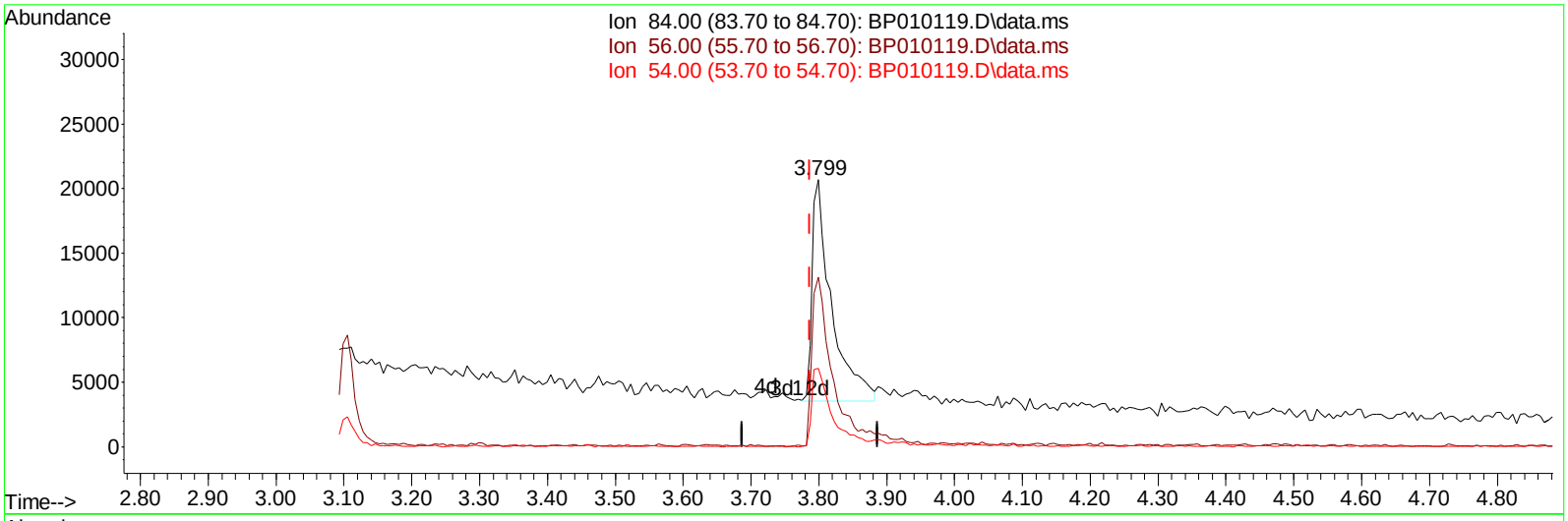

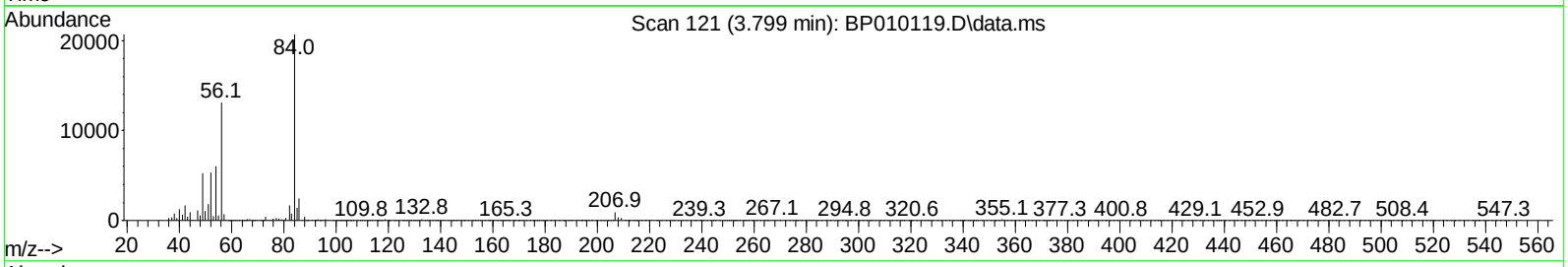

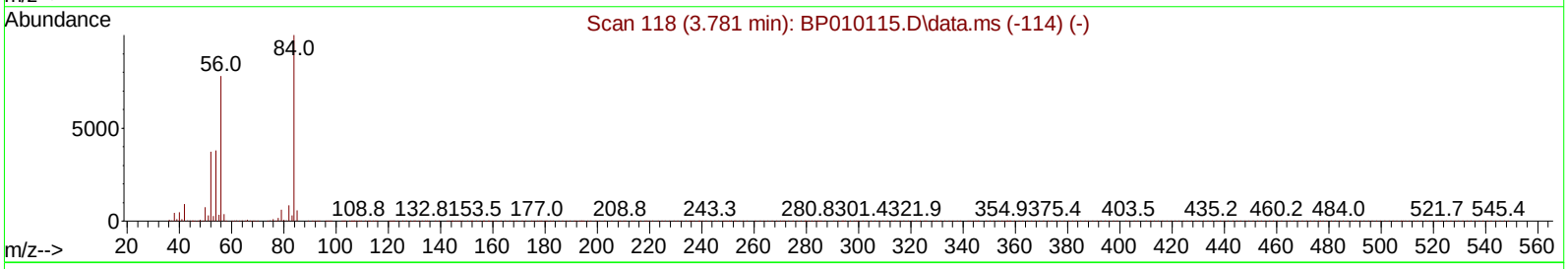

TIC: BP010119.D\data.ms

(4) Pyridine-d5 (S)

3.799min (+ 0.012) 3.02 ng/ul m

response 33852

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	63.47#
54.00	24.80	29.32
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	7.934	152	173749	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	756507	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	500611	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	1019456	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	733444	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	555277	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	15860m	3.980	ng/uL	0.00
4) Pyridine-d5	3.799	84	33852m	3.020	ng/ul	0.01
7) Phenol-d5	7.093	99	305404	19.324	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	211165	22.267	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.464	132	254758	21.766	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	269965	20.617	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	136016	22.606	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	146819	22.343	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	259504	20.286	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	347964	18.970	ng/ul	-0.01
46) Dimethylphthalate-d6	13.975	166	1019011	26.167	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1079391	22.902	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	107126	14.943	ng/ul	0.00
60) Fluorene-d10	15.557	176	852002	25.434	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	95112	14.755	ng/ul	0.00
73) Anthracene-d10	17.416	188	1390437	28.192	ng/ul	0.00
81) Pyrene-d10	19.651	212	1245142	29.851	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	459049	15.769	ng/ul	0.00

Target Compounds Qvalue

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

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