Quantitation Report (Qedit)

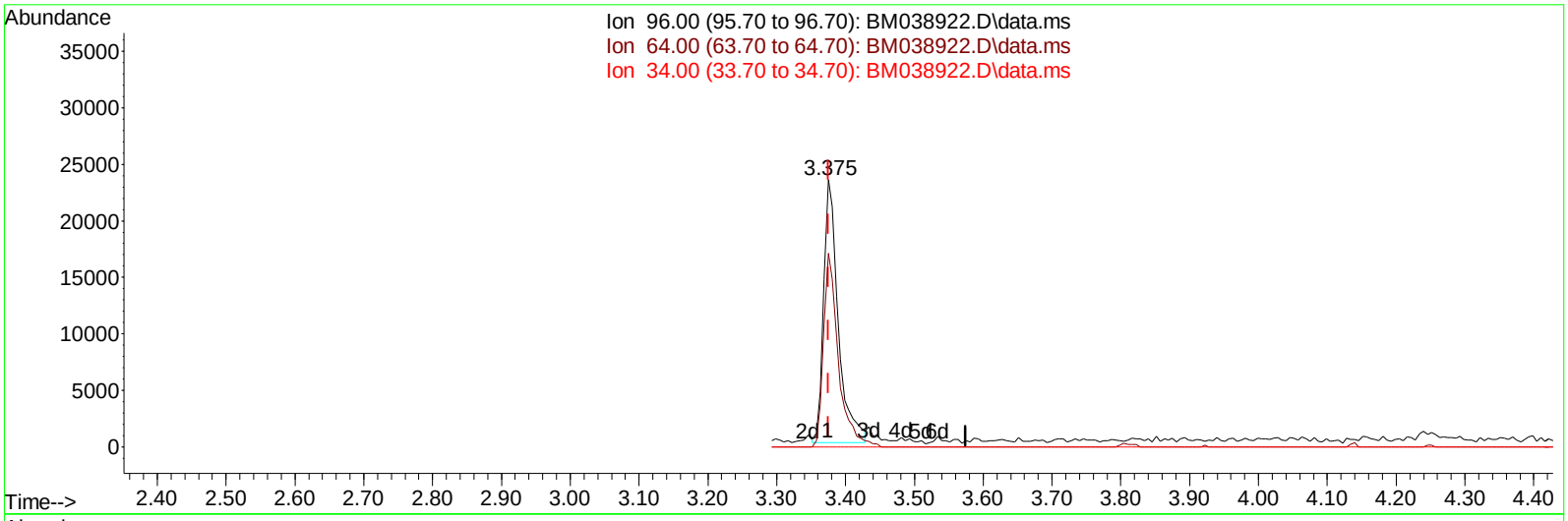
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030723\
Data File : BM038922.D
Acq On : 08 Mar 2023 08:07
Operator : CG/JU
Sample : 01746-04
Misc :
ALS Vial : 35 Sample Multiplier: 1

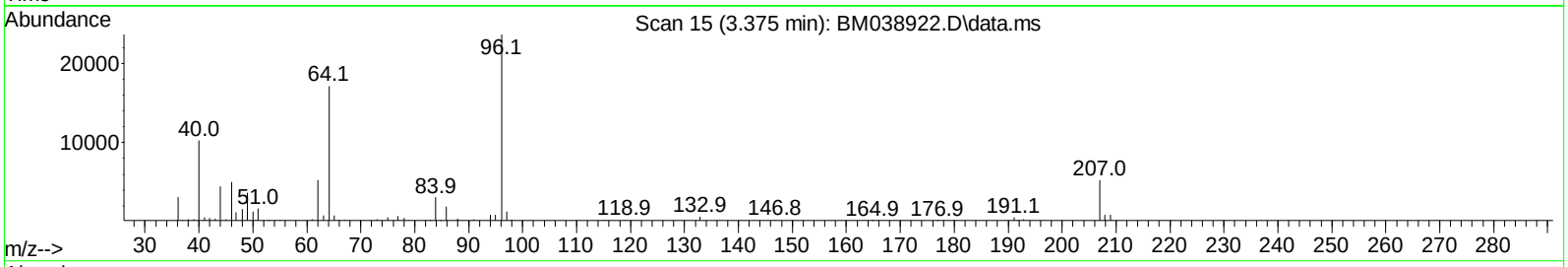
Quant Time: Mar 08 11:56:03 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 08 04:25:35 2023
Response via : Initial Calibration

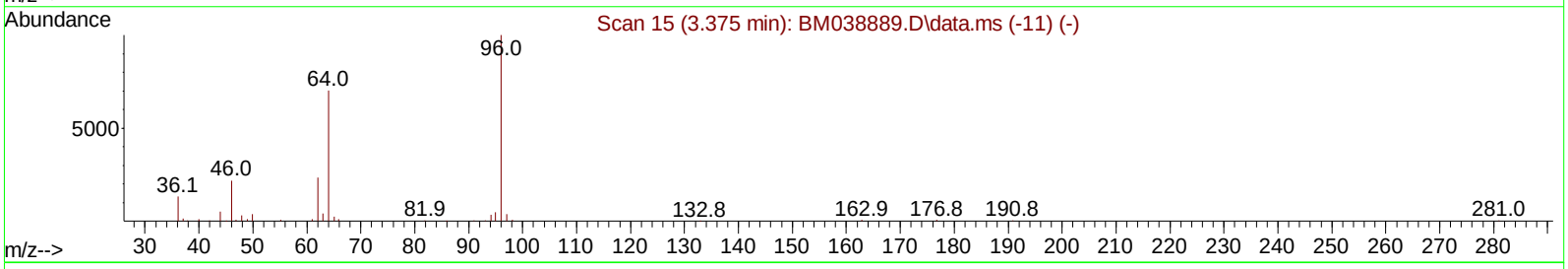
Instrument :
BNA_M
ClientSampleId :
DCAC8

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/08/2023
Supervised By :Jagrut Upadhyay 03/08/2023







TIC: BM038922.D\data.ms

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

3.375min (-0.000) 3.86 ng/uL

response 34205

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.20	72.44
34.00	0.00	0.00
0.00	0.00	0.00

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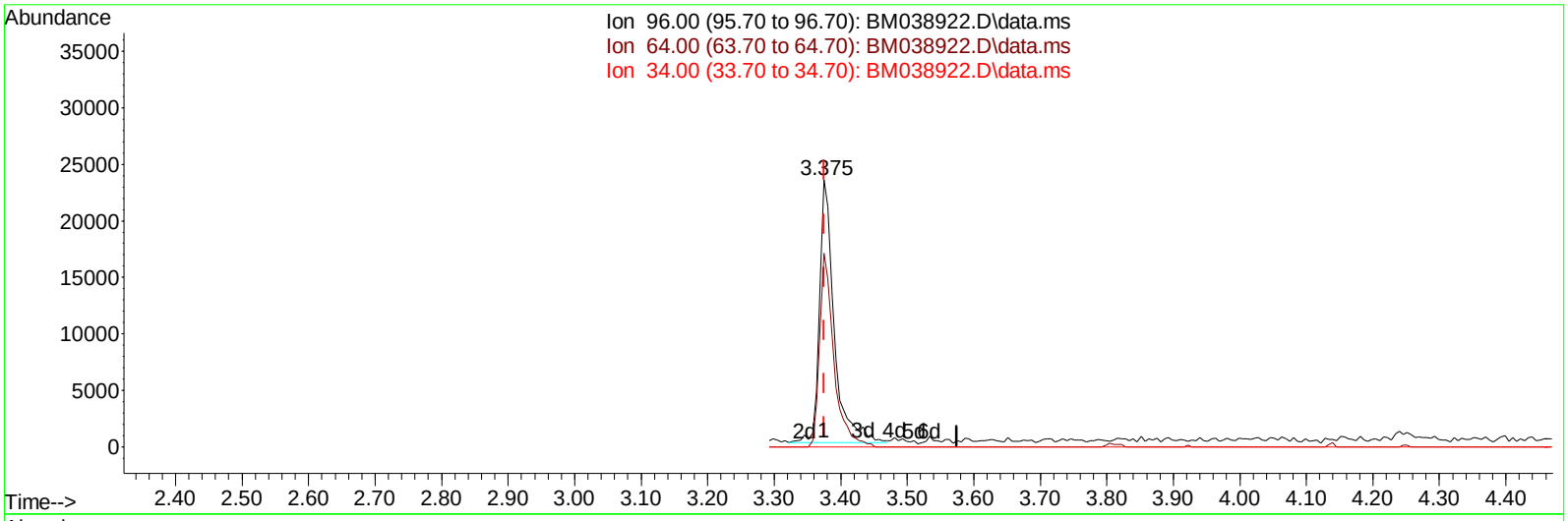

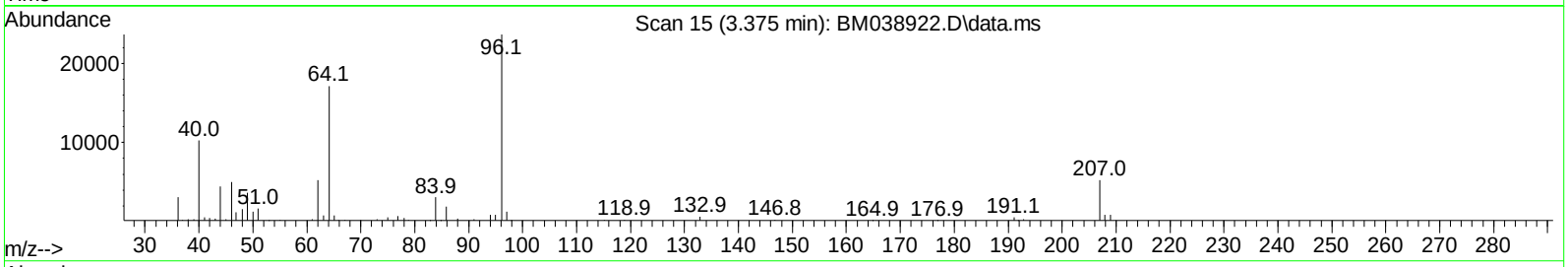

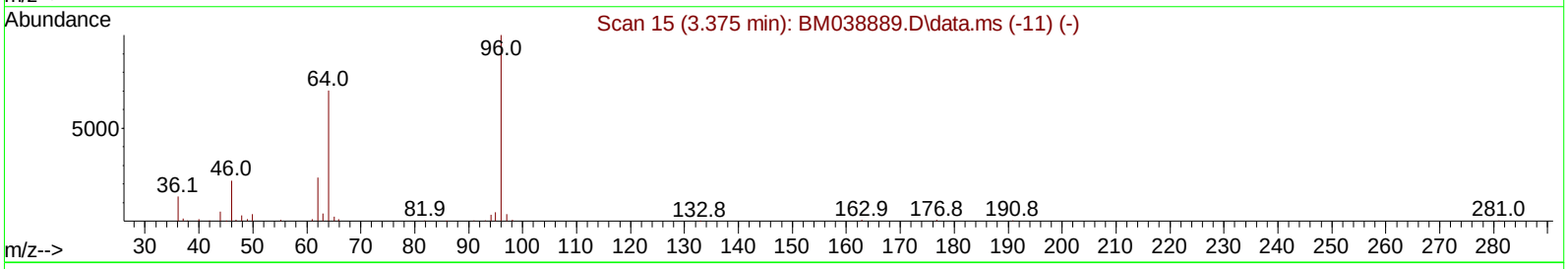

TIC: BM038922.D\data.ms

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

3.375min (-0.000) 4.08 ng/uL m

response 36155

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.20	72.44
34.00	0.00	0.00
0.00	0.00	0.00

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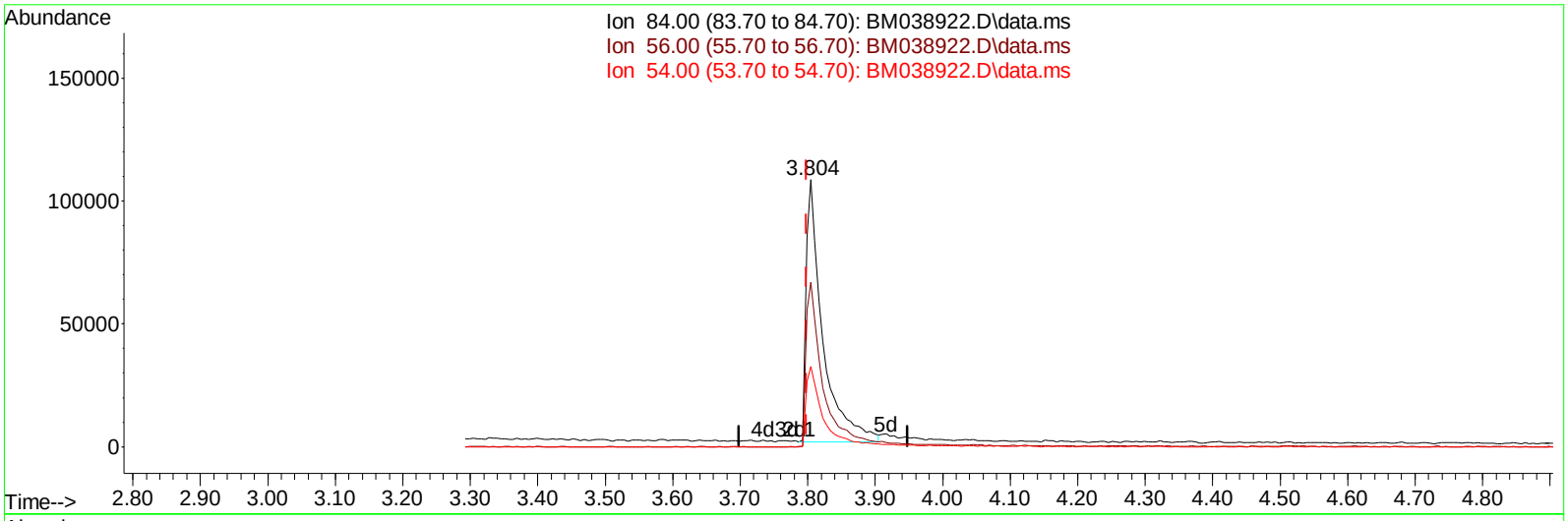

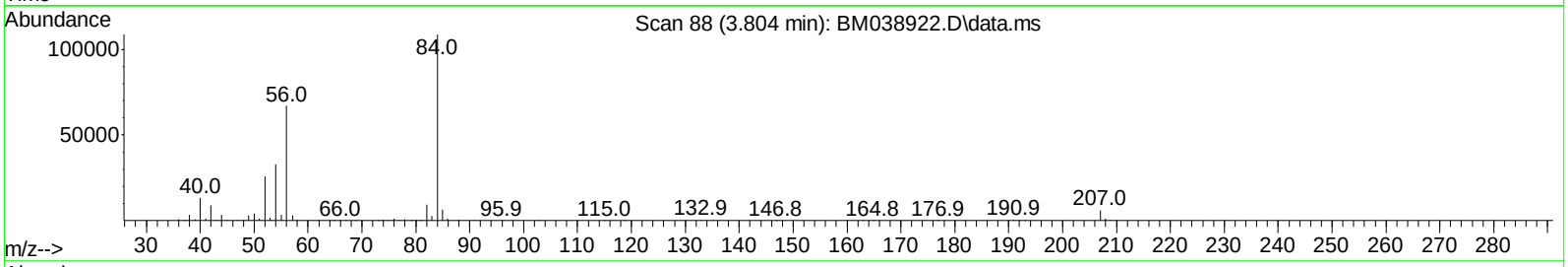

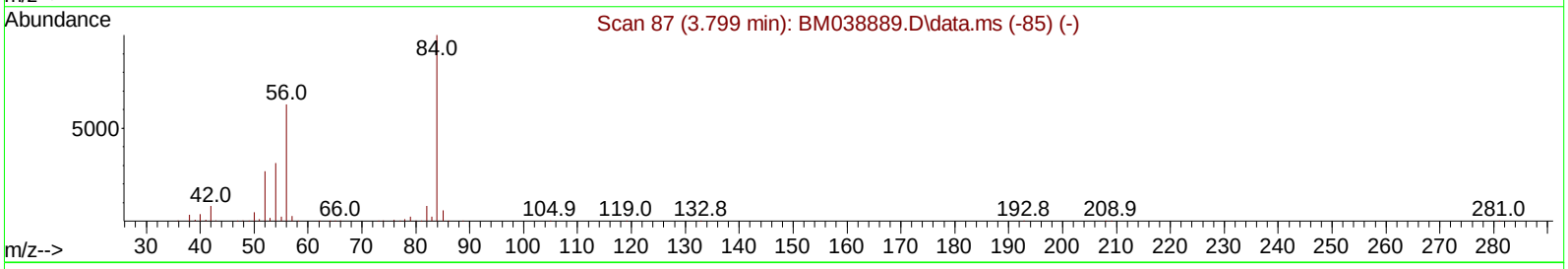

TIC: BM038922.D\data.ms

(4) Pyridine-d5 (S)

3.804min (+ 0.006) 7.43 ng/ul

response 181136

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.20	61.74
54.00	31.50	30.15
0.00	0.00	0.00

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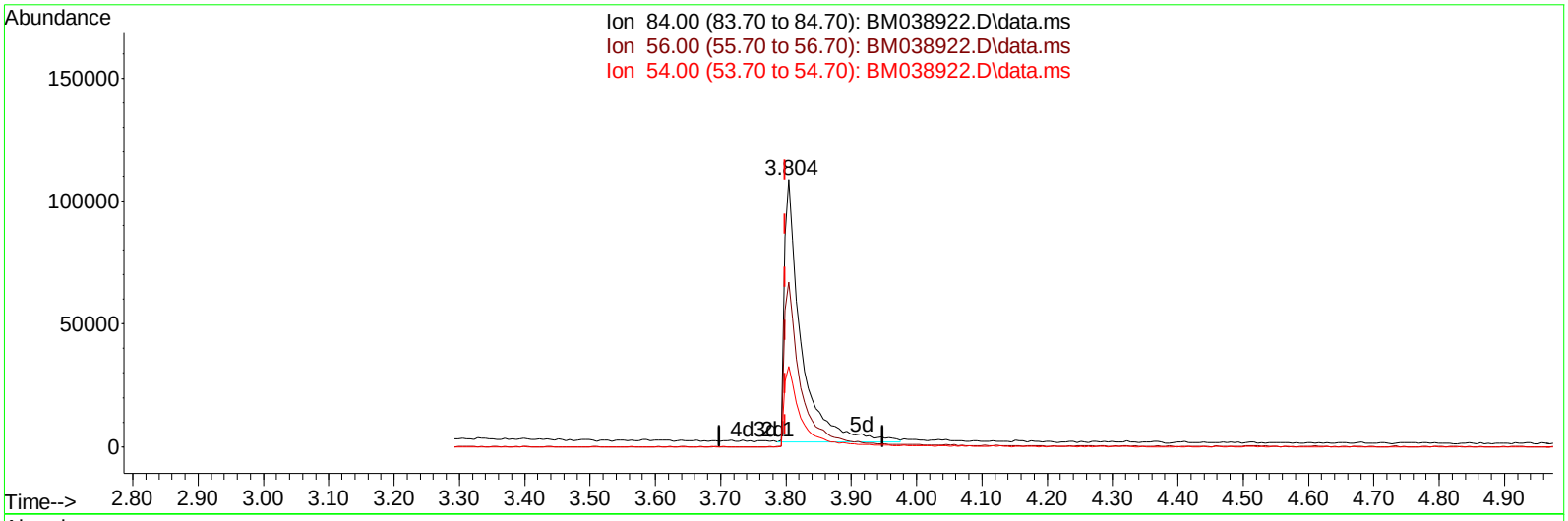

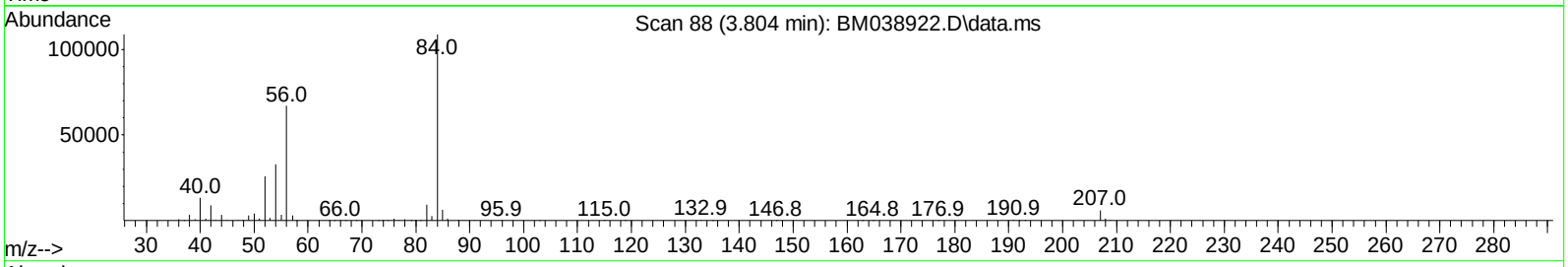

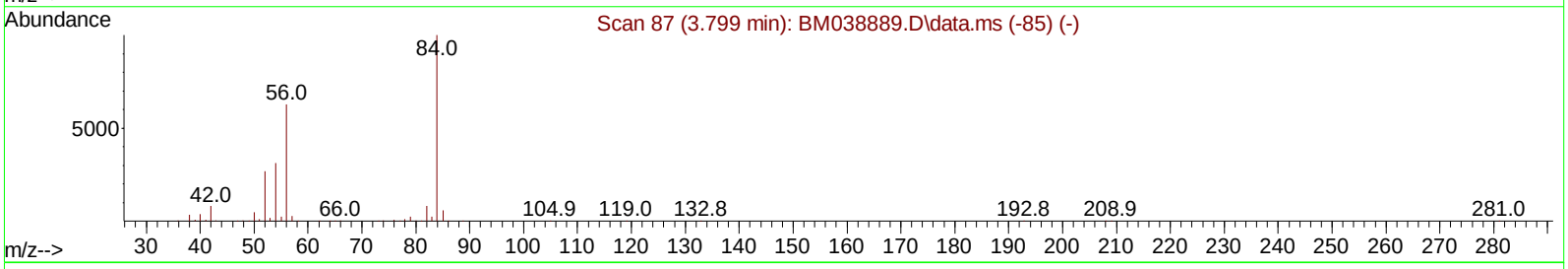

TIC: BM038922.D\data.ms

(4) Pyridine-d5 (S)

3.804min (+ 0.006) 7.77 ng/ul m

response 189286

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.20	61.74
54.00	31.50	30.15
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	330985	20.000	ng/ul	0.00
20) Naphthalene-d8	10.816	136	1494758	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	939937	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	1991767	20.000	ng/ul	0.00
79) Chrysene-d12	21.556	240	1876449	20.000	ng/ul	0.00
88) Perylene-d12	24.009	264	2123687	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	36155m	4.079	ng/uL	0.00
4) Pyridine-d5	3.804	84	189286m	7.767	ng/ul	0.00
7) Phenol-d5	7.151	99	153242	4.938	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.322	67	530636	26.763	ng/ul	0.00
11) 2-Chlorophenol-d4	7.528	132	488746	21.137	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	297071	12.152	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	334276	31.855	ng/ul	0.00
24) 2-Nitrophenol-d4	9.892	143	333153	33.882	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	555554	24.315	ng/ul	0.00
31) 4-Chloroaniline-d4	10.945	131	828469	21.753	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	2072482	28.468	ng/ul	0.00
49) Acenaphthylene-d8	14.327	160	2410570	28.330	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	51645	3.544	ng/ul	0.00
60) Fluorene-d10	15.627	176	1887894	29.404	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	266979	25.783	ng/ul	0.00
73) Anthracene-d10	17.480	188	3184076	32.420	ng/ul	0.00
81) Pyrene-d10	19.768	212	3687559	35.730	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.850	264	3654048	34.143	ng/ul	0.00
Target Compounds						Qvalue

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

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