Quantitation Report (Qedit)

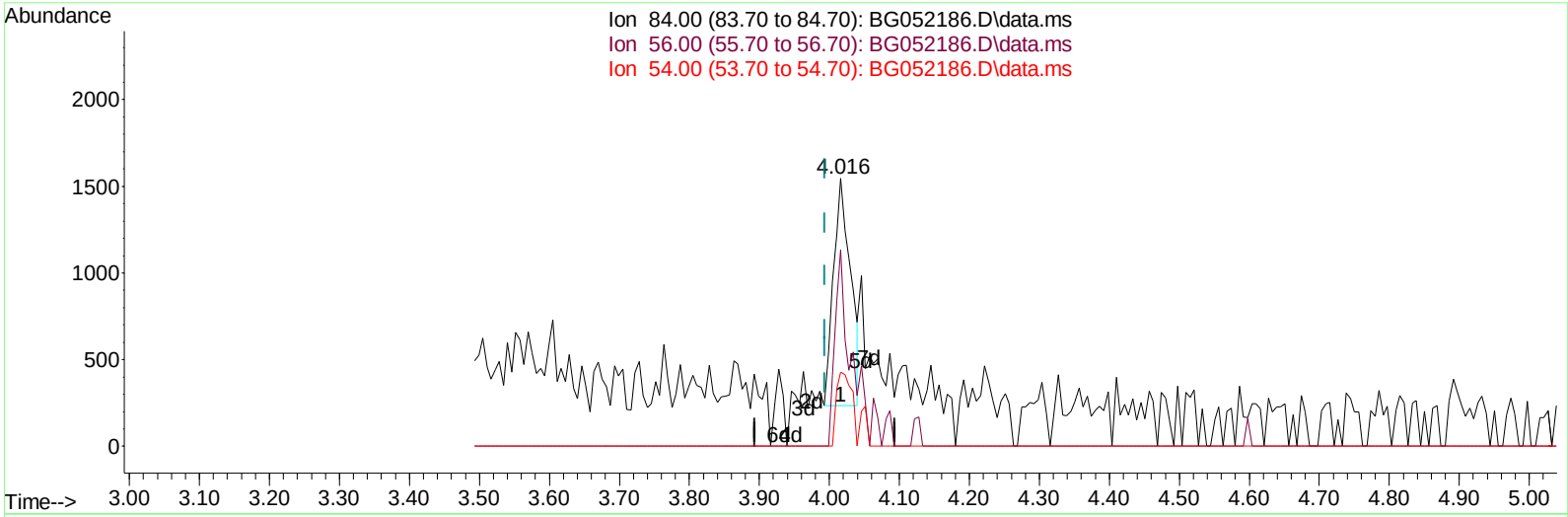
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052186.D
Acq On : 27 Jan 2022 15:27
Operator : CG/JU
Sample : N1279-06
Misc :
ALS Vial : 44 Sample Multiplier: 1

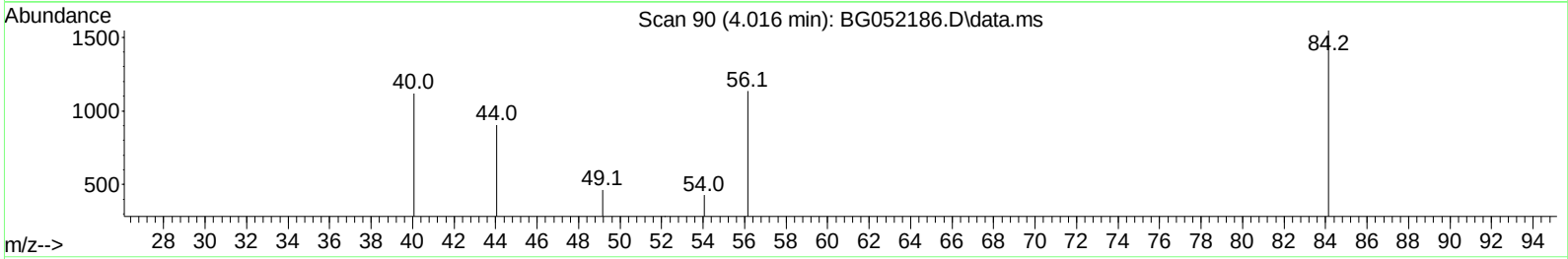
Quant Time: Jan 27 23:59:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 26 23:58:52 2022
Response via : Initial Calibration

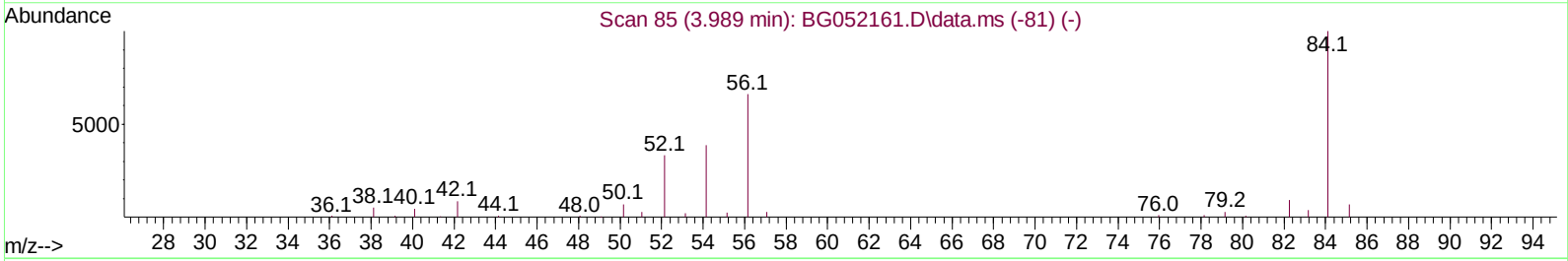
Instrument :
BNA_G
ClientSampleId :
C0VD6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/28/2022
Supervised By :Yogesh Patel 01/28/2022







TIC: BG052186.D\data.ms

(4) Pyridine-d5 (S)

4.016min (+ 0.022) 1.16 ng/ul

response 2236

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	73.33
54.00	31.50	27.70
0.00	0.00	0.00

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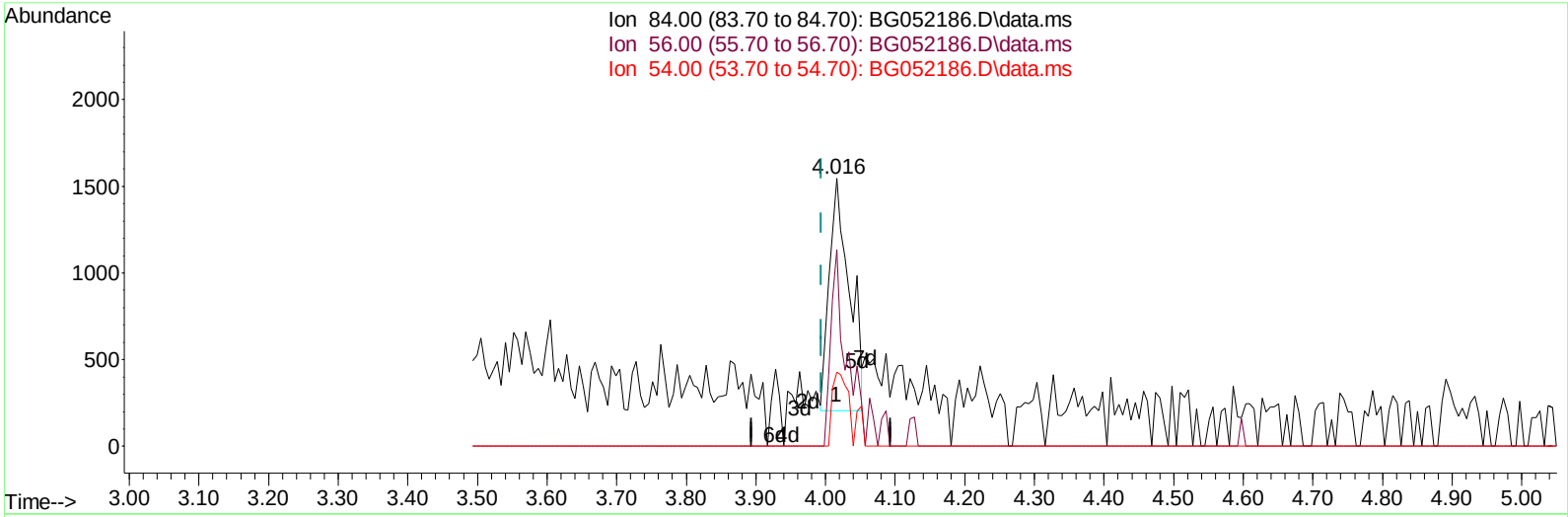

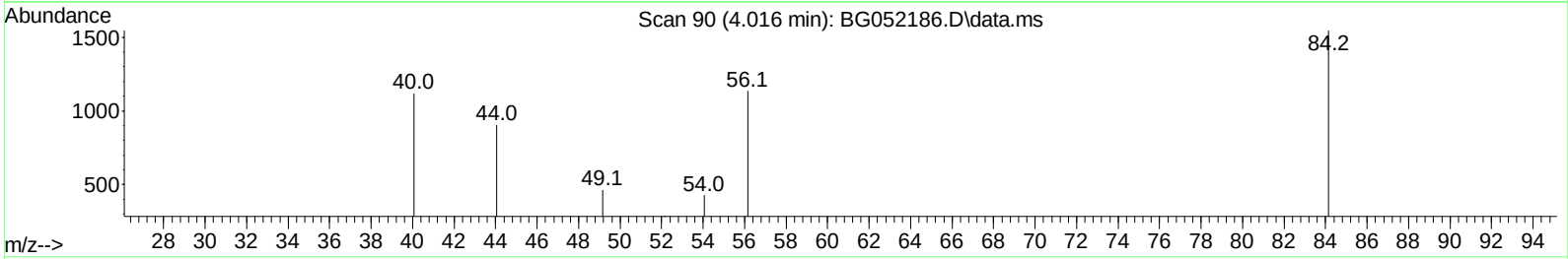

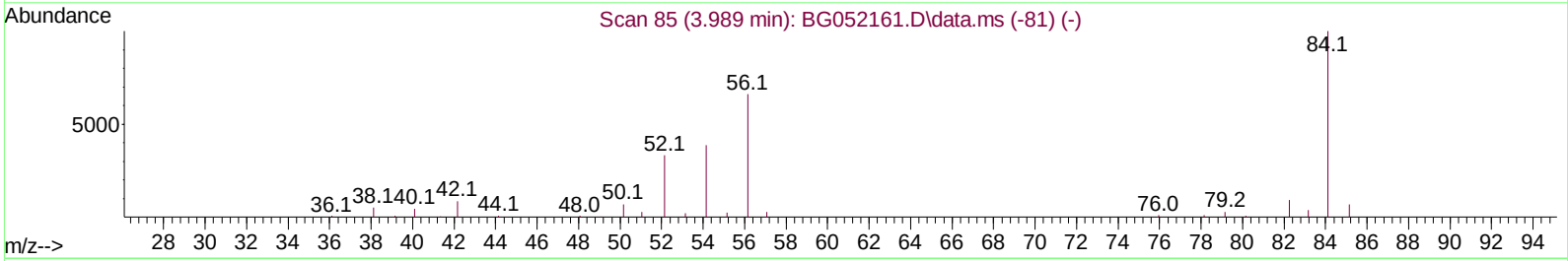

TIC: BG052186.D\data.ms

(4) Pyridine-d5 (S)

4.016min (+ 0.022) 1.39 ng/ul m

response 2672

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	73.33
54.00	31.50	27.70
0.00	0.00	0.00

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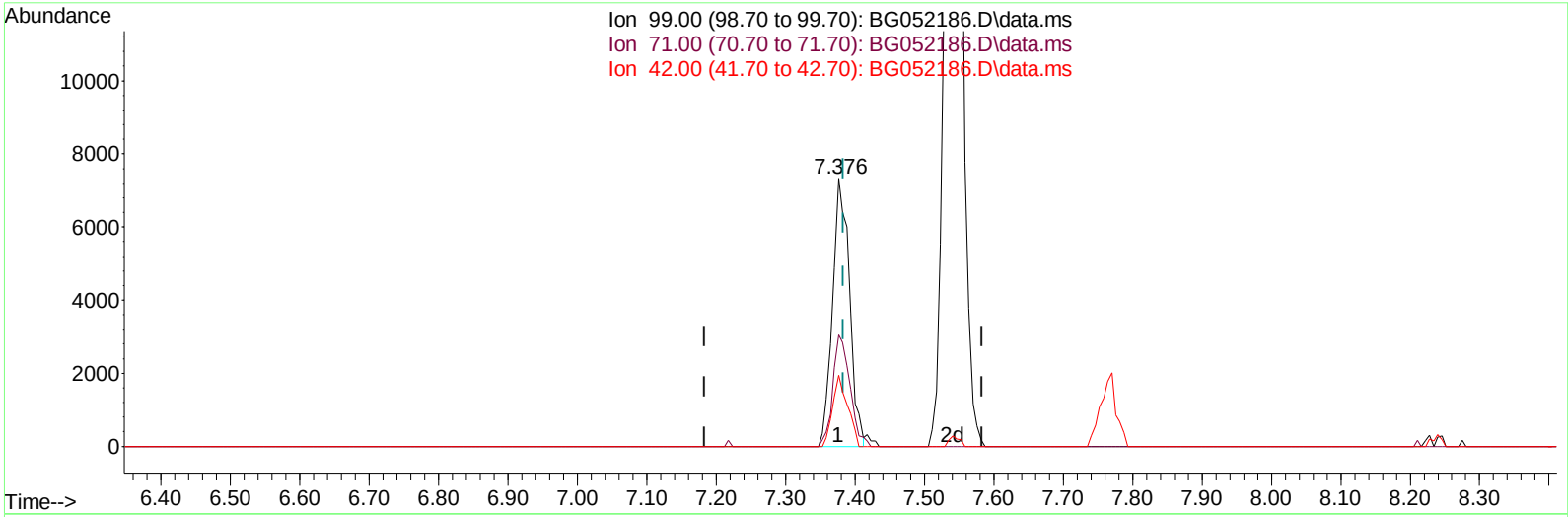

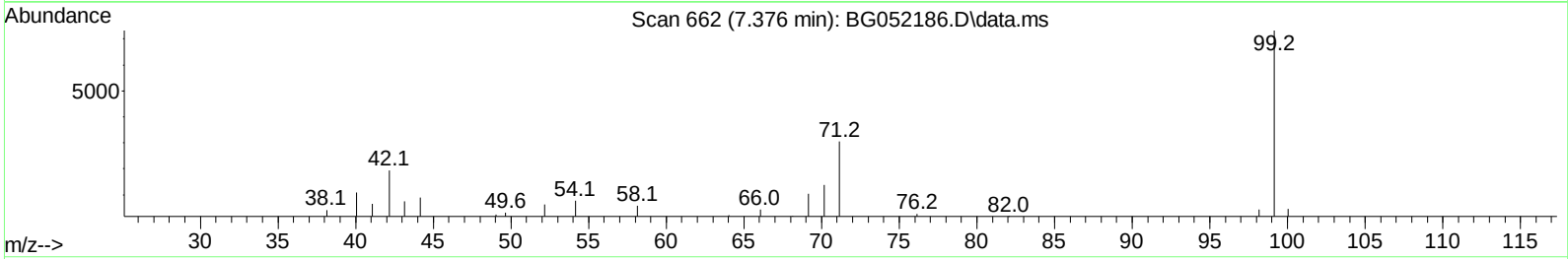

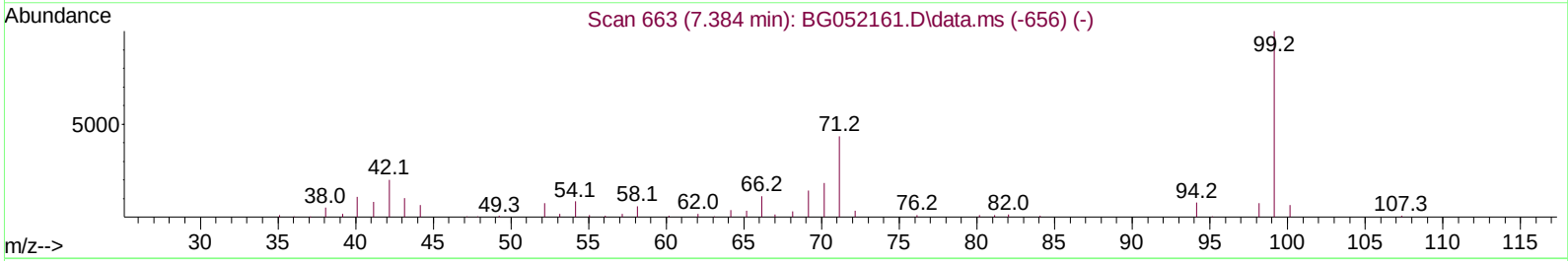

TIC: BG052186.D\data.ms

(7) Phenol-d5 (S)

7.376min (-0.007) 5.72 ng/ul

response 12346

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	38.50	41.74
42.00	22.60	26.44
0.00	0.00	0.00

Quantitation Report (Qedit)

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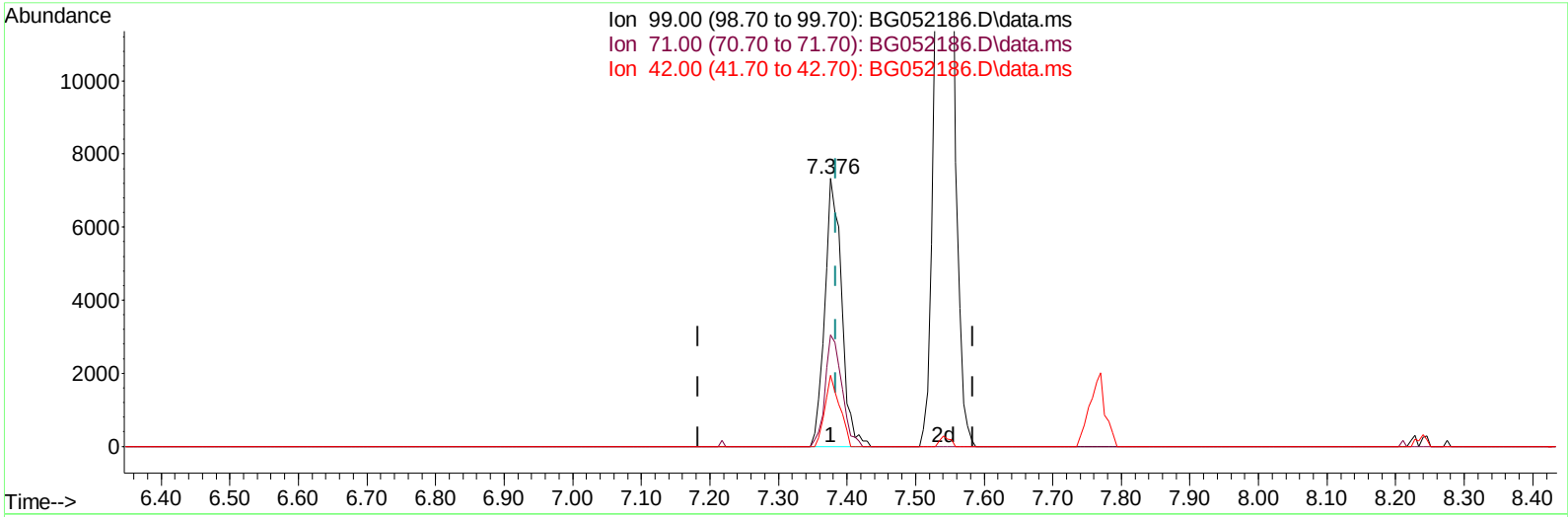

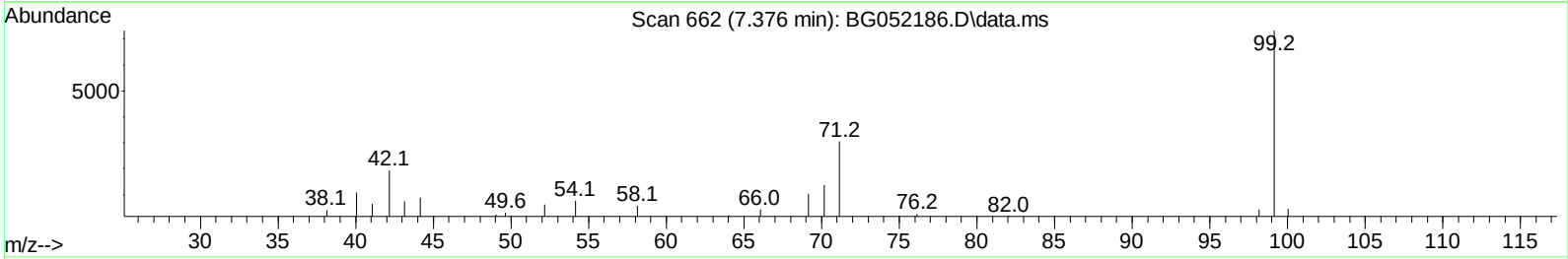

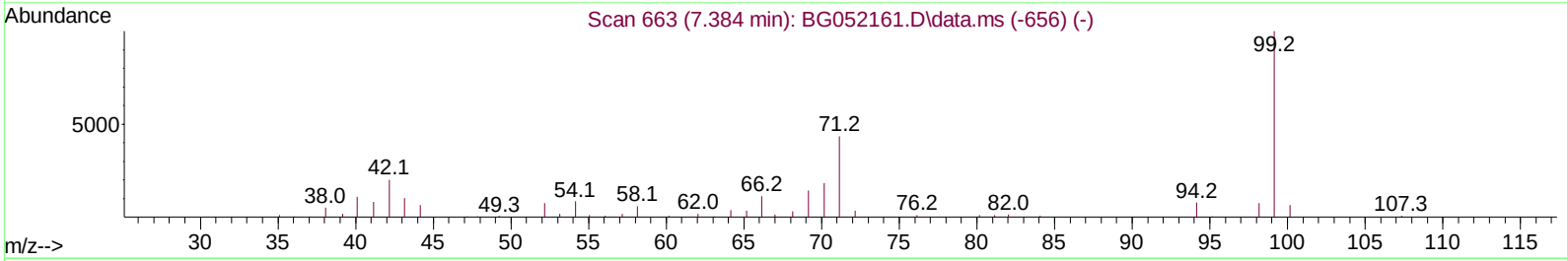

TIC: BG052186.D\data.ms

(7) Phenol-d5 (S)

7.376min (-0.007) 5.82 ng/ul m

response 12568

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	38.50	41.74
42.00	22.60	26.44
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	8.234	152	22072	20.000	ng/ul		0.00
20) Naphthalene-d8	11.077	136	92467	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.884	164	69039	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.627	188	174153	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.945	240	176060	20.000	ng/ul	#	0.00
88) Perylene-d12	25.422	264	177052	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.570	96	2733	4.425	ng/uL		0.00
4) Pyridine-d5	4.016	84	2672m	1.392	ng/ul		0.02
7) Phenol-d5	7.376	99	12568m	5.818	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	41344	30.148	ng/ul		0.00
11) 2-Chlorophenol-d4	7.764	132	35375	24.733	ng/ul		0.00
15) 4-Methylphenol-d8	8.945	113	24243	14.436	ng/ul		0.00
21) Nitrobenzene-d5	9.409	128	25930	35.489	ng/ul		0.00
24) 2-Nitrophenol-d4	10.137	143	26790	33.153	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.689	165	46631	29.523	ng/ul		0.00
31) 4-Chloroaniline-d4	11.200	131	41130	18.698	ng/ul		0.00
46) Dimethylphthalate-d6	14.267	166	195714	34.296	ng/ul		0.00
49) Acenaphthylene-d8	14.572	160	235395	34.149	ng/ul		0.00
54) 4-Nitrophenol-d4	15.066	143	4507	4.724	ng/ul		0.00
60) Fluorene-d10	15.871	176	173706	34.912	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.982	200	26363	23.996	ng/ul		0.00
73) Anthracene-d10	17.727	188	326213	39.486	ng/ul		0.00
81) Pyrene-d10	20.006	212	402217	39.695	ng/ul		0.00
92) Benzo(a)pyrene-d12	25.187	264	383569	41.152	ng/ul		0.00
Target Compounds							Qvalue

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

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