Quantitation Report (Qedit)

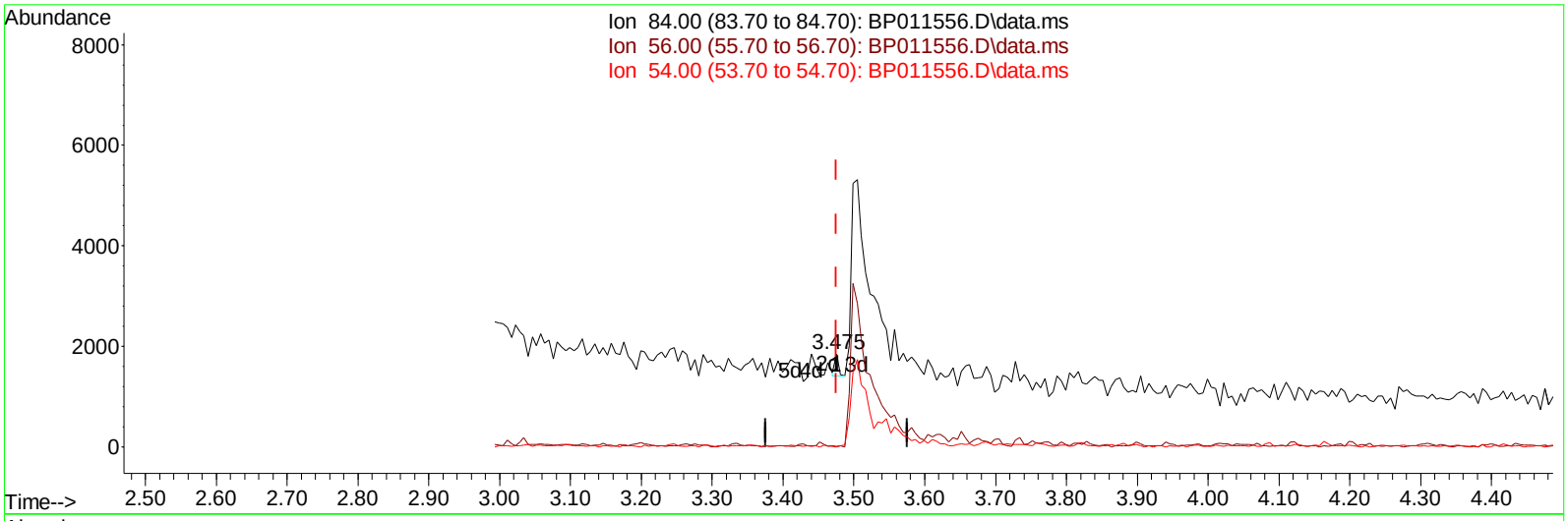
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082422\
Data File : BP011556.D
Acq On : 25 Aug 2022 09:14
Operator : CG/JU
Sample : N4268-01DL 5X
Misc :
ALS Vial : 12 Sample Multiplier: 1

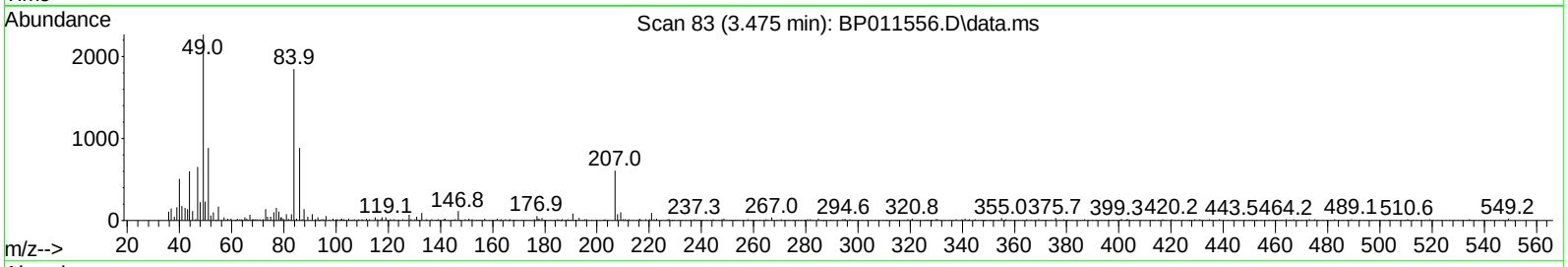
Quant Time: Aug 26 01:01:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 25 02:14:54 2022
Response via : Initial Calibration

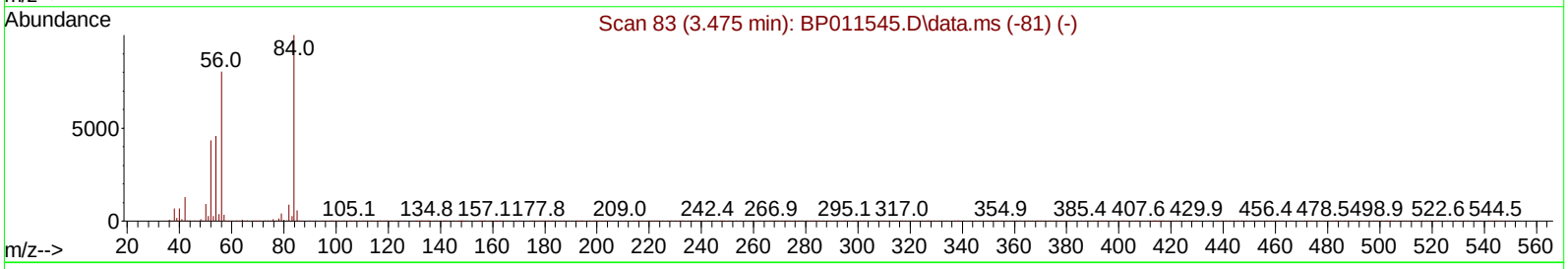
Instrument :
BNA_P
ClientSampleId :
C0AB7DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/26/2022
Supervised By :mohammad ahmed 08/27/2022







TIC: BP011556.D\data.ms

(4) Pyridine-d5 (S)

3.475min (-0.000) 0.04 ng/ul

response 168

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.90	0.59#
54.00	42.10	0.00#
0.00	0.00	0.00

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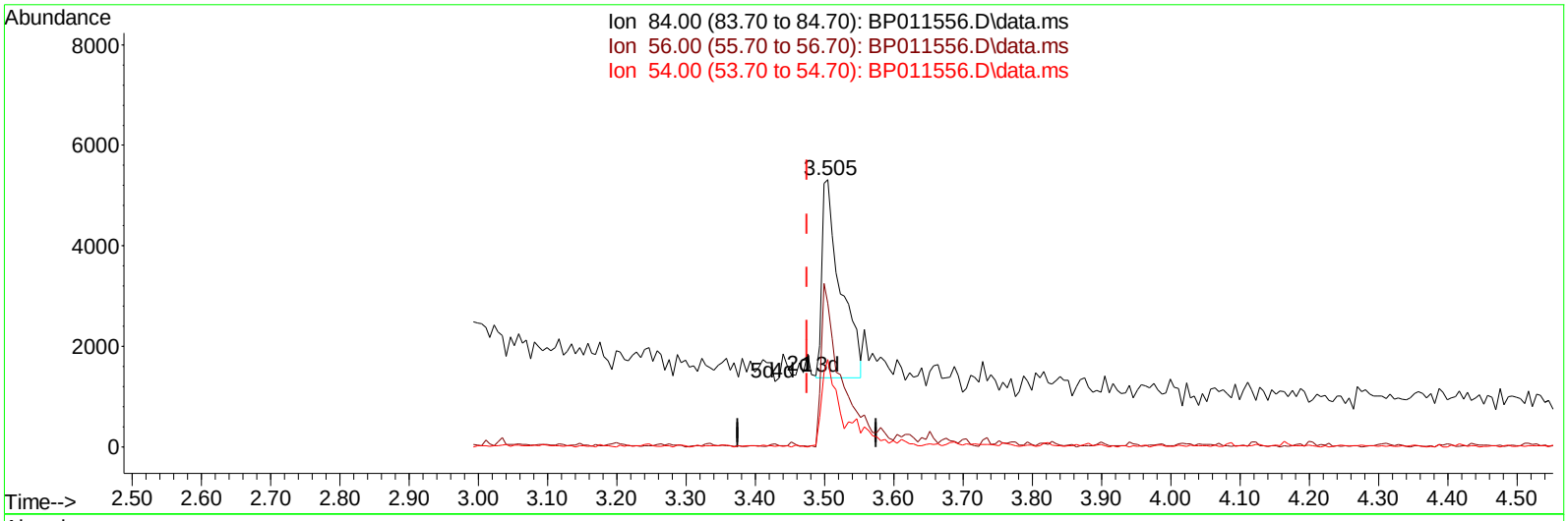

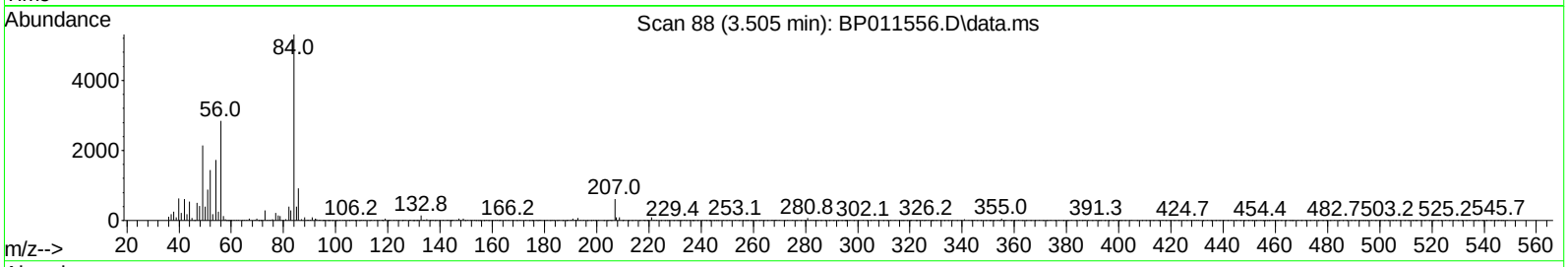

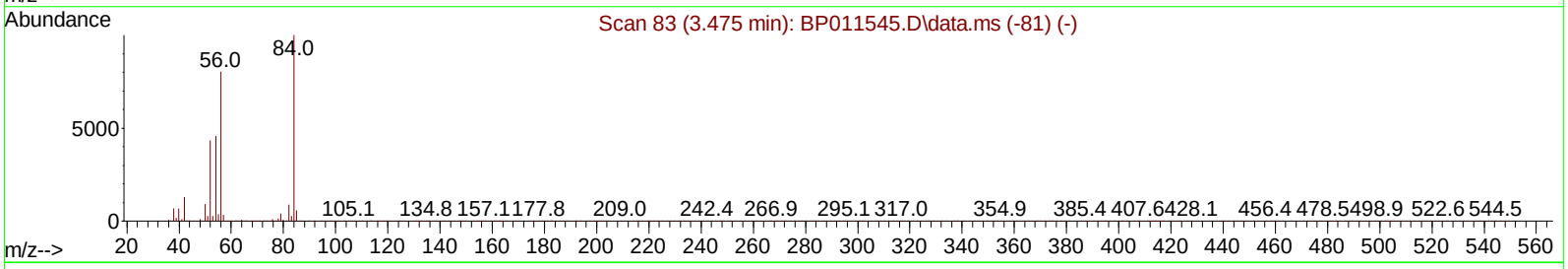

TIC: BP011556.D\data.ms

(4) Pyridine-d5 (S)

3.505min (+ 0.029) 1.60 ng/ul m

response 7272

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.90	53.54#
54.00	42.10	32.57#
0.00	0.00	0.00

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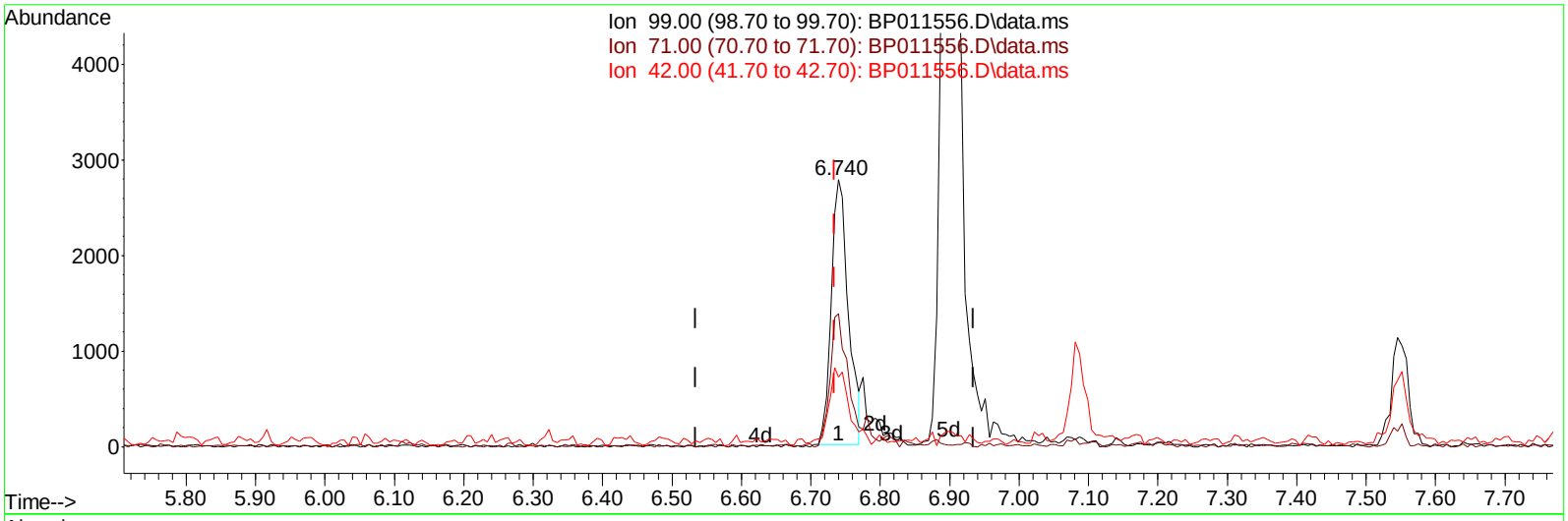

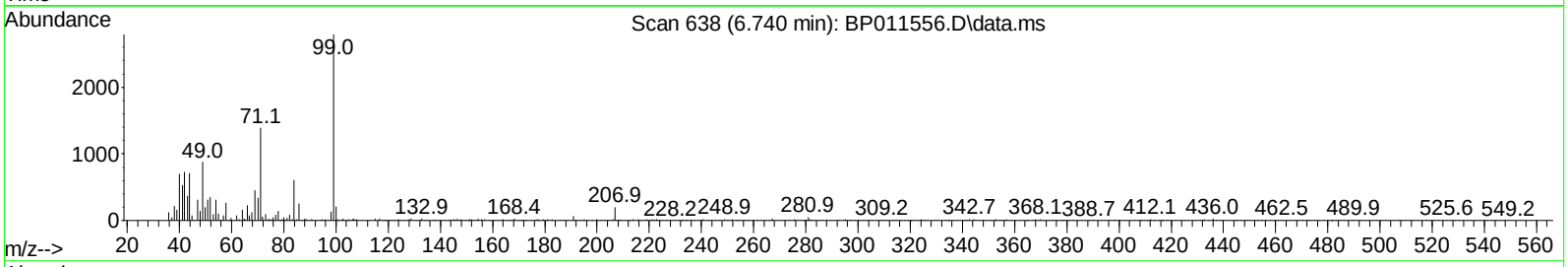

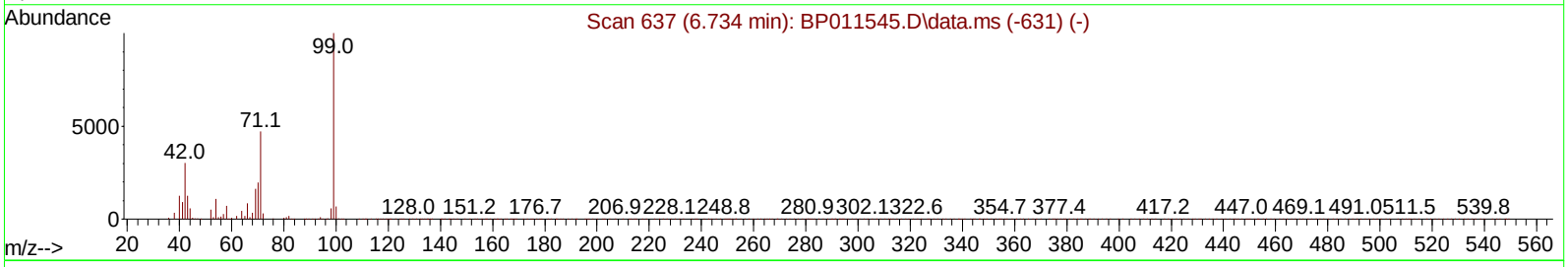

TIC: BP011556.D\data.ms

(7) Phenol-d5 (S)

6.740min (+ 0.006) 0.95 ng/ul

response 4805

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.20	49.79
42.00	25.30	25.97
0.00	0.00	0.00

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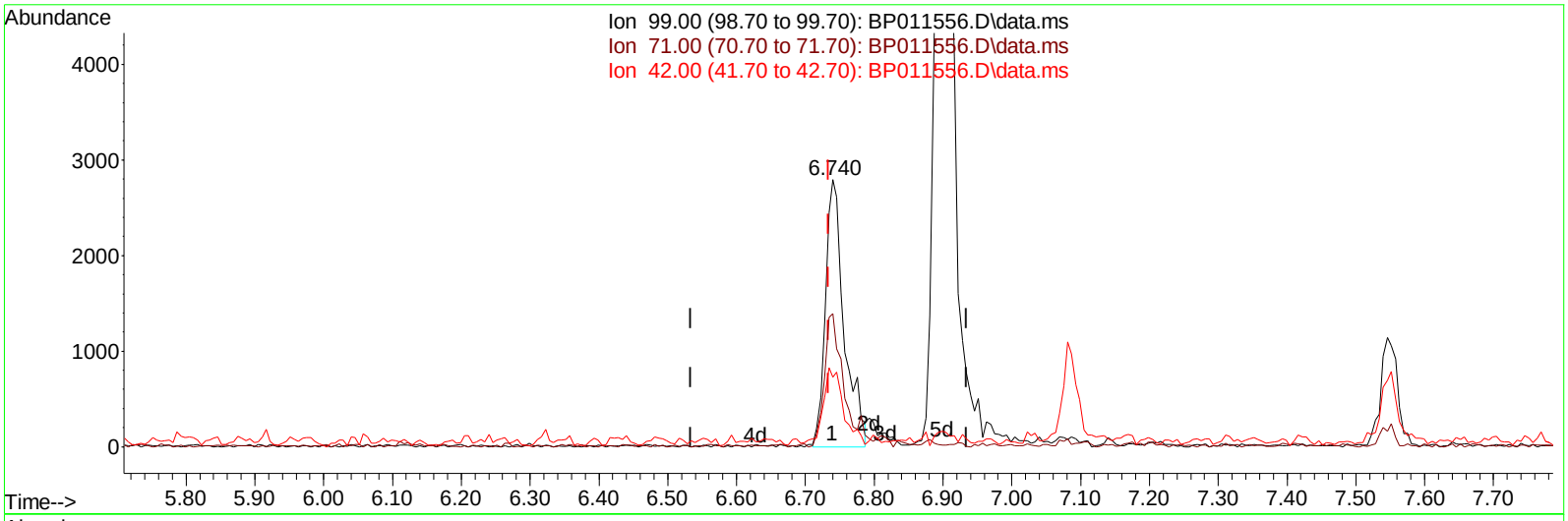

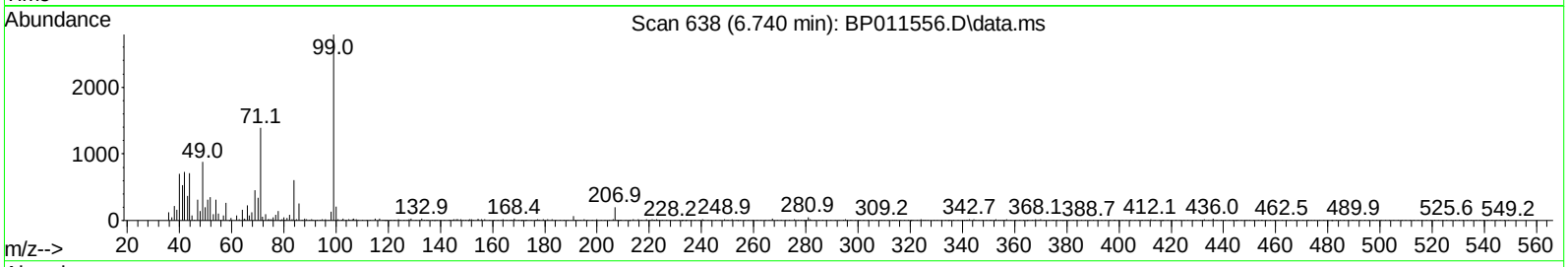

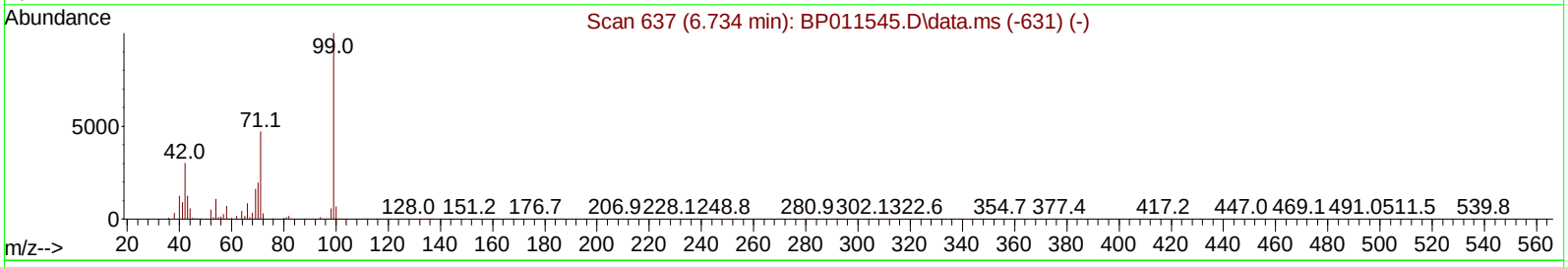

TIC: BP011556.D\data.ms

(7) Phenol-d5 (S)

6.740min (+ 0.006) 1.05 ng/ul m

response 5326

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.20	49.79
42.00	25.30	25.97
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	7.552	152	55191	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	238271	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.222	164	140217	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.992	188	293339	20.000	ng/ul	0.00
79) Chrysene-d12	21.127	240	299643	20.000	ng/ul	0.00
88) Perylene-d12	23.415	264	302527	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.505	84	7272m	1.597	ng/ul	0.03
7) Phenol-d5	6.740	99	5326m	1.050	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.899	67	18512	5.618	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	15483	4.151	ng/ul	0.00
15) 4-Methylphenol-d8	8.275	113	9835	2.423	ng/ul	0.00
21) Nitrobenzene-d5	8.716	128	9464	5.517	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	8521	4.881	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	16310	5.061	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	17958	3.348	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	76435	7.707	ng/ul	0.00
49) Acenaphthylene-d8	13.910	160	74531	6.762	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.228	176	63969	7.667	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	6466	4.079	ng/ul	0.00
73) Anthracene-d10	17.092	188	101849	7.840	ng/ul	0.00
81) Pyrene-d10	19.357	212	121924	8.042	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.263	264	112437	7.640	ng/ul	0.00
Target Compounds						Qvalue
52) Acenaphthene	14.287	153	407393	46.449	ng/ul	99
61) Fluorene	15.281	166	138198	13.967	ng/ul	98
74) Anthracene	17.128	178	21645	1.351	ng/ul	96
80) Fluoranthene	19.033	202	78490	4.210	ng/ul	98
82) Pyrene	19.386	202	43076	2.199	ng/ul	99

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

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