Quantitation Report (Qedit)

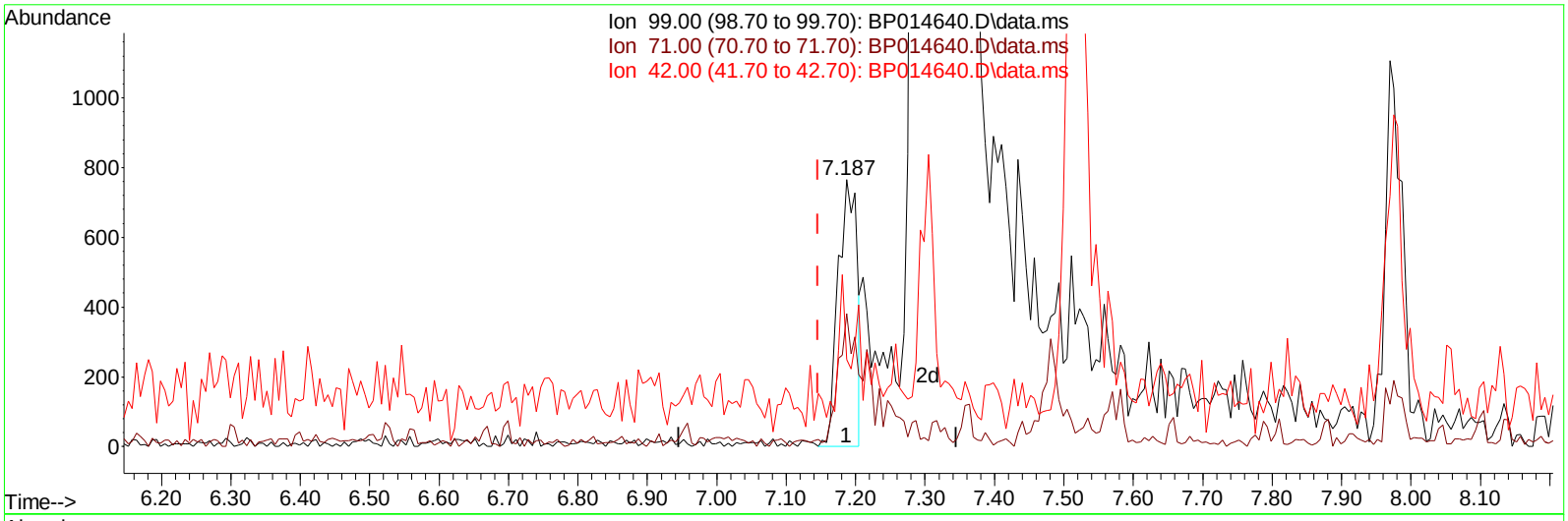
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014640.D
Acq On : 10 Apr 2023 22:46
Operator : CG/JU
Sample : 02194-13
Misc :
ALS Vial : 16 Sample Multiplier: 1

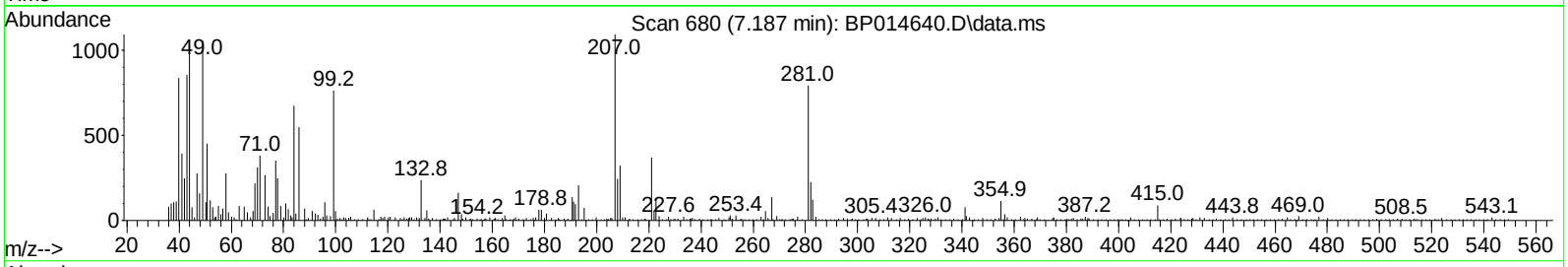
Quant Time: Apr 11 01:25:12 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration

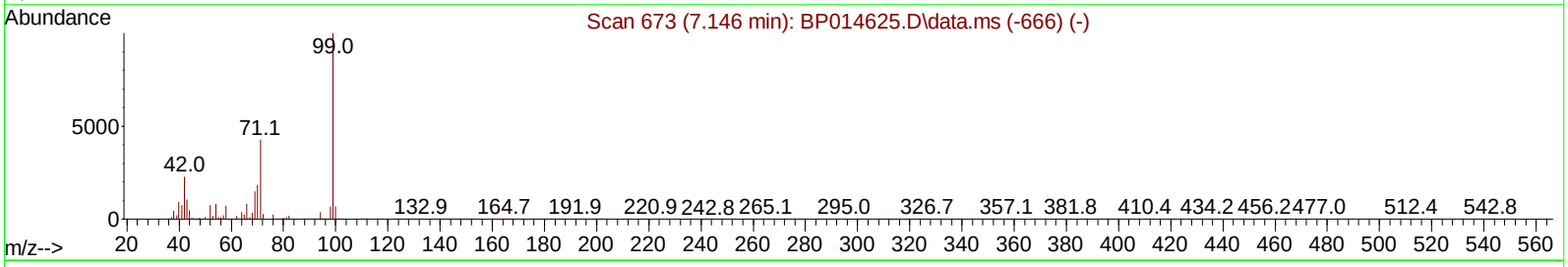
Instrument :
BNA_P
ClientSampleId :
BHDX0

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/11/2023
Supervised By :Jagrut Upadhyay 04/11/2023







TIC: BP014640.D\data.ms

(7) Phenol-d5 (S)

7.187min (+ 0.041) 0.19 ng/ul

response 1457

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	49.93#
42.00	18.90	32.55#
0.00	0.00	0.00

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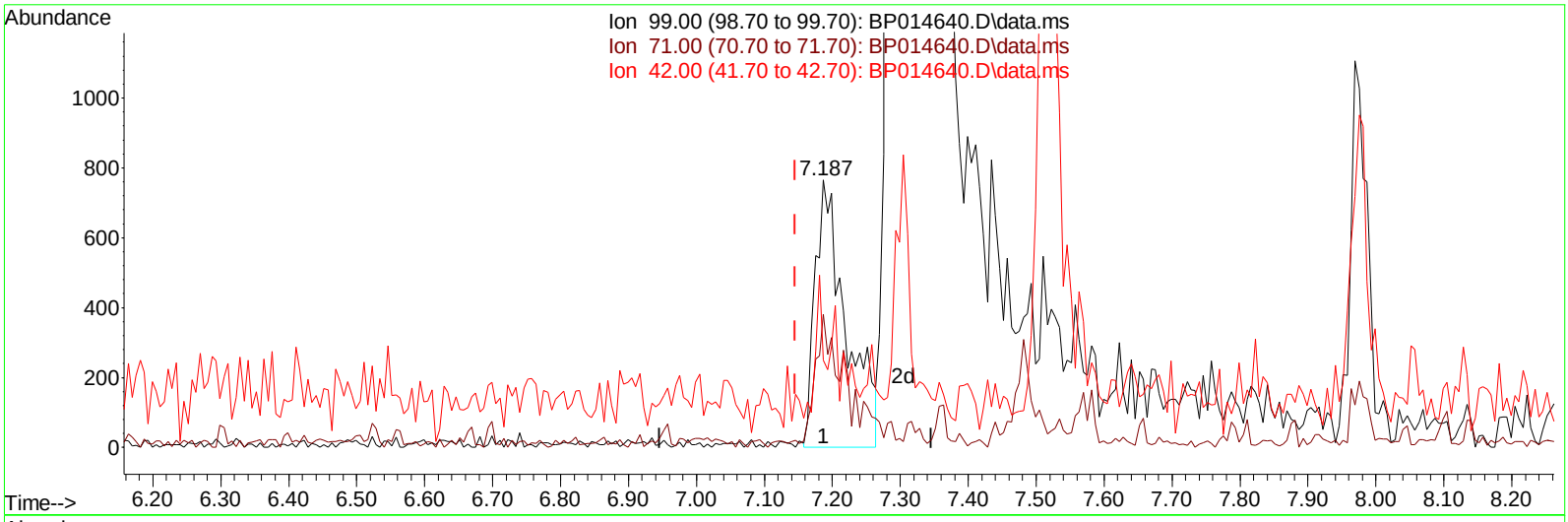

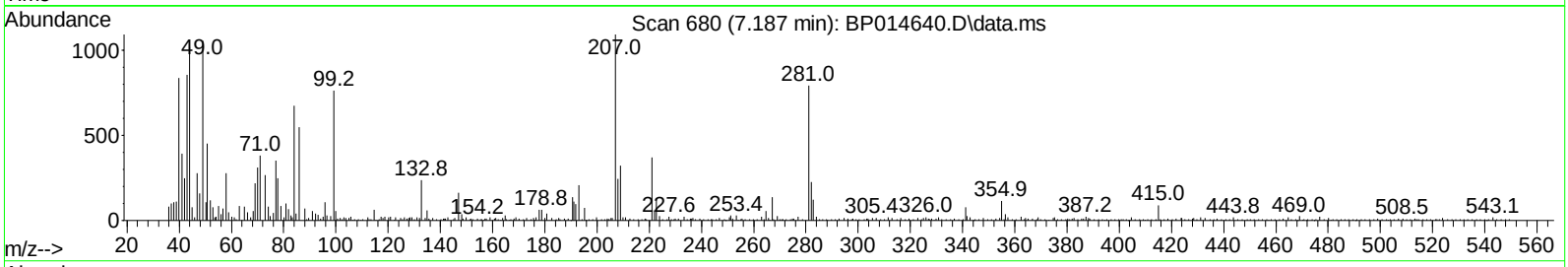

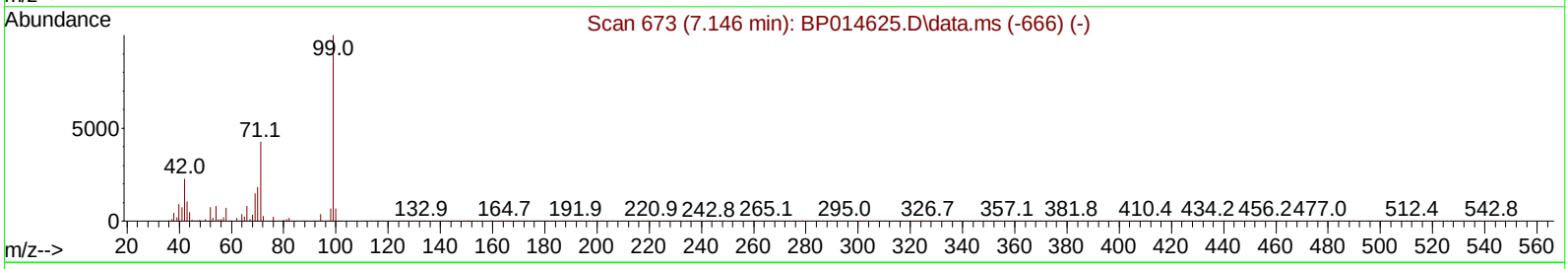

TIC: BP014640.D\data.ms

(7) Phenol-d5 (S)

7.187min (+ 0.041) 0.32 ng/ul m

response 2418

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	40.30	49.93#
42.00	18.90	32.55#
0.00	0.00	0.00

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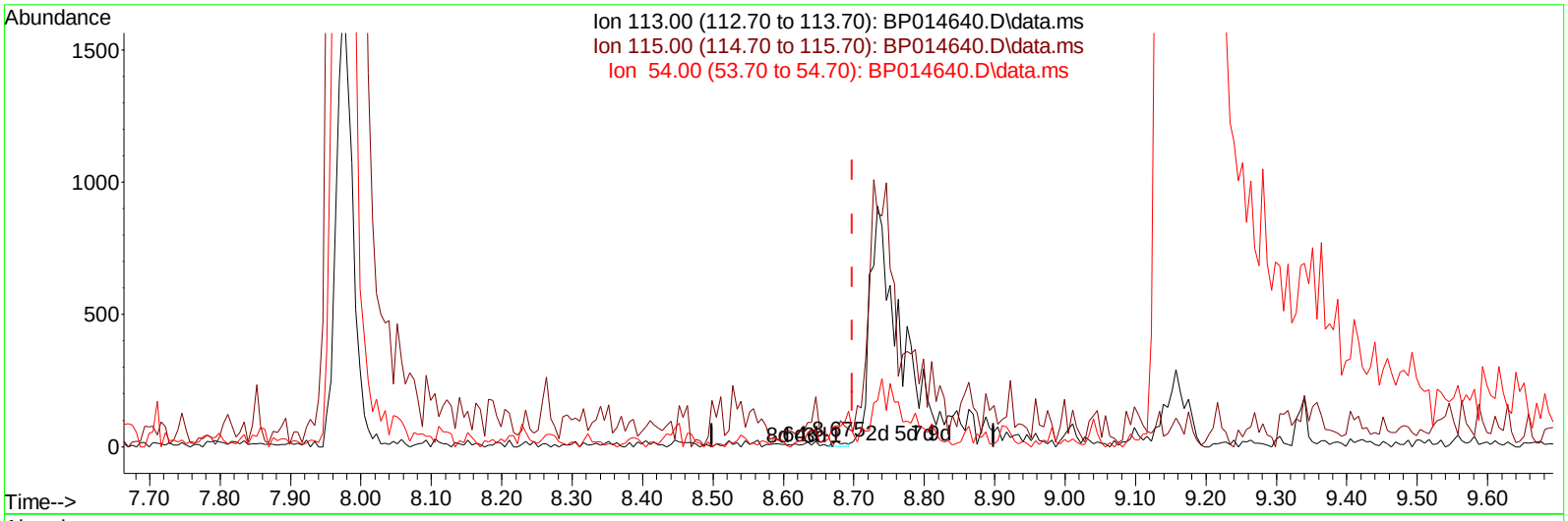

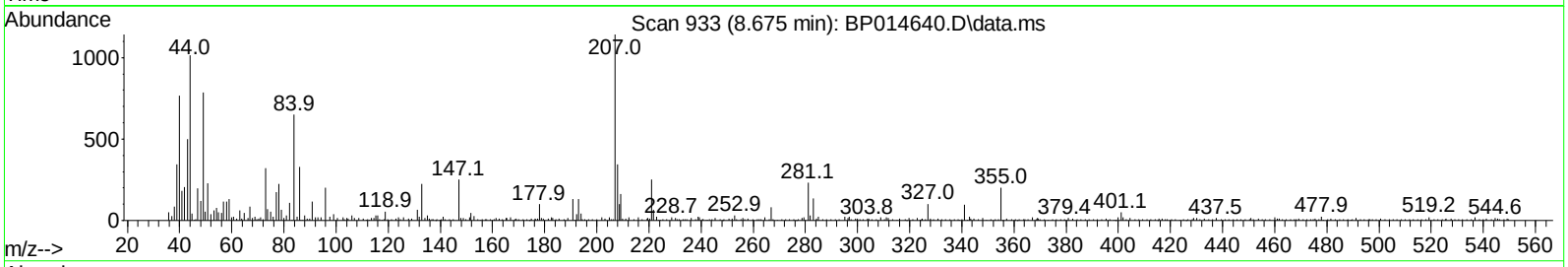

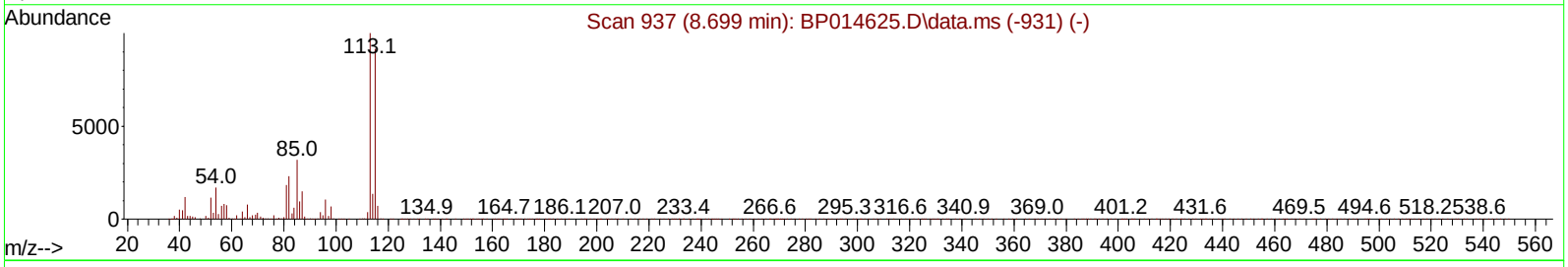

TIC: BP014640.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.675min (-0.023) 0.00 ng/ul

response 23

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	89.90	160.00#
54.00	15.10	395.00#
0.00	0.00	0.00

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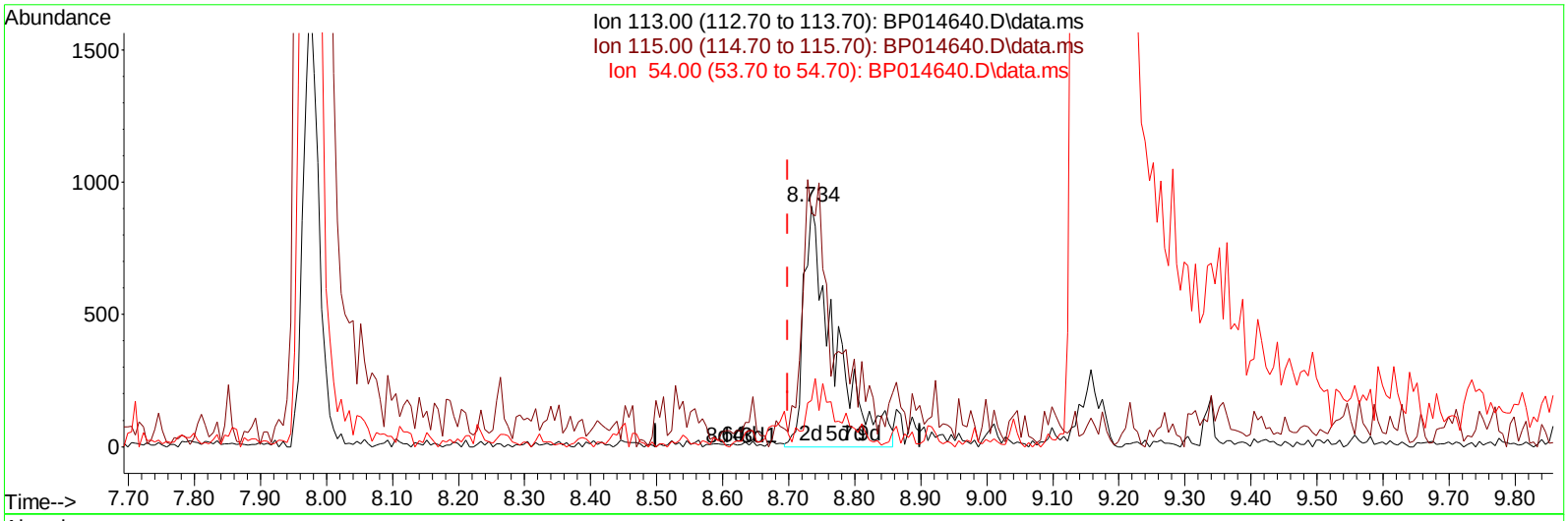

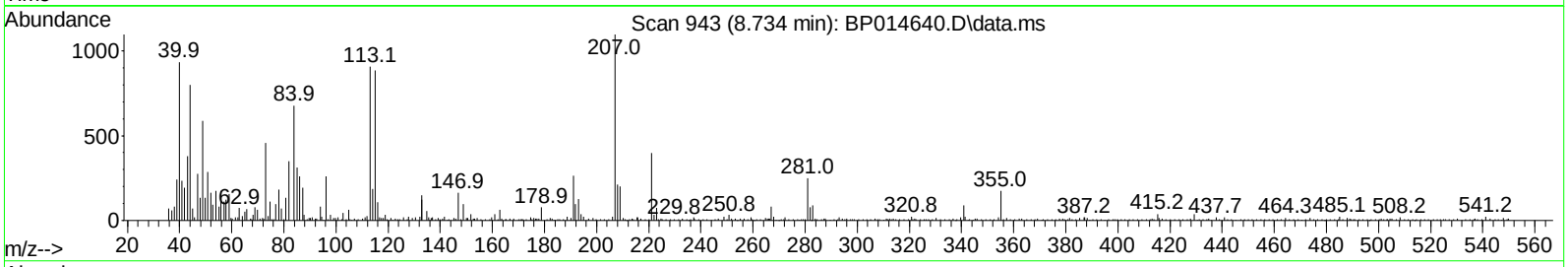

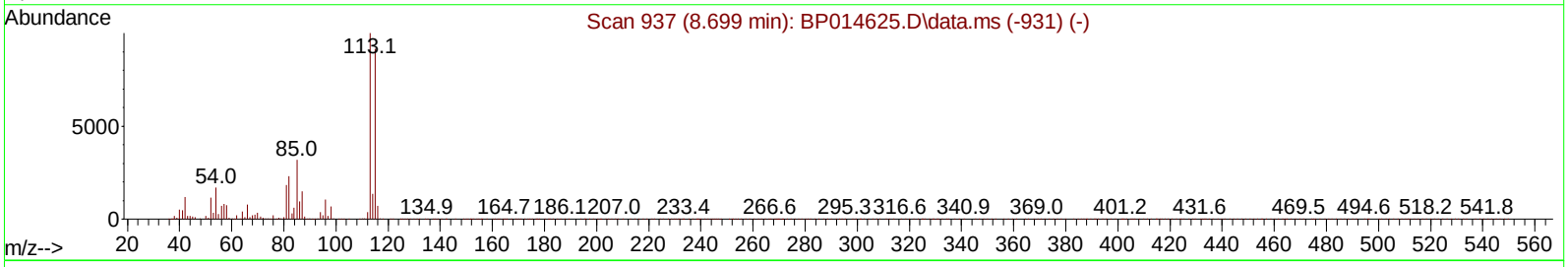

TIC: BP014640.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.734min (+ 0.035) 0.48 ng/ul m

response 2943

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	89.90	97.25
54.00	15.10	19.25#
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	7.975	152	82142	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	333751	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	208439	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	447216	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	499181	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	597999	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	11806	5.574	ng/uL	0.00
4) Pyridine-d5	3.811	84	44479	8.194	ng/ul	0.02
7) Phenol-d5	7.187	99	2418m	0.319	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.305	67	156780	32.256	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	35972	6.570	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	2943m	0.482	ng/ul	0.04
21) Nitrobenzene-d5	9.158	128	78218	29.031	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	17997	5.833	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	51275	9.126	ng/ul	0.02
31) 4-Chloroaniline-d4	10.952	131	141437	19.111	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	408988	24.200	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	601366	32.068	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.622	176	448568	31.827	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.492	188	804373	37.165	ng/ul	0.00
81) Pyrene-d10	19.722	212	1057223	31.531	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	1215835	38.268	ng/ul	0.00
Target Compounds						Qvalue

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

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