Quantitation Report (Qedit)

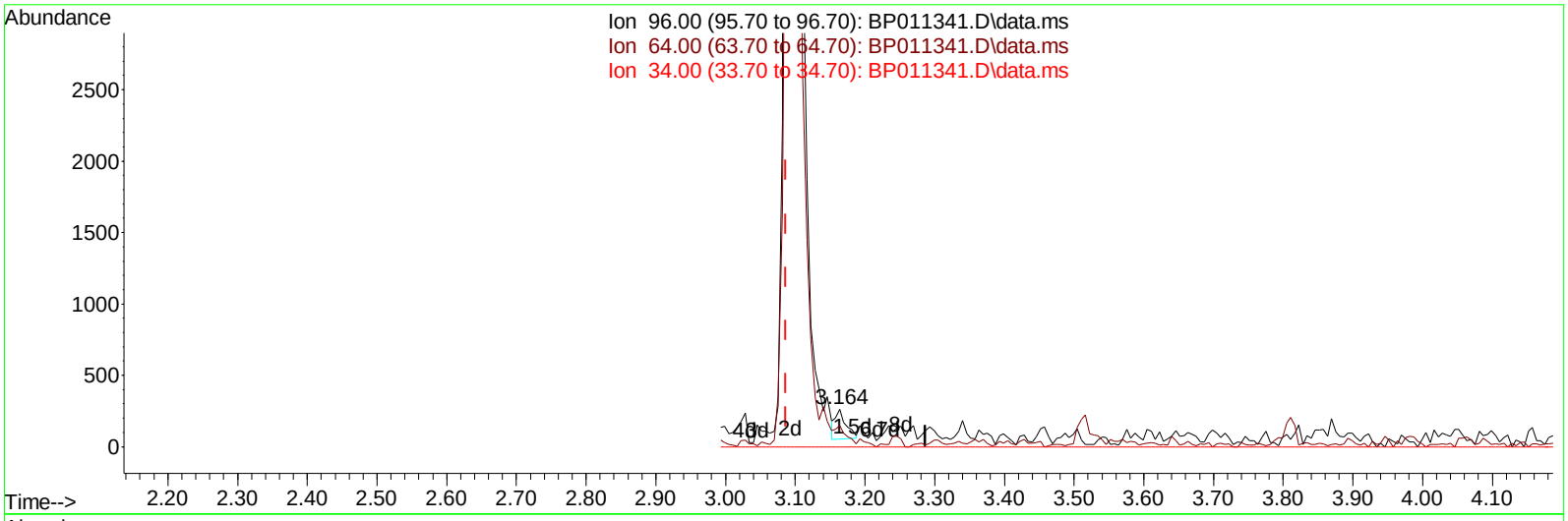
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011341.D
Acq On : 09 Aug 2022 21:54
Operator : CG/JU
Sample : N3900-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

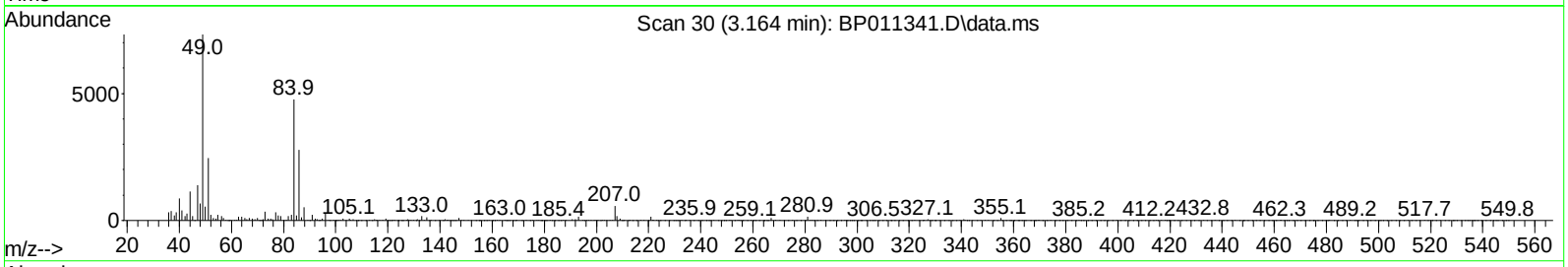
Quant Time: Aug 09 23:16:50 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

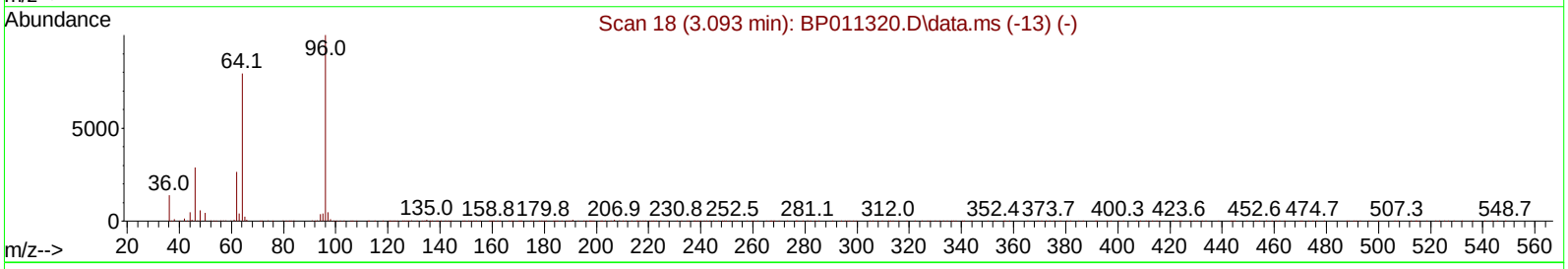
Instrument :
BNA_P
ClientSampleId :
C0C86

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011341.D\data.ms

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

3.164min (+ 0.077) 0.05 ng/uL

response 223

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	56.06
34.00	0.00	0.00
0.00	0.00	0.00

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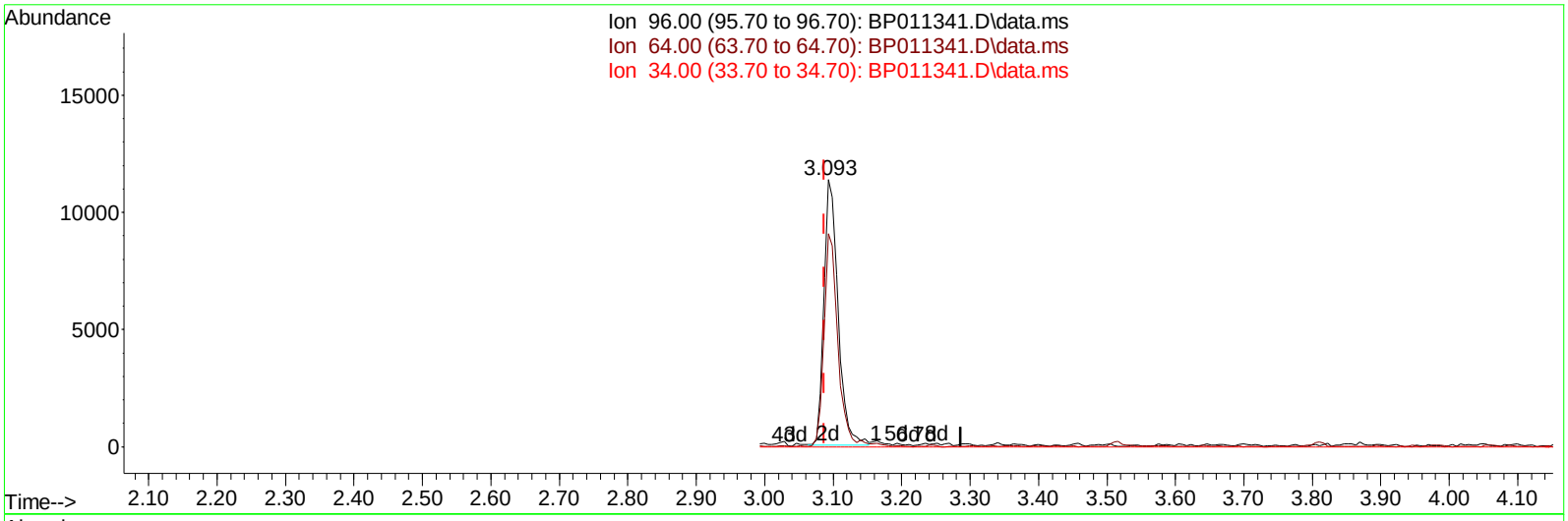

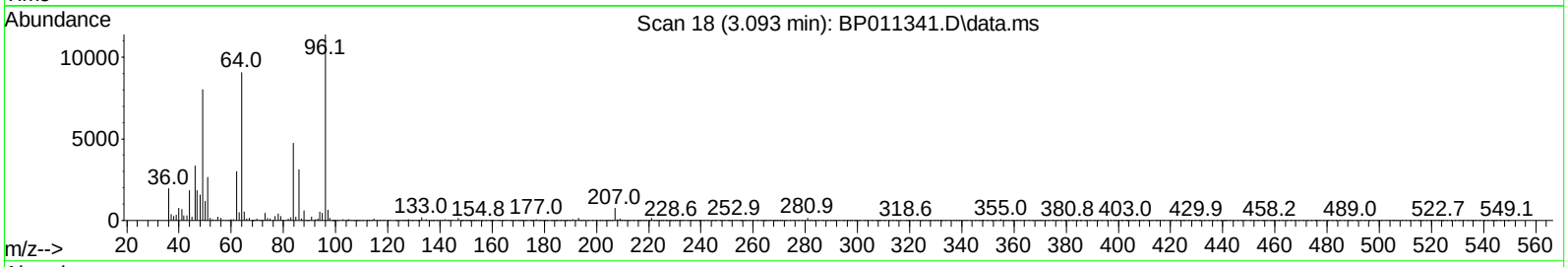

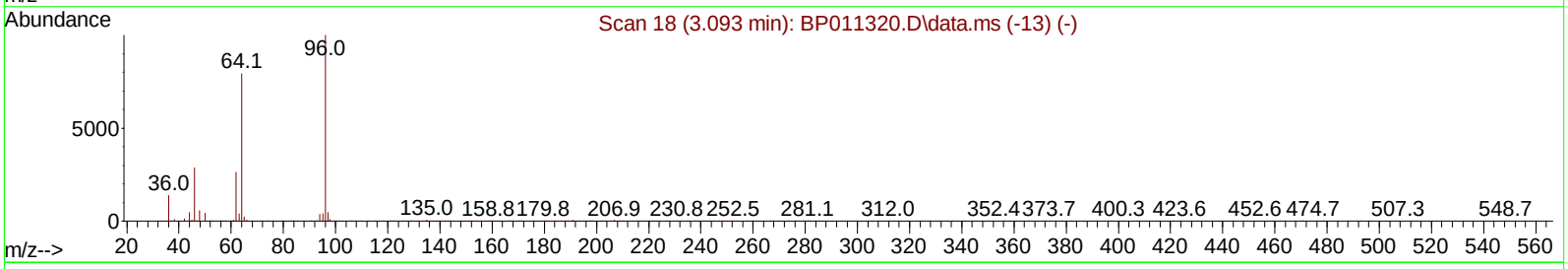

TIC: BP011341.D\data.ms

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

3.093min (+ 0.006) 3.89 ng/uL m

response 16402

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	79.77#
34.00	0.00	0.00
0.00	0.00	0.00

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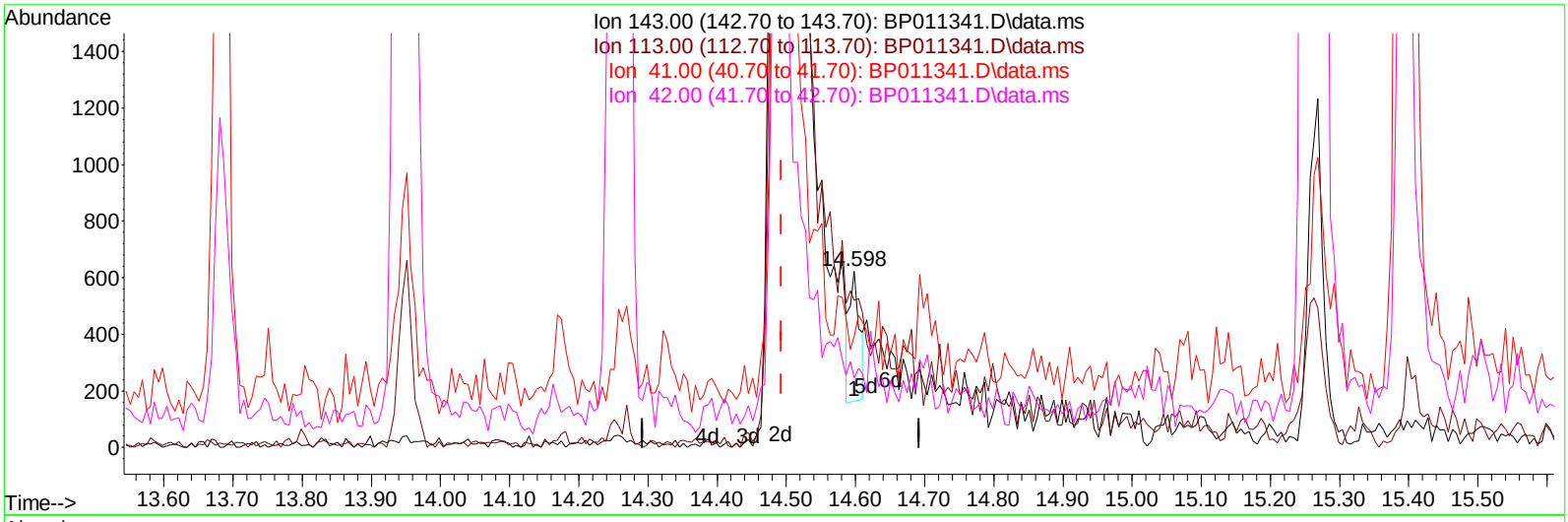

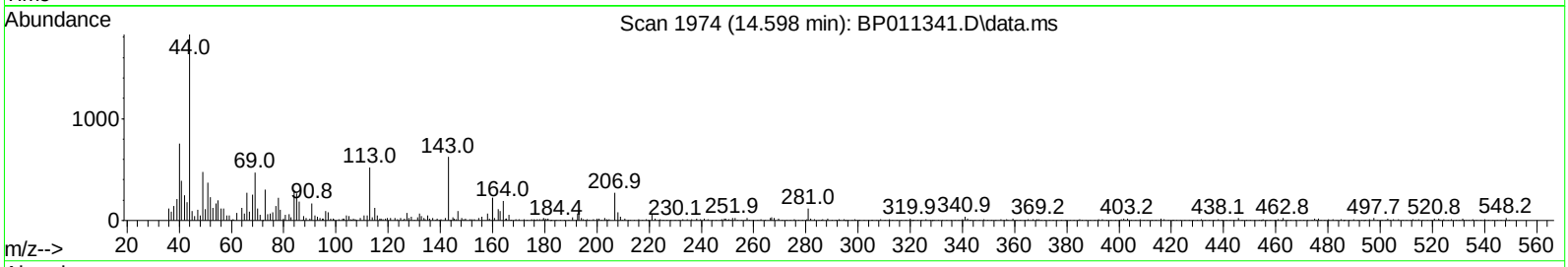

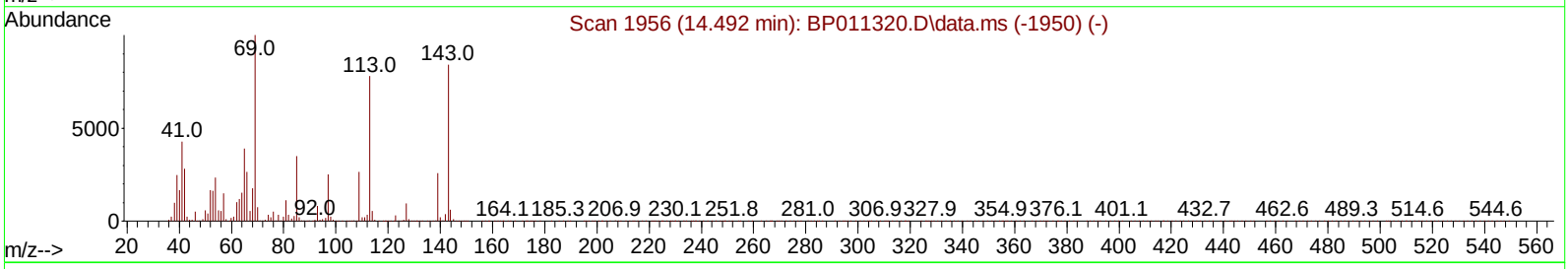

TIC: BP011341.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.598min (+ 0.106) 0.11 ng/ul

response 456

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	83.63
41.00	54.40	63.08
42.00	35.50	39.49

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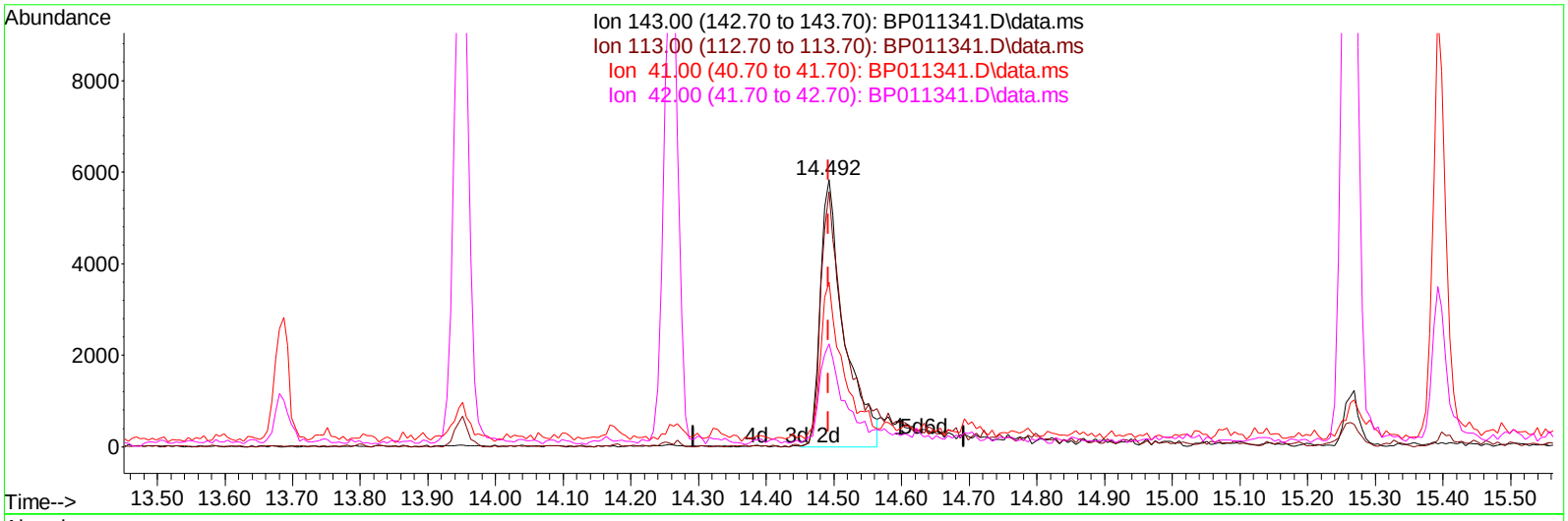

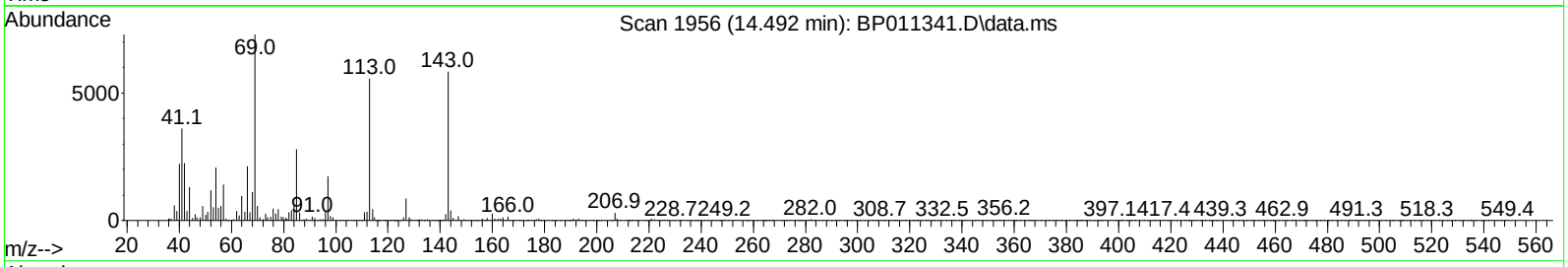

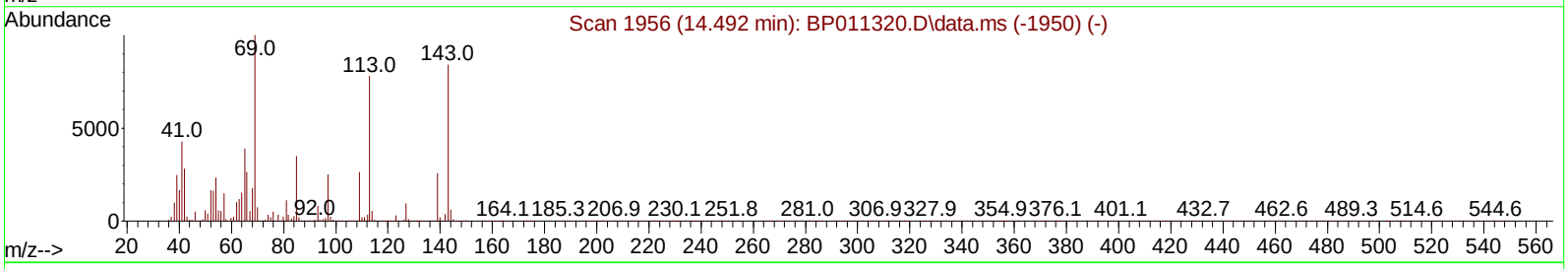

TIC: BP011341.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 3.54 ng/ul m

response 14188

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	95.59#
41.00	54.40	61.72
42.00	35.50	38.55

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	129312	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	547692	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	316016	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	614358	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	565282	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	583316	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	16402m	3.888	ng/uL	0.00
4) Pyridine-d5	3.511	84	77209	7.256	ng/ul	0.01
7) Phenol-d5	6.775	99	67517	5.900	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	247361	30.485	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	216627	24.126	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	134628	14.983	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	133595	35.841	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	125914	36.896	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	204640	29.694	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	308229	25.582	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	827445	37.737	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	932519	35.330	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	14188m	3.543	ng/ul	0.00
60) Fluorene-d10	15.263	176	686746	36.623	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	96603	33.853	ng/ul	0.00
73) Anthracene-d10	17.133	188	1153630	41.777	ng/ul	0.00
81) Pyrene-d10	19.392	212	1301957	41.805	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.315	264	1236654	42.039	ng/ul	-0.01
Target Compounds						Qvalue

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

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