Quantitation Report (Qedit)

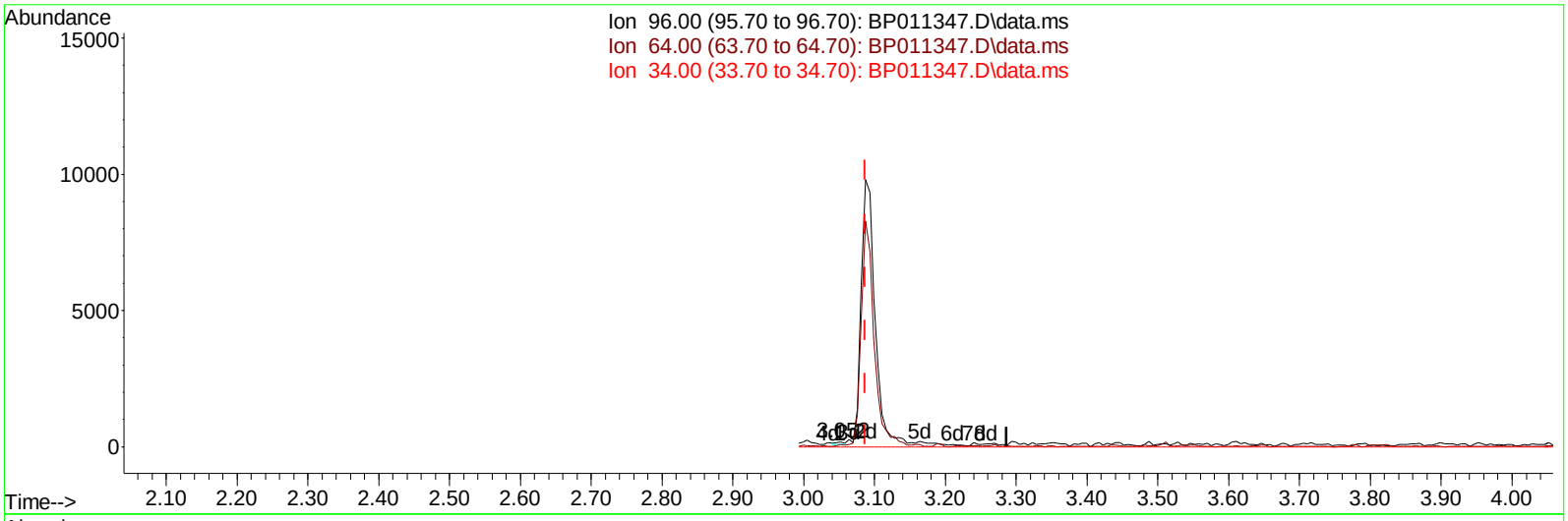
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011347.D
Acq On : 10 Aug 2022 01:23
Operator : CG/JU
Sample : N3900-15
Misc :
ALS Vial : 28 Sample Multiplier: 1

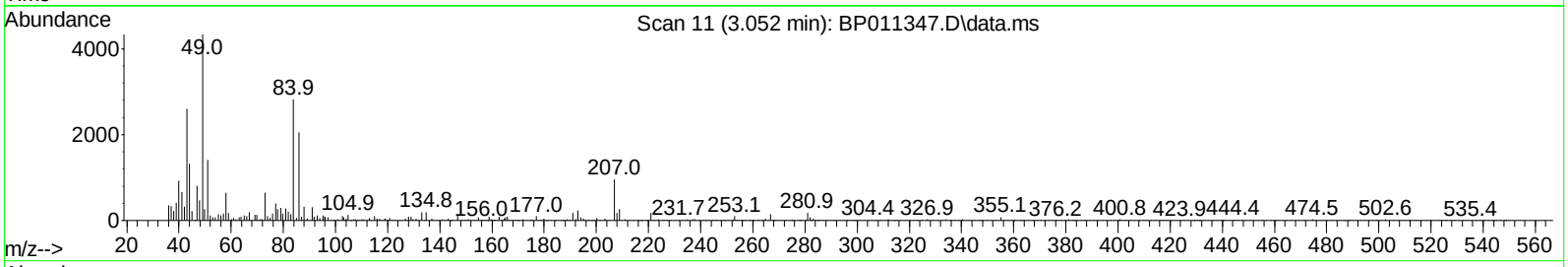
Quant Time: Aug 10 02:19:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

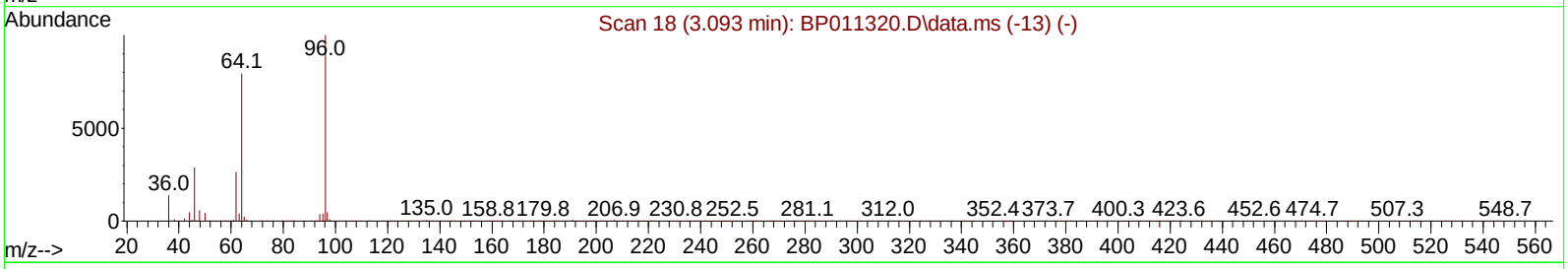
Instrument :
BNA_P
ClientSampleId :
C0CA8

Manual IntegrationsAPPROVED

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







TIC: BP011347.D\data.ms

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

3.052min (-0.035) 0.02 ng/uL

response 74

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	49.42
34.00	0.00	0.00
0.00	0.00	0.00

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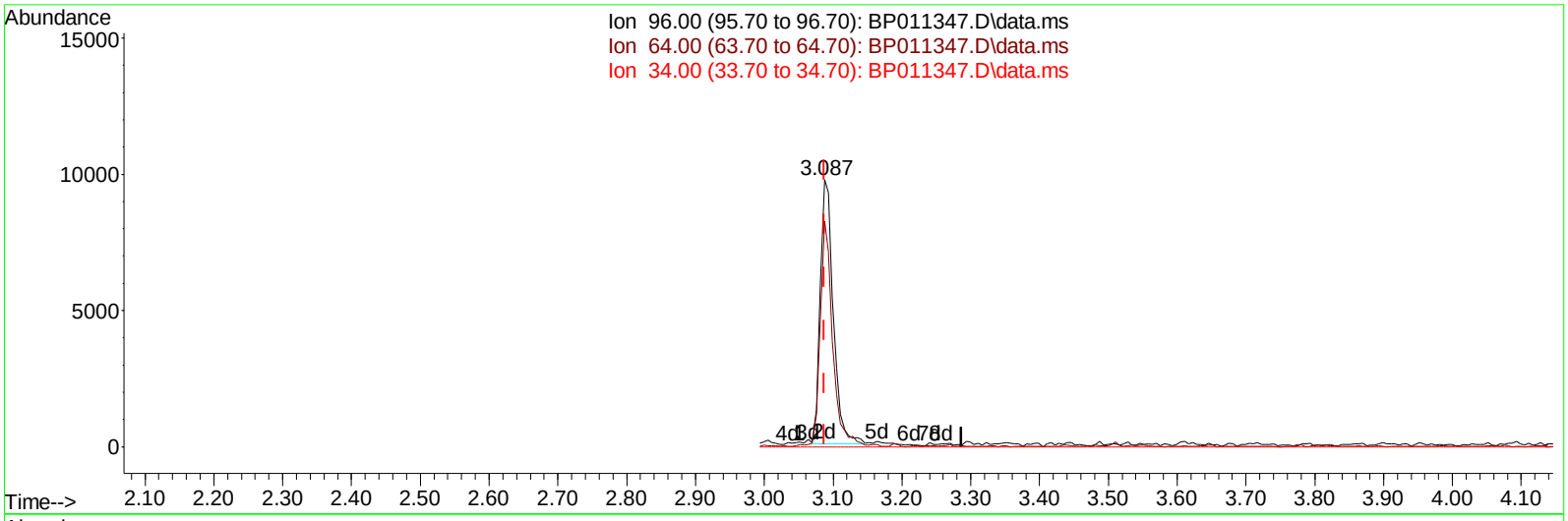

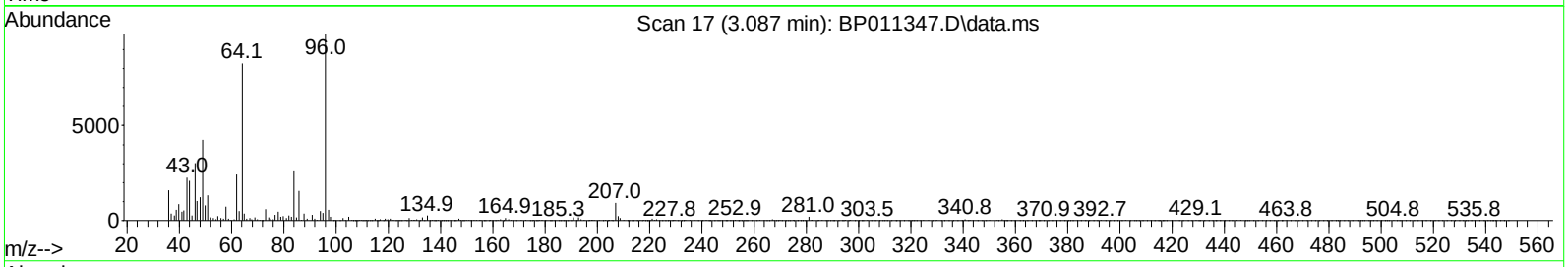

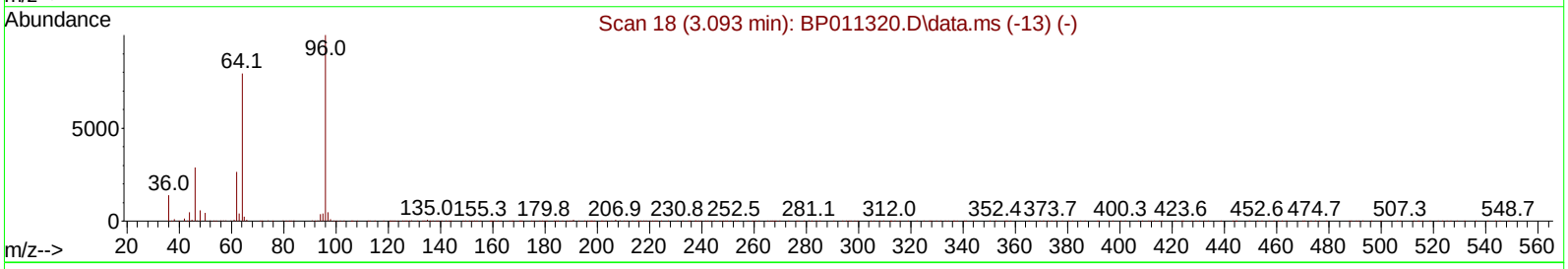

TIC: BP011347.D\data.ms

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

3.087min (+ 0.000) 3.84 ng/uL m

response 12937

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	84.49#
34.00	0.00	0.00
0.00	0.00	0.00

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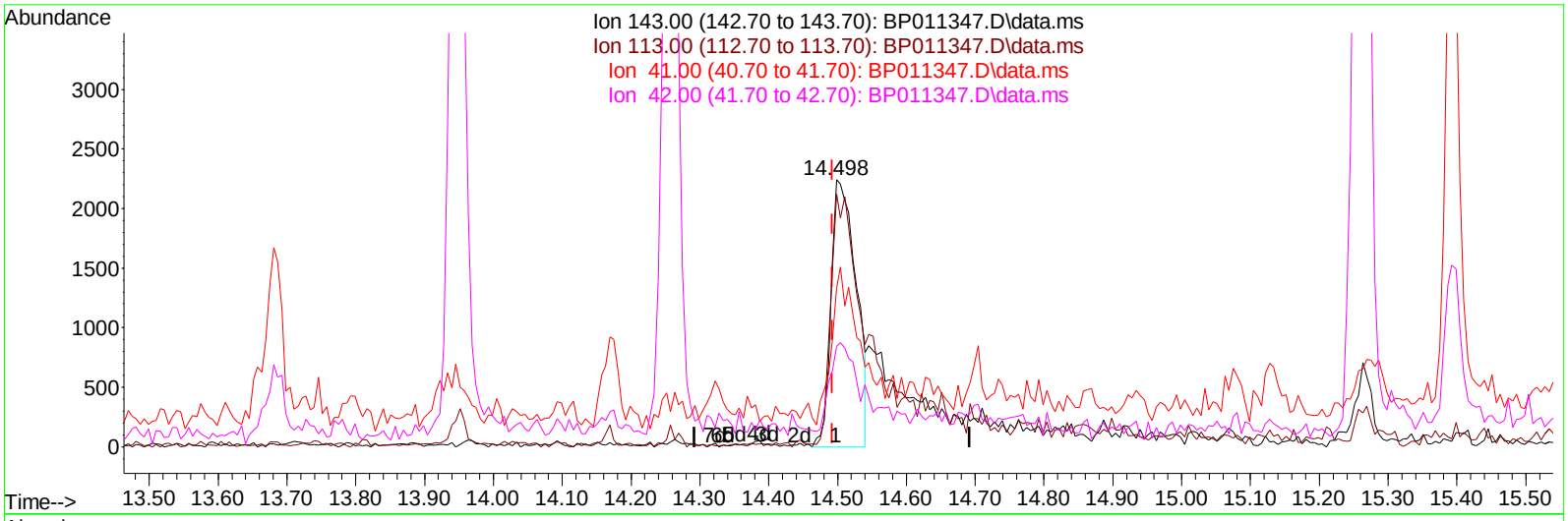

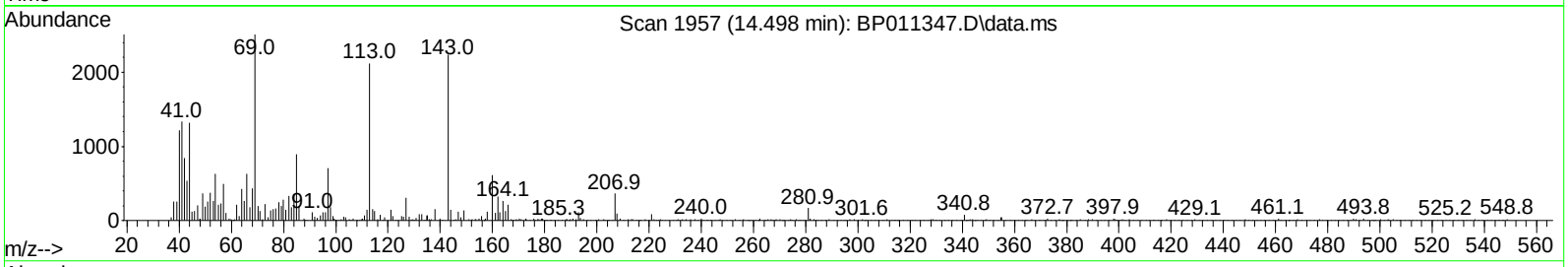

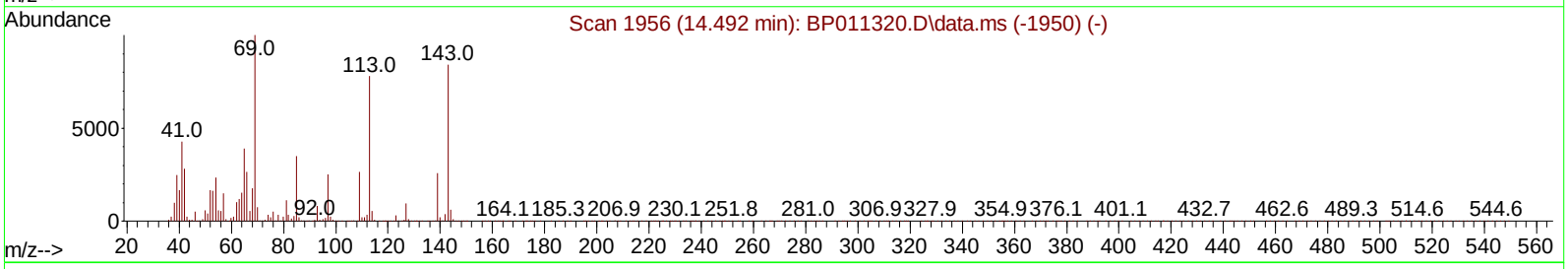

TIC: BP011347.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 1.96 ng/ul

response 5619

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.69#
41.00	54.40	59.53
42.00	35.50	37.71

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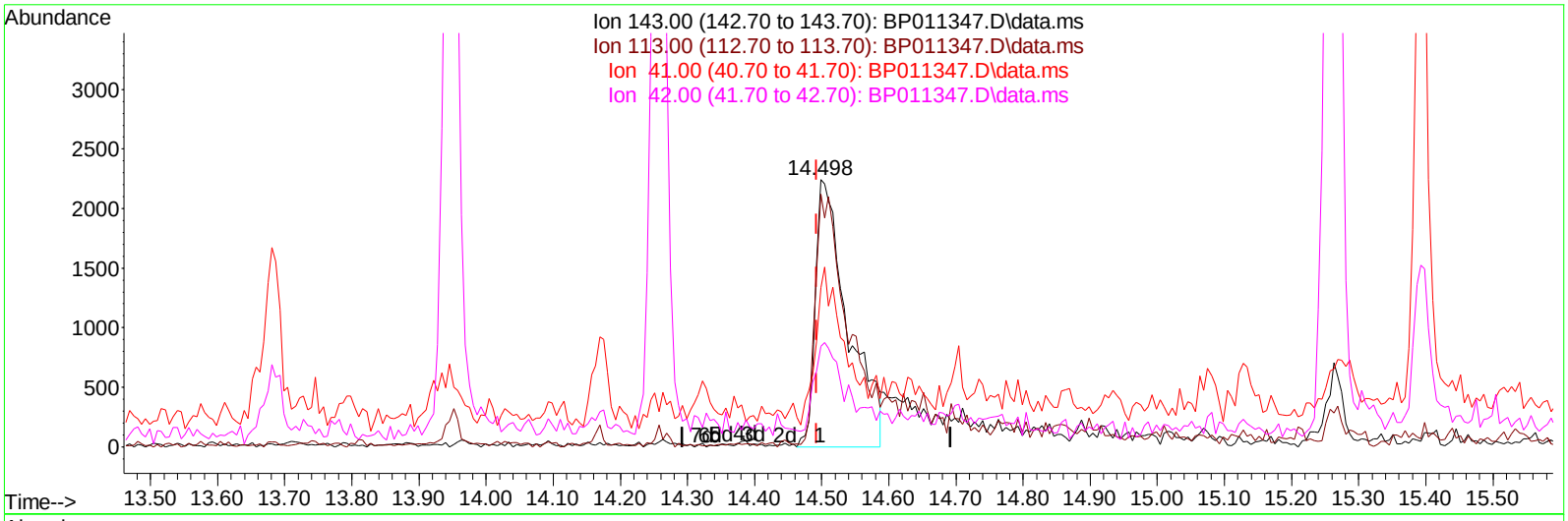

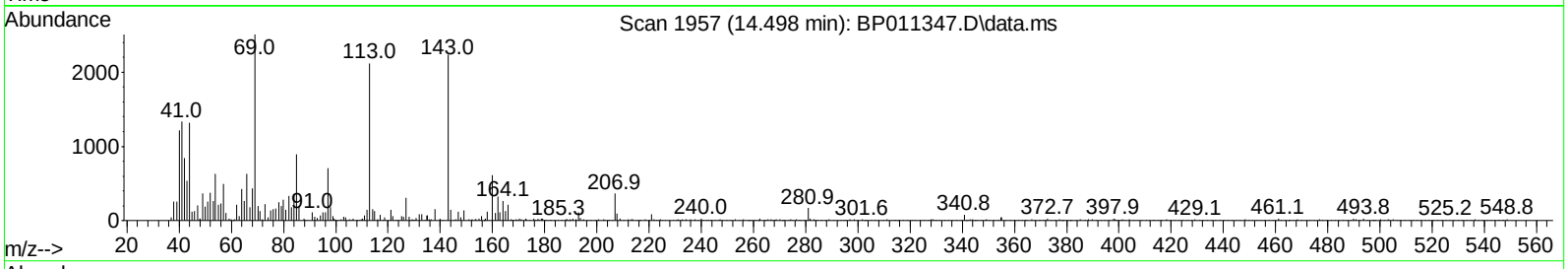

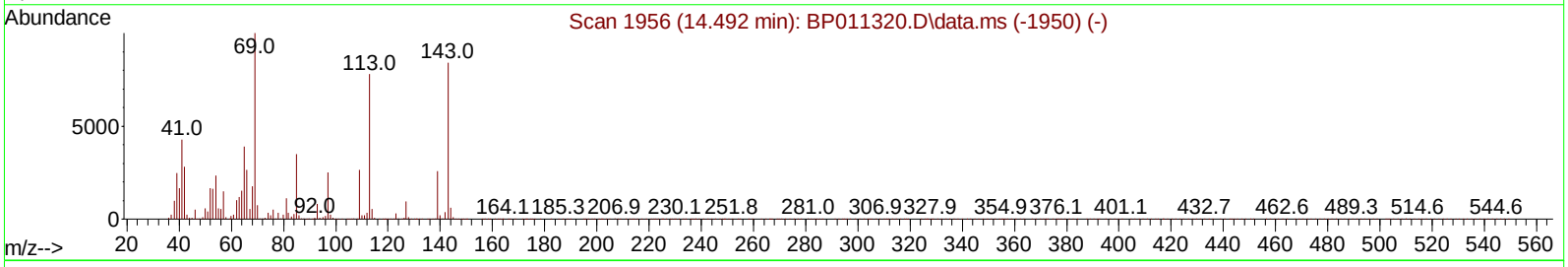

TIC: BP011347.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.62 ng/ul m

response 7504

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.69#
41.00	54.40	59.53
42.00	35.50	37.71

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	103328	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	413641	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	226130	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	444632	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	445696	20.000	ng/ul	-0.01
88) Perylene-d12	23.463	264	458821	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	12937m	3.838	ng/uL	0.00
4) Pyridine-d5	3.505	84	58405	6.869	ng/ul	0.00
7) Phenol-d5	6.775	99	43647	4.773	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	161257	24.871	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	142379	19.844	ng/ul	0.00
15) 4-Methylphenol-d8	8.311	113	81824	11.396	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	87002	30.905	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	79940	31.016	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	124900	23.996	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	156886	17.241	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	480429	30.620	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	557783	29.533	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	7504m	2.619	ng/ul	0.00
60) Fluorene-d10	15.263	176	398194	29.676	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.393	200	50726	24.562	ng/ul	0.00
73) Anthracene-d10	17.128	188	681761	34.113	ng/ul	-0.01
81) Pyrene-d10	19.386	212	831139	33.848	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.310	264	781803	33.788	ng/ul	-0.02
Target Compounds						Qvalue

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

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