Quantitation Report (Qedit)

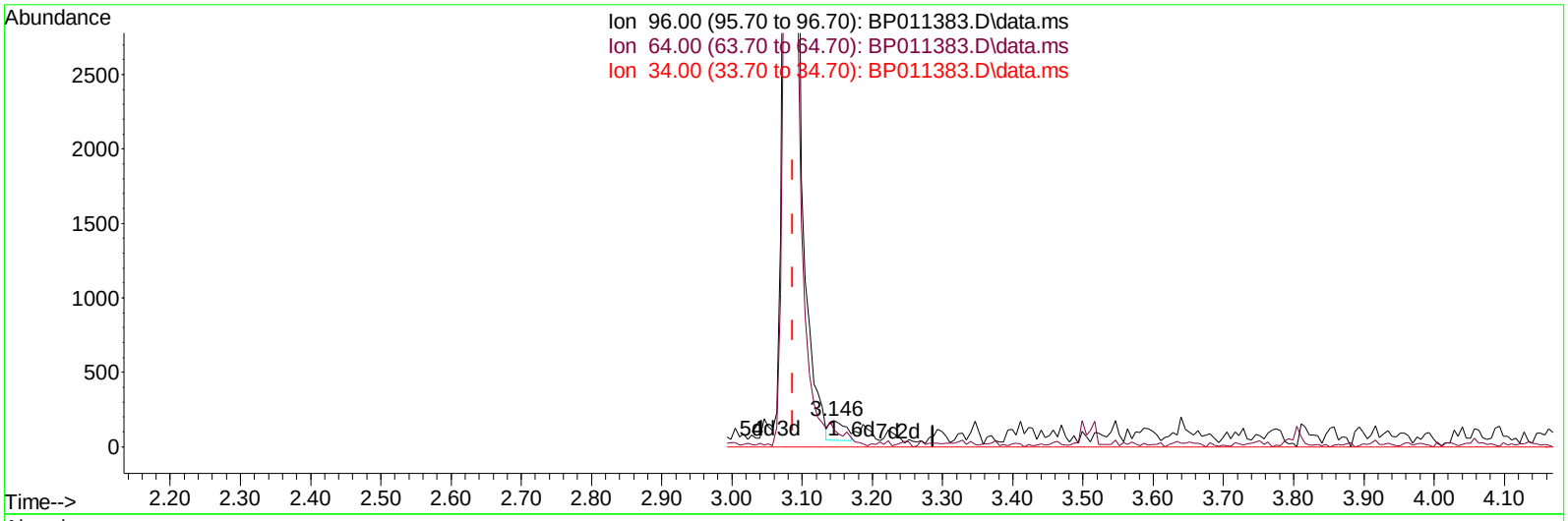
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011383.D
Acq On : 11 Aug 2022 12:44
Operator : CG/JU
Sample : N4128-04
Misc :
ALS Vial : 6 Sample Multiplier: 1

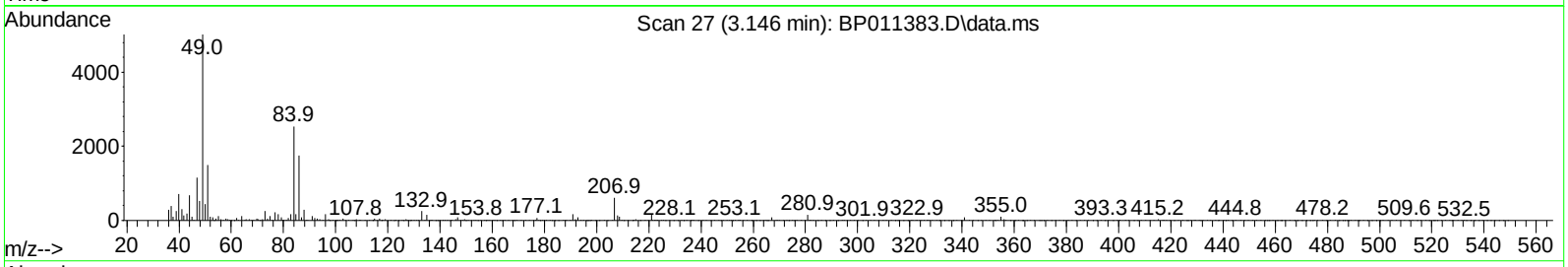
Quant Time: Aug 11 14:12:12 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

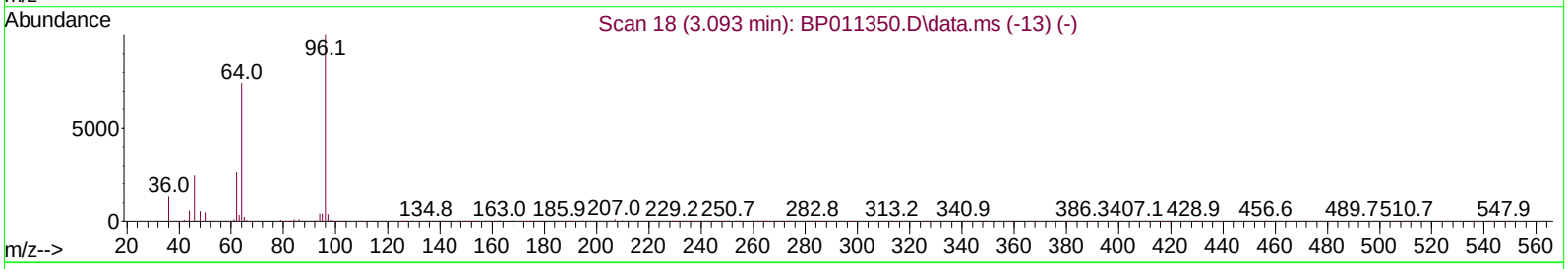
Instrument :
BNA_P
ClientSampleId :
BGC11

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/11/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011383.D\data.ms

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

3.146min (+ 0.059) 0.09 ng/uL

response 216

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	63.48
34.00	0.00	0.00
0.00	0.00	0.00

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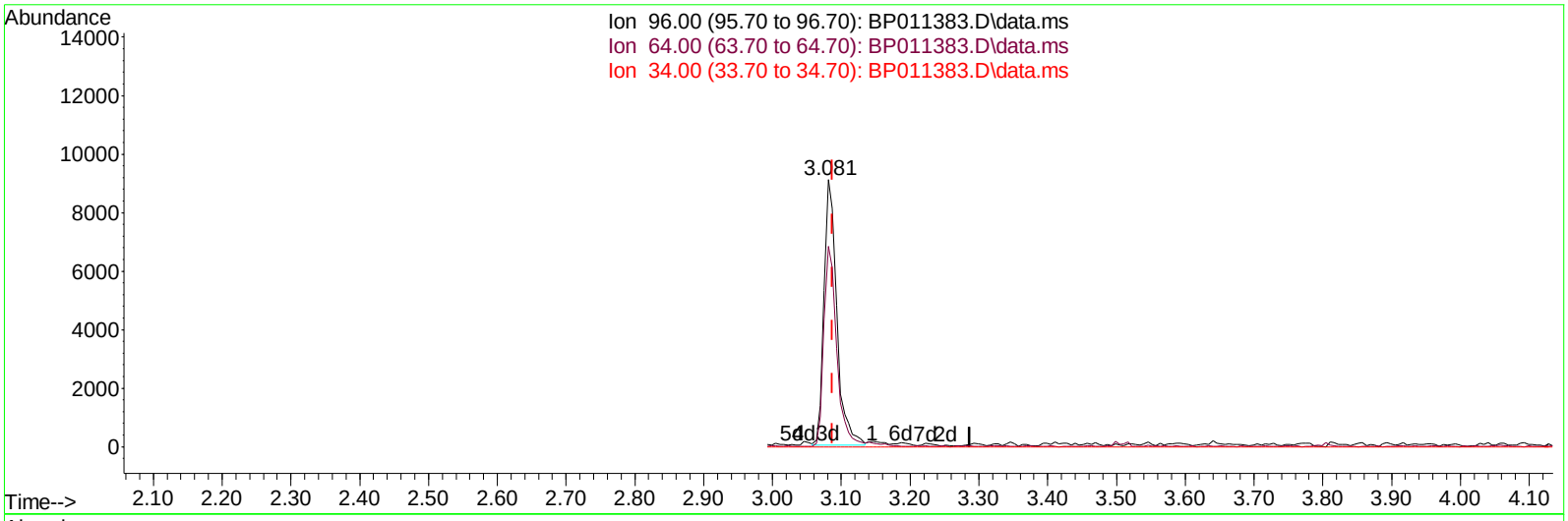

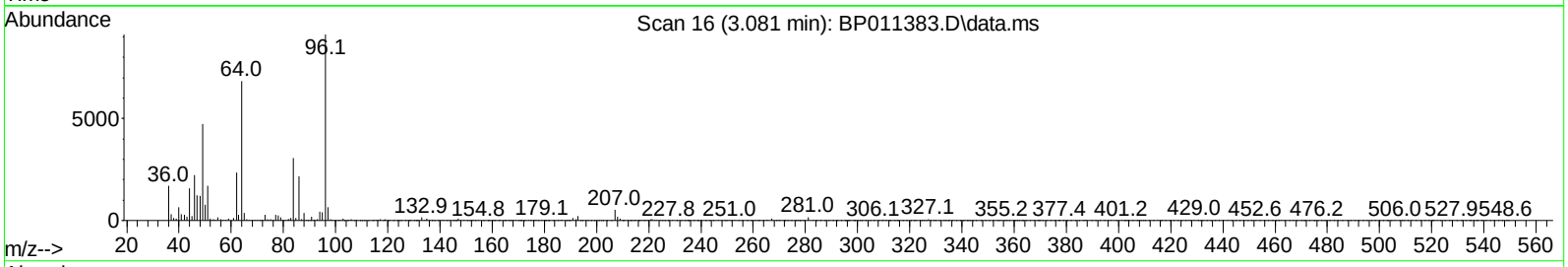

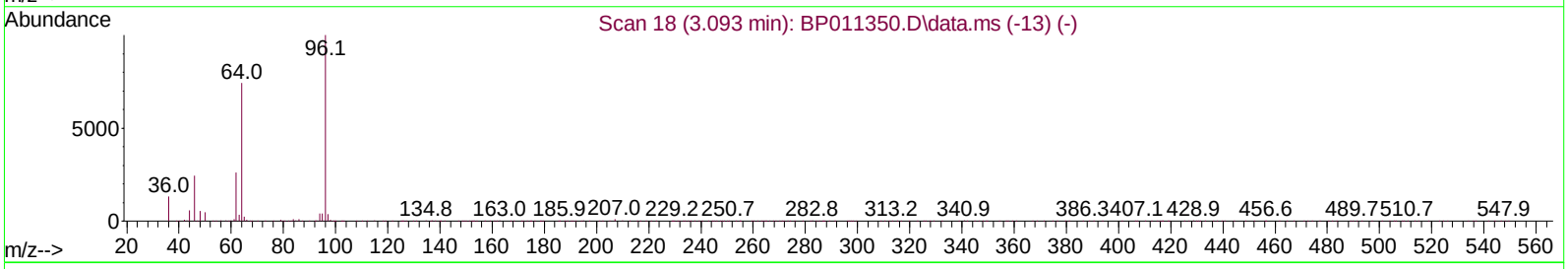

TIC: BP011383.D\data.ms

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

3.081min (-0.006) 4.76 ng/uL m

response 11690

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	74.98#
34.00	0.00	0.00
0.00	0.00	0.00

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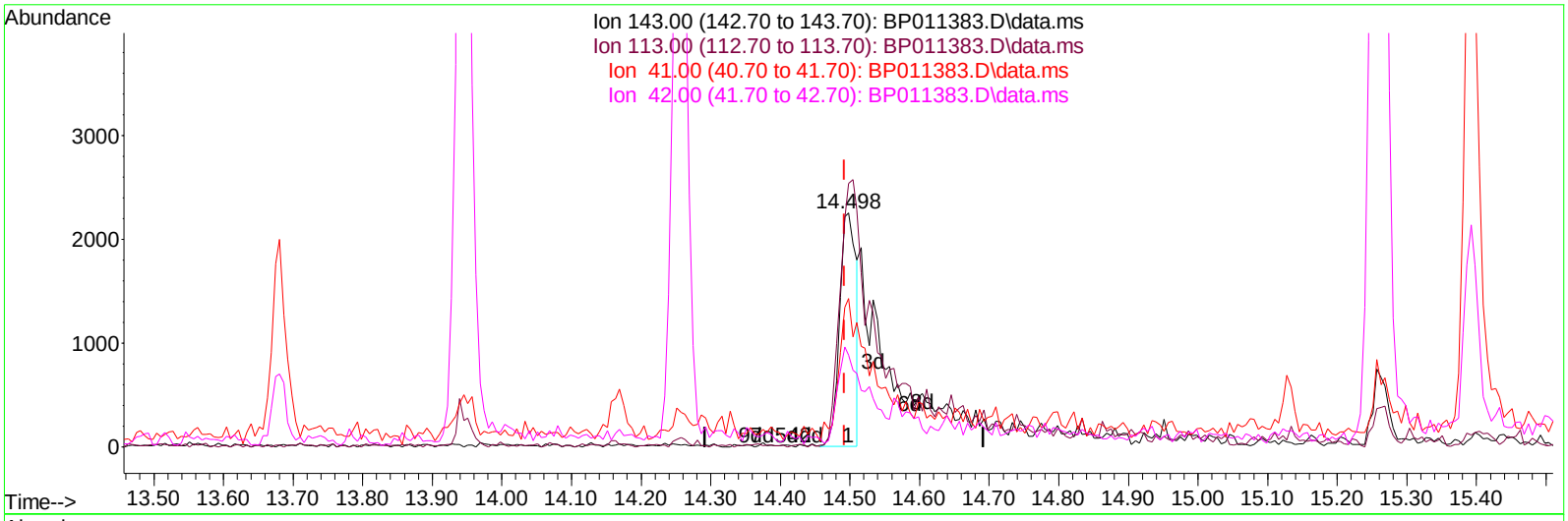

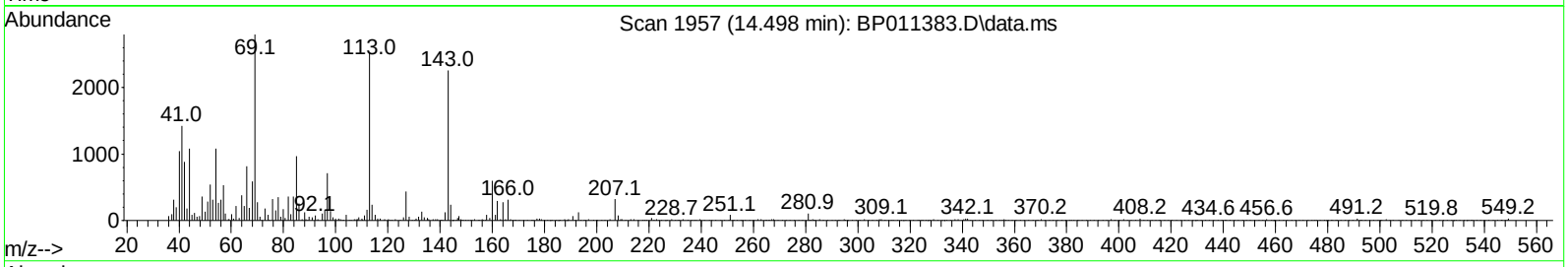

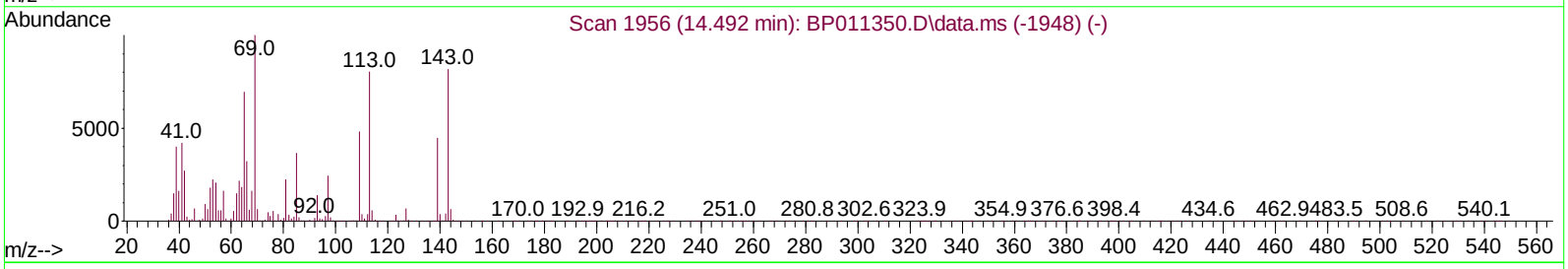

TIC: BP011383.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 1.64 ng/ul

response 3734

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	112.46#
41.00	54.40	63.24
42.00	35.50	39.11

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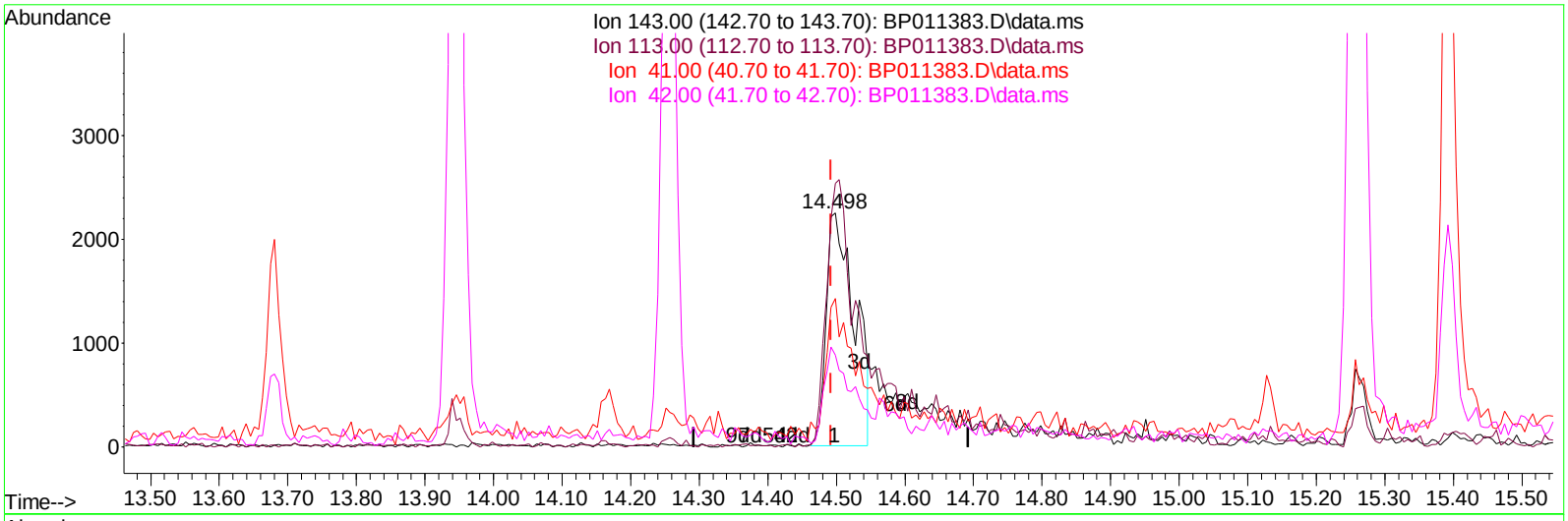

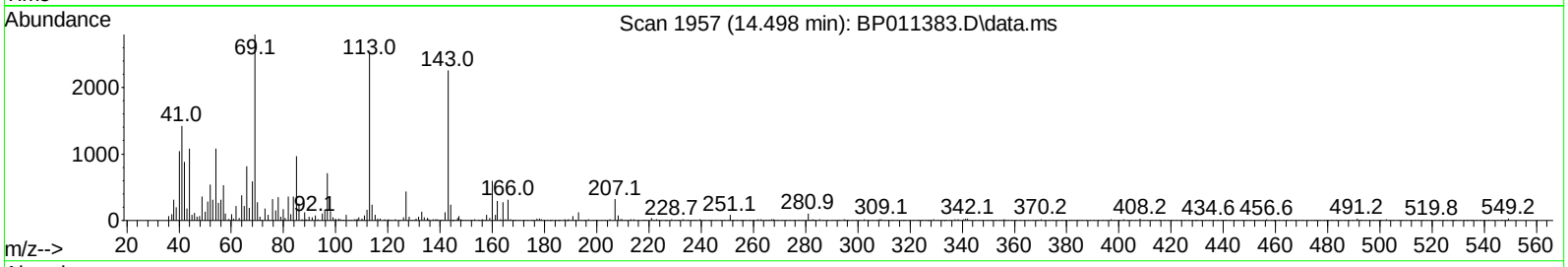

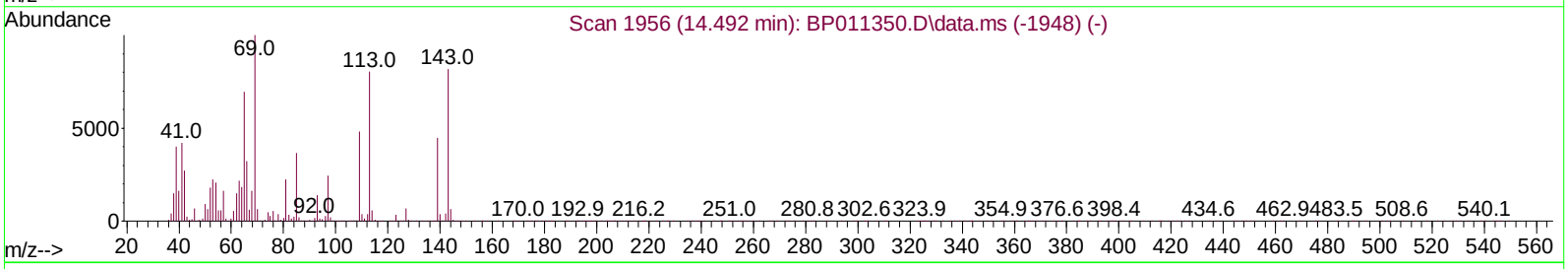

TIC: BP011383.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.006) 2.80 ng/ul m

response 6366

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	112.46#
41.00	54.40	63.24
42.00	35.50	39.11

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	75224	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.369	136	314707	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	179132	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	371755	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	358914	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	389523	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11690m	4.764	ng/uL	0.00
4) Pyridine-d5	3.499	84	53268	8.606	ng/ul	0.00
7) Phenol-d5	6.763	99	44443	6.676	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	155714	32.989	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.116	132	138156	26.450	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	83002	15.879	ng/ul	-0.01
21) Nitrobenzene-d5	8.746	128	83728	39.092	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	81292	41.456	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.998	165	123086	31.082	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.528	131	174370	25.186	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	497610	40.036	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	574921	38.427	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	6366m	2.804	ng/ul	0.00
60) Fluorene-d10	15.263	176	414877	39.031	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.392	200	61192	35.438	ng/ul	0.00
73) Anthracene-d10	17.128	188	715096	42.796	ng/ul	-0.01
81) Pyrene-d10	19.392	212	847078	42.838	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	830388	42.272	ng/ul	0.00
Target Compounds						Qvalue

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

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