Quantitation Report (Qedit)

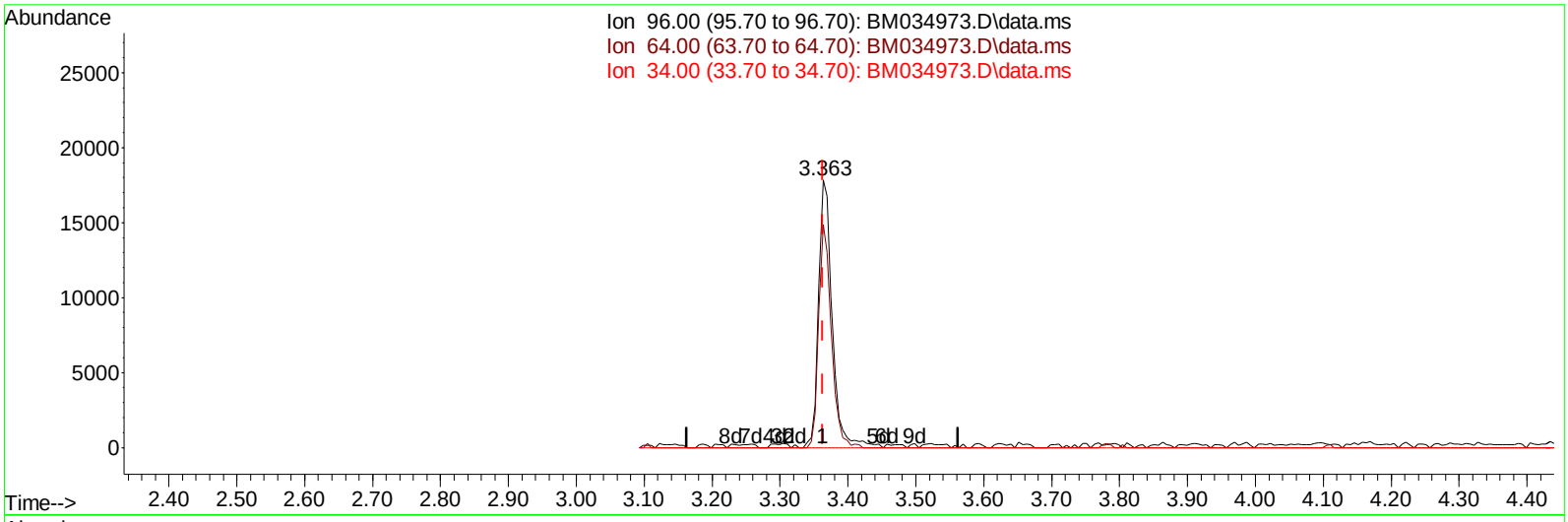
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051022\
Data File : BM034973.D
Acq On : 10 May 2022 21:01
Operator : CG/JU
Sample : N2603-07
Misc :
ALS Vial : 17 Sample Multiplier: 1

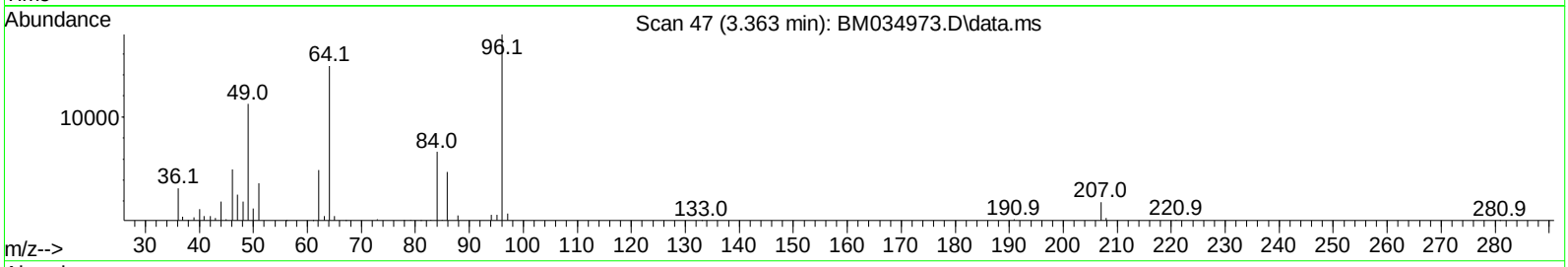
Quant Time: May 11 02:50:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 10 16:40:09 2022
Response via : Initial Calibration

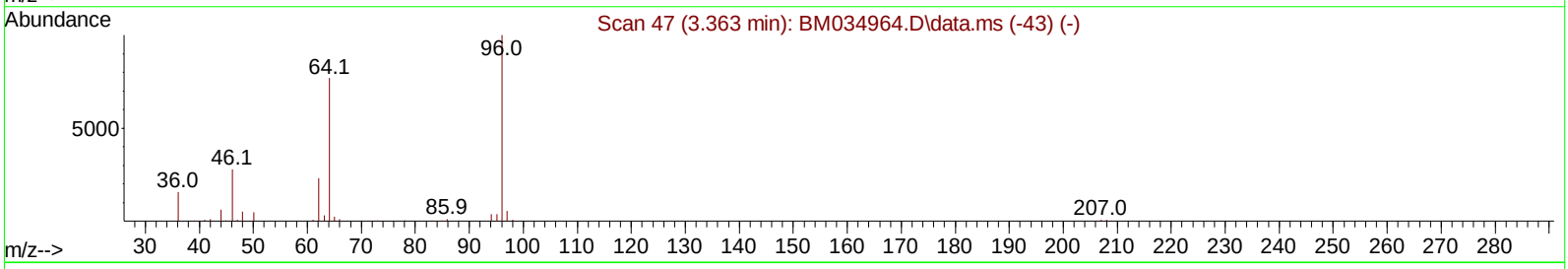
Instrument :
BNA_M
ClientSampleId :
EW488

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/11/2022
Supervised By :mohammad ahmed 05/12/2022







TIC: BM034973.D\data.ms

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

3.363min (+ 0.000) 4.98 ng/uL

response 24885

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.40	83.29
34.00	0.00	0.00
0.00	0.00	0.00

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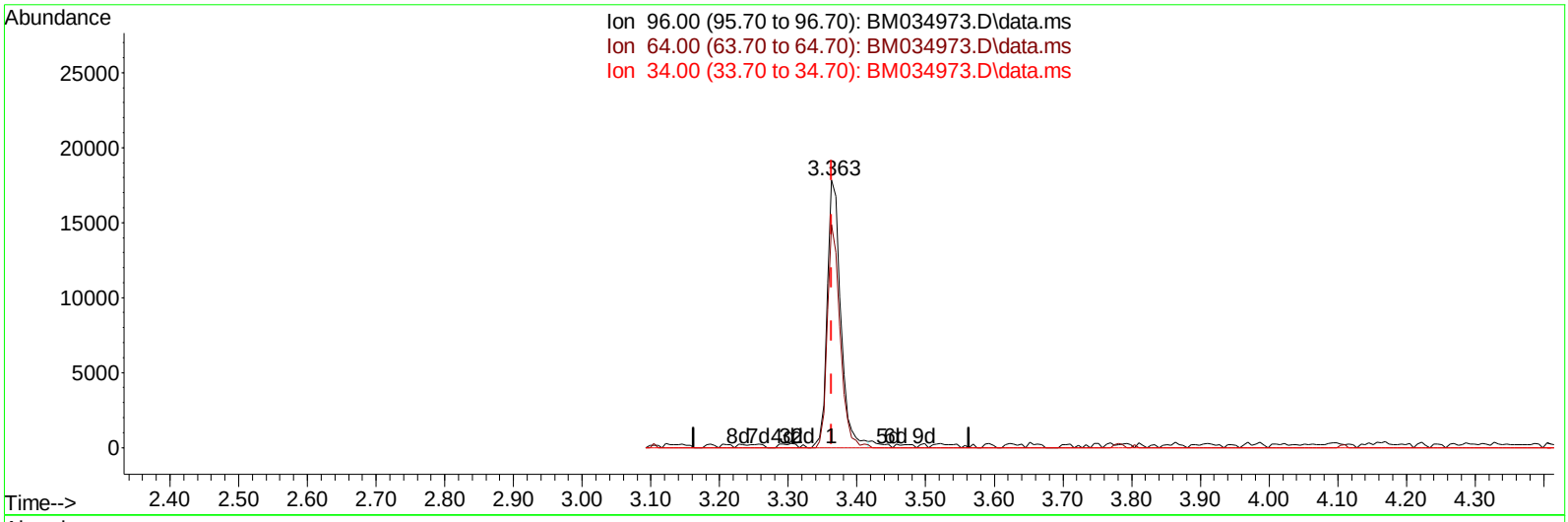

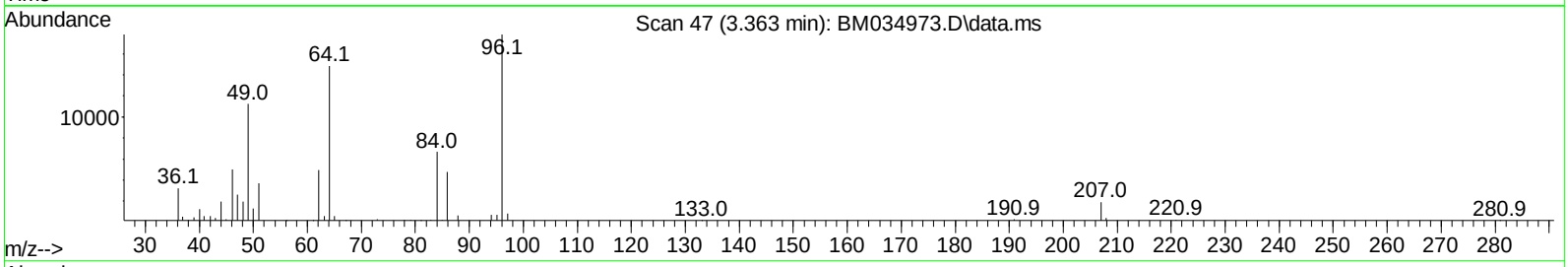

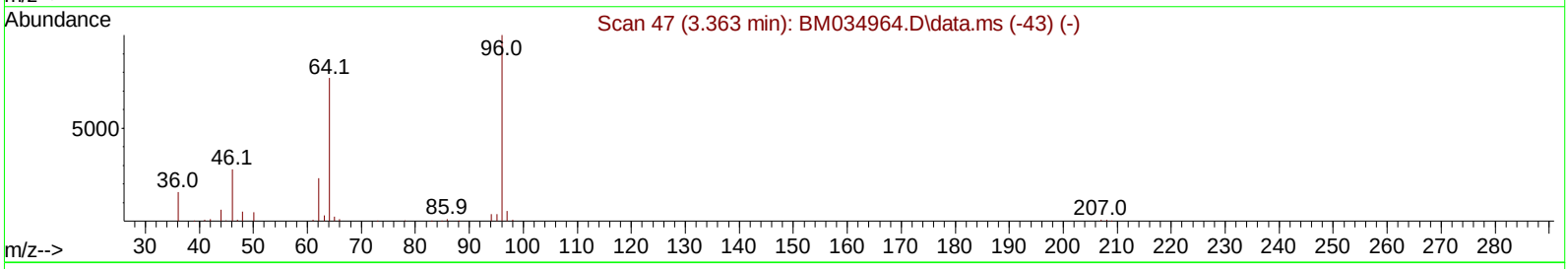

TIC: BM034973.D\data.ms

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

3.363min (+ 0.000) 4.90 ng/uL m

response 24481

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.40	83.29
34.00	0.00	0.00
0.00	0.00	0.00

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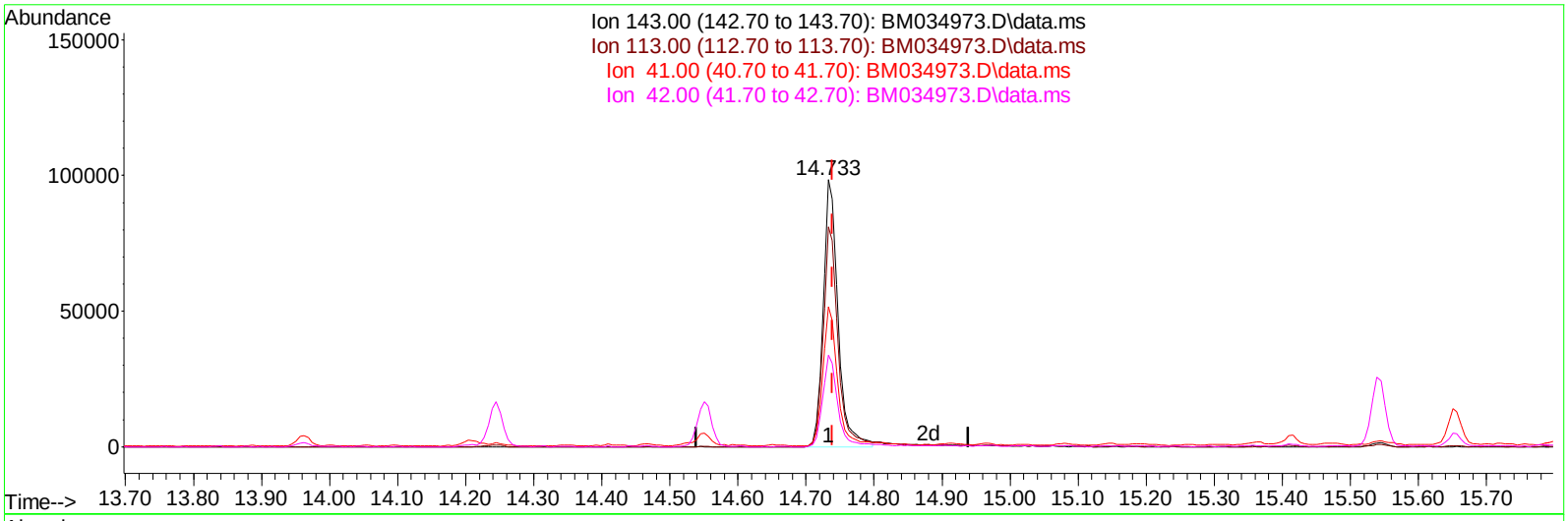

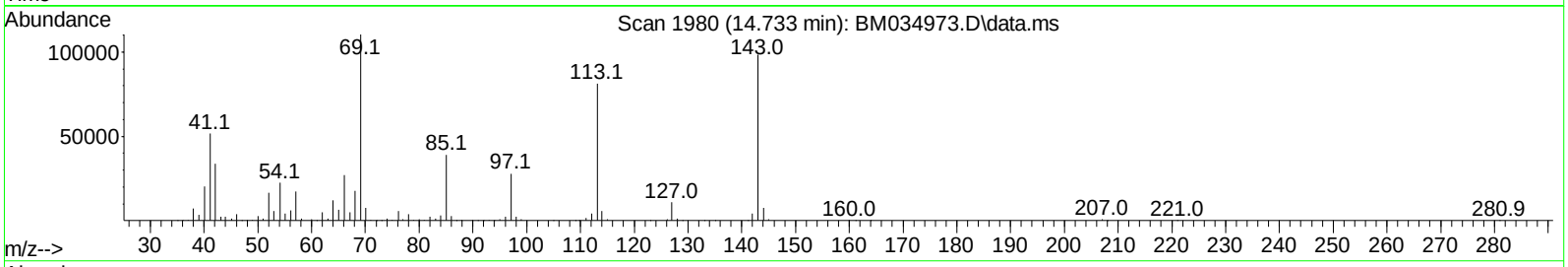

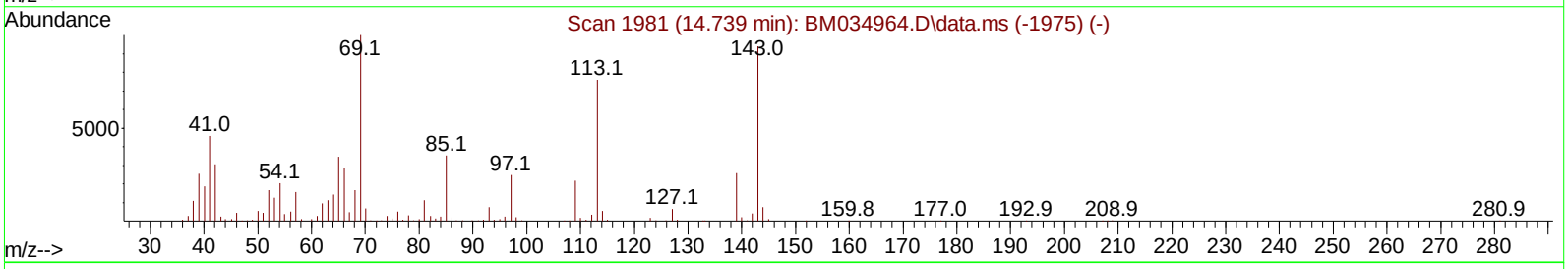

TIC: BM034973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (-0.006) 19.01 ng/ul

response 148330

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	82.51
41.00	42.20	52.48#
42.00	28.30	34.35#

Quantitation Report (Qedit)

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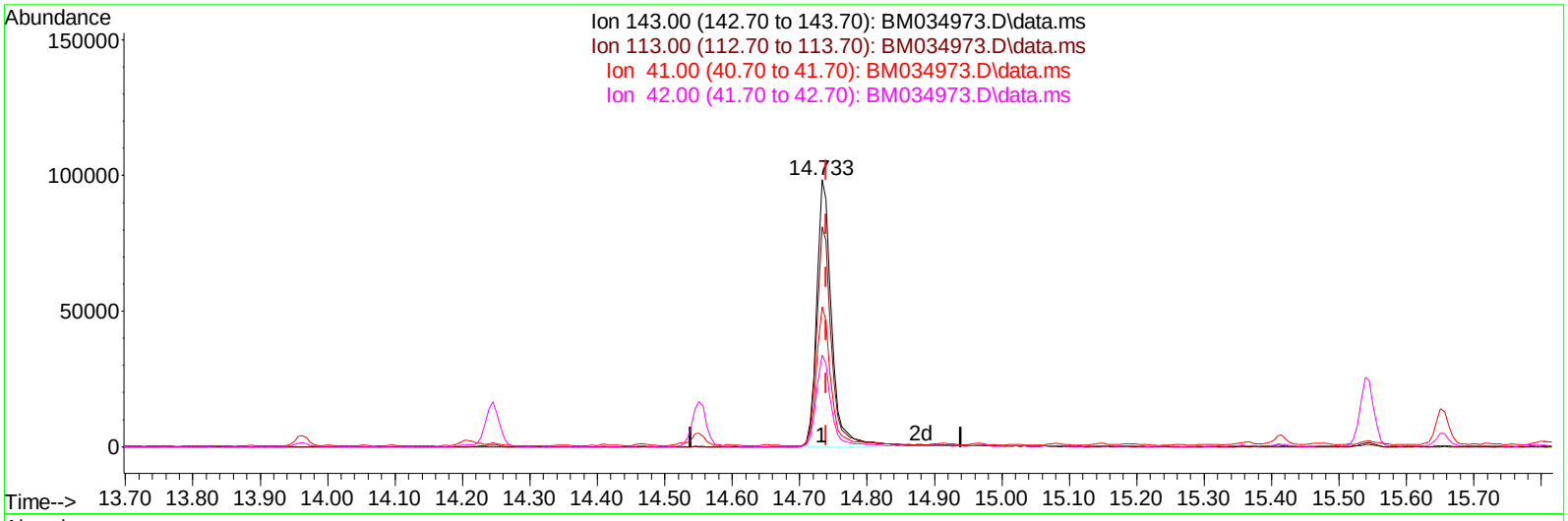

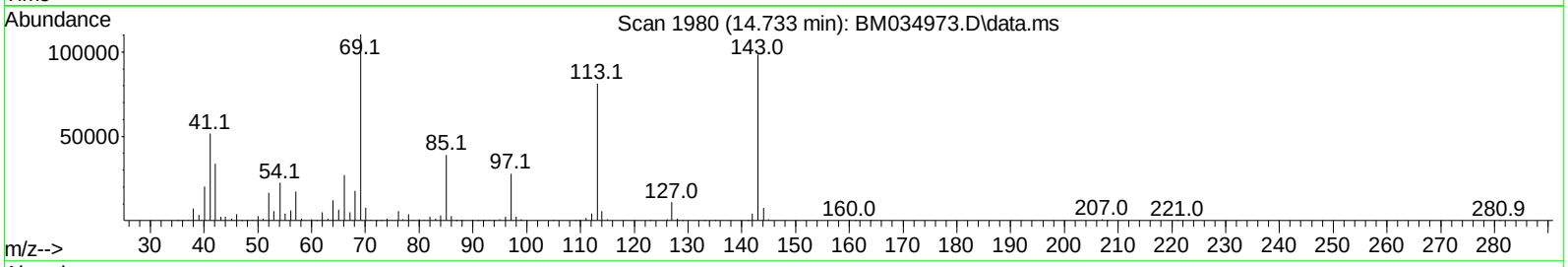

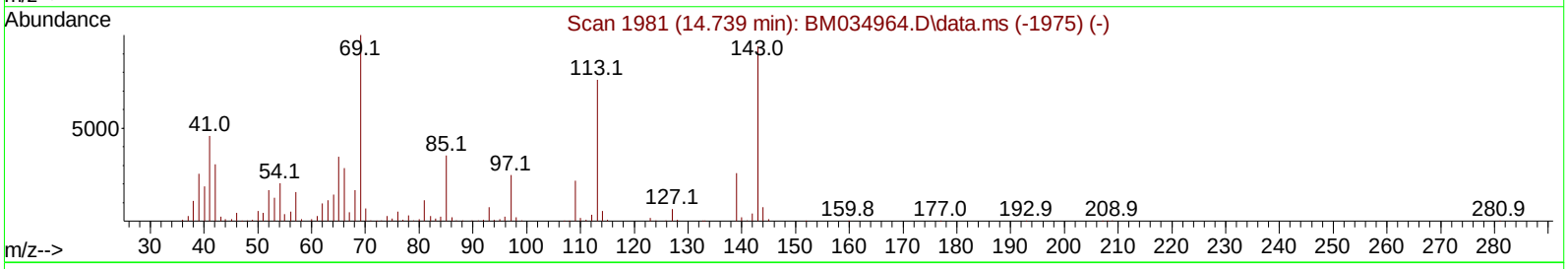

TIC: BM034973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (-0.006) 19.22 ng/ul m

response 149982

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	82.51
41.00	42.20	52.48#
42.00	28.30	34.35#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	209089	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	925763	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	605562	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	1190939	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240	836709	20.000	ng/ul	0.00
88) Perylene-d12	23.850	264	766346	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	24481m	4.898	ng/uL	0.00
4) Pyridine-d5	3.781	84	104262	7.569	ng/ul	0.00
7) Phenol-d5	7.081	99	420223	23.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.251	67	310085	27.312	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	379538	26.701	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	343805	23.313	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	193059	27.209	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	208980	27.307	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	361949	25.271	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	320205	15.702	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1297508	28.411	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1498860	25.865	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	149982m	19.220	ng/ul	0.00
60) Fluorene-d10	15.545	176	1101704	28.594	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	155538	20.725	ng/ul	0.00
73) Anthracene-d10	17.392	188	1625554	28.055	ng/ul	0.00
81) Pyrene-d10	19.680	212	1611447	30.843	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.697	264	1159044	28.939	ng/ul	0.00
Target Compounds						Qvalue
16) Acetophenone	8.745	105	27664	1.208	ng/ul#	92

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

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