Quantitation Report (Qedit)

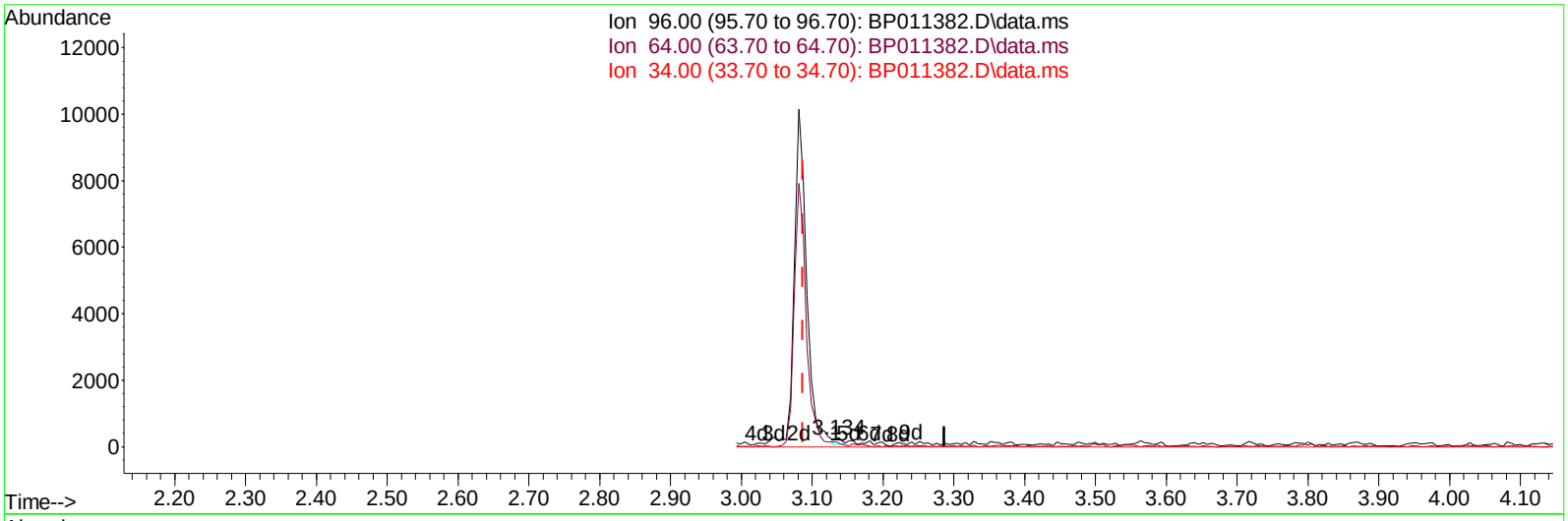
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011382.D
Acq On : 11 Aug 2022 12:09
Operator : CG/JU
Sample : N4128-07
Misc :
ALS Vial : 5 Sample Multiplier: 1

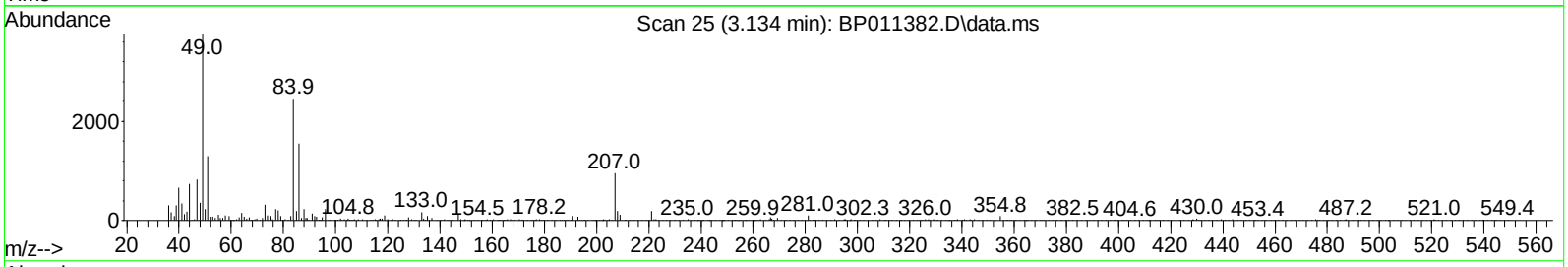
Quant Time: Aug 11 14:07:43 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

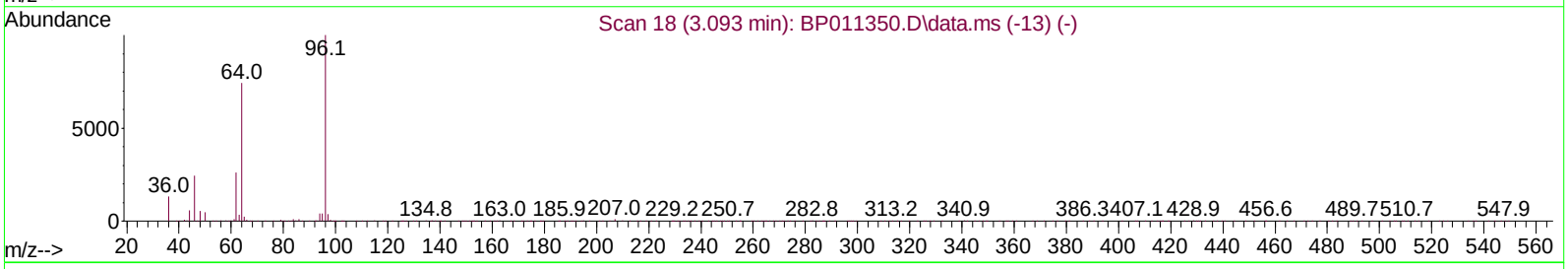
Instrument :
BNA_P
ClientSampleId :
BGBX3

Manual IntegrationsAPPROVED

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







TIC: BP011382.D\data.ms

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

3.134min (+ 0.047) 0.05 ng/uL

response 131

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	64.38
34.00	0.00	0.00
0.00	0.00	0.00

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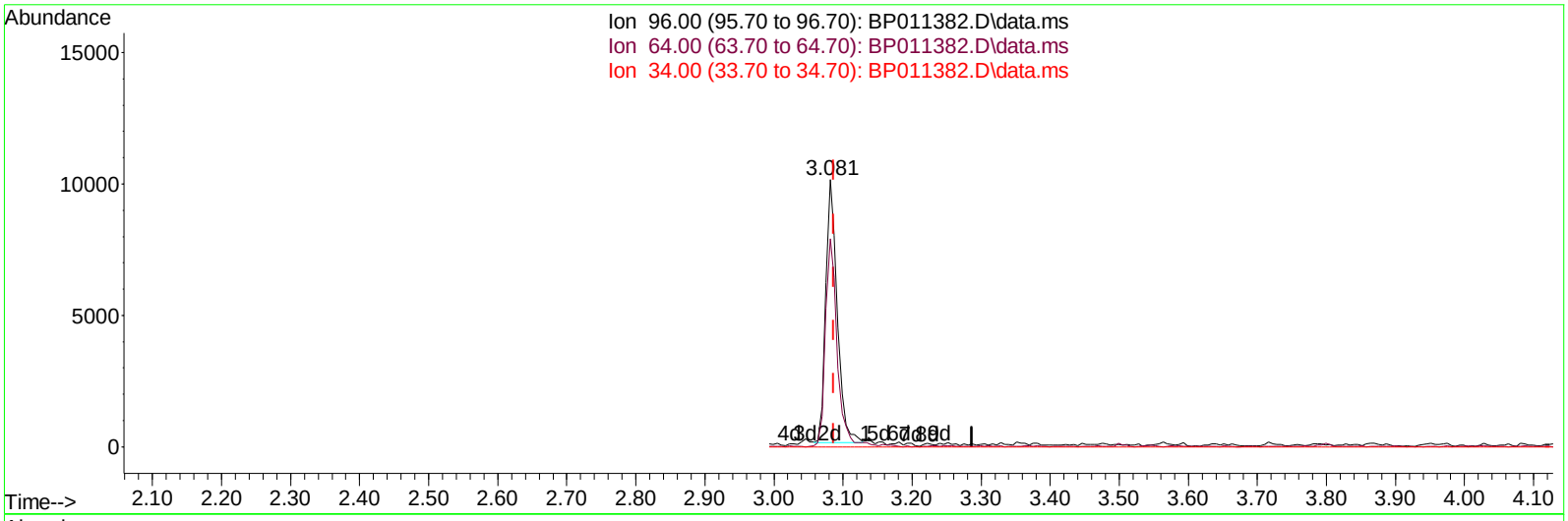

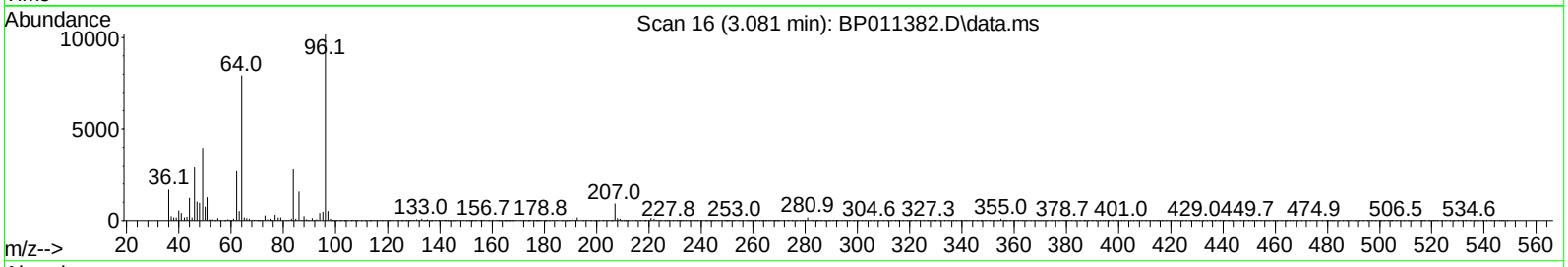

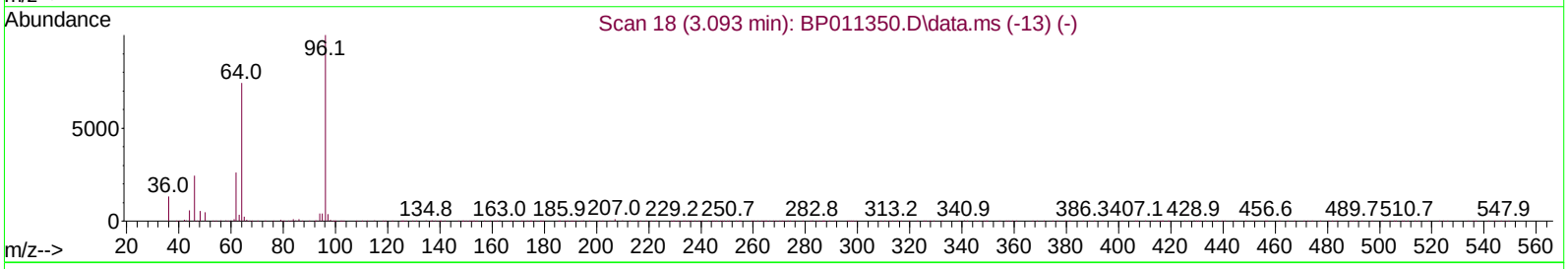

TIC: BP011382.D\data.ms

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

3.081min (-0.006) 4.26 ng/uL m

response 11644

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	77.92#
34.00	0.00	0.00
0.00	0.00	0.00

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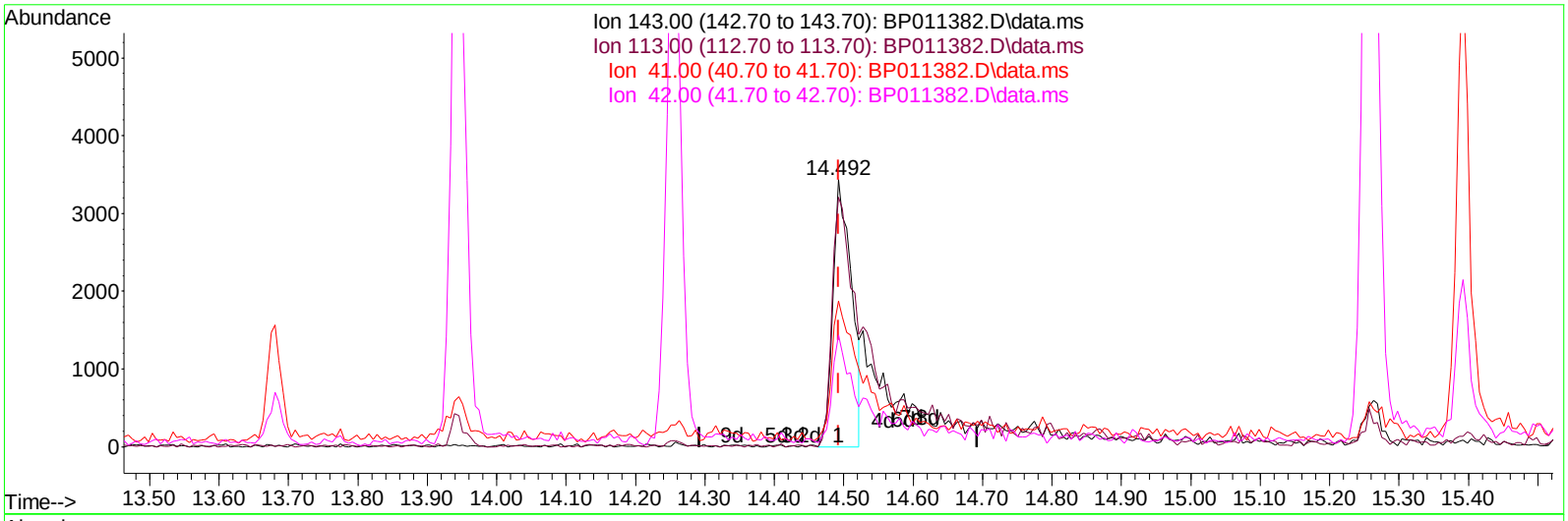

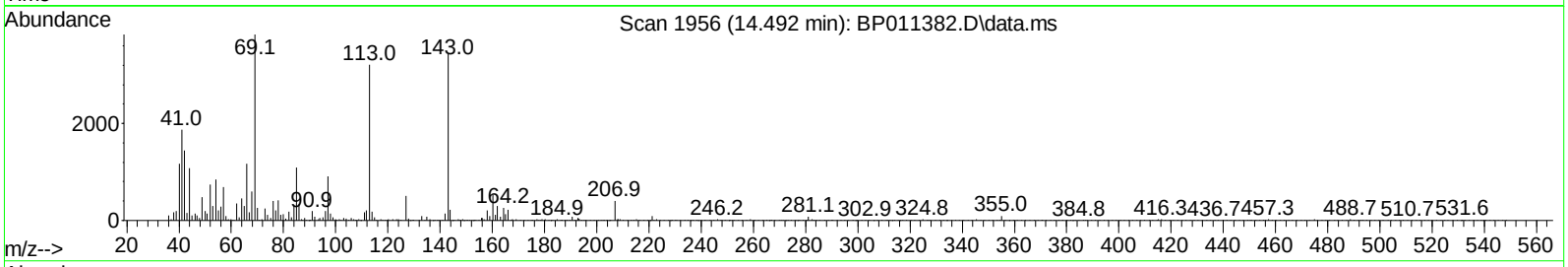

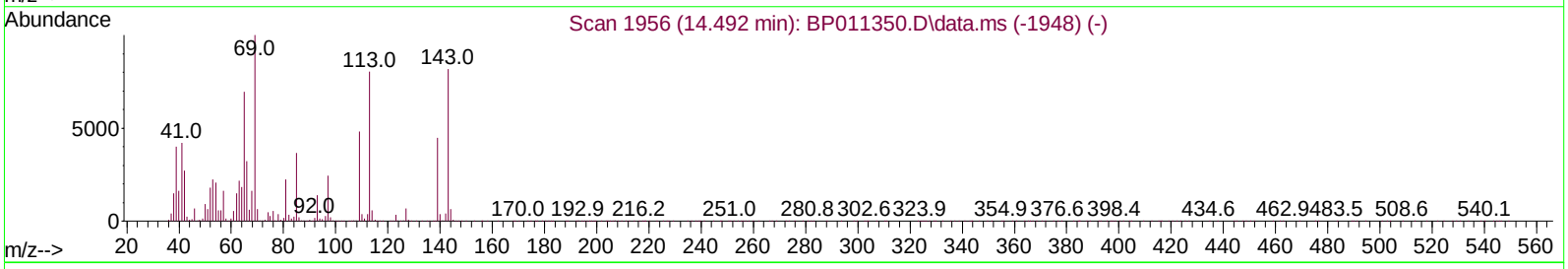

TIC: BP011382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 2.63 ng/ul

response 6499

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.50#
41.00	54.40	54.66
42.00	35.50	41.96

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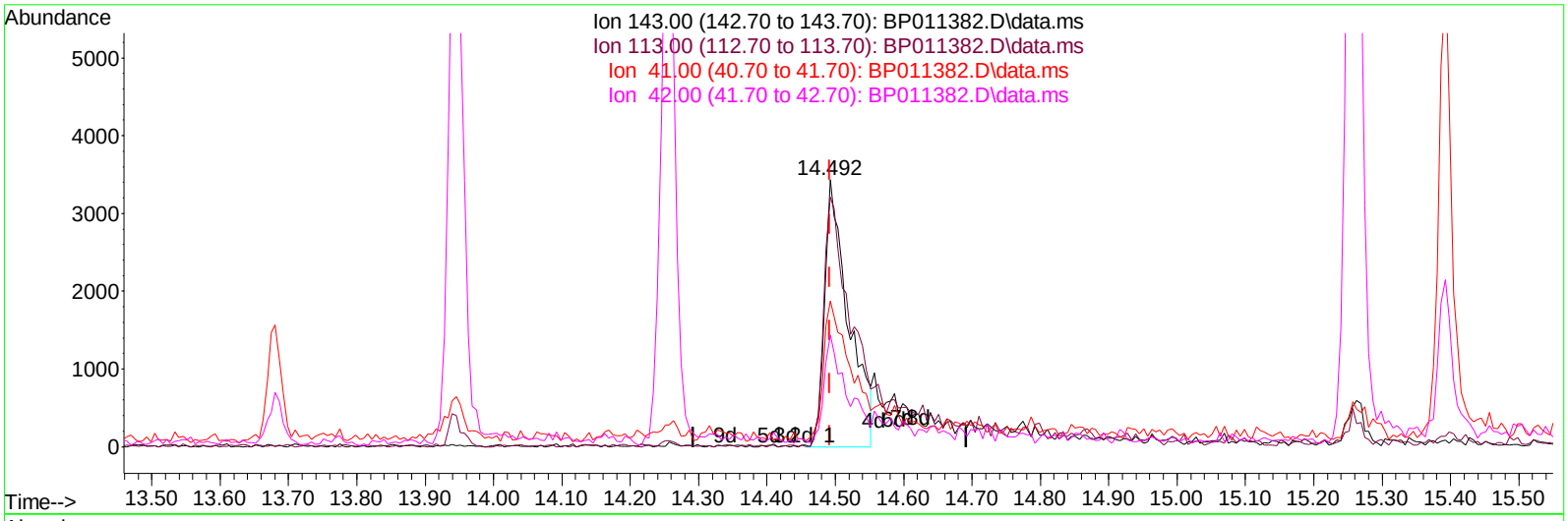

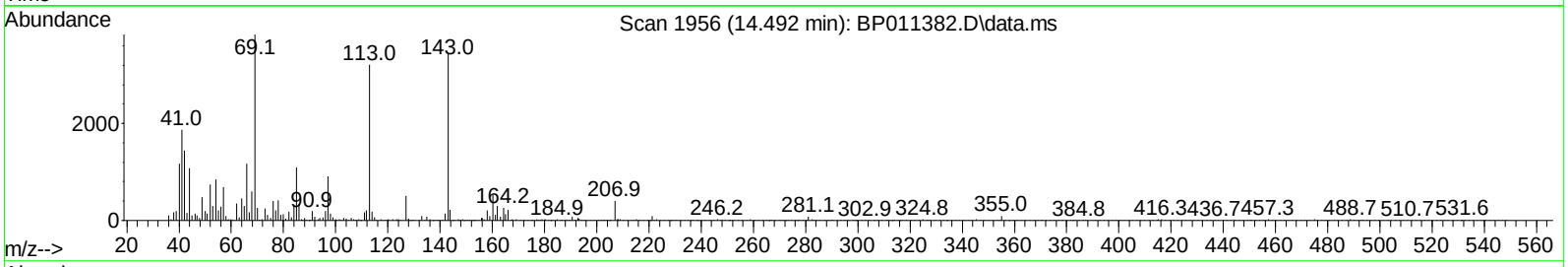

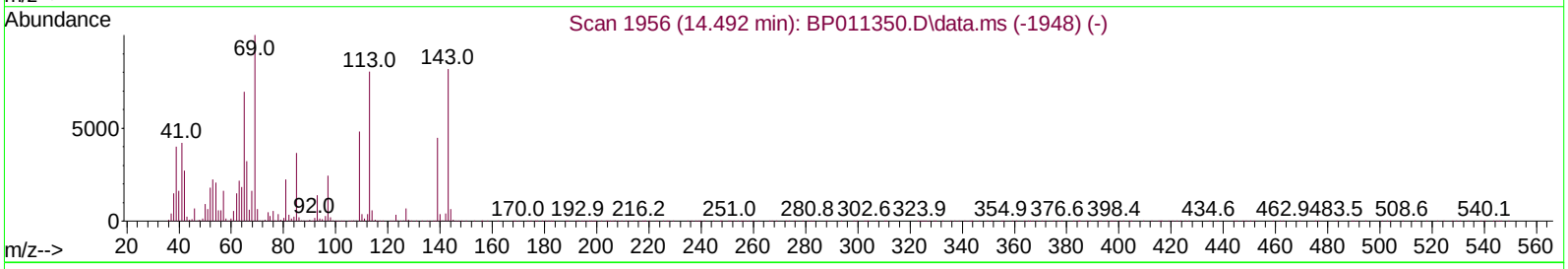

TIC: BP011382.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 3.37 ng/ul m

response 8351

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.50#
41.00	54.40	54.66
42.00	35.50	41.96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	83701	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.369	136	345597	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	195277	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	406966	20.000	ng/ul	-0.01
79) Chrysene-d12	21.163	240	398587	20.000	ng/ul	# 0.00
88) Perylene-d12	23.468	264	426532	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11644m	4.265	ng/uL	0.00
4) Pyridine-d5	3.499	84	38882	5.646	ng/ul	0.00
7) Phenol-d5	6.763	99	49196	6.642	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	149640	28.491	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.110	132	139816	24.057	ng/ul	-0.01
15) 4-Methylphenol-d8	8.299	113	89186	15.334	ng/ul	-0.02
21) Nitrobenzene-d5	8.746	128	80385	34.177	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	77835	36.145	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.998	165	122607	28.194	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.522	131	162033	21.313	ng/ul	-0.01
46) Dimethylphthalate-d6	13.681	166	488748	36.072	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	549061	33.664	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	8351m	3.375	ng/ul	0.00
60) Fluorene-d10	15.257	176	407148	35.137	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.392	200	60802	32.165	ng/ul	0.00
73) Anthracene-d10	17.128	188	713290	38.994	ng/ul	-0.01
81) Pyrene-d10	19.392	212	829677	37.782	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	836131	38.871	ng/ul	0.00
Target Compounds						Qvalue

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

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