Quantitation Report (Qedit)

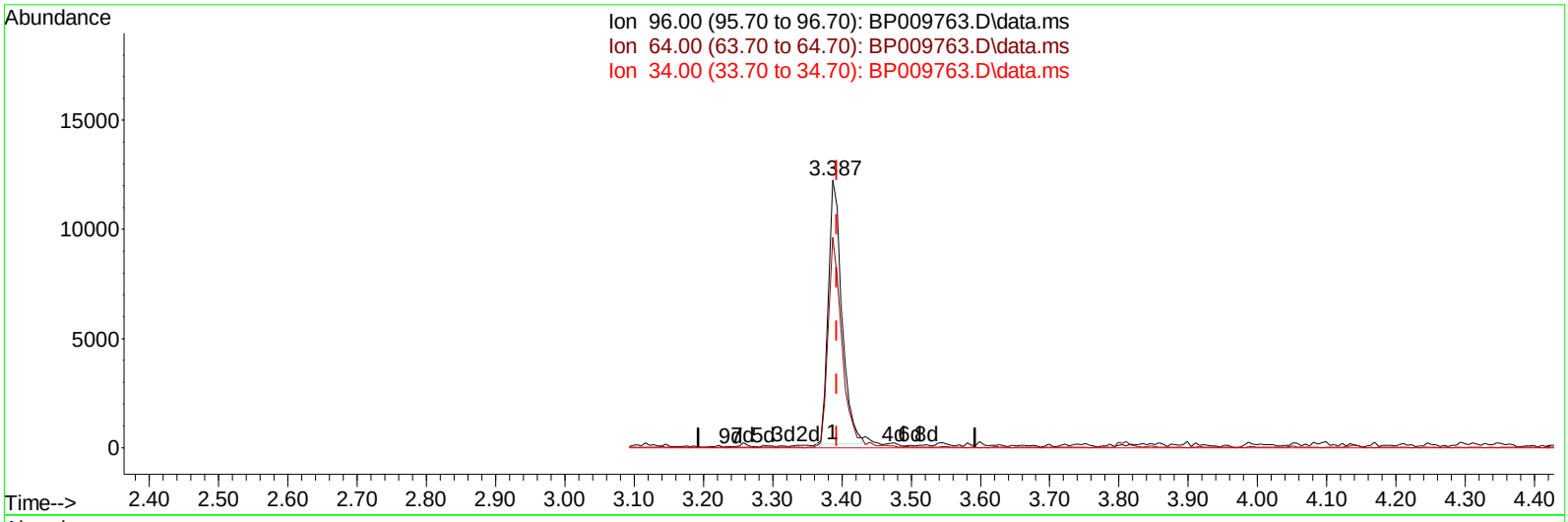
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009763.D
Acq On : 11 Apr 2022 22:08
Operator : CG/JU
Sample : N2338-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

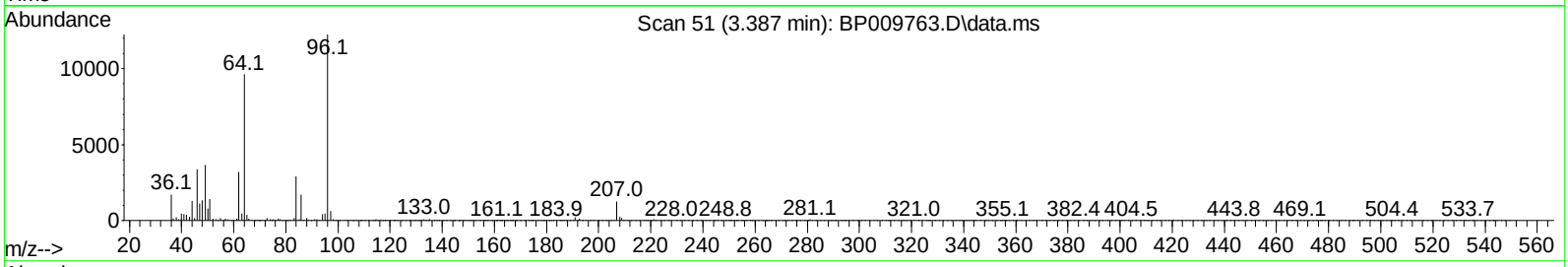
Quant Time: Apr 12 01:53:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 07 19:16:52 2022
Response via : Initial Calibration

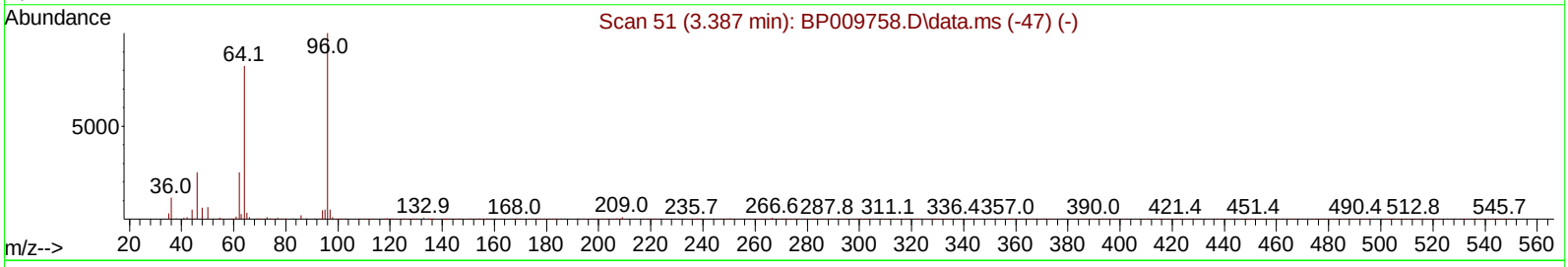
Instrument :
BNA_P
ClientSampleId :
C0J68

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/13/2022







TIC: BP009763.D\data.ms

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

3.387min (-0.006) 5.88 ng/uL

response 16477

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.57#
34.00	0.00	0.00
0.00	0.00	0.00

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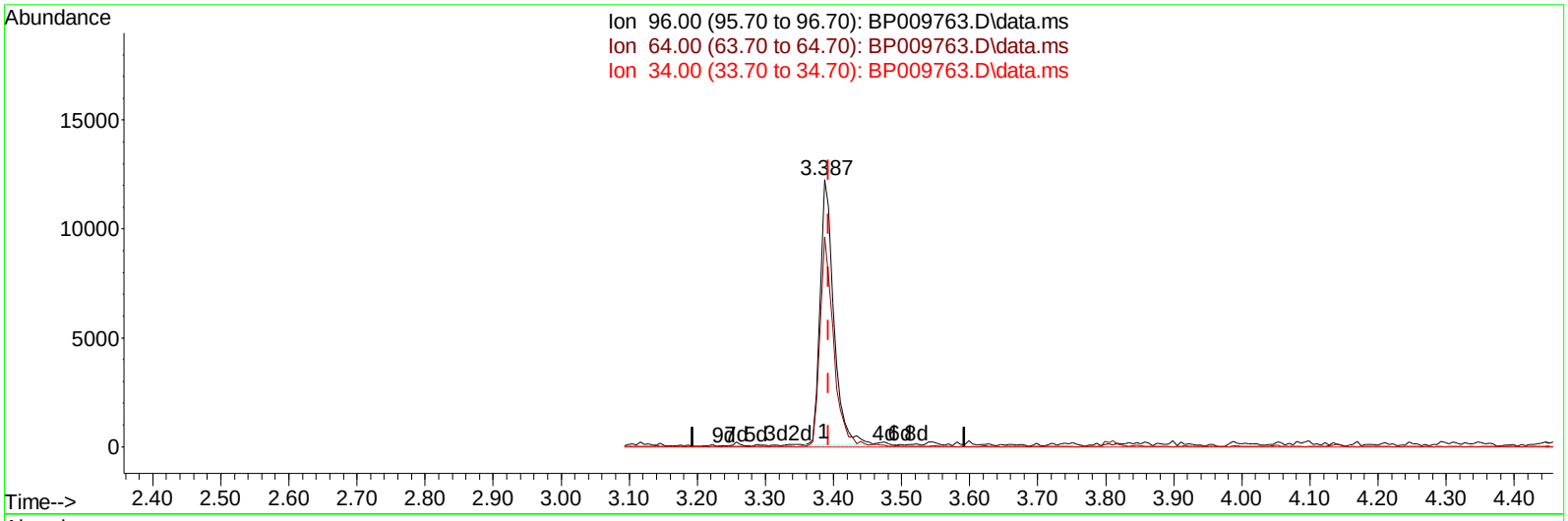

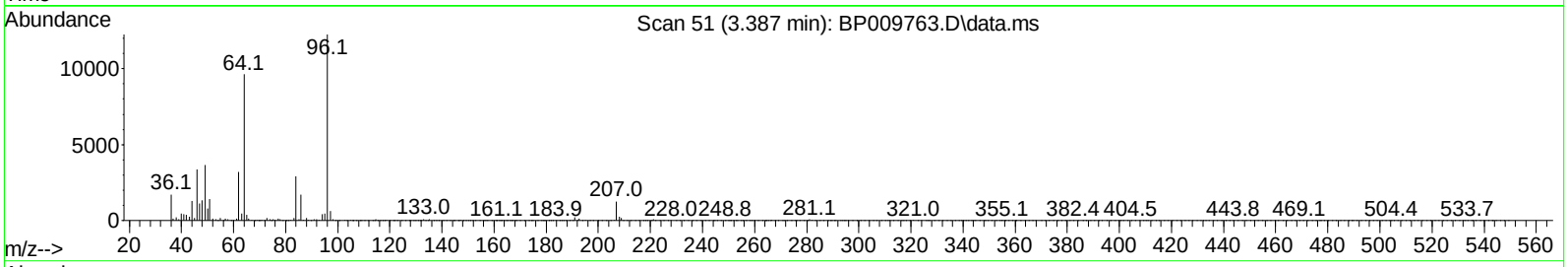

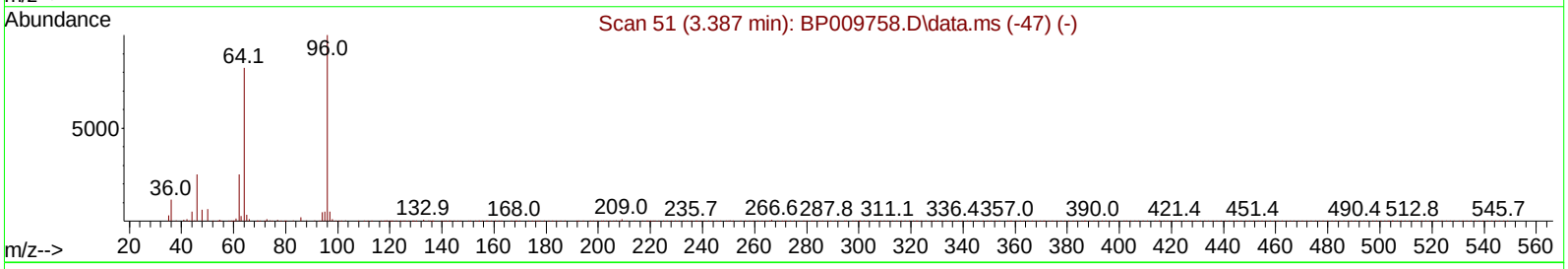

TIC: BP009763.D\data.ms

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

3.387min (-0.006) 6.15 ng/uL m

response 17228

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.57#
34.00	0.00	0.00
0.00	0.00	0.00

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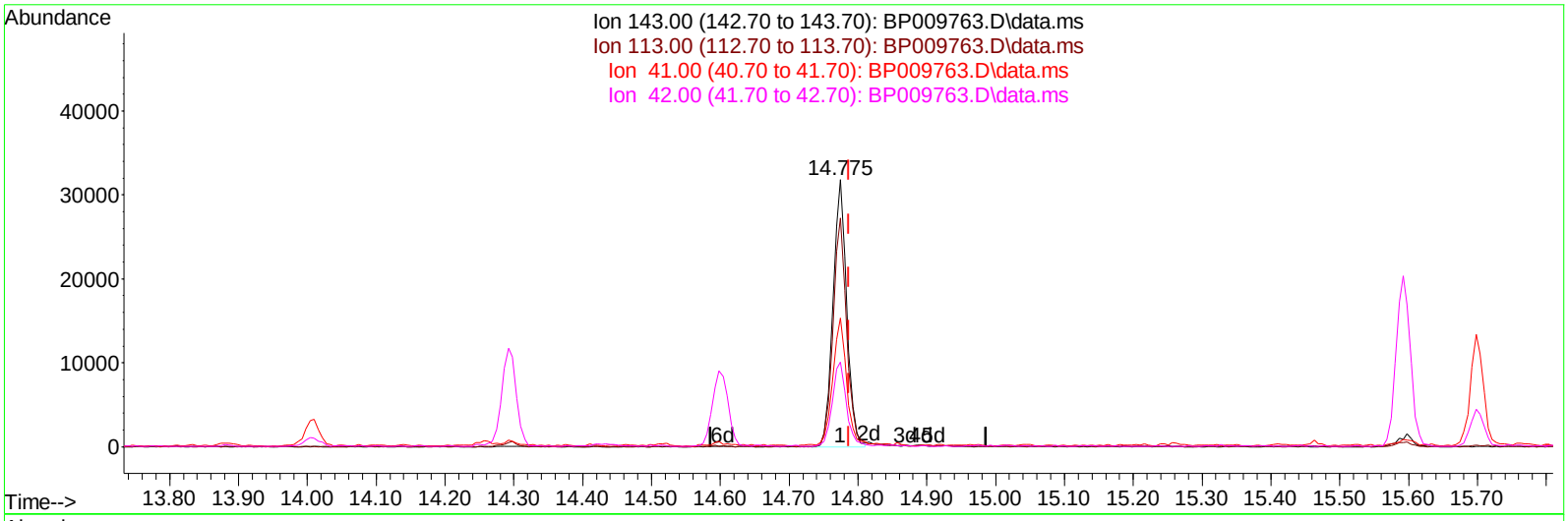

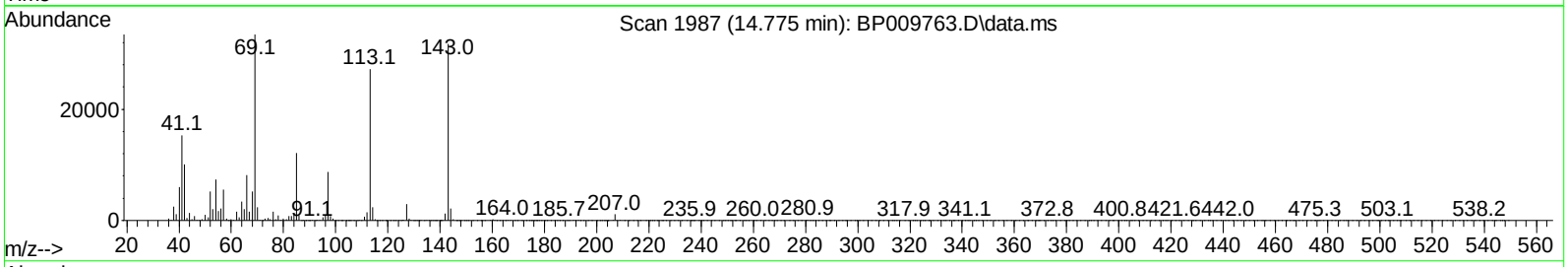

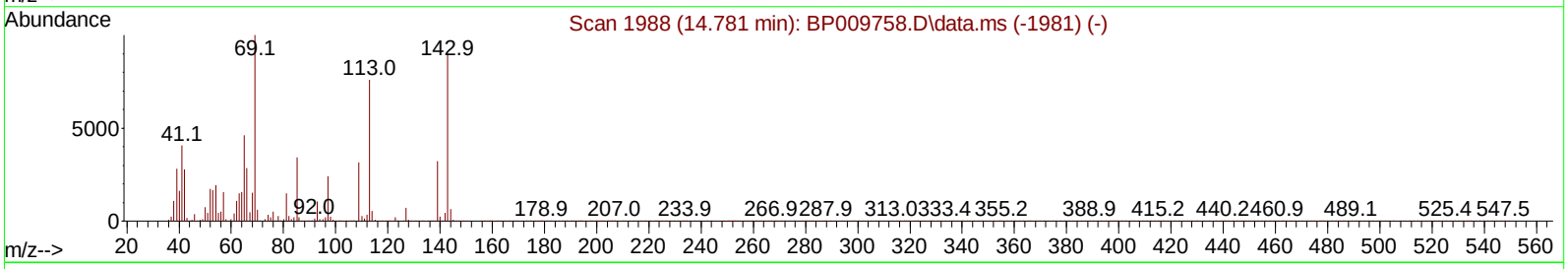

TIC: BP009763.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 7.70 ng/ul

response 43443

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.74#
41.00	23.20	48.26#
42.00	18.00	31.79#

Quantitation Report (Qedit)

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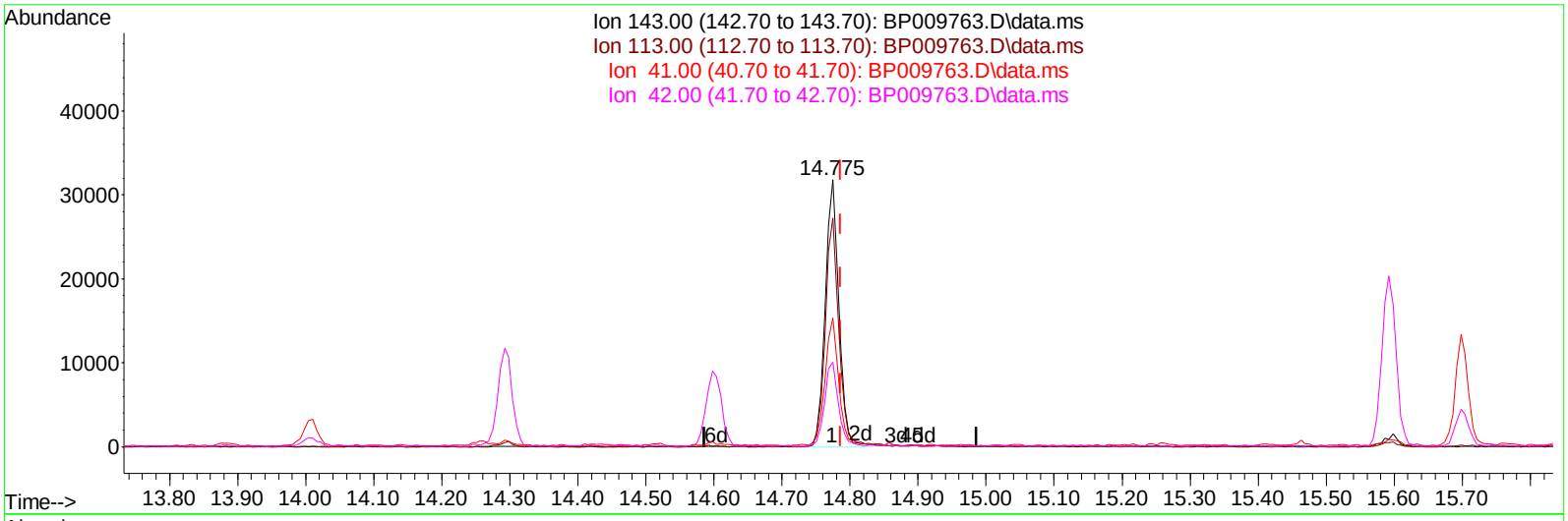

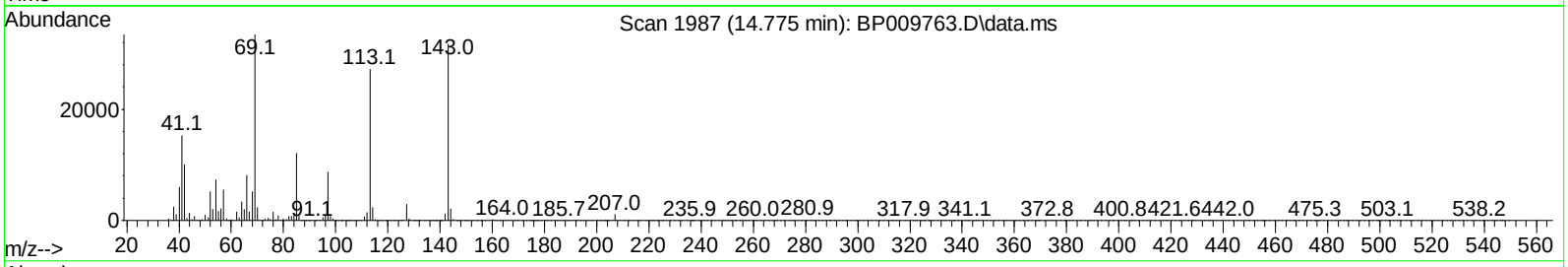

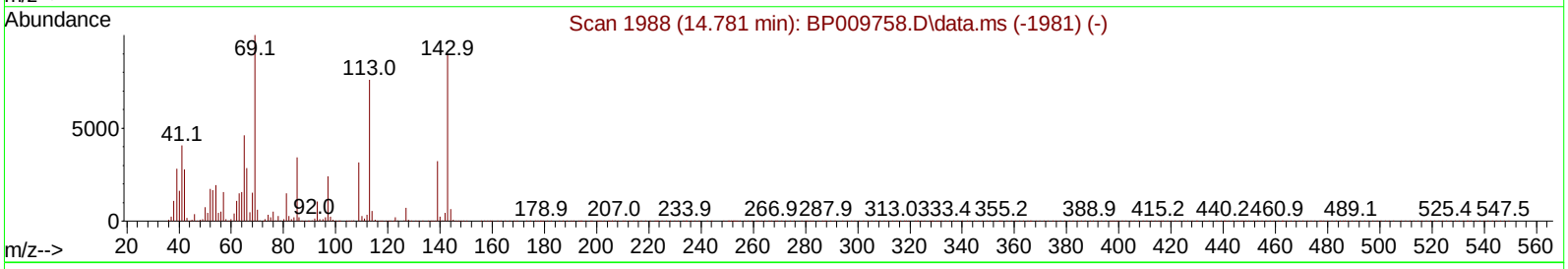

TIC: BP009763.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 7.79 ng/ul m

response 43984

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.74#
41.00	23.20	48.26#
42.00	18.00	31.79#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	124993	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	557685	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.598	164	396201	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	886114	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	926256	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	867044	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17228m	6.151	ng/uL	0.00
4) Pyridine-d5	3.805	84	91812	11.890	ng/ul	0.00
7) Phenol-d5	7.122	99	77810	7.091	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	229859	35.185	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.498	132	216440	26.148	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	154467	16.862	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	146240	36.363	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.851	143	154451	35.030	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.392	165	265149	28.738	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.904	131	387402	29.481	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	1156404	37.157	ng/ul	-0.01
49) Acenaphthylene-d8	14.292	160	1235215	33.501	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	43984m	7.793	ng/ul	-0.01
60) Fluorene-d10	15.592	176	1002542	38.293	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	186904	35.605	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1736222	40.923	ng/ul	-0.01
81) Pyrene-d10	19.680	212	2081194	41.140	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	1900141	42.083	ng/ul	-0.02

Target Compounds Qvalue

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

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