Quantitation Report (Qedit)

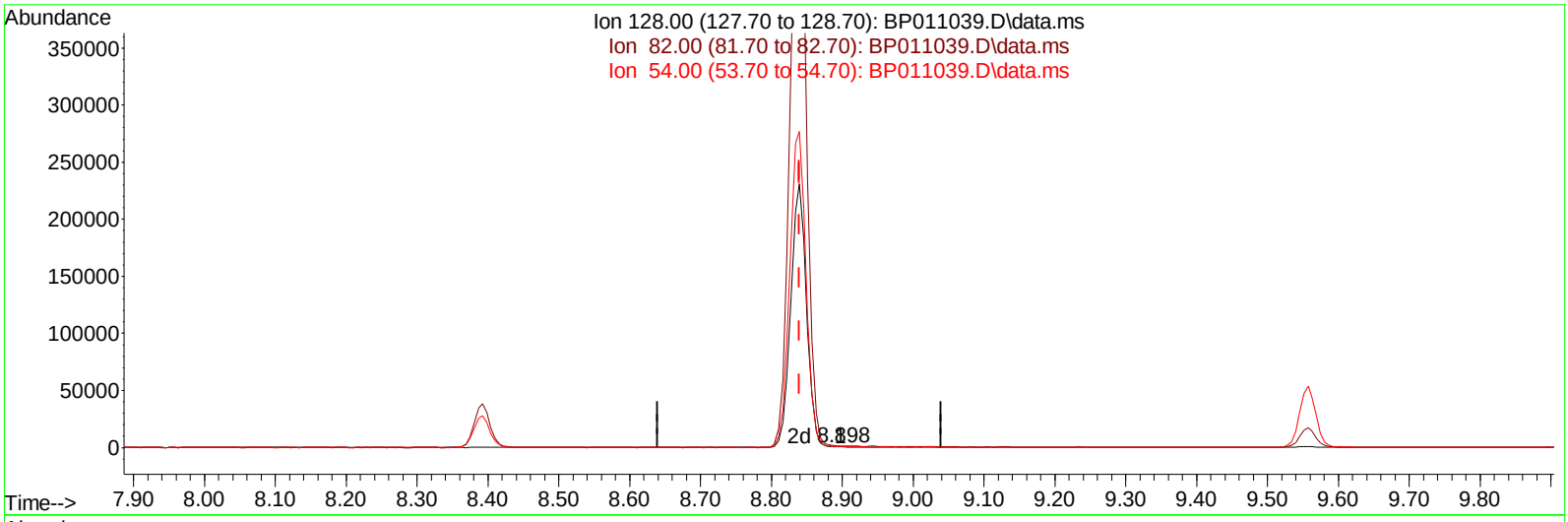
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011039.D
Acq On : 19 Jul 2022 19:00
Operator : CG/JU
Sample : N3753-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

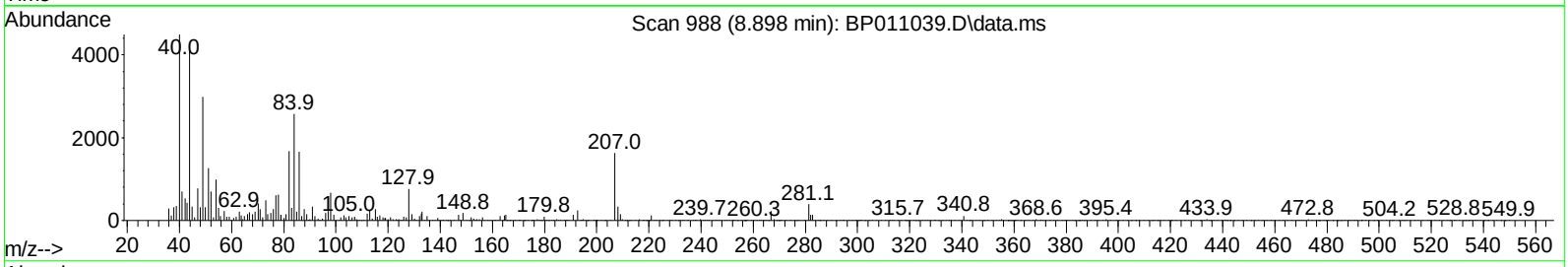
Quant Time: Jul 19 23:41:29 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 19 03:25:54 2022
Response via : Initial Calibration

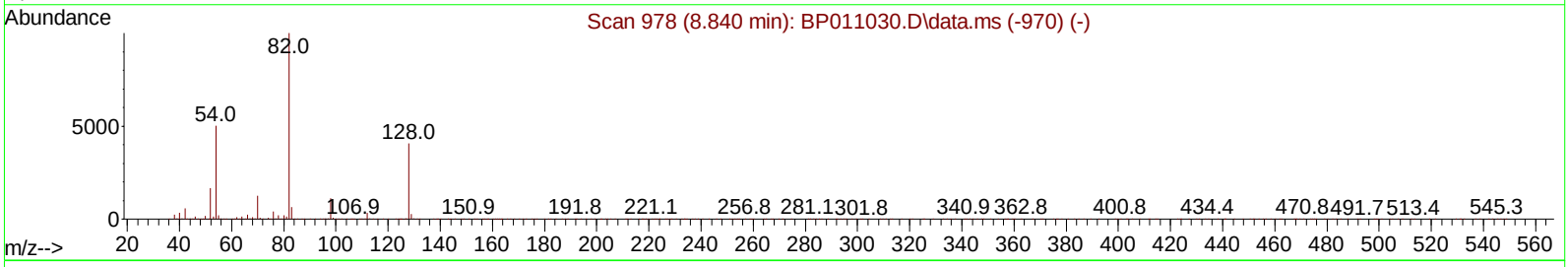
Instrument :
BNA_P
ClientSampleId :
GBHQ2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/20/2022
Supervised By :mohammad ahmed 07/21/2022







TIC: BP011039.D\data.ms

(21) Nitrobenzene-d5 (S)

8.898min (+ 0.059) 0.04 ng/ul

response 425

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	219.95
54.00	155.90	129.27
0.00	0.00	0.00

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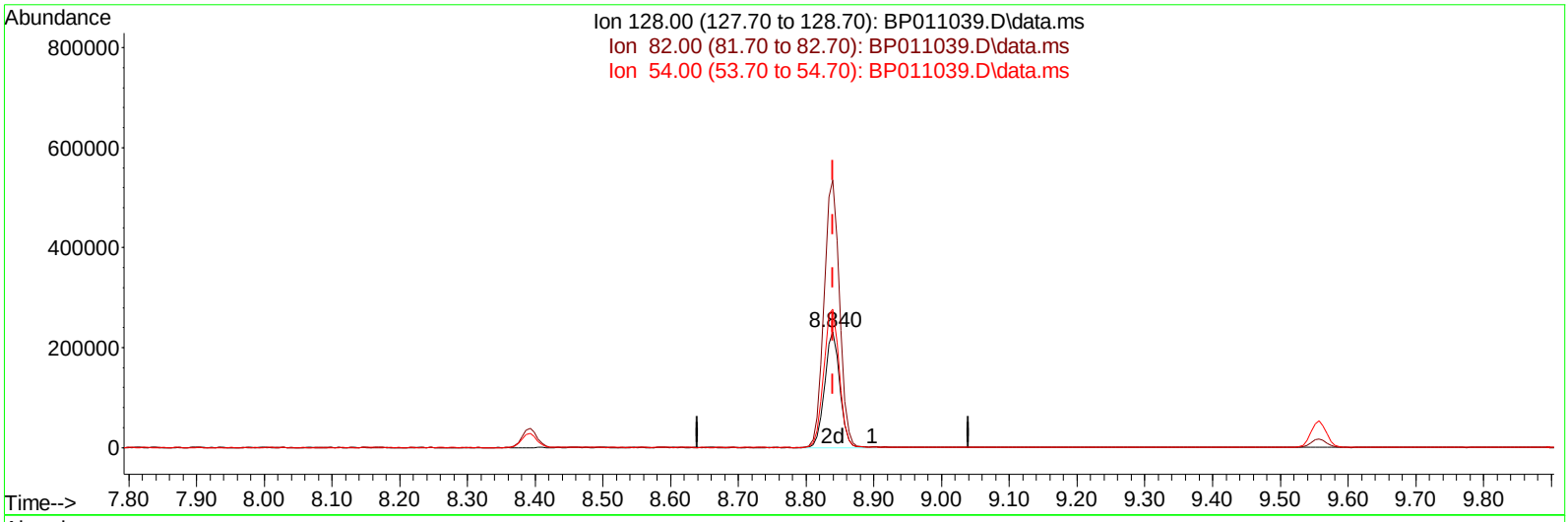

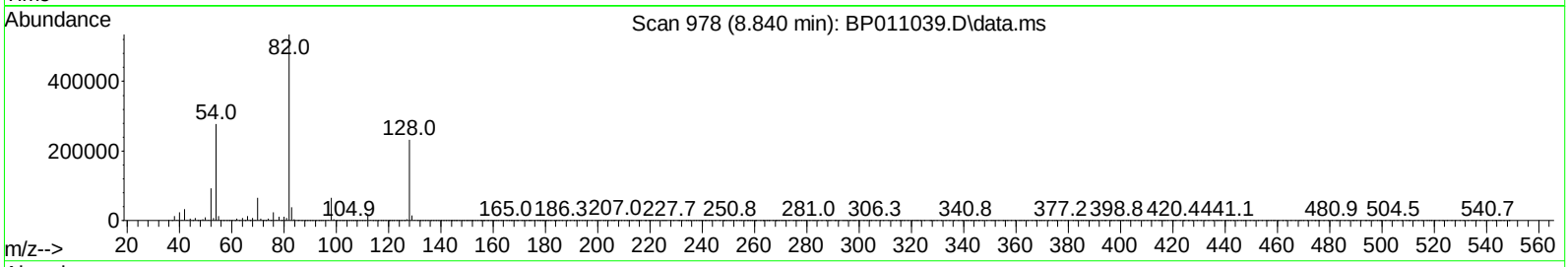

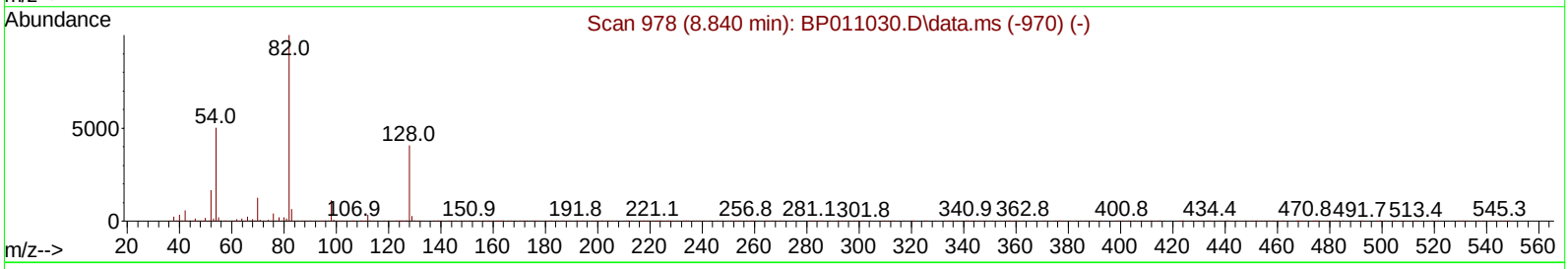

TIC: BP011039.D\data.ms

(21) Nitrobenzene-d5 (S)

8.840min (-0.000) 37.49 ng/ul m

response 359121

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	249.70	231.23
54.00	155.90	119.81#
0.00	0.00	0.00

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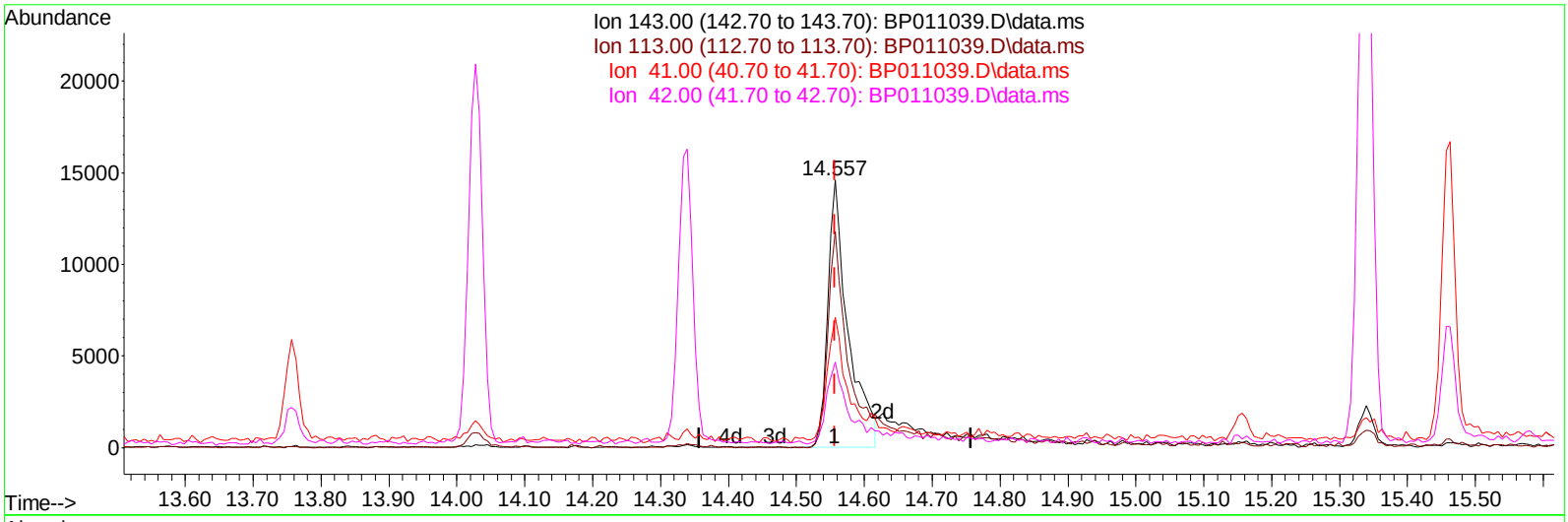

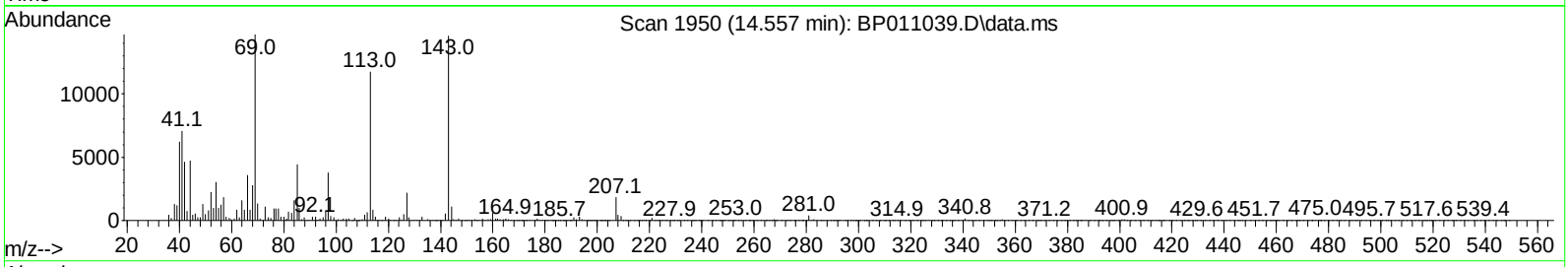

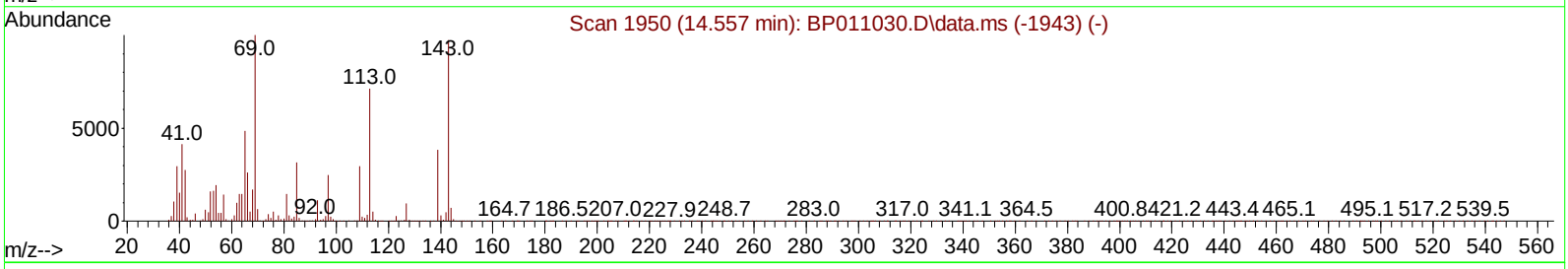

TIC: BP011039.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.000) 3.00 ng/ul

response 30172

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	80.68
41.00	59.00	48.61
42.00	41.10	31.93#

Quantitation Report (Qedit)

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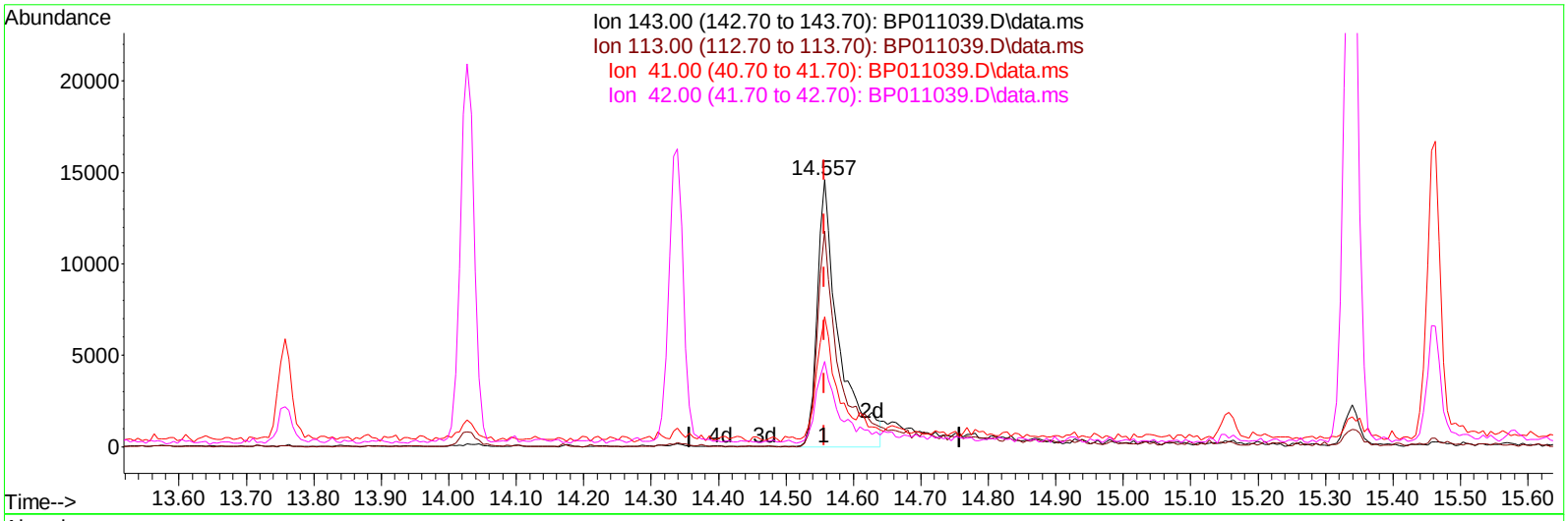

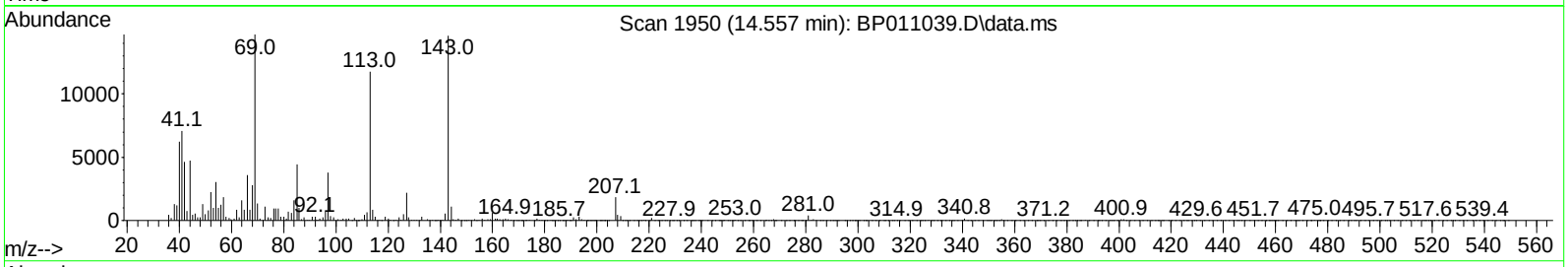

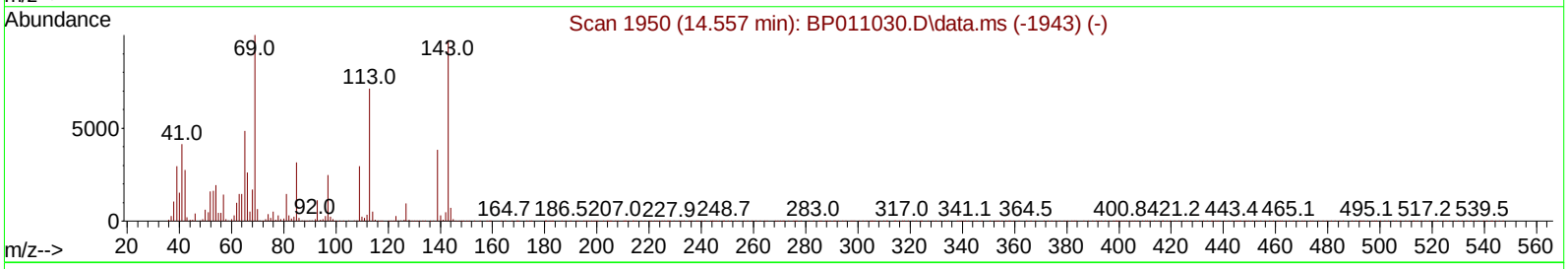

TIC: BP011039.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.000) 3.23 ng/ul m

response 32538

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	80.68
41.00	59.00	48.61
42.00	41.10	31.93#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	331347	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	1314607	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.339	164	727148	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1451798	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1416556	20.000	ng/ul	# 0.00
88) Perylene-d12	23.580	264	1440190	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	40541	4.411	ng/uL	0.00
4) Pyridine-d5	3.593	84	266754	11.229	ng/ul	0.00
7) Phenol-d5	6.857	99	162109	5.960	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	609083	34.825	ng/ul	0.00
11) 2-Chlorophenol-d4	7.210	132	555492	25.275	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	322025	14.789	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	359121m	37.486	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	313341	33.054	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.092	165	516616	29.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	841089	29.429	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	2017173	40.589	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160	2367657	40.378	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	32538m	3.235	ng/ul	0.00
60) Fluorene-d10	15.339	176	1741686	41.230	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	207432	28.874	ng/ul	0.00
73) Anthracene-d10	17.204	188	2827905	45.819	ng/ul	0.00
81) Pyrene-d10	19.463	212	3262096	43.847	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.427	264	3220386	46.557	ng/ul	0.00
Target Compounds						Qvalue

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

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