Quantitation Report (Qedit)

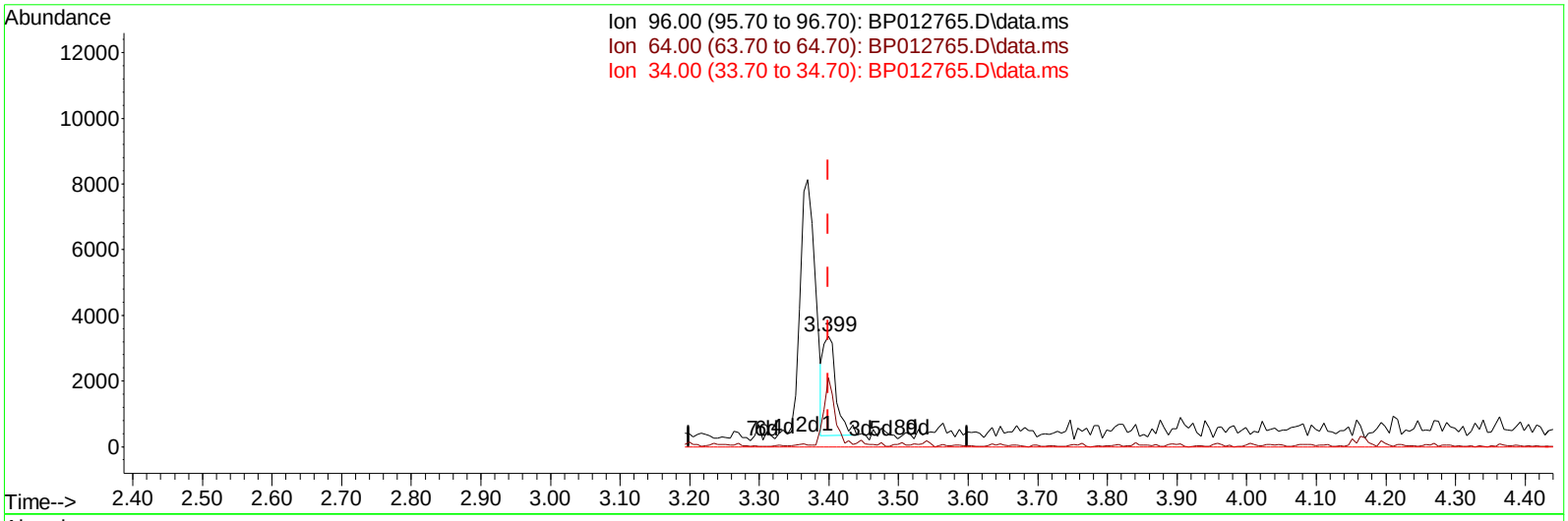
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
Data File : BP012765.D
Acq On : 26 Nov 2022 13:36
Operator : CG/JU
Sample : N5592-05DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

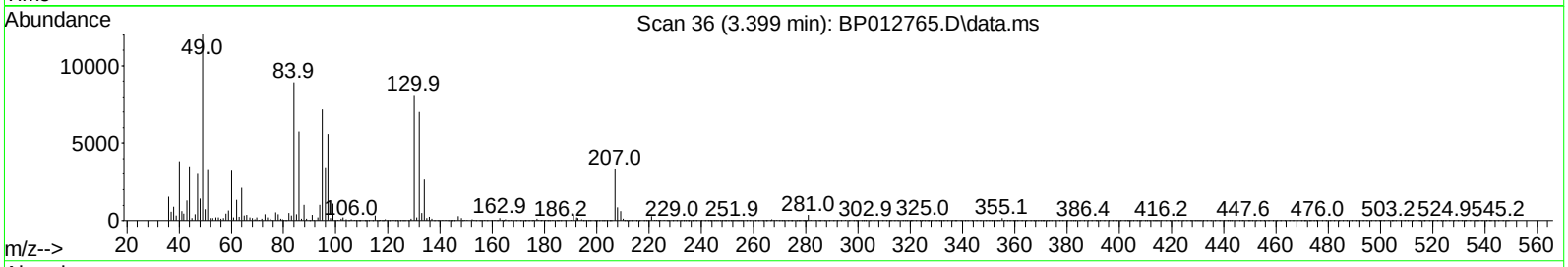
Quant Time: Nov 27 21:51:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 25 15:59:22 2022
Response via : Initial Calibration

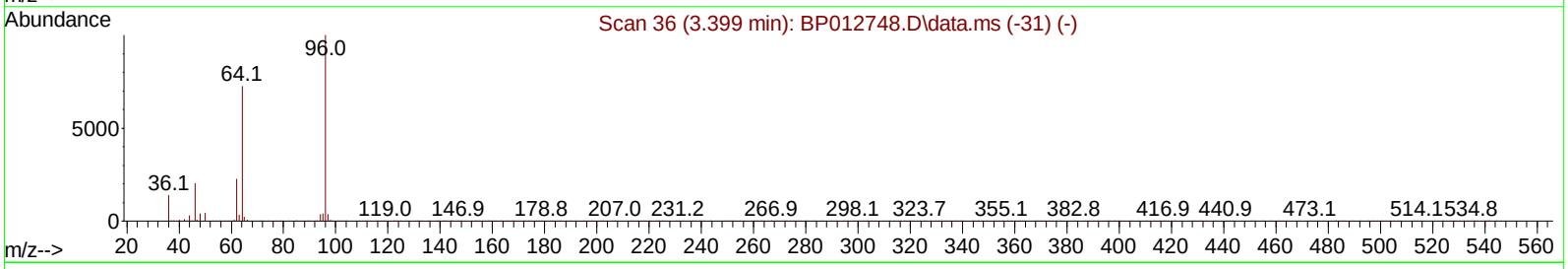
Instrument :
BNA_P
ClientSampleId :
H0AB0DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/28/2022
Supervised By :mohammad ahmed 11/30/2022







TIC: BP012765.D\data.ms

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

3.399min (+ 0.000) 0.33 ng/uL

response 3912

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	62.38
34.00	0.00	0.00
0.00	0.00	0.00

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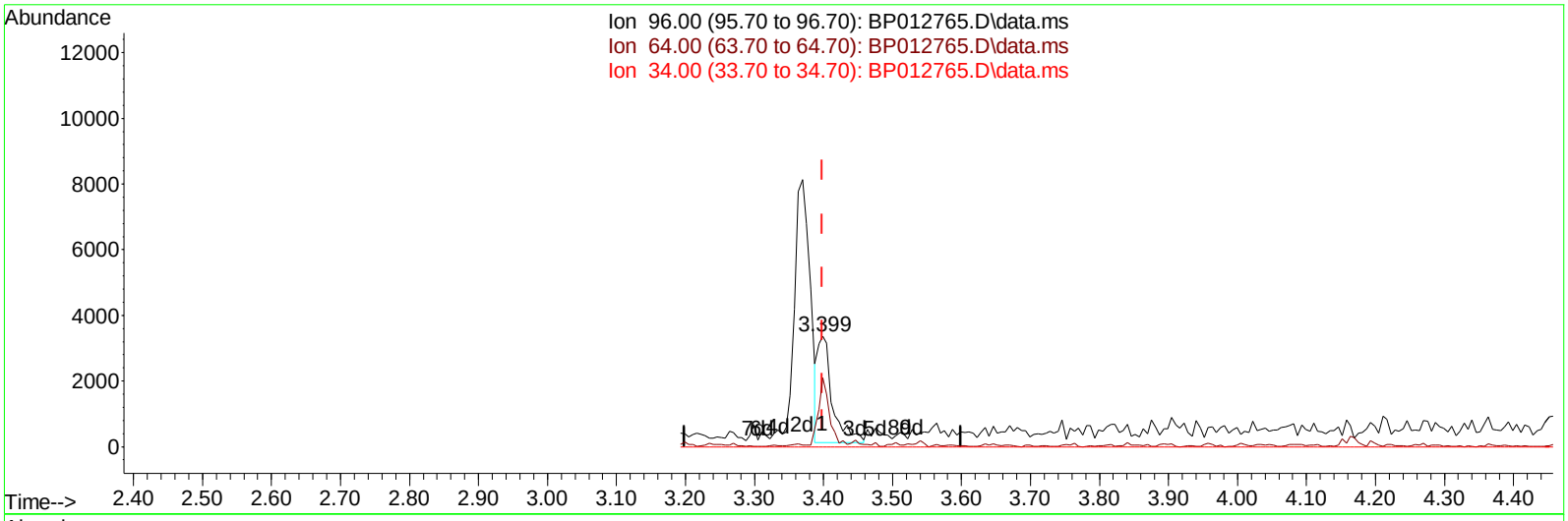

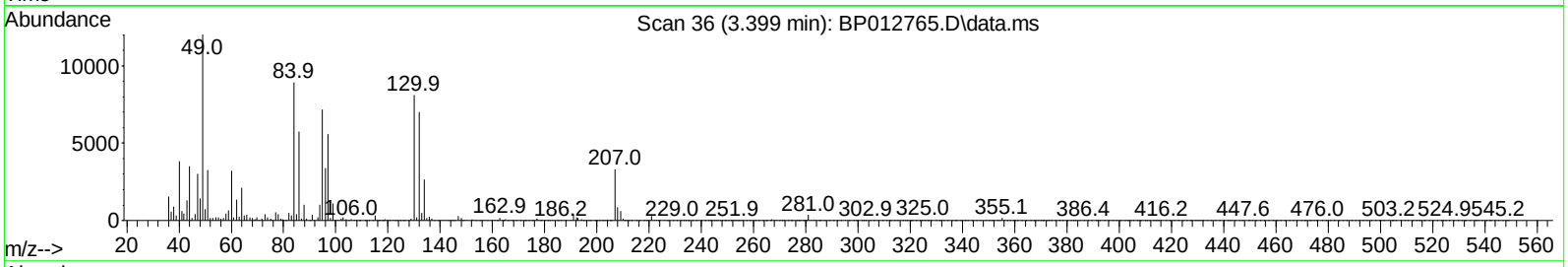

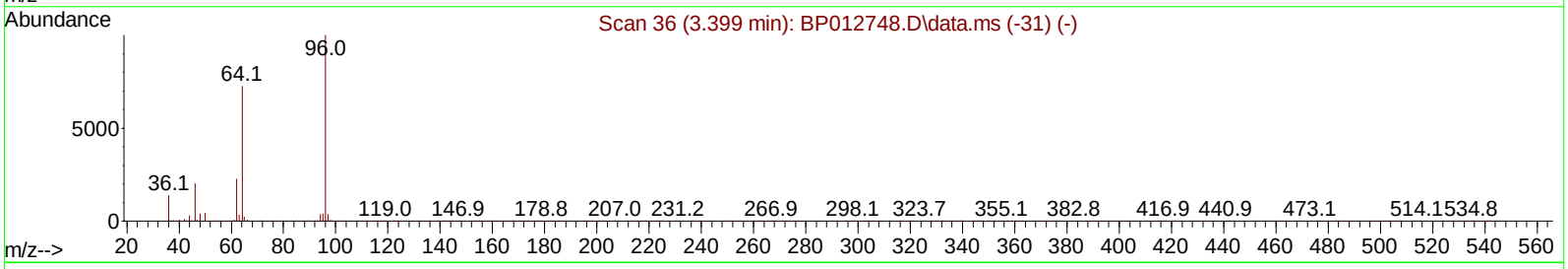

TIC: BP012765.D\data.ms

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

3.399min (+ 0.000) 0.40 ng/uL m

response 4778

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	62.38
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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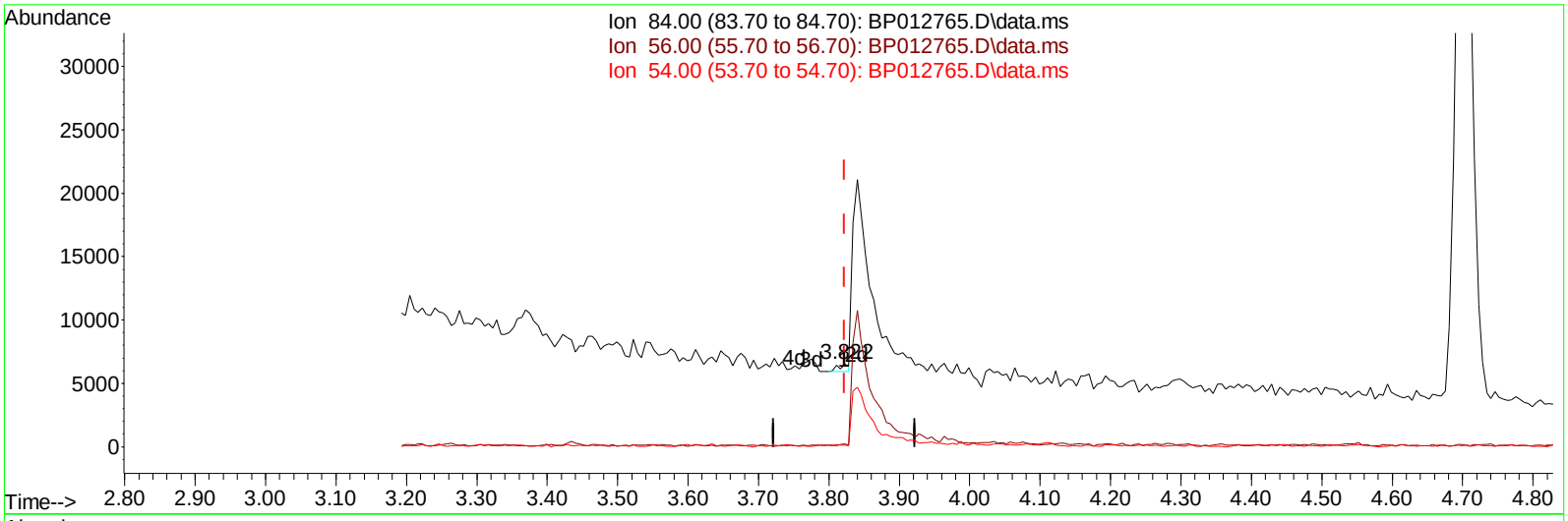

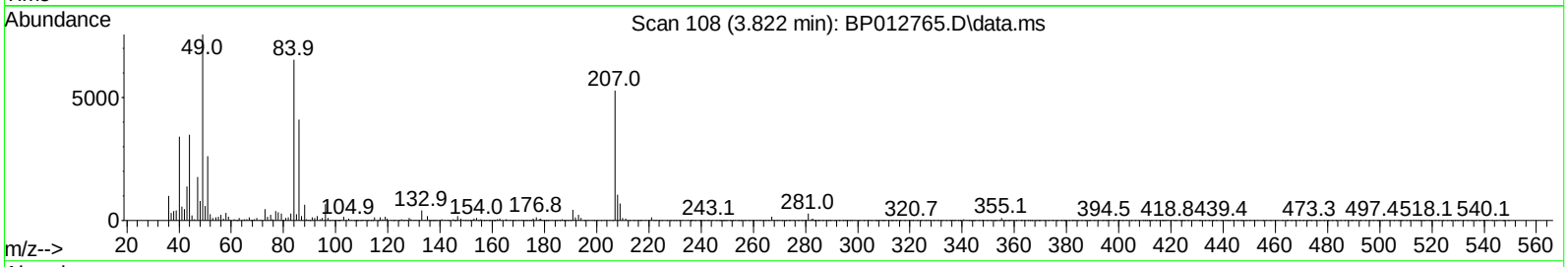

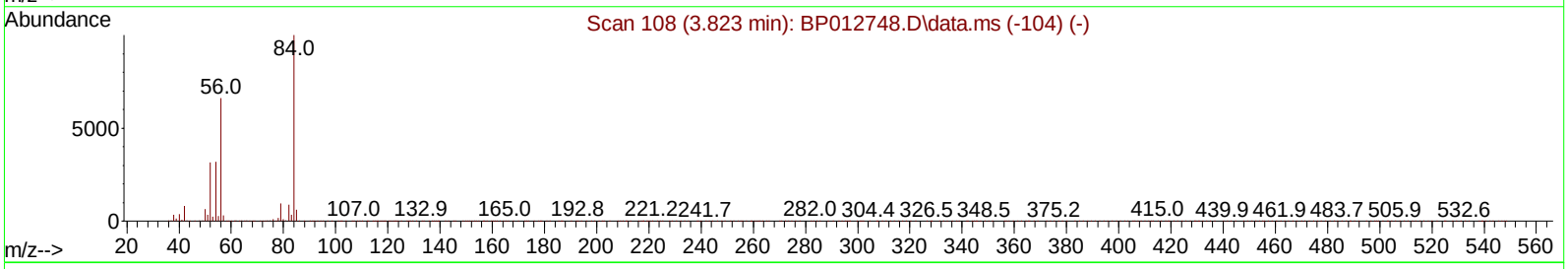

TIC: BP012765.D\data.ms

(4) Pyridine-d5 (S)

3.822min (+ 0.000) 0.02 ng/ul

response 659

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	3.65#
54.00	31.30	2.09#
0.00	0.00	0.00

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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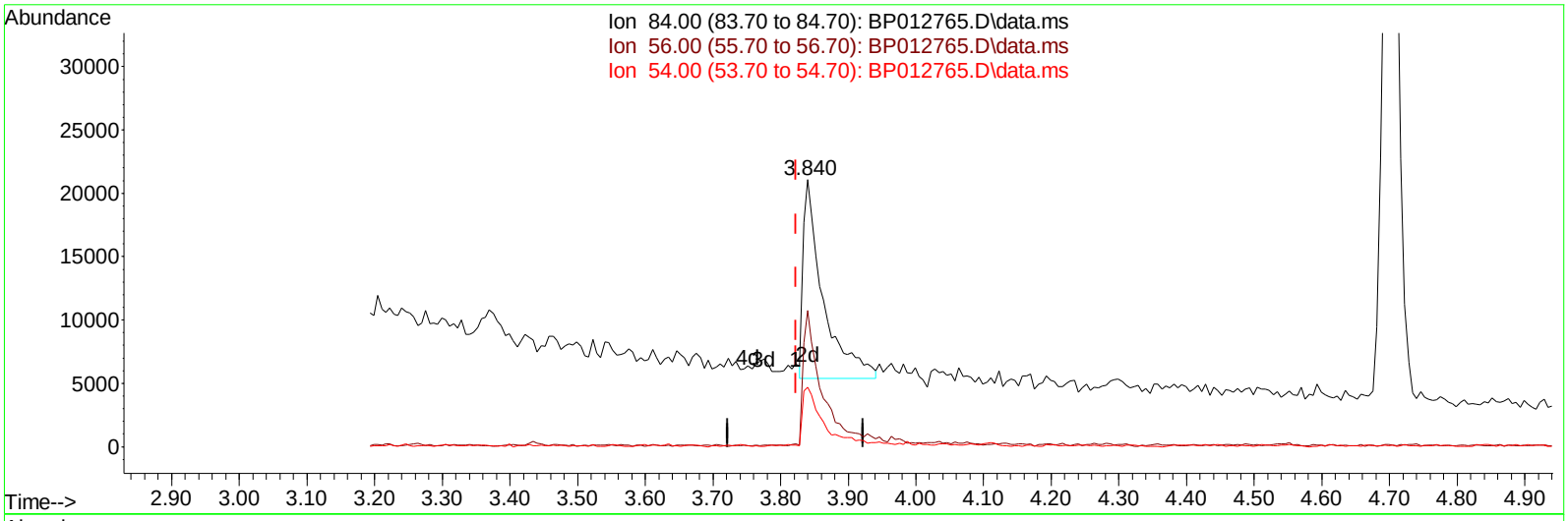

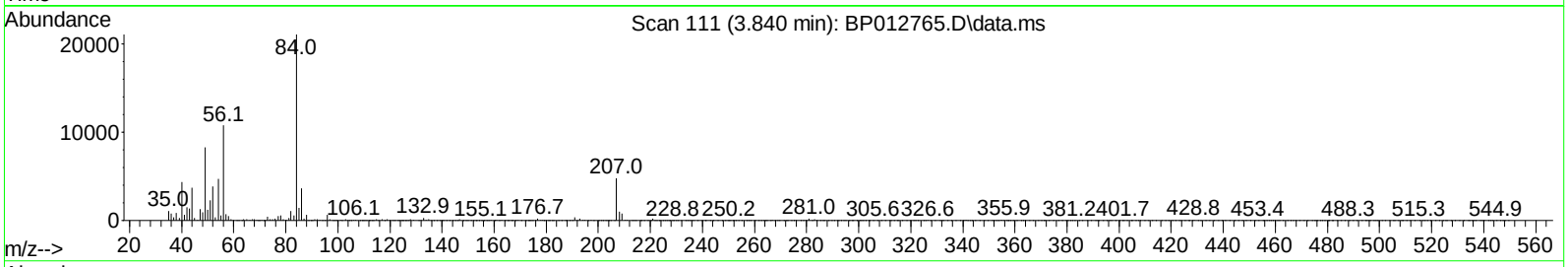

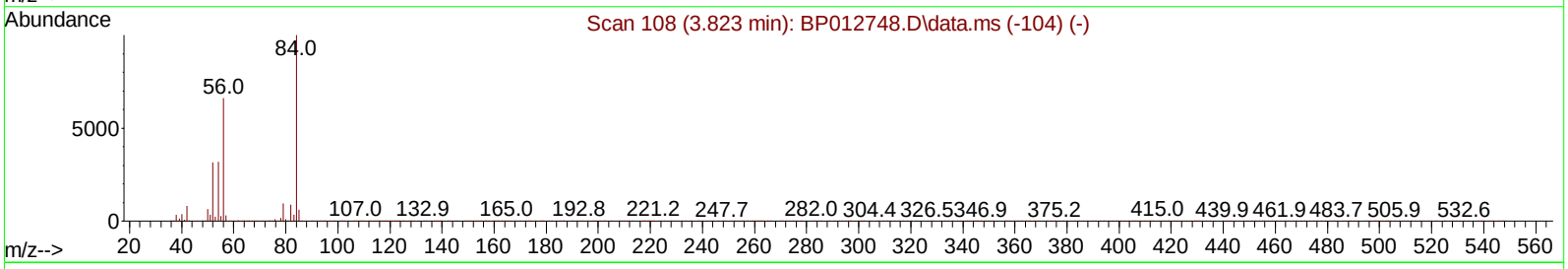

TIC: BP012765.D\data.ms

(4) Pyridine-d5 (S)

3.840min (+ 0.018) 0.89 ng/ul m

response 31849

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	64.40	51.11#
54.00	31.30	22.35#
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/28/2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	511529	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	2232558	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.657	164	1562920	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	3433828	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	2829351	20.000	ng/ul	-0.01
88) Perylene-d12	24.010	264	2202474	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	4778m	0.400	ng/uL	0.00
4) Pyridine-d5	3.840	84	31849m	0.892	ng/ul	0.02
7) Phenol-d5	7.163	99	26380	0.628	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	68351	2.639	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	61855	1.931	ng/ul	0.00
15) 4-Methylphenol-d8	8.722	113	43812	1.294	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	27525	2.014	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.910	143	26405	1.820	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.452	165	71717	2.118	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	63459	1.369	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	363811	3.382	ng/ul	-0.01
49) Acenaphthylene-d8	14.351	160	325259	2.591	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.645	176	317187	3.367	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.510	188	539224	3.567	ng/ul	0.00
81) Pyrene-d10	19.733	212	612222	3.724	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	339242	3.136	ng/ul	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.434	88	156156	12.180	ng/uL	98

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

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