Quantitation Report (Qedit)

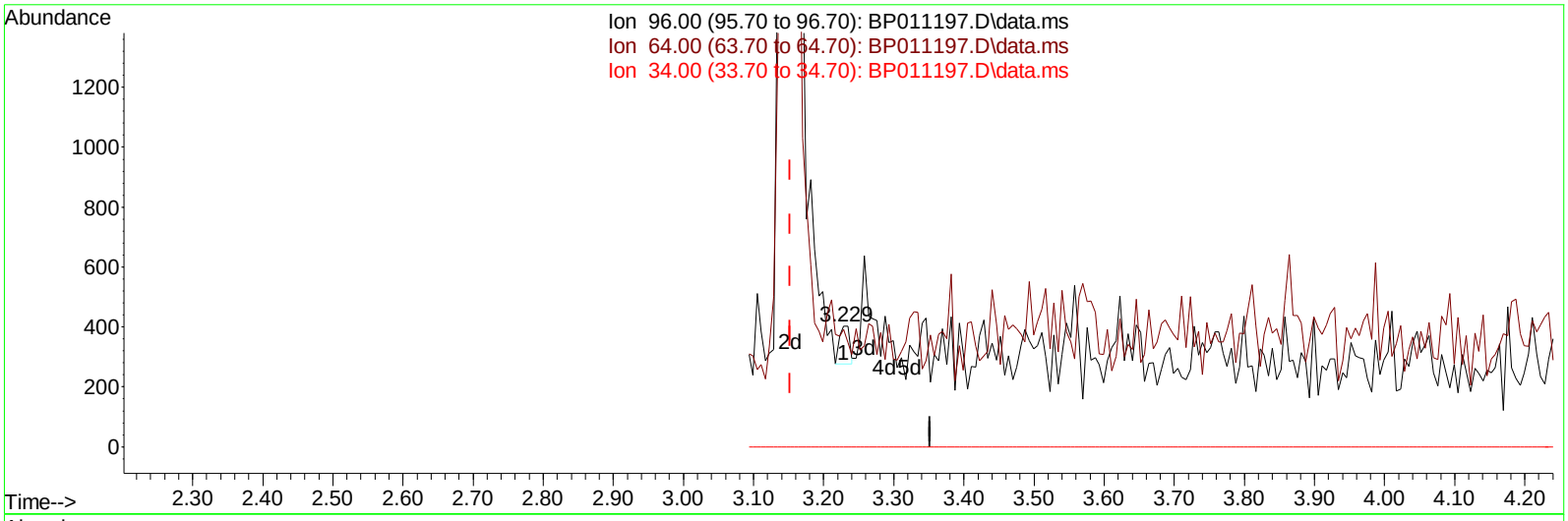
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
Data File : BP011197.D
Acq On : 01 Aug 2022 13:37
Operator : CG/JU
Sample : N3790-15
Misc :
ALS Vial : 6 Sample Multiplier: 1

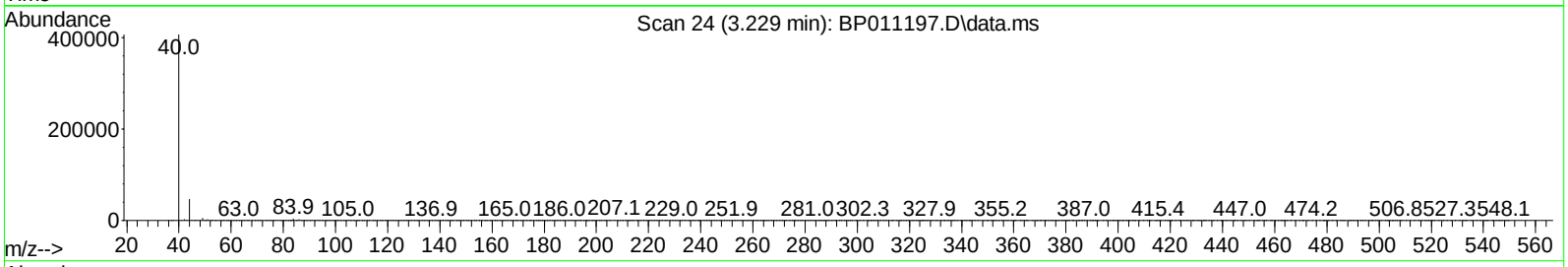
Quant Time: Aug 01 22:59:44 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 28 16:05:22 2022
Response via : Initial Calibration

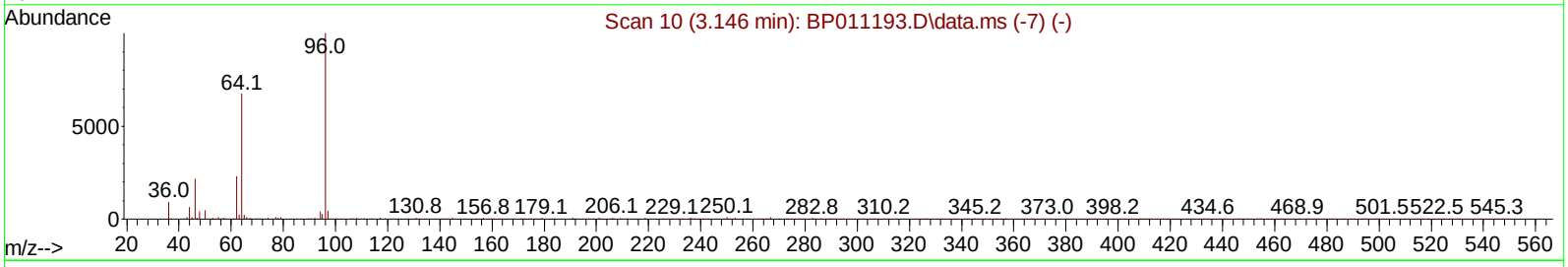
Instrument :
BNA_P
ClientSampleId :
DBUK9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/02/2022
Supervised By :mohammad ahmed 08/03/2022







TIC: BP011197.D\data.ms

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

3.229min (+ 0.077) 0.02 ng/uL

response 129

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	97.76
34.00	0.00	0.00
0.00	0.00	0.00

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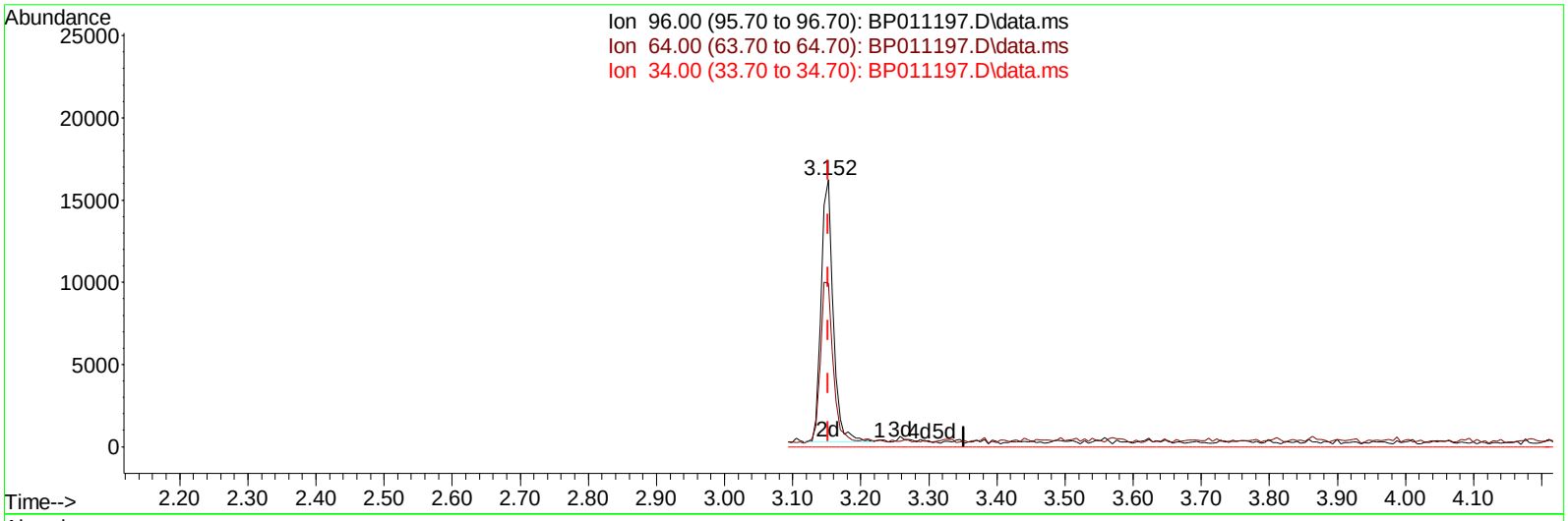

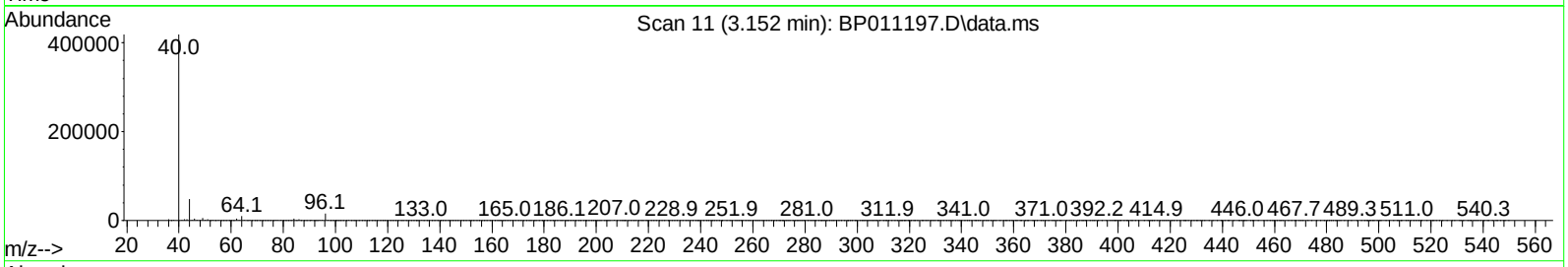

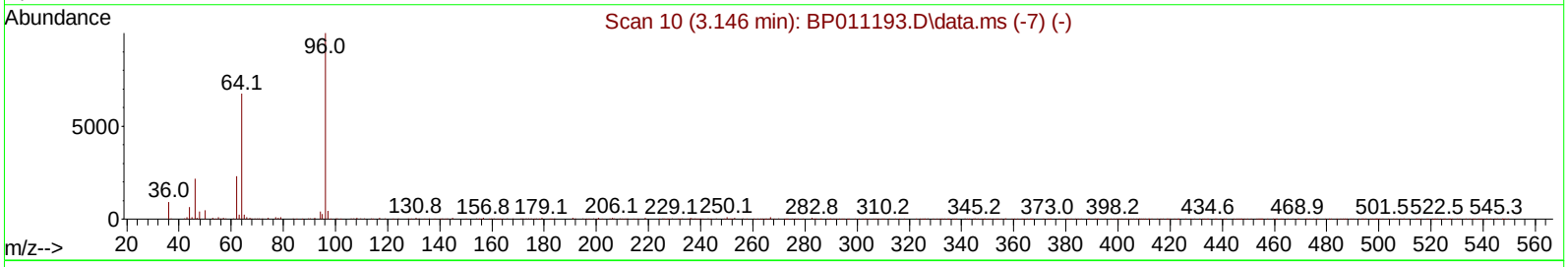

TIC: BP011197.D\data.ms

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

3.152min (+ 0.000) 2.42 ng/uL m

response 19648

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	61.14#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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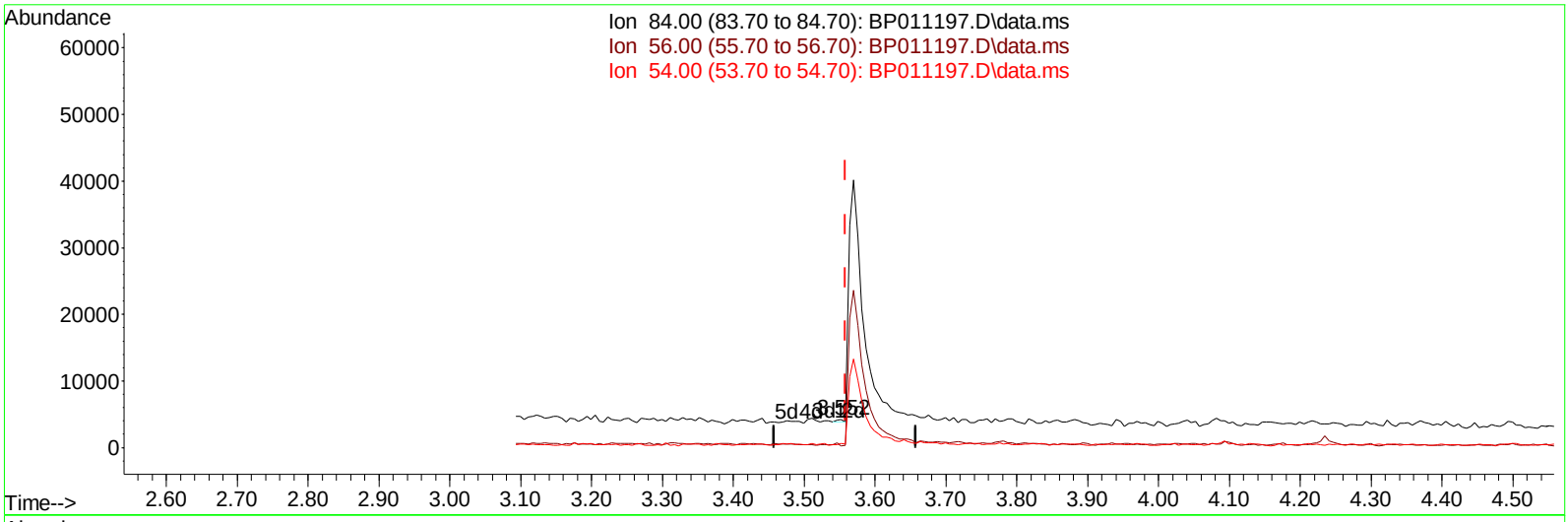

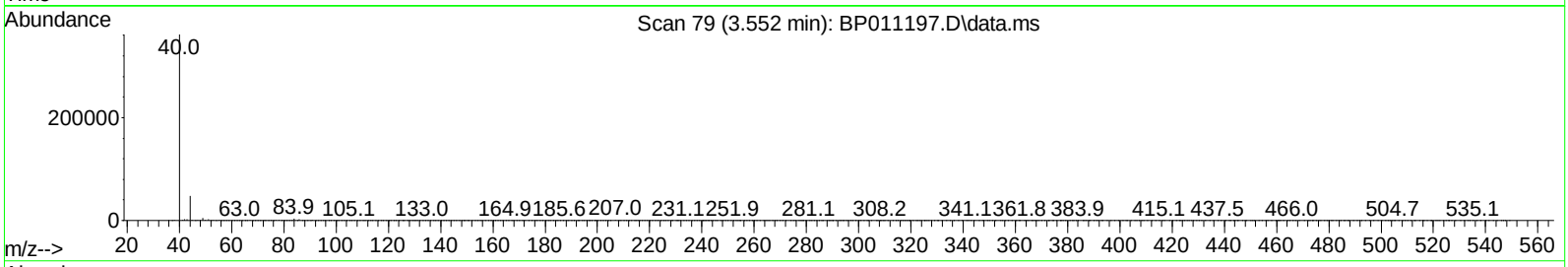

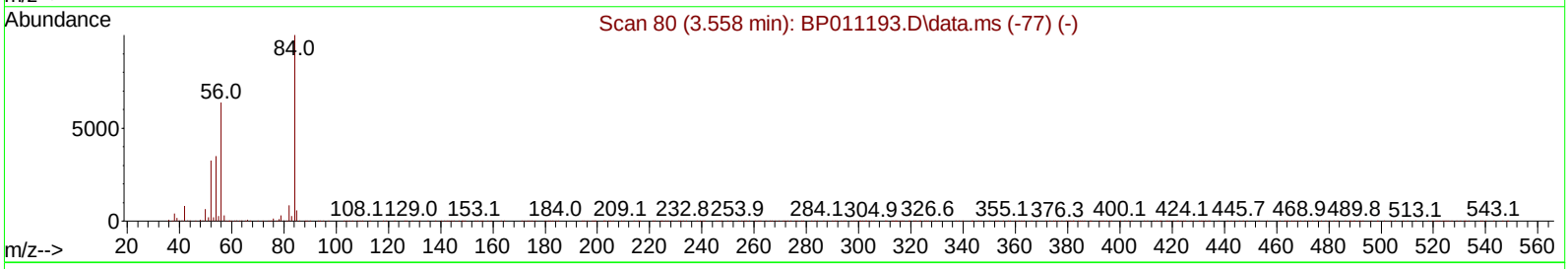

TIC: BP011197.D\data.ms

(4) Pyridine-d5 (S)

3.552min (-0.006) 0.01 ng/ul

response 246

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	6.56#
54.00	44.80	13.86#
0.00	0.00	0.00

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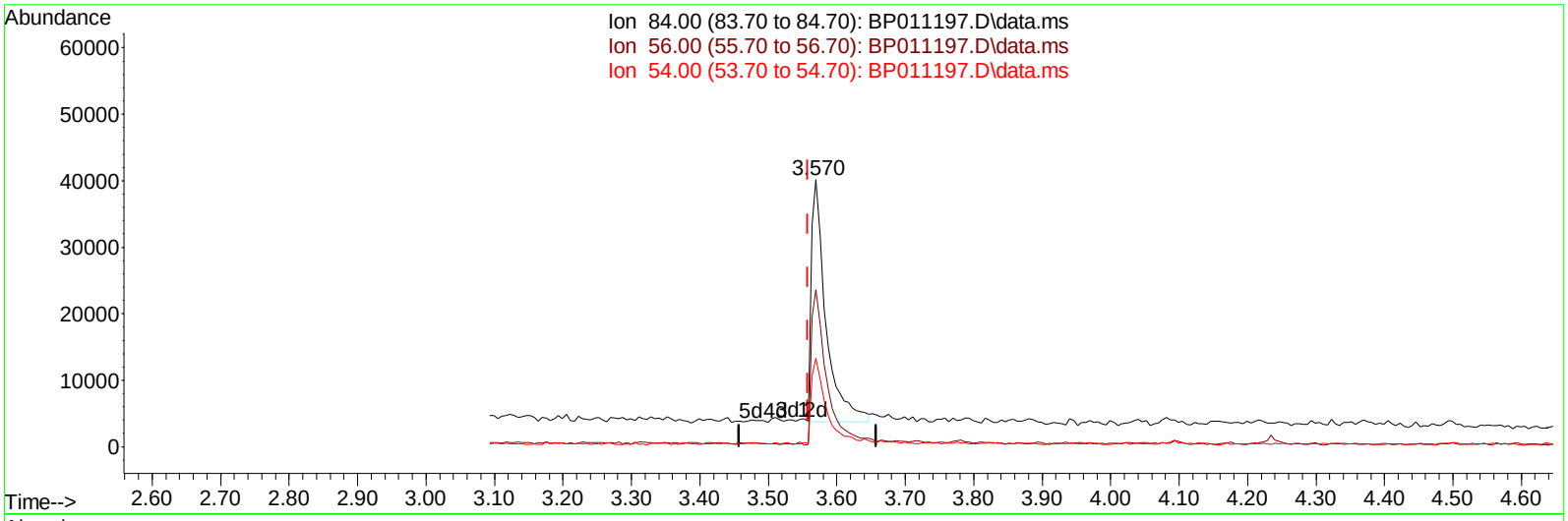

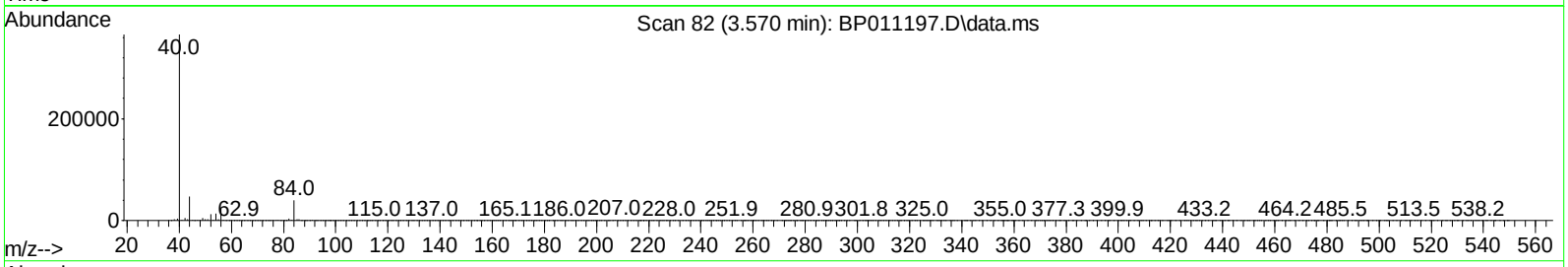

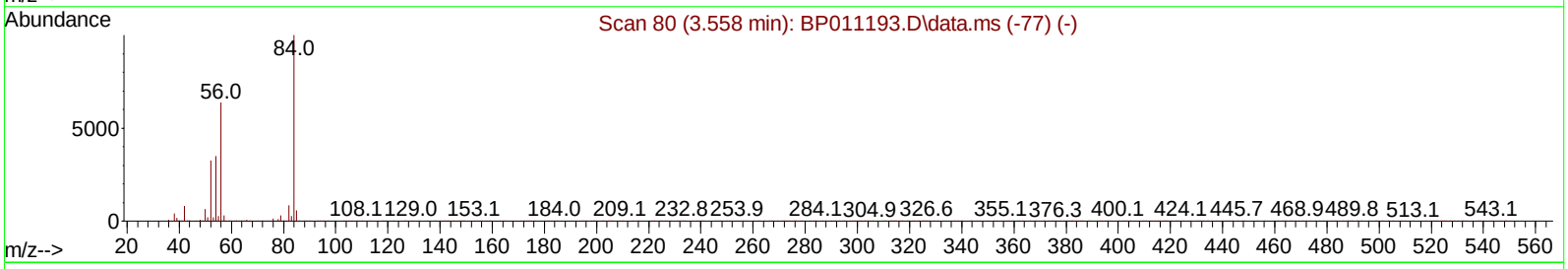

TIC: BP011197.D\data.ms

(4) Pyridine-d5 (S)

3.570min (+ 0.012) 2.35 ng/ul m

response 53980

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	58.88#
54.00	44.80	33.11#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	371209	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	1594083	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	745404	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	1295333	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	1329163	20.000	ng/ul	# 0.00
88) Perylene-d12	23.557	264	1412462	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	19648m	2.424	ng/uL	0.00
4) Pyridine-d5	3.570	84	53980m	2.354	ng/ul	0.01
7) Phenol-d5	6.828	99	638132	19.695	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	486300	21.226	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	529038	20.719	ng/ul	0.00
15) 4-Methylphenol-d8	8.364	113	435747	16.440	ng/ul	0.00
21) Nitrobenzene-d5	8.811	128	267983	21.631	ng/ul	0.00
24) 2-Nitrophenol-d4	9.528	143	250540	21.113	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	412201	19.710	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	291099	8.586	ng/ul	0.00
46) Dimethylphthalate-d6	13.734	166	1165477	22.949	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	1463253	22.928	ng/ul	0.00
54) 4-Nitrophenol-d4	14.540	143	115324	12.624	ng/ul	0.00
60) Fluorene-d10	15.316	176	937897	21.978	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.446	200	82020	12.561	ng/ul	0.00
73) Anthracene-d10	17.187	188	1156065	20.255	ng/ul	0.00
81) Pyrene-d10	19.445	212	1545816	21.643	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.404	264	1629330	23.224	ng/ul	0.00
Target Compounds						Qvalue

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

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