Quantitation Report (Qedit)

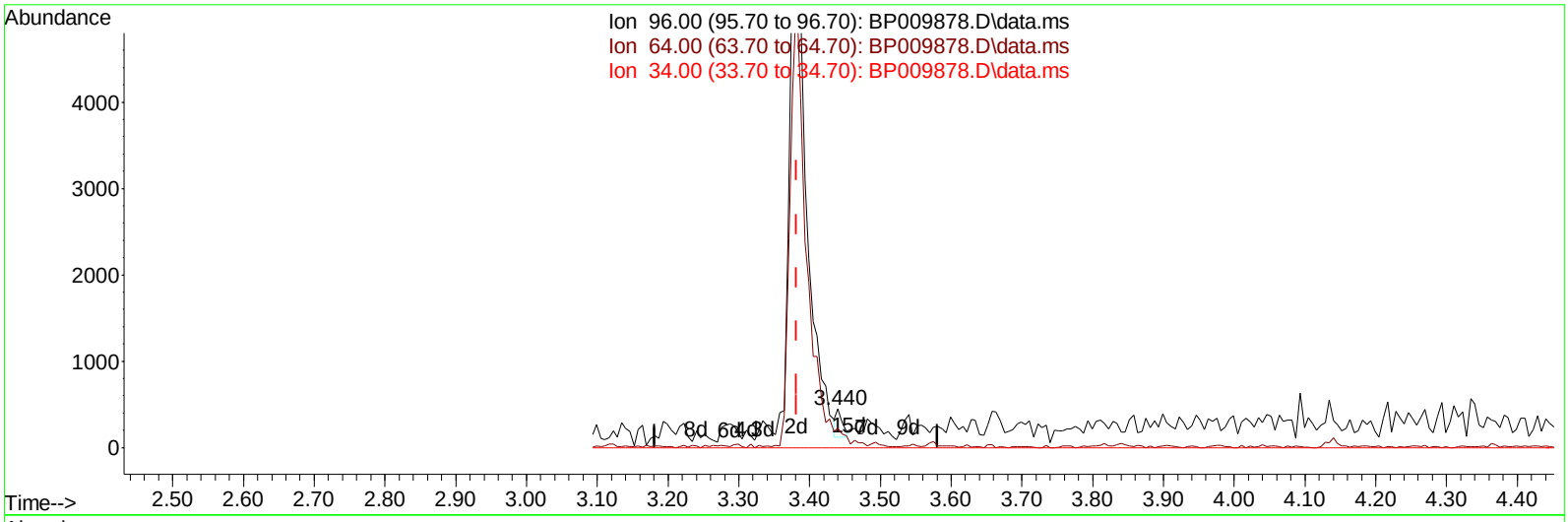
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
Data File : BP009878.D
Acq On : 16 Apr 2022 01:10
Operator : CG/JU
Sample : N2421-03
Misc :
ALS Vial : 29 Sample Multiplier: 1

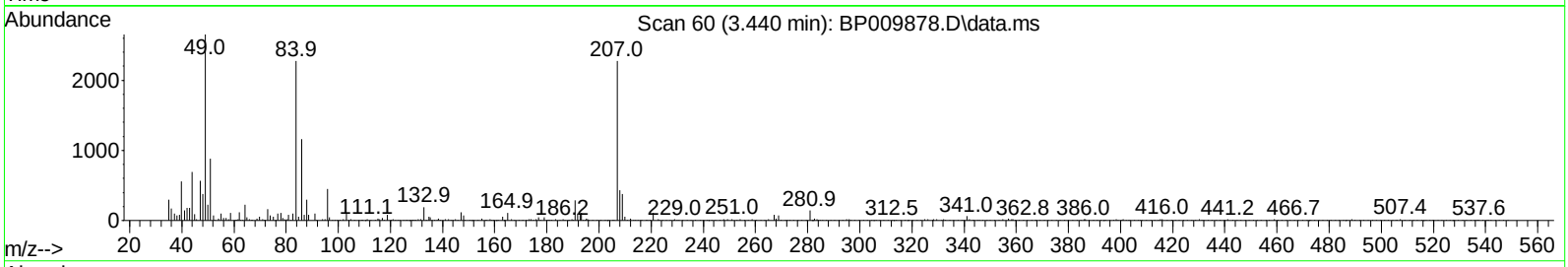
Quant Time: Apr 18 01:07:04 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 18 00:59:30 2022
Response via : Initial Calibration

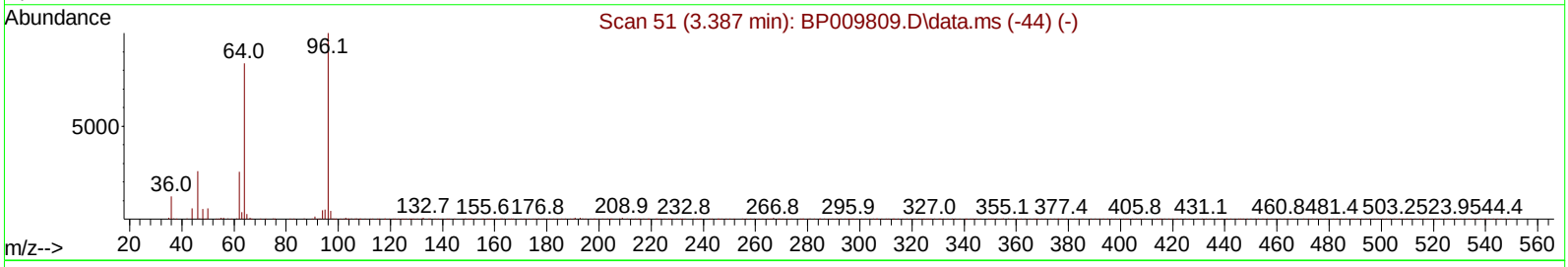
Instrument :
BNA_P
ClientSampleId :
C0J72

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022
Supervised By :Yogesh Patel 04/18/2022







TIC: BP009878.D\data.ms

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

3.440min (+ 0.059) 0.09 ng/uL

response 195

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.40
34.00	0.00	0.00
0.00	0.00	0.00

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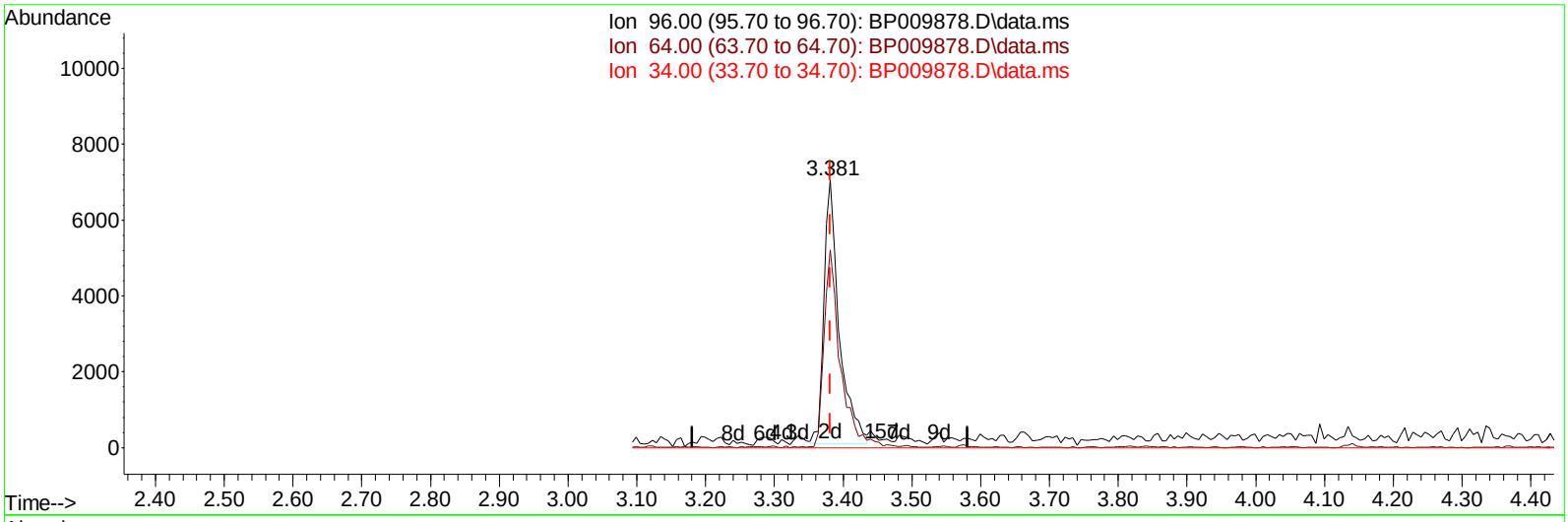

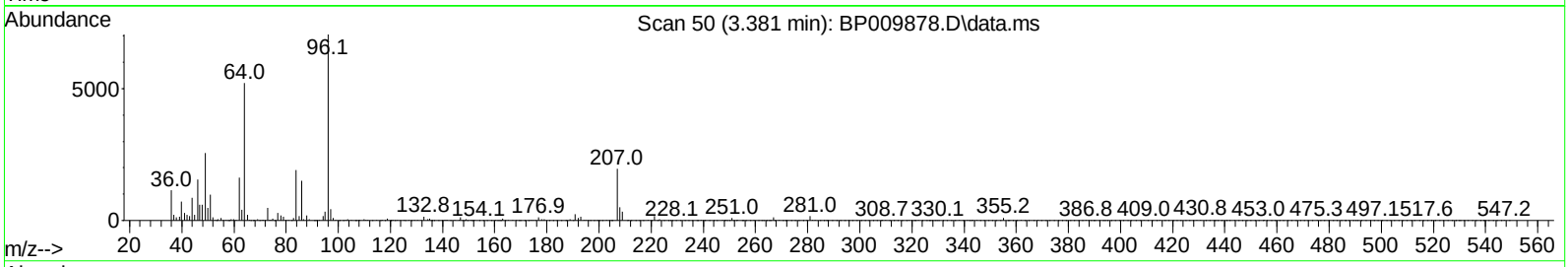

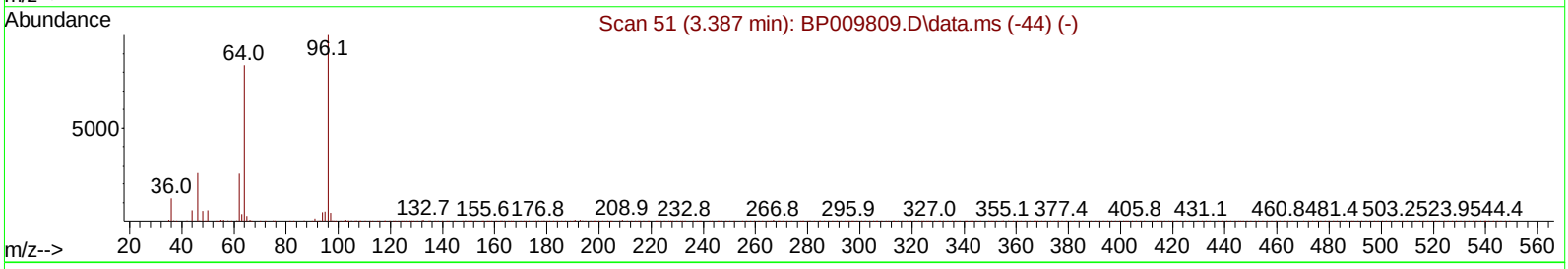

TIC: BP009878.D\data.ms

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

3.381min (+ 0.000) 4.98 ng/uL m

response 10641

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.86#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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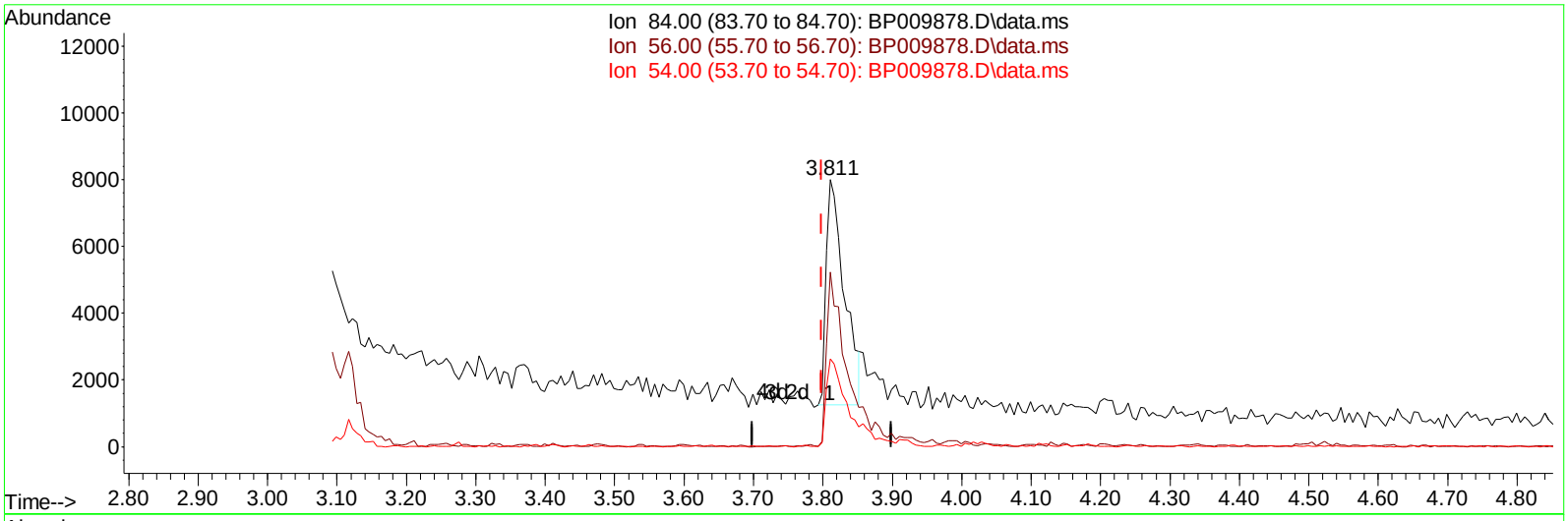

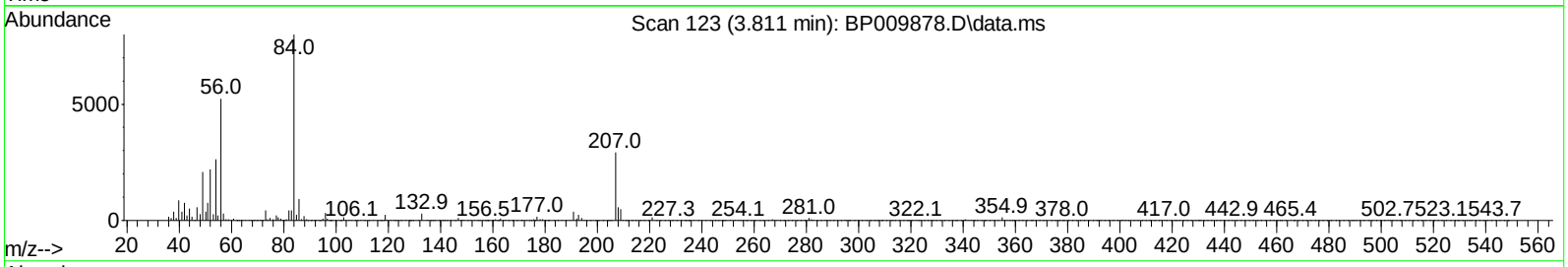

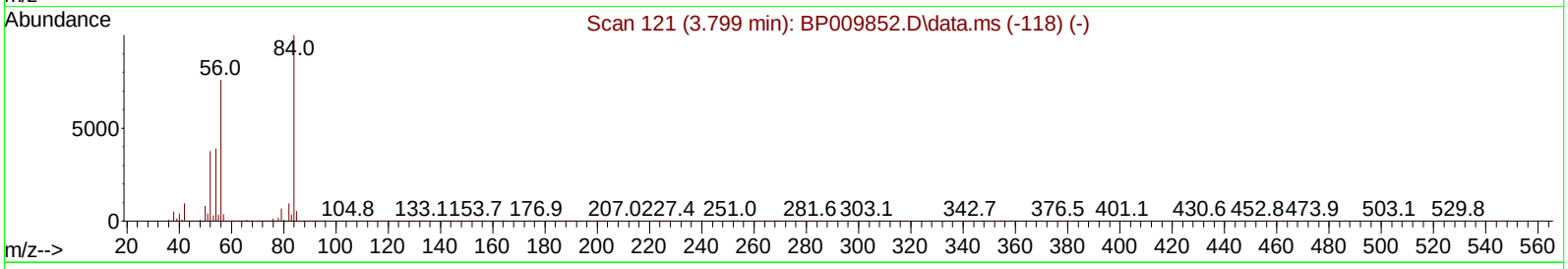

TIC: BP009878.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.012) 2.12 ng/ul

response 12462

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	65.44#
54.00	24.80	32.95#
0.00	0.00	0.00

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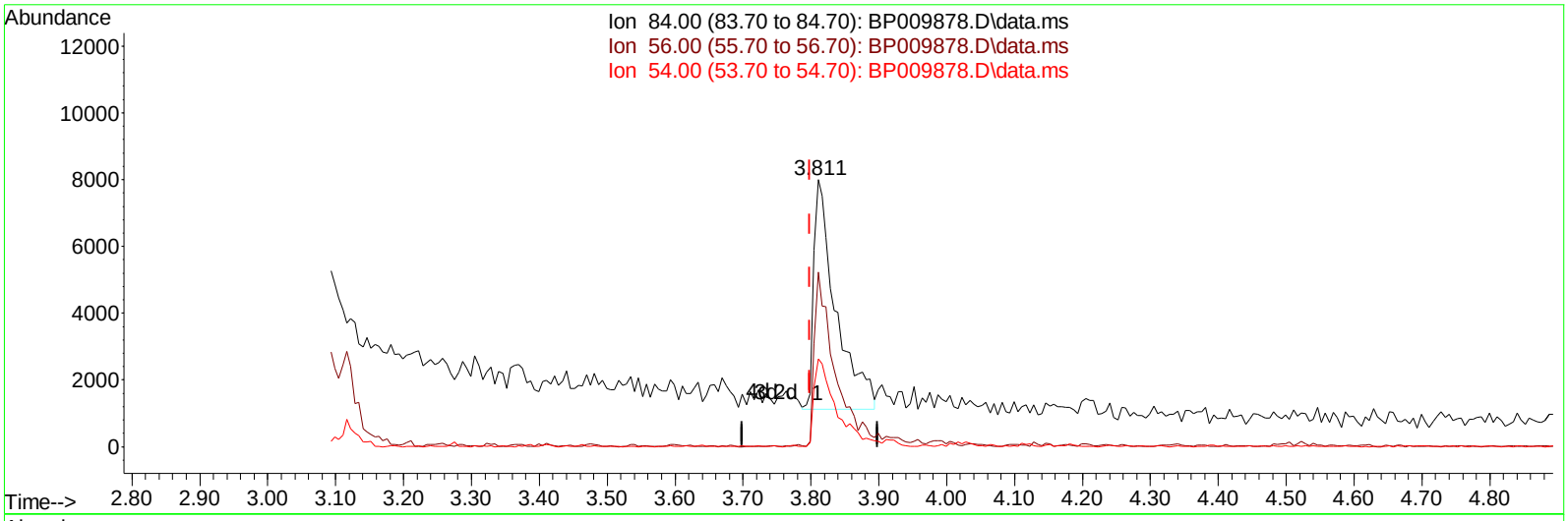

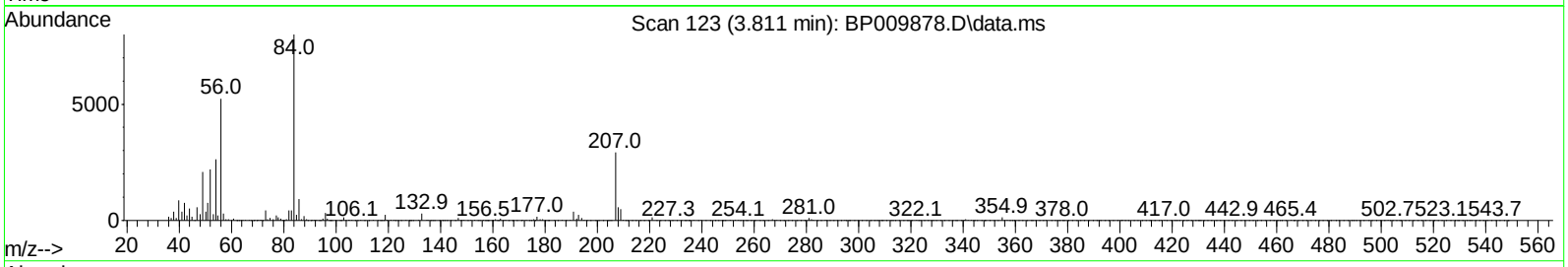

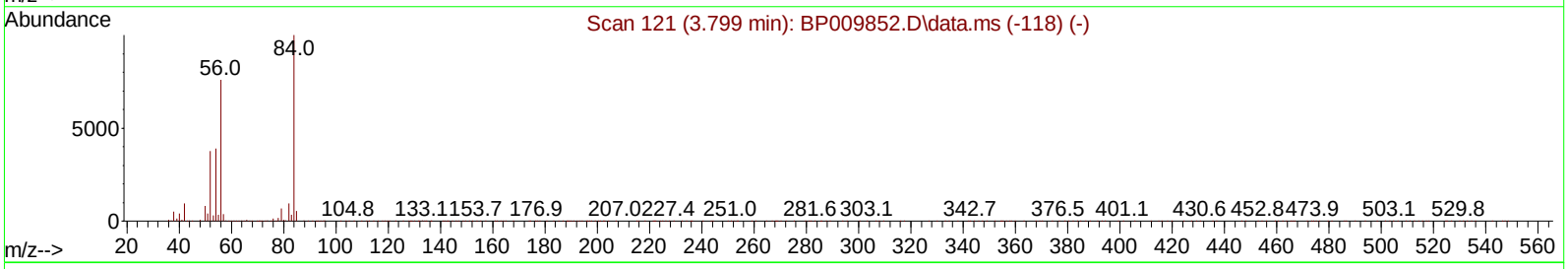

TIC: BP009878.D\data.ms

(4) Pyridine-d5 (S)

3.811min (+ 0.012) 2.62 ng/ul m

response 15446

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	65.44#
54.00	24.80	32.95#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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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	7.958	152	95335	20.000	ng/ul		0.00
20) Naphthalene-d8	10.763	136	445088	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.587	164	332642	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.339	188	758873	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.416	240	861820	20.000	ng/ul	#	0.00
88) Perylene-d12	23.880	264	844737	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.381	96	10641m	4.981	ng/uL		0.00
4) Pyridine-d5	3.811	84	15446m	2.623	ng/ul		0.01
7) Phenol-d5	7.111	99	70239	8.392	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	188599	37.850	ng/ul		0.00
11) 2-Chlorophenol-d4	7.487	132	185224	29.338	ng/ul		0.00
15) 4-Methylphenol-d8	8.658	113	139815	20.011	ng/ul		0.00
21) Nitrobenzene-d5	9.116	128	125379	39.063	ng/ul		0.00
24) 2-Nitrophenol-d4	9.840	143	131189	37.281	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.375	165	236748	32.151	ng/ul		0.00
31) 4-Chloroaniline-d4	10.893	131	105566	10.066	ng/ul		0.00
46) Dimethylphthalate-d6	13.993	166	1020863	39.069	ng/ul		0.00
49) Acenaphthylene-d8	14.281	160	1096651	35.426	ng/ul		0.00
54) 4-Nitrophenol-d4	14.763	143	40366	8.518	ng/ul		0.00
60) Fluorene-d10	15.581	176	893456	40.647	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	171748	38.204	ng/ul		0.00
73) Anthracene-d10	17.439	188	1649533	45.399	ng/ul		0.00
81) Pyrene-d10	19.669	212	2086304	44.325	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.727	264	2107673	47.912	ng/ul		0.00
Target Compounds							Qvalue

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

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