Quantitation Report (Qedit)

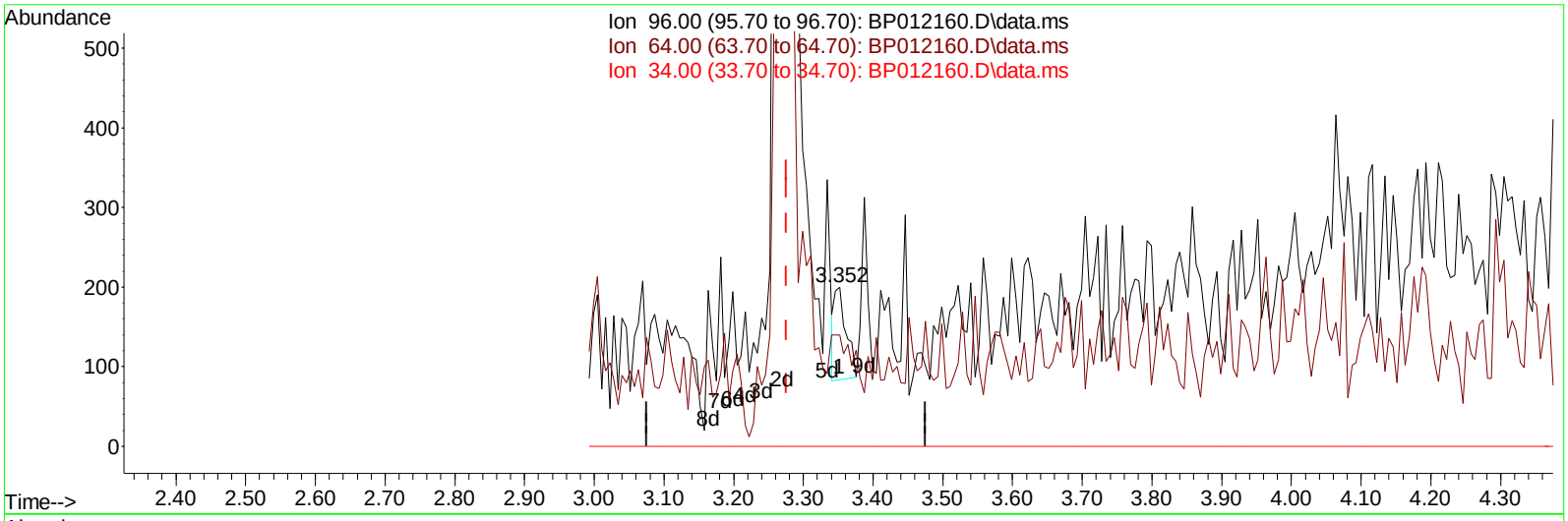
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
Data File : BP012160.D
Acq On : 15 Oct 2022 04:26
Operator : CG/JU
Sample : N4984-18
Misc :
ALS Vial : 35 Sample Multiplier: 1

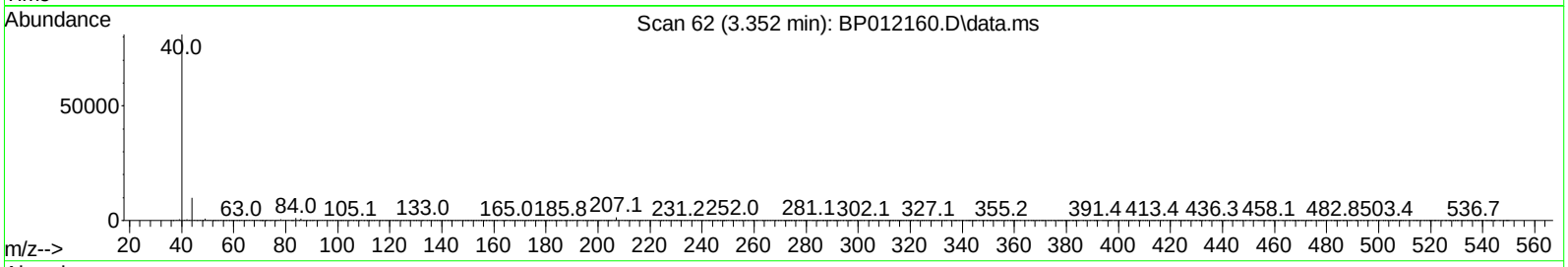
Quant Time: Oct 15 05:03:15 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 12 23:03:21 2022
Response via : Initial Calibration

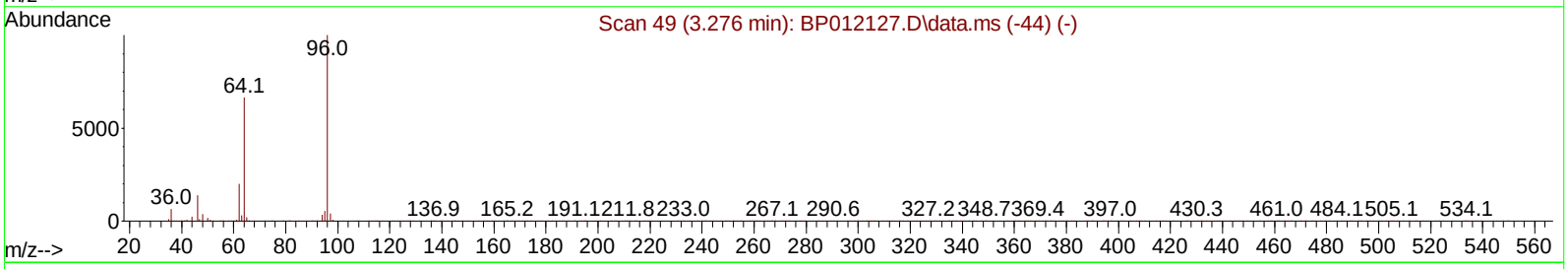
Instrument :
BNA_P
ClientSampleId :
BGNB9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/15/2022
Supervised By :mohammad ahmed 10/16/2022







TIC: BP012160.D\data.ms

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

3.352min (+ 0.076) 0.03 ng/uL

response 138

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	70.00
34.00	0.00	0.00
0.00	0.00	0.00

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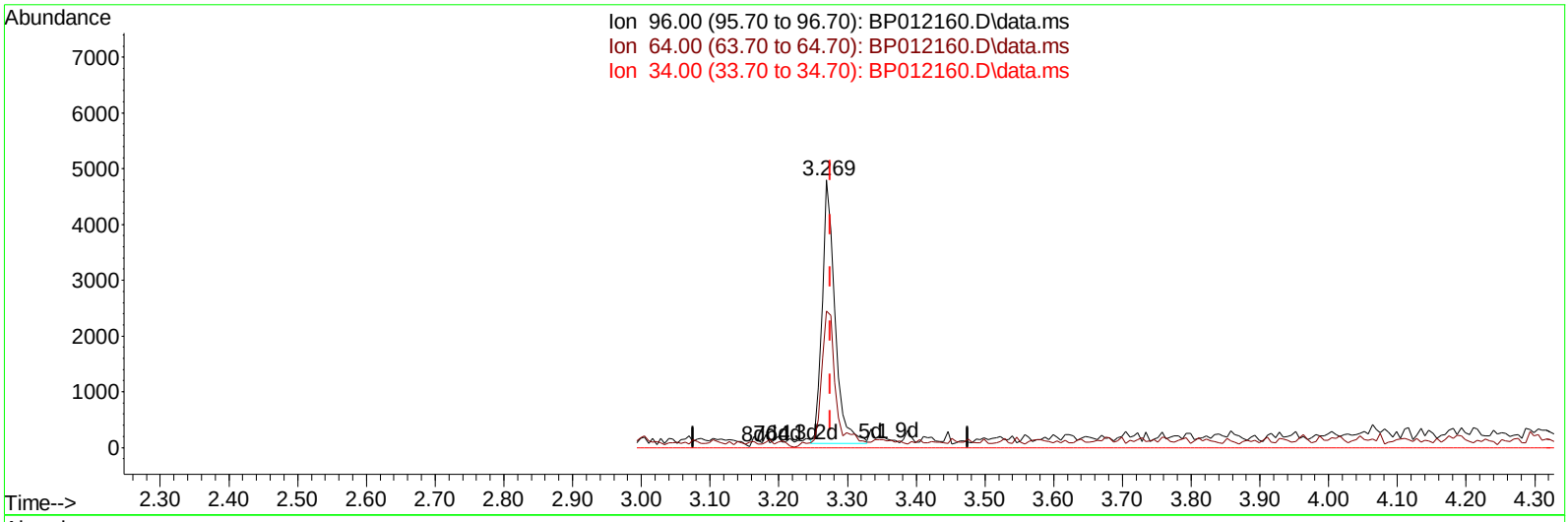

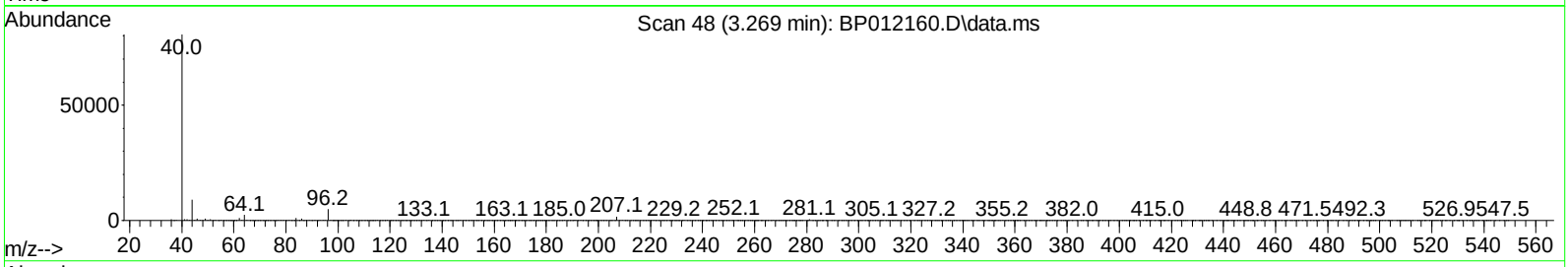

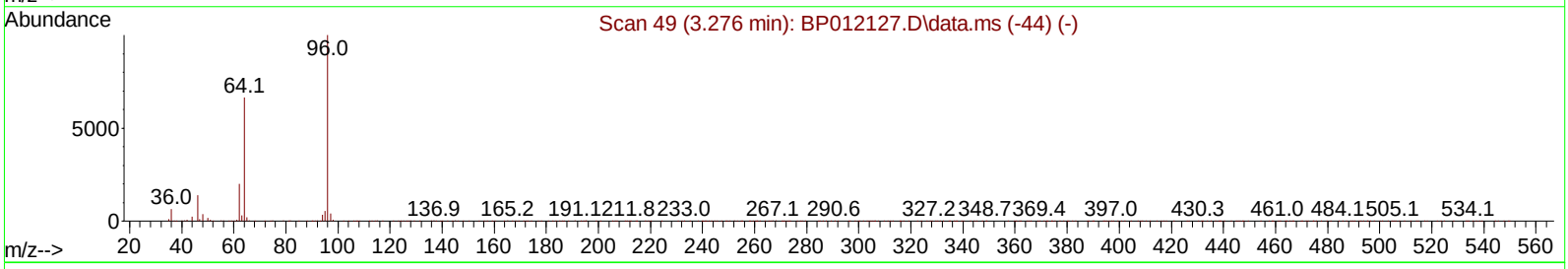

TIC: BP012160.D\data.ms

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

3.269min (-0.006) 1.14 ng/uL m

response 6114

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.10	51.01#
34.00	0.00	0.00
0.00	0.00	0.00

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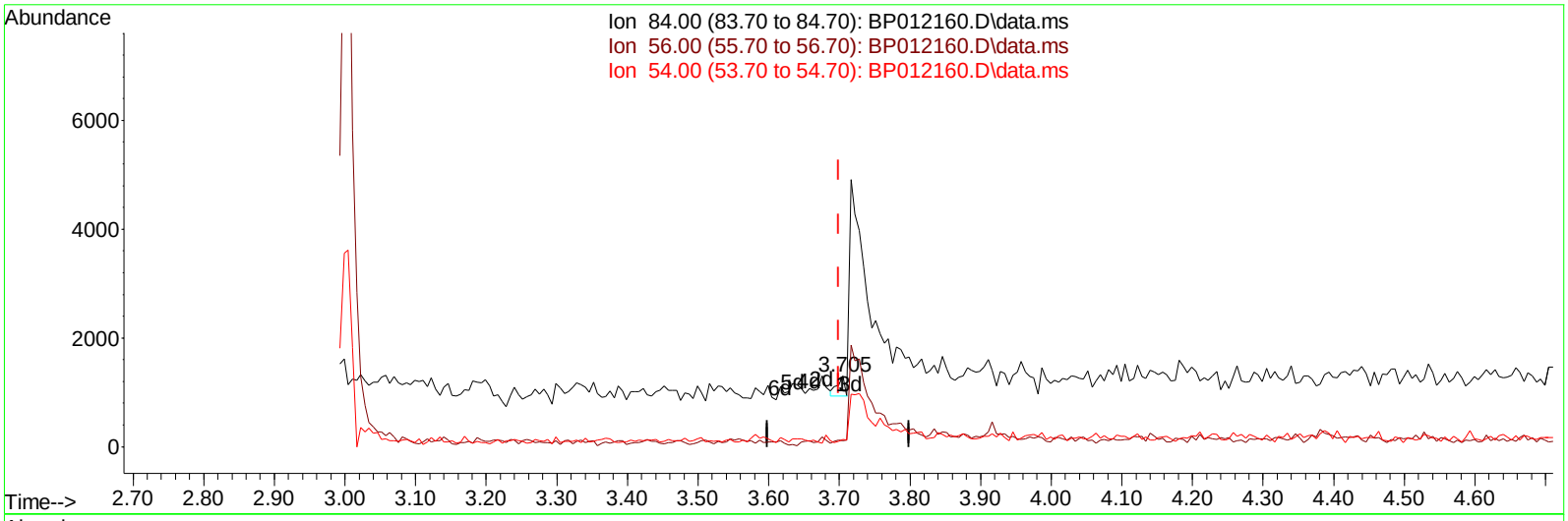

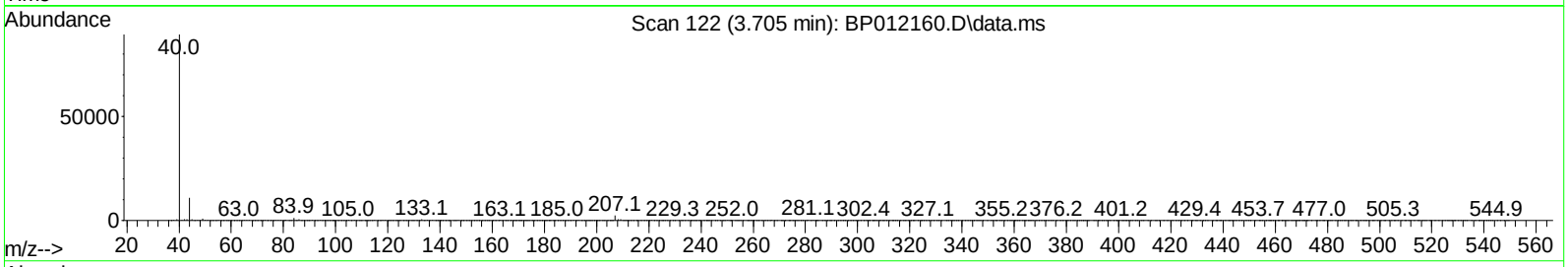

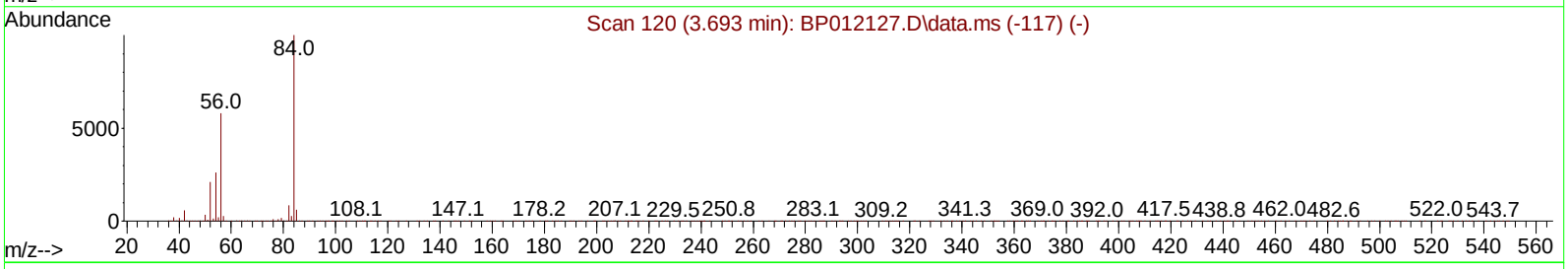

TIC: BP012160.D\data.ms

(4) Pyridine-d5 (S)

3.705min (+ 0.006) 0.02 ng/ul

response 264

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	8.56#
54.00	26.00	10.49#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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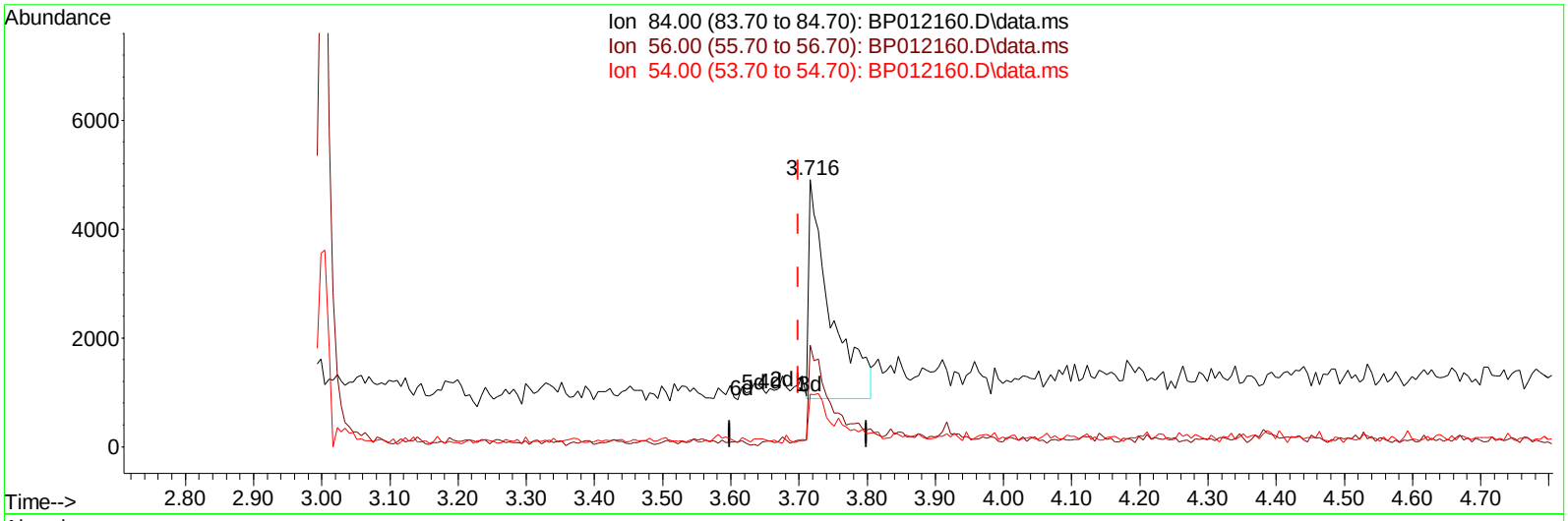

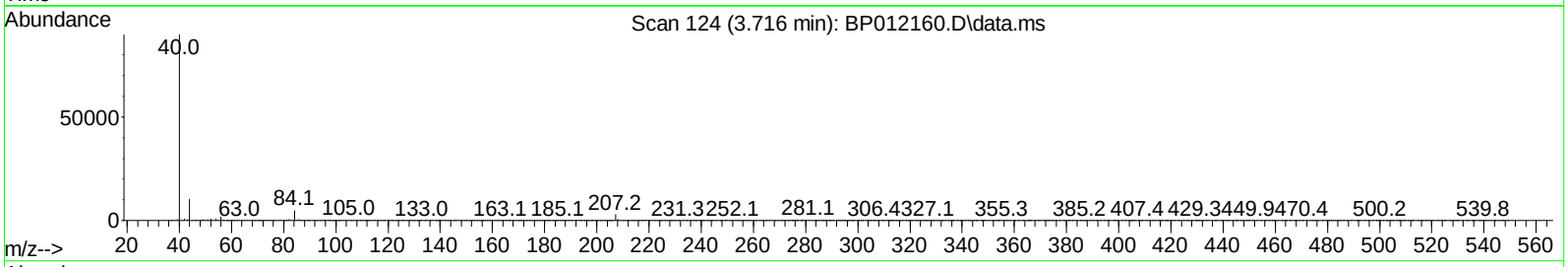

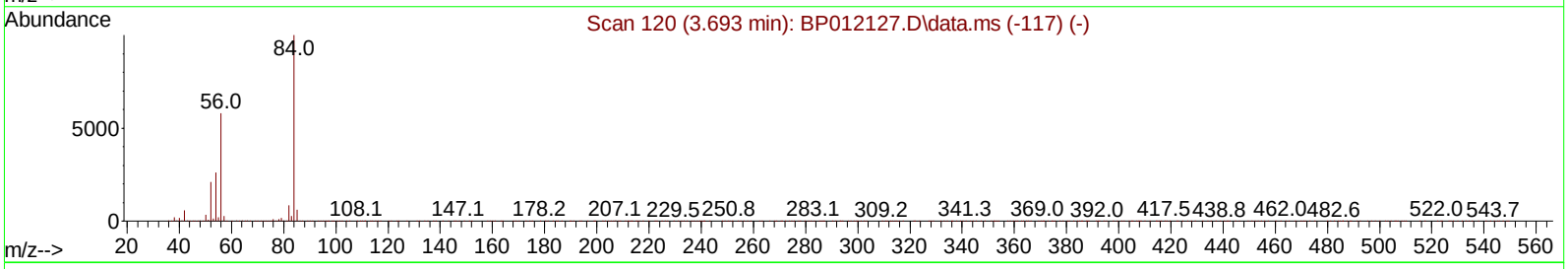

TIC: BP012160.D\data.ms

(4) Pyridine-d5 (S)

3.716min (+ 0.018) 0.62 ng/ul m

response 8920

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	38.11#
54.00	26.00	19.65#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Reviewed By :Yogesh Patel 10/15/2022
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	211669	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.646	136	874597	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.486	164	464653	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.245	188	911568	20.000	ng/ul	-0.01
79) Chrysene-d12	21.339	240	975339	20.000	ng/ul	# 0.00
88) Perylene-d12	23.751	264	1052749	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	6114m	1.139	ng/uL	0.00
4) Pyridine-d5	3.716	84	8920m	0.625	ng/ul	0.02
7) Phenol-d5	7.010	99	373541	21.942	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	227631	23.557	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.369	132	297428	22.011	ng/ul	-0.01
15) 4-Methylphenol-d8	8.552	113	291887	22.323	ng/ul	-0.01
21) Nitrobenzene-d5	9.010	128	136407	22.084	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.728	143	133629	21.336	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	10.269	165	248069	20.437	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.787	131	321938	17.998	ng/ul	-0.02
46) Dimethylphthalate-d6	13.904	166	687638	23.367	ng/ul	-0.01
49) Acenaphthylene-d8	14.181	160	862203	23.694	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.704	143	63443	10.982	ng/ul	0.00
60) Fluorene-d10	15.486	176	636628	23.912	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.610	200	33121	6.881	ng/ul	-0.01
73) Anthracene-d10	17.345	188	1011687	25.558	ng/ul	-0.01
81) Pyrene-d10	19.586	212	1196315	23.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.598	264	1299165	25.386	ng/ul	-0.01
Target Compounds						Qvalue
8) Phenol	7.034	94	18918	1.059	ng/ul#	83

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

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