Quantitation Report (Qedit)

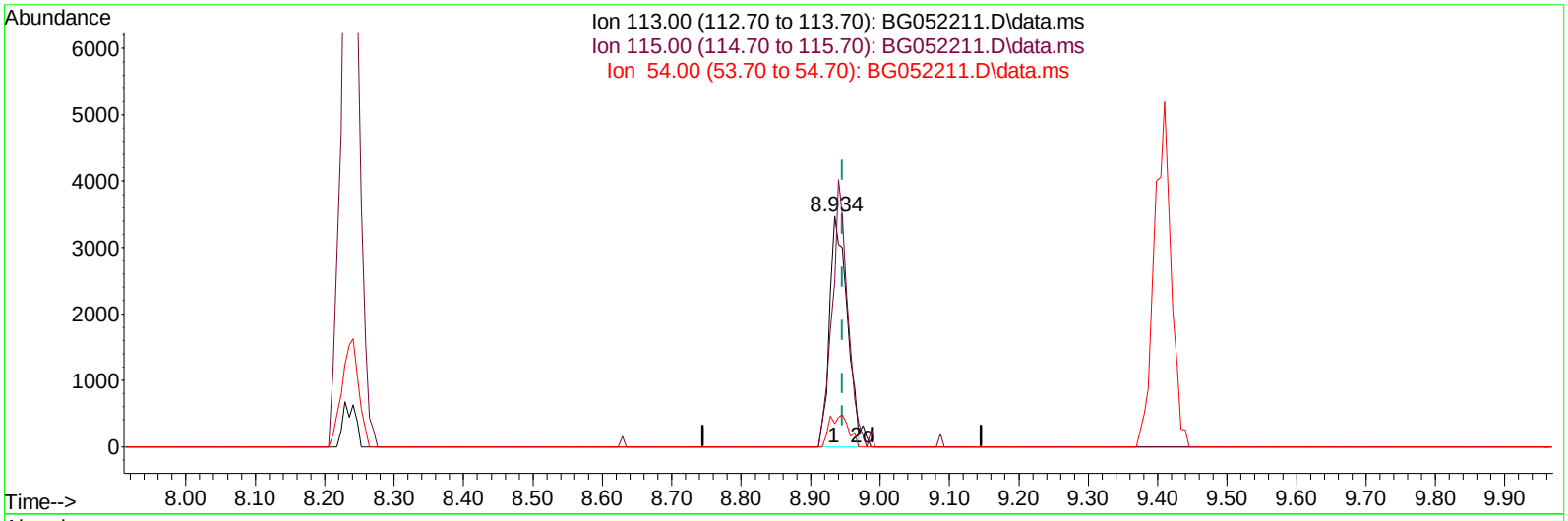
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
Data File : BG052211.D
Acq On : 28 Jan 2022 11:05
Operator : CG/JU
Sample : N1273-06DL 5X
Misc :
ALS Vial : 71 Sample Multiplier: 1

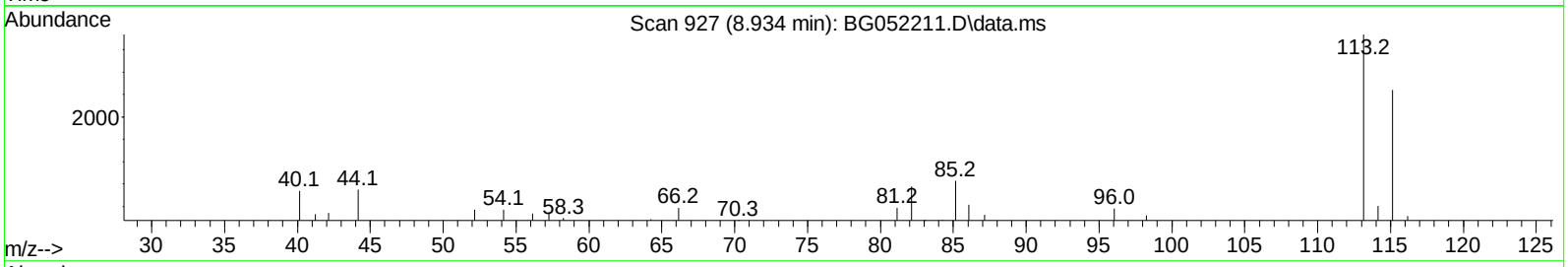
Quant Time: Jan 29 00:50:43 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 26 23:58:52 2022
Response via : Initial Calibration

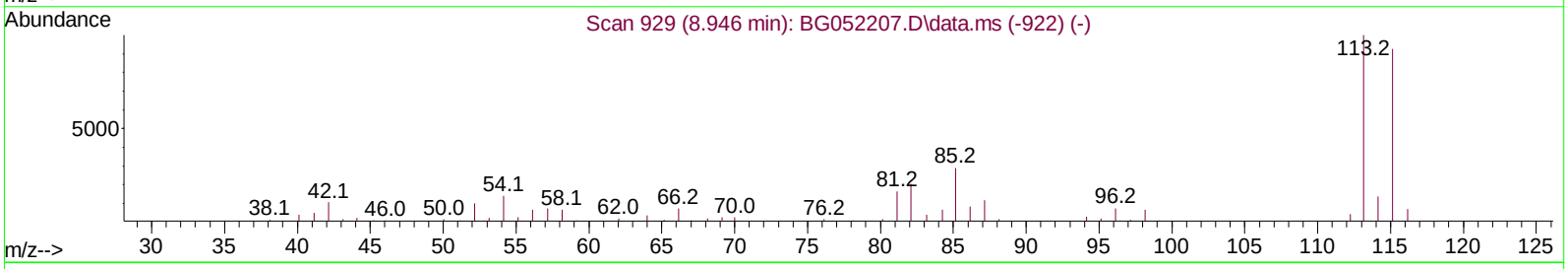
Instrument :
BNA_G
ClientSampleId :
C0HK5DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/31/2022
Supervised By :mohammad ahmed 01/31/2022







TIC: BG052211.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.934min (-0.012) 2.90 ng/ul

response 6174

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	88.40	71.38
54.00	13.00	10.02#
0.00	0.00	0.00

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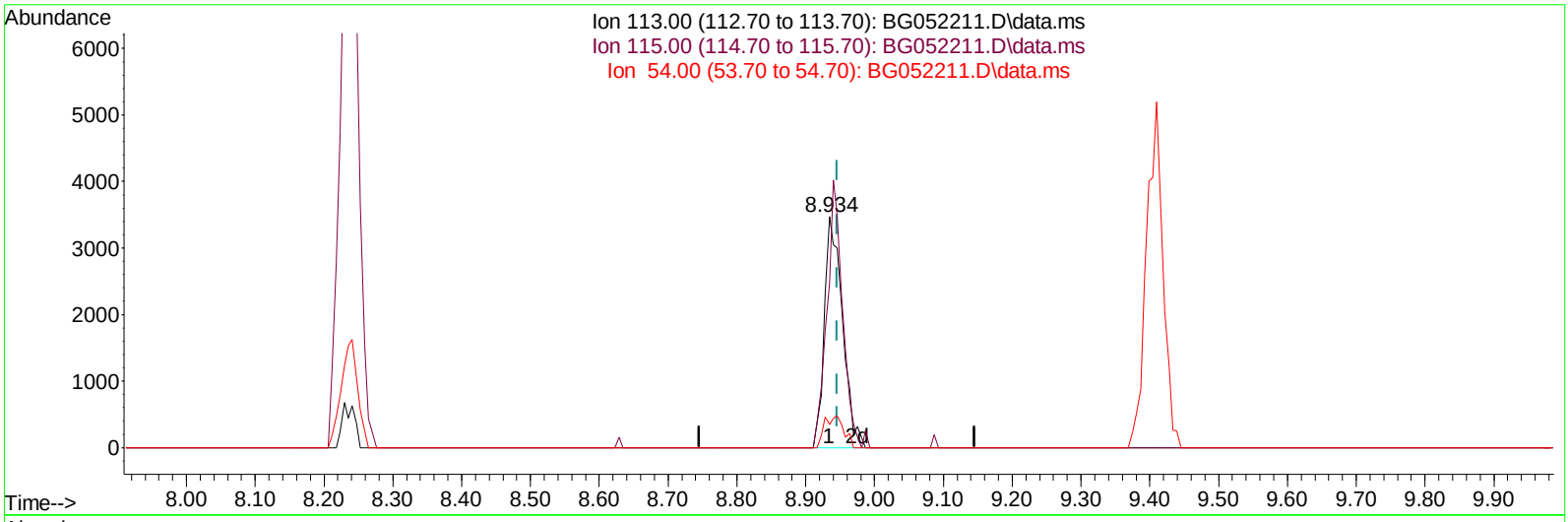

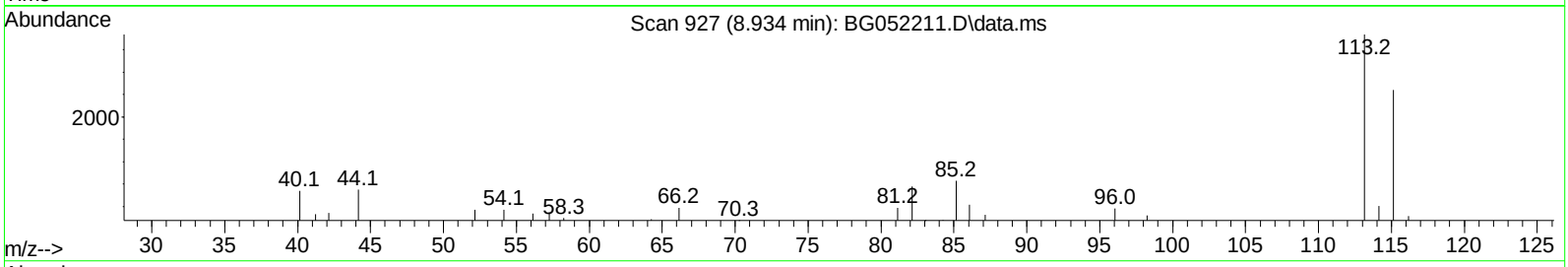

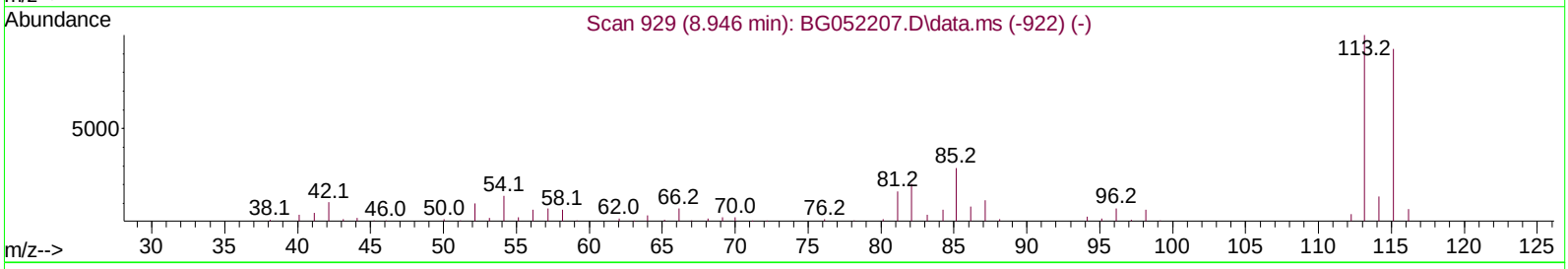

TIC: BG052211.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.934min (-0.012) 2.98 ng/ul m

response 6339

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	88.40	71.38
54.00	13.00	10.02#
0.00	0.00	0.00

Quantitation Report (Qedit)

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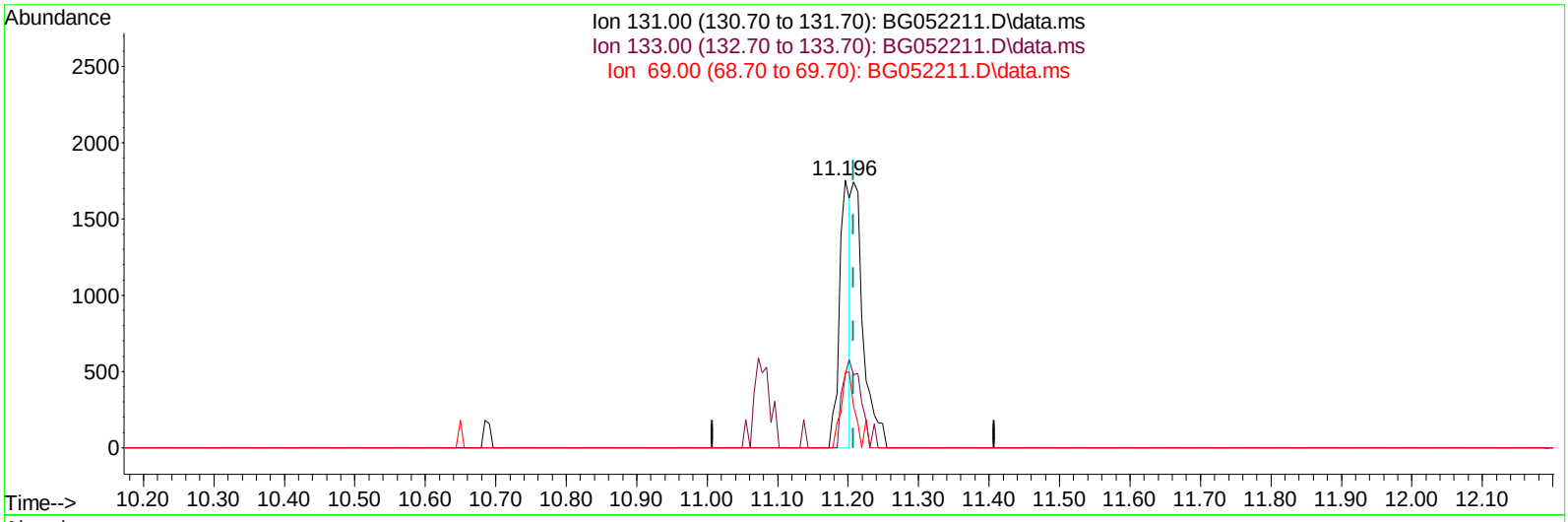

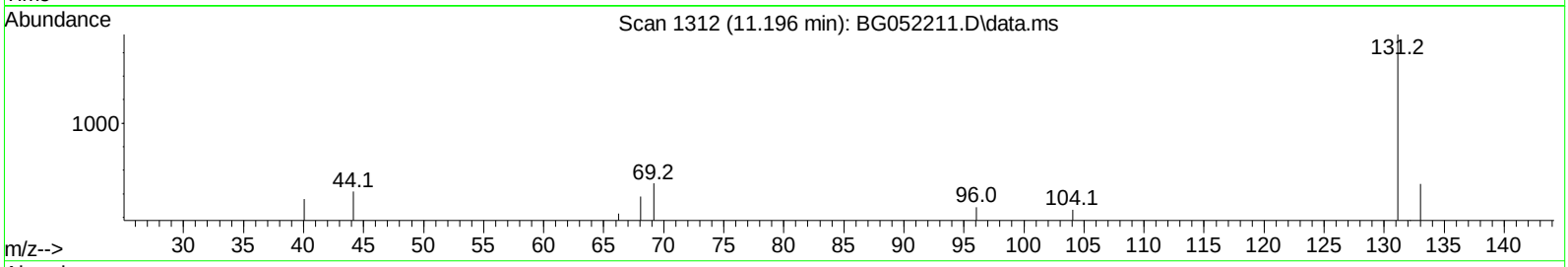

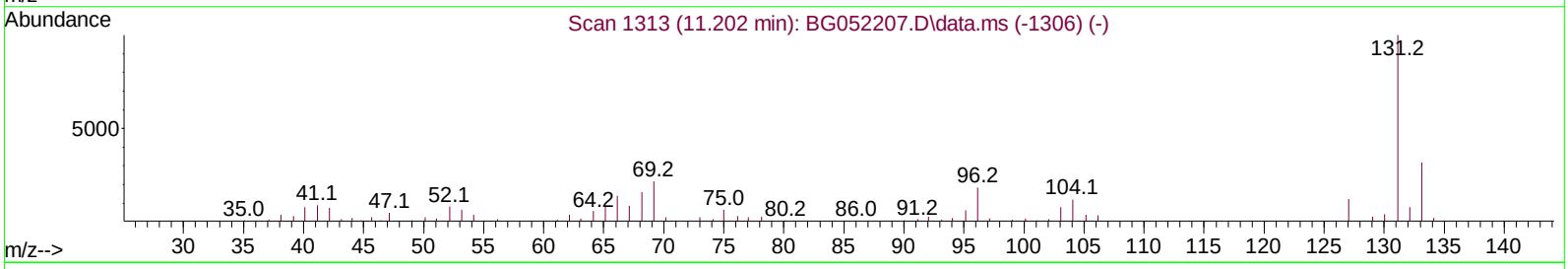

TIC: BG052211.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.196min (-0.012) 0.66 ng/ul

response 1890

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.20	27.59
69.00	21.30	27.94#
0.00	0.00	0.00

Quantitation Report (Qedit)

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ALS Vial : 71 Sample Multiplier: 1

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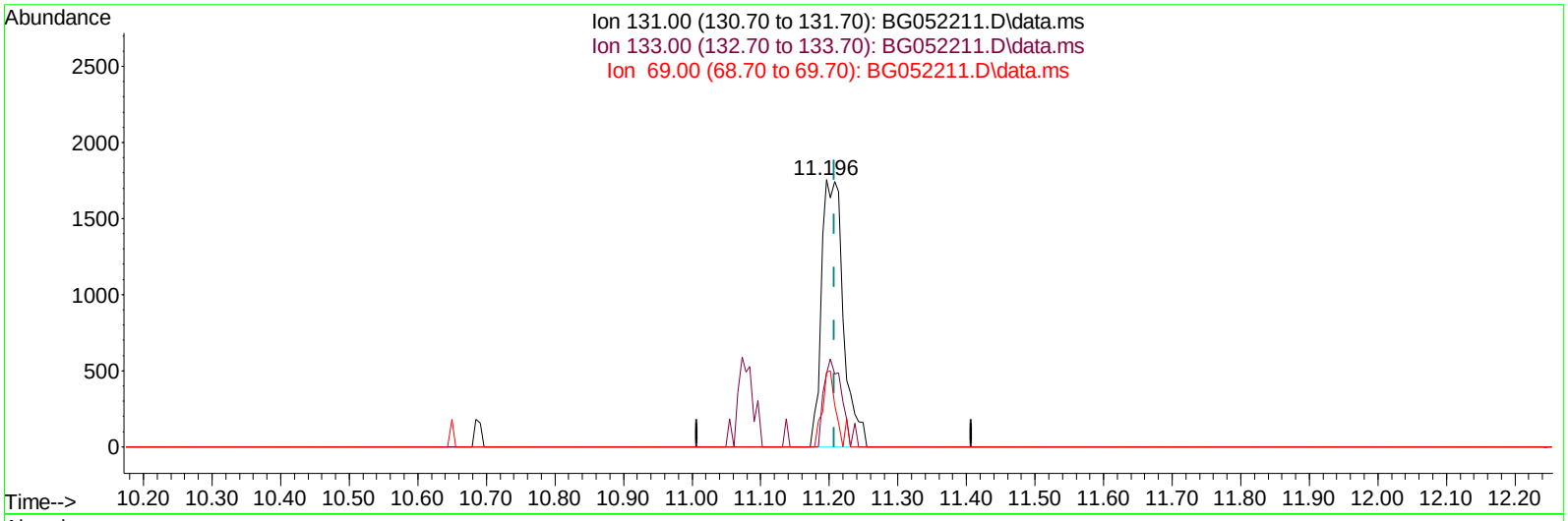

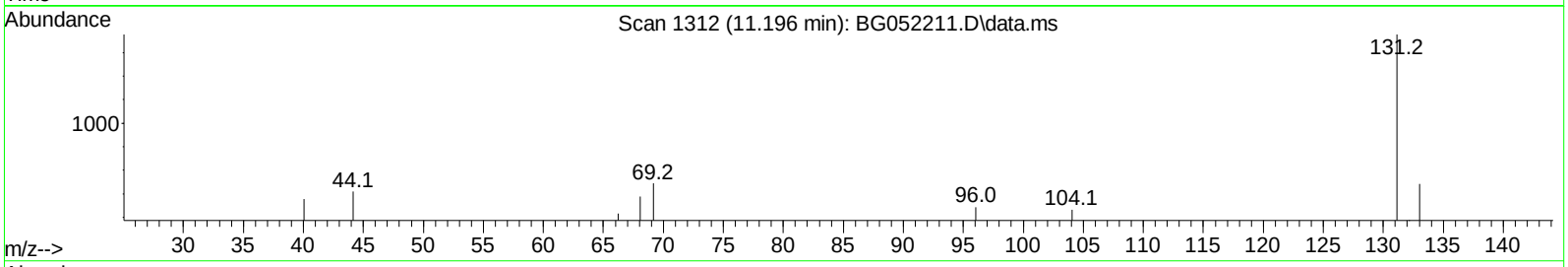

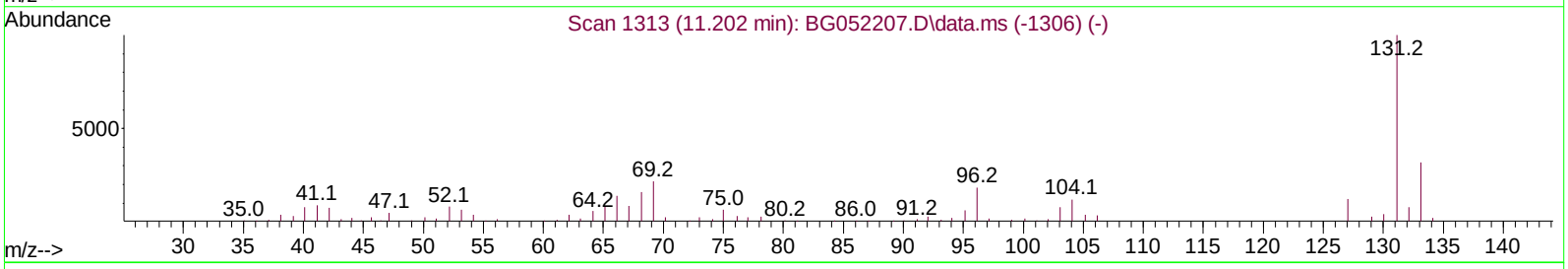

TIC: BG052211.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.196min (-0.012) 1.35 ng/ul m

response 3864

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.20	27.59
69.00	21.30	27.94#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.235	152	27988	20.000	ng/ul	0.00
20) Naphthalene-d8	11.073	136	120289	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.879	164	87653	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.628	188	213224	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	222101	20.000	ng/ul	# 0.00
88) Perylene-d12	25.430	264	233214	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	794	1.014	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.383	99	3633	1.326	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	10605	6.099	ng/ul	0.00
11) 2-Chlorophenol-d4	7.765	132	9104	5.020	ng/ul	0.00
15) 4-Methylphenol-d8	8.934	113	6339m	2.977	ng/ul	-0.01
21) Nitrobenzene-d5	9.404	128	5932	6.241	ng/ul	0.00
24) 2-Nitrophenol-d4	10.139	143	5628	5.354	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	11565	5.628	ng/ul	0.00
31) 4-Chloroaniline-d4	11.196	131	3864m	1.350	ng/ul	-0.01
46) Dimethylphthalate-d6	14.268	166	50308	6.944	ng/ul	0.00
49) Acenaphthylene-d8	14.574	160	59071	6.750	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.872	176	44146	6.989	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.728	188	78612	7.772	ng/ul	0.00
81) Pyrene-d10	20.008	212	96470	7.547	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.183	264	89223	7.267	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.600	88	15693	17.365	ng/uL#	90

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

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