Quantitation Report (Qedit)

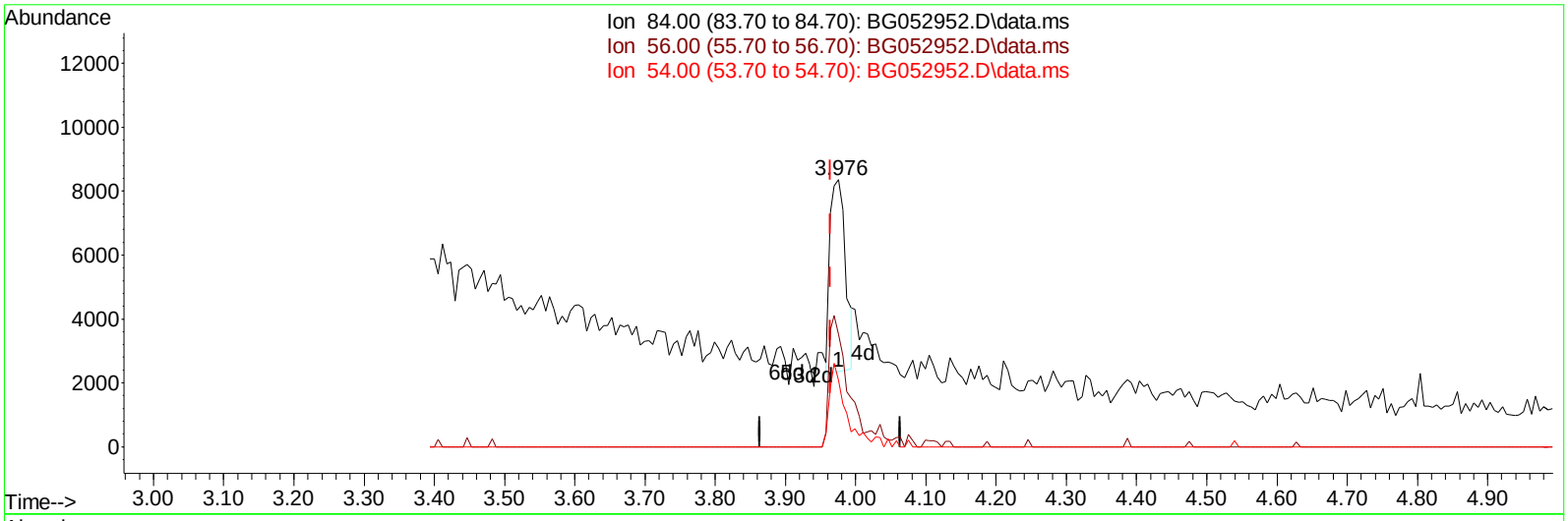
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033022\
Data File : BG052952.D
Acq On : 30 Mar 2022 16:47
Operator : CG/JU
Sample : N2068-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

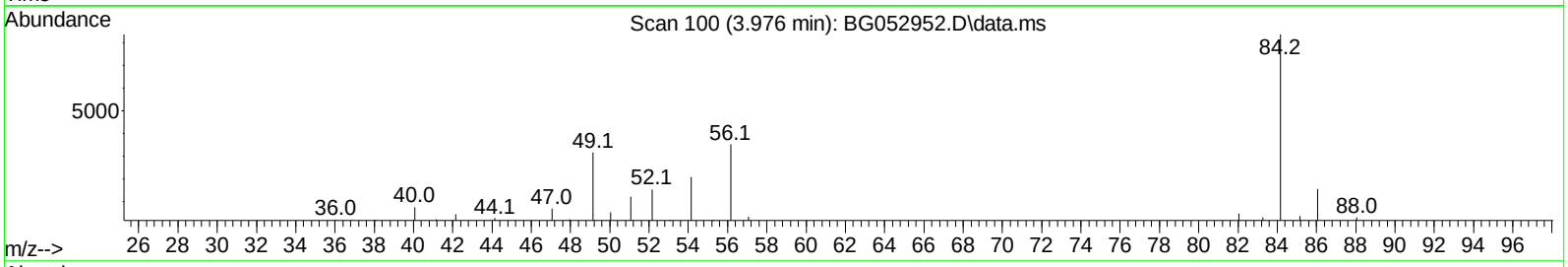
Quant Time: Mar 31 00:52:52 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 31 00:50:49 2022
Response via : Initial Calibration

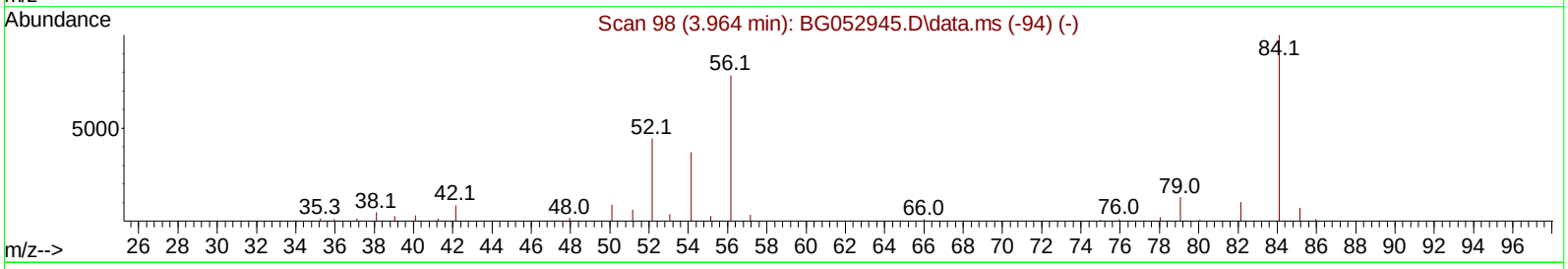
Instrument :
BNA_G
ClientSampleId :
YAYX2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/31/2022
Supervised By :Yogesh Patel 04/06/2022







TIC: BG052952.D\data.ms

(4) Pyridine-d5 (S)

3.976min (+ 0.011) 5.69 ng/ul

response 9207

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	42.14#
54.00	28.70	25.00
0.00	0.00	0.00

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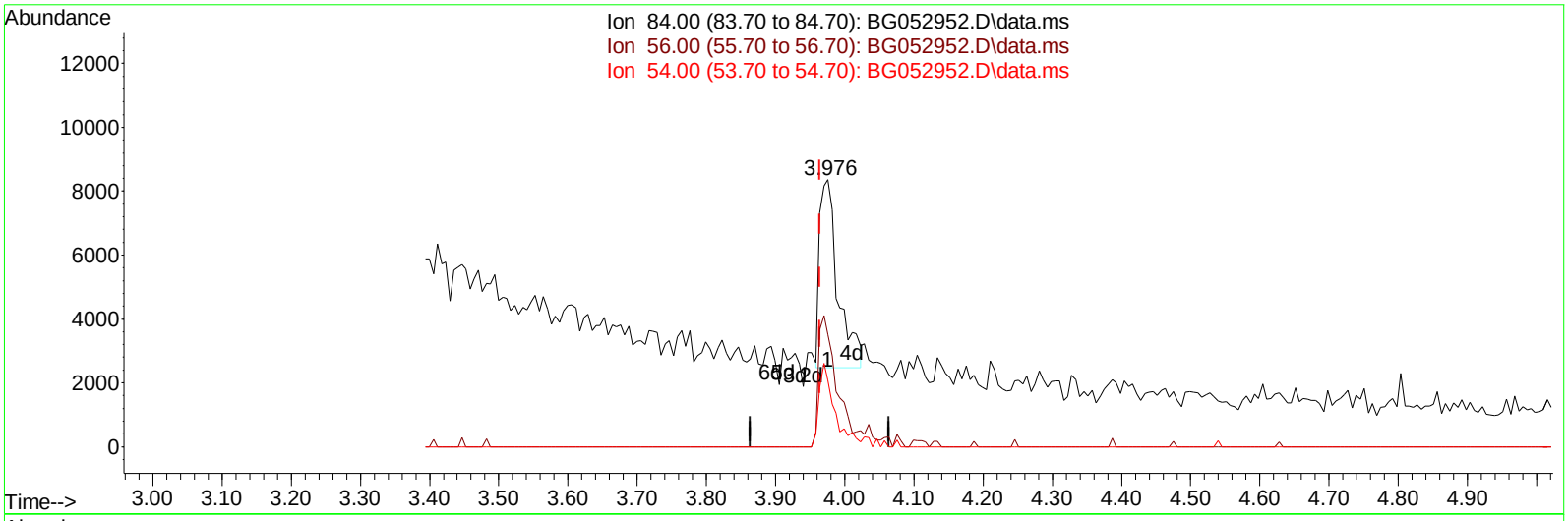

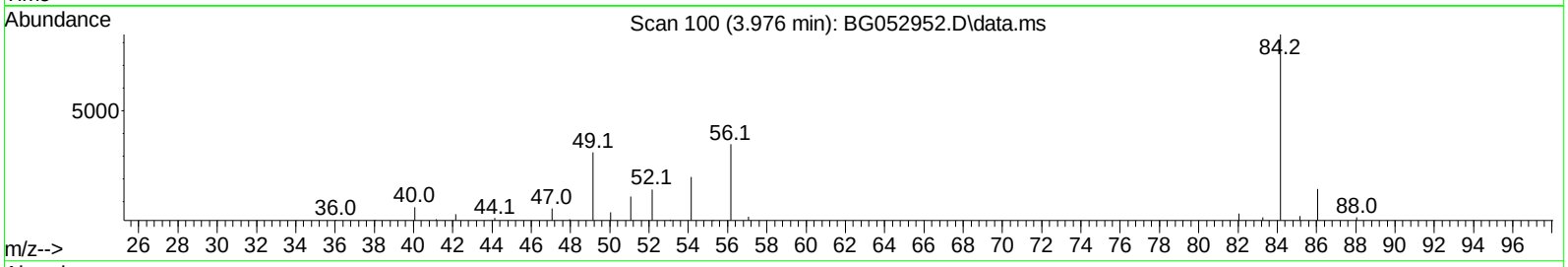

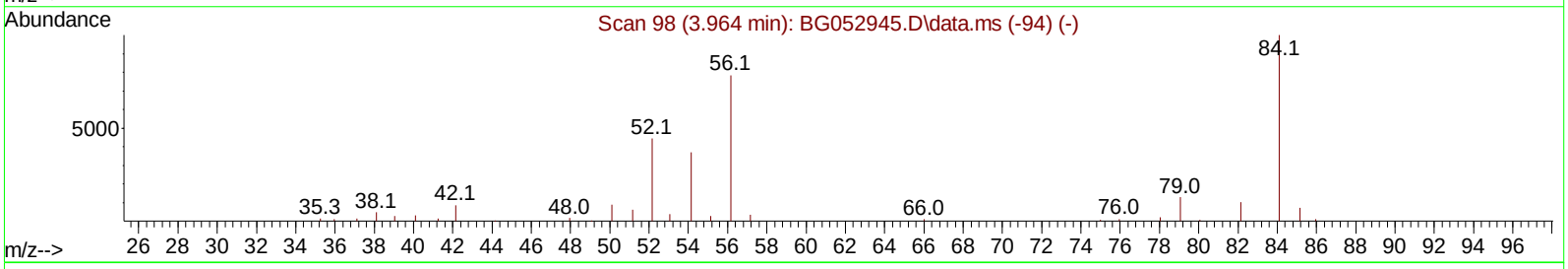

TIC: BG052952.D\data.ms

(4) Pyridine-d5 (S)

3.976min (+ 0.011) 6.75 ng/ul m

response 10915

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	42.14#
54.00	28.70	25.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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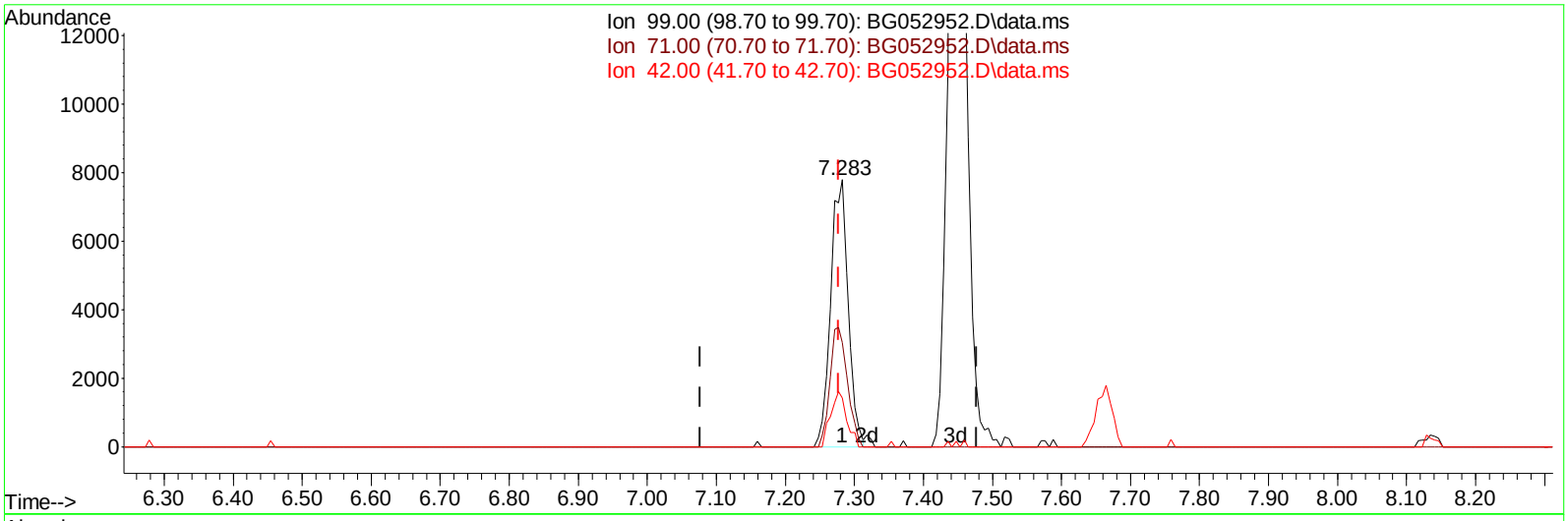

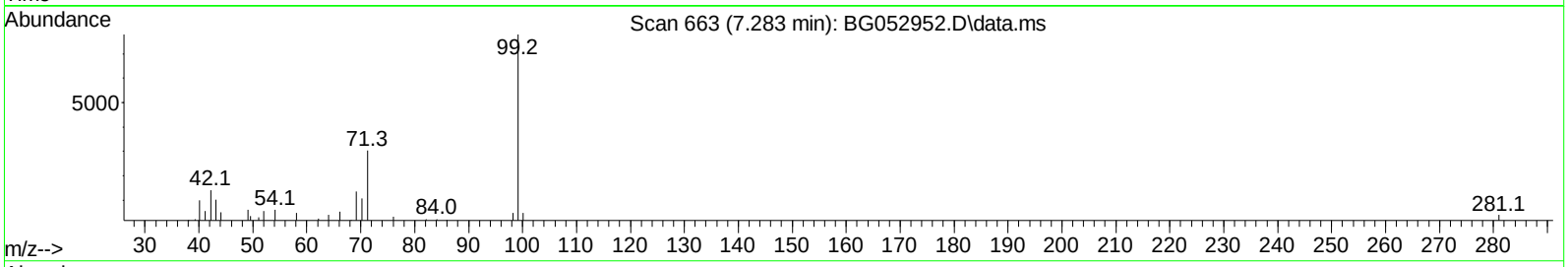

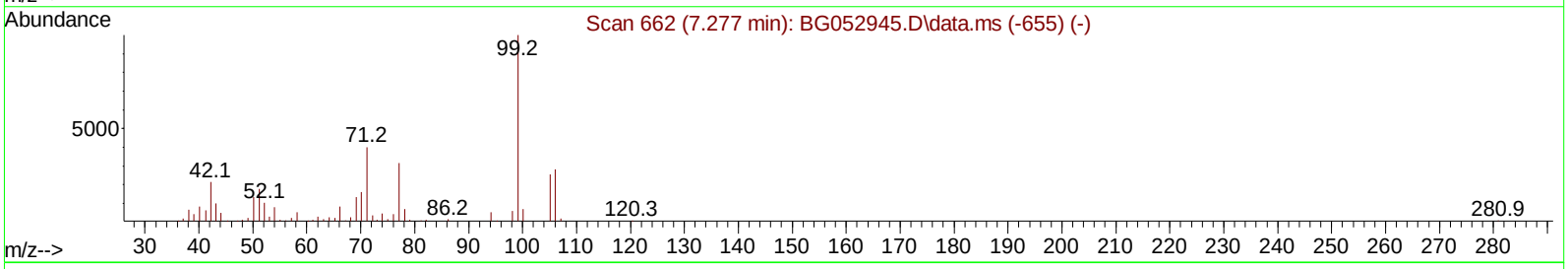

TIC: BG052952.D\data.ms

(7) Phenol-d5 (S)

7.283min (+ 0.006) 6.72 ng/ul

response 13778

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	36.40	39.14
42.00	19.00	18.24
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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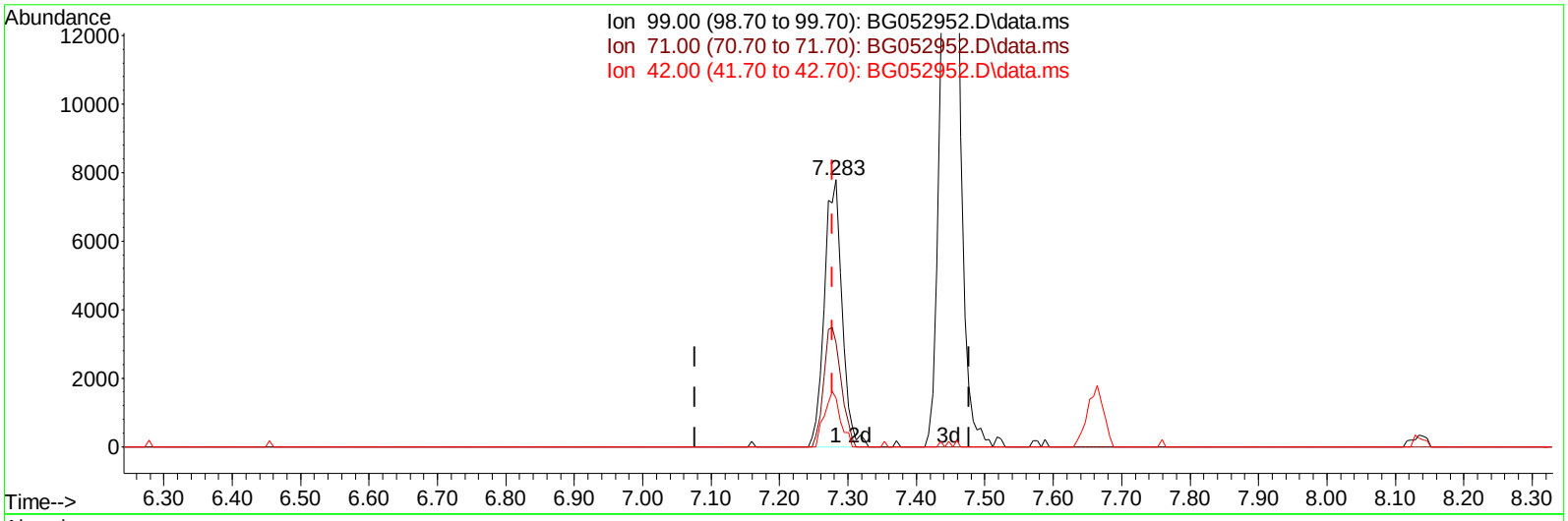
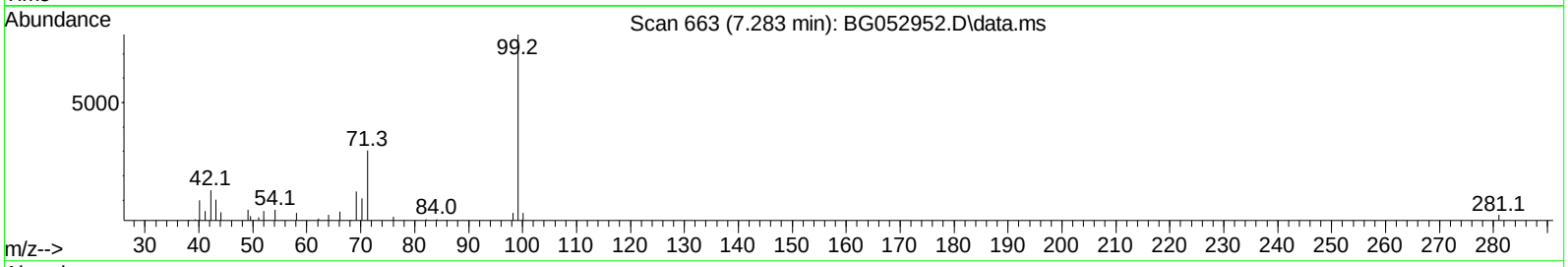
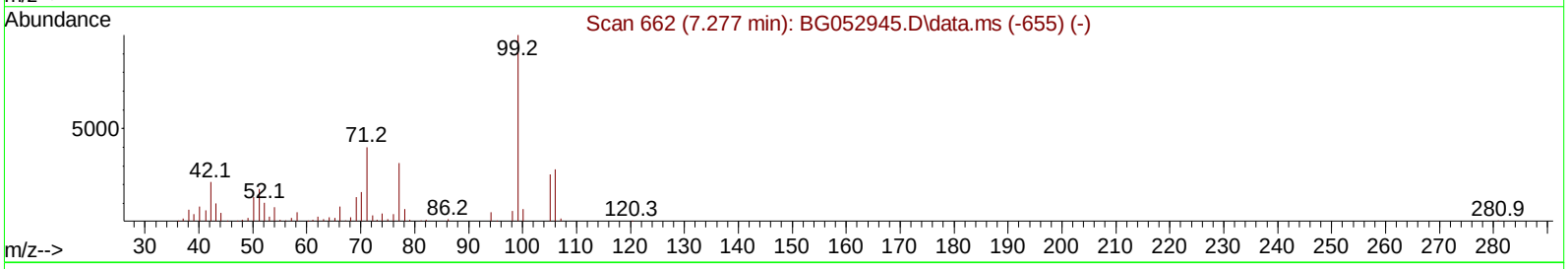
Supervised By :Yogesh Patel 04/06/2022







TIC: BG052952.D\data.ms

(7) Phenol-d5 (S)

7.283min (+ 0.006) 6.82 ng/ul m

response 13973

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	36.40	39.14
42.00	19.00	18.24
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 04/06/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.135	152	22883	20.000	ng/ul		0.00
20) Naphthalene-d8	10.943	136	99663	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.749	164	83861	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.493	188	220697	20.000	ng/ul		0.00
79) Chrysene-d12	21.775	240	243938	20.000	ng/ul	#	0.00
88) Perylene-d12	25.082	264	249880	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.553	96	2997	4.626	ng/uL		0.00
4) Pyridine-d5	3.976	84	10915m	6.749	ng/ul		0.01
7) Phenol-d5	7.283	99	13973m	6.818	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.447	67	39170	31.292	ng/ul		0.00
11) 2-Chlorophenol-d4	7.665	132	32593	23.420	ng/ul		0.00
15) 4-Methylphenol-d8	8.828	113	24964	16.815	ng/ul		0.00
21) Nitrobenzene-d5	9.292	128	22898	31.917	ng/ul		0.00
24) 2-Nitrophenol-d4	10.014	143	24469	32.337	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.555	165	43081	26.170	ng/ul		0.00
31) 4-Chloroaniline-d4	11.072	131	59244	27.230	ng/ul		0.00
46) Dimethylphthalate-d6	14.144	166	211918	32.893	ng/ul		0.00
49) Acenaphthylene-d8	14.444	160	223606	28.562	ng/ul		0.00
54) 4-Nitrophenol-d4	14.920	143	8657	7.992	ng/ul		0.00
60) Fluorene-d10	15.742	176	188154	33.261	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.848	200	39415	35.094	ng/ul		0.00
73) Anthracene-d10	17.592	188	364257	35.444	ng/ul		0.00
81) Pyrene-d10	19.872	212	465888	34.140	ng/ul		0.00
92) Benzo(a)pyrene-d12	24.859	264	452434	35.042	ng/ul		0.00

Target Compounds Qvalue

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

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