Quantitation Report (Qedit)

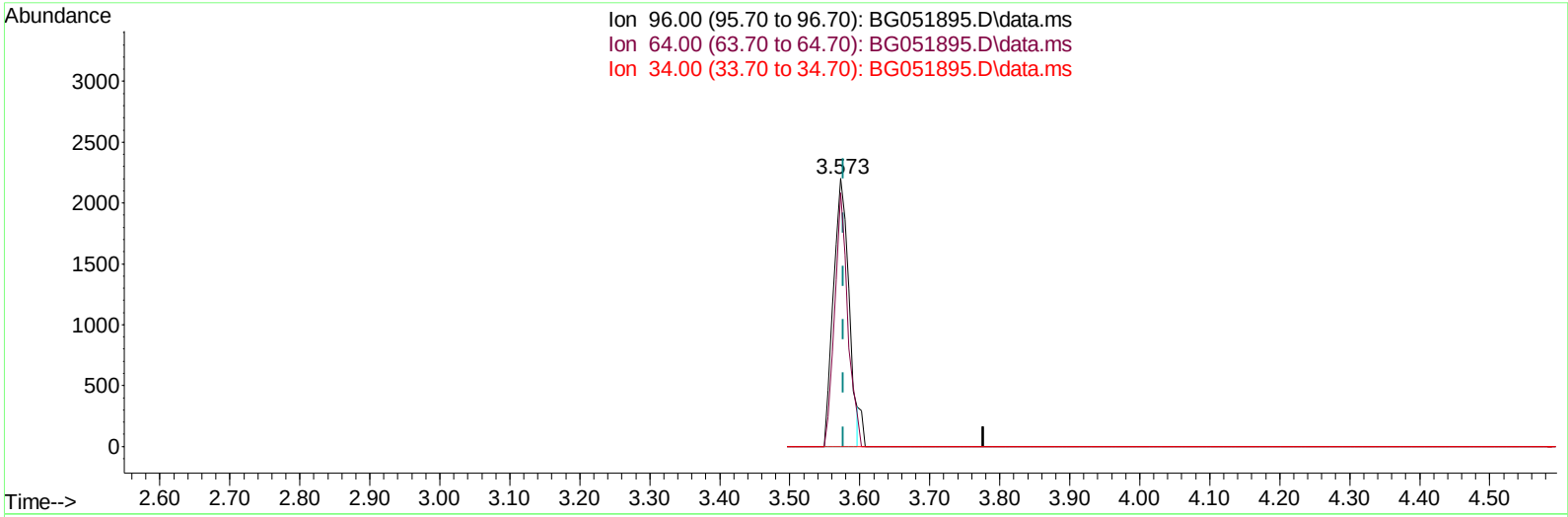
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010622\
Data File : BG051895.D
Acq On : 6 Jan 2022 16:25
Operator : CG/JU
Sample : N1026-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

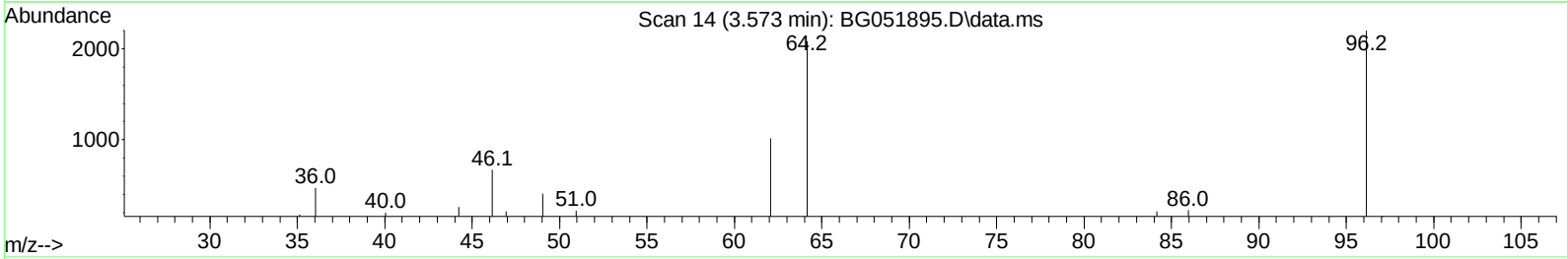
Quant Time: Jan 06 23:56:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 06 14:43:02 2022
Response via : Initial Calibration

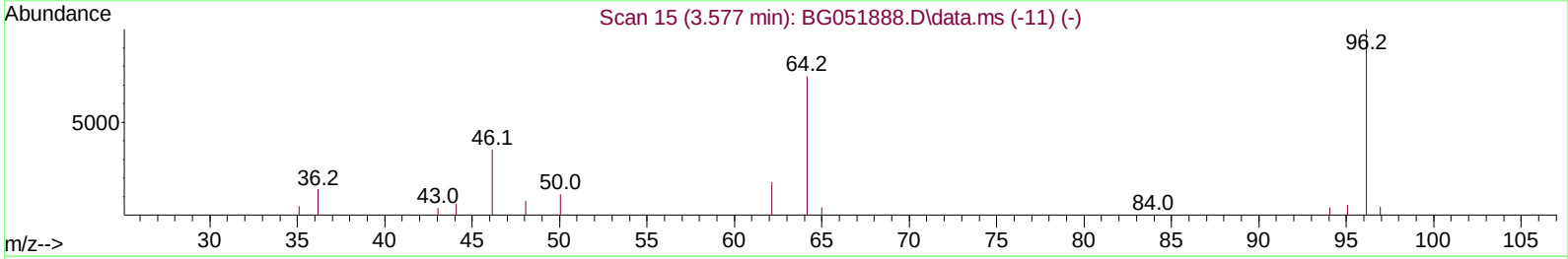
Instrument :
BNA_G
ClientSampleId :
BGLH0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/07/2022
Supervised By :mohammad ahmed 01/12/2022







TIC: BG051895.D\data.ms

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

3.573min (-0.004) 5.34 ng/uL

response 3323

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	94.73#
34.00	0.00	0.00
0.00	0.00	0.00

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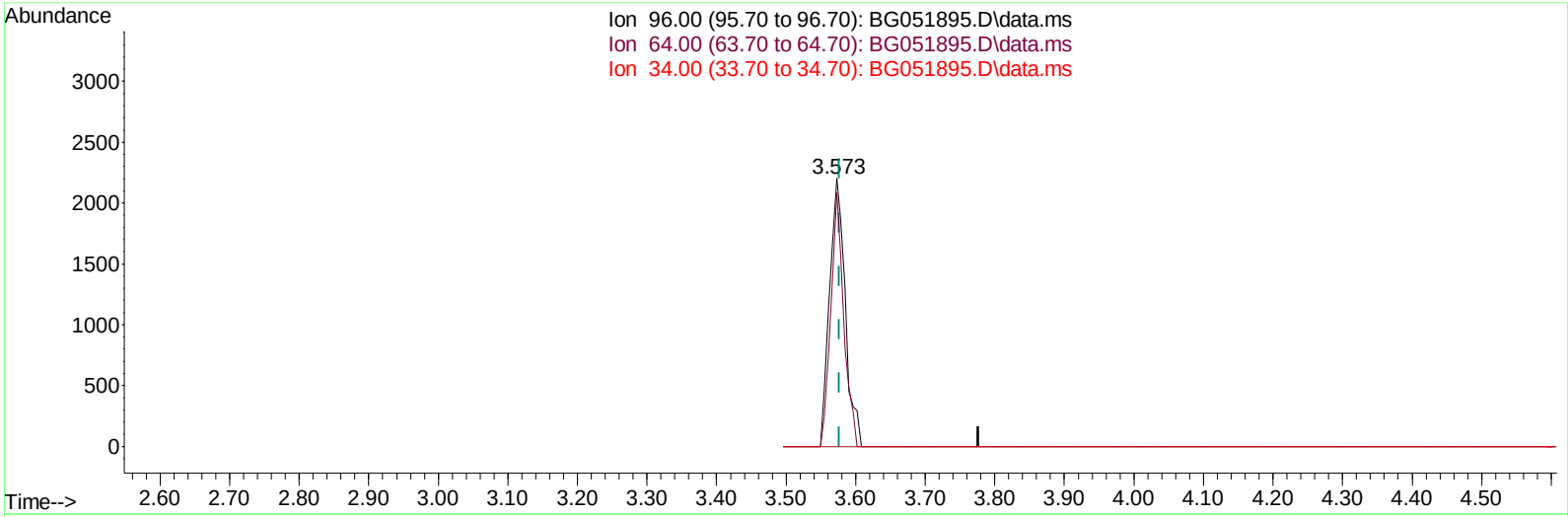

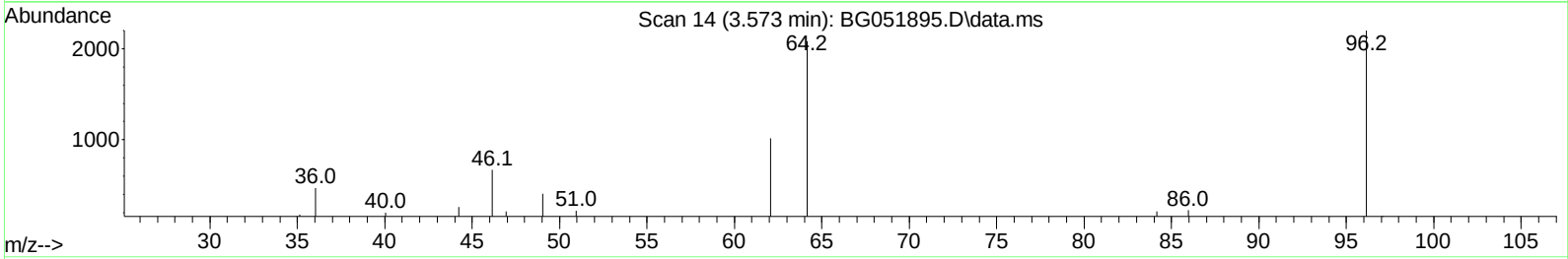

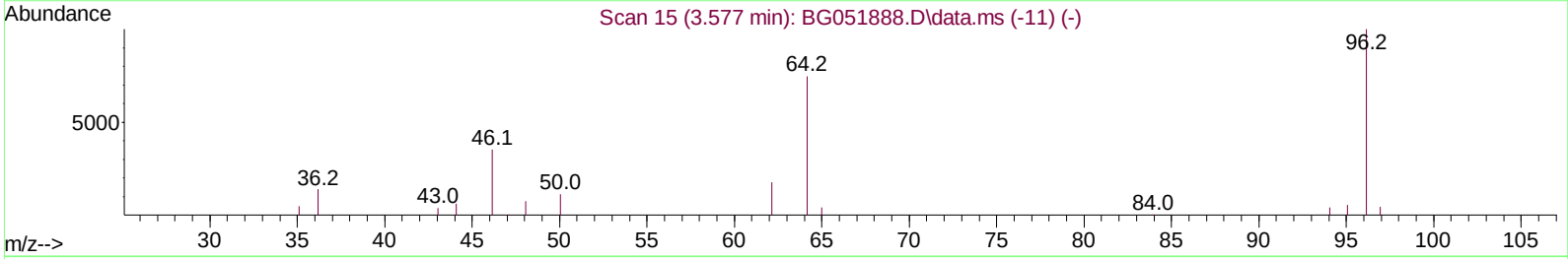

TIC: BG051895.D\data.ms

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

3.573min (-0.004) 5.51 ng/uL m

response 3426

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	94.73#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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ALS Vial : 11 Sample Multiplier: 1

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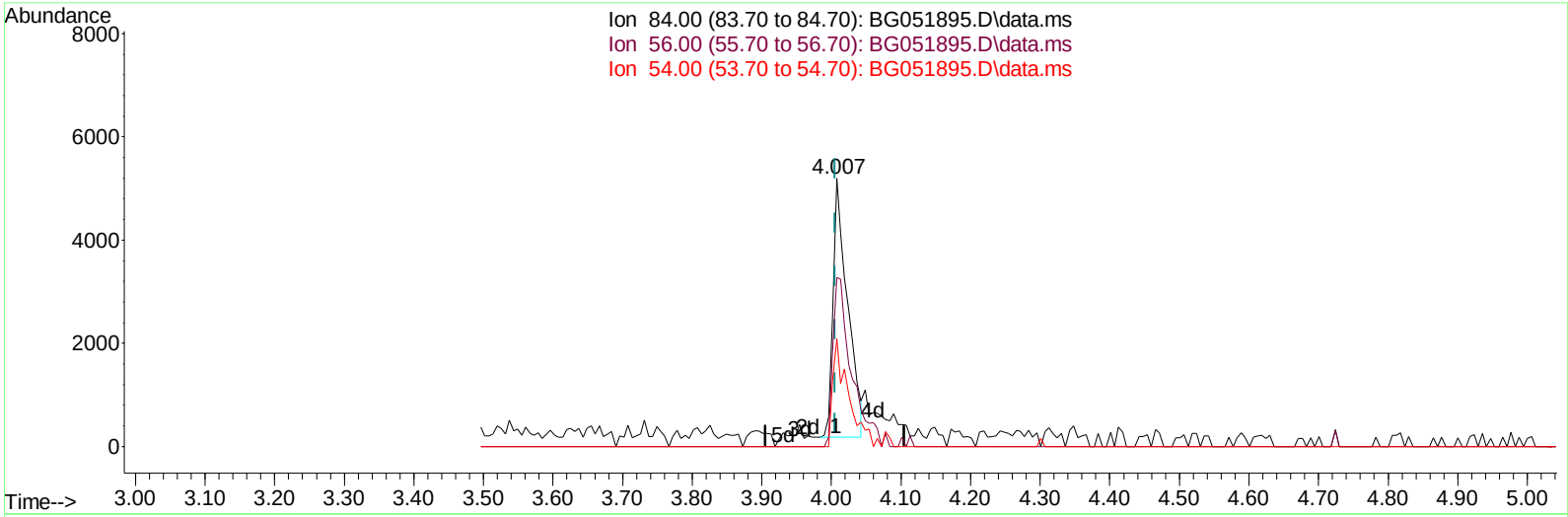

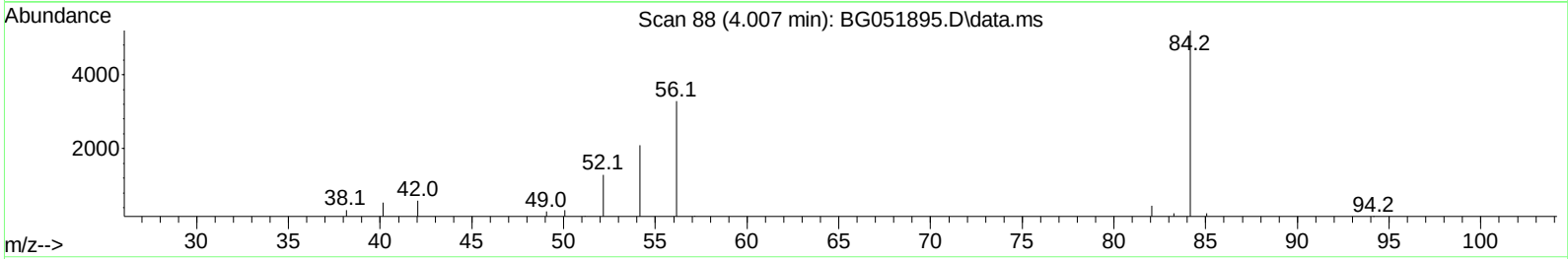

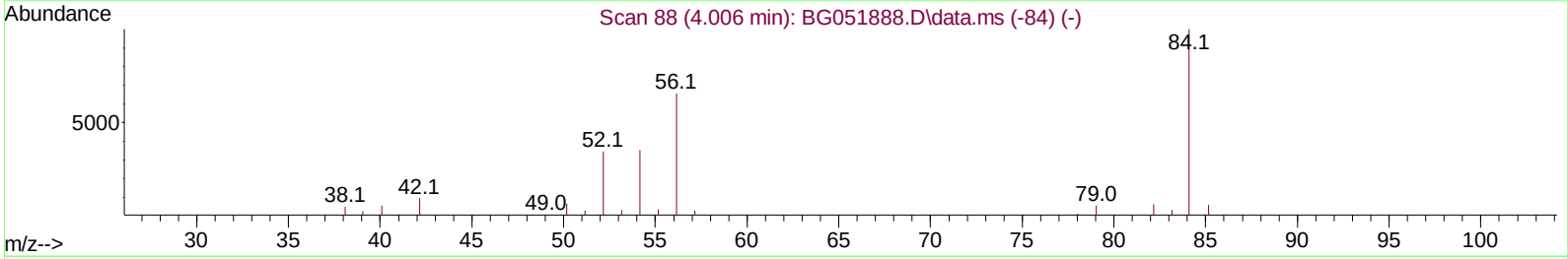

TIC: BG051895.D\data.ms

(4) Pyridine-d5 (S)

4.007min (+ 0.002) 3.80 ng/ul

response 7343

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	63.18
54.00	31.50	40.27#
0.00	0.00	0.00

Quantitation Report (Qedit)

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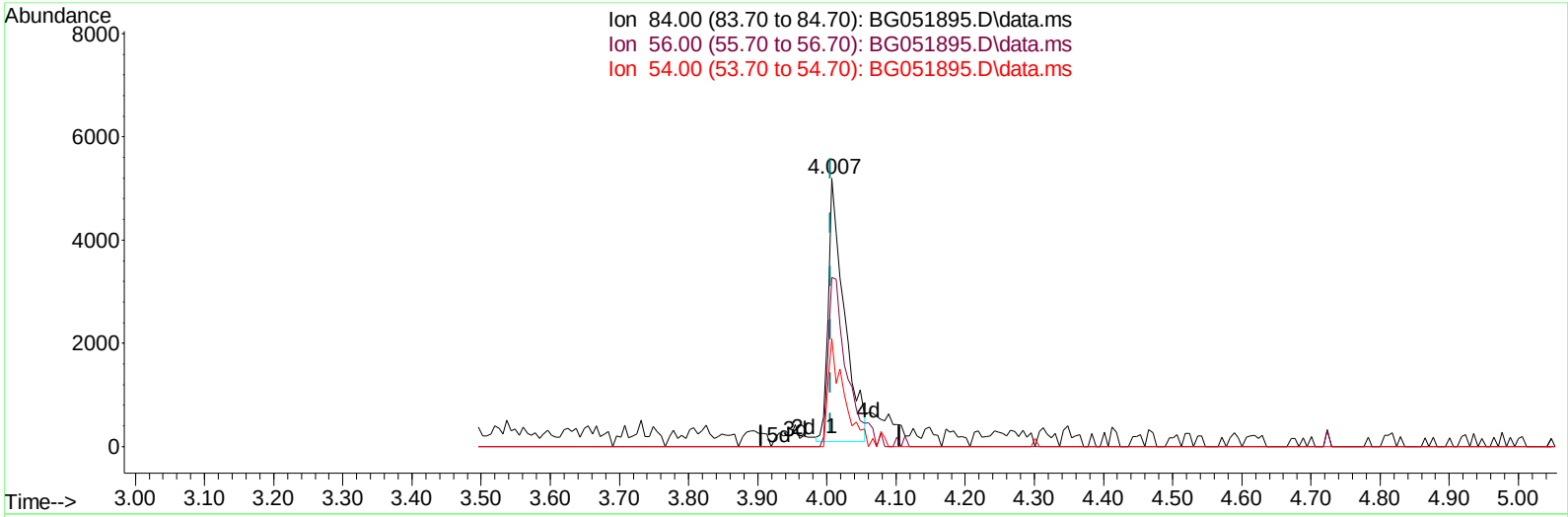

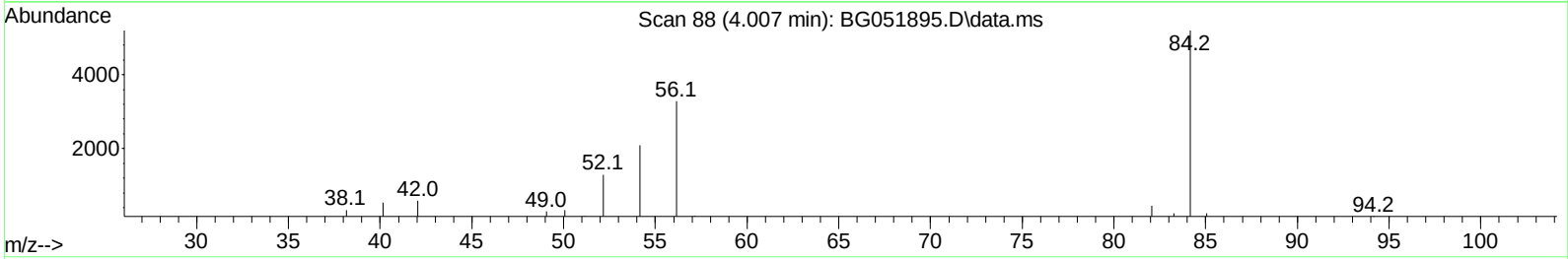

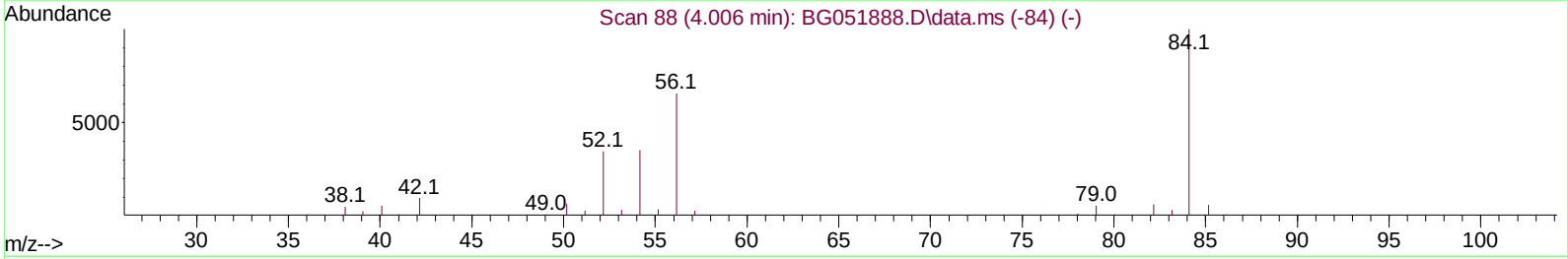

TIC: BG051895.D\data.ms

(4) Pyridine-d5 (S)

4.007min (+ 0.002) 4.23 ng/ul m

response 8188

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	63.18
54.00	31.50	40.27#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.249	152	22228	20.000	ng/ul	0.00
20) Naphthalene-d8	11.086	136	98090	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.893	164	72242	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.636	188	186880	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.954	240	209343	20.000	ng/ul	# 0.00
88) Perylene-d12	25.443	264	211025	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.573	96	3426m	5.507	ng/uL	0.00
4) Pyridine-d5	4.007	84	8188m	4.234	ng/ul	0.00
7) Phenol-d5	7.385	99	14870	6.835	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	45937	33.263	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.773	132	36915	25.629	ng/ul	0.00
15) 4-Methylphenol-d8	8.948	113	26725	15.803	ng/ul	0.00
21) Nitrobenzene-d5	9.418	128	25690	33.145	ng/ul	0.00
24) 2-Nitrophenol-d4	10.146	143	27879	32.523	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.698	165	47927	28.604	ng/ul	0.00
31) 4-Chloroaniline-d4	11.209	131	53444	22.904	ng/ul	0.00
46) Dimethylphthalate-d6	14.282	166	200522	33.581	ng/ul	0.00
49) Acenaphthylene-d8	14.587	160	238598	33.079	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	7738	7.751	ng/ul	0.00
60) Fluorene-d10	15.879	176	177928	34.176	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.979	200	38938	33.028	ng/ul	0.00
73) Anthracene-d10	17.736	188	340994	38.464	ng/ul	0.00
81) Pyrene-d10	20.015	212	453077	37.606	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.202	264	426480	38.389	ng/ul	0.00

Target Compounds Qvalue

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

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