Quantitation Report (Qedit)

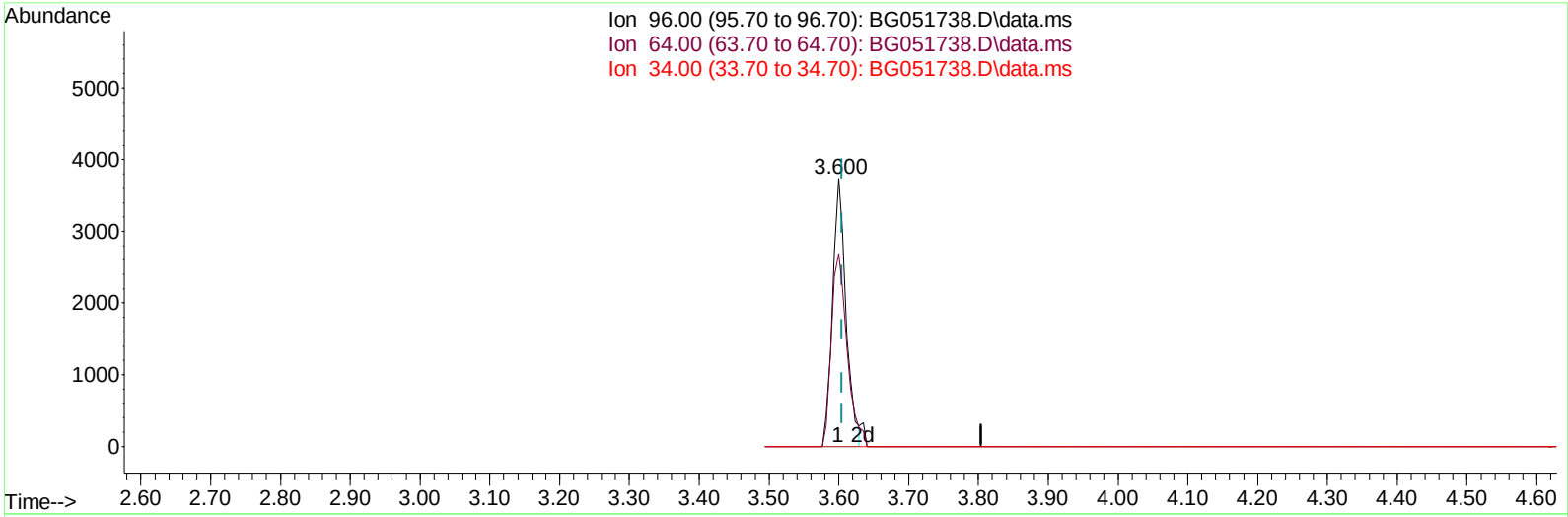
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
Data File : BG051738.D
Acq On : 22 Dec 2021 6:24
Operator : CG/JU
Sample : M5090-12
Misc :
ALS Vial : 66 Sample Multiplier: 1

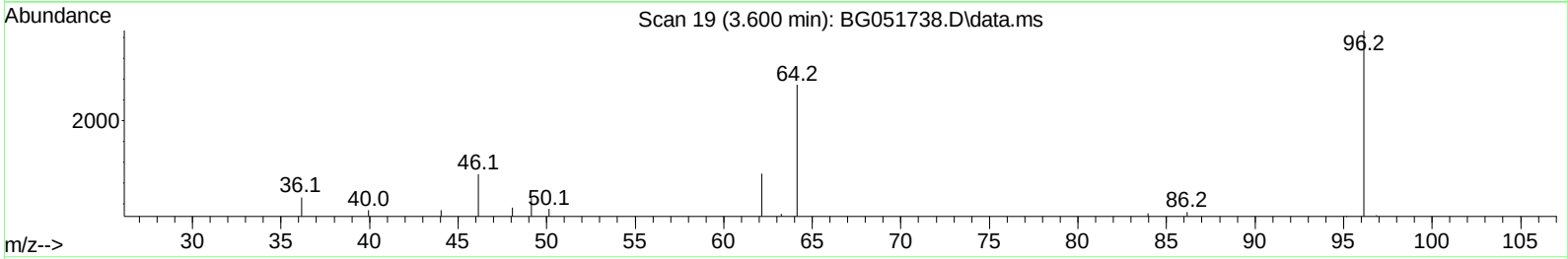
Quant Time: Dec 22 07:24:38 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 14 14:19:57 2021
Response via : Initial Calibration

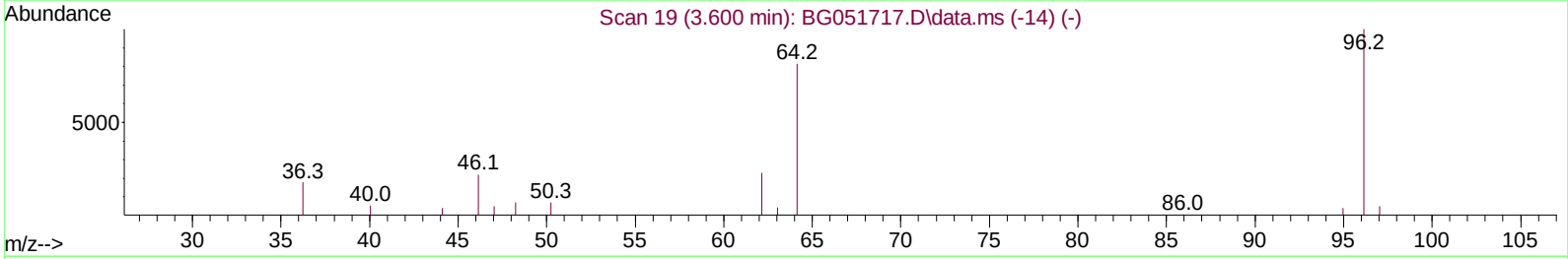
Instrument :
BNA_G
ClientSampleId :
BG3E2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/22/2021
Supervised By :mohammad ahmed 12/28/2021







TIC: BG051738.D\data.ms

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

3.600min (-0.005) 5.44 ng/uL

response 5016

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.03
34.00	0.00	0.00
0.00	0.00	0.00

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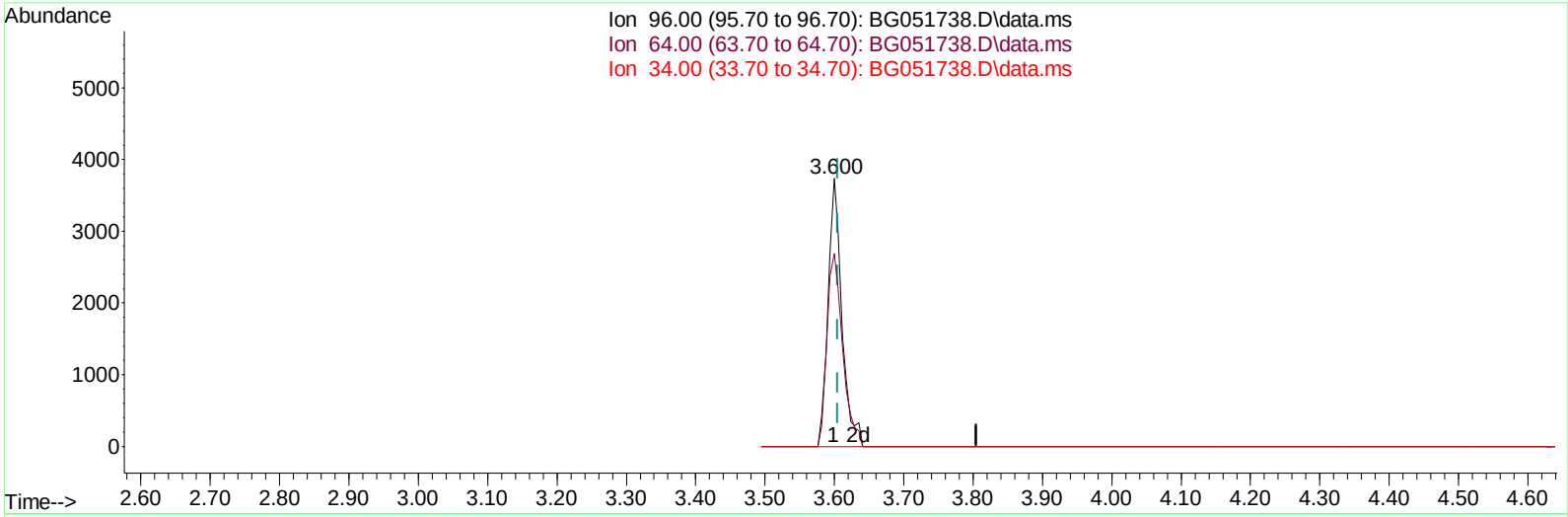

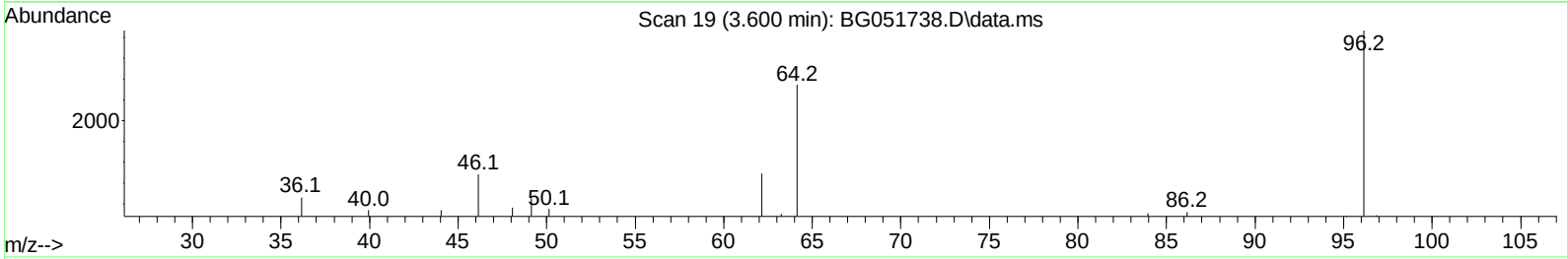

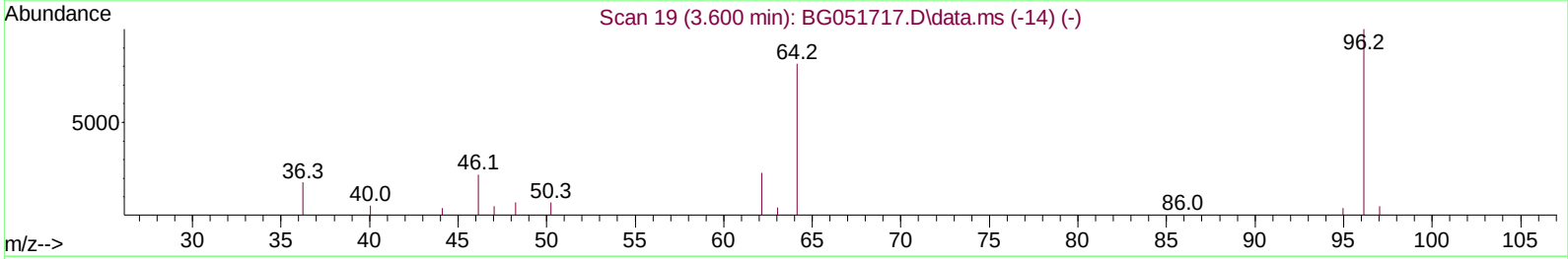

TIC: BG051738.D\data.ms

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

3.600min (-0.005) 5.57 ng/uL m

response 5131

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	72.03
34.00	0.00	0.00
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	8.276	152	38281	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.113	136	168859	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.908	164	122453	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.657	188	283281	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.975	240	273020	20.000	ng/ul	#-0.01
88) Perylene-d12	25.476	264	263712	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	5131m	5.566	ng/uL	0.00
4) Pyridine-d5	4.028	84	21478	7.712	ng/ul	-0.02
7) Phenol-d5	7.406	99	28053	8.207	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.577	67	76982	37.886	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.800	132	70045	29.465	ng/ul	-0.01
15) 4-Methylphenol-d8	8.969	113	49772	18.839	ng/ul	-0.01
21) Nitrobenzene-d5	9.445	128	46478	37.543	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.173	143	48723	36.002	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	91979	31.340	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.230	131	112102	29.011	ng/ul	-0.01
46) Dimethylphthalate-d6	14.297	166	347044	35.409	ng/ul	-0.02
49) Acenaphthylene-d8	14.602	160	418734	34.400	ng/ul	-0.01
54) 4-Nitrophenol-d4	15.066	143	8696	5.456	ng/ul	-0.01
60) Fluorene-d10	15.901	176	308485	35.131	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.995	200	30143	20.423	ng/ul	-0.02
73) Anthracene-d10	17.757	188	517602	38.303	ng/ul	0.00
81) Pyrene-d10	20.030	212	608709	37.045	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.241	264	559588	40.329	ng/ul	-0.01
Target Compounds						Qvalue
16) Acetophenone	9.098	105	6792	1.637	ng/ul	96
35) 4-Chloro-3-methylphenol	12.358	107	16154	5.163	ng/ul	95

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

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