Quantitation Report (Qedit)

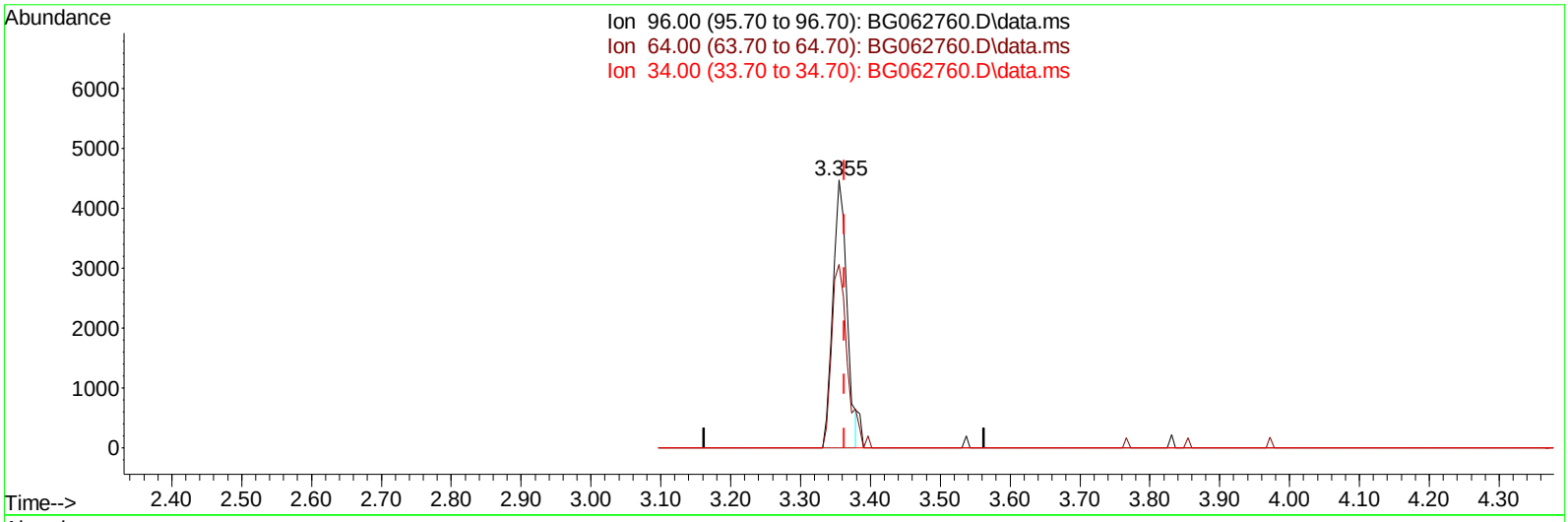
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
Data File : BG062760.D
Acq On : 12 Sep 2024 12:17
Operator : JU/RC
Sample : P3922-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

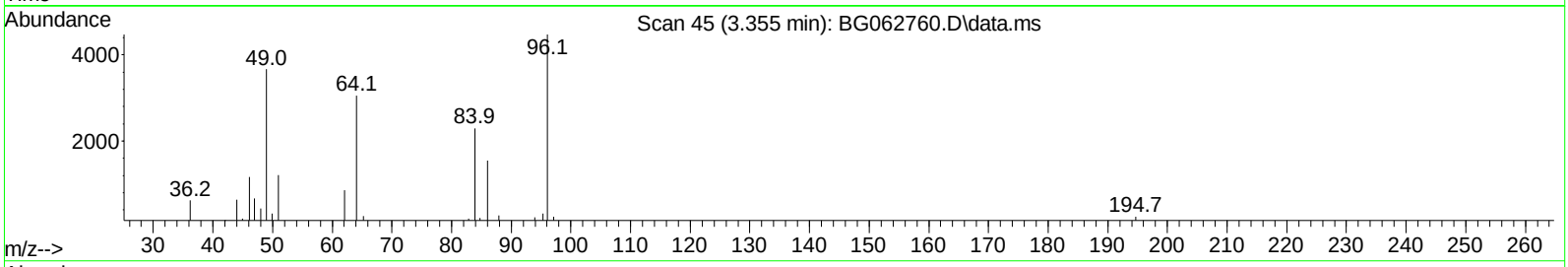
Quant Time: Sep 12 14:42:39 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
Response via : Initial Calibration

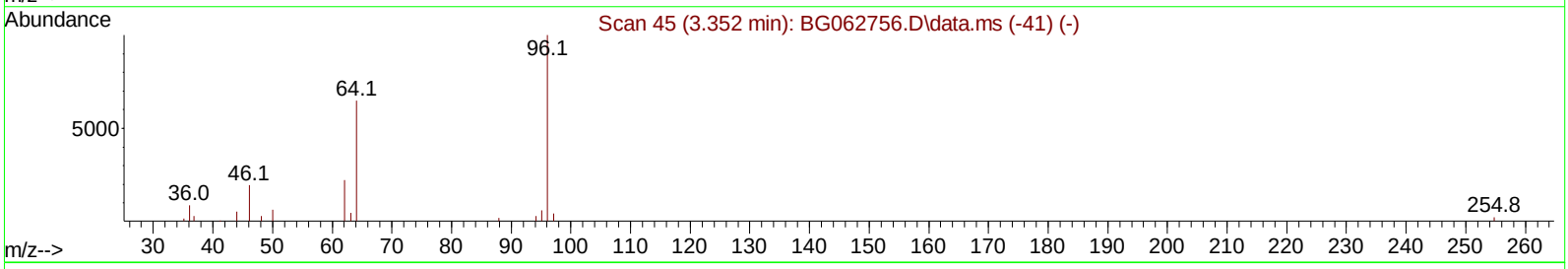
Instrument :
BNA_G
ClientSampleId :
C0AN7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/13/2024
Supervised By :mohammad ahmed 09/14/2024







TIC: BG062760.D\data.ms

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

3.355min (-0.008) 5.72 ng/uL

response 6120

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	68.40
34.00	0.00	0.00
0.00	0.00	0.00

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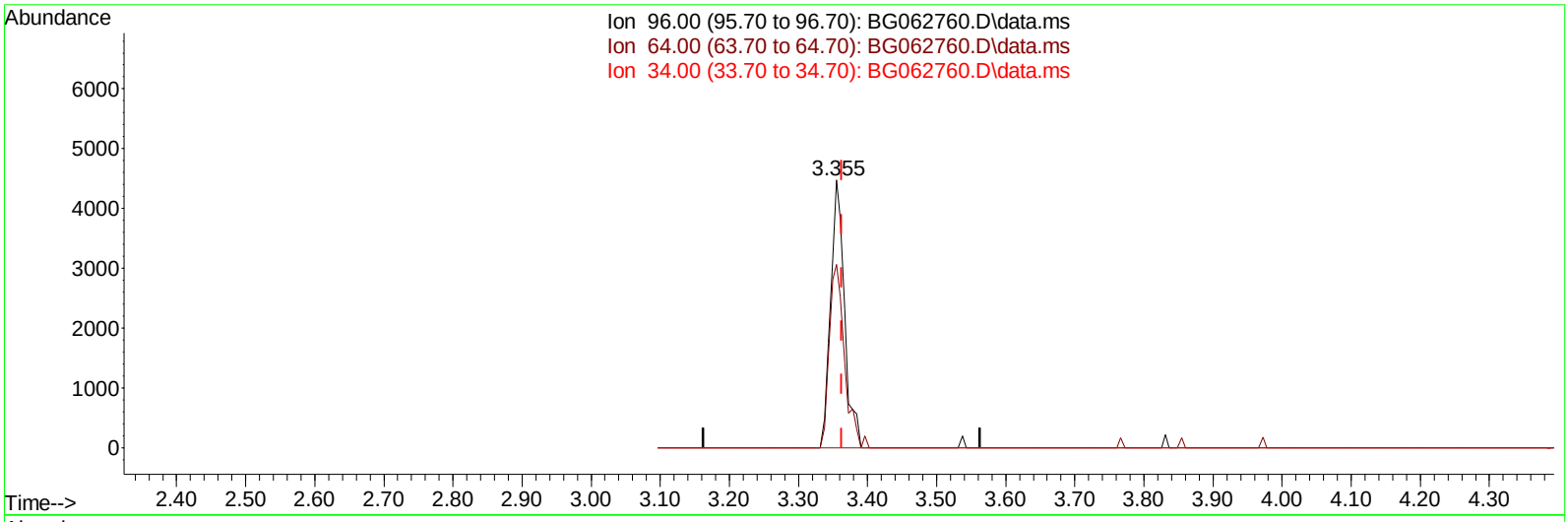

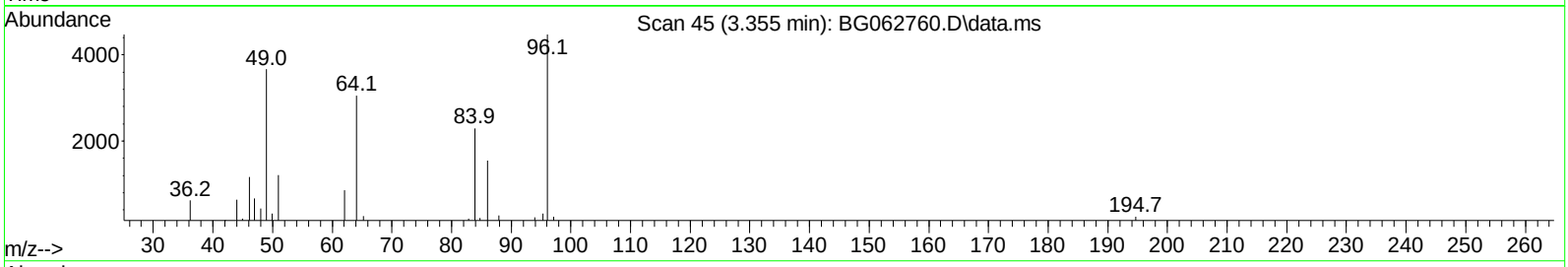

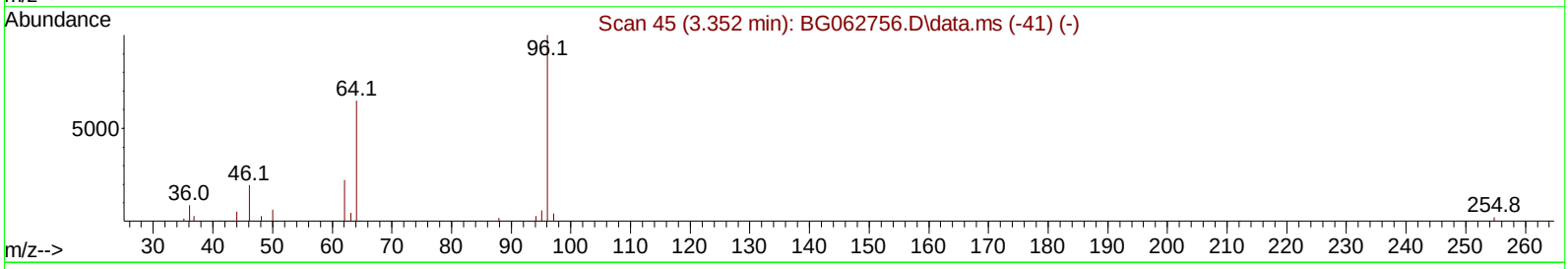

TIC: BG062760.D\data.ms

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

3.355min (-0.008) 5.90 ng/uL m

response 6320

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	68.40
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Acq On : 12 Sep 2024 12:17
Operator : JU/RC
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Misc :
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Quant Time: Sep 12 14:44:20 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.897	152	48399	20.000	ng/ul	0.00
20) Naphthalene-d8	10.694	136	264993	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.530	164	212882	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.274	188	542183	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	544090	20.000	ng/ul	0.00
88) Perylene-d12	24.583	264	599910	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.355	96	6320m	5.903	ng/uL	0.00
4) Pyridine-d5	3.760	84	47730	14.005	ng/ul	0.00
7) Phenol-d5	7.063	99	46229	10.893	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.227	67	53248	20.927	ng/ul	0.00
11) 2-Chlorophenol-d4	7.427	132	72918	23.588	ng/ul	0.00
15) 4-Methylphenol-d8	8.596	113	82169	23.233	ng/ul	-0.01
21) Nitrobenzene-d5	9.048	128	38792	20.482	ng/ul	0.00
24) 2-Nitrophenol-d4	9.771	143	50411	24.338	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.312	165	116837	25.901	ng/ul	0.00
31) 4-Chloroaniline-d4	10.823	131	107659	17.363	ng/ul	0.00
46) Dimethylphthalate-d6	13.937	166	404685	24.686	ng/ul	0.00
49) Acenaphthylene-d8	14.219	160	422362	23.198	ng/ul	0.00
54) 4-Nitrophenol-d4	14.730	143	26222	9.416	ng/ul	0.00
60) Fluorene-d10	15.523	176	357098	24.173	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.635	200	71195	22.828	ng/ul	0.00
73) Anthracene-d10	17.374	188	672529	26.283	ng/ul	0.00
81) Pyrene-d10	19.665	212	813312	26.564	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.378	264	859112	27.121	ng/ul	0.00

Target Compounds Qvalue

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

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