

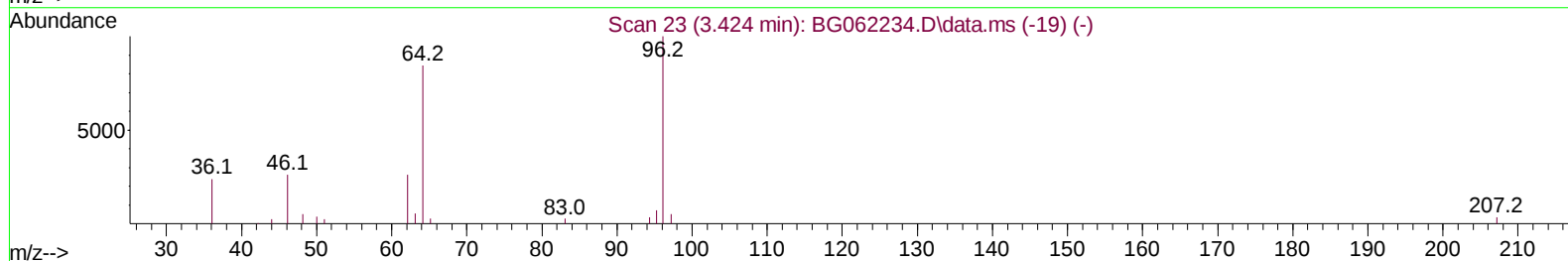
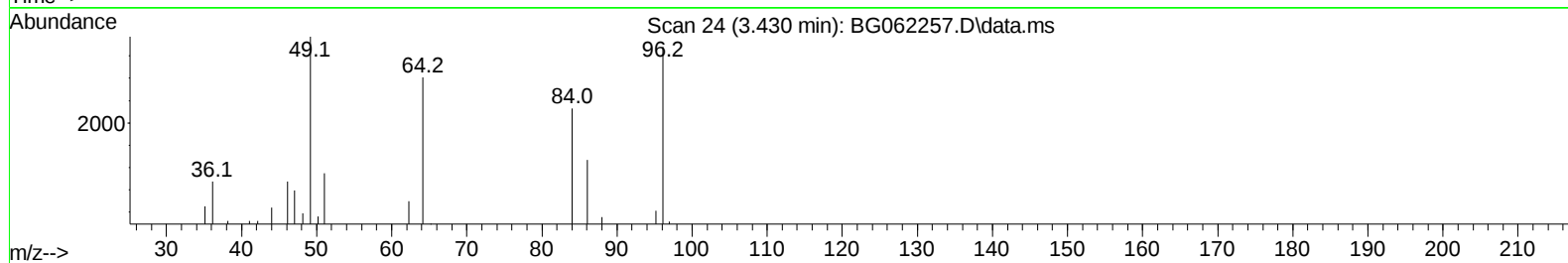
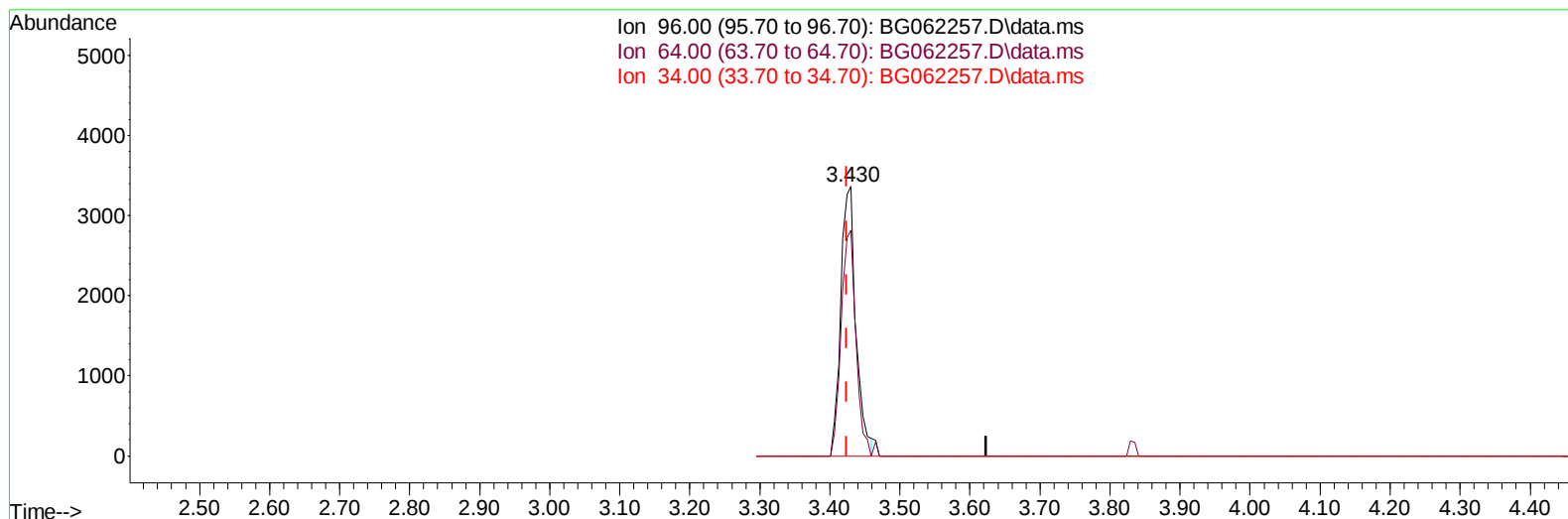
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080524\
Data File : BG062257.D
Acq On : 6 Aug 2024 5:03
Operator : MA/JU
Sample : P3361-20
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E20Q0

Manual IntegrationsAPPROVED

Quant Time: Aug 06 05:49:54 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 06 02:21:27 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024
Supervised By :mohammad ahmed 08/08/2024



TIC: BG062257.D\data.ms

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

3.430min (+ 0.006) 3.27 ng/uL

response 5144

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	84.50	83.94
34.00	0.00	0.00
0.00	0.00	0.00

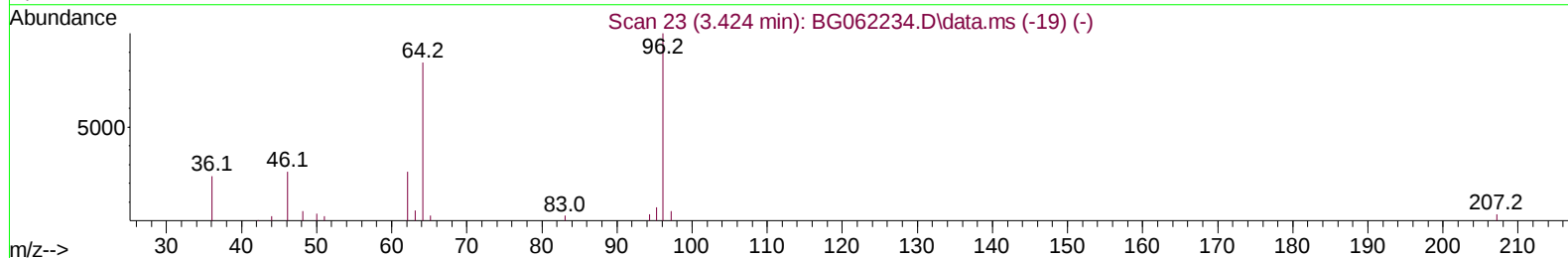
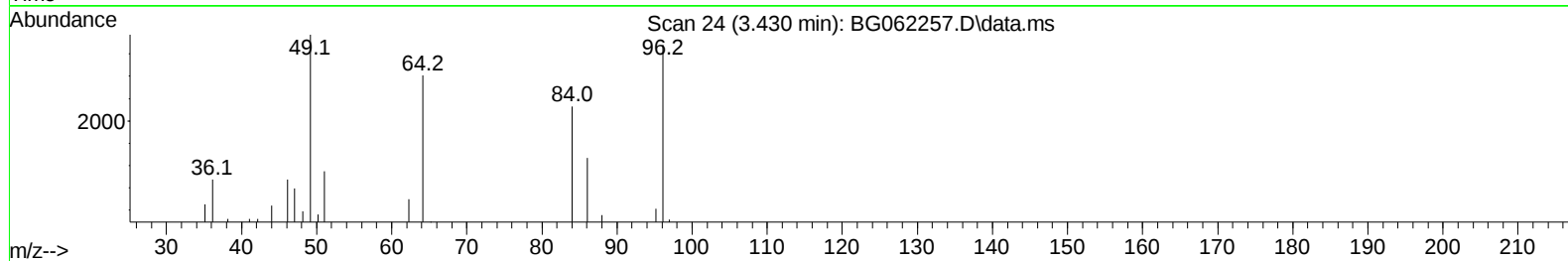
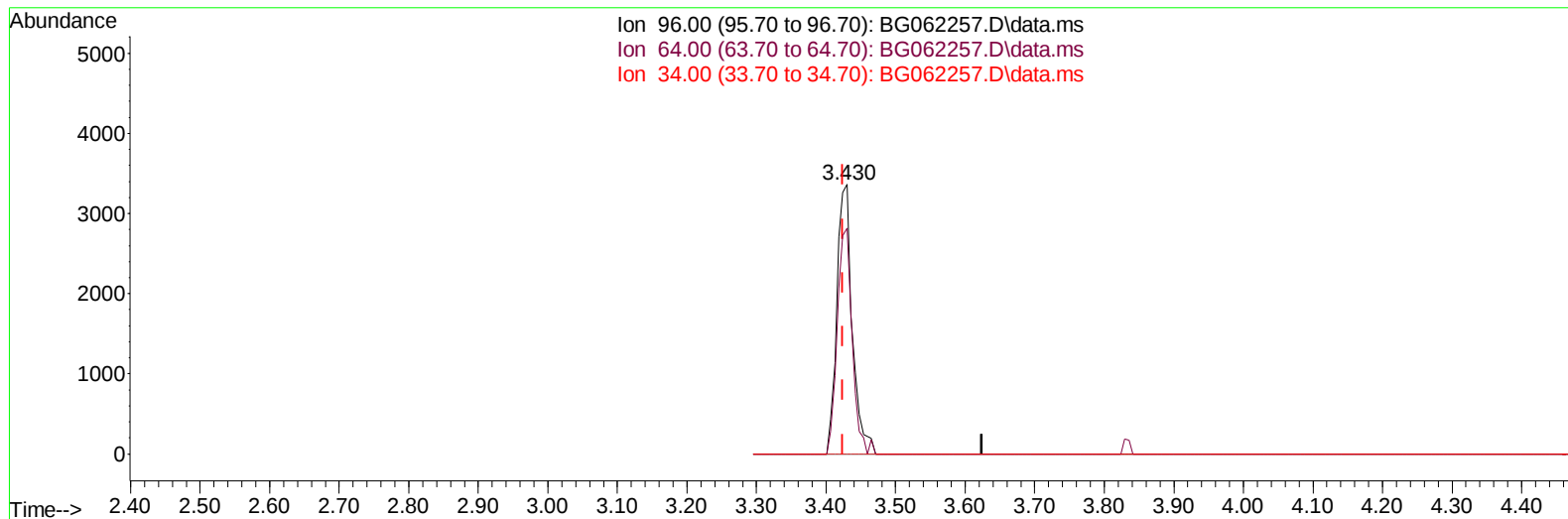
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TIC: BG062257.D\data.ms

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

3.430min (+ 0.006) 3.31 ng/uL m

response 5211

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	84.50	83.94
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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 ClientSampleId :
 E20Q0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/06/2024
 Supervised By :mohammad ahmed 08/08/2024

Quant Time: Aug 06 05:58:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 06 02:21:27 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.965	152	67791	20.000	ng/ul	0.00
20) Naphthalene-d8	10.755	136	318379	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.580	164	257735	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.323	188	574042	20.000	ng/ul	0.00
79) Chrysene-d12	21.564	240	560595	20.000	ng/ul	-0.01
88) Perylene-d12	24.660	264	646352	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.430	96	5211m	3.314	ng/uL	0.00
4) Pyridine-d5	3.829	84	83103	16.802	ng/ul	0.00
7) Phenol-d5	7.113	99	117417	18.147	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.290	67	74377	18.673	ng/ul	0.00
11) 2-Chlorophenol-d4	7.489	132	88402	19.069	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	101092	18.757	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	42908	21.856	ng/ul	0.00
24) 2-Nitrophenol-d4	9.833	143	43462	23.534	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.368	165	100792	18.791	ng/ul	0.00
31) 4-Chloroaniline-d4	10.885	131	144534	18.578	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	364602	18.335	ng/ul	0.00
49) Acenaphthylene-d8	14.274	160	419200	18.593	ng/ul	0.00
54) 4-Nitrophenol-d4	14.750	143	42620	15.309	ng/ul	-0.01
60) Fluorene-d10	15.572	176	332519	18.630	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.678	200	24860	14.623	ng/ul	0.00
73) Anthracene-d10	17.423	188	512062	18.495	ng/ul	0.00
81) Pyrene-d10	19.702	212	632049	18.969	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.449	264	656122	18.989	ng/ul	-0.02

Target Compounds Qvalue

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

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