

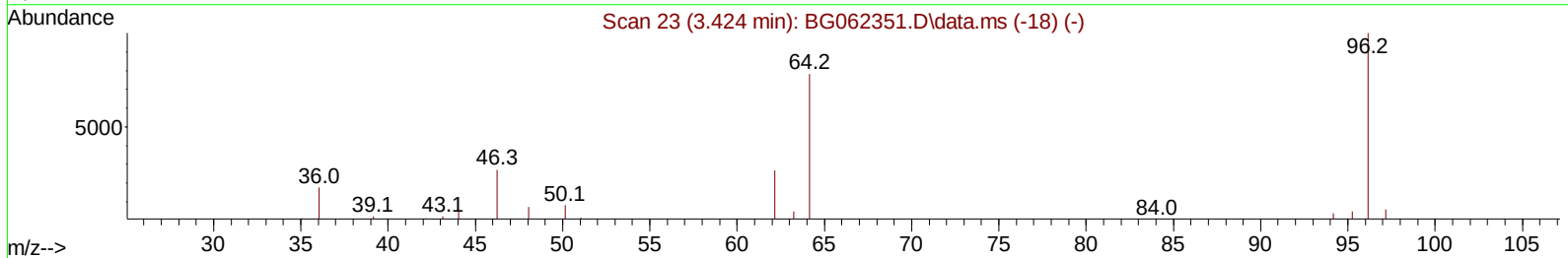
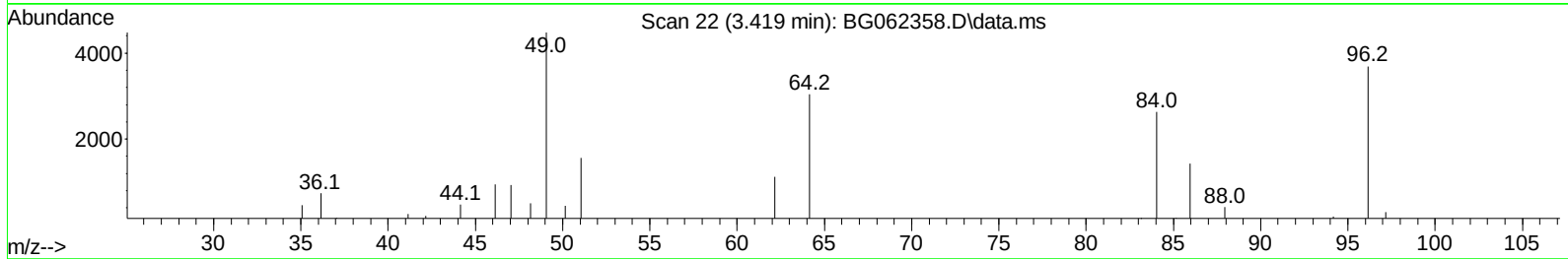
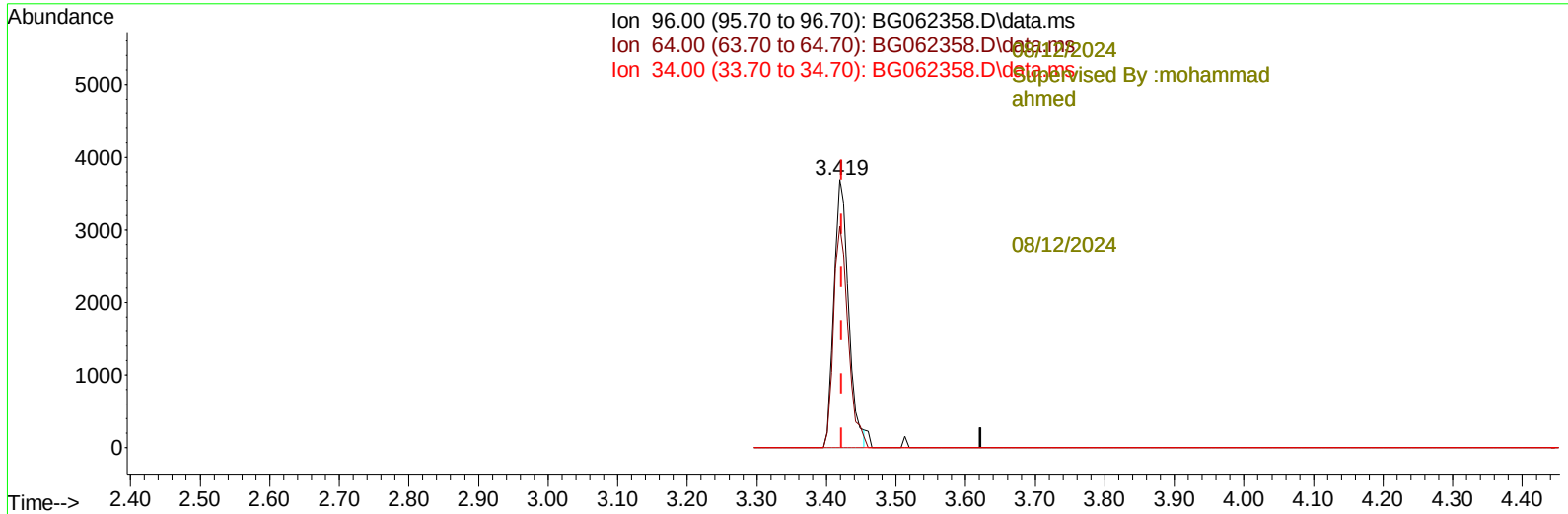
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062358.D
Acq On : 9 Aug 2024 13:16
Operator : MA/JU
Sample : P3480-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E0AD5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/12/2024
Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 09 14:51:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 09 01:00:03 2024
Response via : Initial Calibration



TIC: BG062358.D\data.ms

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

3.419min (-0.004) 4.88 ng/uL

response 5458

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	84.50	82.63
34.00	0.00	0.00
0.00	0.00	0.00

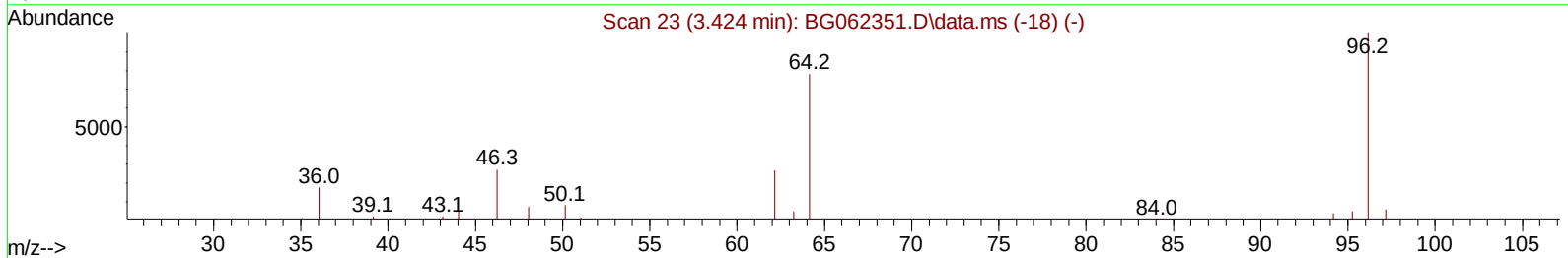
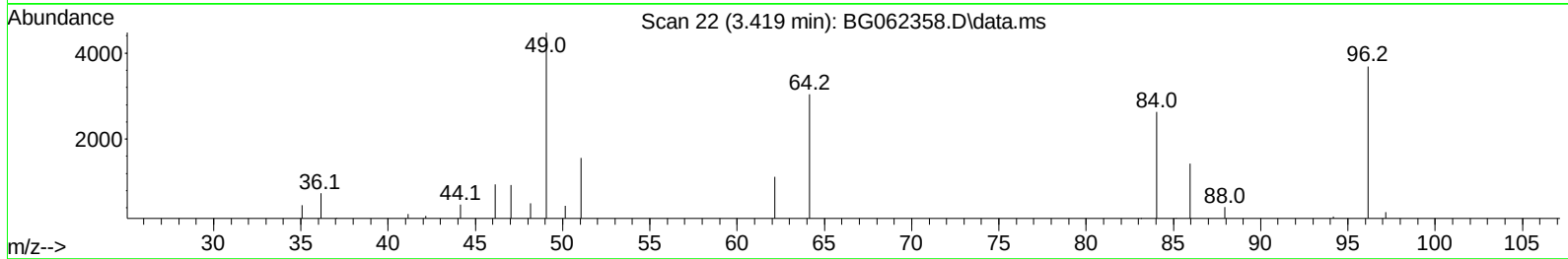
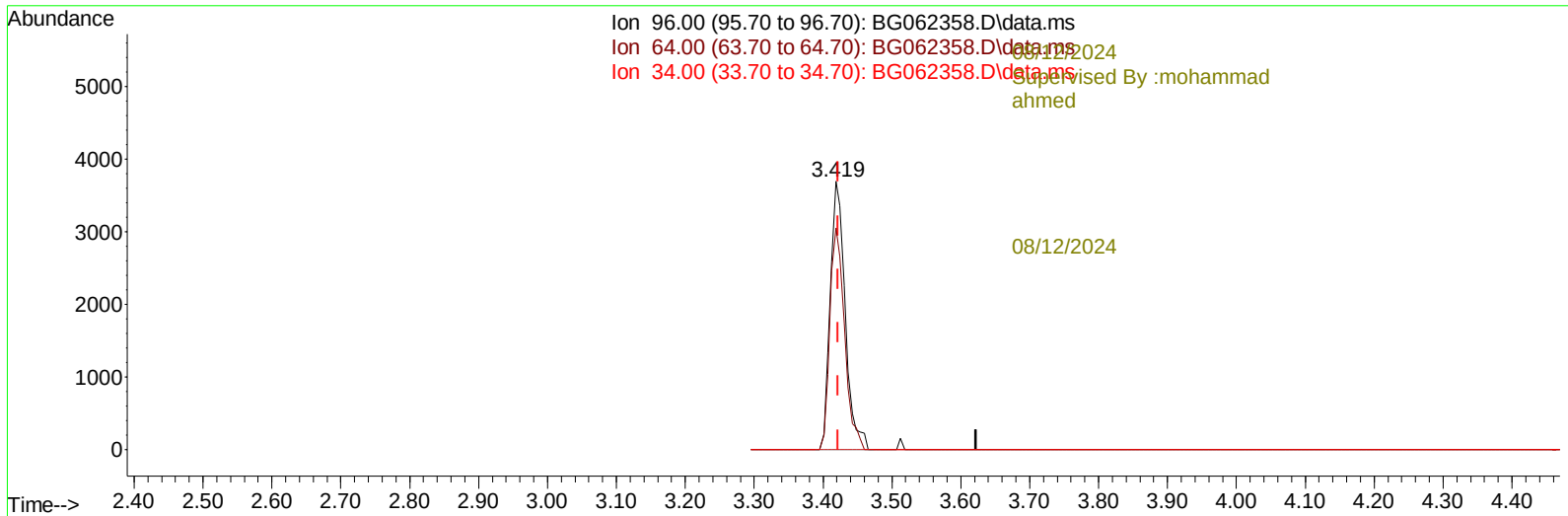
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TIC: BG062358.D\data.ms

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

3.419min (-0.004) 4.95 ng/uL m

response 5538

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	84.50	82.63
34.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/12/2024
 Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 10 01:39:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:00:03 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----08/12/2024							
Supervised By :mohammad							
ahmed							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.960	152		48223	20.000	ng/ul	0.00
20) Naphthalene-d8	10.750	136		247289	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.574	164		197751	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.312	188		430205	20.000	ng/ul	0.00
79) Chrysene-d12	21.553	240		425196	20.000	ng/ul	0.00
88) Perylene-d12	24.643	264		493018	20.000	ng/ul	0.00
-----08/12/2024							
Supervised By :mohammad							
ahmed							
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.419	96		5538m	4.952	ng/uL	0.00
4) Pyridine-d5	3.830	84		21603	6.140	ng/ul	0.00
7) Phenol-d5	7.108	99		44005	9.561	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.284	67		95790	33.808	ng/ul	0.00
11) 2-Chlorophenol-d4	7.490	132		105508	31.994	ng/ul	0.00
15) 4-Methylphenol-d8	8.647	113		85925	22.412	ng/ul	0.00
21) Nitrobenzene-d5	9.111	128		61092	40.065	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143		67300	46.918	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.368	165		142929	34.307	ng/ul	0.00
31) 4-Chloroaniline-d4	10.879	131		52724	8.725	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166		485672	31.832	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160		535059	30.930	ng/ul	0.00
54) 4-Nitrophenol-d4	14.750	143		17053	7.984	ng/ul	0.00
60) Fluorene-d10	15.567	176		445479	32.529	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.673	200		60123	47.189	ng/ul	0.00
73) Anthracene-d10	17.412	188		705614	34.006	ng/ul	0.00
81) Pyrene-d10	19.697	212		872884	34.539	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.431	264		940099	35.670	ng/ul	0.00

Target Compounds Qvalue

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

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