

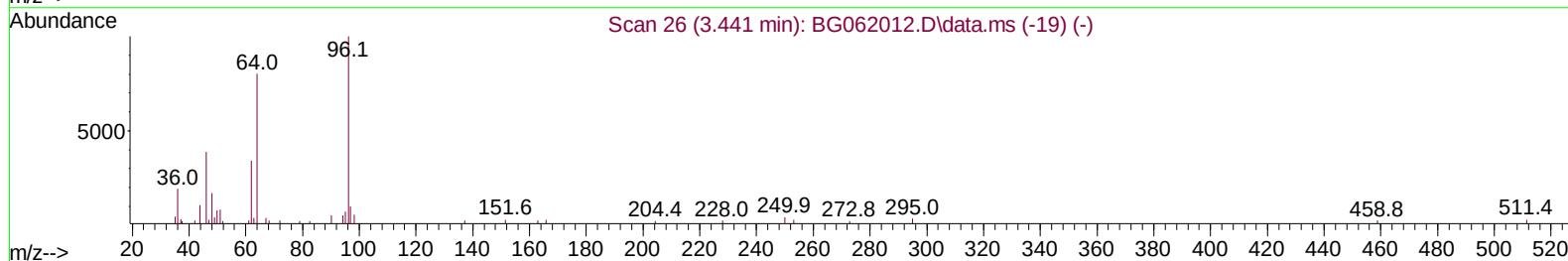
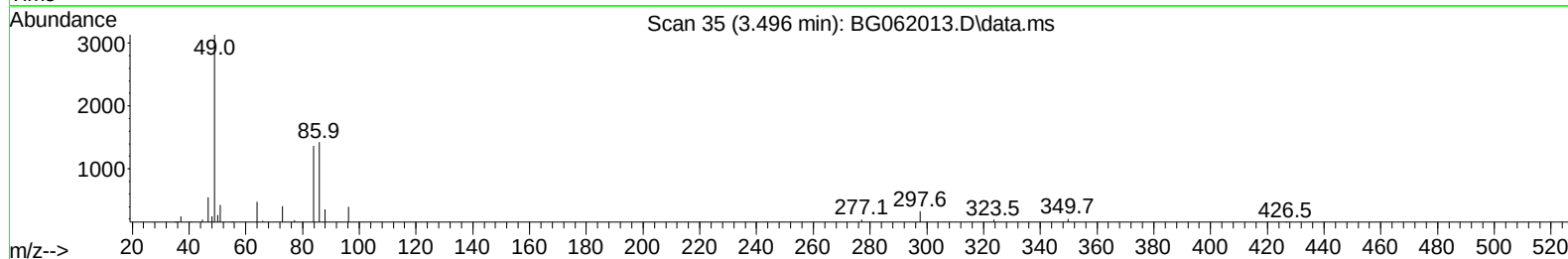
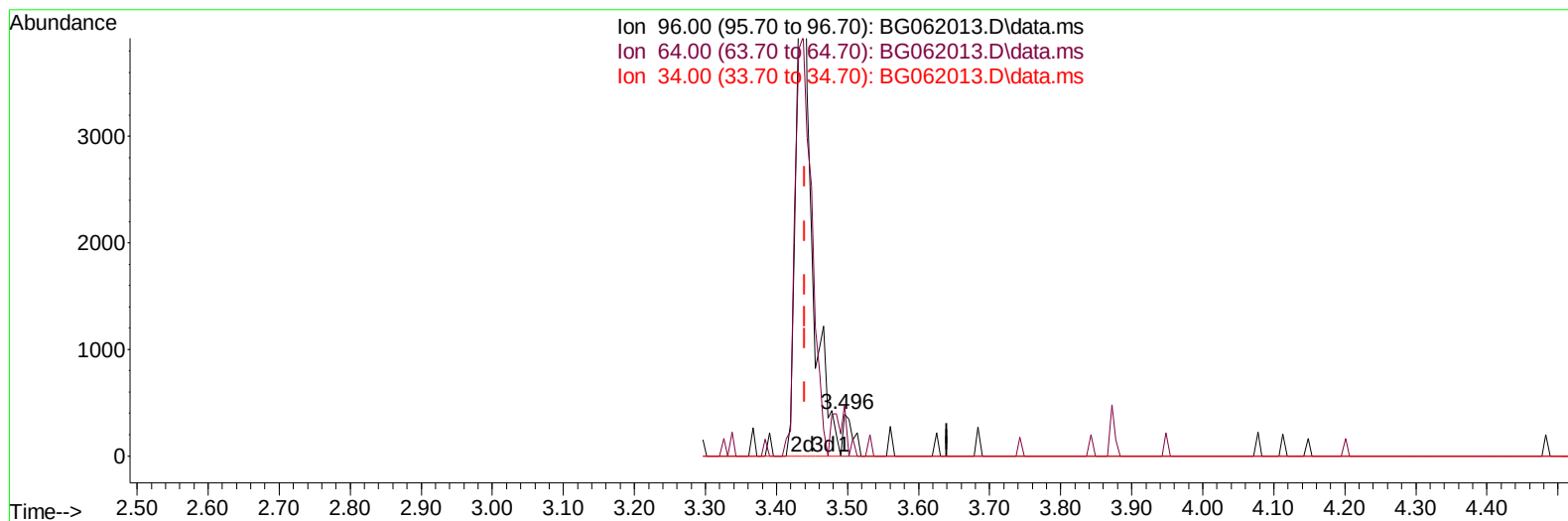
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062013.D
Acq On : 11 Jul 2024 17:02
Operator : MA/JU
Sample : PB161883BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK883

Manual IntegrationsAPPROVED

Quant Time: Jul 12 04:50:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 11 12:11:32 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/12/2024
Supervised By :mohammad ahmed 07/15/2024



TIC: BG062013.D\data.ms

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

3.496min (+ 0.056) 0.31 ng/uL

response 392

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	121.88
34.00	0.00	0.00
0.00	0.00	0.00

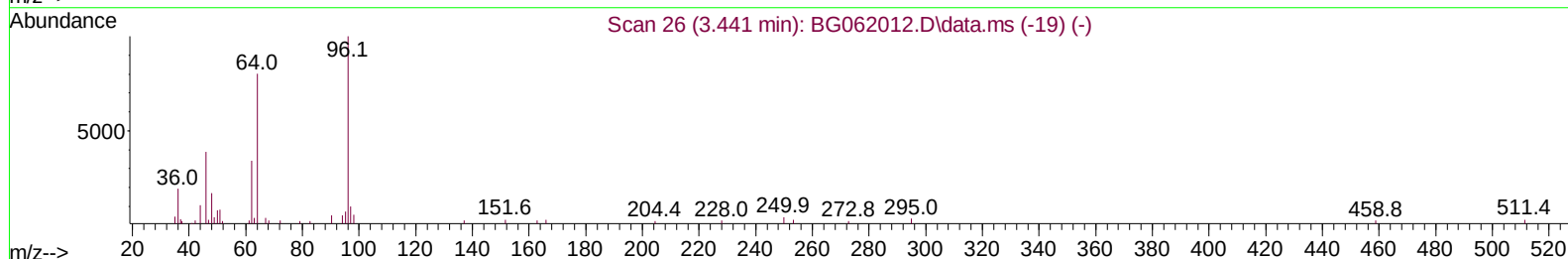
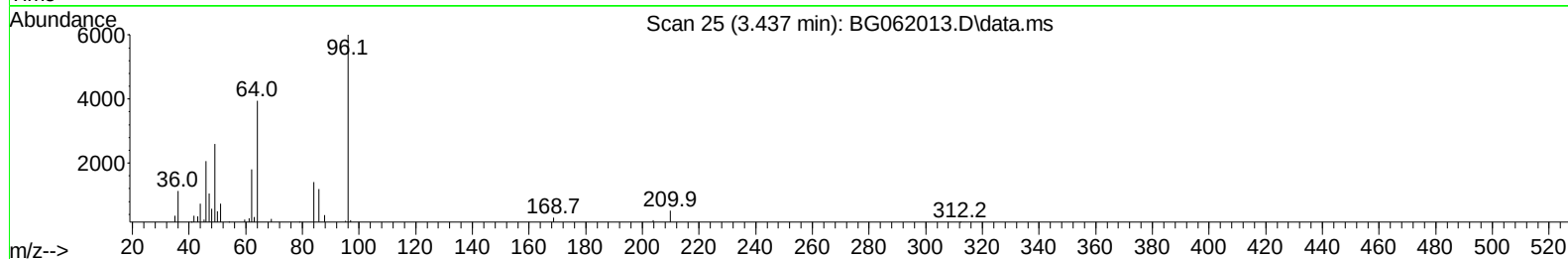
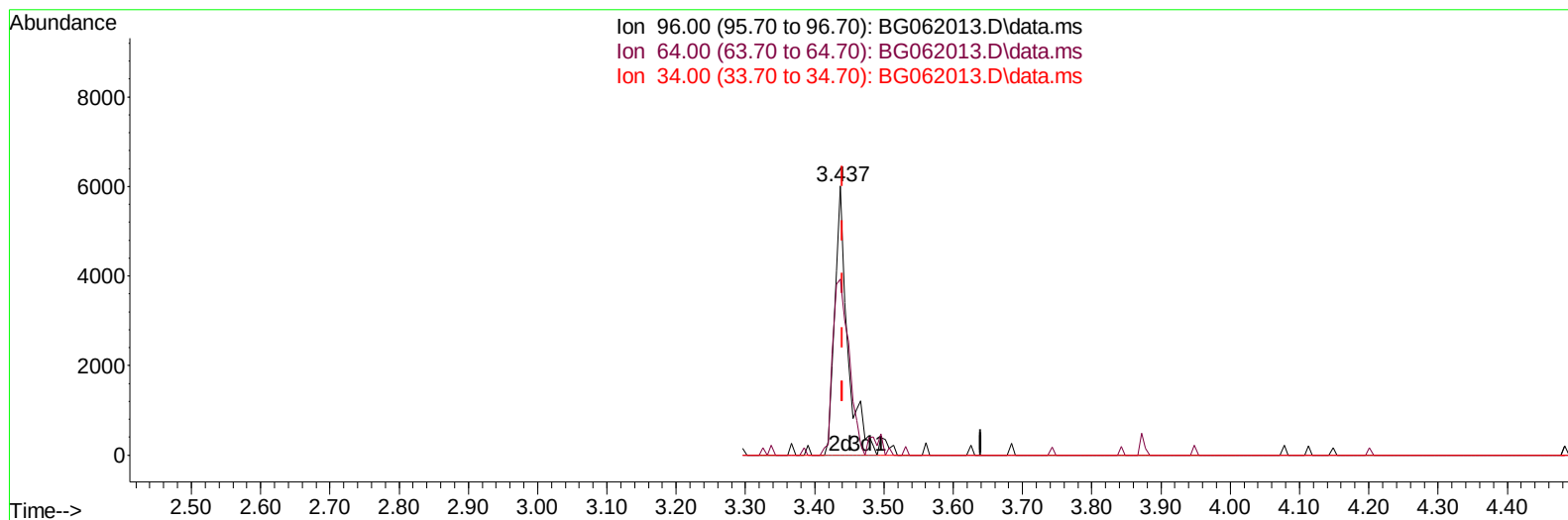
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(3) 1,4-Dioxane-d8 (S)

3.437min (-0.003) 6.11 ng/uL m

response 7746

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	65.55#
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Jul 12 04:52:18 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.038	152	47656	20.000	ng/ul	0.00
20) Naphthalene-d8	10.852	136	212329	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.671	164	140103	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.415	188	363338	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.669	240	381018	20.000	ng/ul	0.00
88) Perylene-d12	24.877	264	456523	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.437	96	7746m	6.108	ng/uL	0.00
4) Pyridine-d5	3.854	84	102414	26.264	ng/ul	0.00
7) Phenol-d5	7.192	99	136305	26.136	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.362	67	90574	27.424	ng/ul	0.00
11) 2-Chlorophenol-d4	7.568	132	103996	29.070	ng/ul	0.00
15) 4-Methylphenol-d8	8.743	113	110765	26.975	ng/ul	0.00
21) Nitrobenzene-d5	9.201	128	52528	32.528	ng/ul	0.00
24) 2-Nitrophenol-d4	9.924	143	62257	33.760	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.464	165	110285	31.005	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	138877	26.893	ng/ul	0.00
46) Dimethylphthalate-d6	14.072	166	363690	31.575	ng/ul	0.00
49) Acenaphthylene-d8	14.366	160	380970	31.631	ng/ul	0.00
54) 4-Nitrophenol-d4	14.871	143	56220	30.564	ng/ul	0.00
60) Fluorene-d10	15.658	176	312592	32.237	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.776	200	54538	28.581	ng/ul	0.00
73) Anthracene-d10	17.515	188	526177	31.018	ng/ul	0.00
81) Pyrene-d10	19.794	212	708268	31.603	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.659	264	739373	32.656	ng/ul	0.00

Target Compounds Qvalue

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

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