

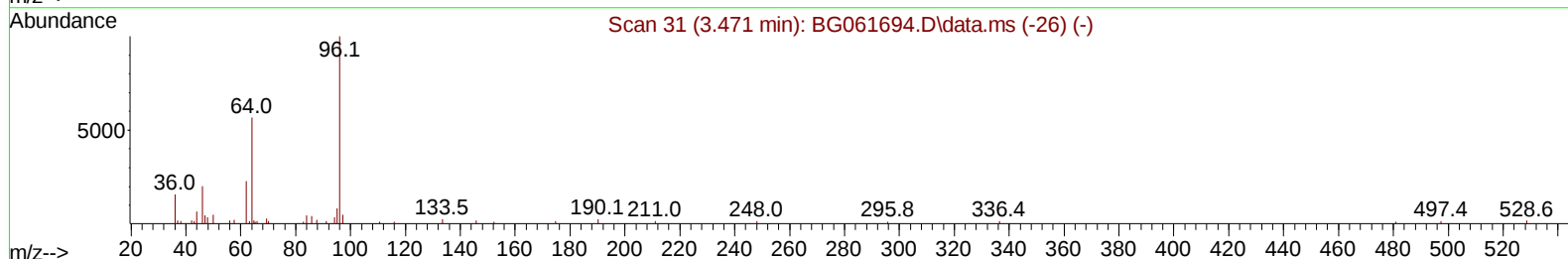
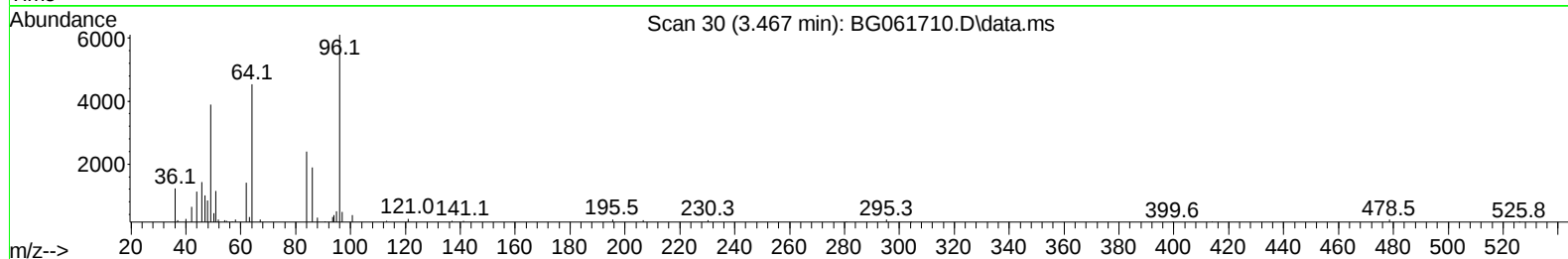
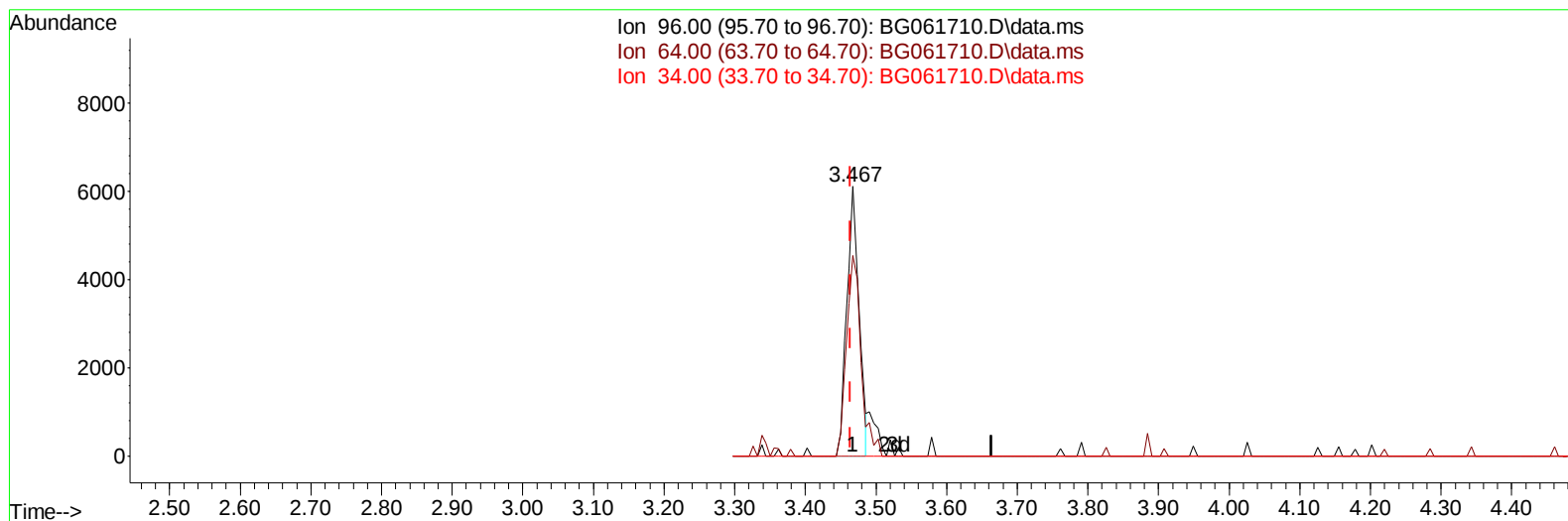
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061710.D
Acq On : 25 Jun 2024 20:00
Operator : MA/JU
Sample : P2873-06
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0SH9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/26/2024
Supervised By :mohammad ahmed 06/27/2024

Quant Time: Jun 25 23:07:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 21 05:03:55 2024
Response via : Initial Calibration



TIC: BG061710.D\data.ms

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

3.467min (+ 0.003) 4.53 ng/uL

response 7401

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	74.30
34.00	0.00	0.00
0.00	0.00	0.00

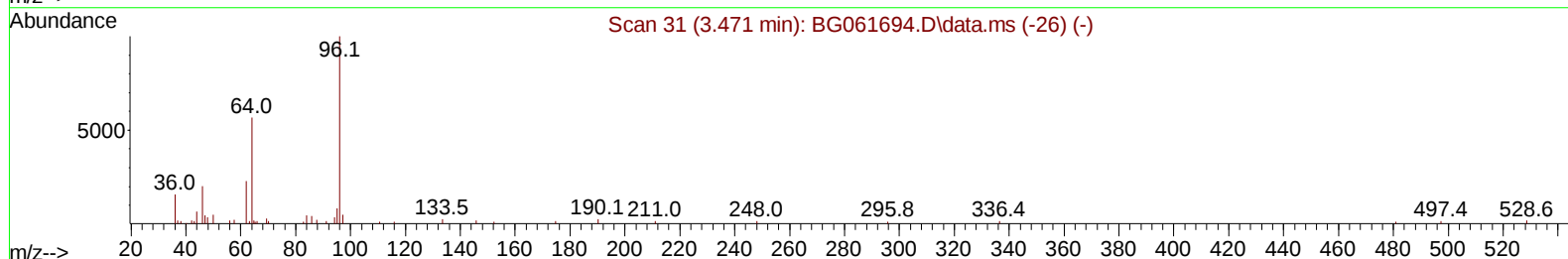
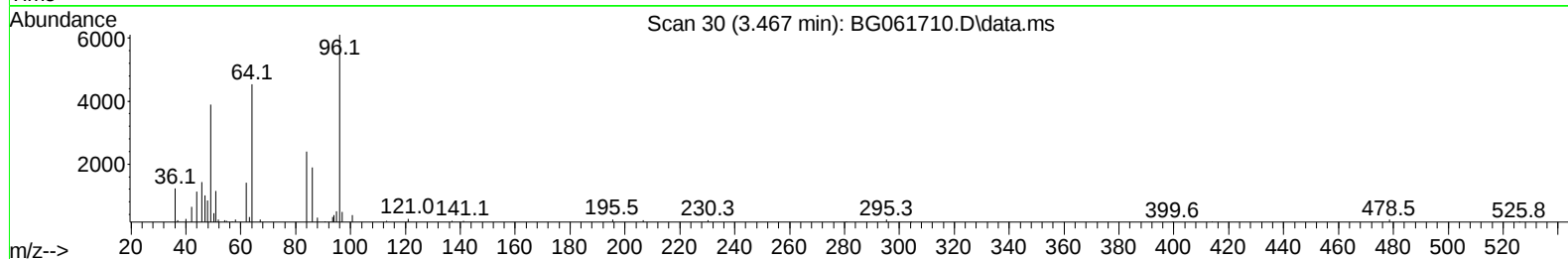
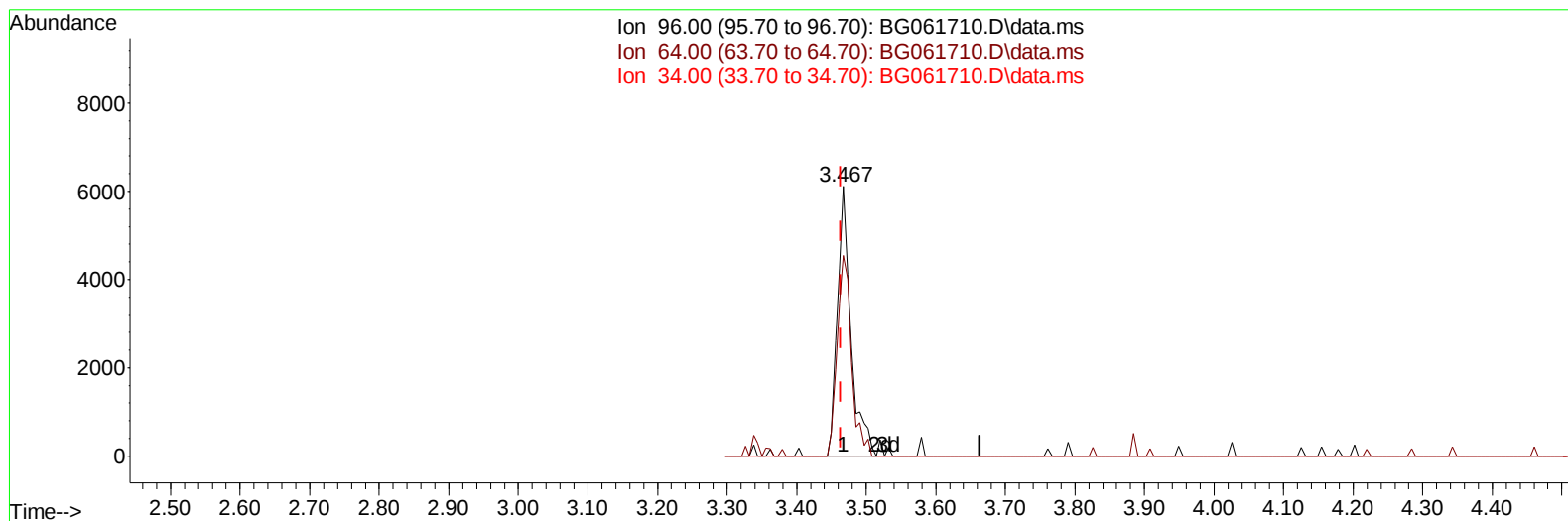
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TIC: BG061710.D\data.ms

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

3.467min (+ 0.003) 5.09 ng/uL m

response 8315

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	74.30
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.074	152	59006	20.000	ng/ul	0.00
20) Naphthalene-d8	10.882	136	311258	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.701	164	241707	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	588127	20.000	ng/ul	0.00
79) Chrysene-d12	21.705	240	501406	20.000	ng/ul	0.00
88) Perylene-d12	24.936	264	592269	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.467	96	8315m	5.091	ng/uL	0.00
4) Pyridine-d5	3.885	84	113721	23.027	ng/ul	0.00
7) Phenol-d5	7.216	99	164866	25.737	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	98948	24.779	ng/ul	0.00
11) 2-Chlorophenol-d4	7.598	132	111111	25.277	ng/ul	0.00
15) 4-Methylphenol-d8	8.767	113	133508	26.960	ng/ul	0.00
21) Nitrobenzene-d5	9.237	128	59281	28.050	ng/ul	0.00
24) 2-Nitrophenol-d4	9.960	143	71430	33.371	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.494	165	139554	27.534	ng/ul	0.00
31) 4-Chloroaniline-d4	11.017	131	197757	25.771	ng/ul	0.00
46) Dimethylphthalate-d6	14.102	166	506053	27.207	ng/ul	0.00
49) Acenaphthylene-d8	14.396	160	575544	28.235	ng/ul	0.00
54) 4-Nitrophenol-d4	14.878	143	84671	26.506	ng/ul	0.00
60) Fluorene-d10	15.688	176	474770	29.156	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.794	200	63636	25.304	ng/ul	0.00
73) Anthracene-d10	17.539	188	768743	28.636	ng/ul	0.00
81) Pyrene-d10	19.819	212	829002	29.599	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.713	264	830552	29.361	ng/ul	0.00

Target Compounds Qvalue

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

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