

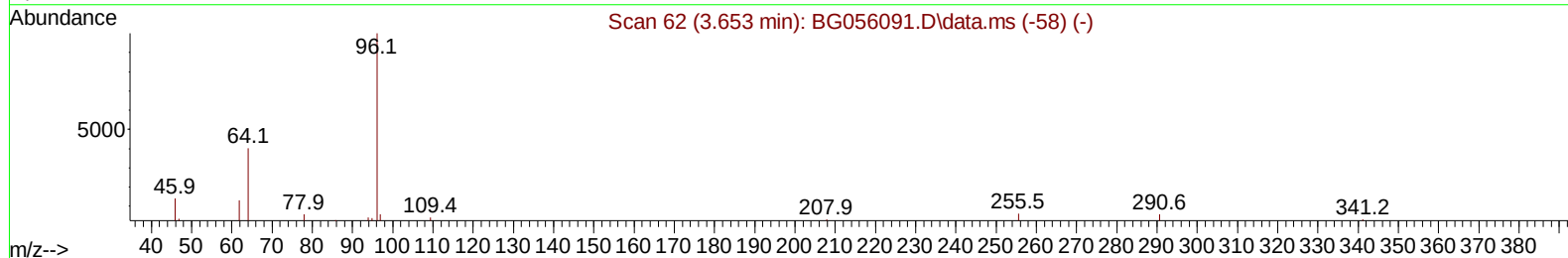
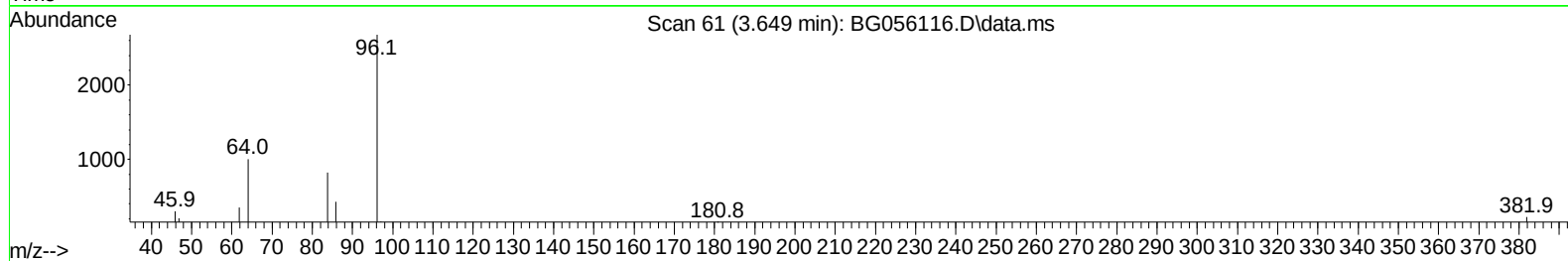
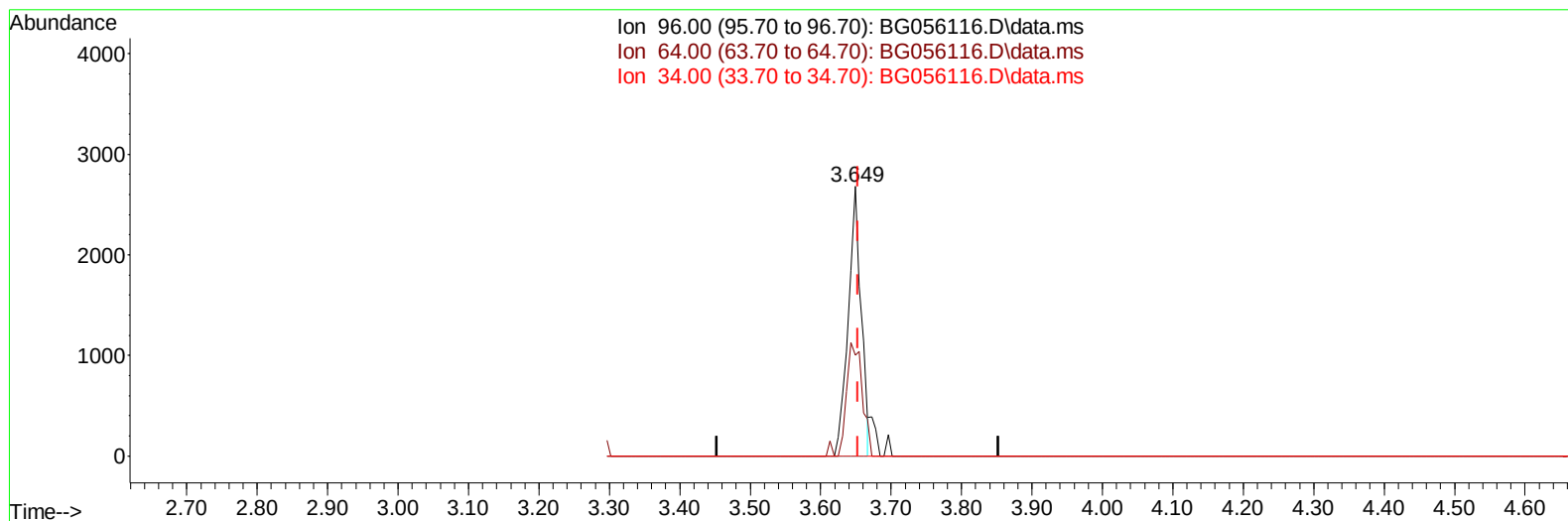
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056116.D
 Acq On : 25 Dec 2022 10:02
 Operator : CG/JU
 Sample : N6017-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYG0

Manual IntegrationsAPPROVED

Quant Time: Dec 25 22:37:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 22:54:38 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/26/2022
 Supervised By :Sohil Jodhani 12/26/2022



TIC: BG056116.D\data.ms

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

3.649min (-0.004) 3.49 ng/uL

response 3378

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.20	37.46
34.00	0.00	0.00
0.00	0.00	0.00

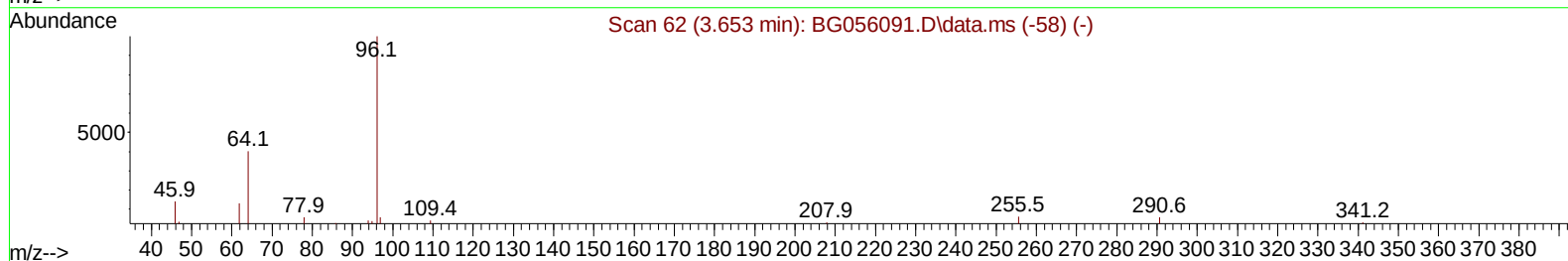
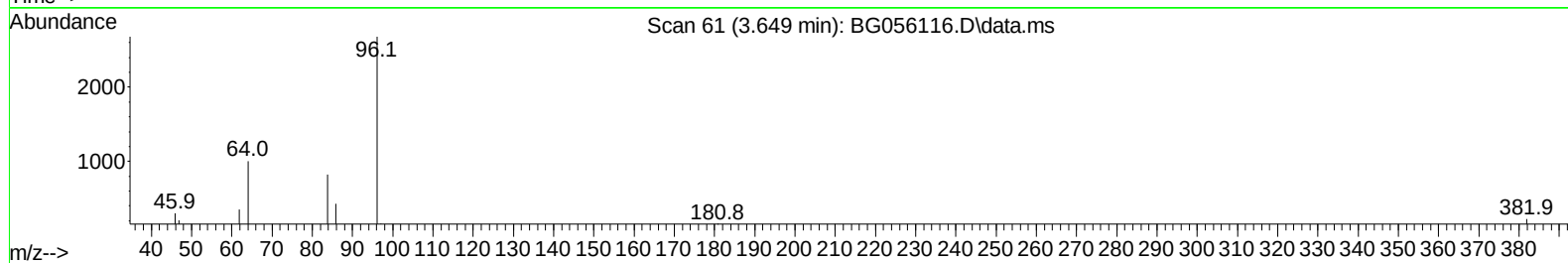
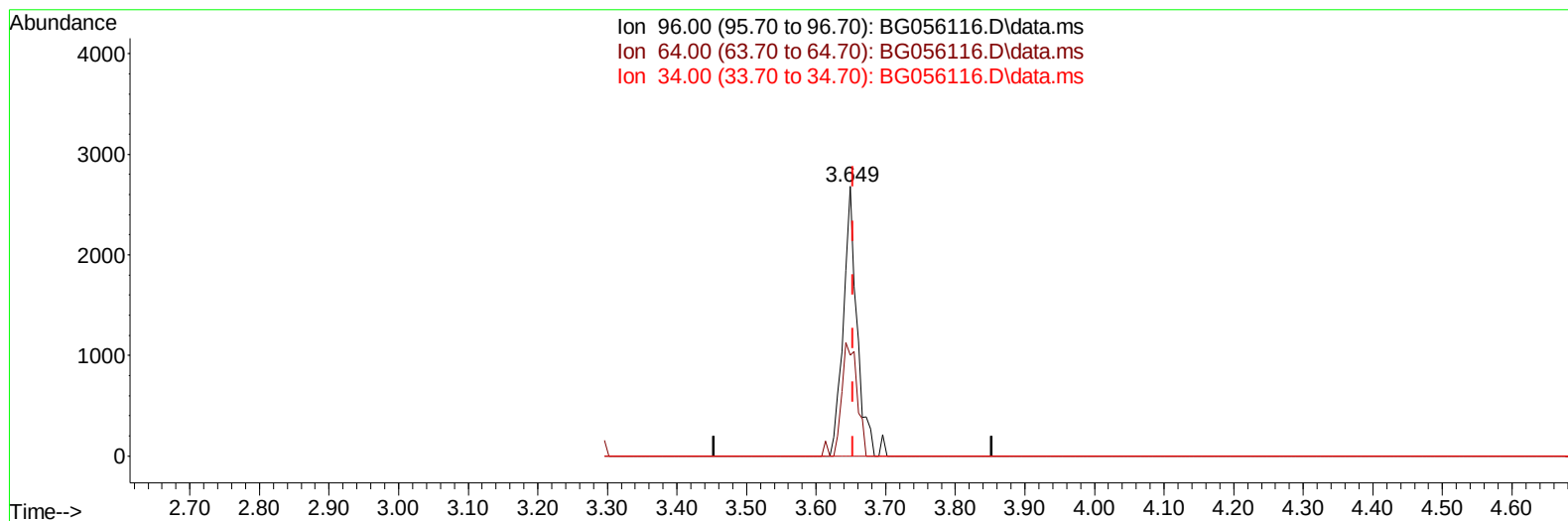
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TIC: BG056116.D\data.ms

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

3.649min (-0.004) 3.73 ng/uL m

response 3608

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.20	37.46
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	8.349	152	40523	20.000	ng/ul	0.00
20) Naphthalene-d8	11.175	136	167384	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.953	164	147727	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.685	188	367398	20.000	ng/ul	0.00
79) Chrysene-d12	21.986	240	358130	20.000	ng/ul	0.00
88) Perylene-d12	25.458	264	377711	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.649	96	3608m	3.727	ng/uL	0.00
4) Pyridine-d5	4.083	84	37567	11.902	ng/ul	0.00
7) Phenol-d5	7.474	99	73312	19.944	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.650	67	32609	22.508	ng/ul	0.00
11) 2-Chlorophenol-d4	7.873	132	53950	24.833	ng/ul	0.00
15) 4-Methylphenol-d8	9.031	113	59127	21.373	ng/ul	0.00
21) Nitrobenzene-d5	9.518	128	28006	22.972	ng/ul	0.00
24) 2-Nitrophenol-d4	10.247	143	39277	24.659	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.782	165	80051	22.616	ng/ul	0.00
31) 4-Chloroaniline-d4	11.299	131	72634	18.143	ng/ul	0.00
46) Dimethylphthalate-d6	14.342	166	270543	23.662	ng/ul	0.00
49) Acenaphthylene-d8	14.653	160	308744	23.911	ng/ul	0.00
54) 4-Nitrophenol-d4	15.112	143	27618	16.047	ng/ul	0.00
60) Fluorene-d10	15.934	176	259647	23.007	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.040	200	50732	21.480	ng/ul	0.00
73) Anthracene-d10	17.785	188	402209	23.690	ng/ul	0.00
81) Pyrene-d10	20.053	212	479811	24.259	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.218	264	468590	24.408	ng/ul	0.00

Target Compounds Qvalue

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

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