

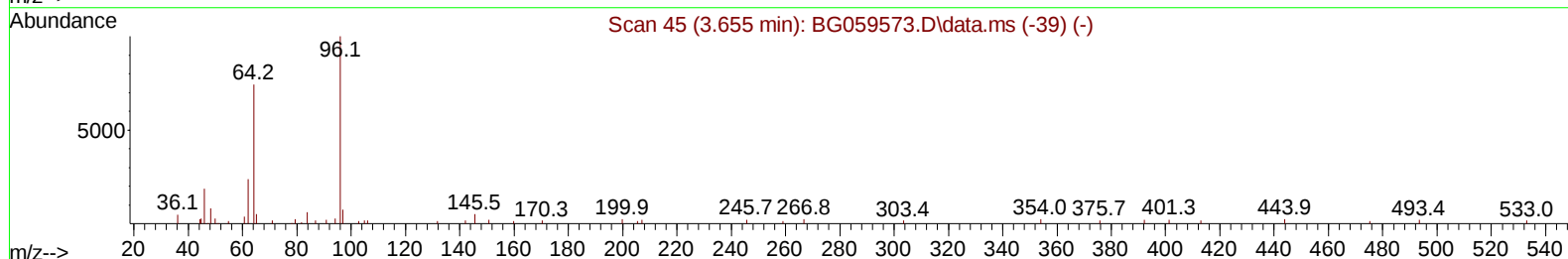
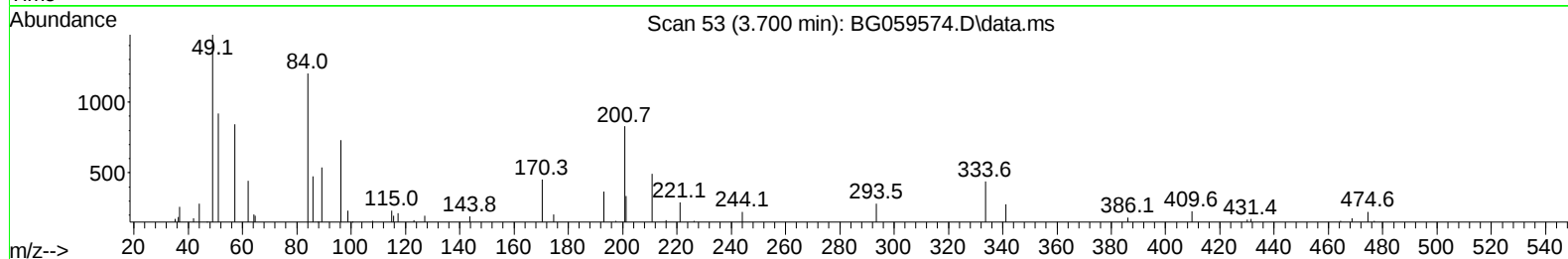
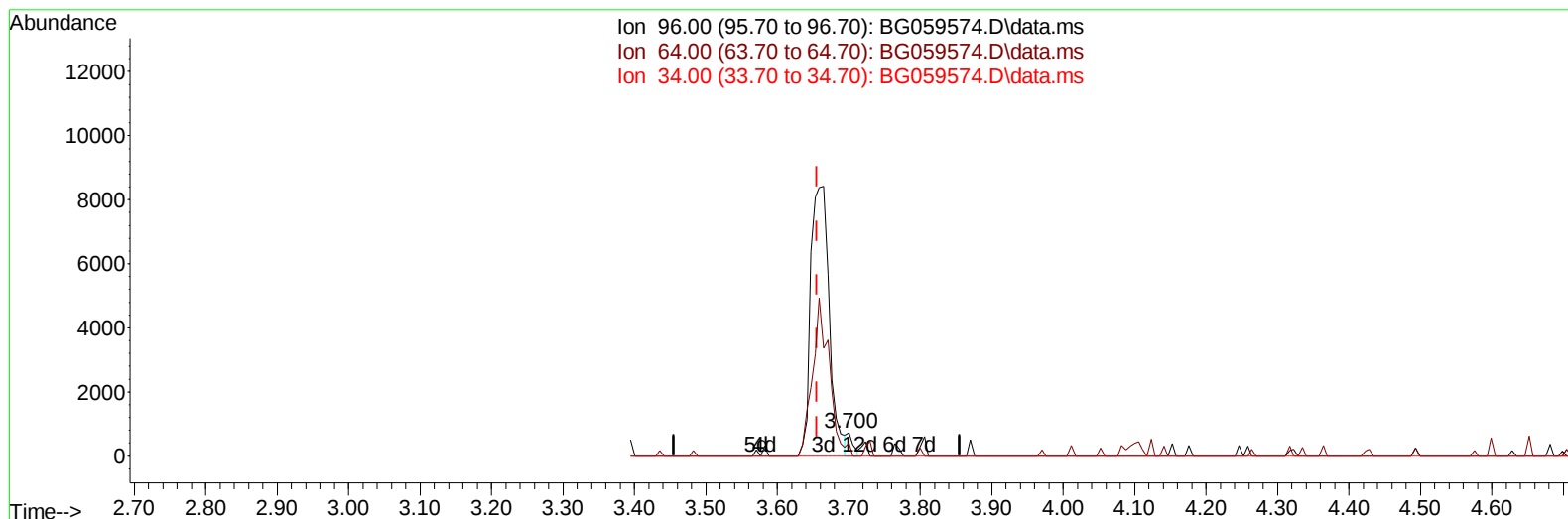
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110123\
 Data File : BG059574.D
 Acq On : 1 Nov 2023 11:16
 Operator : MA/JU
 Sample : PB156564BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK564

Manual IntegrationsAPPROVED

Quant Time: Nov 02 02:30:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 28 03:56:42 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/03/2023
 Supervised By :mohammad ahmed 11/03/2023



TIC: BG059574.D\data.ms

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

3.700min (+ 0.045) 0.18 ng/uL

response 442

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	52.90	55.40
34.00	0.00	0.00
0.00	0.00	0.00

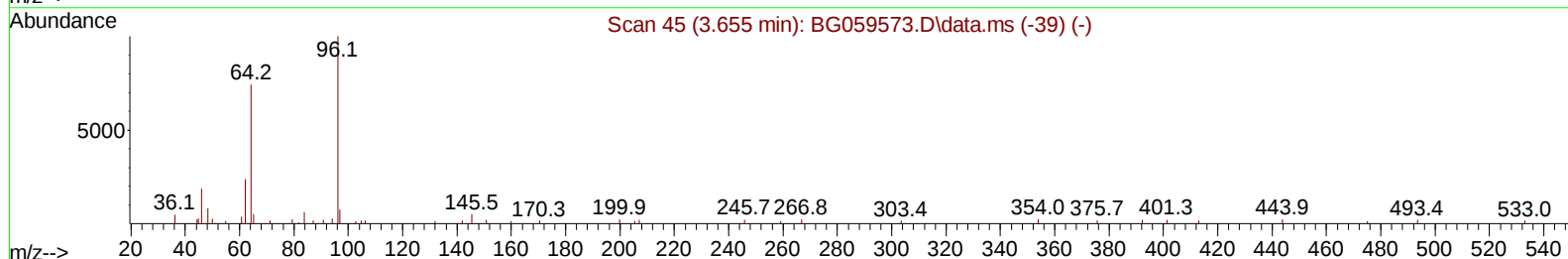
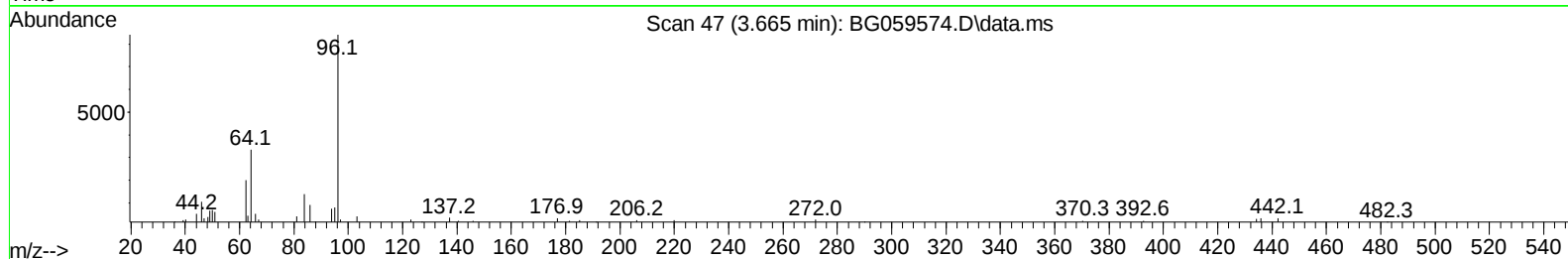
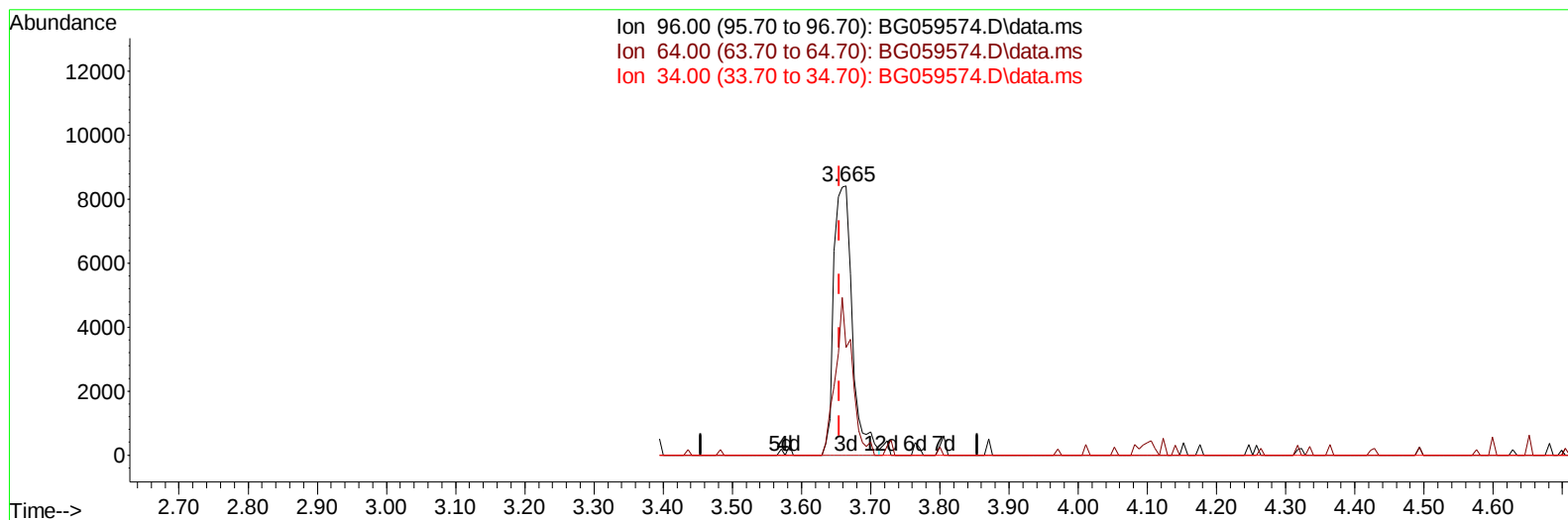
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TIC: BG059574.D\data.ms

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

3.665min (+ 0.010) 6.23 ng/uL m

response 15702

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	52.90	39.87#
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	8.359	152	94241	20.000	ng/ul	0.00
20) Naphthalene-d8	11.209	136	421683	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.987	164	273492	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.736	188	547714	20.000	ng/ul	0.00
79) Chrysene-d12	22.073	240	446351	20.000	ng/ul	0.00
88) Perylene-d12	25.674	264	489152	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.665	96	15702m	6.234	ng/uL	0.00
4) Pyridine-d5	4.094	84	210513	29.908	ng/ul	0.00
7) Phenol-d5	7.466	99	274929	28.389	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.672	67	121670	27.353	ng/ul	0.00
11) 2-Chlorophenol-d4	7.872	132	225817	28.742	ng/ul	0.00
15) 4-Methylphenol-d8	9.041	113	234763	30.555	ng/ul	0.00
21) Nitrobenzene-d5	9.558	128	122106	40.725	ng/ul	0.00
24) 2-Nitrophenol-d4	10.275	143	129920	42.209	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.797	165	211459	31.181	ng/ul	0.00
31) 4-Chloroaniline-d4	11.350	131	347987	31.171	ng/ul	0.00
46) Dimethylphthalate-d6	14.382	166	736216	29.713	ng/ul	0.00
49) Acenaphthylene-d8	14.687	160	887639	30.526	ng/ul	0.00
54) 4-Nitrophenol-d4	15.151	143	135176	31.813	ng/ul	0.00
60) Fluorene-d10	15.974	176	621955	29.372	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.074	200	90125	34.746	ng/ul	0.00
73) Anthracene-d10	17.836	188	960972	32.385	ng/ul	0.00
81) Pyrene-d10	20.110	212	985989	31.780	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.427	264	932098	32.360	ng/ul	-0.01

Target Compounds Qvalue

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

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