

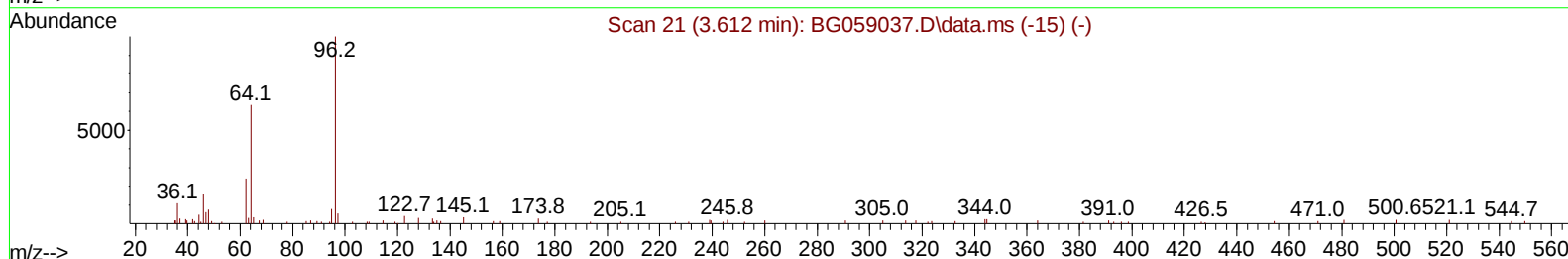
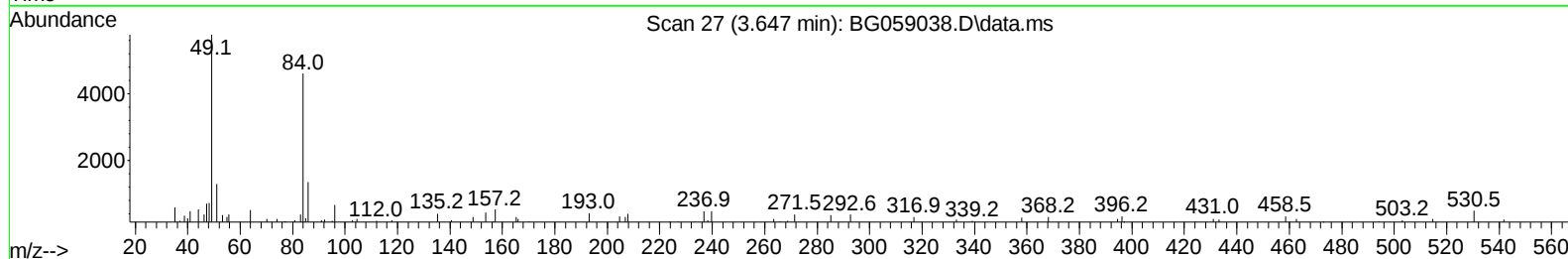
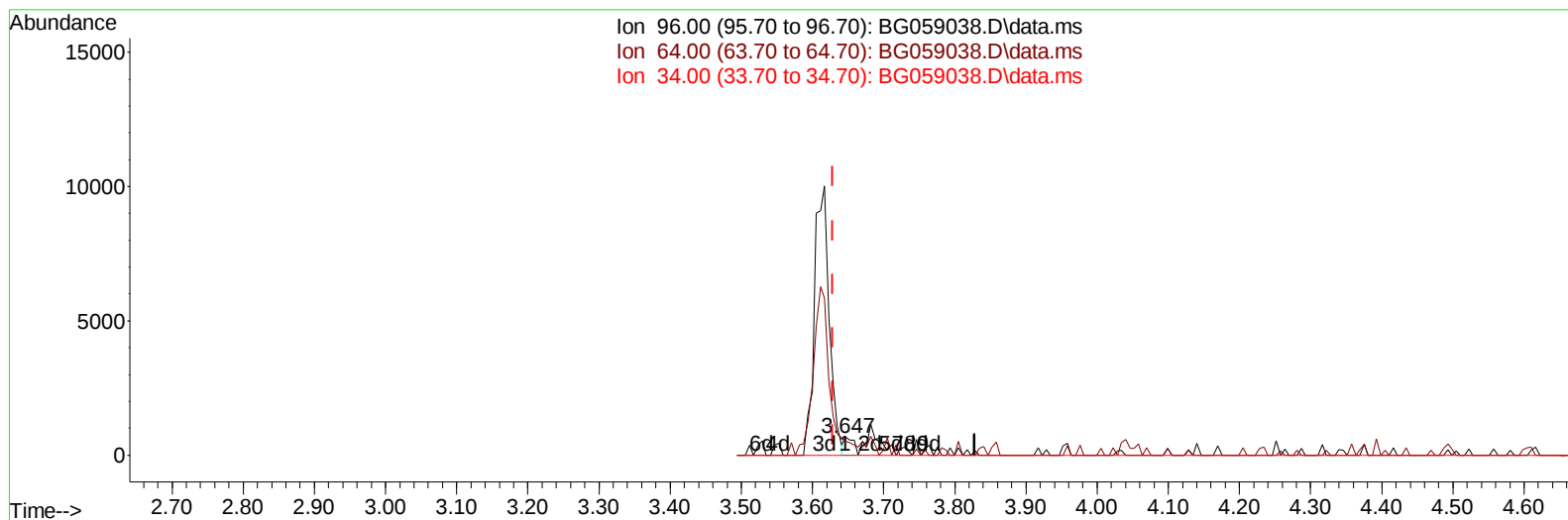
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091323\
 Data File : BG059038.D
 Acq On : 13 Sep 2023 13:13
 Operator : MA/JU
 Sample : PB155387BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK387

Manual IntegrationsAPPROVED

Quant Time: Sep 14 01:04:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:54:19 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/14/2023
 Supervised By :mohammad ahmed 09/14/2023



TIC: BG059038.D\data.ms

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

3.647min (+ 0.018) 0.25 ng/uL

response 627

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.70	78.20
34.00	0.00	0.00
0.00	0.00	0.00

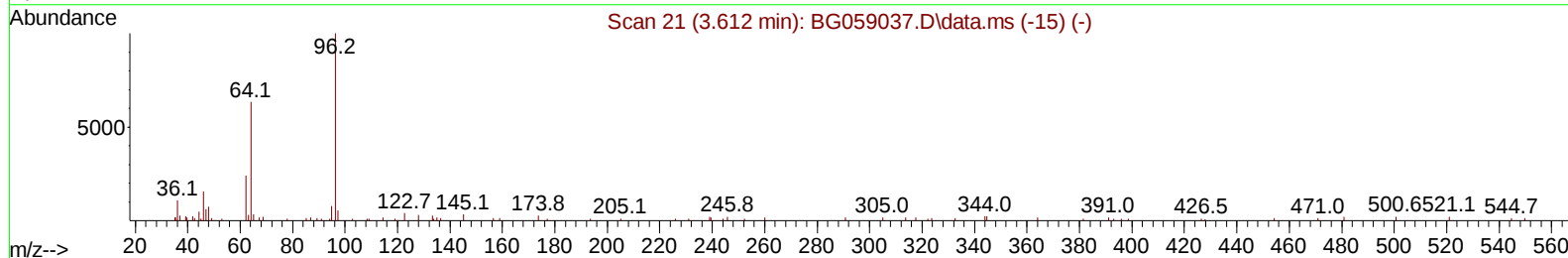
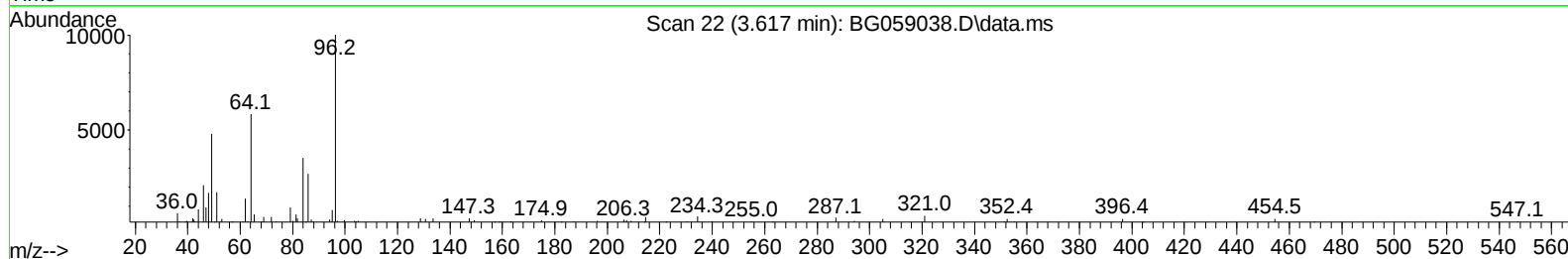
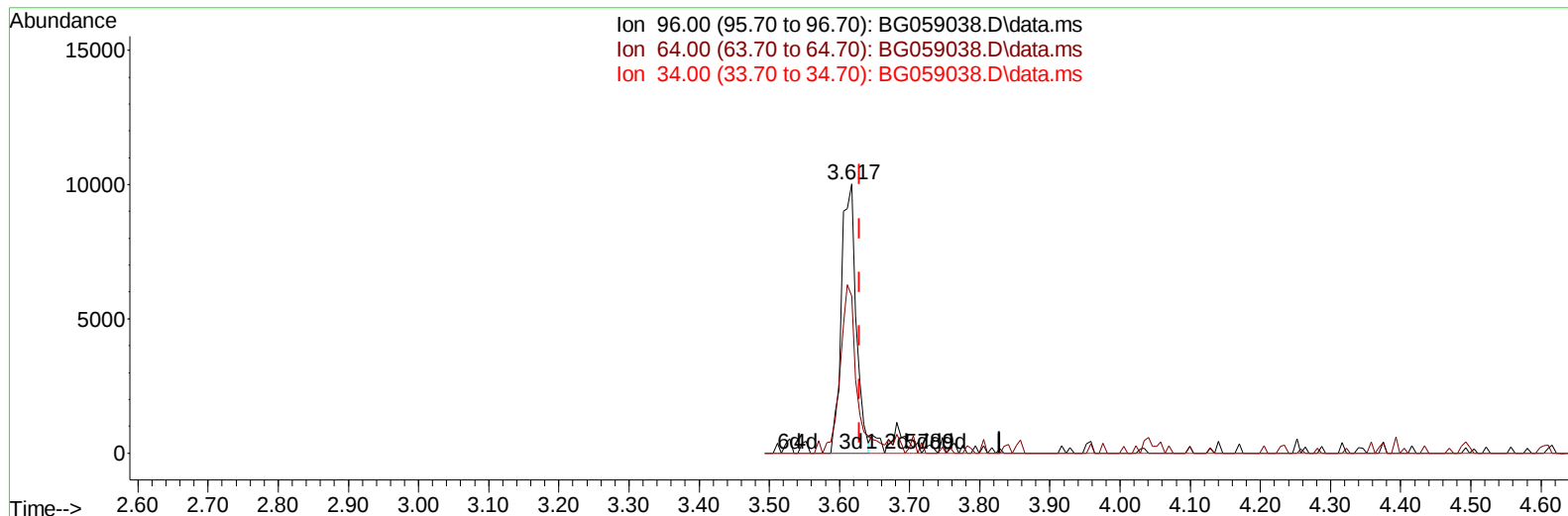
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(3) 1,4-Dioxane-d8 (S)

3.617min (-0.011) 5.83 ng/uL m

response 14522

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.70	58.29#
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Sep 14 01:17:24 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.306	152	105845	20.000	ng/ul	0.00
20) Naphthalene-d8	11.155	136	477604	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.939	164	360402	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.689	188	979912	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.013	240	833003	20.000	ng/ul	0.00
88) Perylene-d12	25.580	264	946252	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.617	96	14522m	5.832	ng/uL	-0.01
4) Pyridine-d5	4.046	84	173314	21.983	ng/ul	-0.01
7) Phenol-d5	7.419	99	284156	26.623	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.624	67	177648	28.384	ng/ul	0.00
11) 2-Chlorophenol-d4	7.818	132	194910	28.473	ng/ul	0.00
15) 4-Methylphenol-d8	8.987	113	228713	27.736	ng/ul	0.00
21) Nitrobenzene-d5	9.504	128	106046	30.611	ng/ul	0.00
24) 2-Nitrophenol-d4	10.227	143	117689	30.960	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.744	165	230131	28.485	ng/ul	0.00
31) 4-Chloroaniline-d4	11.291	131	144716	13.320	ng/ul	0.00
46) Dimethylphthalate-d6	14.334	166	947652	33.898	ng/ul	0.00
49) Acenaphthylene-d8	14.640	160	932525	29.955	ng/ul	0.00
54) 4-Nitrophenol-d4	15.110	143	143944	30.998	ng/ul	0.00
60) Fluorene-d10	15.926	176	782262	31.531	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.032	200	149177	31.014	ng/ul	0.00
73) Anthracene-d10	17.789	188	1371048	31.518	ng/ul	0.00
81) Pyrene-d10	20.063	212	1505795	34.125	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.333	264	1430673	30.843	ng/ul	0.01

Target Compounds Qvalue

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

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