

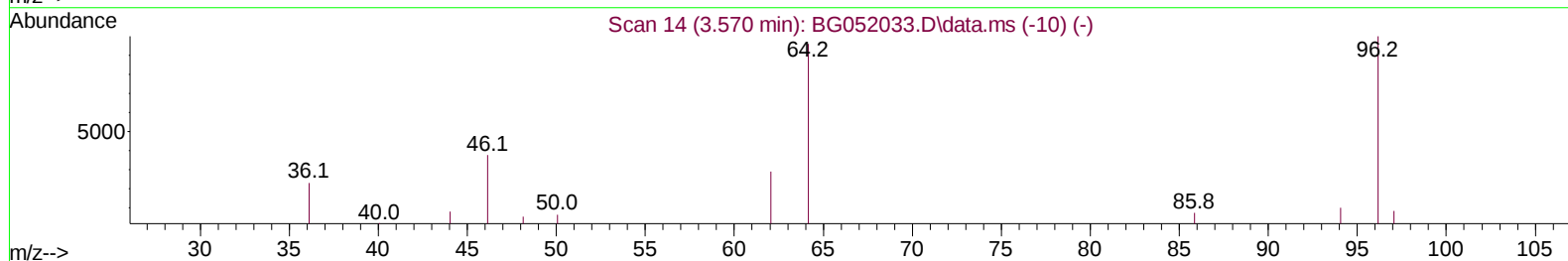
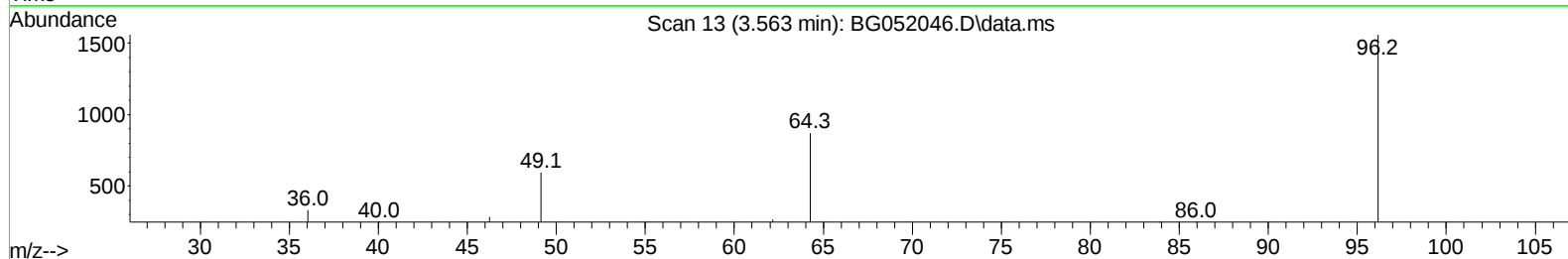
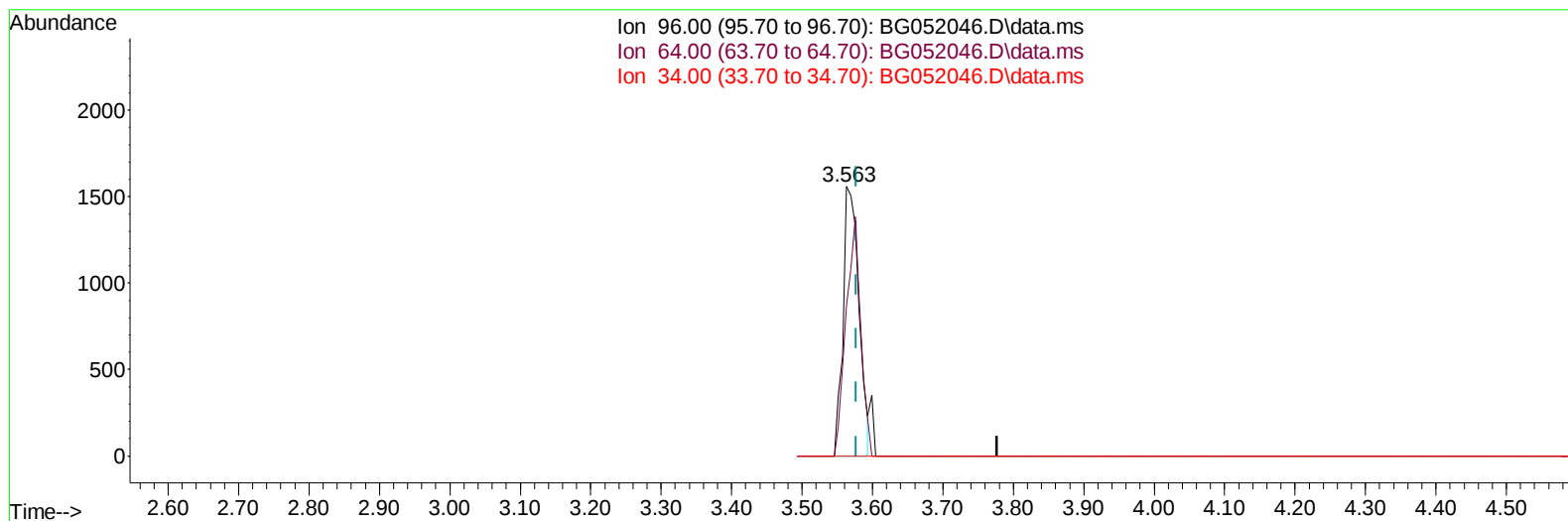
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052046.D
 Acq On : 18 Jan 2022 10:28
 Operator : CG/JU
 Sample : N1169-18
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLY5

Manual IntegrationsAPPROVED

Quant Time: Jan 18 11:12:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/18/2022
 Supervised By :mohammad ahmed 01/21/2022



TIC: BG052046.D\data.ms

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

3.563min (-0.014) 3.65 ng/uL

response 2399

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	55.87#
34.00	0.00	0.00
0.00	0.00	0.00

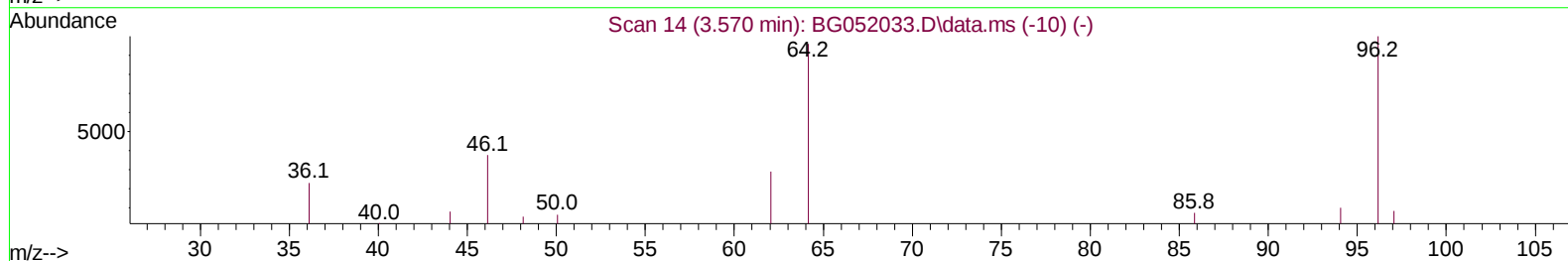
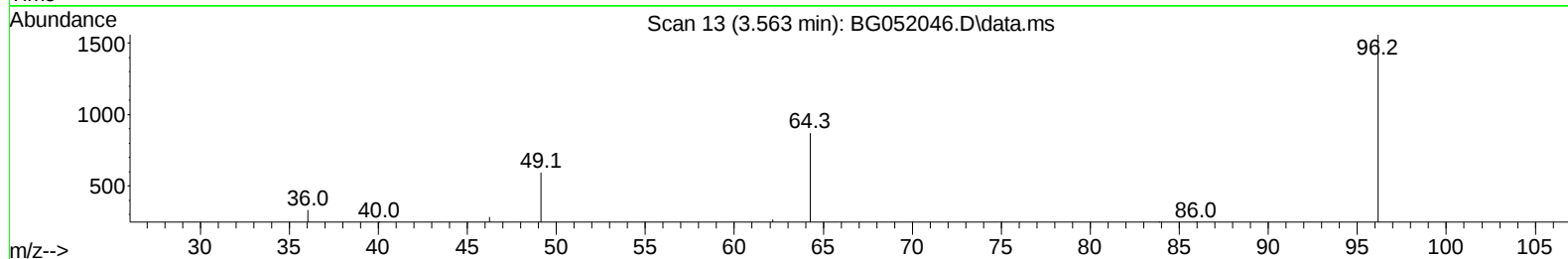
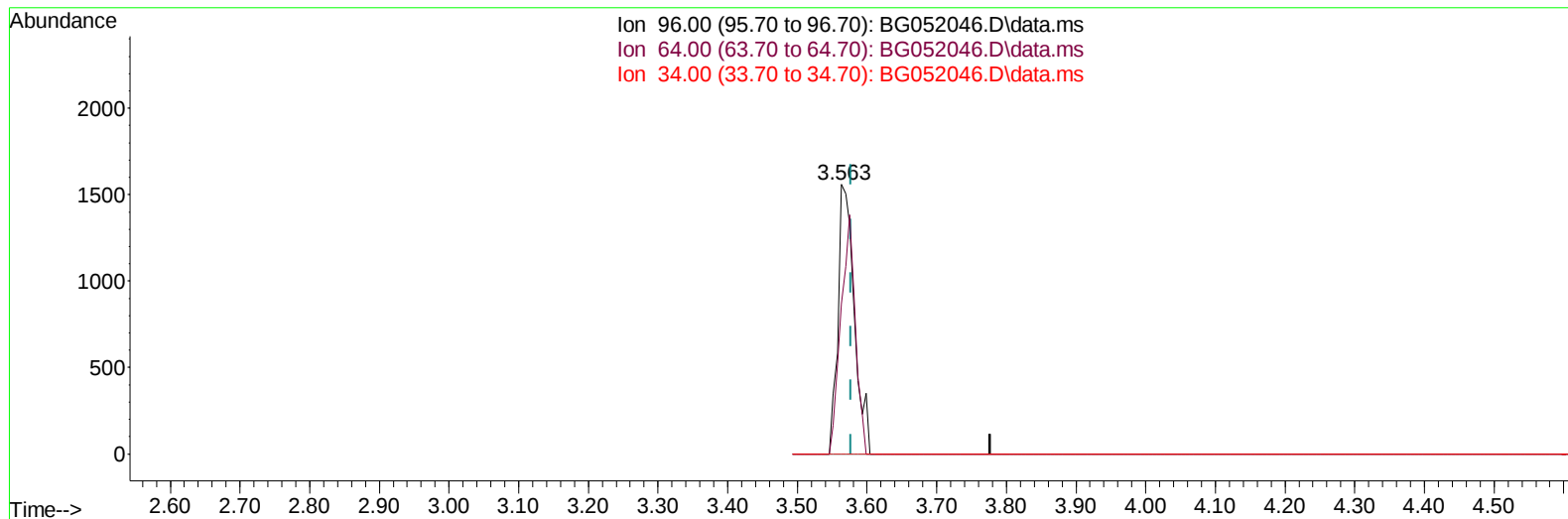
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TIC: BG052046.D\data.ms

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

3.563min (-0.014) 3.84 ng/uL m

response 2523

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	55.87#
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	8.245	152	23458	20.000	ng/ul	0.00
20) Naphthalene-d8	11.083	136	102235	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.889	164	77619	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.639	188	192580	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.956	240	191082	20.000	ng/ul	# 0.00
88) Perylene-d12	25.446	264	190678	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.563	96	2523m	3.843	ng/uL	-0.01
4) Pyridine-d5	3.998	84	15184	7.440	ng/ul	0.00
7) Phenol-d5	7.382	99	51980	22.640	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.546	67	30278	20.775	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.769	132	34378	22.616	ng/ul	0.00
15) 4-Methylphenol-d8	8.944	113	40341	22.603	ng/ul	0.00
21) Nitrobenzene-d5	9.414	128	18727	23.182	ng/ul	0.00
24) 2-Nitrophenol-d4	10.143	143	21351	23.897	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.695	165	40394	23.130	ng/ul	0.00
31) 4-Chloroaniline-d4	11.206	131	52391	21.542	ng/ul	0.00
46) Dimethylphthalate-d6	14.278	166	154303	24.050	ng/ul	0.00
49) Acenaphthylene-d8	14.584	160	191349	24.690	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143	20607	19.211	ng/ul	0.00
60) Fluorene-d10	15.876	176	138773	24.808	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.988	200	21403	17.617	ng/ul	0.00
73) Anthracene-d10	17.738	188	238608	26.119	ng/ul	0.00
81) Pyrene-d10	20.018	212	279700	25.434	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.205	264	255738	25.477	ng/ul	0.01

Target Compounds Qvalue

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

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