

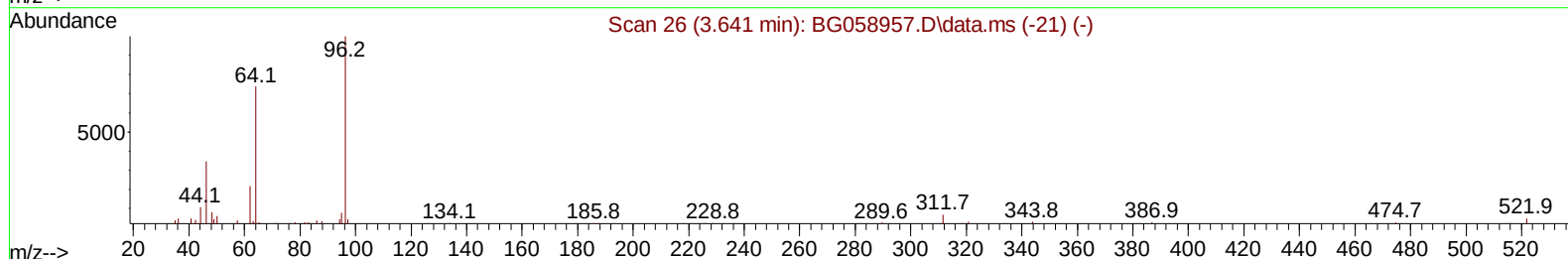
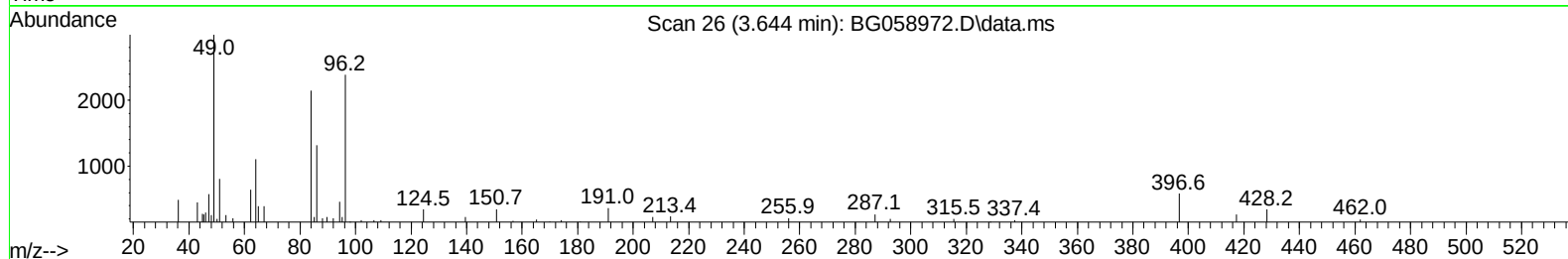
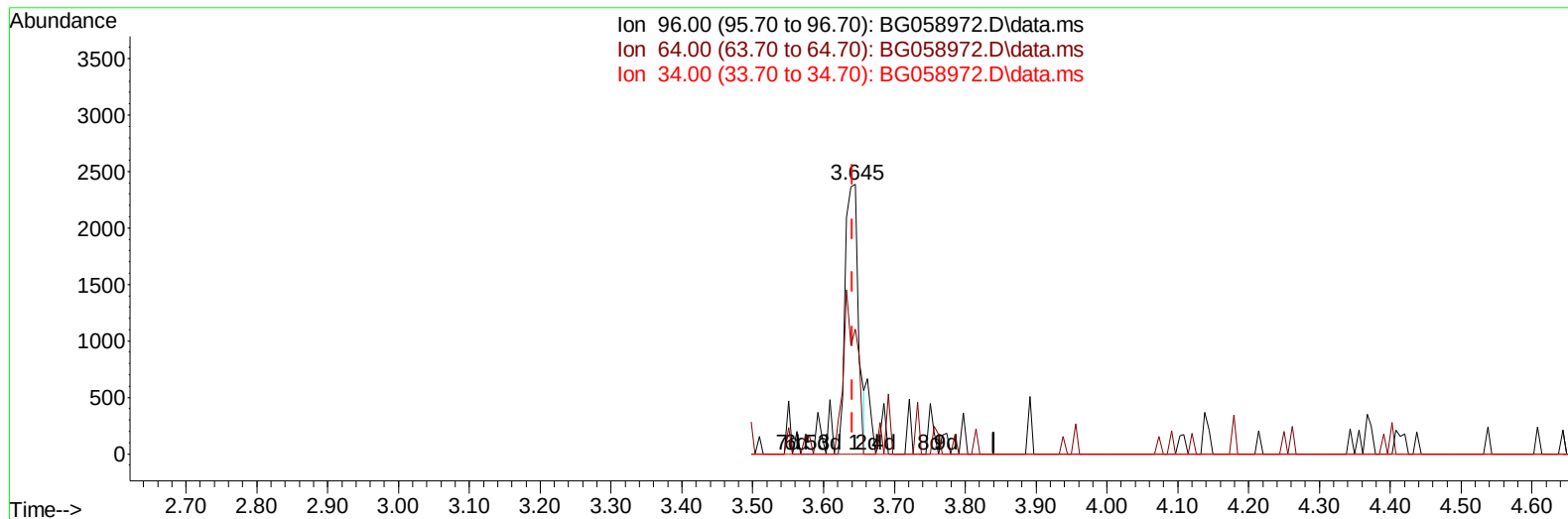
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG058972.D
Acq On : 11 Sep 2023 6:33
Operator : MA/JU
Sample : 04195-05
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
YBY19

Manual IntegrationsAPPROVED

Quant Time: Sep 11 07:20:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Sep 11 00:07:05 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/11/2023
Supervised By :mohammad ahmed 09/13/2023



TIC: BG058972.D\data.ms

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

3.644min (+ 0.003) 1.99 ng/uL

response 3075

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.70	46.29#
34.00	0.00	0.00
0.00	0.00	0.00

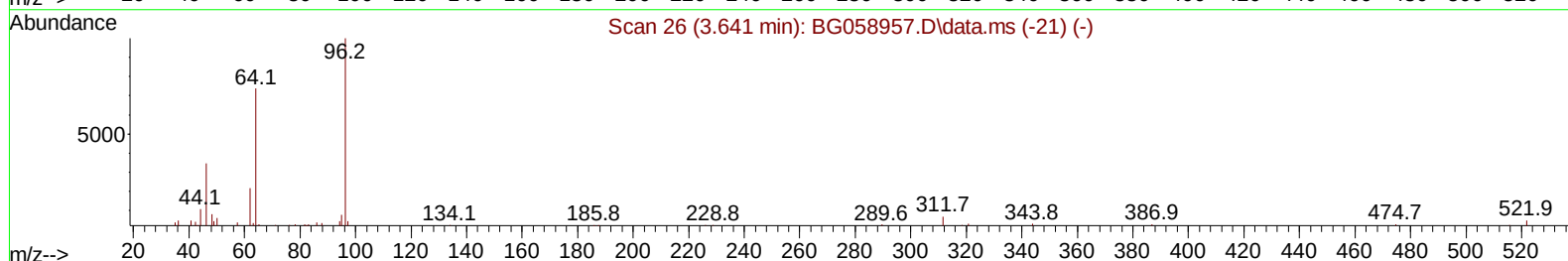
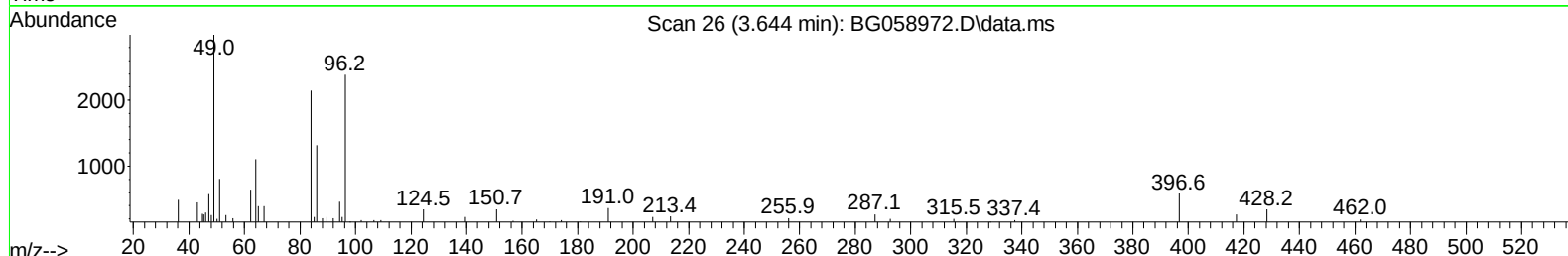
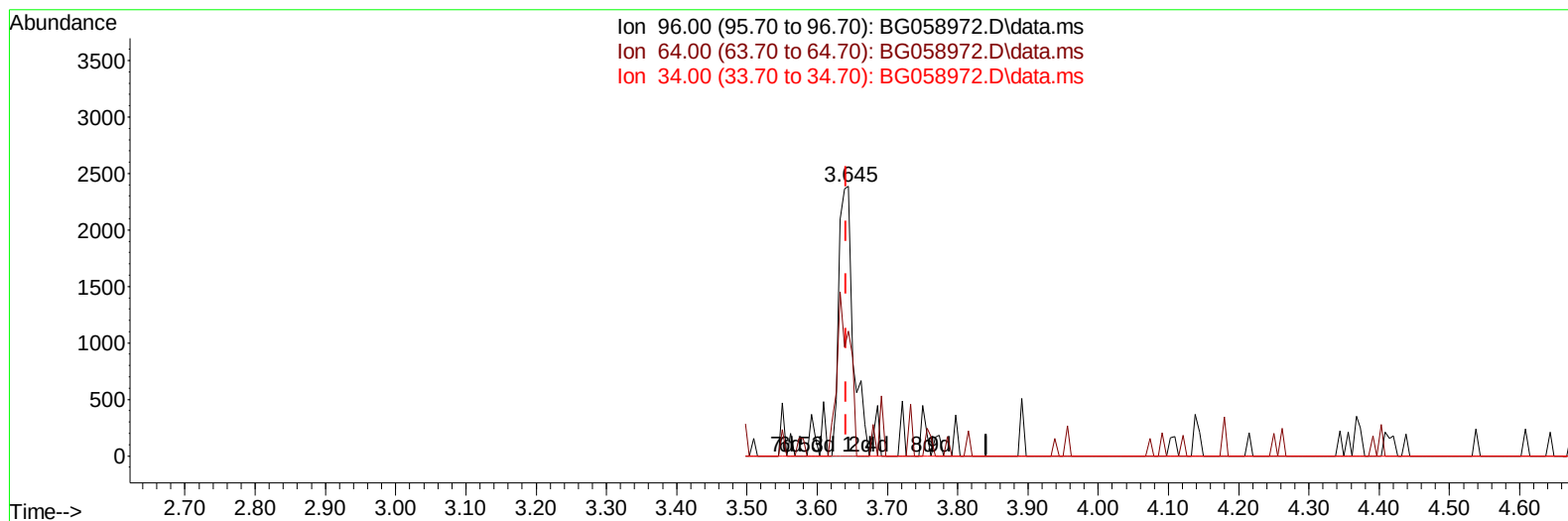
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TIC: BG058972.D\data.ms

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

3.644min (+ 0.003) 2.20 ng/uL m

response 3409

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.70	46.29#
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	8.327	152	65830	20.000	ng/ul	0.00
20) Naphthalene-d8	11.165	136	309885	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.949	164	240475	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.699	188	637787	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.017	240	640216	20.000	ng/ul	0.00
88) Perylene-d12	25.584	264	722476	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.644	96	3409m	2.201	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.434	99	75802	11.419	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.640	67	51470	13.223	ng/ul	0.00
11) 2-Chlorophenol-d4	7.840	132	50412	11.841	ng/ul	0.00
15) 4-Methylphenol-d8	9.003	113	54704	10.666	ng/ul	0.00
21) Nitrobenzene-d5	9.520	128	30499	13.569	ng/ul	0.00
24) 2-Nitrophenol-d4	10.237	143	24335	9.866	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.760	165	61817	11.793	ng/ul	0.00
31) 4-Chloroaniline-d4	11.300	131	21137	2.999	ng/ul	0.00
46) Dimethylphthalate-d6	14.344	166	168243	9.019	ng/ul	0.00
49) Acenaphthylene-d8	14.649	160	270922	13.043	ng/ul	0.00
54) 4-Nitrophenol-d4	15.108	143	20557	6.635	ng/ul	0.00
60) Fluorene-d10	15.936	176	244843	14.791	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.030	200	20119	6.426	ng/ul	0.00
73) Anthracene-d10	17.799	188	417780	14.756	ng/ul	0.00
81) Pyrene-d10	20.066	212	507396	14.961	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.331	264	477198	13.474	ng/ul	0.00

Target Compounds Qvalue

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

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