

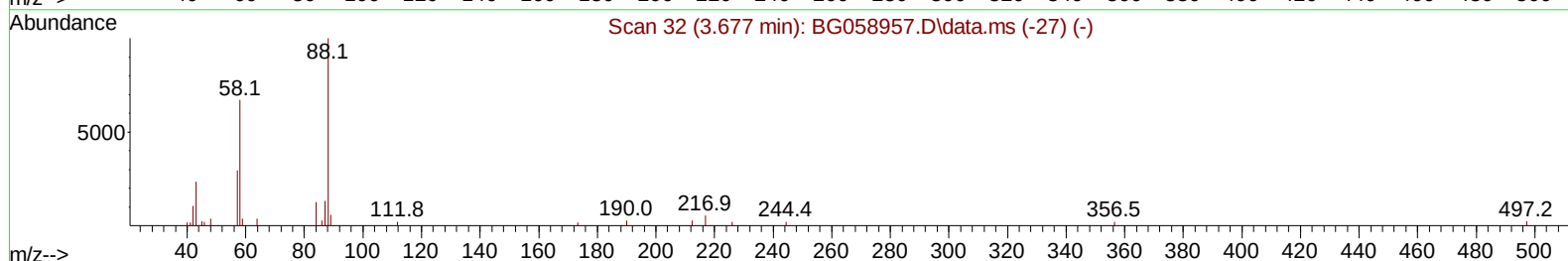
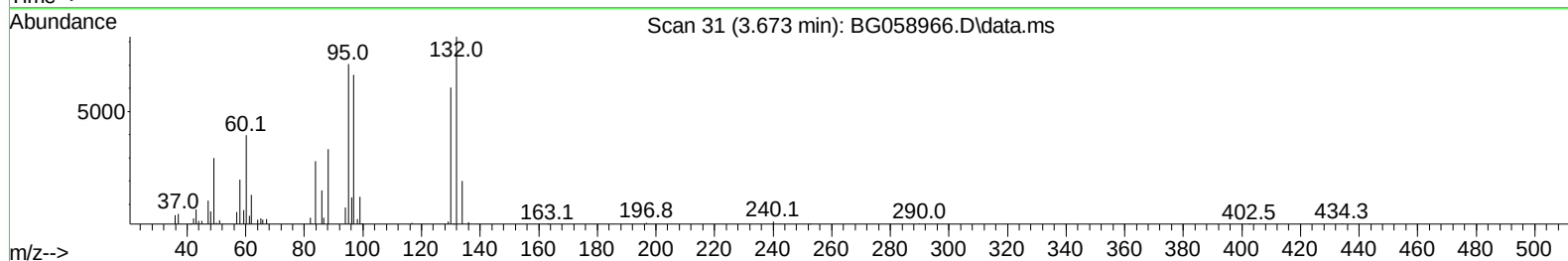
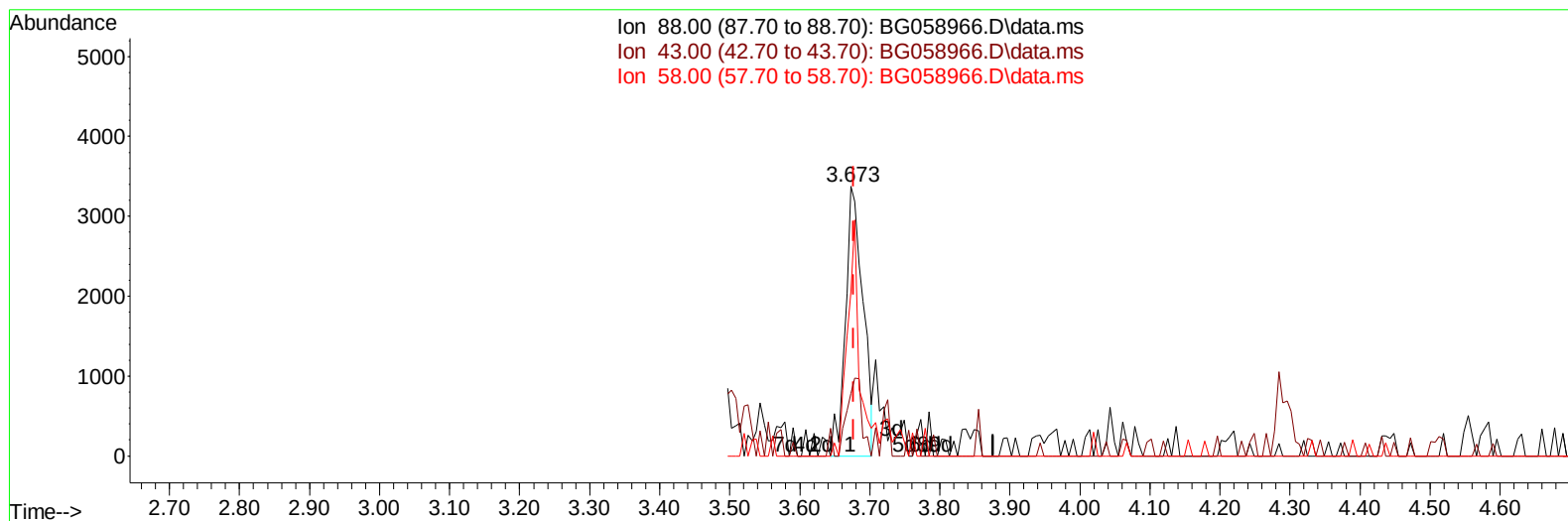
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG058966.D
Acq On : 11 Sep 2023 2:25
Operator : MA/JU
Sample : 04247-17
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BBJ77

Manual IntegrationsAPPROVED

Quant Time: Sep 11 03:20:05 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: BG058966.D\data.ms

(2) 1,4-Dioxane

3.673min (-0.004) 3.83 ng/uL

response 5937

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	23.30	23.27
58.00	67.30	61.04
0.00	0.00	0.00

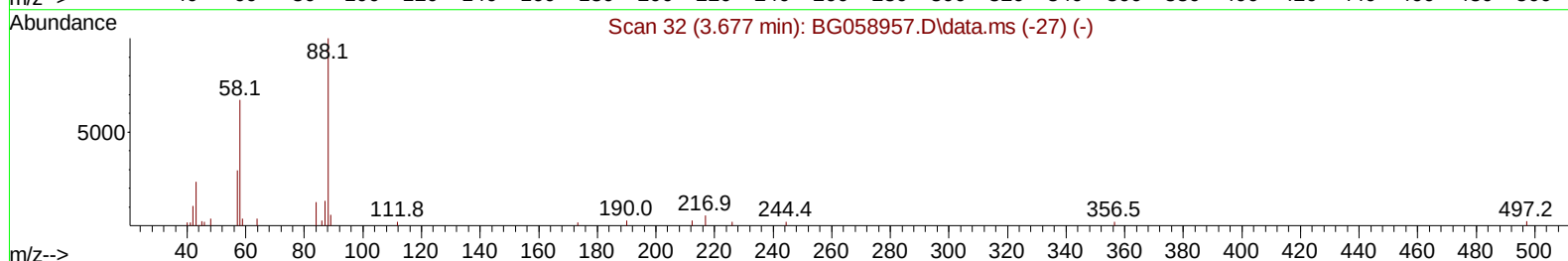
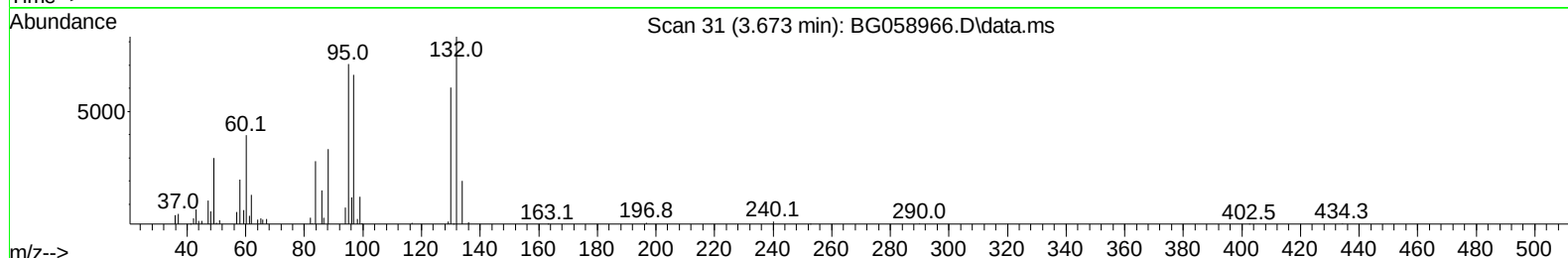
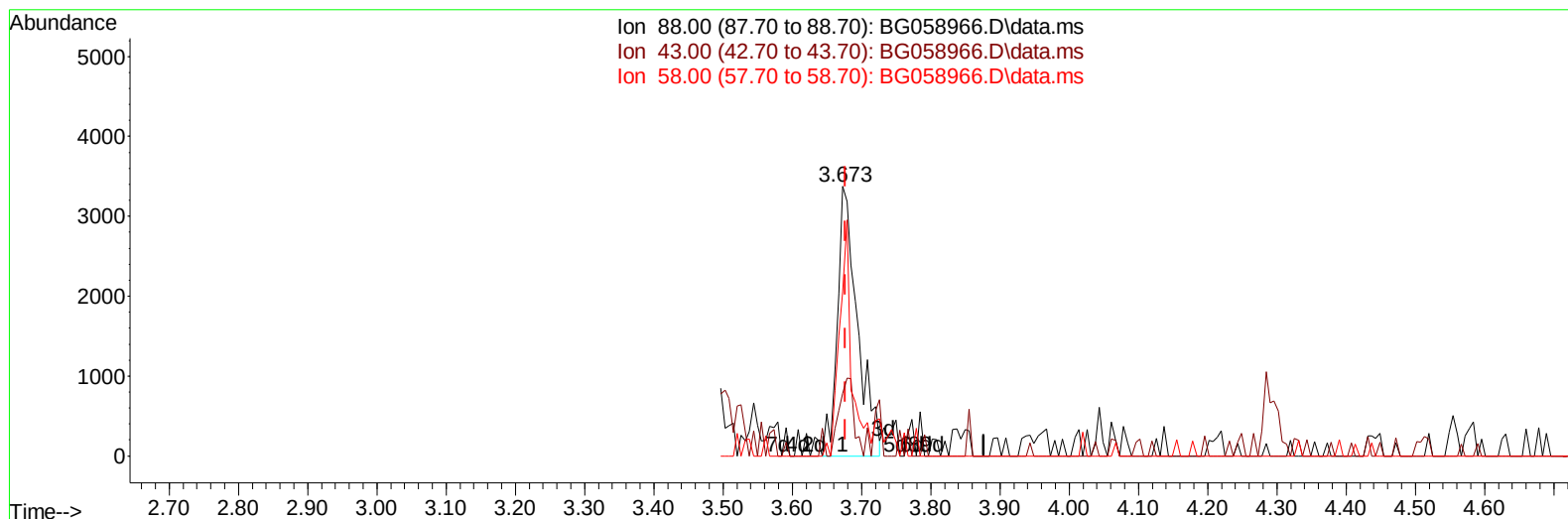
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TIC: BG058966.D\data.ms

(2) 1,4-Dioxane

3.673min (-0.004) 4.41 ng/uL m

response 6844

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	23.30	23.27
58.00	67.30	61.04
0.00	0.00	0.00

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Quant Time: Sep 11 03:21:36 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.326	152	51197	20.000	ng/ul	0.00
20) Naphthalene-d8	11.170	136	271619	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.948	164	216190	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.698	188	579454	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.016	240	563405	20.000	ng/ul	0.00
88) Perylene-d12	25.589	264	659799	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.638	96	11147	9.254	ng/uL	0.00
4) Pyridine-d5	4.073	84	42762	11.213	ng/ul	0.00
7) Phenol-d5	7.433	99	38694	7.495	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.645	67	105074	34.709	ng/ul	0.00
11) 2-Chlorophenol-d4	7.839	132	89896	27.149	ng/ul	0.00
15) 4-Methylphenol-d8	9.002	113	68941	17.284	ng/ul	0.00
21) Nitrobenzene-d5	9.519	128	61392	31.160	ng/ul	0.00
24) 2-Nitrophenol-d4	10.236	143	67910	31.413	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.759	165	119152	25.933	ng/ul	0.00
31) 4-Chloroaniline-d4	11.311	131	154799	25.054	ng/ul	0.00
46) Dimethylphthalate-d6	14.343	166	501149	29.884	ng/ul	0.00
49) Acenaphthylene-d8	14.648	160	530293	28.397	ng/ul	0.00
54) 4-Nitrophenol-d4	15.107	143	19066	6.845	ng/ul	0.00
60) Fluorene-d10	15.935	176	447675	30.081	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.035	200	77069	27.095	ng/ul	0.00
73) Anthracene-d10	17.798	188	797840	31.016	ng/ul	0.00
81) Pyrene-d10	20.066	212	927460	31.076	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.342	264	992500	30.686	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.673	88	6844m	4.410	ng/uL	Qvalue

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

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