

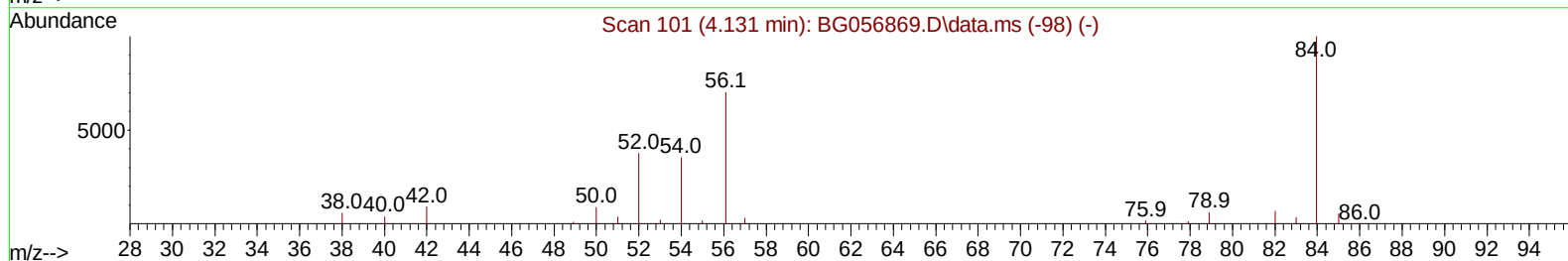
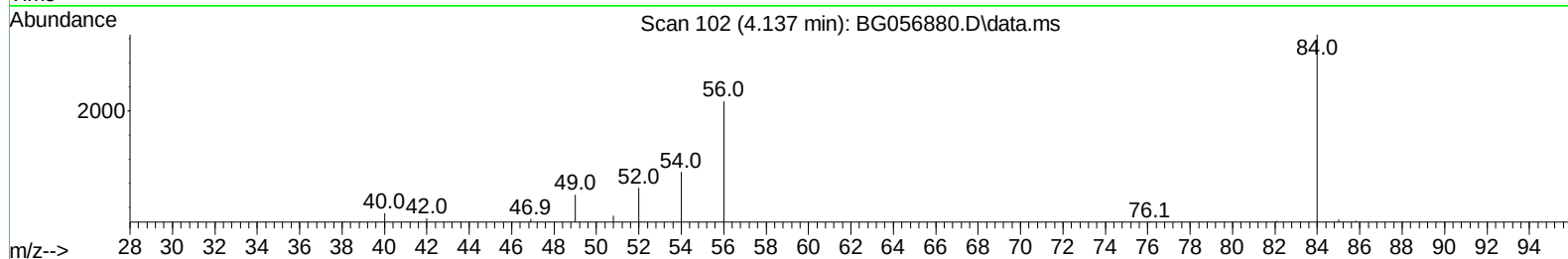
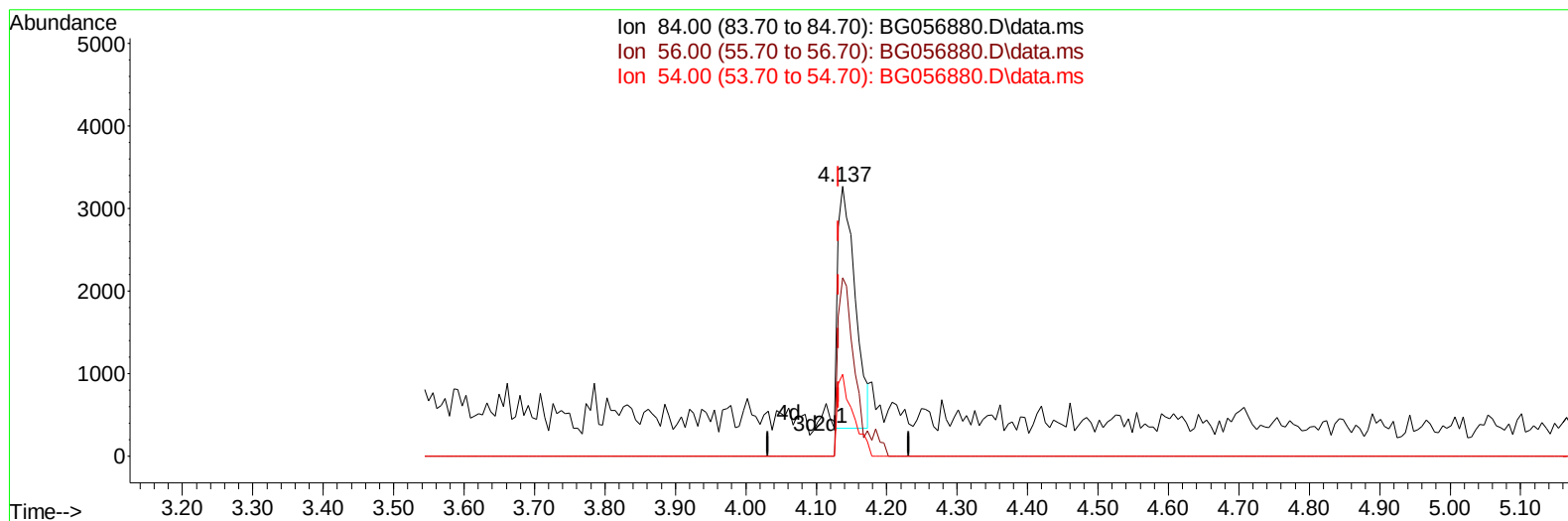
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030823\
 Data File : BG056880.D
 Acq On : 8 Mar 2023 13:36
 Operator : CG/JU
 Sample : 01784-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCAD2

Manual IntegrationsAPPROVED

Quant Time: Mar 08 14:17:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 07 23:54:44 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/09/2023
 Supervised By :Jagrut Upadhyay 03/09/2023



TIC: BG056880.D\data.ms

(4) Pyridine-d5 (S)

4.137min (+ 0.006) 4.54 ng/u1

response 4935

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.50	66.23
54.00	34.90	30.28
0.00	0.00	0.00

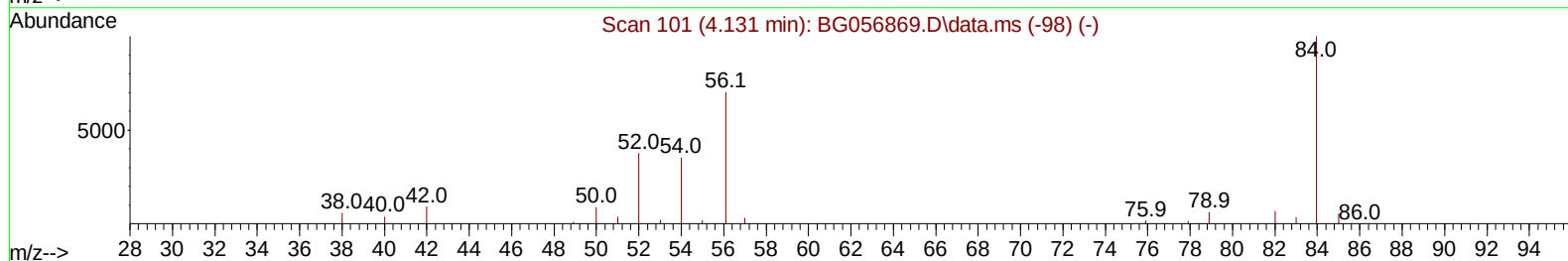
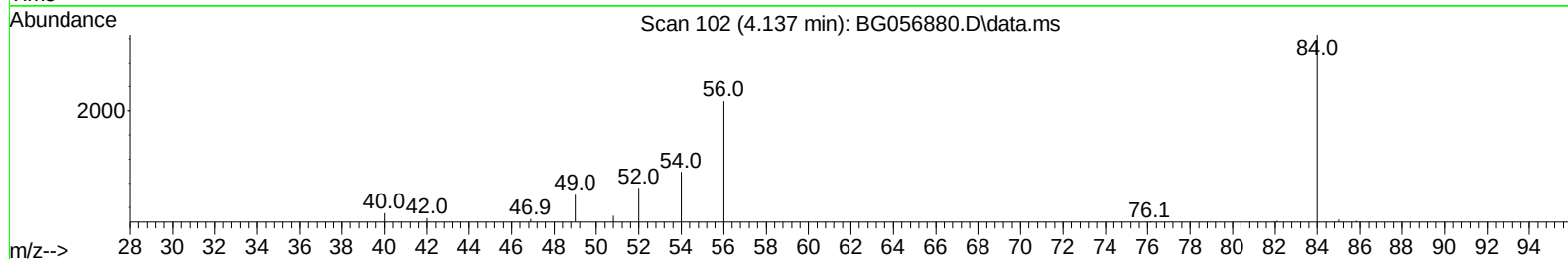
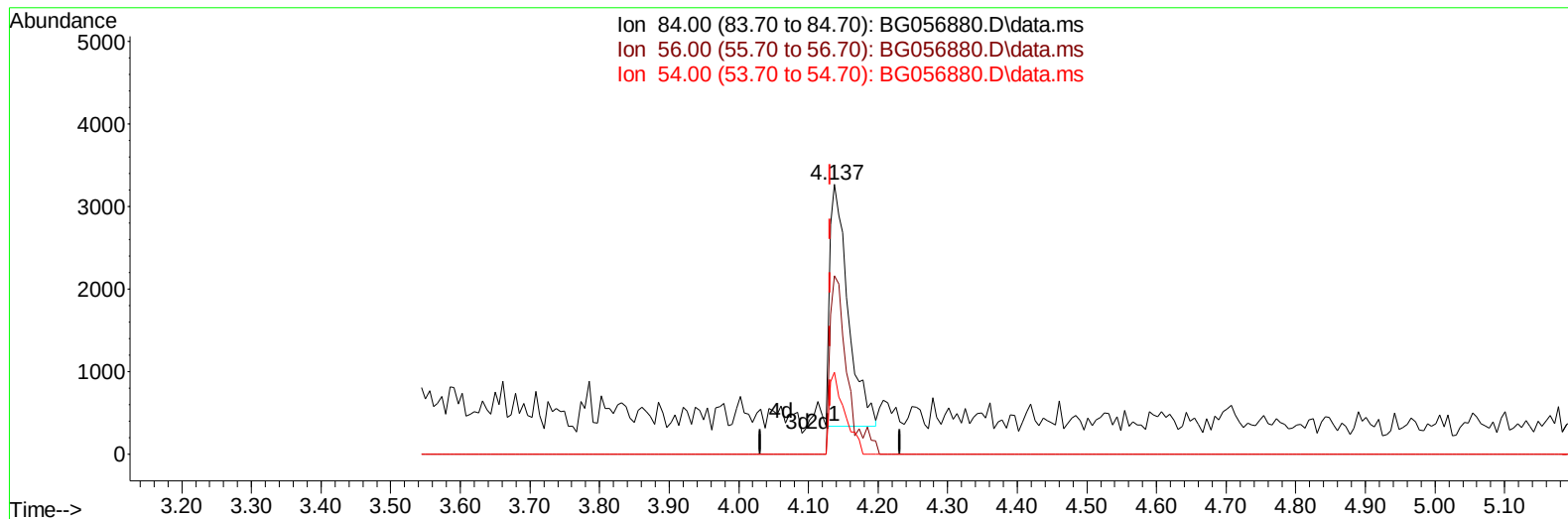
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TIC: BG056880.D\data.ms

(4) Pyridine-d5 (S)

4.137min (+ 0.006) 4.92 ng/ul m

response 5346

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.50	66.23
54.00	34.90	30.28
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	8.421	152	12949	20.000	ng/ul	0.00
20) Naphthalene-d8	11.258	136	52203	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.031	164	44561	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.774	188	120625	20.000	ng/ul	0.00
79) Chrysene-d12	22.099	240	123733	20.000	ng/ul	-0.01
88) Perylene-d12	25.648	264	132048	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.691	96	1506	4.510	ng/uL	0.00
4) Pyridine-d5	4.137	84	5346m	4.922	ng/ul	0.00
7) Phenol-d5	7.539	99	8520	6.420	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.716	67	24777	29.722	ng/ul	0.00
11) 2-Chlorophenol-d4	7.939	132	19704	23.722	ng/ul	0.00
15) 4-Methylphenol-d8	9.108	113	15640	15.618	ng/ul	0.00
21) Nitrobenzene-d5	9.590	128	13811	31.898	ng/ul	0.00
24) 2-Nitrophenol-d4	10.324	143	16177	31.582	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.865	165	26626	26.836	ng/ul	0.00
31) 4-Chloroaniline-d4	11.382	131	27917	21.354	ng/ul	0.00
46) Dimethylphthalate-d6	14.414	166	125210	32.850	ng/ul	0.00
49) Acenaphthylene-d8	14.731	160	129671	30.984	ng/ul	0.00
54) 4-Nitrophenol-d4	15.189	143	5210	8.067	ng/ul	-0.01
60) Fluorene-d10	16.018	176	112894	31.941	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.118	200	27825	31.918	ng/ul	0.00
73) Anthracene-d10	17.868	188	216841	37.194	ng/ul	-0.01
81) Pyrene-d10	20.136	212	277047	36.760	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.407	264	258960	35.815	ng/ul	-0.01

Target Compounds Qvalue

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

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