

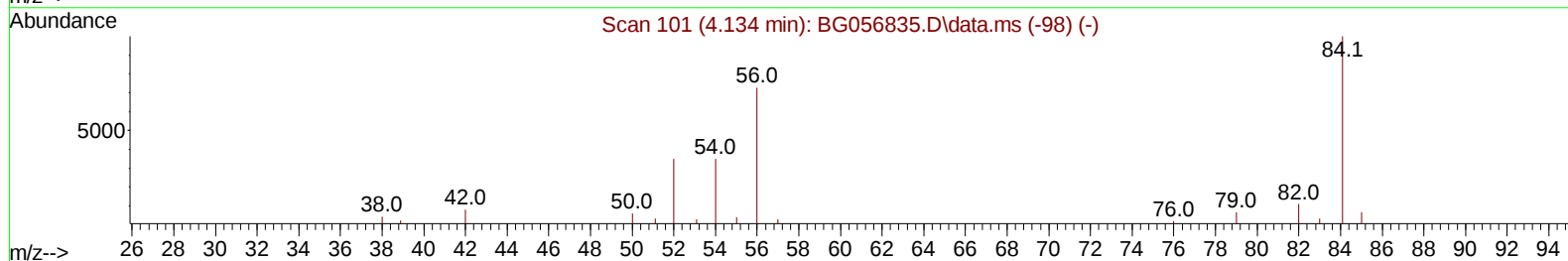
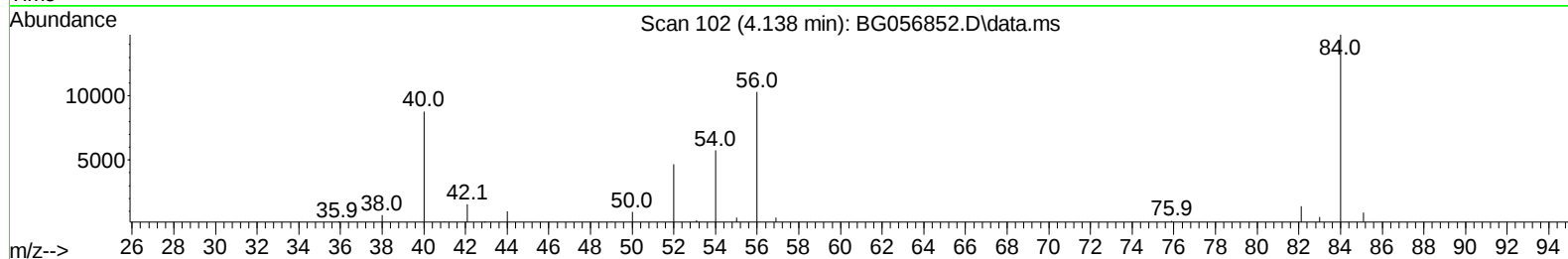
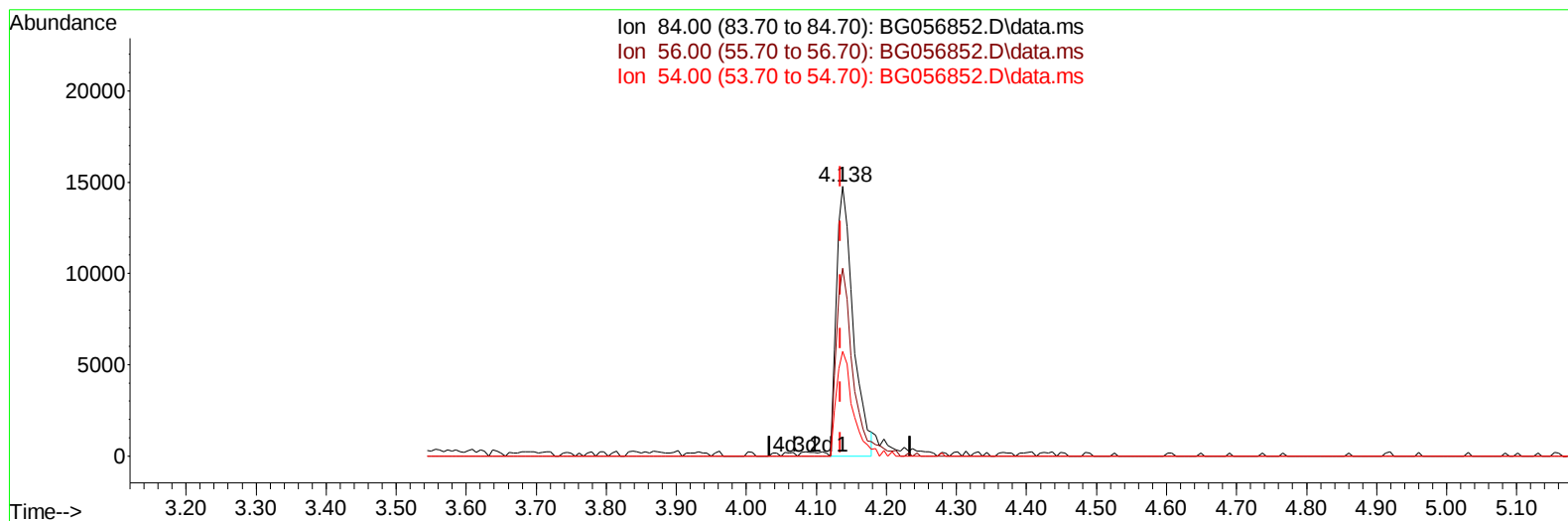
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056852.D
Acq On : 6 Mar 2023 22:14
Operator : CG/JU
Sample : 01712-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX9

Manual IntegrationsAPPROVED

Quant Time: Mar 07 00:55:31 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 02 00:12:28 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/07/2023
Supervised By :Jagrut Upadhyay 03/07/2023



TIC: BG056852.D\data.ms

(4) Pyridine-d5 (S)

4.138min (+ 0.004) 19.35 ng/u1

response 24665

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	67.40	69.78
54.00	38.20	38.85
0.00	0.00	0.00

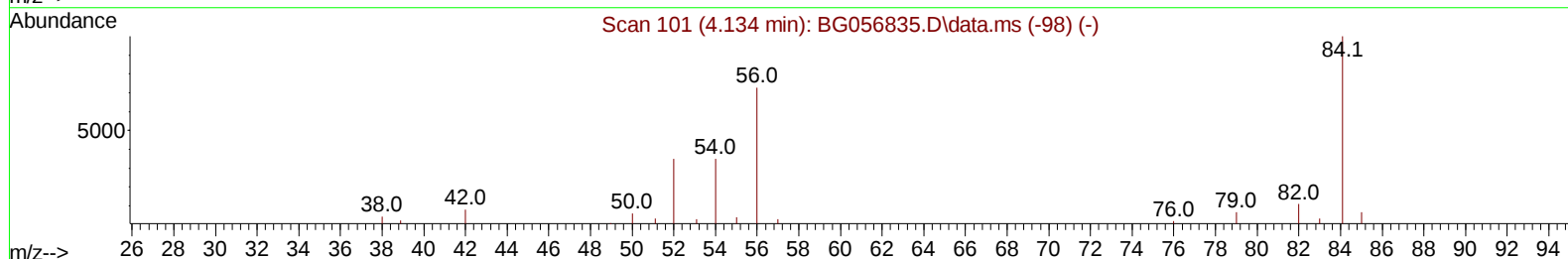
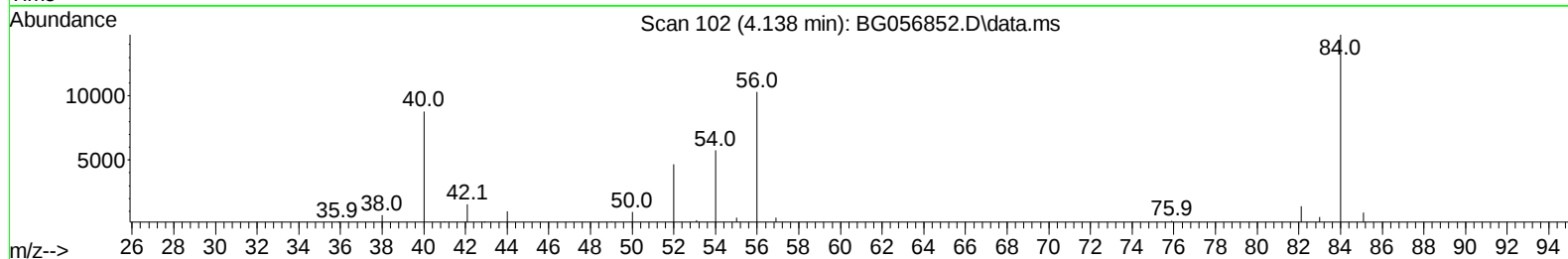
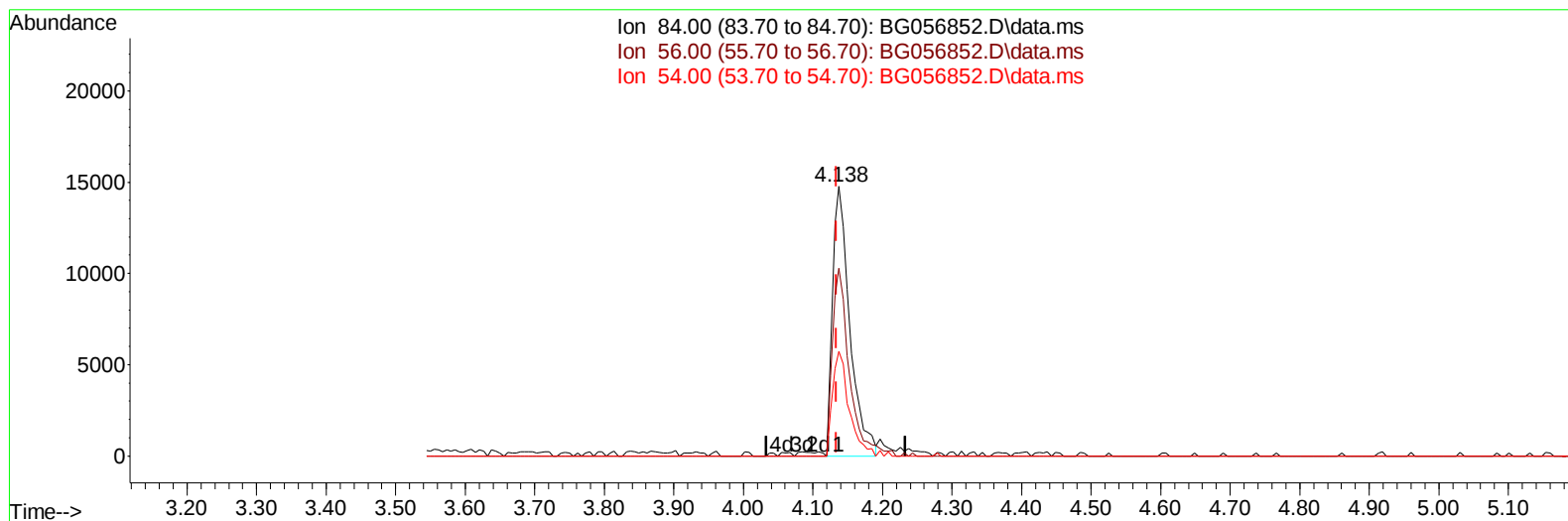
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TIC: BG056852.D\data.ms

(4) Pyridine-d5 (S)

4.138min (+ 0.004) 19.82 ng/ul m

response 25255

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	67.40	69.78
54.00	38.20	38.85
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	17514	20.000	ng/ul	0.00
20) Naphthalene-d8	11.253	136	78683	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.019	164	58152	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.757	188	128747	20.000	ng/ul	0.00
79) Chrysene-d12	22.064	240	116232	20.000	ng/ul	0.00
88) Perylene-d12	25.589	264	126747	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.703	96	1635	5.179	ng/uL	0.00
4) Pyridine-d5	4.138	84	25255m	19.817	ng/ul	0.00
7) Phenol-d5	7.539	99	39679	22.640	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.716	67	26292	24.593	ng/ul	0.00
11) 2-Chlorophenol-d4	7.939	132	27403	23.545	ng/ul	0.00
15) 4-Methylphenol-d8	9.108	113	32116	23.982	ng/ul	0.00
21) Nitrobenzene-d5	9.596	128	15902	24.116	ng/ul	0.00
24) 2-Nitrophenol-d4	10.318	143	19152	24.400	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.859	165	35714	24.152	ng/ul	0.00
31) 4-Chloroaniline-d4	11.376	131	39899	20.688	ng/ul	0.00
46) Dimethylphthalate-d6	14.408	166	115351	24.585	ng/ul	0.00
49) Acenaphthylene-d8	14.719	160	135496	25.409	ng/ul	0.00
54) 4-Nitrophenol-d4	15.178	143	14121	17.457	ng/ul	0.00
60) Fluorene-d10	16.006	176	106581	24.557	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.106	200	16553	19.217	ng/ul	0.00
73) Anthracene-d10	17.851	188	162786	26.062	ng/ul	0.00
81) Pyrene-d10	20.113	212	192488	24.473	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.348	264	180563	26.393	ng/ul	0.00

Target Compounds Qvalue

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

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