

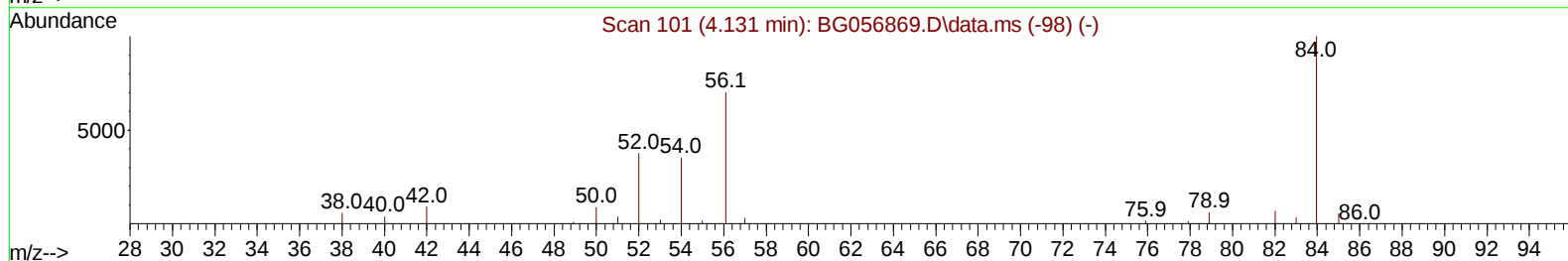
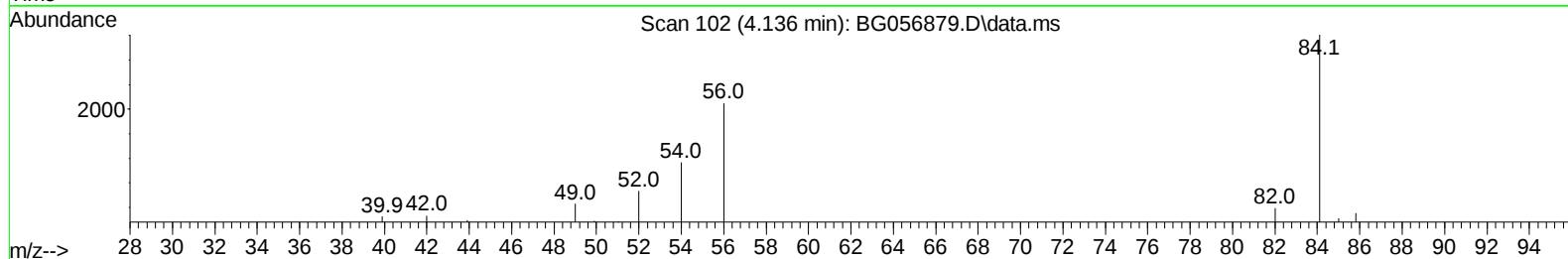
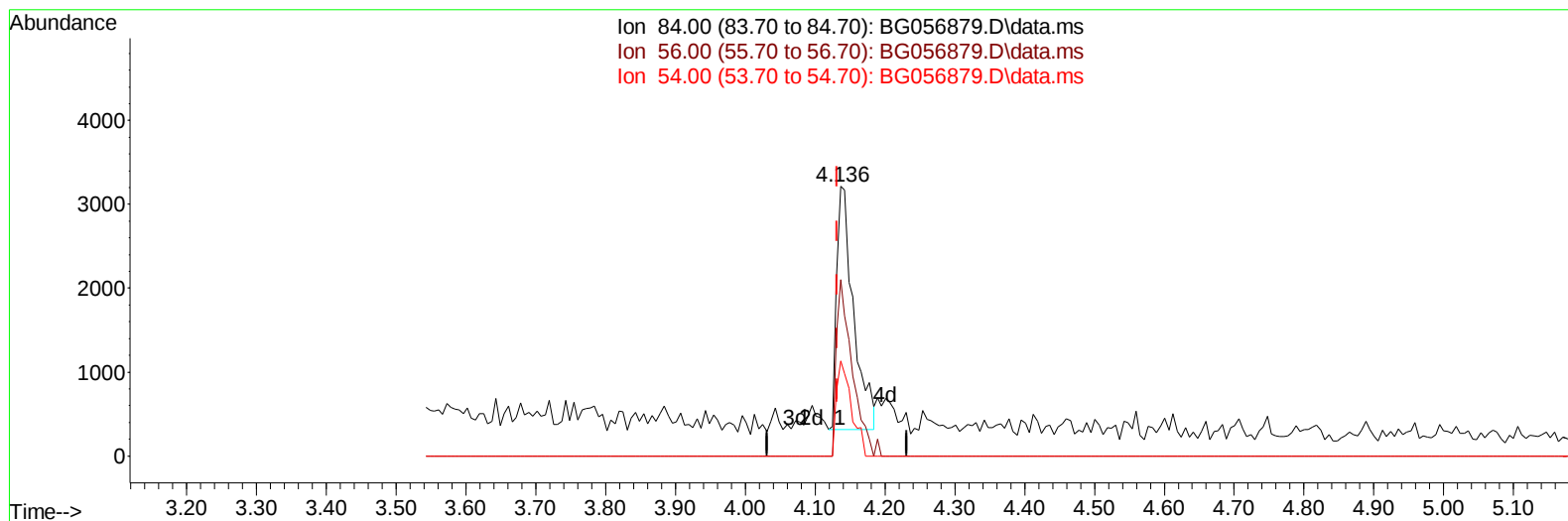
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030823\
 Data File : BG056879.D
 Acq On : 8 Mar 2023 12:55
 Operator : CG/JU
 Sample : 01784-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCAD1

Manual IntegrationsAPPROVED

Quant Time: Mar 08 14:17:38 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: BG056879.D\data.ms

(4) Pyridine-d5 (S)

4.136min (+ 0.005) 4.54 ng/u1

response 4813

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.50	65.35
54.00	34.90	35.21
0.00	0.00	0.00

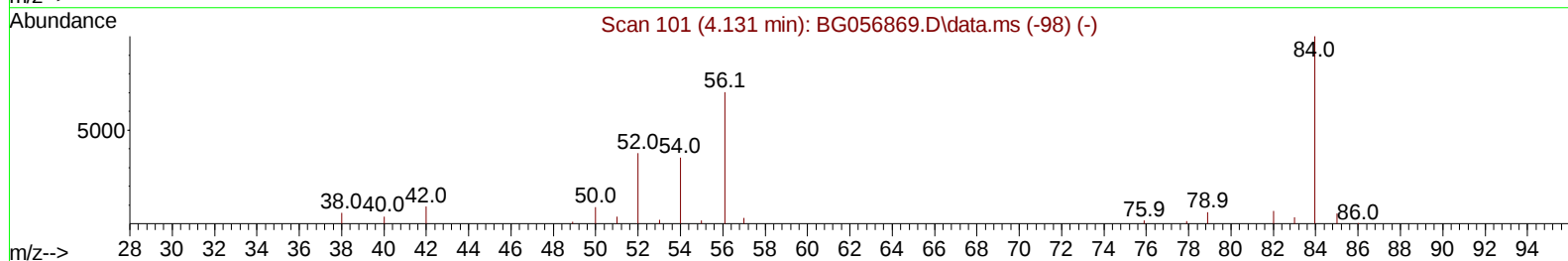
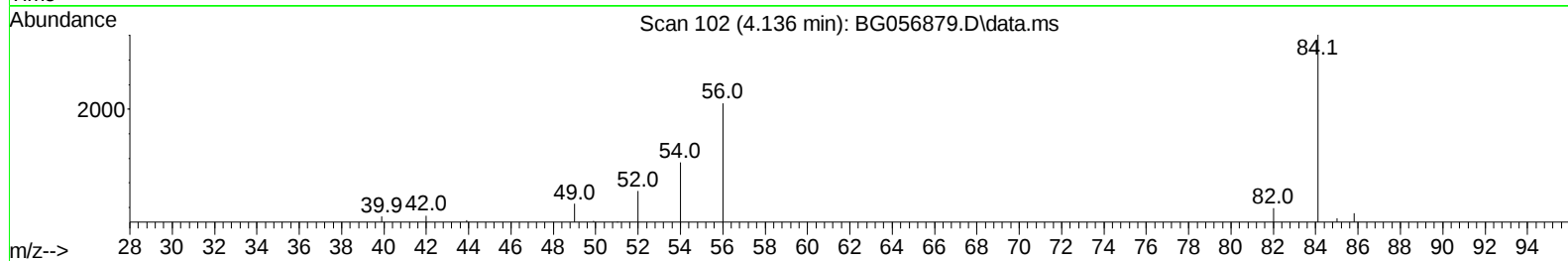
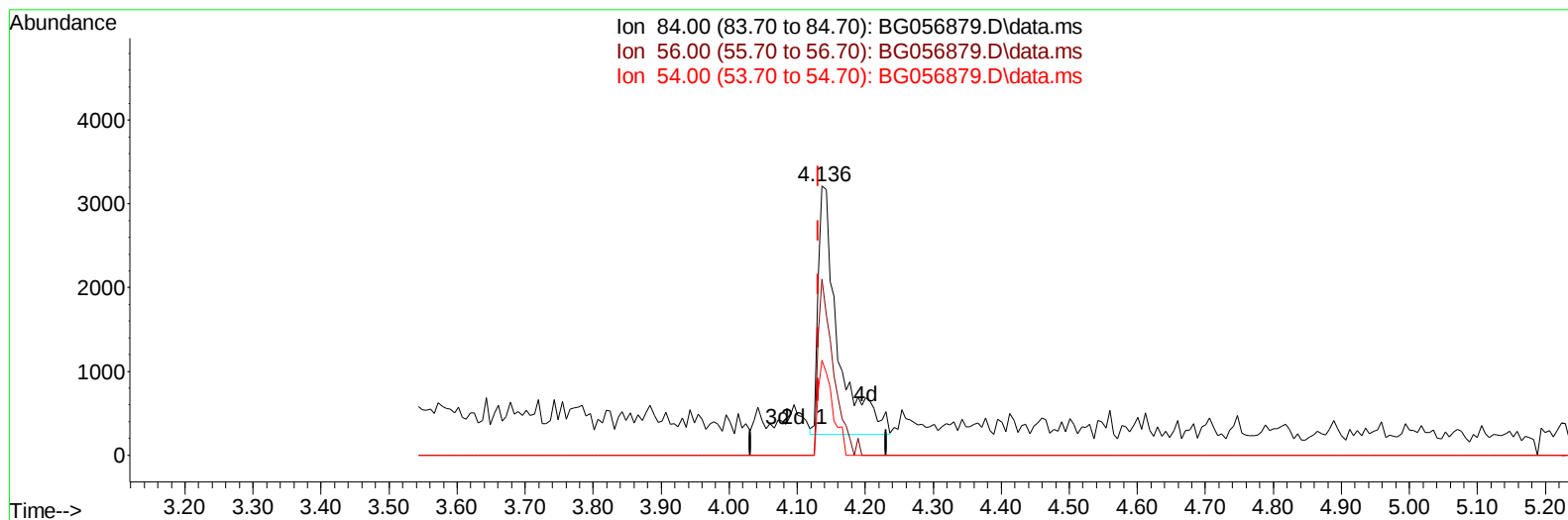
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(4) Pyridine-d5 (S)

4.136min (+ 0.005) 5.64 ng/u1 m

response 5974

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	69.50	65.35
54.00	34.90	35.21
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.414	152	12637	20.000	ng/ul	0.00
20) Naphthalene-d8	11.258	136	53366	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.035	164	43715	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.773	188	123823	20.000	ng/ul	0.00
79) Chrysene-d12	22.098	240	123912	20.000	ng/ul	-0.01
88) Perylene-d12	25.647	264	133271	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.690	96	1661	5.097	ng/uL	0.00
4) Pyridine-d5	4.136	84	5974m	5.635	ng/ul	0.00
7) Phenol-d5	7.538	99	8645	6.675	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.720	67	24941	30.657	ng/ul	0.00
11) 2-Chlorophenol-d4	7.938	132	20285	25.025	ng/ul	0.00
15) 4-Methylphenol-d8	9.101	113	15577	15.939	ng/ul	0.00
21) Nitrobenzene-d5	9.595	128	14165	32.002	ng/ul	0.00
24) 2-Nitrophenol-d4	10.323	143	16149	30.841	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.864	165	27183	26.801	ng/ul	0.00
31) 4-Chloroaniline-d4	11.381	131	27991	20.944	ng/ul	0.00
46) Dimethylphthalate-d6	14.419	166	124499	33.295	ng/ul	0.00
49) Acenaphthylene-d8	14.730	160	130617	31.814	ng/ul	0.00
54) 4-Nitrophenol-d4	15.188	143	5049	7.969	ng/ul	-0.01
60) Fluorene-d10	16.017	176	111484	32.152	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.116	200	26646	29.776	ng/ul	0.00
73) Anthracene-d10	17.873	188	213924	35.746	ng/ul	0.00
81) Pyrene-d10	20.135	212	268325	35.551	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.406	264	252832	34.647	ng/ul	-0.01

Target Compounds Qvalue

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

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