

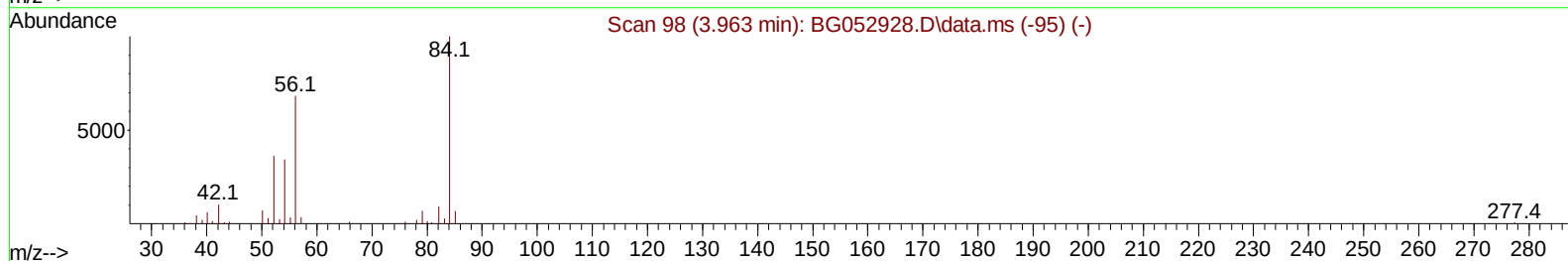
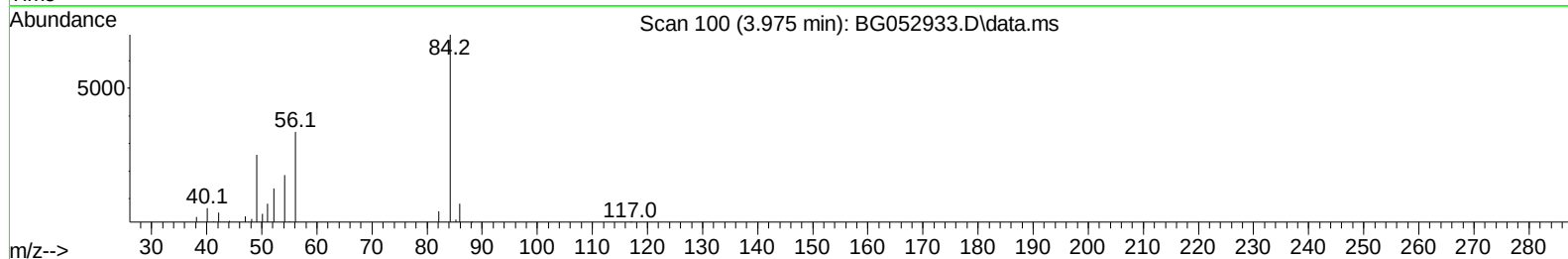
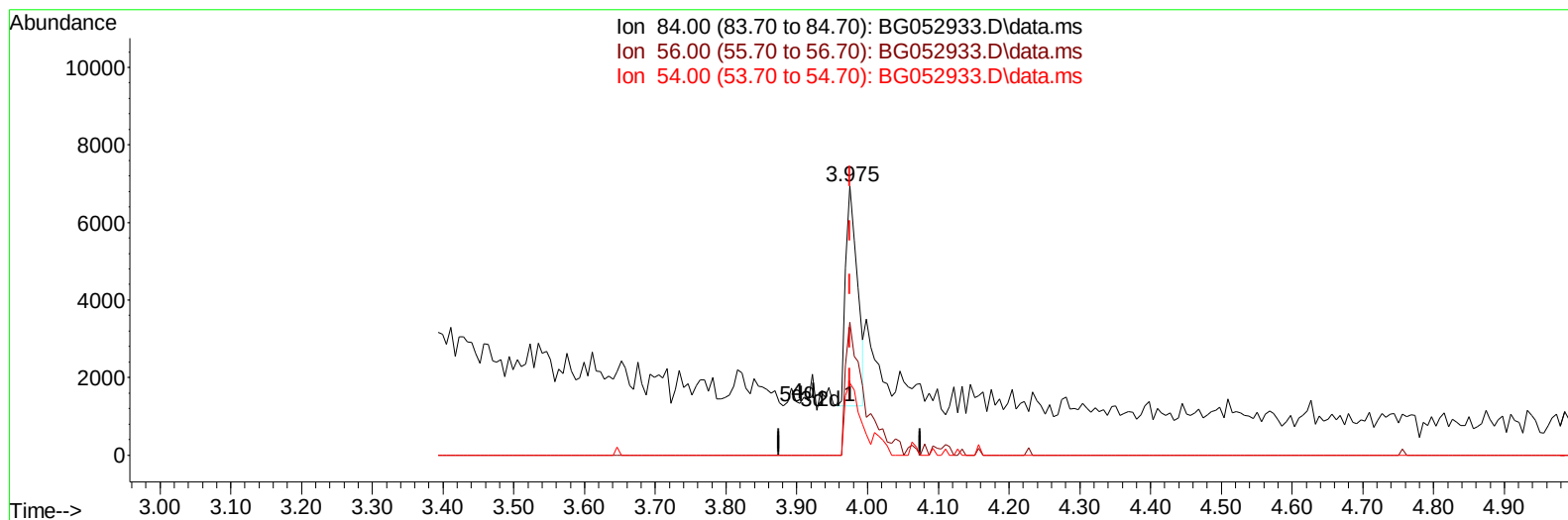
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
Data File : BG052933.D
Acq On : 29 Mar 2022 19:01
Operator : CG/JU
Sample : N2097-20
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF0

Manual IntegrationsAPPROVED

Quant Time: Mar 30 01:10:38 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 23 13:16:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/30/2022
Supervised By :Yogesh Patel 04/02/2022



TIC: BG052933.D\data.ms

(4) Pyridine-d5 (S)

3.975min (-0.000) 3.87 ng/u1

response 6485

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	49.47#
54.00	28.70	26.83
0.00	0.00	0.00

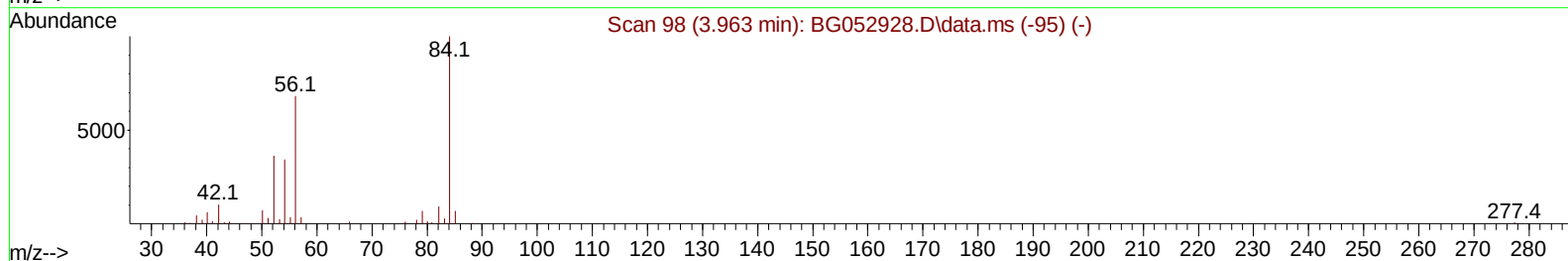
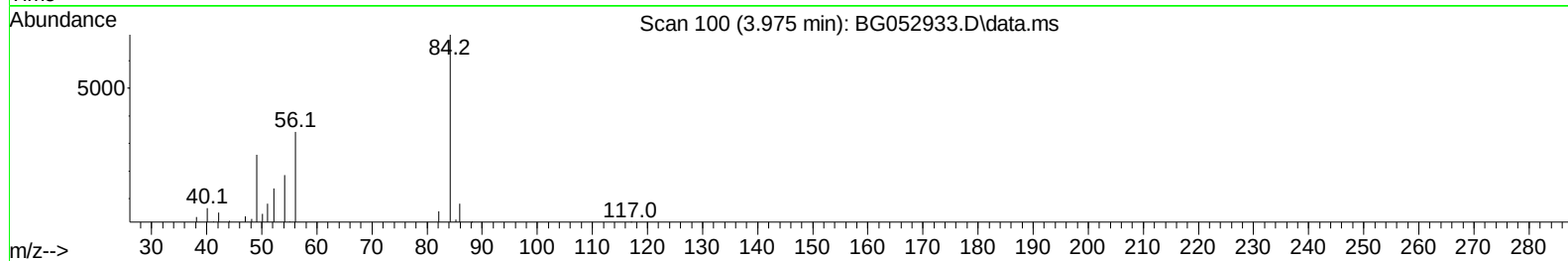
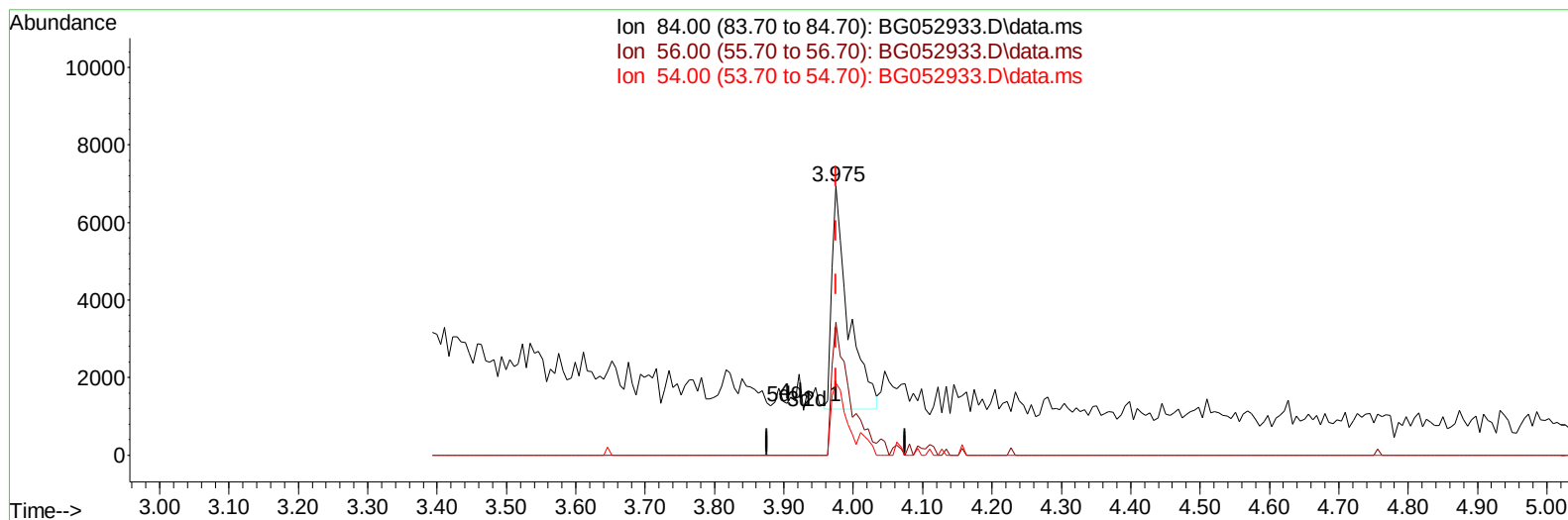
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TIC: BG052933.D\data.ms

(4) Pyridine-d5 (S)

3.975min (-0.000) 5.64 ng/ul m

response 9444

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	49.47#
54.00	28.70	26.83
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	8.134	152	23702	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.942	136	109211	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.754	164	89751	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	222395	20.000	ng/ul	0.00
79) Chrysene-d12	21.780	240	245670	20.000	ng/ul	#-0.01
88) Perylene-d12	25.093	264	244713	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.552	96	2760	4.113	ng/uL	0.00
4) Pyridine-d5	3.975	84	9444m	5.638	ng/ul	0.00
7) Phenol-d5	7.282	99	15752	7.420	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.453	67	44233	34.116	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	36426	25.270	ng/ul	-0.01
15) 4-Methylphenol-d8	8.827	113	28752	18.697	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	26011	33.087	ng/ul	0.00
24) 2-Nitrophenol-d4	10.020	143	28628	34.526	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	51510	28.555	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	42183	17.693	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	240359	34.860	ng/ul	-0.01
49) Acenaphthylene-d8	14.449	160	256807	30.651	ng/ul	0.00
54) 4-Nitrophenol-d4	14.925	143	8232	7.101	ng/ul	0.00
60) Fluorene-d10	15.741	176	209395	34.587	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.847	200	39918	35.271	ng/ul	-0.01
73) Anthracene-d10	17.598	188	372396	35.959	ng/ul	0.00
81) Pyrene-d10	19.877	212	451148	32.827	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.864	264	458016	36.223	ng/ul	-0.01

Target Compounds Qvalue

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

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