

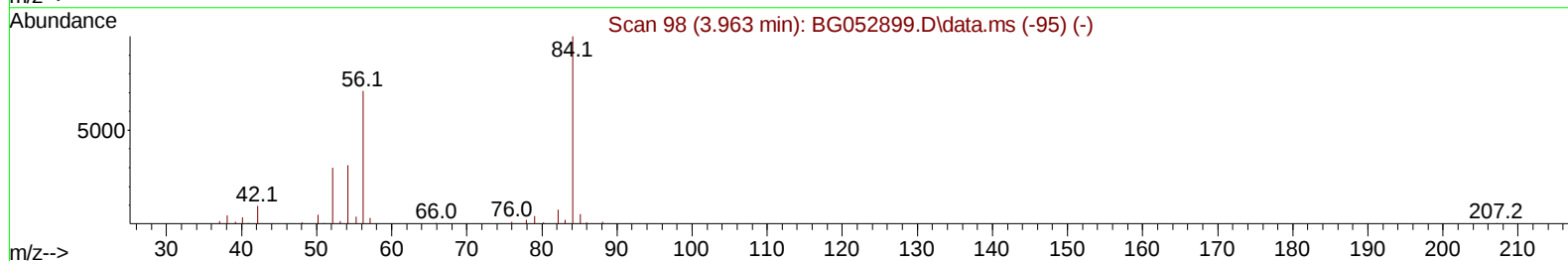
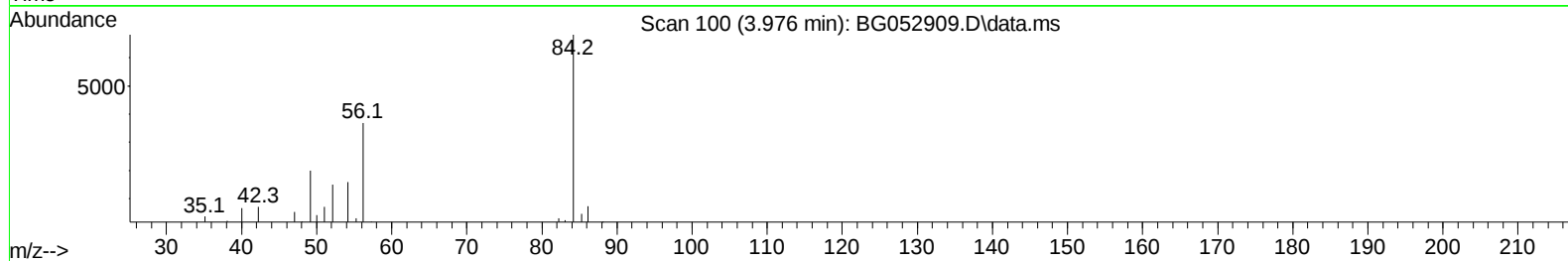
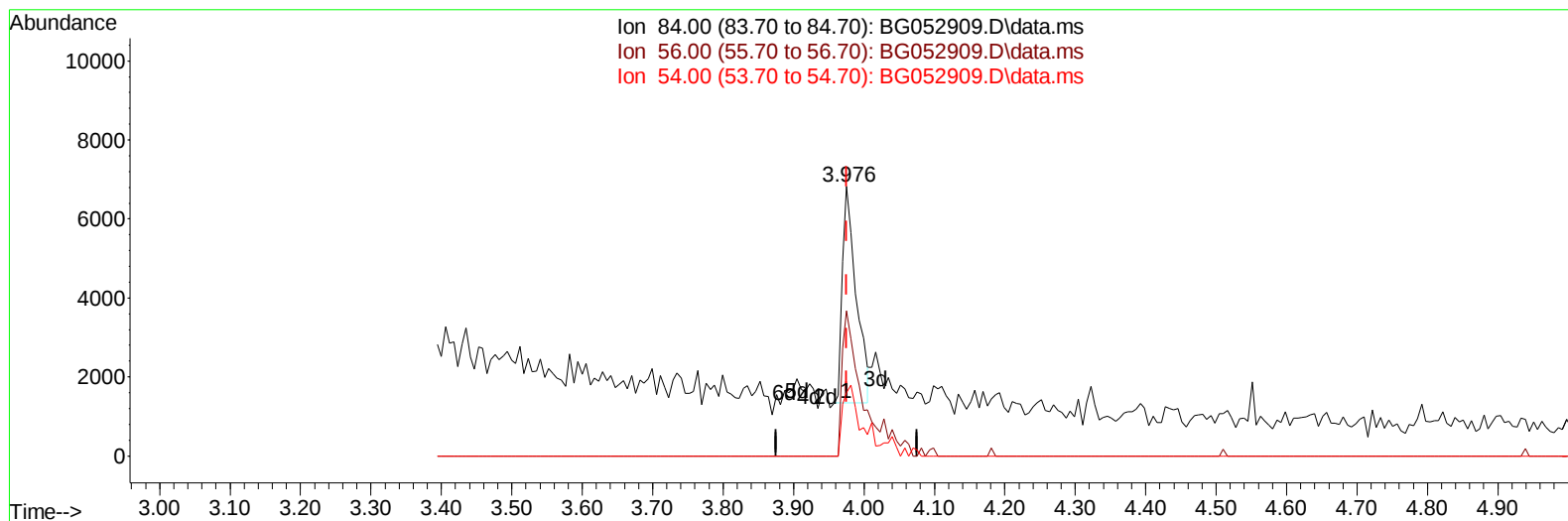
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
 Data File : BG052909.D
 Acq On : 29 Mar 2022 2:59
 Operator : CG/JU
 Sample : N2100-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AF9

Manual IntegrationsAPPROVED

Quant Time: Mar 29 05:39:32 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 03/29/2022
 Supervised By :mohammad ahmed 03/29/2022



TIC: BG052909.D\data.ms

(4) Pyridine-d5 (S)

3.976min (+ 0.000) 4.01 ng/u1

response 7392

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	53.98
54.00	28.70	23.37
0.00	0.00	0.00

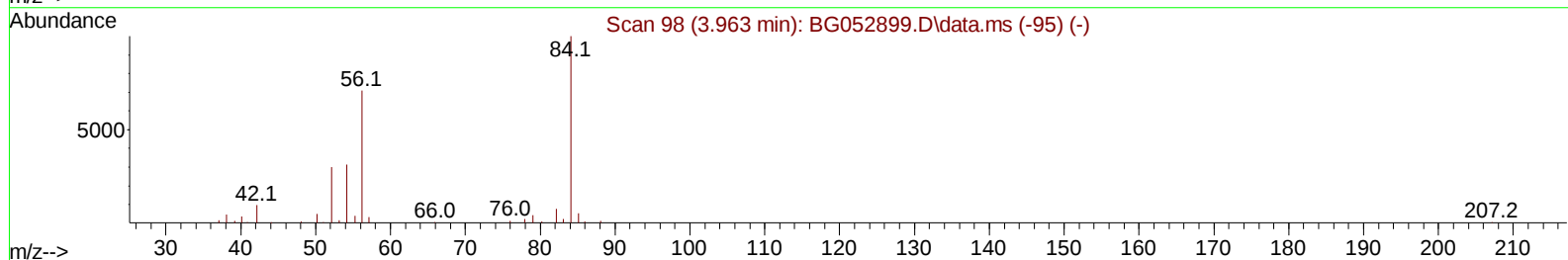
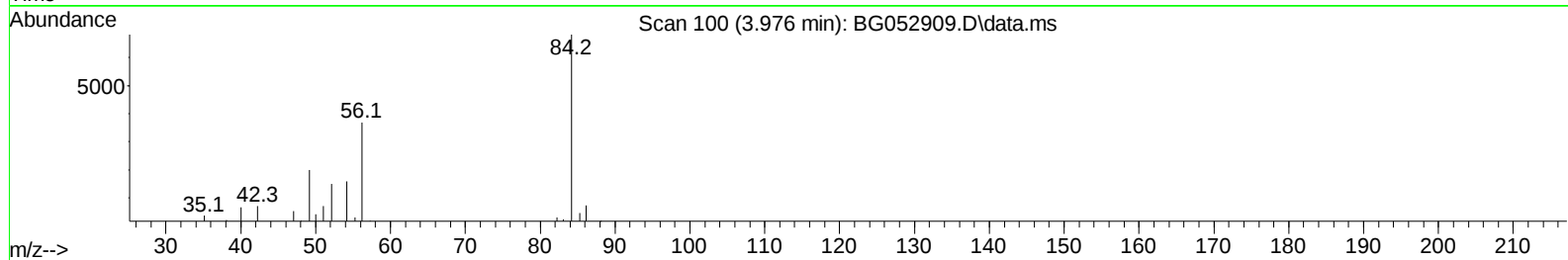
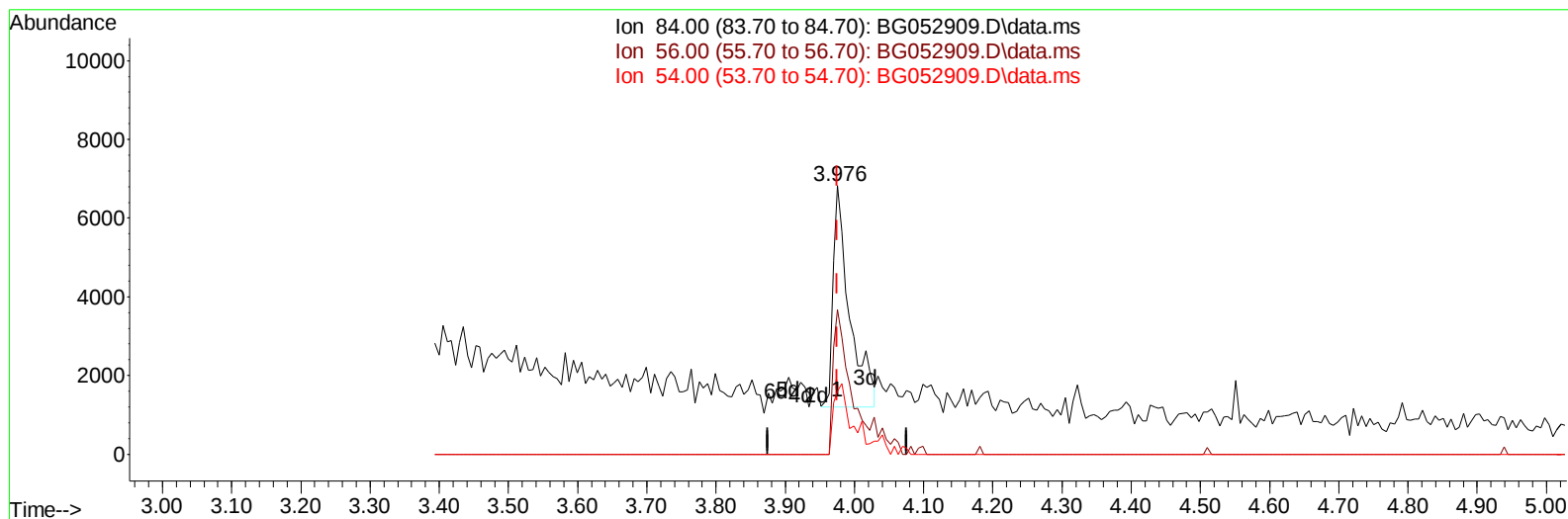
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TIC: BG052909.D\data.ms

(4) Pyridine-d5 (S)

3.976min (+ 0.000) 5.02 ng/u1 m

response 9253

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	53.98
54.00	28.70	23.37
0.00	0.00	0.00

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 Misc :
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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	8.135	152	26088	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.948	136	120205	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.755	164	97242	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	239284	20.000	ng/ul	0.00
79) Chrysene-d12	21.781	240	259798	20.000	ng/ul	-0.01
88) Perylene-d12	25.094	264	261275	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.553	96	3608	4.885	ng/uL	0.00
4) Pyridine-d5	3.976	84	9253m	5.018	ng/ul	0.00
7) Phenol-d5	7.277	99	18625	7.971	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.447	67	53063	37.183	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.665	132	44291	27.916	ng/ul	-0.01
15) 4-Methylphenol-d8	8.828	113	32325	19.098	ng/ul	0.00
21) Nitrobenzene-d5	9.292	128	32313	37.344	ng/ul	0.00
24) 2-Nitrophenol-d4	10.020	143	35147	38.511	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.561	165	61392	30.920	ng/ul	0.00
31) 4-Chloroaniline-d4	11.072	131	74297	28.313	ng/ul	0.00
46) Dimethylphthalate-d6	14.150	166	274636	36.762	ng/ul	-0.01
49) Acenaphthylene-d8	14.450	160	308852	34.023	ng/ul	0.00
54) 4-Nitrophenol-d4	14.925	143	8545	6.803	ng/ul	0.00
60) Fluorene-d10	15.742	176	247797	37.777	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.848	200	44514	36.555	ng/ul	-0.01
73) Anthracene-d10	17.598	188	431147	38.694	ng/ul	0.00
81) Pyrene-d10	19.878	212	528201	36.343	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.871	264	531852	39.396	ng/ul	0.00

Target Compounds Qvalue

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

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