

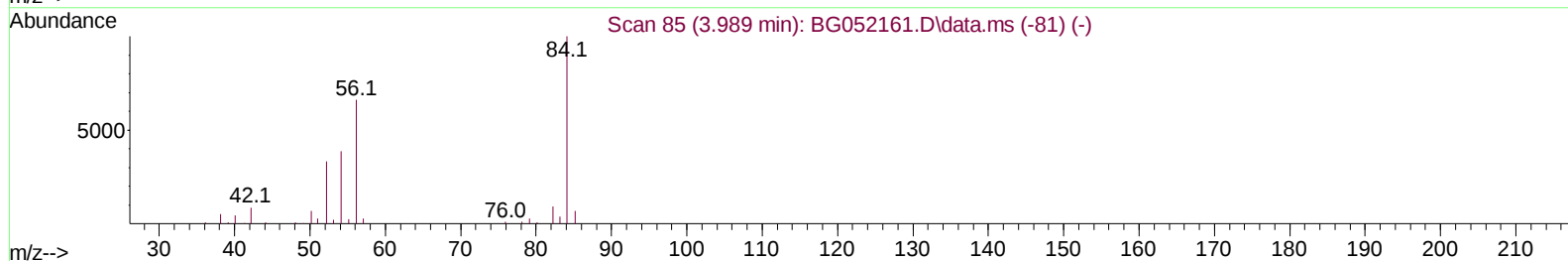
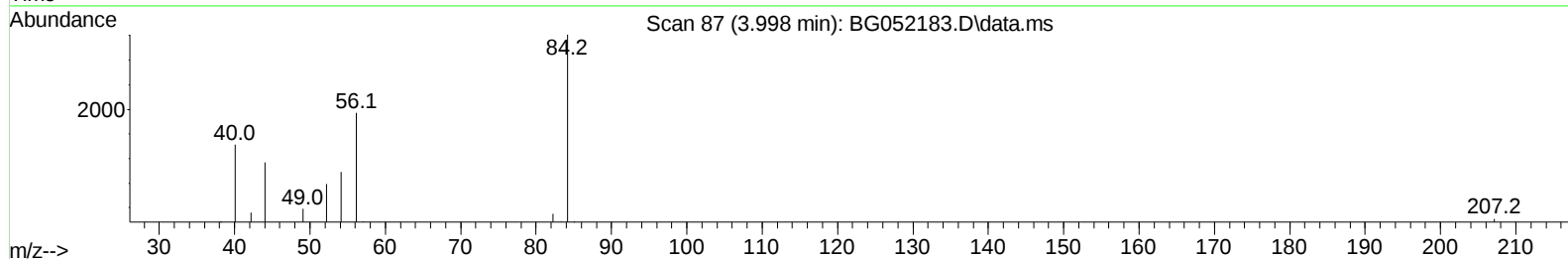
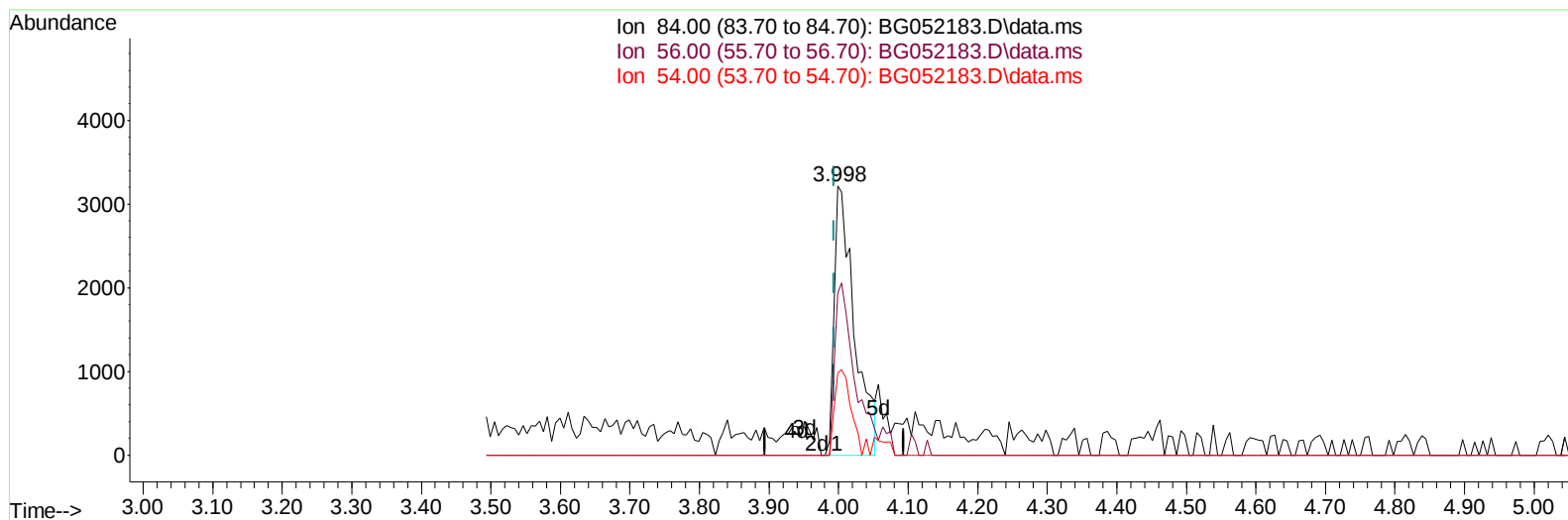
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052183.D
 Acq On : 27 Jan 2022 13:34
 Operator : CG/JU
 Sample : N1279-03
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0VD5

Manual IntegrationsAPPROVED

Quant Time: Jan 27 23:58:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/28/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052183.D\data.ms

(4) Pyridine-d5 (S)

3.998min (+ 0.005) 3.32 ng/u1

response 6474

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	60.31
54.00	31.50	30.54
0.00	0.00	0.00

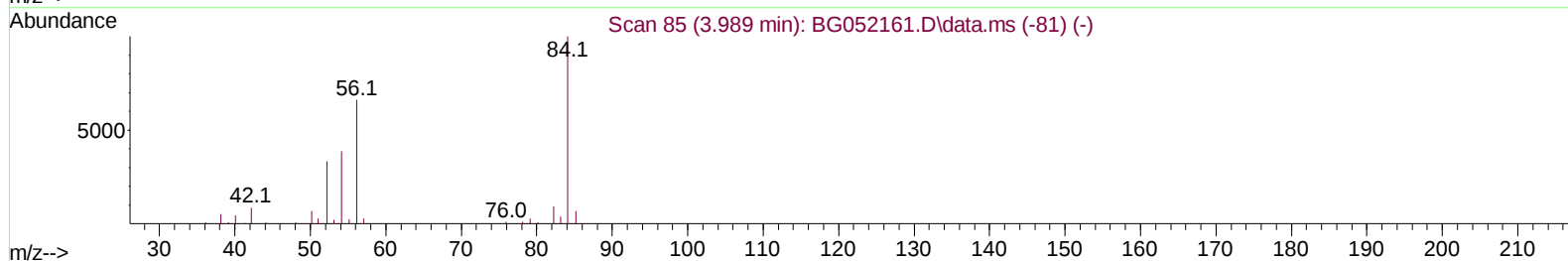
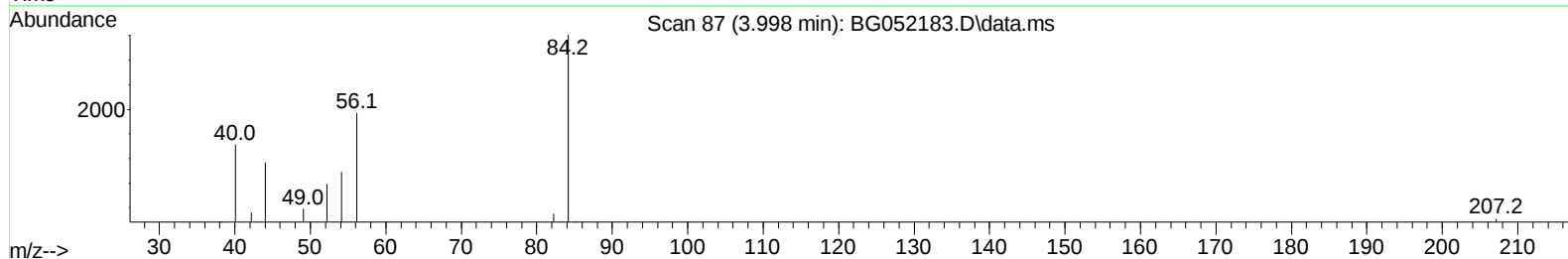
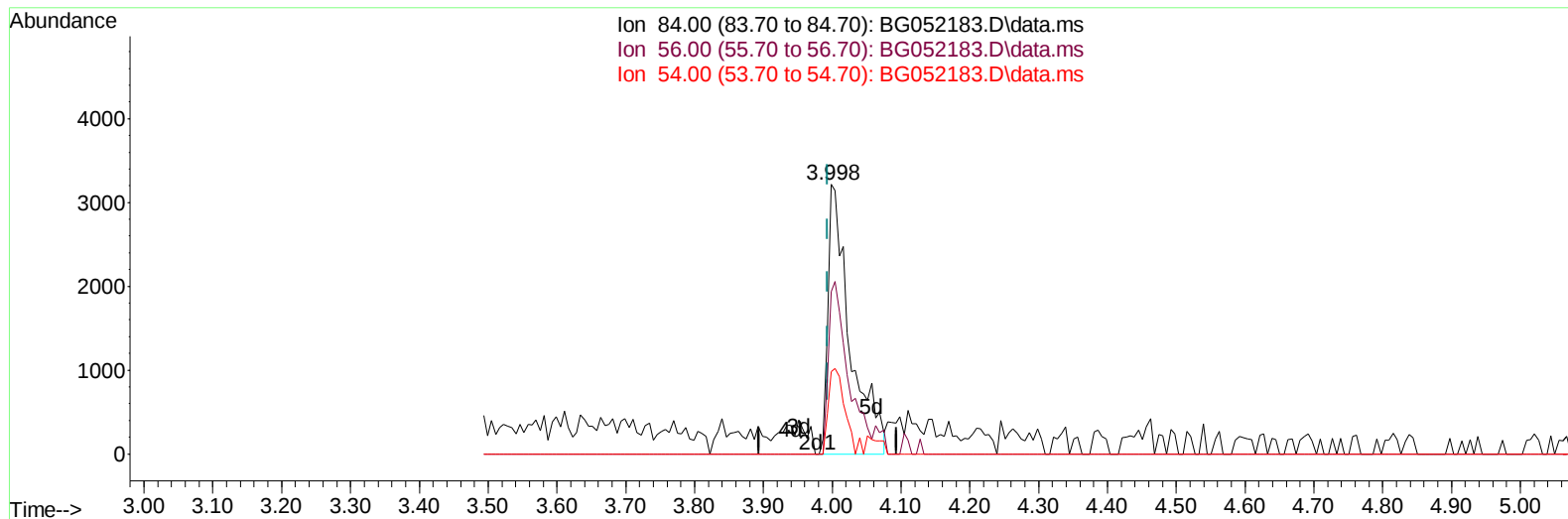
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TIC: BG052183.D\data.ms

(4) Pyridine-d5 (S)

3.998min (+ 0.005) 3.68 ng/u1 m

response 7187

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	60.31
54.00	31.50	30.54
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	8.234	152	22447	20.000	ng/ul	0.00
20) Naphthalene-d8	11.077	136	97119	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.878	164	72053	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	186973	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.945	240	188655	20.000	ng/ul	# 0.00
88) Perylene-d12	25.422	264	189852	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.564	96	2537	4.039	ng/uL	0.00
4) Pyridine-d5	3.998	84	7187m	3.680	ng/ul	0.00
7) Phenol-d5	7.382	99	12587	5.729	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	40328	28.916	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	34572	23.768	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	23800	13.936	ng/ul	0.00
21) Nitrobenzene-d5	9.409	128	24380	31.769	ng/ul	0.00
24) 2-Nitrophenol-d4	10.137	143	26386	31.089	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.689	165	46908	28.275	ng/ul	0.00
31) 4-Chloroaniline-d4	11.200	131	47409	20.521	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	192410	32.307	ng/ul	0.00
49) Acenaphthylene-d8	14.578	160	232052	32.255	ng/ul	0.00
54) 4-Nitrophenol-d4	15.066	143	4583	4.603	ng/ul	0.00
60) Fluorene-d10	15.871	176	174709	33.645	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	28181	23.892	ng/ul	0.00
73) Anthracene-d10	17.727	188	337021	37.997	ng/ul	0.00
81) Pyrene-d10	20.006	212	410471	37.805	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.182	264	397103	39.732	ng/ul	0.00

Target Compounds Qvalue

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

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