

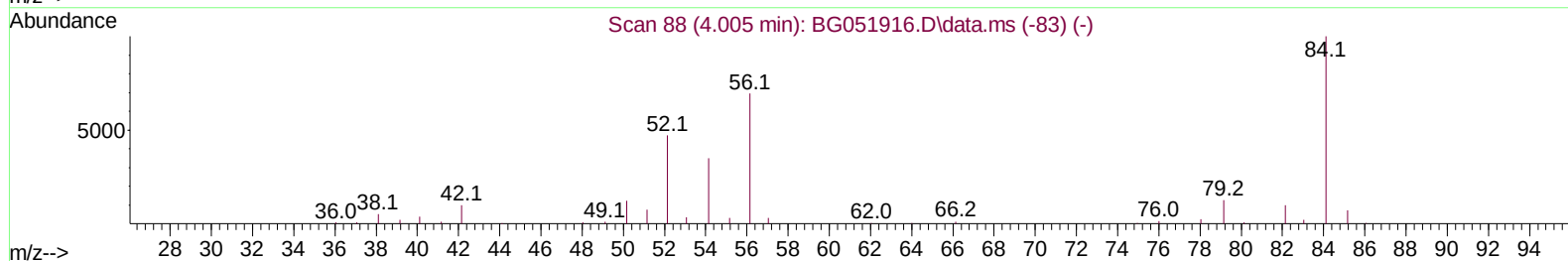
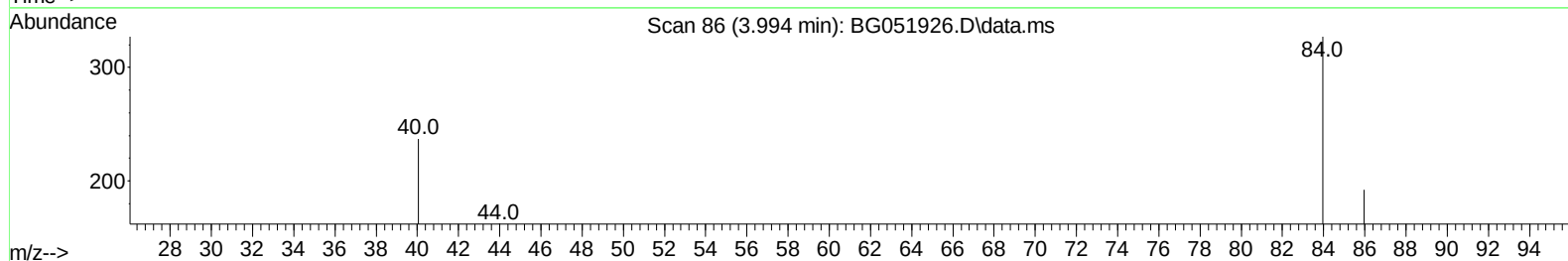
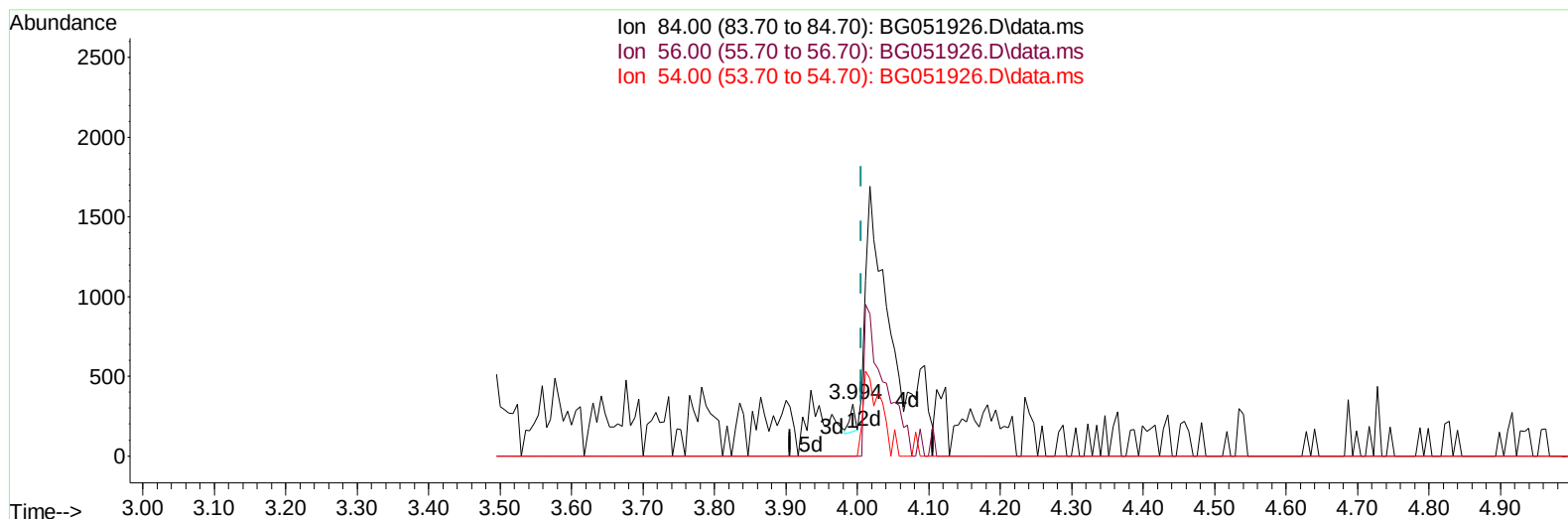
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010622\
 Data File : BG051926.D
 Acq On : 7 Jan 2022 22:13
 Operator : CG/JU
 Sample : N1026-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLG7

Manual IntegrationsAPPROVED

Quant Time: Jan 07 23:15:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/10/2022
 Supervised By :mohammad ahmed 01/12/2022



TIC: BG051926.D\data.ms

(4) Pyridine-d5 (S)

3.994min (-0.012) 0.05 ng/u1

response	89	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	0.00#
54.00	31.50	0.00#
0.00	0.00	0.00

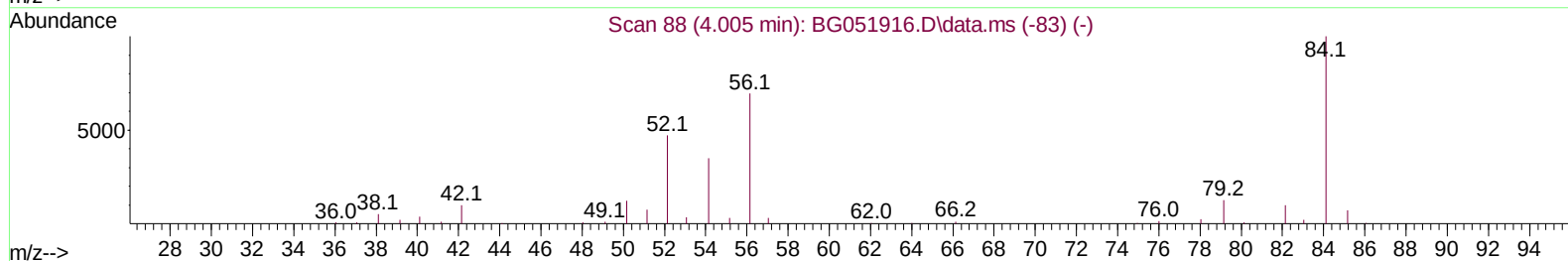
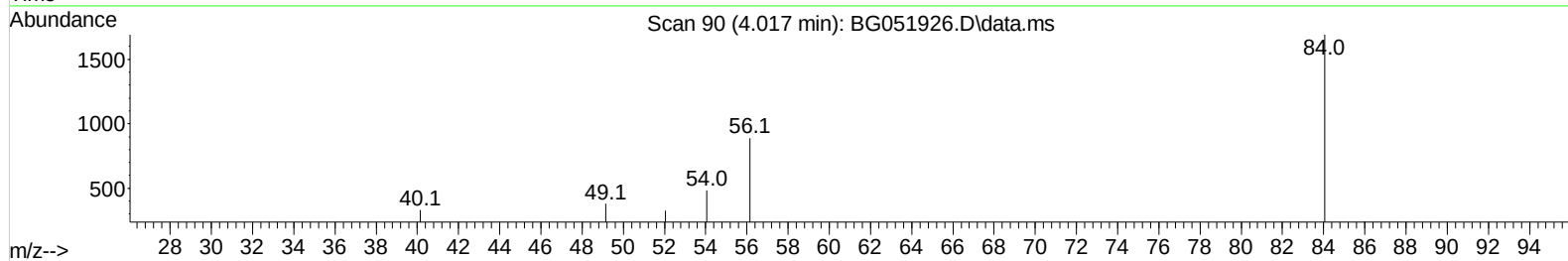
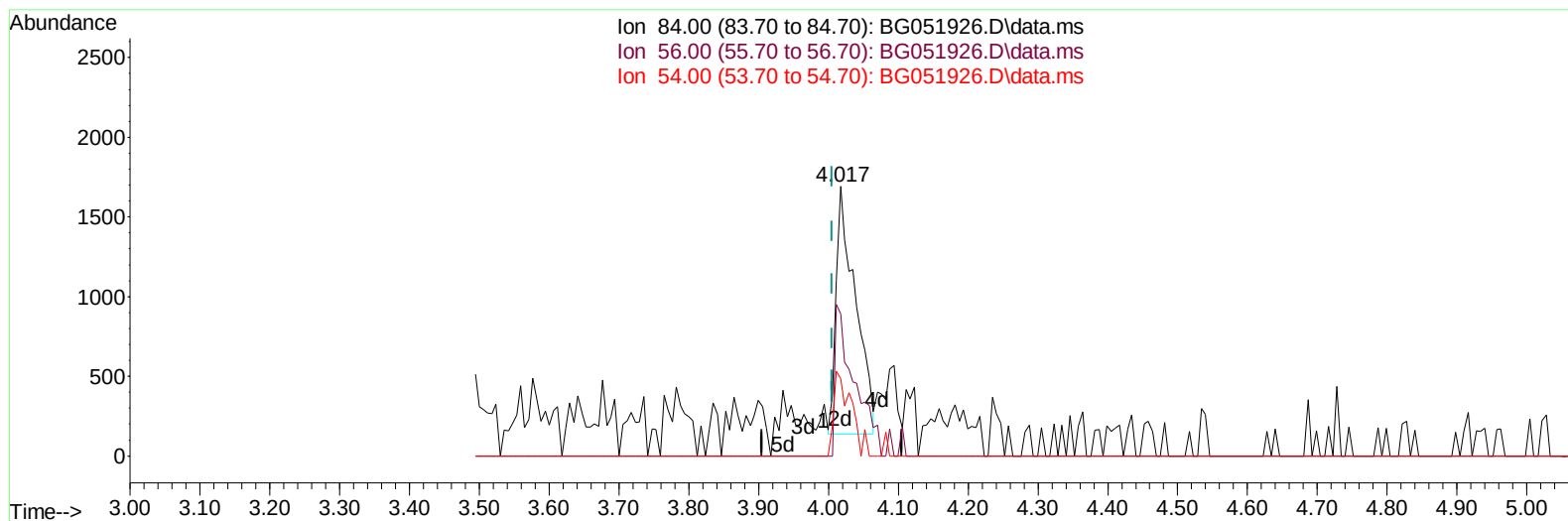
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(4) Pyridine-d5 (S)

4.017min (+ 0.012) 1.68 ng/ul m

response 2985

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	52.63#
54.00	31.50	28.74
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	8.247	152	20413	20.000	ng/ul	0.00
20) Naphthalene-d8	11.084	136	84336	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.885	164	63282	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.634	188	172906	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.952	240	200541	20.000	ng/ul	# 0.00
88) Perylene-d12	25.441	264	197161	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.571	96	4083	7.147	ng/uL	0.00
4) Pyridine-d5	4.017	84	2985m	1.681	ng/ul	0.01
7) Phenol-d5	7.383	99	34387	17.212	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.548	67	26213	20.668	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.771	132	24894	18.820	ng/ul	0.00
15) 4-Methylphenol-d8	8.946	113	27707	17.840	ng/ul	0.00
21) Nitrobenzene-d5	9.416	128	14635	21.961	ng/ul	0.00
24) 2-Nitrophenol-d4	10.150	143	14605	19.816	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.696	165	25306	17.566	ng/ul	0.00
31) 4-Chloroaniline-d4	11.207	131	33976	16.935	ng/ul	0.00
46) Dimethylphthalate-d6	14.280	166	119235	22.795	ng/ul	0.00
49) Acenaphthylene-d8	14.585	160	114818	18.172	ng/ul	0.00
54) 4-Nitrophenol-d4	15.061	143	15778	18.042	ng/ul	0.00
60) Fluorene-d10	15.878	176	91903	20.152	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.983	200	16009	14.677	ng/ul	0.00
73) Anthracene-d10	17.734	188	166173	20.259	ng/ul	0.00
81) Pyrene-d10	20.013	212	222406	19.270	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.200	264	192669	18.563	ng/ul	0.00

Target Compounds Qvalue

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

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