

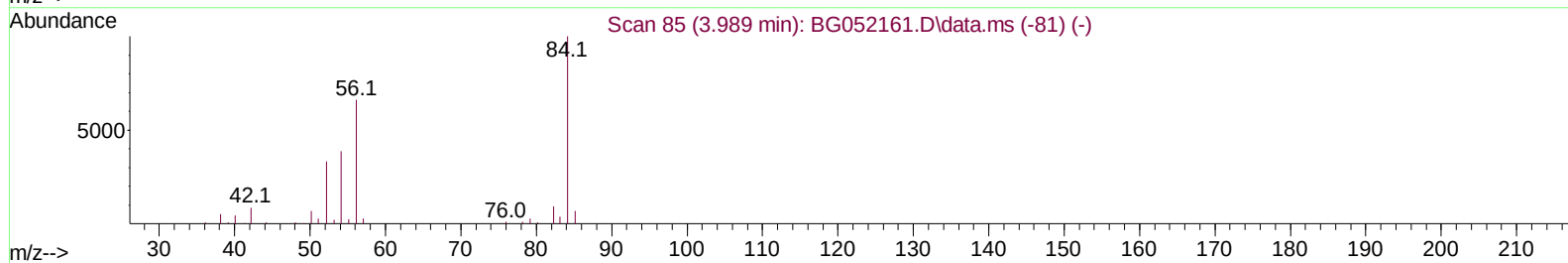
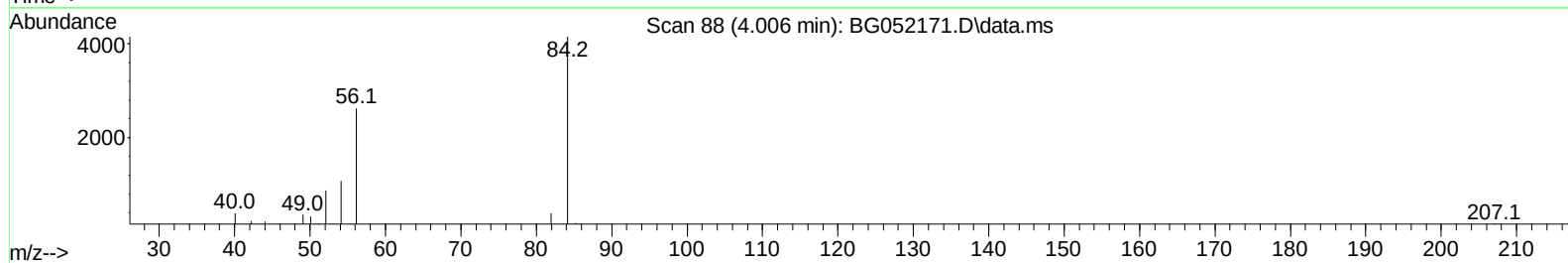
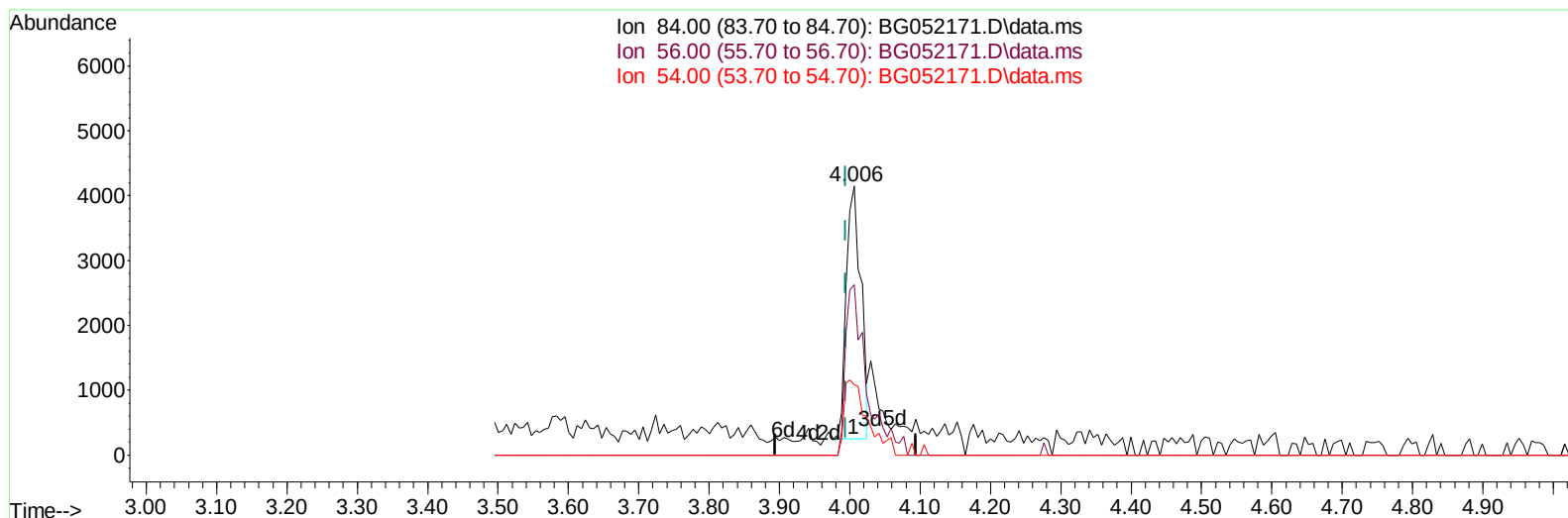
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052171.D
 Acq On : 27 Jan 2022 5:27
 Operator : CG/JU
 Sample : N1230-12
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q55

Manual IntegrationsAPPROVED

Quant Time: Jan 27 06:40:53 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 01/27/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052171.D\data.ms

(4) Pyridine-d5 (S)

4.006min (+ 0.012) 3.01 ng/u1

response 5669

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	63.28
54.00	31.50	26.01
0.00	0.00	0.00

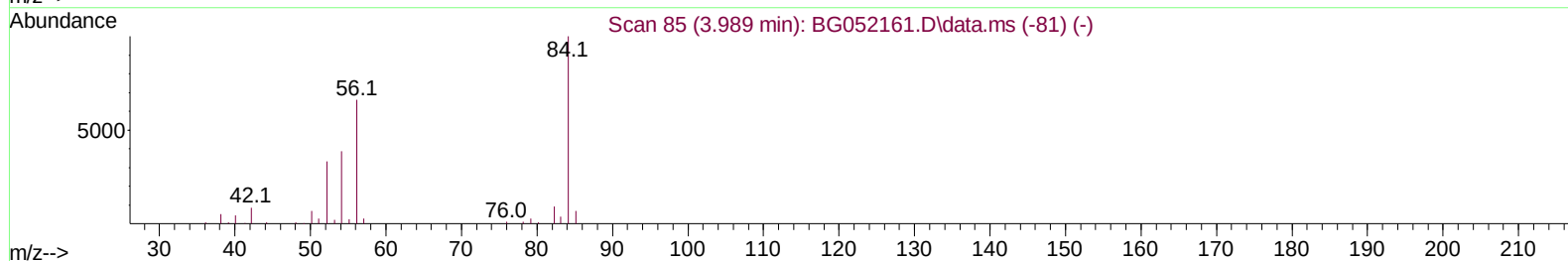
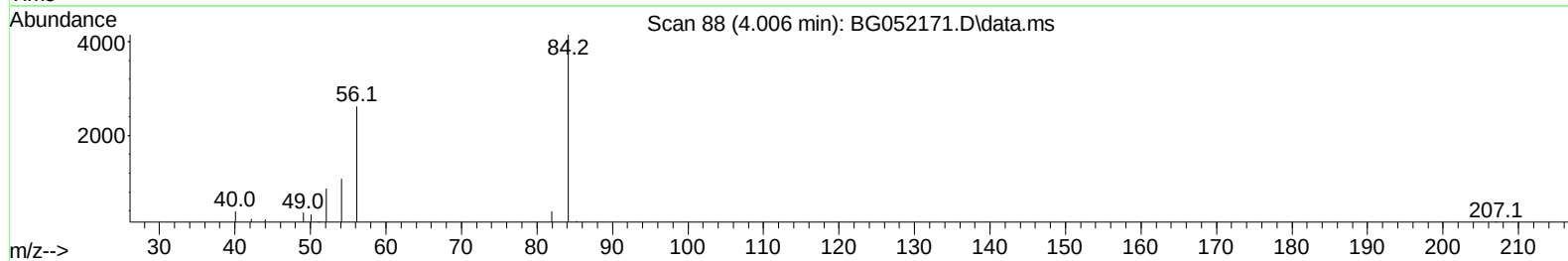
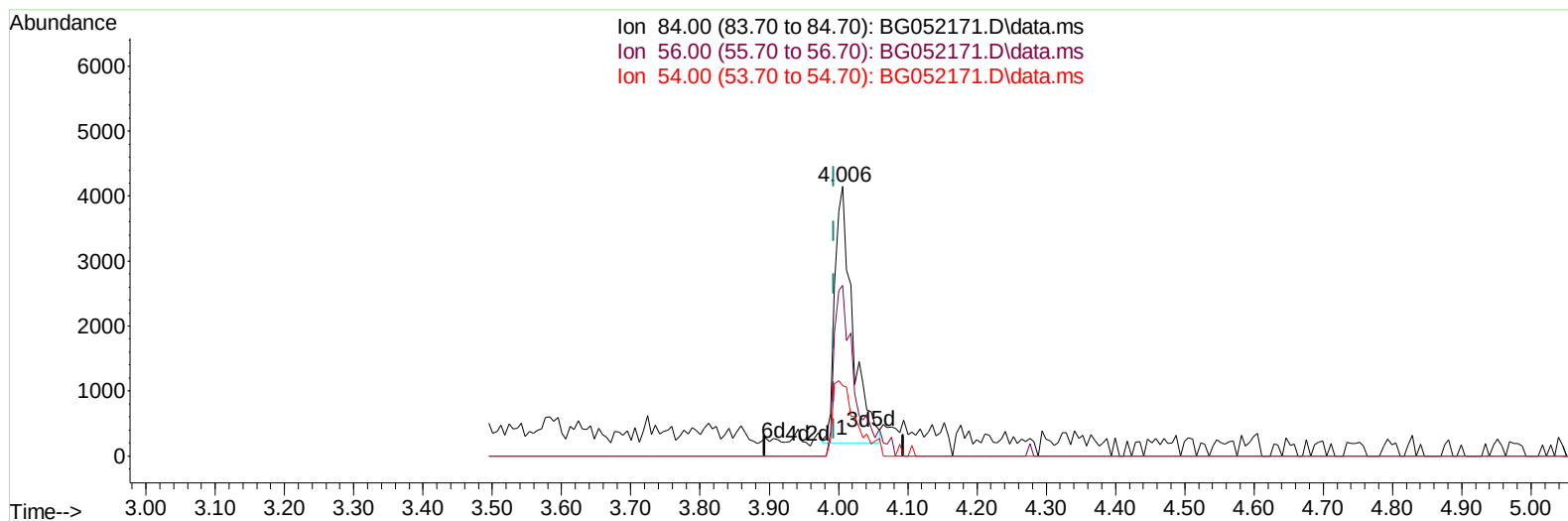
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TIC: BG052171.D\data.ms

(4) Pyridine-d5 (S)

4.006min (+ 0.012) 3.76 ng/ul m

response 7065

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	63.28
54.00	31.50	26.01
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	8.236	152	21627	20.000	ng/ul	0.00
20) Naphthalene-d8	11.079	136	91454	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.879	164	69032	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.629	188	173947	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.946	240	178868	20.000	ng/ul	# 0.00
88) Perylene-d12	25.430	264	176898	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	2432	4.018	ng/uL	0.00
4) Pyridine-d5	4.006	84	7065m	3.755	ng/ul	0.01
7) Phenol-d5	7.384	99	15902	7.513	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	41925	31.201	ng/ul	0.00
11) 2-Chlorophenol-d4	7.766	132	38923	27.774	ng/ul	0.00
15) 4-Methylphenol-d8	8.940	113	28346	17.227	ng/ul	0.00
21) Nitrobenzene-d5	9.405	128	25662	35.511	ng/ul	0.00
24) 2-Nitrophenol-d4	10.139	143	27854	34.851	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	49834	31.900	ng/ul	0.00
31) 4-Chloroaniline-d4	11.202	131	41133	18.907	ng/ul	0.00
46) Dimethylphthalate-d6	14.269	166	205195	35.961	ng/ul	0.00
49) Acenaphthylene-d8	14.574	160	244055	35.408	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	6798	7.126	ng/ul	0.00
60) Fluorene-d10	15.872	176	181080	36.398	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.978	200	30283	27.597	ng/ul	0.00
73) Anthracene-d10	17.729	188	328515	39.812	ng/ul	0.00
81) Pyrene-d10	20.008	212	405827	39.423	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.183	264	374816	40.248	ng/ul	0.00

Target Compounds Qvalue

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

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