

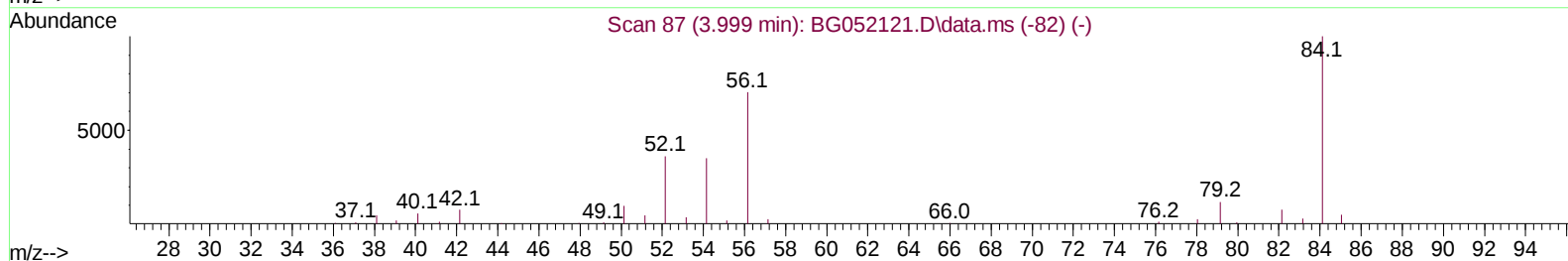
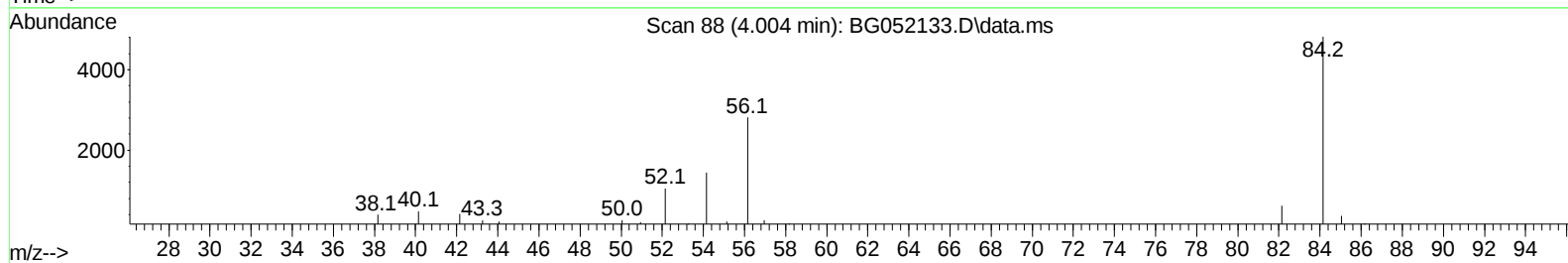
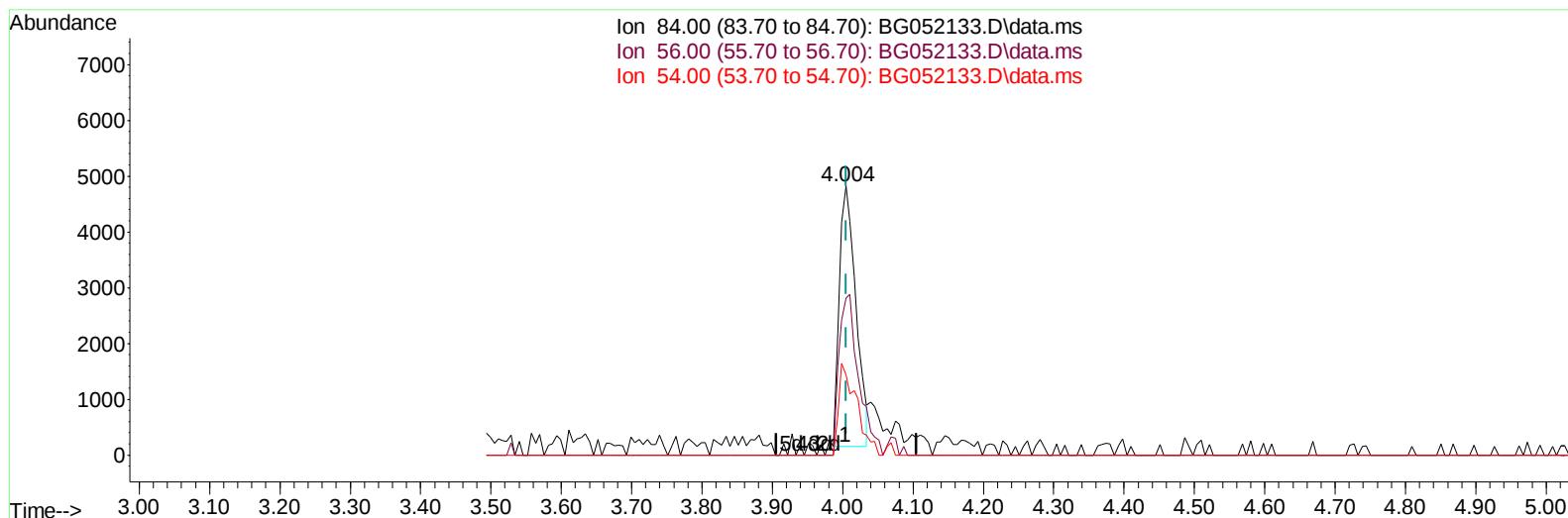
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012422\
 Data File : BG052133.D
 Acq On : 24 Jan 2022 19:22
 Operator : CG/JU
 Sample : N1199-21
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q02

Manual IntegrationsAPPROVED

Quant Time: Jan 25 00:00:58 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 01/25/2022
 Supervised By :mohammad ahmed 01/31/2022



TIC: BG052133.D\data.ms

(4) Pyridine-d5 (S)

4.004min (-0.001) 3.15 ng/u1

response 7651

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	58.41
54.00	31.50	29.81
0.00	0.00	0.00

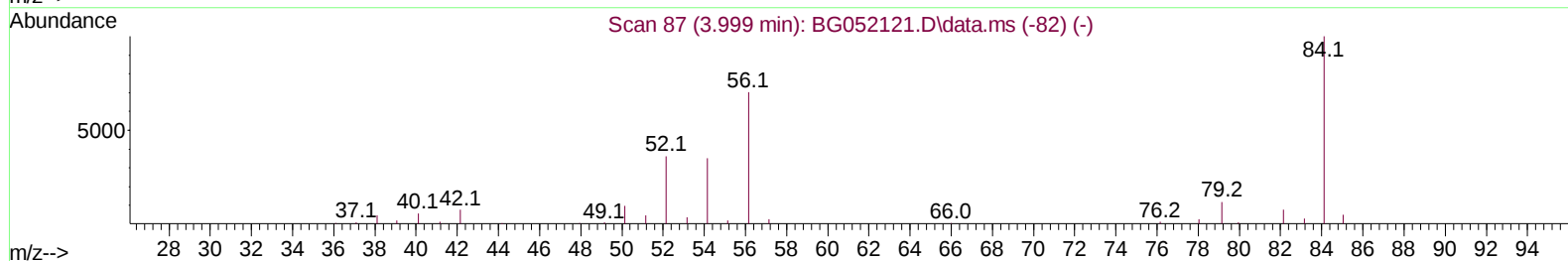
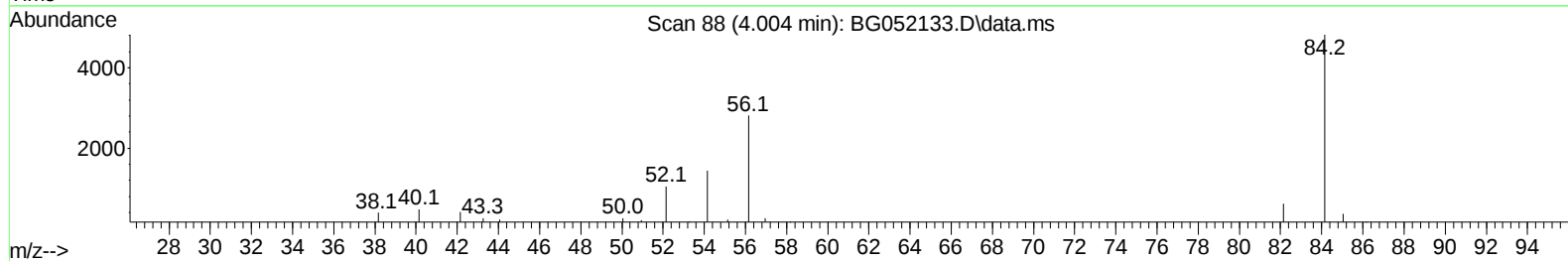
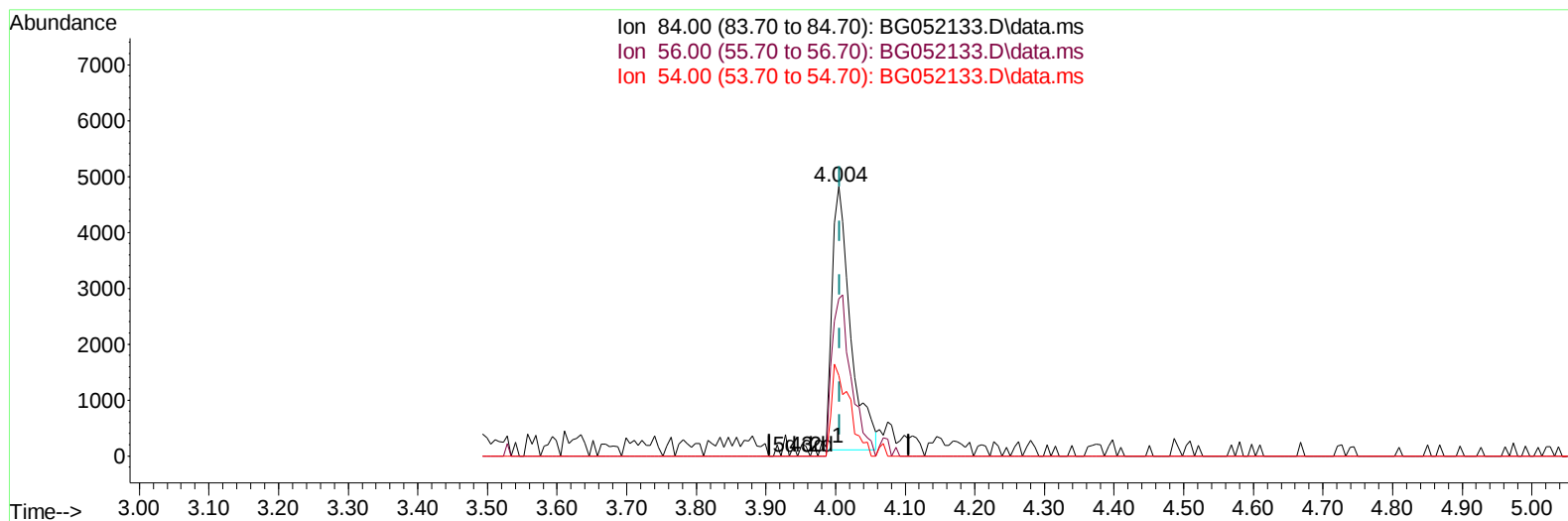
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(4) Pyridine-d5 (S)

4.004min (-0.001) 3.56 ng/ul m

response 8643

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	58.41
54.00	31.50	29.81
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	8.240	152	27904	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.083	136	112860	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.890	164	83073	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.639	188	186006	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.956	240	180663	20.000	ng/ul	# 0.00
88) Perylene-d12	25.446	264	178044	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.570	96	2898	3.711	ng/uL	0.00
4) Pyridine-d5	4.004	84	8643m	3.560	ng/ul	0.00
7) Phenol-d5	7.388	99	18225	6.673	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	44165	25.475	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.770	132	43165	23.872	ng/ul	0.00
15) 4-Methylphenol-d8	8.945	113	32249	15.190	ng/ul	0.00
21) Nitrobenzene-d5	9.415	128	27968	31.362	ng/ul	0.00
24) 2-Nitrophenol-d4	10.149	143	31243	31.677	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.695	165	60031	31.139	ng/ul	0.00
31) 4-Chloroaniline-d4	11.206	131	40538	15.099	ng/ul	0.00
46) Dimethylphthalate-d6	14.279	166	217139	31.622	ng/ul	0.00
49) Acenaphthylene-d8	14.584	160	272434	32.845	ng/ul	0.00
54) 4-Nitrophenol-d4	15.072	143	7308	6.366	ng/ul	0.00
60) Fluorene-d10	15.876	176	195120	32.591	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.982	200	32480	27.680	ng/ul	0.00
73) Anthracene-d10	17.739	188	334946	37.960	ng/ul	0.00
81) Pyrene-d10	20.018	212	389926	37.502	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.205	264	358316	38.228	ng/ul	0.01

Target Compounds Qvalue

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

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