

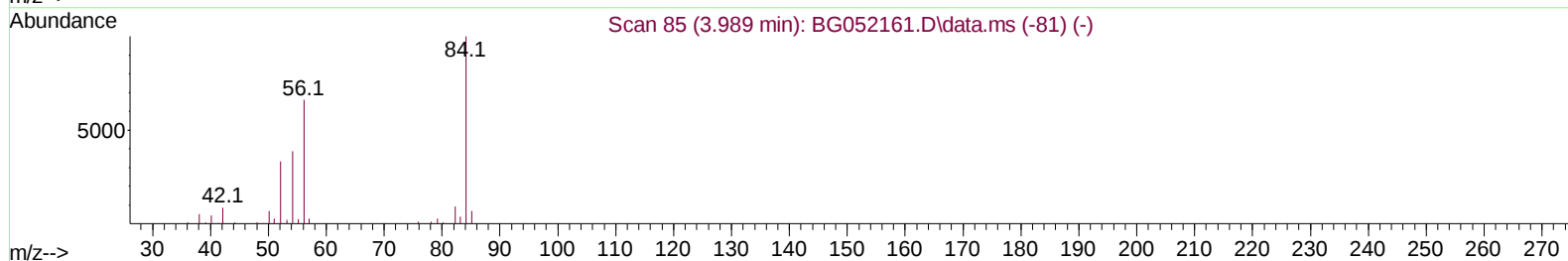
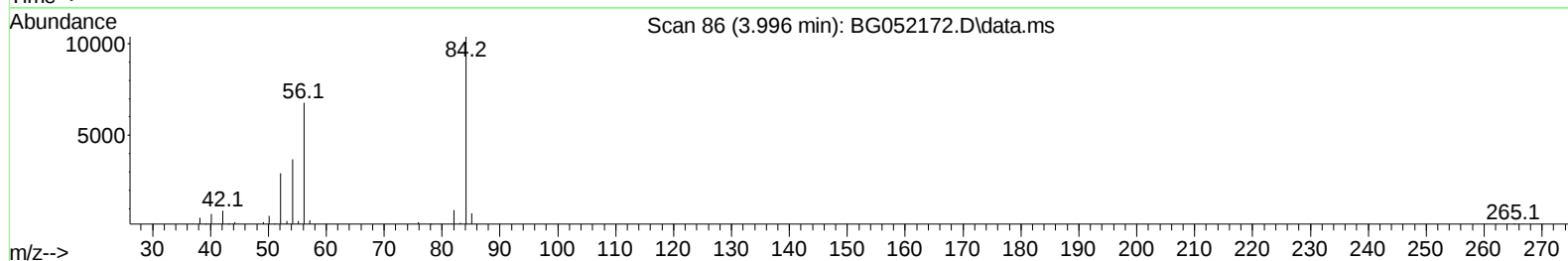
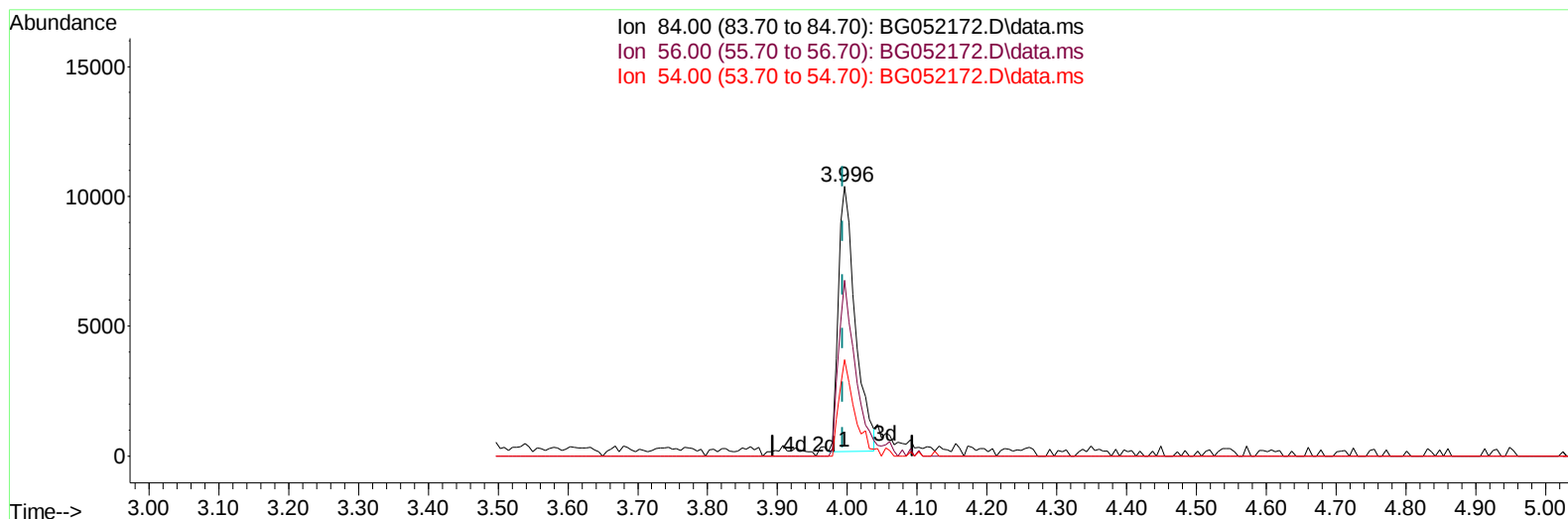
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052172.D
 Acq On : 27 Jan 2022 6:04
 Operator : CG/JU
 Sample : N1230-19
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q21

Manual IntegrationsAPPROVED

Quant Time: Jan 27 06:41:05 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: BG052172.D\data.ms

(4) Pyridine-d5 (S)

3.996min (+ 0.002) 8.92 ng/u1

response 17114

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	65.15
54.00	31.50	35.66
0.00	0.00	0.00

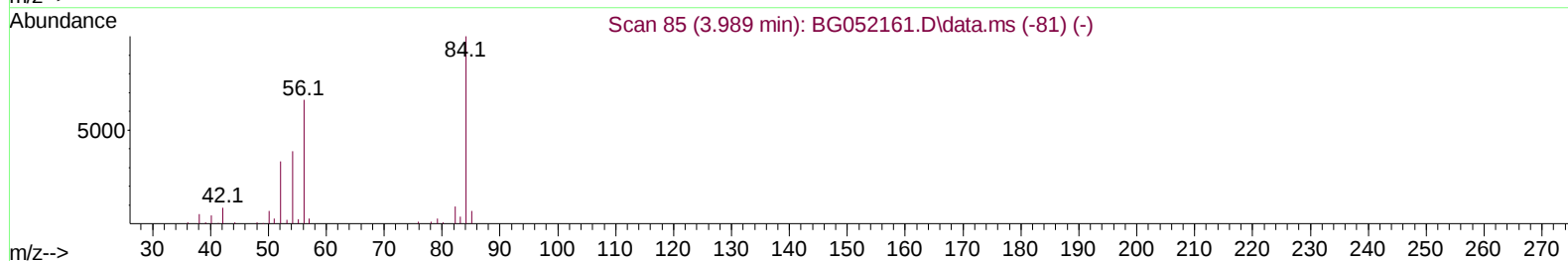
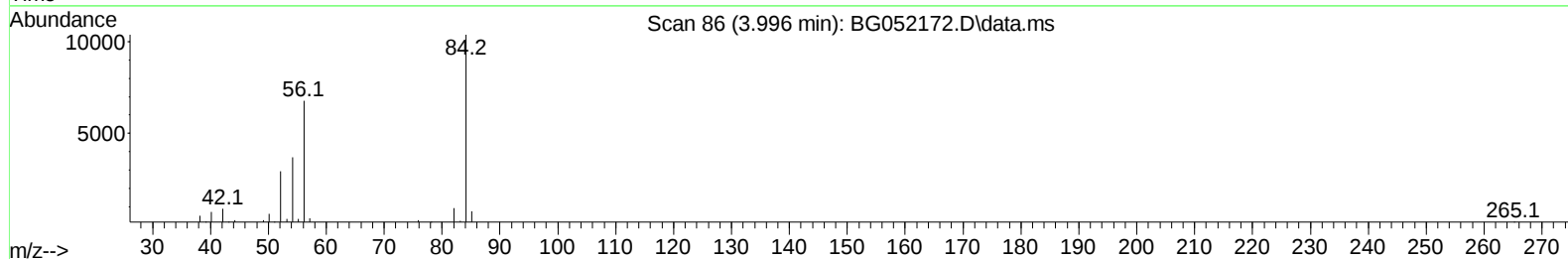
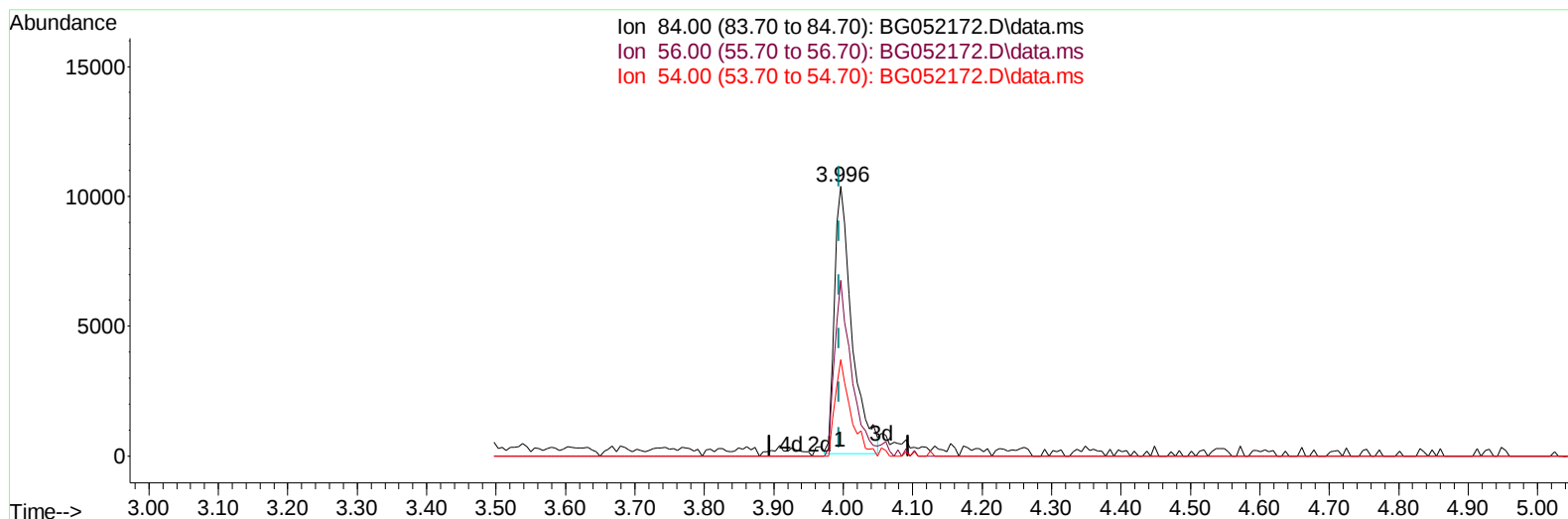
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TIC: BG052172.D\data.ms

(4) Pyridine-d5 (S)

3.996min (+ 0.002) 9.42 ng/u1 m

response 18084

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	65.15
54.00	31.50	35.66
0.00	0.00	0.00

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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	8.238	152	22057	20.000	ng/ul	0.00
20) Naphthalene-d8	11.075	136	92733	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.882	164	69610	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.631	188	171970	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.943	240	185270	20.000	ng/ul	# 0.00
88) Perylene-d12	25.420	264	180787	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.568	96	2762	4.474	ng/uL	0.00
4) Pyridine-d5	3.996	84	18084m	9.424	ng/ul	0.00
7) Phenol-d5	7.380	99	19259	8.921	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.539	67	43385	31.658	ng/ul	0.00
11) 2-Chlorophenol-d4	7.762	132	42077	29.439	ng/ul	0.00
15) 4-Methylphenol-d8	8.943	113	32812	19.552	ng/ul	0.00
21) Nitrobenzene-d5	9.407	128	26509	36.177	ng/ul	0.00
24) 2-Nitrophenol-d4	10.141	143	29554	36.468	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.687	165	51833	32.722	ng/ul	0.00
31) 4-Chloroaniline-d4	11.198	131	54517	24.713	ng/ul	0.00
46) Dimethylphthalate-d6	14.271	166	208328	36.207	ng/ul	0.00
49) Acenaphthylene-d8	14.576	160	244300	35.150	ng/ul	0.00
54) 4-Nitrophenol-d4	15.064	143	8432	8.765	ng/ul	0.00
60) Fluorene-d10	15.869	176	184488	36.775	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.980	200	28812	26.558	ng/ul	0.00
73) Anthracene-d10	17.731	188	322814	39.571	ng/ul	0.00
81) Pyrene-d10	20.010	212	403974	37.887	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.185	264	383504	40.295	ng/ul	0.00

Target Compounds Qvalue

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

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