

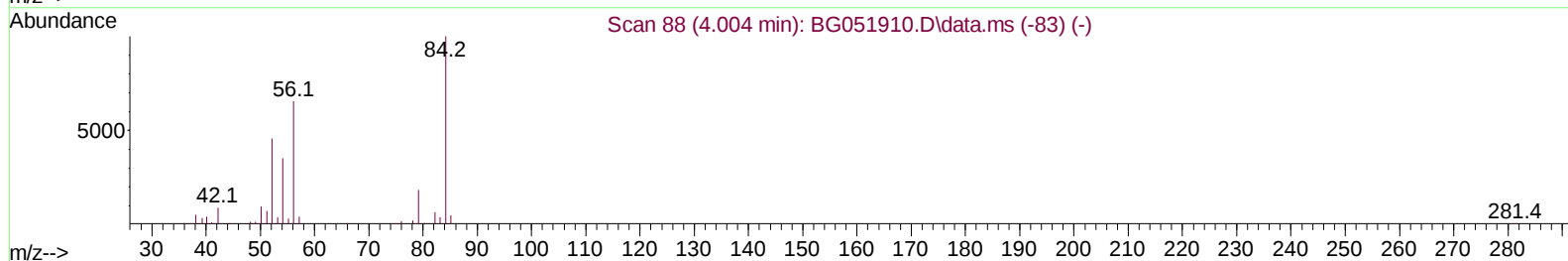
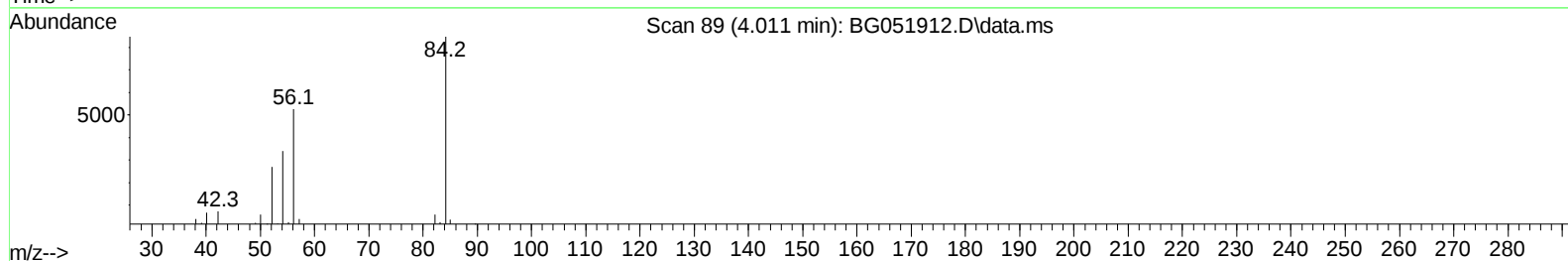
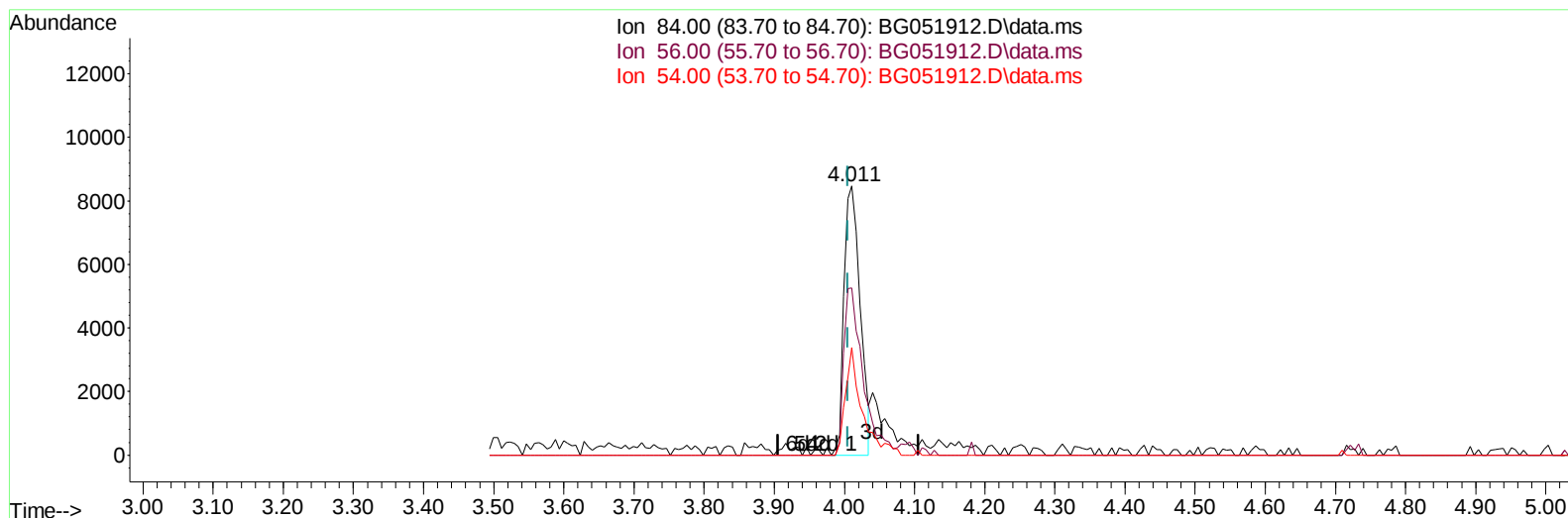
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010622\
Data File : BG051912.D
Acq On : 7 Jan 2022 13:00
Operator : CG/JU
Sample : N1027-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0J34

Manual IntegrationsAPPROVED

Quant Time: Jan 07 22:52:54 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: BG051912.D\data.ms

(4) Pyridine-d5 (S)

4.011min (+ 0.005) 6.72 ng/u1

response 13574

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	62.11
54.00	31.50	40.08#
0.00	0.00	0.00

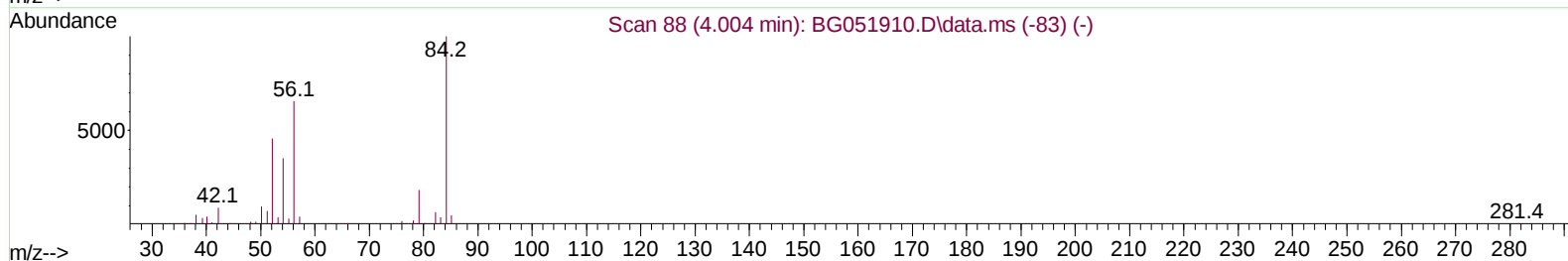
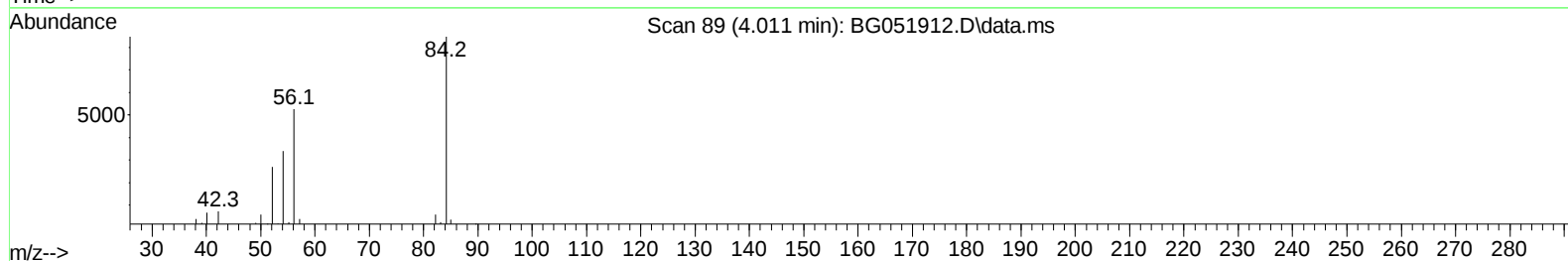
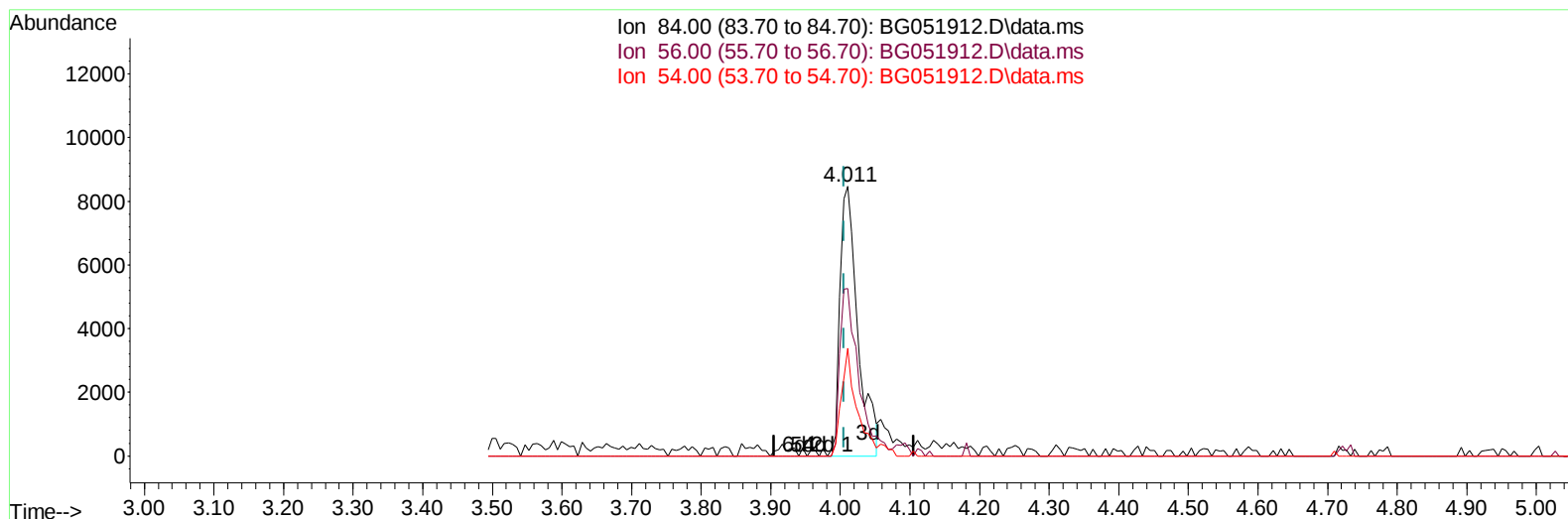
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TIC: BG051912.D\data.ms

(4) Pyridine-d5 (S)

4.011min (+ 0.005) 7.53 ng/u1 m

response 15202

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	62.11
54.00	31.50	40.08#
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	8.246	152	23220	20.000	ng/ul	0.00
20) Naphthalene-d8	11.089	136	98568	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.890	164	71763	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.639	188	185275	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.957	240	195159	20.000	ng/ul	# 0.00
88) Perylene-d12	25.446	264	197963	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.576	96	3859	5.939	ng/uL	0.00
4) Pyridine-d5	4.011	84	15202m	7.526	ng/ul	0.00
7) Phenol-d5	7.383	99	13998	6.159	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.553	67	46970	32.558	ng/ul	0.00
11) 2-Chlorophenol-d4	7.776	132	37454	24.892	ng/ul	0.00
15) 4-Methylphenol-d8	8.951	113	24754	14.012	ng/ul	0.00
21) Nitrobenzene-d5	9.421	128	26707	34.290	ng/ul	0.00
24) 2-Nitrophenol-d4	10.155	143	6752	7.838	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.696	165	46538	27.640	ng/ul	0.00
31) 4-Chloroaniline-d4	11.213	131	64036	27.310	ng/ul	0.00
46) Dimethylphthalate-d6	14.279	166	210693	35.519	ng/ul	0.00
49) Acenaphthylene-d8	14.585	160	244682	34.148	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	15.883	176	182324	35.254	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/ul	
73) Anthracene-d10	17.739	188	331220	37.686	ng/ul	0.00
81) Pyrene-d10	20.012	212	420490	37.438	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.205	264	377698	36.242	ng/ul	0.01

Target Compounds Qvalue

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

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