

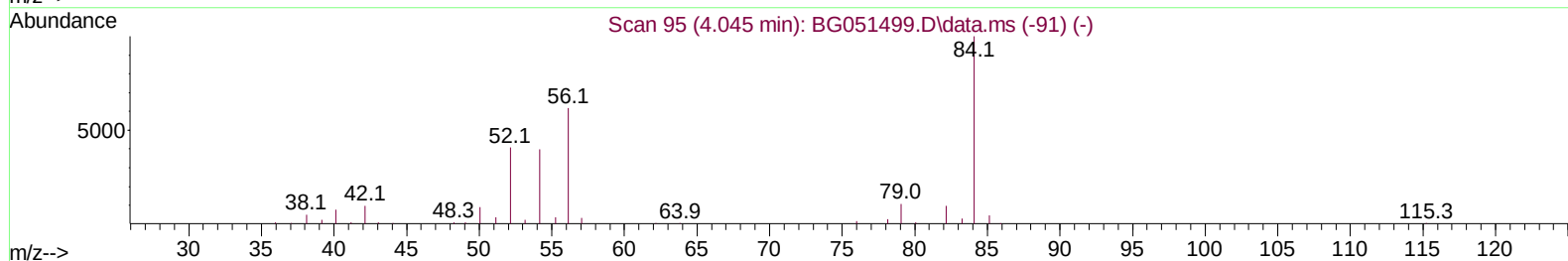
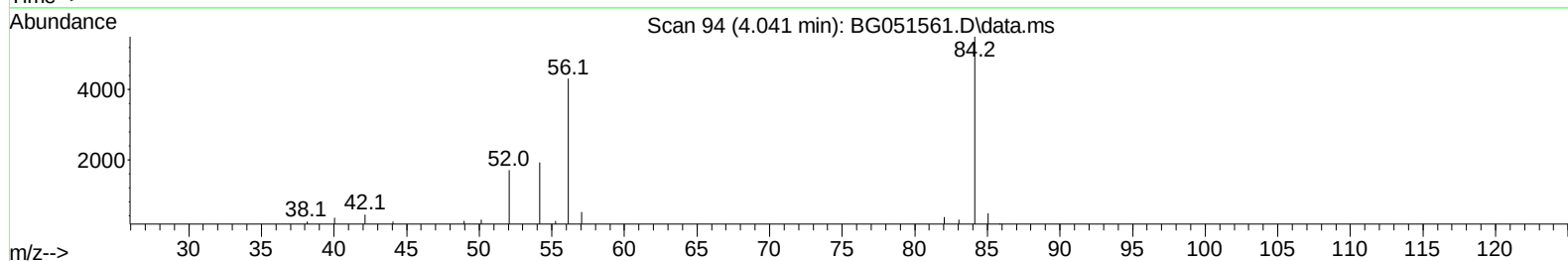
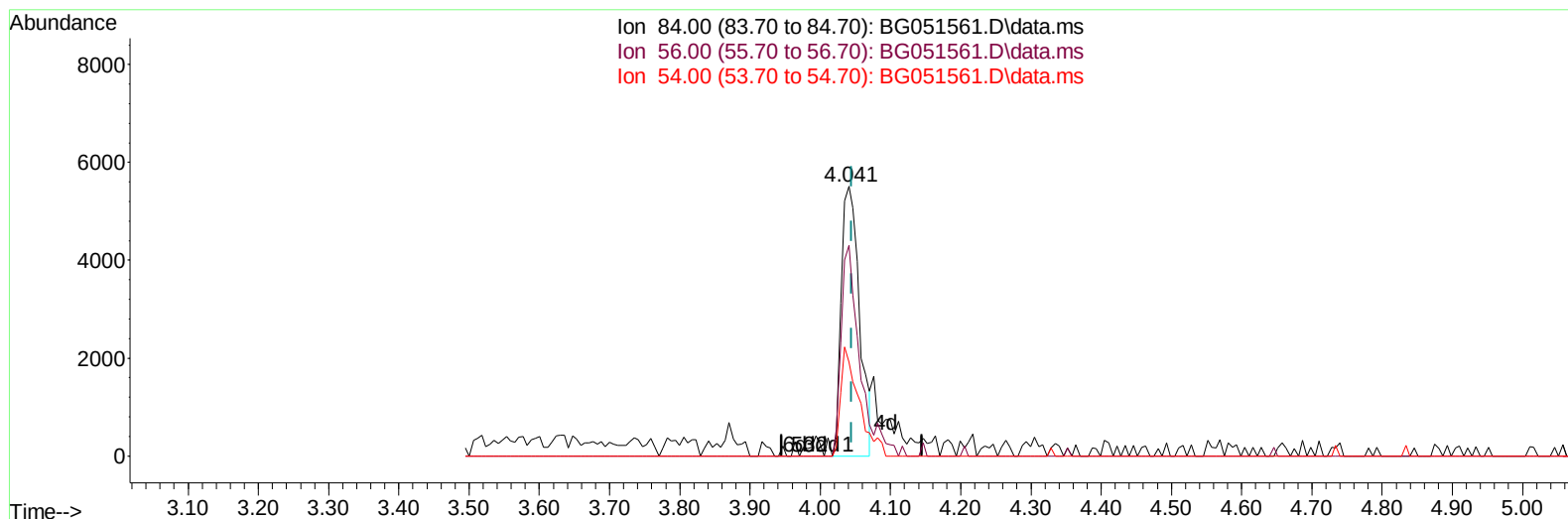
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121421\
Data File : BG051561.D
Acq On : 16 Dec 2021 7:57
Operator : CG/JU
Sample : M4971-13
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BG3D0

Manual IntegrationsAPPROVED

Quant Time: Dec 16 08:41:02 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 14 14:19:57 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/16/2021
Supervised By :Yogesh Patel 12/23/2021



TIC: BG051561.D\data.ms

(4) Pyridine-d5 (S)

4.041min (-0.005) 4.38 ng/u1

response 9967

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	78.21
54.00	31.50	34.84
0.00	0.00	0.00

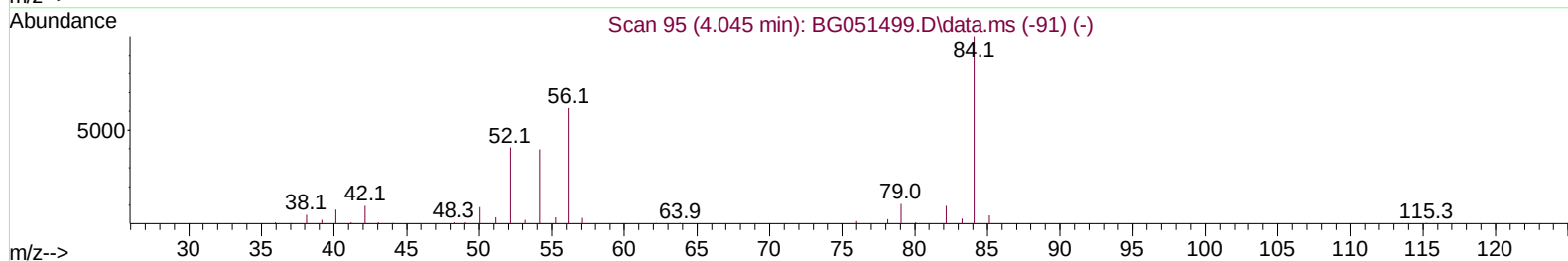
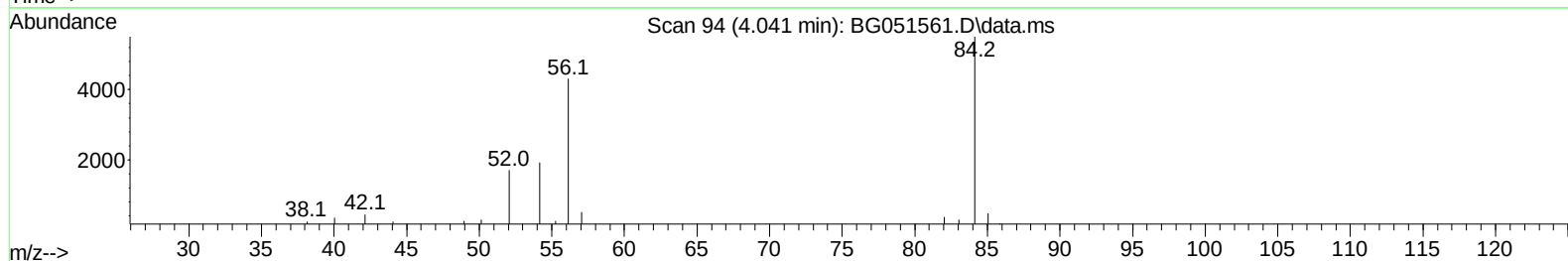
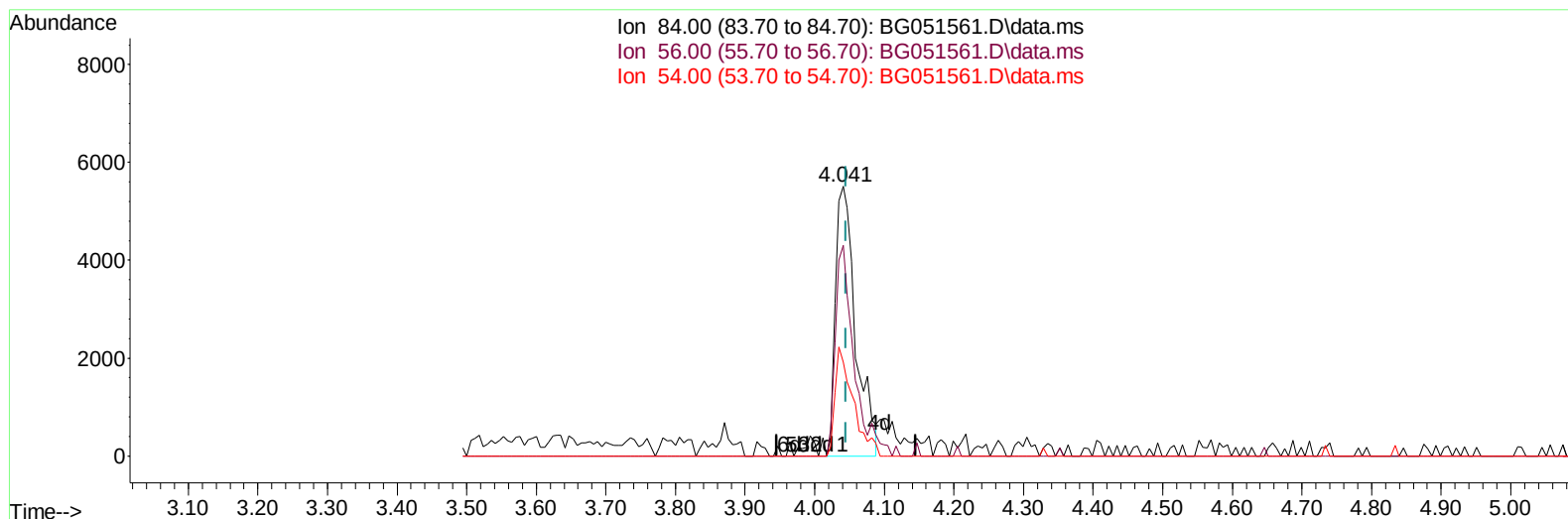
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(4) Pyridine-d5 (S)

4.041min (-0.005) 4.84 ng/ul m

response 11027

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	78.21
54.00	31.50	34.84
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	8.276	152	31305	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.113	136	128723	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.914	164	93209	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.658	188	225928	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.975	240	231922	20.000	ng/ul	#-0.01
88) Perylene-d12	25.471	264	230842	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.606	96	4126	5.474	ng/uL	0.00
4) Pyridine-d5	4.041	84	11027m	4.842	ng/ul	0.00
7) Phenol-d5	7.407	99	22727	8.131	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.583	67	55259	33.256	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.800	132	53846	27.699	ng/ul	-0.01
15) 4-Methylphenol-d8	8.969	113	40053	18.538	ng/ul	-0.01
21) Nitrobenzene-d5	9.445	128	32692	34.641	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.179	143	35130	34.051	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.720	165	68020	30.403	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.237	131	56208	19.082	ng/ul	0.00
46) Dimethylphthalate-d6	14.303	166	261187	35.010	ng/ul	-0.01
49) Acenaphthylene-d8	14.609	160	308176	33.261	ng/ul	0.00
54) 4-Nitrophenol-d4	15.067	143	10683	8.806	ng/ul	-0.01
60) Fluorene-d10	15.901	176	229801	34.381	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.995	200	45143	38.351	ng/ul	-0.02
73) Anthracene-d10	17.757	188	408167	37.872	ng/ul	0.00
81) Pyrene-d10	20.031	212	503066	36.041	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.236	264	470599	38.745	ng/ul	-0.02

Target Compounds Qvalue

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

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