

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
Data File : BG051654.D
Acq On : 19 Dec 2021 8:21
Operator : CG/JU
Sample : M5154-07
Misc :
ALS Vial : 71 Sample Multiplier: 1

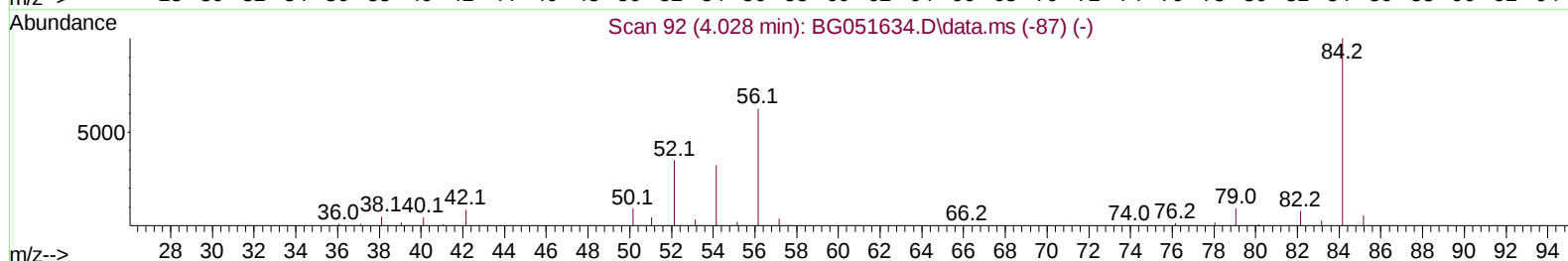
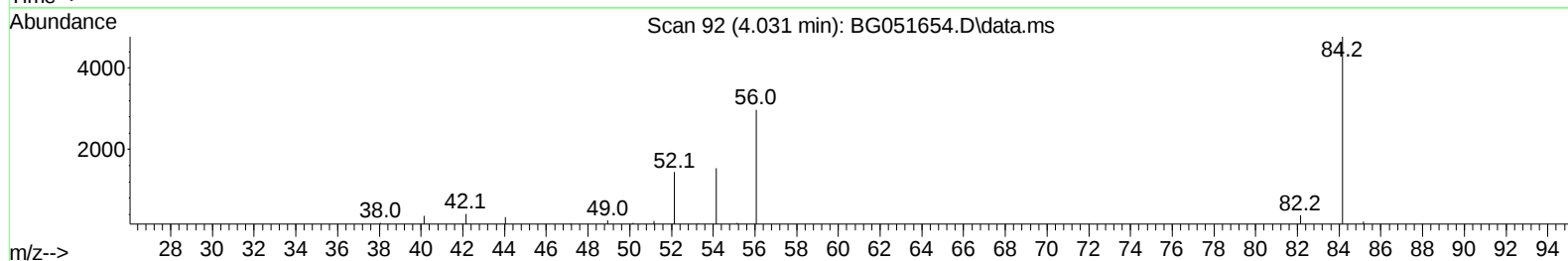
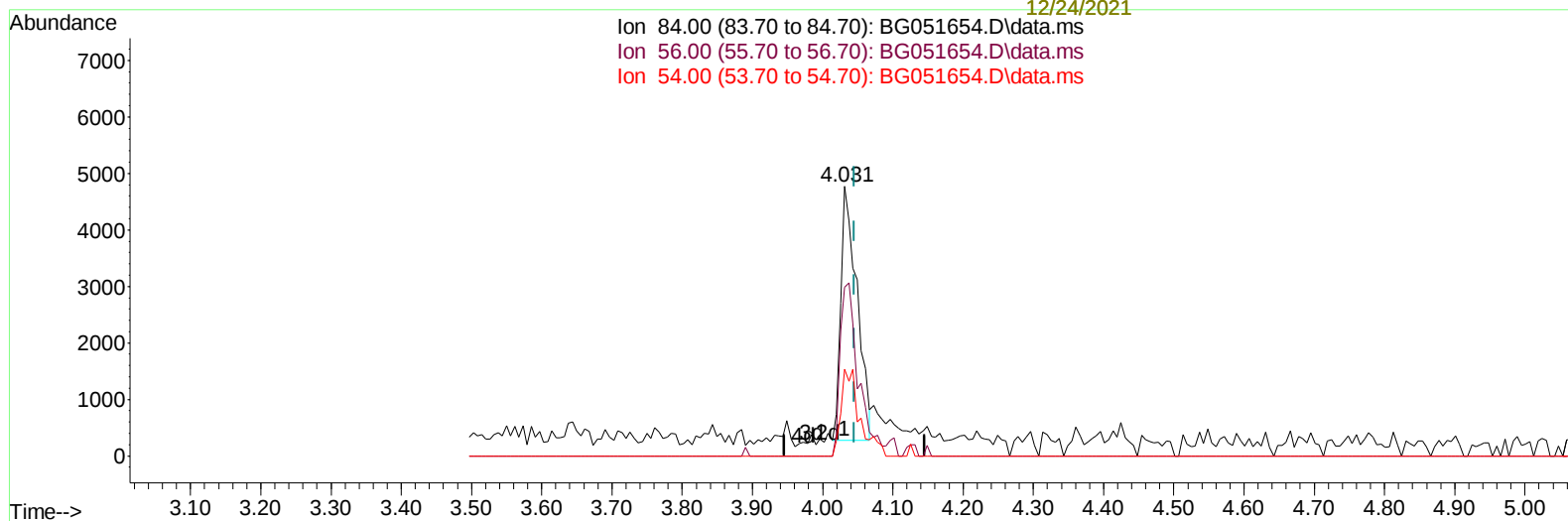
Instrument :
BNA_G
ClientSampleId :
BGL87

Manual IntegrationsAPPROVED

Quant Time: Dec 20 00:10:18 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
12/20/2021

Supervised By :mohammad
ahmed
12/24/2021



TIC: BG051654.D\data.ms

(4) Pyridine-d5 (S)

4.031min (-0.014) 2.18 ng/u1

response 7229

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	62.51
54.00	31.50	32.12
0.00	0.00	0.00

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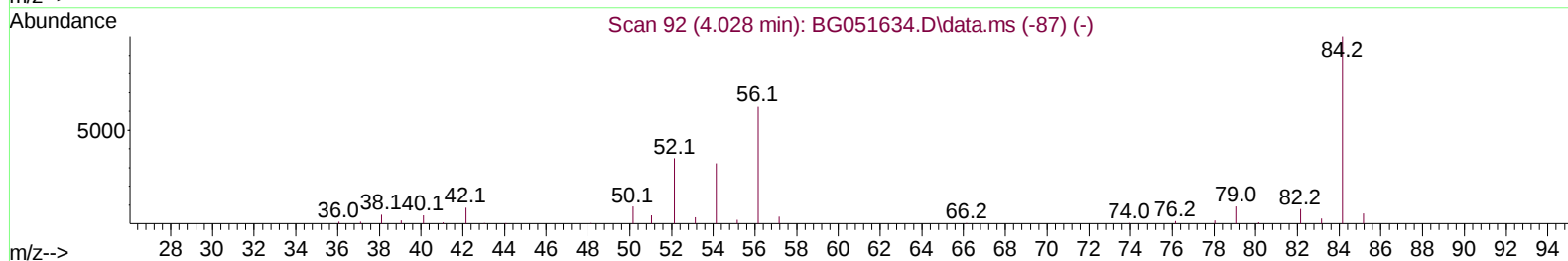
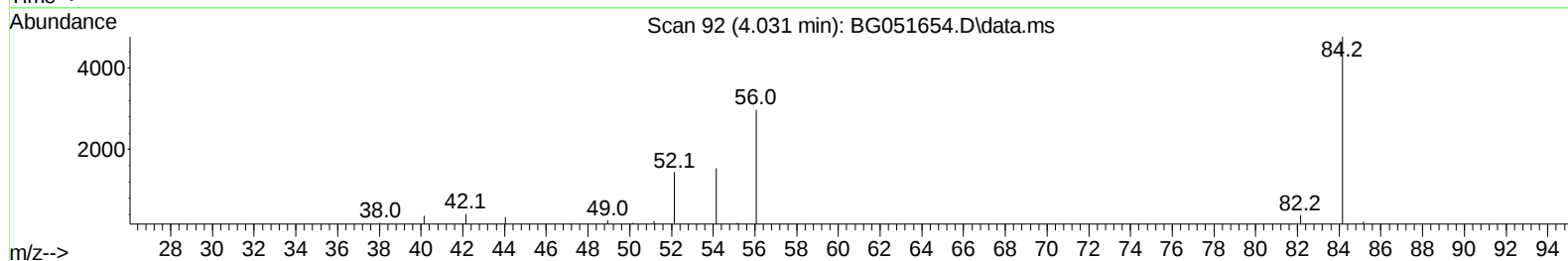
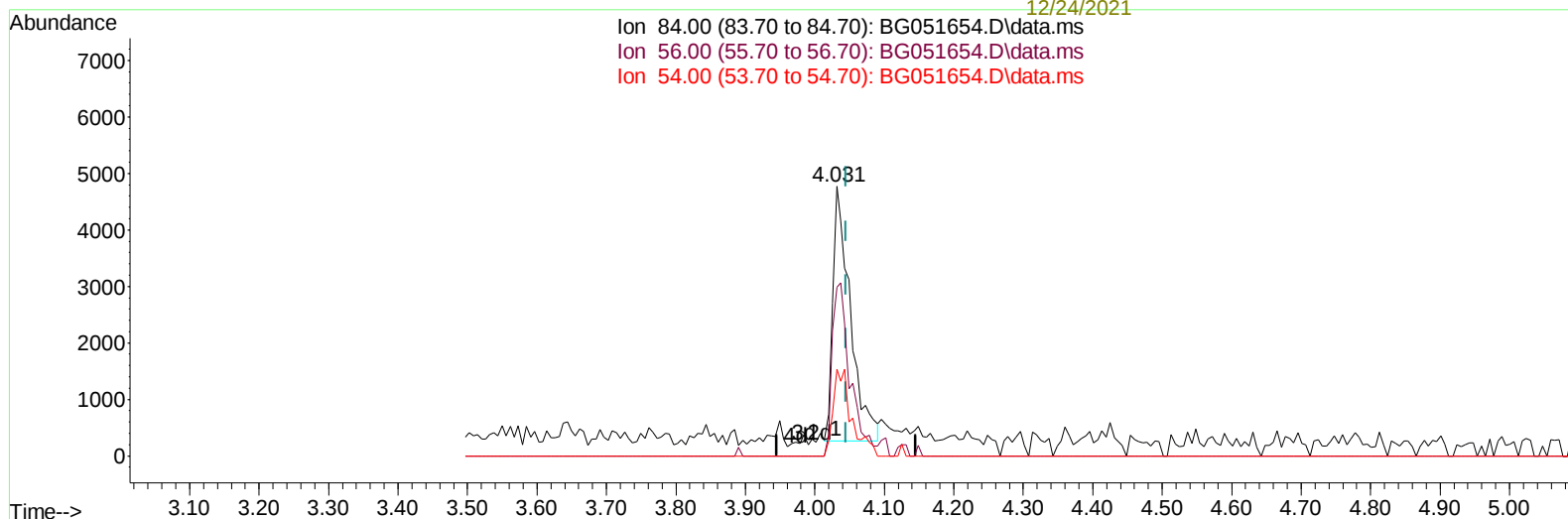
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TIC: BG051654.D\data.ms

(4) Pyridine-d5 (S)

4.031min (-0.014) 2.39 ng/ul m

response 7919

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	68.00	62.51
54.00	31.50	32.12
0.00	0.00	0.00

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Quant Time: Dec 10 00:10:18 2021
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.273	152		45618	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.110	136		190412	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.905	164		137247	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.654	188		310837	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.972	240		299576	20.000	ng/ul	#-0.01
88) Perylene-d12	25.467	264		305243	20.000	ng/ul	-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.597	96		4518	4.113	ng/uL	0.00
4) Pyridine-d5	4.031	84		7919m	2.386	ng/ul	-0.01
7) Phenol-d5	7.403	99		96369	23.660	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.580	67		57639	23.804	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.797	132		68439	24.159	ng/ul	-0.01
15) 4-Methylphenol-d8	8.966	113		76224	24.210	ng/ul	-0.01
21) Nitrobenzene-d5	9.442	128		35739	25.601	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.176	143		38939	25.516	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.716	165		78881	23.835	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.228	131		87540	20.091	ng/ul	-0.01
46) Dimethylphthalate-d6	14.300	166		251185	22.866	ng/ul	-0.01
49) Acenaphthylene-d8	14.605	160		320454	23.488	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143		33862	18.955	ng/ul	0.00
60) Fluorene-d10	15.898	176		225393	22.901	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.998	200		27537	17.004	ng/ul	-0.01
73) Anthracene-d10	17.754	188		354140	23.883	ng/ul	0.00
81) Pyrene-d10	20.027	212		401305	22.258	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.232	264		382535	23.818	ng/ul	-0.02

Target Compounds Qvalue

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

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