

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
 Data File : BG056104.D
 Acq On : 25 Dec 2022 1:15
 Operator : CG/JU
 Sample : N0017-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBYF2

Manual IntegrationsAPPROVED

Quant Time: Dec 25 01:52:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG122422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 24 22:54:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut

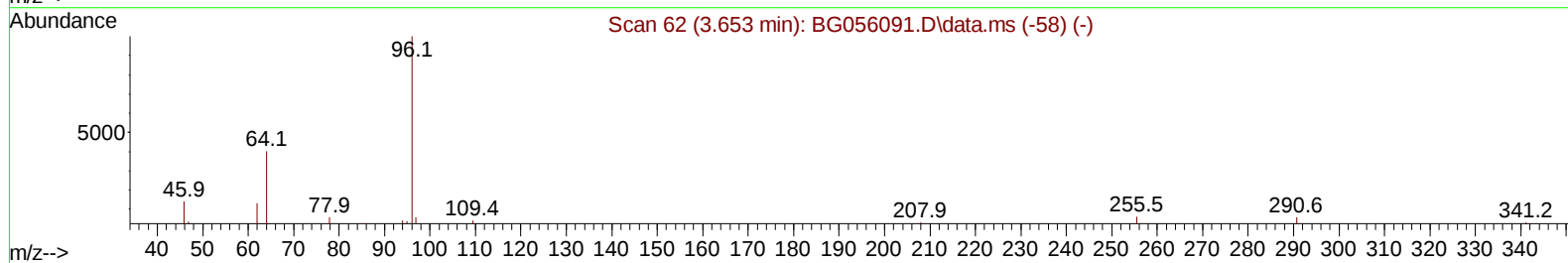
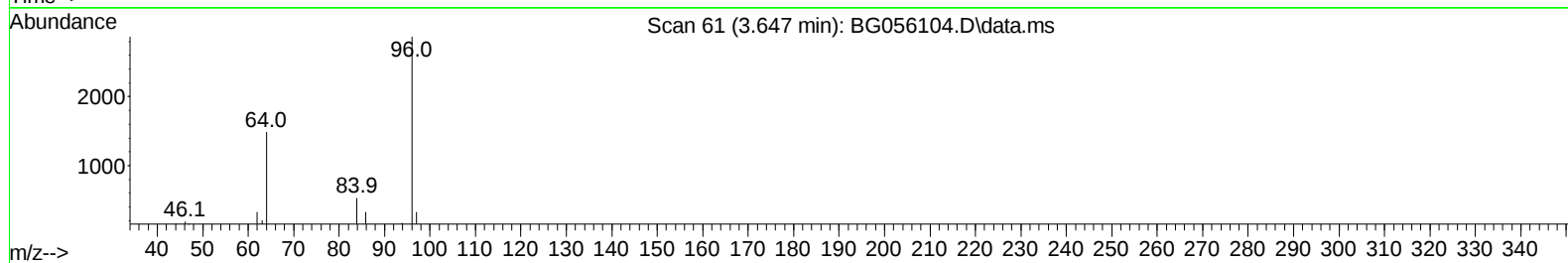
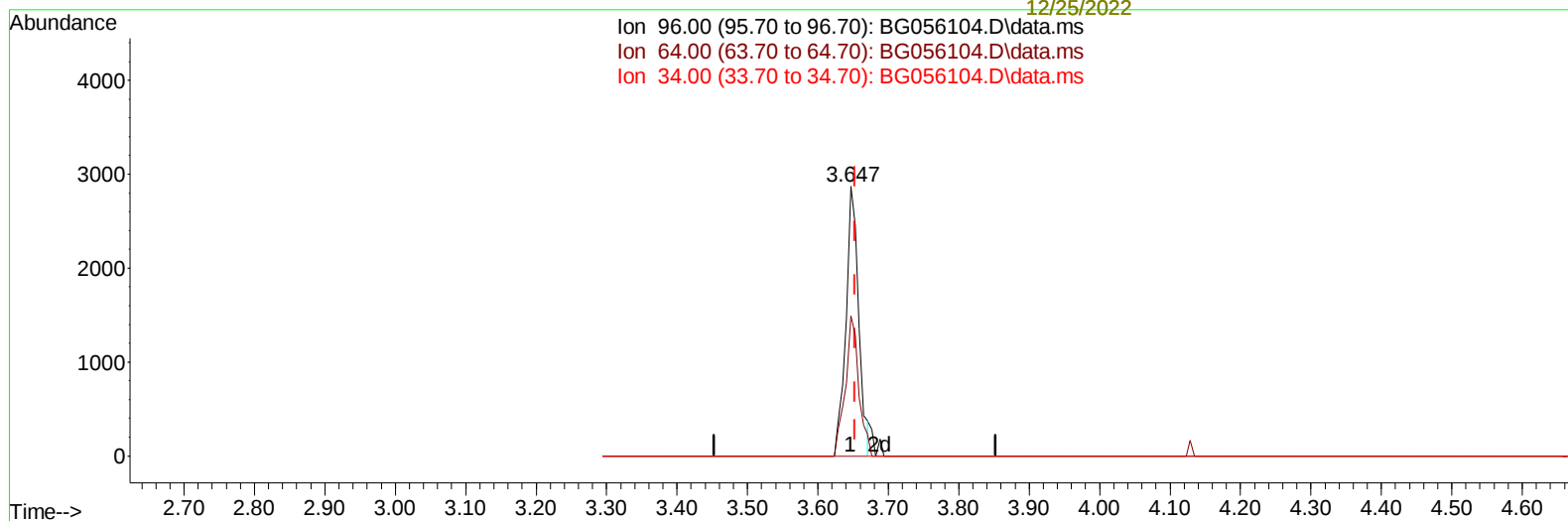
Upadhyay

12/25/2022

Supervised By :Yogesh

Patel

12/25/2022



TIC: BG056104.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.647min (-0.006) 3.80 ng/uL

response 3555

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.20	52.00#
34.00	0.00	0.00
0.00	0.00	0.00

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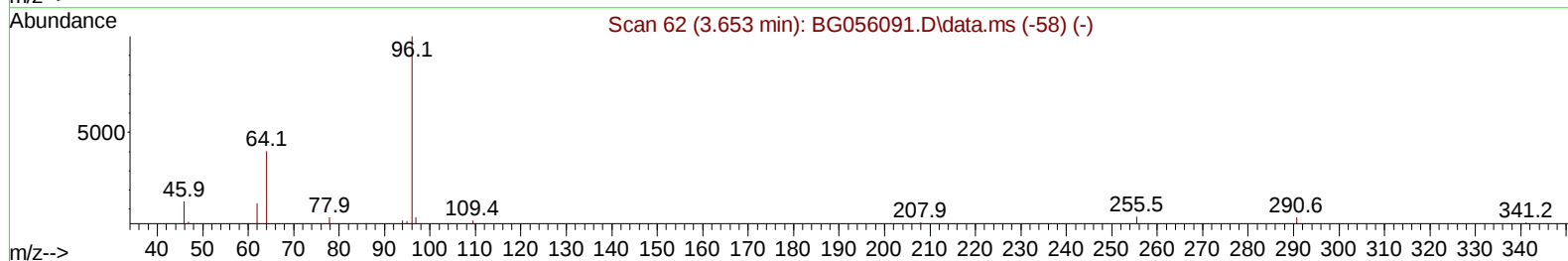
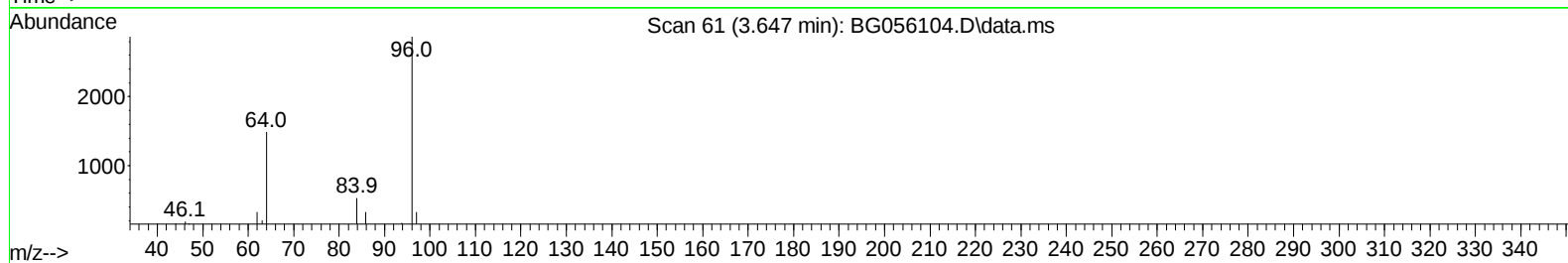
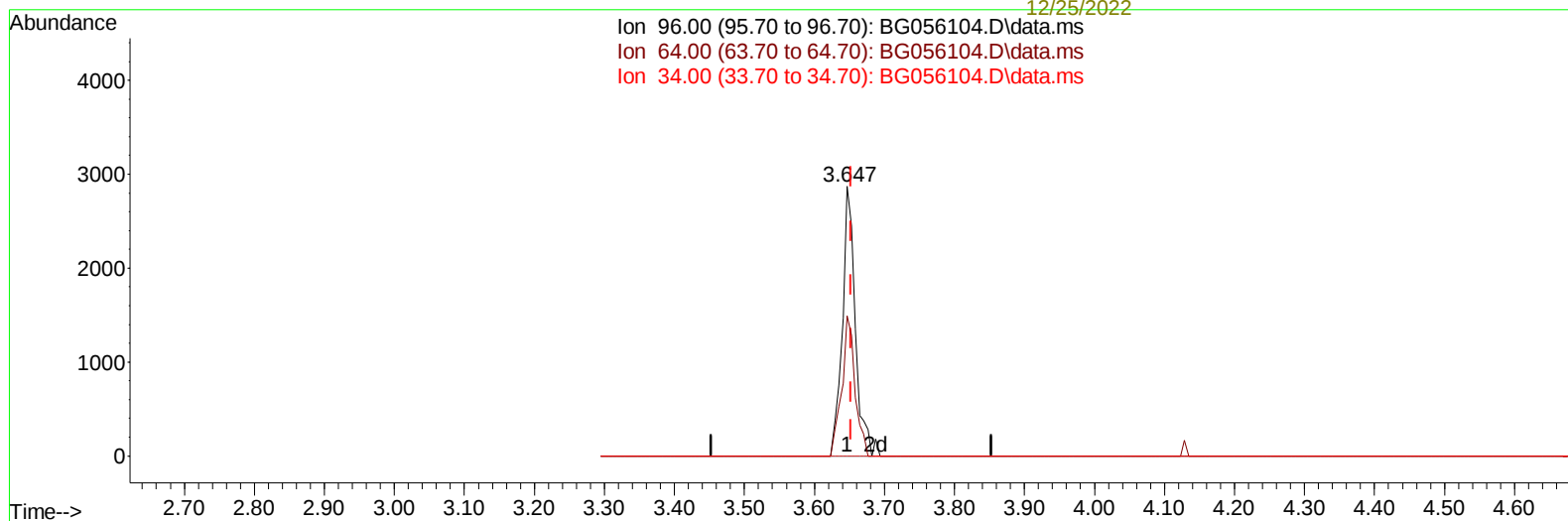
Upadhyay

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TIC: BG056104.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.647min (-0.006) 3.90 ng/uL m

response 3654

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	40.20	52.00#
34.00	0.00	0.00
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	8.347	152	39177	20.000	ng/ul	0.00
20) Naphthalene-d8	11.179	136	157916	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.951	164	136875	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.689	188	353916	20.000	ng/ul	0.00
79) Chrysene-d12	21.990	240	345423	20.000	ng/ul	# 0.00
88) Perylene-d12	25.462	264	370521	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.647	96	3654m	3.904	ng/uL	0.00
4) Pyridine-d5	4.082	84	59380	19.459	ng/ul	0.00
7) Phenol-d5	7.472	99	81999	23.074	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.654	67	32859	23.460	ng/ul	0.00
11) 2-Chlorophenol-d4	7.865	132	52540	25.015	ng/ul	0.00
15) 4-Methylphenol-d8	9.035	113	65664	24.551	ng/ul	0.00
21) Nitrobenzene-d5	9.516	128	27723	24.103	ng/ul	0.00
24) 2-Nitrophenol-d4	10.245	143	36595	24.352	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.780	165	79074	23.680	ng/ul	0.00
31) 4-Chloroaniline-d4	11.303	131	64207	16.999	ng/ul	0.00
46) Dimethylphthalate-d6	14.340	166	256415	24.205	ng/ul	0.00
49) Acenaphthylene-d8	14.652	160	288771	24.138	ng/ul	0.00
54) 4-Nitrophenol-d4	15.110	143	36644	22.979	ng/ul	0.00
60) Fluorene-d10	15.938	176	247360	23.656	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.038	200	49121	21.591	ng/ul	0.00
73) Anthracene-d10	17.789	188	384302	23.498	ng/ul	0.00
81) Pyrene-d10	20.051	212	466594	24.459	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.221	264	452030	24.002	ng/ul	0.00

Target Compounds Qvalue

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

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