Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056103.D
Acq On : 25 Dec 2022 00:34
Operator : CG/JU
Sample : N6017-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Dec 25 01:52:47 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

Instrument :
BNA_G
ClientSampleId :
DBYF1

Manual IntegrationsAPPROVED

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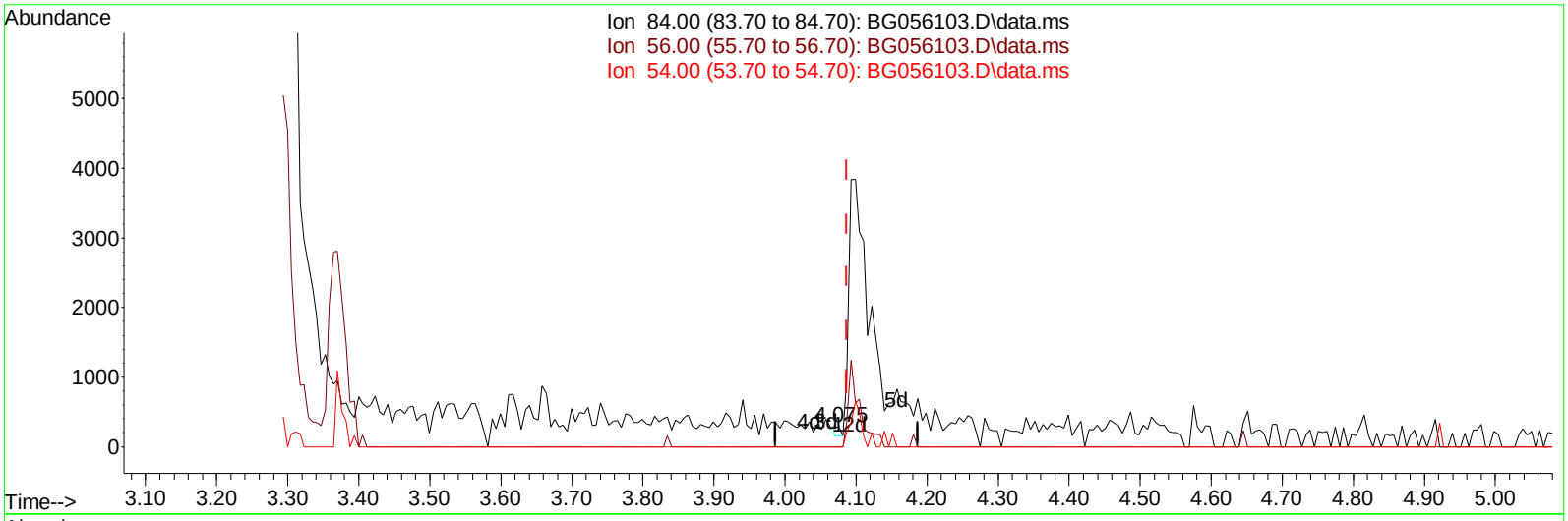
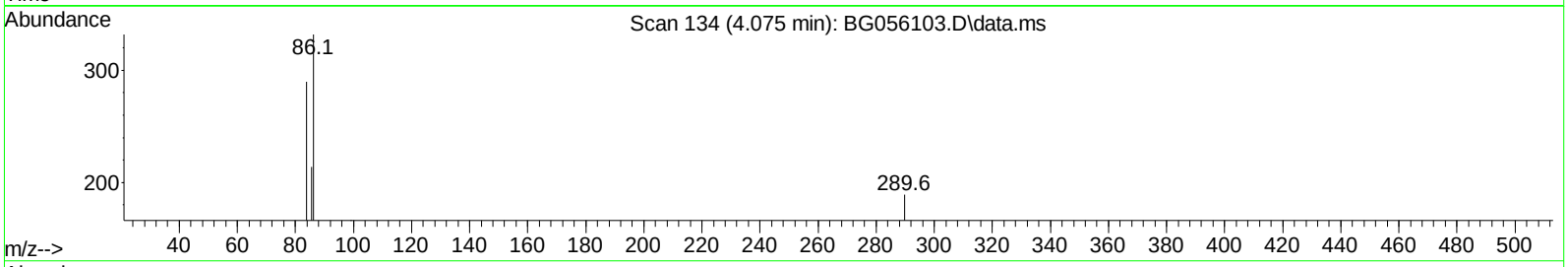
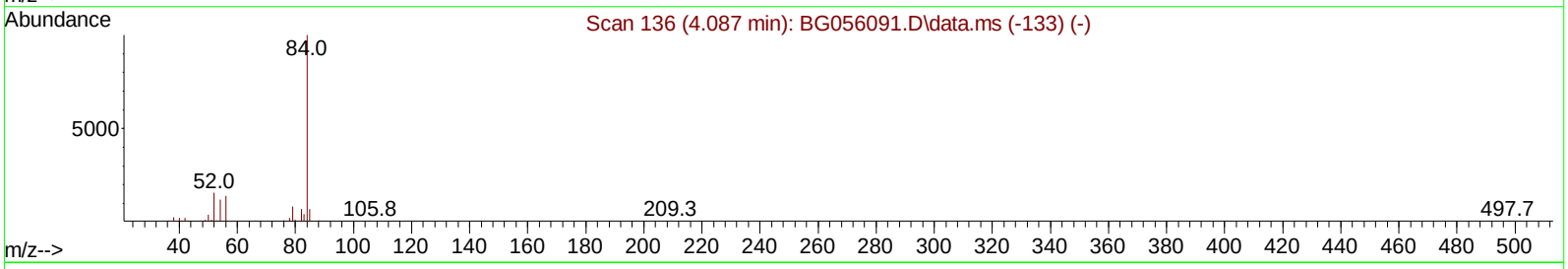
Supervised By :Yogesh Patel 12/25/2022







TIC: BG056103.D\data.ms

(4) Pyridine-d5 (S)

4.075min (-0.012) 0.02 ng/ul

response 47

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	13.80	0.00#
54.00	11.70	0.00#
0.00	0.00	0.00

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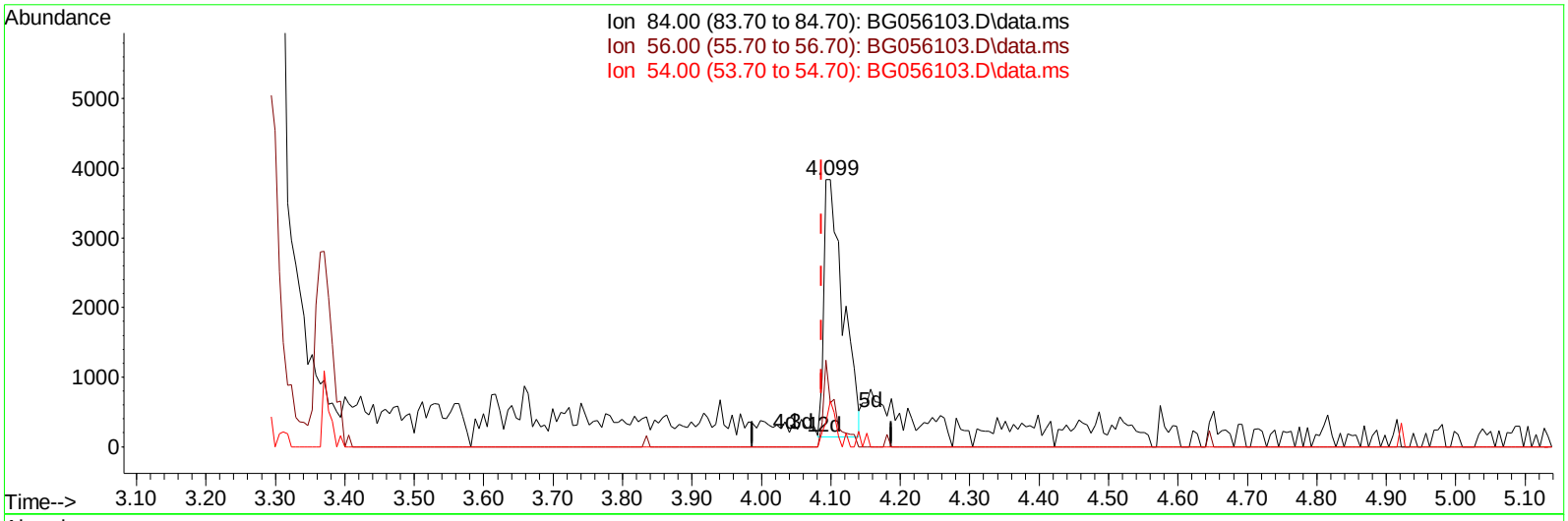

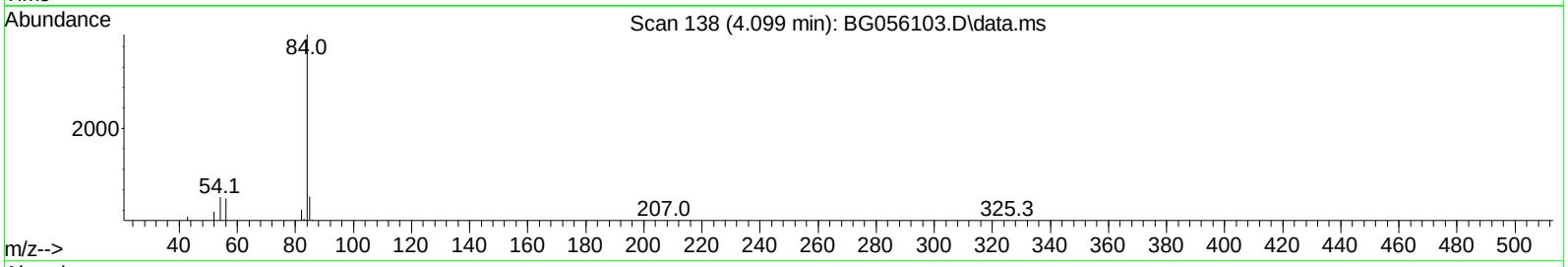

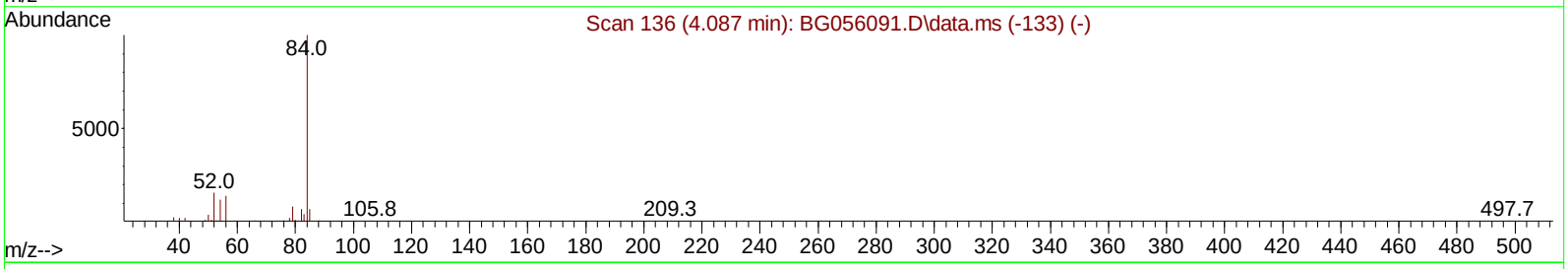

TIC: BG056103.D\data.ms

(4) Pyridine-d5 (S)

4.099min (+ 0.011) 2.42 ng/ul m

response 7057

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	13.80	16.32
54.00	11.70	17.03#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.341	152	37475	20.000	ng/ul	0.00
20) Naphthalene-d8	11.179	136	149612	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.951	164	134282	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.689	188	356630	20.000	ng/ul	0.00
79) Chrysene-d12	21.990	240	349897	20.000	ng/ul	0.00
88) Perylene-d12	25.456	264	374289	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.652	96	2471	2.760	ng/uL	0.00
4) Pyridine-d5	4.099	84	7057m	2.418	ng/ul	0.01
7) Phenol-d5	7.471	99	59005	17.358	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.653	67	22982	17.153	ng/ul	0.00
11) 2-Chlorophenol-d4	7.871	132	37205	18.518	ng/ul	0.00
15) 4-Methylphenol-d8	9.034	113	42009	16.420	ng/ul	0.00
21) Nitrobenzene-d5	9.516	128	18651	17.116	ng/ul	0.00
24) 2-Nitrophenol-d4	10.245	143	24646	17.311	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.785	165	55591	17.571	ng/ul	0.00
31) 4-Chloroaniline-d4	11.296	131	23506	6.569	ng/ul	0.00
46) Dimethylphthalate-d6	14.340	166	190164	18.298	ng/ul	0.00
49) Acenaphthylene-d8	14.651	160	206696	17.611	ng/ul	0.00
54) 4-Nitrophenol-d4	15.109	143	28301	18.090	ng/ul	0.00
60) Fluorene-d10	15.938	176	187516	18.279	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.032	200	36098	15.746	ng/ul	-0.01
73) Anthracene-d10	17.789	188	313776	19.039	ng/ul	0.00
81) Pyrene-d10	20.051	212	399482	20.673	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.215	264	378022	19.870	ng/ul	0.00

Target Compounds Qvalue

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

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