Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122422\
Data File : BG056102.D
Acq On : 24 Dec 2022 23:54
Operator : CG/JU
Sample : N6017-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 25 01:52:40 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 :
DBYF0

Manual IntegrationsAPPROVED

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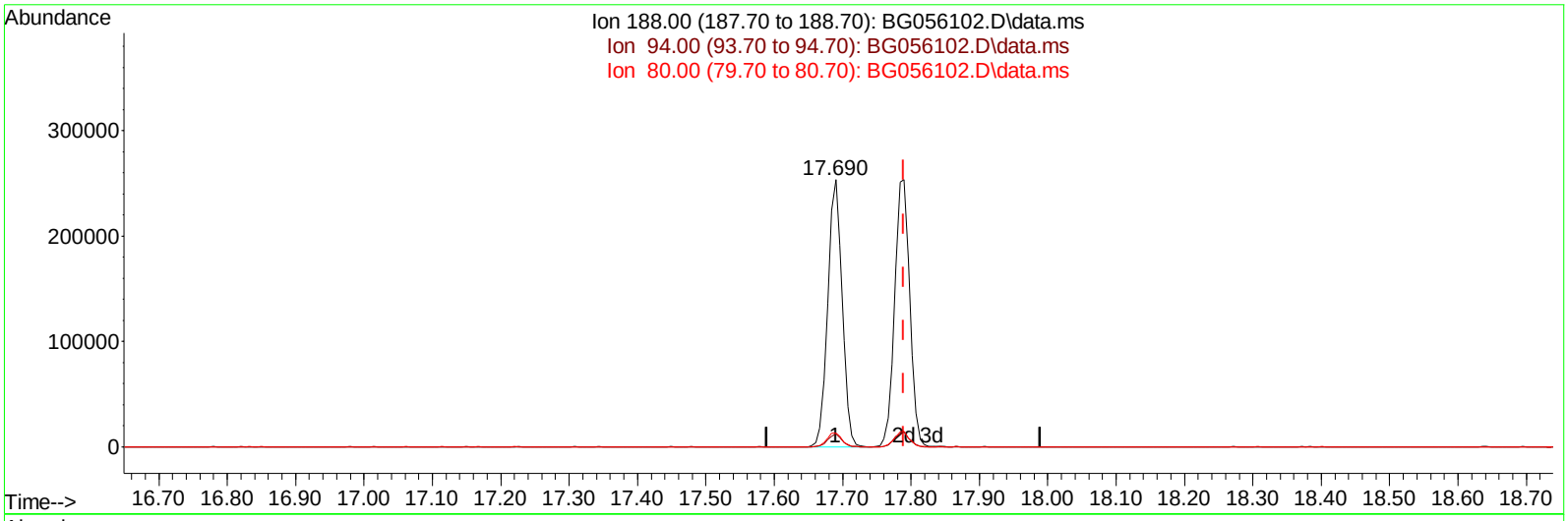
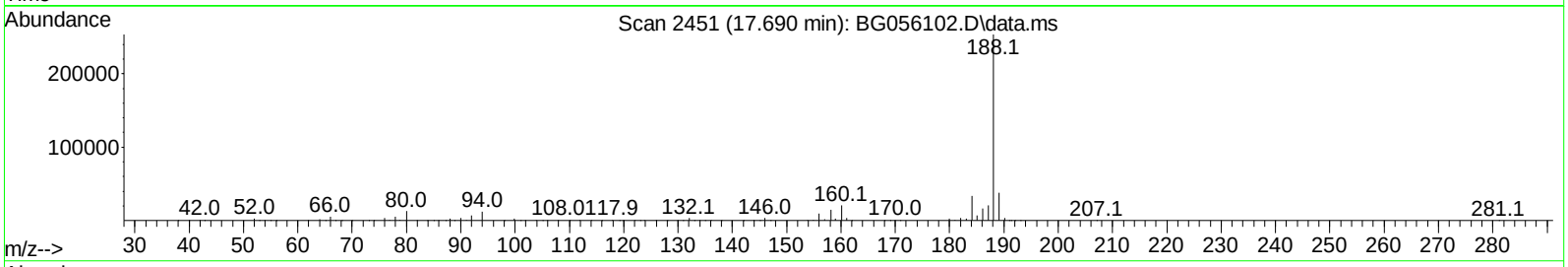
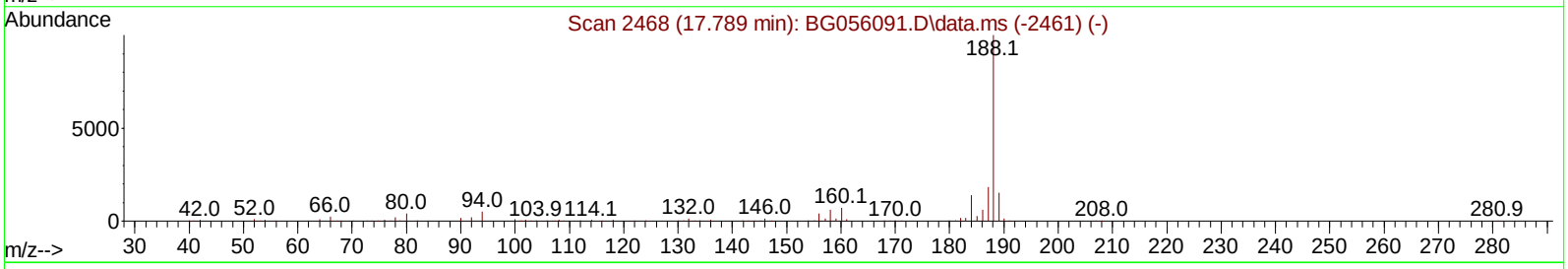
Supervised By :Yogesh Patel 12/25/2022







TIC: BG056102.D\data.ms

(73) Anthracene-d10 (S)

17.690min (-0.099) 21.64 ng/ul

response 368101

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.10	4.93
80.00	4.20	5.00
0.00	0.00	0.00

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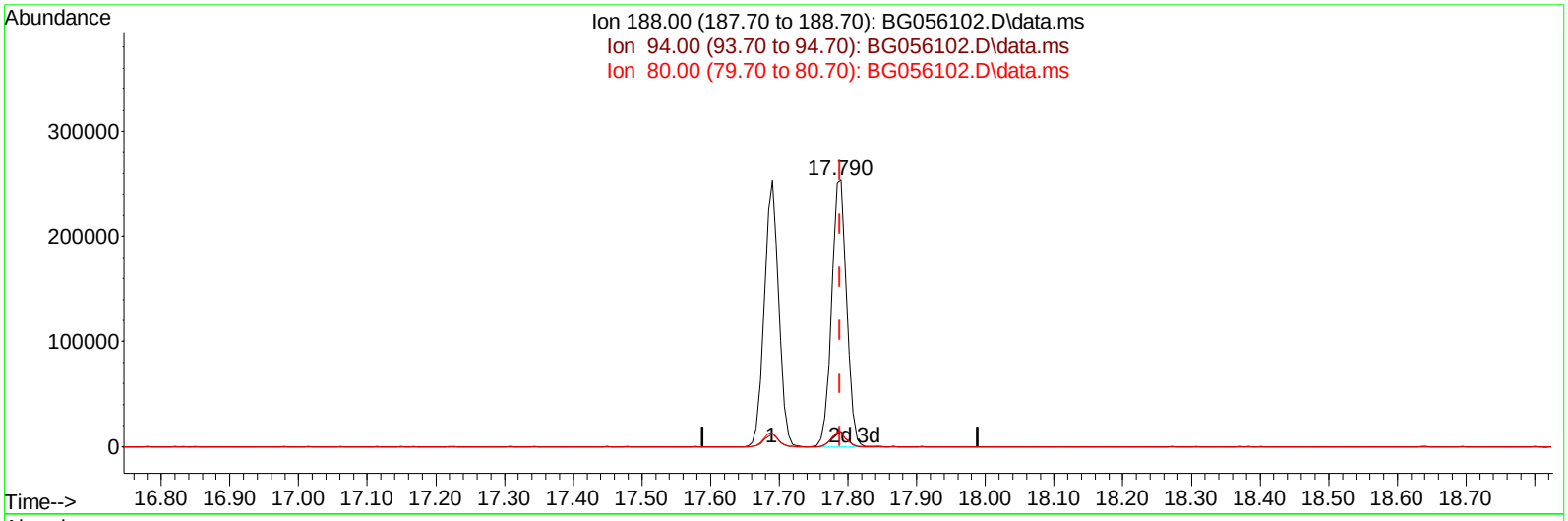

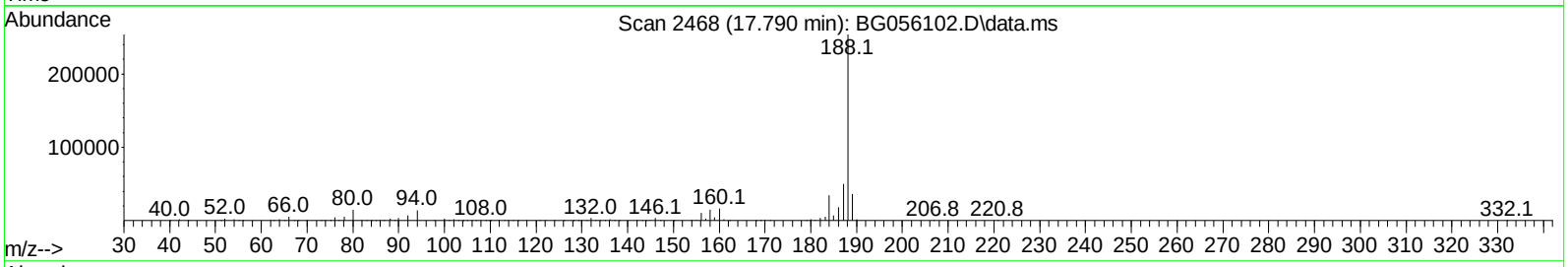

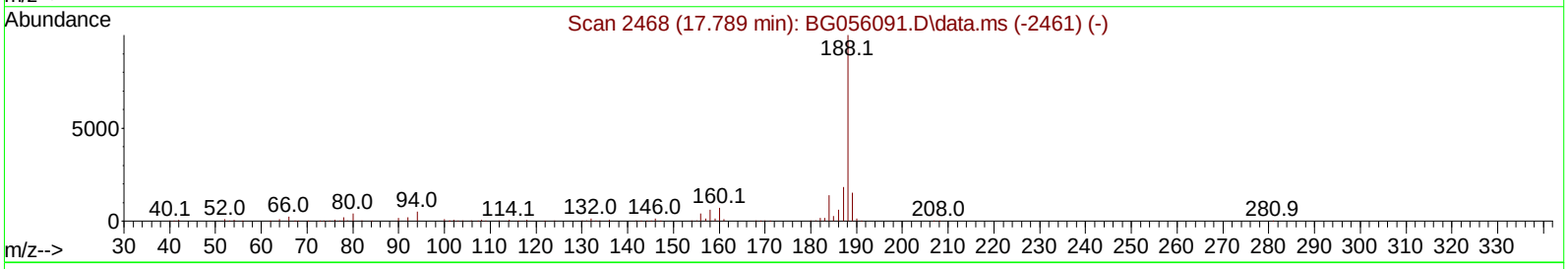

TIC: BG056102.D\data.ms

(73) Anthracene-d10 (S)

17.790min (+ 0.001) 22.70 ng/ul m

response 386124

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	5.10	5.49
80.00	4.20	5.73#
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	8.342	152	37528	20.000	ng/ul	0.00
20) Naphthalene-d8	11.180	136	151544	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.952	164	134270	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.690	188	368101	20.000	ng/ul	0.00
79) Chrysene-d12	21.991	240	368421	20.000	ng/ul	0.00
88) Perylene-d12	25.457	264	389338	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.648	96	3835	4.278	ng/uL	0.00
4) Pyridine-d5	4.082	84	55247	18.900	ng/ul	0.00
7) Phenol-d5	7.473	99	74618	21.920	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.655	67	30894	23.026	ng/ul	0.00
11) 2-Chlorophenol-d4	7.866	132	50643	25.171	ng/ul	0.00
15) 4-Methylphenol-d8	9.035	113	56008	21.861	ng/ul	0.00
21) Nitrobenzene-d5	9.517	128	26357	23.879	ng/ul	0.00
24) 2-Nitrophenol-d4	10.246	143	32741	22.704	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.780	165	70211	21.910	ng/ul	0.00
31) 4-Chloroaniline-d4	11.297	131	69952	19.299	ng/ul	0.00
46) Dimethylphthalate-d6	14.341	166	245196	23.595	ng/ul	0.00
49) Acenaphthylene-d8	14.652	160	267632	22.805	ng/ul	0.00
54) 4-Nitrophenol-d4	15.111	143	34435	22.013	ng/ul	0.00
60) Fluorene-d10	15.939	176	233883	22.801	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.033	200	45530	19.241	ng/ul	-0.01
73) Anthracene-d10	17.790	188	386124m	22.699	ng/ul	0.00
81) Pyrene-d10	20.052	212	466867	22.945	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.216	264	460345	23.262	ng/ul	0.00

Target Compounds Qvalue

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

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