Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062822\
Data File : BM035729.D
Acq On : 28 Jun 2022 18:10
Operator : CG/JU
Sample : N3401-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Jun 29 04:27:29 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 29 04:24:18 2022
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
C0AE3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/29/2022
Supervised By :mohammad ahmed 07/01/2022

06/29/2022
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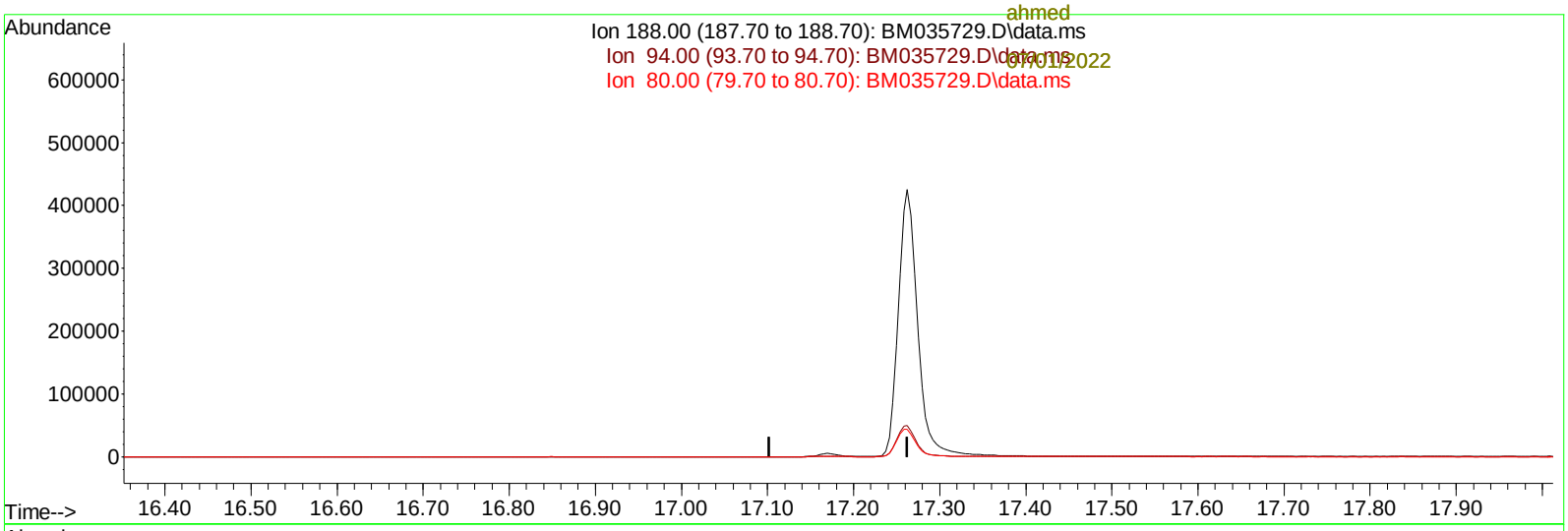
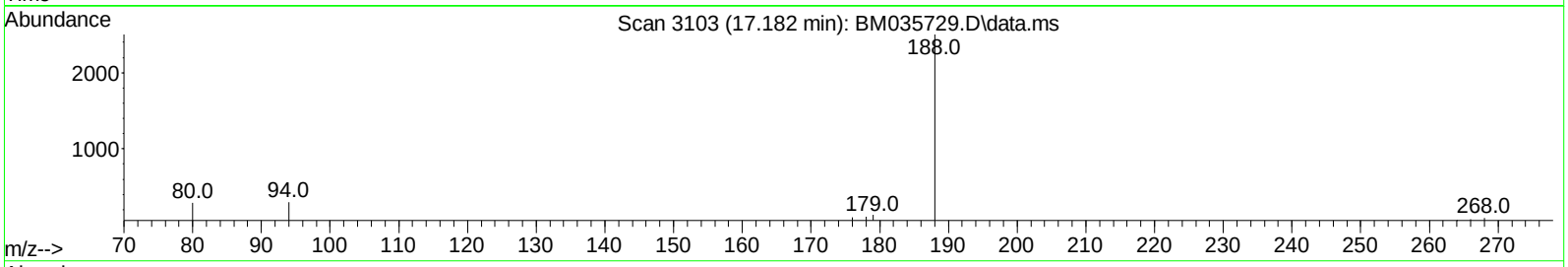
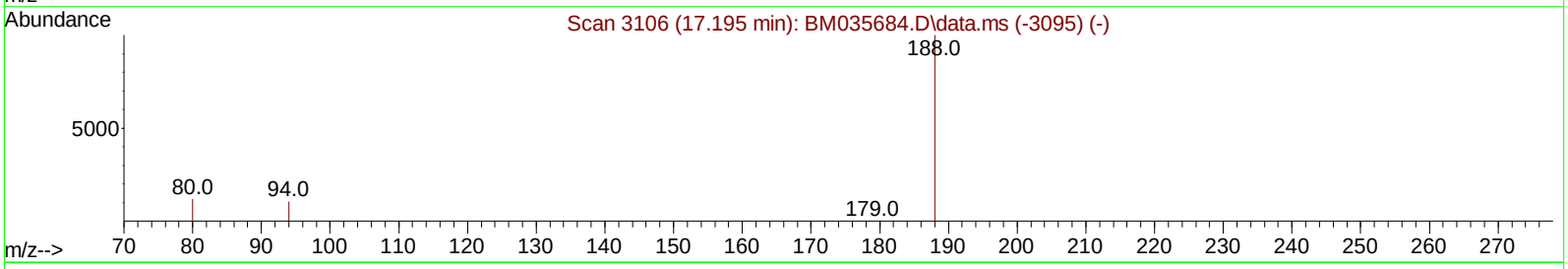
07/01/2022







TIC: BM035729.D\data.ms

(13) Phenanthrene-d10 (I)

17.182min (-17.182) 0.00 ng/ul

response 0

Ion	Exp%	Act%
188.00	100.00	0.00
94.00	11.40	0.00#
80.00	12.70	0.00#
0.00	0.00	0.00

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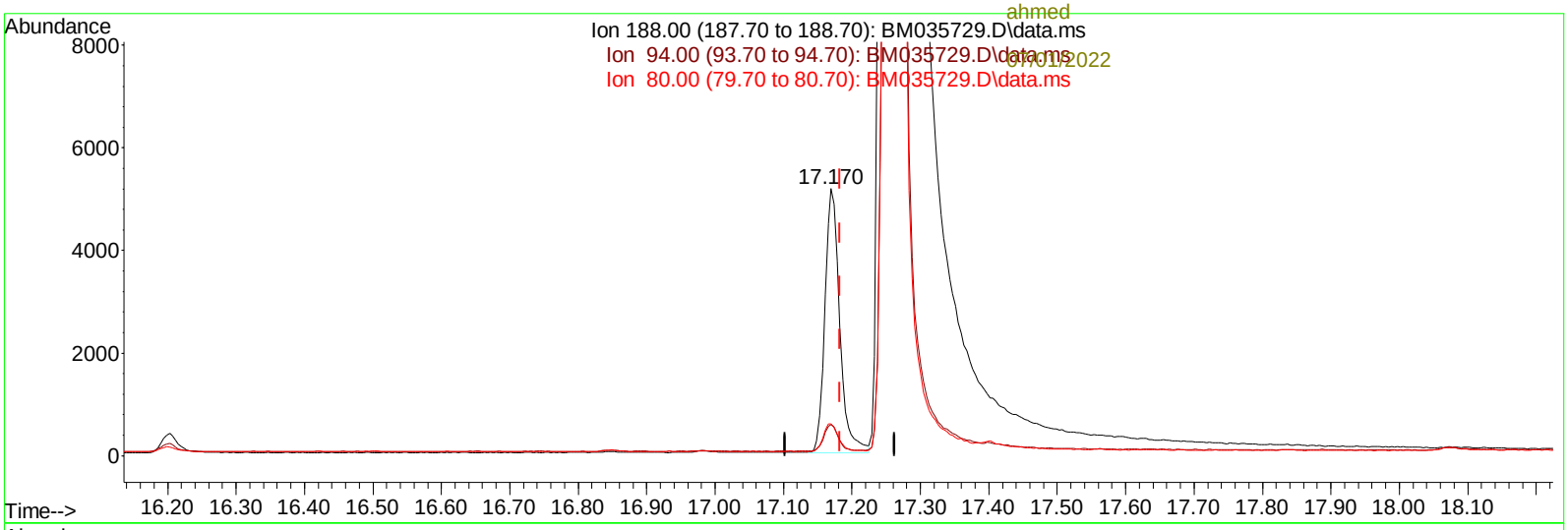

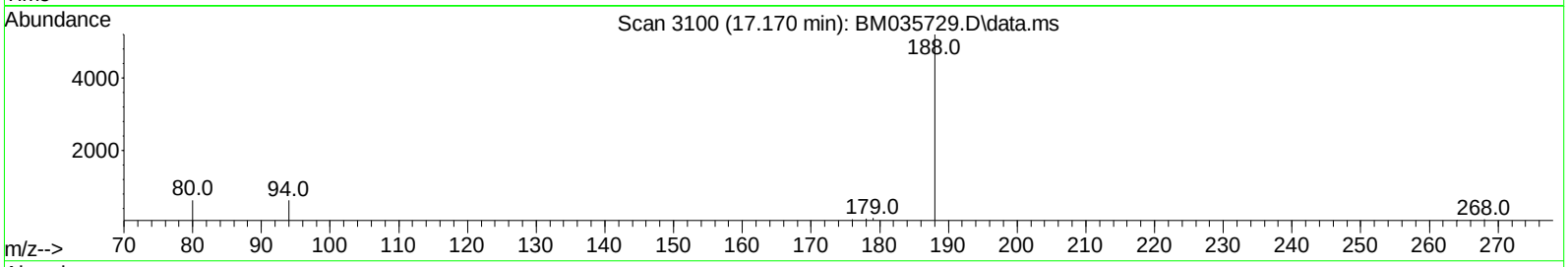

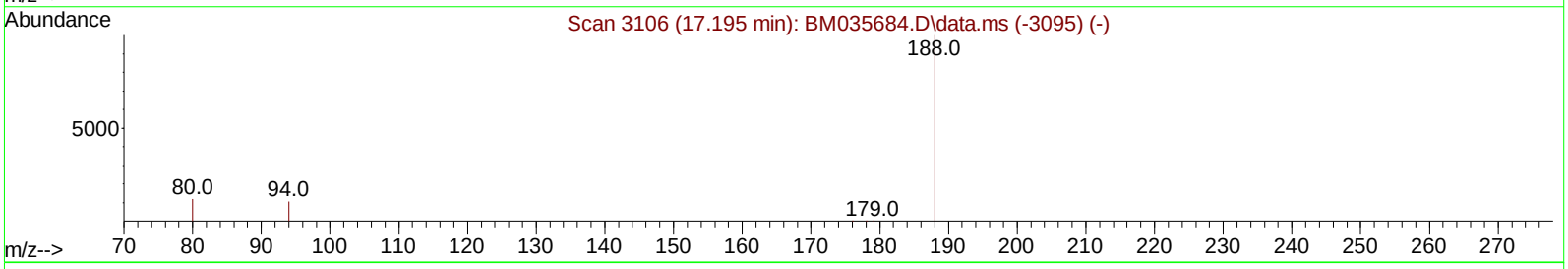

TIC: BM035729.D\data.ms

(13) Phenanthrene-d10 (I)

17.170min (-0.013) 0.40 ng/ul m

response 7759

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	11.40	11.78
80.00	12.70	12.05
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Data File : BM035729.D
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Sample : N3401-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Jun 29 04:27:29 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062522.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.795	152	2441	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.578	136	7731	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.422	164	4106	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.170	188	7759m	0.400	ng/ul	-0.01
17) Chrysene-d12	21.380	240	6461	0.400	ng/ul	# 0.00
23) Perylene-d12	23.692	264	8069	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.209	96	14444	4.360	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.217	152	3506	0.290	ng/ul	0.02
18) Fluoranthene-d10	19.205	212	7366	0.338	ng/ul	0.00
Target Compounds						Qvalue
21) Benzo(a)anthracene	21.377	228	690	0.031	ng/ul#	22
22) Chrysene	21.424	228	1038	0.036	ng/ul#	42
24) Benzo(b)fluoranthene	22.987	252	1435	0.041	ng/ul#	1
25) Benzo(k)fluoranthene	23.034	252	1292	0.037	ng/ul#	1

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

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