

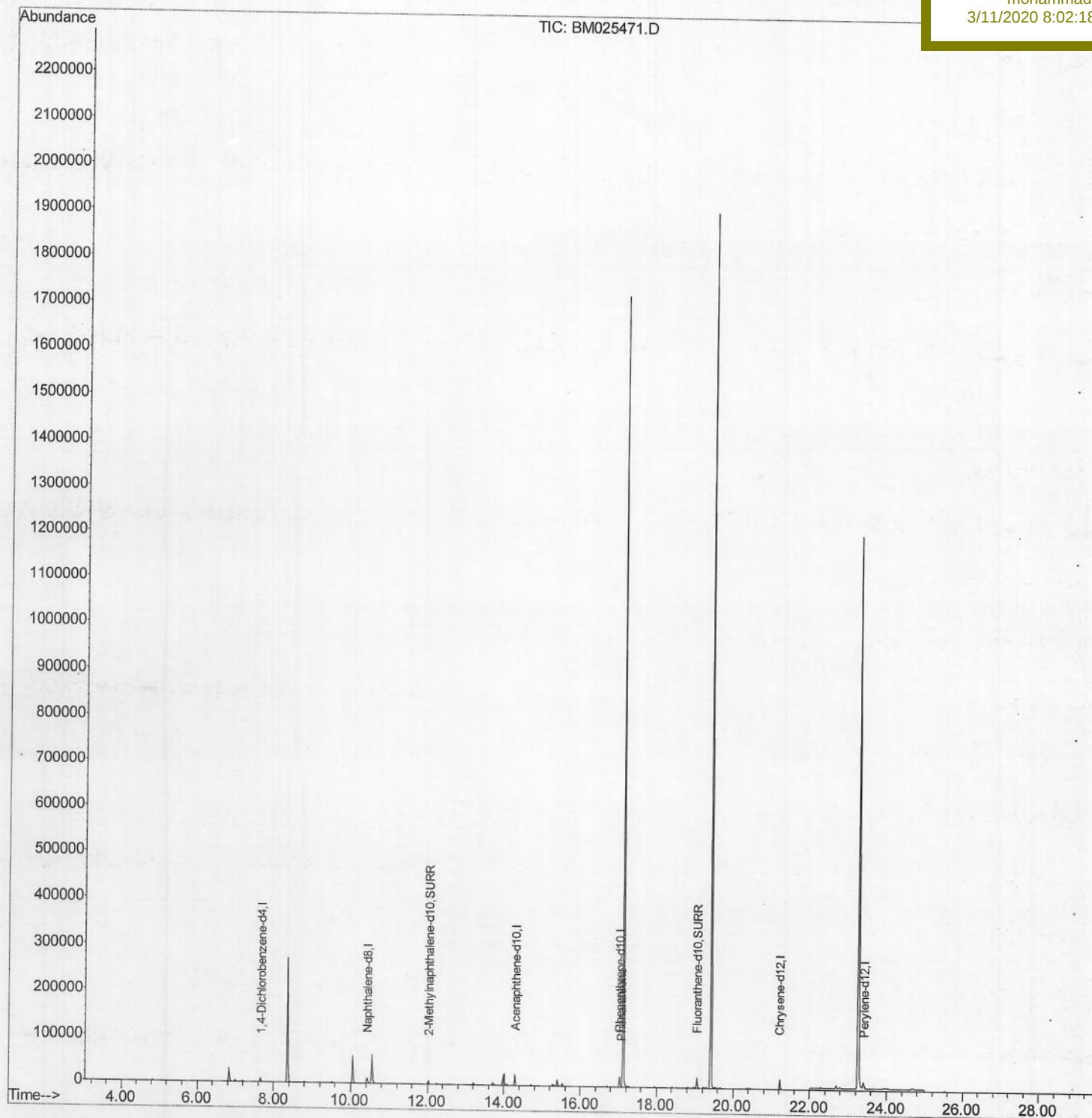
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030920\
Data File : BM025471.D
Acq On : 11 Mar 2020 02:20
Operator : CG/JU
Sample : L1617-11
Misc :
ALS Vial : 65 Sample Multiplier: 1

Quant Time: Mar 11 03:46:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 10 23:48:01 2020
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
DBDS5

Manual Integrations
APPROVED

mohammad
3/11/2020 8:02:18 AM



Quantitation Report (Qedit)

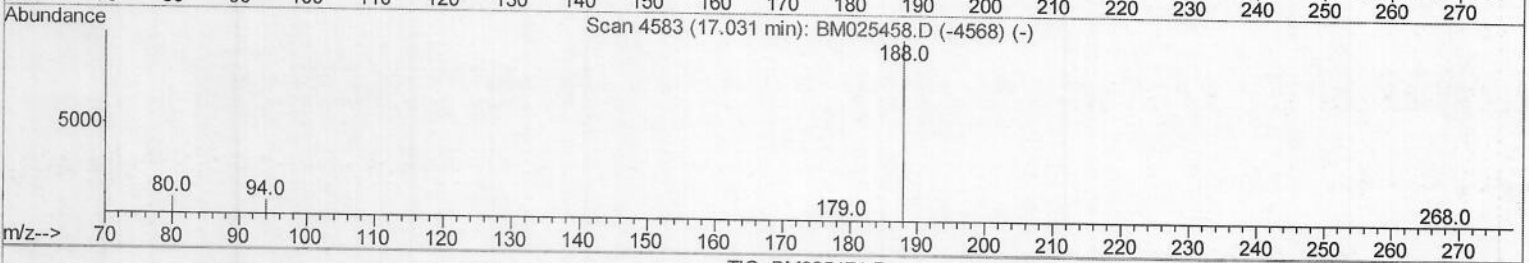
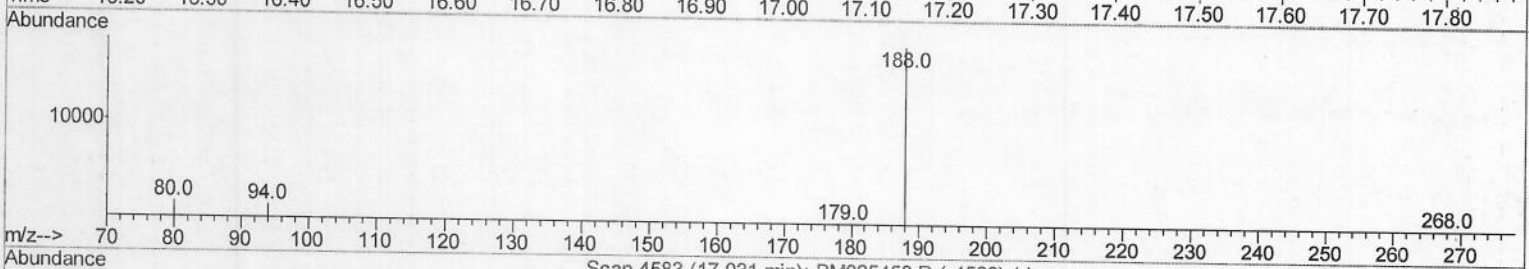
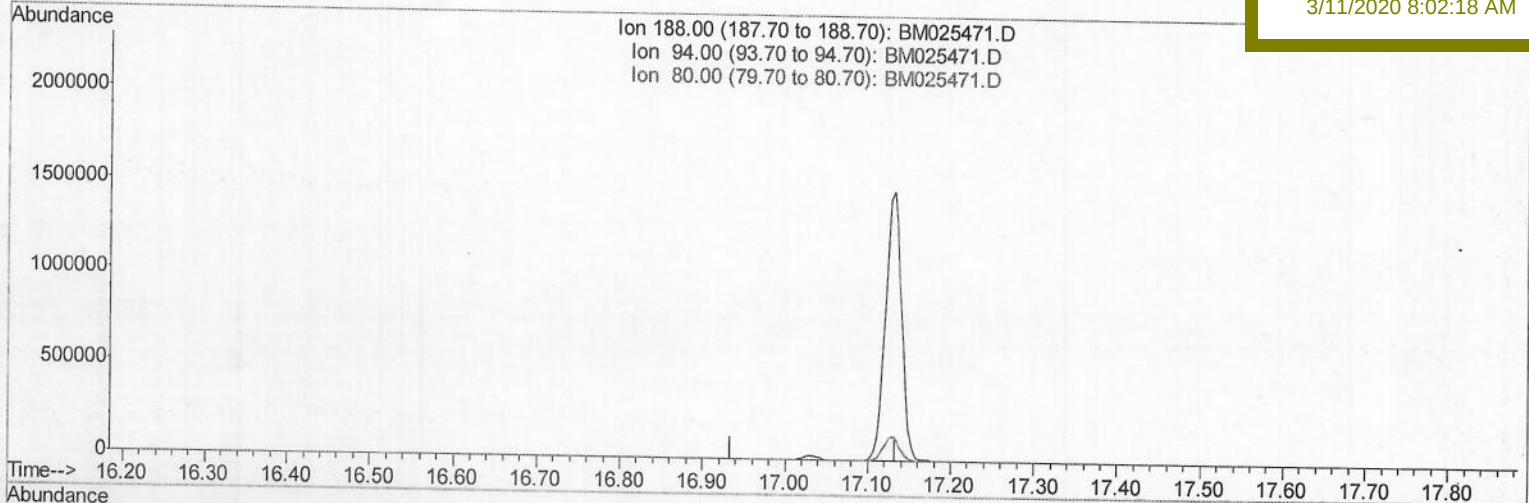
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030920\
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TIC: BM025471.D

(10) Phenanthrene-d10 (I)

17.034min (-17.034) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	7.10	0.00#
80.00	8.70	0.00#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030920\

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Acq On : 11 Mar 2020 02:20

Operator : CG/JU

Sample : L1617-11

Misc :

ALS Vial : 65 Sample Multiplier: 1

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Mar 10 23:48:01 2020

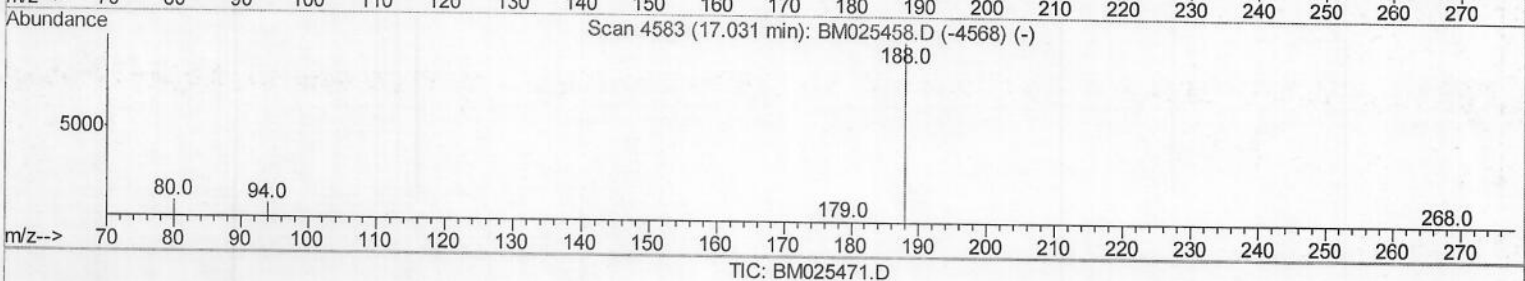
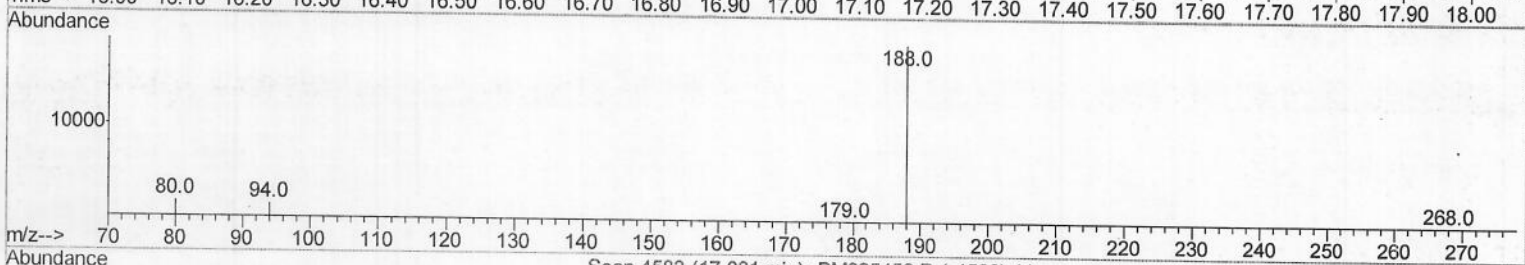
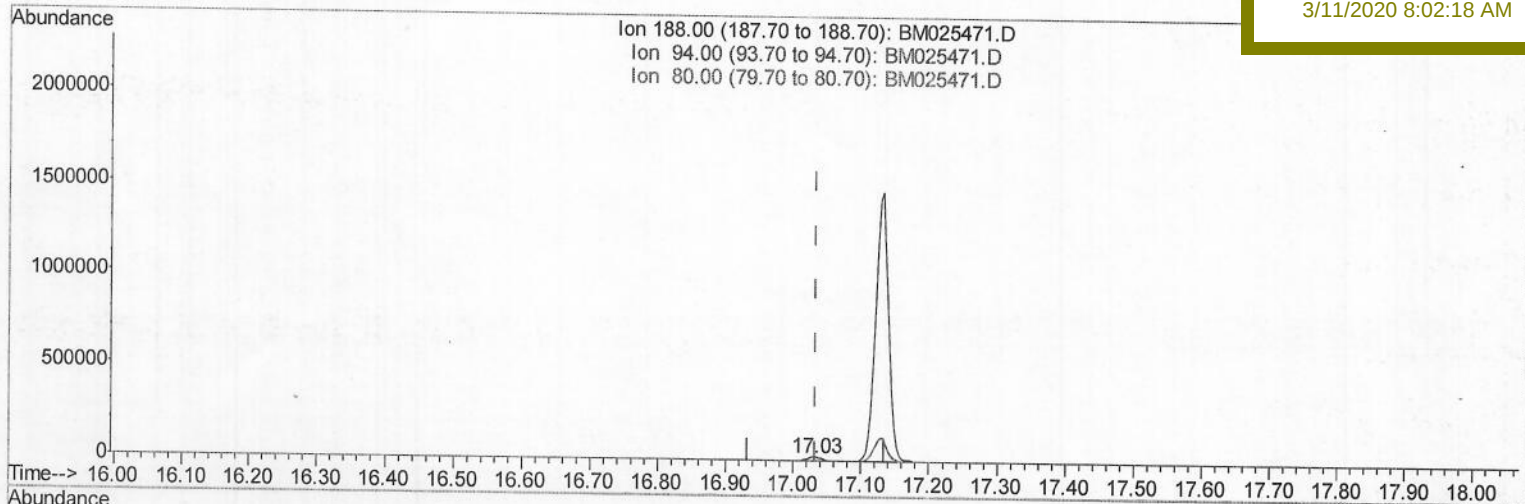
Response via : Initial Calibration

Instrument :

BNA_M

Client Sampled :

DBDS5

Manual Integrations
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(10) Phenanthrene-d10 (I)

17.031min (-0.003) 0.40ng/ul m15V 03/11/20

response 26282

Ion	Exp%	Act%
188.00	100	100
94.00	7.10	8.03
80.00	8.70	9.30
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030920\
 Data File : BM025471.D
 Acq On : 11 Mar 2020 02:20
 Operator : CG/JU
 Sample : L1617-11
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Quant Time: Mar 11 03:46:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 23:48:01 2020
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 DBDS5

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3674	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16471	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	12084	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	26282m)	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	19422	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	18247	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10739	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	26931	0.37	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.07	178	2084	0.025	ng/ul#	Ovalue 89

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