

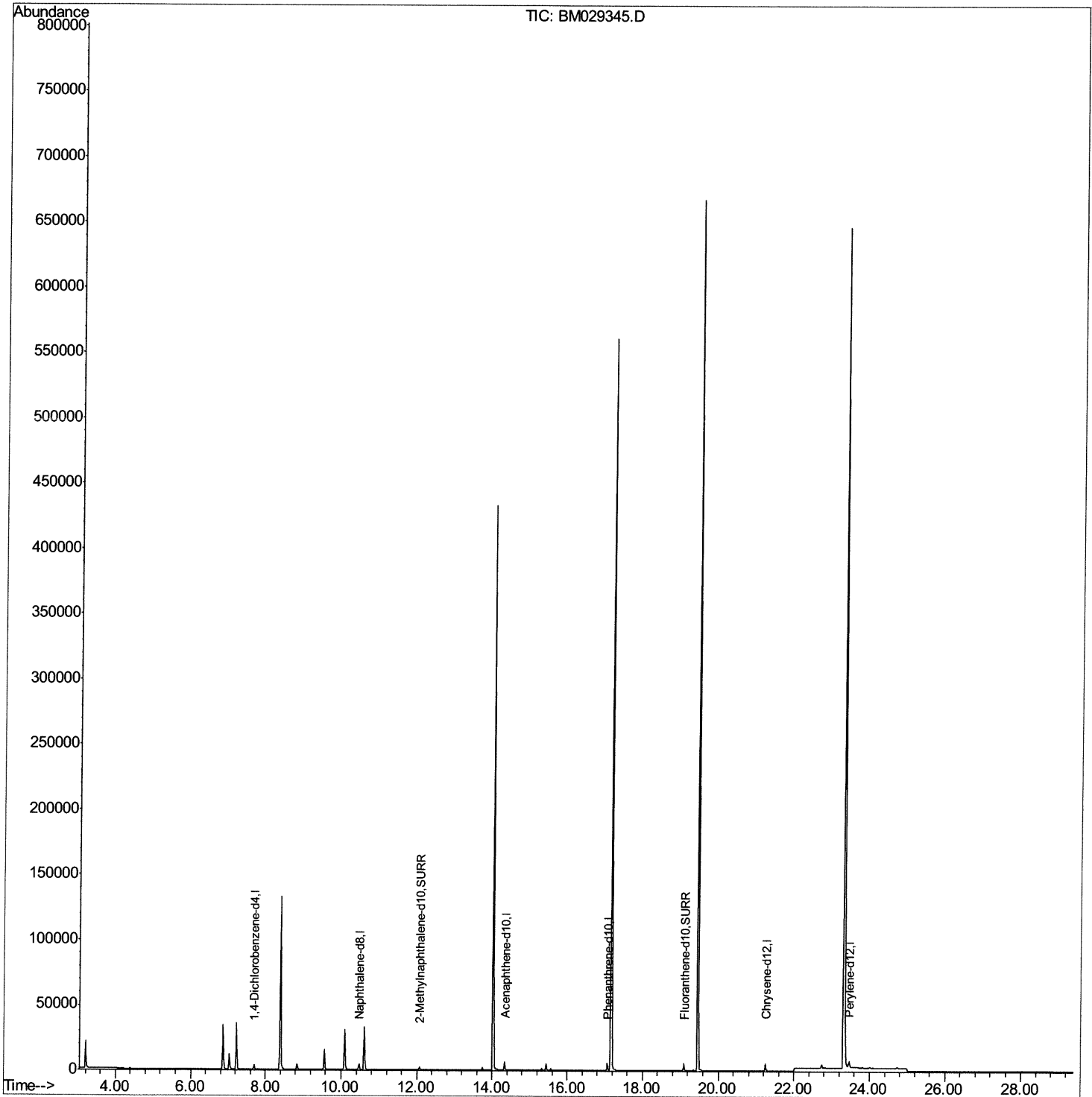
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
Data File : BM029345.D
Acq On : 05 Apr 2021 17:09
Operator : CG/JU
Sample : PB135391BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK91

Manual Integrations
APPROVED

mohammad
4/6/2021 3:23:27 PM

Quant Time: Apr 05 17:43:00 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 05 15:51:17 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

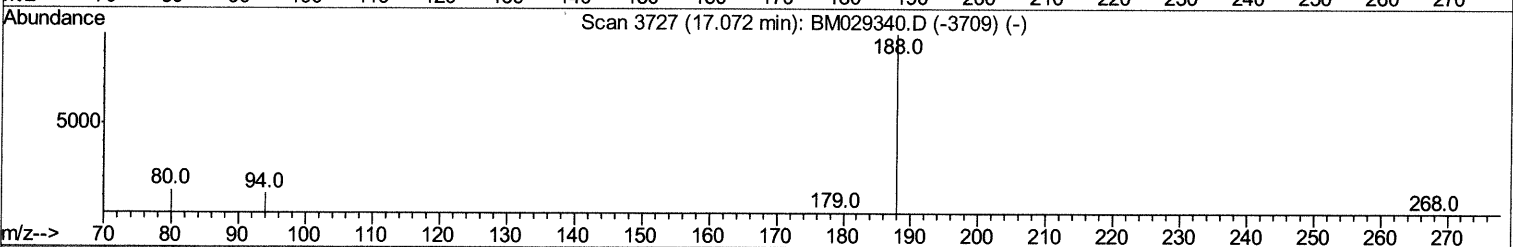
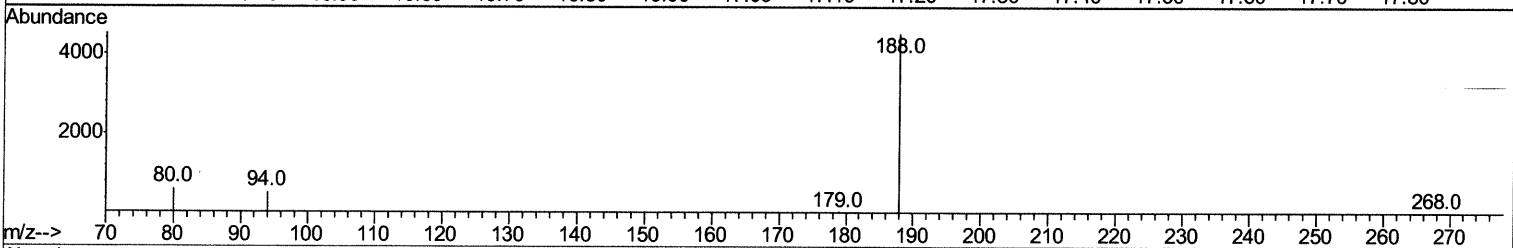
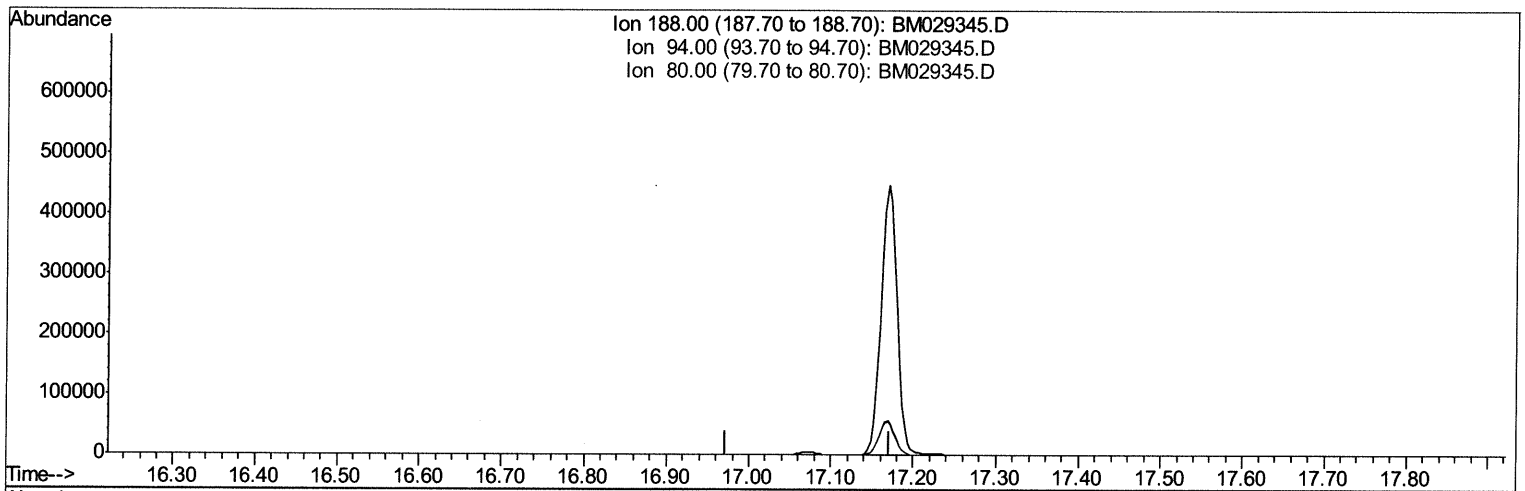
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Quant Time: Apr 05 17:41:18 2021
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TIC: BM029345.D

(10) Phenanthrene-d10 (I)

17.072min (-17.072) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	12.30	0.00#
80.00	13.70	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

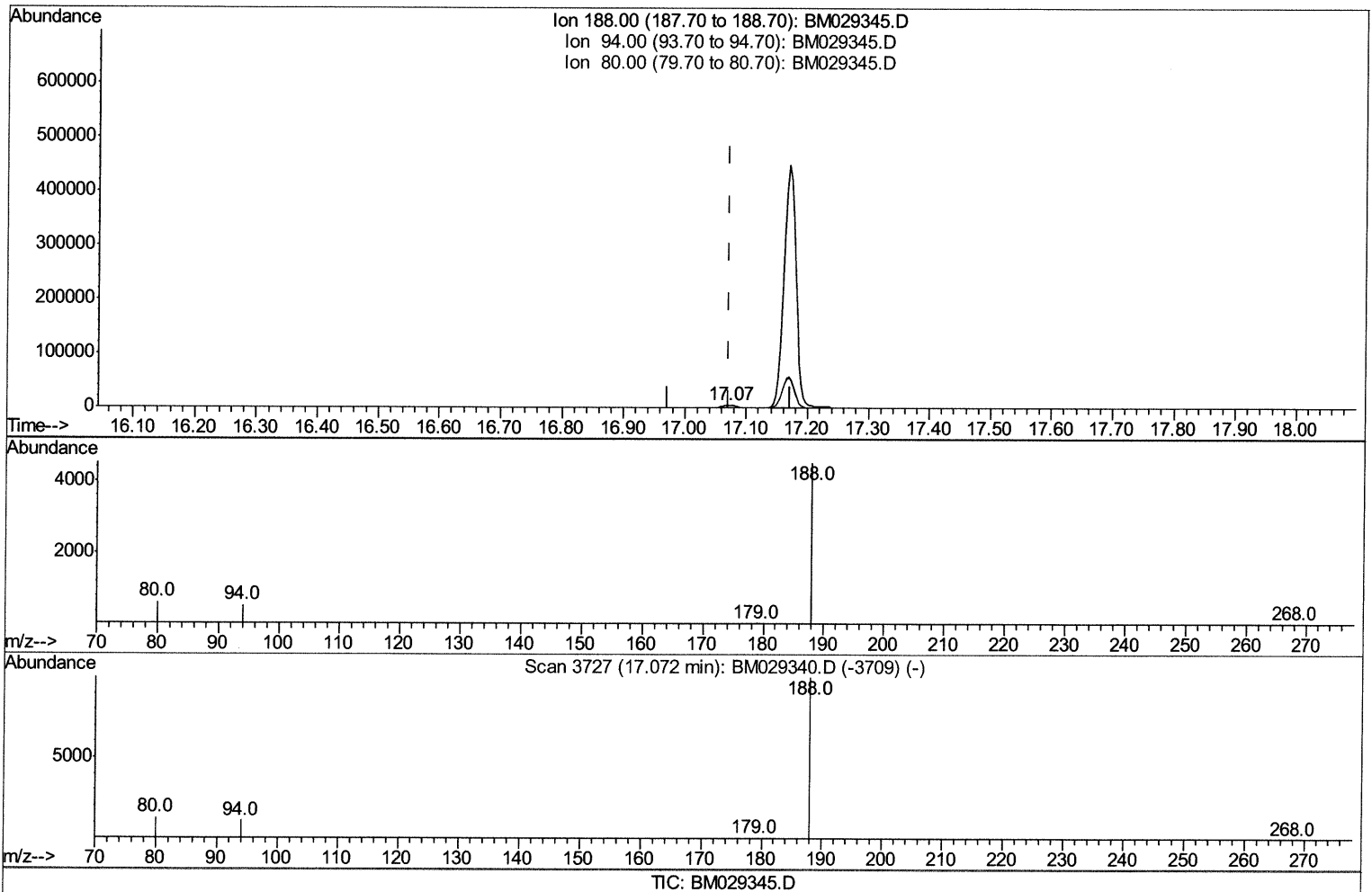
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 Sample : PB135391BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
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(10) Phenanthrene-d10 (I)

17.072min (-0.000) 0.40ng/ul m 04/07/21 JU

response 6203

Ion	Exp%	Act%
188.00	100	100
94.00	12.30	11.97
80.00	13.70	13.94
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
Data File : BM029345.D
Acq On : 05 Apr 2021 17:09
Operator : CG/JU
Sample : PB135391BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK91

Manual Integrations
APPROVED

mohammad
4/6/2021 3:23:27 PM

Quant Time: Apr 05 17:43:00 2021
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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1701	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	5782	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3411	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	6203m>	0.40	ng/ul>	0.00
16) Chrysene-d12	21.25	240	5931	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	7734	0.40	ng/ul	0.00

04/07/21 JU

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	12.07	152	3150	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	6018	0.38	ng/ul	0.00

Target Compounds Ovalue

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