

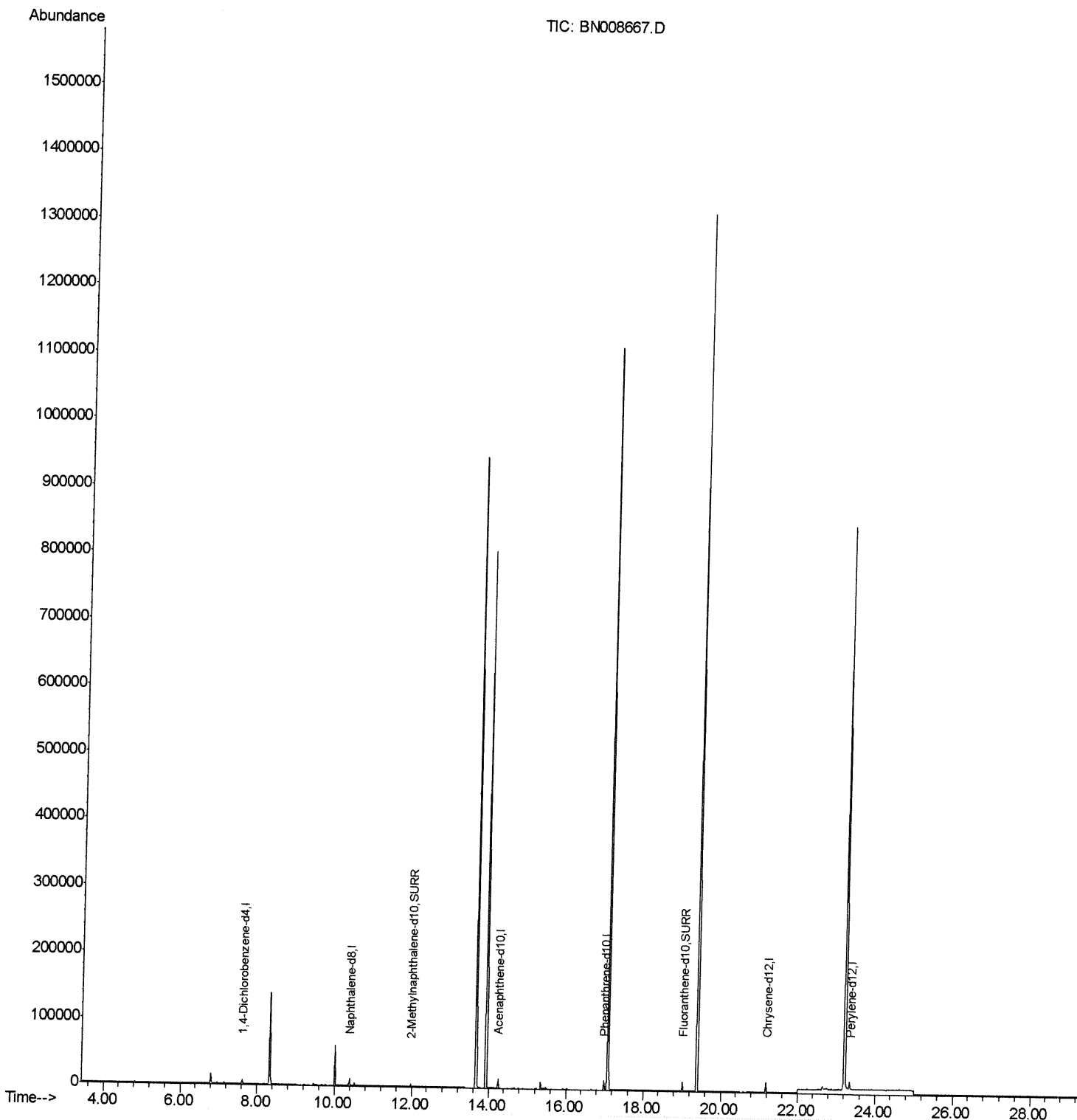
Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
Data File : BN008667.D
Acq On : 20 Nov 2019 19:08
Operator : JU
Sample : K5815-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A5460

Manual Integrations
APPROVED

mohammad
11/21/2019 1:06:19 PM

Quant Time: Nov 21 01:04:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 20 04:45:41 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

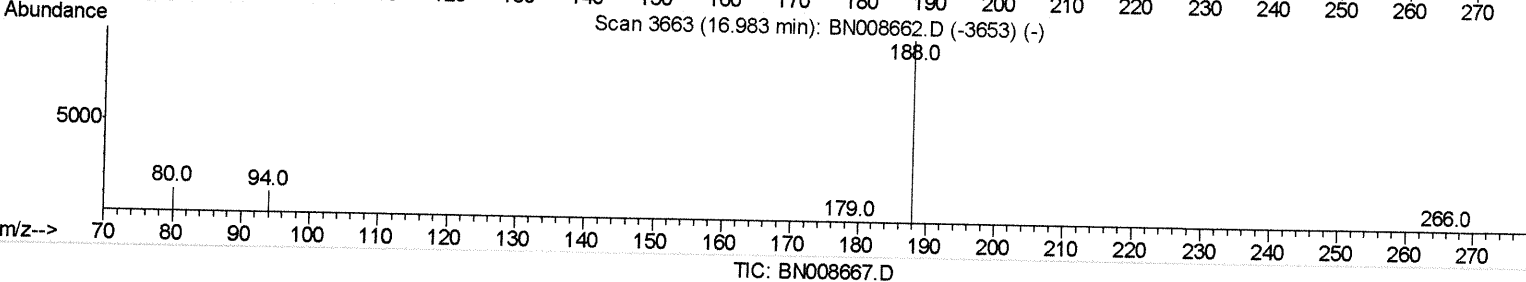
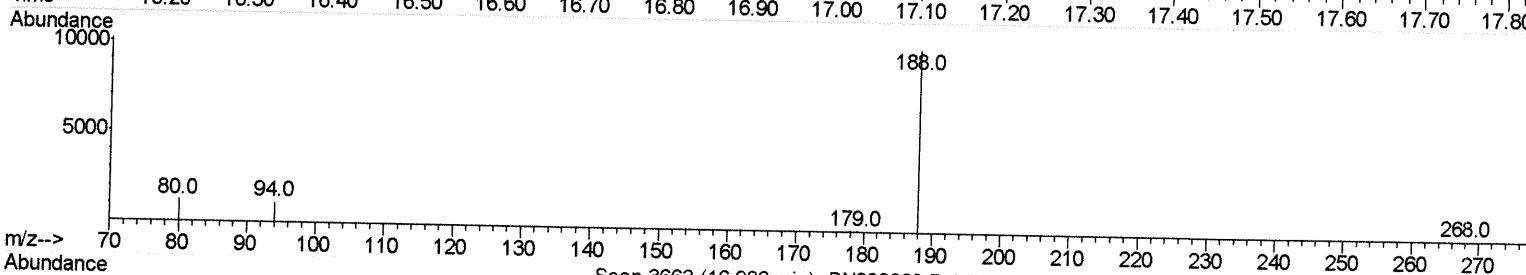
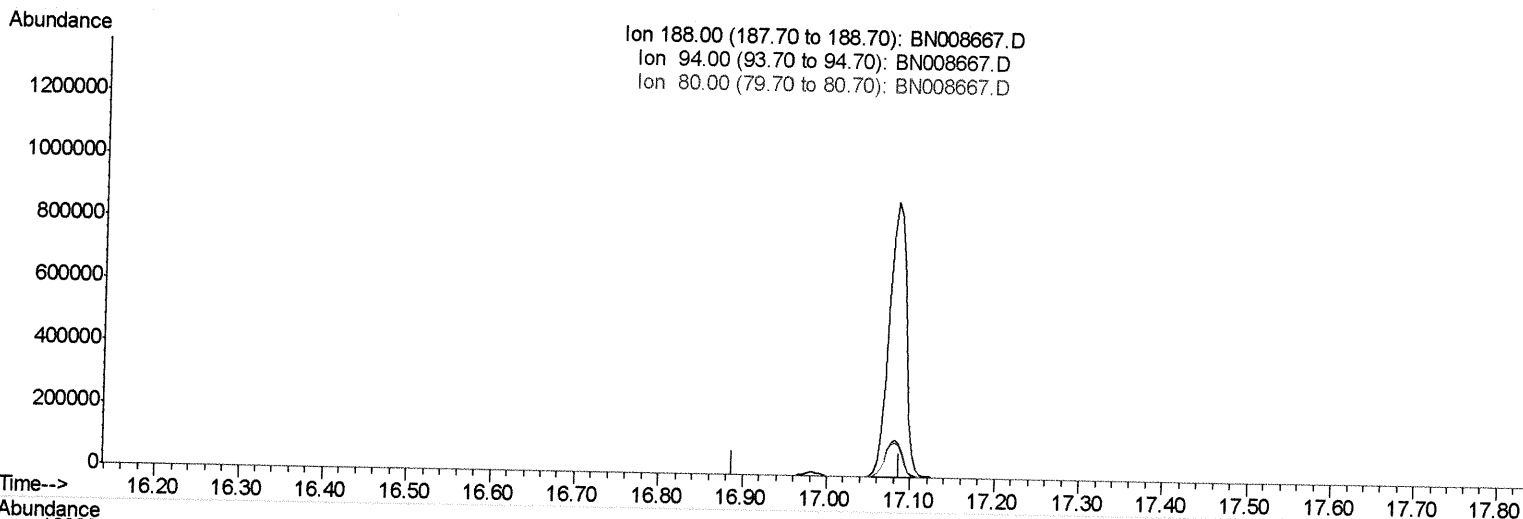
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(10) Phenanthrene-d10 (I)

16.987min (-16.987) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	10.60	0.00#
80.00	11.80	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

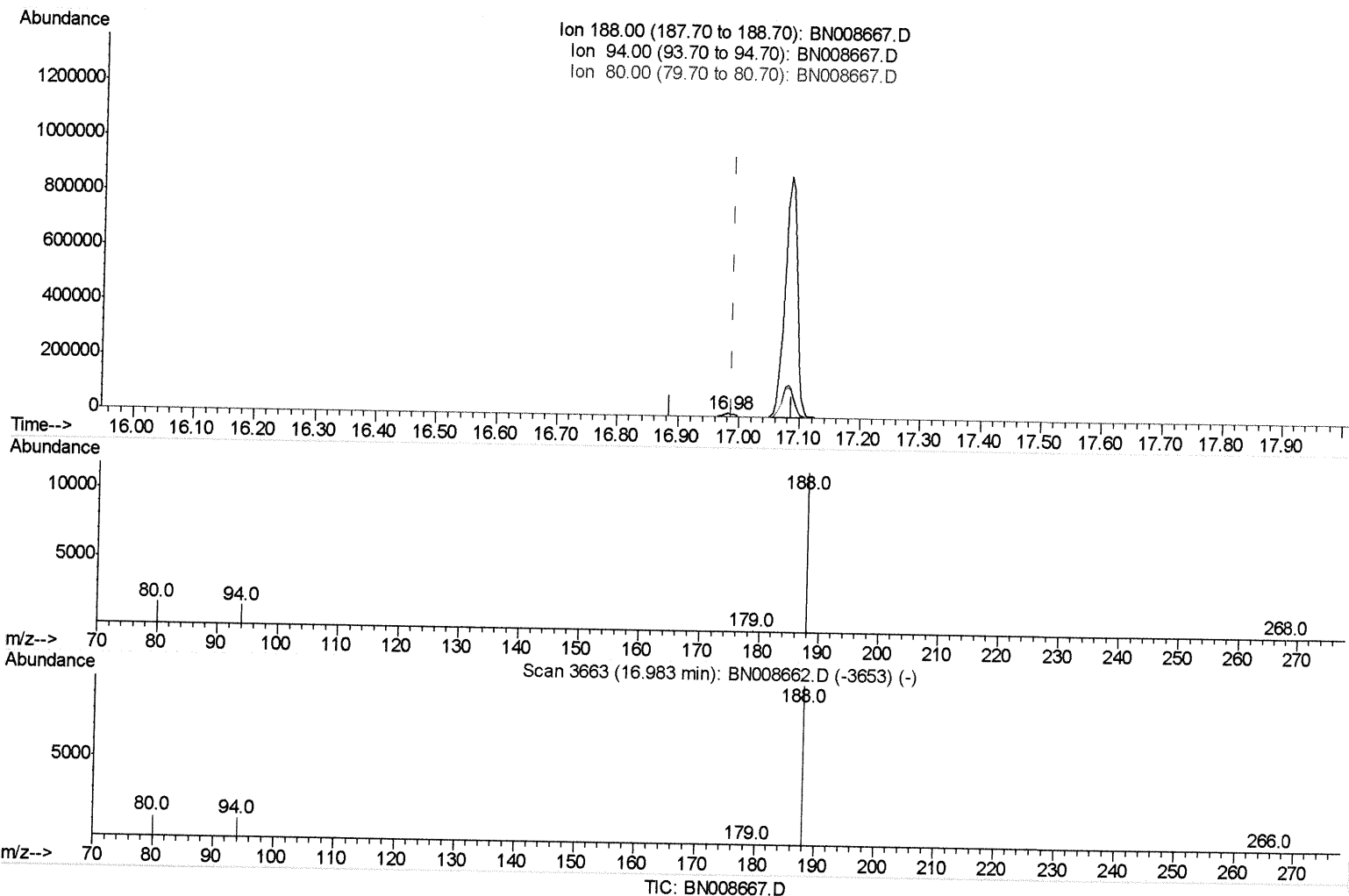
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 Data File : BN008667.D
 Acq On : 20 Nov 2019 19:08
 Operator : JU
 Sample : K5815-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5460

Manual Integrations
 APPROVED

mohammad
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Quant Time: Nov 21 00:43:11 2019
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 Response via : Initial Calibration



(10) Phenanthrene-d10 (I)

16.983min (-0.004) 0.40ng/ul m JU 11/25/19

response 16239

Ion	Exp%	Act%
188.00	100	100
94.00	10.60	11.80
80.00	11.80	13.18
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008667.D
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 Operator : JU
 Sample : K5815-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A5460

Manual Integrations
 APPROVED

mohammad
 11/21/2019 1:06:19 PM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3825	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	14341	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7923	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	16239m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13608	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	13903	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	6872	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	15784	0.32	ng/ul	0.00

Target Compounds

Ovalue

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