

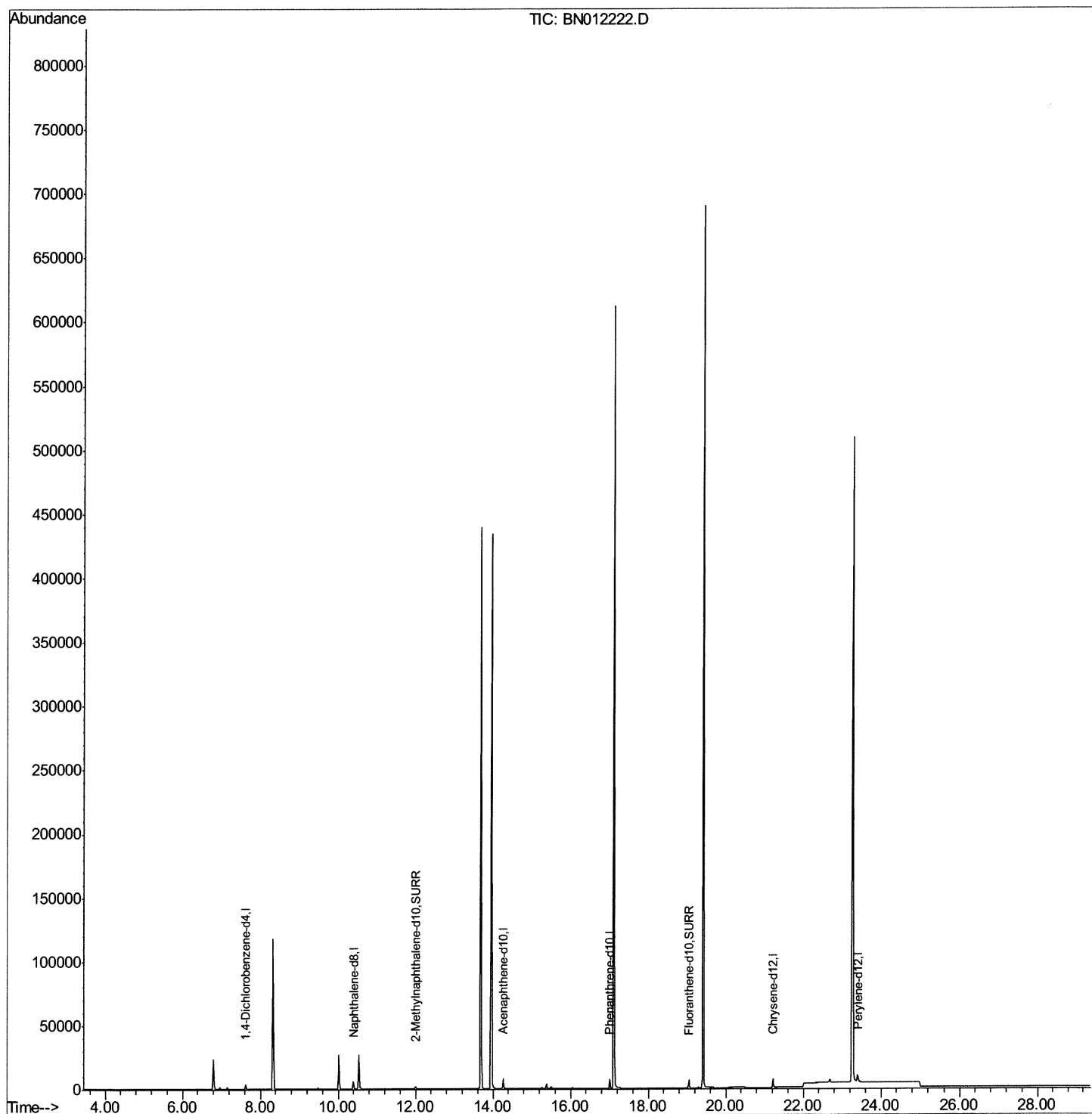
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101420\
Data File : BN012222.D
Acq On : 15 Oct 2020 09:35
Operator : CG/JU
Sample : PB132151BL
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK51

Manual Integrations
APPROVED

mohammad
10/19/2020 1:19:07 PM

Quant Time: Oct 15 10:50:59 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 15 10:47:40 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

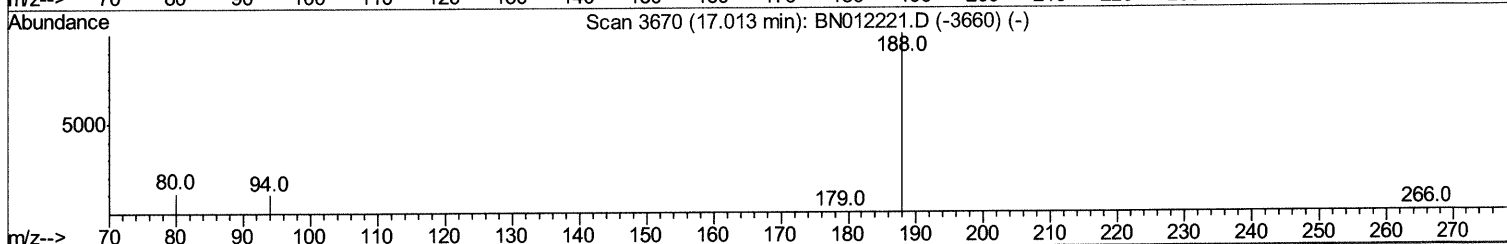
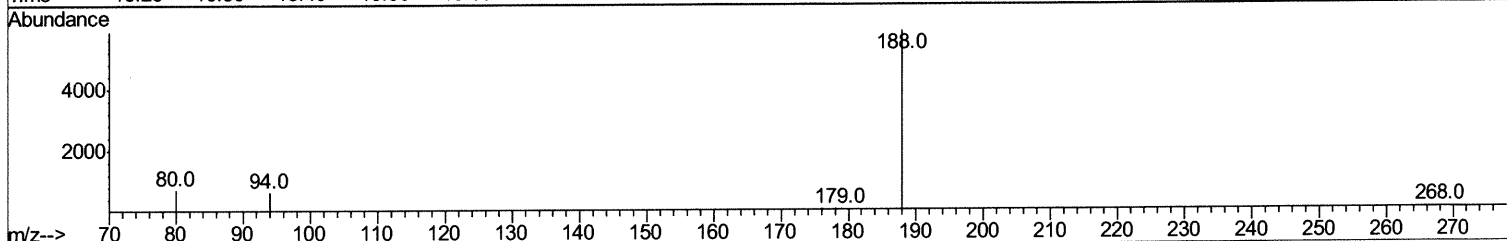
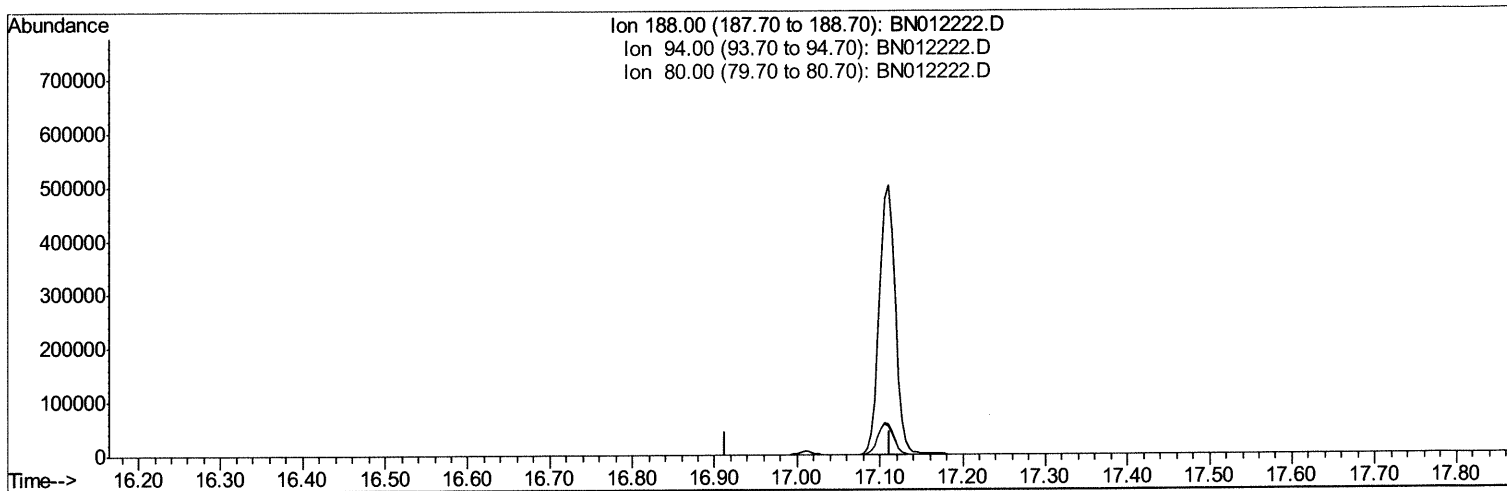
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 10/19/2020 1:19:07 PM

Quant Time: Oct 15 10:49:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 10:47:40 2020
 Response via : Initial Calibration



TIC: BN012222.D

(10) Phenanthrene-d10 (I)
 17.013min (-17.013) 0.00ng/ul
 response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	11.20	0.00#
80.00	12.20	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

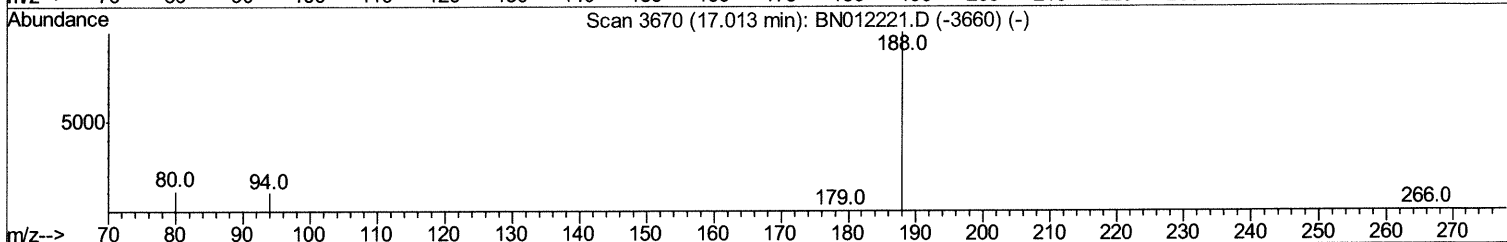
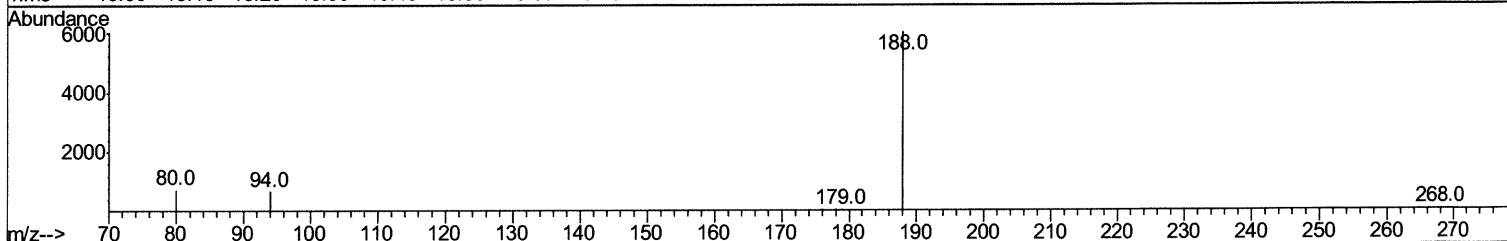
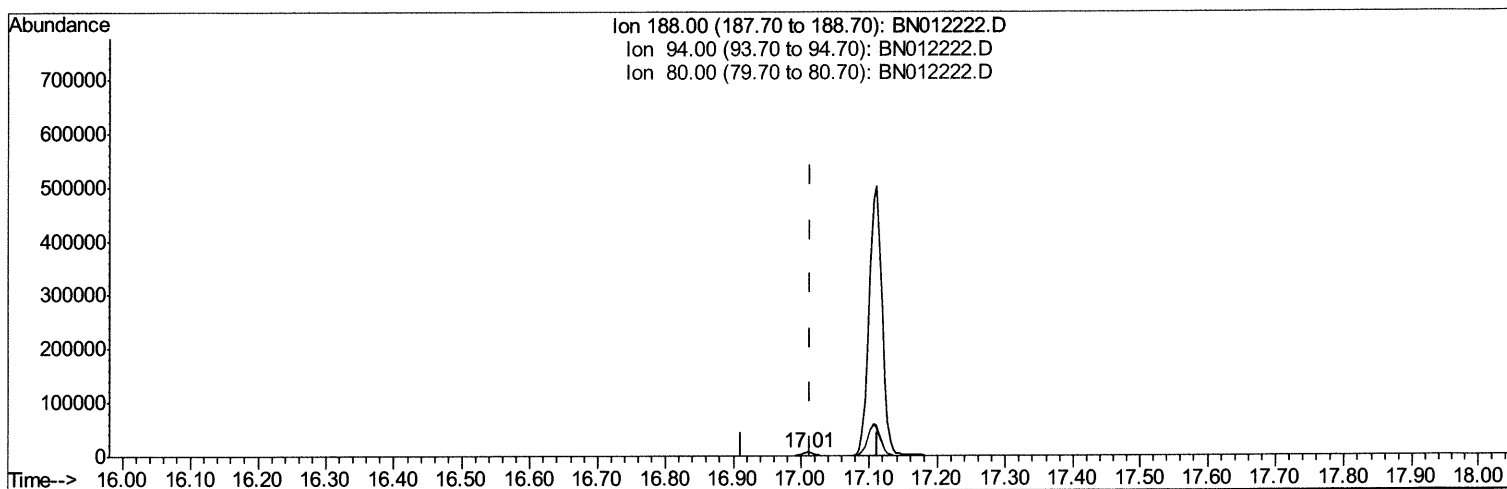
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Response via : Initial Calibration



TIC: BN012222.D

(10) Phenanthrene-d10 (I)

17.009min (-0.004) 0.40ng/ul m 10/20/20 JU

response 8630

Ion	Exp%	Act%
188.00	100	100
94.00	11.20	11.48
80.00	12.20	12.63
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101420\
Data File : BN012222.D
Acq On : 15 Oct 2020 09:35
Operator : CG/JU
Sample : PB132151BL
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Manual Integrations
APPROVED

mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1768	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7401	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4308	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8630m>	0.40	ng/ul>	0.00
16) Chrysene-d12	21.21	240	6859	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	7521	0.40	ng/ul	0.00

10/20/20 JU

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	12.00	152	3191	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	7167	0.30	ng/ul	0.00

Target Compounds Ovalue

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