

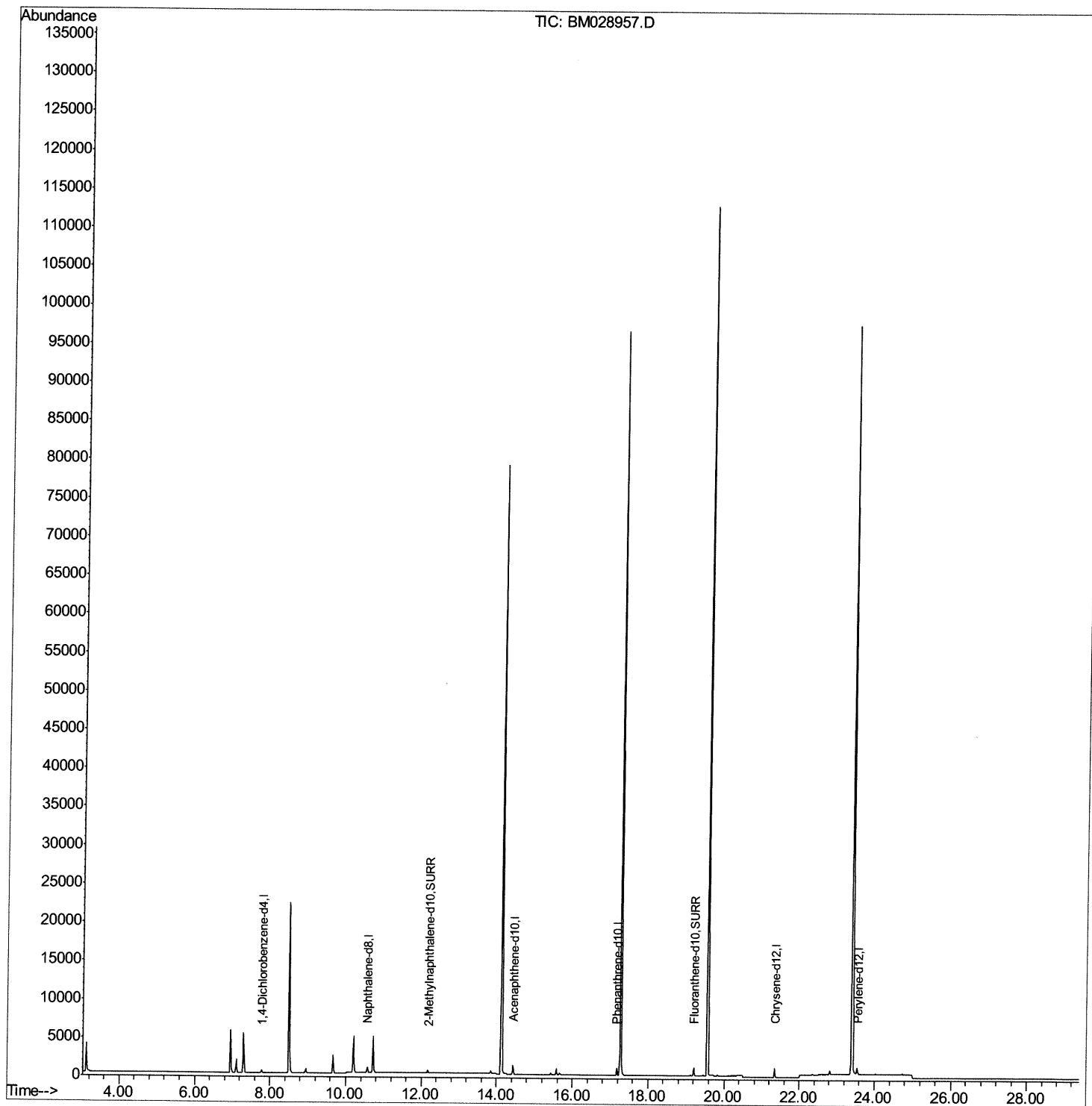
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030121\
Data File : BM028957.D
Acq On : 01 Mar 2021 10:23
Operator : CG/JU
Sample : PB134806BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK06

Manual Integrations
APPROVED

mohammad
3/1/2021 4:33:14 PM

Quant Time: Mar 01 10:55:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 01 10:25:15 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

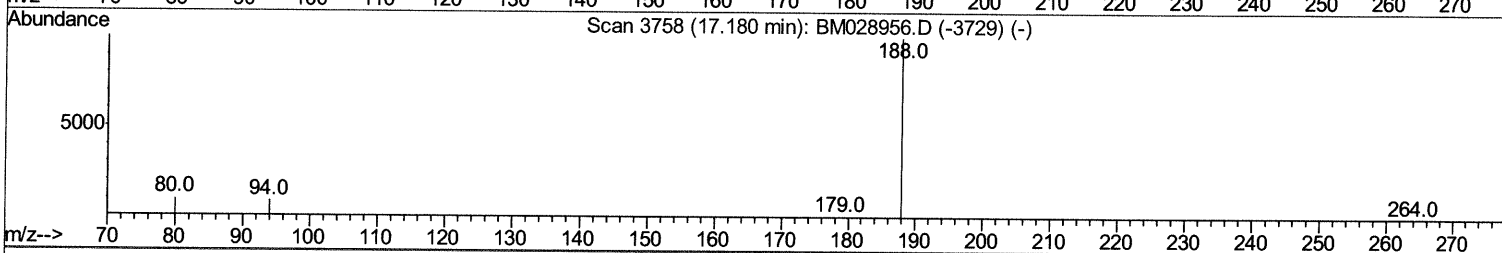
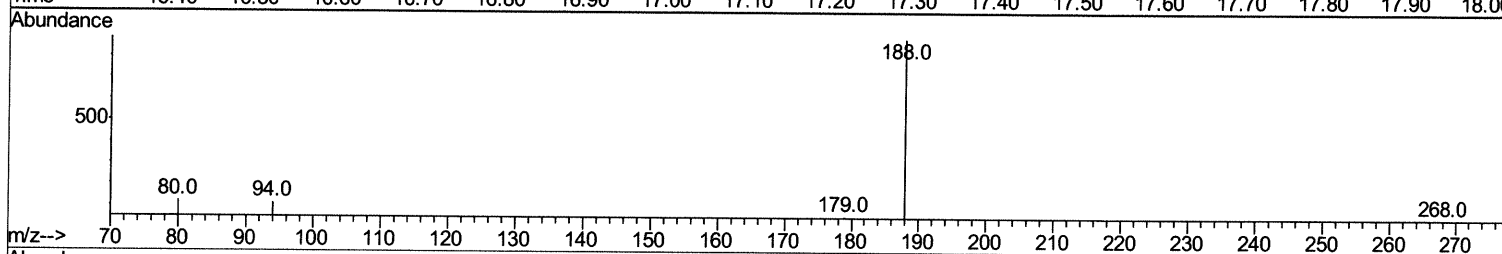
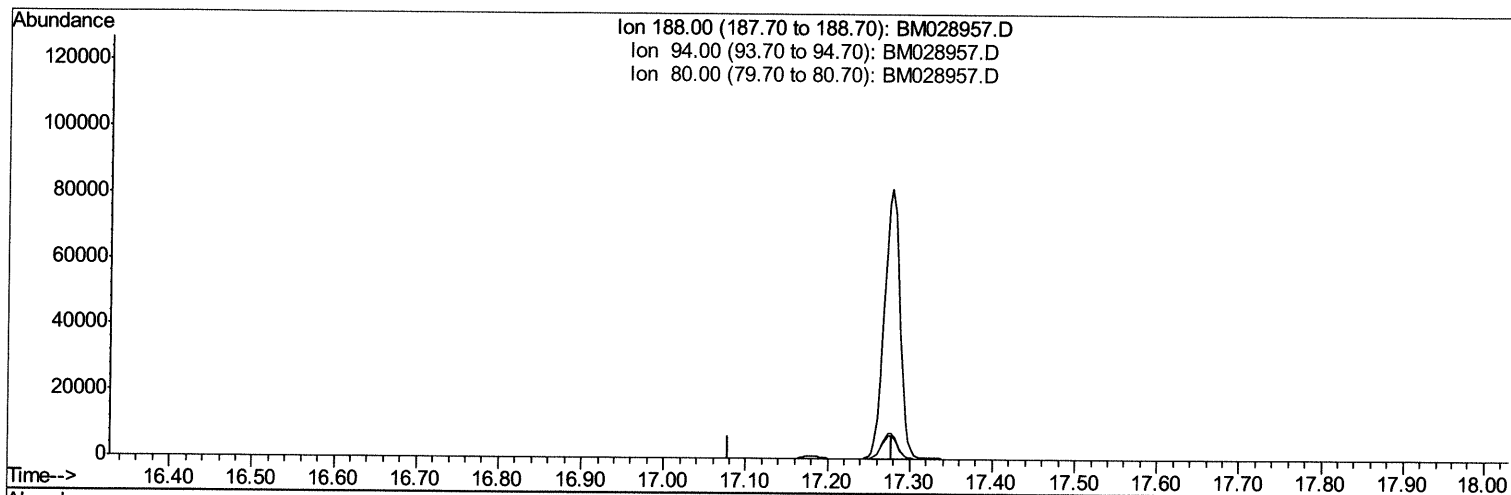
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Quant Time: Mar 01 10:53:40 2021
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TIC: BM028957.D

(10) Phenanthrene-d10 (I)

17.180min (-17.180) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	13.50	0.00#
80.00	14.00	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

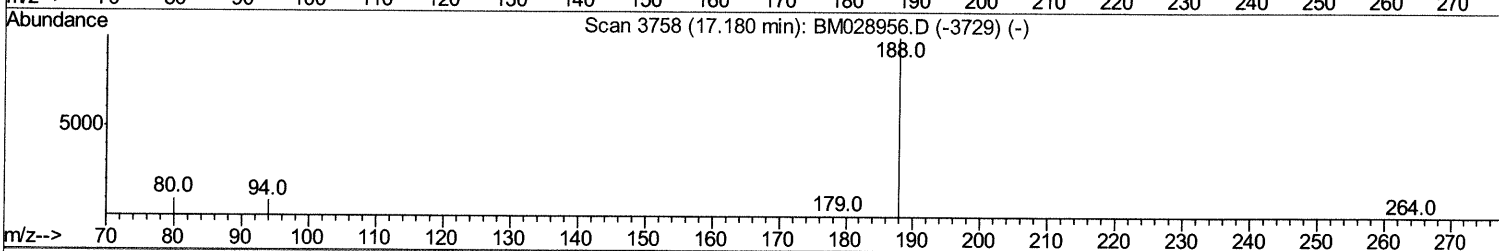
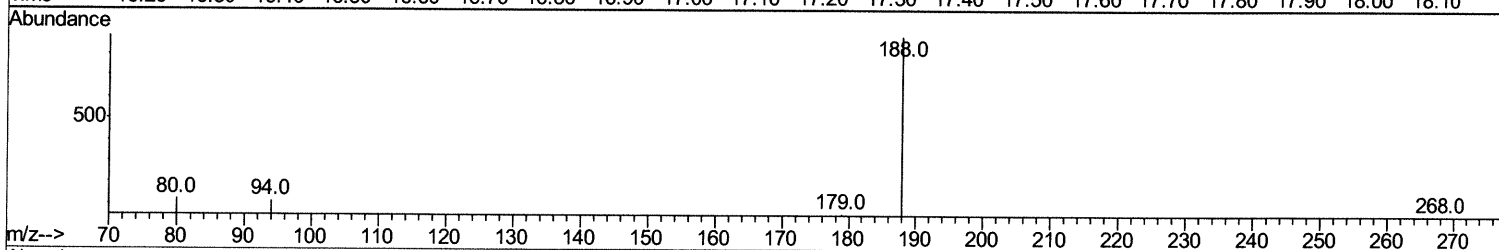
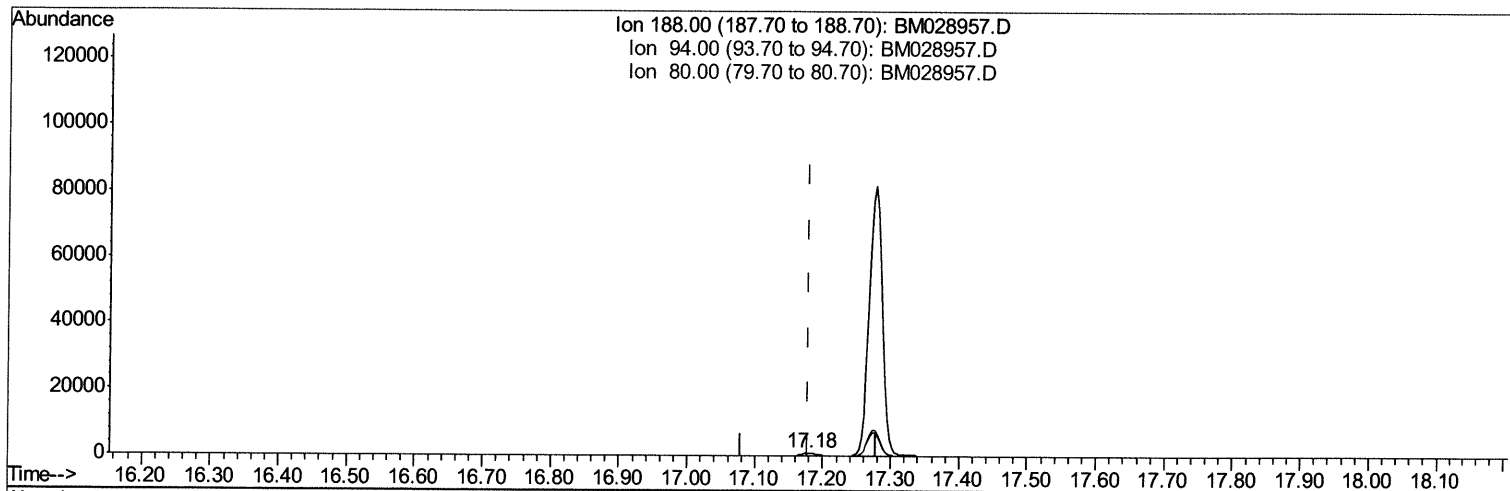
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BNA_M
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Manual Integrations
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3/1/2021 4:33:14 PM

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Response via : Initial Calibration



TIC: BM028957.D

(10) Phenanthrene-d10 (I)

17.180min (+0.000) 0.40ng/ul m 03/02/21 JU

response 1295

Ion	Exp%	Act%
188.00	100	100
94.00	13.50	12.64
80.00	14.00	13.66
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030121\
 Data File : BM028957.D
 Acq On : 01 Mar 2021 10:23
 Operator : CG/JU
 Sample : PB134806BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK06

Manual Integrations
 APPROVED

mohammad
 3/1/2021 4:33:14 PM

Quant Time: Mar 01 10:55:15 2021
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 QLast Update : Mon Mar 01 10:25:15 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	269	0.40	ng/ul	0.00
2) Naphthalene-d8	10.58	136	1079	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.44	164	626	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.18	188	1295m >	0.40	ng/ul >	0.00
16) Chrysene-d12	21.35	240	1095	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	1182	0.40	ng/ul	0.00

03/02/21

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.18	152	587	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.20	212	1153	0.34	ng/ul	0.00

Target Compounds	Ovalue

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