

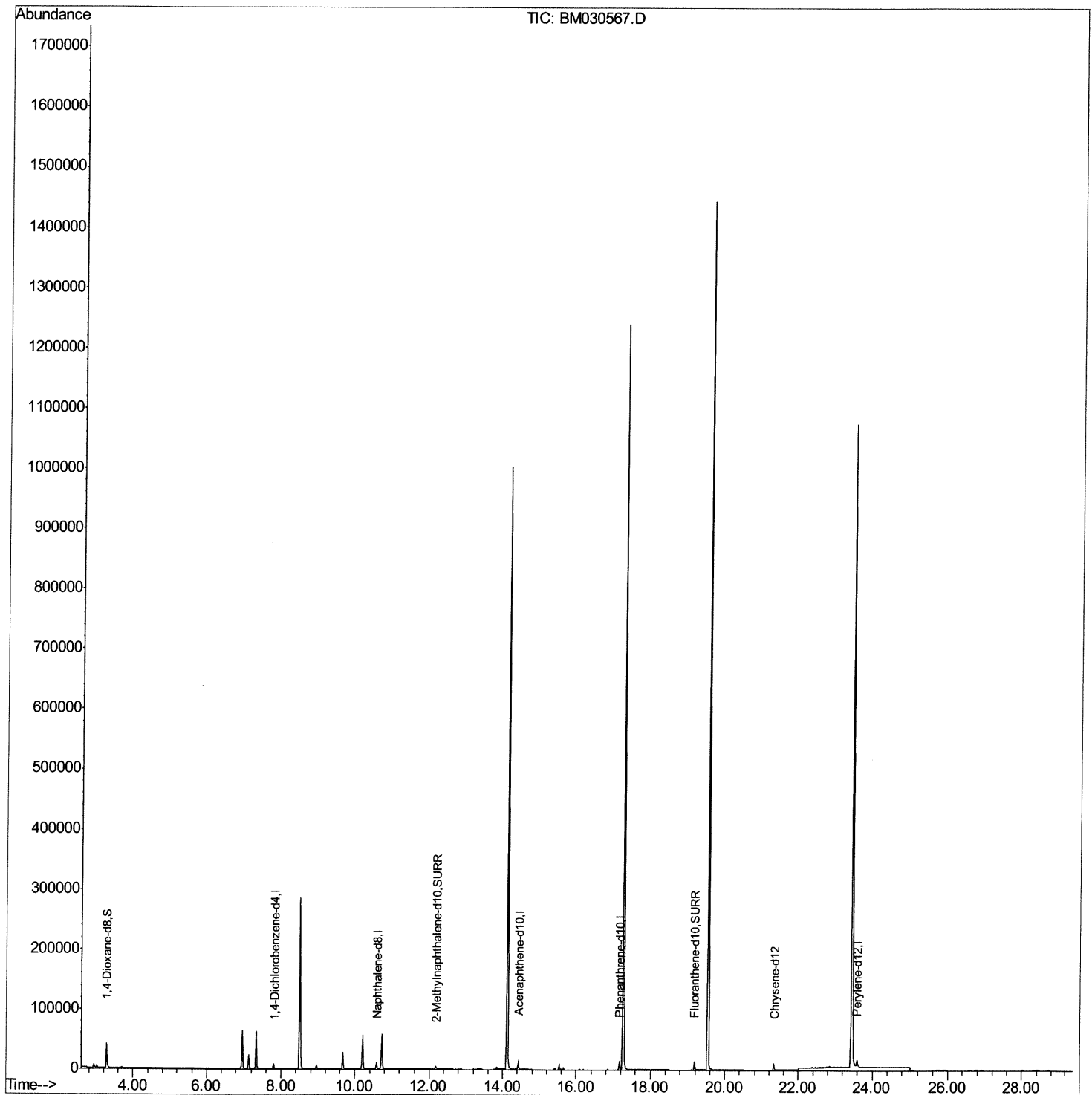
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062421\
Data File : BM030567.D
Acq On : 24 Jun 2021 09:31
Operator : CG/JU
Sample : PB137202BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK202

Manual Integrations
APPROVED

mohammad
6/24/2021 2:55:42 PM

Quant Time: Jun 24 10:24:09 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 24 01:55:26 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

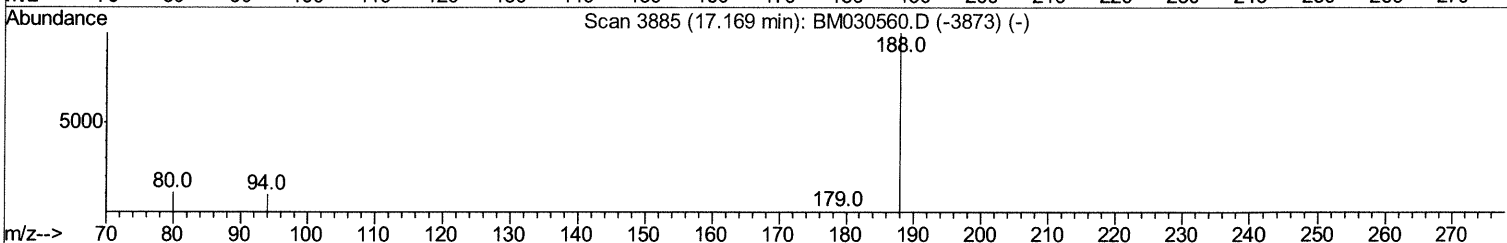
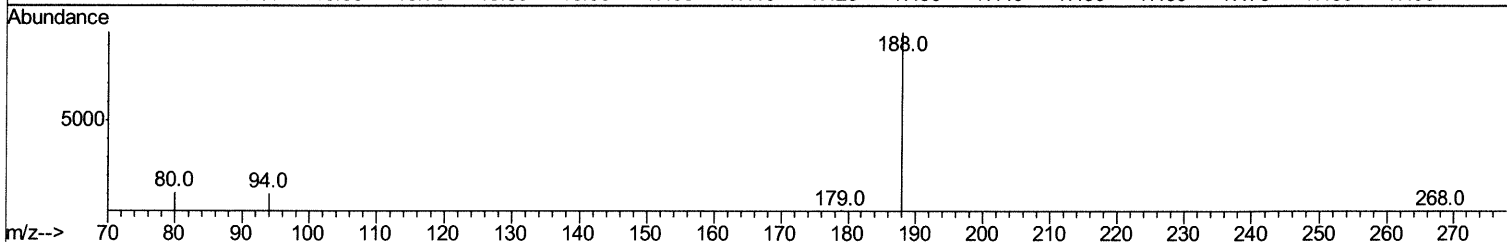
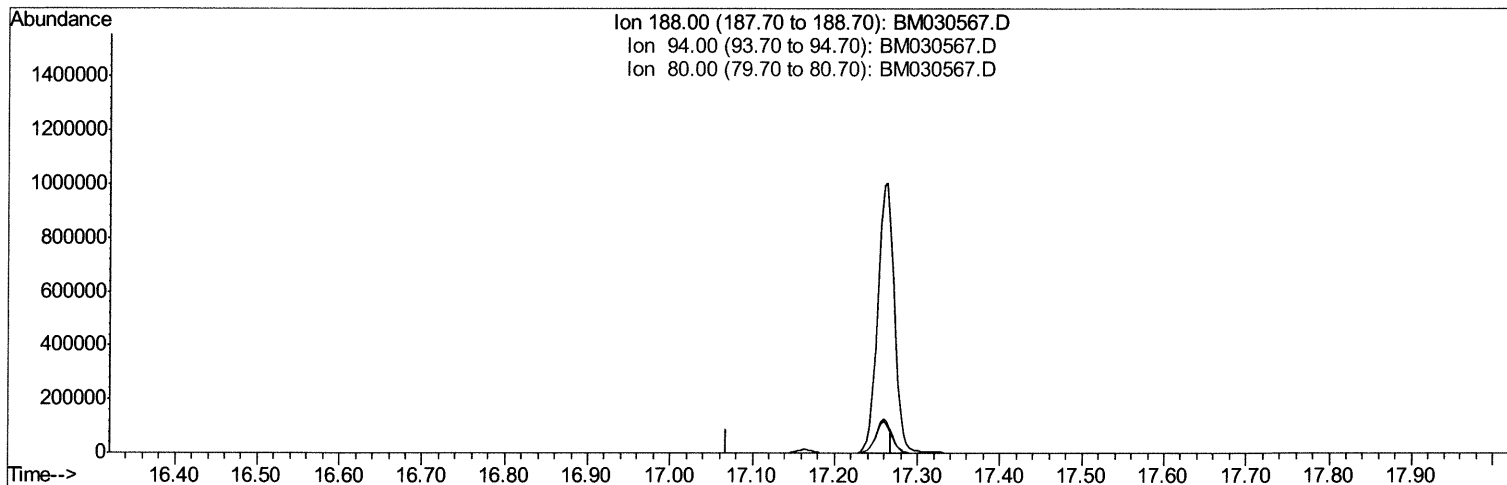
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Quant Time: Jun 24 10:12:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 24 01:55:26 2021
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TIC: BM030567.D

(13) Phenanthrene-d10 (I)

17.169min (-17.169) 0.00ng/ul

response 0

Ion	Exp%	Act%
188.00	100	0.00
94.00	10.30	0.00#
80.00	11.60	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

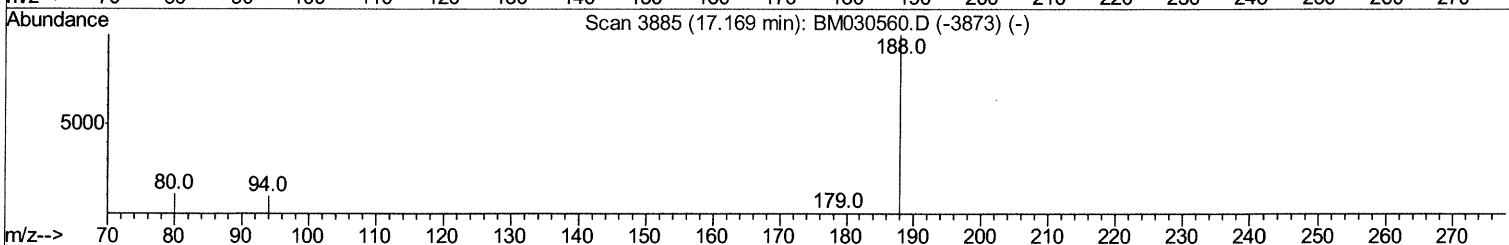
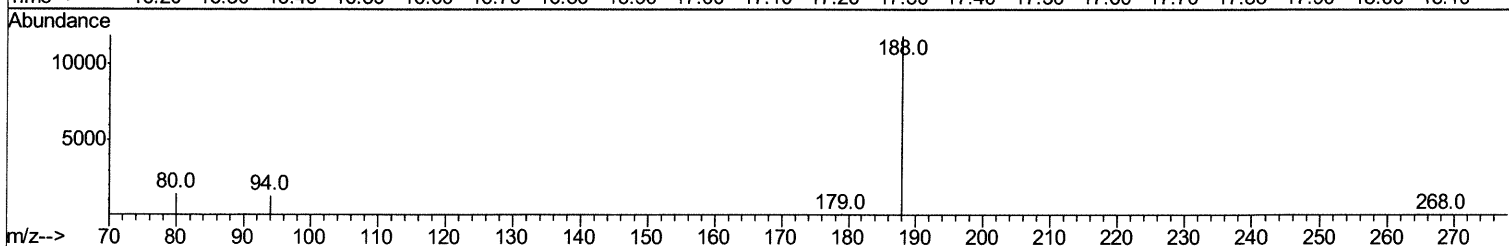
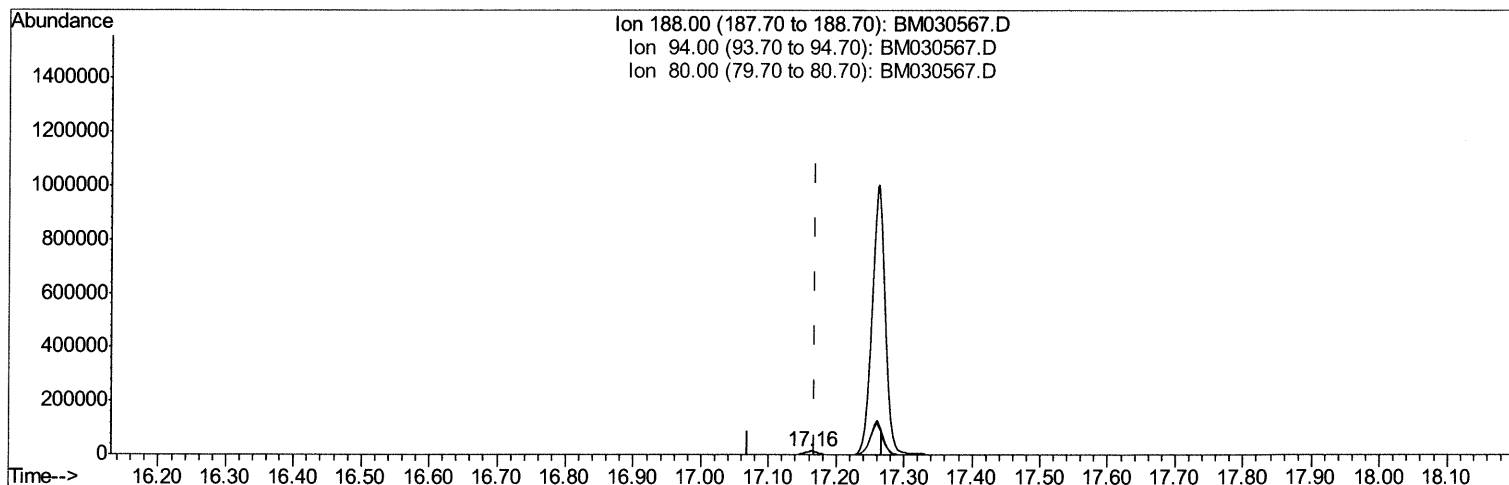
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Manual Integrations
 APPROVED

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 Response via : Initial Calibration



TIC: BM030567.D

(13) Phenanthrene-d10 (I)

17.162min (-0.007) 0.40ng/ul m

06/28/21 JU

response 17089

Ion	Exp%	Act%
188.00	100	100
94.00	10.30	10.91
80.00	11.60	12.02
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062421\
Data File : BM030567.D
Acq On : 24 Jun 2021 09:31
Operator : CG/JU
Sample : PB137202BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK202

Manual Integrations
APPROVED

mohammad
6/24/2021 2:55:42 PM

Quant Time: Jun 24 10:24:09 2021
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	3941	0.40	ng/ul	0.00
4) Naphthalene-d8	10.59	136	15218	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.43	164	9146	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.16	188	17089m	0.40	ng/ul	0.00
17) Chrysene-d12	21.33	240	12888	0.40	ng/ul	0.00
23) Perylene-d12	23.59	264	16712	0.40	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	25731	6.10	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.18	152	8174	0.34	ng/ul	0.00
18) Fluoranthene-d10	19.19	212	15793	0.40	ng/ul	0.00

Target Compounds Ovalue

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

06/28/21JU