

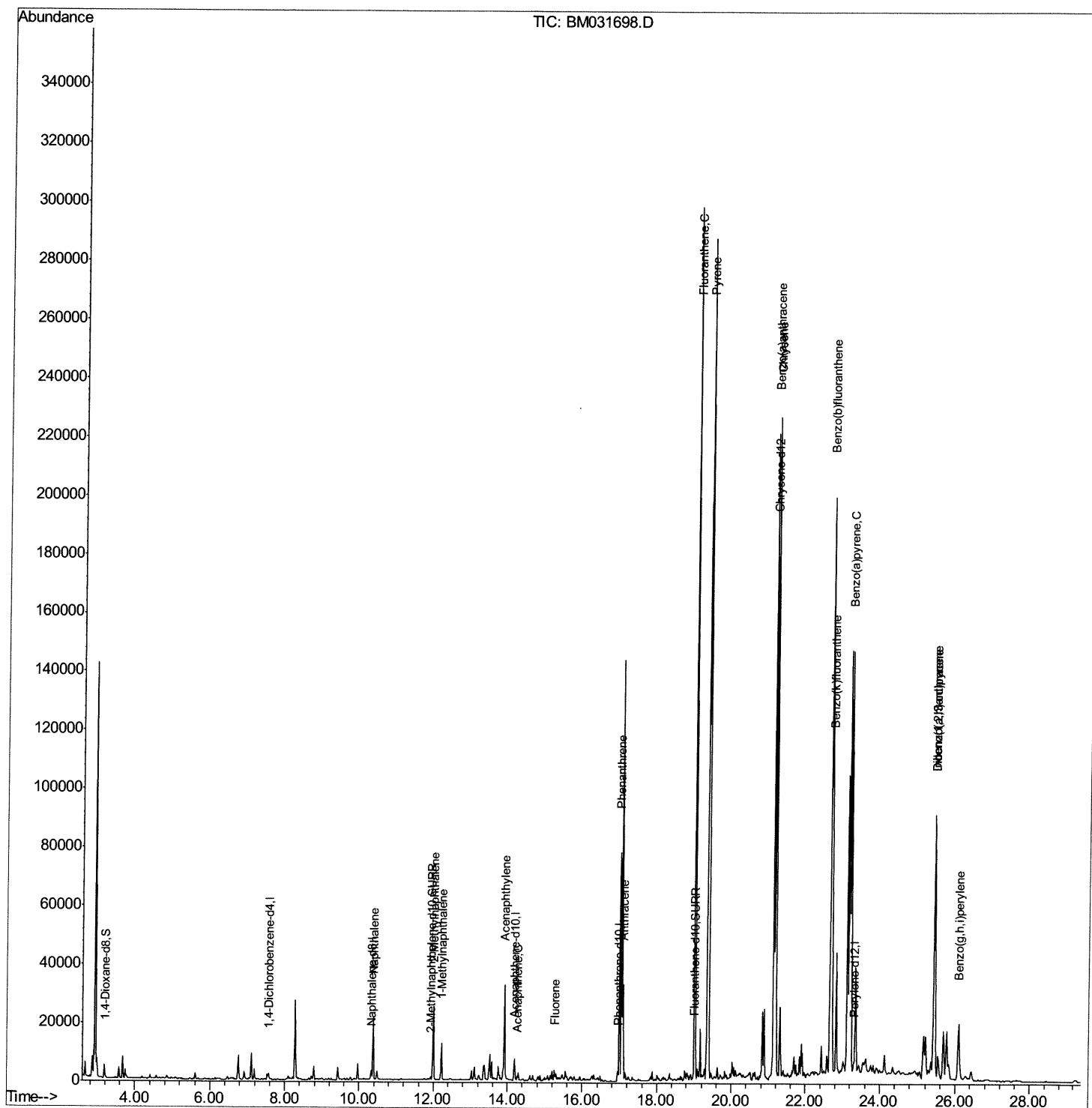
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
Data File : BM031698.D  
Acq On : 13 Aug 2021 04:41  
Operator : CG/JU  
Sample : M3182-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C00C8

Manual Integrations  
APPROVED

mohammad  
8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:16:36 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Aug 12 17:28:09 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

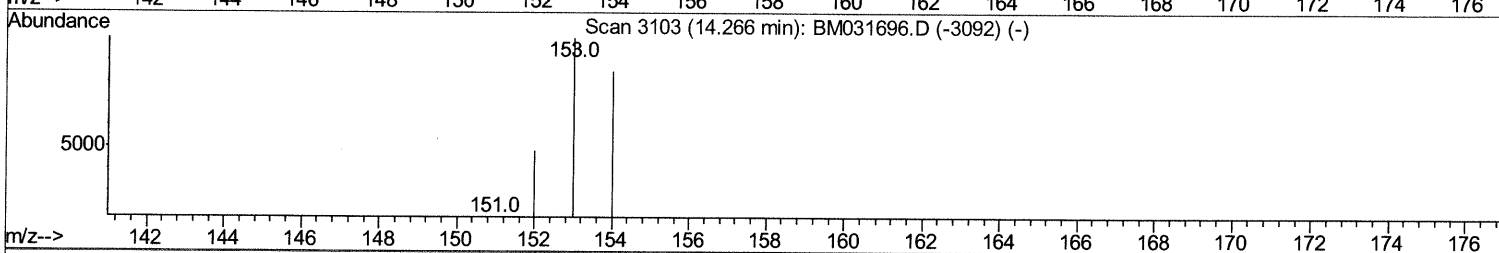
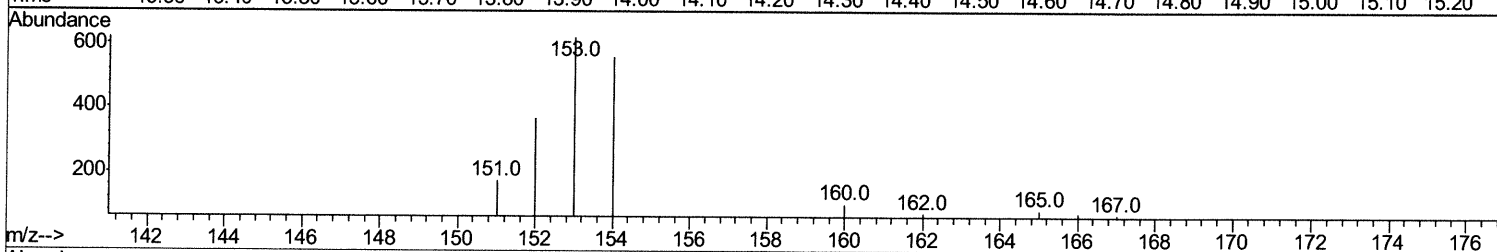
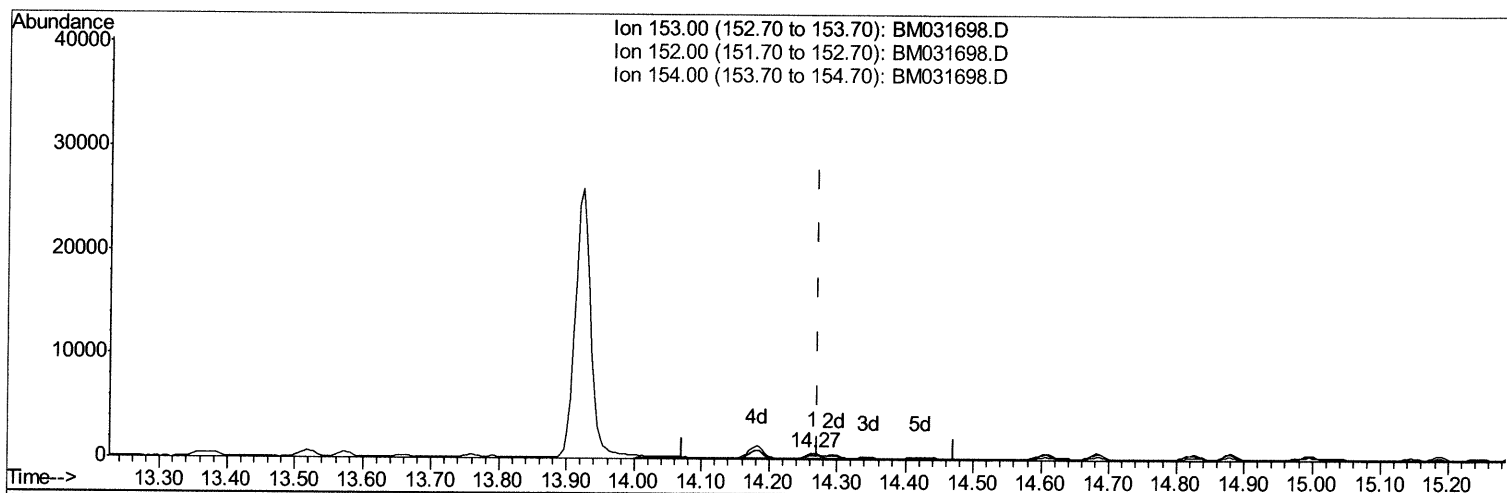
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TIC: BM031698.D

(11) Acenaphthene (C)

14.265min (-0.007) 0.09ng/ul

response 849

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 153.00 | 100   | 100   |
| 152.00 | 50.90 | 59.03 |
| 154.00 | 87.70 | 90.32 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

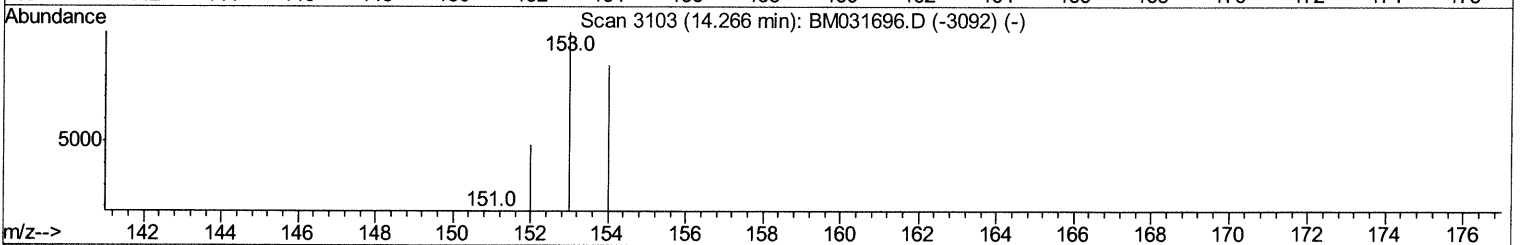
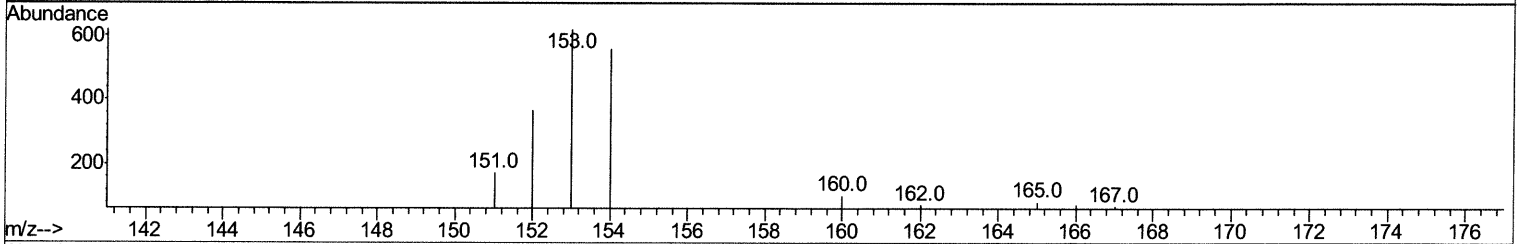
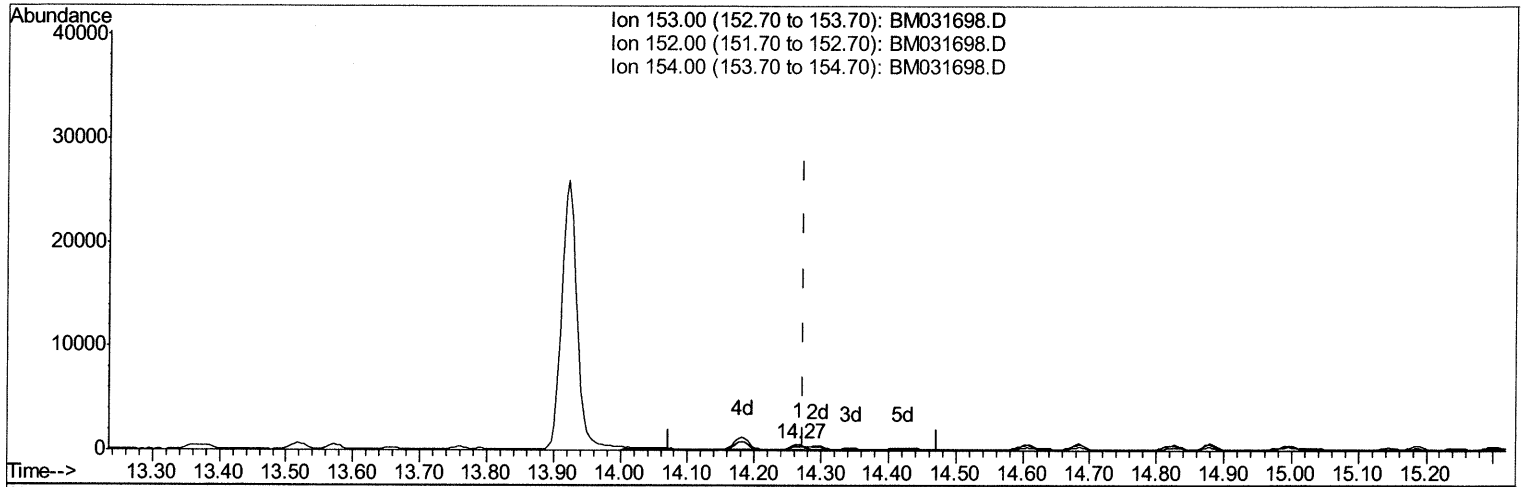
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
 Data File : BM031698.D  
 Acq On : 13 Aug 2021 04:41  
 Operator : CG/JU  
 Sample : M3182-10  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 C00C8

**Manual Integrations**  
**APPROVED**

mohammad  
 8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:13:10 2021  
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TIC: BM031698.D

(11) Acenaphthene (C)

14.265min (-0.007) 0.14ng/ul m *08/16/21JU*

response 1261

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 153.00 | 100   | 100   |
| 152.00 | 50.90 | 59.03 |
| 154.00 | 87.70 | 90.32 |
| 0.00   | 0.00  | 0.00  |

# Quantitation Report (Qedit)

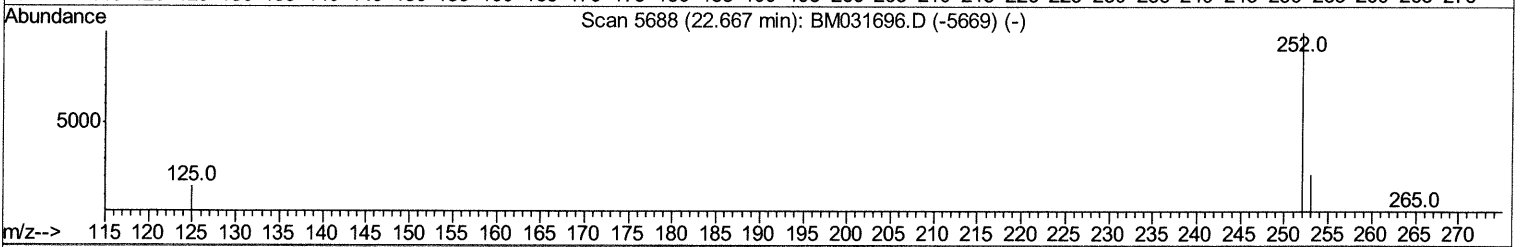
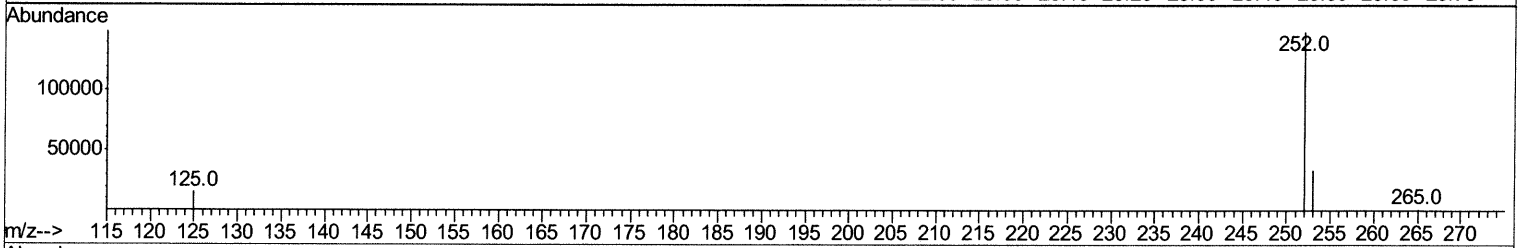
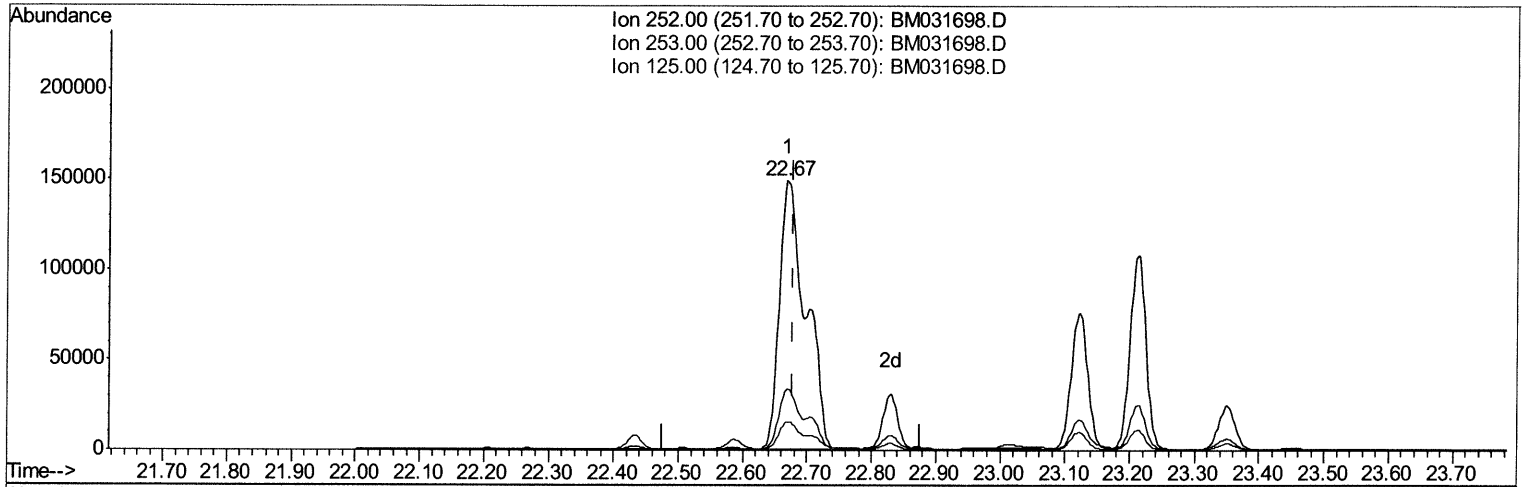
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
 Data File : BM031698.D  
 Acq On : 13 Aug 2021 04:41  
 Operator : CG/JU  
 Sample : M3182-10  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C00C8

Manual Integrations  
 APPROVED

mohammad  
 8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:14:32 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M  
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TIC: BM031698.D

(24) Benzo(b)fluoranthene

22.667min (-0.009) 21.63ng/ul

response 420162

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 26.50 | 22.49 |
| 125.00 | 11.40 | 10.32 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

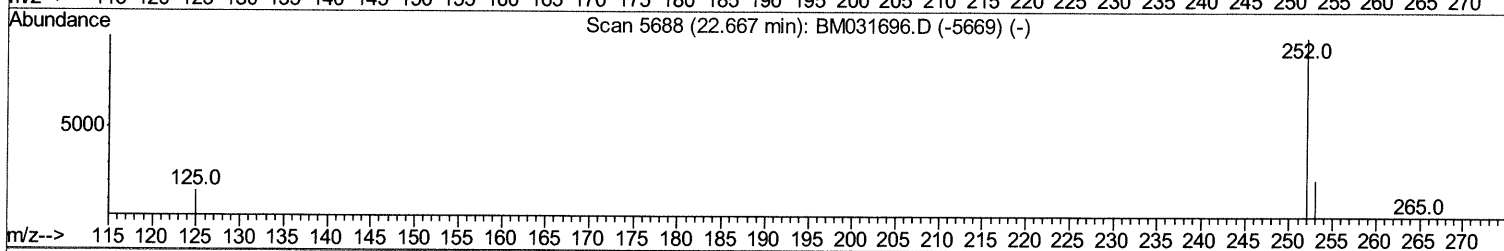
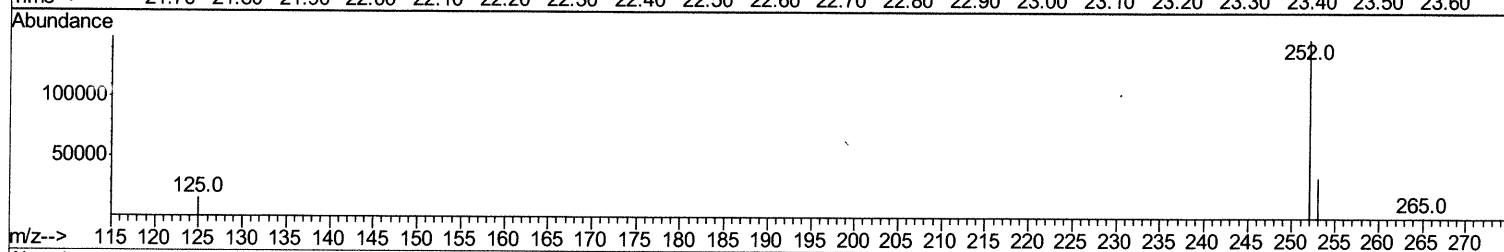
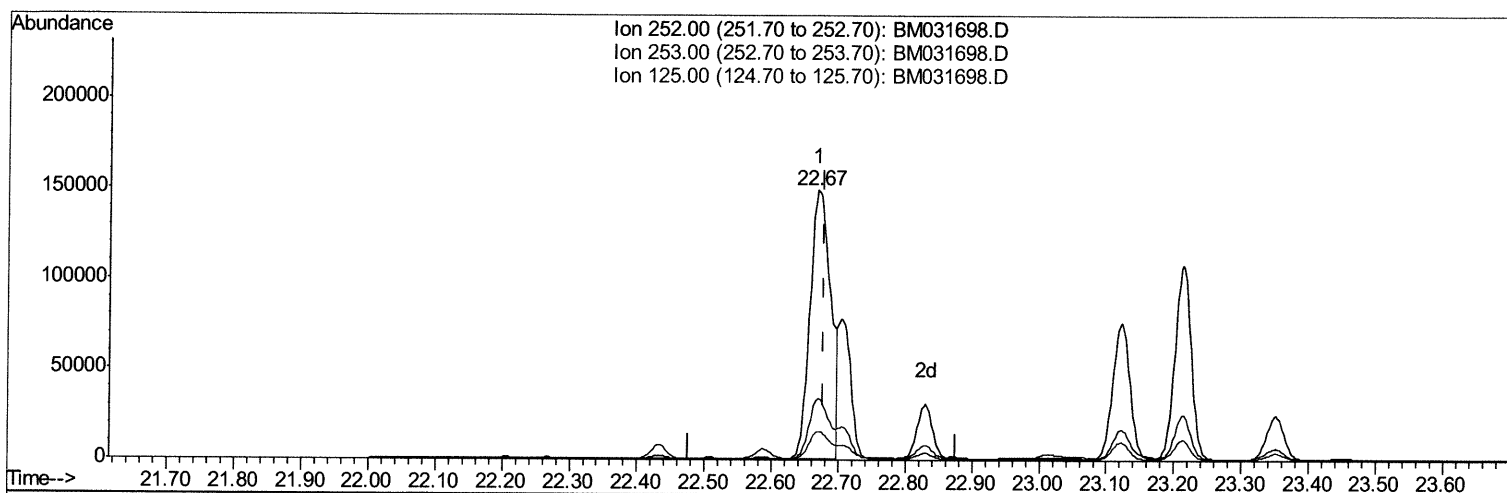
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
Data File : BM031698.D  
Acq On : 13 Aug 2021 04:41  
Operator : CG/JU  
Sample : M3182-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
C00C8

Manual Integrations  
APPROVED

mohammad  
8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:14:32 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Aug 12 17:28:09 2021  
Response via : Initial Calibration



TIC: BM031698.D

(24) Benzo(b)fluoranthene

22.667min (-0.009) 16.30ng/ul m

08/16/21 JU

response 316643

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 26.50 | 22.49 |
| 125.00 | 11.40 | 10.32 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

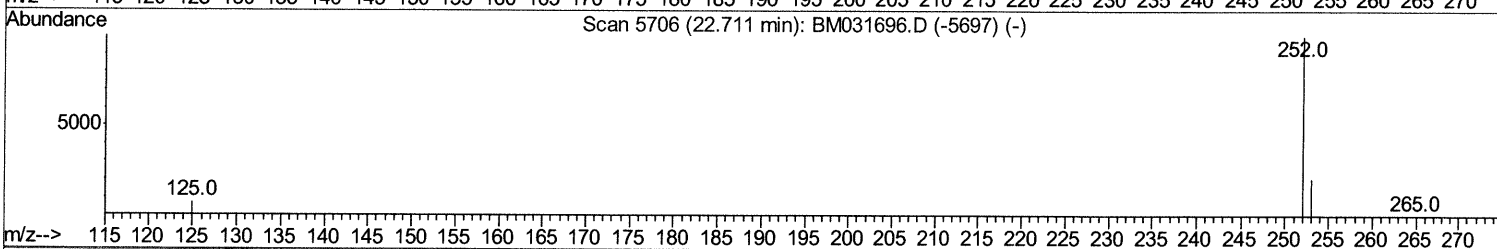
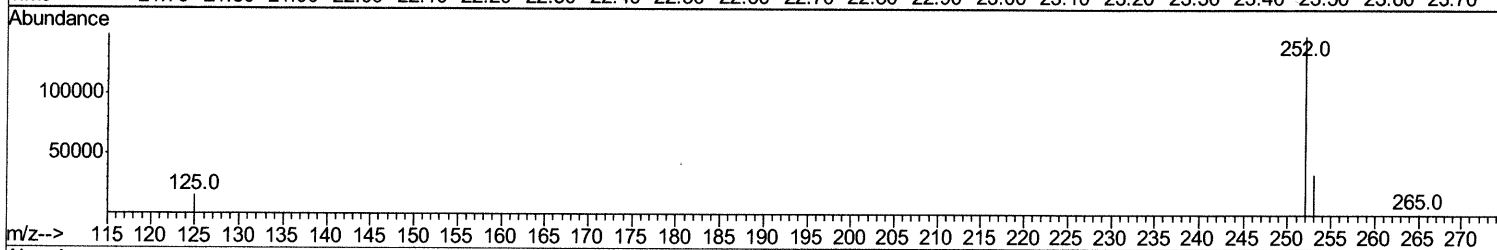
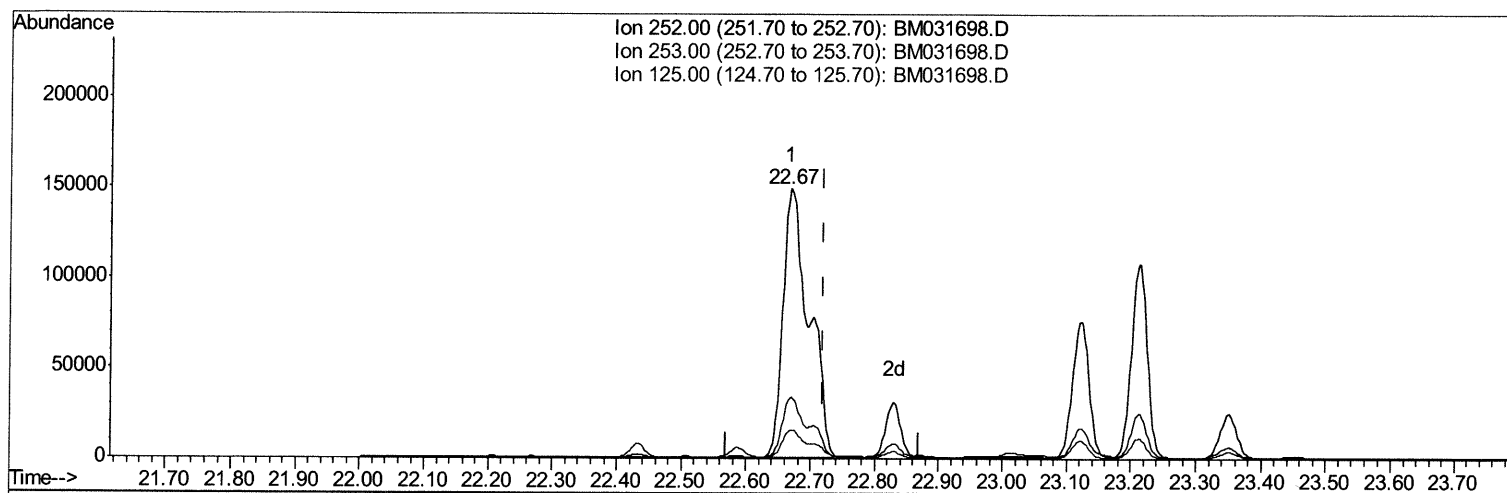
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
Data File : BM031698.D  
Acq On : 13 Aug 2021 04:41  
Operator : CG/JU  
Sample : M3182-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C00C8

Manual Integrations  
APPROVED

mohammad  
8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:15:26 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Aug 12 17:28:09 2021  
Response via : Initial Calibration



TIC: BM031698.D

(25) Benzo(k)fluoranthene

22.667min (-0.053) 20.93ng/ul

response 420218

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 26.60 | 22.49 |
| 125.00 | 11.00 | 10.32 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

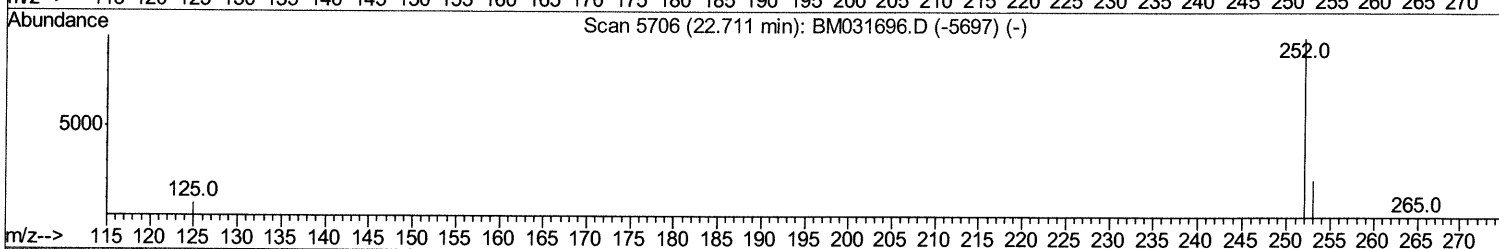
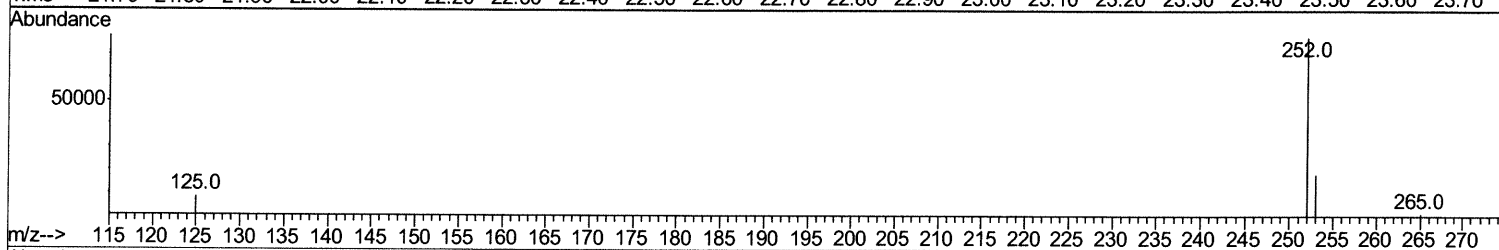
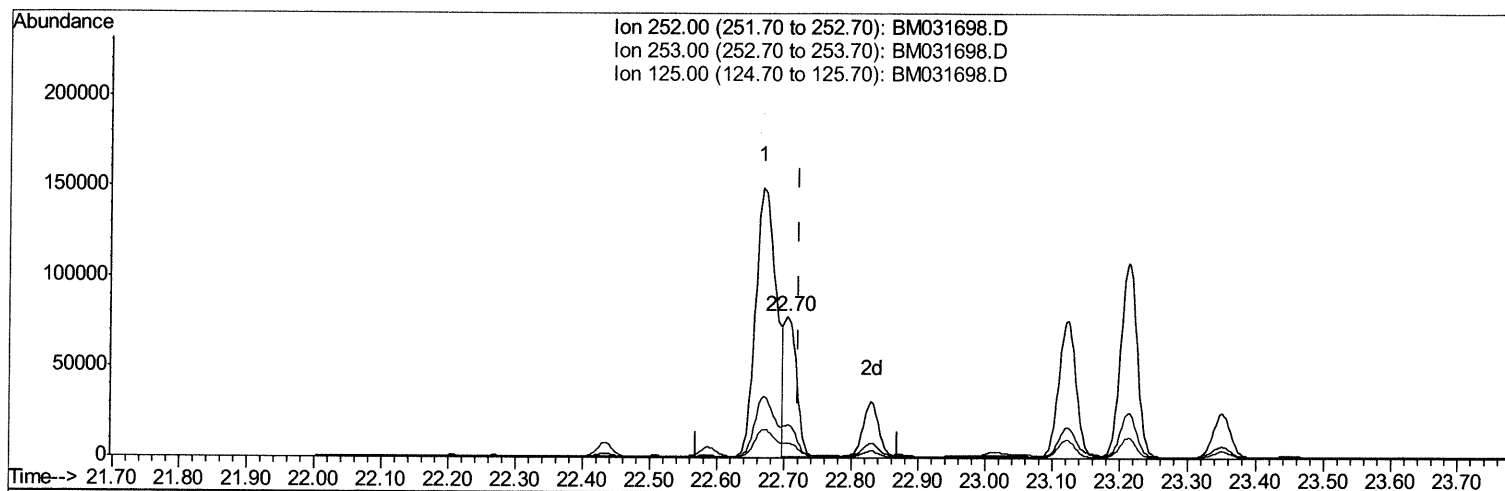
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
Data File : BM031698.D  
Acq On : 13 Aug 2021 04:41  
Operator : CG/JU  
Sample : M3182-10  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C00C8

Manual Integrations  
APPROVED

mohammad  
8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:15:26 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM072821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Aug 12 17:28:09 2021  
Response via : Initial Calibration



TIC: BM031698.D

(25) Benzo(k)fluoranthene

22.704min (-0.017) 5.11ng/ul m

08/16/21 JU

response 102615

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 26.60 | 23.64 |
| 125.00 | 11.00 | 10.50 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM081221\  
 Data File : BM031698.D  
 Acq On : 13 Aug 2021 04:41  
 Operator : CG/JU  
 Sample : M3182-10  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C00C8

Manual Integrations  
 APPROVED

mohammad  
 8/13/2021 5:15:01 PM

Quant Time: Aug 13 05:16:36 2021  
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| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.57  | 152  | 1090     | 0.40 | ng/ul | 0.00     |
| 4) Naphthalene-d8         | 10.32 | 136  | 3983     | 0.40 | ng/ul | 0.00     |
| 9) Acenaphthene-d10       | 14.20 | 164  | 2548     | 0.40 | ng/ul | 0.00     |
| 13) Phenanthrene-d10      | 16.95 | 188  | 4592     | 0.40 | ng/ul | -0.01    |
| 17) Chrysene-d12          | 21.15 | 240  | 4034     | 0.40 | ng/ul | 0.00     |
| 23) Perylene-d12          | 23.30 | 264  | 4312     | 0.40 | ng/ul | -0.02    |

## System Monitoring Compounds

|                            |       |     |      |      |       |       |
|----------------------------|-------|-----|------|------|-------|-------|
| 3) 1,4-Dioxane-d8          | 3.18  | 96  | 2945 | 2.43 | ng/ul | 0.00  |
| 6) 2-Methylnaphthalene-d10 | 11.92 | 152 | 1300 | 0.19 | ng/ul | -0.01 |
| 18) Fluoranthene-d10       | 18.99 | 212 | 2749 | 0.20 | ng/ul | -0.01 |

## Target Compounds

|                            |       |     |         |        | Ovalue |               |
|----------------------------|-------|-----|---------|--------|--------|---------------|
| 5) Naphthalene             | 10.37 | 128 | 28105   | 2.369  | ng/ul  | 97            |
| 7) 2-Methylnaphthalene     | 12.00 | 142 | 17600   | 2.171  | ng/ul  | 99            |
| 8) 1-Methylnaphthalene     | 12.22 | 142 | 8490    | 1.060  | ng/ul  | 99            |
| 10) Acenaphthylene         | 13.92 | 152 | 39363   | 3.627  | ng/ul  | 97            |
| 11) Acenaphthene           | 14.27 | 153 | 1261m   | 0.138  | ng/ul  | > 08/16/21 JU |
| 12) Fluorene               | 15.26 | 166 | 2657    | 0.253  | ng/ul# | 87            |
| 15) Phenanthrene           | 16.99 | 178 | 79459   | 5.449  | ng/ul  | 97            |
| 16) Anthracene             | 17.09 | 178 | 32773   | 2.401  | ng/ul  | 98            |
| 19) Fluoranthene           | 19.02 | 202 | 246658  | 14.319 | ng/ul  | 97            |
| 20) Pyrene                 | 19.38 | 202 | 238355  | 13.643 | ng/ul  | 97            |
| 21) Benzo(a)anthracene     | 21.13 | 228 | 180274  | 11.067 | ng/ul  | 96            |
| 22) Chrysene               | 21.19 | 228 | 215570  | 12.791 | ng/ul  | 98            |
| 24) Benzo(b)fluoranthene   | 22.67 | 252 | 316643m | 16.298 | ng/ul  | > 08/16/21 JU |
| 25) Benzo(k)fluoranthene   | 22.70 | 252 | 102615m | 5.112  | ng/ul  | >             |
| 26) Benzo(a)pyrene         | 23.21 | 252 | 184364  | 10.836 | ng/ul# | 89            |
| 27) Indeno(1,2,3-cd)pyrene | 25.43 | 276 | 136005  | 6.184  | ng/ul# | 99            |
| 28) Dibenzo(a,h)anthracene | 25.44 | 278 | 33467   | 1.888  | ng/ul  | 98            |
| 29) Benzo(a,h,i)perylene   | 26.09 | 276 | 38005   | 2.026  | ng/ul  | 98            |

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