

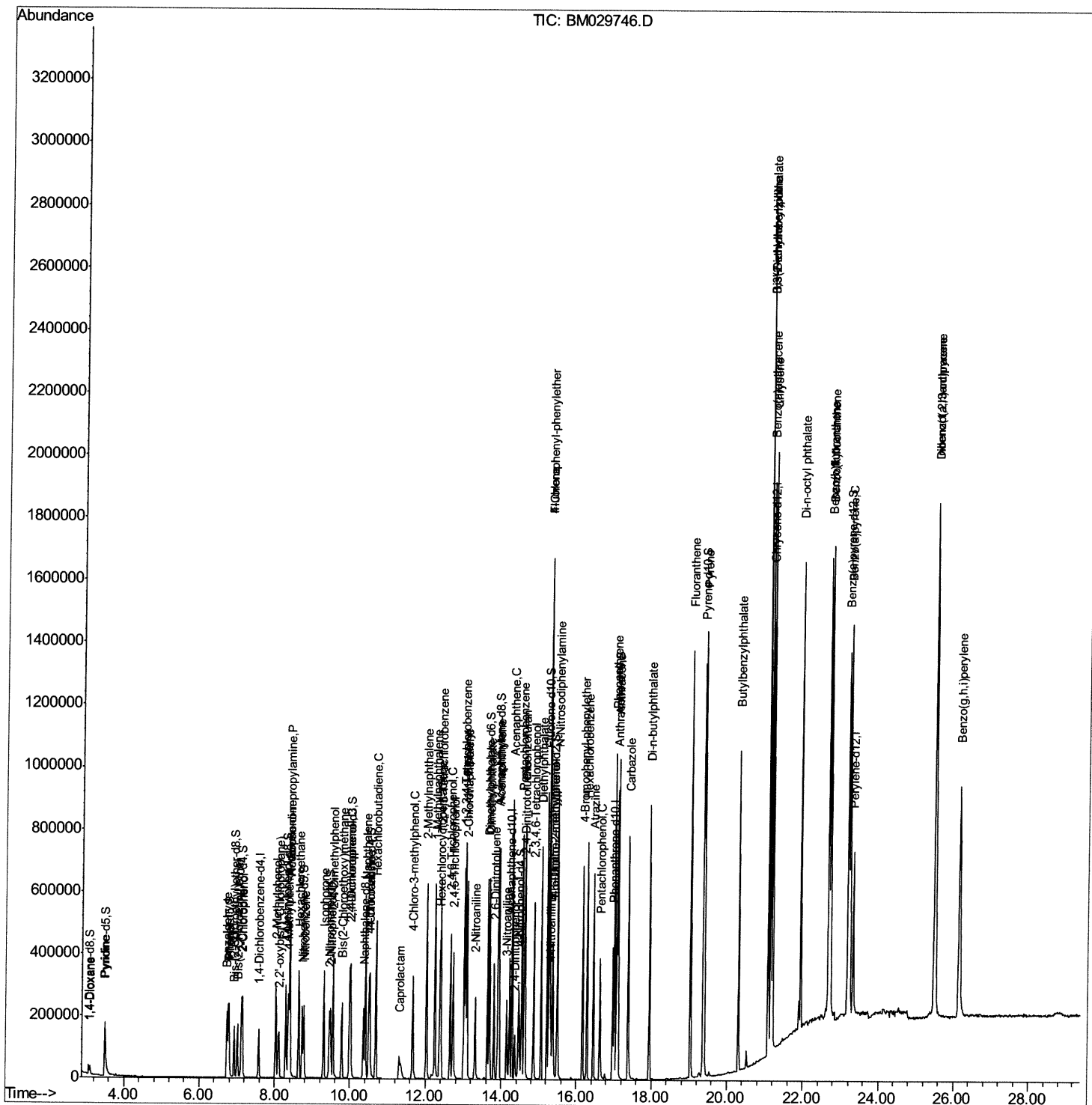
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Data Path   : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
Data File  : BM029746.D  
Acq On     : 01 May 2021   14:00  
Operator   : CG/JU  
Sample     : SSTD04012  
Misc       :  
ALS Vial   : 5      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**ClientSampleId :**  
SSTD040012

**Manual Integrations**  
**APPROVED**

mohammad  
5/3/2021 12:24:21 PM

Quant Time: May 01 14:35:46 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050121.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat May 01 14:02:00 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

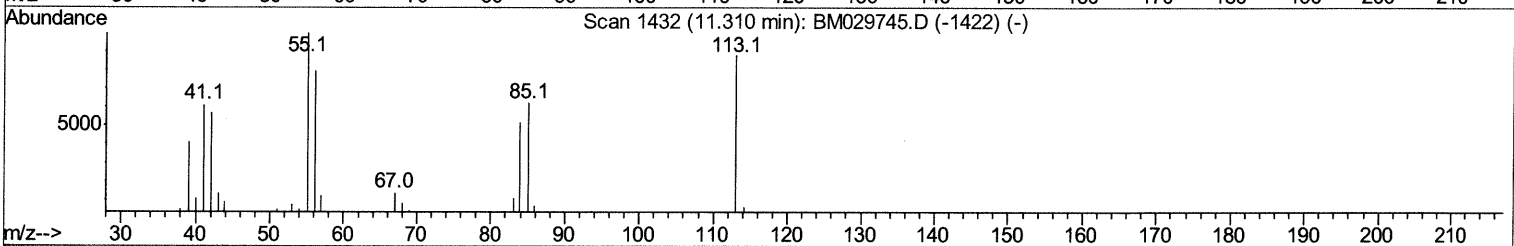
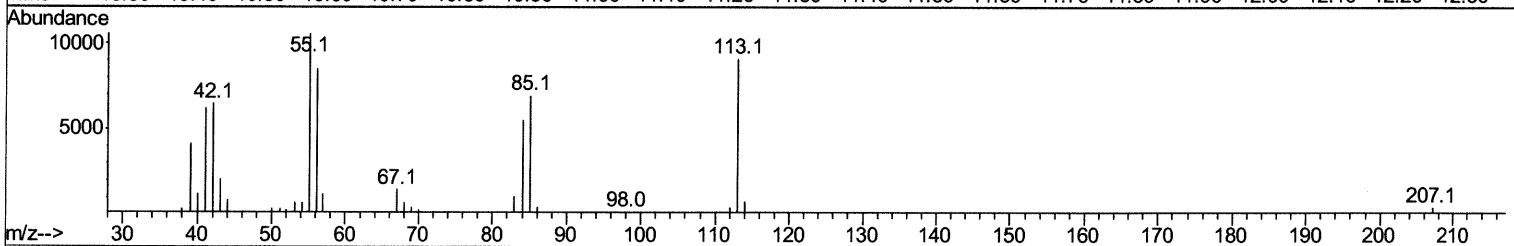
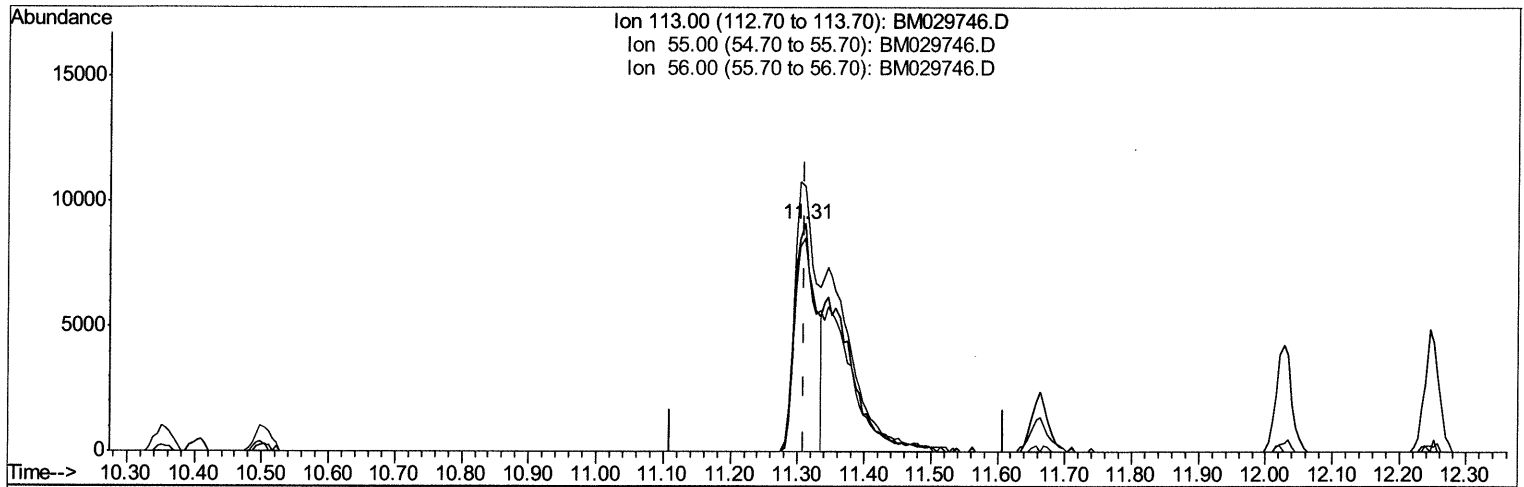
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
 Data File : BM029746.D  
 Acq On : 01 May 2021 14:00  
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**Instrument :**  
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**Manual Integrations**  
**APPROVED**

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 5/3/2021 12:24:21 PM

Quant Time: May 01 14:33:49 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050121.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 01 14:02:00 2021  
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TIC: BM029746.D

(34) Caprolactam

11.310min (-0.000) 24.67ng/ul

response 19318

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 113.50 | 116.03 |
| 56.00  | 90.40  | 93.56  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

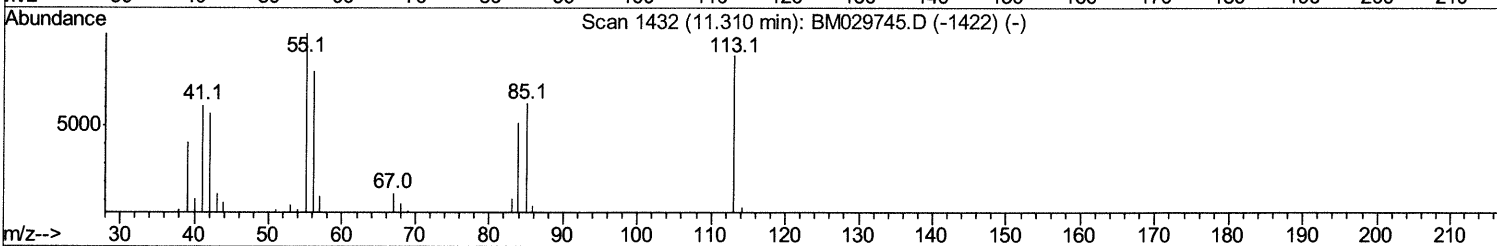
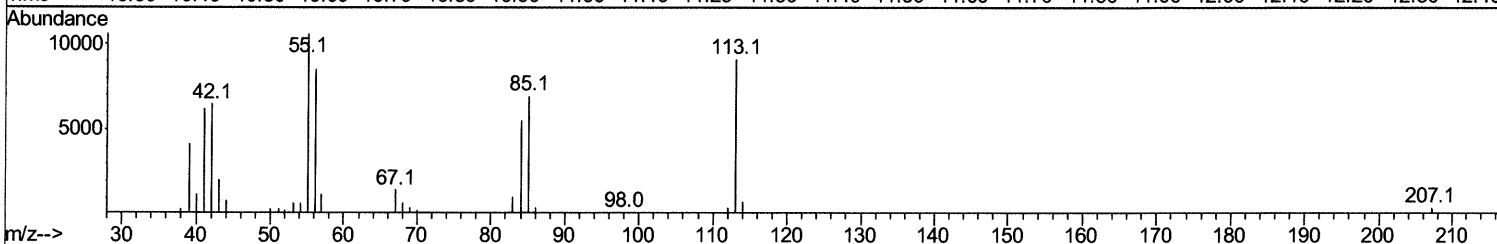
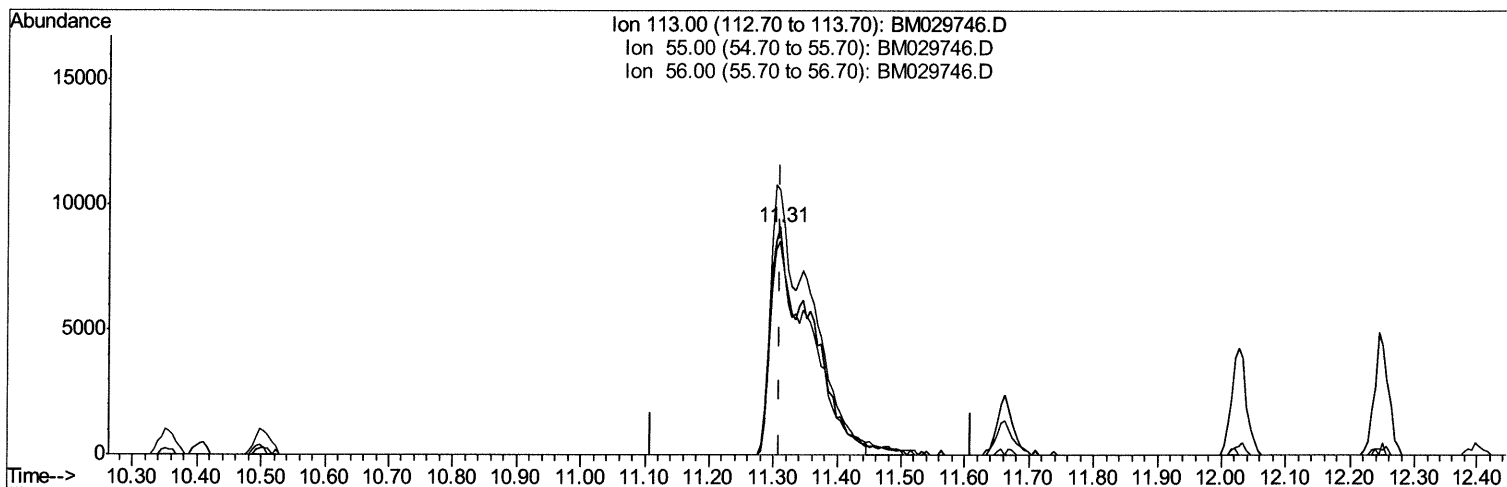
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
 Data File : BM029746.D  
 Acq On : 01 May 2021 14:00  
 Operator : CG/JU  
 Sample : SSTD04012  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD040012

Manual Integrations  
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mohammad  
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 Response via : Initial Calibration



TIC: BM029746.D

(34) Caprolactam

11.310min (-0.000) 48.77ng/ul m 05/04/21 JU

response 38192

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 113.50 | 116.03 |
| 56.00  | 90.40  | 93.56  |
| 0.00   | 0.00   | 0.00   |

## Quantitation Report (Qedit)

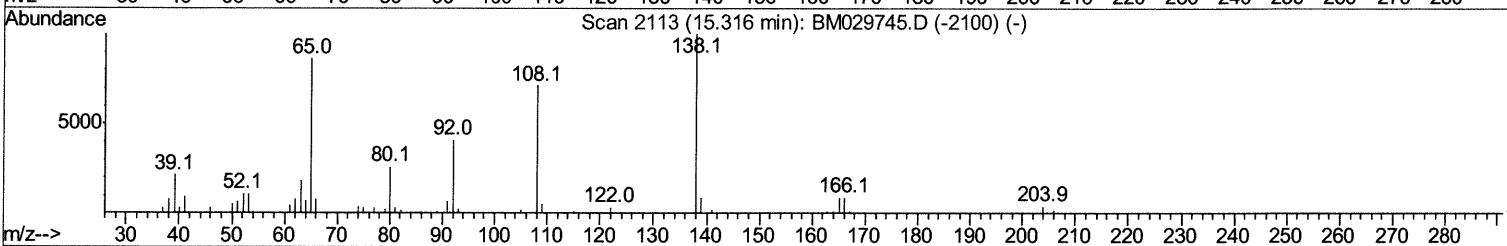
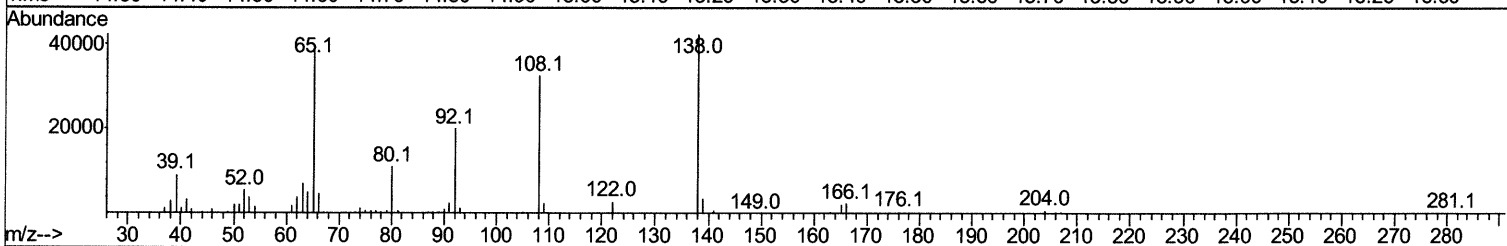
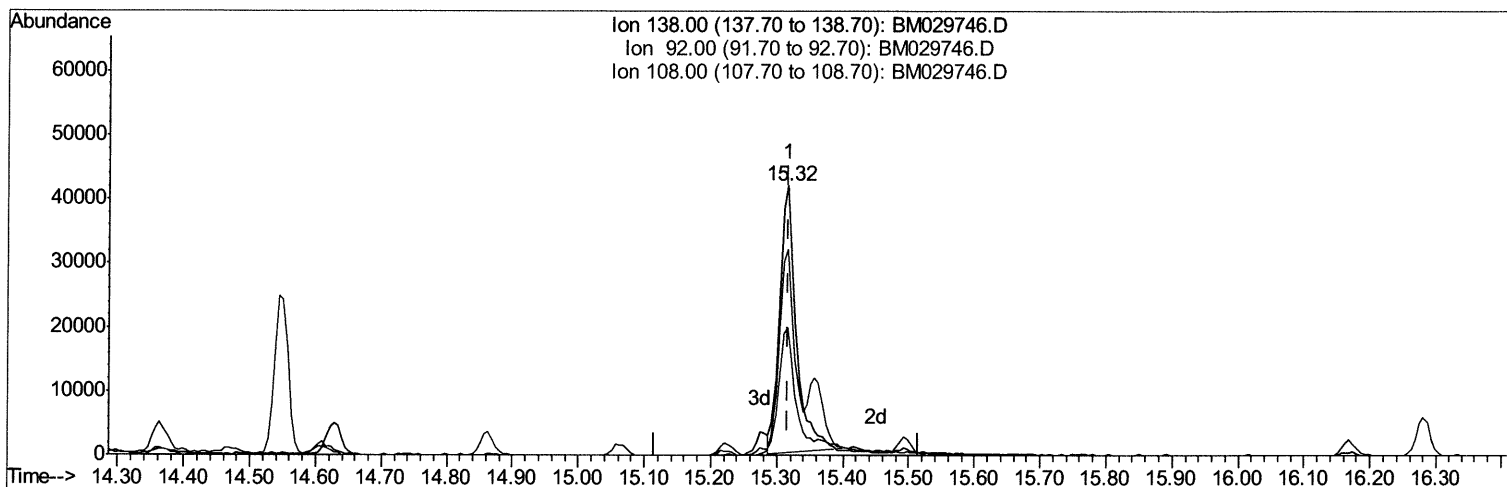
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
Data File : BM029746.D  
Acq On : 01 May 2021 14:00  
Operator : CG/JU  
Sample : SSTD04012  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD040012

Manual Integrations  
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QLast Update : Sat May 01 14:02:00 2021  
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TIC: BM029746.D

(63) 4-Nitroaniline

15.315min (-0.000) 38.25ng/ul

response 74395

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 42.50 | 47.42 |
| 108.00 | 71.00 | 76.55 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

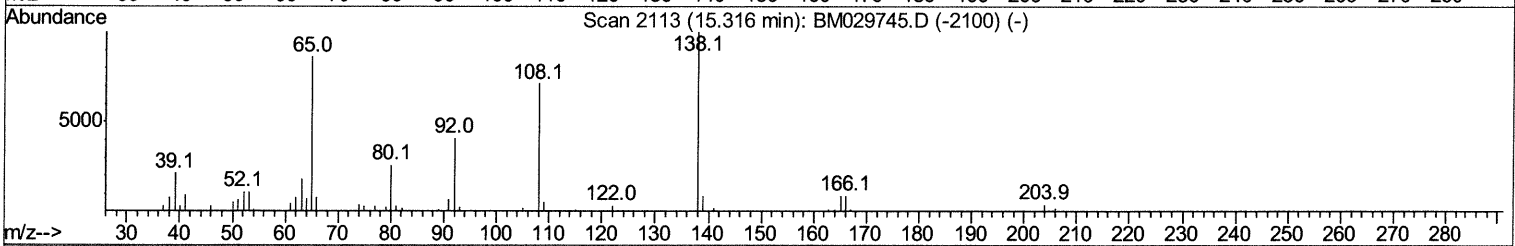
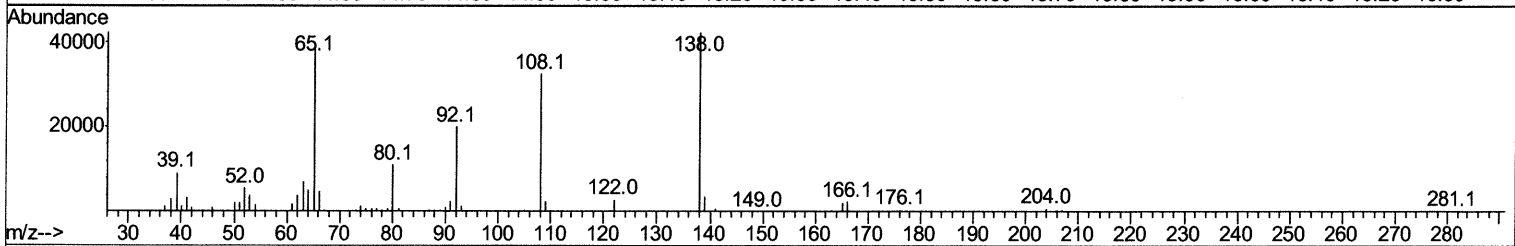
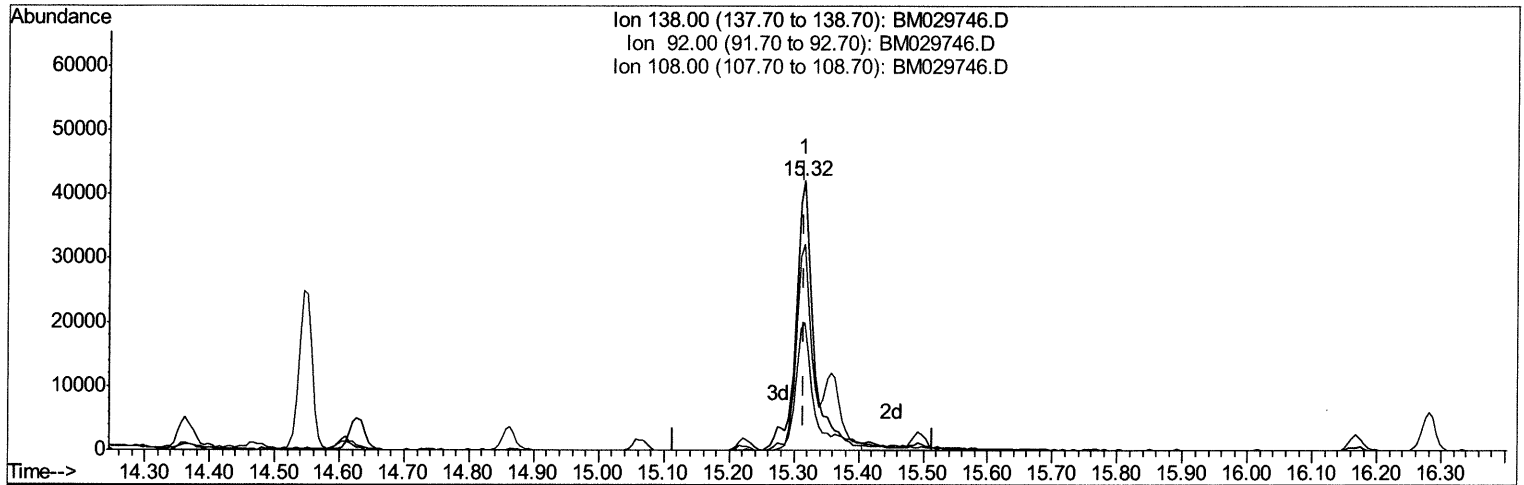
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
Data File : BM029746.D  
Acq On : 01 May 2021 14:00  
Operator : CG/JU  
Sample : SSTD04012  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD040012

Manual Integrations  
APPROVED

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Quant Time: May 01 14:33:49 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050121.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat May 01 14:02:00 2021  
Response via : Initial Calibration



TIC: BM029746.D

(63) 4-Nitroaniline

15.315min (-0.000) 43.11ng/ul m 05/04/21 JU

response 83848

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 138.00 | 100   | 100   |
| 92.00  | 42.50 | 47.42 |
| 108.00 | 71.00 | 76.55 |
| 0.00   | 0.00  | 0.00  |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
 Data File : BM029746.D  
 Acq On : 01 May 2021 14:00  
 Operator : CG/JU  
 Sample : SST040012  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST040012

Manual Integrations  
 APPROVED

mohammad  
 5/3/2021 12:24:21 PM

Quant Time: May 01 14:35:46 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050121.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 01 14:02:00 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.58  | 152  | 50409    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.36 | 136  | 191600   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.23 | 164  | 140296   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.97 | 188  | 284641   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.16 | 240  | 361415   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.32 | 264  | 380520   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.08  | 96  | 15472  | 12.06 | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.49  | 84  | 83988  | 25.06 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.76  | 99  | 130340 | 32.74 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.92  | 67  | 69929  | 30.39 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.12  | 132 | 123650 | 39.91 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.30  | 113 | 117815 | 38.13 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.73  | 128 | 61327  | 50.32 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.45  | 143 | 77158  | 66.41 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 9.99  | 165 | 158085 | 51.93 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.50 | 131 | 166464 | 38.65 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.64 | 166 | 460753 | 43.55 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 13.92 | 160 | 565949 | 45.24 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.46 | 143 | 59532  | 34.55 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.22 | 176 | 385987 | 42.39 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.36 | 200 | 90636  | 67.54 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.07 | 188 | 595499 | 43.81 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.37 | 212 | 732435 | 43.42 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.19 | 264 | 924296 | 45.63 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        |        | Ovalue      |
|--------------------------------|-------|-----|---------|--------|--------|-------------|
| 2) 1,4-Dioxane                 | 3.11  | 88  | 15277   | 11.866 | ng/uL  | 97          |
| 5) Pyridine                    | 3.50  | 79  | 90844   | 26.380 | ng/ul  | 96          |
| 6) Benzaldehyde                | 6.73  | 77  | 77567   | 36.308 | ng/ul  | 97          |
| 8) Phenol                      | 6.79  | 94  | 127842  | 30.555 | ng/ul  | 98          |
| 10) Bis(2-Chloroethyl)ether    | 7.02  | 93  | 101298  | 32.765 | ng/ul  | 98          |
| 12) 2-Chlorophenol             | 7.15  | 128 | 120880  | 36.567 | ng/ul  | 99          |
| 13) 2-Methylphenol             | 8.03  | 108 | 110166  | 35.696 | ng/ul  | 96          |
| 14) 2,2'-oxybis(1-Chloropropan | 8.12  | 45  | 107815  | 26.607 | ng/ul  | 98          |
| 16) Acetophenone               | 8.40  | 105 | 177198  | 35.088 | ng/ul  | 97          |
| 17) N-Nitroso-di-n-propylamine | 8.39  | 70  | 90928   | 37.853 | ng/ul  | 99          |
| 18) 4-Methylphenol             | 8.36  | 108 | 121097  | 36.953 | ng/ul  | 100         |
| 19) Hexachloroethane           | 8.64  | 117 | 56694   | 40.042 | ng/ul  | 95          |
| 22) Nitrobenzene               | 8.77  | 77  | 132910  | 37.407 | ng/ul  | 98          |
| 23) Isophorone                 | 9.30  | 82  | 251840  | 39.672 | ng/ul  | 99          |
| 25) 2-Nitrophenol              | 9.48  | 139 | 79688   | 59.535 | ng/ul  | 98          |
| 26) 2,4-Dimethylphenol         | 9.55  | 107 | 142843  | 39.269 | ng/ul  | 99          |
| 27) Bis(2-Chloroethoxy)methane | 9.79  | 93  | 147473  | 38.888 | ng/ul  | 96          |
| 29) 2,4-Dichlorophenol         | 10.02 | 162 | 146221  | 47.455 | ng/ul  | 94          |
| 30) Naphthalene                | 10.40 | 128 | 411435  | 39.638 | ng/ul  | 98          |
| 32) 4-Chloroaniline            | 10.53 | 127 | 164054  | 37.453 | ng/ul  | 96          |
| 33) Hexachlorobutadiene        | 10.69 | 225 | 119959  | 54.217 | ng/ul  | 97          |
| 34) Caprolactam                | 11.31 | 113 | 38192m> | 48.768 | ng/ul> | 95/04/21 JU |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\  
 Data File : BM029746.D  
 Acq On : 01 May 2021 14:00  
 Operator : CG/JU  
 Sample : SST040012  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST040012

Manual Integrations  
 APPROVED

mohammad  
 5/3/2021 12:24:21 PM

Quant Time: May 01 14:35:46 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM050121.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat May 01 14:02:00 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min)    |
|--------------------------------|-------|------|----------|--------|--------|-------------|
| 35) 4-Chloro-3-methylphenol    | 11.66 | 107  | 124626   | 41.636 | ng/ul  | 99          |
| 36) 2-Methylnaphthalene        | 12.03 | 142  | 322795   | 44.021 | ng/ul  | 98          |
| 37) 1-Methylnaphthalene        | 12.25 | 142  | 310966   | 42.952 | ng/ul  | 97          |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.40 | 216  | 222372   | 45.873 | ng/ul  | 99          |
| 40) Hexachlorocyclopentadiene  | 12.38 | 237  | 109648   | 34.434 | ng/ul  | 95          |
| 41) 2,4,6-Trichlorophenol      | 12.65 | 196  | 133971   | 50.261 | ng/ul  | 95          |
| 42) 2,4,5-Trichlorophenol      | 12.73 | 196  | 140182   | 49.386 | ng/ul  | 98          |
| 43) 1,1'-Biphenyl              | 13.06 | 154  | 438546   | 38.397 | ng/ul  | 99          |
| 44) 2-Chloronaphthalene        | 13.09 | 162  | 349170   | 38.725 | ng/ul  | 99          |
| 45) 2-Nitroaniline             | 13.31 | 65   | 74593    | 38.561 | ng/ul  | 96          |
| 47) Dimethylphthalate          | 13.69 | 163  | 443202   | 40.924 | ng/ul  | 100         |
| 48) 2,6-Dinitrotoluene         | 13.82 | 165  | 92731    | 55.018 | ng/ul  | 96          |
| 50) Acenaphthylene             | 13.94 | 152  | 529801   | 40.954 | ng/ul  | 99          |
| 51) 3-Nitroaniline             | 14.14 | 138  | 81766    | 43.868 | ng/ul  | 99          |
| 52) Acenaphthene               | 14.29 | 153  | 358586   | 37.328 | ng/ul  | 98          |
| 53) 2,4-Dinitrophenol          | 14.36 | 184  | 48021    | 53.095 | ng/ul  | 94          |
| 55) 4-Nitrophenol              | 14.47 | 109  | 44259    | 26.995 | ng/ul  | 87          |
| 56) Dibenzofuran               | 14.63 | 168  | 540028   | 40.452 | ng/ul  | 100         |
| 57) 2,4-Dinitrotoluene         | 14.61 | 165  | 133670   | 55.578 | ng/ul  | 99          |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.86 | 232  | 133518   | 58.700 | ng/ul  | 99          |
| 59) Diethylphthalate           | 15.06 | 149  | 429353   | 40.267 | ng/ul  | 99          |
| 61) Fluorene                   | 15.28 | 166  | 450096   | 40.296 | ng/ul  | 99          |
| 62) 4-Chlorophenyl-phenylether | 15.28 | 204  | 252938   | 45.115 | ng/ul  | 97          |
| 63) 4-Nitroaniline             | 15.32 | 138  | 83848m   | 43.105 | ng/ul  | 05/04/21 JU |
| 66) 4,6-Dinitro-2-methylphenol | 15.37 | 198  | 86856    | 61.535 | ng/ul# | 95          |
| 67) N-Nitrosodiphenylamine     | 15.49 | 169  | 390941   | 42.995 | ng/ul  | 99          |
| 68) 4-Bromophenyl-phenylether  | 16.17 | 248  | 156361   | 51.622 | ng/ul  | 97          |
| 69) Hexachlorobenzene          | 16.28 | 284  | 168052   | 49.805 | ng/ul  | 100         |
| 70) Atrazine                   | 16.45 | 200  | 149500   | 42.919 | ng/ul  | 99          |
| 71) Pentachlorophenol          | 16.63 | 266  | 88253    | 46.060 | ng/ul  | 98          |
| 72) Phenanthrene               | 17.02 | 178  | 667239   | 39.706 | ng/ul  | 100         |
| 74) Anthracene                 | 17.10 | 178  | 690480   | 39.987 | ng/ul  | 99          |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.02 | 216  | 231398   | 50.339 | ng/uL  | 99          |
| 76) Pentachlorobenzene         | 14.55 | 250  | 216096   | 48.152 | ng/uL  | 97          |
| 77) Carbazole                  | 17.38 | 167  | 590851   | 37.175 | ng/ul  | 99          |
| 78) Di-n-butylphthalate        | 17.95 | 149  | 680373   | 41.444 | ng/ul  | 99          |
| 80) Fluoranthene               | 19.03 | 202  | 872833   | 40.277 | ng/ul  | 99          |
| 82) Pyrene                     | 19.40 | 202  | 902266   | 39.720 | ng/ul  | 99          |
| 83) Butylbenzylphthalate       | 20.30 | 149  | 328293   | 41.128 | ng/ul  | 98          |
| 84) 3,3'-Dichlorobenzidine     | 21.08 | 252  | 390034   | 52.112 | ng/ul  | 100         |
| 85) Benzo(a)anthracene         | 21.14 | 228  | 978856   | 41.983 | ng/ul  | 100         |
| 86) Bis(2-ethylhexyl)phthalate | 21.08 | 149  | 552687   | 42.138 | ng/ul  | 99          |
| 87) Chrysene                   | 21.19 | 228  | 981033   | 43.036 | ng/ul  | 100         |
| 89) Di-n-octyl phthalate       | 21.94 | 149  | 975604   | 40.192 | ng/ul  | 100         |
| 90) Benzo(b)fluoranthene       | 22.68 | 252  | 1037299  | 42.180 | ng/ul  | 100         |
| 91) Benzo(k)fluoranthene       | 22.72 | 252  | 1061363  | 43.194 | ng/ul  | 99          |
| 93) Benzo(a)pyrene             | 23.23 | 252  | 932865   | 41.862 | ng/ul  | 99          |
| 94) Indeno(1,2,3-cd)pyrene     | 25.50 | 276  | 1187343  | 39.623 | ng/ul  | 100         |
| 95) Dibenzo(a,h)anthracene     | 25.51 | 278  | 998900   | 39.509 | ng/ul  | 99          |
| 96) Benzo(g,h,i)perylene       | 26.16 | 276  | 966953   | 39.343 | ng/ul  | 98          |

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Sample : SST04012  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION  
QLast Update : Sat May 01 14:02:00 2021  
Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

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