

(QT Reviewed)

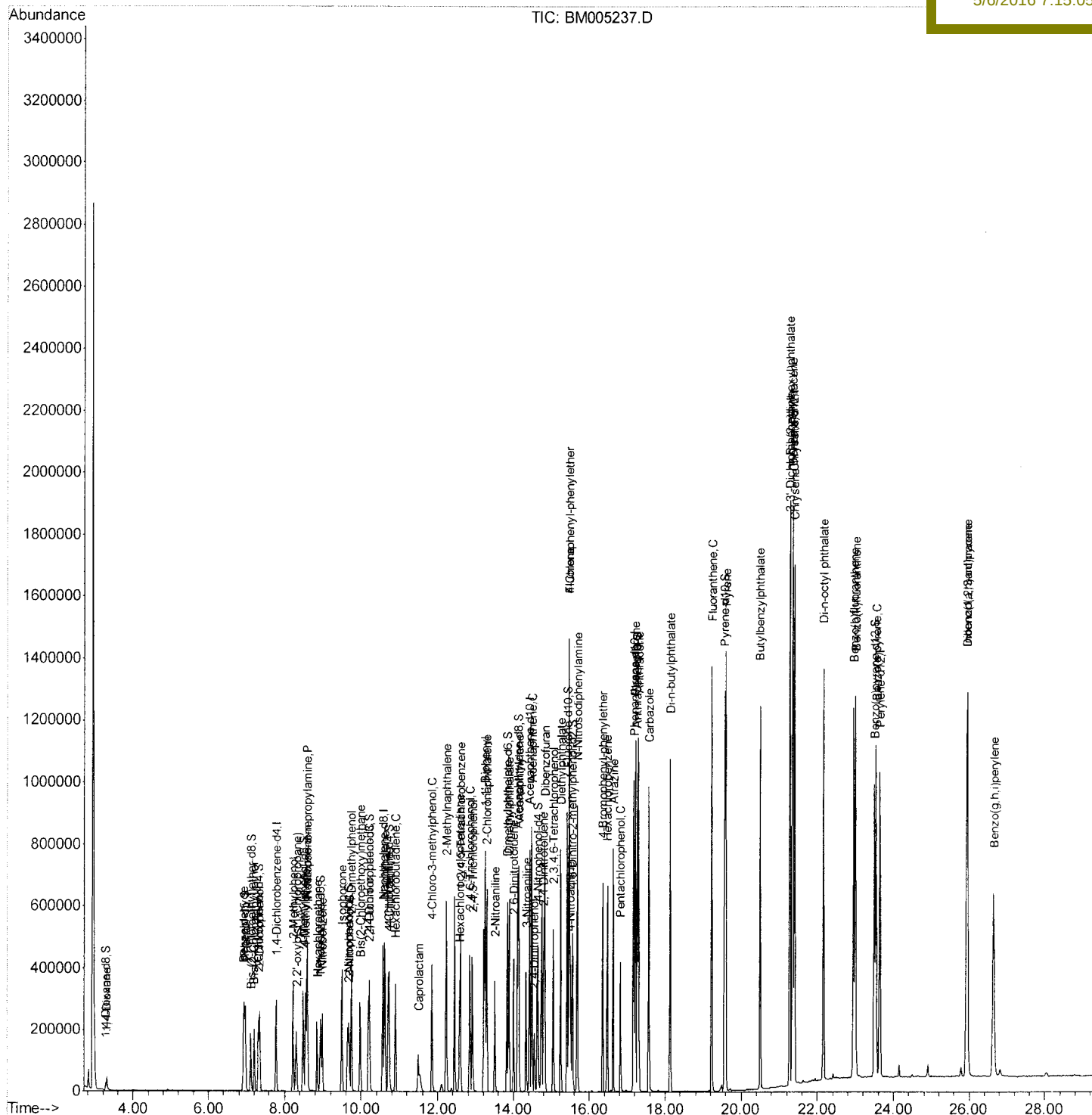
```
Data Path   : Z:\HPCHEM1\BNA_M\Data\BM050516\  
Data File  : BM005237.D  
Acq On     : 05 May 2016   16:30  
Operator   : UM/SJ  
Sample     : SSTDCCC020  
Misc       :  
ALS Vial   : 5      Sample Multiplier: 1
```

Quant Time: May 05 17:08:34 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu May 05 16:29:44 2016  
Response via : Initial Calibration

**Instrument :**  
BNA\_M  
**LabSampleId :**  
SSTD02046

## Manual Integrations APPROVED

sohil  
5/6/2016 7:15:05 PM



# Quantitation Report (Qedit)

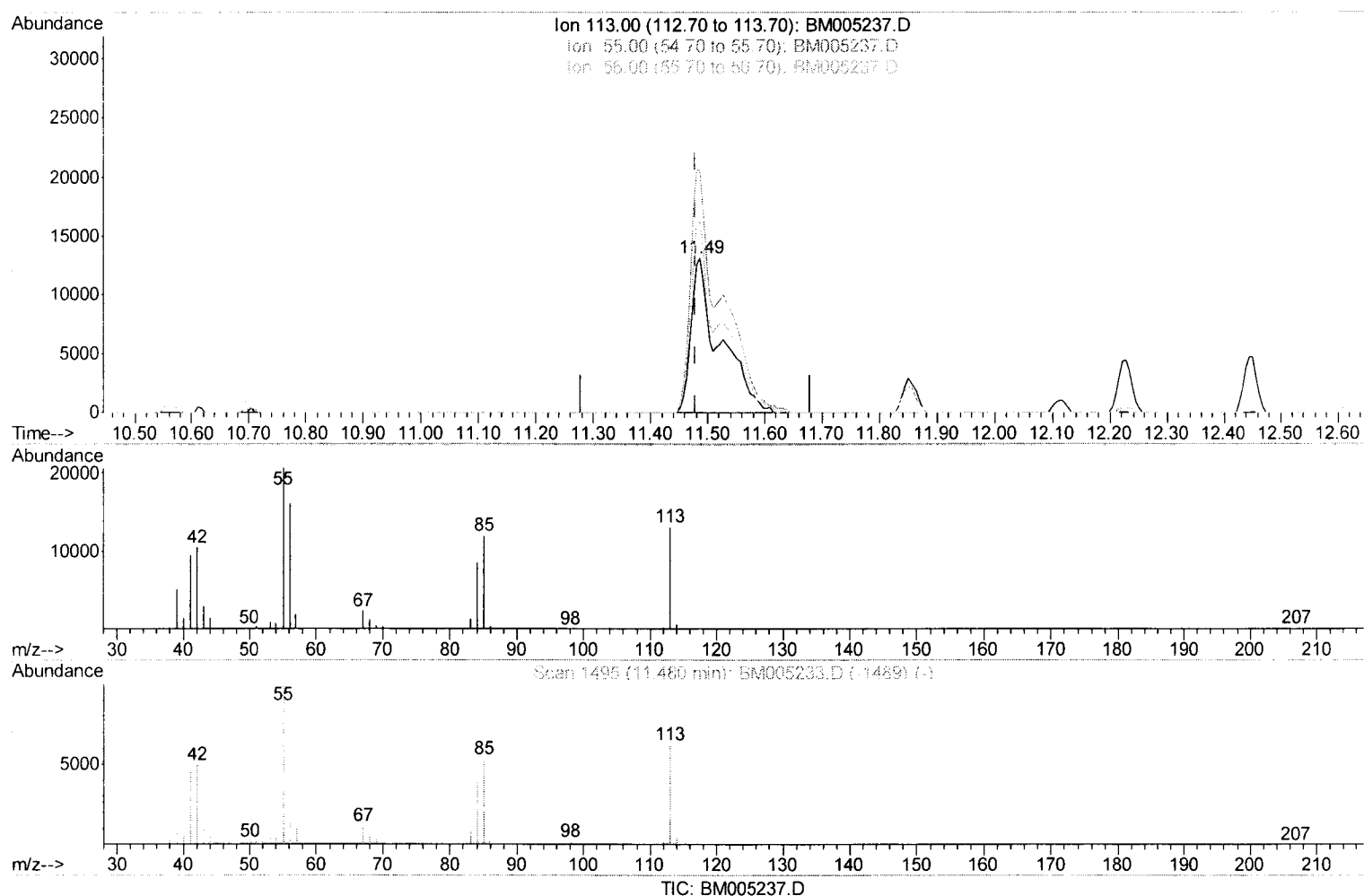
Data Path : Z:\HPCHEM1\BNA\_M\Data\BM050516\  
 Data File : BM005237.D  
 Acq On : 05 May 2016 16:30  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02046

Manual Integrations  
 APPROVED

sohil  
 5/6/2016 7:15:05 PM

Quant Time: May 05 17:04:09 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 05 16:29:44 2016  
 Response via : Initial Calibration



(32) Caprolactam

11.486min (+0.006) 20.39ng/ul m

response 46098

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 168.20 | 156.85 |
| 56.00  | 120.80 | 123.53 |
| 0.00   | 0.00   | 0.00   |

U.M  
 05/07/16

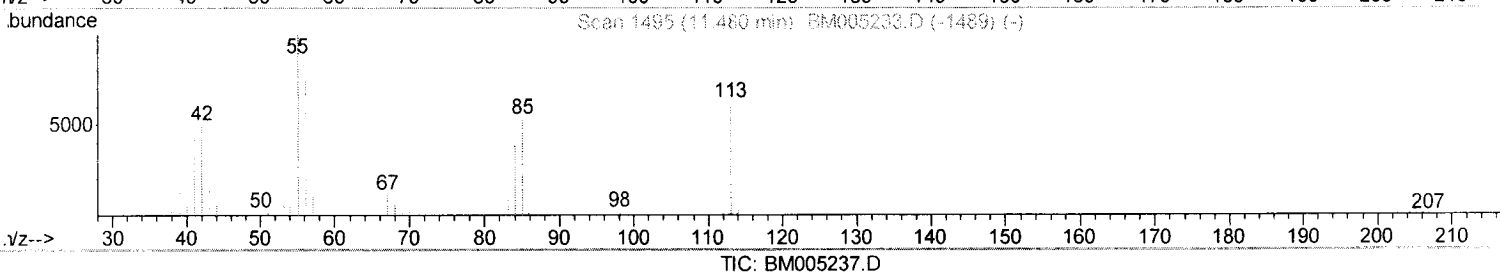
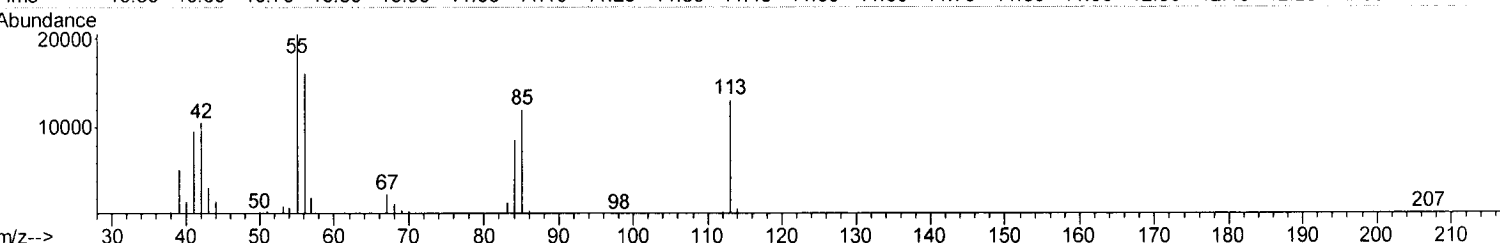
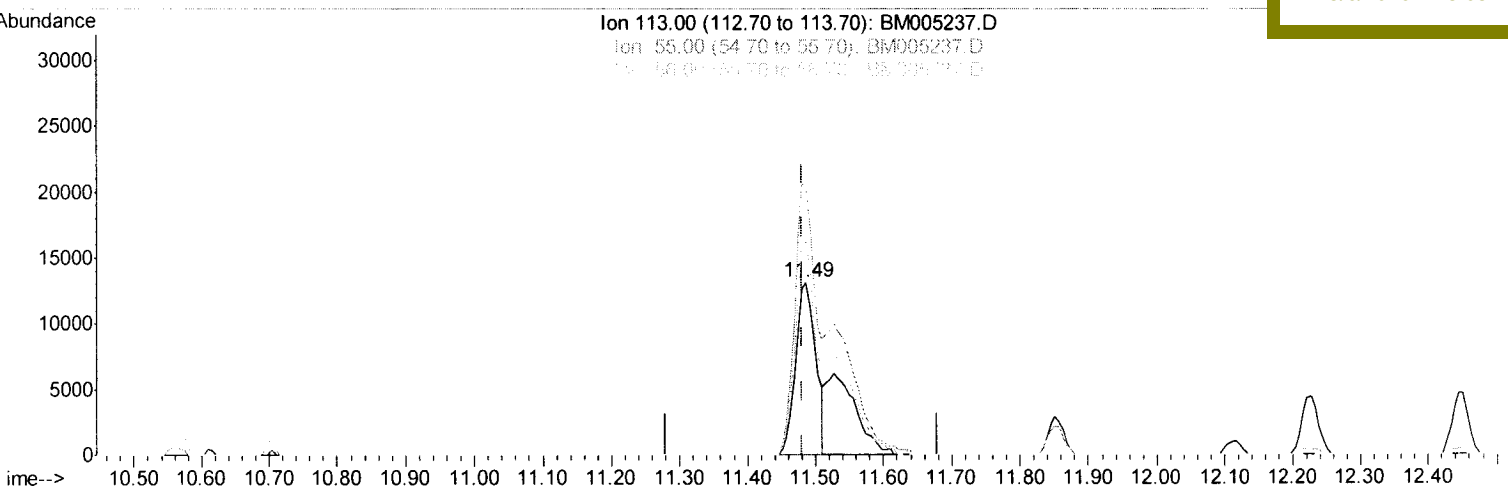
Data Path : Z:\HPCHEM1\BNA\_M\Data\BM050516\  
Data File : BM005237.D  
Acq On : 05 May 2016 16:30  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
LabSampleId :  
SSTD02046

Manual Integrations  
APPROVED

sohil  
5/6/2016 7:15:05 PM

Quant Time: May 05 17:04:09 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu May 05 16:29:44 2016  
Response via : Initial Calibration



(32) Caprolactam

11.486min (+0.006) 11.97ng/ul

response 27046

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 168.20 | 156.85 |
| 56.00  | 120.80 | 123.53 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM050516\  
 Data File : BM005237.D  
 Acq On : 05 May 2016 16:30  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02046

Quant Time: May 05 17:08:34 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 05 16:29:44 2016  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

sohil  
 5/6/2016 7:15:05 PM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.77  | 152  | 81056    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.56 | 136  | 394308   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.42 | 164  | 258750   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.17 | 188  | 629657   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.36 | 240  | 713059   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.64 | 264  | 715760   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.27  | 96  | 13343  | 7.74  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.95  | 99  | 146259 | 19.89 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.11  | 67  | 85030  | 20.27 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.31  | 132 | 111702 | 20.12 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.48  | 113 | 123114 | 20.26 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.93  | 128 | 58028  | 20.61 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.65  | 143 | 66852  | 20.97 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.19 | 165 | 121903 | 20.58 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.70 | 131 | 169926 | 23.83 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.83 | 166 | 424505 | 20.47 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.11 | 160 | 502910 | 20.67 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 77304  | 20.42 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.41 | 176 | 369219 | 20.62 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.54 | 200 | 75918  | 21.43 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.26 | 188 | 587737 | 21.12 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.56 | 212 | 674839 | 20.50 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.49 | 264 | 665634 | 21.01 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |             | Qvalue |
|--------------------------------|-------|-----|--------|-------------|--------|
| 2) 1,4-Dioxane                 | 3.30  | 88  | 24073  | 7.66 ng/uL# | 94     |
| 4) Benzaldehyde                | 6.92  | 77  | 92142  | 25.87 ng/ul | 99     |
| 6) Phenol                      | 6.97  | 94  | 152257 | 20.04 ng/ul | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.20  | 93  | 115485 | 20.19 ng/ul | 98     |
| 10) 2-Chlorophenol             | 7.34  | 128 | 114817 | 20.22 ng/ul | 98     |
| 11) 2-Methylphenol             | 8.22  | 108 | 118792 | 20.23 ng/ul | 100    |
| 12) 2,2'-oxybis(1-Chloropropan | 8.30  | 45  | 157125 | 20.25 ng/ul | 99     |
| 14) Acetophenone               | 8.59  | 105 | 191709 | 21.29 ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.58  | 70  | 99896  | 21.24 ng/ul | 98     |
| 16) 4-Methylphenol             | 8.55  | 108 | 133515 | 20.49 ng/ul | 97     |
| 17) Hexachloroethane           | 8.85  | 117 | 43197  | 20.21 ng/ul | 95     |
| 20) Nitrobenzene               | 8.97  | 77  | 144047 | 20.44 ng/ul | 97     |
| 21) Isophorone                 | 9.49  | 82  | 284668 | 20.80 ng/ul | 98     |
| 23) 2-Nitrophenol              | 9.68  | 139 | 72061  | 20.98 ng/ul | 98     |
| 24) 2,4-Dimethylphenol         | 9.74  | 107 | 150363 | 20.64 ng/ul | 99     |
| 25) Bis(2-Chloroethoxy)methane | 9.97  | 93  | 169708 | 19.91 ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 10.21 | 162 | 124938 | 20.59 ng/ul | 98     |
| 28) Naphthalene                | 10.61 | 128 | 405261 | 20.43 ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.73 | 127 | 173867 | 23.91 ng/ul | 97     |
| 31) Hexachlorobutadiene        | 10.89 | 225 | 75491  | 20.94 ng/ul | 98     |
| 32) Caprolactam                | 11.49 | 113 | 46098m | 20.39 ng/ul | 98     |
| 33) 4-Chloro-3-methylphenol    | 11.85 | 107 | 150776 | 20.73 ng/ul | 99     |
| 34) 2-Methylnaphthalene        | 12.23 | 142 | 307806 | 20.74 ng/ul | 97     |

UM  
 05/07/16

## Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM050516\  
 Data File : BM005237.D  
 Acq On : 05 May 2016 16:30  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02046

Manual Integrations  
 APPROVED

sohil  
 5/6/2016 7:15:05 PM

Quant Time: May 05 17:08:34 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM050516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 05 16:29:44 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev (Min) |
|--------------------------------|-------|------|----------|-------|--------|-----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.60 | 216  | 158108   | 20.09 | ng/ul  | 99        |
| 37) Hexachlorocyclopentadiene  | 12.57 | 237  | 78610    | 19.55 | ng/ul  | 94        |
| 38) 2,4,6-Trichlorophenol      | 12.85 | 196  | 106033   | 20.23 | ng/ul  | 94        |
| 39) 2,4,5-Trichlorophenol      | 12.92 | 196  | 118536   | 20.38 | ng/ul  | 99        |
| 40) 1,1'-Biphenyl              | 13.24 | 154  | 414175   | 20.21 | ng/ul  | 100       |
| 41) 2-Chloronaphthalene        | 13.29 | 162  | 315745   | 20.37 | ng/ul  | 99        |
| 42) 2-Nitroaniline             | 13.50 | 65   | 100222   | 21.00 | ng/ul  | 98        |
| 44) Dimethylphthalate          | 13.87 | 163  | 423266   | 20.42 | ng/ul  | 100       |
| 45) 2,6-Dinitrotoluene         | 14.00 | 165  | 87513    | 21.41 | ng/ul# | 90        |
| 47) Acenaphthylene             | 14.14 | 152  | 530012   | 20.59 | ng/ul  | 100       |
| 48) 3-Nitroaniline             | 14.33 | 138  | 96358    | 22.12 | ng/ul  | 97        |
| 49) Acenaphthene               | 14.48 | 153  | 348080   | 20.52 | ng/ul  | 99        |
| 50) 2,4-Dinitrophenol          | 14.54 | 184  | 42842    | 18.22 | ng/ul  | 97        |
| 52) 4-Nitrophenol              | 14.64 | 109  | 64387    | 20.46 | ng/ul  | 99        |
| 53) Dibenzofuran               | 14.82 | 168  | 510128   | 20.73 | ng/ul  | 99        |
| 54) 2,4-Dinitrotoluene         | 14.79 | 165  | 132726   | 21.78 | ng/ul  | 98        |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.04 | 232  | 104102   | 21.05 | ng/ul# | 96        |
| 56) Diethylphthalate           | 15.24 | 149  | 433809   | 20.72 | ng/ul  | 99        |
| 58) Fluorene                   | 15.47 | 166  | 405577   | 21.08 | ng/ul  | 99        |
| 59) 4-Chlorophenyl-phenylether | 15.46 | 204  | 201283   | 21.11 | ng/ul  | 99        |
| 60) 4-Nitroaniline             | 15.50 | 138  | 102213   | 22.14 | ng/ul  | 97        |
| 63) 4,6-Dinitro-2-methylphenol | 15.55 | 198  | 80016    | 21.50 | ng/ul# | 93        |
| 64) N-Nitrosodiphenylamine     | 15.67 | 169  | 364805   | 21.16 | ng/ul  | 99        |
| 65) 4-Bromophenyl-phenylether  | 16.36 | 248  | 128369   | 21.26 | ng/ul  | 96        |
| 66) Hexachlorobenzene          | 16.47 | 284  | 140994   | 20.82 | ng/ul  | 98        |
| 67) Atrazine                   | 16.63 | 200  | 139850   | 22.23 | ng/ul  | 98        |
| 68) Pentachlorophenol          | 16.82 | 266  | 76837    | 20.42 | ng/ul  | 97        |
| 69) Phenanthrene               | 17.21 | 178  | 693710   | 20.95 | ng/ul  | 99        |
| 71) Anthracene                 | 17.30 | 178  | 700885   | 21.23 | ng/ul  | 100       |
| 72) Carbazole                  | 17.57 | 167  | 645400   | 22.22 | ng/ul  | 99        |
| 73) Di-n-butylphthalate        | 18.13 | 149  | 756058   | 21.34 | ng/ul  | 100       |
| 74) Fluoranthene               | 19.23 | 202  | 828479   | 22.59 | ng/ul  | 99        |
| 77) Pyrene                     | 19.59 | 202  | 854796   | 20.67 | ng/ul  | 99        |
| 78) Butylbenzylphthalate       | 20.49 | 149  | 363925   | 21.19 | ng/ul  | 97        |
| 79) 3,3'-Dichlorobenzidine     | 21.28 | 252  | 290257   | 22.64 | ng/ul  | 99        |
| 80) Benzo(a)anthracene         | 21.34 | 228  | 839251   | 20.46 | ng/ul  | 99        |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 511729   | 21.45 | ng/ul# | 98        |
| 82) Chrysene                   | 21.40 | 228  | 805841   | 20.71 | ng/ul  | 99        |
| 84) Di-n-octyl phthalate       | 22.16 | 149  | 921734   | 22.05 | ng/ul  | 98        |
| 85) Benzo(b)fluoranthene       | 22.95 | 252  | 877953   | 20.98 | ng/ul  | 99        |
| 86) Benzo(k)fluoranthene       | 23.00 | 252  | 825569   | 20.96 | ng/ul  | 99        |
| 88) Benzo(a)pyrene             | 23.54 | 252  | 822555   | 20.79 | ng/ul  | 100       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.94 | 276  | 906227   | 20.99 | ng/ul  | 98        |
| 90) Dibenzo(a,h)anthracene     | 25.95 | 278  | 767717   | 21.25 | ng/ul  | 99        |
| 91) Benzo(g,h,i)perylene       | 26.64 | 276  | 761331   | 20.82 | ng/ul  | 99        |

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