

(QT Reviewed)

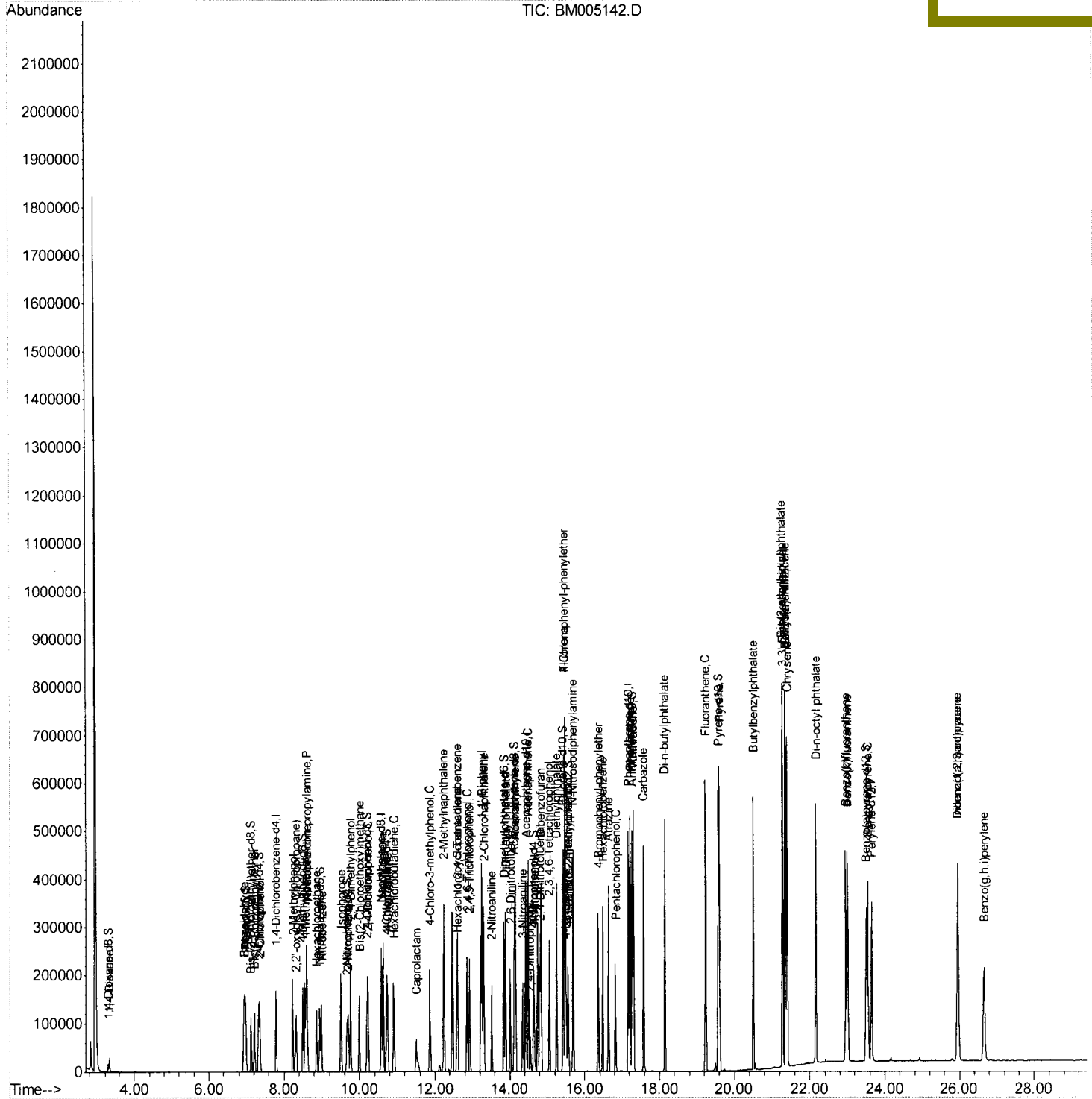
```
Data Path   : Z:\HPCHEM1\BNA_M\DATA\BM043016\  
Data File  : BM005142.D  
Acq On     : 29 Apr 2016   14:49  
Operator   : UM/SJ  
Sample     : SSTDCCC020  
Misc       :  
ALS Vial   : 2      Sample Multiplier: 1
```

Quant Time: Apr 30 03:04:52 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Apr 29 02:20:37 2016  
Response via : Initial Calibration

**Instrument :**  
BNA\_M  
**LabSampleId :**  
SSTD02054

## Manual Integrations

sohil  
4/30/2016 11:27:29 AM



# Quantitation Report (Qedit)

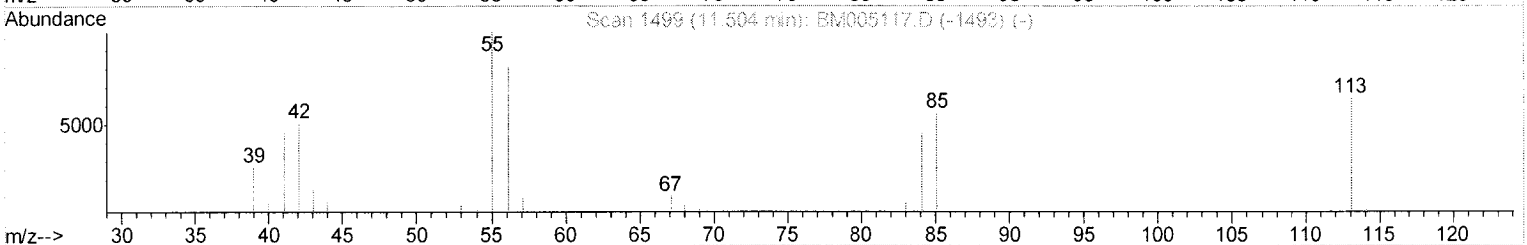
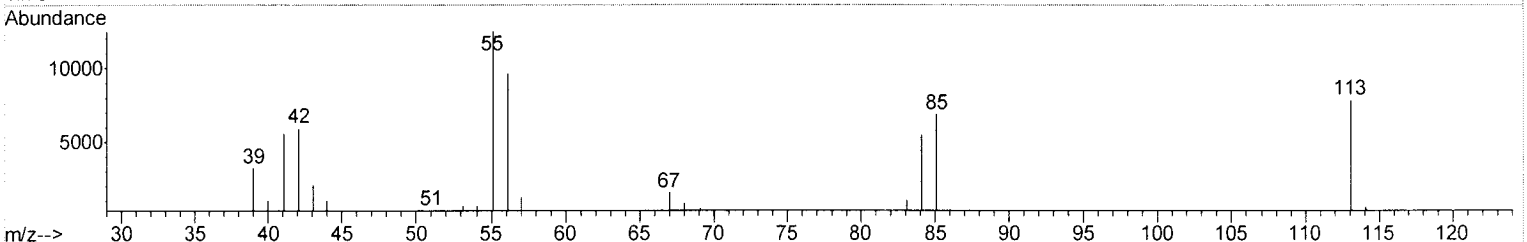
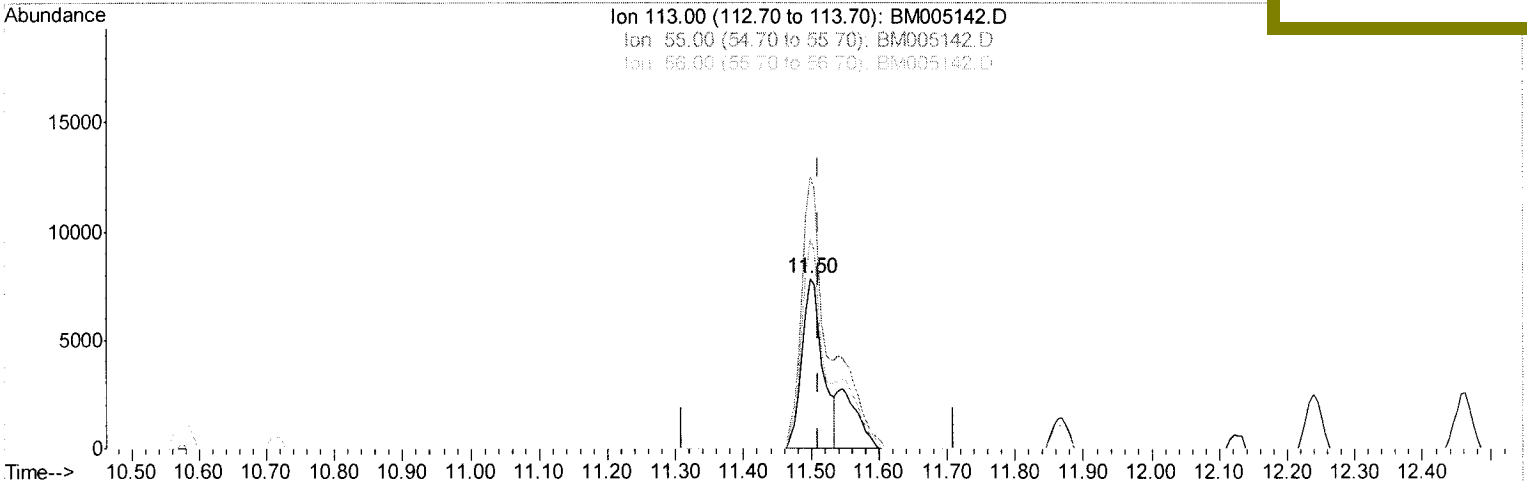
Data Path : Z:\HPCHEM1\BNA\_M\Data\BM043016\  
 Data File : BM005142.D  
 Acq On : 29 Apr 2016 14:49  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA M  
 LabSampleId :  
 SSTD02054

Quant Time: Apr 29 15:19:43 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

sohil  
 4/30/2016 11:27:29 AM



TIC: BM005142.D

(32) Caprolactam

11.498min (-0.012) 14.53ng/ul

response 16559

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 146.20 | 159.05 |
| 56.00  | 109.60 | 122.68 |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

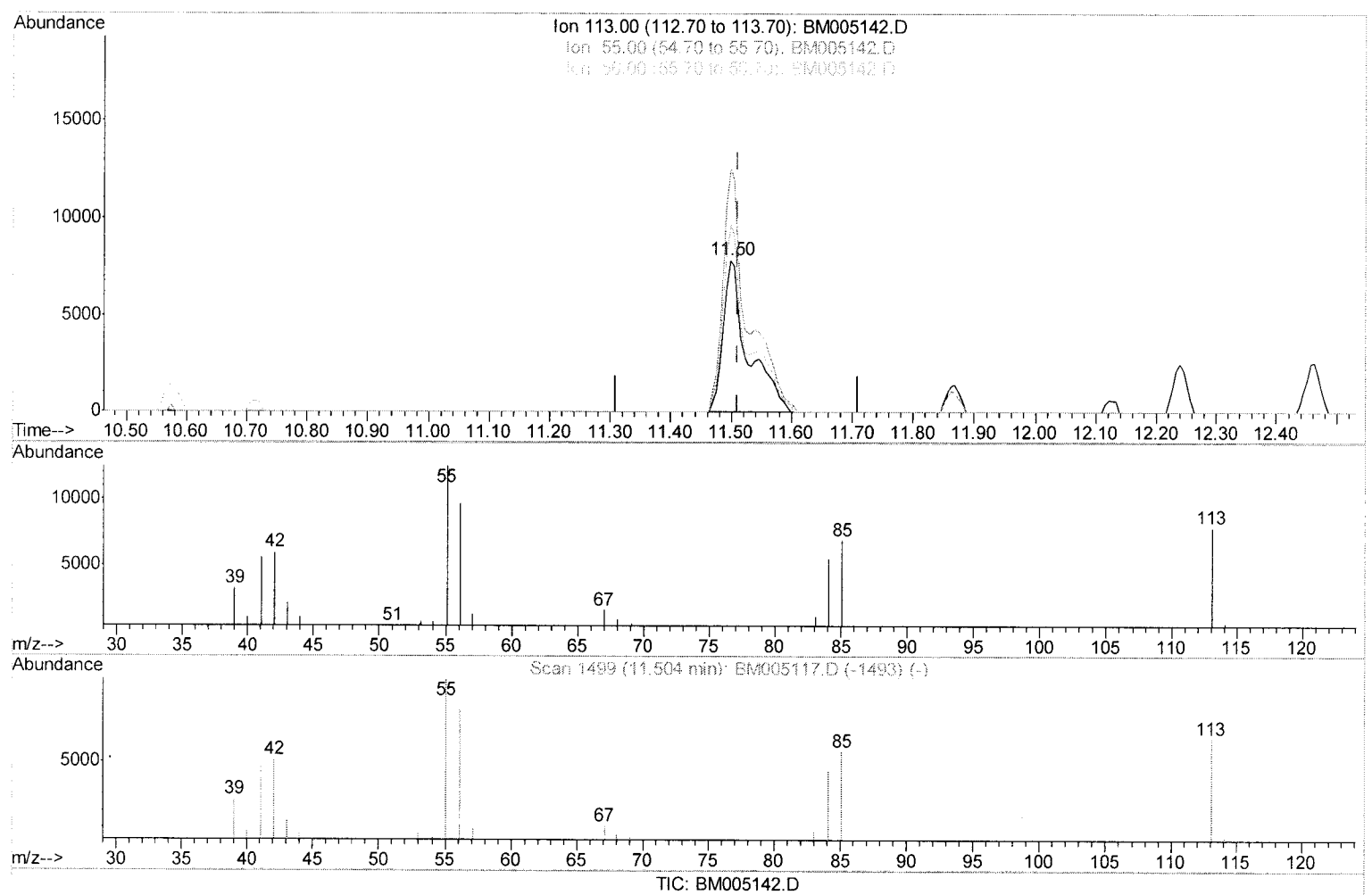
Data Path : Z:\HPCHEM1\BNA\_M\Data\BM043016\  
 Data File : BM005142.D  
 Acq On : 29 Apr 2016 14:49  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

sohil  
 4/30/2016 11:27:29 AM

Quant Time: Apr 29 15:19:43 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration



(32) Caprolactam

11.498min (-0.012) 19.57ng/ul m

response 22296

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 146.20 | 159.05 |
| 56.00  | 109.60 | 122.68 |
| 0.00   | 0.00   | 0.00   |

U.M  
 05/01/2016

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
 Data File : BM005142.D  
 Acq On : 29 Apr 2016 14:49  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

sohil  
 4/30/2016 11:27:29 AM

Quant Time: Apr 30 03:04:52 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev (Min) |
|---------------------------|-------|------|----------|-------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.79  | 152  | 47074    | 20.00 | ng/ul | 0.00      |
| 18) Naphthalene-d8        | 10.57 | 136  | 218361   | 20.00 | ng/ul | 0.00      |
| 35) Acenaphthene-d10      | 14.43 | 164  | 132689   | 20.00 | ng/ul | 0.00      |
| 61) Phenanthrene-d10      | 17.17 | 188  | 305284   | 20.00 | ng/ul | 0.00      |
| 75) Chrysene-d12          | 21.37 | 240  | 292386   | 20.00 | ng/ul | 0.00      |
| 83) Perylene-d12          | 23.64 | 264  | 243737   | 20.00 | ng/ul | -0.01     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96  | 7984   | 8.62  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 83881  | 21.54 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 49478  | 22.33 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.32  | 132 | 65826  | 21.37 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.49  | 113 | 69256  | 20.98 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.95  | 128 | 31882  | 21.08 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.66  | 143 | 35826  | 21.23 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.20 | 165 | 68587  | 21.40 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 90139  | 23.99 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.84 | 166 | 219861 | 20.98 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.12 | 160 | 259669 | 20.81 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 35806  | 19.55 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.42 | 176 | 186436 | 20.25 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.54 | 200 | 31593  | 18.19 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.27 | 188 | 284091 | 20.75 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.57 | 212 | 305270 | 22.93 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.50 | 264 | 229515 | 21.02 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |       | Qvalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.32  | 88  | 14646  | 8.52  | ng/uL | 95     |
| 4) Benzaldehyde                | 6.93  | 77  | 50097  | 25.91 | ng/ul | 98     |
| 6) Phenol                      | 6.99  | 94  | 88086  | 21.70 | ng/ul | 99     |
| 8) Bis(2-Chloroethyl)ether     | 7.22  | 93  | 68244  | 21.93 | ng/ul | 99     |
| 10) 2-Chlorophenol             | 7.35  | 128 | 67589  | 21.35 | ng/ul | 95     |
| 11) 2-Methylphenol             | 8.23  | 108 | 67388  | 21.27 | ng/ul | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.32  | 45  | 89706  | 22.29 | ng/ul | 99     |
| 14) Acetophenone               | 8.60  | 105 | 106423 | 21.42 | ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.59  | 70  | 54272  | 21.81 | ng/ul | 99     |
| 16) 4-Methylphenol             | 8.56  | 108 | 75391  | 21.38 | ng/ul | 99     |
| 17) Hexachloroethane           | 8.86  | 117 | 25859  | 20.51 | ng/ul | 89     |
| 20) Nitrobenzene               | 8.99  | 77  | 78910  | 21.20 | ng/ul | 96     |
| 21) Isophorone                 | 9.50  | 82  | 153102 | 21.66 | ng/ul | 97     |
| 23) 2-Nitrophenol              | 9.70  | 139 | 39450  | 21.51 | ng/ul | 97     |
| 24) 2,4-Dimethylphenol         | 9.76  | 107 | 81740  | 20.70 | ng/ul | 99     |
| 25) Bis(2-Chloroethoxy)methane | 9.99  | 93  | 95111  | 21.65 | ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 10.23 | 162 | 69027  | 20.95 | ng/ul | 99     |
| 28) Naphthalene                | 10.63 | 128 | 223601 | 20.30 | ng/ul | 100    |
| 30) 4-Chloroaniline            | 10.74 | 127 | 91222  | 23.85 | ng/ul | 99     |
| 31) Hexachlorobutadiene        | 10.91 | 225 | 41289  | 18.99 | ng/ul | 97     |
| 32) Caprolactam                | 11.50 | 113 | 22296m | 19.57 | ng/ul | 97     |
| 33) 4-Chloro-3-methylphenol    | 11.86 | 107 | 81261  | 21.57 | ng/ul | 97     |
| 34) 2-Methylnaphthalene        | 12.24 | 142 | 167376 | 20.53 | ng/ul | 99     |

U.M.  
 05/01/2016

## Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
 Data File : BM005142.D  
 Acq On : 29 Apr 2016 14:49  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA M  
 LabSampleId :  
 SSTD02054

Manual Integrations  
 APPROVED

sohil  
 4/30/2016 11:27:29 AM

Quant Time: Apr 30 03:04:52 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.61 | 216  | 86445    | 20.85 | ng/ul | 99       |
| 37) Hexachlorocyclopentadiene  | 12.59 | 237  | 44947    | 18.27 | ng/ul | 96       |
| 38) 2,4,6-Trichlorophenol      | 12.86 | 196  | 57166    | 21.68 | ng/ul | 98       |
| 39) 2,4,5-Trichlorophenol      | 12.93 | 196  | 63475    | 21.49 | ng/ul | 99       |
| 40) 1,1'-Biphenyl              | 13.26 | 154  | 224052   | 21.35 | ng/ul | 99       |
| 41) 2-Chloronaphthalene        | 13.30 | 162  | 167247   | 20.95 | ng/ul | 100      |
| 42) 2-Nitroaniline             | 13.51 | 65   | 48999    | 22.63 | ng/ul | 96       |
| 44) Dimethylphthalate          | 13.89 | 163  | 219284   | 20.86 | ng/ul | 100      |
| 45) 2,6-Dinitrotoluene         | 14.01 | 165  | 43270    | 21.32 | ng/ul | 94       |
| 47) Acenaphthylene             | 14.15 | 152  | 274969   | 20.82 | ng/ul | 99       |
| 48) 3-Nitroaniline             | 14.34 | 138  | 45801    | 22.18 | ng/ul | 100      |
| 49) Acenaphthene               | 14.49 | 153  | 178597   | 20.56 | ng/ul | 100      |
| 50) 2,4-Dinitrophenol          | 14.54 | 184  | 16973    | 14.33 | ng/ul | 98       |
| 52) 4-Nitrophenol              | 14.65 | 109  | 29837    | 18.22 | ng/ul | 97       |
| 53) Dibenzofuran               | 14.83 | 168  | 258703   | 20.19 | ng/ul | 99       |
| 54) 2,4-Dinitrotoluene         | 14.80 | 165  | 63568    | 20.61 | ng/ul | 95       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.06 | 232  | 52473    | 20.46 | ng/ul | 98       |
| 56) Diethylphthalate           | 15.25 | 149  | 219808   | 20.72 | ng/ul | 99       |
| 58) Fluorene                   | 15.48 | 166  | 203053   | 20.42 | ng/ul | 99       |
| 59) 4-Chlorophenyl-phenylether | 15.47 | 204  | 101488   | 20.22 | ng/ul | 98       |
| 60) 4-Nitroaniline             | 15.50 | 138  | 45395    | 20.38 | ng/ul | 98       |
| 63) 4,6-Dinitro-2-methylphenol | 15.56 | 198  | 33808    | 18.49 | ng/ul | 97       |
| 64) N-Nitrosodiphenylamine     | 15.69 | 169  | 180642   | 21.53 | ng/ul | 99       |
| 65) 4-Bromophenyl-phenylether  | 16.37 | 248  | 62989    | 21.14 | ng/ul | 100      |
| 66) Hexachlorobenzene          | 16.48 | 284  | 69799    | 20.56 | ng/ul | 99       |
| 67) Atrazine                   | 16.64 | 200  | 67018    | 21.67 | ng/ul | 98       |
| 68) Pentachlorophenol          | 16.83 | 266  | 37854    | 19.34 | ng/ul | 98       |
| 69) Phenanthrene               | 17.22 | 178  | 335838   | 20.67 | ng/ul | 99       |
| 71) Anthracene                 | 17.31 | 178  | 338766   | 20.68 | ng/ul | 99       |
| 72) Carbazole                  | 17.58 | 167  | 307736   | 21.91 | ng/ul | 99       |
| 73) Di-n-butylphthalate        | 18.14 | 149  | 367648   | 22.36 | ng/ul | 100      |
| 74) Fluoranthene               | 19.24 | 202  | 381998   | 21.36 | ng/ul | 99       |
| 77) Pyrene                     | 19.60 | 202  | 385542   | 22.83 | ng/ul | 98       |
| 78) Butylbenzylphthalate       | 20.50 | 149  | 156452   | 24.08 | ng/ul | 98       |
| 79) 3,3'-Dichlorobenzidine     | 21.29 | 252  | 111995   | 21.42 | ng/ul | 99       |
| 80) Benzo(a)anthracene         | 21.35 | 228  | 352585   | 21.07 | ng/ul | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 215842   | 23.66 | ng/ul | 100      |
| 82) Chrysene                   | 21.40 | 228  | 329964   | 20.60 | ng/ul | 100      |
| 84) Di-n-octyl phthalate       | 22.16 | 149  | 347744   | 22.95 | ng/ul | 99       |
| 85) Benzo(b)fluoranthene       | 22.96 | 252  | 306306   | 20.86 | ng/ul | 99       |
| 86) Benzo(k)fluoranthene       | 23.00 | 252  | 300350   | 21.04 | ng/ul | 100      |
| 88) Benzo(a)pyrene             | 23.54 | 252  | 285887   | 20.85 | ng/ul | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.95 | 276  | 294263   | 22.50 | ng/ul | 98       |
| 90) Dibenzo(a,h)anthracene     | 25.96 | 278  | 245981   | 22.51 | ng/ul | 99       |
| 91) Benzo(g,h,i)perylene       | 26.66 | 276  | 245558   | 22.95 | ng/ul | 98       |

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