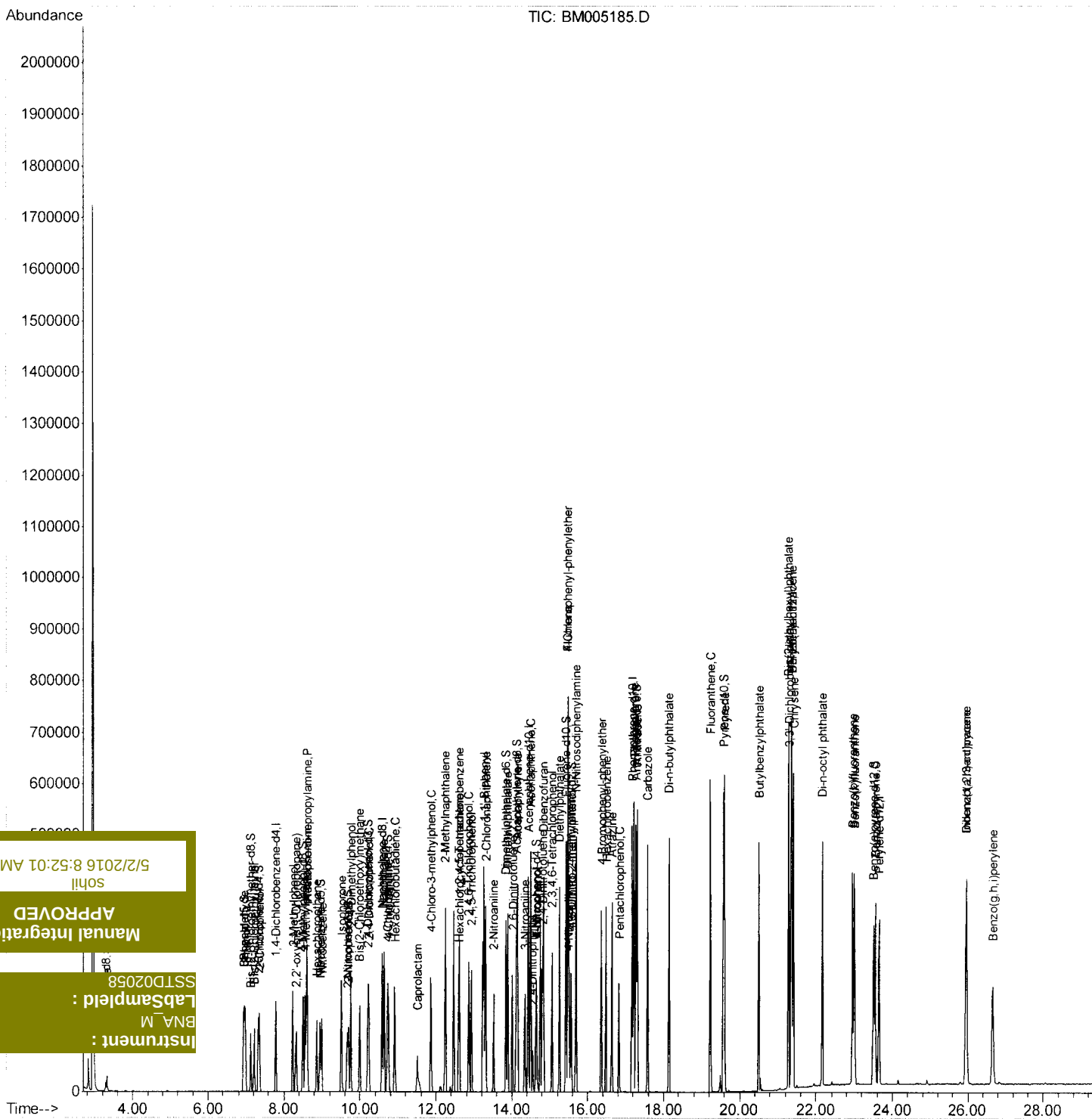


(QT Reviewed)

```
Data Path   : Z:\HPCHEM1\BNA_M\DATA\BM043016\  
Data File  : BM005185.D  
Acq On     : 30 Apr 2016   22:01  
Operator   : UM/SJ  
Sample     : SSTDCCC020EC  
Misc       :  
ALS Vial   : 2      Sample Multiplier: 1
```

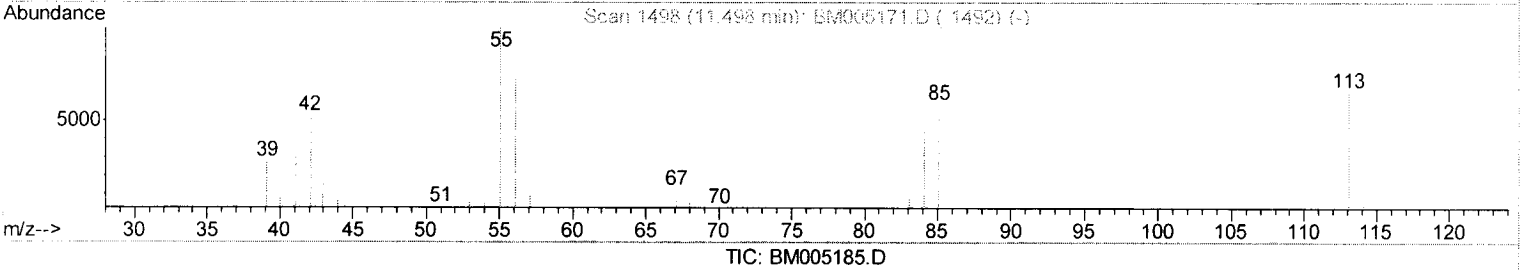
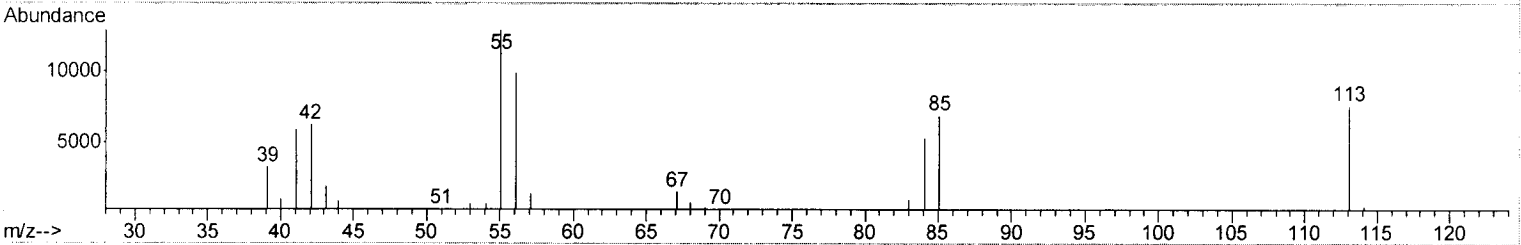
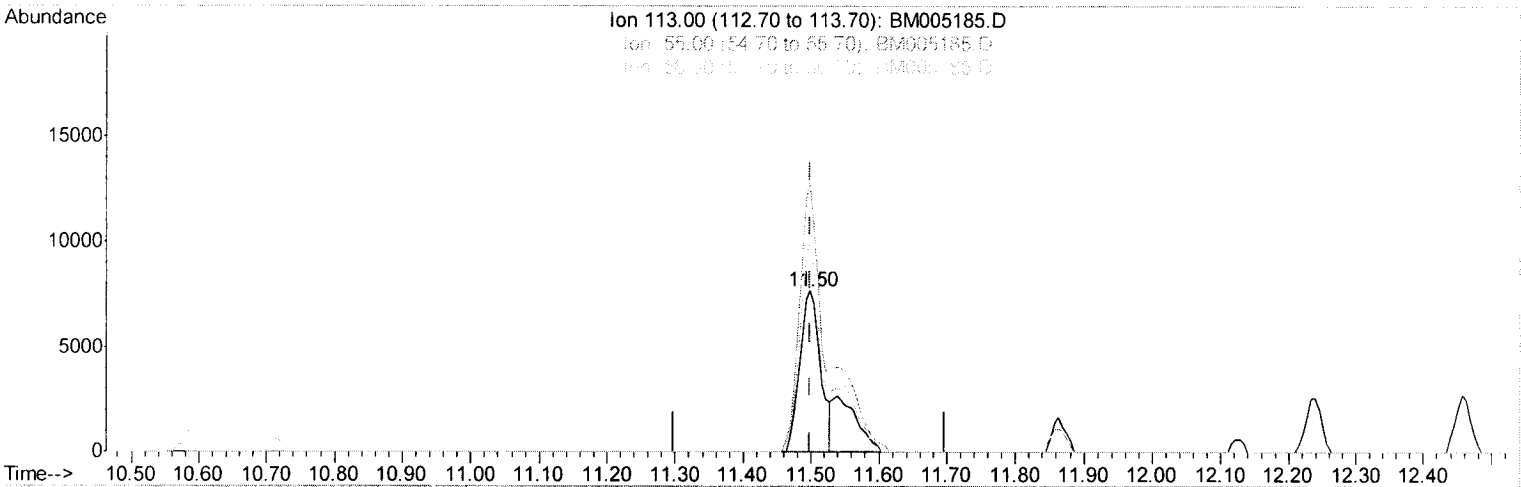
Quant Time: May 01 04:53:27 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun May 01 04:14:40 2016  
Response via : Initial Calibration



# Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
 Data File : BM005185.D  
 Acq On : 30 Apr 2016 22:01  
 Operator : UM/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 01 04:16:59 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun May 01 04:14:40 2016  
 Response via : Initial Calibration



(32) Caprolactam

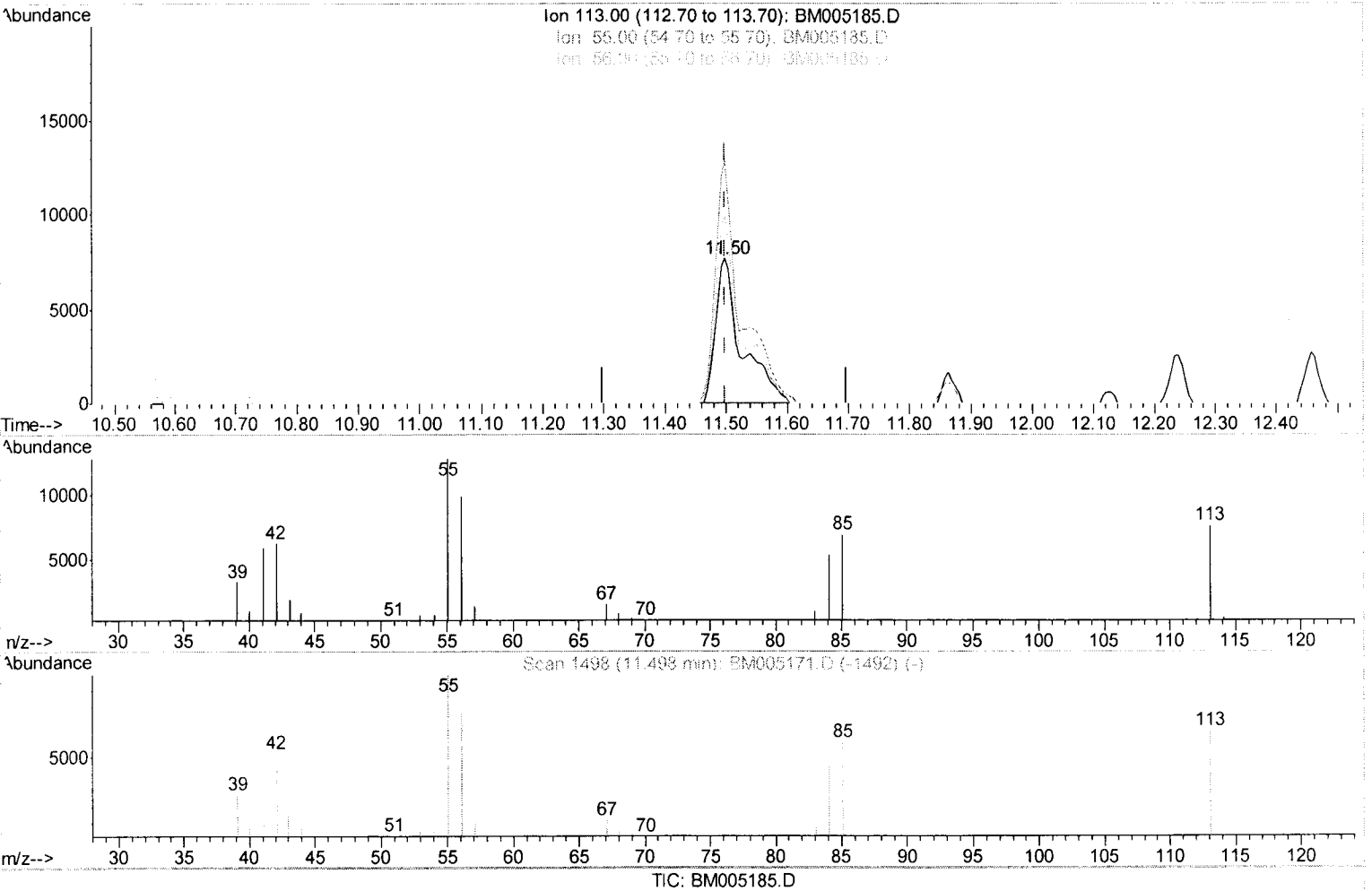
11.498min (+0.000) 13.72ng/ul

|                     |        |
|---------------------|--------|
| 5/2/2016 8:52:01 AM |        |
| Sohil               | Act%   |
| APPROVED            | 100    |
| Manual Integrations | 168.13 |
| 55.00 146.20        |        |
| SSTD02058           | 129.89 |
| LabSampled :        |        |
| BNA_M               | 0.00   |
| Instrument :        |        |

# Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
 Data File : BM005185.D  
 Acq On : 30 Apr 2016 22:01  
 Operator : UM/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 01 04:16:59 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun May 01 04:14:40 2016  
 Response via : Initial Calibration



(32) Caprolactam

11.498min (+0.000) 19.24ng/ul m

response 22808

Ion Exp% Act%

*U.A*  
*05/04/16*

00  
 68.13  
 29.89  
 00

Manual Integrations  
 APPROVED  
 5/2/2016 8:52:01 AM  
 Sohil

SSTD02058  
 LabSampled :  
 BNA\_M  
 Instrument :

SOM02.2-EPA-BM042516.M Sun May 01 04:53:25 2016

## Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
Data File : BM005185.D  
Acq On : 30 Apr 2016 22:01  
Operator : UM/SJ  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 01 04:53:27 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun May 01 04:14:40 2016  
Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.79  | 152  | 48270    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.57 | 136  | 227164   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.43 | 164  | 140520   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.17 | 188  | 311934   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.36 | 240  | 268708   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.64 | 264  | 239257   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96  | 7916   | 8.33  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 85416  | 21.39 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 50011  | 22.01 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.32  | 132 | 67785  | 21.46 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.49  | 113 | 70879  | 20.94 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.95  | 128 | 33692  | 21.41 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.66  | 143 | 37789  | 21.53 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.20 | 165 | 71276  | 21.37 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 94718  | 24.23 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.83 | 166 | 228715 | 20.61 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.12 | 160 | 276058 | 20.89 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 36463  | 18.80 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.42 | 176 | 195671 | 20.06 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.54 | 200 | 33809  | 19.06 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.27 | 188 | 291000 | 20.80 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.57 | 212 | 292211 | 23.89 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.50 | 264 | 218440 | 20.38 | ng/ul | 0.00 |

## Target Compounds

| Target Compounds               | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|--------------------------------|-------|------|----------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.32  | 88   | 14428    | 8.19  | ng/uL | 95     |
| 4) Benzaldehyde                | 6.93  | 77   | 51828    | 26.14 | ng/ul | 97     |
| 6) Phenol                      | 6.98  | 94   | 89238    | 21.44 | ng/ul | 98     |
| 8) Bis-(2-Chloroethyl)ether    | 7.21  | 93   | 69318    | 21.72 | ng/ul | 98     |
| 10) 2-Chlorophenol             | 7.35  | 128  | 68834    | 21.20 | ng/ul | 97     |
| 11) 2-Methylphenol             | 8.23  | 108  | 69190    | 21.30 | ng/ul | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.32  | 45   | 92716    | 22.46 | ng/ul | 97     |
| 14) Acetophenone               | 8.60  | 105  | 111517   | 21.89 | ng/ul | 98     |
| 15) N-Nitroso-di-n-propylamine | 8.59  | 70   | 56424    | 22.11 | ng/ul | 98     |
| 16) 4-Methylphenol             | 8.56  | 108  | 76803    | 21.24 | ng/ul | 95     |
| 17) 2,4-Dichlorophenol         | 8.86  | 117  | 26850    | 20.77 | ng/ul | 93     |
| 18) 2,6-Dichlorophenol         | 8.99  | 77   | 82559    | 21.33 | ng/ul | 99     |
| 19) 2-Nitrophenol              | 9.50  | 82   | 161422   | 21.95 | ng/ul | 98     |
| 20) 4-Nitrophenol              | 9.69  | 139  | 40632    | 21.30 | ng/ul | 91     |
| 21) 2-Methylphenol             | 9.75  | 107  | 85215    | 20.74 | ng/ul | 99     |
| 22) 2,4-Dichlorophenol         | 9.99  | 93   | 98537    | 21.56 | ng/ul | 99     |
| 23) 2,6-Dichlorophenol         | 10.23 | 162  | 72041    | 21.02 | ng/ul | 99     |
| 24) 2,4-Dichlorophenol         | 10.63 | 128  | 234618   | 20.48 | ng/ul | 100    |
| 25) 2,6-Dichlorophenol         | 10.74 | 127  | 96405    | 24.23 | ng/ul | 100    |
| 26) 2,4-Dichlorophenol         | 10.90 | 225  | 44291    | 19.58 | ng/ul | 98     |
| 27) 2,4-Dichlorophenol         | 11.50 | 113  | 22808m   | 19.24 | ng/ul | 98     |
| 28) 2,4-Dichlorophenol         | 11.86 | 107  | 83028    | 21.18 | ng/ul | 98     |
| 29) 2,4-Dichlorophenol         | 12.24 | 142  | 173925   | 20.51 | ng/ul | 100    |

5/2/2016 8:52:01 AM  
Sohil  
APPROVED  
Manual Integrations  
SSTDD02058  
Lab Sampled :  
BNA\_M  
Instrument :

UM  
05/04/16

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM043016\  
 Data File : BM005185.D  
 Acq On : 30 Apr 2016 22:01  
 Operator : UM/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02058

Manual Integrations  
 APPROVED

sohil  
 5/2/2016 8:52:01 AM

Quant Time: May 01 04:53:27 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun May 01 04:14:40 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.61 | 216  | 90111    | 20.53 | ng/ul | 99       |
| 37) Hexachlorocyclopentadiene  | 12.59 | 237  | 46141    | 17.71 | ng/ul | 95       |
| 38) 2,4,6-Trichlorophenol      | 12.85 | 196  | 59599    | 21.34 | ng/ul | 100      |
| 39) 2,4,5-Trichlorophenol      | 12.93 | 196  | 66301    | 21.20 | ng/ul | 98       |
| 40) 1,1'-Biphenyl              | 13.26 | 154  | 233276   | 20.99 | ng/ul | 99       |
| 41) 2-Chloronaphthalene        | 13.30 | 162  | 177390   | 20.98 | ng/ul | 100      |
| 42) 2-Nitroaniline             | 13.51 | 65   | 52394    | 22.85 | ng/ul | 98       |
| 44) Dimethylphthalate          | 13.88 | 163  | 226885   | 20.38 | ng/ul | 100      |
| 45) 2,6-Dinitrotoluene         | 14.01 | 165  | 46994    | 21.86 | ng/ul | 97       |
| 47) Acenaphthylene             | 14.15 | 152  | 291636   | 20.86 | ng/ul | 99       |
| 48) 3-Nitroaniline             | 14.34 | 138  | 49258    | 22.53 | ng/ul | 99       |
| 49) Acenaphthene               | 14.49 | 153  | 189469   | 20.59 | ng/ul | 99       |
| 50) 2,4-Dinitrophenol          | 14.55 | 184  | 18024    | 14.37 | ng/ul | 99       |
| 52) 4-Nitrophenol              | 14.65 | 109  | 30822    | 17.77 | ng/ul | 96       |
| 53) Dibenzofuran               | 14.83 | 168  | 275117   | 20.28 | ng/ul | 100      |
| 54) 2,4-Dinitrotoluene         | 14.80 | 165  | 67819    | 20.77 | ng/ul | 97       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.06 | 232  | 54147    | 19.94 | ng/ul | 99       |
| 56) Diethylphthalate           | 15.25 | 149  | 230298   | 20.50 | ng/ul | 100      |
| 58) Fluorene                   | 15.48 | 166  | 213927   | 20.31 | ng/ul | 99       |
| 59) 4-Chlorophenyl-phenylether | 15.47 | 204  | 106879   | 20.11 | ng/ul | 98       |
| 60) 4-Nitroaniline             | 15.50 | 138  | 48016    | 20.36 | ng/ul | 98       |
| 63) 4,6-Dinitro-2-methylphenol | 15.56 | 198  | 35857    | 19.19 | ng/ul | 95       |
| 64) N-Nitrosodiphenylamine     | 15.69 | 169  | 189767   | 22.13 | ng/ul | 99       |
| 65) 4-Bromophenyl-phenylether  | 16.36 | 248  | 66000    | 21.68 | ng/ul | 98       |
| 66) Hexachlorobenzene          | 16.48 | 284  | 71757    | 20.69 | ng/ul | 97       |
| 67) Atrazine                   | 16.64 | 200  | 68294    | 21.62 | ng/ul | 98       |
| 68) Pentachlorophenol          | 16.83 | 266  | 37026    | 18.51 | ng/ul | 99       |
| 69) Phenanthrene               | 17.22 | 178  | 343713   | 20.70 | ng/ul | 100      |
| 71) Anthracene                 | 17.31 | 178  | 346559   | 20.71 | ng/ul | 100      |
| 72) Carbazole                  | 17.58 | 167  | 306717   | 21.37 | ng/ul | 100      |
| 73) Di-n-butylphthalate        | 18.14 | 149  | 364634   | 21.70 | ng/ul | 100      |
| 74) Fluoranthene               | 19.23 | 202  | 365119   | 19.98 | ng/ul | 100      |
| 77) Pyrene                     | 19.60 | 202  | 368160   | 23.72 | ng/ul | 99       |
| 78) Butylbenzylphthalate       | 20.50 | 149  | 142499   | 23.87 | ng/ul | 97       |
| 79) 3,3'-Dichlorobenzidine     | 21.29 | 252  | 99632    | 20.73 | ng/ul | 98       |
| 80) Benzo(a)anthracene         | 21.35 | 228  | 317949   | 20.67 | ng/ul | 100      |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 192184   | 22.92 | ng/ul | 98       |
| 82) Chrysene                   | 21.40 | 228  | 302773   | 20.57 | ng/ul | 100      |
| 84) Di-n-octyl phthalate       | 22.16 | 149  | 308962   | 20.77 | ng/ul | 98       |
| 85) Benzo(b)fluoranthene       | 22.96 | 252  | 293807   | 20.39 | ng/ul | 99       |
| 86) Benzo(k)fluoranthene       | 23.00 | 252  | 273997   | 19.55 | ng/ul | 100      |
| 88) Benzo(a)pyrene             | 23.54 | 252  | 272415   | 20.24 | ng/ul | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.95 | 276  | 290368   | 22.62 | ng/ul | 99       |
| 90) Dibenzo(a,h)anthracene     | 25.96 | 278  | 244988   | 22.84 | ng/ul | 98       |
| 91) Benzo(g,h,i)perylene       | 26.66 | 276  | 241460   | 22.99 | ng/ul | 99       |

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