



# Quantitation Report (Qedit)

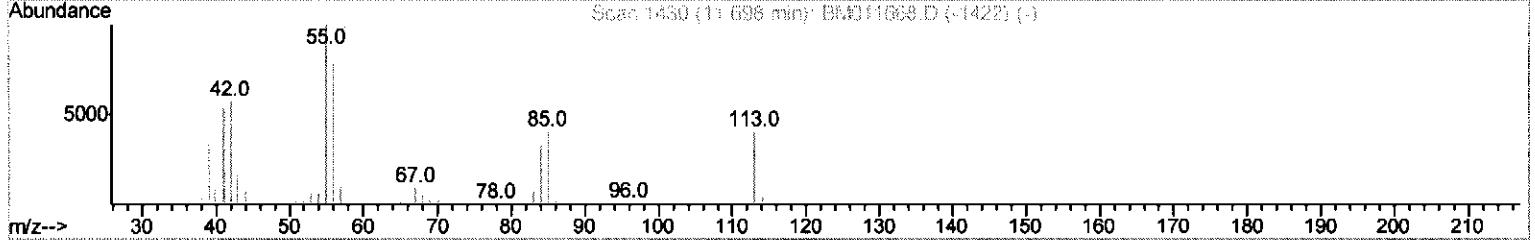
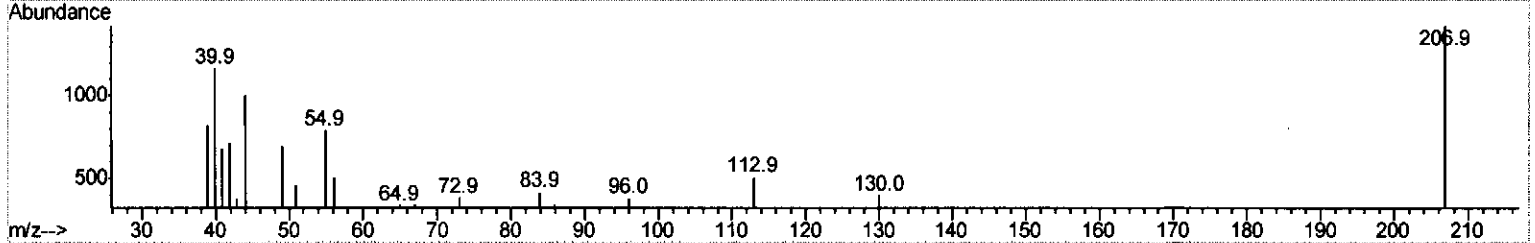
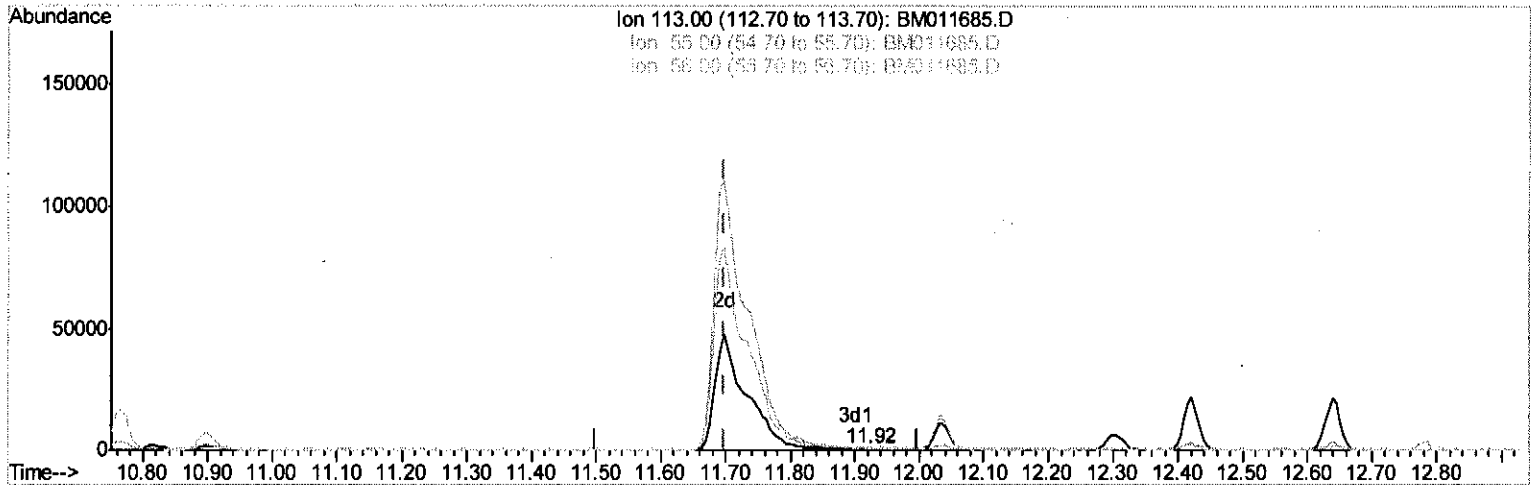
Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\  
 Data File : BM011685.D  
 Acq On : 22 Sep 2017 10:22  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampled :  
 SSTD02034

Manual Integrations  
 APPROVED

Sohil  
 9/25/2017 6:23:19 PM

Quant Time: Sep 22 23:37:00 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 22 08:34:40 2017  
 Response via : Initial Calibration



TIC: BM011685.D

(32) Caprolactam

11.922min (+0.223) 0.02ng/ul

response 179

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 164.60 | 156.13 |
| 56.00  | 115.90 | 98.81  |
| 0.00   | 0.00   | 0.00   |

# Quantitation Report (Qedit)

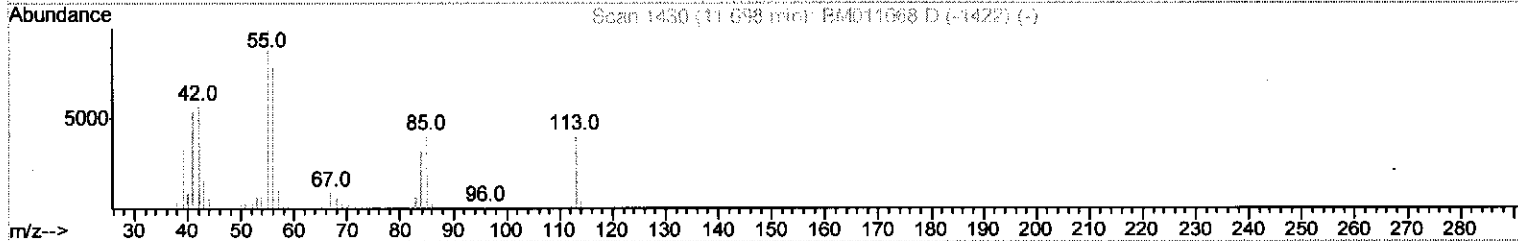
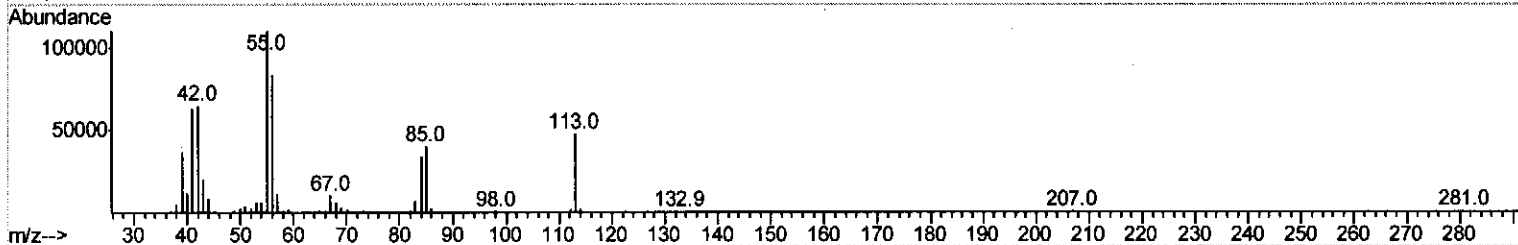
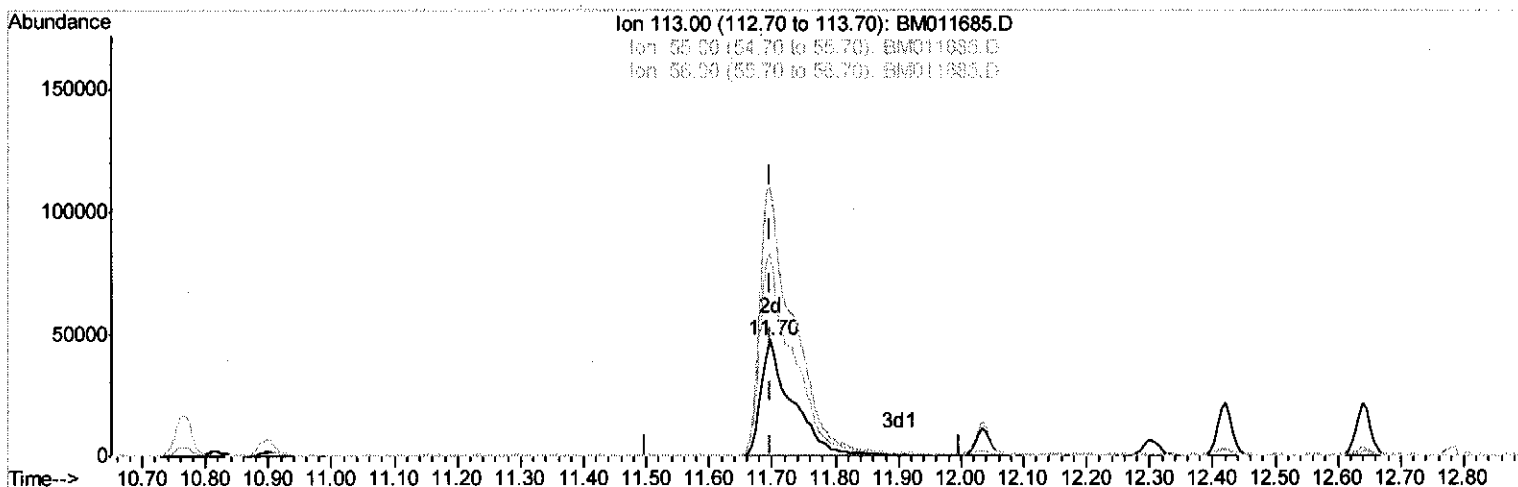
Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\  
 Data File : BM011685.D  
 Acq On : 22 Sep 2017 10:22  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Manual Integrations  
 APPROVED

Sohil  
 9/25/2017 6:23:19 PM

Quant Time: Sep 22 23:37:00 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 22 08:34:40 2017  
 Response via : Initial Calibration



TIC: BM011685.D

(32) Caprolactam

11.698min (-0.000) 18.65ng/ul mju 10/20/17

response 160952

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100    | 100     |
| 55.00  | 164.60 | 231.73# |
| 56.00  | 115.90 | 175.38# |
| 0.00   | 0.00   | 0.00    |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\  
 Data File : BM011685.D  
 Acq On : 22 Sep 2017 10:22  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Manual Integrations  
 APPROVED

Sohil  
 9/25/2017 6:23:19 PM

Quant Time: Sep 22 23:38:54 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 22 08:34:40 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.96  | 152  | 391412   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.77 | 136  | 1789551  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.59 | 164  | 979656   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.33 | 188  | 2068058  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.50 | 240  | 1730016  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.87 | 264  | 1320973  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.39  | 96  | 63968   | 7.35  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.11  | 99  | 741190  | 19.72 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.28  | 67  | 540481  | 21.34 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.49  | 132 | 585542  | 20.67 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.66  | 113 | 577444  | 19.68 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.12  | 128 | 289410  | 21.30 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.85  | 143 | 307874  | 21.80 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.38 | 165 | 571275  | 20.10 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.90 | 131 | 741008  | 23.62 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.99 | 166 | 1626576 | 19.64 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.29 | 160 | 2056565 | 20.53 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.77 | 143 | 279604  | 17.96 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.58 | 176 | 1407974 | 19.63 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.70 | 200 | 227859  | 18.94 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.43 | 188 | 2073311 | 20.36 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.72 | 212 | 2135775 | 26.37 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.71 | 264 | 1280571 | 20.34 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue |     |
|--------------------------------|-------|-----|---------|--------|--------|-----|
| 2) 1,4-Dioxane                 | 3.42  | 88  | 70257   | 7.366  | ng/uL# | 59  |
| 4) Benzaldehyde                | 7.10  | 77  | 450896  | 20.995 | ng/ul  | 98  |
| 6) Phenol                      | 7.14  | 94  | 749158  | 20.081 | ng/ul# | 91  |
| 8) Bis(2-Chloroethyl)ether     | 7.37  | 93  | 588552  | 20.038 | ng/ul# | 81  |
| 10) 2-Chlorophenol             | 7.52  | 128 | 563368  | 20.389 | ng/ul# | 89  |
| 11) 2-Methylphenol             | 8.39  | 108 | 557516  | 19.709 | ng/ul  | 93  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.49  | 45  | 1158950 | 24.344 | ng/ul# | 90  |
| 14) Acetophenone               | 8.79  | 105 | 911745  | 20.329 | ng/ul# | 89  |
| 15) N-Nitroso-di-n-propylamine | 8.77  | 70  | 528694  | 21.968 | ng/ul# | 71  |
| 16) 4-Methylphenol             | 8.72  | 108 | 611203  | 19.747 | ng/ul  | 97  |
| 17) Hexachloroethane           | 9.04  | 117 | 236502  | 22.390 | ng/ul# | 80  |
| 20) Nitrobenzene               | 9.17  | 77  | 738162  | 23.108 | ng/ul  | 98  |
| 21) Isophorone                 | 9.69  | 82  | 1364156 | 21.210 | ng/ul  | 96  |
| 23) 2-Nitrophenol              | 9.88  | 139 | 313380  | 21.120 | ng/ul  | 91  |
| 24) 2,4-Dimethylphenol         | 9.93  | 107 | 665554  | 20.963 | ng/ul  | 96  |
| 25) Bis(2-Chloroethoxy)methane | 10.17 | 93  | 821640  | 19.592 | ng/ul  | 97  |
| 27) 2,4-Dichlorophenol         | 10.41 | 162 | 545710  | 19.910 | ng/ul  | 98  |
| 28) Naphthalene                | 10.82 | 128 | 1830876 | 19.730 | ng/ul  | 98  |
| 30) 4-Chloroaniline            | 10.92 | 127 | 703666  | 23.606 | ng/ul  | 99  |
| 31) Hexachlorobutadiene        | 11.09 | 225 | 357760  | 22.998 | ng/ul  | 97  |
| 32) Caprolactam                | 11.70 | 113 | 160952m | 18.654 | ng/ul  | 97  |
| 33) 4-Chloro-3-methylphenol    | 12.03 | 107 | 585420  | 20.146 | ng/ul  | 97  |
| 34) 2-Methylnaphthalene        | 12.42 | 142 | 1310725 | 19.020 | ng/ul  | 100 |

74 10/3/17

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\  
 Data File : BM011685.D  
 Acq On : 22 Sep 2017 10:22  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02034

Manual Integrations  
 APPROVED

Sohil  
 9/25/2017 6:23:19 PM

Quant Time: Sep 22 23:38:54 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 22 08:34:40 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.64 | 142  | 1244340  | 18.958 | ng/ul  | 96       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.79 | 216  | 657745   | 22.634 | ng/ul# | 94       |
| 38) Hexachlorocyclopentadiene  | 12.76 | 237  | 318037   | 19.525 | ng/ul# | 99       |
| 39) 2,4,6-Trichlorophenol      | 13.03 | 196  | 438865   | 22.181 | ng/ul  | 98       |
| 40) 2,4,5-Trichlorophenol      | 13.09 | 196  | 439270   | 21.896 | ng/ul  | 97       |
| 41) 1,1'-Biphenyl              | 13.43 | 154  | 1643125  | 20.156 | ng/ul# | 95       |
| 42) 2-Chloronaphthalene        | 13.47 | 162  | 1265077  | 20.624 | ng/ul  | 98       |
| 43) 2-Nitroaniline             | 13.67 | 65   | 465513   | 24.435 | ng/ul# | 78       |
| 45) Dimethylphthalate          | 14.04 | 163  | 1553716  | 19.594 | ng/ul# | 98       |
| 46) 2,6-Dinitrotoluene         | 14.17 | 165  | 304364   | 21.734 | ng/ul# | 88       |
| 48) Acenaphthylene             | 14.32 | 152  | 2051244  | 20.347 | ng/ul  | 99       |
| 49) 3-Nitroaniline             | 14.50 | 138  | 326967   | 18.935 | ng/ul  | 96       |
| 50) Acenaphthene               | 14.66 | 153  | 1349705  | 19.814 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 14.70 | 184  | 147860   | 17.089 | ng/ul# | 63       |
| 53) 4-Nitrophenol              | 14.79 | 109  | 229345   | 22.927 | ng/ul# | 78       |
| 54) Dibenzofuran               | 14.99 | 168  | 1865751  | 19.581 | ng/ul  | 100      |
| 55) 2,4-Dinitrotoluene         | 14.95 | 165  | 423995   | 20.596 | ng/ul# | 66       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.22 | 232  | 391512   | 21.413 | ng/ul# | 83       |
| 57) Diethylphthalate           | 15.40 | 149  | 1590402  | 19.273 | ng/ul  | 96       |
| 59) Fluorene                   | 15.64 | 166  | 1542471  | 19.316 | ng/ul  | 100      |
| 60) 4-Chlorophenyl-phenylether | 15.63 | 204  | 775351   | 20.723 | ng/ul  | 92       |
| 61) 4-Nitroaniline             | 15.66 | 138  | 346204   | 18.716 | ng/ul  | 96       |
| 64) 4,6-Dinitro-2-methylphenol | 15.72 | 198  | 231200   | 18.750 | ng/ul# | 88       |
| 65) N-Nitrosodiphenylamine     | 15.84 | 169  | 1255465  | 19.930 | ng/ul  | 98       |
| 66) 4-Bromophenyl-phenylether  | 16.52 | 248  | 462096   | 21.877 | ng/ul  | 95       |
| 67) Hexachlorobenzene          | 16.64 | 284  | 505271   | 21.731 | ng/ul# | 88       |
| 68) Atrazine                   | 16.79 | 200  | 450441   | 20.638 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 16.98 | 266  | 302545   | 20.265 | ng/ul  | 98       |
| 70) Phenanthrene               | 17.37 | 178  | 2276243  | 19.747 | ng/ul  | 99       |
| 72) Anthracene                 | 17.47 | 178  | 2379251  | 20.115 | ng/ul  | 98       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.39 | 216  | 649807   | 23.580 | ng/uL  | 99       |
| 74) Pentachlorobenzene         | 14.91 | 250  | 643733   | 23.032 | ng/uL  | 98       |
| 75) Carbazole                  | 17.73 | 167  | 2031646  | 20.133 | ng/ul  | 97       |
| 76) Di-n-butylphthalate        | 18.28 | 149  | 2748647  | 20.737 | ng/ul  | 100      |
| 77) Fluoranthene               | 19.39 | 202  | 2589567  | 21.841 | ng/ul# | 94       |
| 80) Pvrene                     | 19.74 | 202  | 2661704  | 26.370 | ng/ul# | 89       |
| 81) Butylbenzylphthalate       | 20.63 | 149  | 1210897  | 26.086 | ng/ul  | 90       |
| 82) 3,3'-Dichlorobenzidine     | 21.42 | 252  | 586059   | 18.654 | ng/ul# | 99       |
| 83) Benzo(a)anthracene         | 21.49 | 228  | 2094394  | 21.987 | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.39 | 149  | 1824064  | 25.588 | ng/ul# | 95       |
| 85) Chrysene                   | 21.54 | 228  | 1852973  | 21.457 | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.30 | 149  | 2899525  | 30.928 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.15 | 252  | 1613596  | 20.305 | ng/ul# | 97       |
| 89) Benzo(k)fluoranthene       | 23.19 | 252  | 1626485  | 21.311 | ng/ul# | 95       |
| 91) Benzo(a)pyrene             | 23.76 | 252  | 1512925  | 20.170 | ng/ul# | 97       |
| 92) Indeno(1,2,3-cd)pyrene     | 26.30 | 276  | 1699308  | 20.594 | ng/ul# | 89       |
| 93) Dibenzo(a,h)anthracene     | 26.31 | 278  | 1450936  | 20.738 | ng/ul# | 96       |
| 94) Benzo(a,h,i)perylene       | 27.05 | 276  | 1395230  | 20.880 | ng/ul# | 94       |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092217\  
Data File : BM011685.D  
Acq On : 22 Sep 2017 10:22  
Operator : SJ/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
LabSampleId :  
SSTD02034

Manual Integrations  
APPROVED

Sohil  
9/25/2017 6:23:19 PM

Quant Time: Sep 22 23:38:54 2017  
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Sep 22 08:34:40 2017  
Response via : Initial Calibration

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

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