

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042920\  
 Data File : BN010666.D  
 Acq On : 29 Apr 2020 21:55  
 Operator : CG/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 SSTD02090

Quant Time: Apr 30 04:13:41 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 30 03:53:20 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.58  | 152  | 535495   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.33 | 136  | 2338879  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.20 | 164  | 1504389  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.96 | 188  | 3365470  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.16 | 240  | 2928451  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.32 | 264  | 2662189  | 20.00 | ng/ul | -0.01    |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.20  | 96  | 78108   | 7.63  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 6.77  | 99  | 836734  | 19.79 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.93  | 67  | 483885  | 22.63 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.12  | 132 | 695285  | 19.29 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.29  | 113 | 695004  | 19.50 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.72  | 128 | 351543  | 19.51 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.43  | 143 | 395498  | 19.69 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 9.97  | 165 | 772080  | 19.61 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 10.47 | 131 | 1022108 | 22.64 | ng/ul | 0.00  |
| 44) Dimethylphthalate-d6       | 13.63 | 166 | 2221994 | 19.91 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 13.89 | 160 | 2631523 | 19.19 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 14.42 | 143 | 390019  | 17.95 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 15.20 | 176 | 1915432 | 19.12 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.33 | 200 | 330310  | 18.79 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 17.05 | 188 | 3032500 | 19.45 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 19.36 | 212 | 3310170 | 20.02 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 23.19 | 264 | 2582371 | 18.70 | ng/ul | -0.02 |

## Target Compounds

|                                |       |     |         |       |        | Qvalue |
|--------------------------------|-------|-----|---------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.23  | 88  | 81592   | 7.54  | ng/uL# | 92     |
| 4) Benzaldehyde                | 6.73  | 77  | 562041  | 25.60 | ng/ul  | 98     |
| 6) Phenol                      | 6.80  | 94  | 861789  | 19.69 | ng/ul  | 98     |
| 8) Bis(2-Chloroethyl)ether     | 7.02  | 93  | 656720  | 19.49 | ng/ul  | 96     |
| 10) 2-Chlorophenol             | 7.15  | 128 | 722332  | 19.70 | ng/ul  | 98     |
| 11) 2-Methylphenol             | 8.02  | 108 | 668629  | 19.30 | ng/ul  | 96     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.11  | 45  | 448247  | 20.92 | ng/ul  | 99     |
| 14) Acetophenone               | 8.39  | 105 | 1117752 | 20.74 | ng/ul  | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.39  | 70  | 526612  | 20.40 | ng/ul  | 96     |
| 16) 4-Methylphenol             | 8.35  | 108 | 729375  | 19.68 | ng/ul  | 100    |
| 17) Hexachloroethane           | 8.64  | 117 | 285742  | 18.79 | ng/ul  | 87     |
| 20) Nitrobenzene               | 8.76  | 77  | 834390  | 19.82 | ng/ul  | 98     |
| 21) Isophorone                 | 9.29  | 82  | 1452591 | 18.71 | ng/ul  | 99     |
| 23) 2-Nitrophenol              | 9.46  | 139 | 437357  | 20.14 | ng/ul  | 96     |
| 24) 2,4-Dimethylphenol         | 9.53  | 107 | 831270  | 19.41 | ng/ul  | 97     |
| 25) Bis(2-Chloroethoxy)methane | 9.77  | 93  | 934074  | 19.78 | ng/ul  | 98     |
| 27) 2,4-Dichlorophenol         | 9.99  | 162 | 747947  | 19.28 | ng/ul  | 98     |
| 28) Naphthalene                | 10.39 | 128 | 2373113 | 18.88 | ng/ul  | 100    |
| 30) 4-Chloroaniline            | 10.50 | 127 | 1056005 | 23.56 | ng/ul  | 100    |
| 31) Hexachlorobutadiene        | 10.68 | 225 | 494222  | 18.30 | ng/ul  | 99     |
| 32) Caprolactam                | 11.28 | 113 | 214947  | 17.46 | ng/ul  | 99     |
| 33) 4-Chloro-3-methylphenol    | 11.63 | 107 | 796765  | 19.60 | ng/ul  | 97     |
| 34) 2-Methylnaphthalene        | 12.00 | 142 | 1786995 | 20.12 | ng/ul  | 97     |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042920\  
 Data File : BN010666.D  
 Acq On : 29 Apr 2020 21:55  
 Operator : CG/JU  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD02090

Quant Time: Apr 30 04:13:41 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 30 03:53:20 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.23 | 142  | 1677155  | 18.86 | ng/uL  | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.38 | 216  | 958631   | 19.11 | ng/uL  | 93       |
| 38) Hexachlorocyclopentadiene  | 12.37 | 237  | 374669   | 15.35 | ng/uL  | 96       |
| 39) 2,4,6-Trichlorophenol      | 12.63 | 196  | 612888   | 18.86 | ng/uL  | 98       |
| 40) 2,4,5-Trichlorophenol      | 12.70 | 196  | 665911   | 19.05 | ng/uL  | 98       |
| 41) 1,1'-Biphenyl              | 13.03 | 154  | 2271506  | 19.66 | ng/uL  | 99       |
| 42) 2-Chloronaphthalene        | 13.07 | 162  | 1785747  | 19.12 | ng/uL  | 98       |
| 43) 2-Nitroaniline             | 13.27 | 65   | 452923   | 20.91 | ng/uL  | 92       |
| 45) Dimethylphthalate          | 13.67 | 163  | 2194677  | 18.94 | ng/uL  | 100      |
| 46) 2,6-Dinitrotoluene         | 13.79 | 165  | 484405   | 19.66 | ng/uL  | 100      |
| 48) Acenaphthylene             | 13.92 | 152  | 2813094  | 19.44 | ng/uL  | 99       |
| 49) 3-Nitroaniline             | 14.11 | 138  | 466935   | 19.91 | ng/uL  | 90       |
| 50) Acenaphthene               | 14.27 | 153  | 1884344  | 19.03 | ng/uL  | 97       |
| 51) 2,4-Dinitrophenol          | 14.32 | 184  | 196286   | 16.81 | ng/uL  | 97       |
| 53) 4-Nitrophenol              | 14.44 | 109  | 318625   | 18.11 | ng/uL  | 95       |
| 54) Dibenzofuran               | 14.61 | 168  | 2681482  | 19.34 | ng/uL  | 100      |
| 55) 2,4-Dinitrotoluene         | 14.58 | 165  | 675470   | 19.24 | ng/uL  | 98       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.84 | 232  | 589863   | 19.31 | ng/uL  | 97       |
| 57) Diethylphthalate           | 15.05 | 149  | 2175929  | 18.68 | ng/uL  | 99       |
| 59) Fluorene                   | 15.26 | 166  | 2198218  | 19.34 | ng/uL  | 97       |
| 60) 4-Chlorophenyl-phenylether | 15.26 | 204  | 1108987  | 19.08 | ng/uL  | 99       |
| 61) 4-Nitroaniline             | 15.28 | 138  | 494833   | 18.63 | ng/uL  | 93       |
| 64) 4,6-Dinitro-2-methylphenol | 15.35 | 198  | 363629   | 19.23 | ng/uL# | 92       |
| 65) N-Nitrosodiphenylamine     | 15.47 | 169  | 1914846  | 19.64 | ng/uL  | 97       |
| 66) 4-Bromophenyl-phenylether  | 16.16 | 248  | 687326   | 18.71 | ng/uL  | 96       |
| 67) Hexachlorobenzene          | 16.27 | 284  | 762763   | 18.25 | ng/uL  | 97       |
| 68) Atrazine                   | 16.44 | 200  | 608862   | 15.81 | ng/uL  | 99       |
| 69) Pentachlorophenol          | 16.62 | 266  | 438663   | 16.44 | ng/uL  | 95       |
| 70) Phenanthrene               | 17.00 | 178  | 3543546  | 18.96 | ng/uL  | 98       |
| 72) Anthracene                 | 17.09 | 178  | 3655904  | 19.30 | ng/uL  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.00 | 216  | 937136   | 17.56 | ng/uL  | 97       |
| 74) Pentachlorobenzene         | 14.53 | 250  | 896489   | 17.23 | ng/uL  | 95       |
| 75) Carbazole                  | 17.36 | 167  | 3181733  | 20.45 | ng/uL  | 99       |
| 76) Di-n-butylphthalate        | 17.95 | 149  | 3829351  | 19.53 | ng/uL  | 99       |
| 77) Fluoranthene               | 19.02 | 202  | 4189008  | 20.63 | ng/uL  | 96       |
| 80) Pyrene                     | 19.39 | 202  | 4239103  | 20.03 | ng/uL  | 97       |
| 81) Butylbenzylphthalate       | 20.31 | 149  | 1730162  | 18.96 | ng/uL  | 95       |
| 82) 3,3'-Dichlorobenzidine     | 21.09 | 252  | 1097424  | 17.39 | ng/uL  | 97       |
| 83) Benzo(a)anthracene         | 21.14 | 228  | 3783937  | 18.63 | ng/uL  | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 21.10 | 149  | 2562187  | 19.44 | ng/uL  | 99       |
| 85) Chrysene                   | 21.20 | 228  | 3438689  | 18.00 | ng/uL  | 95       |
| 87) Di-n-octyl phthalate       | 21.97 | 149  | 4095680  | 21.63 | ng/uL  | 100      |
| 88) Benzo(b)fluoranthene       | 22.68 | 252  | 3362502  | 19.36 | ng/uL  | 98       |
| 89) Benzo(k)fluoranthene       | 22.73 | 252  | 3144476  | 18.24 | ng/uL  | 98       |
| 91) Benzo(a)pyrene             | 23.23 | 252  | 3081989  | 19.97 | ng/uL  | 97       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.46 | 276  | 3405976  | 19.61 | ng/uL  | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.47 | 278  | 2851915  | 19.55 | ng/uL  | 99       |
| 94) Benzo(g,h,i)perylene       | 26.11 | 276  | 2766861  | 19.25 | ng/uL  | 98       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042920\  
Data File : BN010666.D  
Acq On : 29 Apr 2020 21:55  
Operator : CG/JU  
Sample : SSTDCCC020EC  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD02090

Quant Time: Apr 30 04:13:41 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Apr 30 03:53:20 2020  
Response via : Initial Calibration

| Internal Standards                                                      | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                   |      |      |          |      |       |          |
| (#)= qualifier out of range (m)= manual integration (+)= signals summed |      |      |          |      |       |          |

