

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042020\  
 Data File : BN010555.D  
 Acq On : 20 Apr 2020 11:09  
 Operator : CG/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 SSTD02080

Manual Integrations  
 APPROVED

mohammad  
 4/20/2020 2:27:43 PM

Quant Time: Apr 20 11:54:18 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 20 11:50:51 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.59  | 152  | 351904   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.35 | 136  | 1603759  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.22 | 164  | 1074008  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.97 | 188  | 2544652  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.18 | 240  | 2853452  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.34 | 264  | 2674472  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.20  | 96  | 51017   | 7.59  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.78  | 99  | 556770  | 20.04 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.94  | 67  | 318028  | 22.63 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.13  | 132 | 471053  | 19.89 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.30  | 113 | 469910  | 20.06 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.73  | 128 | 231294  | 18.72 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 258083  | 18.74 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.98  | 165 | 522193  | 19.34 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.49 | 131 | 721713  | 23.31 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.64 | 166 | 1548286 | 19.43 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.91 | 160 | 1897679 | 19.38 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.43 | 143 | 294540  | 18.98 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.22 | 176 | 1392604 | 19.47 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.35 | 200 | 236882  | 17.82 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.07 | 188 | 2309457 | 19.59 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.37 | 212 | 2816469 | 17.49 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.21 | 264 | 2678797 | 19.31 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue |     |
|--------------------------------|-------|-----|---------|--------|--------|-----|
| 2) 1,4-Dioxane                 | 3.23  | 88  | 53647   | 7.540  | ng/uL  | 96  |
| 4) Benzaldehyde                | 6.75  | 77  | 362601  | 25.131 | ng/ul  | 93  |
| 6) Phenol                      | 6.80  | 94  | 576463  | 20.043 | ng/ul  | 98  |
| 8) Bis(2-Chloroethyl)ether     | 7.03  | 93  | 424349  | 19.161 | ng/ul  | 95  |
| 10) 2-Chlorophenol             | 7.17  | 128 | 485971  | 20.163 | ng/ul  | 95  |
| 11) 2-Methylphenol             | 8.03  | 108 | 445277  | 19.556 | ng/ul  | 96  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.12  | 45  | 290293  | 20.615 | ng/ul  | 98  |
| 14) Acetophenone               | 8.40  | 105 | 744638  | 21.024 | ng/ul  | 99  |
| 15) N-Nitroso-di-n-propylamine | 8.40  | 70  | 345901  | 20.389 | ng/ul  | 97  |
| 16) 4-Methylphenol             | 8.36  | 108 | 493958  | 20.280 | ng/ul  | 96  |
| 17) Hexachloroethane           | 8.66  | 117 | 184596  | 18.467 | ng/ul  | 92  |
| 20) Nitrobenzene               | 8.78  | 77  | 559677  | 19.390 | ng/ul  | 99  |
| 21) Isophorone                 | 9.30  | 82  | 978469  | 18.378 | ng/ul  | 98  |
| 23) 2-Nitrophenol              | 9.47  | 139 | 279653  | 18.776 | ng/ul  | 96  |
| 24) 2,4-Dimethylphenol         | 9.55  | 107 | 560328  | 19.076 | ng/ul  | 99  |
| 25) Bis(2-Chloroethoxy)methane | 9.78  | 93  | 613094  | 18.930 | ng/ul  | 97  |
| 27) 2,4-Dichlorophenol         | 10.00 | 162 | 521222  | 19.595 | ng/ul  | 99  |
| 28) Naphthalene                | 10.40 | 128 | 1636266 | 18.984 | ng/ul  | 99  |
| 30) 4-Chloroaniline            | 10.51 | 127 | 734803  | 23.904 | ng/ul  | 100 |
| 31) Hexachlorobutadiene        | 10.70 | 225 | 345222  | 18.644 | ng/ul  | 91  |
| 32) Caprolactam                | 11.27 | 113 | 148276  | 17.567 | ng/ul  | 97  |
| 33) 4-Chloro-3-methylphenol    | 11.65 | 107 | 544963  | 19.551 | ng/ul  | 97  |
| 34) 2-Methylnaphthalene        | 12.02 | 142 | 1220614 | 20.038 | ng/ul  | 98  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042020\  
 Data File : BN010555.D  
 Acq On : 20 Apr 2020 11:09  
 Operator : CG/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 SSTD02080

Manual Integrations  
 APPROVED

mohammad  
 4/20/2020 2:27:43 PM

Quant Time: Apr 20 11:54:18 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Apr 20 11:50:51 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.24 | 142  | 1157693  | 18.989 | ng/uL  | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.40 | 216  | 672095   | 18.770 | ng/uL  | 96       |
| 38) Hexachlorocyclopentadiene  | 12.39 | 237  | 266662   | 15.305 | ng/uL  | 95       |
| 39) 2,4,6-Trichlorophenol      | 12.64 | 196  | 428967   | 18.487 | ng/uL  | 97       |
| 40) 2,4,5-Trichlorophenol      | 12.71 | 196  | 473560   | 18.977 | ng/uL  | 98       |
| 41) 1,1'-Biphenyl              | 13.05 | 154  | 1595823  | 19.347 | ng/uL  | 99       |
| 42) 2-Chloronaphthalene        | 13.09 | 162  | 1281865  | 19.225 | ng/uL  | 100      |
| 43) 2-Nitroaniline             | 13.29 | 65   | 310397   | 20.068 | ng/uL  | 97       |
| 45) Dimethylphthalate          | 13.69 | 163  | 1548393  | 18.715 | ng/uL  | 100      |
| 46) 2,6-Dinitrotoluene         | 13.80 | 165  | 350410   | 19.916 | ng/uL  | 99       |
| 48) Acenaphthylene             | 13.94 | 152  | 2030738  | 19.656 | ng/uL  | 99       |
| 49) 3-Nitroaniline             | 14.12 | 138  | 338674   | 20.231 | ng/uL  | 94       |
| 50) Acenaphthene               | 14.29 | 153  | 1365704  | 19.317 | ng/uL  | 97       |
| 51) 2,4-Dinitrophenol          | 14.33 | 184  | 137098   | 16.450 | ng/uL  | 93       |
| 53) 4-Nitrophenol              | 14.45 | 109  | 251402   | 20.016 | ng/uL  | 98       |
| 54) Dibenzofuran               | 14.62 | 168  | 1995003  | 20.157 | ng/uL  | 99       |
| 55) 2,4-Dinitrotoluene         | 14.59 | 165  | 488565   | 19.492 | ng/uL# | 98       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.86 | 232  | 432572m  | 19.837 | ng/uL  |          |
| 57) Diethylphthalate           | 15.06 | 149  | 1542577  | 18.545 | ng/uL  | 100      |
| 59) Fluorene                   | 15.27 | 166  | 1559038  | 19.212 | ng/uL  | 97       |
| 60) 4-Chlorophenyl-phenylether | 15.27 | 204  | 802952   | 19.348 | ng/uL  | 98       |
| 61) 4-Nitroaniline             | 15.29 | 138  | 412900   | 21.776 | ng/uL  | 100      |
| 64) 4,6-Dinitro-2-methylphenol | 15.36 | 198  | 261160   | 18.269 | ng/uL# | 89       |
| 65) N-Nitrosodiphenylamine     | 15.49 | 169  | 1417862  | 19.234 | ng/uL  | 97       |
| 66) 4-Bromophenyl-phenylether  | 16.17 | 248  | 484767   | 17.449 | ng/uL  | 99       |
| 67) Hexachlorobenzene          | 16.29 | 284  | 557495   | 17.641 | ng/uL  | 99       |
| 68) Atrazine                   | 16.45 | 200  | 455150   | 15.631 | ng/uL  | 98       |
| 69) Pentachlorophenol          | 16.63 | 266  | 354102   | 17.552 | ng/uL  | 96       |
| 70) Phenanthrene               | 17.01 | 178  | 2752803  | 19.480 | ng/uL  | 98       |
| 72) Anthracene                 | 17.10 | 178  | 2745789  | 19.168 | ng/uL  | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.01 | 216  | 676528   | 16.763 | ng/uL  | 96       |
| 74) Pentachlorobenzene         | 14.55 | 250  | 646902   | 16.443 | ng/uL  | 98       |
| 75) Carbazole                  | 17.37 | 167  | 2551675  | 21.686 | ng/uL  | 99       |
| 76) Di-n-butylphthalate        | 17.96 | 149  | 2720427  | 18.354 | ng/uL  | 100      |
| 77) Fluoranthene               | 19.03 | 202  | 3451314  | 22.478 | ng/uL  | 97       |
| 80) Pvrene                     | 19.40 | 202  | 3625406  | 17.584 | ng/uL  | 96       |
| 81) Butylbenzylphthalate       | 20.33 | 149  | 1411183  | 15.869 | ng/uL  | 98       |
| 82) 3,3'-Dichlorobenzidine     | 21.10 | 252  | 1000621  | 16.270 | ng/uL  | 95       |
| 83) Benzo(a)anthracene         | 21.16 | 228  | 3626558  | 18.321 | ng/uL  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.12 | 149  | 2103475  | 16.377 | ng/uL  | 98       |
| 85) Chrysene                   | 21.22 | 228  | 3394346  | 18.237 | ng/uL  | 99       |
| 87) Di-n-octyl phthalate       | 21.99 | 149  | 3681425  | 19.352 | ng/uL  | 100      |
| 88) Benzo(b)fluoranthene       | 22.70 | 252  | 3388316  | 19.415 | ng/uL  | 96       |
| 89) Benzo(k)fluoranthene       | 22.75 | 252  | 3237104  | 18.696 | ng/uL  | 97       |
| 91) Benzo(a)pyrene             | 23.26 | 252  | 3111417  | 20.067 | ng/uL  | 98       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.49 | 276  | 3309601  | 18.963 | ng/uL  | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.50 | 278  | 2789613  | 19.031 | ng/uL  | 98       |
| 94) Benzo(a,h,i)perylene       | 26.14 | 276  | 2755825  | 19.080 | ng/uL  | 96       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN042020\  
Data File : BN010555.D  
Acq On : 20 Apr 2020 11:09  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
SSTD02080

Manual Integrations  
APPROVED

mohammad  
4/20/2020 2:27:43 PM

Quant Time: Apr 20 11:54:18 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040920MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Apr 20 11:50:51 2020  
Response via : Initial Calibration

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

