

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF111522\  
 Data File : BF131237.D  
 Acq On : 15 Nov 2022 17:41  
 Operator : CG\JU  
 Sample : SSTDICCC040  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICCC040

Quant Time: Nov 16 00:27:21 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF111522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 15 22:44:23 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Christian  
 Giraldo

11/16/2022  
 Supervised By :Jagrut  
 Upadhyay

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.975  | 152  | 102349   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.269  | 136  | 367123   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 10.033 | 164  | 201928   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.527 | 188  | 365201   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 14.186 | 240  | 256429   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 15.727 | 264  | 193495   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.575  | 112  | 461236   | 76.638 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.592  | 99   | 583411   | 74.148 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.545  | 82   | 544397   | 66.906 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.828 | 330  | 158087   | 75.753 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.351  | 172  | 1059235  | 73.227 | ng    | 0.00     |
| 79) Terphenyl-d14             | 13.127 | 244  | 1197935  | 81.908 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.787  | 88   | 107220   | 36.194 | ng    | 100      |
| 3) Pyridine                   | 3.557  | 79   | 301573   | 37.075 | ng    | 100      |
| 4) n-Nitrosodimethylamine     | 3.528  | 42   | 178241   | 37.986 | ng    | 100      |
| 6) Aniline                    | 6.640  | 93   | 371118m  | 38.422 | ng    |          |
| 8) 2-Chlorophenol             | 6.757  | 128  | 257746   | 37.227 | ng    | 100      |
| 9) Benzaldehyde               | 6.528  | 77   | 188021   | 36.722 | ng    | 100      |
| 10) Phenol                    | 6.604  | 94   | 325192   | 35.355 | ng    | 100      |
| 11) bis(2-Chloroethyl)ether   | 6.710  | 93   | 253917   | 38.586 | ng    | 100      |
| 12) 1,3-Dichlorobenzene       | 6.916  | 146  | 292265   | 37.959 | ng    | 100      |
| 13) 1,4-Dichlorobenzene       | 6.992  | 146  | 295384   | 37.767 | ng    | 100      |
| 14) 1,2-Dichlorobenzene       | 7.151  | 146  | 271567   | 37.485 | ng    | 100      |
| 15) Benzyl Alcohol            | 7.116  | 79   | 259920   | 43.192 | ng    | 100      |
| 16) 2,2'-oxybis(1-Chloropr... | 7.251  | 45   | 487557   | 44.188 | ng    | 100      |
| 17) 2-Methylphenol            | 7.216  | 107  | 223106   | 38.528 | ng    | 100      |
| 18) Hexachloroethane          | 7.492  | 117  | 108306   | 37.138 | ng    | 100      |
| 19) n-Nitroso-di-n-propyla... | 7.392  | 70   | 201209   | 38.948 | ng    | 100      |
| 20) 3+4-Methylphenols         | 7.375  | 107  | 289945   | 40.075 | ng    | 100      |
| 22) Acetophenone              | 7.387  | 105  | 368956   | 40.159 | ng    | 100      |
| 24) Nitrobenzene              | 7.563  | 77   | 291227   | 36.948 | ng    | 100      |
| 25) Isophorone                | 7.804  | 82   | 510800   | 40.600 | ng    | 100      |
| 26) 2-Nitrophenol             | 7.881  | 139  | 105032   | 30.032 | ng    | 100      |
| 27) 2,4-Dimethylphenol        | 7.910  | 122  | 212225   | 39.569 | ng    | 100      |
| 28) bis(2-Chloroethoxy)met... | 8.010  | 93   | 286681   | 40.208 | ng    | 100      |
| 29) 2,4-Dichlorophenol        | 8.116  | 162  | 230224   | 40.151 | ng    | 100      |
| 30) 1,2,4-Trichlorobenzene    | 8.204  | 180  | 264717   | 42.812 | ng    | 100      |
| 31) Naphthalene               | 8.292  | 128  | 775218   | 38.646 | ng    | 100      |
| 32) Benzoic acid              | 8.022  | 122  | 140882   | 41.612 | ng    | 100      |
| 33) 4-Chloroaniline           | 8.334  | 127  | 303326   | 37.710 | ng    | 100      |
| 34) Hexachlorobutadiene       | 8.404  | 225  | 170963   | 40.465 | ng    | 100      |
| 35) Caprolactam               | 8.716  | 113  | 67582    | 38.447 | ng    | 100      |
| 36) 4-Chloro-3-methylphenol   | 8.810  | 107  | 238509   | 37.588 | ng    | 100      |
| 37) 2-Methylnaphthalene       | 8.986  | 142  | 510263   | 39.461 | ng    | 100      |
| 38) 1-Methylnaphthalene       | 9.086  | 142  | 499375   | 39.434 | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 9.145  | 216  | 255930   | 38.813 | ng    | 100      |
| 41) Hexachlorocyclopentadiene | 9.133  | 237  | 140792   | 54.723 | ng    | 100      |
| 43) 2,4,6-Trichlorophenol     | 9.257  | 196  | 175877   | 40.643 | ng    | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF111522\  
 Data File : BF131237.D  
 Acq On : 15 Nov 2022 17:41  
 Operator : CG\JU  
 Sample : SSTDICCC040  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICCC040

Quant Time: Nov 16 00:27:21 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF111522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 15 22:44:23 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Christian  
 Giraldo

11/16/2022  
 Supervised By :Jagrut  
 Upadhyay

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 9.292  | 196  | 182494   | 37.693 | ng    | 100      |
| 46) 1,1'-Biphenyl             | 9.451  | 154  | 608200   | 38.955 | ng    | 100      |
| 47) 2-Chloronaphthalene       | 9.475  | 162  | 489497   | 38.636 | ng    | 100      |
| 48) 2-Nitroaniline            | 9.569  | 65   | 152460   | 33.064 | ng    | 100      |
| 49) Acenaphthylene            | 9.898  | 152  | 761205   | 39.320 | ng    | 100      |
| 50) Dimethylphthalate         | 9.751  | 163  | 570941   | 39.749 | ng    | 100      |
| 51) 2,6-Dinitrotoluene        | 9.816  | 165  | 111717   | 35.391 | ng    | 100      |
| 52) Acenaphthene              | 10.069 | 154  | 472282   | 40.898 | ng    | 100      |
| 53) 3-Nitroaniline            | 9.986  | 138  | 117811   | 32.644 | ng    | 100      |
| 54) 2,4-Dinitrophenol         | 10.086 | 184  | 38026    | 22.950 | ng    | 100      |
| 55) Dibenzofuran              | 10.239 | 168  | 670869   | 39.229 | ng    | 100      |
| 56) 4-Nitrophenol             | 10.128 | 139  | 92682    | 38.585 | ng    | 100      |
| 57) 2,4-Dinitrotoluene        | 10.216 | 165  | 134378   | 32.024 | ng    | 100      |
| 58) Fluorene                  | 10.586 | 166  | 526616   | 37.914 | ng    | 100      |
| 59) 2,3,4,6-Tetrachlorophenol | 10.351 | 232  | 159732   | 42.906 | ng    | 100      |
| 60) Diethylphthalate          | 10.457 | 149  | 542100   | 39.188 | ng    | 100      |
| 61) 4-Chlorophenyl-phenyle... | 10.575 | 204  | 265081   | 41.009 | ng    | 100      |
| 62) 4-Nitroaniline            | 10.604 | 138  | 116653   | 31.985 | ng    | 100      |
| 63) Azobenzene                | 10.739 | 77   | 566434   | 39.120 | ng    | 100      |
| 65) 4,6-Dinitro-2-methylph... | 10.628 | 198  | 59620    | 25.391 | ng    | 100      |
| 66) n-Nitrosodiphenylamine    | 10.698 | 169  | 455045   | 37.063 | ng    | 100      |
| 67) 4-Bromophenyl-phenylether | 11.069 | 248  | 172054   | 41.664 | ng    | 100      |
| 68) Hexachlorobenzene         | 11.133 | 284  | 184986   | 40.194 | ng    | 100      |
| 69) Atrazine                  | 11.222 | 200  | 151668   | 39.906 | ng    | 100      |
| 70) Pentachlorophenol         | 11.322 | 266  | 113859   | 56.429 | ng    | 100      |
| 71) Phenanthrene              | 11.557 | 178  | 787685   | 39.216 | ng    | 100      |
| 72) Anthracene                | 11.610 | 178  | 778546   | 39.937 | ng    | 100      |
| 73) Carbazole                 | 11.763 | 167  | 698216   | 40.276 | ng    | 100      |
| 74) Di-n-butylphthalate       | 12.092 | 149  | 829196   | 42.409 | ng    | 100      |
| 75) Fluoranthene              | 12.751 | 202  | 877435   | 43.641 | ng    | 100      |
| 77) Benzidine                 | 12.869 | 184  | 235718   | 31.159 | ng    | 100      |
| 78) Pyrene                    | 12.980 | 202  | 891274   | 40.512 | ng    | 100      |
| 80) Butylbenzylphthalate      | 13.604 | 149  | 318377   | 40.420 | ng    | 100      |
| 81) Benzo(a)anthracene        | 14.174 | 228  | 707234   | 38.586 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 14.139 | 252  | 212488   | 40.618 | ng    | 100      |
| 83) Chrysene                  | 14.216 | 228  | 695110   | 40.157 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 14.163 | 149  | 389390   | 45.795 | ng    | 100      |
| 85) Di-n-octyl phthalate      | 14.798 | 149  | 539113   | 49.998 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 17.315 | 276  | 542352   | 40.321 | ng    | 100      |
| 88) Benzo(b)fluoranthene      | 15.268 | 252  | 544708   | 39.817 | ng    | 100      |
| 89) Benzo(k)fluoranthene      | 15.304 | 252  | 523055   | 39.779 | ng    | 100      |
| 90) Benzo(a)pyrene            | 15.662 | 252  | 436749   | 39.821 | ng    | 100      |
| 91) Dibenzo(a,h)anthracene    | 17.345 | 278  | 445170   | 40.765 | ng    | 100      |
| 92) Benzo(g,h,i)perylene      | 17.798 | 276  | 448902   | 39.454 | ng    | 100      |

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

## Manual Integrations

11/16/2022  
Supervised By :Jagrut  
Upadhyay

11/16/2022

