

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG100522\  
 Data File : BG055075.D  
 Acq On : 5 Oct 2022 12:13  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: Oct 06 03:36:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 04 07:12:22 2022  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                    | Amount | Calc.   | %Dev   | Area% | Dev(min) |
|------|-----------------------------|--------|---------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 20.000 | 20.000  | 0.0    | 113   | 0.00     |
| 2    | 1,4-Dioxane                 | 40.000 | 37.730  | 5.7    | 110   | 0.00     |
| 3    | Pyridine                    | 40.000 | 39.110  | 2.2    | 111   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 40.000 | 37.788  | 5.5    | 104   | 0.00     |
| 5 S  | 2-Fluorophenol              | 80.000 | 92.813  | -16.0  | 126   | 0.00     |
| 6    | Aniline                     | 40.000 | 40.230  | -0.6   | 110   | 0.00     |
| 7 S  | Phenol-d6                   | 80.000 | 91.343  | -14.2  | 121   | 0.00     |
| 8    | 2-Chlorophenol              | 40.000 | 47.289  | -18.2  | 127   | 0.00     |
| 9    | Benzaldehyde                | 40.000 | 37.427  | 6.4    | 108   | 0.00     |
| 10 C | Phenol                      | 40.000 | 44.809  | -12.0  | 119   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 39.014  | 2.5    | 109   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 40.258  | -0.6   | 115   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 39.890  | 0.3    | 113   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 39.770  | 0.6    | 114   | 0.00     |
| 15   | Benzyl Alcohol              | 40.000 | 46.054  | -15.1  | 122   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 37.590  | 6.0    | 103   | 0.00     |
| 17   | 2-Methylphenol              | 40.000 | 43.586  | -9.0   | 117   | 0.00     |
| 18   | Hexachloroethane            | 40.000 | 43.280  | -8.2   | 119   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 40.168  | -0.4   | 111   | 0.00     |
| 20   | 3+4-Methylphenols           | 40.000 | 46.760  | -16.9  | 124   | 0.00     |
| 21 I | Naphthalene-d8              | 20.000 | 20.000  | 0.0    | 109   | 0.00     |
| 22   | Acetophenone                | 40.000 | 40.157  | -0.4   | 108   | 0.00     |
| 23 S | Nitrobenzene-d5             | 80.000 | 90.654  | -13.3  | 116   | 0.00     |
| 24   | Nitrobenzene                | 40.000 | 43.831  | -9.6   | 112   | 0.00     |
| 25   | Isophorone                  | 40.000 | 43.354  | -8.4   | 113   | 0.00     |
| 26 C | 2-Nitrophenol               | 40.000 | 59.386  | -48.5# | 155   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 40.000 | 47.746  | -19.4  | 123   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 41.959  | -4.9   | 110   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 45.063  | -12.7  | 131   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 40.980  | -2.4   | 109   | 0.00     |
| 31   | Naphthalene                 | 40.000 | 40.828  | -2.1   | 109   | 0.00     |
| 32   | Benzoic acid                | 40.000 | 55.601  | -39.0# | 171   | 0.00     |
| 33   | 4-Chloroaniline             | 40.000 | 42.588  | -6.5   | 113   | 0.00     |
| 34 C | Hexachlorobutadiene         | 40.000 | 41.696  | -4.2   | 111   | 0.00     |
| 35   | Caprolactam                 | 40.000 | 55.546  | -38.9# | 143   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 45.424  | -13.6  | 125   | 0.00     |
| 37   | 2-Methylnaphthalene         | 40.000 | 40.822  | -2.1   | 109   | 0.00     |
| 38   | 1-Methylnaphthalene         | 40.000 | 41.082  | -2.7   | 110   | 0.00     |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000  | 0.0    | 112   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 39.744  | 0.6    | 110   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 47.076  | -17.7  | 130   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 104.127 | -30.2# | 149   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 46.821  | -17.1  | 142   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 46.656  | -16.6  | 138   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 78.881  | 1.4    | 110   | 0.00     |
| 46   | 1,1'-Biphenyl               | 40.000 | 39.999  | 0.0    | 109   | 0.00     |
| 47   | 2-Chloronaphthalene         | 40.000 | 40.001  | -0.0   | 110   | 0.00     |

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                   | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|----------------------------|--------|--------|--------|-------|----------|
| 48   | 2-Nitroaniline             | 40.000 | 47.069 | -17.7  | 142   | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 39.909 | 0.2    | 109   | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 46.982 | -17.5  | 126   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 45.467 | -13.7  | 134   | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 41.213 | -3.0   | 115   | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 45.611 | -14.0  | 132   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 51.376 | -28.4# | 144   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 41.110 | -2.8   | 113   | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 49.152 | -22.9  | 149   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 46.632 | -16.6  | 138   | 0.00     |
| 58   | Fluorene                   | 40.000 | 40.775 | -1.9   | 113   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 48.792 | -22.0  | 149   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 48.711 | -21.8  | 131   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 40.809 | -2.0   | 113   | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 48.807 | -22.0  | 129   | 0.00     |
| 63   | Azobenzene                 | 40.000 | 39.318 | 1.7    | 111   | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 118   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 51.907 | -29.8# | 155   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 41.258 | -3.1   | 117   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 42.246 | -5.6   | 119   | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 41.081 | -2.7   | 117   | 0.00     |
| 69   | Atrazine                   | 40.000 | 50.363 | -25.9# | 135   | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 49.724 | -24.3# | 158   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 40.615 | -1.5   | 117   | 0.00     |
| 72   | Anthracene                 | 40.000 | 40.698 | -1.7   | 116   | 0.00     |
| 73   | Carbazole                  | 40.000 | 43.092 | -7.7   | 122   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 51.409 | -28.5# | 139   | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 40.609 | -1.5   | 117   | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 113   | 0.00     |
| 77   | Benzydine                  | 40.000 | 54.591 | -36.5# | 139   | 0.00     |
| 78   | Pyrene                     | 40.000 | 42.004 | -5.0   | 117   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 83.597 | -4.5   | 116   | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 49.630 | -24.1  | 152   | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 41.428 | -3.6   | 113   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 48.493 | -21.2  | 127   | 0.00     |
| 83   | Chrysene                   | 40.000 | 40.443 | -1.1   | 112   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 47.020 | -17.6  | 140   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 40.000 | 47.289 | -18.2  | 143   | 0.00     |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 112   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 41.479 | -3.7   | 113   | 0.00     |
| 88   | Benzo(b)fluoranthene       | 40.000 | 41.202 | -3.0   | 111   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 40.000 | 40.538 | -1.3   | 111   | 0.00     |
| 90 C | Benzo(a)pyrene             | 40.000 | 41.166 | -2.9   | 112   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 42.344 | -5.9   | 117   | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 41.340 | -3.4   | 113   | 0.00     |

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Max. RRF Dev : 25% Max. Rel. Area : 150%

| Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|----------|--------|-------|------|-------|----------|
|----------|--------|-------|------|-------|----------|

(#) = Out of Range

SPCC's out = 0 CCC's out = 2