

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022119\  
 Data File : BG039572.D  
 Acq On : 21 Feb 2019 13:11  
 Operator : JU/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: Feb 21 16:58:10 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 21 16:54:20 2019  
 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    | 152   | -0.02    |
| 2    | 1,4-Dioxane                  | 40.000 | 34.589 | 13.5   | 134   | -0.02    |
| 3    | Pyridine                     | 40.000 | 40.266 | -0.7   | 146   | 0.00     |
| 4    | n-Nitrosodimethylamine       | 40.000 | 36.831 | 7.9    | 141   | 0.00     |
| 5 S  | 2-Fluorophenol               | 80.000 | 79.907 | 0.1    | 154   | 0.00     |
| 6    | Aniline                      | 40.000 | 41.113 | -2.8   | 160   | 0.00     |
| 7 S  | Phenol-d6                    | 80.000 | 86.851 | -8.6   | 165   | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 42.919 | -7.3   | 163   | -0.02    |
| 9    | Benzaldehyde                 | 40.000 | 45.659 | -14.1  | 168   | -0.01    |
| 10 C | Phenol                       | 40.000 | 43.280 | -8.2   | 168   | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 41.913 | -4.8   | 162   | -0.02    |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 39.685 | 0.8    | 154   | -0.02    |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 40.709 | -1.8   | 157   | -0.02    |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 39.966 | 0.1    | 156   | -0.02    |
| 15   | Benzyl Alcohol               | 40.000 | 43.527 | -8.8   | 164   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 35.689 | 10.8   | 141   | -0.02    |
| 17   | 2-Methylphenol               | 40.000 | 44.270 | -10.7  | 171   | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 40.910 | -2.3   | 159   | -0.02    |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 41.543 | -3.9   | 157   | -0.02    |
| 20   | 3+4-Methylphenols            | 40.000 | 45.391 | -13.5  | 170   | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0    | 158   | -0.02    |
| 22   | Acetophenone                 | 40.000 | 39.169 | 2.1    | 162   | -0.02    |
| 23 S | Nitrobenzene-d5              | 80.000 | 94.856 | -18.6  | 193   | -0.01    |
| 24   | Nitrobenzene                 | 40.000 | 46.002 | -15.0  | 189   | -0.01    |
| 25   | Isophorone                   | 40.000 | 39.085 | 2.3    | 159   | -0.02    |
| 26 C | 2-Nitrophenol                | 40.000 | 55.732 | -39.3# | 262   | -0.02    |
| 27   | 2,4-Dimethylphenol           | 40.000 | 42.096 | -5.2   | 172   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 40.426 | -1.1   | 162   | -0.02    |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 44.766 | -11.9  | 181   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 40.887 | -2.2   | 165   | -0.02    |
| 31   | Naphthalene                  | 40.000 | 38.480 | 3.8    | 160   | -0.02    |
| 32   | Benzoic acid                 | 40.000 | 52.086 | -30.2# | 245   | 0.03     |
| 33   | 4-Chloroaniline              | 40.000 | 41.920 | -4.8   | 171   | -0.02    |
| 34 C | Hexachlorobutadiene          | 40.000 | 40.762 | -1.9   | 165   | -0.02    |
| 35   | Caprolactam                  | 40.000 | 45.198 | -13.0  | 174   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 44.959 | -12.4  | 178   | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 39.261 | 1.8    | 163   | -0.02    |
| 38   | 1-Methylnaphthalene          | 40.000 | 39.552 | 1.1    | 164   | -0.02    |
| 39 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0    | 169   | -0.02    |
| 40   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 39.876 | 0.3    | 173   | -0.02    |
| 41 P | Hexachlorocyclopentadiene    | 40.000 | 43.470 | -8.7   | 187   | -0.02    |
| 42 S | 2,4,6-Tribromophenol         | 80.000 | 91.302 | -14.1  | 195   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol        | 40.000 | 45.294 | -13.2  | 189   | 0.00     |
| 44   | 2,4,5-Trichlorophenol        | 40.000 | 45.707 | -14.3  | 186   | 0.00     |

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

|      | Compound                   | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|----------------------------|--------|--------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 80.000 | 73.842 | 7.7    | 163   | -0.02    |
| 46   | 1,1'-Biphenyl              | 40.000 | 37.673 | 5.8    | 168   | -0.02    |
| 47   | 2-Chloronaphthalene        | 40.000 | 39.152 | 2.1    | 173   | -0.02    |
| 48   | 2-Nitroaniline             | 40.000 | 48.202 | -20.5  | 234   | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 38.170 | 4.6    | 167   | -0.02    |
| 50   | Dimethylphthalate          | 40.000 | 38.578 | 3.6    | 169   | -0.02    |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 47.391 | -18.5  | 219   | -0.01    |
| 52 C | Acenaphthene               | 40.000 | 37.569 | 6.1    | 162   | -0.02    |
| 53   | 3-Nitroaniline             | 40.000 | 46.850 | -17.1  | 217   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 55.254 | -38.1# | 276   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 38.783 | 3.0    | 171   | -0.02    |
| 56 P | 4-Nitrophenol              | 40.000 | 42.099 | -5.2   | 192   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 52.189 | -30.5# | 255   | 0.00     |
| 58   | Fluorene                   | 40.000 | 38.425 | 3.9    | 168   | -0.02    |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 44.647 | -11.6  | 186   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 37.841 | 5.4    | 168   | -0.02    |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 40.746 | -1.9   | 181   | -0.02    |
| 62   | 4-Nitroaniline             | 40.000 | 46.659 | -16.6  | 211   | 0.00     |
| 63   | Azobenzene                 | 40.000 | 35.880 | 10.3   | 159   | -0.02    |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 167   | -0.02    |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 56.189 | -40.5# | 286   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 39.097 | 2.3    | 170   | -0.02    |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 41.615 | -4.0   | 182   | -0.02    |
| 68   | Hexachlorobenzene          | 40.000 | 41.754 | -4.4   | 183   | -0.02    |
| 69   | Atrazine                   | 40.000 | 34.417 | 14.0   | 147   | -0.02    |
| 70 C | Pentachlorophenol          | 40.000 | 40.173 | -0.4   | 169   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 37.448 | 6.4    | 165   | -0.02    |
| 72   | Anthracene                 | 40.000 | 38.091 | 4.8    | 167   | -0.02    |
| 73   | Carbazole                  | 40.000 | 37.226 | 6.9    | 164   | -0.02    |
| 74   | Di-n-butylphthalate        | 40.000 | 37.325 | 6.7    | 163   | -0.03    |
| 75 C | Fluoranthene               | 40.000 | 37.545 | 6.1    | 167   | -0.02    |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 172   | -0.02    |
| 77   | Benzidine                  | 40.000 | 33.440 | 16.4   | 152   | -0.02    |
| 78   | Pyrene                     | 40.000 | 37.407 | 6.5    | 167   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 71.580 | 10.5   | 161   | -0.02    |
| 80   | Butylbenzylphthalate       | 40.000 | 39.982 | 0.0    | 174   | -0.02    |
| 81   | Benzo(a)anthracene         | 40.000 | 38.063 | 4.8    | 171   | -0.02    |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 40.572 | -1.4   | 179   | -0.02    |
| 83   | Chrysene                   | 40.000 | 38.304 | 4.2    | 170   | -0.02    |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 39.279 | 1.8    | 172   | -0.04    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 39.423 | 1.4    | 172   | -0.06    |
| 86   | Indeno(1,2,3-cd)pyrene     | 40.000 | 37.477 | 6.3    | 164   | -0.06    |
| 87 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 165   | -0.03    |

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|      | Compound               | Amount | Calc.  | %Dev | Area | % Dev(min) |
|------|------------------------|--------|--------|------|------|------------|
| 88   | Benzo(b)fluoranthene   | 40.000 | 40.322 | -0.8 | 170  | -0.03      |
| 89   | Benzo(k)fluoranthene   | 40.000 | 38.618 | 3.5  | 167  | -0.03      |
| 90 C | Benzo(a)pyrene         | 40.000 | 40.739 | -1.8 | 173  | -0.03      |
| 91   | Dibenzo(a,h)anthracene | 40.000 | 37.569 | 6.1  | 161  | -0.06      |
| 92   | Benzo(q,h,i)perylene   | 40.000 | 37.199 | 7.0  | 158  | -0.05      |

(#) = Out of Range

SPCC's out = 0 CCC's out = 1