

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG112125\  
 Data File : BG064838.D  
 Acq On : 21 Nov 2025 10:09  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: Nov 21 10:44:03 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG11225.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 12 17:10:56 2025  
 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   | 161   | 0.00     |
| 2    | 1,4-Dioxane                 | 40.000 | 40.340 | -0.9  | 154   | 0.00     |
| 3    | Pyridine                    | 40.000 | 41.933 | -4.8  | 164   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 40.000 | 38.165 | 4.6   | 151   | 0.00     |
| 5 S  | 2-Fluorophenol              | 80.000 | 80.214 | -0.3  | 159   | 0.00     |
| 6    | Aniline                     | 40.000 | 41.129 | -2.8  | 160   | 0.00     |
| 7 S  | Phenol-d6                   | 80.000 | 82.416 | -3.0  | 160   | 0.00     |
| 8    | 2-Chlorophenol              | 40.000 | 41.957 | -4.9  | 163   | 0.00     |
| 9    | Benzaldehyde                | 40.000 | 38.162 | 4.6   | 149   | 0.00     |
| 10 C | Phenol                      | 40.000 | 41.496 | -3.7  | 162   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 41.685 | -4.2  | 161   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 40.089 | -0.2  | 157   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 39.107 | 2.2   | 155   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 40.280 | -0.7  | 157   | 0.00     |
| 15   | Benzyl Alcohol              | 40.000 | 40.410 | -1.0  | 156   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 36.020 | 9.9   | 143   | -0.01    |
| 17   | 2-Methylphenol              | 40.000 | 40.659 | -1.6  | 158   | 0.00     |
| 18   | Hexachloroethane            | 40.000 | 38.940 | 2.7   | 156   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 40.914 | -2.3  | 155   | 0.00     |
| 20   | 3+4-Methylphenols           | 40.000 | 40.632 | -1.6  | 160   | 0.00     |
| 21 I | Naphthalene-d8              | 20.000 | 20.000 | 0.0   | 162   | 0.00     |
| 22   | Acetophenone                | 40.000 | 39.743 | 0.6   | 157   | 0.00     |
| 23 S | Nitrobenzene-d5             | 80.000 | 78.605 | 1.7   | 156   | 0.00     |
| 24   | Nitrobenzene                | 40.000 | 39.848 | 0.4   | 160   | 0.00     |
| 25   | Isophorone                  | 40.000 | 40.438 | -1.1  | 165   | 0.00     |
| 26 C | 2-Nitrophenol               | 40.000 | 43.822 | -9.6  | 169   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 40.000 | 40.675 | -1.7  | 163   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 41.970 | -4.9  | 170   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 43.024 | -7.6  | 168   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 39.455 | 1.4   | 160   | 0.00     |
| 31   | Naphthalene                 | 40.000 | 39.895 | 0.3   | 161   | 0.00     |
| 32   | Benzoic acid                | 40.000 | 44.069 | -10.2 | 199   | 0.02     |
| 33   | 4-Chloroaniline             | 40.000 | 42.075 | -5.2  | 164   | 0.00     |
| 34 C | Hexachlorobutadiene         | 40.000 | 39.256 | 1.9   | 158   | 0.00     |
| 35   | Caprolactam                 | 40.000 | 42.726 | -6.8  | 176   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 41.794 | -4.5  | 170   | 0.00     |
| 37   | 2-Methylnaphthalene         | 40.000 | 40.000 | 0.0   | 163   | 0.00     |
| 38   | 1-Methylnaphthalene         | 40.000 | 40.128 | -0.3  | 165   | 0.00     |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000 | 0.0   | 170   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 38.242 | 4.4   | 162   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 35.207 | 12.0  | 149   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 80.094 | -0.1  | 173   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 39.781 | 0.5   | 166   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 40.392 | -1.0  | 169   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 75.696 | 5.4   | 163   | 0.00     |
| 46   | 1,1'-Biphenyl               | 40.000 | 38.492 | 3.8   | 165   | 0.00     |
| 47   | 2-Chloronaphthalene         | 40.000 | 39.054 | 2.4   | 168   | 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 | 40.493 | -1.2  | 169   | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 39.039 | 2.4   | 167   | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 38.824 | 2.9   | 169   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 40.624 | -1.6  | 172   | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 40.574 | -1.4  | 172   | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 42.162 | -5.4  | 176   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 41.306 | -3.3  | 192   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 38.918 | 2.7   | 170   | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 42.970 | -7.4  | 181   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 41.542 | -3.9  | 174   | 0.00     |
| 58   | Fluorene                   | 40.000 | 38.141 | 4.6   | 166   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 41.598 | -4.0  | 179   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 39.399 | 1.5   | 170   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 38.995 | 2.5   | 170   | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 42.352 | -5.9  | 175   | 0.00     |
| 63   | Azobenzene                 | 40.000 | 38.871 | 2.8   | 169   | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0   | 167   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 45.377 | -13.4 | 180   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 41.699 | -4.2  | 169   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 40.880 | -2.2  | 166   | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 39.824 | 0.4   | 165   | 0.00     |
| 69   | Atrazine                   | 40.000 | 39.546 | 1.1   | 165   | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 43.693 | -9.2  | 199   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 39.653 | 0.9   | 165   | 0.00     |
| 72   | Anthracene                 | 40.000 | 40.053 | -0.1  | 168   | 0.00     |
| 73   | Carbazole                  | 40.000 | 39.436 | 1.4   | 163   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 40.319 | -0.8  | 167   | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 37.627 | 5.9   | 157   | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0   | 138   | 0.00     |
| 77   | Benzydine                  | 40.000 | 32.834 | 17.9  | 121   | 0.00     |
| 78   | Pyrene                     | 40.000 | 44.812 | -12.0 | 156   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 85.792 | -7.2  | 149   | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 42.880 | -7.2  | 150   | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 39.439 | 1.4   | 135   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 36.569 | 8.6   | 123   | 0.00     |
| 83   | Chrysene                   | 40.000 | 39.672 | 0.8   | 135   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 41.320 | -3.3  | 141   | -0.01    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 39.130 | 2.2   | 130   | -0.02    |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0   | 119   | -0.02    |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 41.015 | -2.5  | 120   | -0.01    |
| 88   | Benzo(b)fluoranthene       | 40.000 | 39.833 | 0.4   | 117   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 40.000 | 39.697 | 0.8   | 121   | 0.00     |
| 90 C | Benzo(a)pyrene             | 40.000 | 39.831 | 0.4   | 118   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 41.097 | -2.7  | 120   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 41.370 | -3.4  | 121   | -0.02    |

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QLast Update : Wed Nov 12 17:10:56 2025  
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) |
|----------|--------|-------|------|-------|----------|
|----------|--------|-------|------|-------|----------|

(#) = Out of Range

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