

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081425\  
 Data File : BF143397.D  
 Acq On : 14 Aug 2025 22:56  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Aug 15 01:35:30 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF081225.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 12 13:25:39 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    | 91    | 0.00     |
| 2    | 1,4-Dioxane                 | 40.000 | 34.533 | 13.7   | 79    | -0.03    |
| 3    | Pyridine                    | 40.000 | 34.717 | 13.2   | 78    | -0.02    |
| 4    | n-Nitrosodimethylamine      | 40.000 | 34.645 | 13.4   | 77    | -0.02    |
| 5 S  | 2-Fluorophenol              | 80.000 | 73.205 | 8.5    | 82    | 0.00     |
| 6    | Aniline                     | 40.000 | 34.171 | 14.6   | 77    | 0.00     |
| 7 S  | Phenol-d6                   | 80.000 | 70.261 | 12.2   | 79    | 0.00     |
| 8    | 2-Chlorophenol              | 40.000 | 37.322 | 6.7    | 83    | 0.00     |
| 9    | Benzaldehyde                | 40.000 | 32.651 | 18.4   | 72    | 0.00     |
| 10 C | Phenol                      | 40.000 | 35.396 | 11.5   | 79    | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 35.405 | 11.5   | 80    | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 35.839 | 10.4   | 81    | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 35.936 | 10.2   | 83    | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 35.880 | 10.3   | 82    | 0.00     |
| 15   | Benzyl Alcohol              | 40.000 | 36.604 | 8.5    | 81    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 33.764 | 15.6   | 77    | 0.00     |
| 17   | 2-Methylphenol              | 40.000 | 35.456 | 11.4   | 79    | 0.00     |
| 18   | Hexachloroethane            | 40.000 | 38.942 | 2.6    | 86    | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 33.566 | 16.1   | 77    | 0.00     |
| 20   | 3+4-Methylphenols           | 40.000 | 35.190 | 12.0   | 78    | 0.00     |
| 21 I | Naphthalene-d8              | 20.000 | 20.000 | 0.0    | 86    | 0.00     |
| 22   | Acetophenone                | 40.000 | 34.852 | 12.9   | 77    | 0.00     |
| 23 S | Nitrobenzene-d5             | 80.000 | 77.681 | 2.9    | 80    | 0.00     |
| 24   | Nitrobenzene                | 40.000 | 37.898 | 5.3    | 79    | 0.00     |
| 25   | Isophorone                  | 40.000 | 34.512 | 13.7   | 75    | 0.00     |
| 26 C | 2-Nitrophenol               | 40.000 | 45.217 | -13.0  | 107   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 40.000 | 37.878 | 5.3    | 80    | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 35.741 | 10.6   | 78    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 38.587 | 3.5    | 81    | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 37.101 | 7.2    | 81    | 0.00     |
| 31   | Naphthalene                 | 40.000 | 35.988 | 10.0   | 79    | 0.00     |
| 32   | Benzoic acid                | 40.000 | 40.763 | -1.9   | 91    | 0.00     |
| 33   | 4-Chloroaniline             | 40.000 | 35.380 | 11.5   | 76    | 0.00     |
| 34 C | Hexachlorobutadiene         | 40.000 | 38.848 | 2.9    | 84    | 0.00     |
| 35   | Caprolactam                 | 40.000 | 36.300 | 9.3    | 75    | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 35.367 | 11.6   | 74    | 0.00     |
| 37   | 2-Methylnaphthalene         | 40.000 | 35.329 | 11.7   | 77    | 0.00     |
| 38   | 1-Methylnaphthalene         | 40.000 | 34.624 | 13.4   | 76    | 0.00     |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000 | 0.0    | 80    | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 38.695 | 3.3    | 78    | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 51.058 | -27.6# | 112   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 80.730 | -0.9   | 74    | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 41.879 | -4.7   | 78    | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 39.313 | 1.7    | 73    | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 74.656 | 6.7    | 78    | 0.00     |
| 46   | 1,1'-Biphenyl               | 40.000 | 37.371 | 6.6    | 76    | 0.00     |
| 47   | 2-Chloronaphthalene         | 40.000 | 37.462 | 6.3    | 77    | 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 | 38.817 | 3.0    | 80    | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 36.573 | 8.6    | 74    | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 37.666 | 5.8    | 75    | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 40.788 | -2.0   | 84    | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 36.423 | 8.9    | 74    | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 41.020 | -2.6   | 76    | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 35.584 | 11.0   | 69    | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 36.113 | 9.7    | 74    | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 30.288 | 24.3   | 59    | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 39.688 | 0.8    | 80    | 0.00     |
| 58   | Fluorene                   | 40.000 | 35.081 | 12.3   | 73    | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 39.774 | 0.6    | 74    | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 36.866 | 7.8    | 72    | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 36.239 | 9.4    | 73    | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 36.245 | 9.4    | 68    | 0.00     |
| 63   | Azobenzene                 | 40.000 | 33.836 | 15.4   | 68    | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 73    | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 43.140 | -7.9   | 88    | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 38.245 | 4.4    | 70    | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 39.924 | 0.2    | 72    | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 38.074 | 4.8    | 70    | 0.00     |
| 69   | Atrazine                   | 40.000 | 40.779 | -1.9   | 72    | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 32.879 | 17.8   | 59    | 0.00     |
| 71   | Phenanthrene               | 40.000 | 35.962 | 10.1   | 66    | 0.00     |
| 72   | Anthracene                 | 40.000 | 36.392 | 9.0    | 68    | 0.00     |
| 73   | Carbazole                  | 40.000 | 34.117 | 14.7   | 63    | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 42.631 | -6.6   | 72    | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 34.718 | 13.2   | 66    | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 85    | 0.00     |
| 77   | Benzydine                  | 40.000 | 37.244 | 6.9    | 73    | 0.00     |
| 78   | Pyrene                     | 40.000 | 30.918 | 22.7   | 66    | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 63.756 | 20.3   | 70    | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 50.287 | -25.7# | 117   | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 35.746 | 10.6   | 77    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 44.526 | -11.3  | 89    | 0.00     |
| 83   | Chrysene                   | 40.000 | 36.741 | 8.1    | 79    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 51.120 | -27.8# | 121   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 40.000 | 50.655 | -26.6# | 119   | 0.01     |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 105   | 0.01     |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 37.381 | 6.5    | 97    | 0.02     |
| 88   | Benzo(b)fluoranthene       | 40.000 | 37.954 | 5.1    | 98    | 0.00     |
| 89   | Benzo(k)fluoranthene       | 40.000 | 32.831 | 17.9   | 89    | 0.01     |
| 90 C | Benzo(a)pyrene             | 40.000 | 36.515 | 8.7    | 96    | 0.01     |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 38.021 | 4.9    | 98    | 0.02     |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 37.096 | 7.3    | 96    | 0.02     |

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|----------|--------|-------|------|-------|----------|
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(#) = Out of Range                      SPCC's out = 0    CCC's out = 1