

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060525\  
 Data File : BF142625.D  
 Acq On : 05 Jun 2025 09:37  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jun 05 11:34:22 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 20 16:26:47 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  | 125   | 0.00     |
| 2    | 1,4-Dioxane                 | 40.000 | 40.410 | -1.0 | 131   | 0.00     |
| 3    | Pyridine                    | 40.000 | 40.947 | -2.4 | 133   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 40.000 | 38.290 | 4.3  | 122   | -0.02    |
| 5 S  | 2-Fluorophenol              | 80.000 | 77.896 | 2.6  | 129   | 0.00     |
| 6    | Aniline                     | 40.000 | 39.471 | 1.3  | 129   | -0.01    |
| 7 S  | Phenol-d6                   | 80.000 | 77.312 | 3.4  | 127   | 0.00     |
| 8    | 2-Chlorophenol              | 40.000 | 39.178 | 2.1  | 127   | 0.00     |
| 9    | Benzaldehyde                | 40.000 | 37.107 | 7.2  | 122   | 0.00     |
| 10 C | Phenol                      | 40.000 | 38.996 | 2.5  | 128   | -0.01    |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 39.048 | 2.4  | 128   | -0.01    |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 38.592 | 3.5  | 126   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 38.687 | 3.3  | 125   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 38.413 | 4.0  | 126   | 0.00     |
| 15   | Benzyl Alcohol              | 40.000 | 38.121 | 4.7  | 124   | -0.01    |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 38.613 | 3.5  | 126   | -0.01    |
| 17   | 2-Methylphenol              | 40.000 | 38.723 | 3.2  | 125   | 0.00     |
| 18   | Hexachloroethane            | 40.000 | 38.760 | 3.1  | 126   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 36.981 | 7.5  | 122   | -0.02    |
| 20   | 3+4-Methylphenols           | 40.000 | 37.528 | 6.2  | 122   | 0.00     |
| 21 I | Naphthalene-d8              | 20.000 | 20.000 | 0.0  | 127   | 0.00     |
| 22   | Acetophenone                | 40.000 | 37.065 | 7.3  | 122   | -0.01    |
| 23 S | Nitrobenzene-d5             | 80.000 | 75.377 | 5.8  | 123   | -0.01    |
| 24   | Nitrobenzene                | 40.000 | 37.704 | 5.7  | 123   | -0.01    |
| 25   | Isophorone                  | 40.000 | 36.930 | 7.7  | 121   | -0.01    |
| 26 C | 2-Nitrophenol               | 40.000 | 39.397 | 1.5  | 126   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 40.000 | 38.580 | 3.6  | 126   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 37.793 | 5.5  | 126   | -0.01    |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 38.280 | 4.3  | 125   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 37.811 | 5.5  | 125   | 0.00     |
| 31   | Naphthalene                 | 40.000 | 37.509 | 6.2  | 125   | 0.00     |
| 32   | Benzoic acid                | 40.000 | 36.415 | 9.0  | 117   | -0.04    |
| 33   | 4-Chloroaniline             | 40.000 | 38.402 | 4.0  | 125   | 0.00     |
| 34 C | Hexachlorobutadiene         | 40.000 | 36.666 | 8.3  | 120   | 0.00     |
| 35   | Caprolactam                 | 40.000 | 38.036 | 4.9  | 122   | -0.02    |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 36.714 | 8.2  | 120   | -0.01    |
| 37   | 2-Methylnaphthalene         | 40.000 | 37.375 | 6.6  | 125   | 0.00     |
| 38   | 1-Methylnaphthalene         | 40.000 | 37.001 | 7.5  | 122   | 0.00     |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000 | 0.0  | 123   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 38.259 | 4.4  | 123   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 34.650 | 13.4 | 105   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 73.469 | 8.2  | 119   | -0.01    |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 38.593 | 3.5  | 124   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 37.970 | 5.1  | 121   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 71.781 | 10.3 | 119   | -0.01    |
| 46   | 1,1'-Biphenyl               | 40.000 | 37.345 | 6.6  | 121   | -0.01    |
| 47   | 2-Chloronaphthalene         | 40.000 | 37.561 | 6.1  | 121   | 0.00     |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060525\  
 Data File : BF142625.D  
 Acq On : 05 Jun 2025 09:37  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jun 05 11:34:22 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 20 16:26:47 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) |
|------|----------------------------|--------|--------|-------|-------|----------|
| 48   | 2-Nitroaniline             | 40.000 | 38.573 | 3.6   | 122   | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 37.526 | 6.2   | 121   | -0.01    |
| 50   | Dimethylphthalate          | 40.000 | 37.211 | 7.0   | 121   | -0.01    |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 38.373 | 4.1   | 120   | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 37.134 | 7.2   | 121   | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 38.576 | 3.6   | 124   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 35.863 | 10.3  | 109   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 37.075 | 7.3   | 121   | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 35.358 | 11.6  | 110   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 39.006 | 2.5   | 123   | -0.01    |
| 58   | Fluorene                   | 40.000 | 37.066 | 7.3   | 122   | -0.01    |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 37.830 | 5.4   | 119   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 36.575 | 8.6   | 118   | -0.01    |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 37.050 | 7.4   | 121   | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 37.923 | 5.2   | 119   | -0.01    |
| 63   | Azobenzene                 | 40.000 | 37.006 | 7.5   | 119   | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0   | 121   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 40.095 | -0.2  | 117   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 37.790 | 5.5   | 121   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 37.992 | 5.0   | 122   | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 38.159 | 4.6   | 121   | 0.00     |
| 69   | Atrazine                   | 40.000 | 36.706 | 8.2   | 114   | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 35.392 | 11.5  | 108   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 37.001 | 7.5   | 118   | 0.00     |
| 72   | Anthracene                 | 40.000 | 37.793 | 5.5   | 120   | 0.00     |
| 73   | Carbazole                  | 40.000 | 36.874 | 7.8   | 117   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 35.120 | 12.2  | 111   | -0.01    |
| 75 C | Fluoranthene               | 40.000 | 36.208 | 9.5   | 117   | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0   | 125   | 0.00     |
| 77   | Benzydine                  | 40.000 | 39.060 | 2.3   | 114   | 0.00     |
| 78   | Pyrene                     | 40.000 | 36.770 | 8.1   | 115   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 69.729 | 12.8  | 112   | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 39.243 | 1.9   | 120   | -0.01    |
| 81   | Benzo(a)anthracene         | 40.000 | 37.407 | 6.5   | 122   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 44.828 | -12.1 | 140   | 0.00     |
| 83   | Chrysene                   | 40.000 | 38.373 | 4.1   | 126   | -0.01    |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 45.647 | -14.1 | 140   | -0.01    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 44.953 | -12.4 | 143   | -0.01    |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0   | 136   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 34.708 | 13.2  | 123   | -0.01    |
| 88   | Benzo(b)fluoranthene       | 40.000 | 39.903 | 0.2   | 143   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 40.000 | 33.683 | 15.8  | 122   | -0.01    |
| 90 C | Benzo(a)pyrene             | 40.000 | 37.849 | 5.4   | 133   | -0.01    |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 34.758 | 13.1  | 122   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 34.574 | 13.6  | 122   | -0.02    |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060525\  
Data File : BF142625.D  
Acq On : 05 Jun 2025 09:37  
Operator : RC/JU  
Sample : SSTDCCC040  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_F  
LabSampleId :  
SSTDCCC040

Quant Time: Jun 05 11:34:22 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue May 20 16:26:47 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