

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF022425\  
 Data File : BF141734.D  
 Acq On : 24 Feb 2025 09:52  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Feb 24 11:50:52 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF020625.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 06 16:58:59 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    | 118   | -0.01    |
| 2    | 1,4-Dioxane                 | 40.000 | 39.955 | 0.1    | 117   | -0.02    |
| 3    | Pyridine                    | 40.000 | 42.593 | -6.5   | 121   | -0.02    |
| 4    | n-Nitrosodimethylamine      | 40.000 | 36.168 | 9.6    | 105   | -0.04    |
| 5 S  | 2-Fluorophenol              | 80.000 | 82.619 | -3.3   | 120   | -0.01    |
| 6    | Aniline                     | 40.000 | 38.415 | 4.0    | 111   | -0.02    |
| 7 S  | Phenol-d6                   | 80.000 | 80.912 | -1.1   | 119   | -0.01    |
| 8    | 2-Chlorophenol              | 40.000 | 41.675 | -4.2   | 122   | -0.01    |
| 9    | Benzaldehyde                | 40.000 | 35.392 | 11.5   | 108   | 0.00     |
| 10 C | Phenol                      | 40.000 | 39.970 | 0.1    | 117   | -0.02    |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 41.397 | -3.5   | 121   | -0.01    |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 41.469 | -3.7   | 123   | -0.01    |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 41.118 | -2.8   | 122   | -0.01    |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 41.510 | -3.8   | 123   | -0.01    |
| 15   | Benzyl Alcohol              | 40.000 | 38.748 | 3.1    | 111   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 36.997 | 7.5    | 108   | -0.01    |
| 17   | 2-Methylphenol              | 40.000 | 40.348 | -0.9   | 119   | -0.01    |
| 18   | Hexachloroethane            | 40.000 | 41.091 | -2.7   | 120   | -0.01    |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 39.600 | 1.0    | 116   | -0.02    |
| 20   | 3+4-Methylphenols           | 40.000 | 40.106 | -0.3   | 118   | -0.02    |
| 21 I | Naphthalene-d8              | 20.000 | 20.000 | 0.0    | 118   | -0.01    |
| 22   | Acetophenone                | 40.000 | 39.603 | 1.0    | 117   | -0.02    |
| 23 S | Nitrobenzene-d5             | 80.000 | 79.027 | 1.2    | 116   | -0.01    |
| 24   | Nitrobenzene                | 40.000 | 39.250 | 1.9    | 115   | -0.01    |
| 25   | Isophorone                  | 40.000 | 39.691 | 0.8    | 117   | -0.01    |
| 26 C | 2-Nitrophenol               | 40.000 | 41.819 | -4.5   | 119   | -0.01    |
| 27   | 2,4-Dimethylphenol          | 40.000 | 40.869 | -2.2   | 117   | -0.01    |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 41.074 | -2.7   | 120   | -0.01    |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 41.080 | -2.7   | 120   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 41.242 | -3.1   | 122   | -0.01    |
| 31   | Naphthalene                 | 40.000 | 40.202 | -0.5   | 119   | -0.01    |
| 32   | Benzoic acid                | 40.000 | 36.197 | 9.5    | 105   | -0.02    |
| 33   | 4-Chloroaniline             | 40.000 | 40.108 | -0.3   | 118   | -0.02    |
| 34 C | Hexachlorobutadiene         | 40.000 | 40.820 | -2.1   | 120   | -0.01    |
| 35   | Caprolactam                 | 40.000 | 37.900 | 5.3    | 107   | -0.03    |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 39.830 | 0.4    | 116   | -0.01    |
| 37   | 2-Methylnaphthalene         | 40.000 | 40.317 | -0.8   | 120   | -0.01    |
| 38   | 1-Methylnaphthalene         | 40.000 | 39.889 | 0.3    | 117   | -0.01    |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000 | 0.0    | 115   | -0.02    |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 41.402 | -3.5   | 118   | -0.01    |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 53.675 | -34.2# | 146   | -0.01    |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 77.011 | 3.7    | 111   | -0.02    |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 40.669 | -1.7   | 116   | -0.01    |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 40.093 | -0.2   | 114   | -0.01    |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 79.837 | 0.2    | 118   | -0.01    |
| 46   | 1,1'-Biphenyl               | 40.000 | 40.676 | -1.7   | 118   | -0.02    |
| 47   | 2-Chloronaphthalene         | 40.000 | 40.362 | -0.9   | 118   | -0.02    |

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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 | 37.784 | 5.5    | 107   | -0.02    |
| 49   | Acenaphthylene             | 40.000 | 39.638 | 0.9    | 115   | -0.01    |
| 50   | Dimethylphthalate          | 40.000 | 39.993 | 0.0    | 116   | -0.02    |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 40.054 | -0.1   | 114   | -0.01    |
| 52 C | Acenaphthene               | 40.000 | 38.999 | 2.5    | 112   | -0.02    |
| 53   | 3-Nitroaniline             | 40.000 | 37.105 | 7.2    | 107   | -0.01    |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 40.016 | -0.0   | 115   | -0.01    |
| 55   | Dibenzofuran               | 40.000 | 38.888 | 2.8    | 113   | -0.02    |
| 56 P | 4-Nitrophenol              | 40.000 | 52.281 | -30.7# | 143   | -0.01    |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 38.253 | 4.4    | 110   | -0.01    |
| 58   | Fluorene                   | 40.000 | 38.813 | 3.0    | 113   | -0.02    |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 39.459 | 1.4    | 114   | -0.02    |
| 60   | Diethylphthalate           | 40.000 | 38.552 | 3.6    | 112   | -0.02    |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 39.386 | 1.5    | 114   | -0.02    |
| 62   | 4-Nitroaniline             | 40.000 | 33.916 | 15.2   | 95    | -0.02    |
| 63   | Azobenzene                 | 40.000 | 37.762 | 5.6    | 109   | -0.02    |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 104   | -0.02    |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 39.807 | 0.5    | 101   | -0.02    |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 42.595 | -6.5   | 113   | -0.02    |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 42.845 | -7.1   | 112   | -0.02    |
| 68   | Hexachlorobenzene          | 40.000 | 42.563 | -6.4   | 112   | -0.02    |
| 69   | Atrazine                   | 40.000 | 34.134 | 14.7   | 92    | -0.02    |
| 70 C | Pentachlorophenol          | 40.000 | 40.338 | -0.8   | 100   | -0.01    |
| 71   | Phenanthrene               | 40.000 | 39.886 | 0.3    | 105   | -0.02    |
| 72   | Anthracene                 | 40.000 | 40.029 | -0.1   | 106   | -0.02    |
| 73   | Carbazole                  | 40.000 | 37.234 | 6.9    | 97    | -0.02    |
| 74   | Di-n-butylphthalate        | 40.000 | 38.383 | 4.0    | 100   | -0.02    |
| 75 C | Fluoranthene               | 40.000 | 34.718 | 13.2   | 91    | -0.02    |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 83    | -0.02    |
| 77   | Benzidine                  | 40.000 | 35.808 | 10.5   | 62    | -0.02    |
| 78   | Pyrene                     | 40.000 | 44.341 | -10.9  | 89    | -0.02    |
| 79 S | Terphenyl-d14              | 80.000 | 87.051 | -8.8   | 89    | -0.02    |
| 80   | Butylbenzylphthalate       | 40.000 | 41.738 | -4.3   | 84    | -0.02    |
| 81   | Benzo(a)anthracene         | 40.000 | 40.133 | -0.3   | 82    | -0.02    |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 43.296 | -8.2   | 90    | -0.02    |
| 83   | Chrysene                   | 40.000 | 40.808 | -2.0   | 86    | -0.03    |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 45.839 | -14.6  | 91    | -0.03    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 54.897 | -37.2# | 112   | -0.04    |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 116   | -0.04    |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 47.835 | -19.6  | 133   | -0.06    |
| 88   | Benzo(b)fluoranthene       | 40.000 | 34.508 | 13.7   | 101   | -0.04    |
| 89   | Benzo(k)fluoranthene       | 40.000 | 33.351 | 16.6   | 97    | -0.04    |
| 90 C | Benzo(a)pyrene             | 40.000 | 39.873 | 0.3    | 116   | -0.04    |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 47.430 | -18.6  | 131   | -0.06    |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 48.159 | -20.4  | 133   | -0.07    |

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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 = 1