

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF100825\  
 Data File : BF143887.D  
 Acq On : 08 Oct 2025 10:13  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Oct 08 10:41:51 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF100625.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Oct 06 15:29:47 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 101   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.583 | 0.664 | -13.9 | 111   | 0.00     |
| 3    | Pyridine                    | 1.697 | 1.869 | -10.1 | 107   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.893 | 0.976 | -9.3  | 109   | -0.03    |
| 5 S  | 2-Fluorophenol              | 1.259 | 1.370 | -8.8  | 106   | 0.00     |
| 6    | Aniline                     | 2.250 | 2.430 | -8.0  | 105   | 0.00     |
| 7 S  | Phenol-d6                   | 1.501 | 1.603 | -6.8  | 105   | -0.01    |
| 8    | 2-Chlorophenol              | 1.231 | 1.334 | -8.4  | 105   | 0.00     |
| 9    | Benzaldehyde                | 1.158 | 1.366 | -18.0 | 104   | 0.00     |
| 10 C | Phenol                      | 1.795 | 2.034 | -13.3 | 112   | -0.02    |
| 11   | bis(2-Chloroethyl)ether     | 1.386 | 1.486 | -7.2  | 104   | -0.01    |
| 12   | 1,3-Dichlorobenzene         | 1.474 | 1.532 | -3.9  | 103   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.454 | 1.548 | -6.5  | 105   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.392 | 1.475 | -6.0  | 104   | 0.00     |
| 15   | Benzyl Alcohol              | 1.125 | 1.221 | -8.5  | 106   | -0.01    |
| 16   | 2,2'-oxybis(1-Chloropropane | 3.156 | 3.374 | -6.9  | 106   | 0.00     |
| 17   | 2-Methylphenol              | 1.017 | 1.097 | -7.9  | 106   | 0.00     |
| 18   | Hexachloroethane            | 0.485 | 0.525 | -8.2  | 106   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.028 | 1.050 | -2.1  | 104   | -0.02    |
| 20   | 3+4-Methylphenols           | 1.179 | 1.286 | -9.1  | 103   | -0.02    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 101   | 0.00     |
| 22   | Acetophenone                | 0.423 | 0.443 | -4.7  | 104   | -0.01    |
| 23 S | Nitrobenzene-d5             | 0.329 | 0.359 | -9.1  | 105   | -0.01    |
| 24   | Nitrobenzene                | 0.366 | 0.402 | -9.8  | 104   | -0.01    |
| 25   | Isophorone                  | 0.694 | 0.729 | -5.0  | 105   | -0.02    |
| 26 C | 2-Nitrophenol               | 0.157 | 0.171 | -8.9  | 100   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.282 | 0.311 | -10.3 | 107   | -0.01    |
| 28   | bis(2-Chloroethoxy)methane  | 0.438 | 0.454 | -3.7  | 104   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.294 | 0.312 | -6.1  | 103   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.289 | 0.303 | -4.8  | 103   | 0.00     |
| 31   | Naphthalene                 | 0.910 | 0.956 | -5.1  | 104   | 0.00     |
| 32   | Benzoic acid                | 0.195 | 0.198 | -1.5  | 100   | -0.05    |
| 33   | 4-Chloroaniline             | 0.341 | 0.360 | -5.6  | 103   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.201 | 0.216 | -7.5  | 106   | 0.00     |
| 35   | Caprolactam                 | 0.093 | 0.102 | -9.7  | 108   | -0.04    |
| 36 C | 4-Chloro-3-methylphenol     | 0.282 | 0.302 | -7.1  | 107   | -0.01    |
| 37   | 2-Methylnaphthalene         | 0.606 | 0.626 | -3.3  | 103   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.588 | 0.604 | -2.7  | 101   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.553 | 0.573 | -3.6  | 101   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.229 | 0.244 | -6.6  | 99    | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.199 | 0.213 | -7.0  | 107   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.404 | 0.428 | -5.9  | 100   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.427 | 0.435 | -1.9  | 99    | -0.01    |
| 45 S | 2-Fluorobiphenyl            | 1.260 | 1.197 | 5.0   | 101   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.324 | 1.365 | -3.1  | 103   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.104 | 1.147 | -3.9  | 105   | 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                   | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|----------------------------|-------|-------|-------|-------|----------|
| 48   | 2-Nitroaniline             | 0.349 | 0.394 | -12.9 | 104   | 0.00     |
| 49   | Acenaphthylene             | 1.552 | 1.636 | -5.4  | 105   | 0.00     |
| 50   | Dimethylphthalate          | 1.270 | 1.326 | -4.4  | 107   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.236 | 0.266 | -12.7 | 105   | -0.01    |
| 52 C | Acenaphthene               | 1.017 | 1.047 | -2.9  | 103   | 0.00     |
| 53   | 3-Nitroaniline             | 0.288 | 0.320 | -11.1 | 108   | -0.01    |
| 54 P | 2,4-Dinitrophenol          | 0.118 | 0.115 | 2.5   | 100   | 0.00     |
| 55   | Dibenzofuran               | 1.436 | 1.486 | -3.5  | 103   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.218 | 0.245 | -12.4 | 109   | -0.01    |
| 57   | 2,4-Dinitrotoluene         | 0.320 | 0.367 | -14.7 | 104   | 0.00     |
| 58   | Fluorene                   | 1.083 | 1.100 | -1.6  | 104   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.302 | 0.328 | -8.6  | 105   | 0.00     |
| 60   | Diethylphthalate           | 1.180 | 1.260 | -6.8  | 107   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.573 | 0.587 | -2.4  | 102   | 0.00     |
| 62   | 4-Nitroaniline             | 0.277 | 0.308 | -11.2 | 109   | -0.02    |
| 63   | Azobenzene                 | 1.233 | 1.300 | -5.4  | 105   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.101 | 0.104 | -3.0  | 100   | -0.02    |
| 66 c | n-Nitrosodiphenylamine     | 0.628 | 0.645 | -2.7  | 104   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.221 | 0.221 | 0.0   | 102   | 0.00     |
| 68   | Hexachlorobenzene          | 0.256 | 0.262 | -2.3  | 104   | 0.00     |
| 69   | Atrazine                   | 0.190 | 0.196 | -3.2  | 107   | -0.01    |
| 70 C | Pentachlorophenol          | 0.147 | 0.152 | -3.4  | 103   | 0.00     |
| 71   | Phenanthrene               | 0.970 | 0.981 | -1.1  | 107   | 0.00     |
| 72   | Anthracene                 | 0.980 | 1.010 | -3.1  | 105   | 0.00     |
| 73   | Carbazole                  | 0.868 | 0.895 | -3.1  | 108   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.005 | 1.079 | -7.4  | 109   | 0.00     |
| 75 C | Fluoranthene               | 1.002 | 1.052 | -5.0  | 108   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 115   | 0.00     |
| 77   | Benzydine                  | 0.694 | 0.756 | -8.9  | 114   | 0.00     |
| 78   | Pyrene                     | 1.667 | 1.756 | -5.3  | 108   | 0.00     |
| 79 S | Terphenyl-d14              | 1.170 | 1.254 | -7.2  | 113   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.542 | 0.605 | -11.6 | 112   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.309 | 1.342 | -2.5  | 110   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.399 | 0.417 | -4.5  | 109   | 0.00     |
| 83   | Chrysene                   | 1.211 | 1.288 | -6.4  | 115   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.653 | 0.714 | -9.3  | 111   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.072 | 1.129 | -5.3  | 109   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.225 | 1.337 | -9.1  | 107   | -0.01    |
| 88   | Benzo(b)fluoranthene       | 1.149 | 1.192 | -3.7  | 105   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.063 | 1.203 | -13.2 | 121   | 0.00     |
| 90 C | Benzo(a)pyrene             | 0.996 | 1.095 | -9.9  | 111   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 0.986 | 1.080 | -9.5  | 107   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 0.986 | 1.074 | -8.9  | 107   | -0.02    |

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| Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|----------|-------|------|------|-------|----------|
|----------|-------|------|------|-------|----------|

(#) = Out of Range

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