

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051925\  
 Data File : BP024686.D  
 Acq On : 19 May 2025 12:48  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 19 14:38:00 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP051325.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 13 16:21:39 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 123   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.549 | 0.489 | 10.9   | 119   | 0.00     |
| 3    | Pyridine                    | 1.218 | 1.176 | 3.4    | 121   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.538 | 0.499 | 7.2    | 119   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.169 | 1.150 | 1.6    | 128   | 0.00     |
| 6    | Aniline                     | 1.927 | 1.866 | 3.2    | 120   | 0.00     |
| 7 S  | Phenol-d6                   | 1.492 | 1.492 | 0.0    | 125   | 0.01     |
| 8    | 2-Chlorophenol              | 1.327 | 1.332 | -0.4   | 126   | 0.00     |
| 9    | Benzaldehyde                | 0.966 | 0.927 | 4.0    | 119   | -0.01    |
| 10 C | Phenol                      | 1.540 | 1.515 | 1.6    | 124   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.266 | 1.186 | 6.3    | 120   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.527 | 1.462 | 4.3    | 125   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.536 | 1.497 | 2.5    | 125   | -0.01    |
| 14   | 1,2-Dichlorobenzene         | 1.497 | 1.436 | 4.1    | 124   | 0.00     |
| 15   | Benzyl Alcohol              | 1.134 | 1.139 | -0.4   | 122   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.516 | 1.413 | 6.8    | 119   | -0.01    |
| 17   | 2-Methylphenol              | 1.105 | 1.105 | 0.0    | 124   | 0.00     |
| 18   | Hexachloroethane            | 0.589 | 0.544 | 7.6    | 122   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.032 | 0.994 | 3.7    | 117   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.503 | 1.498 | 0.3    | 122   | 0.01     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 121   | 0.00     |
| 22   | Acetophenone                | 0.500 | 0.474 | 5.2    | 119   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.396 | 0.371 | 6.3    | 117   | 0.00     |
| 24   | Nitrobenzene                | 0.348 | 0.326 | 6.3    | 117   | 0.00     |
| 25   | Isophorone                  | 0.687 | 0.647 | 5.8    | 116   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.171 | 0.181 | -5.8   | 127   | -0.01    |
| 27   | 2,4-Dimethylphenol          | 0.301 | 0.303 | -0.7   | 123   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.410 | 0.382 | 6.8    | 115   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.292 | 0.294 | -0.7   | 122   | 0.01     |
| 30   | 1,2,4-Trichlorobenzene      | 0.329 | 0.317 | 3.6    | 125   | 0.00     |
| 31   | Naphthalene                 | 1.036 | 0.989 | 4.5    | 122   | 0.00     |
| 32   | Benzoic acid                | 0.181 | 0.183 | -1.1   | 115   | 0.02     |
| 33   | 4-Chloroaniline             | 0.418 | 0.426 | -1.9   | 123   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.213 | 0.203 | 4.7    | 124   | 0.00     |
| 35   | Caprolactam                 | 0.106 | 0.109 | -2.8   | 120   | 0.01     |
| 36 C | 4-Chloro-3-methylphenol     | 0.356 | 0.352 | 1.1    | 119   | 0.02     |
| 37   | 2-Methylnaphthalene         | 0.671 | 0.648 | 3.4    | 122   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.718 | 0.689 | 4.0    | 120   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 118   | -0.01    |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.581 | 0.567 | 2.4    | 121   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.352 | 0.524 | -48.9# | 173#  | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.339 | 0.350 | -3.2   | 128   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.376 | 0.389 | -3.5   | 123   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.419 | 0.420 | -0.2   | 122   | 0.02     |
| 45 S | 2-Fluorobiphenyl            | 1.527 | 1.471 | 3.7    | 121   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.478 | 1.423 | 3.7    | 121   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.127 | 1.099 | 2.5    | 121   | 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             | 0.334 | 0.330 | 1.2  | 116   | 0.00     |
| 49   | Acenaphthylene             | 1.886 | 1.814 | 3.8  | 118   | 0.00     |
| 50   | Dimethylphthalate          | 1.524 | 1.431 | 6.1  | 116   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.322 | 0.314 | 2.5  | 118   | -0.01    |
| 52 C | Acenaphthene               | 1.084 | 1.034 | 4.6  | 118   | 0.00     |
| 53   | 3-Nitroaniline             | 0.329 | 0.334 | -1.5 | 115   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.169 | 0.173 | -2.4 | 118   | 0.00     |
| 55   | Dibenzofuran               | 1.730 | 1.657 | 4.2  | 119   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.250 | 0.219 | 12.4 | 100   | 0.04     |
| 57   | 2,4-Dinitrotoluene         | 0.454 | 0.465 | -2.4 | 120   | 0.00     |
| 58   | Fluorene                   | 1.410 | 1.400 | 0.7  | 123   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.388 | 0.388 | 0.0  | 121   | 0.00     |
| 60   | Diethylphthalate           | 1.541 | 1.495 | 3.0  | 122   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.703 | 0.692 | 1.6  | 124   | 0.00     |
| 62   | 4-Nitroaniline             | 0.317 | 0.343 | -8.2 | 122   | 0.00     |
| 63   | Azobenzene                 | 1.318 | 1.225 | 7.1  | 115   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 122   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.129 | 0.132 | -2.3 | 125   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.614 | 0.585 | 4.7  | 120   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.236 | 0.231 | 2.1  | 125   | 0.00     |
| 68   | Hexachlorobenzene          | 0.299 | 0.295 | 1.3  | 128   | 0.00     |
| 69   | Atrazine                   | 0.225 | 0.223 | 0.9  | 125   | 0.00     |
| 70 C | Pentachlorophenol          | 0.168 | 0.164 | 2.4  | 117   | 0.00     |
| 71   | Phenanthrene               | 1.112 | 1.047 | 5.8  | 122   | 0.00     |
| 72   | Anthracene                 | 1.119 | 1.084 | 3.1  | 123   | 0.00     |
| 73   | Carbazole                  | 1.012 | 0.995 | 1.7  | 121   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.302 | 1.256 | 3.5  | 122   | 0.00     |
| 75 C | Fluoranthene               | 1.300 | 1.236 | 4.9  | 121   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 108   | 0.00     |
| 77   | Benzidine                  | 0.545 | 0.596 | -9.4 | 115   | 0.01     |
| 78   | Pyrene                     | 1.198 | 1.261 | -5.3 | 118   | 0.00     |
| 79 S | Terphenyl-d14              | 1.166 | 1.259 | -8.0 | 120   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.533 | 0.570 | -6.9 | 116   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.256 | 1.234 | 1.8  | 110   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.466 | 0.455 | 2.4  | 102   | 0.00     |
| 83   | Chrysene                   | 1.190 | 1.150 | 3.4  | 109   | 0.01     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.783 | 0.792 | -1.1 | 111   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.281 | 1.160 | 9.4  | 98    | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 84    | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.460 | 1.430 | 2.1  | 84    | 0.01     |
| 88   | Benzo(b)fluoranthene       | 1.145 | 1.167 | -1.9 | 89    | 0.01     |
| 89   | Benzo(k)fluoranthene       | 1.180 | 1.204 | -2.0 | 89    | 0.01     |
| 90 C | Benzo(a)pyrene             | 1.099 | 1.092 | 0.6  | 86    | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.188 | 1.163 | 2.1  | 84    | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 1.181 | 1.155 | 2.2  | 84    | 0.02     |

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