

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF013025\  
 Data File : BF141317.D  
 Acq On : 30 Jan 2025 17:04  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jan 30 17:29:09 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF012925.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 29 17:41:25 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  | 111   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.573 | 0.545 | 4.9  | 113   | 0.00     |
| 3    | Pyridine                    | 1.380 | 1.337 | 3.1  | 113   | -0.02    |
| 4    | n-Nitrosodimethylamine      | 0.789 | 0.773 | 2.0  | 115   | -0.03    |
| 5 S  | 2-Fluorophenol              | 1.303 | 1.218 | 6.5  | 113   | 0.00     |
| 6    | Aniline                     | 1.403 | 1.365 | 2.7  | 112   | 0.00     |
| 7 S  | Phenol-d6                   | 1.657 | 1.555 | 6.2  | 114   | -0.01    |
| 8    | 2-Chlorophenol              | 1.396 | 1.321 | 5.4  | 113   | -0.01    |
| 9    | Benzaldehyde                | 0.960 | 0.869 | 9.5  | 108   | 0.00     |
| 10 C | Phenol                      | 1.757 | 1.669 | 5.0  | 113   | -0.01    |
| 11   | bis(2-Chloroethyl)ether     | 1.347 | 1.282 | 4.8  | 116   | -0.01    |
| 12   | 1,3-Dichlorobenzene         | 1.540 | 1.459 | 5.3  | 114   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.547 | 1.470 | 5.0  | 114   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.448 | 1.347 | 7.0  | 112   | 0.00     |
| 15   | Benzyl Alcohol              | 1.133 | 1.089 | 3.9  | 112   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.342 | 2.249 | 4.0  | 115   | 0.00     |
| 17   | 2-Methylphenol              | 1.135 | 1.076 | 5.2  | 115   | 0.00     |
| 18   | Hexachloroethane            | 0.571 | 0.547 | 4.2  | 115   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.985 | 0.907 | 7.9  | 113   | -0.01    |
| 20   | 3+4-Methylphenols           | 1.393 | 1.308 | 6.1  | 113   | -0.01    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 113   | 0.00     |
| 22   | Acetophenone                | 0.475 | 0.434 | 8.6  | 112   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.379 | 0.357 | 5.8  | 113   | 0.00     |
| 24   | Nitrobenzene                | 0.393 | 0.377 | 4.1  | 114   | -0.01    |
| 25   | Isophorone                  | 0.656 | 0.621 | 5.3  | 114   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.185 | 0.179 | 3.2  | 114   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.236 | 0.226 | 4.2  | 113   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.418 | 0.393 | 6.0  | 114   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.289 | 0.273 | 5.5  | 113   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.330 | 0.307 | 7.0  | 114   | 0.00     |
| 31   | Naphthalene                 | 1.057 | 0.966 | 8.6  | 111   | 0.00     |
| 32   | Benzoic acid                | 0.217 | 0.232 | -6.9 | 118   | -0.03    |
| 33   | 4-Chloroaniline             | 0.358 | 0.318 | 11.2 | 106   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.194 | 0.180 | 7.2  | 113   | 0.00     |
| 35   | Caprolactam                 | 0.094 | 0.090 | 4.3  | 115   | -0.02    |
| 36 C | 4-Chloro-3-methylphenol     | 0.310 | 0.294 | 5.2  | 115   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.672 | 0.614 | 8.6  | 111   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.660 | 0.612 | 7.3  | 114   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 112   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.569 | 0.535 | 6.0  | 113   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.168 | 0.179 | -6.5 | 117   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.183 | 0.174 | 4.9  | 113   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.372 | 0.366 | 1.6  | 113   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.405 | 0.377 | 6.9  | 112   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.299 | 1.172 | 9.8  | 111   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.560 | 1.440 | 7.7  | 111   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.216 | 1.123 | 7.6  | 112   | 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             | 0.382 | 0.370 | 3.1   | 114   | 0.00     |
| 49   | Acenaphthylene             | 1.702 | 1.581 | 7.1   | 111   | 0.00     |
| 50   | Dimethylphthalate          | 1.351 | 1.255 | 7.1   | 114   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.314 | 0.300 | 4.5   | 114   | 0.00     |
| 52 C | Acenaphthene               | 1.216 | 1.139 | 6.3   | 113   | 0.00     |
| 53   | 3-Nitroaniline             | 0.307 | 0.290 | 5.5   | 108   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.146 | 0.155 | -6.2  | 119   | 0.00     |
| 55   | Dibenzofuran               | 1.627 | 1.501 | 7.7   | 111   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.225 | 0.225 | 0.0   | 113   | -0.01    |
| 57   | 2,4-Dinitrotoluene         | 0.406 | 0.389 | 4.2   | 113   | 0.00     |
| 58   | Fluorene                   | 1.262 | 1.151 | 8.8   | 111   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.319 | 0.307 | 3.8   | 113   | 0.00     |
| 60   | Diethylphthalate           | 1.310 | 1.231 | 6.0   | 113   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.617 | 0.569 | 7.8   | 113   | 0.00     |
| 62   | 4-Nitroaniline             | 0.302 | 0.290 | 4.0   | 113   | -0.01    |
| 63   | Azobenzene                 | 1.311 | 1.233 | 5.9   | 113   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 115   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.118 | 0.134 | -13.6 | 122   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.647 | 0.615 | 4.9   | 112   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.224 | 0.214 | 4.5   | 113   | 0.00     |
| 68   | Hexachlorobenzene          | 0.238 | 0.226 | 5.0   | 113   | 0.00     |
| 69   | Atrazine                   | 0.216 | 0.195 | 9.7   | 105   | 0.00     |
| 70 C | Pentachlorophenol          | 0.139 | 0.150 | -7.9  | 120   | 0.00     |
| 71   | Phenanthrene               | 1.075 | 1.019 | 5.2   | 114   | 0.00     |
| 72   | Anthracene                 | 1.053 | 1.004 | 4.7   | 114   | 0.00     |
| 73   | Carbazole                  | 0.934 | 0.891 | 4.6   | 113   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.132 | 1.125 | 0.6   | 119   | 0.00     |
| 75 C | Fluoranthene               | 1.062 | 1.032 | 2.8   | 117   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 120   | 0.00     |
| 77   | Benzydine                  | 0.641 | 0.693 | -8.1  | 114   | 0.00     |
| 78   | Pyrene                     | 1.920 | 1.754 | 8.6   | 120   | 0.00     |
| 79 S | Terphenyl-d14              | 1.297 | 1.159 | 10.6  | 118   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.671 | 0.665 | 0.9   | 128   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.380 | 1.213 | 12.1  | 115   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.439 | 0.414 | 5.7   | 119   | 0.00     |
| 83   | Chrysene                   | 1.265 | 1.178 | 6.9   | 125   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.851 | 0.849 | 0.2   | 130   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.321 | 1.367 | -3.5  | 134   | -0.02    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 109   | -0.02    |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.291 | 1.261 | 2.3   | 110   | -0.03    |
| 88   | Benzo(b)fluoranthene       | 1.258 | 1.253 | 0.4   | 123   | -0.02    |
| 89   | Benzo(k)fluoranthene       | 1.114 | 1.025 | 8.0   | 104   | -0.02    |
| 90 C | Benzo(a)pyrene             | 1.063 | 1.019 | 4.1   | 110   | -0.02    |
| 91   | Dibenzo(a,h)anthracene     | 1.039 | 1.014 | 2.4   | 109   | -0.04    |
| 92   | Benzo(g,h,i)perylene       | 1.080 | 1.072 | 0.7   | 113   | -0.04    |

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|----------|-------|------|------|-------|----------|

(#) = Out of Range

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