

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF021925\  
 Data File : BF141693.D  
 Acq On : 20 Feb 2025 00:05  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Feb 20 02:06:49 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 94    | -0.01    |
| 2    | 1,4-Dioxane                 | 0.513 | 0.564 | -9.9  | 103   | -0.01    |
| 3    | Pyridine                    | 1.222 | 1.368 | -11.9 | 102   | -0.02    |
| 4    | n-Nitrosodimethylamine      | 0.809 | 0.767 | 5.2   | 88    | -0.04    |
| 5 S  | 2-Fluorophenol              | 1.276 | 1.393 | -9.2  | 102   | -0.01    |
| 6    | Aniline                     | 1.513 | 1.473 | 2.6   | 90    | -0.02    |
| 7 S  | Phenol-d6                   | 1.612 | 1.669 | -3.5  | 98    | -0.02    |
| 8    | 2-Chlorophenol              | 1.378 | 1.463 | -6.2  | 100   | -0.02    |
| 9    | Benzaldehyde                | 0.857 | 0.820 | 4.3   | 94    | 0.00     |
| 10 C | Phenol                      | 1.678 | 1.743 | -3.9  | 98    | -0.02    |
| 11   | bis(2-Chloroethyl)ether     | 1.261 | 1.337 | -6.0  | 100   | -0.01    |
| 12   | 1,3-Dichlorobenzene         | 1.455 | 1.554 | -6.8  | 102   | -0.01    |
| 13 C | 1,4-Dichlorobenzene         | 1.487 | 1.584 | -6.5  | 102   | -0.01    |
| 14   | 1,2-Dichlorobenzene         | 1.380 | 1.475 | -6.9  | 101   | -0.01    |
| 15   | Benzyl Alcohol              | 1.236 | 1.233 | 0.2   | 92    | -0.01    |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.158 | 2.096 | 2.9   | 91    | -0.01    |
| 17   | 2-Methylphenol              | 1.079 | 1.118 | -3.6  | 98    | -0.01    |
| 18   | Hexachloroethane            | 0.550 | 0.588 | -6.9  | 100   | -0.01    |
| 19 P | n-Nitroso-di-n-propylamine  | 0.964 | 0.973 | -0.9  | 95    | -0.02    |
| 20   | 3+4-Methylphenols           | 1.371 | 1.421 | -3.6  | 98    | -0.02    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 95    | -0.01    |
| 22   | Acetophenone                | 0.498 | 0.505 | -1.4  | 97    | -0.02    |
| 23 S | Nitrobenzene-d5             | 0.383 | 0.384 | -0.3  | 95    | -0.02    |
| 24   | Nitrobenzene                | 0.374 | 0.380 | -1.6  | 96    | -0.01    |
| 25   | Isophorone                  | 0.611 | 0.619 | -1.3  | 96    | -0.02    |
| 26 C | 2-Nitrophenol               | 0.186 | 0.198 | -6.5  | 97    | -0.01    |
| 27   | 2,4-Dimethylphenol          | 0.217 | 0.224 | -3.2  | 95    | -0.01    |
| 28   | bis(2-Chloroethoxy)methane  | 0.392 | 0.412 | -5.1  | 99    | -0.01    |
| 29 C | 2,4-Dichlorophenol          | 0.294 | 0.307 | -4.4  | 98    | -0.01    |
| 30   | 1,2,4-Trichlorobenzene      | 0.320 | 0.332 | -3.8  | 100   | -0.01    |
| 31   | Naphthalene                 | 1.041 | 1.078 | -3.6  | 99    | -0.01    |
| 32   | Benzoic acid                | 0.226 | 0.189 | 16.4  | 78    | -0.04    |
| 33   | 4-Chloroaniline             | 0.363 | 0.361 | 0.6   | 94    | -0.02    |
| 34 C | Hexachlorobutadiene         | 0.192 | 0.203 | -5.7  | 100   | -0.01    |
| 35   | Caprolactam                 | 0.089 | 0.088 | 1.1   | 90    | -0.04    |
| 36 C | 4-Chloro-3-methylphenol     | 0.325 | 0.323 | 0.6   | 93    | -0.01    |
| 37   | 2-Methylnaphthalene         | 0.677 | 0.691 | -2.1  | 98    | -0.01    |
| 38   | 1-Methylnaphthalene         | 0.656 | 0.661 | -0.8  | 96    | -0.01    |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 91    | -0.02    |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.579 | 0.618 | -6.7  | 96    | -0.01    |
| 41 P | Hexachlorocyclopentadiene   | 0.192 | 0.142 | 26.0# | 63    | -0.02    |
| 42 S | 2,4,6-Tribromophenol        | 0.201 | 0.198 | 1.5   | 90    | -0.02    |
| 43 C | 2,4,6-Trichlorophenol       | 0.374 | 0.390 | -4.3  | 94    | -0.01    |
| 44   | 2,4,5-Trichlorophenol       | 0.405 | 0.421 | -4.0  | 93    | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.298 | 1.372 | -5.7  | 98    | -0.01    |
| 46   | 1,1'-Biphenyl               | 1.551 | 1.641 | -5.8  | 96    | -0.02    |
| 47   | 2-Chloronaphthalene         | 1.147 | 1.205 | -5.1  | 96    | -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                   | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|----------------------------|-------|-------|--------|-------|----------|
| 48   | 2-Nitroaniline             | 0.385 | 0.380 | 1.3    | 89    | -0.02    |
| 49   | Acenaphthylene             | 1.705 | 1.777 | -4.2   | 95    | -0.02    |
| 50   | Dimethylphthalate          | 1.358 | 1.429 | -5.2   | 96    | -0.02    |
| 51   | 2,6-Dinitrotoluene         | 0.297 | 0.309 | -4.0   | 93    | -0.02    |
| 52 C | Acenaphthene               | 1.147 | 1.161 | -1.2   | 92    | -0.02    |
| 53   | 3-Nitroaniline             | 0.319 | 0.308 | 3.4    | 88    | -0.02    |
| 54 P | 2,4-Dinitrophenol          | 0.176 | 0.123 | 30.1#  | 63    | -0.01    |
| 55   | Dibenzofuran               | 1.683 | 1.718 | -2.1   | 93    | -0.02    |
| 56 P | 4-Nitrophenol              | 0.237 | 0.181 | 23.6   | 66    | -0.01    |
| 57   | 2,4-Dinitrotoluene         | 0.395 | 0.398 | -0.8   | 91    | -0.02    |
| 58   | Fluorene                   | 1.301 | 1.330 | -2.2   | 94    | -0.02    |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.349 | 0.354 | -1.4   | 92    | -0.02    |
| 60   | Diethylphthalate           | 1.336 | 1.371 | -2.6   | 94    | -0.02    |
| 61   | 4-Chlorophenyl-phenylether | 0.626 | 0.661 | -5.6   | 96    | -0.02    |
| 62   | 4-Nitroaniline             | 0.324 | 0.296 | 8.6    | 80    | -0.03    |
| 63   | Azobenzene                 | 1.328 | 1.325 | 0.2    | 91    | -0.02    |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 85    | -0.02    |
| 65   | 4,6-Dinitro-2-methylphenol | 0.139 | 0.124 | 10.8   | 74    | -0.02    |
| 66 c | n-Nitrosodiphenylamine     | 0.640 | 0.701 | -9.5   | 95    | -0.02    |
| 67   | 4-Bromophenyl-phenylether  | 0.224 | 0.248 | -10.7  | 95    | -0.02    |
| 68   | Hexachlorobenzene          | 0.230 | 0.251 | -9.1   | 93    | -0.02    |
| 69   | Atrazine                   | 0.192 | 0.184 | 4.2    | 84    | -0.02    |
| 70 C | Pentachlorophenol          | 0.147 | 0.114 | 22.4#  | 62    | -0.02    |
| 71   | Phenanthrene               | 1.101 | 1.134 | -3.0   | 88    | -0.02    |
| 72   | Anthracene                 | 1.107 | 1.131 | -2.2   | 88    | -0.02    |
| 73   | Carbazole                  | 0.996 | 0.966 | 3.0    | 83    | -0.02    |
| 74   | Di-n-butylphthalate        | 1.150 | 1.206 | -4.9   | 89    | -0.02    |
| 75 C | Fluoranthene               | 1.167 | 1.058 | 9.3    | 78    | -0.02    |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 71    | -0.02    |
| 77   | Benzydine                  | 0.441 | 0.425 | 3.6    | 57    | -0.02    |
| 78   | Pyrene                     | 1.703 | 1.869 | -9.7   | 76    | -0.02    |
| 79 S | Terphenyl-d14              | 1.185 | 1.319 | -11.3  | 78    | -0.02    |
| 80   | Butylbenzylphthalate       | 0.590 | 0.690 | -16.9  | 80    | -0.02    |
| 81   | Benzo(a)anthracene         | 1.329 | 1.386 | -4.3   | 73    | -0.02    |
| 82   | 3,3'-Dichlorobenzidine     | 0.381 | 0.430 | -12.9  | 81    | -0.02    |
| 83   | Chrysene                   | 1.228 | 1.304 | -6.2   | 77    | -0.03    |
| 84   | Bis(2-ethylhexyl)phthalate | 0.665 | 0.899 | -35.2# | 92    | -0.02    |
| 85 c | Di-n-octyl phthalate       | 0.949 | 1.358 | -43.1# | 99    | -0.04    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 97    | -0.04    |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.236 | 1.449 | -17.2  | 109   | -0.06    |
| 88   | Benzo(b)fluoranthene       | 1.343 | 1.215 | 9.5    | 88    | -0.04    |
| 89   | Benzo(k)fluoranthene       | 1.147 | 1.183 | -3.1   | 101   | -0.04    |
| 90 C | Benzo(a)pyrene             | 1.087 | 1.111 | -2.2   | 100   | -0.04    |
| 91   | Dibenzo(a,h)anthracene     | 1.001 | 1.182 | -18.1  | 109   | -0.06    |
| 92   | Benzo(g,h,i)perylene       | 1.048 | 1.209 | -15.4  | 107   | -0.07    |

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

(#) = Out of Range

SPCC's out = 0 CCC's out = 2