

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\  
 Data File : BF084544.D  
 Acq On : 19 Jan 2016 9:30  
 Operator : SJ/UM  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jan 19 17:20:45 2016  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jan 19 17:05:27 2016  
 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                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 86    | -0.03    |
| 2    | 1,4-Dioxane                 | 0.586 | 0.579 | 1.2   | 81    | -0.05    |
| 3    | Pyridine                    | 1.633 | 1.397 | 14.5  | 67    | -0.08    |
| 4    | n-Nitrosodimethylamine      | 0.838 | 0.664 | 20.8  | 64    | -0.06    |
| 5 S  | 2-Fluorophenol              | 1.236 | 1.139 | 7.8   | 84    | -0.05    |
| 6    | Aniline                     | 2.272 | 2.094 | 7.8   | 77    | -0.03    |
| 7 S  | Phenol-d6                   | 1.553 | 1.481 | 4.6   | 81    | -0.03    |
| 8    | 2-Chlorophenol              | 1.403 | 1.360 | 3.1   | 84    | -0.05    |
| 9    | Benzaldehyde                | 1.011 | 0.883 | 12.7  | 75    | -0.05    |
| 10 C | Phenol                      | 1.766 | 1.561 | 11.6  | 70    | -0.05    |
| 11   | bis(2-Chloroethyl)ether     | 1.368 | 1.362 | 0.4   | 89    | -0.05    |
| 12   | 1,3-Dichlorobenzene         | 1.447 | 1.476 | -2.0  | 90    | -0.03    |
| 13 C | 1,4-Dichlorobenzene         | 1.451 | 1.433 | 1.2   | 81    | -0.03    |
| 14   | 1,2-Dichlorobenzene         | 1.390 | 1.349 | 2.9   | 89    | -0.03    |
| 15   | Benzyl Alcohol              | 1.306 | 1.091 | 16.5  | 68    | -0.03    |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.491 | 2.063 | 17.2  | 73    | -0.03    |
| 17   | 2-Methylphenol              | 1.176 | 1.184 | -0.7  | 81    | -0.03    |
| 18   | Hexachloroethane            | 0.580 | 0.589 | -1.6  | 84    | -0.03    |
| 19 P | n-Nitroso-di-n-propylamine  | 1.051 | 0.995 | 5.3   | 83    | -0.03    |
| 20   | 3+4-Methylphenols           | 1.382 | 1.234 | 10.7  | 82    | -0.03    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 84    | -0.03    |
| 22   | Acetophenone                | 0.481 | 0.487 | -1.2  | 80    | -0.03    |
| 23 S | Nitrobenzene-d5             | 0.359 | 0.323 | 10.0  | 77    | -0.03    |
| 24   | Nitrobenzene                | 0.364 | 0.401 | -10.2 | 83    | -0.03    |
| 25   | Isophorone                  | 0.687 | 0.673 | 2.0   | 82    | -0.03    |
| 26 C | 2-Nitrophenol               | 0.166 | 0.169 | -1.8  | 87    | -0.05    |
| 27   | 2,4-Dimethylphenol          | 0.336 | 0.309 | 8.0   | 72    | -0.03    |
| 28   | bis(2-Chloroethoxy)methane  | 0.420 | 0.419 | 0.2   | 90    | -0.03    |
| 29 C | 2,4-Dichlorophenol          | 0.280 | 0.292 | -4.3  | 89    | -0.03    |
| 30   | 1,2,4-Trichlorobenzene      | 0.289 | 0.305 | -5.5  | 85    | -0.03    |
| 31   | Naphthalene                 | 0.907 | 0.975 | -7.5  | 91    | -0.03    |
| 32   | Benzoic acid                | 0.198 | 0.212 | -7.1  | 86    | 0.00     |
| 33   | 4-Chloroaniline             | 0.416 | 0.394 | 5.3   | 82    | -0.05    |
| 34 C | Hexachlorobutadiene         | 0.166 | 0.161 | 3.0   | 81    | -0.03    |
| 35   | Caprolactam                 | 0.094 | 0.102 | -8.5  | 80    | -0.01    |
| 36 C | 4-Chloro-3-methylphenol     | 0.313 | 0.300 | 4.2   | 84    | -0.03    |
| 37   | 2-Methylnaphthalene         | 0.651 | 0.592 | 9.1   | 79    | -0.03    |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 78    | -0.03    |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.575 | 0.658 | -14.4 | 89    | -0.03    |
| 40 P | Hexachlorocyclopentadiene   | 0.347 | 0.336 | 3.2   | 74    | -0.03    |
| 41 S | 2,4,6-Tribromophenol        | 0.202 | 0.219 | -8.4  | 80    | -0.03    |
| 42 C | 2,4,6-Trichlorophenol       | 0.430 | 0.435 | -1.2  | 81    | -0.03    |
| 43   | 2,4,5-Trichlorophenol       | 0.412 | 0.442 | -7.3  | 81    | -0.03    |
| 44 S | 2-Fluorobiphenyl            | 1.317 | 1.159 | 12.0  | 78    | -0.03    |

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|------|----------------------------|-------|-------|---------|-------|----------|
| 45   | 1,1'-Biphenyl              | 1.590 | 1.644 | -3.4    | 87    | -0.03    |
| 46   | 2-Chloronaphthalene        | 1.249 | 1.242 | 0.6     | 86    | -0.03    |
| 47   | 2-Nitroaniline             | 0.412 | 0.501 | -21.6   | 99    | -0.03    |
| 48   | Acenaphthylene             | 1.845 | 1.790 | 3.0     | 72    | -0.03    |
| 49   | Dimethylphthalate          | 1.430 | 1.432 | -0.1    | 75    | -0.03    |
| 50   | 2,6-Dinitrotoluene         | 0.314 | 0.308 | 1.9     | 69    | -0.03    |
| 51 C | Acenaphthene               | 1.237 | 1.289 | -4.2    | 76    | -0.03    |
| 52   | 3-Nitroaniline             | 0.389 | 0.403 | -3.6    | 75    | -0.03    |
| 53 P | 2,4-Dinitrophenol          | 0.093 | 0.205 | -120.4# | 121   | -0.03    |
| 54   | Dibenzofuran               | 1.812 | 1.621 | 10.5    | 73    | -0.03    |
| 55 P | 4-Nitrophenol              | 0.282 | 0.313 | -11.0   | 72    | -0.03    |
| 56   | 2,4-Dinitrotoluene         | 0.397 | 0.412 | -3.8    | 69    | -0.03    |
| 57   | Fluorene                   | 1.400 | 1.181 | 15.6    | 70    | -0.03    |
| 58   | 2,3,4,6-Tetrachlorophenol  | 0.352 | 0.381 | -8.2    | 80    | -0.03    |
| 59   | Diethylphthalate           | 1.410 | 1.527 | -8.3    | 83    | -0.03    |
| 60   | 4-Chlorophenyl-phenylether | 0.564 | 0.655 | -16.1   | 91    | -0.03    |
| 61   | 4-Nitroaniline             | 0.346 | 0.409 | -18.2   | 94    | -0.02    |
| 62   | Azobenzene                 | 1.546 | 1.646 | -6.5    | 82    | -0.03    |
| 63 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0     | 85    | -0.03    |
| 64   | 4,6-Dinitro-2-methylphenol | 0.109 | 0.128 | -17.4   | 96    | -0.03    |
| 65 c | n-Nitrosodiphenylamine     | 0.639 | 0.597 | 6.6     | 79    | -0.03    |
| 66   | 4-Bromophenyl-phenylether  | 0.215 | 0.216 | -0.5    | 80    | -0.03    |
| 67   | Hexachlorobenzene          | 0.242 | 0.232 | 4.1     | 86    | -0.03    |
| 68   | Atrazine                   | 0.209 | 0.207 | 1.0     | 74    | -0.03    |
| 69 C | Pentachlorophenol          | 0.126 | 0.123 | 2.4     | 73    | -0.03    |
| 70   | Phenanthrene               | 1.046 | 0.961 | 8.1     | 71    | -0.03    |
| 71   | Anthracene                 | 1.026 | 1.128 | -9.9    | 97    | -0.03    |
| 72   | Carbazole                  | 1.003 | 0.992 | 1.1     | 80    | -0.05    |
| 73   | Di-n-butylphthalate        | 1.111 | 1.214 | -9.3    | 90    | -0.03    |
| 74 C | Fluoranthene               | 1.093 | 1.108 | -1.4    | 89    | -0.03    |
| 75 I | Chrysene-d12               | 1.000 | 1.000 | 0.0     | 86    | -0.03    |
| 76   | Benzidine                  | 0.717 | 0.715 | 0.3     | 85    | -0.03    |
| 77   | Pyrene                     | 1.569 | 1.450 | 7.6     | 90    | -0.03    |
| 78 S | Terphenyl-d14              | 0.896 | 0.781 | 12.8    | 87    | -0.03    |
| 79   | Butylbenzylphthalate       | 0.679 | 0.636 | 6.3     | 79    | -0.03    |
| 80   | Benzo(a)anthracene         | 1.174 | 1.080 | 8.0     | 84    | -0.03    |
| 81   | 3,3'-Dichlorobenzidine     | 0.410 | 0.457 | -11.5   | 95    | -0.03    |
| 82   | Chrysene                   | 1.128 | 1.035 | 8.2     | 80    | -0.03    |
| 83   | Bis(2-ethylhexyl)phthalate | 0.858 | 0.780 | 9.1     | 82    | -0.03    |
| 84 c | Di-n-octyl phthalate       | 1.449 | 1.273 | 12.1    | 72    | -0.03    |
| 85   | Indeno(1,2,3-cd)pyrene     | 0.895 | 0.842 | 5.9     | 83    | -0.07    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0     | 86    | -0.05    |
| 87   | Benzo(b)fluoranthene       | 1.308 | 1.280 | 2.1     | 80    | -0.05    |

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|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(k)fluoranthene   | 1.070 | 1.160 | -8.4 | 93    | -0.05    |
| 89 C | Benzo(a)pyrene         | 1.110 | 1.163 | -4.8 | 86    | -0.05    |
| 90   | Dibenzo(a,h)anthracene | 0.955 | 0.878 | 8.1  | 82    | -0.07    |
| 91   | Benzo(a,h,i)perylene   | 0.989 | 0.893 | 9.7  | 83    | -0.08    |

(#) = Out of Range

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