

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF120816\  
 Data File : BF091503.D  
 Acq On : 8 Dec 2016 14:59  
 Operator : UM/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Dec 09 00:16:15 2016  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF112916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 29 13:33:46 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 67    | -0.05    |
| 2    | 1,4-Dioxane                 | 0.554 | 0.513 | 7.4    | 64    | -0.11    |
| 3    | Pyridine                    | 1.567 | 1.472 | 6.1    | 64    | -0.12    |
| 4    | n-Nitrosodimethylamine      | 0.822 | 0.843 | -2.6   | 71    | -0.12    |
| 5 S  | 2-Fluorophenol              | 1.224 | 1.149 | 6.1    | 64    | -0.06    |
| 6    | Aniline                     | 2.002 | 2.178 | -8.8   | 75    | -0.06    |
| 7 S  | Phenol-d6                   | 1.507 | 1.694 | -12.4  | 81    | -0.04    |
| 8    | 2-Chlorophenol              | 1.342 | 1.545 | -15.1  | 79    | -0.05    |
| 9    | Benzaldehyde                | 0.969 | 0.919 | 5.2    | 64    | -0.05    |
| 10 C | Phenol                      | 1.773 | 2.157 | -21.7# | 85    | -0.05    |
| 11   | bis(2-Chloroethyl)ether     | 1.267 | 1.455 | -14.8  | 84    | -0.05    |
| 12   | 1,3-Dichlorobenzene         | 1.493 | 1.526 | -2.2   | 75    | -0.06    |
| 13 C | 1,4-Dichlorobenzene         | 1.485 | 1.498 | -0.9   | 72    | -0.06    |
| 14   | 1,2-Dichlorobenzene         | 1.383 | 1.496 | -8.2   | 76    | -0.05    |
| 15   | Benzyl Alcohol              | 1.206 | 1.160 | 3.8    | 63    | -0.06    |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.724 | 2.775 | -1.9   | 70    | -0.06    |
| 17   | 2-Methylphenol              | 1.061 | 1.036 | 2.4    | 65    | -0.05    |
| 18   | Hexachloroethane            | 0.527 | 0.567 | -7.6   | 72    | -0.06    |
| 19 P | n-Nitroso-di-n-propylamine  | 0.949 | 1.039 | -9.5   | 82    | -0.05    |
| 20   | 3+4-Methylphenols           | 1.295 | 1.321 | -2.0   | 77    | -0.05    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 67    | -0.06    |
| 22   | Acetophenone                | 0.474 | 0.520 | -9.7   | 71    | -0.05    |
| 23 S | Nitrobenzene-d5             | 0.357 | 0.368 | -3.1   | 72    | -0.06    |
| 24   | Nitrobenzene                | 0.365 | 0.440 | -20.5  | 80    | -0.05    |
| 25   | Isophorone                  | 0.632 | 0.661 | -4.6   | 69    | -0.06    |
| 26 C | 2-Nitrophenol               | 0.167 | 0.192 | -15.0  | 82    | -0.05    |
| 27   | 2,4-Dimethylphenol          | 0.311 | 0.294 | 5.5    | 60    | -0.05    |
| 28   | bis(2-Chloroethoxy)methane  | 0.381 | 0.425 | -11.5  | 82    | -0.05    |
| 29 C | 2,4-Dichlorophenol          | 0.274 | 0.320 | -16.8  | 83    | -0.05    |
| 30   | 1,2,4-Trichlorobenzene      | 0.307 | 0.339 | -10.4  | 74    | -0.06    |
| 31   | Naphthalene                 | 0.951 | 1.003 | -5.5   | 72    | -0.06    |
| 32   | Benzoic acid                | 0.230 | 0.268 | -16.5  | 73    | -0.05    |
| 33   | 4-Chloroaniline             | 0.406 | 0.448 | -10.3  | 79    | -0.05    |
| 34 C | Hexachlorobutadiene         | 0.189 | 0.236 | -24.9# | 82    | -0.06    |
| 35   | Caprolactam                 | 0.079 | 0.088 | -11.4  | 73    | -0.06    |
| 36 C | 4-Chloro-3-methylphenol     | 0.296 | 0.373 | -26.0# | 87    | -0.05    |
| 37   | 2-Methylnaphthalene         | 0.646 | 0.761 | -17.8  | 80    | -0.05    |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 75    | -0.06    |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.564 | 0.570 | -1.1   | 86    | -0.06    |
| 40 P | Hexachlorocyclopentadiene   | 0.261 | 0.272 | -4.2   | 80    | -0.06    |
| 41 S | 2,4,6-Tribromophenol        | 0.189 | 0.204 | -7.9   | 87    | -0.06    |
| 42 C | 2,4,6-Trichlorophenol       | 0.366 | 0.411 | -12.3  | 85    | -0.05    |
| 43   | 2,4,5-Trichlorophenol       | 0.367 | 0.385 | -4.9   | 78    | -0.05    |
| 44 S | 2-Fluorobiphenyl            | 1.105 | 1.244 | -12.6  | 84    | -0.05    |

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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) |
|------|----------------------------|-------|-------|--------|-------|----------|
| 45   | 1,1'-Biphenyl              | 1.438 | 1.392 | 3.2    | 84    | -0.06    |
| 46   | 2-Chloronaphthalene        | 1.071 | 1.042 | 2.7    | 83    | -0.06    |
| 47   | 2-Nitroaniline             | 0.385 | 0.390 | -1.3   | 86    | -0.06    |
| 48   | Acenaphthylene             | 1.686 | 1.637 | 2.9    | 77    | -0.06    |
| 49   | Dimethylphthalate          | 1.211 | 1.361 | -12.4  | 85    | -0.06    |
| 50   | 2,6-Dinitrotoluene         | 0.271 | 0.329 | -21.4  | 89    | -0.05    |
| 51 C | Acenaphthene               | 1.137 | 1.186 | -4.3   | 78    | -0.05    |
| 52   | 3-Nitroaniline             | 0.313 | 0.305 | 2.6    | 87    | -0.06    |
| 53 P | 2,4-Dinitrophenol          | 0.118 | 0.146 | -23.7  | 86    | -0.06    |
| 54   | Dibenzofuran               | 1.523 | 1.536 | -0.9   | 79    | -0.06    |
| 55 P | 4-Nitrophenol              | 0.226 | 0.216 | 4.4    | 75    | -0.05    |
| 56   | 2,4-Dinitrotoluene         | 0.351 | 0.438 | -24.8  | 91    | -0.05    |
| 57   | Fluorene                   | 1.169 | 1.178 | -0.8   | 82    | -0.06    |
| 58   | 2,3,4,6-Tetrachlorophenol  | 0.314 | 0.363 | -15.6  | 86    | -0.05    |
| 59   | Diethylphthalate           | 1.195 | 1.323 | -10.7  | 85    | -0.06    |
| 60   | 4-Chlorophenyl-phenylether | 0.586 | 0.622 | -6.1   | 89    | -0.06    |
| 61   | 4-Nitroaniline             | 0.294 | 0.313 | -6.5   | 78    | -0.05    |
| 62   | Azobenzene                 | 1.220 | 1.398 | -14.6  | 83    | -0.06    |
| 63 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 75    | -0.06    |
| 64   | 4,6-Dinitro-2-methylphenol | 0.110 | 0.141 | -28.2# | 98    | -0.06    |
| 65 c | n-Nitrosodiphenylamine     | 0.606 | 0.604 | 0.3    | 80    | -0.06    |
| 66   | 4-Bromophenyl-phenylether  | 0.215 | 0.245 | -14.0  | 82    | -0.06    |
| 67   | Hexachlorobenzene          | 0.232 | 0.234 | -0.9   | 83    | -0.06    |
| 68   | Atrazine                   | 0.196 | 0.197 | -0.5   | 86    | -0.06    |
| 69 C | Pentachlorophenol          | 0.137 | 0.158 | -15.3  | 80    | -0.06    |
| 70   | Phenanthrene               | 1.039 | 1.021 | 1.7    | 81    | -0.06    |
| 71   | Anthracene                 | 1.031 | 1.154 | -11.9  | 80    | -0.06    |
| 72   | Carbazole                  | 0.936 | 0.947 | -1.2   | 82    | -0.06    |
| 73   | Di-n-butylphthalate        | 1.061 | 1.160 | -9.3   | 79    | -0.06    |
| 74 C | Fluoranthene               | 0.984 | 1.193 | -21.2# | 86    | -0.06    |
| 75 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 89    | -0.07    |
| 76   | Benzidine                  | 0.587 | 0.363 | 38.2#  | 52    | -0.06    |
| 77   | Pyrene                     | 1.518 | 1.543 | -1.6   | 84    | -0.06    |
| 78 S | Terphenyl-d14              | 0.920 | 1.030 | -12.0  | 93    | -0.06    |
| 79   | Butylbenzylphthalate       | 0.568 | 0.601 | -5.8   | 95    | -0.07    |
| 80   | Benzo(a)anthracene         | 1.170 | 1.335 | -14.1  | 103   | -0.06    |
| 81   | 3,3'-Dichlorobenzidine     | 0.412 | 0.436 | -5.8   | 99    | -0.06    |
| 82   | Chrysene                   | 1.157 | 1.359 | -17.5  | 101   | -0.06    |
| 83   | Bis(2-ethylhexyl)phthalate | 0.720 | 0.854 | -18.6  | 102   | -0.06    |
| 84 c | Di-n-octyl phthalate       | 1.090 | 1.269 | -16.4  | 101   | -0.06    |
| 85   | Indeno(1,2,3-cd)pyrene     | 1.060 | 0.984 | 7.2    | 79    | -0.13    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 78    | -0.09    |
| 87   | Benzo(b)fluoranthene       | 1.243 | 1.459 | -17.4  | 83    | -0.07    |

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| 88   | Benzo(k)fluoranthene   | 1.045 | 1.130 | -8.1 | 101   | -0.07    |
| 89 C | Benzo(a)pyrene         | 1.116 | 1.203 | -7.8 | 88    | -0.09    |
| 90   | Dibenzo(a,h)anthracene | 1.033 | 1.102 | -6.7 | 81    | -0.13    |
| 91   | Benzo(g,h,i)perylene   | 1.065 | 1.118 | -5.0 | 78    | -0.13    |

(#) = Out of Range

SPCC's out = 0 CCC's out = 4