

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF081715\  
 Data File : BF081041.D  
 Acq On : 18 Aug 2015 11:06  
 Operator : UM/IZ  
 Sample : SSTDICC040  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Aug 18 13:05:20 2015  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081715.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 17 17:57:25 2015  
 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   | 97    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.627 | 0.663  | -5.7  | 100   | 0.00     |
| 3    | Pyridine                    | 1.769 | 1.811  | -2.4  | 88    | -0.02    |
| 4    | n-Nitrosodimethylamine      | 0.985 | 1.080  | -9.6  | 104   | 0.02     |
| 5 S  | 2-Fluorophenol              | 1.330 | 1.367  | -2.8  | 103   | 0.01     |
| 6    | Aniline                     | 2.523 | 2.604  | -3.2  | 102   | 0.01     |
| 7 S  | Phenol-d6                   | 1.735 | 1.888  | -8.8  | 100   | 0.02     |
| 8    | 2-Chlorophenol              | 1.455 | 1.501  | -3.2  | 92    | 0.01     |
| 9    | Benzaldehyde                | 1.118 | 1.209  | -8.1  | 98    | 0.01     |
| 10 C | Phenol                      | 2.049 | 2.263  | -10.4 | 96    | 0.02     |
| 11   | bis(2-Chloroethyl)ether     | 1.572 | 1.699  | -8.1  | 103   | 0.01     |
| 12   | 1,3-Dichlorobenzene         | 1.564 | 1.705  | -9.0  | 105   | 0.01     |
| 13 C | 1,4-Dichlorobenzene         | 1.549 | 1.736  | -12.1 | 109   | 0.01     |
| 14   | 1,2-Dichlorobenzene         | 1.449 | 1.408  | 2.8   | 87    | 0.00     |
| 15   | Benzyl Alcohol              | 1.404 | 1.545  | -10.0 | 114   | 0.01     |
| 16   | 2,2'-oxybis(1-Chloropropane | 3.089 | 3.147  | -1.9  | 95    | 0.01     |
| 17   | 2-Methylphenol              | 1.249 | 1.226  | 1.8   | 92    | 0.01     |
| 18   | Hexachloroethane            | 0.626 | 0.497  | 20.6  | 74    | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.202 | 1.265  | -5.2  | 95    | 0.01     |
| 20   | 3+4-Methylphenols           | 1.585 | 1.576  | 0.6   | 97    | 0.01     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000  | 0.0   | 101   | 0.01     |
| 22   | Acetophenone                | 0.525 | 0.519  | 1.1   | 93    | 0.01     |
| 23 S | Nitrobenzene-d5             | 0.438 | 0.425  | 3.0   | 99    | 0.01     |
| 24   | Nitrobenzene                | 0.458 | 0.491  | -7.2  | 102   | 0.01     |
| 25   | Isophorone                  | 0.816 | 0.772  | 5.4   | 94    | 0.01     |
| 26 C | 2-Nitrophenol               | 0.190 | 0.170  | 10.5  | 94    | 0.01     |
| 27   | 2,4-Dimethylphenol          | 0.337 | 0.335  | 0.6   | 91    | 0.01     |
| 28   | bis(2-Chloroethoxy)methane  | 0.482 | 0.445  | 7.7   | 93    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.284 | 0.287  | -1.1  | 96    | 0.01     |
| 30   | 1,2,4-Trichlorobenzene      | 0.315 | 0.301  | 4.4   | 91    | 0.00     |
| 31   | Naphthalene                 | 1.115 | 1.173  | -5.2  | 104   | 0.01     |
| 32   | Benzoic acid                | 0.189 | 0.174  | 7.9   | 108   | 0.06     |
| 33   | 4-Chloroaniline             | 0.470 | 0.503  | -7.0  | 115   | 0.01     |
| 34 C | Hexachlorobutadiene         | 0.178 | 0.191  | -7.3  | 106   | 0.01     |
| 35   | Caprolactam                 | 0.099 | 0.095  | 4.0   | 91    | 0.05     |
| 36 C | 4-Chloro-3-methylphenol     | 0.338 | 0.342  | -1.2  | 100   | 0.02     |
| 37   | 2-Methylnaphthalene         | 0.704 | 0.625  | 11.2  | 86    | 0.00     |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000  | 0.0   | 94    | 0.01     |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.645 | 0.716  | -11.0 | 104   | 0.01     |
| 40 P | Hexachlorocyclopentadiene   | 0.334 | 0.041# | 87.7# | 11#   | 0.01     |
| 41 S | 2,4,6-Tribromophenol        | 0.173 | 0.151  | 12.7  | 85    | 0.00     |
| 42 C | 2,4,6-Trichlorophenol       | 0.456 | 0.496  | -8.8  | 102   | 0.01     |
| 43   | 2,4,5-Trichlorophenol       | 0.456 | 0.450  | 1.3   | 83    | 0.01     |
| 44 S | 2-Fluorobiphenyl            | 1.526 | 1.447  | 5.2   | 99    | 0.00     |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF081715\  
 Data File : BF081041.D  
 Acq On : 18 Aug 2015 11:06  
 Operator : UM/IZ  
 Sample : SSTDCCCC040  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCCC040

Quant Time: Aug 18 13:05:20 2015  
 Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081715.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 17 17:57:25 2015  
 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) |
|------|----------------------------|-------|--------|--------|-------|----------|
| 45   | 1,1'-Biphenyl              | 1.762 | 1.922  | -9.1   | 92    | 0.01     |
| 46   | 2-Chloronaphthalene        | 1.415 | 1.490  | -5.3   | 101   | 0.01     |
| 47   | 2-Nitroaniline             | 0.579 | 0.578  | 0.2    | 92    | 0.01     |
| 48   | Acenaphthylene             | 2.532 | 2.514  | 0.7    | 99    | 0.01     |
| 49   | Dimethylphthalate          | 1.647 | 1.750  | -6.3   | 93    | 0.01     |
| 50   | 2,6-Dinitrotoluene         | 0.354 | 0.340  | 4.0    | 83    | 0.01     |
| 51 C | Acenaphthene               | 1.516 | 1.450  | 4.4    | 92    | 0.01     |
| 52   | 3-Nitroaniline             | 0.443 | 0.412  | 7.0    | 89    | 0.01     |
| 53 P | 2,4-Dinitrophenol          | 0.167 | 0.019# | 88.6#  | 10#   | 0.01     |
| 54   | Dibenzofuran               | 2.031 | 1.939  | 4.5    | 95    | 0.01     |
| 55 P | 4-Nitrophenol              | 0.311 | 0.289  | 7.1    | 91    | 0.01     |
| 56   | 2,4-Dinitrotoluene         | 0.469 | 0.417  | 11.1   | 81    | 0.01     |
| 57   | Fluorene                   | 1.562 | 1.355  | 13.3   | 86    | 0.00     |
| 58   | 2,3,4,6-Tetrachlorophenol  | 0.353 | 0.316  | 10.5   | 76    | 0.00     |
| 59   | Diethylphthalate           | 1.743 | 1.803  | -3.4   | 94    | 0.01     |
| 60   | 4-Chlorophenyl-phenylether | 0.661 | 0.623  | 5.7    | 87    | 0.00     |
| 61   | 4-Nitroaniline             | 0.452 | 0.436  | 3.5    | 92    | 0.01     |
| 62   | Azobenzene                 | 2.009 | 1.955  | 2.7    | 85    | 0.01     |
| 63 I | Phenanthrene-d10           | 1.000 | 1.000  | 0.0    | 79    | 0.01     |
| 64   | 4,6-Dinitro-2-methylphenol | 0.119 | 0.025  | 79.0#  | 15#   | 0.01     |
| 65 c | n-Nitrosodiphenylamine     | 0.714 | 0.837  | -17.2  | 89    | 0.01     |
| 66   | 4-Bromophenyl-phenylether  | 0.196 | 0.228  | -16.3  | 94    | 0.01     |
| 67   | Hexachlorobenzene          | 0.199 | 0.211  | -6.0   | 78    | 0.01     |
| 68   | Atrazine                   | 0.199 | 0.170  | 14.6   | 68    | 0.01     |
| 69 C | Pentachlorophenol          | 0.119 | 0.127  | -6.7   | 82    | 0.01     |
| 70   | Phenanthrene               | 1.185 | 1.247  | -5.2   | 85    | 0.01     |
| 71   | Anthracene                 | 1.153 | 1.155  | -0.2   | 77    | 0.01     |
| 72   | Carbazole                  | 1.110 | 1.105  | 0.5    | 70    | 0.01     |
| 73   | Di-n-butylphthalate        | 1.344 | 1.407  | -4.7   | 79    | 0.00     |
| 74 C | Fluoranthene               | 1.279 | 1.325  | -3.6   | 79    | 0.01     |
| 75 I | Chrysene-d12               | 1.000 | 1.000  | 0.0    | 82    | 0.01     |
| 76   | Benzidine                  | 0.756 | 0.927  | -22.6  | 82    | 0.01     |
| 77   | Pyrene                     | 1.626 | 1.596  | 1.8    | 85    | 0.01     |
| 78 S | Terphenyl-d14              | 0.933 | 0.849  | 9.0    | 71    | 0.01     |
| 79   | Butylbenzylphthalate       | 0.699 | 0.740  | -5.9   | 78    | 0.00     |
| 80   | Benzo(a)anthracene         | 1.341 | 1.300  | 3.1    | 75    | 0.01     |
| 81   | 3,3'-Dichlorobenzidine     | 0.434 | 0.507  | -16.8  | 96    | 0.01     |
| 82   | Chrysene                   | 1.209 | 1.119  | 7.4    | 70    | 0.01     |
| 83   | Bis(2-ethylhexyl)phthalate | 0.895 | 1.000  | -11.7  | 87    | 0.01     |
| 84 c | Di-n-octyl phthalate       | 1.361 | 1.778  | -30.6# | 99    | 0.01     |
| 85   | Indeno(1,2,3-cd)pyrene     | 0.849 | 0.769  | 9.4    | 71    | 0.02     |
| 86 I | Perylene-d12               | 1.000 | 1.000  | 0.0    | 73    | 0.01     |
| 87   | Benzo(b)fluoranthene       | 1.415 | 1.379  | 2.5    | 65    | 0.01     |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF081715\  
Data File : BF081041.D  
Acq On : 18 Aug 2015 11:06  
Operator : UM/IZ  
Sample : SSTDICCC040  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_F  
LabSampleId :  
SSTDCCC040

Quant Time: Aug 18 13:05:20 2015  
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081715.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Aug 17 17:57:25 2015  
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) |
|------|------------------------|-------|-------|-------|-------|----------|
| 88   | Benzo(k)fluoranthene   | 1.318 | 1.503 | -14.0 | 90    | 0.01     |
| 89 C | Benzo(a)pyrene         | 1.233 | 1.313 | -6.5  | 77    | 0.02     |
| 90   | Dibenzo(a,h)anthracene | 0.960 | 0.987 | -2.8  | 73    | 0.02     |
| 91   | Benzo(g,h,i)perylene   | 0.974 | 0.952 | 2.3   | 70    | 0.03     |

(#) = Out of Range

SPCC's out = 2 CCC's out = 1