

Data Path : Z:\HPCHEM1\BNA F\Data\BF120717\  
 Data File : BF101225.D  
 Acq On : 8 Dec 2017 1:11  
 Operator : SJ/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Dec 08 02:33:34 2017  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Dec 04 13:07:49 2017  
 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   | 88    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.500 | 0.505  | -1.0  | 86    | -0.02    |
| 3    | Pyridine                    | 1.297 | 1.383  | -6.6  | 84    | -0.02    |
| 4    | n-Nitrosodimethylamine      | 0.634 | 0.704  | -11.0 | 94    | -0.02    |
| 5 S  | 2-Fluorophenol              | 1.210 | 1.259  | -4.0  | 89    | 0.00     |
| 6    | Aniline                     | 1.911 | 1.866  | 2.4   | 84    | 0.00     |
| 7 S  | Phenol-d6                   | 1.469 | 1.467  | 0.1   | 86    | 0.00     |
| 8    | 2-Chlorophenol              | 1.320 | 1.302  | 1.4   | 86    | 0.00     |
| 9    | Benzaldehyde                | 0.899 | 0.832  | 7.5   | 84    | 0.00     |
| 10 C | Phenol                      | 1.634 | 1.585  | 3.0   | 84    | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.226 | 1.187  | 3.2   | 84    | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.537 | 1.538  | -0.1  | 86    | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.591 | 1.573  | 1.1   | 86    | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.484 | 1.442  | 2.8   | 86    | 0.00     |
| 15   | Benzyl Alcohol              | 1.135 | 1.043  | 8.1   | 78    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.666 | 1.559  | 6.4   | 79    | 0.00     |
| 17   | 2-Methylphenol              | 1.066 | 1.028  | 3.6   | 84    | 0.00     |
| 18   | Hexachloroethane            | 0.511 | 0.385  | 24.7  | 65    | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.990 | 0.935  | 5.6   | 83    | 0.00     |
| 20   | 3+4-Methylphenols           | 1.390 | 1.310  | 5.8   | 81    | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000  | 0.0   | 79    | 0.00     |
| 22   | Acetophenone                | 0.483 | 0.503  | -4.1  | 81    | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.335 | 0.342  | -2.1  | 79    | 0.00     |
| 24   | Nitrobenzene                | 0.329 | 0.332  | -0.9  | 79    | 0.00     |
| 25   | Isophorone                  | 0.573 | 0.565  | 1.4   | 77    | 0.00     |
| 26 C | 2-Nitrophenol               | 0.180 | 0.123  | 31.7# | 53    | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.261 | 0.265  | -1.5  | 79    | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.385 | 0.390  | -1.3  | 79    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.284 | 0.282  | 0.7   | 76    | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.312 | 0.317  | -1.6  | 80    | 0.00     |
| 31   | Naphthalene                 | 0.986 | 0.958  | 2.8   | 76    | 0.00     |
| 32   | Benzoic acid                | 0.203 | 0.128  | 36.9# | 51    | -0.02    |
| 33   | 4-Chloroaniline             | 0.396 | 0.389  | 1.8   | 75    | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.192 | 0.198  | -3.1  | 81    | 0.00     |
| 35   | Caprolactam                 | 0.089 | 0.073  | 18.0  | 62    | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.294 | 0.265  | 9.9   | 70    | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.670 | 0.610  | 9.0   | 71    | 0.00     |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000  | 0.0   | 62    | 0.00     |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.727 | 0.838  | -15.3 | 72    | 0.00     |
| 40 P | Hexachlorocyclopentadiene   | 0.215 | 0.012# | 94.4# | 3#    | 0.00     |
| 41 S | 2,4,6-Tribromophenol        | 0.240 | 0.205  | 14.6  | 52    | 0.00     |
| 42 C | 2,4,6-Trichlorophenol       | 0.426 | 0.448  | -5.2  | 64    | 0.00     |
| 43   | 2,4,5-Trichlorophenol       | 0.455 | 0.479  | -5.3  | 64    | 0.00     |
| 44 S | 2-Fluorobiphenyl            | 1.462 | 1.705  | -16.6 | 72    | 0.00     |

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 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                   | AvqRF | CCRF   | %Dev   | Area% | Dev(min) |
|------|----------------------------|-------|--------|--------|-------|----------|
| 45   | 1,1'-Biphenyl              | 1.832 | 2.026  | -10.6  | 69    | 0.00     |
| 46   | 2-Chloronaphthalene        | 1.301 | 1.403  | -7.8   | 67    | 0.00     |
| 47   | 2-Nitroaniline             | 0.385 | 0.419  | -8.8   | 67    | 0.00     |
| 48   | Acenaphthylene             | 2.139 | 2.171  | -1.5   | 63    | 0.00     |
| 49   | Dimethylphthalate          | 1.613 | 1.747  | -8.3   | 67    | -0.01    |
| 50   | 2,6-Dinitrotoluene         | 0.340 | 0.286  | 15.9   | 52    | 0.00     |
| 51 C | Acenaphthene               | 1.281 | 1.276  | 0.4    | 63    | 0.00     |
| 52   | 3-Nitroaniline             | 0.382 | 0.401  | -5.0   | 64    | 0.00     |
| 53 P | 2,4-Dinitrophenol          | 0.155 | 0.000# | 100.0# | 0#    | -9.96#   |
| 54   | Dibenzofuran               | 1.845 | 1.814  | 1.7    | 61    | 0.00     |
| 55 P | 4-Nitrophenol              | 0.258 | 0.178  | 31.0#  | 41#   | 0.00     |
| 56   | 2,4-Dinitrotoluene         | 0.455 | 0.319  | 29.9#  | 42#   | 0.00     |
| 57   | Fluorene                   | 1.500 | 1.442  | 3.9    | 59    | 0.00     |
| 58   | 2,3,4,6-Tetrachlorophenol  | 0.365 | 0.328  | 10.1   | 55    | 0.00     |
| 59   | Diethylphthalate           | 1.591 | 1.645  | -3.4   | 63    | 0.00     |
| 60   | 4-Chlorophenyl-phenylether | 0.741 | 0.755  | -1.9   | 63    | 0.00     |
| 61   | 4-Nitroaniline             | 0.397 | 0.401  | -1.0   | 62    | 0.00     |
| 62   | Azobenzene                 | 1.350 | 1.235  | 8.5    | 56    | 0.00     |
| 63 I | Phenanthrene-d10           | 1.000 | 1.000  | 0.0    | 52    | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol | 0.134 | 0.008  | 94.0#  | 3#    | 0.00     |
| 65 c | n-Nitrosodiphenylamine     | 0.706 | 0.793  | -12.3  | 58    | -0.01    |
| 66   | 4-Bromophenyl-phenylether  | 0.224 | 0.245  | -9.4   | 56    | 0.00     |
| 67   | Hexachlorobenzene          | 0.240 | 0.244  | -1.7   | 53    | 0.00     |
| 68   | Atrazine                   | 0.216 | 0.147  | 31.9#  | 35#   | -0.01    |
| 69 C | Pentachlorophenol          | 0.132 | 0.103  | 22.0#  | 38#   | 0.00     |
| 70   | Phenanthrene               | 1.101 | 1.094  | 0.6    | 52    | 0.00     |
| 71   | Anthracene                 | 1.115 | 1.109  | 0.5    | 51    | 0.00     |
| 72   | Carbazole                  | 1.018 | 1.002  | 1.6    | 51    | 0.00     |
| 73   | Di-n-butylphthalate        | 1.277 | 1.325  | -3.8   | 53    | 0.00     |
| 74 C | Fluoranthene               | 1.221 | 1.146  | 6.1    | 49#   | 0.00     |
| 75 I | Chrysene-d12               | 1.000 | 1.000  | 0.0    | 47#   | 0.00     |
| 76   | Benzidine                  | 0.650 | 0.866  | -33.2# | 63    | 0.00     |
| 77   | Pyrene                     | 1.380 | 1.466  | -6.2   | 50#   | 0.00     |
| 78 S | Terphenyl-d14              | 0.914 | 1.016  | -11.2  | 53    | 0.00     |
| 79   | Butylbenzylphthalate       | 0.608 | 0.651  | -7.1   | 49#   | 0.00     |
| 80   | Benzo(a)anthracene         | 1.263 | 1.272  | -0.7   | 47#   | 0.00     |
| 81   | 3,3'-Dichlorobenzidine     | 0.437 | 0.487  | -11.4  | 52    | 0.00     |
| 82   | Chrysene                   | 1.173 | 1.165  | 0.7    | 47#   | 0.00     |
| 83   | Bis(2-ethylhexyl)phthalate | 0.868 | 0.939  | -8.2   | 50#   | 0.00     |
| 84 c | Di-n-octyl phthalate       | 1.444 | 1.518  | -5.1   | 48#   | 0.00     |
| 85   | Indeno(1,2,3-cd)pyrene     | 1.043 | 0.894  | 14.3   | 41#   | 0.02     |
| 86 I | Perylene-d12               | 1.000 | 1.000  | 0.0    | 40#   | 0.00     |
| 87   | Benzo(b)fluoranthene       | 1.198 | 1.245  | -3.9   | 42#   | 0.00     |

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|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(k)fluoranthene   | 1.180 | 1.256 | -6.4 | 41#   | 0.00     |
| 89 C | Benzo(a)pyrene         | 1.089 | 1.106 | -1.6 | 40#   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 0.960 | 0.984 | -2.5 | 41#   | 0.01     |
| 91   | Benzo(a,h,i)perylene   | 0.905 | 0.915 | -1.1 | 41#   | 0.01     |

(#) = Out of Range

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