

Data Path : Z:\HPCHEM1\BNA F\Data\BF051316\  
 Data File : BF086976.D  
 Acq On : 13 May 2016 17:39  
 Operator : UM/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 13 23:01:49 2016  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri May 13 14:55:23 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   | 99    | -0.08    |
| 2    | 1,4-Dioxane                 | 0.580 | 0.583 | -0.5  | 99    | -0.15    |
| 3    | Pyridine                    | 1.417 | 1.657 | -16.9 | 115   | -0.19    |
| 4    | n-Nitrosodimethylamine      | 1.028 | 1.111 | -8.1  | 102   | -0.18    |
| 5 S  | 2-Fluorophenol              | 1.181 | 1.264 | -7.0  | 104   | -0.09    |
| 6    | Aniline                     | 1.909 | 1.936 | -1.4  | 101   | -0.08    |
| 7 S  | Phenol-d6                   | 1.578 | 1.505 | 4.6   | 94    | -0.08    |
| 8    | 2-Chlorophenol              | 1.425 | 1.401 | 1.7   | 97    | -0.08    |
| 9    | Benzaldehyde                | 0.914 | 0.857 | 6.2   | 96    | -0.08    |
| 10 C | Phenol                      | 1.876 | 1.821 | 2.9   | 93    | -0.08    |
| 11   | bis(2-Chloroethyl)ether     | 1.463 | 1.535 | -4.9  | 97    | -0.08    |
| 12   | 1,3-Dichlorobenzene         | 1.542 | 1.470 | 4.7   | 95    | -0.09    |
| 13 C | 1,4-Dichlorobenzene         | 1.573 | 1.472 | 6.4   | 98    | -0.08    |
| 14   | 1,2-Dichlorobenzene         | 1.371 | 1.411 | -2.9  | 98    | -0.08    |
| 15   | Benzyl Alcohol              | 1.090 | 1.111 | -1.9  | 99    | -0.08    |
| 16   | 2,2'-oxybis(1-Chloropropane | 3.180 | 3.085 | 3.0   | 101   | -0.08    |
| 17   | 2-Methylphenol              | 1.134 | 1.087 | 4.1   | 92    | -0.08    |
| 18   | Hexachloroethane            | 0.592 | 0.590 | 0.3   | 100   | -0.09    |
| 19 P | n-Nitroso-di-n-propylamine  | 1.103 | 1.081 | 2.0   | 90    | -0.08    |
| 20   | 3+4-Methylphenols           | 1.481 | 1.470 | 0.7   | 94    | -0.08    |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 100   | -0.08    |
| 22   | Acetophenone                | 0.472 | 0.461 | 2.3   | 99    | -0.08    |
| 23 S | Nitrobenzene-d5             | 0.398 | 0.414 | -4.0  | 104   | -0.07    |
| 24   | Nitrobenzene                | 0.410 | 0.383 | 6.6   | 93    | -0.08    |
| 25   | Isophorone                  | 0.739 | 0.721 | 2.4   | 103   | -0.07    |
| 26 C | 2-Nitrophenol               | 0.180 | 0.195 | -8.3  | 103   | -0.08    |
| 27   | 2,4-Dimethylphenol          | 0.307 | 0.317 | -3.3  | 112   | -0.07    |
| 28   | bis(2-Chloroethoxy)methane  | 0.399 | 0.367 | 8.0   | 89    | -0.08    |
| 29 C | 2,4-Dichlorophenol          | 0.270 | 0.278 | -3.0  | 104   | -0.08    |
| 30   | 1,2,4-Trichlorobenzene      | 0.302 | 0.298 | 1.3   | 99    | -0.08    |
| 31   | Naphthalene                 | 1.002 | 1.008 | -0.6  | 102   | -0.08    |
| 32   | Benzoic acid                | 0.180 | 0.172 | 4.4   | 97    | -0.07    |
| 33   | 4-Chloroaniline             | 0.389 | 0.370 | 4.9   | 94    | -0.08    |
| 34 C | Hexachlorobutadiene         | 0.175 | 0.162 | 7.4   | 92    | -0.08    |
| 35   | Caprolactam                 | 0.092 | 0.099 | -7.6  | 114   | -0.07    |
| 36 C | 4-Chloro-3-methylphenol     | 0.327 | 0.305 | 6.7   | 94    | -0.08    |
| 37   | 2-Methylnaphthalene         | 0.586 | 0.571 | 2.6   | 95    | -0.08    |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 101   | -0.08    |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.582 | 0.598 | -2.7  | 107   | -0.08    |
| 40 P | Hexachlorocyclopentadiene   | 0.276 | 0.223 | 19.2  | 80    | -0.08    |
| 41 S | 2,4,6-Tribromophenol        | 0.212 | 0.215 | -1.4  | 106   | -0.08    |
| 42 C | 2,4,6-Trichlorophenol       | 0.389 | 0.387 | 0.5   | 105   | -0.08    |
| 43   | 2,4,5-Trichlorophenol       | 0.402 | 0.380 | 5.5   | 96    | -0.07    |
| 44 S | 2-Fluorobiphenyl            | 1.190 | 1.158 | 2.7   | 111   | -0.08    |

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

|      | Compound                   | AvqRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|----------------------------|-------|-------|--------|-------|----------|
| 45   | 1,1'-Biphenyl              | 1.591 | 1.588 | 0.2    | 98    | -0.08    |
| 46   | 2-Chloronaphthalene        | 1.246 | 1.281 | -2.8   | 101   | -0.08    |
| 47   | 2-Nitroaniline             | 0.456 | 0.463 | -1.5   | 109   | -0.08    |
| 48   | Acenaphthylene             | 1.833 | 1.955 | -6.7   | 104   | -0.08    |
| 49   | Dimethylphthalate          | 1.526 | 1.519 | 0.5    | 108   | -0.08    |
| 50   | 2,6-Dinitrotoluene         | 0.321 | 0.361 | -12.5  | 121   | -0.07    |
| 51 C | Acenaphthene               | 1.140 | 1.227 | -7.6   | 102   | -0.08    |
| 52   | 3-Nitroaniline             | 0.338 | 0.415 | -22.8  | 128   | -0.07    |
| 53 P | 2,4-Dinitrophenol          | 0.134 | 0.163 | -21.6  | 121   | -0.08    |
| 54   | Dibenzofuran               | 1.464 | 1.562 | -6.7   | 102   | -0.08    |
| 55 P | 4-Nitrophenol              | 0.201 | 0.248 | -23.4  | 118   | -0.08    |
| 56   | 2,4-Dinitrotoluene         | 0.397 | 0.384 | 3.3    | 97    | -0.08    |
| 57   | Fluorene                   | 1.248 | 1.357 | -8.7   | 106   | -0.08    |
| 58   | 2,3,4,6-Tetrachlorophenol  | 0.303 | 0.319 | -5.3   | 101   | -0.08    |
| 59   | Diethylphthalate           | 1.487 | 1.349 | 9.3    | 92    | -0.08    |
| 60   | 4-Chlorophenyl-phenylether | 0.635 | 0.535 | 15.7   | 99    | -0.08    |
| 61   | 4-Nitroaniline             | 0.325 | 0.411 | -26.5# | 120   | -0.07    |
| 62   | Azobenzene                 | 1.604 | 1.509 | 5.9    | 98    | -0.08    |
| 63 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 108   | -0.08    |
| 64   | 4,6-Dinitro-2-methylphenol | 0.124 | 0.133 | -7.3   | 119   | -0.07    |
| 65 c | n-Nitrosodiphenylamine     | 0.614 | 0.560 | 8.8    | 103   | -0.08    |
| 66   | 4-Bromophenyl-phenylether  | 0.203 | 0.210 | -3.4   | 105   | -0.08    |
| 67   | Hexachlorobenzene          | 0.237 | 0.213 | 10.1   | 106   | -0.08    |
| 68   | Atrazine                   | 0.205 | 0.200 | 2.4    | 97    | -0.08    |
| 69 C | Pentachlorophenol          | 0.110 | 0.124 | -12.7  | 112   | -0.08    |
| 70   | Phenanthrene               | 1.067 | 0.949 | 11.1   | 105   | -0.08    |
| 71   | Anthracene                 | 1.092 | 1.157 | -6.0   | 110   | -0.08    |
| 72   | Carbazole                  | 0.938 | 0.950 | -1.3   | 112   | -0.08    |
| 73   | Di-n-butylphthalate        | 1.146 | 1.053 | 8.1    | 98    | -0.08    |
| 74 C | Fluoranthene               | 1.098 | 1.117 | -1.7   | 108   | -0.09    |
| 75 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 116   | -0.08    |
| 76   | Benzidine                  | 0.510 | 0.612 | -20.0  | 127   | -0.08    |
| 77   | Pyrene                     | 1.531 | 1.375 | 10.2   | 108   | -0.08    |
| 78 S | Terphenyl-d14              | 0.943 | 0.946 | -0.3   | 117   | -0.08    |
| 79   | Butylbenzylphthalate       | 0.648 | 0.650 | -0.3   | 121   | -0.08    |
| 80   | Benzo(a)anthracene         | 1.123 | 1.171 | -4.3   | 125   | -0.08    |
| 81   | 3,3'-Dichlorobenzidine     | 0.713 | 0.746 | -4.6   | 115   | -0.08    |
| 82   | Chrysene                   | 1.142 | 1.154 | -1.1   | 131   | -0.08    |
| 83   | Bis(2-ethylhexyl)phthalate | 0.779 | 0.772 | 0.9    | 116   | -0.08    |
| 84 c | Di-n-octyl phthalate       | 1.290 | 1.393 | -8.0   | 127   | -0.07    |
| 85   | Indeno(1,2,3-cd)pyrene     | 0.950 | 0.961 | -1.2   | 117   | -0.13    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 121   | -0.09    |
| 87   | Benzo(b)fluoranthene       | 1.300 | 1.461 | -12.4  | 148   | -0.08    |

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|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(k)fluoranthene   | 1.064 | 0.945 | 11.2 | 105   | -0.08    |
| 89 C | Benzo(a)pyrene         | 1.121 | 1.078 | 3.8  | 119   | -0.09    |
| 90   | Dibenzo(a,h)anthracene | 0.945 | 0.932 | 1.4  | 115   | -0.13    |
| 91   | Benzo(a,h,i)perylene   | 0.960 | 0.976 | -1.7 | 118   | -0.15    |

(#) = Out of Range

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