

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050119\  
 Data File : BF114058.D  
 Acq On : 1 May 2019 10:15  
 Operator : JU/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 01 14:23:29 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 23 03:32:35 2019  
 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   | 117   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.551 | 0.501 | 9.1   | 106   | 0.02     |
| 3    | Pyridine                    | 1.505 | 1.458 | 3.1   | 113   | 0.01     |
| 4    | n-Nitrosodimethylamine      | 0.515 | 0.501 | 2.7   | 114   | 0.02     |
| 5 S  | 2-Fluorophenol              | 1.169 | 1.128 | 3.5   | 111   | 0.00     |
| 6    | Aniline                     | 2.015 | 2.008 | 0.3   | 116   | 0.00     |
| 7 S  | Phenol-d6                   | 1.467 | 1.460 | 0.5   | 115   | 0.01     |
| 8    | 2-Chlorophenol              | 1.335 | 1.342 | -0.5  | 116   | 0.00     |
| 9    | Benzaldehyde                | 0.859 | 0.792 | 7.8   | 109   | 0.00     |
| 10 C | Phenol                      | 1.754 | 1.718 | 2.1   | 113   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.320 | 1.313 | 0.5   | 116   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.512 | 1.506 | 0.4   | 115   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.521 | 1.497 | 1.6   | 113   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.398 | 1.391 | 0.5   | 113   | 0.00     |
| 15   | Benzyl Alcohol              | 0.988 | 1.165 | -17.9 | 135   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.541 | 1.508 | 2.1   | 113   | 0.00     |
| 17   | 2-Methylphenol              | 1.170 | 1.100 | 6.0   | 108   | 0.00     |
| 18   | Hexachloroethane            | 0.533 | 0.535 | -0.4  | 116   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.950 | 0.909 | 4.3   | 112   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.322 | 1.291 | 2.3   | 110   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 116   | 0.00     |
| 22   | Acetophenone                | 0.445 | 0.434 | 2.5   | 112   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.324 | 0.341 | -5.2  | 119   | 0.00     |
| 24   | Nitrobenzene                | 0.358 | 0.368 | -2.8  | 116   | 0.00     |
| 25   | Isophorone                  | 0.663 | 0.644 | 2.9   | 114   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.163 | 0.190 | -16.6 | 132   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.266 | 0.261 | 1.9   | 111   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.422 | 0.419 | 0.7   | 114   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.304 | 0.309 | -1.6  | 116   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.339 | 0.343 | -1.2  | 116   | 0.00     |
| 31   | Naphthalene                 | 0.950 | 0.949 | 0.1   | 113   | 0.00     |
| 32   | Benzoic acid                | 0.180 | 0.198 | -10.0 | 121   | 0.03     |
| 33   | 4-Chloroaniline             | 0.402 | 0.407 | -1.2  | 118   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.211 | 0.210 | 0.5   | 114   | 0.00     |
| 35   | Caprolactam                 | 0.097 | 0.095 | 2.1   | 115   | 0.01     |
| 36 C | 4-Chloro-3-methylphenol     | 0.301 | 0.300 | 0.3   | 114   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.641 | 0.630 | 1.7   | 112   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.618 | 0.608 | 1.6   | 113   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 115   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.650 | 0.654 | -0.6  | 114   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.362 | 0.318 | 12.2  | 98    | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.244 | 0.241 | 1.2   | 110   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.413 | 0.421 | -1.9  | 116   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.463 | 0.463 | 0.0   | 114   | 0.00     |

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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                   | AvqRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|----------------------------|-------|-------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.279 | 1.192 | 6.8    | 108   | 0.00     |
| 46   | 1,1'-Biphenyl              | 1.528 | 1.507 | 1.4    | 112   | 0.00     |
| 47   | 2-Chloronaphthalene        | 1.242 | 1.223 | 1.5    | 112   | 0.00     |
| 48   | 2-Nitroaniline             | 0.326 | 0.350 | -7.4   | 120   | 0.00     |
| 49   | Acenaphthylene             | 1.945 | 1.909 | 1.9    | 111   | 0.00     |
| 50   | Dimethylphthalate          | 1.446 | 1.435 | 0.8    | 115   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.310 | 0.332 | -7.1   | 119   | 0.00     |
| 52 C | Acenaphthene               | 1.150 | 1.149 | 0.1    | 114   | 0.00     |
| 53   | 3-Nitroaniline             | 0.326 | 0.348 | -6.7   | 120   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.112 | 0.152 | -35.7# | 163#  | 0.00     |
| 55   | Dibenzofuran               | 1.722 | 1.673 | 2.8    | 109   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.258 | 0.273 | -5.8   | 119   | 0.01     |
| 57   | 2,4-Dinitrotoluene         | 0.392 | 0.438 | -11.7  | 127   | 0.00     |
| 58   | Fluorene                   | 1.296 | 1.275 | 1.6    | 111   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.407 | 0.397 | 2.5    | 111   | 0.00     |
| 60   | Diethylphthalate           | 1.374 | 1.322 | 3.8    | 109   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.687 | 0.673 | 2.0    | 110   | 0.00     |
| 62   | 4-Nitroaniline             | 0.318 | 0.354 | -11.3  | 127   | 0.00     |
| 63   | Azobenzene                 | 1.331 | 1.264 | 5.0    | 107   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 115   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.083 | 0.109 | -31.3# | 152#  | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.600 | 0.588 | 2.0    | 112   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.242 | 0.236 | 2.5    | 111   | 0.00     |
| 68   | Hexachlorobenzene          | 0.266 | 0.260 | 2.3    | 110   | 0.00     |
| 69   | Atrazine                   | 0.195 | 0.182 | 6.7    | 107   | 0.00     |
| 70 C | Pentachlorophenol          | 0.150 | 0.152 | -1.3   | 114   | 0.00     |
| 71   | Phenanthrene               | 1.005 | 0.975 | 3.0    | 110   | 0.00     |
| 72   | Anthracene                 | 1.015 | 0.996 | 1.9    | 111   | 0.00     |
| 73   | Carbazole                  | 0.906 | 0.916 | -1.1   | 114   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.065 | 1.044 | 2.0    | 110   | 0.00     |
| 75 C | Fluoranthene               | 1.097 | 1.066 | 2.8    | 113   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 120   | 0.00     |
| 77   | Benzidine                  | 0.596 | 0.576 | 3.4    | 125   | 0.00     |
| 78   | Pyrene                     | 1.399 | 1.373 | 1.9    | 115   | 0.00     |
| 79 S | Terphenyl-d14              | 0.976 | 0.925 | 5.2    | 111   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.550 | 0.576 | -4.7   | 124   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.178 | 1.306 | -10.9  | 132   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.432 | 0.509 | -17.8  | 136   | 0.00     |
| 83   | Chrysene                   | 1.148 | 1.235 | -7.6   | 128   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.700 | 0.791 | -13.0  | 136   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.096 | 1.292 | -17.9  | 141   | -0.02    |
| 86   | Indeno(1,2,3-cd)pyrene     | 1.087 | 1.149 | -5.7   | 136   | 0.00     |
| 87 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 150   | 0.00     |

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|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.265 | 1.280 | -1.2 | 152#  | -0.01    |
| 89   | Benzo(k)fluoranthene   | 1.088 | 1.028 | 5.5  | 137   | 0.00     |
| 90 C | Benzo(a)pyrene         | 1.094 | 1.078 | 1.5  | 146   | 0.00     |
| 91   | Dibenzo(a,h)anthracene | 1.027 | 0.928 | 9.6  | 136   | 0.00     |
| 92   | Benzo(q,h,i)perylene   | 1.037 | 0.905 | 12.7 | 133   | 0.00     |

(#) = Out of Range

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