

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091619\  
 Data File : BF117073.D  
 Acq On : 16 Sep 2019 21:05  
 Operator : HP/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Sep 17 11:45:20 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Sep 13 16:23:17 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    | 64    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.536 | 0.509  | 5.0    | 64    | -0.04    |
| 3    | Pyridine                    | 1.467 | 1.391  | 5.2    | 62    | -0.02    |
| 4    | n-Nitrosodimethylamine      | 0.504 | 0.472  | 6.3    | 64    | -0.04    |
| 5 S  | 2-Fluorophenol              | 1.281 | 1.248  | 2.6    | 64    | 0.00     |
| 6    | Aniline                     | 2.132 | 2.072  | 2.8    | 65    | 0.00     |
| 7 S  | Phenol-d6                   | 1.656 | 1.609  | 2.8    | 65    | 0.01     |
| 8    | 2-Chlorophenol              | 1.382 | 1.331  | 3.7    | 64    | 0.00     |
| 9    | Benzaldehyde                | 0.904 | 0.926  | -2.4   | 69    | 0.00     |
| 10 C | Phenol                      | 1.825 | 1.788  | 2.0    | 65    | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.341 | 1.248  | 6.9    | 64    | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.551 | 1.525  | 1.7    | 66    | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.583 | 1.513  | 4.4    | 63    | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.473 | 1.429  | 3.0    | 65    | 0.00     |
| 15   | Benzyl Alcohol              | 1.270 | 1.194  | 6.0    | 62    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.325 | 1.201  | 9.4    | 61    | 0.00     |
| 17   | 2-Methylphenol              | 1.158 | 1.118  | 3.5    | 66    | 0.00     |
| 18   | Hexachloroethane            | 0.609 | 0.368  | 39.6#  | 41#   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.008 | 0.954  | 5.4    | 65    | 0.00     |
| 20   | 3+4-Methylphenols           | 1.443 | 1.484  | -2.8   | 70    | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000  | 0.0    | 64    | 0.00     |
| 22   | Acetophenone                | 0.508 | 0.499  | 1.8    | 67    | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.381 | 0.369  | 3.1    | 65    | 0.00     |
| 24   | Nitrobenzene                | 0.376 | 0.361  | 4.0    | 65    | 0.00     |
| 25   | Isophorone                  | 0.674 | 0.620  | 8.0    | 64    | 0.00     |
| 26 C | 2-Nitrophenol               | 0.161 | 0.088  | 45.3#  | 36#   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.256 | 0.237  | 7.4    | 63    | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.415 | 0.386  | 7.0    | 64    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.268 | 0.242  | 9.7    | 60    | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.290 | 0.270  | 6.9    | 63    | 0.00     |
| 31   | Naphthalene                 | 0.962 | 0.905  | 5.9    | 63    | 0.00     |
| 32   | Benzoic acid                | 0.180 | 0.092  | 48.9#  | 36#   | -0.03    |
| 33   | 4-Chloroaniline             | 0.427 | 0.397  | 7.0    | 62    | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.164 | 0.154  | 6.1    | 63    | 0.00     |
| 35   | Caprolactam                 | 0.091 | 0.080  | 12.1   | 60    | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.313 | 0.280  | 10.5   | 61    | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.648 | 0.590  | 9.0    | 62    | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.607 | 0.556  | 8.4    | 62    | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000  | 0.0    | 57    | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.516 | 0.531  | -2.9   | 62    | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.194 | 0.000# | 100.0# | 0#    | -8.97#   |
| 42 S | 2,4,6-Tribromophenol        | 0.167 | 0.140  | 16.2   | 49#   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.350 | 0.326  | 6.9    | 55    | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.374 | 0.338  | 9.6    | 55    | 0.01     |

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|      | Compound                   | AvqRF | CCRF   | %Dev   | Area% | Dev(min) |
|------|----------------------------|-------|--------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.279 | 1.336  | -4.5   | 63    | 0.00     |
| 46   | 1,1'-Biphenyl              | 1.569 | 1.614  | -2.9   | 62    | 0.00     |
| 47   | 2-Chloronaphthalene        | 1.238 | 1.229  | 0.7    | 60    | 0.00     |
| 48   | 2-Nitroaniline             | 0.397 | 0.431  | -8.6   | 66    | 0.00     |
| 49   | Acenaphthylene             | 1.855 | 1.774  | 4.4    | 57    | 0.00     |
| 50   | Dimethylphthalate          | 1.416 | 1.411  | 0.4    | 60    | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.288 | 0.234  | 18.7   | 49#   | 0.00     |
| 52 C | Acenaphthene               | 1.099 | 1.041  | 5.3    | 56    | 0.00     |
| 53   | 3-Nitroaniline             | 0.364 | 0.397  | -9.1   | 65    | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.102 | 0.000# | 100.0# | 0#    | -9.91#   |
| 55   | Dibenzofuran               | 1.622 | 1.523  | 6.1    | 56    | 0.00     |
| 56 P | 4-Nitrophenol              | 0.236 | 0.154  | 34.7#  | 38#   | 0.01     |
| 57   | 2,4-Dinitrotoluene         | 0.374 | 0.272  | 27.3#  | 42#   | 0.00     |
| 58   | Fluorene                   | 1.307 | 1.176  | 10.0   | 54    | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.287 | 0.248  | 13.6   | 50    | 0.00     |
| 60   | Diethylphthalate           | 1.393 | 1.345  | 3.4    | 58    | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.565 | 0.542  | 4.1    | 56    | 0.00     |
| 62   | 4-Nitroaniline             | 0.378 | 0.393  | -4.0   | 61    | 0.00     |
| 63   | Azobenzene                 | 1.423 | 1.267  | 11.0   | 53    | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000  | 0.0    | 48#   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.084 | 0.009  | 89.3#  | 5#    | -0.28    |
| 66 c | n-Nitrosodiphenylamine     | 0.691 | 0.718  | -3.9   | 54    | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.204 | 0.206  | -1.0   | 52    | 0.00     |
| 68   | Hexachlorobenzene          | 0.223 | 0.214  | 4.0    | 50    | 0.00     |
| 69   | Atrazine                   | 0.170 | 0.162  | 4.7    | 46#   | 0.00     |
| 70 C | Pentachlorophenol          | 0.105 | 0.091  | 13.3   | 45#   | 0.00     |
| 71   | Phenanthrene               | 1.136 | 1.095  | 3.6    | 49#   | 0.00     |
| 72   | Anthracene                 | 1.160 | 1.133  | 2.3    | 50#   | 0.00     |
| 73   | Carbazole                  | 1.101 | 1.078  | 2.1    | 50#   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.310 | 1.318  | -0.6   | 51    | 0.00     |
| 75 C | Fluoranthene               | 1.167 | 1.174  | -0.6   | 51    | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000  | 0.0    | 48#   | 0.00     |
| 77   | Benzidine                  | 0.644 | 0.829  | -28.7# | 65    | 0.00     |
| 78   | Pyrene                     | 1.678 | 1.612  | 3.9    | 52    | 0.00     |
| 79 S | Terphenyl-d14              | 1.070 | 1.079  | -0.8   | 54    | 0.00     |
| 80   | Butylbenzylphthalate       | 0.793 | 0.790  | 0.4    | 53    | 0.00     |
| 81   | Benzo(a)anthracene         | 1.336 | 1.261  | 5.6    | 49#   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.431 | 0.467  | -8.4   | 56    | 0.00     |
| 83   | Chrysene                   | 1.303 | 1.234  | 5.3    | 49#   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 1.022 | 1.054  | -3.1   | 55    | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.609 | 1.746  | -8.5   | 56    | 0.00     |
| 86   | Indeno(1,2,3-cd)pyrene     | 0.951 | 0.729  | 23.3   | 44#   | 0.00     |
| 87 I | Perylene-d12               | 1.000 | 1.000  | 0.0    | 36#   | 0.00     |

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|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.211 | 1.109 | 8.4  | 37#   | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.148 | 1.236 | -7.7 | 40#   | 0.00     |
| 90 C | Benzo(a)pyrene         | 1.063 | 1.005 | 5.5  | 37#   | 0.00     |
| 91   | Dibenzo(a,h)anthracene | 0.847 | 0.866 | -2.2 | 44#   | 0.00     |
| 92   | Benzo(g,h,i)perylene   | 0.844 | 0.846 | -0.2 | 44#   | 0.00     |

(#) = Out of Range

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