

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040620\  
 Data File : BF119811.D  
 Acq On : 7 Apr 2020 10:21  
 Operator : MA/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Apr 07 11:16:37 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 06 11:15:16 2020  
 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                     | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|------------------------------|--------|--------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4       | 20.000 | 20.000 | 0.0    | 74    | 0.00     |
| 2    | 1,4-Dioxane                  | 40.000 | 35.868 | 10.3   | 64    | 0.00     |
| 3    | Pyridine                     | 40.000 | 34.912 | 12.7   | 62    | 0.00     |
| 4    | n-Nitrosodimethylamine       | 40.000 | 33.499 | 16.3   | 59    | 0.00     |
| 5 S  | 2-Fluorophenol               | 80.000 | 78.304 | 2.1    | 69    | 0.00     |
| 6    | Aniline                      | 40.000 | 36.912 | 7.7    | 64    | 0.00     |
| 7 S  | Phenol-d6                    | 80.000 | 78.448 | 1.9    | 68    | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 40.333 | -0.8   | 70    | 0.00     |
| 9    | Benzaldehyde                 | 40.000 | 34.184 | 14.5   | 61    | 0.00     |
| 10 C | Phenol                       | 40.000 | 41.145 | -2.9   | 72    | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 38.499 | 3.8    | 68    | 0.00     |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 40.337 | -0.8   | 71    | 0.00     |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 40.586 | -1.5   | 72    | 0.00     |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 40.909 | -2.3   | 71    | 0.00     |
| 15   | Benzyl Alcohol               | 40.000 | 38.088 | 4.8    | 65    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 34.266 | 14.3   | 60    | 0.00     |
| 17   | 2-Methylphenol               | 40.000 | 39.165 | 2.1    | 69    | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 40.828 | -2.1   | 72    | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 37.554 | 6.1    | 66    | 0.00     |
| 20   | 3+4-Methylphenols            | 40.000 | 39.127 | 2.2    | 68    | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0    | 72    | 0.00     |
| 22   | Acetophenone                 | 40.000 | 39.413 | 1.5    | 69    | 0.00     |
| 23 S | Nitrobenzene-d5              | 80.000 | 84.287 | -5.4   | 73    | 0.00     |
| 24   | Nitrobenzene                 | 40.000 | 39.675 | 0.8    | 69    | 0.00     |
| 25   | Isophorone                   | 40.000 | 38.116 | 4.7    | 67    | 0.00     |
| 26 C | 2-Nitrophenol                | 40.000 | 51.011 | -27.5# | 88    | 0.00     |
| 27   | 2,4-Dimethylphenol           | 40.000 | 36.826 | 7.9    | 64    | 0.00     |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 38.395 | 4.0    | 67    | 0.00     |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 40.638 | -1.6   | 71    | 0.00     |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 40.964 | -2.4   | 72    | 0.00     |
| 31   | Naphthalene                  | 40.000 | 41.021 | -2.6   | 71    | 0.00     |
| 32   | Benzoic acid                 | 40.000 | 38.044 | 4.9    | 67    | 0.00     |
| 33   | 4-Chloroaniline              | 40.000 | 38.573 | 3.6    | 66    | 0.00     |
| 34 C | Hexachlorobutadiene          | 40.000 | 43.407 | -8.5   | 76    | 0.00     |
| 35   | Caprolactam                  | 40.000 | 37.577 | 6.1    | 64    | 0.00     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 39.345 | 1.6    | 68    | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 40.602 | -1.5   | 71    | 0.00     |
| 38   | 1-Methylnaphthalene          | 40.000 | 40.217 | -0.5   | 69    | 0.00     |
| 39 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0    | 72    | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 42.646 | -6.6   | 74    | 0.00     |
| 41 P | Hexachlorocyclopentadiene    | 40.000 | 36.698 | 8.3    | 64    | 0.00     |
| 42 S | 2,4,6-Tribromophenol         | 80.000 | 91.214 | -14.0  | 79    | 0.00     |
| 43 C | 2,4,6-Trichlorophenol        | 40.000 | 41.518 | -3.8   | 73    | 0.00     |
| 44   | 2,4,5-Trichlorophenol        | 40.000 | 41.686 | -4.2   | 72    | 0.00     |

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

|      | Compound                   | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|----------------------------|--------|--------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 80.000 | 83.145 | -3.9   | 75    | 0.00     |
| 46   | 1,1'-Biphenyl              | 40.000 | 42.234 | -5.6   | 73    | 0.00     |
| 47   | 2-Chloronaphthalene        | 40.000 | 41.101 | -2.8   | 71    | 0.00     |
| 48   | 2-Nitroaniline             | 40.000 | 42.064 | -5.2   | 71    | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 40.390 | -1.0   | 69    | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 41.003 | -2.5   | 71    | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 43.617 | -9.0   | 74    | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 40.681 | -1.7   | 71    | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 41.051 | -2.6   | 70    | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 42.058 | -5.1   | 82    | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 40.526 | -1.3   | 70    | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 37.694 | 5.8    | 63    | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 45.737 | -14.3  | 78    | 0.00     |
| 58   | Fluorene                   | 40.000 | 41.549 | -3.9   | 71    | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 41.523 | -3.8   | 70    | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 41.002 | -2.5   | 71    | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 42.550 | -6.4   | 74    | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 41.351 | -3.4   | 70    | 0.00     |
| 63   | Azobenzene                 | 40.000 | 39.030 | 2.4    | 68    | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 71    | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 50.688 | -26.7# | 87    | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 39.890 | 0.3    | 69    | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 42.240 | -5.6   | 72    | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 42.181 | -5.5   | 73    | 0.00     |
| 69   | Atrazine                   | 40.000 | 40.480 | -1.2   | 70    | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 39.961 | 0.1    | 69    | 0.00     |
| 71   | Phenanthrene               | 40.000 | 40.183 | -0.5   | 69    | 0.00     |
| 72   | Anthracene                 | 40.000 | 40.170 | -0.4   | 69    | 0.00     |
| 73   | Carbazole                  | 40.000 | 39.308 | 1.7    | 67    | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 42.822 | -7.1   | 74    | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 40.044 | -0.1   | 68    | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 72    | 0.00     |
| 77   | Benzidine                  | 40.000 | 29.234 | 26.9#  | 52    | 0.00     |
| 78   | Pyrene                     | 40.000 | 38.245 | 4.4    | 69    | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 81.885 | -2.4   | 74    | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 41.533 | -3.8   | 75    | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 39.607 | 1.0    | 69    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 38.532 | 3.7    | 66    | 0.00     |
| 83   | Chrysene                   | 40.000 | 38.617 | 3.5    | 67    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 45.003 | -12.5  | 81    | 0.00     |
| 85 c | Di-n-octyl phthalate       | 40.000 | 46.049 | -15.1  | 79    | 0.00     |
| 86   | Indeno(1,2,3-cd)pyrene     | 40.000 | 41.619 | -4.0   | 77    | 0.00     |
| 87 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 73    | 0.00     |

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|      | Compound               | Amount | Calc.  | %Dev | Area% | Dev(min) |
|------|------------------------|--------|--------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 40.000 | 38.426 | 3.9  | 70    | 0.00     |
| 89   | Benzo(k)fluoranthene   | 40.000 | 40.401 | -1.0 | 71    | 0.00     |
| 90 C | Benzo(a)pyrene         | 40.000 | 39.728 | 0.7  | 71    | 0.00     |
| 91   | Dibenzo(a,h)anthracene | 40.000 | 40.465 | -1.2 | 77    | 0.00     |
| 92   | Benzo(g,h,i)perylene   | 40.000 | 39.397 | 1.5  | 77    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 1