

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082719\  
 Data File : BF116588.D  
 Acq On : 27 Aug 2019 17:05  
 Operator : HP/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Aug 28 06:36:31 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 23 13:38:31 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                     | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|------------------------------|--------|--------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4       | 20.000 | 20.000 | 0.0    | 62    | 0.00     |
| 2    | 1,4-Dioxane                  | 40.000 | 33.481 | 16.3   | 54    | -0.04    |
| 3    | Pyridine                     | 40.000 | 32.961 | 17.6   | 56    | -0.04    |
| 4    | n-Nitrosodimethylamine       | 40.000 | 36.474 | 8.8    | 57    | -0.04    |
| 5 S  | 2-Fluorophenol               | 80.000 | 69.461 | 13.2   | 57    | 0.00     |
| 6    | Aniline                      | 40.000 | 36.216 | 9.5    | 57    | 0.00     |
| 7 S  | Phenol-d6                    | 80.000 | 73.524 | 8.1    | 58    | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 36.231 | 9.4    | 60    | 0.00     |
| 9    | Benzaldehyde                 | 40.000 | 34.747 | 13.1   | 53    | 0.00     |
| 10 C | Phenol                       | 40.000 | 37.784 | 5.5    | 61    | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 35.719 | 10.7   | 57    | 0.00     |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 39.027 | 2.4    | 61    | 0.00     |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 42.002 | -5.0   | 63    | 0.00     |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 44.785 | -12.0  | 67    | 0.00     |
| 15   | Benzyl Alcohol               | 40.000 | 39.024 | 2.4    | 62    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 39.113 | 2.2    | 63    | 0.00     |
| 17   | 2-Methylphenol               | 40.000 | 44.661 | -11.7  | 69    | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 48.856 | -22.1  | 76    | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 41.886 | -4.7   | 70    | 0.00     |
| 20   | 3+4-Methylphenols            | 40.000 | 48.850 | -22.1  | 73    | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0    | 77    | 0.00     |
| 22   | Acetophenone                 | 40.000 | 36.202 | 9.5    | 73    | 0.00     |
| 23 S | Nitrobenzene-d5              | 80.000 | 76.523 | 4.3    | 76    | 0.00     |
| 24   | Nitrobenzene                 | 40.000 | 39.176 | 2.1    | 77    | 0.00     |
| 25   | Isophorone                   | 40.000 | 37.377 | 6.6    | 79    | 0.00     |
| 26 C | 2-Nitrophenol                | 40.000 | 51.313 | -28.3# | 100   | 0.00     |
| 27   | 2,4-Dimethylphenol           | 40.000 | 37.745 | 5.6    | 79    | 0.00     |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 37.816 | 5.5    | 79    | 0.00     |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 41.806 | -4.5   | 83    | 0.00     |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 42.997 | -7.5   | 82    | 0.00     |
| 31   | Naphthalene                  | 40.000 | 41.026 | -2.6   | 80    | 0.00     |
| 32   | Benzoic acid                 | 40.000 | 53.151 | -32.9# | 98    | 0.00     |
| 33   | 4-Chloroaniline              | 40.000 | 38.870 | 2.8    | 76    | 0.00     |
| 34 C | Hexachlorobutadiene          | 40.000 | 42.432 | -6.1   | 83    | 0.00     |
| 35   | Caprolactam                  | 40.000 | 34.617 | 13.5   | 71    | 0.00     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 37.382 | 6.5    | 76    | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 33.917 | 15.2   | 69    | 0.00     |
| 38   | 1-Methylnaphthalene          | 40.000 | 33.605 | 16.0   | 67    | 0.00     |
| 39 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0    | 78    | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 37.519 | 6.2    | 73    | 0.00     |
| 41 P | Hexachlorocyclopentadiene    | 40.000 | 43.117 | -7.8   | 86    | 0.00     |
| 42 S | 2,4,6-Tribromophenol         | 80.000 | 90.738 | -13.4  | 87    | 0.00     |
| 43 C | 2,4,6-Trichlorophenol        | 40.000 | 38.919 | 2.7    | 76    | 0.00     |
| 44   | 2,4,5-Trichlorophenol        | 40.000 | 33.011 | 17.5   | 65    | 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 | 76.118 | 4.9    | 75    | 0.00     |
| 46   | 1,1'-Biphenyl              | 40.000 | 35.203 | 12.0   | 70    | 0.00     |
| 47   | 2-Chloronaphthalene        | 40.000 | 31.108 | 22.2   | 62    | 0.00     |
| 48   | 2-Nitroaniline             | 40.000 | 33.124 | 17.2   | 66    | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 39.211 | 2.0    | 76    | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 37.415 | 6.5    | 75    | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 41.582 | -4.0   | 81    | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 41.895 | -4.7   | 82    | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 42.422 | -6.1   | 81    | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 57.834 | -44.6# | 135   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 40.643 | -1.6   | 80    | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 43.108 | -7.8   | 82    | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 44.322 | -10.8  | 87    | 0.00     |
| 58   | Fluorene                   | 40.000 | 33.004 | 17.5   | 66    | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 43.023 | -7.6   | 82    | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 42.386 | -6.0   | 84    | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 38.091 | 4.8    | 77    | -0.01    |
| 62   | 4-Nitroaniline             | 40.000 | 34.972 | 12.6   | 66    | 0.00     |
| 63   | Azobenzene                 | 40.000 | 38.966 | 2.6    | 76    | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 60    | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 56.766 | -41.9# | 99    | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 50.667 | -26.7# | 77    | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 52.514 | -31.3# | 81    | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 50.832 | -27.1# | 79    | 0.00     |
| 69   | Atrazine                   | 40.000 | 49.689 | -24.2  | 75    | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 42.881 | -7.2   | 65    | 0.00     |
| 71   | Phenanthrene               | 40.000 | 41.434 | -3.6   | 62    | 0.00     |
| 72   | Anthracene                 | 40.000 | 42.196 | -5.5   | 63    | 0.00     |
| 73   | Carbazole                  | 40.000 | 42.115 | -5.3   | 63    | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 52.664 | -31.7# | 82    | -0.01    |
| 75 C | Fluoranthene               | 40.000 | 56.285 | -40.7# | 84    | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 78    | 0.00     |
| 77   | Benzidine                  | 40.000 | 40.677 | -1.7   | 78    | 0.00     |
| 78   | Pyrene                     | 40.000 | 40.760 | -1.9   | 82    | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 84.721 | -5.9   | 88    | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 43.534 | -8.8   | 89    | -0.01    |
| 81   | Benzo(a)anthracene         | 40.000 | 41.288 | -3.2   | 82    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 40.720 | -1.8   | 80    | -0.01    |
| 83   | Chrysene                   | 40.000 | 40.534 | -1.3   | 81    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 42.427 | -6.1   | 90    | -0.01    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 39.802 | 0.5    | 85    | -0.03    |
| 86   | Indeno(1,2,3-cd)pyrene     | 40.000 | 33.635 | 15.9   | 81    | -0.03    |
| 87 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 51    | -0.02    |

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|------|------------------------|--------|--------|--------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 40.000 | 58.582 | -46.5# | 73    | -0.02    |
| 89   | Benzo(k)fluoranthene   | 40.000 | 55.319 | -38.3# | 67    | -0.02    |
| 90 C | Benzo(a)pyrene         | 40.000 | 43.301 | -8.3   | 56    | -0.02    |
| 91   | Dibenzo(a,h)anthracene | 40.000 | 57.493 | -43.7# | 81    | -0.02    |
| 92   | Benzo(q,h,i)perylene   | 40.000 | 57.333 | -43.3# | 83    | -0.02    |

(#) = Out of Range

SPCC's out = 0 CCC's out = 3