

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM052318\  
 Data File : BM015287.D  
 Acq On : 23 May 2018 22:01  
 Operator : SJ/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 24 02:17:25 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM052218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 22 15:47:28 2018  
 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  | 110   | 0.00     |
| 2    | 1,4-Dioxane                  | 40.000 | 39.131 | 2.2  | 105   | 0.00     |
| 3    | Pyridine                     | 40.000 | 41.329 | -3.3 | 109   | -0.01    |
| 4    | n-Nitrosodimethylamine       | 40.000 | 42.733 | -6.8 | 113   | 0.00     |
| 5 S  | 2-Fluorophenol               | 80.000 | 82.565 | -3.2 | 110   | 0.00     |
| 6    | Aniline                      | 40.000 | 41.913 | -4.8 | 113   | 0.00     |
| 7 S  | Phenol-d6                    | 80.000 | 83.883 | -4.9 | 112   | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 41.575 | -3.9 | 111   | 0.00     |
| 9    | Benzaldehyde                 | 40.000 | 41.617 | -4.0 | 115   | 0.00     |
| 10 C | Phenol                       | 40.000 | 41.393 | -3.5 | 111   | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 41.916 | -4.8 | 113   | 0.00     |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 40.602 | -1.5 | 111   | 0.00     |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 40.793 | -2.0 | 111   | 0.00     |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 41.065 | -2.7 | 112   | 0.00     |
| 15   | Benzyl Alcohol               | 40.000 | 43.452 | -8.6 | 116   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 42.633 | -6.6 | 118   | 0.00     |
| 17   | 2-Methylphenol               | 40.000 | 42.248 | -5.6 | 113   | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 42.571 | -6.4 | 114   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 43.928 | -9.8 | 118   | 0.00     |
| 20   | 3+4-Methylphenols            | 40.000 | 42.323 | -5.8 | 114   | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0  | 115   | 0.00     |
| 22   | Acetophenone                 | 40.000 | 41.081 | -2.7 | 116   | 0.00     |
| 23 S | Nitrobenzene-d5              | 80.000 | 83.283 | -4.1 | 116   | 0.00     |
| 24   | Nitrobenzene                 | 40.000 | 40.931 | -2.3 | 114   | 0.00     |
| 25   | Isophorone                   | 40.000 | 43.261 | -8.2 | 119   | 0.00     |
| 26 C | 2-Nitrophenol                | 40.000 | 41.610 | -4.0 | 117   | 0.00     |
| 27   | 2,4-Dimethylphenol           | 40.000 | 41.638 | -4.1 | 116   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 42.262 | -5.7 | 119   | 0.00     |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 41.993 | -5.0 | 115   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 40.773 | -1.9 | 115   | 0.00     |
| 31   | Naphthalene                  | 40.000 | 40.346 | -0.9 | 115   | 0.00     |
| 32   | Benzoic acid                 | 40.000 | 41.398 | -3.5 | 119   | 0.02     |
| 33   | 4-Chloroaniline              | 40.000 | 41.242 | -3.1 | 115   | 0.00     |
| 34 C | Hexachlorobutadiene          | 40.000 | 41.401 | -3.5 | 116   | 0.00     |
| 35   | Caprolactam                  | 40.000 | 40.485 | -1.2 | 112   | 0.01     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 41.852 | -4.6 | 116   | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 41.294 | -3.2 | 116   | 0.00     |
| 38 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0  | 116   | 0.00     |
| 39   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 41.316 | -3.3 | 118   | 0.00     |
| 40 P | Hexachlorocyclopentadiene    | 40.000 | 41.427 | -3.6 | 123   | 0.00     |
| 41 S | 2,4,6-Tribromophenol         | 80.000 | 84.120 | -5.2 | 115   | 0.00     |
| 42 C | 2,4,6-Trichlorophenol        | 40.000 | 42.716 | -6.8 | 117   | 0.00     |
| 43   | 2,4,5-Trichlorophenol        | 40.000 | 42.329 | -5.8 | 117   | 0.00     |
| 44 S | 2-Fluorobiphenyl             | 80.000 | 82.465 | -3.1 | 118   | 0.00     |

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

|      | Compound                   | Amount | Calc.  | %Dev  | Area% | Dev(min) |
|------|----------------------------|--------|--------|-------|-------|----------|
| 45   | 1,1'-Biphenyl              | 40.000 | 40.879 | -2.2  | 117   | 0.00     |
| 46   | 2-Chloronaphthalene        | 40.000 | 40.994 | -2.5  | 117   | 0.00     |
| 47   | 2-Nitroaniline             | 40.000 | 41.117 | -2.8  | 116   | 0.00     |
| 48   | Acenaphthylene             | 40.000 | 41.577 | -3.9  | 116   | 0.00     |
| 49   | Dimethylphthalate          | 40.000 | 40.906 | -2.3  | 116   | 0.00     |
| 50   | 2,6-Dinitrotoluene         | 40.000 | 42.058 | -5.1  | 115   | 0.00     |
| 51 C | Acenaphthene               | 40.000 | 40.431 | -1.1  | 116   | 0.00     |
| 52   | 3-Nitroaniline             | 40.000 | 40.052 | -0.1  | 112   | 0.00     |
| 53 P | 2,4-Dinitrophenol          | 40.000 | 37.916 | 5.2   | 113   | 0.00     |
| 54   | Dibenzofuran               | 40.000 | 40.087 | -0.2  | 116   | 0.00     |
| 55 P | 4-Nitrophenol              | 40.000 | 39.006 | 2.5   | 109   | 0.00     |
| 56   | 2,4-Dinitrotoluene         | 40.000 | 39.983 | 0.0   | 112   | 0.00     |
| 57   | Fluorene                   | 40.000 | 40.166 | -0.4  | 114   | 0.00     |
| 58   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 41.532 | -3.8  | 115   | 0.00     |
| 59   | Diethylphthalate           | 40.000 | 40.765 | -1.9  | 115   | 0.00     |
| 60   | 4-Chlorophenyl-phenylether | 40.000 | 40.761 | -1.9  | 117   | 0.00     |
| 61   | 4-Nitroaniline             | 40.000 | 39.171 | 2.1   | 110   | 0.00     |
| 62   | Azobenzene                 | 40.000 | 42.500 | -6.3  | 117   | 0.00     |
| 63 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0   | 111   | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol | 40.000 | 41.187 | -3.0  | 112   | 0.00     |
| 65 c | n-Nitrosodiphenylamine     | 40.000 | 42.297 | -5.7  | 114   | 0.00     |
| 66   | 4-Bromophenyl-phenylether  | 40.000 | 43.066 | -7.7  | 116   | 0.00     |
| 67   | Hexachlorobenzene          | 40.000 | 41.963 | -4.9  | 115   | 0.00     |
| 68   | Atrazine                   | 40.000 | 42.543 | -6.4  | 111   | 0.00     |
| 69 C | Pentachlorophenol          | 40.000 | 41.615 | -4.0  | 112   | 0.00     |
| 70   | Phenanthrene               | 40.000 | 40.454 | -1.1  | 111   | 0.00     |
| 71   | Anthracene                 | 40.000 | 41.325 | -3.3  | 111   | 0.00     |
| 72   | Carbazole                  | 40.000 | 40.406 | -1.0  | 107   | 0.00     |
| 73   | Di-n-butylphthalate        | 40.000 | 43.370 | -8.4  | 112   | 0.00     |
| 74 C | Fluoranthene               | 40.000 | 39.628 | 0.9   | 106   | 0.00     |
| 75 I | Chrysene-d12               | 20.000 | 20.000 | 0.0   | 101   | 0.00     |
| 76   | Benzidine                  | 40.000 | 44.300 | -10.7 | 102   | 0.00     |
| 77   | Pyrene                     | 40.000 | 42.287 | -5.7  | 105   | 0.00     |
| 78 S | Terphenyl-d14              | 80.000 | 84.986 | -6.2  | 106   | 0.00     |
| 79   | Butylbenzylphthalate       | 40.000 | 44.639 | -11.6 | 105   | 0.00     |
| 80   | Benzo(a)anthracene         | 40.000 | 40.784 | -2.0  | 101   | 0.00     |
| 81   | 3,3'-Dichlorobenzidine     | 40.000 | 41.561 | -3.9  | 99    | 0.00     |
| 82   | Chrysene                   | 40.000 | 40.480 | -1.2  | 100   | 0.00     |
| 83   | Bis(2-ethylhexyl)phthalate | 40.000 | 44.758 | -11.9 | 104   | 0.00     |
| 84 c | Di-n-octyl phthalate       | 40.000 | 42.035 | -5.1  | 102   | 0.00     |
| 85   | Indeno(1,2,3-cd)pyrene     | 40.000 | 40.302 | -0.8  | 98    | 0.00     |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0   | 98    | 0.00     |
| 87   | Benzo(b)fluoranthene       | 40.000 | 41.171 | -2.9  | 99    | 0.00     |

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|------|------------------------|--------|--------|------|-------|----------|
| 88   | Benzo(k)fluoranthene   | 40.000 | 40.727 | -1.8 | 99    | 0.00     |
| 89 C | Benzo(a)pyrene         | 40.000 | 41.404 | -3.5 | 99    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 40.000 | 41.007 | -2.5 | 99    | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 40.000 | 40.514 | -1.3 | 98    | 0.00     |

(#) = Out of Range

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