# Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051117\
Data File : BF095098.D
Acq On : 12 May 2017 00:57
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_F
**LabSampleId :** SSTDCCC040

Quant Time: May 12 04:40:54 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 2017
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 | 99 | -0.03 |
| 2 | | 1,4-Dioxane | 40.000 | 38.905 | 2.7 | 101 | -0.08 |
| 3 | | Pyridine | 40.000 | 38.661 | 3.3 | 98 | -0.07 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 35.522 | 11.2 | 91 | -0.07 |
| 5 | S | 2-Fluorophenol | 80.000 | 78.503 | 1.9 | 98 | -0.04 |
| 6 | | Aniline | 40.000 | 41.288 | -3.2 | 97 | -0.03 |
| 7 | S | Phenol-d6 | 80.000 | 83.241 | -4.1 | 103 | -0.03 |
| 8 | | 2-Chlorophenol | 40.000 | 40.032 | -0.1 | 100 | -0.02 |
| 9 | | Benzaldehyde | 40.000 | 40.382 | -1.0 | 98 | -0.02 |
| 10 | C | Phenol | 40.000 | 44.088 | -10.2 | 109 | -0.03 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 40.824 | -2.1 | 101 | -0.03 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 39.795 | 0.5 | 106 | -0.04 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 40.597 | -1.5 | 100 | -0.04 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 40.515 | -1.3 | 101 | -0.04 |
| 15 | | Benzyl Alcohol | 40.000 | 45.208 | -13.0 | 107 | -0.02 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 40.454 | -1.1 | 103 | -0.02 |
| 17 | | 2-Methylphenol | 40.000 | 44.004 | -10.0 | 113 | -0.02 |
| 18 | | Hexachloroethane | 40.000 | 39.923 | 0.2 | 98 | -0.04 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 40.557 | -1.4 | 103 | -0.02 |
| 20 | | 3+4-Methylphenols | 40.000 | 40.771 | -1.9 | 101 | -0.02 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 104 | -0.03 |
| 22 | | Acetophenone | 40.000 | 38.746 | 3.1 | 102 | -0.02 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 79.412 | 0.7 | 103 | -0.02 |
| 24 | | Nitrobenzene | 40.000 | 39.430 | 1.4 | 105 | -0.02 |
| 25 | | Isophorone | 40.000 | 38.582 | 3.5 | 101 | -0.02 |
| 26 | C | 2-Nitrophenol | 40.000 | 38.863 | 2.8 | 99 | -0.02 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 38.733 | 3.2 | 105 | -0.02 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 39.960 | 0.1 | 105 | -0.02 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 39.932 | 0.2 | 108 | -0.01 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.718 | 0.7 | 106 | -0.02 |
| 31 | | Naphthalene | 40.000 | 39.358 | 1.6 | 105 | -0.03 |
| 32 | | Benzoic acid | 40.000 | 13.281 | 66.8# | 27 | -0.05 |
| 33 | | 4-Chloroaniline | 40.000 | 41.231 | -3.1 | 108 | -0.02 |
| 34 | C | Hexachlorobutadiene | 40.000 | 38.089 | 4.8 | 102 | -0.04 |
| 35 | | Caprolactam | 40.000 | 37.698 | 5.8 | 95 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 40.906 | -2.3 | 107 | -0.01 |
| 37 | | 2-Methylnaphthalene | 40.000 | 39.839 | 0.4 | 106 | -0.03 |
| 38 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 107 | -0.03 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 40.453 | -1.1 | 103 | -0.03 |
| 40 | P | Hexachlorocyclopentadiene | 40.000 | 34.756 | 13.1 | 89 | -0.04 |
| 41 | S | 2,4,6-Tribromophenol | 80.000 | 78.025 | 2.5 | 98 | -0.03 |
| 42 | C | 2,4,6-Trichlorophenol | 40.000 | 40.309 | -0.8 | 103 | -0.02 |
| 43 | | 2,4,5-Trichlorophenol | 40.000 | 36.838 | 7.9 | 94 | 0.00 |
| 44 | S | 2-Fluorobiphenyl | 80.000 | 77.058 | 3.7 | 101 | -0.03 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 40.000 | 40.909 | -2.3 | 104 | -0.03 |
| 46 | | 2-Chloronaphthalene | 40.000 | 40.631 | -1.6 | 106 | -0.02 |
| 47 | | 2-Nitroaniline | 40.000 | 42.542 | -6.4 | 112 | -0.02 |
| 48 | | Acenaphthylene | 40.000 | 38.176 | 4.6 | 100 | -0.03 |
| 49 | | Dimethylphthalate | 40.000 | 41.183 | -3.0 | 107 | -0.03 |
| 50 | | 2,6-Dinitrotoluene | 40.000 | 39.392 | 1.5 | 101 | -0.02 |
| 51 | C | Acenaphthene | 40.000 | 39.431 | 1.4 | 103 | -0.04 |
| 52 | | 3-Nitroaniline | 40.000 | 43.271 | -8.2 | 110 | -0.01 |
| 53 | P | 2,4-Dinitrophenol | 40.000 | 7.079 | 82.3# | 3 | 0.05 |
| 54 | | Dibenzofuran | 40.000 | 41.619 | -4.0 | 110 | -0.03 |
| 55 | P | 4-Nitrophenol | 40.000 | 31.799 | 20.5 | 78 | 0.02 |
| 56 | | 2,4-Dinitrotoluene | 40.000 | 43.108 | -7.8 | 109 | -0.01 |
| 57 | | Fluorene | 40.000 | 40.190 | -0.5 | 108 | -0.03 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 38.784 | 3.0 | 102 | -0.02 |
| 59 | | Diethylphthalate | 40.000 | 40.895 | -2.2 | 110 | -0.03 |
| 60 | | 4-Chlorophenyl-phenylether | 40.000 | 41.616 | -4.0 | 108 | -0.04 |
| 61 | | 4-Nitroaniline | 40.000 | 43.483 | -8.7 | 115 | -0.01 |
| 62 | | Azobenzene | 40.000 | 39.533 | 1.2 | 101 | -0.04 |
| 63 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 106 | -0.03 |
| 64 | | 4,6-Dinitro-2-methylphenol | 40.000 | 8.014 | 80.0# | 21 | 0.00 |
| 65 | c | n-Nitrosodiphenylamine | 40.000 | 40.478 | -1.2 | 108 | -0.03 |
| 66 | | 4-Bromophenyl-phenylether | 40.000 | 39.851 | 0.4 | 103 | -0.04 |
| 67 | | Hexachlorobenzene | 40.000 | 40.001 | -0.0 | 106 | -0.03 |
| 68 | | Atrazine | 40.000 | 41.004 | -2.5 | 114 | -0.02 |
| 69 | C | Pentachlorophenol | 40.000 | 24.146 | 39.6# | 64 | -0.01 |
| 70 | | Phenanthrene | 40.000 | 39.390 | 1.5 | 107 | -0.03 |
| 71 | | Anthracene | 40.000 | 40.214 | -0.5 | 108 | -0.03 |
| 72 | | Carbazole | 40.000 | 40.114 | -0.3 | 109 | -0.02 |
| 73 | | Di-n-butylphthalate | 40.000 | 43.497 | -8.7 | 114 | -0.03 |
| 74 | C | Fluoranthene | 40.000 | 38.363 | 4.1 | 101 | -0.03 |
| 75 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 89 | -0.02 |
| 76 | | Benzidine | 40.000 | 21.341 | 46.6# | 39 | -0.01 |
| 77 | | Pyrene | 40.000 | 46.271 | -15.7 | 101 | -0.02 |
| 78 | S | Terphenyl-d14 | 80.000 | 92.139 | -15.2 | 103 | -0.03 |
| 79 | | Butylbenzylphthalate | 40.000 | 46.726 | -16.8 | 101 | -0.03 |
| 80 | | Benzo(a)anthracene | 40.000 | 39.178 | 2.1 | 87 | -0.02 |
| 81 | | 3,3'-Dichlorobenzidine | 40.000 | 33.645 | 15.9 | 72 | -0.03 |
| 82 | | Chrysene | 40.000 | 39.817 | 0.5 | 87 | -0.03 |
| 83 | | Bis(2-ethylhexyl)phthalate | 40.000 | 46.920 | -17.3 | 101 | -0.04 |
| 84 | c | Di-n-octyl phthalate | 40.000 | 43.657 | -9.1 | 94 | -0.04 |
| 85 | | Indeno(1,2,3-cd)pyrene | 40.000 | 47.748 | -19.4 | 108 | -0.02 |
| 86 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 95 | -0.02 |
| 87 | | Benzo(b)fluoranthene | 40.000 | 37.765 | 5.6 | 83 | -0.02 |

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| | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 40.000 | 40.147 | -0.4 | 92 | -0.02 |
| 89 C | Benzo(a)pyrene | 40.000 | 38.363 | 4.1 | 87 | -0.02 |
| 90 | Dibenzo(a,h)anthracene | 40.000 | 45.708 | -14.3 | 106 | -0.03 |
| 91 | Benzo(g,h,i)perylene | 40.000 | 49.975 | -24.9 | 119 | -0.02 |

(#) = Out of Range SPCC's out = 0 CCC's out = 1