# Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA G\Data\BG120817\
Data File : BG031433.D
Acq On : 8 Dec 2017 23:28
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_G
**LabSampleId :** SSTDCCC040

Quant Time: Dec 11 10:10:07 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 06 17:58:37 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 | 87 | 0.00 |
| 2 | | 1,4-Dioxane | 40.000 | 30.041 | 24.9 | 66 | 0.00 |
| 3 | | Pyridine | 40.000 | 31.993 | 20.0 | 66 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 40.550 | -1.4 | 85 | 0.00 |
| 5 | S | 2-Fluorophenol | 80.000 | 75.893 | 5.1 | 80 | 0.00 |
| 6 | | Aniline | 40.000 | 41.213 | -3.0 | 87 | 0.00 |
| 7 | S | Phenol-d6 | 80.000 | 82.996 | -3.7 | 86 | 0.00 |
| 8 | | 2-Chlorophenol | 40.000 | 40.926 | -2.3 | 86 | 0.00 |
| 9 | | Benzaldehyde | 40.000 | 34.947 | 12.6 | 74 | 0.00 |
| 10 | C | Phenol | 40.000 | 41.725 | -4.3 | 87 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 40.658 | -1.6 | 86 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 39.195 | 2.0 | 84 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.767 | 0.6 | 85 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 39.666 | 0.8 | 84 | 0.00 |
| 15 | | Benzyl Alcohol | 40.000 | 40.296 | -0.7 | 84 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 43.888 | -9.7 | 94 | 0.00 |
| 17 | | 2-Methylphenol | 40.000 | 41.418 | -3.5 | 86 | 0.00 |
| 18 | | Hexachloroethane | 40.000 | 40.779 | -1.9 | 86 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 43.985 | -10.0 | 91 | 0.00 |
| 20 | | 3+4-Methylphenols | 40.000 | 41.680 | -4.2 | 87 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 90 | 0.00 |
| 22 | | Acetophenone | 40.000 | 40.594 | -1.5 | 89 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 82.375 | -3.0 | 91 | 0.00 |
| 24 | | Nitrobenzene | 40.000 | 41.114 | -2.8 | 90 | 0.00 |
| 25 | | Isophorone | 40.000 | 41.753 | -4.4 | 92 | 0.00 |
| 26 | C | 2-Nitrophenol | 40.000 | 40.865 | -2.2 | 90 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 39.804 | 0.5 | 87 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 41.020 | -2.6 | 90 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 40.698 | -1.7 | 87 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.512 | 1.2 | 87 | 0.00 |
| 31 | | Naphthalene | 40.000 | 39.257 | 1.9 | 88 | 0.00 |
| 32 | | Benzoic acid | 40.000 | 24.624 | 38.4# | 51 | -0.01 |
| 33 | | 4-Chloroaniline | 40.000 | 42.079 | -5.2 | 91 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 38.142 | 4.6 | 86 | 0.00 |
| 35 | | Caprolactam | 40.000 | 43.945 | -9.9 | 100 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 43.211 | -8.0 | 95 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 41.028 | -2.6 | 92 | 0.00 |
| 38 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 101 | 0.00 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 37.223 | 6.9 | 93 | 0.00 |
| 40 | P | Hexachlorocyclopentadiene | 40.000 | 38.074 | 4.8 | 102 | 0.00 |
| 41 | S | 2,4,6-Tribromophenol | 80.000 | 64.098 | 19.9 | 81 | 0.00 |
| 42 | C | 2,4,6-Trichlorophenol | 40.000 | 37.612 | 6.0 | 92 | 0.00 |
| 43 | | 2,4,5-Trichlorophenol | 40.000 | 38.086 | 4.8 | 95 | 0.00 |
| 44 | S | 2-Fluorobiphenyl | 80.000 | 74.344 | 7.1 | 94 | 0.00 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 40.000 | 37.529 | 6.2 | 95 | 0.00 |
| 46 | | 2-Chloronaphthalene | 40.000 | 37.958 | 5.1 | 95 | 0.00 |
| 47 | | 2-Nitroaniline | 40.000 | 42.246 | -5.6 | 105 | 0.00 |
| 48 | | Acenaphthylene | 40.000 | 38.665 | 3.3 | 97 | 0.00 |
| 49 | | Dimethylphthalate | 40.000 | 39.978 | 0.1 | 101 | 0.00 |
| 50 | | 2,6-Dinitrotoluene | 40.000 | 41.317 | -3.3 | 101 | 0.00 |
| 51 | C | Acenaphthene | 40.000 | 39.024 | 2.4 | 98 | 0.00 |
| 52 | | 3-Nitroaniline | 40.000 | 40.816 | -2.0 | 100 | 0.00 |
| 53 | P | 2,4-Dinitrophenol | 40.000 | 16.031 | 59.9# | 30 | 0.00 |
| 54 | | Dibenzofuran | 40.000 | 39.539 | 1.2 | 98 | 0.00 |
| 55 | P | 4-Nitrophenol | 40.000 | 38.175 | 4.6 | 95 | 0.00 |
| 56 | | 2,4-Dinitrotoluene | 40.000 | 41.989 | -5.0 | 104 | 0.00 |
| 57 | | Fluorene | 40.000 | 40.123 | -0.3 | 102 | 0.00 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 38.494 | 3.8 | 97 | 0.00 |
| 59 | | Diethylphthalate | 40.000 | 40.654 | -1.6 | 104 | 0.00 |
| 60 | | 4-Chlorophenyl-phenylether | 40.000 | 40.159 | -0.4 | 101 | 0.00 |
| 61 | | 4-Nitroaniline | 40.000 | 41.627 | -4.1 | 105 | 0.00 |
| 62 | | Azobenzene | 40.000 | 43.031 | -7.6 | 109 | 0.00 |
| 63 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 108 | 0.00 |
| 64 | | 4,6-Dinitro-2-methylphenol | 40.000 | 26.967 | 32.6# | 71 | 0.00 |
| 65 | c | n-Nitrosodiphenylamine | 40.000 | 38.063 | 4.8 | 102 | 0.00 |
| 66 | | 4-Bromophenyl-phenylether | 40.000 | 36.131 | 9.7 | 95 | 0.00 |
| 67 | | Hexachlorobenzene | 40.000 | 33.541 | 16.1 | 89 | 0.00 |
| 68 | | Atrazine | 40.000 | 37.325 | 6.7 | 102 | 0.00 |
| 69 | C | Pentachlorophenol | 40.000 | 34.233 | 14.4 | 92 | 0.00 |
| 70 | | Phenanthrene | 40.000 | 39.197 | 2.0 | 105 | 0.00 |
| 71 | | Anthracene | 40.000 | 39.001 | 2.5 | 105 | 0.00 |
| 72 | | Carbazole | 40.000 | 38.872 | 2.8 | 104 | 0.00 |
| 73 | | Di-n-butylphthalate | 40.000 | 38.196 | 4.5 | 104 | 0.00 |
| 74 | C | Fluoranthene | 40.000 | 39.969 | 0.1 | 108 | 0.00 |
| 75 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 106 | 0.00 |
| 76 | | Benzidine | 40.000 | 30.755 | 23.1 | 78 | 0.00 |
| 77 | | Pyrene | 40.000 | 42.531 | -6.3 | 110 | 0.00 |
| 78 | S | Terphenyl-d14 | 80.000 | 76.666 | 4.2 | 100 | 0.00 |
| 79 | | Butylbenzylphthalate | 40.000 | 40.308 | -0.8 | 105 | 0.00 |
| 80 | | Benzo(a)anthracene | 40.000 | 39.380 | 1.5 | 103 | 0.00 |
| 81 | | 3,3'-Dichlorobenzidine | 40.000 | 36.683 | 8.3 | 95 | 0.00 |
| 82 | | Chrysene | 40.000 | 40.639 | -1.6 | 106 | 0.00 |
| 83 | | Bis(2-ethylhexyl)phthalate | 40.000 | 40.159 | -0.4 | 107 | 0.00 |
| 84 | c | Di-n-octyl phthalate | 40.000 | 41.165 | -2.9 | 109 | 0.00 |
| 85 | | Indeno(1,2,3-cd)pyrene | 40.000 | 36.757 | 8.1 | 96 | 0.00 |
| 86 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 105 | 0.00 |
| 87 | | Benzo(b)fluoranthene | 40.000 | 39.962 | 0.1 | 104 | 0.00 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 40.000 | 39.168 | 2.1 | 102 | 0.00 |
| 89 C | Benzo(a)pyrene | 40.000 | 39.231 | 1.9 | 103 | 0.00 |
| 90 | Dibenzo(a,h)anthracene | 40.000 | 36.535 | 8.7 | 95 | 0.00 |
| 91 | Benzo(g,h,i)perylene | 40.000 | 37.211 | 7.0 | 97 | 0.00 |

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