# Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071316\
Data File : BG023080.D
Acq On : 13 Jul 2016 21:51
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 3 Sample Multiplier: 1

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

Quant Time: Jul 14 04:36:48 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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 | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 1 | I | 1,4-Dichlorobenzene-d4 | 1.000 | 1.000 | 0.0 | 100 | 0.00 |
| 2 | | 1,4-Dioxane | 0.475 | 0.402 | 15.4 | 88 | 0.00 |
| 3 | | Pyridine | 1.627 | 1.469 | 9.7 | 91 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 0.698 | 0.641 | 8.2 | 91 | 0.00 |
| 5 | S | 2-Fluorophenol | 1.223 | 1.161 | 5.1 | 97 | 0.00 |
| 6 | | Aniline | 2.654 | 2.475 | 6.7 | 93 | 0.00 |
| 7 | S | Phenol-d6 | 1.915 | 1.798 | 6.1 | 92 | 0.00 |
| 8 | | 2-Chlorophenol | 1.301 | 1.217 | 6.5 | 94 | 0.00 |
| 9 | | Benzaldehyde | 1.280 | 1.163 | 9.1 | 93 | 0.00 |
| 10 | C | Phenol | 1.971 | 1.878 | 4.7 | 95 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 1.562 | 1.453 | 7.0 | 94 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 1.593 | 1.488 | 6.6 | 97 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 1.595 | 1.529 | 4.1 | 97 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 1.585 | 1.441 | 9.1 | 95 | 0.00 |
| 15 | | Benzyl Alcohol | 1.754 | 1.637 | 6.7 | 91 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 1.908 | 1.851 | 3.0 | 98 | 0.00 |
| 17 | | 2-Methylphenol | 1.283 | 1.198 | 6.6 | 93 | 0.00 |
| 18 | | Hexachloroethane | 0.673 | 0.640 | 4.9 | 103 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 1.727 | 1.564 | 9.4 | 88 | 0.00 |
| 20 | | 3+4-Methylphenols | 1.789 | 1.653 | 7.6 | 87 | 0.00 |
| 21 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 95 | 0.00 |
| 22 | | Acetophenone | 0.600 | 0.542 | 9.7 | 86 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 0.467 | 0.440 | 5.8 | 90 | 0.00 |
| 24 | | Nitrobenzene | 0.496 | 0.474 | 4.4 | 93 | 0.00 |
| 25 | | Isophorone | 0.939 | 0.840 | 10.5 | 83 | 0.00 |
| 26 | C | 2-Nitrophenol | 0.195 | 0.186 | 4.6 | 86 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 0.337 | 0.293 | 13.1 | 83 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 0.524 | 0.494 | 5.7 | 89 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 0.359 | 0.333 | 7.2 | 86 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 0.411 | 0.380 | 7.5 | 89 | 0.00 |
| 31 | | Naphthalene | 1.018 | 0.940 | 7.7 | 89 | 0.00 |
| 32 | | Benzoic acid | 0.306 | 0.273 | 10.8 | 80 | 0.02 |
| 33 | | 4-Chloroaniline | 0.498 | 0.459 | 7.8 | 88 | 0.00 |
| 34 | C | Hexachlorobutadiene | 0.334 | 0.302 | 9.6 | 88 | 0.00 |
| 35 | | Caprolactam | 0.148 | 0.123 | 16.9 | 79 | 0.01 |
| 36 | C | 4-Chloro-3-methylphenol | 0.491 | 0.443 | 9.8 | 84 | 0.01 |
| 37 | | 2-Methylnaphthalene | 0.805 | 0.743 | 7.7 | 86 | 0.00 |
| 38 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 87 | 0.00 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 0.862 | 0.787 | 8.7 | 81 | 0.00 |
| 40 | P | Hexachlorocyclopentadiene | 0.496 | 0.526 | -6.0 | 93 | 0.00 |
| 41 | S | 2,4,6-Tribromophenol | 0.359 | 0.296 | 17.5 | 74 | 0.00 |
| 42 | C | 2,4,6-Trichlorophenol | 0.495 | 0.457 | 7.7 | 80 | 0.01 |
| 43 | | 2,4,5-Trichlorophenol | 0.509 | 0.469 | 7.9 | 83 | 0.01 |
| 44 | S | 2-Fluorobiphenyl | 1.270 | 1.198 | 5.7 | 85 | 0.00 |

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| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 1.501 | 1.436 | 4.3 | 84 | 0.00 |
| 46 | | 2-Chloronaphthalene | 1.217 | 1.177 | 3.3 | 85 | 0.00 |
| 47 | | 2-Nitroaniline | 0.519 | 0.511 | 1.5 | 87 | 0.01 |
| 48 | | Acenaphthylene | 1.826 | 1.718 | 5.9 | 83 | 0.00 |
| 49 | | Dimethylphthalate | 1.634 | 1.463 | 10.5 | 81 | 0.00 |
| 50 | | 2,6-Dinitrotoluene | 0.367 | 0.334 | 9.0 | 81 | 0.01 |
| 51 | C | Acenaphthene | 1.193 | 1.130 | 5.3 | 85 | 0.00 |
| 52 | | 3-Nitroaniline | 0.366 | 0.344 | 6.0 | 83 | 0.01 |
| 53 | P | 2,4-Dinitrophenol | 0.252 | 0.273 | -8.3 | 92 | 0.00 |
| 54 | | Dibenzofuran | 1.786 | 1.645 | 7.9 | 83 | 0.00 |
| 55 | P | 4-Nitrophenol | 0.307 | 0.259 | 15.6 | 74 | 0.02 |
| 56 | | 2,4-Dinitrotoluene | 0.535 | 0.492 | 8.0 | 82 | 0.00 |
| 57 | | Fluorene | 1.537 | 1.419 | 7.7 | 84 | 0.00 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 0.509 | 0.440 | 13.6 | 76 | 0.00 |
| 59 | | Diethylphthalate | 1.662 | 1.505 | 9.4 | 81 | 0.00 |
| 60 | | 4-Chlorophenyl-phenylether | 0.936 | 0.816 | 12.8 | 77 | 0.00 |
| 61 | | 4-Nitroaniline | 0.396 | 0.370 | 6.6 | 84 | 0.00 |
| 62 | | Azobenzene | 1.590 | 1.554 | 2.3 | 88 | 0.00 |
| 63 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 87 | 0.00 |
| 64 | | 4,6-Dinitro-2-methylphenol | 0.149 | 0.130 | 12.8 | 73 | 0.00 |
| 65 | c | n-Nitrosodiphenylamine | 0.576 | 0.547 | 5.0 | 82 | 0.00 |
| 66 | | 4-Bromophenyl-phenylether | 0.260 | 0.225 | 13.5 | 75 | 0.00 |
| 67 | | Hexachlorobenzene | 0.283 | 0.242 | 14.5 | 74 | 0.00 |
| 68 | | Atrazine | 0.256 | 0.237 | 7.4 | 82 | 0.00 |
| 69 | C | Pentachlorophenol | 0.183 | 0.154 | 15.8 | 69 | 0.00 |
| 70 | | Phenanthrene | 0.980 | 0.914 | 6.7 | 83 | 0.00 |
| 71 | | Anthracene | 0.992 | 0.928 | 6.5 | 83 | 0.00 |
| 72 | | Carbazole | 0.858 | 0.809 | 5.7 | 84 | 0.00 |
| 73 | | Di-n-butylphthalate | 0.954 | 0.958 | -0.4 | 86 | 0.00 |
| 74 | C | Fluoranthene | 1.203 | 1.067 | 11.3 | 82 | 0.00 |
| 75 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 82 | 0.02 |
| 76 | | Benzidine | 0.573 | 0.365 | 36.3# | 50 | 0.00 |
| 77 | | Pyrene | 1.124 | 1.080 | 3.9 | 83 | 0.00 |
| 78 | S | Terphenyl-d14 | 0.646 | 0.660 | -2.2 | 84 | 0.00 |
| 79 | | Butylbenzylphthalate | 0.445 | 0.453 | -1.8 | 84 | 0.01 |
| 80 | | Benzo(a)anthracene | 1.119 | 1.055 | 5.7 | 79 | 0.02 |
| 81 | | 3,3'-Dichlorobenzidine | 0.491 | 0.445 | 9.4 | 73 | 0.02 |
| 82 | | Chrysene | 1.050 | 0.997 | 5.0 | 80 | 0.03 |
| 83 | | Bis(2-ethylhexyl)phthalate | 0.618 | 0.614 | 0.6 | 83 | 0.03 |
| 84 | c | Di-n-octyl phthalate | 1.031 | 1.032 | -0.1 | 82 | 0.04 |
| 85 | | Indeno(1,2,3-cd)pyrene | 1.338 | 1.127 | 15.8 | 71 | 0.11 |
| 86 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 79 | 0.06 |
| 87 | | Benzo(b)fluoranthene | 1.154 | 1.083 | 6.2 | 76 | 0.05 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 1.095 | 1.043 | 4.7 | 75 | 0.06 |
| 89 C | Benzo(a)pyrene | 1.106 | 1.029 | 7.0 | 75 | 0.06 |
| 90 | Dibenzo(a,h)anthracene | 1.098 | 0.971 | 11.6 | 71 | 0.10 |
| 91 | Benzo(g,h,i)perylene | 1.099 | 0.948 | 13.7 | 69 | 0.13 |

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