# Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
Data File : BF080360.D
Acq On : 6 Jul 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

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

Quant Time: Jul 07 05:39:19 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 07 03:05:40 2015
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 | 98 | 0.00 |
| 2 | | 1,4-Dioxane | 0.516 | 0.539 | -4.5 | 101 | -0.01 |
| 3 | | Pyridine | 1.312 | 1.414 | -7.8 | 111 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 0.617 | 0.692 | -12.2 | 110 | -0.01 |
| 5 | S | 2-Fluorophenol | 1.092 | 1.089 | 0.3 | 98 | 0.00 |
| 6 | | Aniline | 2.004 | 2.035 | -1.5 | 99 | 0.00 |
| 7 | S | Phenol-d6 | 1.448 | 1.414 | 2.3 | 99 | 0.00 |
| 8 | | 2-Chlorophenol | 1.266 | 1.270 | -0.3 | 98 | 0.00 |
| 9 | | Benzaldehyde | 0.805 | 0.827 | -2.7 | 101 | 0.00 |
| 10 | C | Phenol | 1.570 | 1.501 | 4.4 | 98 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 1.184 | 1.152 | 2.7 | 100 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 1.554 | 1.536 | 1.2 | 96 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 1.451 | 1.484 | -2.3 | 101 | -0.01 |
| 14 | | 1,2-Dichlorobenzene | 1.469 | 1.475 | -0.4 | 99 | 0.00 |
| 15 | | Benzyl Alcohol | 1.120 | 1.171 | -4.6 | 99 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 1.915 | 1.917 | -0.1 | 100 | 0.00 |
| 17 | | 2-Methylphenol | 0.981 | 0.981 | 0.0 | 98 | 0.00 |
| 18 | | Hexachloroethane | 0.466 | 0.464 | 0.4 | 96 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 0.939 | 0.945 | -0.6 | 99 | 0.00 |
| 20 | | 3+4-Methylphenols | 1.368 | 1.415 | -3.4 | 99 | 0.00 |
| 21 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 101 | 0.00 |
| 22 | | Acetophenone | 0.461 | 0.464 | -0.7 | 97 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 0.343 | 0.348 | -1.5 | 97 | 0.00 |
| 24 | | Nitrobenzene | 0.344 | 0.339 | 1.5 | 100 | 0.00 |
| 25 | | Isophorone | 0.660 | 0.670 | -1.5 | 101 | 0.00 |
| 26 | C | 2-Nitrophenol | 0.149 | 0.152 | -2.0 | 97 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 0.293 | 0.290 | 1.0 | 100 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 0.410 | 0.406 | 1.0 | 98 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 0.260 | 0.270 | -3.8 | 103 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 0.309 | 0.299 | 3.2 | 97 | 0.00 |
| 31 | | Naphthalene | 1.026 | 1.023 | 0.3 | 99 | 0.00 |
| 32 | | Benzoic acid | 0.143 | 0.171 | -19.6 | 109 | 0.00 |
| 33 | | 4-Chloroaniline | 0.409 | 0.413 | -1.0 | 103 | 0.00 |
| 34 | C | Hexachlorobutadiene | 0.188 | 0.184 | 2.1 | 97 | 0.00 |
| 35 | | Caprolactam | 0.080 | 0.081 | -1.3 | 100 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 0.269 | 0.261 | 3.0 | 99 | 0.00 |
| 37 | | 2-Methylnaphthalene | 0.654 | 0.642 | 1.8 | 97 | 0.00 |
| 38 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 103 | 0.00 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 0.634 | 0.616 | 2.8 | 98 | 0.00 |
| 40 | P | Hexachlorocyclopentadiene | 0.223 | 0.220 | 1.3 | 97 | 0.00 |
| 41 | S | 2,4,6-Tribromophenol | 0.209 | 0.212 | -1.4 | 99 | 0.00 |
| 42 | C | 2,4,6-Trichlorophenol | 0.418 | 0.415 | 0.7 | 96 | 0.00 |
| 43 | | 2,4,5-Trichlorophenol | 0.487 | 0.488 | -0.2 | 100 | 0.00 |
| 44 | S | 2-Fluorobiphenyl | 1.482 | 1.467 | 1.0 | 100 | 0.00 |

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| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 1.775 | 1.748 | 1.5 | 100 | 0.00 |
| 46 | | 2-Chloronaphthalene | 1.354 | 1.348 | 0.4 | 99 | 0.00 |
| 47 | | 2-Nitroaniline | 0.379 | 0.396 | -4.5 | 101 | 0.00 |
| 48 | | Acenaphthylene | 2.253 | 2.193 | 2.7 | 99 | 0.00 |
| 49 | | Dimethylphthalate | 1.606 | 1.456 | 9.3 | 99 | 0.00 |
| 50 | | 2,6-Dinitrotoluene | 0.336 | 0.343 | -2.1 | 100 | 0.00 |
| 51 | C | Acenaphthene | 1.349 | 1.367 | -1.3 | 102 | 0.00 |
| 52 | | 3-Nitroaniline | 0.378 | 0.384 | -1.6 | 97 | 0.00 |
| 53 | P | 2,4-Dinitrophenol | 0.120 | 0.118 | 1.7 | 92 | 0.00 |
| 54 | | Dibenzofuran | 1.918 | 1.922 | -0.2 | 103 | 0.00 |
| 55 | P | 4-Nitrophenol | 0.277 | 0.299 | -7.9 | 99 | 0.00 |
| 56 | | 2,4-Dinitrotoluene | 0.432 | 0.434 | -0.5 | 100 | 0.00 |
| 57 | | Fluorene | 1.651 | 1.650 | 0.1 | 99 | 0.00 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 0.390 | 0.395 | -1.3 | 99 | 0.00 |
| 59 | | Diethylphthalate | 1.604 | 1.579 | 1.6 | 98 | 0.00 |
| 60 | | 4-Chlorophenyl-phenylether | 0.731 | 0.727 | 0.5 | 99 | 0.00 |
| 61 | | 4-Nitroaniline | 0.367 | 0.397 | -8.2 | 100 | 0.00 |
| 62 | | Azobenzene | 1.523 | 1.513 | 0.7 | 98 | 0.00 |
| 63 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 100 | 0.00 |
| 64 | | 4,6-Dinitro-2-methylphenol | 0.106 | 0.105 | 0.9 | 94 | 0.00 |
| 65 | c | n-Nitrosodiphenylamine | 0.637 | 0.647 | -1.6 | 100 | 0.00 |
| 66 | | 4-Bromophenyl-phenylether | 0.229 | 0.227 | 0.9 | 98 | 0.00 |
| 67 | | Hexachlorobenzene | 0.240 | 0.246 | -2.5 | 101 | 0.00 |
| 68 | | Atrazine | 0.190 | 0.200 | -5.3 | 105 | 0.00 |
| 69 | C | Pentachlorophenol | 0.119 | 0.100 | 16.0 | 88 | 0.00 |
| 70 | | Phenanthrene | 1.200 | 1.186 | 1.2 | 104 | 0.00 |
| 71 | | Anthracene | 1.176 | 1.171 | 0.4 | 97 | 0.00 |
| 72 | | Carbazole | 1.097 | 1.136 | -3.6 | 105 | 0.00 |
| 73 | | Di-n-butylphthalate | 1.255 | 1.238 | 1.4 | 98 | 0.00 |
| 74 | C | Fluoranthene | 1.289 | 1.247 | 3.3 | 98 | -0.01 |
| 75 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 100 | -0.01 |
| 76 | | Benzidine | 0.507 | 0.550 | -8.5 | 102 | 0.00 |
| 77 | | Pyrene | 1.210 | 1.274 | -5.3 | 103 | 0.00 |
| 78 | S | Terphenyl-d14 | 0.857 | 0.914 | -6.7 | 106 | 0.00 |
| 79 | | Butylbenzylphthalate | 0.483 | 0.530 | -9.7 | 100 | -0.01 |
| 80 | | Benzo(a)anthracene | 1.205 | 1.221 | -1.3 | 99 | -0.01 |
| 81 | | 3,3'-Dichlorobenzidine | 0.400 | 0.417 | -4.2 | 97 | -0.01 |
| 82 | | Chrysene | 1.128 | 1.159 | -2.7 | 98 | -0.01 |
| 83 | | Bis(2-ethylhexyl)phthalate | 0.736 | 0.794 | -7.9 | 97 | -0.01 |
| 84 | c | Di-n-octyl phthalate | 1.215 | 1.299 | -6.9 | 97 | -0.01 |
| 85 | | Indeno(1,2,3-cd)pyrene | 1.326 | 1.354 | -2.1 | 97 | 0.00 |
| 86 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 97 | 0.00 |
| 87 | | Benzo(b)fluoranthene | 1.257 | 1.213 | 3.5 | 87 | 0.00 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 1.190 | 1.301 | -9.3 | 112 | 0.00 |
| 89 C | Benzo(a)pyrene | 1.149 | 1.187 | -3.3 | 98 | -0.01 |
| 90 | Dibenzo(a,h)anthracene | 1.132 | 1.157 | -2.2 | 97 | 0.00 |
| 91 | Benzo(g,h,i)perylene | 1.148 | 1.178 | -2.6 | 97 | 0.00 |

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