# Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
Data File : BF077394.D
Acq On : 20 Feb 2015 21:05
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SSTDCCC040

Quant Time: Feb 21 01:38:14 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 20 23:55:25 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 | 116 | 0.00 |
| 2 | | 1,4-Dioxane | 0.508 | 0.523 | -3.0 | 121 | 0.00 |
| 3 | | Pyridine | 1.454 | 1.160 | 20.2 | 87 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 0.632 | 0.653 | -3.3 | 124 | 0.00 |
| 5 | S | 2-Fluorophenol | 1.198 | 1.198 | 0.0 | 115 | 0.00 |
| 6 | | Aniline | 2.129 | 2.127 | 0.1 | 112 | 0.00 |
| 7 | S | Phenol-d6 | 1.540 | 1.482 | 3.8 | 108 | 0.00 |
| 8 | | 2-Chlorophenol | 1.413 | 1.420 | -0.5 | 115 | 0.00 |
| 9 | | Benzaldehyde | 0.935 | 0.875 | 6.4 | 110 | -0.01 |
| 10 | C | Phenol | 1.828 | 1.840 | -0.7 | 110 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 1.313 | 1.246 | 5.1 | 110 | -0.01 |
| 12 | | 1,3-Dichlorobenzene | 1.539 | 1.542 | -0.2 | 114 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 1.566 | 1.565 | 0.1 | 113 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 1.489 | 1.511 | -1.5 | 114 | 0.00 |
| 15 | | Benzyl Alcohol | 1.171 | 1.128 | 3.7 | 109 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 1.877 | 1.860 | 0.9 | 111 | -0.01 |
| 17 | | 2-Methylphenol | 1.105 | 1.141 | -3.3 | 111 | 0.00 |
| 18 | | Hexachloroethane | 0.523 | 0.521 | 0.4 | 113 | -0.01 |
| 19 | P | n-Nitroso-di-n-propylamine | 0.978 | 1.035 | -5.8 | 117 | 0.00 |
| 20 | | 3+4-Methylphenols | 1.455 | 1.437 | 1.2 | 110 | -0.01 |
| 21 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 111 | 0.00 |
| 22 | | Acetophenone | 0.470 | 0.469 | 0.2 | 113 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 0.322 | 0.324 | -0.6 | 108 | -0.01 |
| 24 | | Nitrobenzene | 0.338 | 0.343 | -1.5 | 110 | -0.01 |
| 25 | | Isophorone | 0.638 | 0.652 | -2.2 | 106 | -0.01 |
| 26 | C | 2-Nitrophenol | 0.139 | 0.157 | -12.9 | 113 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 0.310 | 0.324 | -4.5 | 113 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 0.403 | 0.438 | -8.7 | 117 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 0.291 | 0.310 | -6.5 | 114 | -0.01 |
| 30 | | 1,2,4-Trichlorobenzene | 0.320 | 0.337 | -5.3 | 114 | 0.00 |
| 31 | | Naphthalene | 1.000 | 1.023 | -2.3 | 114 | 0.00 |
| 32 | | Benzoic acid | 0.151 | 0.176 | -16.6 | 118 | 0.00 |
| 33 | | 4-Chloroaniline | 0.445 | 0.439 | 1.3 | 104 | -0.01 |
| 34 | C | Hexachlorobutadiene | 0.183 | 0.195 | -6.6 | 116 | 0.00 |
| 35 | | Caprolactam | 0.089 | 0.094 | -5.6 | 111 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 0.293 | 0.315 | -7.5 | 112 | 0.00 |
| 37 | | 2-Methylnaphthalene | 0.710 | 0.719 | -1.3 | 106 | -0.01 |
| 38 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 112 | -0.01 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 0.574 | 0.570 | 0.7 | 107 | -0.01 |
| 40 | P | Hexachlorocyclopentadiene | 0.308 | 0.312 | -1.3 | 103 | -0.01 |
| 41 | S | 2,4,6-Tribromophenol | 0.193 | 0.202 | -4.7 | 109 | 0.00 |
| 42 | C | 2,4,6-Trichlorophenol | 0.380 | 0.405 | -6.6 | 110 | 0.00 |
| 43 | | 2,4,5-Trichlorophenol | 0.406 | 0.408 | -0.5 | 106 | 0.00 |
| 44 | S | 2-Fluorobiphenyl | 1.283 | 1.313 | -2.3 | 114 | 0.00 |

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| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 1.515 | 1.484 | 2.0 | 106 | -0.01 |
| 46 | | 2-Chloronaphthalene | 1.188 | 1.165 | 1.9 | 109 | 0.00 |
| 47 | | 2-Nitroaniline | 0.293 | 0.310 | -5.8 | 113 | 0.00 |
| 48 | | Acenaphthylene | 1.881 | 1.880 | 0.1 | 110 | 0.00 |
| 49 | | Dimethylphthalate | 1.406 | 1.366 | 2.8 | 107 | 0.00 |
| 50 | | 2,6-Dinitrotoluene | 0.282 | 0.290 | -2.8 | 102 | -0.01 |
| 51 | C | Acenaphthene | 1.123 | 1.143 | -1.8 | 110 | 0.00 |
| 52 | | 3-Nitroaniline | 0.325 | 0.358 | -10.2 | 113 | 0.00 |
| 53 | P | 2,4-Dinitrophenol | 0.086 | 0.079 | 8.1 | 98 | -0.01 |
| 54 | | Dibenzofuran | 1.733 | 1.766 | -1.9 | 111 | 0.00 |
| 55 | P | 4-Nitrophenol | 0.261 | 0.258 | 1.1 | 99 | -0.01 |
| 56 | | 2,4-Dinitrotoluene | 0.381 | 0.399 | -4.7 | 103 | 0.00 |
| 57 | | Fluorene | 1.386 | 1.388 | -0.1 | 108 | 0.00 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 0.327 | 0.342 | -4.6 | 107 | 0.00 |
| 59 | | Diethylphthalate | 1.386 | 1.383 | 0.2 | 108 | 0.00 |
| 60 | | 4-Chlorophenyl-phenylether | 0.668 | 0.656 | 1.8 | 106 | 0.00 |
| 61 | | 4-Nitroaniline | 0.312 | 0.330 | -5.8 | 111 | 0.00 |
| 62 | | Azobenzene | 1.315 | 1.347 | -2.4 | 108 | 0.00 |
| 63 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 106 | 0.00 |
| 64 | | 4,6-Dinitro-2-methylphenol | 0.075 | 0.069 | 8.0 | 93 | -0.01 |
| 65 | c | n-Nitrosodiphenylamine | 0.664 | 0.678 | -2.1 | 104 | -0.01 |
| 66 | | 4-Bromophenyl-phenylether | 0.234 | 0.237 | -1.3 | 108 | 0.00 |
| 67 | | Hexachlorobenzene | 0.251 | 0.255 | -1.6 | 110 | 0.00 |
| 68 | | Atrazine | 0.216 | 0.194 | 10.2 | 101 | 0.00 |
| 69 | C | Pentachlorophenol | 0.132 | 0.138 | -4.5 | 104 | 0.00 |
| 70 | | Phenanthrene | 1.159 | 1.128 | 2.7 | 104 | -0.01 |
| 71 | | Anthracene | 1.150 | 1.147 | 0.3 | 108 | 0.00 |
| 72 | | Carbazole | 1.094 | 1.151 | -5.2 | 107 | 0.00 |
| 73 | | Di-n-butylphthalate | 1.284 | 1.271 | 1.0 | 99 | 0.00 |
| 74 | C | Fluoranthene | 1.266 | 1.287 | -1.7 | 106 | 0.00 |
| 75 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 98 | 0.01 |
| 76 | | Benzidine | 0.676 | 0.613 | 9.3 | 96 | 0.00 |
| 77 | | Pyrene | 1.223 | 1.244 | -1.7 | 101 | -0.01 |
| 78 | S | Terphenyl-d14 | 0.856 | 0.872 | -1.9 | 100 | 0.00 |
| 79 | | Butylbenzylphthalate | 0.503 | 0.554 | -10.1 | 103 | 0.00 |
| 80 | | Benzo(a)anthracene | 1.172 | 1.170 | 0.2 | 100 | 0.01 |
| 81 | | 3,3'-Dichlorobenzidine | 0.430 | 0.456 | -6.0 | 102 | 0.01 |
| 82 | | Chrysene | 1.101 | 1.094 | 0.6 | 101 | 0.01 |
| 83 | | Bis(2-ethylhexyl)phthalate | 0.807 | 0.827 | -2.5 | 99 | 0.01 |
| 84 | c | Di-n-octyl phthalate | 1.348 | 1.355 | -0.5 | 94 | 0.05 |
| 85 | | Indeno(1,2,3-cd)pyrene | 1.255 | 1.290 | -2.8 | 101 | 0.09 |
| 86 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 99 | 0.07 |
| 87 | | Benzo(b)fluoranthene | 1.163 | 1.174 | -0.9 | 94 | 0.06 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 1.140 | 1.163 | -2.0 | 108 | 0.07 |
| 89 C | Benzo(a)pyrene | 1.108 | 1.137 | -2.6 | 98 | 0.08 |
| 90 | Dibenzo(a,h)anthracene | 1.056 | 1.093 | -3.5 | 100 | 0.09 |
| 91 | Benzo(g,h,i)perylene | 1.066 | 1.113 | -4.4 | 101 | 0.09 |

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