# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012320\
Data File : BF118677.D
Acq On : 23 Jan 2020 15:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Jan 24 02:41:26 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 20 14:45:03 2020
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 | 79 | 0.00 |
| 2 | | 1,4-Dioxane | 0.522 | 0.517 | 1.0 | 75 | -0.02 |
| 3 | | Pyridine | 1.470 | 1.466 | 0.3 | 75 | -0.02 |
| 4 | | n-Nitrosodimethylamine | 0.630 | 0.539 | 14.4 | 64 | -0.02 |
| 5 | S | 2-Fluorophenol | 1.077 | 1.093 | -1.5 | 76 | 0.00 |
| 6 | | Aniline | 1.843 | 1.846 | -0.2 | 75 | 0.00 |
| 7 | S | Phenol-d6 | 1.400 | 1.396 | 0.3 | 74 | 0.00 |
| 8 | | 2-Chlorophenol | 1.209 | 1.215 | -0.5 | 75 | 0.00 |
| 9 | | Benzaldehyde | 0.851 | 0.847 | 0.5 | 77 | 0.00 |
| 10 | C | Phenol | 1.595 | 1.549 | 2.9 | 73 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 1.183 | 1.188 | -0.4 | 76 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 1.415 | 1.469 | -3.8 | 78 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 1.420 | 1.471 | -3.6 | 77 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 1.335 | 1.372 | -2.8 | 76 | 0.00 |
| 15 | | Benzyl Alcohol | 1.110 | 1.084 | 2.3 | 74 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 1.648 | 1.542 | 6.4 | 70 | -0.01 |
| 17 | | 2-Methylphenol | 1.007 | 1.010 | -0.3 | 76 | 0.00 |
| 18 | | Hexachloroethane | 0.515 | 0.499 | 3.1 | 73 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 0.945 | 0.851 | 9.9 | 69 | 0.00 |
| 20 | | 3+4-Methylphenols | 1.229 | 1.204 | 2.0 | 72 | 0.00 |
| 21 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 78 | 0.00 |
| 22 | | Acetophenone | 0.474 | 0.460 | 3.0 | 72 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 0.358 | 0.332 | 7.3 | 70 | 0.00 |
| 24 | | Nitrobenzene | 0.366 | 0.341 | 6.8 | 70 | 0.00 |
| 25 | | Isophorone | 0.652 | 0.629 | 3.5 | 74 | 0.00 |
| 26 | C | 2-Nitrophenol | 0.156 | 0.148 | 5.1 | 73 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 0.270 | 0.254 | 5.9 | 71 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 0.420 | 0.414 | 1.4 | 75 | -0.01 |
| 29 | C | 2,4-Dichlorophenol | 0.298 | 0.302 | -1.3 | 76 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 0.346 | 0.354 | -2.3 | 77 | 0.00 |
| 31 | | Naphthalene | 0.903 | 0.960 | -6.3 | 78 | 0.00 |
| 32 | | Benzoic acid | 0.202 | 0.176 | 12.9 | 67 | 0.00 |
| 33 | | 4-Chloroaniline | 0.418 | 0.434 | -3.8 | 77 | 0.00 |
| 34 | C | Hexachlorobutadiene | 0.210 | 0.212 | -1.0 | 76 | -0.01 |
| 35 | | Caprolactam | 0.098 | 0.100 | -2.0 | 81 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 0.294 | 0.297 | -1.0 | 76 | 0.00 |
| 37 | | 2-Methylnaphthalene | 0.638 | 0.665 | -4.2 | 77 | 0.00 |
| 38 | | 1-Methylnaphthalene | 0.606 | 0.622 | -2.6 | 76 | 0.00 |
| 39 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 79 | -0.01 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 0.626 | 0.635 | -1.4 | 75 | 0.00 |
| 41 | P | Hexachlorocyclopentadiene | 0.316 | 0.253 | 19.9 | 60 | 0.00 |
| 42 | S | 2,4,6-Tribromophenol | 0.231 | 0.228 | 1.3 | 74 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 0.415 | 0.404 | 2.7 | 74 | 0.00 |
| 44 | | 2,4,5-Trichlorophenol | 0.424 | 0.426 | -0.5 | 76 | 0.00 |

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|---|---|---|---|---|---|---|---|
| 45 | S | 2-Fluorobiphenyl | 1.217 | 1.144 | 6.0 | 76 | 0.00 |
| 46 | | 1,1'-Biphenyl | 1.378 | 1.459 | -5.9 | 77 | 0.00 |
| 47 | | 2-Chloronaphthalene | 1.162 | 1.216 | -4.6 | 77 | 0.00 |
| 48 | | 2-Nitroaniline | 0.357 | 0.320 | 10.4 | 68 | 0.00 |
| 49 | | Acenaphthylene | 1.634 | 1.750 | -7.1 | 79 | 0.00 |
| 50 | | Dimethylphthalate | 1.303 | 1.376 | -5.6 | 80 | 0.00 |
| 51 | | 2,6-Dinitrotoluene | 0.292 | 0.287 | 1.7 | 75 | 0.00 |
| 52 | C | Acenaphthene | 1.093 | 1.118 | -2.3 | 76 | 0.00 |
| 53 | | 3-Nitroaniline | 0.333 | 0.330 | 0.9 | 75 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 0.116 | 0.111 | 4.3 | 78 | 0.00 |
| 55 | | Dibenzofuran | 1.515 | 1.599 | -5.5 | 77 | 0.00 |
| 56 | P | 4-Nitrophenol | 0.252 | 0.246 | 2.4 | 74 | 0.00 |
| 57 | | 2,4-Dinitrotoluene | 0.368 | 0.369 | -0.3 | 78 | 0.00 |
| 58 | | Fluorene | 1.189 | 1.211 | -1.9 | 75 | -0.01 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 0.373 | 0.373 | 0.0 | 75 | 0.00 |
| 60 | | Diethylphthalate | 1.263 | 1.314 | -4.0 | 78 | -0.01 |
| 61 | | 4-Chlorophenyl-phenylether | 0.647 | 0.652 | -0.8 | 73 | 0.00 |
| 62 | | 4-Nitroaniline | 0.342 | 0.346 | -1.2 | 78 | -0.01 |
| 63 | | Azobenzene | 1.224 | 1.213 | 0.9 | 74 | 0.00 |
| 64 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 82 | -0.01 |
| 65 | | 4,6-Dinitro-2-methylphenol | 0.088 | 0.088 | 0.0 | 82 | 0.00 |
| 66 | c | n-Nitrosodiphenylamine | 0.603 | 0.602 | 0.2 | 77 | 0.00 |
| 67 | | 4-Bromophenyl-phenylether | 0.247 | 0.246 | 0.4 | 78 | -0.01 |
| 68 | | Hexachlorobenzene | 0.280 | 0.279 | 0.4 | 78 | 0.00 |
| 69 | | Atrazine | 0.209 | 0.207 | 1.0 | 78 | 0.00 |
| 70 | C | Pentachlorophenol | 0.156 | 0.148 | 5.1 | 75 | 0.00 |
| 71 | | Phenanthrene | 0.965 | 1.002 | -3.8 | 79 | 0.00 |
| 72 | | Anthracene | 0.955 | 0.996 | -4.3 | 79 | 0.00 |
| 73 | | Carbazole | 0.876 | 0.913 | -4.2 | 79 | 0.00 |
| 74 | | Di-n-butylphthalate | 0.973 | 1.042 | -7.1 | 82 | -0.01 |
| 75 | C | Fluoranthene | 1.018 | 1.086 | -6.7 | 80 | 0.00 |
| 76 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 81 | 0.00 |
| 77 | | Benzidine | 0.534 | 0.653 | -22.3 | 93 | 0.00 |
| 78 | | Pyrene | 1.358 | 1.404 | -3.4 | 80 | 0.00 |
| 79 | S | Terphenyl-d14 | 0.917 | 0.935 | -2.0 | 79 | 0.00 |
| 80 | | Butylbenzylphthalate | 0.556 | 0.582 | -4.7 | 83 | -0.01 |
| 81 | | Benzo(a)anthracene | 1.203 | 1.246 | -3.6 | 79 | -0.01 |
| 82 | | 3,3'-Dichlorobenzidine | 0.429 | 0.498 | -16.1 | 87 | -0.01 |
| 83 | | Chrysene | 1.153 | 1.213 | -5.2 | 81 | 0.00 |
| 84 | | Bis(2-ethylhexyl)phthalate | 0.679 | 0.728 | -7.2 | 84 | -0.02 |
| 85 | c | Di-n-octyl phthalate | 1.005 | 1.177 | -17.1 | 94 | -0.02 |
| 86 | | Indeno(1,2,3-cd)pyrene | 1.002 | 1.144 | -14.2 | 99 | 0.00 |
| 87 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 100 | 0.00 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(b)fluoranthene | 1.257 | 1.220 | 2.9 | 89 | 0.00 |
| 89 | Benzo(k)fluoranthene | 1.165 | 1.186 | -1.8 | 98 | 0.00 |
| 90 C | Benzo(a)pyrene | 1.084 | 1.102 | -1.7 | 97 | 0.00 |
| 91 | Dibenzo(a,h)anthracene | 0.959 | 0.924 | 3.6 | 96 | 0.00 |
| 92 | Benzo(g,h,i)perylene | 0.931 | 0.901 | 3.2 | 97 | -0.01 |

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