# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112119\
Data File : BF117900.D
Acq On : 21 Nov 2019 9:46
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Nov 21 15:14:43 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 13 18:36:07 2019
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 | 88 | 0.00 |
| 2 | | 1,4-Dioxane | 40.000 | 45.386 | -13.5 | 102 | 0.00 |
| 3 | | Pyridine | 40.000 | 33.809 | 15.5 | 76 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 34.753 | 13.1 | 79 | 0.00 |
| 5 | S | 2-Fluorophenol | 80.000 | 77.116 | 3.6 | 89 | 0.00 |
| 6 | | Aniline | 40.000 | 39.813 | 0.5 | 89 | 0.00 |
| 7 | S | Phenol-d6 | 80.000 | 78.891 | 1.4 | 91 | 0.00 |
| 8 | | 2-Chlorophenol | 40.000 | 41.008 | -2.5 | 92 | 0.00 |
| 9 | | Benzaldehyde | 40.000 | 37.641 | 5.9 | 84 | 0.00 |
| 10 | C | Phenol | 40.000 | 43.329 | -8.3 | 98 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 39.699 | 0.8 | 89 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 40.427 | -1.1 | 91 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.959 | 0.1 | 89 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 40.225 | -0.6 | 92 | 0.00 |
| 15 | | Benzyl Alcohol | 40.000 | 42.885 | -7.2 | 95 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 37.236 | 6.9 | 81 | 0.00 |
| 17 | | 2-Methylphenol | 40.000 | 40.528 | -1.3 | 90 | 0.00 |
| 18 | | Hexachloroethane | 40.000 | 41.431 | -3.6 | 92 | -0.01 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 40.421 | -1.1 | 92 | 0.00 |
| 20 | | 3+4-Methylphenols | 40.000 | 39.338 | 1.7 | 89 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 90 | 0.00 |
| 22 | | Acetophenone | 40.000 | 38.751 | 3.1 | 91 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 80.290 | -0.4 | 93 | 0.00 |
| 24 | | Nitrobenzene | 40.000 | 40.256 | -0.6 | 93 | 0.00 |
| 25 | | Isophorone | 40.000 | 38.162 | 4.6 | 91 | 0.00 |
| 26 | C | 2-Nitrophenol | 40.000 | 43.074 | -7.7 | 99 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 37.788 | 5.5 | 87 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 38.716 | 3.2 | 91 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 38.946 | 2.6 | 91 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.048 | 2.4 | 91 | 0.00 |
| 31 | | Naphthalene | 40.000 | 40.192 | -0.5 | 93 | 0.00 |
| 32 | | Benzoic acid | 40.000 | 40.495 | -1.2 | 95 | 0.02 |
| 33 | | 4-Chloroaniline | 40.000 | 39.614 | 1.0 | 93 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 40.217 | -0.5 | 93 | 0.00 |
| 35 | | Caprolactam | 40.000 | 39.755 | 0.6 | 94 | 0.01 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 40.249 | -0.6 | 95 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 39.729 | 0.7 | 92 | 0.00 |
| 38 | | 1-Methylnaphthalene | 40.000 | 40.029 | -0.1 | 93 | 0.00 |
| 39 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 93 | 0.00 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 38.849 | 2.9 | 93 | 0.00 |
| 41 | P | Hexachlorocyclopentadiene | 40.000 | 36.709 | 8.2 | 88 | -0.01 |
| 42 | S | 2,4,6-Tribromophenol | 80.000 | 81.543 | -1.9 | 99 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 40.000 | 38.716 | 3.2 | 92 | 0.00 |
| 44 | | 2,4,5-Trichlorophenol | 40.000 | 38.736 | 3.2 | 92 | 0.00 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | S | 2-Fluorobiphenyl | 80.000 | 82.575 | -3.2 | 93 | 0.00 |
| 46 | | 1,1'-Biphenyl | 40.000 | 39.389 | 1.5 | 93 | 0.00 |
| 47 | | 2-Chloronaphthalene | 40.000 | 38.965 | 2.6 | 93 | 0.00 |
| 48 | | 2-Nitroaniline | 40.000 | 40.435 | -1.1 | 96 | 0.00 |
| 49 | | Acenaphthylene | 40.000 | 39.794 | 0.5 | 94 | 0.00 |
| 50 | | Dimethylphthalate | 40.000 | 39.711 | 0.7 | 96 | 0.00 |
| 51 | | 2,6-Dinitrotoluene | 40.000 | 39.925 | 0.2 | 95 | 0.00 |
| 52 | C | Acenaphthene | 40.000 | 39.125 | 2.2 | 94 | 0.00 |
| 53 | | 3-Nitroaniline | 40.000 | 40.472 | -1.2 | 96 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 40.000 | 42.067 | -5.2 | 110 | 0.00 |
| 55 | | Dibenzofuran | 40.000 | 39.042 | 2.4 | 93 | 0.00 |
| 56 | P | 4-Nitrophenol | 40.000 | 40.383 | -1.0 | 96 | 0.01 |
| 57 | | 2,4-Dinitrotoluene | 40.000 | 42.370 | -5.9 | 102 | 0.00 |
| 58 | | Fluorene | 40.000 | 39.189 | 2.0 | 97 | 0.00 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 38.339 | 4.2 | 92 | 0.00 |
| 60 | | Diethylphthalate | 40.000 | 40.230 | -0.6 | 96 | 0.00 |
| 61 | | 4-Chlorophenyl-phenylether | 40.000 | 39.686 | 0.8 | 95 | 0.00 |
| 62 | | 4-Nitroaniline | 40.000 | 42.008 | -5.0 | 105 | 0.00 |
| 63 | | Azobenzene | 40.000 | 38.905 | 2.7 | 93 | 0.00 |
| 64 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 96 | 0.00 |
| 65 | | 4,6-Dinitro-2-methylphenol | 40.000 | 44.763 | -11.9 | 112 | 0.00 |
| 66 | c | n-Nitrosodiphenylamine | 40.000 | 35.869 | 10.3 | 89 | 0.00 |
| 67 | | 4-Bromophenyl-phenylether | 40.000 | 39.111 | 2.2 | 97 | 0.00 |
| 68 | | Hexachlorobenzene | 40.000 | 37.538 | 6.2 | 92 | 0.00 |
| 69 | | Atrazine | 40.000 | 39.357 | 1.6 | 98 | 0.00 |
| 70 | C | Pentachlorophenol | 40.000 | 37.266 | 6.8 | 92 | 0.00 |
| 71 | | Phenanthrene | 40.000 | 37.232 | 6.9 | 95 | 0.00 |
| 72 | | Anthracene | 40.000 | 37.870 | 5.3 | 97 | 0.00 |
| 73 | | Carbazole | 40.000 | 38.762 | 3.1 | 96 | 0.00 |
| 74 | | Di-n-butylphthalate | 40.000 | 39.705 | 0.7 | 102 | 0.00 |
| 75 | C | Fluoranthene | 40.000 | 42.468 | -6.2 | 101 | 0.00 |
| 76 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 106 | 0.00 |
| 77 | | Benzidine | 40.000 | 34.824 | 12.9 | 92 | 0.00 |
| 78 | | Pyrene | 40.000 | 37.203 | 7.0 | 102 | 0.00 |
| 79 | S | Terphenyl-d14 | 80.000 | 75.618 | 5.5 | 104 | 0.00 |
| 80 | | Butylbenzylphthalate | 40.000 | 40.516 | -1.3 | 113 | 0.00 |
| 81 | | Benzo(a)anthracene | 40.000 | 39.068 | 2.3 | 108 | 0.00 |
| 82 | | 3,3'-Dichlorobenzidine | 40.000 | 39.919 | 0.2 | 109 | 0.00 |
| 83 | | Chrysene | 40.000 | 38.444 | 3.9 | 109 | 0.00 |
| 84 | | Bis(2-ethylhexyl)phthalate | 40.000 | 42.752 | -6.9 | 122 | 0.00 |
| 85 | c | Di-n-octyl phthalate | 40.000 | 43.615 | -9.0 | 125 | 0.00 |
| 86 | | Indeno(1,2,3-cd)pyrene | 40.000 | 36.662 | 8.3 | 113 | 0.01 |
| 87 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 115 | 0.00 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(b)fluoranthene | 40.000 | 39.392 | 1.5 | 114 | 0.00 |
| 89 | Benzo(k)fluoranthene | 40.000 | 36.933 | 7.7 | 112 | 0.00 |
| 90 C | Benzo(a)pyrene | 40.000 | 38.063 | 4.8 | 113 | 0.00 |
| 91 | Dibenzo(a,h)anthracene | 40.000 | 37.346 | 6.6 | 115 | 0.01 |
| 92 | Benzo(g,h,i)perylene | 40.000 | 36.262 | 9.3 | 112 | 0.02 |

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