# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102618\
Data File : BF110207.D
Acq On : 26 Oct 2018 14:03
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Oct 26 17:20:51 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 23 15:06:18 2018
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 | 69 | -0.01 |
| 2 | | 1,4-Dioxane | 40.000 | 38.185 | 4.5 | 67 | -0.02 |
| 3 | | Pyridine | 40.000 | 38.857 | 2.9 | 64 | -0.01 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 39.090 | 2.3 | 66 | -0.02 |
| 5 | S | 2-Fluorophenol | 80.000 | 78.772 | 1.5 | 68 | -0.01 |
| 6 | | Aniline | 40.000 | 39.489 | 1.3 | 68 | -0.01 |
| 7 | S | Phenol-d6 | 80.000 | 78.722 | 1.6 | 68 | -0.01 |
| 8 | | 2-Chlorophenol | 40.000 | 39.915 | 0.2 | 68 | -0.02 |
| 9 | | Benzaldehyde | 40.000 | 37.444 | 6.4 | 64 | -0.01 |
| 10 | C | Phenol | 40.000 | 39.164 | 2.1 | 68 | -0.01 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 38.616 | 3.5 | 67 | -0.01 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 39.552 | 1.1 | 69 | -0.01 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.363 | 1.6 | 69 | -0.01 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 39.637 | 0.9 | 69 | -0.02 |
| 15 | | Benzyl Alcohol | 40.000 | 40.873 | -2.2 | 69 | -0.02 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 39.350 | 1.6 | 68 | -0.01 |
| 17 | | 2-Methylphenol | 40.000 | 39.156 | 2.1 | 67 | -0.01 |
| 18 | | Hexachloroethane | 40.000 | 40.074 | -0.2 | 70 | -0.02 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 38.762 | 3.1 | 67 | -0.02 |
| 20 | | 3+4-Methylphenols | 40.000 | 39.514 | 1.2 | 68 | -0.01 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 69 | -0.02 |
| 22 | | Acetophenone | 40.000 | 39.614 | 1.0 | 69 | -0.01 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 80.776 | -1.0 | 68 | -0.02 |
| 24 | | Nitrobenzene | 40.000 | 40.746 | -1.9 | 69 | -0.02 |
| 25 | | Isophorone | 40.000 | 38.794 | 3.0 | 66 | -0.02 |
| 26 | C | 2-Nitrophenol | 40.000 | 43.065 | -7.7 | 70 | -0.01 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 39.527 | 1.2 | 68 | -0.02 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 38.951 | 2.6 | 68 | -0.02 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 40.313 | -0.8 | 68 | -0.01 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.847 | 0.4 | 68 | -0.01 |
| 31 | | Naphthalene | 40.000 | 39.709 | 0.7 | 69 | -0.02 |
| 32 | | Benzoic acid | 40.000 | 40.236 | -0.6 | 67 | 0.00 |
| 33 | | 4-Chloroaniline | 40.000 | 39.632 | 0.9 | 69 | -0.01 |
| 34 | C | Hexachlorobutadiene | 40.000 | 40.398 | -1.0 | 71 | -0.01 |
| 35 | | Caprolactam | 40.000 | 40.171 | -0.4 | 67 | -0.01 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 39.597 | 1.0 | 67 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 39.896 | 0.3 | 69 | -0.01 |
| 38 | | 1-Methylnaphthalene | 40.000 | 39.125 | 2.2 | 68 | -0.01 |
| 39 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 70 | -0.01 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 39.240 | 1.9 | 69 | -0.01 |
| 41 | P | Hexachlorocyclopentadiene | 40.000 | 38.454 | 3.9 | 63 | -0.01 |
| 42 | S | 2,4,6-Tribromophenol | 80.000 | 80.195 | -0.2 | 69 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 40.000 | 40.011 | -0.0 | 68 | -0.01 |
| 44 | | 2,4,5-Trichlorophenol | 40.000 | 39.667 | 0.8 | 68 | -0.01 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | S | 2-Fluorobiphenyl | 80.000 | 80.363 | -0.5 | 70 | -0.01 |
| 46 | | 1,1'-Biphenyl | 40.000 | 39.243 | 1.9 | 70 | -0.01 |
| 47 | | 2-Chloronaphthalene | 40.000 | 39.137 | 2.2 | 68 | -0.02 |
| 48 | | 2-Nitroaniline | 40.000 | 41.296 | -3.2 | 70 | -0.01 |
| 49 | | Acenaphthylene | 40.000 | 39.317 | 1.7 | 69 | -0.02 |
| 50 | | Dimethylphthalate | 40.000 | 38.914 | 2.7 | 69 | -0.01 |
| 51 | | 2,6-Dinitrotoluene | 40.000 | 41.813 | -4.5 | 70 | -0.01 |
| 52 | C | Acenaphthene | 40.000 | 39.537 | 1.2 | 70 | -0.01 |
| 53 | | 3-Nitroaniline | 40.000 | 40.046 | -0.1 | 67 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 40.000 | 41.349 | -3.4 | 75 | 0.00 |
| 55 | | Dibenzofuran | 40.000 | 39.211 | 2.0 | 69 | -0.01 |
| 56 | P | 4-Nitrophenol | 40.000 | 40.884 | -2.2 | 66 | 0.00 |
| 57 | | 2,4-Dinitrotoluene | 40.000 | 42.874 | -7.2 | 71 | -0.01 |
| 58 | | Fluorene | 40.000 | 39.028 | 2.4 | 69 | -0.01 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 39.424 | 1.4 | 67 | -0.01 |
| 60 | | Diethylphthalate | 40.000 | 39.250 | 1.9 | 69 | -0.01 |
| 61 | | 4-Chlorophenyl-phenylether | 40.000 | 39.204 | 2.0 | 70 | -0.01 |
| 62 | | 4-Nitroaniline | 40.000 | 40.580 | -1.4 | 69 | 0.00 |
| 63 | | Azobenzene | 40.000 | 39.366 | 1.6 | 69 | -0.02 |
| 64 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 69 | -0.01 |
| 65 | | 4,6-Dinitro-2-methylphenol | 40.000 | 42.328 | -5.8 | 70 | -0.01 |
| 66 | c | n-Nitrosodiphenylamine | 40.000 | 39.266 | 1.8 | 68 | -0.02 |
| 67 | | 4-Bromophenyl-phenylether | 40.000 | 40.193 | -0.5 | 69 | -0.01 |
| 68 | | Hexachlorobenzene | 40.000 | 39.829 | 0.4 | 69 | -0.01 |
| 69 | | Atrazine | 40.000 | 40.804 | -2.0 | 70 | -0.01 |
| 70 | C | Pentachlorophenol | 40.000 | 42.815 | -7.0 | 66 | -0.01 |
| 71 | | Phenanthrene | 40.000 | 39.216 | 2.0 | 69 | -0.01 |
| 72 | | Anthracene | 40.000 | 39.681 | 0.8 | 69 | -0.01 |
| 73 | | Carbazole | 40.000 | 40.076 | -0.2 | 70 | 0.00 |
| 74 | | Di-n-butylphthalate | 40.000 | 40.930 | -2.3 | 70 | -0.02 |
| 75 | C | Fluoranthene | 40.000 | 40.102 | -0.3 | 70 | -0.01 |
| 76 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 70 | 0.00 |
| 77 | | Benzidine | 40.000 | 51.455 | -28.6# | 77 | -0.01 |
| 78 | | Pyrene | 40.000 | 39.951 | 0.1 | 71 | -0.01 |
| 79 | S | Terphenyl-d14 | 80.000 | 78.666 | 1.7 | 71 | -0.01 |
| 80 | | Butylbenzylphthalate | 40.000 | 41.437 | -3.6 | 71 | -0.01 |
| 81 | | Benzo(a)anthracene | 40.000 | 39.662 | 0.8 | 69 | 0.00 |
| 82 | | 3,3'-Dichlorobenzidine | 40.000 | 40.150 | -0.4 | 69 | 0.00 |
| 83 | | Chrysene | 40.000 | 39.480 | 1.3 | 70 | -0.01 |
| 84 | | Bis(2-ethylhexyl)phthalate | 40.000 | 39.412 | 1.5 | 68 | -0.02 |
| 85 | c | Di-n-octyl phthalate | 40.000 | 39.798 | 0.5 | 69 | -0.02 |
| 86 | | Indeno(1,2,3-cd)pyrene | 40.000 | 32.181 | 19.5 | 62 | -0.01 |
| 87 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 66 | -0.01 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(b)fluoranthene | 40.000 | 40.769 | -1.9 | 66 | 0.00 |
| 89 | Benzo(k)fluoranthene | 40.000 | 41.408 | -3.5 | 68 | -0.01 |
| 90 C | Benzo(a)pyrene | 40.000 | 39.695 | 0.8 | 65 | -0.01 |
| 91 | Dibenzo(a,h)anthracene | 40.000 | 36.469 | 8.8 | 62 | -0.02 |
| 92 | Benzo(g,h,i)perylene | 40.000 | 36.505 | 8.7 | 62 | -0.01 |

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