# Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_F\Data\BF012318\
Data File : BF102334.D
Acq On : 23 Jan 2018 9:56
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Jan 23 17:14:53 2018
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 23 00:43:41 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 | 102 | 0.00 |
| 2 | | 1,4-Dioxane | 40.000 | 39.759 | 0.6 | 100 | 0.01 |
| 3 | | Pyridine | 40.000 | 40.875 | -2.2 | 100 | 0.01 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 40.948 | -2.4 | 100 | 0.01 |
| 5 | S | 2-Fluorophenol | 80.000 | 80.482 | -0.6 | 101 | 0.00 |
| 6 | | Aniline | 40.000 | 39.689 | 0.8 | 100 | 0.00 |
| 7 | S | Phenol-d6 | 80.000 | 79.991 | 0.0 | 101 | 0.01 |
| 8 | | 2-Chlorophenol | 40.000 | 41.319 | -3.3 | 103 | 0.00 |
| 9 | | Benzaldehyde | 40.000 | 40.436 | -1.1 | 101 | 0.00 |
| 10 | C | Phenol | 40.000 | 38.932 | 2.7 | 98 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 41.045 | -2.6 | 102 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 40.260 | -0.6 | 102 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 40.286 | -0.7 | 102 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 41.116 | -2.8 | 103 | 0.00 |
| 15 | | Benzyl Alcohol | 40.000 | 40.135 | -0.3 | 101 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 40.313 | -0.8 | 100 | 0.00 |
| 17 | | 2-Methylphenol | 40.000 | 40.182 | -0.5 | 99 | 0.00 |
| 18 | | Hexachloroethane | 40.000 | 41.311 | -3.3 | 103 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 40.432 | -1.1 | 102 | 0.00 |
| 20 | | 3+4-Methylphenols | 40.000 | 41.157 | -2.9 | 100 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 104 | 0.00 |
| 22 | | Acetophenone | 40.000 | 39.076 | 2.3 | 103 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 79.245 | 0.9 | 102 | 0.00 |
| 24 | | Nitrobenzene | 40.000 | 39.720 | 0.7 | 101 | 0.00 |
| 25 | | Isophorone | 40.000 | 40.308 | -0.8 | 104 | 0.00 |
| 26 | C | 2-Nitrophenol | 40.000 | 41.235 | -3.1 | 103 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 39.835 | 0.4 | 102 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 40.117 | -0.3 | 103 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 40.183 | -0.5 | 103 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.725 | 0.7 | 103 | 0.00 |
| 31 | | Naphthalene | 40.000 | 39.410 | 1.5 | 103 | 0.00 |
| 32 | | Benzoic acid | 40.000 | 36.282 | 9.3 | 92 | 0.01 |
| 33 | | 4-Chloroaniline | 40.000 | 39.777 | 0.6 | 103 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 40.658 | -1.6 | 103 | 0.00 |
| 35 | | Caprolactam | 40.000 | 39.581 | 1.0 | 100 | 0.01 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 40.038 | -0.1 | 102 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 39.018 | 2.5 | 102 | 0.00 |
| 38 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 103 | 0.00 |
| 39 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 40.087 | -0.2 | 103 | 0.00 |
| 40 | P | Hexachlorocyclopentadiene | 40.000 | 41.209 | -3.0 | 99 | 0.00 |
| 41 | S | 2,4,6-Tribromophenol | 80.000 | 77.826 | 2.7 | 101 | 0.00 |
| 42 | C | 2,4,6-Trichlorophenol | 40.000 | 40.481 | -1.2 | 103 | 0.00 |
| 43 | | 2,4,5-Trichlorophenol | 40.000 | 41.219 | -3.0 | 104 | 0.00 |
| 44 | S | 2-Fluorobiphenyl | 80.000 | 79.942 | 0.1 | 102 | 0.00 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 1,1'-Biphenyl | 40.000 | 39.878 | 0.3 | 104 | 0.00 |
| 46 | | 2-Chloronaphthalene | 40.000 | 39.565 | 1.1 | 102 | 0.00 |
| 47 | | 2-Nitroaniline | 40.000 | 41.261 | -3.2 | 106 | 0.00 |
| 48 | | Acenaphthylene | 40.000 | 39.176 | 2.1 | 101 | 0.00 |
| 49 | | Dimethylphthalate | 40.000 | 39.952 | 0.1 | 104 | 0.00 |
| 50 | | 2,6-Dinitrotoluene | 40.000 | 39.690 | 0.8 | 99 | 0.00 |
| 51 | C | Acenaphthene | 40.000 | 39.357 | 1.6 | 103 | 0.00 |
| 52 | | 3-Nitroaniline | 40.000 | 39.556 | 1.1 | 101 | 0.00 |
| 53 | P | 2,4-Dinitrophenol | 40.000 | 39.427 | 1.4 | 98 | 0.00 |
| 54 | | Dibenzofuran | 40.000 | 40.695 | -1.7 | 102 | 0.00 |
| 55 | P | 4-Nitrophenol | 40.000 | 39.440 | 1.4 | 95 | 0.01 |
| 56 | | 2,4-Dinitrotoluene | 40.000 | 39.556 | 1.1 | 100 | 0.00 |
| 57 | | Fluorene | 40.000 | 41.010 | -2.5 | 105 | 0.00 |
| 58 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 40.623 | -1.6 | 106 | 0.00 |
| 59 | | Diethylphthalate | 40.000 | 39.581 | 1.0 | 103 | 0.00 |
| 60 | | 4-Chlorophenyl-phenylether | 40.000 | 40.951 | -2.4 | 103 | 0.00 |
| 61 | | 4-Nitroaniline | 40.000 | 39.943 | 0.1 | 102 | 0.00 |
| 62 | | Azobenzene | 40.000 | 39.582 | 1.0 | 102 | 0.00 |
| 63 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 100 | 0.00 |
| 64 | | 4,6-Dinitro-2-methylphenol | 40.000 | 43.246 | -8.1 | 103 | 0.01 |
| 65 | c | n-Nitrosodiphenylamine | 40.000 | 40.610 | -1.5 | 103 | 0.00 |
| 66 | | 4-Bromophenyl-phenylether | 40.000 | 41.733 | -4.3 | 104 | 0.00 |
| 67 | | Hexachlorobenzene | 40.000 | 40.821 | -2.1 | 102 | 0.00 |
| 68 | | Atrazine | 40.000 | 40.374 | -0.9 | 102 | 0.00 |
| 69 | C | Pentachlorophenol | 40.000 | 40.077 | -0.2 | 92 | 0.00 |
| 70 | | Phenanthrene | 40.000 | 39.477 | 1.3 | 101 | 0.00 |
| 71 | | Anthracene | 40.000 | 39.670 | 0.8 | 101 | 0.00 |
| 72 | | Carbazole | 40.000 | 39.510 | 1.2 | 101 | 0.00 |
| 73 | | Di-n-butylphthalate | 40.000 | 40.053 | -0.1 | 100 | 0.00 |
| 74 | C | Fluoranthene | 40.000 | 39.255 | 1.9 | 100 | 0.01 |
| 75 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 99 | 0.01 |
| 76 | | Benzidine | 40.000 | 37.328 | 6.7 | 99 | 0.00 |
| 77 | | Pyrene | 40.000 | 39.981 | 0.0 | 99 | 0.01 |
| 78 | S | Terphenyl-d14 | 80.000 | 78.420 | 2.0 | 100 | 0.00 |
| 79 | | Butylbenzylphthalate | 40.000 | 41.103 | -2.8 | 99 | 0.00 |
| 80 | | Benzo(a)anthracene | 40.000 | 39.738 | 0.7 | 100 | 0.01 |
| 81 | | 3,3'-Dichlorobenzidine | 40.000 | 41.164 | -2.9 | 101 | 0.00 |
| 82 | | Chrysene | 40.000 | 39.429 | 1.4 | 99 | 0.00 |
| 83 | | Bis(2-ethylhexyl)phthalate | 40.000 | 41.236 | -3.1 | 100 | 0.00 |
| 84 | c | Di-n-octyl phthalate | 40.000 | 40.754 | -1.9 | 100 | 0.01 |
| 85 | | Indeno(1,2,3-cd)pyrene | 40.000 | 38.369 | 4.1 | 102 | 0.02 |
| 86 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 96 | 0.01 |
| 87 | | Benzo(b)fluoranthene | 40.000 | 42.239 | -5.6 | 99 | 0.01 |

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|---|---|---|---|---|---|---|
| 88 | Benzo(k)fluoranthene | 40.000 | 38.092 | 4.8 | 95 | 0.01 |
| 89 C | Benzo(a)pyrene | 40.000 | 40.341 | -0.9 | 97 | 0.01 |
| 90 | Dibenzo(a,h)anthracene | 40.000 | 39.399 | 1.5 | 100 | 0.02 |
| 91 | Benzo(g,h,i)perylene | 40.000 | 40.081 | -0.2 | 102 | 0.02 |

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