# Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040519\
Data File : BF113621.D
Acq On : 5 Apr 2019 17:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Apr 05 23:41:32 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 04 04:05:31 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 | 104 | 0.00 |
| 2 | | 1,4-Dioxane | 40.000 | 38.118 | 4.7 | 99 | 0.00 |
| 3 | | Pyridine | 40.000 | 40.225 | -0.6 | 102 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 40.416 | -1.0 | 103 | 0.01 |
| 5 | S | 2-Fluorophenol | 80.000 | 80.151 | -0.2 | 102 | 0.00 |
| 6 | | Aniline | 40.000 | 41.035 | -2.6 | 101 | 0.00 |
| 7 | S | Phenol-d6 | 80.000 | 79.923 | 0.1 | 102 | 0.00 |
| 8 | | 2-Chlorophenol | 40.000 | 40.385 | -1.0 | 103 | 0.00 |
| 9 | | Benzaldehyde | 40.000 | 41.588 | -4.0 | 100 | 0.00 |
| 10 | C | Phenol | 40.000 | 40.785 | -2.0 | 104 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 39.625 | 0.9 | 101 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 39.870 | 0.3 | 102 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.867 | 0.3 | 101 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 40.384 | -1.0 | 102 | 0.00 |
| 15 | | Benzyl Alcohol | 40.000 | 42.241 | -5.6 | 103 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 38.835 | 2.9 | 98 | 0.00 |
| 17 | | 2-Methylphenol | 40.000 | 40.140 | -0.4 | 102 | 0.00 |
| 18 | | Hexachloroethane | 40.000 | 39.270 | 1.8 | 100 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 40.137 | -0.3 | 101 | 0.00 |
| 20 | | 3+4-Methylphenols | 40.000 | 42.186 | -5.5 | 105 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 103 | 0.00 |
| 22 | | Acetophenone | 40.000 | 40.555 | -1.4 | 104 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 80.487 | -0.6 | 103 | 0.00 |
| 24 | | Nitrobenzene | 40.000 | 39.956 | 0.1 | 102 | 0.00 |
| 25 | | Isophorone | 40.000 | 40.001 | -0.0 | 104 | 0.00 |
| 26 | C | 2-Nitrophenol | 40.000 | 41.735 | -4.3 | 106 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 39.296 | 1.8 | 101 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 39.755 | 0.6 | 101 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 41.475 | -3.7 | 105 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.771 | 0.6 | 101 | 0.00 |
| 31 | | Naphthalene | 40.000 | 40.077 | -0.2 | 102 | 0.00 |
| 32 | | Benzoic acid | 40.000 | 34.694 | 13.3 | 88 | 0.01 |
| 33 | | 4-Chloroaniline | 40.000 | 42.875 | -7.2 | 106 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 40.163 | -0.4 | 102 | 0.00 |
| 35 | | Caprolactam | 40.000 | 43.691 | -9.2 | 108 | 0.01 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 41.729 | -4.3 | 105 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 40.330 | -0.8 | 102 | 0.00 |
| 38 | | 1-Methylnaphthalene | 40.000 | 40.244 | -0.6 | 102 | 0.00 |
| 39 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 104 | 0.00 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 38.926 | 2.7 | 101 | 0.00 |
| 41 | P | Hexachlorocyclopentadiene | 40.000 | 35.274 | 11.8 | 90 | 0.00 |
| 42 | S | 2,4,6-Tribromophenol | 80.000 | 79.434 | 0.7 | 100 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 40.000 | 39.967 | 0.1 | 104 | 0.00 |
| 44 | | 2,4,5-Trichlorophenol | 40.000 | 40.058 | -0.1 | 104 | 0.00 |

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040519\
Data File : BF113621.D
Acq On : 5 Apr 2019 17:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Apr 05 23:41:32 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 04 04:05:31 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) |
|---|---|---|---|---|---|---|---|
| 45 | S | 2-Fluorobiphenyl | 80.000 | 78.769 | 1.5 | 97 | 0.00 |
| 46 | | 1,1'-Biphenyl | 40.000 | 39.300 | 1.8 | 102 | 0.00 |
| 47 | | 2-Chloronaphthalene | 40.000 | 39.358 | 1.6 | 103 | 0.00 |
| 48 | | 2-Nitroaniline | 40.000 | 41.530 | -3.8 | 108 | 0.00 |
| 49 | | Acenaphthylene | 40.000 | 39.080 | 2.3 | 100 | 0.00 |
| 50 | | Dimethylphthalate | 40.000 | 39.420 | 1.4 | 102 | 0.00 |
| 51 | | 2,6-Dinitrotoluene | 40.000 | 40.665 | -1.7 | 104 | 0.00 |
| 52 | C | Acenaphthene | 40.000 | 39.397 | 1.5 | 102 | 0.00 |
| 53 | | 3-Nitroaniline | 40.000 | 41.634 | -4.1 | 107 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 40.000 | 38.106 | 4.7 | 93 | 0.00 |
| 55 | | Dibenzofuran | 40.000 | 39.374 | 1.6 | 101 | 0.00 |
| 56 | P | 4-Nitrophenol | 40.000 | 37.293 | 6.8 | 95 | 0.00 |
| 57 | | 2,4-Dinitrotoluene | 40.000 | 41.397 | -3.5 | 105 | 0.00 |
| 58 | | Fluorene | 40.000 | 40.117 | -0.3 | 104 | 0.00 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 40.558 | -1.4 | 104 | 0.00 |
| 60 | | Diethylphthalate | 40.000 | 39.452 | 1.4 | 101 | 0.00 |
| 61 | | 4-Chlorophenyl-phenylether | 40.000 | 40.053 | -0.1 | 102 | 0.00 |
| 62 | | 4-Nitroaniline | 40.000 | 43.121 | -7.8 | 110 | 0.00 |
| 63 | | Azobenzene | 40.000 | 39.233 | 1.9 | 101 | -0.01 |
| 64 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 105 | 0.00 |
| 65 | | 4,6-Dinitro-2-methylphenol | 40.000 | 40.700 | -1.8 | 107 | 0.00 |
| 66 | c | n-Nitrosodiphenylamine | 40.000 | 39.319 | 1.7 | 103 | 0.00 |
| 67 | | 4-Bromophenyl-phenylether | 40.000 | 39.305 | 1.7 | 102 | 0.00 |
| 68 | | Hexachlorobenzene | 40.000 | 39.568 | 1.1 | 102 | 0.00 |
| 69 | | Atrazine | 40.000 | 39.565 | 1.1 | 102 | 0.00 |
| 70 | C | Pentachlorophenol | 40.000 | 40.546 | -1.4 | 107 | -0.01 |
| 71 | | Phenanthrene | 40.000 | 39.685 | 0.8 | 102 | -0.01 |
| 72 | | Anthracene | 40.000 | 39.550 | 1.1 | 102 | -0.01 |
| 73 | | Carbazole | 40.000 | 40.839 | -2.1 | 105 | 0.00 |
| 74 | | Di-n-butylphthalate | 40.000 | 39.639 | 0.9 | 101 | -0.02 |
| 75 | C | Fluoranthene | 40.000 | 40.827 | -2.1 | 104 | 0.00 |
| 76 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 107 | -0.01 |
| 77 | | Benzidine | 40.000 | 38.582 | 3.5 | 105 | -0.01 |
| 78 | | Pyrene | 40.000 | 39.104 | 2.2 | 106 | -0.01 |
| 79 | S | Terphenyl-d14 | 80.000 | 76.186 | 4.8 | 103 | -0.01 |
| 80 | | Butylbenzylphthalate | 40.000 | 40.006 | -0.0 | 106 | -0.02 |
| 81 | | Benzo(a)anthracene | 40.000 | 41.094 | -2.7 | 109 | -0.02 |
| 82 | | 3,3'-Dichlorobenzidine | 40.000 | 41.515 | -3.8 | 108 | -0.01 |
| 83 | | Chrysene | 40.000 | 38.905 | 2.7 | 105 | -0.01 |
| 84 | | Bis(2-ethylhexyl)phthalate | 40.000 | 40.993 | -2.5 | 108 | -0.02 |
| 85 | c | Di-n-octyl phthalate | 40.000 | 39.005 | 2.5 | 105 | -0.02 |
| 86 | | Indeno(1,2,3-cd)pyrene | 40.000 | 35.046 | 12.4 | 108 | -0.02 |
| 87 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 105 | -0.02 |

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040519\
Data File : BF113621.D
Acq On : 5 Apr 2019 17:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: Apr 05 23:41:32 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 04 04:05:31 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) |
|---|---|---|---|---|---|---|
| 88 | Benzo(b)fluoranthene | 40.000 | 44.159 | -10.4 | 117 | -0.02 |
| 89 | Benzo(k)fluoranthene | 40.000 | 38.274 | 4.3 | 94 | -0.02 |
| 90 C | Benzo(a)pyrene | 40.000 | 40.689 | -1.7 | 106 | -0.02 |
| 91 | Dibenzo(a,h)anthracene | 40.000 | 39.123 | 2.2 | 107 | -0.02 |
| 92 | Benzo(g,h,i)perylene | 40.000 | 39.080 | 2.3 | 109 | -0.02 |

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