# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021819\
Data File : BM018834.D
Acq On : 18 Feb 2019 12:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_M
**LabSampleId :** SSTDCCC040

Quant Time: Feb 18 15:35:08 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 11 16:02:12 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 | 100 | -0.01 |
| 2 | | 1,4-Dioxane | 40.000 | 37.826 | 5.4 | 94 | 0.00 |
| 3 | | Pyridine | 40.000 | 41.779 | -4.4 | 95 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 42.893 | -7.2 | 102 | 0.00 |
| 5 | S | 2-Fluorophenol | 80.000 | 78.880 | 1.4 | 97 | -0.01 |
| 6 | | Aniline | 40.000 | 41.563 | -3.9 | 101 | -0.01 |
| 7 | S | Phenol-d6 | 80.000 | 84.016 | -5.0 | 102 | 0.00 |
| 8 | | 2-Chlorophenol | 40.000 | 41.103 | -2.8 | 101 | -0.01 |
| 9 | | Benzaldehyde | 40.000 | 44.600 | -11.5 | 108 | -0.01 |
| 10 | C | Phenol | 40.000 | 41.816 | -4.5 | 102 | -0.01 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 41.687 | -4.2 | 102 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 39.917 | 0.2 | 99 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.945 | 0.1 | 100 | -0.01 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 40.242 | -0.6 | 100 | -0.01 |
| 15 | | Benzyl Alcohol | 40.000 | 43.297 | -8.2 | 103 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 42.578 | -6.4 | 107 | -0.01 |
| 17 | | 2-Methylphenol | 40.000 | 42.082 | -5.2 | 102 | 0.00 |
| 18 | | Hexachloroethane | 40.000 | 39.758 | 0.6 | 98 | -0.01 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 42.946 | -7.4 | 103 | -0.01 |
| 20 | | 3+4-Methylphenols | 40.000 | 42.834 | -7.1 | 103 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 103 | -0.01 |
| 22 | | Acetophenone | 40.000 | 40.662 | -1.7 | 105 | -0.01 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 82.281 | -2.9 | 104 | -0.01 |
| 24 | | Nitrobenzene | 40.000 | 40.470 | -1.2 | 103 | -0.01 |
| 25 | | Isophorone | 40.000 | 41.440 | -3.6 | 104 | -0.01 |
| 26 | C | 2-Nitrophenol | 40.000 | 41.610 | -4.0 | 103 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 40.943 | -2.4 | 104 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 40.581 | -1.5 | 103 | -0.01 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 41.449 | -3.6 | 104 | -0.01 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 39.288 | 1.8 | 102 | 0.00 |
| 31 | | Naphthalene | 40.000 | 39.644 | 0.9 | 103 | -0.01 |
| 32 | | Benzoic acid | 40.000 | 34.463 | 13.8 | 89 | 0.00 |
| 33 | | 4-Chloroaniline | 40.000 | 41.680 | -4.2 | 105 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 38.122 | 4.7 | 100 | -0.01 |
| 35 | | Caprolactam | 40.000 | 42.698 | -6.7 | 107 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 42.547 | -6.4 | 107 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 40.925 | -2.3 | 106 | 0.00 |
| 38 | | 1-Methylnaphthalene | 40.000 | 40.863 | -2.2 | 106 | -0.01 |
| 39 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 107 | 0.00 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 38.223 | 4.4 | 102 | -0.01 |
| 41 | P | Hexachlorocyclopentadiene | 40.000 | 33.727 | 15.7 | 89 | 0.00 |
| 42 | S | 2,4,6-Tribromophenol | 80.000 | 83.436 | -4.3 | 109 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 40.000 | 40.273 | -0.7 | 104 | 0.00 |
| 44 | | 2,4,5-Trichlorophenol | 40.000 | 40.561 | -1.4 | 105 | -0.01 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | S | 2-Fluorobiphenyl | 80.000 | 78.386 | 2.0 | 106 | -0.01 |
| 46 | | 1,1'-Biphenyl | 40.000 | 39.778 | 0.6 | 107 | 0.00 |
| 47 | | 2-Chloronaphthalene | 40.000 | 39.720 | 0.7 | 107 | -0.01 |
| 48 | | 2-Nitroaniline | 40.000 | 43.773 | -9.4 | 111 | 0.00 |
| 49 | | Acenaphthylene | 40.000 | 40.874 | -2.2 | 108 | 0.00 |
| 50 | | Dimethylphthalate | 40.000 | 40.270 | -0.7 | 108 | -0.01 |
| 51 | | 2,6-Dinitrotoluene | 40.000 | 41.628 | -4.1 | 107 | 0.00 |
| 52 | C | Acenaphthene | 40.000 | 39.745 | 0.6 | 106 | 0.00 |
| 53 | | 3-Nitroaniline | 40.000 | 40.772 | -1.9 | 108 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 40.000 | 40.052 | -0.1 | 112 | 0.00 |
| 55 | | Dibenzofuran | 40.000 | 40.090 | -0.2 | 108 | -0.01 |
| 56 | P | 4-Nitrophenol | 40.000 | 36.495 | 8.8 | 96 | 0.00 |
| 57 | | 2,4-Dinitrotoluene | 40.000 | 42.904 | -7.3 | 109 | 0.00 |
| 58 | | Fluorene | 40.000 | 40.692 | -1.7 | 109 | 0.00 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 41.040 | -2.6 | 107 | 0.00 |
| 60 | | Diethylphthalate | 40.000 | 40.698 | -1.7 | 109 | 0.00 |
| 61 | | 4-Chlorophenyl-phenylether | 40.000 | 39.698 | 0.8 | 106 | 0.00 |
| 62 | | 4-Nitroaniline | 40.000 | 39.206 | 2.0 | 103 | 0.00 |
| 63 | | Azobenzene | 40.000 | 42.275 | -5.7 | 111 | 0.00 |
| 64 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 110 | 0.00 |
| 65 | | 4,6-Dinitro-2-methylphenol | 40.000 | 36.375 | 9.1 | 99 | 0.00 |
| 66 | c | n-Nitrosodiphenylamine | 40.000 | 40.120 | -0.3 | 109 | 0.00 |
| 67 | | 4-Bromophenyl-phenylether | 40.000 | 39.138 | 2.2 | 107 | -0.01 |
| 68 | | Hexachlorobenzene | 40.000 | 39.020 | 2.4 | 108 | 0.00 |
| 69 | | Atrazine | 40.000 | 38.760 | 3.1 | 106 | 0.00 |
| 70 | C | Pentachlorophenol | 40.000 | 40.946 | -2.4 | 106 | -0.01 |
| 71 | | Phenanthrene | 40.000 | 39.841 | 0.4 | 111 | 0.00 |
| 72 | | Anthracene | 40.000 | 40.763 | -1.9 | 111 | -0.01 |
| 73 | | Carbazole | 40.000 | 40.823 | -2.1 | 111 | 0.00 |
| 74 | | Di-n-butylphthalate | 40.000 | 41.739 | -4.3 | 112 | 0.00 |
| 75 | C | Fluoranthene | 40.000 | 40.371 | -0.9 | 110 | 0.00 |
| 76 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 107 | 0.00 |
| 77 | | Benzidine | 40.000 | 33.898 | 15.3 | 74 | 0.00 |
| 78 | | Pyrene | 40.000 | 41.326 | -3.3 | 110 | 0.00 |
| 79 | S | Terphenyl-d14 | 80.000 | 85.519 | -6.9 | 111 | 0.00 |
| 80 | | Butylbenzylphthalate | 40.000 | 43.327 | -8.3 | 111 | 0.00 |
| 81 | | Benzo(a)anthracene | 40.000 | 40.222 | -0.6 | 107 | 0.00 |
| 82 | | 3,3'-Dichlorobenzidine | 40.000 | 38.913 | 2.7 | 101 | 0.00 |
| 83 | | Chrysene | 40.000 | 40.278 | -0.7 | 107 | 0.00 |
| 84 | | Bis(2-ethylhexyl)phthalate | 40.000 | 43.464 | -8.7 | 112 | 0.00 |
| 85 | c | Di-n-octyl phthalate | 40.000 | 43.528 | -8.8 | 111 | 0.00 |
| 86 | | Indeno(1,2,3-cd)pyrene | 40.000 | 34.188 | 14.5 | 84 | -0.01 |
| 87 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 96 | 0.00 |

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| | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|
| 88 | Benzo(b)fluoranthene | 40.000 | 42.326 | -5.8 | 102 | 0.00 |
| 89 | Benzo(k)fluoranthene | 40.000 | 40.932 | -2.3 | 99 | 0.00 |
| 90 C | Benzo(a)pyrene | 40.000 | 40.587 | -1.5 | 97 | -0.01 |
| 91 | Dibenzo(a,h)anthracene | 40.000 | 36.983 | 7.5 | 84 | -0.01 |
| 92 | Benzo(g,h,i)perylene | 40.000 | 36.218 | 9.5 | 81 | -0.02 |

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