# Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050622\
Data File : BG053469.D
Acq On : 6 May 2022 10:12
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :**
BNA_G
**LabSampleId :**
SSTDCCC040

Quant Time: May 06 23:30:23 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 05 02:13:53 2022
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 | 93 | 0.00 |
| 2 | | 1,4-Dioxane | 40.000 | 39.303 | 1.7 | 92 | 0.00 |
| 3 | | Pyridine | 40.000 | 43.309 | -8.3 | 96 | 0.00 |
| 4 | | n-Nitrosodimethylamine | 40.000 | 41.249 | -3.1 | 93 | 0.00 |
| 5 | S | 2-Fluorophenol | 80.000 | 80.818 | -1.0 | 93 | 0.00 |
| 6 | | Aniline | 40.000 | 43.195 | -8.0 | 97 | 0.00 |
| 7 | S | Phenol-d6 | 80.000 | 83.557 | -4.4 | 95 | 0.00 |
| 8 | | 2-Chlorophenol | 40.000 | 40.236 | -0.6 | 93 | 0.00 |
| 9 | | Benzaldehyde | 40.000 | 39.203 | 2.0 | 96 | 0.00 |
| 10 | C | Phenol | 40.000 | 41.679 | -4.2 | 93 | 0.00 |
| 11 | | bis(2-Chloroethyl)ether | 40.000 | 40.186 | -0.5 | 90 | 0.00 |
| 12 | | 1,3-Dichlorobenzene | 40.000 | 38.824 | 2.9 | 90 | 0.00 |
| 13 | C | 1,4-Dichlorobenzene | 40.000 | 39.258 | 1.9 | 93 | 0.00 |
| 14 | | 1,2-Dichlorobenzene | 40.000 | 41.047 | -2.6 | 95 | 0.00 |
| 15 | | Benzyl Alcohol | 40.000 | 42.662 | -6.7 | 96 | 0.00 |
| 16 | | 2,2'-oxybis(1-Chloropropane | 40.000 | 41.077 | -2.7 | 94 | 0.00 |
| 17 | | 2-Methylphenol | 40.000 | 42.374 | -5.9 | 97 | -0.01 |
| 18 | | Hexachloroethane | 40.000 | 38.019 | 5.0 | 90 | 0.00 |
| 19 | P | n-Nitroso-di-n-propylamine | 40.000 | 42.565 | -6.4 | 97 | 0.00 |
| 20 | | 3+4-Methylphenols | 40.000 | 41.984 | -5.0 | 94 | 0.00 |
| 21 | I | Naphthalene-d8 | 20.000 | 20.000 | 0.0 | 95 | 0.00 |
| 22 | | Acetophenone | 40.000 | 40.139 | -0.3 | 96 | 0.00 |
| 23 | S | Nitrobenzene-d5 | 80.000 | 82.247 | -2.8 | 95 | 0.00 |
| 24 | | Nitrobenzene | 40.000 | 41.350 | -3.4 | 95 | 0.00 |
| 25 | | Isophorone | 40.000 | 42.736 | -6.8 | 98 | 0.00 |
| 26 | C | 2-Nitrophenol | 40.000 | 41.955 | -4.9 | 97 | 0.00 |
| 27 | | 2,4-Dimethylphenol | 40.000 | 40.774 | -1.9 | 94 | 0.00 |
| 28 | | bis(2-Chloroethoxy)methane | 40.000 | 40.853 | -2.1 | 96 | 0.00 |
| 29 | C | 2,4-Dichlorophenol | 40.000 | 42.285 | -5.7 | 95 | 0.00 |
| 30 | | 1,2,4-Trichlorobenzene | 40.000 | 41.360 | -3.4 | 98 | 0.00 |
| 31 | | Naphthalene | 40.000 | 41.005 | -2.5 | 97 | 0.00 |
| 32 | | Benzoic acid | 40.000 | 37.342 | 6.6 | 91 | 0.00 |
| 33 | | 4-Chloroaniline | 40.000 | 43.252 | -8.1 | 98 | 0.00 |
| 34 | C | Hexachlorobutadiene | 40.000 | 39.231 | 1.9 | 92 | 0.00 |
| 35 | | Caprolactam | 40.000 | 46.149 | -15.4 | 107 | 0.00 |
| 36 | C | 4-Chloro-3-methylphenol | 40.000 | 43.904 | -9.8 | 99 | 0.00 |
| 37 | | 2-Methylnaphthalene | 40.000 | 42.223 | -5.6 | 98 | 0.00 |
| 38 | | 1-Methylnaphthalene | 40.000 | 41.645 | -4.1 | 100 | 0.00 |
| 39 | I | Acenaphthene-d10 | 20.000 | 20.000 | 0.0 | 101 | 0.00 |
| 40 | | 1,2,4,5-Tetrachlorobenzene | 40.000 | 39.926 | 0.2 | 99 | 0.00 |
| 41 | P | Hexachlorocyclopentadiene | 40.000 | 36.260 | 9.4 | 94 | 0.00 |
| 42 | S | 2,4,6-Tribromophenol | 80.000 | 84.602 | -5.8 | 102 | 0.00 |
| 43 | C | 2,4,6-Trichlorophenol | 40.000 | 41.681 | -4.2 | 99 | 0.00 |
| 44 | | 2,4,5-Trichlorophenol | 40.000 | 41.698 | -4.2 | 99 | 0.00 |
| 45 | S | 2-Fluorobiphenyl | 80.000 | 78.709 | 1.6 | 99 | 0.00 |
| 46 | | 1,1'-Biphenyl | 40.000 | 40.104 | -0.3 | 101 | 0.00 |
| 47 | | 2-Chloronaphthalene | 40.000 | 39.106 | 2.2 | 99 | 0.00 |

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| | | Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 48 | | 2-Nitroaniline | 40.000 | 42.683 | -6.7 | 103 | 0.00 |
| 49 | | Acenaphthylene | 40.000 | 40.701 | -1.8 | 100 | 0.00 |
| 50 | | Dimethylphthalate | 40.000 | 40.901 | -2.3 | 101 | 0.00 |
| 51 | | 2,6-Dinitrotoluene | 40.000 | 41.349 | -3.4 | 100 | 0.00 |
| 52 | C | Acenaphthene | 40.000 | 40.578 | -1.4 | 100 | 0.00 |
| 53 | | 3-Nitroaniline | 40.000 | 43.349 | -8.4 | 104 | 0.00 |
| 54 | P | 2,4-Dinitrophenol | 40.000 | 40.793 | -2.0 | 105 | 0.00 |
| 55 | | Dibenzofuran | 40.000 | 40.215 | -0.5 | 102 | 0.00 |
| 56 | P | 4-Nitrophenol | 40.000 | 40.560 | -1.4 | 97 | 0.00 |
| 57 | | 2,4-Dinitrotoluene | 40.000 | 41.919 | -4.8 | 101 | 0.00 |
| 58 | | Fluorene | 40.000 | 40.741 | -1.9 | 101 | 0.00 |
| 59 | | 2,3,4,6-Tetrachlorophenol | 40.000 | 40.025 | -0.1 | 97 | 0.00 |
| 60 | | Diethylphthalate | 40.000 | 40.781 | -2.0 | 100 | 0.00 |
| 61 | | 4-Chlorophenyl-phenylether | 40.000 | 41.368 | -3.4 | 103 | 0.00 |
| 62 | | 4-Nitroaniline | 40.000 | 42.277 | -5.7 | 103 | 0.00 |
| 63 | | Azobenzene | 40.000 | 40.967 | -2.4 | 103 | 0.00 |
| 64 | I | Phenanthrene-d10 | 20.000 | 20.000 | 0.0 | 104 | 0.00 |
| 65 | | 4,6-Dinitro-2-methylphenol | 40.000 | 43.620 | -9.0 | 105 | 0.00 |
| 66 | c | n-Nitrosodiphenylamine | 40.000 | 39.947 | 0.1 | 101 | 0.00 |
| 67 | | 4-Bromophenyl-phenylether | 40.000 | 40.101 | -0.3 | 101 | 0.00 |
| 68 | | Hexachlorobenzene | 40.000 | 38.713 | 3.2 | 100 | 0.00 |
| 69 | | Atrazine | 40.000 | 39.069 | 2.3 | 101 | 0.00 |
| 70 | C | Pentachlorophenol | 40.000 | 38.689 | 3.3 | 102 | 0.00 |
| 71 | | Phenanthrene | 40.000 | 39.903 | 0.2 | 103 | 0.00 |
| 72 | | Anthracene | 40.000 | 39.831 | 0.4 | 104 | 0.00 |
| 73 | | Carbazole | 40.000 | 39.807 | 0.5 | 103 | 0.00 |
| 74 | | Di-n-butylphthalate | 40.000 | 39.629 | 0.9 | 101 | 0.00 |
| 75 | C | Fluoranthene | 40.000 | 39.165 | 2.1 | 101 | 0.00 |
| 76 | I | Chrysene-d12 | 20.000 | 20.000 | 0.0 | 99 | 0.00 |
| 77 | | Benzidine | 40.000 | 45.745 | -14.4 | 115 | 0.00 |
| 78 | | Pyrene | 40.000 | 41.474 | -3.7 | 101 | 0.00 |
| 79 | S | Terphenyl-d14 | 80.000 | 81.042 | -1.3 | 100 | 0.00 |
| 80 | | Butylbenzylphthalate | 40.000 | 41.990 | -5.0 | 100 | 0.00 |
| 81 | | Benzo(a)anthracene | 40.000 | 40.746 | -1.9 | 101 | 0.00 |
| 82 | | 3,3'-Dichlorobenzidine | 40.000 | 41.752 | -4.4 | 100 | 0.00 |
| 83 | | Chrysene | 40.000 | 40.985 | -2.5 | 101 | 0.00 |
| 84 | | Bis(2-ethylhexyl)phthalate | 40.000 | 42.235 | -5.6 | 102 | 0.00 |
| 85 | c | Di-n-octyl phthalate | 40.000 | 42.138 | -5.3 | 101 | 0.00 |
| 86 | I | Perylene-d12 | 20.000 | 20.000 | 0.0 | 100 | -0.01 |
| 87 | | Indeno(1,2,3-cd)pyrene | 40.000 | 41.054 | -2.6 | 101 | -0.01 |
| 88 | | Benzo(b)fluoranthene | 40.000 | 40.380 | -1.0 | 98 | -0.01 |
| 89 | | Benzo(k)fluoranthene | 40.000 | 40.648 | -1.6 | 103 | 0.00 |
| 90 | C | Benzo(a)pyrene | 40.000 | 40.606 | -1.5 | 100 | 0.00 |
| 91 | | Dibenzo(a,h)anthracene | 40.000 | 40.548 | -1.4 | 101 | -0.02 |
| 92 | | Benzo(g,h,i)perylene | 40.000 | 40.705 | -1.8 | 101 | -0.01 |

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|---|---|---|---|---|---|

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