# Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081216\
Data File : BM007047.D
Acq On : 12 Aug 2016 17:28
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_M
**LabSampleId :** SSTD02060

Quant Time: Aug 15 15:53:53 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 12 23:03:04 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 1 | I | 1,4-Dichlorobenzene-d4 | 1.000 | 1.000 | 0.0 | 120 | 0.00 |
| 2 | | 1,4-Dioxane | 0.476 | 0.507 | -6.5 | 131 | 0.00 |
| 3 | S | 1,4-Dioxane-d8 | 0.458 | 0.528 | -15.3 | 136 | 0.00 |
| 4 | | Benzaldehyde | 0.974 | 1.104 | -13.3 | 126 | 0.00 |
| 5 | S | Phenol-d5 | 1.620 | 1.617 | 0.2 | 121 | 0.00 |
| 6 | | Phenol | 1.707 | 1.686 | 1.2 | 120 | 0.00 |
| 7 | S | Bis-(2-Chloroethyl)ether-d8 | 0.929 | 0.935 | -0.6 | 120 | 0.00 |
| 8 | | Bis(2-Chloroethyl)ether | 1.259 | 1.275 | -1.3 | 120 | 0.00 |
| 9 | S | 2-Chlorophenol-d4 | 1.290 | 1.331 | -3.2 | 124 | 0.00 |
| 10 | | 2-Chlorophenol | 1.320 | 1.342 | -1.7 | 120 | 0.00 |
| 11 | | 2-Methylphenol | 1.283 | 1.234 | 3.8 | 118 | 0.00 |
| 12 | | 2,2'-oxybis(1-Chloropropane | 1.443 | 1.432 | 0.8 | 121 | 0.00 |
| 13 | S | 4-Methylphenol-d8 | 1.333 | 1.262 | 5.3 | 117 | 0.00 |
| 14 | | Acetophenone | 2.053 | 1.957 | 4.7 | 115 | 0.00 |
| 15 | P | N-Nitroso-di-n-propylamine | 1.068 | 1.006 | 5.8 | 113 | 0.00 |
| 16 | | 4-Methylphenol | 1.401 | 1.324 | 5.5 | 117 | 0.00 |
| 17 | | Hexachloroethane | 0.551 | 0.560 | -1.6 | 122 | 0.00 |
| 18 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 118 | 0.00 |
| 19 | S | Nitrobenzene-d5 | 0.145 | 0.151 | -4.1 | 117 | 0.00 |
| 20 | | Nitrobenzene | 0.374 | 0.380 | -1.6 | 116 | 0.00 |
| 21 | | Isophorone | 0.700 | 0.696 | 0.6 | 114 | 0.00 |
| 22 | S | 2-Nitrophenol-d4 | 0.155 | 0.164 | -5.8 | 120 | 0.00 |
| 23 | C | 2-Nitrophenol | 0.170 | 0.183 | -7.6 | 121 | 0.00 |
| 24 | | 2,4-Dimethylphenol | 0.389 | 0.390 | -0.3 | 114 | 0.00 |
| 25 | | Bis(2-Chloroethoxy)methane | 0.420 | 0.423 | -0.7 | 116 | 0.00 |
| 26 | S | 2,4-Dichlorophenol-d3 | 0.325 | 0.332 | -2.2 | 118 | 0.00 |
| 27 | C | 2,4-Dichlorophenol | 0.324 | 0.328 | -1.2 | 116 | 0.00 |
| 28 | | Naphthalene | 0.991 | 0.974 | 1.7 | 113 | 0.00 |
| 29 | S | 4-Chloroaniline-d4 | 0.387 | 0.388 | -0.3 | 110 | 0.00 |
| 30 | | 4-Chloroaniline | 0.397 | 0.391 | 1.5 | 107 | 0.00 |
| 31 | C | Hexachlorobutadiene | 0.239 | 0.244 | -2.1 | 117 | 0.00 |
| 32 | | Caprolactam | 0.110 | 0.104 | 5.5 | 112 | 0.00 |
| 33 | C | 4-Chloro-3-methylphenol | 0.354 | 0.351 | 0.8 | 114 | 0.00 |
| 34 | | 2-Methylnaphthalene | 0.764 | 0.741 | 3.0 | 111 | 0.00 |
| 35 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 108 | 0.00 |
| 36 | | 1,2,4,5-Tetrachlorobenzene | 0.707 | 0.757 | -7.1 | 111 | 0.00 |
| 37 | | Hexachlorocyclopentadiene | 0.450 | 0.475 | -5.6 | 110 | 0.00 |
| 38 | C | 2,4,6-Trichlorophenol | 0.454 | 0.489 | -7.7 | 112 | 0.00 |
| 39 | | 2,4,5-Trichlorophenol | 0.475 | 0.505 | -6.3 | 110 | 0.00 |
| 40 | | 1,1'-Biphenyl | 1.605 | 1.682 | -4.8 | 110 | 0.00 |
| 41 | | 2-Chloronaphthalene | 1.259 | 1.311 | -4.1 | 109 | 0.00 |
| 42 | | 2-Nitroaniline | 0.361 | 0.382 | -5.8 | 110 | 0.00 |
| 43 | S | Dimethylphthalate-d6 | 1.653 | 1.658 | -0.3 | 106 | 0.00 |
| 44 | | Dimethylphthalate | 1.652 | 1.654 | -0.1 | 106 | 0.00 |

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| | | Compound | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 45 | | 2,6-Dinitrotoluene | 0.321 | 0.328 | -2.2 | 108 | 0.00 |
| 46 | S | Acenaphthylene-d8 | 1.985 | 2.060 | -3.8 | 108 | 0.00 |
| 47 | | Acenaphthylene | 2.043 | 2.083 | -2.0 | 106 | 0.00 |
| 48 | | 3-Nitroaniline | 0.317 | 0.296 | 6.6 | 94 | 0.00 |
| 49 | C | Acenaphthene | 1.327 | 1.324 | 0.2 | 105 | 0.00 |
| 50 | | 2,4-Dinitrophenol | 0.190 | 0.157 | 17.4 | 101 | 0.00 |
| 51 | S | 4-Nitrophenol-d4 | 0.296 | 0.290 | 2.0 | 105 | 0.00 |
| 52 | | 4-Nitrophenol | 0.301 | 0.286 | 5.0 | 102 | 0.00 |
| 53 | | Dibenzofuran | 1.969 | 1.949 | 1.0 | 105 | 0.00 |
| 54 | | 2,4-Dinitrotoluene | 0.474 | 0.478 | -0.8 | 106 | 0.00 |
| 55 | | 2,3,4,6-Tetrachlorophenol | 0.459 | 0.467 | -1.7 | 106 | 0.00 |
| 56 | | Diethylphthalate | 1.742 | 1.674 | 3.9 | 104 | 0.00 |
| 57 | S | Fluorene-d10 | 1.445 | 1.446 | -0.1 | 106 | 0.00 |
| 58 | | Fluorene | 1.588 | 1.569 | 1.2 | 103 | 0.00 |
| 59 | | 4-Chlorophenyl-phenylether | 0.838 | 0.826 | 1.4 | 103 | 0.00 |
| 60 | | 4-Nitroaniline | 0.372 | 0.280 | 24.7# | 78 | 0.00 |
| 61 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 103 | 0.00 |
| 62 | S | 4,6-Dinitro-2-methylphenol- | 0.111 | 0.100 | 9.9 | 99 | 0.00 |
| 63 | | 4,6-Dinitro-2-methylphenol | 0.120 | 0.110 | 8.3 | 97 | 0.00 |
| 64 | | N-Nitrosodiphenylamine | 0.566 | 0.588 | -3.9 | 102 | 0.00 |
| 65 | | 4-Bromophenyl-phenylether | 0.227 | 0.239 | -5.3 | 107 | 0.00 |
| 66 | | Hexachlorobenzene | 0.256 | 0.268 | -4.7 | 105 | 0.00 |
| 67 | | Atrazine | 0.228 | 0.238 | -4.4 | 102 | 0.00 |
| 68 | C | Pentachlorophenol | 0.158 | 0.168 | -6.3 | 111 | 0.00 |
| 69 | | Phenanthrene | 1.076 | 1.084 | -0.7 | 101 | 0.00 |
| 70 | S | Anthracene-d10 | 0.927 | 0.956 | -3.1 | 103 | 0.00 |
| 71 | | Anthracene | 1.104 | 1.118 | -1.3 | 100 | 0.00 |
| 72 | | Carbazole | 0.951 | 0.987 | -3.8 | 99 | 0.00 |
| 73 | | Di-n-butylphthalate | 1.163 | 1.258 | -8.2 | 107 | 0.00 |
| 74 | C | Fluoranthene | 1.305 | 1.393 | -6.7 | 100 | 0.00 |
| 75 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 94 | 0.00 |
| 76 | S | Pyrene-d10 | 0.844 | 0.896 | -6.2 | 99 | 0.00 |
| 77 | | Pyrene | 1.106 | 1.151 | -4.1 | 97 | 0.00 |
| 78 | | Butylbenzylphthalate | 0.430 | 0.487 | -13.3 | 104 | 0.00 |
| 79 | | 3,3'-Dichlorobenzidine | 0.386 | 0.325 | 15.8 | 68 | 0.00 |
| 80 | | Benzo(a)anthracene | 1.182 | 1.211 | -2.5 | 95 | 0.00 |
| 81 | | Bis(2-ethylhexyl)phthalate | 0.636 | 0.713 | -12.1 | 102 | 0.00 |
| 82 | | Chrysene | 1.080 | 1.103 | -2.1 | 94 | 0.00 |
| 83 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 95 | 0.00 |
| 84 | | Di-n-octyl phthalate | 1.032 | 1.140 | -10.5 | 99 | 0.00 |
| 85 | | Benzo(b)fluoranthene | 1.205 | 1.214 | -0.7 | 93 | 0.00 |
| 86 | | Benzo(k)fluoranthene | 1.131 | 1.136 | -0.4 | 94 | 0.00 |
| 87 | S | Benzo(a)pyrene-d12 | 0.915 | 0.932 | -1.9 | 95 | 0.00 |

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|---|---|---|---|---|---|---|
| 88 C | Benzo(a)pyrene | 1.142 | 1.160 | -1.6 | 94 | 0.00 |
| 89 | Indeno(1,2,3-cd)pyrene | 1.400 | 1.411 | -0.8 | 93 | 0.00 |
| 90 | Dibenzo(a,h)anthracene | 1.175 | 1.192 | -1.4 | 93 | 0.00 |
| 91 | Benzo(g,h,i)perylene | 1.183 | 1.189 | -0.5 | 95 | 0.00 |

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