# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102919\
Data File : BE100698.D
Acq On : 29 Oct 2019 19:03
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

**Instrument :** BNA_E
**LabSampleId :** SSTDCCC0.4

Quant Time: Oct 29 19:46:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 23 16:49:33 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 | AvgRF | CCRF | %Dev | Area% | Dev(min) |
|---|---|---|---|---|---|---|---|
| 1 | I | 1,4-Dichlorobenzene-d4 | 1.000 | 1.000 | 0.0 | 71 | 0.00 |
| 2 | | 1,4-Dioxane | 0.714 | 0.772 | -8.1 | 78 | 0.00 |
| 3 | | n-Nitrosodimethylamine | 0.755 | 0.785 | -4.0 | 80 | 0.00 |
| 4 | S | 2-Fluorophenol | 1.235 | 1.322 | -7.0 | 80 | 0.00 |
| 5 | S | Phenol-d6 | 1.742 | 1.984 | -13.9 | 86 | 0.00 |
| 6 | | bis(2-Chloroethyl)ether | 1.430 | 1.660 | -16.1 | 88 | 0.00 |
| 7 | I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 79 | 0.00 |
| 8 | S | Nitrobenzene-d5 | 0.248 | 0.334 | -34.7# | 121 | 0.00 |
| 9 | | Naphthalene | 4.206 | 4.693 | -11.6 | 90 | 0.00 |
| 10 | | Hexachlorobutadiene | 0.183 | 0.203 | -10.9 | 91 | 0.00 |
| 11 | SURR | 2-Methylnaphthalene-d10 | 0.631 | 0.703 | -11.4 | 91 | 0.00 |
| 12 | | 2-Methylnaphthalene | 0.717 | 0.806 | -12.4 | 92 | 0.00 |
| 13 | I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 81 | 0.00 |
| 14 | S | 2,4,6-Tribromophenol | 0.106 | 0.144 | -35.8# | 126 | 0.00 |
| 15 | S | 2-Fluorobiphenyl | 1.457 | 1.558 | -6.9 | 89 | 0.00 |
| 16 | | Acenaphthylene | 1.721 | 1.759 | -2.2 | 86 | -0.01 |
| 17 | | Acenaphthene | 1.112 | 1.236 | -11.2 | 94 | 0.00 |
| 18 | | Fluorene | 1.477 | 1.650 | -11.7 | 94 | 0.00 |
| 19 | I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 88 | -0.01 |
| 20 | | 4-Bromophenyl-phenylether | 0.204 | 0.223 | -9.3 | 100 | 0.00 |
| 21 | | Hexachlorobenzene | 0.194 | 0.209 | -7.7 | 100 | -0.01 |
| 22 | | Pentachlorophenol | 0.062 | 0.055 | 11.3 | 87 | 0.00 |
| 23 | | Phenanthrene | 1.100 | 1.237 | -12.5 | 98 | -0.01 |
| 24 | | Anthracene | 1.021 | 1.101 | -7.8 | 99 | 0.00 |
| 25 | SURR | Fluoranthene-d10 | 5.162 | 5.566 | -7.8 | 99 | 0.00 |
| 26 | | Fluoranthene | 1.436 | 1.568 | -9.2 | 97 | 0.00 |
| 27 | I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 76 | 0.00 |
| 28 | | Pyrene | 1.118 | 1.379 | -23.3 | 94 | 0.00 |
| 29 | S | Terphenyl-d14 | 0.875 | 1.032 | -17.9 | 95 | 0.00 |
| 30 | | Benzo(a)anthracene | 1.305 | 1.454 | -11.4 | 86 | -0.01 |
| 31 | | Chrysene | 1.304 | 1.445 | -10.8 | 85 | 0.00 |
| 32 | | Bis(2-ethylhexyl)phthalate | 2.729 | 2.767 | -1.4 | 77 | -0.01 |
| 33 | | Indeno(1,2,3-cd)pyrene | 1.603 | 1.701 | -6.1 | 84 | -0.01 |
| 34 | I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 72 | -0.01 |
| 35 | | Benzo(b)fluoranthene | 1.166 | 1.226 | -5.1 | 82 | 0.00 |
| 36 | | Benzo(k)fluoranthene | 1.200 | 1.323 | -10.3 | 85 | -0.01 |
| 37 | C | Benzo(a)pyrene | 1.086 | 1.121 | -3.2 | 80 | -0.01 |
| 38 | | Dibenzo(a,h)anthracene | 1.122 | 1.195 | -6.5 | 85 | -0.01 |
| 39 | | Benzo(g,h,i)perylene | 1.129 | 1.218 | -7.9 | 83 | -0.02 |

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

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