# Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE050119\
Data File : BE099454.D
Acq On : 1 May 2019 9:03
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

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

Quant Time: May 02 07:32:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 29 18:36:43 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 | 0.400 | 0.400 | 0.0 | 107 | 0.00 |
| 2 | | 1,4-Dioxane | 0.400 | 0.357 | 10.8 | 111 | 0.00 |
| 3 | | n-Nitrosodimethylamine | 0.400 | 0.365 | 8.8 | 110 | 0.00 |
| 4 | S | 2-Fluorophenol | 0.400 | 0.426 | -6.5 | 119 | 0.00 |
| 5 | S | Phenol-d6 | 0.400 | 0.498 | -24.5 | 136 | 0.00 |
| 6 | | bis(2-Chloroethyl)ether | 0.400 | 0.395 | 1.3 | 108 | 0.00 |
| 7 | I | Naphthalene-d8 | 0.400 | 0.400 | 0.0 | 130 | 0.00 |
| 8 | S | Nitrobenzene-d5 | 0.400 | 0.392 | 2.0 | 127 | 0.00 |
| 9 | | Naphthalene | 0.400 | 0.413 | -3.2 | 132 | 0.00 |
| 10 | | Hexachlorobutadiene | 0.400 | 0.395 | 1.3 | 126 | 0.00 |
| 11 | SURR | 2-Methylnaphthalene-d10 | 0.400 | 0.427 | -6.7 | 140 | 0.00 |
| 12 | | 2-Methylnaphthalene | 0.400 | 0.433 | -8.2 | 141 | 0.00 |
| 13 | I | Acenaphthene-d10 | 0.400 | 0.400 | 0.0 | 140 | 0.00 |
| 14 | S | 2,4,6-Tribromophenol | 0.400 | 0.409 | -2.2 | 148 | 0.00 |
| 15 | S | 2-Fluorobiphenyl | 0.400 | 0.397 | 0.8 | 136 | -0.02 |
| 16 | | Acenaphthylene | 0.400 | 0.406 | -1.5 | 142 | 0.00 |
| 17 | | Acenaphthene | 0.400 | 0.409 | -2.2 | 144 | 0.00 |
| 18 | | Fluorene | 0.400 | 0.400 | 0.0 | 139 | 0.00 |
| 19 | I | Phenanthrene-d10 | 0.400 | 0.400 | 0.0 | 133 | 0.00 |
| 20 | | 4-Bromophenyl-phenylether | 0.400 | 0.403 | -0.8 | 132 | 0.00 |
| 21 | | Hexachlorobenzene | 0.400 | 0.421 | -5.2 | 139 | 0.00 |
| 22 | | Pentachlorophenol | 0.400 | 0.515 | -28.7# | 157 | 0.00 |
| 23 | | Phenanthrene | 0.400 | 0.395 | 1.3 | 129 | 0.00 |
| 24 | | Anthracene | 0.400 | 0.392 | 2.0 | 131 | 0.00 |
| 25 | SURR | Fluoranthene-d10 | 0.400 | 0.362 | 9.5 | 120 | 0.00 |
| 26 | | Fluoranthene | 0.400 | 0.367 | 8.3 | 121 | 0.00 |
| 27 | I | Chrysene-d12 | 0.400 | 0.400 | 0.0 | 107 | 0.00 |
| 28 | | Pyrene | 0.400 | 0.500 | -25.0 | 118 | 0.00 |
| 29 | S | Terphenyl-d14 | 0.400 | 0.484 | -21.0 | 115 | 0.00 |
| 30 | | Benzo(a)anthracene | 0.400 | 0.461 | -15.3 | 109 | -0.01 |
| 31 | | Chrysene | 0.400 | 0.450 | -12.5 | 106 | 0.00 |
| 32 | | Bis(2-ethylhexyl)phthalate | 0.400 | 0.424 | -6.0 | 99 | -0.01 |
| 33 | | Indeno(1,2,3-cd)pyrene | 0.400 | 0.449 | -12.2 | 105 | -0.01 |
| 34 | I | Perylene-d12 | 0.400 | 0.400 | 0.0 | 104 | 0.00 |
| 35 | | Benzo(b)fluoranthene | 0.400 | 0.424 | -6.0 | 109 | 0.00 |
| 36 | | Benzo(k)fluoranthene | 0.400 | 0.411 | -2.7 | 102 | 0.00 |
| 37 | C | Benzo(a)pyrene | 0.400 | 0.419 | -4.7 | 106 | 0.00 |
| 38 | | Dibenzo(a,h)anthracene | 0.400 | 0.411 | -2.7 | 104 | 0.00 |
| 39 | | Benzo(g,h,i)perylene | 0.400 | 0.417 | -4.2 | 105 | -0.01 |

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