Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081016\
Data File : BE092261.D
Acq On : 10 Aug 2016 18:58
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc : confirmation run
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 11 01:28:49 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 11:12:32 2016
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	5.000	5.000	0.0	147	0.00
2		1,4-Dioxane	0.400	0.376	6.0	136	0.00
3		n-Nitrosodimethylamine	0.400	0.483	-20.7	165	-0.01
4	S	2-Fluorophenol	0.400	0.413	-3.2	167	0.00
5	S	Phenol-d6	0.400	0.387	3.3	182	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.428	-7.0	174	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	147	0.00
8	S	Nitrobenzene-d5	0.400	0.376	6.0	163	-0.02
9		Naphthalene	0.400	0.398	0.5	146	-0.02
10		Hexachlorobutadiene	0.400	0.393	1.8	145	-0.02
11		2-Methylnaphthalene	0.400	0.401	-0.3	150	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	140	0.00
13	S	2,4,6-Tribromophenol	0.400	0.514	-28.5#	165	0.00
14	S	2-Fluorobiphenyl	0.400	0.403	-0.8	145	-0.01
15		Acenaphthylene	0.400	0.401	-0.3	147	0.00
16		Acenaphthene	0.400	0.398	0.5	142	0.00
17		Fluorene	0.400	0.401	-0.3	144	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	141	0.00
19		Hexachlorobenzene	0.400	0.381	4.8	135	0.00
20		Pentachlorophenol	0.400	0.451	-12.7	188	0.00
21		Phenanthrene	0.400	0.373	6.8	137	0.00
22		Anthracene	0.400	0.383	4.3	146	0.00
23		Fluoranthene	0.400	0.388	3.0	144	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	141	-0.02
25		Pyrene	0.400	0.412	-3.0	153	0.00
26	S	Terphenyl-d14	0.400	0.414	-3.5	158	0.00
27		Benzo(a)anthracene	0.400	0.417	-4.2	158	-0.02
28		Chrysene	0.400	0.375	6.3	130	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.516	-29.0#	166	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	161	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	148	0.00
32		Benzo(b)fluoranthene	0.400	0.400	0.0	160	0.00
33		Benzo(k)fluoranthene	0.400	0.406	-1.5	149	0.00
34	C	Benzo(a)pyrene	0.400	0.406	-1.5	164	0.00
35		Dibenzo(a,h)anthracene	0.400	0.401	-0.3	165	0.00
36		Benzo(g,h,i)perylene	0.400	0.394	1.5	153	-0.02

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