Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111716\
Data File : BE092557.D
Acq On : 17 Nov 2016 9:16
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 18 13:13:45 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 15:43:18 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	106	-0.02
2		1,4-Dioxane	0.400	0.441	-10.2	130	0.00
3		n-Nitrosodimethylamine	0.400	0.498	-24.5	131	-0.01
4	S	2-Fluorophenol	0.400	0.404	-1.0	107	0.00
5	S	Phenol-d6	0.400	0.451	-12.7	125	0.00
6		bis(2-Chloroethyl)ether	0.400	0.352	12.0	106	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.400	0.407	-1.7	108	-0.02
9		Naphthalene	0.400	0.409	-2.2	108	-0.02
10		Hexachlorobutadiene	0.400	0.405	-1.3	106	0.00
11		2-Methylnaphthalene	0.400	0.398	0.5	111	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	114	0.00
13	S	2,4,6-Tribromophenol	0.400	0.402	-0.5	122	-0.01
14	S	2-Fluorobiphenyl	0.400	0.386	3.5	109	0.00
15		Acenaphthylene	0.400	0.384	4.0	114	0.00
16		Acenaphthene	0.400	0.395	1.3	113	0.00
17		Fluorene	0.400	0.401	-0.3	115	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	106	0.00
19		Hexachlorobenzene	0.400	0.415	-3.7	123	0.00
20		Pentachlorophenol	0.400	0.504	-26.0#	147	0.00
21		Phenanthrene	0.400	0.411	-2.7	113	-0.02
22		Anthracene	0.400	0.393	1.8	111	0.00
23		Fluoranthene	0.400	0.410	-2.5	116	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
25		Pyrene	0.400	0.386	3.5	118	0.00
26	S	Terphenyl-d14	0.400	0.376	6.0	116	-0.02
27		Benzo(a)anthracene	0.400	0.461	-15.3	126	0.00
28		Chrysene	0.400	0.417	-4.2	114	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.533	-33.3#	153	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	109	0.00
31	I	Perylene-d12	5.000	5.000	0.0	106	-0.01
32		Benzo(b)fluoranthene	0.400	0.382	4.5	106	-0.01
33		Benzo(k)fluoranthene	0.400	0.400	0.0	113	-0.01
34	C	Benzo(a)pyrene	0.400	0.390	2.5	112	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.382	4.5	109	-0.02
36		Benzo(g,h,i)perylene	0.400	0.387	3.3	107	-0.02

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