Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111816\
Data File : BE092570.D
Acq On : 18 Nov 2016 17:35
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 21 05:39:35 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	105	-0.02
2		1,4-Dioxane	0.400	0.386	3.5	117	0.00
3		n-Nitrosodimethylamine	0.400	0.444	-11.0	109	0.00
4	S	2-Fluorophenol	0.400	0.412	-3.0	109	0.00
5	S	Phenol-d6	0.400	0.428	-7.0	116	0.00
6		bis(2-Chloroethyl)ether	0.400	0.348	13.0	105	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	107	0.00
8	S	Nitrobenzene-d5	0.400	0.410	-2.5	111	-0.02
9		Naphthalene	0.400	0.398	0.5	107	-0.02
10		Hexachlorobutadiene	0.400	0.406	-1.5	108	0.00
11		2-Methylnaphthalene	0.400	0.396	1.0	112	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	114	0.00
13	S	2,4,6-Tribromophenol	0.400	0.407	-1.7	124	0.00
14	S	2-Fluorobiphenyl	0.400	0.394	1.5	111	0.00
15		Acenaphthylene	0.400	0.384	4.0	114	0.00
16		Acenaphthene	0.400	0.395	1.3	112	0.00
17		Fluorene	0.400	0.398	0.5	114	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
19		Hexachlorobenzene	0.400	0.372	7.0	111	0.00
20		Pentachlorophenol	0.400	0.498	-24.5	146	0.00
21		Phenanthrene	0.400	0.408	-2.0	113	0.00
22		Anthracene	0.400	0.395	1.3	112	0.00
23		Fluoranthene	0.400	0.405	-1.3	115	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	105	0.00
25		Pyrene	0.400	0.400	0.0	113	0.00
26	S	Terphenyl-d14	0.400	0.389	2.8	110	0.00
27		Benzo(a)anthracene	0.400	0.478	-19.5	121	0.00
28		Chrysene	0.400	0.464	-16.0	118	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.461	-15.3	116	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.428	-7.0	105	0.00
31	I	Perylene-d12	5.000	5.000	0.0	103	-0.01
32		Benzo(b)fluoranthene	0.400	0.374	6.5	100	-0.01
33		Benzo(k)fluoranthene	0.400	0.400	0.0	109	-0.01
34	C	Benzo(a)pyrene	0.400	0.386	3.5	108	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.375	6.3	104	-0.02
36		Benzo(g,h,i)perylene	0.400	0.378	5.5	102	0.00

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