Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE112316\
Data File : BE092572.D
Acq On : 23 Nov 2016 9:43
Operator : UM/SJ
Sample : SSTDICCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 23 16:01:19 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	104	-0.02
2		1,4-Dioxane	0.400	0.382	4.5	114	0.00
3		n-Nitrosodimethylamine	0.400	0.435	-8.7	103	0.00
4	S	2-Fluorophenol	0.400	0.414	-3.5	108	0.00
5	S	Phenol-d6	0.400	0.417	-4.2	111	0.00
6		bis(2-Chloroethyl)ether	0.400	0.331	17.3	98	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
8	S	Nitrobenzene-d5	0.400	0.390	2.5	100	0.00
9		Naphthalene	0.400	0.393	1.8	102	-0.02
10		Hexachlorobutadiene	0.400	0.402	-0.5	102	0.00
11		2-Methylnaphthalene	0.400	0.384	4.0	104	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	0.400	0.385	3.8	110	0.00
14	S	2-Fluorobiphenyl	0.400	0.383	4.3	102	0.00
15		Acenaphthylene	0.400	0.377	5.8	105	0.00
16		Acenaphthene	0.400	0.390	2.5	104	0.00
17		Fluorene	0.400	0.389	2.8	105	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
19		Hexachlorobenzene	0.400	0.379	5.3	107	0.00
20		Pentachlorophenol	0.400	0.458	-14.5	125	0.00
21		Phenanthrene	0.400	0.398	0.5	104	-0.02
22		Anthracene	0.400	0.389	2.8	104	0.00
23		Fluoranthene	0.400	0.397	0.8	107	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	103	0.00
25		Pyrene	0.400	0.375	6.3	104	0.00
26	S	Terphenyl-d14	0.400	0.376	6.0	105	0.00
27		Benzo(a)anthracene	0.400	0.449	-12.2	110	0.00
28		Chrysene	0.400	0.422	-5.5	105	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.470	-17.5	117	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	95	0.00
31	I	Perylene-d12	5.000	5.000	0.0	99	-0.01
32		Benzo(b)fluoranthene	0.400	0.397	0.8	103	0.00
33		Benzo(k)fluoranthene	0.400	0.359	10.3	95	0.00
34	C	Benzo(a)pyrene	0.400	0.368	8.0	99	0.00
35		Dibenzo(a,h)anthracene	0.400	0.358	10.5	95	-0.02
36		Benzo(g,h,i)perylene	0.400	0.363	9.3	94	0.00

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