Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110816\
Data File : BE092517.D
Acq On : 8 Nov 2016 20:15
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 09 09:58:25 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE110816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 16:27:22 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	112	-0.02
2		1,4-Dioxane	0.400	0.390	2.5	127	0.00
3		n-Nitrosodimethylamine	0.400	0.369	7.8	93	0.01
4	S	2-Fluorophenol	0.400	0.369	7.8	112	0.00
5	S	Phenol-d6	0.400	0.330	17.5	105	0.00
6		bis(2-Chloroethyl)ether	0.400	0.335	16.3	102	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	113	0.00
8	S	Nitrobenzene-d5	0.400	0.354	11.5	115	0.00
9		Naphthalene	0.400	0.387	3.3	108	0.00
10		Hexachlorobutadiene	0.400	0.394	1.5	108	-0.02
11		2-Methylnaphthalene	0.400	0.377	5.8	111	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	112	0.00
13	S	2,4,6-Tribromophenol	0.400	0.480	-20.0	117	0.00
14	S	2-Fluorobiphenyl	0.400	0.408	-2.0	118	0.00
15		Acenaphthylene	0.400	0.391	2.3	117	-0.02
16		Acenaphthene	0.400	0.392	2.0	114	0.00
17		Fluorene	0.400	0.394	1.5	116	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
19		Hexachlorobenzene	0.400	0.394	1.5	115	0.00
20		Pentachlorophenol	0.400	0.316	21.0	100	0.00
21		Phenanthrene	0.400	0.413	-3.2	116	0.00
22		Anthracene	0.400	0.398	0.5	112	0.00
23		Fluoranthene	0.400	0.387	3.3	108	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	108	0.00
25		Pyrene	0.400	0.380	5.0	111	0.00
26	S	Terphenyl-d14	0.400	0.389	2.8	114	0.00
27		Benzo(a)anthracene	0.400	0.383	4.3	112	0.00
28		Chrysene	0.400	0.348	13.0	99	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.393	1.8	99	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.375	6.3	100	0.00
31	I	Perylene-d12	5.000	5.000	0.0	101	-0.01
32		Benzo(b)fluoranthene	0.400	0.393	1.8	97	0.00
33		Benzo(k)fluoranthene	0.400	0.319	20.3	89	0.00
34	C	Benzo(a)pyrene	0.400	0.371	7.3	97	0.00
35		Dibenzo(a,h)anthracene	0.400	0.379	5.3	98	-0.02
36		Benzo(g,h,i)perylene	0.400	0.387	3.3	101	0.00

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