Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081016\
Data File : BE092245.D
Acq On : 10 Aug 2016 9:04
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 10 18:51:38 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	103	0.00
2		1,4-Dioxane	0.400	0.422	-5.5	105	0.00
3		n-Nitrosodimethylamine	0.400	0.441	-10.2	101	0.00
4	S	2-Fluorophenol	0.400	0.350	12.5	99	0.00
5	S	Phenol-d6	0.400	0.376	6.0	124	0.00
6		bis(2-Chloroethyl)ether	0.400	0.353	11.8	101	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	101	0.00
8	S	Nitrobenzene-d5	0.400	0.394	1.5	118	-0.02
9		Naphthalene	0.400	0.394	1.5	100	0.00
10		Hexachlorobutadiene	0.400	0.414	-3.5	105	0.00
11		2-Methylnaphthalene	0.400	0.392	2.0	101	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	0.400	0.470	-17.5	99	0.00
14	S	2-Fluorobiphenyl	0.400	0.394	1.5	101	0.00
15		Acenaphthylene	0.400	0.386	3.5	100	0.00
16		Acenaphthene	0.400	0.401	-0.3	102	0.00
17		Fluorene	0.400	0.390	2.5	100	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	100	0.00
19		Hexachlorobenzene	0.400	0.405	-1.3	102	0.00
20		Pentachlorophenol	0.400	0.365	8.8	94	0.02
21		Phenanthrene	0.400	0.386	3.5	101	0.00
22		Anthracene	0.400	0.359	10.3	98	0.00
23		Fluoranthene	0.400	0.373	6.8	99	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
25		Pyrene	0.400	0.380	5.0	100	0.00
26	S	Terphenyl-d14	0.400	0.377	5.8	102	0.02
27		Benzo(a)anthracene	0.400	0.386	3.5	104	0.00
28		Chrysene	0.400	0.445	-11.2	110	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.457	-14.2	93	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.358	10.5	99	0.00
31	I	Perylene-d12	5.000	5.000	0.0	101	0.00
32		Benzo(b)fluoranthene	0.400	0.403	-0.8	111	0.00
33		Benzo(k)fluoranthene	0.400	0.354	11.5	89	0.01
34	C	Benzo(a)pyrene	0.400	0.361	9.8	100	0.00
35		Dibenzo(a,h)anthracene	0.400	0.354	11.5	100	0.00
36		Benzo(g,h,i)perylene	0.400	0.370	7.5	99	0.00

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