Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE033116\
Data File : BE091564.D
Acq On : 30 Mar 2016 21:22
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 31 00:55:58 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE033016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 30 18:45:59 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	110	0.00
2		1,4-Dioxane	0.400	0.370	7.5	104	-0.01
3		n-Nitrosodimethylamine	0.400	0.387	3.3	104	-0.03
4	S	2-Fluorophenol	0.400	0.376	6.0	103	-0.01
5	S	Phenol-d6	0.400	0.367	8.3	104	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	109	0.00
7	S	Nitrobenzene-d5	0.400	0.370	7.5	103	-0.02
8		Naphthalene	0.400	0.393	1.8	104	-0.02
9		2-Methylnaphthalene	0.400	0.386	3.5	104	-0.02
10	I	Acenaphthene-d10	5.000	5.000	0.0	104	-0.03
11	S	2,4,6-Tribromophenol	0.400	0.295	26.3#	65	-0.01
12	S	2-Fluorobiphenyl	0.400	0.405	-1.3	103	-0.02
13		Acenaphthylene	0.400	0.405	-1.3	117	-0.01
14		Acenaphthene	0.400	0.395	1.3	103	-0.01
15		Fluorene	0.400	0.388	3.0	101	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	102	-0.01
17		Hexachlorobenzene	0.400	0.395	1.3	101	-0.02
18		Pentachlorophenol	0.400	0.379	5.3	111	-0.02
19		Phenanthrene	0.400	0.394	1.5	99	-0.01
20		Anthracene	0.400	0.373	6.8	97	-0.01
21		Fluoranthene	0.400	0.377	5.8	97	-0.02
22	I	Chrysene-d12	5.000	5.000	0.0	110	-0.02
23		Pyrene	0.400	0.353	11.8	96	0.00
24	S	Terphenyl-d14	0.400	0.357	10.8	96	-0.02
25		Benzo(a)anthracene	0.400	0.376	6.0	101	-0.02
26		Chrysene	0.400	0.366	8.5	98	-0.02
27		Indeno(1,2,3-cd)pyrene	0.400	0.353	11.8	98	-0.06
28	I	Perylene-d12	5.000	5.000	0.0	102	-0.03
29		Benzo(b)fluoranthene	0.400	0.399	0.3	101	-0.02
30		Benzo(k)fluoranthene	0.400	0.364	9.0	95	-0.02
31	C	Benzo(a)pyrene	0.400	0.374	6.5	97	-0.03
32		Dibenzo(a,h)anthracene	0.400	0.378	5.5	97	-0.05
33		Benzo(g,h,i)perylene	0.400	0.378	5.5	97	-0.05

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