Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040116\
Data File : BE091590.D
Acq On : 31 Mar 2016 19:54
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc : LOQ WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 01 01:11:41 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
2		1,4-Dioxane	0.698	0.658	5.7	90	-0.01
3		n-Nitrosodimethylamine	0.908	0.790	13.0	85	-0.03
4	S	2-Fluorophenol	1.309	1.241	5.2	94	-0.01
5	S	Phenol-d6	1.770	1.719	2.9	100	-0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
7	S	Nitrobenzene-d5	0.336	0.309	8.0	100	-0.02
8		Naphthalene	1.033	1.002	3.0	100	-0.02
9		2-Methylnaphthalene	0.675	0.662	1.9	103	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	107	-0.03
11	S	2,4,6-Tribromophenol	0.222	0.153	31.1#	63	-0.01
12	S	2-Fluorobiphenyl	1.421	1.371	3.5	101	-0.02
13		Acenaphthylene	2.016	2.165	-7.4	128	-0.01
14		Acenaphthene	1.237	1.211	2.1	106	-0.01
15		Fluorene	1.591	1.549	2.6	104	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	105	-0.01
17		Hexachlorobenzene	0.184	0.183	0.5	105	-0.02
18		Pentachlorophenol	0.103	0.081	21.4	95	-0.02
19		Phenanthrene	1.171	1.137	2.9	101	-0.01
20		Anthracene	1.080	1.049	2.9	104	-0.01
21		Fluoranthene	1.380	1.339	3.0	102	-0.02
22	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
23		Pyrene	1.020	1.042	-2.2	101	0.00
24	S	Terphenyl-d14	0.634	0.646	-1.9	99	-0.02
25		Benzo(a)anthracene	1.174	1.223	-4.2	102	-0.02
26		Chrysene	1.004	1.040	-3.6	100	0.00
27		Indeno(1,2,3-cd)pyrene	1.157	1.196	-3.4	104	-0.05
28	I	Perylene-d12	1.000	1.000	0.0	102	-0.03
29		Benzo(b)fluoranthene	1.213	1.230	-1.4	102	-0.02
30		Benzo(k)fluoranthene	1.174	1.149	2.1	102	-0.02
31	C	Benzo(a)pyrene	1.108	1.156	-4.3	107	-0.03
32		Dibenzo(a,h)anthracene	1.088	1.092	-0.4	103	-0.05
33		Benzo(g,h,i)perylene	1.111	1.113	-0.2	102	-0.05

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