Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032816\
Data File : BE091551.D
Acq On : 28 Mar 2016 23:31
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 29 01:34:43 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 25 19:00:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	91	0.00
2		1,4-Dioxane	0.669	0.698	-4.3	90	-0.02
3		n-Nitrosodimethylamine	0.581	0.664	-14.3	105	-0.02
4	S	2-Fluorophenol	0.975	1.151	-18.1	115	-0.02
5	S	Phenol-d6	1.330	1.576	-18.5	120	-0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
7	S	Nitrobenzene-d5	0.217	0.250	-15.2	133	-0.01
8		Naphthalene	1.007	1.008	-0.1	100	-0.02
9		2-Methylnaphthalene	0.629	0.657	-4.5	109	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	108	-0.01
11	S	2,4,6-Tribromophenol	0.085	0.119	-40.0#	175#	-0.01
12	S	2-Fluorobiphenyl	1.408	1.367	2.9	106	-0.01
13		Acenaphthylene	1.752	1.910	-9.0	132	0.00
14		Acenaphthene	1.237	1.274	-3.0	111	-0.01
15		Fluorene	1.511	1.503	0.5	113	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
17		Hexachlorobenzene	0.176	0.179	-1.7	113	-0.02
18		Pentachlorophenol	0.047	0.064	-36.2#	176#	-0.02
19		Phenanthrene	1.145	1.138	0.6	105	0.00
20		Anthracene	0.948	0.997	-5.2	123	0.00
21		Fluoranthene	1.071	1.199	-12.0	129	-0.02
22	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
23		Pyrene	1.181	1.196	-1.3	125	0.00
24	S	Terphenyl-d14	0.822	0.820	0.2	120	0.00
25		Benzo(a)anthracene	1.154	1.238	-7.3	132	0.00
26		Chrysene	1.516	1.502	0.9	110	0.00
27		Indeno(1,2,3-cd)pyrene	1.179	1.325	-12.4	142	-0.02
28	I	Perylene-d12	1.000	1.000	0.0	116	-0.01
29		Benzo(b)fluoranthene	1.111	1.152	-3.7	136	-0.01
30		Benzo(k)fluoranthene	1.107	1.162	-5.0	127	0.00
31	C	Benzo(a)pyrene	0.926	1.112	-20.1#	164#	-0.01
32		Dibenzo(a,h)anthracene	0.984	1.036	-5.3	136	-0.02
33		Benzo(g,h,i)perylene	1.029	1.071	-4.1	132	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1