Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040216\
Data File : BE091604.D
Acq On : 1 Apr 2016 23:45
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 02 01:10:46 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	116	0.00
2		1,4-Dioxane	0.698	0.590	15.5	94	-0.02
3		n-Nitrosodimethylamine	0.908	0.768	15.4	96	-0.03
4	S	2-Fluorophenol	1.309	1.202	8.2	106	-0.02
5	S	Phenol-d6	1.770	1.673	5.5	113	-0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
7	S	Nitrobenzene-d5	0.336	0.303	9.8	114	-0.02
8		Naphthalene	1.033	0.996	3.6	115	-0.02
9		2-Methylnaphthalene	0.675	0.642	4.9	116	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	120	-0.03
11	S	2,4,6-Tribromophenol	0.222	0.139	37.4#	64	-0.01
12	S	2-Fluorobiphenyl	1.421	1.369	3.7	113	-0.02
13		Acenaphthylene	2.016	1.940	3.8	129	-0.01
14		Acenaphthene	1.237	1.180	4.6	115	-0.01
15		Fluorene	1.591	1.478	7.1	112	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	115	-0.01
17		Hexachlorobenzene	0.184	0.183	0.5	114	-0.02
18		Pentachlorophenol	0.103	0.071	31.1#	90	-0.02
19		Phenanthrene	1.171	1.130	3.5	110	-0.01
20		Anthracene	1.080	1.035	4.2	113	-0.01
21		Fluoranthene	1.380	1.314	4.8	110	-0.02
22	I	Chrysene-d12	1.000	1.000	0.0	105	-0.02
23		Pyrene	1.020	1.102	-8.0	112	0.00
24	S	Terphenyl-d14	0.634	0.665	-4.9	107	0.00
25		Benzo(a)anthracene	1.174	1.171	0.3	102	-0.02
26		Chrysene	1.004	1.057	-5.3	107	0.00
27		Indeno(1,2,3-cd)pyrene	1.157	1.218	-5.3	111	-0.05
28	I	Perylene-d12	1.000	1.000	0.0	106	-0.03
29		Benzo(b)fluoranthene	1.213	1.190	1.9	103	-0.02
30		Benzo(k)fluoranthene	1.174	1.193	-1.6	110	-0.02
31	C	Benzo(a)pyrene	1.108	1.148	-3.6	112	-0.02
32		Dibenzo(a,h)anthracene	1.088	1.117	-2.7	110	-0.05
33		Benzo(g,h,i)perylene	1.111	1.144	-3.0	109	-0.05

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