Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE040216\
Data File : BE091592.D
Acq On : 1 Apr 2016 16:05
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 01 23:46:32 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	103	-0.02
2		1,4-Dioxane	0.698	0.620	11.2	88	-0.02
3		n-Nitrosodimethylamine	0.908	0.800	11.9	90	-0.03
4	S	2-Fluorophenol	1.309	1.205	7.9	95	-0.02
5	S	Phenol-d6	1.770	1.643	7.2	99	-0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
7	S	Nitrobenzene-d5	0.336	0.294	12.5	97	-0.02
8		Naphthalene	1.033	1.006	2.6	102	-0.02
9		2-Methylnaphthalene	0.675	0.646	4.3	102	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	108	-0.03
11	S	2,4,6-Tribromophenol	0.222	0.140	36.9#	58	-0.01
12	S	2-Fluorobiphenyl	1.421	1.364	4.0	102	-0.02
13		Acenaphthylene	2.016	1.948	3.4	117	-0.01
14		Acenaphthene	1.237	1.196	3.3	106	-0.01
15		Fluorene	1.591	1.509	5.2	103	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	110	-0.01
17		Hexachlorobenzene	0.184	0.179	2.7	107	-0.02
18		Pentachlorophenol	0.103	0.071	31.1#	86	-0.02
19		Phenanthrene	1.171	1.106	5.6	103	-0.01
20		Anthracene	1.080	0.971	10.1	101	-0.01
21		Fluoranthene	1.380	1.212	12.2	97	-0.02
22	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
23		Pyrene	1.020	1.002	1.8	97	0.00
24	S	Terphenyl-d14	0.634	0.603	4.9	93	0.00
25		Benzo(a)anthracene	1.174	1.115	5.0	93	-0.02
26		Chrysene	1.004	1.008	-0.4	98	0.00
27		Indeno(1,2,3-cd)pyrene	1.157	1.165	-0.7	102	-0.05
28	I	Perylene-d12	1.000	1.000	0.0	102	-0.02
29		Benzo(b)fluoranthene	1.213	1.140	6.0	95	-0.02
30		Benzo(k)fluoranthene	1.174	1.143	2.6	102	-0.02
31	C	Benzo(a)pyrene	1.108	1.069	3.5	100	-0.03
32		Dibenzo(a,h)anthracene	1.088	1.056	2.9	100	-0.05
33		Benzo(g,h,i)perylene	1.111	1.089	2.0	100	-0.05

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