Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040116\
Data File : BE091578.D
Acq On : 31 Mar 2016 10:15
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 01 00:49:40 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	105	0.00
2		1,4-Dioxane	0.698	0.675	3.3	97	-0.01
3		n-Nitrosodimethylamine	0.908	0.807	11.1	92	-0.03
4	S	2-Fluorophenol	1.309	1.248	4.7	99	-0.01
5	S	Phenol-d6	1.770	1.673	5.5	102	-0.02
6	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
7	S	Nitrobenzene-d5	0.336	0.305	9.2	98	-0.02
8		Naphthalene	1.033	1.042	-0.9	103	-0.02
9		2-Methylnaphthalene	0.675	0.658	2.5	102	-0.01
10	I	Acenaphthene-d10	1.000	1.000	0.0	105	-0.03
11	S	2,4,6-Tribromophenol	0.222	0.162	27.0#	66	-0.01
12	S	2-Fluorobiphenyl	1.421	1.407	1.0	102	-0.02
13		Acenaphthylene	2.016	1.999	0.8	117	-0.01
14		Acenaphthene	1.237	1.218	1.5	105	-0.01
15		Fluorene	1.591	1.548	2.7	103	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.01
17		Hexachlorobenzene	0.184	0.180	2.2	104	-0.02
18		Pentachlorophenol	0.103	0.096	6.8	114	-0.02
19		Phenanthrene	1.171	1.147	2.0	104	-0.01
20		Anthracene	1.080	1.006	6.9	102	-0.01
21		Fluoranthene	1.380	1.335	3.3	104	-0.02
22	I	Chrysene-d12	1.000	1.000	0.0	114	-0.02
23		Pyrene	1.020	0.939	7.9	104	0.00
24	S	Terphenyl-d14	0.634	0.595	6.2	104	-0.02
25		Benzo(a)anthracene	1.174	1.080	8.0	103	-0.02
26		Chrysene	1.004	0.934	7.0	103	-0.02
27		Indeno(1,2,3-cd)pyrene	1.157	1.043	9.9	103	-0.06
28	I	Perylene-d12	1.000	1.000	0.0	107	-0.03
29		Benzo(b)fluoranthene	1.213	1.206	0.6	106	-0.02
30		Benzo(k)fluoranthene	1.174	1.109	5.5	103	-0.02
31	C	Benzo(a)pyrene	1.108	1.056	4.7	104	-0.03
32		Dibenzo(a,h)anthracene	1.088	1.046	3.9	104	-0.05
33		Benzo(g,h,i)perylene	1.111	1.076	3.2	104	-0.05

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