Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089953.D
Acq On : 26 Jun 2015 21:25
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 27 01:01:53 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 2015
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	86	0.00
2		1,4-Dioxane	0.675	0.617	8.6	93	0.00
3		n-Nitrosodimethylamine	0.899	0.816	9.2	86	0.00
4	S	2-Fluorophenol	0.903	0.966	-7.0	98	0.00
5	S	Phenol-d6	1.383	2.219	-60.4#	149	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
7	S	Nitrobenzene-d5	0.372	0.397	-6.7	105	0.00
8		Nitrobenzene	0.396	0.416	-5.1	104	0.00
9		Naphthalene	1.112	1.220	-9.7	104	0.00
10		Hexachlorobutadiene	0.176	0.175	0.6	94	0.00
11		2-Methylnaphthalene	0.761	0.943	-23.9	120	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
13	S	2,4,6-Tribromophenol	0.064	0.115	-79.7#	189#	0.00
14	S	2-Fluorobiphenyl	1.457	1.891	-29.8#	127	0.00
15		Acenaphthylene	8.767	11.263	-28.5#	121	0.00
16		Acenaphthene	1.288	1.691	-31.3#	122	0.00
17		Fluorene	1.750	2.277	-30.1#	124	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
19		4-Bromophenyl-phenylether	0.204	0.256	-25.5#	125	-0.02
20		Hexachlorobenzene	0.894	1.138	-27.3#	123	0.00
21		Pentachlorophenol	0.033	0.062	-87.9#	211#	-0.02
22		Phenanthrene	1.254	1.580	-26.0#	123	-0.02
23		Anthracene	1.164	1.498	-28.7#	124	0.00
24		Fluoranthene	1.361	1.780	-30.8#	126	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
26		Pyrene	1.266	1.603	-26.6#	126	0.00
27	S	Terphenyl-d14	0.863	1.089	-26.2#	131	0.00
28		Benzo(a)anthracene	1.195	1.593	-33.3#	138	0.00
29		Chrysene	1.306	1.676	-28.3#	135	0.00
30		Bis(2-ethylhexyl)phthalate	0.102	0.156	-52.9#	172#	0.00
31		Indeno(1,2,3-cd)pyrene	0.960	1.379	-43.6#	144	0.00
32	I	Perylene-d12	1.000	1.000	0.0	95	0.00
33		Benzo(b)fluoranthene	1.058	1.468	-38.8#	139	0.00
34		Benzo(k)fluoranthene	1.199	1.677	-39.9#	142	0.00
35	C	Benzo(a)pyrene	0.940	1.322	-40.6#	144	0.00
36		Dibenzo(a,h)anthracene	0.875	1.260	-44.0#	148	0.00
37		Benzo(g,h,i)perylene	0.941	1.329	-41.2#	144	0.00

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