Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE033017\
Data File : BE092876.D
Acq On : 30 Mar 2017 15:51
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 31 01:34:13 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 29 17:47:56 2017
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	0.668	0.707	-5.8	108	-0.01
3		n-Nitrosodimethylamine	0.691	0.703	-1.7	115	0.00
4	S	2-Fluorophenol	1.164	1.014	12.9	93	0.00
5	S	Phenol-d6	1.708	1.537	10.0	101	0.00
6		bis(2-Chloroethyl)ether	1.625	1.681	-3.4	110	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.287	0.284	1.0	116	0.00
9		Naphthalene	1.095	1.129	-3.1	109	0.00
10		Hexachlorobutadiene	0.147	0.145	1.4	105	0.00
11		2-Methylnaphthalene	0.695	0.685	1.4	106	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	0.078	0.063	19.2	97	0.00
14	S	2-Fluorobiphenyl	1.648	1.675	-1.6	108	-0.01
15		Acenaphthylene	7.666	6.538	14.7	102	0.00
16		Acenaphthene	1.367	1.302	4.8	106	0.00
17		Fluorene	1.629	1.565	3.9	104	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.02
19		Hexachlorobenzene	0.185	0.197	-6.5	116	0.00
20		Pentachlorophenol	0.048	0.036	25.0#	95	0.00
21		Phenanthrene	1.217	1.415	-16.3	121	0.00
22		Anthracene	1.106	1.139	-3.0	119	0.00
23		Fluoranthene	5.597	4.877	12.9	101	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	102	-0.02
25		Pyrene	4.972	4.529	8.9	99	0.00
26	S	Terphenyl-d14	0.864	0.807	6.6	99	-0.02
27		Benzo(a)anthracene	1.185	1.053	11.1	99	0.00
28		Chrysene	1.250	1.297	-3.8	107	0.00
29		Bis(2-ethylhexyl)phthalate	0.362	0.224	38.1#	80	-0.02
30		Indeno(1,2,3-cd)pyrene	4.702	3.853	18.1	89	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	101	0.00
32		Benzo(b)fluoranthene	1.331	1.099	17.4	87	-0.01
33		Benzo(k)fluoranthene	1.192	1.097	8.0	99	-0.01
34	C	Benzo(a)pyrene	1.100	0.869	21.0#	87	0.00
35		Dibenzo(a,h)anthracene	1.170	0.994	15.0	91	-0.02
36		Benzo(g,h,i)perylene	4.851	4.089	15.7	88	0.00

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