Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be102116\
Data File : BE092464.D
Acq On : 21 Oct 2016 11:51
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 21 17:26:19 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	170#	-0.02
2		1,4-Dioxane	0.635	0.604	4.9	160#	-0.01
3		n-Nitrosodimethylamine	0.588	0.614	-4.4	196#	-0.01
4	S	2-Fluorophenol	0.999	0.858	14.1	156#	0.00
5	S	Phenol-d6	1.388	1.156	16.7	163#	0.00
6		bis(2-Chloroethyl)ether	1.222	1.093	10.6	195#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	164#	0.00
8	S	Nitrobenzene-d5	0.299	0.275	8.0	178#	0.00
9		Naphthalene	1.082	1.073	0.8	163#	-0.02
10		Hexachlorobutadiene	0.170	0.164	3.5	159#	-0.02
11		2-Methylnaphthalene	0.716	0.696	2.8	162#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	153#	0.00
13	S	2,4,6-Tribromophenol	0.144	0.106	26.4#	132	0.00
14	S	2-Fluorobiphenyl	1.184	1.250	-5.6	164#	0.00
15		Acenaphthylene	7.240	7.080	2.2	153#	0.00
16		Acenaphthene	1.134	1.157	-2.0	151#	0.00
17		Fluorene	1.438	1.398	2.8	151#	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	136	0.02
19		Hexachlorobenzene	0.214	0.235	-9.8	159#	0.00
20		Pentachlorophenol	0.050	0.045	10.0	145	0.00
21		Phenanthrene	1.149	1.322	-15.1	165#	0.00
22		Anthracene	1.105	1.206	-9.1	163#	0.02
23		Fluoranthene	5.460	5.490	-0.5	148	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	142	0.00
25		Pyrene	4.333	4.834	-11.6	165#	0.00
26	S	Terphenyl-d14	0.620	0.634	-2.3	147	0.02
27		Benzo(a)anthracene	4.873	5.165	-6.0	154#	0.00
28		Chrysene	4.462	4.937	-10.6	147	0.00
29		Bis(2-ethylhexyl)phthalate	0.420	0.399	5.0	133	0.00
30		Indeno(1,2,3-cd)pyrene	4.734	4.358	7.9	133	0.04
31	I	Perylene-d12	1.000	1.000	0.0	136	0.02
32		Benzo(b)fluoranthene	1.211	1.234	-1.9	139	0.02
33		Benzo(k)fluoranthene	1.267	1.275	-0.6	147	0.02
34	C	Benzo(a)pyrene	1.187	1.180	0.6	143	0.02
35		Dibenzo(a,h)anthracene	1.088	1.036	4.8	135	0.04
36		Benzo(g,h,i)perylene	4.673	4.428	5.2	133	0.04

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