Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE030717\
Data File : BE092805.D
Acq On : 7 Mar 2017 11:06
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 08 04:20:09 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE030617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 06 15:46:01 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	80	-0.02
2		1,4-Dioxane	0.743	0.818	-10.1	102	-0.01
3		n-Nitrosodimethylamine	0.839	0.682	18.7	75	-0.02
4	S	2-Fluorophenol	1.058	0.845	20.1	77	-0.02
5	S	Phenol-d6	1.260	0.946	24.9	74	0.00
6		bis(2-Chloroethyl)ether	1.374	1.573	-14.5	104	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	84	-0.02
8	S	Nitrobenzene-d5	0.306	0.249	18.6	83	-0.02
9		Naphthalene	1.070	1.111	-3.8	95	0.00
10		Hexachlorobutadiene	0.154	0.161	-4.5	92	-0.02
11		2-Methylnaphthalene	0.692	0.660	4.6	90	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.02
13	S	2,4,6-Tribromophenol	0.132	0.101	23.5	75	-0.02
14	S	2-Fluorobiphenyl	1.179	1.128	4.3	93	-0.01
15		Acenaphthylene	7.684	7.338	4.5	96	0.00
16		Acenaphthene	1.076	1.037	3.6	93	-0.02
17		Fluorene	1.463	1.407	3.8	93	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
19		Hexachlorobenzene	0.205	0.202	1.5	101	-0.02
20		Pentachlorophenol	0.069	0.060	13.0	95	-0.11
21		Phenanthrene	1.199	1.153	3.8	95	0.00
22		Anthracene	1.209	1.152	4.7	97	0.00
23		Fluoranthene	5.839	5.644	3.3	96	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	90	-0.02
25		Pyrene	4.529	4.412	2.6	98	0.00
26	S	Terphenyl-d14	0.603	0.590	2.2	96	-0.02
27		Benzo(a)anthracene	5.028	4.519	10.1	100	-0.02
28		Chrysene	4.585	4.953	-8.0	94	-0.02
29		Bis(2-ethylhexyl)phthalate	0.549	0.320	41.7#	75	-0.02
30		Indeno(1,2,3-cd)pyrene	4.965	4.614	7.1	99	-0.03
31	I	Perylene-d12	1.000	1.000	0.0	89	-0.02
32		Benzo(b)fluoranthene	1.332	1.278	4.1	102	-0.02
33		Benzo(k)fluoranthene	1.340	1.410	-5.2	97	-0.02
34	C	Benzo(a)pyrene	1.211	1.194	1.4	101	-0.02
35		Dibenzo(a,h)anthracene	1.157	1.080	6.7	98	-0.02
36		Benzo(g,h,i)perylene	4.951	4.780	3.5	96	-0.03

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