Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE022017\
Data File : BE092753.D
Acq On : 20 Feb 2017 22:52
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 100 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 21 00:54:24 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE021517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 15 17:05:20 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	185#	0.00
2		1,4-Dioxane	0.710	0.653	8.0	172#	0.00
3		n-Nitrosodimethylamine	0.740	0.697	5.8	199#	-0.01
4	S	2-Fluorophenol	0.973	1.175	-20.8	232#	-0.02
5	S	Phenol-d6	1.194	1.755	-47.0#	281#	-0.02
6		bis(2-Chloroethyl)ether	1.321	1.433	-8.5	218#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	195#	-0.02
8	S	Nitrobenzene-d5	0.279	0.324	-16.1	234#	-0.02
9		Naphthalene	1.082	1.105	-2.1	197#	0.00
10		Hexachlorobutadiene	0.159	0.172	-8.2	202#	0.00
11		2-Methylnaphthalene	0.680	0.681	-0.1	200#	-0.02
12	I	Acenaphthene-d10	1.000	1.000	0.0	201#	0.00
13	S	2,4,6-Tribromophenol	0.106	0.137	-29.2#	266#	-0.02
14	S	2-Fluorobiphenyl	1.219	1.298	-6.5	211#	-0.02
15		Acenaphthylene	7.544	7.486	0.8	201#	0.00
16		Acenaphthene	1.096	1.126	-2.7	205#	-0.02
17		Fluorene	1.407	1.395	0.9	201#	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	202#	0.00
19		Hexachlorobenzene	0.193	0.185	4.1	176#	0.00
20		Pentachlorophenol	0.042	0.044	-4.8	246#	-0.09
21		Phenanthrene	1.188	1.092	8.1	176#	0.00
22		Anthracene	1.094	1.091	0.3	206#	0.00
23		Fluoranthene	5.096	4.817	5.5	190#	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	154#	0.00
25		Pyrene	4.761	5.794	-21.7	193#	0.00
26	S	Terphenyl-d14	0.638	0.825	-29.3#	205#	-0.02
27		Benzo(a)anthracene	4.849	5.407	-11.5	175#	0.00
28		Chrysene	5.752	5.965	-3.7	159#	0.00
29		Bis(2-ethylhexyl)phthalate	0.674	1.089	-61.6#	288#	0.00
30		Indeno(1,2,3-cd)pyrene	6.486	6.192	4.5	146	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	158#	-0.01
32		Benzo(b)fluoranthene	1.414	1.290	8.8	148	-0.01
33		Benzo(k)fluoranthene	1.495	1.380	7.7	160#	-0.01
34	C	Benzo(a)pyrene	1.285	1.253	2.5	163#	-0.01
35		Dibenzo(a,h)anthracene	1.355	1.207	10.9	143	-0.02
36		Benzo(g,h,i)perylene	5.728	4.983	13.0	138	-0.02

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