Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE020117\
Data File : BE092707.D
Acq On : 1 Feb 2017 9:58
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 01 12:32:28 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE011617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 17 10:14:01 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	150#	-0.02
2		1,4-Dioxane	0.775	0.740	4.5	136	0.00
3		n-Nitrosodimethylamine	0.902	0.882	2.2	168#	-0.01
4	S	2-Fluorophenol	1.116	1.300	-16.5	206#	-0.01
5	S	Phenol-d6	1.368	1.610	-17.7	215#	-0.02
6		bis(2-Chloroethyl)ether	1.453	1.432	1.4	176#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	145	0.00
8	S	Nitrobenzene-d5	0.325	0.339	-4.3	182#	-0.02
9		Naphthalene	1.044	1.055	-1.1	158#	-0.02
10		Hexachlorobutadiene	0.154	0.164	-6.5	162#	-0.02
11		2-Methylnaphthalene	0.648	0.646	0.3	161#	-0.02
12	I	Acenaphthene-d10	1.000	1.000	0.0	141	0.00
13	S	2,4,6-Tribromophenol	0.126	0.146	-15.9	204#	-0.01
14	S	2-Fluorobiphenyl	1.230	1.282	-4.2	158#	-0.02
15		Acenaphthylene	6.890	7.059	-2.5	163#	-0.02
16		Acenaphthene	1.105	1.125	-1.8	151#	-0.02
17		Fluorene	1.387	1.373	1.0	155#	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	135	0.00
19		Hexachlorobenzene	0.189	0.188	0.5	128	0.00
20		Pentachlorophenol	0.051	0.076	-49.0#	219#	-0.02
21		Phenanthrene	1.104	1.070	3.1	135	0.00
22		Anthracene	1.018	1.080	-6.1	160#	-0.02
23		Fluoranthene	4.895	5.145	-5.1	156#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	125	0.00
25		Pyrene	4.929	5.120	-3.9	156#	-0.02
26	S	Terphenyl-d14	0.707	0.744	-5.2	154#	0.00
27		Benzo(a)anthracene	4.926	5.640	-14.5	167#	0.00
28		Chrysene	5.542	5.201	6.2	127	0.00
29		Bis(2-ethylhexyl)phthalate	0.913	1.541	-68.8#	296#	0.00
30		Indeno(1,2,3-cd)pyrene	5.987	6.840	-14.2	164#	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	136	0.00
32		Benzo(b)fluoranthene	1.210	1.242	-2.6	152#	-0.01
33		Benzo(k)fluoranthene	1.215	1.270	-4.5	165#	-0.01
34	C	Benzo(a)pyrene	1.136	1.195	-5.2	165#	-0.01
35		Dibenzo(a,h)anthracene	1.112	1.188	-6.8	164#	-0.03
36		Benzo(g,h,i)perylene	4.699	4.961	-5.6	159#	-0.02

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