Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE012517\
Data File : BE092684.D
Acq On : 25 Jan 2017 14:19
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 25 17:25:23 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	154#	0.00
2		1,4-Dioxane	0.775	0.684	11.7	129	0.00
3		n-Nitrosodimethylamine	0.902	0.580	35.7#	113	0.02
4	S	2-Fluorophenol	1.116	0.933	16.4	151#	0.00
5	S	Phenol-d6	1.368	1.397	-2.1	191#	-0.02
6		bis(2-Chloroethyl)ether	1.453	1.294	10.9	163#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	179#	0.00
8	S	Nitrobenzene-d5	0.325	0.275	15.4	183#	-0.02
9		Naphthalene	1.044	1.017	2.6	189#	-0.02
10		Hexachlorobutadiene	0.154	0.151	1.9	185#	-0.02
11		2-Methylnaphthalene	0.648	0.642	0.9	198#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	187#	0.00
13	S	2,4,6-Tribromophenol	0.126	0.109	13.5	203#	0.04
14	S	2-Fluorobiphenyl	1.230	1.175	4.5	192#	-0.01
15		Acenaphthylene	6.890	6.863	0.4	210#	-0.02
16		Acenaphthene	1.105	1.114	-0.8	198#	0.00
17		Fluorene	1.387	1.334	3.8	200#	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	182#	0.00
19		Hexachlorobenzene	0.189	0.182	3.7	168#	0.00
20		Pentachlorophenol	0.051	0.000	100.0#	0#	0.02
21		Phenanthrene	1.104	1.136	-2.9	194#	0.00
22		Anthracene	1.018	1.041	-2.3	209#	-0.02
23		Fluoranthene	4.895	4.748	3.0	195#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	154#	0.00
25		Pyrene	4.929	5.115	-3.8	191#	-0.02
26	S	Terphenyl-d14	0.707	0.709	-0.3	180#	0.00
27		Benzo(a)anthracene	4.926	4.555	7.5	165#	0.00
28		Chrysene	5.542	5.396	2.6	161#	0.00
29		Bis(2-ethylhexyl)phthalate	0.913	0.838	8.2	197#	0.00
30		Indeno(1,2,3-cd)pyrene	5.987	5.471	8.6	161#	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	142	0.00
32		Benzo(b)fluoranthene	1.210	1.158	4.3	148	-0.01
33		Benzo(k)fluoranthene	1.215	1.310	-7.8	178#	-0.01
34	C	Benzo(a)pyrene	1.136	1.175	-3.4	170#	0.00
35		Dibenzo(a,h)anthracene	1.112	1.106	0.5	159#	-0.02
36		Benzo(g,h,i)perylene	4.699	4.578	2.6	153#	-0.02

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