Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013117\
Data File : BE092705.D
Acq On : 1 Feb 2017 2:27
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 01 03:24:19 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	195#	0.00
2		1,4-Dioxane	0.775	0.665	14.2	159#	0.01
3		n-Nitrosodimethylamine	0.902	0.625	30.7#	154#	0.01
4	S	2-Fluorophenol	1.116	1.131	-1.3	232#	0.00
5	S	Phenol-d6	1.368	1.583	-15.7	274#	-0.02
6		bis(2-Chloroethyl)ether	1.453	1.242	14.5	199#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	206#	0.00
8	S	Nitrobenzene-d5	0.325	0.301	7.4	231#	-0.02
9		Naphthalene	1.044	1.027	1.6	220#	-0.02
10		Hexachlorobutadiene	0.154	0.161	-4.5	226#	-0.02
11		2-Methylnaphthalene	0.648	0.651	-0.5	232#	-0.02
12	I	Acenaphthene-d10	1.000	1.000	0.0	214#	0.00
13	S	2,4,6-Tribromophenol	0.126	0.132	-4.8	282#	-0.01
14	S	2-Fluorobiphenyl	1.230	1.188	3.4	223#	-0.02
15		Acenaphthylene	6.890	7.051	-2.3	247#	-0.02
16		Acenaphthene	1.105	1.110	-0.5	227#	0.00
17		Fluorene	1.387	1.348	2.8	232#	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	208#	0.00
19		Hexachlorobenzene	0.189	0.183	3.2	193#	0.00
20		Pentachlorophenol	0.051	0.052	-2.0	232#	0.00
21		Phenanthrene	1.104	1.099	0.5	214#	0.00
22		Anthracene	1.018	1.062	-4.3	243#	-0.02
23		Fluoranthene	4.895	4.976	-1.7	233#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	171#	0.00
25		Pyrene	4.929	5.636	-14.3	234#	-0.02
26	S	Terphenyl-d14	0.707	0.777	-9.9	219#	0.00
27		Benzo(a)anthracene	4.926	5.290	-7.4	213#	0.00
28		Chrysene	5.542	5.836	-5.3	193#	0.00
29		Bis(2-ethylhexyl)phthalate	0.913	1.360	-49.0#	356#	0.00
30		Indeno(1,2,3-cd)pyrene	5.987	5.986	0.0	195#	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	167#	0.00
32		Benzo(b)fluoranthene	1.210	1.272	-5.1	191#	-0.01
33		Benzo(k)fluoranthene	1.215	1.290	-6.2	207#	-0.01
34	C	Benzo(a)pyrene	1.136	1.208	-6.3	206#	-0.01
35		Dibenzo(a,h)anthracene	1.112	1.158	-4.1	197#	-0.03
36		Benzo(g,h,i)perylene	4.699	4.776	-1.6	188#	-0.02

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