Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090116\
Data File : BE092339.D
Acq On : 1 Sep 2016 21:27
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Sep 02 11:13:11 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 11:02:43 2016
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	95	0.00
2		1,4-Dioxane	0.717	0.805	-12.3	111	0.00
3		n-Nitrosodimethylamine	0.744	0.662	11.0	90	0.00
4	S	2-Fluorophenol	1.330	1.258	5.4	99	0.00
5	S	Phenol-d6	1.842	1.504	18.3	99	0.00
6		bis(2-Chloroethyl)ether	1.424	1.258	11.7	102	0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8	S	Nitrobenzene-d5	0.365	0.304	16.7	100	0.00
9		Naphthalene	1.106	1.182	-6.9	106	0.00
10		Hexachlorobutadiene	0.175	0.186	-6.3	105	0.00
11		2-Methylnaphthalene	0.728	0.752	-3.3	104	0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	92	-0.02
13	S	2,4,6-Tribromophenol	0.196	0.164	16.3	90	0.00
14	S	2-Fluorobiphenyl	1.526	1.531	-0.3	100	0.00
15		Acenaphthylene	8.673	8.876	-2.3	103	0.00
16		Acenaphthene	1.394	1.473	-5.7	104	0.00
17		Fluorene	1.772	1.815	-2.4	101	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
19		Hexachlorobenzene	0.220	0.234	-6.4	105	0.00
20		Pentachlorophenol	0.070	0.053	24.3	90	0.00
21		Phenanthrene	1.215	1.321	-8.7	104	0.00
22		Anthracene	1.181	1.187	-0.5	101	0.00
23		Fluoranthene	6.088	6.266	-2.9	99	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
25		Pyrene	4.561	4.478	1.8	98	0.00
26	S	Terphenyl-d14	0.689	0.649	5.8	94	0.00
27		Benzo(a)anthracene	5.548	5.613	-1.2	101	0.00
28		Chrysene	4.695	5.294	-12.8	105	0.00
29		Bis(2-ethylhexyl)phthalate	0.534	0.410	23.2	76	0.00
30		Indeno(1,2,3-cd)pyrene	5.479	5.586	-2.0	100	0.00
31	I	Perylene-d12	1.000	1.000	0.0	90	0.00
32		Benzo(b)fluoranthene	1.350	1.283	5.0	100	0.00
33		Benzo(k)fluoranthene	1.317	1.436	-9.0	99	-0.01
34	C	Benzo(a)pyrene	1.282	1.308	-2.0	100	0.00
35		Dibenzo(a,h)anthracene	1.211	1.219	-0.7	100	-0.02
36		Benzo(g,h,i)perylene	5.006	5.150	-2.9	100	0.00

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