Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101716\
Data File : BE092455.D
Acq On : 17 Oct 2016 14:37
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 17 15:39:06 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	163#	-0.02
2		1,4-Dioxane	0.635	0.634	0.2	161#	-0.01
3		n-Nitrosodimethylamine	0.588	0.664	-12.9	204#	-0.02
4	S	2-Fluorophenol	0.999	0.904	9.5	158#	-0.01
5	S	Phenol-d6	1.388	1.207	13.0	163#	-0.02
6		bis(2-Chloroethyl)ether	1.222	1.145	6.3	196#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	153#	0.00
8	S	Nitrobenzene-d5	0.299	0.301	-0.7	183#	-0.02
9		Naphthalene	1.082	1.176	-8.7	167#	-0.02
10		Hexachlorobutadiene	0.170	0.181	-6.5	164#	-0.02
11		2-Methylnaphthalene	0.716	0.752	-5.0	163#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
13	S	2,4,6-Tribromophenol	0.144	0.106	26.4#	123	0.00
14	S	2-Fluorobiphenyl	1.184	1.353	-14.3	165#	0.00
15		Acenaphthylene	7.240	7.413	-2.4	149	0.00
16		Acenaphthene	1.134	1.219	-7.5	148	0.00
17		Fluorene	1.438	1.473	-2.4	148	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.02
19		Hexachlorobenzene	0.214	0.243	-13.6	151#	0.00
20		Pentachlorophenol	0.050	0.043	14.0	127	0.00
21		Phenanthrene	1.149	1.343	-16.9	153#	0.00
22		Anthracene	1.105	1.163	-5.2	144	0.02
23		Fluoranthene	5.460	5.663	-3.7	140	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	126	0.00
25		Pyrene	4.333	5.064	-16.9	153#	0.00
26	S	Terphenyl-d14	0.620	0.689	-11.1	142	0.02
27		Benzo(a)anthracene	4.873	5.064	-3.9	134	0.00
28		Chrysene	4.462	4.977	-11.5	131	0.00
29		Bis(2-ethylhexyl)phthalate	0.420	0.403	4.0	120	0.00
30		Indeno(1,2,3-cd)pyrene	4.734	4.422	6.6	120	0.04
31	I	Perylene-d12	1.000	1.000	0.0	117	0.02
32		Benzo(b)fluoranthene	1.211	1.270	-4.9	123	0.02
33		Benzo(k)fluoranthene	1.267	1.381	-9.0	137	0.02
34	C	Benzo(a)pyrene	1.187	1.241	-4.5	129	0.02
35		Dibenzo(a,h)anthracene	1.088	1.075	1.2	121	0.05
36		Benzo(g,h,i)perylene	4.673	4.648	0.5	120	0.05

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