Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE120916\
Data File : BE092598.D
Acq On : 9 Dec 2016 17:56
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 09 23:35:25 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 15:43:18 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	124	-0.02
2		1,4-Dioxane	0.681	0.816	-19.8	154#	0.00
3		n-Nitrosodimethylamine	0.654	0.607	7.2	134	0.00
4	S	2-Fluorophenol	0.844	0.702	16.8	103	0.00
5	S	Phenol-d6	1.104	1.082	2.0	131	0.00
6		bis(2-Chloroethyl)ether	1.341	1.130	15.7	117	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
8	S	Nitrobenzene-d5	0.269	0.236	12.3	110	-0.02
9		Naphthalene	1.032	1.046	-1.4	125	-0.02
10		Hexachlorobutadiene	0.157	0.162	-3.2	126	0.00
11		2-Methylnaphthalene	0.658	0.653	0.8	129	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
13	S	2,4,6-Tribromophenol	0.112	0.109	2.7	134	0.00
14	S	2-Fluorobiphenyl	1.219	1.219	0.0	128	-0.01
15		Acenaphthylene	6.962	6.814	2.1	132	0.00
16		Acenaphthene	1.136	1.141	-0.4	130	-0.02
17		Fluorene	1.408	1.420	-0.9	131	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	119	0.00
19		Hexachlorobenzene	0.197	0.219	-11.2	148	0.00
20		Pentachlorophenol	0.036	0.040	-11.1	171#	0.00
21		Phenanthrene	1.119	1.214	-8.5	134	-0.02
22		Anthracene	1.053	1.082	-2.8	131	0.00
23		Fluoranthene	5.219	5.323	-2.0	130	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	119	0.00
25		Pyrene	3.975	3.963	0.3	127	0.00
26	S	Terphenyl-d14	0.572	0.558	2.4	126	0.00
27		Benzo(a)anthracene	4.692	4.266	9.1	129	0.00
28		Chrysene	4.250	4.712	-10.9	127	-0.02
29		Bis(2-ethylhexyl)phthalate	0.305	0.342	-12.1	160#	-0.02
30		Indeno(1,2,3-cd)pyrene	4.722	4.286	9.2	114	0.00
31	I	Perylene-d12	1.000	1.000	0.0	114	-0.01
32		Benzo(b)fluoranthene	1.213	1.108	8.7	108	-0.01
33		Benzo(k)fluoranthene	1.270	1.272	-0.2	120	-0.01
34	C	Benzo(a)pyrene	1.122	1.085	3.3	119	-0.01
35		Dibenzo(a,h)anthracene	1.111	1.024	7.8	113	-0.02
36		Benzo(g,h,i)perylene	4.702	4.408	6.3	111	0.00

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