Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA M\DATA\BM112115\
Data File : BM003273.D
Acq On : 22 Nov 2015 06:03
Operator : UM/IZ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.461

Quant Time: Nov 23 01:05:49 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM111815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 04:42:03 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	109	0.00
2	SURR	1,4-Dioxane-d8	0.427	0.407	4.7	103	0.00
3		1,4-Dioxane	0.486	0.462	4.9	104	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
5		Naphthalene	1.027	1.045	-1.8	115	0.00
6	SURR	2-Methylnaphthalene-d10	0.544	0.552	-1.5	114	0.00
7		2-Methylnaphthalene	0.669	0.694	-3.7	116	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
9		Acenaphthylene	1.665	1.948	-17.0	132	0.00
10	C	Acenaphthene	1.406	1.461	-3.9	115	0.00
11		Fluorene	1.521	1.581	-3.9	114	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
13		Pentachlorophenol	0.094	0.105	-11.7	128	0.00
14		Phenanthrene	1.161	1.210	-4.2	114	0.00
15		Anthracene	0.972	1.116	-14.8	128	0.00
16	SURR	Fluoranthene-d10	0.980	1.066	-8.8	122	0.00
17	C	Fluoranthene	1.287	1.390	-8.0	121	0.00
18	I	Chrysene-d12	1.000	1.000	0.0	130	0.00
19		Pyrene	1.371	1.395	-1.8	129	0.00
20		Benzo(a)anthracene	1.160	1.351	-16.5	154#	0.00
21		Chrysene	1.306	1.353	-3.6	131	0.00
22	I	Perylene-d12	1.000	1.000	0.0	151#	0.00
23		Benzo(b)fluoranthene	1.521	1.415	7.0	142	0.00
24		Benzo(k)fluoranthene	1.438	1.388	3.5	140	0.00
25	C	Benzo(a)pyrene	1.353	1.406	-3.9	156#	0.00
26		Indeno(1,2,3-cd)pyrene	1.612	1.591	1.3	147	0.00
27		Dibenzo(a,h)anthracene	1.303	1.280	1.8	147	0.00
28		Benzo(g,h,i)perylene	1.416	1.342	5.2	141	0.00

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