Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061416\
Data File : BM005781.D
Acq On : 14 Jun 2016 10:46
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.466

Quant Time: Jun 14 15:52:41 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM061316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 19:45:38 2016
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	111	0.00
2	SURR	1,4-Dioxane-d8	0.195	0.204	-4.6	108	0.00
3		1,4-Dioxane	0.271	0.269	0.7	109	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.01
5		Naphthalene	0.907	0.993	-9.5	115	-0.01
6	SURR	2-Methylnaphthalene-d10	0.499	0.567	-13.6	115	-0.01
7		2-Methylnaphthalene	0.634	0.727	-14.7	116	-0.01
8	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
9		Acenaphthylene	1.814	1.958	-7.9	112	0.00
10	C	Acenaphthene	1.277	1.383	-8.3	113	0.00
11		Fluorene	1.223	1.328	-8.6	112	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
13		Pentachlorophenol	0.028	0.026	7.1	116	0.00
14		Phenanthrene	0.951	1.024	-7.7	111	0.00
15		Anthracene	1.177	0.964	18.1	83	-0.02
16	SURR	Fluoranthene-d10	0.841	0.888	-5.6	114	0.00
17	C	Fluoranthene	1.199	1.258	-4.9	113	-0.01
18	I	Chrysene-d12	1.000	1.000	0.0	109	-0.01
19		Pyrene	1.375	1.553	-12.9	112	-0.01
20		Benzo(a)anthracene	0.936	1.115	-19.1	126	-0.01
21		Chrysene	1.432	1.488	-3.9	103	-0.01
22	I	Perylene-d12	1.000	1.000	0.0	112	-0.01
23		Benzo(b)fluoranthene	1.118	1.339	-19.8	132	-0.01
24		Benzo(k)fluoranthene	1.392	1.391	0.1	98	0.00
25	C	Benzo(a)pyrene	1.253	1.363	-8.8	111	-0.01
26		Indeno(1,2,3-cd)pyrene	1.437	1.570	-9.3	112	-0.01
27		Dibenzo(a,h)anthracene	1.111	1.199	-7.9	111	-0.01
28		Benzo(g,h,i)perylene	1.270	1.410	-11.0	114	-0.01

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