Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061416\
Data File : BM005783.D
Acq On : 14 Jun 2016 13:25
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.467

Quant Time: Jun 14 16:07:44 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	105	0.00
2	SURR	1,4-Dioxane-d8	0.195	0.214	-9.7	107	0.00
3		1,4-Dioxane	0.271	0.279	-3.0	107	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
5		Naphthalene	0.907	0.992	-9.4	103	-0.01
6	SURR	2-Methylnaphthalene-d10	0.499	0.549	-10.0	99	0.00
7		2-Methylnaphthalene	0.634	0.700	-10.4	100	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
9		Acenaphthylene	1.814	1.915	-5.6	95	0.00
10	C	Acenaphthene	1.277	1.367	-7.0	96	0.00
11		Fluorene	1.223	1.289	-5.4	94	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
13		Pentachlorophenol	0.028	0.027	3.6	105	0.00
14		Phenanthrene	0.951	1.060	-11.5	98	0.00
15		Anthracene	1.177	1.194	-1.4	88	0.00
16	SURR	Fluoranthene-d10	0.841	0.936	-11.3	103	0.00
17	C	Fluoranthene	1.199	1.322	-10.3	102	-0.01
18	I	Chrysene-d12	1.000	1.000	0.0	101	-0.01
19		Pyrene	1.375	1.528	-11.1	103	-0.01
20		Benzo(a)anthracene	0.936	1.006	-7.5	106	-0.01
21		Chrysene	1.432	1.553	-8.4	100	-0.01
22	I	Perylene-d12	1.000	1.000	0.0	104	-0.01
23		Benzo(b)fluoranthene	1.118	1.224	-9.5	112	0.00
24		Benzo(k)fluoranthene	1.392	1.570	-12.8	103	0.00
25	C	Benzo(a)pyrene	1.253	1.359	-8.5	103	0.00
26		Indeno(1,2,3-cd)pyrene	1.437	1.567	-9.0	104	-0.01
27		Dibenzo(a,h)anthracene	1.111	1.201	-8.1	103	0.00
28		Benzo(g,h,i)perylene	1.270	1.395	-9.8	105	0.00

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