Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092815\
Data File : BM002729.D
Acq On : 28 Sep 2015 11:25
Operator : UM/IZ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.487

Quant Time: Sep 29 00:27:47 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM092415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 25 15:05:16 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	119	0.00
2	SURR	1,4-Dioxane-d8	0.413	0.402	2.7	119	0.00
3		1,4-Dioxane	0.479	0.467	2.5	118	-0.02
4	I	Naphthalene-d8	1.000	1.000	0.0	117	-0.01
5		Naphthalene	1.084	1.050	3.1	113	0.00
6	SURR	2-Methylnaphthalene-d10	0.585	0.518	11.5	109	0.00
7		2-Methylnaphthalene	0.698	0.674	3.4	113	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
9		Acenaphthylene	2.063	1.974	4.3	111	-0.02
10	C	Acenaphthene	1.506	1.499	0.5	108	0.00
11		Fluorene	1.717	1.631	5.0	106	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
13		Pentachlorophenol	0.078	0.043	44.9#	60	0.00
14		Phenanthrene	1.302	1.191	8.5	103	0.00
15		Anthracene	1.203	1.149	4.5	103	-0.01
16	SURR	Fluoranthene-d10	0.997	0.987	1.0	106	-0.01
17	C	Fluoranthene	1.416	1.342	5.2	105	0.00
18	I	Chrysene-d12	1.000	1.000	0.0	110	0.00
19		Pyrene	1.559	1.455	6.7	105	0.00
20		Benzo(a)anthracene	1.337	1.300	2.8	107	0.00
21		Chrysene	1.280	1.279	0.1	109	0.00
22	I	Perylene-d12	1.000	1.000	0.0	114	-0.01
23		Benzo(b)fluoranthene	1.442	1.392	3.5	115	0.00
24		Benzo(k)fluoranthene	1.336	1.345	-0.7	110	0.00
25	C	Benzo(a)pyrene	1.328	1.321	0.5	113	0.00
26		Indeno(1,2,3-cd)pyrene	1.481	1.514	-2.2	116	-0.01
27		Dibenzo(a,h)anthracene	1.215	1.232	-1.4	115	-0.01
28		Benzo(g,h,i)perylene	1.256	1.303	-3.7	118	-0.01

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