Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
Data File : BM006344.D
Acq On : 08 Jul 2016 06:44
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.446

Quant Time: Jul 08 08:58:18 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 08 08:04:54 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	89	0.00
2	SURR	1,4-Dioxane-d8	0.169	0.179	-5.9	86	0.00
3		1,4-Dioxane	0.219	0.226	-3.2	86	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
5		Naphthalene	0.933	0.996	-6.8	88	0.00
6	SURR	2-Methylnaphthalene-d10	0.468	0.529	-13.0	93	0.00
7		2-Methylnaphthalene	0.625	0.703	-12.5	92	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.02
9		Acenaphthylene	1.347	1.378	-2.3	91	0.00
10	C	Acenaphthene	1.337	1.437	-7.5	93	0.00
11		Fluorene	1.500	1.607	-7.1	95	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
13		Pentachlorophenol	0.062	0.073	-17.7	122	0.00
14		Phenanthrene	1.106	1.191	-7.7	96	0.00
15		Anthracene	1.025	1.130	-10.2	97	0.00
16	SURR	Fluoranthene-d10	0.843	0.975	-15.7	103	0.00
17	C	Fluoranthene	1.220	1.399	-14.7	103	0.00
18	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
19		Pyrene	1.675	1.746	-4.2	105	0.00
20		Benzo(a)anthracene	1.295	1.462	-12.9	109	0.00
21		Chrysene	1.269	1.442	-13.6	110	0.00
22	I	Perylene-d12	1.000	1.000	0.0	108	0.00
23		Benzo(b)fluoranthene	1.432	1.610	-12.4	116	0.00
24		Benzo(k)fluoranthene	1.332	1.384	-3.9	99	0.00
25	C	Benzo(a)pyrene	1.333	1.452	-8.9	107	0.00
26		Indeno(1,2,3-cd)pyrene	1.479	1.700	-14.9	109	0.00
27		Dibenzo(a,h)anthracene	1.154	1.318	-14.2	109	0.00
28		Benzo(g,h,i)perylene	1.251	1.421	-13.6	108	0.00

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