Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042916\
Data File : BM005135.D
Acq On : 29 Apr 2016 09:15
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.452

Quant Time: Apr 29 13:58:11 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM041816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 19 08:25:48 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	88	0.00
2	SURR	1,4-Dioxane-d8	0.193	0.165	14.5	76	0.00
3		1,4-Dioxane	0.238	0.208	12.6	79	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
5		Naphthalene	0.946	0.960	-1.5	93	0.00
6	SURR	2-Methylnaphthalene-d10	0.476	0.520	-9.2	98	0.00
7		2-Methylnaphthalene	0.630	0.697	-10.6	100	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
9		Acenaphthylene	1.608	1.649	-2.5	111	0.00
10	C	Acenaphthene	1.340	1.368	-2.1	108	0.00
11		Fluorene	1.451	1.583	-9.1	115	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	114	0.00
13		Pentachlorophenol	0.047	0.046	2.1	152#	0.00
14		Phenanthrene	1.099	1.158	-5.4	121	0.00
15		Anthracene	0.951	1.072	-12.7	136	0.00
16	SURR	Fluoranthene-d10	0.815	0.922	-13.1	131	0.00
17	C	Fluoranthene	1.132	1.282	-13.3	132	0.00
18	I	Chrysene-d12	1.000	1.000	0.0	156#	0.00
19		Pyrene	1.631	1.475	9.6	134	0.00
20		Benzo(a)anthracene	1.064	1.117	-5.0	164#	0.00
21		Chrysene	1.282	1.262	1.6	156#	0.00
22	I	Perylene-d12	1.000	1.000	0.0	145	0.00
23		Benzo(b)fluoranthene	1.433	1.467	-2.4	155#	0.00
24		Benzo(k)fluoranthene	1.381	1.419	-2.8	147	0.00
25	C	Benzo(a)pyrene	1.230	1.247	-1.4	152#	0.00
26		Indeno(1,2,3-cd)pyrene	1.273	1.201	5.7	140	0.00
27		Dibenzo(a,h)anthracene	0.972	0.952	2.1	147	0.00
28		Benzo(g,h,i)perylene	1.166	1.048	10.1	132	0.00

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