Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM010815\
Data File : BM000145.D
Acq On : 08 Jan 2015 15:20
Operator : TP/IZ
Sample : SSTD0.422
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jan 08 16:34:16 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM010715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 08 10:41:47 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	99	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	98	-0.01
3		Naphthalene	1.072	1.043	2.7	98	-0.01
4	SURR	2-Methylnaphthalene-d10	0.532	0.520	2.3	98	-0.01
5		2-Methylnaphthalene	0.710	0.694	2.3	98	-0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
7		Acenaphthylene	1.996	1.976	1.0	100	-0.01
8	C	Acenaphthene	1.247	1.228	1.5	100	-0.02
9		Fluorene	1.433	1.449	-1.1	100	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
11		Pentachlorophenol	0.063	0.047	25.4#	97	-0.02
12		Phenanthrene	1.235	1.173	5.0	98	0.00
13		Anthracene	1.159	1.136	2.0	103	0.00
14	SURR	Fluoranthene-d10	0.862	0.858	0.5	102	0.00
15	C	Fluoranthene	1.355	1.343	0.9	102	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	103	-0.01
17		Pyrene	1.669	1.572	5.8	101	0.00
18		Benzo(a)anthracene	1.324	1.279	3.4	101	-0.01
19		Chrysene	1.444	1.404	2.8	104	-0.01
20	I	Perylene-d12	1.000	1.000	0.0	105	-0.01
21		Benzo(b)fluoranthene	1.574	1.571	0.2	115	-0.01
22		Benzo(k)fluoranthene	1.503	1.377	8.4	94	-0.01
23	C	Benzo(a)pyrene	1.425	1.373	3.6	104	-0.01
24		Indeno(1,2,3-cd)pyrene	1.626	1.560	4.1	105	-0.01
25		Dibenzo(a,h)anthracene	1.293	1.256	2.9	106	-0.01
26		Benzo(g,h,i)perylene	1.444	1.380	4.4	103	-0.01

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