Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM010815\
Data File : BM000153.D
Acq On : 09 Jan 2015 01:01
Operator : TP/IZ
Sample : SSTD0.423
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jan 09 03:33:37 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	93	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	93	-0.01
3		Naphthalene	1.072	1.036	3.4	92	-0.01
4	SURR	2-Methylnaphthalene-d10	0.532	0.517	2.8	92	-0.01
5		2-Methylnaphthalene	0.710	0.696	2.0	93	-0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
7		Acenaphthylene	1.996	1.921	3.8	94	-0.01
8	C	Acenaphthene	1.247	1.200	3.8	94	-0.02
9		Fluorene	1.433	1.421	0.8	94	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
11		Pentachlorophenol	0.063	0.044	30.2#	85	-0.02
12		Phenanthrene	1.235	1.179	4.5	92	0.00
13		Anthracene	1.159	1.138	1.8	96	0.00
14	SURR	Fluoranthene-d10	0.862	0.873	-1.3	97	0.00
15	C	Fluoranthene	1.355	1.362	-0.5	97	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	98	-0.01
17		Pyrene	1.669	1.578	5.5	97	0.00
18		Benzo(a)anthracene	1.324	1.260	4.8	95	-0.01
19		Chrysene	1.444	1.383	4.2	97	-0.01
20	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
21		Benzo(b)fluoranthene	1.574	1.438	8.6	99	-0.01
22		Benzo(k)fluoranthene	1.503	1.523	-1.3	98	-0.01
23	C	Benzo(a)pyrene	1.425	1.370	3.9	97	-0.01
24		Indeno(1,2,3-cd)pyrene	1.626	1.579	2.9	100	-0.01
25		Dibenzo(a,h)anthracene	1.293	1.264	2.2	100	-0.01
26		Benzo(g,h,i)perylene	1.444	1.396	3.3	98	-0.01

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