Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021115\
Data File : BM000498.D
Acq On : 12 Feb 2015 04:04
Operator : TP/IZ
Sample : SSTD0.404
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.404

Quant Time: Feb 12 07:34:28 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 11 23:58:48 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	103	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
3		Naphthalene	0.977	0.952	2.6	104	0.00
4	SURR	2-Methylnaphthalene-d10	0.536	0.524	2.2	102	0.00
5		2-Methylnaphthalene	0.683	0.659	3.5	102	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
7		Acenaphthylene	1.597	1.626	-1.8	109	0.00
8	C	Acenaphthene	1.224	1.225	-0.1	105	0.00
9		Fluorene	1.493	1.492	0.1	104	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
11		Pentachlorophenol	0.037	0.048	-29.7#	144	0.00
12		Phenanthrene	1.129	1.078	4.5	105	0.00
13		Anthracene	1.014	0.991	2.3	107	0.00
14	SURR	Fluoranthene-d10	0.886	0.867	2.1	107	0.00
15	C	Fluoranthene	1.303	1.277	2.0	110	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
17		Pyrene	1.468	1.434	2.3	112	0.00
18		Benzo(a)anthracene	1.175	1.127	4.1	114	0.00
19		Chrysene	1.356	1.305	3.8	110	0.00
20	I	Perylene-d12	1.000	1.000	0.0	113	0.00
21		Benzo(b)fluoranthene	1.377	1.414	-2.7	119	0.00
22		Benzo(k)fluoranthene	1.464	1.333	8.9	104	0.00
23	C	Benzo(a)pyrene	1.296	1.271	1.9	114	0.00
24		Indeno(1,2,3-cd)pyrene	1.363	1.243	8.8	110	-0.01
25		Dibenzo(a,h)anthracene	1.078	0.986	8.5	112	0.00
26		Benzo(g,h,i)perylene	1.181	1.067	9.7	109	-0.01

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