Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120914\
Data File : BE089197.D
Acq On : 9 Dec 2014 21:31
Operator : TP/JJ
Sample : SSTD0.437
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTD0.437

Quant Time: Dec 10 01:04:26 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM01.2-EPA-SIM-BE120514.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 05 23:47:26 2014
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	119	-0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	121	-0.02
3		Naphthalene	0.958	0.994	-3.8	120	-0.02
4	SURR	2-Methylnaphthalene-d10	0.463	0.492	-6.3	123	-0.02
5		2-Methylnaphthalene	0.591	0.648	-9.6	126	-0.02
6	I	Acenaphthene-d10	1.000	1.000	0.0	122	-0.03
7		Acenaphthylene	7.001	7.249	-3.5	122	-0.02
8	C	Acenaphthene	4.263	4.489	-5.3	123	-0.02
9		Fluorene	1.162	1.219	-4.9	122	-0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	118	-0.02
11		Pentachlorophenol	0.071	0.051	28.2#	83	-0.01
12		Phenanthrene	4.311	4.570	-6.0	119	-0.02
13		Anthracene	4.248	4.436	-4.4	117	-0.02
14	SURR	Fluoranthene-d10	0.711	0.715	-0.6	113	-0.03
15	C	Fluoranthene	3.890	3.907	-0.4	112	-0.03
16	I	Chrysene-d12	1.000	1.000	0.0	106	-0.03
17		Pyrene	5.537	6.027	-8.8	111	-0.03
18		Benzo(a)anthracene	1.033	1.053	-1.9	103	-0.03
19		Chrysene	1.065	1.114	-4.6	108	-0.03
20	I	Perylene-d12	1.000	1.000	0.0	103	-0.03
21		Benzo(b)fluoranthene	0.989	1.042	-5.4	103	-0.03
22		Benzo(k)fluoranthene	1.160	1.251	-7.8	104	-0.03
23	C	Benzo(a)pyrene	1.040	1.084	-4.2	100	-0.03
24		Indeno(1,2,3-cd)pyrene	3.338	3.353	-0.4	99	-0.04
25		Dibenzo(a,h)anthracene	0.853	0.899	-5.4	102	-0.04
26		Benzo(g,h,i)perylene	4.128	4.182	-1.3	101	-0.05

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