Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120914\
Data File : BE089203.D
Acq On : 10 Dec 2014 1:27
Operator : TP/JJ
Sample : SSTD0.438
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTD0.438

Quant Time: Dec 10 02:16:02 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 : Wed Dec 10 01:02:19 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	123	-0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	124	-0.02
3		Naphthalene	0.958	0.993	-3.7	124	-0.02
4	SURR	2-Methylnaphthalene-d10	0.463	0.491	-6.0	126	-0.02
5		2-Methylnaphthalene	0.591	0.644	-9.0	129	-0.02
6	I	Acenaphthene-d10	1.000	1.000	0.0	122	-0.03
7		Acenaphthylene	7.001	7.422	-6.0	125	-0.02
8	C	Acenaphthene	4.263	4.470	-4.9	123	-0.02
9		Fluorene	1.162	1.220	-5.0	122	-0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	116	-0.03
11		Pentachlorophenol	0.071	0.059	16.9	94	-0.03
12		Phenanthrene	4.311	4.545	-5.4	116	-0.03
13		Anthracene	4.248	4.479	-5.4	116	-0.03
14	SURR	Fluoranthene-d10	0.711	0.697	2.0	108	-0.03
15	C	Fluoranthene	3.890	3.812	2.0	107	-0.03
16	I	Chrysene-d12	1.000	1.000	0.0	106	-0.03
17		Pyrene	5.537	5.800	-4.7	107	-0.03
18		Benzo(a)anthracene	1.033	1.098	-6.3	108	-0.03
19		Chrysene	1.065	1.113	-4.5	108	-0.03
20	I	Perylene-d12	1.000	1.000	0.0	107	-0.03
21		Benzo(b)fluoranthene	0.989	1.071	-8.3	110	-0.03
22		Benzo(k)fluoranthene	1.160	1.248	-7.6	108	-0.03
23	C	Benzo(a)pyrene	1.040	1.130	-8.7	109	-0.03
24		Indeno(1,2,3-cd)pyrene	3.338	3.473	-4.0	107	-0.04
25		Dibenzo(a,h)anthracene	0.853	0.922	-8.1	109	-0.04
26		Benzo(g,h,i)perylene	4.128	4.181	-1.3	106	-0.04

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