Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE083116\
Data File : BE092312.D
Acq On : 31 Aug 2016 9:24
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTD0.473

Quant Time: Aug 31 10:02:21 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 26 11:18:03 2016
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	155#	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	174#	0.00
3		Naphthalene	0.916	0.918	-0.2	168#	0.00
4	SURR	2-Methylnaphthalene-d10	0.527	0.547	-3.8	176#	-0.01
5		2-Methylnaphthalene	0.610	0.635	-4.1	176#	-0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	179#	0.00
7		Acenaphthylene	1.734	1.788	-3.1	177#	0.00
8	C	Acenaphthene	1.212	1.241	-2.4	175#	0.00
9		Fluorene	1.408	1.410	-0.1	170#	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	150	0.00
11		Pentachlorophenol	0.047	0.055	-17.0	194#	0.00
12		Phenanthrene	0.987	1.036	-5.0	148	0.00
13		Anthracene	0.897	1.014	-13.0	165#	-0.02
14	SURR	Fluoranthene-d10	0.992	1.062	-7.1	154#	0.00
15	C	Fluoranthene	1.161	1.217	-4.8	150#	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	135	0.00
17		Pyrene	0.938	1.170	-24.7#	150#	0.00
18		Benzo(a)anthracene	0.858	1.087	-26.7#	154#	0.00
19		Chrysene	1.005	1.121	-11.5	134	0.00
20	I	Perylene-d12	1.000	1.000	0.0	141	0.00
21		Benzo(b)fluoranthene	0.987	1.081	-9.5	141	0.00
22		Benzo(k)fluoranthene	1.156	1.163	-0.6	125	0.00
23	C	Benzo(a)pyrene	1.043	1.112	-6.6	133	0.00
24		Indeno(1,2,3-cd)pyrene	1.132	1.248	-10.2	139	0.00
25		Dibenzo(a,h)anthracene	0.904	1.021	-12.9	143	0.00
26		Benzo(g,h,i)perylene	1.005	1.064	-5.9	134	0.00

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