Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120616\
Data File : BM008194.D
Acq On : 06 Dec 2016 20:40
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.466

Quant Time: Dec 07 06:15:18 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 07 05:56:15 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	74	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
3		Naphthalene	1.119	1.087	2.9	82	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.574	-1.4	86	0.01
5		2-Methylnaphthalene	0.721	0.708	1.8	83	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
7		Acenaphthylene	1.810	1.732	4.3	86	0.00
8	C	Acenaphthene	1.377	1.327	3.6	88	0.00
9		Fluorene	1.583	1.501	5.2	86	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
11		Pentachlorophenol	0.140	0.099	29.3#	60	0.00
12		Phenanthrene	1.246	1.190	4.5	79	0.00
13		Anthracene	1.170	1.287	-10.0	90	0.00
14	SURR	Fluoranthene-d10	1.132	1.018	10.1	74	0.00
15	C	Fluoranthene	1.608	1.316	18.2	69	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	54	0.00
17		Pyrene	1.389	1.763	-26.9#	68	0.00
18		Benzo(a)anthracene	1.228	1.148	6.5	52	0.01
19		Chrysene	1.527	1.530	-0.2	55	0.01
20	I	Perylene-d12	1.000	1.000	0.0	55	0.00
21		Benzo(b)fluoranthene	1.596	1.344	15.8	47	0.00
22		Benzo(k)fluoranthene	1.647	1.753	-6.4	61	0.01
23	C	Benzo(a)pyrene	1.495	1.391	7.0	52	0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.717	7.1	51	0.01
25		Dibenzo(a,h)anthracene	1.481	1.338	9.7	50	0.02
26		Benzo(g,h,i)perylene	1.640	1.539	6.2	51	0.02

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