Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120616\
Data File : BM008181.D
Acq On : 06 Dec 2016 12:52
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.465

Quant Time: Dec 06 13:48:46 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 03 04:54:05 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	79	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
3		Naphthalene	1.119	1.073	4.1	76	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.541	4.4	77	0.00
5		2-Methylnaphthalene	0.721	0.664	7.9	74	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	73	0.00
7		Acenaphthylene	1.810	1.768	2.3	71	0.00
8	C	Acenaphthene	1.377	1.312	4.7	71	0.00
9		Fluorene	1.583	1.407	11.1	66	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	65	0.00
11		Pentachlorophenol	0.140	0.113	19.3	54	0.00
12		Phenanthrene	1.246	1.178	5.5	62	0.00
13		Anthracene	1.170	1.187	-1.5	66	0.00
14	SURR	Fluoranthene-d10	1.132	1.033	8.7	60	0.00
15	C	Fluoranthene	1.608	1.375	14.5	58	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	53	0.00
17		Pyrene	1.389	1.500	-8.0	58	0.00
18		Benzo(a)anthracene	1.228	1.249	-1.7	56	0.00
19		Chrysene	1.527	1.410	7.7	51	0.00
20	I	Perylene-d12	1.000	1.000	0.0	58	0.00
21		Benzo(b)fluoranthene	1.596	1.473	7.7	54	0.00
22		Benzo(k)fluoranthene	1.647	1.462	11.2	53	0.00
23	C	Benzo(a)pyrene	1.495	1.420	5.0	56	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.740	5.8	54	0.00
25		Dibenzo(a,h)anthracene	1.481	1.376	7.1	54	0.00
26		Benzo(g,h,i)perylene	1.640	1.564	4.6	55	0.00

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