Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112116\
Data File : BM007973.D
Acq On : 22 Nov 2016 00:04
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.445

Quant Time: Nov 22 05:01:32 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 22 04:57:47 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	92	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
3		Naphthalene	1.119	1.129	-0.9	96	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.562	0.7	95	0.00
5		2-Methylnaphthalene	0.721	0.721	0.0	96	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
7		Acenaphthylene	1.810	1.829	-1.0	98	0.00
8	C	Acenaphthene	1.377	1.370	0.5	98	0.00
9		Fluorene	1.583	1.587	-0.3	99	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
11		Pentachlorophenol	0.140	0.129	7.9	95	0.00
12		Phenanthrene	1.246	1.243	0.2	101	0.00
13		Anthracene	1.170	1.126	3.8	97	0.00
14	SURR	Fluoranthene-d10	1.132	1.102	2.7	99	0.00
15	C	Fluoranthene	1.608	1.620	-0.7	104	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	93	0.00
17		Pyrene	1.389	1.449	-4.3	97	0.00
18		Benzo(a)anthracene	1.228	1.234	-0.5	96	0.00
19		Chrysene	1.527	1.471	3.7	92	0.00
20	I	Perylene-d12	1.000	1.000	0.0	91	0.00
21		Benzo(b)fluoranthene	1.596	1.658	-3.9	96	0.00
22		Benzo(k)fluoranthene	1.647	1.553	5.7	89	0.00
23	C	Benzo(a)pyrene	1.495	1.499	-0.3	93	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.835	0.7	90	0.00
25		Dibenzo(a,h)anthracene	1.481	1.459	1.5	90	0.00
26		Benzo(g,h,i)perylene	1.640	1.613	1.6	89	0.00

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