Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121216\
Data File : BM008258.D
Acq On : 12 Dec 2016 09:43
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.434

Quant Time: Dec 13 02:28:14 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 12 03:51:49 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	138	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	145	0.00
3		Naphthalene	1.119	1.095	2.1	141	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.568	-0.4	146	0.00
5		2-Methylnaphthalene	0.721	0.712	1.2	143	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
7		Acenaphthylene	1.810	1.763	2.6	139	0.00
8	C	Acenaphthene	1.377	1.302	5.4	137	0.00
9		Fluorene	1.583	1.409	11.0	128	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	123	0.00
11		Pentachlorophenol	0.140	0.121	13.6	110	0.00
12		Phenanthrene	1.246	1.191	4.4	119	0.00
13		Anthracene	1.170	1.143	2.3	121	0.00
14	SURR	Fluoranthene-d10	1.132	1.028	9.2	113	0.00
15	C	Fluoranthene	1.608	1.397	13.1	110	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	90	0.00
17		Pyrene	1.389	1.719	-23.8#	111	0.00
18		Benzo(a)anthracene	1.228	1.436	-16.9	108	0.00
19		Chrysene	1.527	1.359	11.0	82	0.00
20	I	Perylene-d12	1.000	1.000	0.0	97	0.00
21		Benzo(b)fluoranthene	1.596	1.407	11.8	87	0.00
22		Benzo(k)fluoranthene	1.647	1.555	5.6	95	0.00
23	C	Benzo(a)pyrene	1.495	1.452	2.9	96	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.751	5.2	91	0.00
25		Dibenzo(a,h)anthracene	1.481	1.377	7.0	91	0.00
26		Benzo(g,h,i)perylene	1.640	1.512	7.8	89	0.00

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