Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092216\
Data File : BM007520.D
Acq On : 22 Sep 2016 20:28
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.442

Quant Time: Sep 23 12:15:13 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 04:00:23 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	117	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
3		Naphthalene	0.931	0.934	-0.3	117	0.00
4	SURR	2-Methylnaphthalene-d10	0.466	0.464	0.4	115	0.00
5		2-Methylnaphthalene	0.630	0.630	0.0	116	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
7		Acenaphthylene	1.693	1.638	3.2	112	0.00
8	C	Acenaphthene	1.320	1.302	1.4	114	0.00
9		Fluorene	1.618	1.581	2.3	110	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	108	0.00
11		Pentachlorophenol	0.109	0.108	0.9	118	0.00
12		Phenanthrene	1.082	1.090	-0.7	110	0.02
13		Anthracene	0.981	0.978	0.3	109	0.00
14	SURR	Fluoranthene-d10	0.932	0.933	-0.1	109	0.00
15	C	Fluoranthene	1.364	1.377	-1.0	111	-0.01
16	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
17		Pyrene	1.322	1.374	-3.9	115	0.00
18		Benzo(a)anthracene	1.182	1.219	-3.1	116	0.00
19		Chrysene	1.262	1.276	-1.1	109	0.00
20	I	Perylene-d12	1.000	1.000	0.0	112	0.00
21		Benzo(b)fluoranthene	1.335	1.410	-5.6	117	0.00
22		Benzo(k)fluoranthene	1.349	1.340	0.7	112	0.00
23	C	Benzo(a)pyrene	1.241	1.290	-3.9	117	0.00
24		Indeno(1,2,3-cd)pyrene	1.278	1.369	-7.1	122	0.00
25		Dibenzo(a,h)anthracene	0.976	1.079	-10.6	126	0.00
26		Benzo(g,h,i)perylene	1.148	1.212	-5.6	119	0.00

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