Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
Data File : BM008099.D
Acq On : 29 Nov 2016 09:16
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.459

Quant Time: Nov 29 10:25:33 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 05:49:20 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	123	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	144	-0.01
3		Naphthalene	1.119	1.079	3.6	138	-0.01
4	SURR	2-Methylnaphthalene-d10	0.566	0.566	0.0	144	0.00
5		2-Methylnaphthalene	0.721	0.716	0.7	143	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	145	0.00
7		Acenaphthylene	1.810	1.813	-0.2	146	0.00
8	C	Acenaphthene	1.377	1.339	2.8	144	0.00
9		Fluorene	1.583	1.471	7.1	137	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	138	-0.02
11		Pentachlorophenol	0.140	0.131	6.4	133	-0.02
12		Phenanthrene	1.246	1.189	4.6	133	0.00
13		Anthracene	1.170	1.159	0.9	137	0.00
14	SURR	Fluoranthene-d10	1.132	1.118	1.2	138	-0.01
15	C	Fluoranthene	1.608	1.950	-21.3#	172#	-0.01
16	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
17		Pyrene	1.389	1.648	-18.6	131	-0.01
18		Benzo(a)anthracene	1.228	1.397	-13.8	130	-0.01
19		Chrysene	1.527	1.331	12.8	99	0.00
20	I	Perylene-d12	1.000	1.000	0.0	116	-0.01
21		Benzo(b)fluoranthene	1.596	1.478	7.4	110	-0.01
22		Benzo(k)fluoranthene	1.647	1.406	14.6	103	-0.01
23	C	Benzo(a)pyrene	1.495	1.384	7.4	109	-0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.662	10.1	104	-0.01
25		Dibenzo(a,h)anthracene	1.481	1.326	10.5	105	-0.02
26		Benzo(g,h,i)perylene	1.640	1.409	14.1	99	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1