Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122916\
Data File : BM008713.D
Acq On : 29 Dec 2016 09:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.441

Quant Time: Dec 30 00:55:22 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 29 03:29:39 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	99	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	121	0.00
3		Naphthalene	1.125	1.052	6.5	116	0.00
4	SURR	2-Methylnaphthalene-d10	0.613	0.580	5.4	120	0.00
5		2-Methylnaphthalene	0.757	0.649	14.3	107	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
7		Acenaphthylene	1.809	1.987	-9.8	143	0.00
8	C	Acenaphthene	1.361	1.397	-2.6	132	0.00
9		Fluorene	1.553	1.490	4.1	122	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
11		Pentachlorophenol	0.124	0.113	8.9	104	0.00
12		Phenanthrene	1.221	1.256	-2.9	129	0.00
13		Anthracene	1.186	1.233	-4.0	135	0.00
14	SURR	Fluoranthene-d10	1.101	1.261	-14.5	143	0.00
15	C	Fluoranthene	1.498	1.683	-12.3	143	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	129	0.00
17		Pyrene	1.546	1.703	-10.2	144	0.00
18		Benzo(a)anthracene	1.313	1.369	-4.3	142	0.00
19		Chrysene	1.438	1.553	-8.0	137	0.00
20	I	Perylene-d12	1.000	1.000	0.0	134	0.00
21		Benzo(b)fluoranthene	1.576	1.529	3.0	132	0.00
22		Benzo(k)fluoranthene	1.537	1.510	1.8	130	0.00
23	C	Benzo(a)pyrene	1.461	1.502	-2.8	139	0.00
24		Indeno(1,2,3-cd)pyrene	1.741	1.711	1.7	129	0.00
25		Dibenzo(a,h)anthracene	1.391	1.347	3.2	127	0.00
26		Benzo(g,h,i)perylene	1.493	1.477	1.1	129	0.00

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