Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111616\
Data File : BM007925.D
Acq On : 16 Nov 2016 17:28
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.465

Quant Time: Nov 16 17:59:43 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 16:16:41 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	74	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	76	0.00
3		Naphthalene	1.029	1.077	-4.7	77	0.00
4	SURR	2-Methylnaphthalene-d10	0.474	0.505	-6.5	78	0.00
5		2-Methylnaphthalene	0.650	0.686	-5.5	77	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
7		Acenaphthylene	1.882	1.820	3.3	81	0.00
8	C	Acenaphthene	1.432	1.363	4.8	79	0.00
9		Fluorene	1.591	1.511	5.0	80	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	81	0.00
11		Pentachlorophenol	0.094	0.113	-20.2#	107	0.00
12		Phenanthrene	1.117	1.140	-2.1	78	0.00
13		Anthracene	1.020	1.084	-6.3	85	0.00
14	SURR	Fluoranthene-d10	0.905	1.060	-17.1	90	0.00
15	C	Fluoranthene	1.336	1.483	-11.0	85	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	82	0.00
17		Pyrene	1.372	1.495	-9.0	81	0.00
18		Benzo(a)anthracene	1.147	1.260	-9.9	84	0.00
19		Chrysene	1.299	1.399	-7.7	77	0.00
20	I	Perylene-d12	1.000	1.000	0.0	81	0.00
21		Benzo(b)fluoranthene	1.377	1.507	-9.4	77	0.00
22		Benzo(k)fluoranthene	1.428	1.521	-6.5	81	0.00
23	C	Benzo(a)pyrene	1.275	1.400	-9.8	81	0.00
24		Indeno(1,2,3-cd)pyrene	1.547	1.715	-10.9	79	0.00
25		Dibenzo(a,h)anthracene	1.244	1.372	-10.3	79	0.00
26		Benzo(g,h,i)perylene	1.390	1.537	-10.6	78	0.00

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