Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112116\
Data File : BM007958.D
Acq On : 21 Nov 2016 15:10
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.444

Quant Time: Nov 21 15:49:27 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 Nov 21 15:11:37 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	92	0.00
2	I	Naphthalene-d8	0.400	0.400	0.0	93	0.00
3		Naphthalene	0.400	0.391	2.3	91	0.00
4	SURR	2-Methylnaphthalene-d10	0.400	0.386	3.5	90	0.00
5		2-Methylnaphthalene	0.400	0.388	3.0	91	0.00
6	I	Acenaphthene-d10	0.400	0.400	0.0	91	0.00
7		Acenaphthylene	0.400	0.403	-0.8	92	0.00
8	C	Acenaphthene	0.400	0.399	0.3	93	0.00
9		Fluorene	0.400	0.388	3.0	90	0.00
10	I	Phenanthrene-d10	0.400	0.400	0.0	91	0.00
11		Pentachlorophenol	0.400	0.361	9.8	85	0.00
12		Phenanthrene	0.400	0.399	0.3	92	0.00
13		Anthracene	0.400	0.391	2.3	89	0.00
14	SURR	Fluoranthene-d10	0.400	0.389	2.8	90	0.00
15	C	Fluoranthene	0.400	0.379	5.3	89	0.00
16	I	Chrysene-d12	0.400	0.400	0.0	89	0.00
17		Pyrene	0.400	0.396	1.0	88	0.00
18		Benzo(a)anthracene	0.400	0.400	0.0	92	0.00
19		Chrysene	0.400	0.373	6.8	86	0.00
20	I	Perylene-d12	0.400	0.400	0.0	86	0.00
21		Benzo(b)fluoranthene	0.400	0.375	6.3	82	0.00
22		Benzo(k)fluoranthene	0.400	0.411	-2.7	92	0.00
23	C	Benzo(a)pyrene	0.400	0.394	1.5	87	0.00
24		Indeno(1,2,3-cd)pyrene	0.400	0.389	2.8	84	0.00
25		Dibenzo(a,h)anthracene	0.400	0.387	3.3	84	0.00
26		Benzo(g,h,i)perylene	0.400	0.393	1.8	84	0.00

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