Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011516\
Data File : BE091372.D
Acq On : 15 Jan 2016 9:38
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jan 16 05:03:17 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 15 15:14:26 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	91	0.00
2	S	2-Fluorophenol	1.000	1.500	-50.0#	145	-0.01
3	S	Phenol-d6	1.000	1.171	-17.1	135	0.00
4	I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
5	S	Nitrobenzene-d5	1.000	0.982	1.8	98	0.00
6		Naphthalene	1.000	1.012	-1.2	97	0.00
7		2-Methylnaphthalene	1.000	1.009	-0.9	127	0.00
8	I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
9	S	2,4,6-Tribromophenol	1.000	1.032	-3.2	134	0.00
10	S	2-Fluorobiphenyl	1.000	1.029	-2.9	121	0.00
11		Acenaphthylene	1.000	0.994	0.6	127	0.00
12		Acenaphthene	1.000	0.979	2.1	103	0.00
13		Fluorene	1.000	0.967	3.3	114	0.00
14	I	Phenanthrene-d10	5.000	5.000	0.0	93	0.00
15		Phenanthrene	1.000	1.037	-3.7	94	-0.01
16		Anthracene	1.000	1.029	-2.9	112	0.00
17		Fluoranthene	1.000	1.030	-3.0	100	0.00
18	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
19		Pyrene	1.000	0.998	0.2	126	0.00
20	S	Terphenyl-d14	1.000	1.031	-3.1	124	0.00
21		Benzo(a)anthracene	1.000	1.040	-4.0	138	0.00
22		Chrysene	1.000	1.014	-1.4	97	0.00
23		Indeno(1,2,3-cd)pyrene	1.000	0.980	2.0	113	0.00
24	I	Perylene-d12	5.000	5.000	0.0	98	0.00
25		Benzo(b)fluoranthene	1.000	1.029	-2.9	131	0.00
26		Benzo(k)fluoranthene	1.000	1.049	-4.9	120	0.00
27	C	Benzo(a)pyrene	1.000	0.967	3.3	104	0.00
28		Dibenzo(a,h)anthracene	1.000	1.083	-8.3	151	0.00
29		Benzo(g,h,i)perylene	1.000	1.060	-6.0	130	0.00

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