Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021916\
Data File : BE091413.D
Acq On : 19 Feb 2016 13:19
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Feb 19 22:37:56 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 03 00:27:53 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	95	-0.01
2	S	2-Fluorophenol	1.000	1.072	-7.2	127	-0.03
3	S	Phenol-d6	1.000	0.940	6.0	95	-0.02
4	I	Naphthalene-d8	5.000	5.000	0.0	100	-0.02
5	S	Nitrobenzene-d5	1.000	1.113	-11.3	139	-0.02
6		Naphthalene	1.000	0.938	6.2	106	-0.02
7		2-Methylnaphthalene	1.000	1.028	-2.8	120	-0.02
8	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
9	S	2,4,6-Tribromophenol	1.000	1.013	-1.3	145	-0.01
10	S	2-Fluorobiphenyl	1.000	0.957	4.3	109	-0.01
11		Acenaphthylene	1.000	1.039	-3.9	130	0.00
12		Acenaphthene	1.000	0.933	6.7	109	0.00
13		Fluorene	1.000	0.971	2.9	110	0.00
14	I	Phenanthrene-d10	5.000	5.000	0.0	99	-0.01
15		Phenanthrene	1.000	0.949	5.1	106	0.00
16		Anthracene	1.000	1.049	-4.9	115	-0.01
17		Fluoranthene	1.000	1.026	-2.6	121	0.00
18	I	Chrysene-d12	5.000	5.000	0.0	98	0.00
19		Pyrene	1.000	1.094	-9.4	121	0.00
20	S	Terphenyl-d14	1.000	1.065	-6.5	114	0.00
21		Benzo(a)anthracene	1.000	1.098	-9.8	135	0.00
22		Chrysene	1.000	0.995	0.5	107	0.00
23		Indeno(1,2,3-cd)pyrene	1.000	1.125	-12.5	131	0.00
24	I	Perylene-d12	5.000	5.000	0.0	106	0.00
25		Benzo(b)fluoranthene	1.000	0.996	0.4	120	0.00
26		Benzo(k)fluoranthene	1.000	1.056	-5.6	130	0.00
27	C	Benzo(a)pyrene	1.000	1.102	-10.2	139	0.00
28		Dibenzo(a,h)anthracene	1.000	1.110	-11.0	146	0.00
29		Benzo(g,h,i)perylene	1.000	1.051	-5.1	130	0.00

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