Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021716\
Data File : BE091411.D
Acq On : 17 Feb 2016 19:20
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Feb 18 00:00:41 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	97	-0.01
2	S	2-Fluorophenol	1.000	1.134	-13.4	138	-0.04
3	S	Phenol-d6	1.000	1.206	-20.6	131	-0.03
4	I	Naphthalene-d8	5.000	5.000	0.0	102	-0.01
5	S	Nitrobenzene-d5	1.000	1.227	-22.7	162	-0.03
6		Naphthalene	1.000	0.938	6.2	108	-0.02
7		2-Methylnaphthalene	1.000	1.037	-3.7	124	-0.02
8	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
9	S	2,4,6-Tribromophenol	1.000	1.183	-18.3	180	-0.01
10	S	2-Fluorobiphenyl	1.000	0.968	3.2	110	-0.02
11		Acenaphthylene	1.000	1.094	-9.4	137	0.00
12		Acenaphthene	1.000	0.921	7.9	108	0.00
13		Fluorene	1.000	0.964	3.6	109	0.00
14	I	Phenanthrene-d10	5.000	5.000	0.0	96	-0.01
15		Phenanthrene	1.000	0.926	7.4	101	0.00
16		Anthracene	1.000	1.092	-9.2	116	-0.01
17		Fluoranthene	1.000	1.092	-9.2	126	0.00
18	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
19		Pyrene	1.000	1.094	-9.4	124	0.00
20	S	Terphenyl-d14	1.000	1.089	-8.9	120	0.00
21		Benzo(a)anthracene	1.000	1.193	-19.3	152	0.00
22		Chrysene	1.000	0.926	7.4	102	0.00
23		Indeno(1,2,3-cd)pyrene	1.000	1.230	-23.0	148	-0.01
24	I	Perylene-d12	5.000	5.000	0.0	111	0.00
25		Benzo(b)fluoranthene	1.000	1.023	-2.3	129	0.00
26		Benzo(k)fluoranthene	1.000	1.031	-3.1	133	0.00
27	C	Benzo(a)pyrene	1.000	1.120	-12.0	148	0.00
28		Dibenzo(a,h)anthracene	1.000	1.223	-22.3	173	-0.01
29		Benzo(g,h,i)perylene	1.000	1.098	-9.8	143	-0.01

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