Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120914\
Data File : Be089190.d
Acq On : 9 Dec 2014 17:03
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Dec 10 00:56:23 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE120514.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 08 00:50:53 2014
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	116	-0.02
2		1,4-Dioxane	1.000	1.016	-1.6	111	-0.02
3	I	Naphthalene-d8	5.000	5.000	0.0	119	-0.02
4	S	Nitrobenzene-d5	1.000	1.080	-8.0	123	-0.02
5		Naphthalene	1.000	1.042	-4.2	120	-0.02
6		2-Methylnaphthalene	1.000	1.100	-10.0	127	-0.02
7	I	Acenaphthene-d10	5.000	5.000	0.0	124	-0.01
8	S	2-Fluorobiphenyl	1.000	1.066	-6.6	127	-0.01
9		Acenaphthylene	1.000	1.080	-8.0	128	-0.02
10		Acenaphthene	1.000	1.061	-6.1	129	-0.01
11		Fluorene	1.000	1.071	-7.1	127	-0.02
12	I	Phenanthrene-d10	5.000	5.000	0.0	119	-0.02
13		Pentachlorophenol	1.000	0.934	6.6	131	-0.02
14		Phenanthrene	1.000	1.054	-5.4	120	-0.02
15		Anthracene	1.000	1.072	-7.2	121	-0.03
16		Fluoranthene	1.000	1.032	-3.2	115	-0.02
17	I	Chrysene-d12	5.000	5.000	0.0	109	-0.02
18		Pyrene	1.000	1.092	-9.2	114	-0.03
19	S	Terphenyl-d14	1.000	1.093	-9.3	116	-0.02
20		Benzo(a)anthracene	1.000	1.031	-3.1	111	-0.03
21		Chrysene	1.000	1.051	-5.1	109	-0.03
22		Indeno(1,2,3-cd)pyrene	1.000	1.077	-7.7	118	-0.04
23	I	Perylene-d12	5.000	5.000	0.0	109	-0.03
24		Benzo(b)fluoranthene	1.000	1.073	-7.3	112	-0.03
25		Benzo(k)fluoranthene	1.000	1.054	-5.4	110	-0.03
26	C	Benzo(a)pyrene	1.000	1.061	-6.1	114	-0.03
27		Dibenzo(a,h)anthracene	1.000	1.106	-10.6	117	-0.04
28		Benzo(g,h,i)perylene	1.000	1.065	-6.5	111	-0.04

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