Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102014\
Data File : BE088123.D
Acq On : 21 Oct 2014 00:02
Operator : TP/JJ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE102014

Quant Time: Oct 21 03:10:21 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 21 02:11:14 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	118	0.00
2		1,4-Dioxane	1.000	1.070	-7.0	120	0.00
3	I	Naphthalene-d8	5.000	5.000	0.0	118	0.00
4	S	Nitrobenzene-d5	1.000	1.111	-11.1	127	0.00
5		Naphthalene	1.000	1.078	-7.8	123	0.00
6		2-Methylnaphthalene	1.000	1.094	-9.4	124	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	117	0.00
8	S	2-Fluorobiphenyl	1.000	1.106	-10.6	124	0.00
9		Acenaphthylene	1.000	1.118	-11.8	125	0.00
10		Acenaphthene	1.000	1.092	-9.2	126	0.00
11		Fluorene	1.000	1.091	-9.1	123	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	117	0.00
13		Phenanthrene	1.000	1.256	-25.6#	122	0.00
14		Anthracene	1.000	1.074	-7.4	123	0.00
15		Fluoranthene	1.000	1.035	-3.5	121	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	117	0.00
17		Pyrene	1.000	1.035	-3.5	121	0.00
18	S	Terphenyl-d14	1.000	1.060	-6.0	122	0.00
19		Benzo(a)anthracene	1.000	0.996	0.4	120	0.00
20		Chrysene	1.000	1.037	-3.7	121	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	1.030	-3.0	121	0.00
22	I	Perylene-d12	5.000	5.000	0.0	116	0.00
23		Benzo(b)fluoranthene	1.000	1.012	-1.2	117	0.00
24		Benzo(k)fluoranthene	1.000	1.055	-5.5	122	0.00
25	C	Benzo(a)pyrene	1.000	1.039	-3.9	117	0.00
26		Dibenzo(a,h)anthracene	1.000	1.032	-3.2	119	0.00
27		Benzo(g,h,i)perylene	1.000	1.031	-3.1	119	0.00

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