Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be102414\
Data File : BE088216.D
Acq On : 24 Oct 2014 17:09
Operator : TP/JJ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE102414

Quant Time: Oct 27 13:10:08 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102414.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 27 13:04:40 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	108	0.00
2		1,4-Dioxane	1.000	1.107	-10.7	113	-0.01
3	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
4	S	Nitrobenzene-d5	1.000	1.050	-5.0	109	0.00
5		Naphthalene	1.000	1.100	-10.0	112	0.00
6		2-Methylnaphthalene	1.000	1.106	-10.6	110	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
8	S	2-Fluorobiphenyl	1.000	1.106	-10.6	109	0.00
9		Acenaphthylene	1.000	1.132	-13.2	109	0.00
10		Acenaphthene	1.000	1.076	-7.6	107	0.00
11		Fluorene	1.000	1.125	-12.5	108	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
13		Phenanthrene	1.000	1.077	-7.7	106	0.00
14		Anthracene	1.000	1.116	-11.6	107	0.00
15		Fluoranthene	1.000	1.118	-11.8	108	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	109	0.00
17		Pyrene	1.000	1.083	-8.3	109	0.00
18	S	Terphenyl-d14	1.000	1.078	-7.8	110	0.00
19		Benzo(a)anthracene	1.000	1.044	-4.4	112	0.00
20		Chrysene	1.000	1.057	-5.7	112	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	0.982	1.8	116	0.00
22	I	Perylene-d12	5.000	5.000	0.0	112	0.00
23		Benzo(b)fluoranthene	1.000	0.973	2.7	109	0.00
24		Benzo(k)fluoranthene	1.000	1.107	-10.7	117	0.00
25	C	Benzo(a)pyrene	1.000	1.003	-0.3	115	0.00
26		Dibenzo(a,h)anthracene	1.000	1.017	-1.7	119	0.00
27		Benzo(g,h,i)perylene	1.000	1.092	-9.2	119	0.00

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