Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE010715\
Data File : BE089359.D
Acq On : 8 Jan 2015 11:22
Operator : TP/IZ
Sample : G1011-03
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : MW3

Quant Time: Jan 08 16:00:29 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BASE-BE010715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 08 15:21:44 2015
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	94	0.00
2	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
3	S	Nitrobenzene-d5	1.000	6.769	-576.9#	611	0.00
4		Naphthalene	1.000	0.000	100.0#	0	-8.53#
5		2-Methylnaphthalene	1.000	0.010	99.0#	1	0.00
6	I	Acenaphthene-d10	5.000	5.000	0.0	99	0.00
7	S	2-Fluorobiphenyl	1.000	6.445	-544.5#	616	0.00
8		Acenaphthylene	1.000	0.005	99.5#	0	0.00
9	C	Acenaphthene	1.000	0.017	98.3#	2	0.00
10		Fluorene	1.000	0.000	100.0#	0	-11.34#
11	I	Phenanthrene-d10	5.000	5.000	0.0	103	0.00
12		Phenanthrene	1.000	0.017	98.3#	2	0.00
13		Anthracene	1.000	0.000	100.0#	0	-12.84#
14	C	Fluoranthene	1.000	0.007	99.3#	1	0.00
15	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
16		Pyrene	1.000	0.052	94.8#	5	0.00
17	S	Terphenyl-d14	1.000	7.464	-646.4#	752	0.00
18		Benzo(a)anthracene	1.000	0.014	98.6#	1	0.02
19		Chrysene	1.000	0.013	98.7#	1	-0.06
20		Indeno(1,2,3-cd)pyrene	1.000	0.000	100.0#	0	-0.05
21	I	Perylene-d12	5.000	5.000	0.0	97	0.00
22		Benzo(b)fluoranthene	1.000	0.002	99.8#	0	0.01
23		Benzo(k)fluoranthene	1.000	0.002	99.8#	0	0.01
24	C	Benzo(a)pyrene	1.000	0.001	99.9#	0	0.00
25		Dibenzo(a,h)anthracene	1.000	0.001	99.9#	0	0.03
26		Benzo(g,h,i)perylene	1.000	0.001	99.9#	0	0.04

(#) = Out of Range SPCC's out = 0 CCC's out = 3