Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
Data File : BM001658.D
Acq On : 12 Jun 2015 01:13
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM061215

Quant Time: Jun 12 16:46:12 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 16:27:37 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	80	0.00
2		1,4-Dioxane	1.000	1.054	-5.4	84	0.00
3		n-Nitrosodimethylamine	1.000	1.011	-1.1	83	0.00
4	S	2-Fluorophenol	1.000	0.962	3.8	81	0.00
5	S	Phenol-d6	1.000	0.953	4.7	81	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	81	0.00
7	S	Nitrobenzene-d5	1.000	0.953	4.7	82	0.00
8		Nitrobenzene	1.000	0.947	5.3	81	0.00
9		Naphthalene	1.000	0.973	2.7	83	0.00
10		Hexachlorobutadiene	1.000	0.975	2.5	85	0.00
11		2-Methylnaphthalene	1.000	0.983	1.7	85	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	82	0.00
13	S	2,4,6-Tribromophenol	1.000	1.027	-2.7	82	0.00
14	S	2-Fluorobiphenyl	1.000	0.989	1.1	86	0.00
15		Acenaphthylene	1.000	0.975	2.5	83	0.00
16		Acenaphthene	1.000	0.954	4.6	84	0.00
17		Fluorene	1.000	1.001	-0.1	87	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	80	0.00
19		4-Bromophenyl-phenylether	1.000	1.028	-2.8	81	0.00
20		Hexachlorobenzene	1.000	1.132	-13.2	91	0.00
21		Pentachlorophenol	1.000	0.894	10.6	81	0.00
22		Phenanthrene	1.000	0.941	5.9	81	0.00
23		Anthracene	1.000	0.985	1.5	79	0.00
24		Fluoranthene	1.000	1.012	-1.2	84	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	79	0.00
26		Pyrene	1.000	1.002	-0.2	83	0.00
27	S	Terphenyl-d14	1.000	1.003	-0.3	82	0.00
28		Benzo(a)anthracene	1.000	0.984	1.6	82	0.00
29		Chrysene	1.000	0.950	5.0	79	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.963	3.7	81	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.860	14.0	75	0.00
32	I	Perylene-d12	5.000	5.000	0.0	76	0.00
33		Benzo(b)fluoranthene	1.000	0.972	2.8	79	0.00
34		Benzo(k)fluoranthene	1.000	0.985	1.5	78	0.00
35	C	Benzo(a)pyrene	1.000	0.952	4.8	76	0.00
36		Dibenzo(a,h)anthracene	1.000	0.898	10.2	74	0.00
37		Benzo(g,h,i)perylene	1.000	1.003	-0.3	77	0.00

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