Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052215\
Data File : BM001393.D
Acq On : 22 May 2015 21:29
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM052215

Quant Time: May 23 05:00:51 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 23 04:45:39 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	95	0.00
2		1,4-Dioxane	1.000	0.845	15.5	88	0.00
3		n-Nitrosodimethylamine	1.000	0.913	8.7	86	0.00
4	S	2-Fluorophenol	1.000	0.935	6.5	86	0.00
5	S	Phenol-d6	1.000	0.914	8.6	85	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	1.000	0.909	9.1	85	0.00
8		Nitrobenzene	1.000	0.893	10.7	84	0.00
9		Naphthalene	1.000	0.962	3.8	87	0.00
10		Hexachlorobutadiene	1.000	0.956	4.4	86	0.00
11		2-Methylnaphthalene	1.000	0.951	4.9	86	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.790	21.0	74	0.00
14	S	2-Fluorobiphenyl	1.000	0.949	5.1	87	0.00
15		Acenaphthylene	1.000	0.940	6.0	88	0.00
16		Acenaphthene	1.000	0.961	3.9	89	0.00
17		Fluorene	1.000	0.872	12.8	85	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	97	0.00
19		4-Bromophenyl-phenylether	1.000	1.060	-6.0	84	0.00
20		Hexachlorobenzene	1.000	1.023	-2.3	95	0.00
21		Pentachlorophenol	1.000	1.122	-12.2	88	0.00
22		Phenanthrene	1.000	1.043	-4.3	96	0.00
23		Anthracene	1.000	0.929	7.1	81	0.00
24		Fluoranthene	1.000	0.895	10.5	88	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	98	0.00
26		Pyrene	1.000	0.921	7.9	88	0.00
27	S	Terphenyl-d14	1.000	0.910	9.0	87	0.00
28		Benzo(a)anthracene	1.000	0.932	6.8	92	0.00
29		Chrysene	1.000	0.905	9.5	87	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.781	21.9	84	-0.02
31		Indeno(1,2,3-cd)pyrene	1.000	0.922	7.8	87	-0.01
32	I	Perylene-d12	5.000	5.000	0.0	95	0.00
33		Benzo(b)fluoranthene	1.000	0.933	6.7	91	0.00
34		Benzo(k)fluoranthene	1.000	0.936	6.4	85	0.00
35	C	Benzo(a)pyrene	1.000	0.923	7.7	88	0.00
36		Dibenzo(a,h)anthracene	1.000	0.935	6.5	88	-0.01
37		Benzo(g,h,i)perylene	1.000	0.937	6.3	88	-0.01

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