Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM050115\
Data File : BM001147.D
Acq On : 02 May 2015 00:56
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM050115

Quant Time: May 04 17:02:19 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM050115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 04 16:48:07 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	0.876	12.4	80	0.00
3		n-Nitrosodimethylamine	1.000	1.019	-1.9	92	0.00
4	S	2-Fluorophenol	1.000	1.021	-2.1	96	0.00
5	S	Phenol-d6	1.000	1.006	-0.6	95	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	96	0.00
7	S	Nitrobenzene-d5	1.000	0.980	2.0	94	0.00
8		Nitrobenzene	1.000	0.976	2.4	94	0.00
9		Naphthalene	1.000	1.028	-2.8	96	0.00
10		Hexachlorobutadiene	1.000	1.019	-1.9	97	0.00
11		2-Methylnaphthalene	1.000	1.022	-2.2	98	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
13	S	2,4,6-Tribromophenol	1.000	1.244	-24.4	151	0.00
14	S	2-Fluorobiphenyl	1.000	1.027	-2.7	98	0.00
15		Acenaphthylene	1.000	1.006	-0.6	97	0.00
16		Acenaphthene	1.000	1.006	-0.6	99	0.00
17		Fluorene	1.000	1.036	-3.6	97	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	97	0.00
19		Hexachlorobenzene	1.000	1.012	-1.2	100	0.00
20		Pentachlorophenol	1.000	0.846	15.4	88	0.00
21		Phenanthrene	1.000	1.033	-3.3	101	0.00
22		Anthracene	1.000	1.014	-1.4	96	0.00
23		Fluoranthene	1.000	1.017	-1.7	99	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	98	0.00
25		Pyrene	1.000	1.020	-2.0	100	0.00
26	S	Terphenyl-d14	1.000	1.018	-1.8	98	0.00
27		Benzo(a)anthracene	1.000	1.016	-1.6	100	0.00
28		Chrysene	1.000	1.008	-0.8	95	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	1.014	-1.4	96	0.00
30	I	Perylene-d12	5.000	5.000	0.0	98	-0.01
31		Benzo(b)fluoranthene	1.000	0.966	3.4	88	-0.01
32		Benzo(k)fluoranthene	1.000	1.018	-1.8	107	-0.01
33	C	Benzo(a)pyrene	1.000	0.983	1.7	96	0.00
34		Dibenzo(a,h)anthracene	1.000	0.992	0.8	96	-0.01
35		Benzo(g,h,i)perylene	1.000	0.983	1.7	95	0.00

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