Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\Data\BM041015\
Data File : BM000950.D
Acq On : 10 Apr 2015 21:36
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 11 00:10:10 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM040915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 10 23:03:27 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	167	0.00
2		1,4-Dioxane	1.000	0.976	2.4	160	0.00
3		n-Nitrosodimethylamine	1.000	1.002	-0.2	171	0.00
4	S	2-Fluorophenol	1.000	1.001	-0.1	168	0.00
5	S	Phenol-d6	1.000	0.996	0.4	171	0.00
6	C	Phenol	1.000	0.998	0.2	172	-0.02
7		bis(2-Chloroethyl)ether	1.000	0.991	0.9	167	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	165	0.00
9	S	Nitrobenzene-d5	1.000	0.987	1.3	171	0.00
10		Nitrobenzene	1.000	0.971	2.9	175	0.00
11		2,4-Dimethylphenol	1.000	1.007	-0.7	172	0.00
12	C	2,4-Dichlorophenol	1.000	0.973	2.7	171	0.00
13		Naphthalene	1.000	0.975	2.5	165	0.00
14	C	Hexachlorobutadiene	1.000	0.948	5.2	162	0.00
15	C	2-Methylnaphthalene	1.000	0.999	0.1	169	0.00
16		1-Methylnaphthalene	1.000	0.998	0.2	169	-0.01
17	I	Acenaphthene-d10	5.000	5.000	0.0	209	0.01
18	S	2,4,6-Tribromophenol	1.000	0.841	15.9	142	0.01
19	S	2-Fluorobiphenyl	1.000	0.940	6.0	168	0.00
20		Acenaphthylene	1.000	0.840	16.0	139	0.01
21	C	Acenaphthene	1.000	0.948	5.2	184	-0.02
22	P	2,4-Dinitrophenol	1.000	0.939	6.1	210	-0.01
23		Fluorene	1.000	1.015	-1.5	202	-0.02
24	I	Phenanthrene-d10	5.000	5.000	0.0	176	0.01
25		Hexachlorobenzene	1.000	0.956	4.4	160	0.01
26	C	Pentachlorophenol	1.000	0.855	14.5	171	0.00
27		Phenanthrene	1.000	0.974	2.6	162	0.01
28		Anthracene	1.000	0.995	0.5	181	-0.01
29	C	Fluoranthene	1.000	0.958	4.2	163	-0.01
30	I	Chrysene-d12	5.000	5.000	0.0	168	0.00
31		Benzidine	1.000	0.960	4.0	172	0.00
32		Pyrene	1.000	0.982	1.8	163	0.00
33	S	Terphenyl-d14	1.000	1.050	-5.0	180	0.00
34		Benzo(a)anthracene	1.000	0.980	2.0	169	0.00
35		3,3'-Dichlorobenzidine	1.000	0.959	4.1	183	0.00
36		Chrysene	1.000	0.979	2.1	163	0.00
37		Indeno(1,2,3-cd)pyrene	1.000	0.991	0.9	168	0.00
38	I	Perylene-d12	5.000	5.000	0.0	167	0.00
39		Benzo(b)fluoranthene	1.000	0.971	2.9	171	0.00
40		Benzo(k)fluoranthene	1.000	0.984	1.6	161	0.00
41	C	Benzo(a)pyrene	1.000	0.975	2.5	167	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Dibenzo(a,h)anthracene	1.000	0.996	0.4	169	0.00
43	Benzo(g,h,i)perylene	1.000	0.980	2.0	165	0.00

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