Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM120915\
Data File : BM003420.D
Acq On : 09 Dec 2015 19:41
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 10 00:23:29 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM120915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 18:52:47 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	109	0.00
2		1,4-Dioxane	0.467	0.458	1.9	103	0.00
3		Pyridine	1.308	1.327	-1.5	106	0.00
4		n-Nitrosodimethylamine	0.419	0.437	-4.3	108	0.00
5	S	2-Fluorophenol	1.125	1.156	-2.8	106	0.00
6		Aniline	2.056	2.112	-2.7	108	0.00
7	S	Phenol-d6	1.562	1.607	-2.9	107	0.00
8		2-Chlorophenol	1.424	1.464	-2.8	108	0.00
9		Benzaldehyde	0.766	0.787	-2.7	101	0.00
10	C	Phenol	1.657	1.696	-2.4	108	0.00
11		bis(2-Chloroethyl)ether	1.341	1.352	-0.8	106	0.00
12		1,3-Dichlorobenzene	1.565	1.585	-1.3	107	0.00
13	C	1,4-Dichlorobenzene	1.613	1.623	-0.6	106	0.00
14		1,2-Dichlorobenzene	1.541	1.553	-0.8	107	0.00
15		Benzyl Alcohol	1.081	1.138	-5.3	107	0.00
16		2,2'-oxybis(1-Chloropropane	1.433	1.463	-2.1	108	0.00
17		2-Methylphenol	1.159	1.212	-4.6	109	0.00
18		Hexachloroethane	0.554	0.572	-3.2	107	0.00
19	P	n-Nitroso-di-n-propylamine	1.025	1.069	-4.3	108	0.00
20		3+4-Methylphenols	1.603	1.665	-3.9	108	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.490	0.501	-2.2	108	0.00
23	S	Nitrobenzene-d5	0.323	0.335	-3.7	108	0.00
24		Nitrobenzene	0.349	0.360	-3.2	109	0.00
25		Isophorone	0.652	0.673	-3.2	107	0.00
26	C	2-Nitrophenol	0.176	0.180	-2.3	109	0.00
27		2,4-Dimethylphenol	0.307	0.312	-1.6	108	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.390	0.5	106	0.00
29	C	2,4-Dichlorophenol	0.293	0.302	-3.1	108	0.00
30		1,2,4-Trichlorobenzene	0.327	0.329	-0.6	108	0.00
31		Naphthalene	1.047	1.047	0.0	108	0.00
32		Benzoic acid	0.196	0.189	3.6	109	0.01
33		4-Chloroaniline	0.432	0.438	-1.4	108	0.00
34	C	Hexachlorobutadiene	0.194	0.195	-0.5	108	0.00
35		Caprolactam	0.097	0.095	2.1	109	0.00
36	C	4-Chloro-3-methylphenol	0.330	0.338	-2.4	108	0.00
37		2-Methylnaphthalene	0.718	0.711	1.0	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.643	0.651	-1.2	107	0.00
40	P	Hexachlorocyclopentadiene	0.359	0.368	-2.5	108	0.00
41	S	2,4,6-Tribromophenol	0.200	0.203	-1.5	108	0.00
42	C	2,4,6-Trichlorophenol	0.419	0.438	-4.5	107	0.00
43		2,4,5-Trichlorophenol	0.441	0.459	-4.1	109	0.00
44	S	2-Fluorobiphenyl	1.468	1.469	-0.1	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.774	1.777	-0.2	107	0.00
46		2-Chloronaphthalene	1.303	1.321	-1.4	109	0.00
47		2-Nitroaniline	0.323	0.329	-1.9	109	0.00
48		Acenaphthylene	2.172	2.200	-1.3	108	0.00
49		Dimethylphthalate	1.648	1.627	1.3	107	0.00
50		2,6-Dinitrotoluene	0.344	0.354	-2.9	108	0.00
51	C	Acenaphthene	1.259	1.248	0.9	108	0.00
52		3-Nitroaniline	0.352	0.369	-4.8	109	0.00
53	P	2,4-Dinitrophenol	0.165	0.166	-0.6	110	0.00
54		Dibenzofuran	1.818	1.798	1.1	108	0.00
55	P	4-Nitrophenol	0.291	0.289	0.7	110	0.00
56		2,4-Dinitrotoluene	0.446	0.462	-3.6	109	0.00
57		Fluorene	1.507	1.475	2.1	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.339	0.342	-0.9	107	0.00
59		Diethylphthalate	1.618	1.587	1.9	107	0.00
60		4-Chlorophenyl-phenylether	0.762	0.739	3.0	107	0.00
61		4-Nitroaniline	0.359	0.356	0.8	111	0.00
62		Azobenzene	1.269	1.276	-0.6	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.124	-2.5	110	0.00
65	c	n-Nitrosodiphenylamine	0.632	0.653	-3.3	107	0.00
66		4-Bromophenyl-phenylether	0.216	0.223	-3.2	107	0.00
67		Hexachlorobenzene	0.232	0.236	-1.7	108	0.00
68		Atrazine	0.197	0.208	-5.6	107	0.00
69	C	Pentachlorophenol	0.132	0.138	-4.5	110	0.00
70		Phenanthrene	1.122	1.118	0.4	108	0.00
71		Anthracene	1.115	1.126	-1.0	108	0.00
72		Carbazole	0.972	0.993	-2.2	109	0.00
73		Di-n-butylphthalate	1.136	1.201	-5.7	110	0.00
74	C	Fluoranthene	1.157	1.157	0.0	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
76		Benzidine	0.641	0.680	-6.1	109	0.00
77		Pyrene	1.403	1.373	2.1	109	0.00
78	S	Terphenyl-d14	0.848	0.822	3.1	108	0.00
79		Butylbenzylphthalate	0.540	0.550	-1.9	113	0.00
80		Benzo(a)anthracene	1.199	1.194	0.4	108	0.00
81		3,3'-Dichlorobenzidine	0.382	0.417	-9.2	111	0.00
82		Chrysene	1.154	1.149	0.4	110	0.00
83		Bis(2-ethylhexyl)phthalate	0.741	0.764	-3.1	112	0.00
84	c	Di-n-octyl phthalate	1.224	1.281	-4.7	112	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.281	-5.6	109	0.00
86	I	Perylene-d12	1.000	1.000	0.0	112	0.00
87		Benzo(b)fluoranthene	1.216	1.198	1.5	108	0.00

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88	Benzo(k)fluoranthene	1.187	1.192	-0.4	110	0.00
89 C	Benzo(a)pyrene	1.144	1.161	-1.5	110	0.00
90	Dibenzo(a,h)anthracene	1.117	1.145	-2.5	109	-0.01
91	Benzo(g,h,i)perylene	1.128	1.154	-2.3	109	0.00

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