Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061416\
Data File : BG022334.D
Acq On : 14 Jun 2016 14:21
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 15 02:20:56 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 17:40:23 2016
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	141	-0.02
2		1,4-Dioxane	0.496	0.480	3.2	149	-0.02
3		Pyridine	1.340	1.392	-3.9	145	-0.02
4		n-Nitrosodimethylamine	0.300	0.355	-18.3	161#	-0.02
5	S	2-Fluorophenol	1.034	1.167	-12.9	154#	-0.02
6		Aniline	2.066	2.301	-11.4	148	-0.01
7	S	Phenol-d6	1.626	1.795	-10.4	149	-0.01
8		2-Chlorophenol	1.430	1.501	-5.0	145	-0.01
9		Benzaldehyde	0.895	0.977	-9.2	145	-0.01
10	C	Phenol	1.677	1.872	-11.6	152#	-0.01
11		bis(2-Chloroethyl)ether	1.337	1.417	-6.0	146	-0.02
12		1,3-Dichlorobenzene	1.533	1.474	3.8	137	-0.02
13	C	1,4-Dichlorobenzene	1.527	1.528	-0.1	141	-0.01
14		1,2-Dichlorobenzene	1.539	1.485	3.5	138	-0.01
15		Benzyl Alcohol	1.051	1.218	-15.9	162#	-0.01
16		2,2'-oxybis(1-Chloropropane	1.387	1.497	-7.9	154#	-0.01
17		2-Methylphenol	1.251	1.363	-9.0	154#	-0.01
18		Hexachloroethane	0.558	0.560	-0.4	145	-0.01
19	P	n-Nitroso-di-n-propylamine	1.101	1.178	-7.0	143	-0.01
20		3+4-Methylphenols	1.795	1.921	-7.0	151#	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	133	-0.01
22		Acetophenone	0.431	0.467	-8.4	140	-0.01
23	S	Nitrobenzene-d5	0.287	0.322	-12.2	145	-0.01
24		Nitrobenzene	0.288	0.325	-12.8	146	-0.01
25		Isophorone	0.671	0.682	-1.6	137	0.00
26	C	2-Nitrophenol	0.156	0.188	-20.5#	151#	-0.01
27		2,4-Dimethylphenol	0.283	0.294	-3.9	136	-0.01
28		bis(2-Chloroethoxy)methane	0.385	0.401	-4.2	138	-0.01
29	C	2,4-Dichlorophenol	0.262	0.286	-9.2	137	0.00
30		1,2,4-Trichlorobenzene	0.293	0.287	2.0	129	-0.01
31		Naphthalene	0.998	0.988	1.0	136	-0.01
32		Benzoic acid	0.087	0.083	4.6	121	-0.01
33		4-Chloroaniline	0.415	0.442	-6.5	139	-0.01
34	C	Hexachlorobutadiene	0.157	0.147	6.4	130	-0.02
35		Caprolactam	0.144	0.144	0.0	133	0.00
36	C	4-Chloro-3-methylphenol	0.311	0.341	-9.6	142	0.00
37		2-Methylnaphthalene	0.737	0.725	1.6	130	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	123	0.00
39		1,2,4,5-Tetrachlorobenzene	0.515	0.527	-2.3	125	0.00
40	P	Hexachlorocyclopentadiene	0.233	0.171	26.6#	88	-0.01
41	S	2,4,6-Tribromophenol	0.226	0.225	0.4	120	-0.01
42	C	2,4,6-Trichlorophenol	0.345	0.400	-15.9	136	-0.01
43		2,4,5-Trichlorophenol	0.382	0.422	-10.5	135	0.00
44	S	2-Fluorobiphenyl	1.312	1.346	-2.6	125	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.505	1.586	-5.4	125	0.00
46		2-Chloronaphthalene	1.154	1.201	-4.1	124	-0.01
47		2-Nitroaniline	0.275	0.333	-21.1	143	-0.01
48		Acenaphthylene	2.024	2.046	-1.1	124	0.00
49		Dimethylphthalate	1.658	1.599	3.6	121	0.00
50		2,6-Dinitrotoluene	0.360	0.374	-3.9	123	0.00
51	C	Acenaphthene	1.213	1.222	-0.7	125	-0.01
52		3-Nitroaniline	0.400	0.421	-5.2	137	0.00
53	P	2,4-Dinitrophenol	0.108	0.106	1.9	108	0.00
54		Dibenzofuran	1.893	1.885	0.4	122	-0.01
55	P	4-Nitrophenol	0.276	0.289	-4.7	119	0.00
56		2,4-Dinitrotoluene	0.484	0.530	-9.5	127	0.00
57		Fluorene	1.631	1.628	0.2	123	-0.01
58		2,3,4,6-Tetrachlorophenol	0.342	0.350	-2.3	127	0.00
59		Diethylphthalate	1.771	1.698	4.1	121	-0.01
60		4-Chlorophenyl-phenylether	0.738	0.724	1.9	119	0.00
61		4-Nitroaniline	0.424	0.429	-1.2	123	0.00
62		Azobenzene	1.341	1.448	-8.0	133	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	119	0.00
64		4,6-Dinitro-2-methylphenol	0.092	0.096	-4.3	115	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.611	-6.1	122	0.00
66		4-Bromophenyl-phenylether	0.187	0.191	-2.1	117	-0.01
67		Hexachlorobenzene	0.218	0.209	4.1	112	0.00
68		Atrazine	0.213	0.202	5.2	111	0.00
69	C	Pentachlorophenol	0.084	0.076	9.5	106	0.00
70		Phenanthrene	1.086	1.071	1.4	118	0.00
71		Anthracene	1.077	1.069	0.7	116	0.00
72		Carbazole	1.020	1.014	0.6	117	0.00
73		Di-n-butylphthalate	1.334	1.303	2.3	117	0.00
74	C	Fluoranthene	1.249	1.152	7.8	112	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.523	0.524	-0.2	92	0.00
77		Pyrene	1.243	1.321	-6.3	109	0.00
78	S	Terphenyl-d14	0.842	0.903	-7.2	109	0.00
79		Butylbenzylphthalate	0.577	0.664	-15.1	118	0.00
80		Benzo(a)anthracene	1.121	1.127	-0.5	105	0.00
81		3,3'-Dichlorobenzidine	0.315	0.329	-4.4	104	0.00
82		Chrysene	1.091	1.077	1.3	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.848	0.911	-7.4	110	0.00
84	c	Di-n-octyl phthalate	1.408	1.534	-8.9	111	0.00
85		Indeno(1,2,3-cd)pyrene	1.224	1.163	5.0	96	0.00
86	I	Perylene-d12	1.000	1.000	0.0	97	0.00
87		Benzo(b)fluoranthene	1.166	1.185	-1.6	98	0.00

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88	Benzo(k)fluoranthene	1.148	1.160	-1.0	97	-0.01
89 C	Benzo(a)pyrene	1.115	1.148	-3.0	98	0.00
90	Dibenzo(a,h)anthracene	1.050	1.110	-5.7	99	0.00
91	Benzo(g,h,i)perylene	1.093	1.100	-0.6	97	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1