Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022718\
Data File : BG032869.D
Acq On : 28 Feb 2018 00:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 28 03:32:49 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 26 17:50:45 2018
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	128	0.00
2		1,4-Dioxane	0.543	0.550	-1.3	131	0.00
3		Pyridine	1.495	1.604	-7.3	132	0.00
4		n-Nitrosodimethylamine	0.600	0.634	-5.7	133	0.00
5	S	2-Fluorophenol	1.250	1.304	-4.3	130	0.00
6		Aniline	2.359	2.333	1.1	127	0.00
7	S	Phenol-d6	1.742	1.809	-3.8	131	0.00
8		2-Chlorophenol	1.368	1.420	-3.8	131	0.00
9		Benzaldehyde	1.004	0.922	8.2	121	0.00
10	C	Phenol	1.757	1.813	-3.2	132	0.00
11		bis(2-Chloroethyl)ether	1.436	1.412	1.7	126	0.00
12		1,3-Dichlorobenzene	1.603	1.602	0.1	128	0.00
13	C	1,4-Dichlorobenzene	1.613	1.596	1.1	126	0.00
14		1,2-Dichlorobenzene	1.546	1.526	1.3	127	0.00
15		Benzyl Alcohol	1.281	1.303	-1.7	131	0.00
16		2,2'-oxybis(1-Chloropropane	2.469	2.391	3.2	125	0.00
17		2-Methylphenol	1.295	1.320	-1.9	130	0.00
18		Hexachloroethane	0.549	0.572	-4.2	133	0.00
19	P	n-Nitroso-di-n-propylamine	1.230	1.217	1.1	128	0.00
20		3+4-Methylphenols	1.801	1.876	-4.2	134	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
22		Acetophenone	0.505	0.494	2.2	130	0.00
23	S	Nitrobenzene-d5	0.242	0.335	-38.4#	171#	0.00
24		Nitrobenzene	0.289	0.349	-20.8	156#	0.00
25		Isophorone	0.702	0.679	3.3	130	0.00
26	C	2-Nitrophenol	0.103	0.164	-59.2#	224#	0.00
27		2,4-Dimethylphenol	0.305	0.300	1.6	134	0.00
28		bis(2-Chloroethoxy)methane	0.409	0.396	3.2	130	0.00
29	C	2,4-Dichlorophenol	0.277	0.283	-2.2	136	0.00
30		1,2,4-Trichlorobenzene	0.324	0.314	3.1	129	0.00
31		Naphthalene	1.045	1.018	2.6	131	0.00
32		Benzoic acid	0.174	0.229	-31.6#	188#	0.00
33		4-Chloroaniline	0.467	0.453	3.0	130	0.00
34	C	Hexachlorobutadiene	0.182	0.172	5.5	125	0.00
35		Caprolactam	0.111	0.118	-6.3	138	0.00
36	C	4-Chloro-3-methylphenol	0.322	0.336	-4.3	136	0.00
37		2-Methylnaphthalene	0.756	0.741	2.0	131	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
39		1,2,4,5-Tetrachlorobenzene	0.594	0.577	2.9	133	0.00
40	P	Hexachlorocyclopentadiene	0.303	0.304	-0.3	134	0.00
41	S	2,4,6-Tribromophenol	0.210	0.219	-4.3	139	0.00
42	C	2,4,6-Trichlorophenol	0.374	0.397	-6.1	143	0.00
43		2,4,5-Trichlorophenol	0.433	0.463	-6.9	144	0.00
44	S	2-Fluorobiphenyl	1.387	1.325	4.5	129	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.748	1.707	2.3	133	0.00
46		2-Chloronaphthalene	1.306	1.261	3.4	131	0.00
47		2-Nitroaniline	0.282	0.408	-44.7#	197#	0.00
48		Acenaphthylene	2.137	2.067	3.3	130	0.00
49		Dimethylphthalate	1.698	1.603	5.6	128	0.00
50		2,6-Dinitrotoluene	0.244	0.349	-43.0#	191#	0.00
51	C	Acenaphthene	1.331	1.291	3.0	131	0.00
52		3-Nitroaniline	0.308	0.402	-30.5#	175#	0.00
53	P	2,4-Dinitrophenol	0.070	0.068	2.9	138	0.00
54		Dibenzofuran	1.895	1.797	5.2	129	0.00
55	P	4-Nitrophenol	0.233	0.258	-10.7	148	0.00
56		2,4-Dinitrotoluene	0.293	0.465	-58.7#	223#	0.00
57		Fluorene	1.564	1.491	4.7	130	0.00
58		2,3,4,6-Tetrachlorophenol	0.336	0.353	-5.1	140	0.00
59		Diethylphthalate	1.791	1.684	6.0	128	0.00
60		4-Chlorophenyl-phenylether	0.765	0.733	4.2	131	0.00
61		4-Nitroaniline	0.336	0.390	-16.1	153#	0.00
62		Azobenzene	1.602	1.510	5.7	128	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
64		4,6-Dinitro-2-methylphenol	0.054	0.087	-61.1#	219#	0.00
65	c	n-Nitrosodiphenylamine	0.651	0.651	0.0	129	0.00
66		4-Bromophenyl-phenylether	0.211	0.208	1.4	129	0.00
67		Hexachlorobenzene	0.229	0.223	2.6	128	0.00
68		Atrazine	0.214	0.206	3.7	123	0.00
69	C	Pentachlorophenol	0.144	0.147	-2.1	131	0.00
70		Phenanthrene	1.116	1.081	3.1	125	0.00
71		Anthracene	1.120	1.088	2.9	125	0.00
72		Carbazole	1.104	1.041	5.7	121	0.00
73		Di-n-butylphthalate	1.355	1.303	3.8	122	0.00
74	C	Fluoranthene	1.233	1.159	6.0	122	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	127	0.00
76		Benzidine	0.682	0.716	-5.0	137	0.00
77		Pyrene	1.410	1.346	4.5	122	0.00
78	S	Terphenyl-d14	0.890	0.837	6.0	121	0.00
79		Butylbenzylphthalate	0.625	0.677	-8.3	131	0.00
80		Benzo(a)anthracene	1.283	1.259	1.9	124	0.00
81		3,3'-Dichlorobenzidine	0.475	0.480	-1.1	125	0.00
82		Chrysene	1.225	1.216	0.7	126	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	0.980	-5.8	127	-0.01
84	c	Di-n-octyl phthalate	1.411	1.579	-11.9	131	-0.02
85		Indeno(1,2,3-cd)pyrene	1.429	1.432	-0.2	123	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	131	-0.01
87		Benzo(b)fluoranthene	1.183	1.144	3.3	124	-0.02

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88	Benzo(k)fluoranthene	1.175	1.140	3.0	127	-0.01
89 C	Benzo(a)pyrene	1.125	1.102	2.0	127	-0.01
90	Dibenzo(a,h)anthracene	1.132	1.090	3.7	123	-0.02
91	Benzo(g,h,i)perylene	1.116	1.086	2.7	125	-0.02

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