Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG091616\
Data File : BG024010.D
Acq On : 17 Sep 2016 1:46
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 17 04:43:25 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 09 16:24:09 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	153#	-0.04
2		1,4-Dioxane	0.470	0.506	-7.7	150#	-0.03
3		Pyridine	1.412	1.578	-11.8	154#	-0.04
4		n-Nitrosodimethylamine	0.744	0.904	-21.5	186#	-0.03
5	S	2-Fluorophenol	1.139	1.229	-7.9	153#	-0.04
6		Aniline	2.328	2.500	-7.4	159#	-0.04
7	S	Phenol-d6	1.754	1.871	-6.7	154#	-0.04
8		2-Chlorophenol	1.320	1.376	-4.2	152#	-0.04
9		Benzaldehyde	1.246	1.240	0.5	143	-0.04
10	C	Phenol	1.845	1.964	-6.4	151#	-0.04
11		bis(2-Chloroethyl)ether	1.451	1.517	-4.5	165#	-0.04
12		1,3-Dichlorobenzene	1.545	1.551	-0.4	153#	-0.04
13	C	1,4-Dichlorobenzene	1.636	1.618	1.1	156#	-0.04
14		1,2-Dichlorobenzene	1.536	1.547	-0.7	155#	-0.04
15		Benzyl Alcohol	1.546	1.796	-16.2	164#	-0.04
16		2,2'-oxybis(1-Chloropropane	1.945	2.010	-3.3	162#	-0.04
17		2-Methylphenol	1.243	1.318	-6.0	155#	-0.04
18		Hexachloroethane	0.630	0.666	-5.7	159#	-0.04
19	P	n-Nitroso-di-n-propylamine	1.549	1.708	-10.3	167#	-0.04
20		3+4-Methylphenols	1.732	1.840	-6.2	149	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	149	-0.04
22		Acetophenone	0.518	0.527	-1.7	152#	-0.04
23	S	Nitrobenzene-d5	0.448	0.486	-8.5	165#	-0.04
24		Nitrobenzene	0.482	0.516	-7.1	165#	-0.04
25		Isophorone	0.869	0.922	-6.1	160#	-0.04
26	C	2-Nitrophenol	0.183	0.194	-6.0	160#	-0.04
27		2,4-Dimethylphenol	0.311	0.321	-3.2	145	-0.04
28		bis(2-Chloroethoxy)methane	0.439	0.455	-3.6	159#	-0.04
29	C	2,4-Dichlorophenol	0.330	0.346	-4.8	153#	-0.04
30		1,2,4-Trichlorobenzene	0.362	0.371	-2.5	159#	-0.04
31		Naphthalene	1.031	1.000	3.0	151#	-0.04
32		Benzoic acid	0.253	0.259	-2.4	151#	-0.03
33		4-Chloroaniline	0.430	0.439	-2.1	149	-0.04
34	C	Hexachlorobutadiene	0.272	0.287	-5.5	156#	-0.04
35		Caprolactam	0.116	0.117	-0.9	144	-0.04
36	C	4-Chloro-3-methylphenol	0.441	0.468	-6.1	156#	-0.04
37		2-Methylnaphthalene	0.784	0.770	1.8	145	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	144	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.664	0.701	-5.6	150	-0.04
40	P	Hexachlorocyclopentadiene	0.335	0.340	-1.5	148	-0.04
41	S	2,4,6-Tribromophenol	0.360	0.328	8.9	125	-0.03
42	C	2,4,6-Trichlorophenol	0.443	0.469	-5.9	148	-0.04
43		2,4,5-Trichlorophenol	0.448	0.464	-3.6	142	-0.04
44	S	2-Fluorobiphenyl	1.395	1.381	1.0	145	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.609	1.596	0.8	144	-0.04
46		2-Chloronaphthalene	1.181	1.170	0.9	145	-0.04
47		2-Nitroaniline	0.465	0.531	-14.2	167#	-0.03
48		Acenaphthylene	1.969	1.933	1.8	141	-0.03
49		Dimethylphthalate	1.715	1.663	3.0	141	-0.03
50		2,6-Dinitrotoluene	0.362	0.358	1.1	142	-0.03
51	C	Acenaphthene	1.217	1.204	1.1	145	-0.03
52		3-Nitroaniline	0.339	0.354	-4.4	142	-0.03
53	P	2,4-Dinitrophenol	0.226	0.182	19.5	109	-0.03
54		Dibenzofuran	1.900	1.832	3.6	139	-0.03
55	P	4-Nitrophenol	0.269	0.241	10.4	133	-0.03
56		2,4-Dinitrotoluene	0.532	0.520	2.3	139	-0.03
57		Fluorene	1.647	1.585	3.8	141	-0.03
58		2,3,4,6-Tetrachlorophenol	0.455	0.433	4.8	135	-0.04
59		Diethylphthalate	1.831	1.708	6.7	135	-0.03
60		4-Chlorophenyl-phenylether	0.887	0.868	2.1	138	-0.03
61		4-Nitroaniline	0.342	0.316	7.6	131	-0.03
62		Azobenzene	1.773	1.805	-1.8	149	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	132	-0.03
64		4,6-Dinitro-2-methylphenol	0.139	0.134	3.6	126	-0.03
65	c	n-Nitrosodiphenylamine	0.568	0.581	-2.3	135	-0.03
66		4-Bromophenyl-phenylether	0.243	0.249	-2.5	133	-0.03
67		Hexachlorobenzene	0.285	0.278	2.5	128	-0.03
68		Atrazine	0.244	0.235	3.7	124	-0.03
69	C	Pentachlorophenol	0.152	0.136	10.5	112	-0.03
70		Phenanthrene	1.040	1.011	2.8	131	-0.03
71		Anthracene	1.061	1.028	3.1	130	-0.03
72		Carbazole	0.963	0.900	6.5	125	-0.03
73		Di-n-butylphthalate	1.149	1.078	6.2	127	-0.03
74	C	Fluoranthene	1.253	1.193	4.8	125	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	138	-0.04
76		Benzidine	0.619	0.529	14.5	115	-0.03
77		Pyrene	1.230	1.083	12.0	123	-0.03
78	S	Terphenyl-d14	0.878	0.784	10.7	124	-0.03
79		Butylbenzylphthalate	0.474	0.484	-2.1	150	-0.03
80		Benzo(a)anthracene	1.120	1.143	-2.1	144	-0.04
81		3,3'-Dichlorobenzidine	0.448	0.461	-2.9	140	-0.03
82		Chrysene	1.066	1.088	-2.1	145	-0.04
83		Bis(2-ethylhexyl)phthalate	0.649	0.675	-4.0	156#	-0.04
84	c	Di-n-octyl phthalate	1.065	1.165	-9.4	162#	-0.04
85		Indeno(1,2,3-cd)pyrene	1.180	1.391	-17.9	172#	-0.10
86	I	Perylene-d12	1.000	1.000	0.0	153#	-0.06
87		Benzo(b)fluoranthene	1.136	1.171	-3.1	160#	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.127	1.153	-2.3	157#	-0.06
89 C	Benzo(a)pyrene	1.079	1.134	-5.1	164#	-0.06
90	Dibenzo(a,h)anthracene	1.061	1.108	-4.4	162#	-0.10
91	Benzo(g,h,i)perylene	1.020	1.110	-8.8	170#	-0.11

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