Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG090816\
Data File : BG023964.D
Acq On : 8 Sep 2016 11:30
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 08 14:06:13 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 20:05:47 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	124	0.04
2		1,4-Dioxane	0.470	0.471	-0.2	114	0.02
3		Pyridine	1.412	1.634	-15.7	130	0.02
4		n-Nitrosodimethylamine	0.744	0.912	-22.6	152#	0.03
5	S	2-Fluorophenol	1.139	1.215	-6.7	123	0.03
6		Aniline	2.328	2.478	-6.4	128	0.04
7	S	Phenol-d6	1.754	1.922	-9.6	129	0.03
8		2-Chlorophenol	1.320	1.368	-3.6	123	0.04
9		Benzaldehyde	1.246	1.260	-1.1	118	0.03
10	C	Phenol	1.845	1.951	-5.7	122	0.03
11		bis(2-Chloroethyl)ether	1.451	1.452	-0.1	128	0.04
12		1,3-Dichlorobenzene	1.545	1.501	2.8	121	0.04
13	C	1,4-Dichlorobenzene	1.636	1.565	4.3	123	0.04
14		1,2-Dichlorobenzene	1.536	1.526	0.7	124	0.04
15		Benzyl Alcohol	1.546	1.855	-20.0	138	0.03
16		2,2'-oxybis(1-Chloropropane	1.945	1.999	-2.8	131	0.04
17		2-Methylphenol	1.243	1.297	-4.3	124	0.03
18		Hexachloroethane	0.630	0.687	-9.0	134	0.04
19	P	n-Nitroso-di-n-propylamine	1.549	1.786	-15.3	142	0.03
20		3+4-Methylphenols	1.732	1.926	-11.2	127	0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.03
22		Acetophenone	0.518	0.519	-0.2	127	0.03
23	S	Nitrobenzene-d5	0.448	0.474	-5.8	137	0.04
24		Nitrobenzene	0.482	0.508	-5.4	137	0.04
25		Isophorone	0.869	0.918	-5.6	135	0.03
26	C	2-Nitrophenol	0.183	0.199	-8.7	139	0.04
27		2,4-Dimethylphenol	0.311	0.321	-3.2	122	0.04
28		bis(2-Chloroethoxy)methane	0.439	0.440	-0.2	130	0.04
29	C	2,4-Dichlorophenol	0.330	0.349	-5.8	131	0.03
30		1,2,4-Trichlorobenzene	0.362	0.365	-0.8	132	0.04
31		Naphthalene	1.031	1.009	2.1	129	0.03
32		Benzoic acid	0.253	0.232	8.3	115	0.03
33		4-Chloroaniline	0.430	0.451	-4.9	129	0.03
34	C	Hexachlorobutadiene	0.272	0.289	-6.2	133	0.04
35		Caprolactam	0.116	0.126	-8.6	132	0.01
36	C	4-Chloro-3-methylphenol	0.441	0.496	-12.5	140	0.03
37		2-Methylnaphthalene	0.784	0.803	-2.4	128	0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	135	0.03
39		1,2,4,5-Tetrachlorobenzene	0.664	0.651	2.0	130	0.03
40	P	Hexachlorocyclopentadiene	0.335	0.324	3.3	132	0.03
41	S	2,4,6-Tribromophenol	0.360	0.343	4.7	122	0.03
42	C	2,4,6-Trichlorophenol	0.443	0.444	-0.2	132	0.03
43		2,4,5-Trichlorophenol	0.448	0.446	0.4	128	0.03
44	S	2-Fluorobiphenyl	1.395	1.325	5.0	130	0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.609	1.530	4.9	129	0.03
46		2-Chloronaphthalene	1.181	1.107	6.3	129	0.03
47		2-Nitroaniline	0.465	0.518	-11.4	152#	0.03
48		Acenaphthylene	1.969	1.897	3.7	130	0.03
49		Dimethylphthalate	1.715	1.624	5.3	129	0.03
50		2,6-Dinitrotoluene	0.362	0.357	1.4	132	0.03
51	C	Acenaphthene	1.217	1.179	3.1	133	0.03
52		3-Nitroaniline	0.339	0.347	-2.4	131	0.03
53	P	2,4-Dinitrophenol	0.226	0.188	16.8	105	0.03
54		Dibenzofuran	1.900	1.815	4.5	129	0.02
55	P	4-Nitrophenol	0.269	0.247	8.2	128	0.02
56		2,4-Dinitrotoluene	0.532	0.529	0.6	133	0.03
57		Fluorene	1.647	1.598	3.0	133	0.02
58		2,3,4,6-Tetrachlorophenol	0.455	0.467	-2.6	136	0.02
59		Diethylphthalate	1.831	1.745	4.7	129	0.02
60		4-Chlorophenyl-phenylether	0.887	0.863	2.7	128	0.03
61		4-Nitroaniline	0.342	0.334	2.3	129	0.03
62		Azobenzene	1.773	1.807	-1.9	139	0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.02
64		4,6-Dinitro-2-methylphenol	0.139	0.140	-0.7	129	0.03
65	c	n-Nitrosodiphenylamine	0.568	0.568	0.0	129	0.02
66		4-Bromophenyl-phenylether	0.243	0.248	-2.1	130	0.03
67		Hexachlorobenzene	0.285	0.274	3.9	123	0.02
68		Atrazine	0.244	0.240	1.6	124	0.01
69	C	Pentachlorophenol	0.152	0.140	7.9	113	0.02
70		Phenanthrene	1.040	1.031	0.9	131	0.03
71		Anthracene	1.061	1.019	4.0	126	0.02
72		Carbazole	0.963	0.908	5.7	124	0.01
73		Di-n-butylphthalate	1.149	1.104	3.9	127	0.01
74	C	Fluoranthene	1.253	1.193	4.8	122	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	125	0.01
76		Benzidine	0.619	0.559	9.7	110	0.01
77		Pyrene	1.230	1.184	3.7	121	0.01
78	S	Terphenyl-d14	0.878	0.843	4.0	121	0.01
79		Butylbenzylphthalate	0.474	0.479	-1.1	134	0.00
80		Benzo(a)anthracene	1.120	1.156	-3.2	132	0.02
81		3,3'-Dichlorobenzidine	0.448	0.461	-2.9	127	0.02
82		Chrysene	1.066	1.092	-2.4	132	0.02
83		Bis(2-ethylhexyl)phthalate	0.649	0.670	-3.2	140	0.00
84	c	Di-n-octyl phthalate	1.065	1.141	-7.1	143	0.01
85		Indeno(1,2,3-cd)pyrene	1.180	1.366	-15.8	153#	0.04
86	I	Perylene-d12	1.000	1.000	0.0	136	0.03
87		Benzo(b)fluoranthene	1.136	1.172	-3.2	142	0.03

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88	Benzo(k)fluoranthene	1.127	1.137	-0.9	137	0.03
89 C	Benzo(a)pyrene	1.079	1.121	-3.9	144	0.03
90	Dibenzo(a,h)anthracene	1.061	1.096	-3.3	142	0.03
91	Benzo(g,h,i)perylene	1.020	1.113	-9.1	151#	0.03

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