Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG103116\
Data File : BG024573.D
Acq On : 1 Nov 2016 1:51
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 01 12:01:16 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 01 11:50:40 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#	0.00
2		1,4-Dioxane	0.499	0.503	-0.8	161#	0.00
3		Pyridine	1.493	1.421	4.8	151#	0.00
4		n-Nitrosodimethylamine	0.773	0.714	7.6	143	0.00
5	S	2-Fluorophenol	1.220	1.213	0.6	160#	0.00
6		Aniline	2.121	2.146	-1.2	158#	0.00
7	S	Phenol-d6	1.674	1.734	-3.6	163#	0.00
8		2-Chlorophenol	1.241	1.240	0.1	162#	0.00
9		Benzaldehyde	1.034	1.019	1.5	156#	0.00
10	C	Phenol	1.725	1.770	-2.6	163#	0.00
11		bis(2-Chloroethyl)ether	1.296	1.345	-3.8	168#	0.00
12		1,3-Dichlorobenzene	1.536	1.523	0.8	162#	0.00
13	C	1,4-Dichlorobenzene	1.517	1.531	-0.9	162#	0.00
14		1,2-Dichlorobenzene	1.459	1.458	0.1	161#	0.00
15		Benzyl Alcohol	1.557	1.540	1.1	156#	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.296	4.5	152#	0.00
17		2-Methylphenol	1.143	1.157	-1.2	162#	0.00
18		Hexachloroethane	0.583	0.575	1.4	164#	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.191	3.5	149	0.00
20		3+4-Methylphenols	1.535	1.579	-2.9	159#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	150#	0.00
22		Acetophenone	0.514	0.501	2.5	155#	0.00
23	S	Nitrobenzene-d5	0.439	0.442	-0.7	161#	0.00
24		Nitrobenzene	0.489	0.485	0.8	157#	0.00
25		Isophorone	0.817	0.785	3.9	150	0.00
26	C	2-Nitrophenol	0.181	0.197	-8.8	173#	0.00
27		2,4-Dimethylphenol	0.318	0.329	-3.5	160#	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.421	0.5	162#	0.00
29	C	2,4-Dichlorophenol	0.333	0.343	-3.0	161#	0.00
30		1,2,4-Trichlorobenzene	0.395	0.391	1.0	159#	0.00
31		Naphthalene	0.998	0.969	2.9	154#	0.00
32		Benzoic acid	0.264	0.238	9.8	143	0.02
33		4-Chloroaniline	0.433	0.436	-0.7	152#	0.00
34	C	Hexachlorobutadiene	0.309	0.311	-0.6	166#	0.00
35		Caprolactam	0.122	0.116	4.9	147	0.00
36	C	4-Chloro-3-methylphenol	0.402	0.399	0.7	152#	0.00
37		2-Methylnaphthalene	0.728	0.718	1.4	157#	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	137	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.695	-3.0	155#	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.376	1.1	144	0.00
41	S	2,4,6-Tribromophenol	0.299	0.316	-5.7	145	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.439	-8.4	160#	0.00
43		2,4,5-Trichlorophenol	0.438	0.459	-4.8	152#	0.00
44	S	2-Fluorobiphenyl	1.203	1.195	0.7	147	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.390	-0.1	148	0.00
46		2-Chloronaphthalene	1.057	1.082	-2.4	152#	0.00
47		2-Nitroaniline	0.433	0.450	-3.9	141	0.00
48		Acenaphthylene	1.621	1.637	-1.0	145	0.00
49		Dimethylphthalate	1.420	1.421	-0.1	141	0.00
50		2,6-Dinitrotoluene	0.303	0.328	-8.3	147	0.00
51	C	Acenaphthene	1.208	1.120	7.3	134	0.00
52		3-Nitroaniline	0.314	0.327	-4.1	145	0.00
53	P	2,4-Dinitrophenol	0.220	0.238	-8.2	145	0.00
54		Dibenzofuran	1.670	1.652	1.1	142	0.00
55	P	4-Nitrophenol	0.273	0.262	4.0	134	0.00
56		2,4-Dinitrotoluene	0.440	0.471	-7.0	142	0.00
57		Fluorene	1.385	1.367	1.3	139	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.467	-2.9	142	0.00
59		Diethylphthalate	1.489	1.435	3.6	135	0.00
60		4-Chlorophenyl-phenylether	0.777	0.785	-1.0	147	0.00
61		4-Nitroaniline	0.343	0.355	-3.5	144	0.00
62		Azobenzene	1.485	1.402	5.6	133	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.136	1.4	131	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.557	-1.8	139	0.00
66		4-Bromophenyl-phenylether	0.243	0.252	-3.7	143	0.00
67		Hexachlorobenzene	0.268	0.279	-4.1	140	0.00
68		Atrazine	0.235	0.224	4.7	127	0.00
69	C	Pentachlorophenol	0.177	0.167	5.6	122	0.00
70		Phenanthrene	1.019	0.994	2.5	134	0.00
71		Anthracene	1.028	1.003	2.4	133	0.00
72		Carbazole	0.938	0.903	3.7	133	0.00
73		Di-n-butylphthalate	1.081	1.036	4.2	130	0.00
74	C	Fluoranthene	1.289	1.239	3.9	128	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	128	0.00
76		Benzidine	0.471	0.487	-3.4	139	0.00
77		Pyrene	0.978	0.950	2.9	128	0.00
78	S	Terphenyl-d14	0.669	0.600	10.3	122	0.00
79		Butylbenzylphthalate	0.408	0.407	0.2	135	0.00
80		Benzo(a)anthracene	1.099	1.057	3.8	133	0.00
81		3,3'-Dichlorobenzidine	0.443	0.428	3.4	131	0.00
82		Chrysene	1.046	1.007	3.7	131	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.566	2.9	132	0.00
84	c	Di-n-octyl phthalate	0.968	0.971	-0.3	135	0.00
85		Indeno(1,2,3-cd)pyrene	1.444	1.429	1.0	140	0.00
86	I	Perylene-d12	1.000	1.000	0.0	130	0.00
87		Benzo(b)fluoranthene	1.081	1.054	2.5	137	0.00

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88	Benzo(k)fluoranthene	1.044	0.996	4.6	132	0.00
89 C	Benzo(a)pyrene	1.056	1.029	2.6	135	0.00
90	Dibenzo(a,h)anthracene	1.159	1.104	4.7	138	0.00
91	Benzo(g,h,i)perylene	1.158	1.092	5.7	138	0.00

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