Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025496.D
Acq On : 13 Jan 2017 00:14
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 13 02:24:54 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 11:26:22 2017
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	134	0.00
2		1,4-Dioxane	0.469	0.489	-4.3	139	0.00
3		Pyridine	1.361	1.455	-6.9	131	0.00
4		n-Nitrosodimethylamine	0.808	0.914	-13.1	140	0.00
5	S	2-Fluorophenol	1.100	1.205	-9.5	143	0.00
6		Aniline	2.100	2.274	-8.3	138	0.00
7	S	Phenol-d6	1.550	1.727	-11.4	139	0.00
8		2-Chlorophenol	1.241	1.341	-8.1	139	0.00
9		Benzaldehyde	1.134	1.162	-2.5	136	0.00
10	C	Phenol	1.718	1.855	-8.0	136	0.00
11		bis(2-Chloroethyl)ether	1.324	1.419	-7.2	139	0.00
12		1,3-Dichlorobenzene	1.502	1.602	-6.7	136	0.00
13	C	1,4-Dichlorobenzene	1.557	1.597	-2.6	132	0.00
14		1,2-Dichlorobenzene	1.486	1.539	-3.6	135	0.00
15		Benzyl Alcohol	1.479	1.639	-10.8	134	0.00
16		2,2'-oxybis(1-Chloropropane	2.451	2.774	-13.2	145	0.00
17		2-Methylphenol	1.128	1.245	-10.4	140	0.00
18		Hexachloroethane	0.598	0.661	-10.5	136	0.00
19	P	n-Nitroso-di-n-propylamine	1.309	1.399	-6.9	136	0.00
20		3+4-Methylphenols	1.561	1.716	-9.9	138	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	143	0.00
22		Acetophenone	0.556	0.557	-0.2	132	0.00
23	S	Nitrobenzene-d5	0.453	0.469	-3.5	136	0.00
24		Nitrobenzene	0.497	0.507	-2.0	135	0.00
25		Isophorone	0.820	0.851	-3.8	136	0.00
26	C	2-Nitrophenol	0.190	0.203	-6.8	139	0.00
27		2,4-Dimethylphenol	0.312	0.298	4.5	125	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.427	-0.2	139	0.00
29	C	2,4-Dichlorophenol	0.334	0.348	-4.2	137	0.00
30		1,2,4-Trichlorobenzene	0.404	0.401	0.7	135	0.00
31		Naphthalene	0.999	1.019	-2.0	135	0.00
32		Benzoic acid	0.251	0.266	-6.0	137	0.01
33		4-Chloroaniline	0.420	0.444	-5.7	135	0.00
34	C	Hexachlorobutadiene	0.329	0.334	-1.5	136	0.00
35		Caprolactam	0.126	0.122	3.2	129	0.01
36	C	4-Chloro-3-methylphenol	0.412	0.434	-5.3	137	0.00
37		2-Methylnaphthalene	0.740	0.763	-3.1	136	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	137	0.00
39		1,2,4,5-Tetrachlorobenzene	0.660	0.688	-4.2	136	0.00
40	P	Hexachlorocyclopentadiene	0.192	0.222	-15.6	139	0.00
41	S	2,4,6-Tribromophenol	0.322	0.321	0.3	127	0.00
42	C	2,4,6-Trichlorophenol	0.416	0.417	-0.2	130	0.00
43		2,4,5-Trichlorophenol	0.435	0.433	0.5	129	0.00
44	S	2-Fluorobiphenyl	1.235	1.245	-0.8	134	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.345	1.376	-2.3	134	0.00
46		2-Chloronaphthalene	1.051	1.074	-2.2	136	0.00
47		2-Nitroaniline	0.444	0.485	-9.2	138	0.00
48		Acenaphthylene	1.629	1.662	-2.0	135	0.00
49		Dimethylphthalate	1.458	1.475	-1.2	130	0.00
50		2,6-Dinitrotoluene	0.319	0.329	-3.1	134	0.00
51	C	Acenaphthene	1.065	1.086	-2.0	133	0.00
52		3-Nitroaniline	0.306	0.320	-4.6	133	0.00
53	P	2,4-Dinitrophenol	0.183	0.189	-3.3	126	0.00
54		Dibenzofuran	1.684	1.715	-1.8	132	0.00
55	P	4-Nitrophenol	0.229	0.237	-3.5	128	0.00
56		2,4-Dinitrotoluene	0.464	0.493	-6.2	136	0.00
57		Fluorene	1.410	1.442	-2.3	134	0.00
58		2,3,4,6-Tetrachlorophenol	0.466	0.461	1.1	129	0.00
59		Diethylphthalate	1.529	1.547	-1.2	132	0.00
60		4-Chlorophenyl-phenylether	0.799	0.827	-3.5	135	0.00
61		4-Nitroaniline	0.308	0.335	-8.8	128	0.00
62		Azobenzene	1.475	1.566	-6.2	137	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.159	-14.4	132	0.00
65	c	n-Nitrosodiphenylamine	0.541	0.569	-5.2	132	0.00
66		4-Bromophenyl-phenylether	0.247	0.260	-5.3	131	0.00
67		Hexachlorobenzene	0.283	0.293	-3.5	132	0.00
68		Atrazine	0.237	0.187	21.1	99	0.00
69	C	Pentachlorophenol	0.147	0.140	4.8	114	0.00
70		Phenanthrene	1.031	1.064	-3.2	130	0.00
71		Anthracene	1.047	1.082	-3.3	132	0.00
72		Carbazole	0.937	0.953	-1.7	129	0.00
73		Di-n-butylphthalate	1.152	1.169	-1.5	129	0.00
74	C	Fluoranthene	1.361	1.394	-2.4	131	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	142	0.00
76		Benzidine	0.499	0.425	14.8	114	0.00
77		Pyrene	1.045	1.056	-1.1	131	0.00
78	S	Terphenyl-d14	0.754	0.723	4.1	127	0.00
79		Butylbenzylphthalate	0.420	0.437	-4.0	141	0.00
80		Benzo(a)anthracene	1.116	1.145	-2.6	136	0.00
81		3,3'-Dichlorobenzidine	0.433	0.462	-6.7	141	0.00
82		Chrysene	1.070	1.100	-2.8	138	0.00
83		Bis(2-ethylhexyl)phthalate	0.584	0.615	-5.3	142	0.00
84	c	Di-n-octyl phthalate	0.970	1.054	-8.7	144	0.00
85		Indeno(1,2,3-cd)pyrene	1.360	1.488	-9.4	145	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	145	0.00
87		Benzo(b)fluoranthene	1.091	1.124	-3.0	142	0.00

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88	Benzo(k)fluoranthene	1.054	1.093	-3.7	140	0.00
89 C	Benzo(a)pyrene	1.054	1.104	-4.7	143	0.00
90	Dibenzo(a,h)anthracene	1.117	1.180	-5.6	143	0.00
91	Benzo(g,h,i)perylene	1.110	1.172	-5.6	146	-0.01

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