Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_G\DATA\BG072015\
Data File : BG017982.D
Acq On : 20 Jul 2015 10:53
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 20 16:51:43 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 01:50:30 2015
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	138	0.00
2		1,4-Dioxane	0.504	0.424	15.9	123	0.00
3		Pyridine	1.347	1.192	11.5	121	0.00
4		n-Nitrosodimethylamine	0.511	0.441	13.7	124	0.00
5	S	2-Fluorophenol	1.146	1.093	4.6	134	0.00
6		Aniline	2.003	1.880	6.1	133	0.00
7	S	Phenol-d6	1.502	1.427	5.0	135	0.00
8		2-Chlorophenol	1.352	1.313	2.9	139	0.00
9		Benzaldehyde	0.920	0.842	8.5	130	0.00
10	C	Phenol	1.603	1.509	5.9	135	0.00
11		bis(2-Chloroethyl)ether	1.254	1.148	8.5	131	0.00
12		1,3-Dichlorobenzene	1.484	1.426	3.9	137	0.00
13	C	1,4-Dichlorobenzene	1.521	1.469	3.4	138	0.00
14		1,2-Dichlorobenzene	1.454	1.371	5.7	137	0.00
15		Benzyl Alcohol	1.111	1.066	4.1	137	0.00
16		2,2'-oxybis(1-Chloropropane	1.511	1.269	16.0	123	0.00
17		2-Methylphenol	1.118	1.084	3.0	138	0.00
18		Hexachloroethane	0.575	0.541	5.9	135	0.00
19	P	n-Nitroso-di-n-propylamine	1.097	0.984	10.3	128	0.00
20		3+4-Methylphenols	1.513	1.479	2.2	137	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
22		Acetophenone	0.432	0.410	5.1	136	0.00
23	S	Nitrobenzene-d5	0.309	0.293	5.2	134	0.00
24		Nitrobenzene	0.329	0.311	5.5	131	0.00
25		Isophorone	0.648	0.603	6.9	130	0.00
26	C	2-Nitrophenol	0.157	0.178	-13.4	163#	0.00
27		2,4-Dimethylphenol	0.286	0.285	0.3	142	0.00
28		bis(2-Chloroethoxy)methane	0.361	0.338	6.4	131	0.00
29	C	2,4-Dichlorophenol	0.254	0.278	-9.4	149	0.00
30		1,2,4-Trichlorobenzene	0.302	0.301	0.3	144	0.00
31		Naphthalene	0.979	0.924	5.6	134	0.00
32		Benzoic acid	0.110	0.151	-37.3#	218#	0.01
33		4-Chloroaniline	0.436	0.433	0.7	138	0.00
34	C	Hexachlorobutadiene	0.170	0.176	-3.5	145	0.00
35		Caprolactam	0.129	0.128	0.8	141	0.01
36	C	4-Chloro-3-methylphenol	0.289	0.296	-2.4	142	0.00
37		2-Methylnaphthalene	0.703	0.679	3.4	138	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	139	0.00
39		1,2,4,5-Tetrachlorobenzene	0.536	0.537	-0.2	144	0.00
40	P	Hexachlorocyclopentadiene	0.292	0.298	-2.1	149	0.00
41	S	2,4,6-Tribromophenol	0.203	0.224	-10.3	162#	0.00
42	C	2,4,6-Trichlorophenol	0.349	0.372	-6.6	156#	0.00
43		2,4,5-Trichlorophenol	0.363	0.399	-9.9	156#	0.00
44	S	2-Fluorobiphenyl	1.154	1.079	6.5	134	0.00

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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)
45		1,1'-Biphenyl	1.400	1.328	5.1	136	0.00
46		2-Chloronaphthalene	1.135	1.083	4.6	139	0.00
47		2-Nitroaniline	0.305	0.293	3.9	139	0.00
48		Acenaphthylene	1.892	1.786	5.6	134	0.00
49		Dimethylphthalate	1.389	1.320	5.0	135	0.00
50		2,6-Dinitrotoluene	0.315	0.327	-3.8	150	0.00
51	C	Acenaphthene	1.146	1.075	6.2	136	0.00
52		3-Nitroaniline	0.354	0.366	-3.4	144	0.00
53	P	2,4-Dinitrophenol	0.121	0.128	-5.8	159#	0.00
54		Dibenzofuran	1.654	1.551	6.2	136	0.00
55	P	4-Nitrophenol	0.280	0.286	-2.1	147	0.00
56		2,4-Dinitrotoluene	0.422	0.438	-3.8	149	0.00
57		Fluorene	1.421	1.337	5.9	137	0.00
58		2,3,4,6-Tetrachlorophenol	0.313	0.328	-4.8	151#	0.00
59		Diethylphthalate	1.439	1.350	6.2	135	0.00
60		4-Chlorophenyl-phenylether	0.697	0.676	3.0	140	0.00
61		4-Nitroaniline	0.415	0.402	3.1	141	0.00
62		Azobenzene	1.330	1.160	12.8	125	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
64		4,6-Dinitro-2-methylphenol	0.103	0.115	-11.7	166#	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.581	3.3	137	0.00
66		4-Bromophenyl-phenylether	0.200	0.204	-2.0	146	0.00
67		Hexachlorobenzene	0.222	0.223	-0.5	147	0.00
68		Atrazine	0.198	0.195	1.5	138	0.00
69	C	Pentachlorophenol	0.115	0.132	-14.8	164#	0.00
70		Phenanthrene	1.032	0.945	8.4	133	0.00
71		Anthracene	1.004	0.950	5.4	133	0.00
72		Carbazole	0.948	0.882	7.0	133	0.00
73		Di-n-butylphthalate	1.162	1.092	6.0	131	0.00
74	C	Fluoranthene	1.129	1.049	7.1	132	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	136	0.00
76		Benzidine	0.568	0.570	-0.4	139	0.00
77		Pyrene	1.115	1.025	8.1	129	0.00
78	S	Terphenyl-d14	0.794	0.720	9.3	128	0.00
79		Butylbenzylphthalate	0.500	0.511	-2.2	137	0.00
80		Benzo(a)anthracene	1.056	1.000	5.3	133	0.00
81		3,3'-Dichlorobenzidine	0.383	0.419	-9.4	150	0.00
82		Chrysene	1.006	0.946	6.0	133	0.00
83		Bis(2-ethylhexyl)phthalate	0.688	0.687	0.1	134	0.00
84	c	Di-n-octyl phthalate	1.084	1.123	-3.6	137	0.00
85		Indeno(1,2,3-cd)pyrene	1.217	1.219	-0.2	141	0.00
86	I	Perylene-d12	1.000	1.000	0.0	143	0.00
87		Benzo(b)fluoranthene	1.094	1.049	4.1	140	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.093	1.014	7.2	133	0.00
89 C	Benzo(a)pyrene	1.020	0.974	4.5	138	0.00
90	Dibenzo(a,h)anthracene	1.057	1.039	1.7	144	0.00
91	Benzo(g,h,i)perylene	1.019	0.991	2.7	143	0.00

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