Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG052015\
Data File : BG017228.D
Acq On : 21 May 2015 00:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 21 03:39:45 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 21 02:15:39 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	122	-0.01
2		1,4-Dioxane	0.445	0.431	3.1	119	-0.02
3		Pyridine	1.346	1.317	2.2	114	-0.01
4		n-Nitrosodimethylamine	1.021	0.943	7.6	108	-0.02
5	S	2-Fluorophenol	1.127	1.098	2.6	113	-0.02
6		Aniline	2.189	2.089	4.6	110	-0.01
7	S	Phenol-d6	1.582	1.537	2.8	112	-0.01
8		2-Chlorophenol	1.349	1.273	5.6	110	-0.01
9		Benzaldehyde	1.041	0.967	7.1	111	-0.02
10	C	Phenol	1.678	1.659	1.1	116	-0.01
11		bis(2-Chloroethyl)ether	1.258	1.222	2.9	111	-0.01
12		1,3-Dichlorobenzene	1.506	1.388	7.8	111	-0.01
13	C	1,4-Dichlorobenzene	1.520	1.396	8.2	109	-0.01
14		1,2-Dichlorobenzene	1.452	1.326	8.7	108	-0.01
15		Benzyl Alcohol	1.537	1.418	7.7	108	-0.01
16		2,2'-oxybis(1-Chloropropane	2.040	2.018	1.1	118	-0.01
17		2-Methylphenol	1.216	1.145	5.8	113	-0.02
18		Hexachloroethane	0.635	0.605	4.7	113	-0.02
19	P	n-Nitroso-di-n-propylamine	1.378	1.249	9.4	106	-0.01
20		3+4-Methylphenols	1.691	1.548	8.5	106	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	113	-0.01
22		Acetophenone	0.485	0.455	6.2	105	-0.02
23	S	Nitrobenzene-d5	0.428	0.413	3.5	106	-0.01
24		Nitrobenzene	0.419	0.406	3.1	109	-0.01
25		Isophorone	0.790	0.741	6.2	103	-0.02
26	C	2-Nitrophenol	0.187	0.174	7.0	103	-0.02
27		2,4-Dimethylphenol	0.308	0.293	4.9	108	-0.02
28		bis(2-Chloroethoxy)methane	0.384	0.368	4.2	106	-0.01
29	C	2,4-Dichlorophenol	0.291	0.286	1.7	107	-0.01
30		1,2,4-Trichlorobenzene	0.330	0.299	9.4	100	-0.02
31		Naphthalene	0.979	0.932	4.8	105	-0.02
32		Benzoic acid	0.209	0.185	11.5	97	0.00
33		4-Chloroaniline	0.463	0.435	6.0	102	-0.02
34	C	Hexachlorobutadiene	0.209	0.194	7.2	102	-0.02
35		Caprolactam	0.146	0.144	1.4	114	0.00
36	C	4-Chloro-3-methylphenol	0.368	0.365	0.8	109	-0.01
37		2-Methylnaphthalene	0.776	0.697	10.2	103	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	106	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.566	0.562	0.7	102	-0.01
40	P	Hexachlorocyclopentadiene	0.345	0.297	13.9	89	-0.01
41	S	2,4,6-Tribromophenol	0.242	0.243	-0.4	104	0.00
42	C	2,4,6-Trichlorophenol	0.401	0.403	-0.5	101	-0.01
43		2,4,5-Trichlorophenol	0.414	0.436	-5.3	109	-0.01
44	S	2-Fluorobiphenyl	1.363	1.309	4.0	99	-0.01

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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	1.441	1.373	4.7	99	-0.01
46		2-Chloronaphthalene	1.141	1.096	3.9	100	-0.01
47		2-Nitroaniline	0.444	0.467	-5.2	110	-0.01
48		Acenaphthylene	1.953	1.936	0.9	101	-0.01
49		Dimethylphthalate	1.564	1.480	5.4	98	-0.01
50		2,6-Dinitrotoluene	0.347	0.350	-0.9	102	-0.01
51	C	Acenaphthene	1.154	1.134	1.7	103	-0.01
52		3-Nitroaniline	0.384	0.404	-5.2	108	-0.01
53	P	2,4-Dinitrophenol	0.201	0.159	20.9	77	0.00
54		Dibenzofuran	1.715	1.648	3.9	100	-0.01
55	P	4-Nitrophenol	0.310	0.305	1.6	104	-0.01
56		2,4-Dinitrotoluene	0.483	0.513	-6.2	109	0.00
57		Fluorene	1.518	1.473	3.0	99	-0.01
58		2,3,4,6-Tetrachlorophenol	0.380	0.400	-5.3	107	-0.01
59		Diethylphthalate	1.680	1.603	4.6	101	-0.01
60		4-Chlorophenyl-phenylether	0.780	0.745	4.5	98	-0.01
61		4-Nitroaniline	0.429	0.474	-10.5	113	0.00
62		Azobenzene	1.605	1.605	0.0	104	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	-0.01
64		4,6-Dinitro-2-methylphenol	0.137	0.128	6.6	100	0.00
65	c	n-Nitrosodiphenylamine	0.614	0.572	6.8	103	0.00
66		4-Bromophenyl-phenylether	0.218	0.199	8.7	102	0.00
67		Hexachlorobenzene	0.230	0.216	6.1	104	0.00
68		Atrazine	0.225	0.211	6.2	102	-0.01
69	C	Pentachlorophenol	0.144	0.109	24.3#	82	0.00
70		Phenanthrene	1.031	1.007	2.3	107	-0.01
71		Anthracene	1.045	1.015	2.9	108	0.00
72		Carbazole	0.960	0.975	-1.6	112	0.00
73		Di-n-butylphthalate	1.285	1.308	-1.8	113	0.00
74	C	Fluoranthene	1.239	1.259	-1.6	113	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	137	0.00
76		Benzidine	0.653	0.485	25.7#	97	-0.01
77		Pyrene	1.227	1.067	13.0	115	0.00
78	S	Terphenyl-d14	0.960	0.853	11.1	118	-0.01
79		Butylbenzylphthalate	0.558	0.526	5.7	125	0.00
80		Benzo(a)anthracene	1.146	1.075	6.2	124	-0.01
81		3,3'-Dichlorobenzidine	0.444	0.407	8.3	122	0.00
82		Chrysene	1.068	0.997	6.6	124	0.00
83		Bis(2-ethylhexyl)phthalate	0.793	0.747	5.8	123	-0.01
84	c	Di-n-octyl phthalate	1.320	1.275	3.4	128	0.00
85		Indeno(1,2,3-cd)pyrene	1.277	1.238	3.1	128	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	135	-0.01
87		Benzo(b)fluoranthene	1.166	1.058	9.3	123	0.00

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88	Benzo(k)fluoranthene	1.107	1.077	2.7	127	0.00
89 C	Benzo(a)pyrene	1.071	1.020	4.8	127	0.00
90	Dibenzo(a,h)anthracene	1.099	1.049	4.5	128	-0.02
91	Benzo(g,h,i)perylene	1.035	0.985	4.8	127	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 1