Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021816\
Data File : BG020950.D
Acq On : 19 Feb 2016 00:25
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 19 02:49:25 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 11 15:26:53 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	100	0.00
2		1,4-Dioxane	0.432	0.441	-2.1	105	-0.01
3		Pyridine	1.277	1.338	-4.8	101	-0.01
4		n-Nitrosodimethylamine	0.334	0.359	-7.5	105	-0.01
5	S	2-Fluorophenol	1.188	1.237	-4.1	104	-0.01
6		Aniline	2.164	2.298	-6.2	105	-0.01
7	S	Phenol-d6	1.736	1.860	-7.1	106	0.00
8		2-Chlorophenol	1.484	1.582	-6.6	107	-0.01
9		Benzaldehyde	0.869	0.883	-1.6	101	-0.01
10	C	Phenol	1.837	1.954	-6.4	106	0.00
11		bis(2-Chloroethyl)ether	1.467	1.548	-5.5	103	-0.01
12		1,3-Dichlorobenzene	1.569	1.653	-5.4	106	-0.01
13	C	1,4-Dichlorobenzene	1.590	1.678	-5.5	105	-0.01
14		1,2-Dichlorobenzene	1.547	1.644	-6.3	105	-0.01
15		Benzyl Alcohol	1.094	1.172	-7.1	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.530	1.567	-2.4	102	-0.01
17		2-Methylphenol	1.323	1.418	-7.2	107	0.00
18		Hexachloroethane	0.547	0.585	-6.9	106	-0.01
19	P	n-Nitroso-di-n-propylamine	1.120	1.210	-8.0	108	-0.01
20		3+4-Methylphenols	1.844	2.000	-8.5	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	-0.01
22		Acetophenone	0.482	0.494	-2.5	108	-0.01
23	S	Nitrobenzene-d5	0.304	0.305	-0.3	106	-0.01
24		Nitrobenzene	0.317	0.321	-1.3	108	0.00
25		Isophorone	0.641	0.668	-4.2	110	-0.01
26	C	2-Nitrophenol	0.167	0.184	-10.2	113	-0.01
27		2,4-Dimethylphenol	0.320	0.320	0.0	104	-0.01
28		bis(2-Chloroethoxy)methane	0.395	0.406	-2.8	110	-0.01
29	C	2,4-Dichlorophenol	0.287	0.301	-4.9	113	-0.01
30		1,2,4-Trichlorobenzene	0.289	0.293	-1.4	108	0.00
31		Naphthalene	1.035	1.044	-0.9	108	-0.01
32		Benzoic acid	0.256	0.263	-2.7	112	0.00
33		4-Chloroaniline	0.445	0.465	-4.5	112	0.00
34	C	Hexachlorobutadiene	0.147	0.152	-3.4	110	-0.01
35		Caprolactam	0.110	0.118	-7.3	116	0.00
36	C	4-Chloro-3-methylphenol	0.328	0.346	-5.5	113	0.00
37		2-Methylnaphthalene	0.726	0.744	-2.5	111	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	113	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.587	0.586	0.2	112	-0.01
40	P	Hexachlorocyclopentadiene	0.258	0.277	-7.4	117	-0.02
41	S	2,4,6-Tribromophenol	0.197	0.216	-9.6	124	-0.01
42	C	2,4,6-Trichlorophenol	0.394	0.412	-4.6	115	-0.01
43		2,4,5-Trichlorophenol	0.414	0.433	-4.6	117	-0.01
44	S	2-Fluorobiphenyl	1.424	1.447	-1.6	113	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.764	1.783	-1.1	113	-0.01
46		2-Chloronaphthalene	1.263	1.270	-0.6	114	-0.01
47		2-Nitroaniline	0.303	0.313	-3.3	115	-0.01
48		Acenaphthylene	2.196	2.244	-2.2	115	-0.01
49		Dimethylphthalate	1.564	1.646	-5.2	118	-0.01
50		2,6-Dinitrotoluene	0.335	0.360	-7.5	120	-0.01
51	C	Acenaphthene	1.332	1.365	-2.5	115	-0.01
52		3-Nitroaniline	0.389	0.416	-6.9	119	-0.01
53	P	2,4-Dinitrophenol	0.116	0.139	-19.8	135	-0.01
54		Dibenzofuran	1.762	1.796	-1.9	116	-0.01
55	P	4-Nitrophenol	0.294	0.321	-9.2	119	0.00
56		2,4-Dinitrotoluene	0.433	0.483	-11.5	122	-0.01
57		Fluorene	1.622	1.686	-3.9	118	-0.01
58		2,3,4,6-Tetrachlorophenol	0.316	0.339	-7.3	121	-0.01
59		Diethylphthalate	1.608	1.700	-5.7	120	-0.01
60		4-Chlorophenyl-phenylether	0.720	0.756	-5.0	121	-0.01
61		4-Nitroaniline	0.391	0.418	-6.9	119	-0.01
62		Azobenzene	1.264	1.266	-0.2	113	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	121	-0.01
64		4,6-Dinitro-2-methylphenol	0.094	0.106	-12.8	134	-0.01
65	c	n-Nitrosodiphenylamine	0.646	0.644	0.3	118	-0.01
66		4-Bromophenyl-phenylether	0.211	0.216	-2.4	121	-0.02
67		Hexachlorobenzene	0.225	0.229	-1.8	123	-0.01
68		Atrazine	0.196	0.193	1.5	119	-0.01
69	C	Pentachlorophenol	0.129	0.130	-0.8	121	-0.01
70		Phenanthrene	1.139	1.149	-0.9	122	-0.02
71		Anthracene	1.156	1.162	-0.5	121	-0.01
72		Carbazole	1.037	1.020	1.6	118	-0.01
73		Di-n-butylphthalate	1.330	1.306	1.8	117	-0.01
74	C	Fluoranthene	1.230	1.216	1.1	119	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	118	-0.02
76		Benzidine	0.647	0.584	9.7	103	-0.01
77		Pyrene	1.319	1.337	-1.4	120	-0.02
78	S	Terphenyl-d14	0.857	0.870	-1.5	119	-0.02
79		Butylbenzylphthalate	0.632	0.638	-0.9	117	-0.02
80		Benzo(a)anthracene	1.194	1.223	-2.4	122	-0.02
81		3,3'-Dichlorobenzidine	0.443	0.450	-1.6	119	-0.02
82		Chrysene	1.123	1.141	-1.6	120	-0.02
83		Bis(2-ethylhexyl)phthalate	0.936	0.946	-1.1	118	-0.02
84	c	Di-n-octyl phthalate	1.537	1.573	-2.3	119	-0.04
85		Indeno(1,2,3-cd)pyrene	1.291	1.360	-5.3	123	-0.06
86	I	Perylene-d12	1.000	1.000	0.0	122	-0.04
87		Benzo(b)fluoranthene	1.224	1.220	0.3	120	-0.03

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.199	1.227	-2.3	125	-0.03
89 C	Benzo(a)pyrene	1.166	1.182	-1.4	123	-0.04
90	Dibenzo(a,h)anthracene	1.134	1.150	-1.4	123	-0.06
91	Benzo(g,h,i)perylene	1.126	1.153	-2.4	124	-0.07

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