Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021916\
Data File : BG020986.D
Acq On : 20 Feb 2016 00:11
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 20 01:32:58 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	109	0.00
2		1,4-Dioxane	0.432	0.394	8.8	102	-0.01
3		Pyridine	1.277	1.244	2.6	102	-0.01
4		n-Nitrosodimethylamine	0.334	0.317	5.1	101	-0.01
5	S	2-Fluorophenol	1.188	1.172	1.3	107	0.00
6		Aniline	2.164	2.209	-2.1	110	0.00
7	S	Phenol-d6	1.736	1.832	-5.5	113	0.00
8		2-Chlorophenol	1.484	1.546	-4.2	113	-0.01
9		Benzaldehyde	0.869	0.822	5.4	102	0.00
10	C	Phenol	1.837	1.918	-4.4	113	0.00
11		bis(2-Chloroethyl)ether	1.467	1.458	0.6	106	-0.01
12		1,3-Dichlorobenzene	1.569	1.574	-0.3	110	-0.01
13	C	1,4-Dichlorobenzene	1.590	1.636	-2.9	111	-0.01
14		1,2-Dichlorobenzene	1.547	1.591	-2.8	111	-0.01
15		Benzyl Alcohol	1.094	1.124	-2.7	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.530	1.473	3.7	104	-0.01
17		2-Methylphenol	1.323	1.384	-4.6	113	0.00
18		Hexachloroethane	0.547	0.564	-3.1	111	-0.01
19	P	n-Nitroso-di-n-propylamine	1.120	1.153	-2.9	112	-0.01
20		3+4-Methylphenols	1.844	1.974	-7.0	117	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	117	-0.01
22		Acetophenone	0.482	0.472	2.1	113	0.00
23	S	Nitrobenzene-d5	0.304	0.300	1.3	113	0.00
24		Nitrobenzene	0.317	0.310	2.2	114	0.00
25		Isophorone	0.641	0.639	0.3	115	0.00
26	C	2-Nitrophenol	0.167	0.186	-11.4	123	-0.01
27		2,4-Dimethylphenol	0.320	0.313	2.2	111	0.00
28		bis(2-Chloroethoxy)methane	0.395	0.390	1.3	114	-0.01
29	C	2,4-Dichlorophenol	0.287	0.304	-5.9	124	0.00
30		1,2,4-Trichlorobenzene	0.289	0.295	-2.1	119	0.00
31		Naphthalene	1.035	1.042	-0.7	117	-0.01
32		Benzoic acid	0.256	0.273	-6.6	126	0.01
33		4-Chloroaniline	0.445	0.460	-3.4	120	0.00
34	C	Hexachlorobutadiene	0.147	0.149	-1.4	118	-0.01
35		Caprolactam	0.110	0.116	-5.5	124	0.00
36	C	4-Chloro-3-methylphenol	0.328	0.349	-6.4	124	0.00
37		2-Methylnaphthalene	0.726	0.752	-3.6	122	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	124	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.587	0.592	-0.9	124	0.00
40	P	Hexachlorocyclopentadiene	0.258	0.227	12.0	105	-0.01
41	S	2,4,6-Tribromophenol	0.197	0.211	-7.1	132	0.00
42	C	2,4,6-Trichlorophenol	0.394	0.415	-5.3	127	0.00
43		2,4,5-Trichlorophenol	0.414	0.438	-5.8	129	0.00
44	S	2-Fluorobiphenyl	1.424	1.439	-1.1	124	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.764	1.775	-0.6	124	-0.01
46		2-Chloronaphthalene	1.263	1.280	-1.3	125	0.00
47		2-Nitroaniline	0.303	0.309	-2.0	124	0.00
48		Acenaphthylene	2.196	2.228	-1.5	125	0.00
49		Dimethylphthalate	1.564	1.602	-2.4	126	0.00
50		2,6-Dinitrotoluene	0.335	0.356	-6.3	130	0.00
51	C	Acenaphthene	1.332	1.373	-3.1	127	-0.01
52		3-Nitroaniline	0.389	0.396	-1.8	124	0.00
53	P	2,4-Dinitrophenol	0.116	0.144	-24.1	153#	0.00
54		Dibenzofuran	1.762	1.797	-2.0	127	-0.01
55	P	4-Nitrophenol	0.294	0.308	-4.8	125	0.00
56		2,4-Dinitrotoluene	0.433	0.468	-8.1	130	0.00
57		Fluorene	1.622	1.638	-1.0	125	-0.01
58		2,3,4,6-Tetrachlorophenol	0.316	0.330	-4.4	129	0.00
59		Diethylphthalate	1.608	1.638	-1.9	127	0.00
60		4-Chlorophenyl-phenylether	0.720	0.734	-1.9	129	0.00
61		4-Nitroaniline	0.391	0.395	-1.0	123	0.00
62		Azobenzene	1.264	1.228	2.8	120	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
64		4,6-Dinitro-2-methylphenol	0.094	0.109	-16.0	141	0.00
65	c	n-Nitrosodiphenylamine	0.646	0.669	-3.6	126	-0.01
66		4-Bromophenyl-phenylether	0.211	0.224	-6.2	130	-0.01
67		Hexachlorobenzene	0.225	0.236	-4.9	130	-0.01
68		Atrazine	0.196	0.200	-2.0	126	0.00
69	C	Pentachlorophenol	0.129	0.136	-5.4	130	-0.01
70		Phenanthrene	1.139	1.154	-1.3	126	-0.01
71		Anthracene	1.156	1.171	-1.3	125	-0.01
72		Carbazole	1.037	1.047	-1.0	125	0.00
73		Di-n-butylphthalate	1.330	1.338	-0.6	124	-0.01
74	C	Fluoranthene	1.230	1.245	-1.2	125	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	124	-0.02
76		Benzidine	0.647	0.563	13.0	104	0.00
77		Pyrene	1.319	1.350	-2.4	127	-0.01
78	S	Terphenyl-d14	0.857	0.884	-3.2	127	-0.01
79		Butylbenzylphthalate	0.632	0.652	-3.2	125	-0.02
80		Benzo(a)anthracene	1.194	1.221	-2.3	127	-0.01
81		3,3'-Dichlorobenzidine	0.443	0.446	-0.7	123	-0.01
82		Chrysene	1.123	1.142	-1.7	126	-0.01
83		Bis(2-ethylhexyl)phthalate	0.936	0.948	-1.3	124	-0.02
84	c	Di-n-octyl phthalate	1.537	1.572	-2.3	124	-0.03
85		Indeno(1,2,3-cd)pyrene	1.291	1.336	-3.5	126	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	125	-0.03
87		Benzo(b)fluoranthene	1.224	1.262	-3.1	128	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.199	1.202	-0.3	126	-0.02
89 C	Benzo(a)pyrene	1.166	1.187	-1.8	127	-0.03
90	Dibenzo(a,h)anthracene	1.134	1.150	-1.4	127	-0.06
91	Benzo(g,h,i)perylene	1.126	1.135	-0.8	126	-0.06

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