Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022417\
Data File : BG025951.D
Acq On : 25 Feb 2017 3:39
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Feb 25 04:18:44 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Feb 17 15:48:21 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	101	0.00
2		1,4-Dioxane	0.525	0.475	9.5	91	0.00
3		Pyridine	1.564	1.496	4.3	93	-0.01
4		n-Nitrosodimethylamine	0.563	0.535	5.0	93	0.00
5	S	2-Fluorophenol	1.149	1.137	1.0	97	0.00
6		Aniline	2.363	2.403	-1.7	99	0.00
7	S	Phenol-d6	1.700	1.779	-4.6	101	0.00
8		2-Chlorophenol	1.365	1.407	-3.1	101	0.00
9		Benzaldehyde	1.007	0.975	3.2	96	0.00
10	C	Phenol	1.847	1.906	-3.2	99	0.00
11		bis(2-Chloroethyl)ether	1.518	1.468	3.3	95	0.00
12		1,3-Dichlorobenzene	1.578	1.533	2.9	97	0.00
13	C	1,4-Dichlorobenzene	1.599	1.588	0.7	98	0.00
14		1,2-Dichlorobenzene	1.570	1.540	1.9	97	0.00
15		Benzyl Alcohol	1.245	1.327	-6.6	100	0.00
16		2,2'-oxybis(1-Chloropropane	2.203	2.045	7.2	92	0.00
17		2-Methylphenol	1.331	1.419	-6.6	102	0.00
18		Hexachloroethane	0.530	0.519	2.1	98	-0.01
19	P	n-Nitroso-di-n-propylamine	1.242	1.295	-4.3	100	0.00
20		3+4-Methylphenols	1.896	2.042	-7.7	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
22		Acetophenone	0.480	0.481	-0.2	99	0.00
23	S	Nitrobenzene-d5	0.317	0.327	-3.2	101	0.00
24		Nitrobenzene	0.343	0.347	-1.2	100	0.00
25		Isophorone	0.697	0.733	-5.2	103	0.00
26	C	2-Nitrophenol	0.160	0.184	-15.0	109	0.00
27		2,4-Dimethylphenol	0.299	0.314	-5.0	104	0.00
28		bis(2-Chloroethoxy)methane	0.442	0.442	0.0	100	-0.01
29	C	2,4-Dichlorophenol	0.285	0.314	-10.2	107	0.00
30		1,2,4-Trichlorobenzene	0.310	0.312	-0.6	103	0.00
31		Naphthalene	1.034	1.042	-0.8	101	0.00
32		Benzoic acid	0.228	0.278	-21.9	117	0.00
33		4-Chloroaniline	0.449	0.479	-6.7	105	0.00
34	C	Hexachlorobutadiene	0.177	0.176	0.6	101	0.00
35		Caprolactam	0.115	0.142	-23.5	117	0.00
36	C	4-Chloro-3-methylphenol	0.342	0.379	-10.8	108	0.00
37		2-Methylnaphthalene	0.788	0.817	-3.7	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.501	0.490	2.2	104	0.00
40	P	Hexachlorocyclopentadiene	0.274	0.249	9.1	95	-0.01
41	S	2,4,6-Tribromophenol	0.185	0.210	-13.5	120	0.00
42	C	2,4,6-Trichlorophenol	0.324	0.350	-8.0	112	0.00
43		2,4,5-Trichlorophenol	0.368	0.404	-9.8	115	0.00
44	S	2-Fluorobiphenyl	1.145	1.111	3.0	104	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.449	1.421	1.9	105	0.00
46		2-Chloronaphthalene	1.048	1.030	1.7	105	0.00
47		2-Nitroaniline	0.332	0.364	-9.6	110	0.00
48		Acenaphthylene	1.866	1.899	-1.8	108	0.00
49		Dimethylphthalate	1.372	1.387	-1.1	108	0.00
50		2,6-Dinitrotoluene	0.308	0.337	-9.4	111	0.00
51	C	Acenaphthene	1.232	1.255	-1.9	109	0.00
52		3-Nitroaniline	0.345	0.378	-9.6	111	0.00
53	P	2,4-Dinitrophenol	0.161	0.185	-14.9	123	0.00
54		Dibenzofuran	1.643	1.643	0.0	107	0.00
55	P	4-Nitrophenol	0.281	0.308	-9.6	111	0.00
56		2,4-Dinitrotoluene	0.415	0.467	-12.5	113	0.00
57		Fluorene	1.468	1.496	-1.9	109	0.00
58		2,3,4,6-Tetrachlorophenol	0.324	0.359	-10.8	114	0.00
59		Diethylphthalate	1.419	1.446	-1.9	109	0.00
60		4-Chlorophenyl-phenylether	0.680	0.690	-1.5	110	0.00
61		4-Nitroaniline	0.357	0.394	-10.4	115	0.00
62		Azobenzene	1.220	1.230	-0.8	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
64		4,6-Dinitro-2-methylphenol	0.117	0.127	-8.5	116	0.00
65	c	n-Nitrosodiphenylamine	0.607	0.618	-1.8	110	0.00
66		4-Bromophenyl-phenylether	0.210	0.219	-4.3	113	0.00
67		Hexachlorobenzene	0.216	0.225	-4.2	115	0.00
68		Atrazine	0.215	0.225	-4.7	112	0.00
69	C	Pentachlorophenol	0.133	0.148	-11.3	115	0.00
70		Phenanthrene	1.086	1.085	0.1	109	0.00
71		Anthracene	1.107	1.119	-1.1	110	0.00
72		Carbazole	1.112	1.121	-0.8	110	0.00
73		Di-n-butylphthalate	1.092	1.145	-4.9	111	-0.01
74	C	Fluoranthene	1.195	1.187	0.7	109	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	113	-0.01
76		Benzidine	0.631	0.626	0.8	105	0.00
77		Pyrene	1.265	1.236	2.3	109	0.00
78	S	Terphenyl-d14	0.822	0.799	2.8	110	-0.01
79		Butylbenzylphthalate	0.480	0.530	-10.4	117	-0.01
80		Benzo(a)anthracene	1.168	1.173	-0.4	112	0.00
81		3,3'-Dichlorobenzidine	0.416	0.447	-7.5	115	-0.01
82		Chrysene	1.123	1.120	0.3	111	-0.01
83		Bis(2-ethylhexyl)phthalate	0.693	0.742	-7.1	114	-0.03
84	c	Di-n-octyl phthalate	1.145	1.280	-11.8	117	-0.03
85		Indeno(1,2,3-cd)pyrene	1.323	1.396	-5.5	116	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	115	-0.02
87		Benzo(b)fluoranthene	1.198	1.201	-0.3	111	-0.02

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88	Benzo(k)fluoranthene	1.169	1.182	-1.1	114	-0.02
89 C	Benzo(a)pyrene	1.134	1.154	-1.8	114	-0.02
90	Dibenzo(a,h)anthracene	1.139	1.166	-2.4	115	-0.03
91	Benzo(g,h,i)perylene	1.142	1.175	-2.9	116	-0.01

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