Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG090916\
Data File : BG023981.D
Acq On : 9 Sep 2016 19:34
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 10 01:11:23 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 09 16:24:09 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	116	-0.04
2		1,4-Dioxane	0.470	0.390	17.0	88	-0.02
3		Pyridine	1.412	1.564	-10.8	116	-0.03
4		n-Nitrosodimethylamine	0.744	0.885	-19.0	138	-0.03
5	S	2-Fluorophenol	1.139	1.160	-1.8	109	-0.04
6		Aniline	2.328	2.983	-28.1#	143	-0.04
7	S	Phenol-d6	1.754	2.288	-30.4#	143	-0.04
8		2-Chlorophenol	1.320	1.515	-14.8	127	-0.04
9		Benzaldehyde	1.246	1.441	-15.7	126	-0.04
10	C	Phenol	1.845	2.372	-28.6#	139	-0.04
11		bis(2-Chloroethyl)ether	1.451	1.632	-12.5	134	-0.04
12		1,3-Dichlorobenzene	1.545	1.524	1.4	114	-0.04
13	C	1,4-Dichlorobenzene	1.636	1.584	3.2	115	-0.04
14		1,2-Dichlorobenzene	1.536	1.597	-4.0	121	-0.04
15		Benzyl Alcohol	1.546	2.281	-47.5#	158#	-0.04
16		2,2'-oxybis(1-Chloropropane	1.945	2.342	-20.4	143	-0.04
17		2-Methylphenol	1.243	1.684	-35.5#	150	-0.04
18		Hexachloroethane	0.630	0.663	-5.2	120	-0.04
19	P	n-Nitroso-di-n-propylamine	1.549	2.257	-45.7#	167#	-0.04
20		3+4-Methylphenols	1.732	2.483	-43.4#	153#	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	142	-0.04
22		Acetophenone	0.518	0.534	-3.1	147	-0.04
23	S	Nitrobenzene-d5	0.448	0.488	-8.9	158#	-0.04
24		Nitrobenzene	0.482	0.526	-9.1	160#	-0.05
25		Isophorone	0.869	1.014	-16.7	168#	-0.05
26	C	2-Nitrophenol	0.183	0.199	-8.7	157#	-0.04
27		2,4-Dimethylphenol	0.311	0.335	-7.7	144	-0.04
28		bis(2-Chloroethoxy)methane	0.439	0.463	-5.5	155#	-0.04
29	C	2,4-Dichlorophenol	0.330	0.379	-14.8	160#	-0.04
30		1,2,4-Trichlorobenzene	0.362	0.362	0.0	148	-0.04
31		Naphthalene	1.031	1.010	2.0	146	-0.04
32		Benzoic acid	0.253	0.297	-17.4	166#	-0.04
33		4-Chloroaniline	0.430	0.491	-14.2	159#	-0.04
34	C	Hexachlorobutadiene	0.272	0.270	0.7	140	-0.04
35		Caprolactam	0.116	0.155	-33.6#	182#	-0.04
36	C	4-Chloro-3-methylphenol	0.441	0.556	-26.1#	177#	-0.04
37		2-Methylnaphthalene	0.784	0.833	-6.2	150#	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	167#	-0.03
39		1,2,4,5-Tetrachlorobenzene	0.664	0.665	-0.2	165#	-0.03
40	P	Hexachlorocyclopentadiene	0.335	0.250	25.4#	126	-0.04
41	S	2,4,6-Tribromophenol	0.360	0.381	-5.8	168#	-0.03
42	C	2,4,6-Trichlorophenol	0.443	0.454	-2.5	167#	-0.04
43		2,4,5-Trichlorophenol	0.448	0.477	-6.5	169#	-0.04
44	S	2-Fluorobiphenyl	1.395	1.255	10.0	152#	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.609	1.492	7.3	155#	-0.04
46		2-Chloronaphthalene	1.181	1.111	5.9	160#	-0.04
47		2-Nitroaniline	0.465	0.558	-20.0	203#	-0.03
48		Acenaphthylene	1.969	1.907	3.1	161#	-0.03
49		Dimethylphthalate	1.715	1.659	3.3	163#	-0.03
50		2,6-Dinitrotoluene	0.362	0.374	-3.3	172#	-0.03
51	C	Acenaphthene	1.217	1.170	3.9	163#	-0.03
52		3-Nitroaniline	0.339	0.391	-15.3	182#	-0.03
53	P	2,4-Dinitrophenol	0.226	0.239	-5.8	166#	-0.03
54		Dibenzofuran	1.900	1.852	2.5	163#	-0.03
55	P	4-Nitrophenol	0.269	0.321	-19.3	206#	-0.03
56		2,4-Dinitrotoluene	0.532	0.588	-10.5	182#	-0.03
57		Fluorene	1.647	1.656	-0.5	170#	-0.03
58		2,3,4,6-Tetrachlorophenol	0.455	0.496	-9.0	179#	-0.03
59		Diethylphthalate	1.831	1.794	2.0	164#	-0.03
60		4-Chlorophenyl-phenylether	0.887	0.892	-0.6	164#	-0.03
61		4-Nitroaniline	0.342	0.440	-28.7#	211#	-0.03
62		Azobenzene	1.773	1.869	-5.4	178#	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	182#	-0.03
64		4,6-Dinitro-2-methylphenol	0.139	0.145	-4.3	187#	-0.03
65	c	n-Nitrosodiphenylamine	0.568	0.532	6.3	169#	-0.03
66		4-Bromophenyl-phenylether	0.243	0.232	4.5	170#	-0.03
67		Hexachlorobenzene	0.285	0.267	6.3	169#	-0.03
68		Atrazine	0.244	0.257	-5.3	186#	-0.03
69	C	Pentachlorophenol	0.152	0.155	-2.0	176#	-0.03
70		Phenanthrene	1.040	1.015	2.4	181#	-0.03
71		Anthracene	1.061	1.043	1.7	181#	-0.03
72		Carbazole	0.963	1.022	-6.1	196#	-0.03
73		Di-n-butylphthalate	1.149	1.172	-2.0	189#	-0.03
74	C	Fluoranthene	1.253	1.415	-12.9	203#	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	213#	-0.03
76		Benzidine	0.619	0.614	0.8	205#	-0.02
77		Pyrene	1.230	1.165	5.3	203#	-0.03
78	S	Terphenyl-d14	0.878	0.799	9.0	195#	-0.02
79		Butylbenzylphthalate	0.474	0.534	-12.7	255#	-0.02
80		Benzo(a)anthracene	1.120	1.129	-0.8	219#	-0.03
81		3,3'-Dichlorobenzidine	0.448	0.434	3.1	204#	-0.02
82		Chrysene	1.066	1.052	1.3	216#	-0.03
83		Bis(2-ethylhexyl)phthalate	0.649	0.693	-6.8	246#	-0.03
84	c	Di-n-octyl phthalate	1.065	0.989	7.1	212#	-0.03
85		Indeno(1,2,3-cd)pyrene	1.180	1.103	6.5	210#	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	190#	-0.05
87		Benzo(b)fluoranthene	1.136	1.193	-5.0	203#	-0.04

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.127	1.169	-3.7	198#	-0.04
89 C	Benzo(a)pyrene	1.079	1.127	-4.4	203#	-0.05
90	Dibenzo(a,h)anthracene	1.061	1.074	-1.2	196#	-0.08
91	Benzo(g,h,i)perylene	1.020	1.085	-6.4	207#	-0.09

(#) = Out of Range SPCC's out = 0 CCC's out = 2