Table S3. Mean molecular distances between ZmpB subgroups and the ZmpD cluster. Standard errors are in grey.

	***zmpB*_1**	***zmpB*_2**	***zmpB*_3**	***zmpB*_4**	***zmpB*_5**	***zmpB*_6**	***zmpB*_7**	***zmpB*_8**	***zmpB*_9**	***zmpB*_10**	***zmpD*_11**
***zmpB*_1**		0.010	0.009	0.011	0.015	0.014	0.015	0.017	0.018	0.023	0.013
***zmpB*_2**	0.266		0.005	0.012	0.015	0.014	0.013	0.018	0.018	0.019	0.014
***zmpB*_3**	0.225	0.089		0.009	0.016	0.014	0.013	0.019	0.018	0.021	0.014
***zmpB*_4**	0.301	0.353	0.231		0.013	0.014	0.016	0.017	0.020	0.025	0.011
***zmpB*_5**	0.504	0.480	0.508	0.422		0.013	0.015	0.022	0.021	0.027	0.012
***zmpB*_6**	0.440	0.446	0.444	0.439	0.442		0.016	0.023	0.020	0.024	0.013
***zmpB*_7**	0.417	0.416	0.419	0.461	0.521	0.486		0.020	0.018	0.024	0.015
***zmpB*_8**	0.629	0.603	0.612	0.634	0.674	0.679	0.577		0.014	0.022	0.018
***zmpB*_9**	0.619	0.617	0.625	0.627	0.681	0.653	0.574	0.411		0.021	0.017
***zmpB*_10**	0.762	0.726	0.736	0.751	0.811	0.772	0.739	0.681	0.687		0.023
***zmpD*_11**	0.399	0.459	0.458	0.378	0.463	0.461	0.502	0.660	0.660	0.810	