Erratum: Density functional study of the effect of pressure on the ferroelectric GeTe [Phys. Rev. B 73, 214105 (2006)]

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Figure 1, the energy bands of GeTe, did not contain the spin-orbit splittings that it was supposed to. The energy bands are calculated after the self-consistent calculations, so any error in them has no effect on anything else. We thank Manuel Cardona for calling this to our attention and Priya Gopal for calculating the corrected spin-orbit bands.

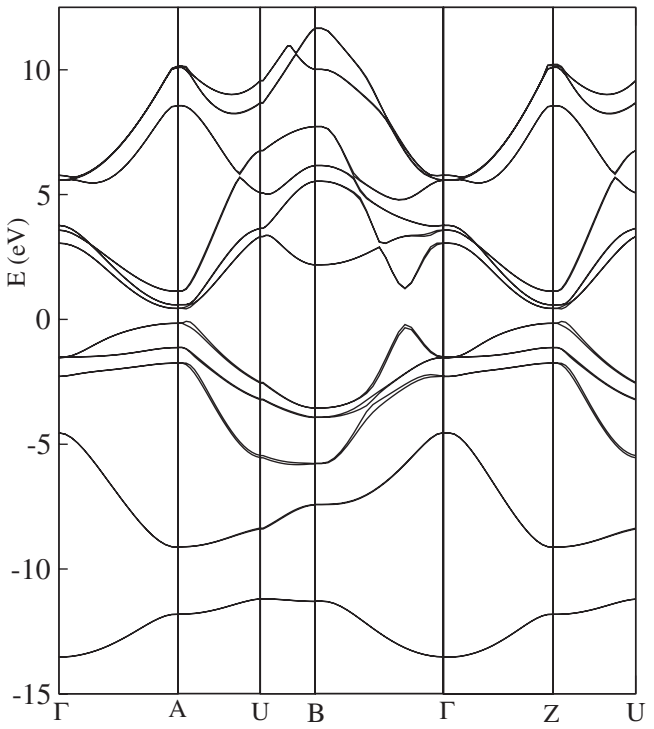

FIG. 1. Energy bands of rhombohedral GeTe. The zero of energy is taken at the top of the valence band. The symmetry points are labeled as described in the text except that, to be consistent with Slater, L has been relabeled A.

