Erratum: Large anomalous Nernst coefficient in an oxide skyrmion crystal Chern insulator **[Phys. Rev. B 98, 205125 (2018)]**

Yo Pierre Mizuta, Hikaru Sawahata, and Fumiyuki Ishii

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In Fig. 5(c) of the original article, incorrectness of thermoelectric conductivities α_{xx} and α_{xy} by a constant factor has been noticed. In this Erratum is the corrected version of Fig. 5. Specifically, the correction made is nothing but multiplication of the originally plotted values by a factor of 600 (note the revision of unit for the horizontal axis).

This correction does not affect any other part of the original article.

FIG. 5. (a) Zoomed-in band structure around the conduction band bottom, and the chemical potential dependence of (b) zero-temperature electric conductivities $\sigma_{xx}(\tau = 10 \text{ fs})$ and σ_{xy} , and of (c) thermoelectric conductivities $\alpha_{xx}(\tau = 10 \text{ fs})$ and α_{xy} at 100 K. The system has two additional electrons as compared to the pristine system, which occupy the lowest two bands in (a), where their respective Chern numbers of -1 are indicated. A Gaussian smearing of 10 meV is employed for (b).