**Erratum: Induced Magnetoelectric Response in *Pnma* Perovskites
[Phys. Rev. Lett. 107, 197603 (2011)]**

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In our Letter, we reported that the space group of the ferroelectric ground state of strained CaMnO_3 is $Pmc2_1$. However, this was a typographical error and the correct space group of our calculated lowest energy structure is $Pmn2_1$, consistent with that reported by Zhao *et al.* in Ref. [1].

[1] H. J. Zhao, M. N. Grisolia, Y. Yang, J. Íñiguez, M. Bibes, X. M. Chen, and L. Bellaiche, *Phys. Rev. B* **92**, 235133 (2015).