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Electronic Structure Calculations for $YBa_2Cu_3O_7$ within the Slave Boson Formalism [Phys. Rev. Lett. 77, 4066 (1996)]

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On page 4067, third paragraph, line 10, $b_0^2 = 0.1$ should be $b_0^2 = 0.05$.
An incorrect figure was printed as Fig. 2 in this article. Below is the proper figure.

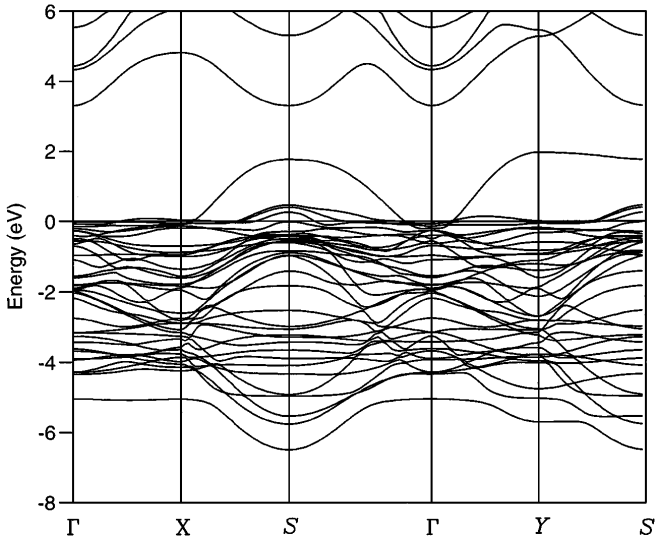

FIG. 2. Band structure of $YBCo_7$, as obtained from LDA + SB self-consistent electronic structure calculations. The energies are referred to the Fermi level.

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