Erratum: "Efficient chain moves for Monte Carlo simulations of a wormlike DNA model: Excluded volume, supercoils, site juxtapositions, knots, and comparisons with random-flight and lattice models" [J. Chem. Phys. 128, 145104 (2008)]

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On page 145104–4 of Ref. 1, starting in the eighth line below Eq. (3), two errors should be corrected in the statement "ν=0.243, 8.2, and 78.8 $e/$Å (e is the electronic charge)…" The values 8.2 and 78.8 should read, respectively, 0.82 and 7.88 instead. These errors are merely typographical. Our simulations were conducted using the correct values of ν. The results of the paper remain unchanged.

[1] Z. Liu and H. S. Chan, J. Chem. Phys. **128**, 145104 (2008).

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