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Erratum: "A quantum propagator for path-integral simulations of rigid molecules" [J. Chem. Phys. 134, 054117 (2011)]

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The captions to Figures 1 and 2 should read $PT = 1500$ K and not $PT = 120$ K.[1] Second, in paragraph 2 of Sec. III (Computational Details) concerning the symmetries of the propagator, the description we gave is somewhat opaque and is better described in a recent publication[2] (which also includes FORTRAN source code to calculate the propagator). Neither the main body of the work nor the conclusions are affected by these changes.

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[1]E. G. Noya, C. Vega, and C. McBride, "A quantum propagator for path-integral simulations of rigid molecules," J. Chem. Phys. **134**, 054117 (2011).
[2]C. McBride, E. G. Noya, and C. Vega, Comput. Phys. Commun. **184**, 885 (2013).

