CORRIGENDUM

Atomistic simulations of spinodal phase separation preceding polymer crystallization

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In the above Letter, the parameters for the nonpolar polymer model used in the simulations were incorrectly specified as being identical to the polyethylene model of Paul *et al.* (ref. 21 of the Letter). However, the potential used in the Letter differed from that of Paul *et al.* in that the two-fold term in the C-C-C-C torsion was changed from $k_{\Phi}^2(1-\cos 2\Phi)$, used in the polyethylene (PE) polymer model of Paul *et al.* to $k_{\Phi}^2(1+\cos 2\Phi)$.