Erratum: Flexible ring polymers in an obstacle environment: Molecular theory of linear viscoelasticity [Phys. Rev. E 74, 021805 (2006)]

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We would like to point out a typographical error in the second term of Eq. (27) in this paper. The actual equation should read as:

$$\frac{\partial \langle X_{p\alpha} X_{p\beta} \rangle}{\partial t} = -2 \frac{k_p}{\zeta_p} \langle X_{p\alpha} X_{p\beta} \rangle + 4 \frac{k_B T}{\zeta_p} \delta_{\alpha\beta} + \kappa_{\alpha\mu} \langle X_{p\mu} X_{p\beta} \rangle + \kappa_{\beta\mu} \langle X_{p\alpha} X_{p\mu} \rangle. \tag{27}$$

This does not modify the results of the paper.

In the second and the third paragraph of Sec. V on melt of rings we have presented two scenarios of modified number of segments between entanglements (N_e^r) and modified entanglement spacing (a_r) , respectively. In presenting these scenarios it was assumed that the structure between entanglements is a fractal structure with $a_r = (N_e^r)^\nu b$. The primitive chain formulation for length of trunk in PPR framework of Sec. II intrinsically assumes and requires a Gaussian structure between entanglements. In order to have a consistent framework the incorrect fractal picture between entanglements and the arguments presented based on it like the expressions for entanglement spacing (a_r) , number of segments between entanglements (N_e^r) , and the plateau modulus ratio in Eq. (44) and Eq. (45) are invalid.

We change the two scenarios presented earlier to those consistent with the Doi-Edwards primitive chain formulation. We change the first scenario to have entanglement spacing (a_r) for rings to be the same as that for linear chains (a_l) and the resulting corrected fractal picture giving the number of segments between entanglements to be $N_e^r = N_e^l N^{1-2\nu}$. We change the second scenario to have the entanglement spacing $a_r \approx 0.6a_l$ and the number of segments between entanglements to be $N_e^r = 0.36N_e^l N^{1-2\nu}$. The calculations presented in the paper are not affected by this change as entanglement spacing (a_r) is a fit parameter in the model.

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