

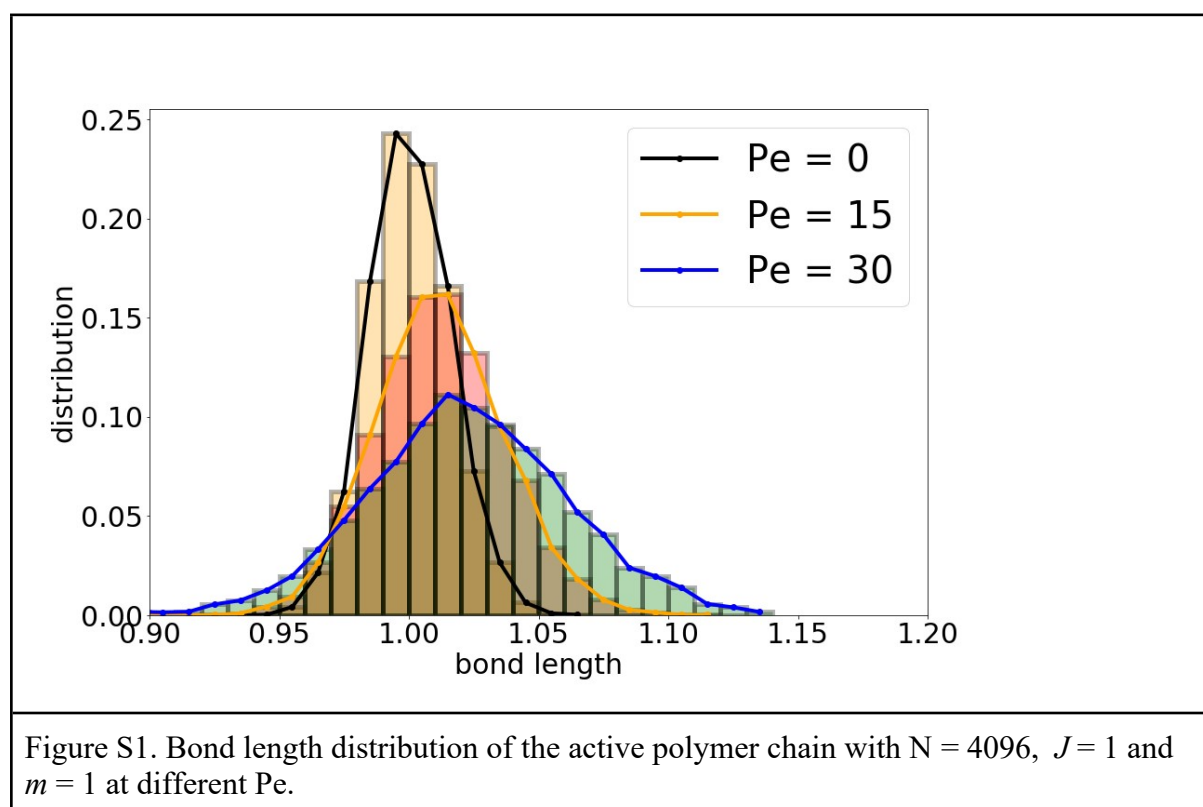
SUPPLEMENTARY INFORMATION

Motility-induced collapse of active Brownian particle polymer chain

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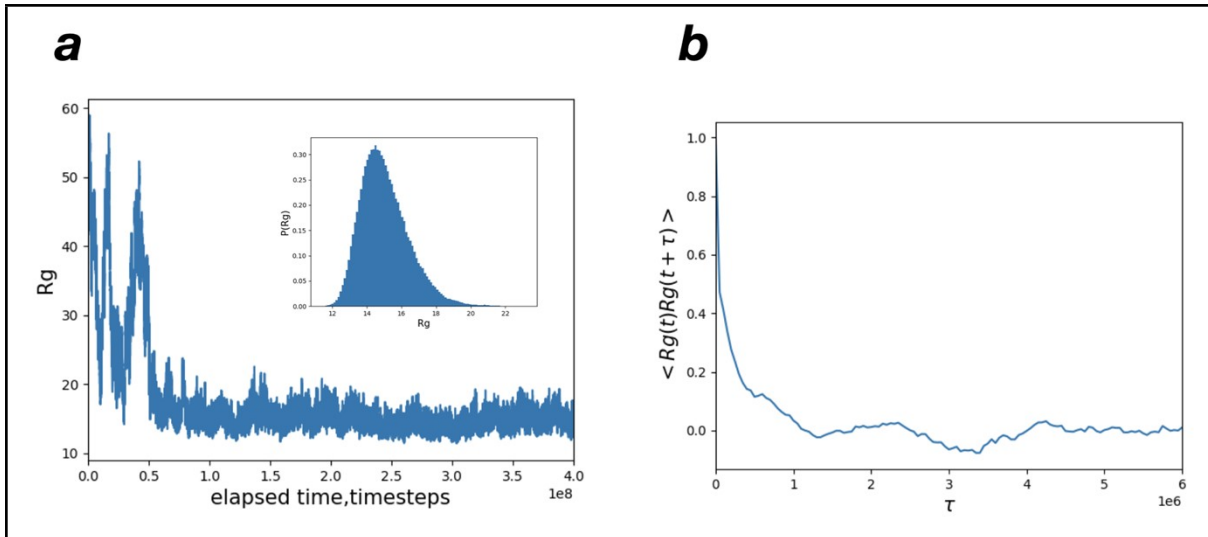


Figure S2. Typical dependence of active polymer size R_g on simulation time with $N = 4096$ and $Pe = 20$. Inset shows density distribution of R_g , obtained within $R_g(t)$ plateau and collected over several independent runs (a). $\langle R_g(t)R_g(t+\tau) \rangle$ autocorrelation function for $N = 4096$ and $Pe = 20$ (b).

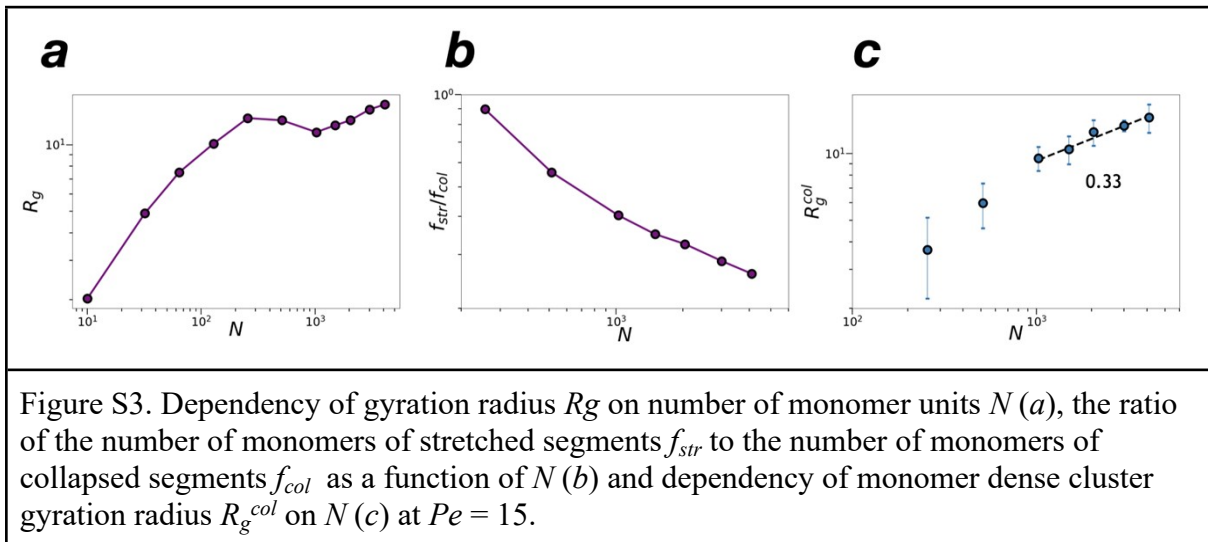


Figure S3. Dependency of gyration radius R_g on number of monomer units N (a), the ratio of the number of monomers of stretched segments f_{str} to the number of monomers of collapsed segments f_{col} as a function of N (b) and dependency of monomer dense cluster gyration radius R_g^{col} on N (c) at $Pe = 15$.

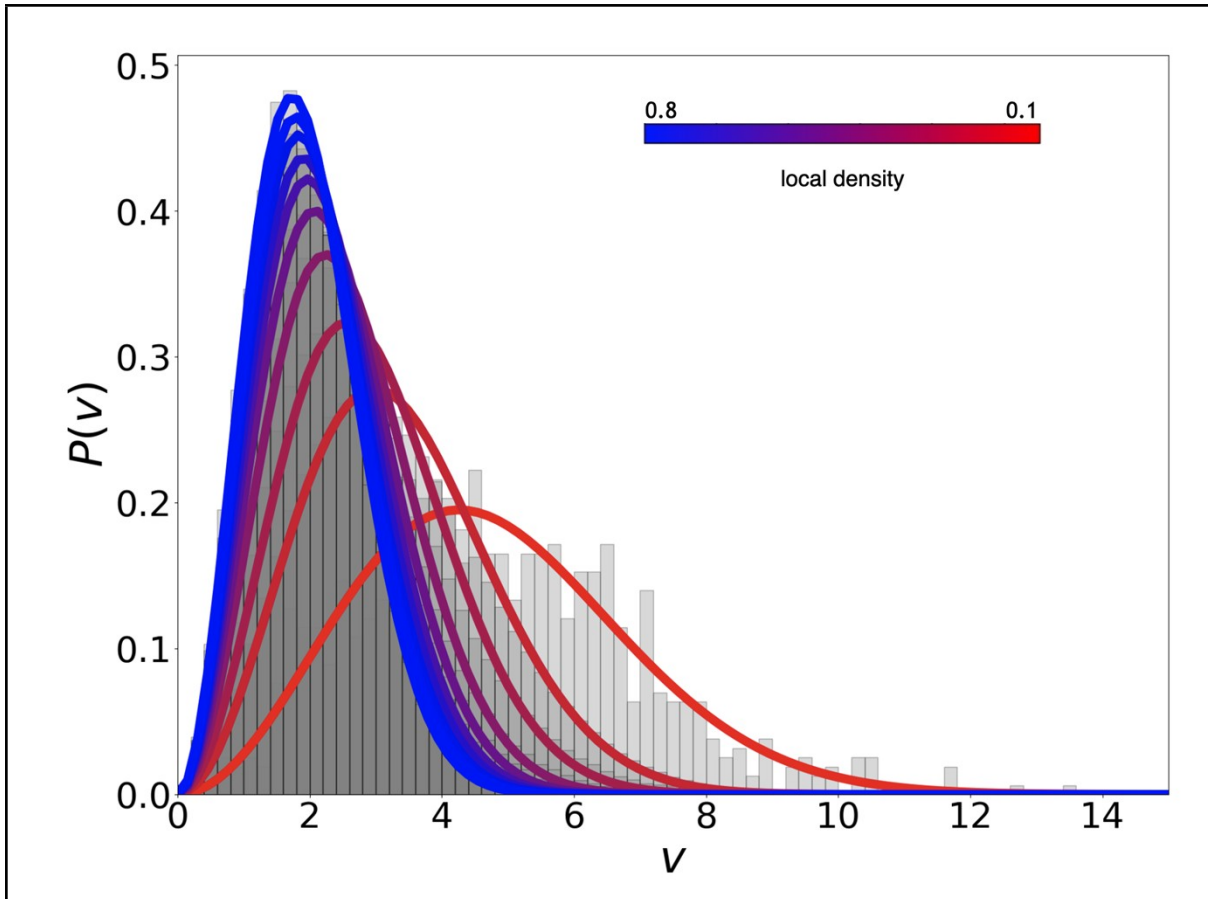


Figure S4. Full set of normalized velocity distributions $P(v)$ of underdamped $N = 4096$ active chain with $Pe = 15$ for different local densities, approximated with Maxwell-Boltzman velocity functions.

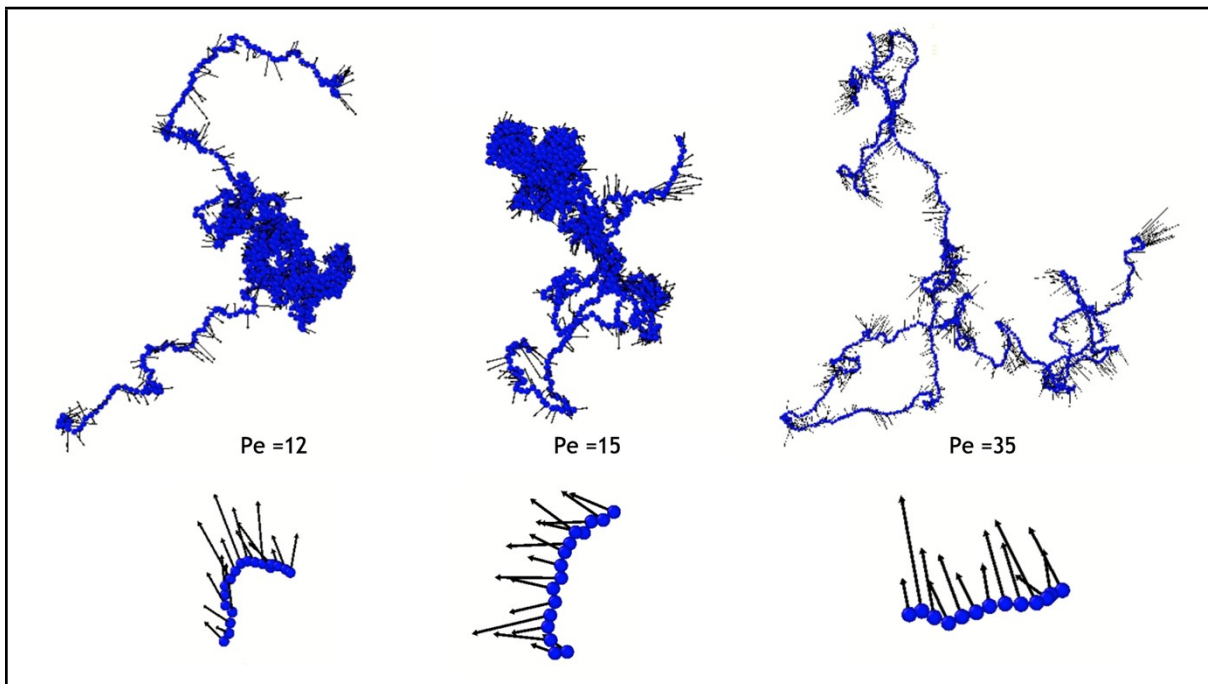


Fig. S5. Typical snapshots of the active chain with $N = 1024$, showing amplitudes and directions of the monomer units velocities (black arrows).

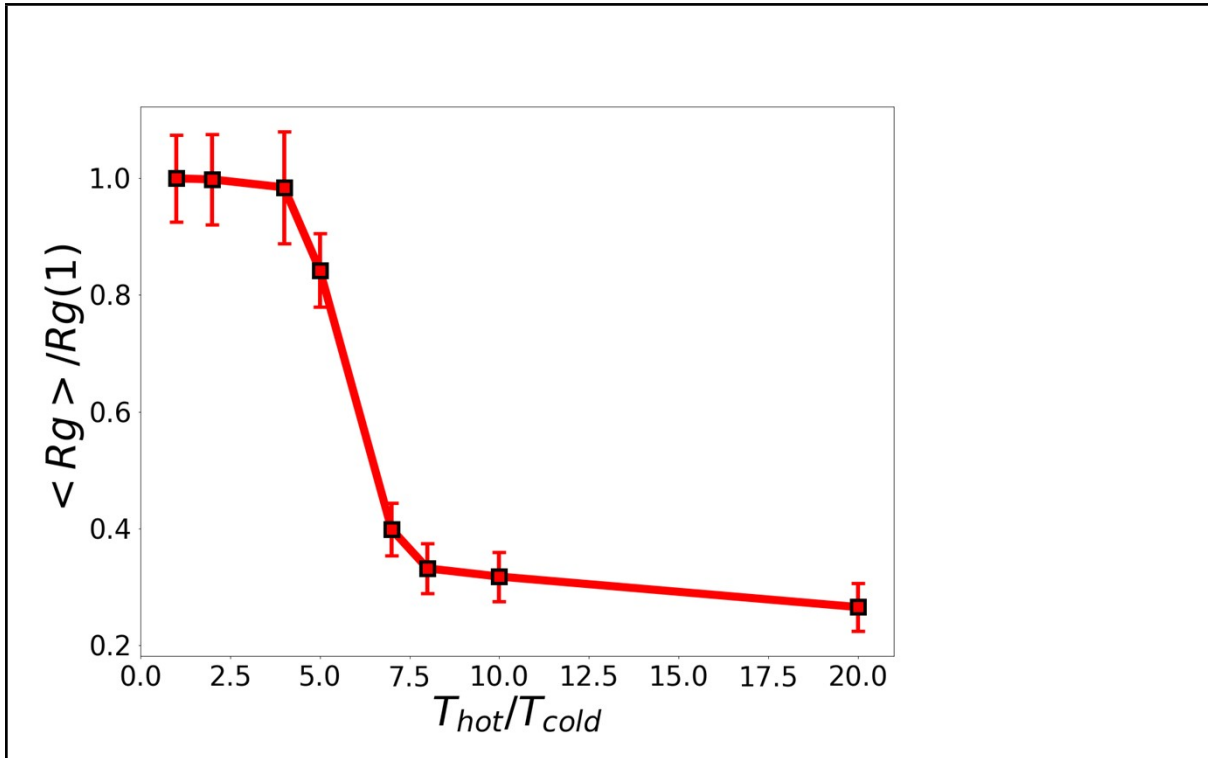


Figure S6. Reduced gyration radius Rg dependence on T_{hot}/T_{cold} for the $[A_{20}B_{80}]_{10}A_{20}$ regularly alternating copolymer with A - hot segments of length 20 and B - cold segments ($T_{cold} = 1$) of length 80.

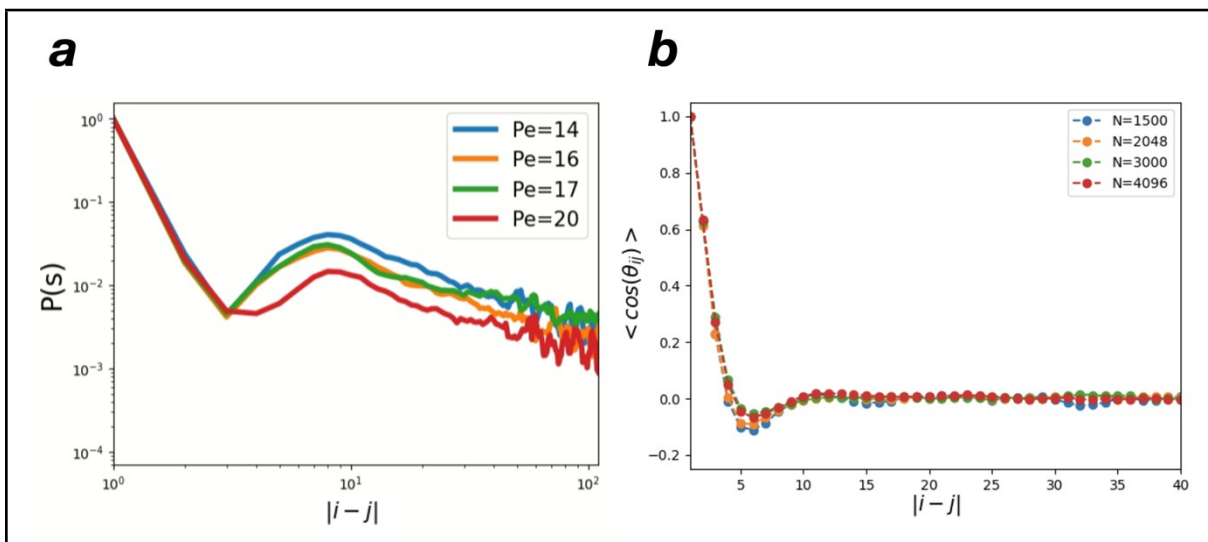


Figure S7. Contact probability P on contour distance along the chain $s = |i-j|$ for collapsed

segments for different Pe within collapsed conformation region of $N = 4096$ active polymer chain (a); Bond-bond autocorrelation function along the chain ($\langle \cos \theta_{ij} \rangle$, where θ_{ij} is angle between) with $N = 4096$ and $Pe = 15$.