Wenlin Zhang

List of Publications by Year in descending order

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Version: 2024-02-01

759233 888059 17 486 12 17 h-index citations g-index papers 17 17 17 699 citing authors docs citations times ranked all docs

#	Article	IF	CITATIONS
1	Molecular Dynamics Simulations of the Effects of Entanglement on Polymer Crystal Nucleation. Macromolecules, 2022, 55, 4899-4906.	4.8	14
2	Molecular Dynamics Simulations of Crystal Nucleation near Interfaces in Incompatible Polymer Blends. Polymers, 2021, 13, 347.	4.5	9
3	Effect of Flow-Induced Nematic Order on Polyethylene Crystal Nucleation. Macromolecules, 2020, 53, 7650-7657.	4.8	24
4	Modeling Intercolloidal Interactions Induced by Adsorption of Mobile Telechelic Polymers onto Particle Surfaces. Macromolecules, 2019, 52, 5357-5365.	4.8	11
5	A metastable nematic precursor accelerates polyethylene oligomer crystallization as determined by atomistic simulations and self-consistent field theory. Journal of Chemical Physics, 2019, 150, 244903.	3.0	22
6	Thermal Fluctuations Lead to Cumulative Disorder and Enhance Charge Transport in Conjugated Polymers. Macromolecular Rapid Communications, 2019, 40, e1900134.	3.9	8
7	Nematic Order Imposes Molecular Weight Effect on Charge Transport in Conjugated Polymers. ACS Central Science, 2018, 4, 413-421.	11.3	16
8	Tension-Induced Nematic Phase Separation in Bidisperse Homopolymer Melts. ACS Central Science, 2018, 4, 1545-1550.	11.3	10
9	Side chain length affects backbone dynamics in poly(3â€alkylthiophene)s. Journal of Polymer Science, Part B: Polymer Physics, 2018, 56, 1193-1202.	2.1	31
10	Direct All-Atom Molecular Dynamics Simulations of the Effects of Short Chain Branching on Polyethylene Oligomer Crystal Nucleation. Macromolecules, 2018, 51, 4762-4769.	4.8	43
11	Predicting Flory-Huggins <mml:math <br="" xmlns:mml="http://www.w3.org/1998/Math/MathML">display="inline"><mml:mrow><mml:mi>i‡</mml:mi></mml:mrow></mml:math> from Simulations. Physical Review Letters, 2017, 119, 017801.	7.8	44
12	Predicting the Flory-Huggins χ Parameter for Polymers with Stiffness Mismatch from Molecular Dynamics Simulations. Polymers, 2016, 8, 241.	4.5	57
13	Using surface-induced ordering to probe the isotropic-to-nematic transition for semiflexible polymers. Soft Matter, 2016, 12, 6141-6147.	2.7	15
14	Surface-Induced Chain Alignment of Semiflexible Polymers. Macromolecules, 2016, 49, 963-971.	4.8	29
15	Molecular Rectification in Conjugated Block Copolymer Photovoltaics. Journal of Physical Chemistry C, 2016, 120, 6978-6988.	3.1	32
16	Predicting Nematic Phases of Semiflexible Polymers. Macromolecules, 2015, 48, 1454-1462.	4.8	43
17	Predicting Chain Dimensions of Semiflexible Polymers from Dihedral Potentials. Macromolecules, 2014, 47, 6453-6461.	4.8	78