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List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Chain-End Modification: A Starting Point for Controlling Polymer Crystal Nucleation. <i>Macromolecules</i> , 2021, 54, 1599-1610.	4.8	8
2	Quasi-continuous melting of model polymer monolayers prompts reinterpretation of polymer melting. <i>Nature Communications</i> , 2021, 12, 1710.	12.8	13
3	Roughening Transition and Quasi-continuous Melting of Monolayers of Ultra-long Alkanes: Why Bulk Polymer Melting Is Strongly First-Order. <i>Macromolecules</i> , 2021, 54, 10135-10149.	4.8	7
4	Property Decoupling across the Embryonic Nucleus-Melt Interface during Polymer Crystal Nucleation. <i>Journal of Physical Chemistry B</i> , 2020, 124, 4793-4804.	2.6	12
5	Monodisperse Polymer Melts Crystallize via Structurally Polydisperse Nanoscale Clusters: Insights from Polyethylene. <i>Polymers</i> , 2020, 12, 447.	4.5	6
6	A coarse-grain model for entangled polyethylene melts and polyethylene crystallization. <i>Journal of Chemical Physics</i> , 2019, 150, 244901.	3.0	23
7	Characterizing key features in the formation of ice and gas hydrate systems. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2019, 377, 20180167.	3.4	22
8	Polymer nucleation under high-driving force, long-chain conditions: Heat release and the separation of time scales. <i>Journal of Chemical Physics</i> , 2019, 150, 114901.	3.0	13
9	Divining the shape of nascent polymer crystal nuclei. <i>Journal of Chemical Physics</i> , 2019, 151, 144901.	3.0	11
10	Does Local Structure Bias How a Crystal Nucleus Evolves?. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 6991-6998.	4.6	19
11	Unraveling Mixed Hydrate Formation: Microscopic Insights into Early Stage Behavior. <i>Journal of Physical Chemistry B</i> , 2016, 120, 13218-13223.	2.6	14
12	Evidence from mixed hydrate nucleation for a funnel model of crystallization. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, 12041-12046.	7.1	47
13	Algorithms for GPU-based molecular dynamics simulations of complex fluids: Applications to water, mixtures, and liquid crystals. <i>Journal of Computational Chemistry</i> , 2015, 36, 1787-1804.	3.3	7