

# Yang Wang

## List of Publications by Year in descending order

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20  
papers

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citations

840776

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888059

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times ranked

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citing authors

#	ARTICLE	IF	CITATIONS
1	Three Tribolayers Self-Generated from SiC Individually Work for Reducing Friction in Different Contact Pressures. <i>Journal of Physical Chemistry C</i> , 2022, 126, 2728-2736.	3.1	9
2	Role of Interfacial Bonding in Tribochemical Wear. <i>Frontiers in Chemistry</i> , 2022, 10, 852371.	3.6	9
3	Chemical-Reaction-Induced deformation of Body-Centered cubic iron in supercritical water leading to high risk of cleavage Fracture: A reactive Molecular dynamics study. <i>Computational Materials Science</i> , 2022, 208, 111354.	3.0	0
4	Non-Empirical Law for Nanoscale Atom-by-Atom Wear. <i>Advanced Science</i> , 2021, 8, 2002827.	11.2	21
5	Cooperative roles of chemical reactions and mechanical friction in chemical mechanical polishing of gallium nitride assisted by OH radicals: tight-binding quantum chemical molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 4075-4084.	2.8	15
6	Role of OH Termination in Mitigating Friction of Diamond-like Carbon under High Load: A Joint Simulation and Experimental Study. <i>Langmuir</i> , 2021, 37, 6292-6300.	3.5	11
7	Selective Wear Behaviors of a Water-Lubricating SiC Surface under Rotating-Contact Conditions Revealed by Large-Scale Reactive Molecular Dynamics Simulations. <i>Journal of Physical Chemistry C</i> , 2021, 125, 14957-14964.	3.1	6
8	Different Etching Mechanisms of Diamond by Oxygen and Hydrogen Plasma: a Reactive Molecular Dynamics Study. <i>Journal of Physical Chemistry C</i> , 2021, 125, 16711-16718.	3.1	16
9	Atom-by-Atom and Sheet-by-Sheet Chemical Mechanical Polishing of Diamond Assisted by OH Radicals: A Tight-Binding Quantum Chemical Molecular Dynamics Simulation Study. <i>ACS Applied Materials &amp; Interfaces</i> , 2021, 13, 41231-41237.	8.0	19
10	Heterogeneous yielding mechanisms of body centered cubic iron for high resistance to chemical reaction-induced deterioration in supercritical water environments: A reactive molecular dynamics study. <i>Scripta Materialia</i> , 2021, 202, 113997.	5.2	3
11	Reactive Molecular Dynamics Simulations of Wear and Tribochemical Reactions of Diamond like Carbon Interfaces with Nanoscale Asperities under H <sub>2</sub> Gas: Implications for Solid Lubricant Coatings. <i>ACS Applied Nano Materials</i> , 2020, 3, 7297-7304.	5.0	21
12	Self-Formed Double Tribolayers Play Collaborative Roles in Achieving Superlow Friction in an Aqueous Environment. <i>Journal of Physical Chemistry C</i> , 2020, 124, 8295-8303.	3.1	24
13	Development of a Transferable ReaxFF Parameter Set for Carbon- and Silicon-Based Solid Systems. <i>Journal of Physical Chemistry C</i> , 2020, 124, 10007-10015.	3.1	22
14	Reactive Molecular Dynamics Simulation on Friction-induced Chemical Reactions of SiC in Water Environments. <i>Journal of Computer Chemistry Japan</i> , 2020, 19, 139-141.	0.1	0
15	Proposal of a new formation mechanism for hydrogenated diamond-like carbon transfer films: Hydrocarbon-emission-induced transfer. <i>Carbon</i> , 2019, 154, 7-12.	10.3	35
16	Prediction of Macroscopic Properties of Diamond-like Carbon from Atomic-Scale Structure. <i>Journal of Physical Chemistry C</i> , 2019, 123, 24609-24614.	3.1	6
17	Triboemission of hydrocarbon molecules from diamond-like carbon friction interface induces atomic-scale wear. <i>Science Advances</i> , 2019, 5, eaax9301.	10.3	70
18	Tribochemical reactions and graphitization of diamond-like carbon against alumina give volcano-type temperature dependence of friction coefficients: A tight-binding quantum chemical molecular dynamics simulation. <i>Carbon</i> , 2018, 133, 350-357.	10.3	52

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19	Tight-Binding Quantum Chemical Molecular Dynamics Study on the Friction and Wear Processes of Diamond-Like Carbon Coatings: Effect of Tensile Stress. ACS Applied Materials & Interfaces, 2017, 9, 34396-34404.	8.0	43
20	Heterogeneous Yielding Mechanisms of Body Center Cubic Iron for High Resistance to Chemical Reaction-Induced Deterioration in Supercritical Water Environments: A Reactive Molecular Dynamics Study. SSRN Electronic Journal, 0, , .	0.4	0