Yunlong Li

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	A review on enhancement of mechanical and tribological properties of polymer composites reinforced by carbon nanotubes and graphene sheet: Molecular dynamics simulations. Composites Part B: Engineering, 2019, 160, 348-361.	12.0	168
2	A comparison study on mechanical properties of polymer composites reinforced by carbon nanotubes and graphene sheet. Composites Part B: Engineering, 2018, 133, 35-41.	12.0	146
3	A molecular dynamics simulation study on enhancement of mechanical and tribological properties of polymer composites by introduction of graphene. Carbon, 2017, 111, 538-545.	10.3	131
4	Enhancement of tribological properties of polymer composites reinforced by functionalized graphene. Composites Part B: Engineering, 2017, 120, 83-91.	12.0	91
5	A study on tribology of nitrile-butadiene rubber composites by incorporation of carbon nanotubes: Molecular dynamics simulations. Carbon, 2016, 100, 145-150.	10.3	75
6	Enhancement of fracture properties of polymer composites reinforced by carbon nanotubes: A molecular dynamics study. Carbon, 2018, 129, 504-509.	10.3	71
7	Molecular dynamics simulations of tribology properties of NBR (Nitrile-Butadiene Rubber) /carbon nanotube composites. Composites Part B: Engineering, 2016, 97, 62-67.	12.0	60
8	Enhanced tribological properties of polymer composites by incorporation of nano-SiO 2 particles: A molecular dynamics simulation study. Computational Materials Science, 2017, 134, 93-99.	3.0	51
9	The effect of sliding velocity on the tribological properties of polymer/carbon nanotube composites. Carbon, 2016, 106, 106-109.	10.3	33
10	A comparative study on enhancement of mechanical and tribological properties of nitrile rubber composites reinforced by different functionalized graphene sheets: Molecular dynamics simulations. Polymer Composites, 2021, 42, 205-219.	4.6	30
11	The interfacial load-transfer enhancement mechanism of amino-functionalised carbon nanotube reinforced epoxy matrix composites: A molecular dynamics study. Composites Science and Technology, 2021, 209, 108790.	7.8	29
12	Effects of carbon nanotubes functionalization on mechanical and tribological properties of nitrile rubber nanocomposites: Molecular dynamics simulations. Computational Materials Science, 2021, 196, 110556.	3.0	24
13	Molecular dynamics study on the reinforcing effect of incorporation of graphene/carbon nanotubes on the mechanical properties of swelling rubber. Polymer Testing, 2021, 102, 107337.	4.8	18
14	Molecular dynamics simulations of mechanical properties of swollen nitrile rubber composites by incorporating carbon nanotubes. Polymer Composites, 2020, 41, 3160-3169.	4.6	14
15	A study on the enhancement of the mechanical properties of weak structural planes based on microbiologically induced calcium carbonate precipitation. Bulletin of Engineering Geology and the Environment, 2020, 79, 4349-4362.	3.5	11
16	A study on effects of stone–thrower–wales defective carbon nanotubes on glass transition temperature of polymer composites using molecular dynamics simulations. Computational Materials Science, 2021, 186, 110005.	3.0	9
17	Molecular Dynamics Simulations of Thermal Properties of Polymer Composites Enhanced by Cross-Linked Graphene Sheets. Acta Mechanica Solida Sinica, 2018, 31, 673-682.	1.9	8
18	Study on Preparation and thermal reflective properties of energy saving pigments with selective solar reflection. IOP Conference Series: Materials Science and Engineering, 2019, 544, 012010.	0.6	5

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19	Tribological properties of swollen nitrile rubber under dry and wet sliding conditions. Materials Research Express, 2020, 7, 015311.	1.6	4
20	A molecular dynamics study on the effect of TSW defective graphene on the glass transition temperature of polymer materials. Polymer Bulletin, 2022, 79, 2205-2218.	3.3	4
21	Driveline Simulation of 2013 Formula Student Electric Racing Vehicle. Applied Mechanics and Materials, 0, 541-542, 424-429.	0.2	3
22	Effect of modified nano/Mg(OH)2 on the flame retardancy and mechanical properties of NBR based on molecular simulation. Modelling and Simulation in Materials Science and Engineering, 0, , .	2.0	0