

Wolfgang Verestek

List of Publications by Year in descending order

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papers

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1684188
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9
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citing authors

| # | ARTICLE | IF | CITATIONS |
|---|--|-----|-----------|
| 1 | Molecular dynamics simulations investigating consecutive nucleation, solidification and grain growth in a twelve-million-atom Fe-system. <i>Journal of Crystal Growth</i> , 2017, 474, 140-145. | 1.5 | 23 |
| 2 | A multiscale modeling on fracture and strength of graphene platelets reinforced epoxy. <i>Engineering Fracture Mechanics</i> , 2020, 235, 107197. | 4.3 | 21 |
| 3 | Molecular dynamics investigations of the strengthening of Al-Cu alloys during thermal ageing. <i>Physical Mesomechanics</i> , 2017, 20, 291-304. | 1.9 | 16 |
| 4 | Microstructural Investigation of Plasma Sprayed Ceramic Coatings Using Peridynamics. <i>Journal of Mechanics</i> , 2020, 36, 183-196. | 1.4 | 9 |
| 5 | Characterization of Cure Behavior in Epoxy Using Molecular Dynamics Simulation Compared with Dielectric Analysis and DSC. <i>Polymers</i> , 2021, 13, 3085. | 4.5 | 9 |
| 6 | Microstructural investigation of plasma sprayed ceramic coatings focusing on the effect of the splat boundary for SOFC sealing applications using peridynamics. <i>Theoretical and Applied Fracture Mechanics</i> , 2021, 112, 102926. | 4.7 | 7 |
| 7 | Introducing a method of constructing realistic closed cell nano-porous iron crystals and MD simulations to investigate the influence of the system size on the stability and the mechanical properties. <i>Computational Materials Science</i> , 2019, 166, 150-154. | 3.0 | 4 |
| 8 | Atomistic-scale modeling of nano-clay-filled shape memory polymers. <i>Computational Materials Science</i> , 2021, 188, 110246. | 3.0 | 4 |
| 9 | Molecular Dynamics Simulation of High-Temperature Creep Behavior of Nickel Polycrystalline Nanopillars. <i>Molecules</i> , 2021, 26, 2606. | 3.8 | 0 |