Gennady B Sushko

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Atomistic simulation of the FEBID-driven growth of iron-based nanostructures. Physical Chemistry Chemical Physics, 2022, 24, 10807-10819.	2.8	2
2	Multiscale modeling of stochastic dynamics processes with <scp>MBN Explorer</scp> . Journal of Computational Chemistry, 2022, 43, 1442-1458.	3.3	1
3	All-atom relativistic molecular dynamics simulations of channeling and radiation processes in oriented crystals. European Physical Journal D, 2021, 75, 1.	1.3	13
4	Irradiation-driven molecular dynamics simulation of the FEBID process for Pt(PF3)4. Beilstein Journal of Nanotechnology, 2021, 12, 1151-1172.	2.8	5
5	Multiscale simulation of the focused electron beam induced deposition process. Scientific Reports, 2020, 10, 20827.	3.3	16
6	Modeling MesoBioNano systems with MBN Studio made easy. Journal of Molecular Graphics and Modelling, 2019, 88, 247-260.	2.4	34
7	Reactive molecular dynamics simulations of organometallic compound W(CO)6 fragmentation,. European Physical Journal D, 2019, 73, 1.	1.3	13
8	Channeling and radiation of 855†MeV electrons and positrons in straight and bent tungsten (1 1 0) crystals. Nuclear Instruments & Methods in Physics Research B, 2018, 424, 26-36.	1.4	12
9	Transonic Panel Flutter in Accelerating or Decelerating Flow Conditions. AIAA Journal, 2018, 56, 997-1010.	2.6	23
10	Molecular dynamics for irradiation driven chemistry: application to the FEBID process*. European Physical Journal D, 2016, 70, 1.	1.3	30
11	Simulation of channeling and radiation of 855 MeV electrons and positrons in a small-amplitude short-period bent crystal. Nuclear Instruments & Methods in Physics Research B, 2016, 387, 41-53.	1.4	21
12	Toward the Exploration of the NiTi Phase Diagram with a Classical Force Field. Journal of Physical Chemistry C, 2016, 120, 25043-25052.	3.1	4
13	Studying chemical reactions in biological systems with MBN Explorer: implementation of molecular mechanics with dynamical topology. European Physical Journal D, 2016, 70, 1.	1.3	33
14	Reconciling simulated melting and ground-state properties of metals with a modified embedded-atom method potential. Journal of Physics Condensed Matter, 2016, 28, 145201.	1.8	9
15	Channeling of ultra-relativistic positrons in bent diamond crystals. St Petersburg Polytechnical University Journal Physics and Mathematics, 2015, 1, 212-218.	0.3	3
16	Electron and positron propagation in straight and periodically bent axial and planar silicon channels. St Petersburg Polytechnical University Journal Physics and Mathematics, 2015, 1, 332-340.	0.3	3
17	A small-amplitude crystalline undulator based on 20ÂGeV electrons and positrons: Simulations. St Petersburg Polytechnical University Journal Physics and Mathematics, 2015, 1, 341-345.	0.3	7
18	Development of collisional data base for elementary processes of electron scattering by atoms and molecules. Nuclear Instruments & Methods in Physics Research B, 2015, 354, 90-95.	1.4	12

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#	Article	IF	CITATIONS
19	Multi-GeV electron and positron channeling in bent silicon crystals. Nuclear Instruments & Methods in Physics Research B, 2015, 355, 39-43.	1.4	18
20	Radiation emission by electrons channeling in bent silicon crystals. European Physical Journal D, 2014, 68, 1.	1.3	17
21	Validation of Classical Force Fields for the Description of Thermo-Mechanical Properties of Transition Metal Materials. Journal of Physical Chemistry A, 2014, 118, 8426-8436.	2.5	8
22	Molecular Dynamics Simulation of Self-Diffusion Processes in Titanium in Bulk Material, on Grain Junctions and on Surface. Journal of Physical Chemistry A, 2014, 118, 6685-6691.	2.5	14
23	Benchmarking of classical force fields by ab initio calculations of atomic clusters: Ti and Ni–Ti case. Computational and Theoretical Chemistry, 2013, 1021, 101-108.	2.5	11
24	Molecular dynamics simulations of the nanoindentation process of titanium crystal. Computational Materials Science, 2013, 76, 20-26.	3.0	65
25	Simulation of ultra-relativistic electrons and positrons channeling in crystals with MBN Explorer. Journal of Computational Physics, 2013, 252, 404-418.	3.8	48
26	Kinetics of liquid-solid phase transition in large nickel clusters. Physical Review B, 2013, 88, .	3.2	12