

Arkady Krasheninnikov

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

239
papers

22,256
citations

68
h-index

147
g-index

259
ext. papers

24,702
ext. citations

7.3
avg, IF

7.28
L-index

#	Paper	IF	Citations
239	Data-Driven Quest for Two-Dimensional Non-van der Waals Materials.. <i>Nano Letters</i> , 2022 ,	11.5	11
238	Physics and theory of defects in 2D materials: the role of reduced dimensionality 2022 , 7-41		1
237	Two-dimensional materials under ion irradiation: from defect production to structure and property engineering 2022 , 259-301		0
236	Compact Laser Devices for Measuring Airborne Microparticle Concentrations and Their Application at the Geophysical Monitoring Center of the Sadovsky Institute of Dynamics of Geospheres, Russian Academy of Sciences. <i>Seismic Instruments</i> , 2022 , 58, 235-243	1.1	
235	Quasi-two-dimensional NaCl crystals encapsulated between graphene sheets and their decomposition under an electron beam. <i>Nanoscale</i> , 2021 , 13, 19626-19633	7.7	1
234	Gas Permeability and Selectivity of a Porous WS ₂ Monolayer. <i>Journal of Physical Chemistry C</i> , 2021 , 125, 25055-25066	3.8	3
233	Boosting the Electrocatalytic Conversion of Nitrogen to Ammonia on Metal-Phthalocyanine-Based Two-Dimensional Conjugated Covalent Organic Frameworks. <i>Journal of the American Chemical Society</i> , 2021 , 143, 19992-20000	16.4	19
232	Polymorphic Phases of Metal Chlorides in the Confined 2D Space of Bilayer Graphene. <i>Advanced Materials</i> , 2021 , 33, e2105898	24	5
231	Photoluminescence line shapes for color centers in silicon carbide from density functional theory calculations. <i>Physical Review B</i> , 2021 , 103,	3.3	7
230	Enhancing Ferromagnetism and Tuning Electronic Properties of CrI Monolayers by Adsorption of Transition-Metal Atoms. <i>ACS Applied Materials & Interfaces</i> , 2021 , 13, 21593-21601	9.5	6
229	Atomistic Simulations of Defect Production in Monolayer and Bulk Hexagonal Boron Nitride under Low- and High-Fluence Ion Irradiation. <i>Nanomaterials</i> , 2021 , 11,	5.4	3
228	Defect Agglomeration and Electron-Beam-Induced Local-Phase Transformations in Single-Layer MoTe ₂ . <i>Journal of Physical Chemistry C</i> , 2021 , 125, 13601-13609	3.8	3
227	Mirror twin boundaries in MoSe monolayers as one dimensional nanotemplates for selective water adsorption. <i>Nanoscale</i> , 2021 , 13, 1038-1047	7.7	5
226	Robust Magnetoelectric Effect in the Decorated Graphene/InSe Heterostructure. <i>ACS Applied Materials & Interfaces</i> , 2021 , 13, 3033-3039	9.5	6
225	Formation of Highly Doped Nanostripes in 2D Transition Metal Dichalcogenides via a Dislocation Climb Mechanism. <i>Advanced Materials</i> , 2021 , 33, e2007819	24	3
224	Controlled generation of luminescent centers in hexagonal boron nitride by irradiation engineering. <i>Science Advances</i> , 2021 , 7,	14.3	12
223	Enhanced Trion Emission in Monolayer MoSe ₂ by Constructing a Type-I Van Der Waals Heterostructure. <i>Advanced Functional Materials</i> , 2021 , 31, 2104960	15.6	2

222	Layer-Dependent Band Gaps of Platinum Dichalcogenides. <i>ACS Nano</i> , 2021 , 15, 13249-13259	16.7	9
221	Tunable electronic properties and enhanced ferromagnetism in CrGeTe monolayer by strain engineering. <i>Nanotechnology</i> , 2021 , 32,	3.4	2
220	Chlorine doping of MoSe flakes by ion implantation. <i>Nanoscale</i> , 2021 , 13, 5834-5846	7.7	8
219	Polymorphic Phases of Metal Chlorides in the Confined 2D Space of Bilayer Graphene (Adv. Mater. 52/2021). <i>Advanced Materials</i> , 2021 , 33, 2170415	24	0
218	Electron-beam-stimulated Atomic Migration Processes in Single-layer MoTe ₂ . <i>Microscopy and Microanalysis</i> , 2020 , 26, 534-537	0.5	
217	Electron Irradiation of Two-dimensional MoS ₂ : Insights into the Influence of Electronic Excitations from First-principle Calculations. <i>Microscopy and Microanalysis</i> , 2020 , 26, 784-785	0.5	0
216	Strain robust spin gapless semiconductors/half-metals in transition metal embedded MoSe monolayer. <i>Journal of Physics Condensed Matter</i> , 2020 , 32, 365305	1.8	4
215	Alkali metals inside bi-layer graphene and MoS ₂ : Insights from first-principles calculations. <i>Nano Energy</i> , 2020 , 75, 104927	17.1	16
214	Enhanced Ferromagnetism and Tunable Magnetism in FeGeTe Monolayer by Strain Engineering. <i>ACS Applied Materials & Interfaces</i> , 2020 , 12, 26367-26373	9.5	32
213	Simulating Raman spectra by combining first-principles and empirical potential approaches with application to defective MoS ₂ . <i>Npj Computational Materials</i> , 2020 , 6,	10.9	9
212	Synergistic electroreduction of carbon dioxide to carbon monoxide on bimetallic layered conjugated metal-organic frameworks. <i>Nature Communications</i> , 2020 , 11, 1409	17.4	166
211	Formation of Defects in Two-Dimensional MoS in the Transmission Electron Microscope at Electron Energies below the Knock-on Threshold: The Role of Electronic Excitations. <i>Nano Letters</i> , 2020 , 20, 2865-2870	11.5	40
210	Revealing the defect-dominated oxygen evolution activity of hematene. <i>Journal of Materials Chemistry A</i> , 2020 , 8, 6709-6716	13	23
209	Band Bending and Valence Band Quantization at Line Defects in MoS. <i>ACS Nano</i> , 2020 , 14, 9176-9187	16.7	8
208	Local vibrational modes of Si vacancy spin qubits in SiC. <i>Physical Review B</i> , 2020 , 101,	3.3	16
207	Megacity Aerosol Pollution and Atmospheric Electric Field Disturbances. <i>Izvestiya - Atmospheric and Oceanic Physics</i> , 2020 , 56, 759-772	1	1
206	Reversible crystalline-to-amorphous phase transformation in monolayer MoS ₂ under grazing ion irradiation. <i>2D Materials</i> , 2020 , 7, 025005	5.9	9
205	Tailoring the Electronic and Magnetic Properties of Hematene by Surface Passivation: Insights from First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 22784-22792	3.8	3

204	Freestanding and Supported MoS Monolayers under Cluster Irradiation: Insights from Molecular Dynamics Simulations. <i>ACS Applied Materials & Interfaces</i> , 2020 , 12, 37454-37463	9.5	10
203	Are two-dimensional materials radiation tolerant?. <i>Nanoscale Horizons</i> , 2020 , 5, 1447-1452	10.8	10
202	Channeling effects in gold nanoclusters under He ion irradiation: insights from molecular dynamics simulations. <i>Nanotechnology</i> , 2020 , 31, 035302	3.4	7
201	Which Transition Metal Atoms Can Be Embedded into Two-Dimensional Molybdenum Dichalcogenides and Add Magnetism?. <i>Nano Letters</i> , 2019 , 19, 4581-4587	11.5	36
200	Widely tunable GaAs bandgap via strain engineering in core/shell nanowires with large lattice mismatch. <i>Nature Communications</i> , 2019 , 10, 2793	17.4	34
199	Electron-Beam-Driven Structure Evolution of Single-Layer MoTe ₂ for Quantum Devices. <i>ACS Applied Nano Materials</i> , 2019 , 2, 3262-3270	5.6	23
198	Tomonaga-Luttinger Liquid in a Box: Electrons Confined within MoS ₂ Mirror-Twin Boundaries. <i>Physical Review X</i> , 2019 , 9,	9.1	21
197	Room-Temperature Ferromagnetism in MoTe ₂ by Post-Growth Incorporation of Vanadium Impurities. <i>Advanced Electronic Materials</i> , 2019 , 5, 1900044	6.4	38
196	Thermal transport in MoS ₂ from molecular dynamics using different empirical potentials. <i>Physical Review B</i> , 2019 , 99,	3.3	31
195	Enhanced sensitivity of MoSe monolayer for gas adsorption induced by electric field. <i>Journal of Physics Condensed Matter</i> , 2019 , 31, 445301	1.8	16
194	Layer Rotation-Angle-Dependent Excitonic Absorption in van der Waals Heterostructures Revealed by Electron Energy Loss Spectroscopy. <i>ACS Nano</i> , 2019 , 13, 9541-9550	16.7	17
193	Nonstoichiometric Phases of Two-Dimensional Transition-Metal Dichalcogenides: From Chalcogen Vacancies to Pure Metal Membranes. <i>Journal of Physical Chemistry Letters</i> , 2019 , 10, 6492-6498	6.4	12
192	Efficient method for calculating Raman spectra of solids with impurities and alloys and its application to two-dimensional transition metal dichalcogenides. <i>Physical Review Materials</i> , 2019 , 3,	3.2	9
191	Perforating Freestanding Molybdenum Disulfide Monolayers with Highly Charged Ions. <i>Journal of Physical Chemistry Letters</i> , 2019 , 10, 904-910	6.4	28
190	Effects of electron beam generated lattice defects on the periodic lattice distortion structure in 1T $\bar{1}$ MoS ₂ and 1T $\bar{1}$ MoSe ₂ thin layers. <i>Physical Review B</i> , 2019 , 99,	3.3	3
189	Metallic Twin Boundaries Boost the Hydrogen Evolution Reaction on the Basal Plane of Molybdenum Selenotellurides. <i>Advanced Energy Materials</i> , 2018 , 8, 1800031	21.8	66
188	Post-Synthesis Modifications of Two-Dimensional MoSe or MoTe by Incorporation of Excess Metal Atoms into the Crystal Structure. <i>ACS Nano</i> , 2018 , 12, 3975-3984	16.7	39
187	Hydrogen-assisted post-growth substitution of tellurium into molybdenum disulfide monolayers with tunable compositions. <i>Nanotechnology</i> , 2018 , 29, 145603	3.4	10

186	MoS ₂ Quantum Dots as Efficient Catalyst Materials for the Oxygen Evolution Reaction. <i>ACS Catalysis</i> , 2018 , 8, 1683-1689	13.1	135
185	Supported Two-Dimensional Materials under Ion Irradiation: The Substrate Governs Defect Production. <i>ACS Applied Materials & Interfaces</i> , 2018 , 10, 30827-30836	9.5	50
184	When defects are not defects. <i>Nature Materials</i> , 2018 , 17, 757-758	27	14
183	Revealing the Atomic Defects of WS ₂ Governing Its Distinct Optical Emissions. <i>Advanced Functional Materials</i> , 2018 , 28, 1704210	15.6	49
182	Reversible superdense ordering of lithium between two graphene sheets. <i>Nature</i> , 2018 , 564, 234-239	50.4	121
181	Observation of charge density waves in free-standing 1T-TaSe ₂ monolayers by transmission electron microscopy. <i>Applied Physics Letters</i> , 2018 , 113, 173103	3.4	14
180	Electron Irradiation-Induced Defects and Phase Transformations in Two-Dimensional Transition Metal Dichalcogenides. <i>Microscopy and Microanalysis</i> , 2018 , 24, 1592-1593	0.5	1
179	1T phase as an efficient hole injection layer to TMDs transistors: a universal approach to achieve p-type contacts. <i>2D Materials</i> , 2018 , 5, 031012	5.9	19
178	Nanostructuring few-layer graphene films with swift heavy ions for electronic application: tuning of electronic and transport properties. <i>Nanoscale</i> , 2018 , 10, 14499-14509	7.7	26
177	Atomic Defects and Doping of Monolayer NbSe. <i>ACS Nano</i> , 2017 , 11, 2894-2904	16.7	46
176	From Permeation to Cluster Arrays: Graphene on Ir(111) Exposed to Carbon Vapor. <i>Nano Letters</i> , 2017 , 17, 3105-3112	11.5	14
175	Two-dimensional MoS ₂ under ion irradiation: from controlled defect production to electronic structure engineering. <i>2D Materials</i> , 2017 , 4, 025078	5.9	99
174	Structural Transformations in Two-Dimensional Transition-Metal Dichalcogenide MoS under an Electron Beam: Insights from First-Principles Calculations. <i>Journal of Physical Chemistry Letters</i> , 2017 , 8, 3061-3067	6.4	68
173	Acoustic gravity waves and the atmospheric electric field perturbations accompanying them. <i>Geomagnetism and Aeronomy</i> , 2017 , 57, 335-346	0.9	4
172	Revisiting Hollandites: Channels Filling by Main-Group Elements Together with Transition Metals in Bi ₂ V ₈ O ₁₆ . <i>Chemistry of Materials</i> , 2017 , 29, 5558-5565	9.6	4
171	Tailoring the optical properties of atomically-thin WS ₂ via ion irradiation. <i>Nanoscale</i> , 2017 , 9, 11027-11034	7.7	62
170	Structural Distortions and Charge Density Waves in Iodine Chains Encapsulated inside Carbon Nanotubes. <i>Nano Letters</i> , 2017 , 17, 3694-3700	11.5	33
169	Engineering the Electronic Properties of Two-Dimensional Transition Metal Dichalcogenides by Introducing Mirror Twin Boundaries. <i>Advanced Electronic Materials</i> , 2017 , 3, 1600468	6.4	59

168	Creating nanoporous graphene with swift heavy ions. <i>Carbon</i> , 2017 , 114, 511-518	10.4	43
167	Charge equilibration times for slow highly charged ions in single layer graphene. <i>Journal of Physics: Conference Series</i> , 2017 , 875, 112001	0.3	
166	Engineering and modifying two-dimensional materials by electron beams. <i>MRS Bulletin</i> , 2017 , 42, 667-676	2	48
165	Interatomic Coulombic Decay: The Mechanism for Rapid Deexcitation of Hollow Atoms. <i>Physical Review Letters</i> , 2017 , 119, 103401	7.4	53
164	Substitutional carbon doping of free-standing and Ru-supported BN sheets: a first-principles study. <i>Journal of Physics Condensed Matter</i> , 2017 , 29, 415301	1.8	4
163	Luminescence of defects in the structural transformation of layered tin dichalcogenides. <i>Applied Physics Letters</i> , 2017 , 111, 262102	3.4	14
162	Vibrational Properties of Metal Phosphorus Trichalcogenides from First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2017 , 121, 27207-27217	3.8	36
161	Towards diluted magnetism in TaAs. <i>Physical Review Materials</i> , 2017 , 1,	3.2	1
160	From Point to Line Defects in Two-Dimensional Transition Metal Dichalcogenides: Insights from Transmission Electron Microscopy and First-Principles Calculations. <i>Carbon Nanostructures</i> , 2017 , 71-85	0.6	1
159	Study of acoustic emission signals during fracture shear deformation. <i>Acoustical Physics</i> , 2016 , 62, 505-513	13	10
158	Vibrational Properties of a Two-Dimensional Silica Kagome Lattice. <i>ACS Nano</i> , 2016 , 10, 10929-10935	16.7	12
157	Mechanical properties and current-carrying capacity of Al reinforced with graphene/BN nanoribbons: a computational study. <i>Nanoscale</i> , 2016 , 8, 20080-20089	7.7	14
156	Electron-Beam Induced Transformations of Layered Tin Dichalcogenides. <i>Nano Letters</i> , 2016 , 16, 4410-6	11.5	82
155	Nanostructured BN-Mg composites: features of interface bonding and mechanical properties. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 965-9	3.6	10
154	Phosphorene under electron beam: from monolayer to one-dimensional chains. <i>Nanoscale</i> , 2016 , 8, 7949-7957	7.57	33
153	Ultrafast electronic response of graphene to a strong and localized electric field. <i>Nature Communications</i> , 2016 , 7, 13948	17.4	91
152	Atomic scale microstructure and properties of Se-deficient two-dimensional MoSe ₂ . <i>ACS Nano</i> , 2015 , 9, 3274-83	16.7	176
151	Solubility of Boron, Carbon, and Nitrogen in Transition Metals: Getting Insight into Trends from First-Principles Calculations. <i>Journal of Physical Chemistry Letters</i> , 2015 , 6, 3263-3268	6.4	40

150	Comment on "Interfacial carbon nanoplatelet formation by ion irradiation of graphene on iridium(111)". <i>ACS Nano</i> , 2015 , 9, 4664-5	16.7	7
149	Line and rotational defects in boron-nitrene: Structure, energetics, and dependence on mechanical strain from first-principles calculations. <i>Physica Status Solidi (B): Basic Research</i> , 2015 , 252, 1725-1730	1.3	6
148	Three-fold rotational defects in two-dimensional transition metal dichalcogenides. <i>Nature Communications</i> , 2015 , 6, 6736	17.4	149
147	Single-Layer ReS ₂ Two-Dimensional Semiconductor with Tunable In-Plane Anisotropy. <i>ACS Nano</i> , 2015 , 9, 11249-57	16.7	286
146	Native defects in bulk and monolayer MoS ₂ from first principles. <i>Physical Review B</i> , 2015 , 91,	3.3	339
145	Xe irradiation of graphene on Ir(111): From trapping to blistering. <i>Physical Review B</i> , 2015 , 92,	3.3	24
144	Silicon and silicon-nitrogen impurities in graphene: Structure, energetics, and effects on electronic transport. <i>Physical Review B</i> , 2015 , 92,	3.3	21
143	Simulations of electromechanical shape transformations of Au nanoparticles. <i>Physica Status Solidi (B): Basic Research</i> , 2015 , 252, 144-148	1.3	2
142	Atomic scale interface design and characterisation. <i>Beilstein Journal of Nanotechnology</i> , 2015 , 6, 1708-113		2
141	Properties of individual dopant atoms in single-layer MoS ₂ : atomic structure, migration, and enhanced reactivity. <i>Advanced Materials</i> , 2014 , 26, 2857-61	24	229
140	Triazine-based graphitic carbon nitride: a two-dimensional semiconductor. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 7450-5	16.4	412
139	Triazine-Based Graphitic Carbon Nitride: a Two-Dimensional Semiconductor. <i>Angewandte Chemie</i> , 2014 , 126, 7580-7585	3.6	125
138	Electronic stopping power from first-principles calculations with account for core electron excitations and projectile ionization. <i>Physical Review B</i> , 2014 , 89,	3.3	59
137	Toward Stronger AlBN Nanotube Composite Materials: Insights into Bonding at the Al/BN Interface from First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2014 , 118, 26894-26901	3.8	21
136	Solid-State Growth of One- and Two-Dimensional Silica Structures on Metal Surfaces. <i>Journal of Physical Chemistry C</i> , 2014 , 118, 21001-21005	3.8	6
135	Semiconductor to Metal to Half-Metal Transition in Pt-Embedded Zigzag Graphene Nanoribbons. <i>Journal of Physical Chemistry C</i> , 2014 , 118, 16133-16139	3.8	16
134	Atomic structure and dynamic behaviour of truly one-dimensional ionic chains inside carbon nanotubes. <i>Nature Materials</i> , 2014 , 13, 1050-4	27	66
133	Coronene encapsulation in single-walled carbon nanotubes: stacked columns, peapods, and nanoribbons. <i>ChemPhysChem</i> , 2014 , 15, 1660-5	3.2	24

132	Fabrication and atomic structure of size-selected, layered MoS ₂ clusters for catalysis. <i>Nanoscale</i> , 2014 , 6, 12463-9	7.7	33
131	Doping Properties and Phase Transition in Single-Layer MoS ₂ . <i>Microscopy and Microanalysis</i> , 2014 , 20, 1750-1751	0.5	1
130	Inorganic Two-Dimensional Materials under Electron Irradiation: Stability, Evolution of the Atomic Structure, and Beam-Mediated Doping. <i>Microscopy and Microanalysis</i> , 2014 , 20, 1756-1757	0.5	
129	Interfacial carbon nanoplatelet formation by ion irradiation of graphene on iridium(111). <i>ACS Nano</i> , 2014 , 8, 12208-18	16.7	23
128	Experimental observation of boron nitride chains. <i>ACS Nano</i> , 2014 , 8, 11950-7	16.7	57
127	Charged Point Defects in the Flatland: Accurate Formation Energy Calculations in Two-Dimensional Materials. <i>Physical Review X</i> , 2014 , 4,	9.1	49
126	Atomic scale study of the life cycle of a dislocation in graphene from birth to annihilation. <i>Nature Communications</i> , 2013 , 4, 2098	17.4	133
125	Strains induced by point defects in graphene on a metal. <i>Physical Review Letters</i> , 2013 , 111, 085501	7.4	43
124	Electronic structure of boron nitride sheets doped with carbon from first-principles calculations. <i>Physical Review B</i> , 2013 , 87,	3.3	129
123	Electronic structures and optical properties of realistic transition metal dichalcogenide heterostructures from first principles. <i>Physical Review B</i> , 2013 , 88,	3.3	342
122	Defects in bilayer silica and graphene: common trends in diverse hexagonal two-dimensional systems. <i>Scientific Reports</i> , 2013 , 3, 3482	4.9	71
121	From point to extended defects in two-dimensional MoS ₂ : Evolution of atomic structure under electron irradiation. <i>Physical Review B</i> , 2013 , 88,	3.3	334
120	Ion-irradiation-induced defects in isotopically-labeled two layered graphene: enhanced in-situ annealing of the damage. <i>Advanced Materials</i> , 2013 , 25, 1004-9	24	66
119	Doped Graphene as a Material for Oxygen Reduction Reaction in Hydrogen Fuel Cells: A Computational Study. <i>ACS Catalysis</i> , 2013 , 3, 159-165	13.1	95
118	Chirality-dependent reactivity of individual single-walled carbon nanotubes. <i>Small</i> , 2013 , 9, 1379-86	11	33
117	Ion impacts on graphene/Ir(111): interface channeling, vacancy funnels, and a nanomesh. <i>Nano Letters</i> , 2013 , 13, 1948-55	11.5	73
116	Dual origin of defect magnetism in graphene and its reversible switching by molecular doping. <i>Nature Communications</i> , 2013 , 4, 2010	17.4	189
115	In situ growth of cellular two-dimensional silicon oxide on metal substrates. <i>ACS Nano</i> , 2013 , 7, 5175-80	16.7	28

114	Two-Dimensional Transition Metal Dichalcogenide Alloys: Stability and Electronic Properties. <i>Journal of Physical Chemistry Letters</i> , 2012 , 3, 3652-6	6.4	241
113	Atomistic description of electron beam damage in nitrogen-doped graphene and single-walled carbon nanotubes. <i>ACS Nano</i> , 2012 , 6, 8837-46	16.7	101
112	Effects of confinement and environment on the electronic structure and exciton binding energy of MoS ₂ from first principles. <i>Physical Review B</i> , 2012 , 86,	3.3	474
111	The Role of Stable and Mobile Carbon Adspecies in Copper-Promoted Graphene Growth. <i>Journal of Physical Chemistry C</i> , 2012 , 116, 5802-5809	3.8	64
110	Two-dimensional transition metal dichalcogenides under electron irradiation: defect production and doping. <i>Physical Review Letters</i> , 2012 , 109, 035503	7.4	768
109	Accurate measurement of electron beam induced displacement cross sections for single-layer graphene. <i>Physical Review Letters</i> , 2012 , 108, 196102	7.4	326
108	Are we van der Waals ready?. <i>Journal of Physics Condensed Matter</i> , 2012 , 24, 424218	1.8	112
107	Stability of graphene edges under electron beam: equilibrium energetics versus dynamic effects. <i>ACS Nano</i> , 2012 , 6, 671-6	16.7	104
106	Gold-embedded zigzag graphene nanoribbons as spin gapless semiconductors. <i>Physical Review B</i> , 2012 , 86,	3.3	38
105	Spin-half paramagnetism in graphene induced by point defects. <i>Nature Physics</i> , 2012 , 8, 199-202	16.2	638
104	van der Waals bonding in layered compounds from advanced density-functional first-principles calculations. <i>Physical Review Letters</i> , 2012 , 108, 235502	7.4	665
103	Direct imaging of a two-dimensional silica glass on graphene. <i>Nano Letters</i> , 2012 , 12, 1081-6	11.5	206
102	Engineering the atomic structure of carbon nanotubes by a focused electron beam: new morphologies at the sub-nanometer scale. <i>ChemPhysChem</i> , 2012 , 13, 2596-600	3.2	15
101	Atom-by-atom observation of grain boundary migration in graphene. <i>Nano Letters</i> , 2012 , 12, 3168-73	11.5	154
100	Ion irradiation tolerance of graphene as studied by atomistic simulations. <i>Applied Physics Letters</i> , 2012 , 100, 233108	3.4	39
99	Strain Fields and Electronic Structure of Vacancy-Type Defects in Graphene from First-Principles Simulation. <i>NATO Science for Peace and Security Series B: Physics and Biophysics</i> , 2012 , 49-59	0.2	
98	Tuning electronic and magnetic properties of zigzag graphene nanoribbons by large-scale bending. <i>Applied Physics Letters</i> , 2012 , 100, 263115	3.4	11
97	Atomistic simulations of the implantation of low-energy boron and nitrogen ions into graphene. <i>Physical Review B</i> , 2011 , 83,	3.3	114

96	Characterization of ion-irradiation-induced defects in multi-walled carbon nanotubes. <i>New Journal of Physics</i> , 2011 , 13, 073004	2.9	41
95	Cutting and controlled modification of graphene with ion beams. <i>Nanotechnology</i> , 2011 , 22, 175306	3.4	119
94	From point defects in graphene to two-dimensional amorphous carbon. <i>Physical Review Letters</i> , 2011 , 106, 105505	7.4	582
93	Structural defects in graphene. <i>ACS Nano</i> , 2011 , 5, 26-41	16.7	2388
92	Stone-Wales-type transformations in carbon nanostructures driven by electron irradiation. <i>Physical Review B</i> , 2011 , 83,	3.3	199
91	Synthesis of graphene nanoribbons encapsulated in single-walled carbon nanotubes. <i>Nano Letters</i> , 2011 , 11, 4352-6	11.5	148
90	Bound and free self-interstitial defects in graphite and bilayer graphene: A computational study. <i>Physical Review B</i> , 2011 , 84,	3.3	31
89	Attractive interaction between transition-metal atom impurities and vacancies in graphene: a first-principles study. <i>Theoretical Chemistry Accounts</i> , 2011 , 129, 625-630	1.9	89
88	Production of defects in hexagonal boron nitride monolayer under ion irradiation. <i>Nuclear Instruments & Methods in Physics Research B</i> , 2011 , 269, 1327-1331	1.2	43
87	Mechanism of swift chemical sputtering: Comparison of Be/C/W dimer bond breaking. <i>Nuclear Instruments & Methods in Physics Research B</i> , 2011 , 269, 1257-1261	1.2	8
86	Berseneva, Krasheninnikov, and Nieminen Reply:. <i>Physical Review Letters</i> , 2011 , 107,	7.4	5
85	Mechanisms of postsynthesis doping of boron nitride nanostructures with carbon from first-principles simulations. <i>Physical Review Letters</i> , 2011 , 107, 035501	7.4	84
84	Boron nitride formation on magnesium studied by ab initio calculations. <i>Physical Review B</i> , 2010 , 81,	3.3	5
83	Response of mechanically strained nanomaterials to irradiation: Insight from atomistic simulations. <i>Physical Review B</i> , 2010 , 82,	3.3	26
82	Effects of ion bombardment on a two-dimensional target: Atomistic simulations of graphene irradiation. <i>Physical Review B</i> , 2010 , 81,	3.3	303
81	A first-principles study on magnetic coupling between carbon adatoms on graphene. <i>New Journal of Physics</i> , 2010 , 12, 113021	2.9	29
80	Migration of gold atoms in graphene ribbons: Role of the edges. <i>Physical Review B</i> , 2010 , 81,	3.3	39
79	Effect of iron nanoparticle geometry on the energetics of carbon interstitials. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2010 , 7, NA-NA		1

78	Defect-induced junctions between single- or double-wall carbon nanotubes and metal crystals. <i>Nanoscale</i> , 2010 , 2, 901-5	7.7	18
77	Migration and localization of metal atoms on strained graphene. <i>Physical Review Letters</i> , 2010 , 105, 196102	7.7	281
76	Electron knock-on damage in hexagonal boron nitride monolayers. <i>Physical Review B</i> , 2010 , 82,	3.3	212
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