Gun-Do Lee

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.

2,593 28 49 g-index

90 2,911 8.1 4.79 ext. papers ext. citations avg, IF L-index

#	Paper	IF	Citations
88	OH molecule-involved formation of point defects in monolayer graphene. <i>Nanotechnology</i> , 2021 , 32, 025704	3.4	
87	Density Functional Theory Study of Edge-Induced Atomic-Scale Structural Phase Transitions of MoS2 Nanocrystals: Implications for a High-Performance Catalyst. <i>ACS Applied Nano Materials</i> , 2021 , 4, 5496-5502	5.6	2
86	Atomically Precise Control of Carbon Insertion into hBN Monolayer Point Vacancies using a Focused Electron Beam Guide. <i>Small</i> , 2021 , 17, e2100693	11	3
85	Quasi-graphitic carbon shell-induced Cu confinement promotes electrocatalytic CO reduction toward C products. <i>Nature Communications</i> , 2021 , 12, 3765	17.4	17
84	Reaction mechanisms of chlorine reduction on hydroxylated alumina in titanium nitride growth: First principles study. <i>Applied Surface Science</i> , 2021 , 550, 149391	6.7	2
83	Two-dimensional iodine-monofluoride epitaxy on WSe2. Npj 2D Materials and Applications, 2021, 5,	8.8	2
82	Phase Engineering of Transition Metal Dichalcogenides via a Thermodynamically Designed Gas-Solid Reaction. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 8430-8439	6.4	
81	Thermodynamically driven self-formation of copper-embedded nitrogen-doped carbon nanofiber catalysts for a cascade electroreduction of carbon dioxide to ethylene. <i>Journal of Materials Chemistry A</i> , 2020 , 8, 11632-11641	13	17
80	Direct observation and catalytic role of mediator atom in 2D materials. <i>Science Advances</i> , 2020 , 6, eaba ²	1942;	5
79	The influence of hydrogen concentration in amorphous carbon films on mechanical properties and fluorine penetration: a density functional theory and molecular dynamics study <i>RSC Advances</i> , 2020 , 10, 6822-6830	3.7	4
78	Thermodynamic insights into interfacial interactions in TiN/amorphous Al2O3 heterostructures: ab initio molecular dynamics and first principles investigation. <i>Inorganic Chemistry Frontiers</i> , 2020 , 7, 4347-	4356	О
77	Scanning Moir Fringe Method: A Superior Approach to Perceive Defects, Interfaces, and Distortion in 2D Materials. <i>ACS Nano</i> , 2020 , 14, 6034-6042	16.7	6
76	Analysis of surface adsorption kinetics of SiH4 and Si2H6 for deposition of a hydrogenated silicon thin film using intermediate pressure SiH4 plasmas. <i>Applied Surface Science</i> , 2019 , 496, 143728	6.7	7
75	Atomic Structure and Dynamics of Epitaxial Platinum Bilayers on Graphene. ACS Nano, 2019, 13, 12162-	126.70	12
74	Atomic Scale Imaging of Reversible Ring Cyclization in Graphene Nanoconstrictions. <i>ACS Nano</i> , 2019 , 13, 2379-2388	16.7	2
73	Effects of nitrogen doping in amorphous carbon layers on the diffusion of fluorine atoms: A first-principles study. <i>Journal of Applied Physics</i> , 2019 , 125, 155701	2.5	8
7 ²	Bonding structure and etching characteristics of amorphous carbon for a hardmask deposited by DC sputtering. <i>Carbon</i> , 2019 , 154, 277-284	10.4	8

(2016-2019)

71	Predictive fabrication of Ni phosphide embedded in carbon nanofibers as active and stable electrocatalysts. <i>Journal of Materials Chemistry A</i> , 2019 , 7, 7451-7458	13	17
70	First principles investigation on energetics, structure, and mechanical properties of amorphous carbon films doped with B, N, and Cl. <i>Scientific Reports</i> , 2019 , 9, 18961	4.9	3
69	Single-Crystalline Nanobelts Composed of Transition Metal Ditellurides. <i>Advanced Materials</i> , 2018 , 30, e1707260	24	15
68	Effects of H and N treatment for BH dosing process on TiN surfaces during atomic layer deposition: an study <i>RSC Advances</i> , 2018 , 8, 21164-21173	3.7	3
67	Overall reaction mechanism for a full atomic layer deposition cycle of W films on TiN surfaces: first-principles study <i>RSC Advances</i> , 2018 , 8, 39039-39046	3.7	3
66	Solid-Phase Epitaxial Growth of an Alumina Layer Having a Stacking-Mismatched Domain Structure of the Intermediate Phase. <i>ACS Applied Materials & Samp; Interfaces</i> , 2018 , 10, 41487-41496	9.5	3
65	The impact of substrate surface defects on the properties of two-dimensional van der Waals heterostructures. <i>Nanoscale</i> , 2018 , 10, 19212-19219	7.7	9
64	Lattice contraction with boron doping in fully strained SiGe epitaxial layers. <i>Japanese Journal of Applied Physics</i> , 2018 , 57, 065504	1.4	4
63	Graphene as a flexible template for controlling magnetic interactions between metal atoms. Journal of Physics Condensed Matter, 2017 , 29, 085001	1.8	1
62	Improved performance of AlGaN-based deep ultraviolet light-emitting diodes with nano-patterned AlN/sapphire substrates. <i>Applied Physics Letters</i> , 2017 , 110, 191103	3.4	63
61	Electrically Driven Reversible Phase Changes in Layered In Se Crystalline Film. <i>Advanced Materials</i> , 2017 , 29, 1703568	24	45
60	Point defects in turbostratic stacked bilayer graphene. <i>Nanoscale</i> , 2017 , 9, 13725-13730	7.7	8
59	Gaseous Nanocarving-Mediated Carbon Framework with Spontaneous Metal Assembly for Structure-Tunable Metal/Carbon Nanofibers. <i>Advanced Materials</i> , 2017 , 29, 1702958	24	10
58	Dissociation reaction of B2H6 on TiN surfaces during atomic layer deposition: first-principles study. <i>RSC Advances</i> , 2017 , 7, 55750-55755	3.7	9
57	In Situ Atomic Level Dynamics of Heterogeneous Nucleation and Growth of Graphene from Inorganic Nanoparticle Seeds. <i>ACS Nano</i> , 2016 , 10, 9397-9410	16.7	8
56	Atomic Structure and Spectroscopy of Single Metal (Cr, V) Substitutional Dopants in Monolayer MoS. <i>ACS Nano</i> , 2016 , 10, 10227-10236	16.7	77
55	Direct imaging of rotating molecules anchored on graphene. <i>Nanoscale</i> , 2016 , 8, 13174-80	7.7	9
54	Elongated Silicon-Carbon Bonds at Graphene Edges. <i>ACS Nano</i> , 2016 , 10, 142-9	16.7	15

53	Electron beam-formed ferromagnetic defects on MoS surface along 1 T phase transition. <i>Scientific Reports</i> , 2016 , 6, 38730	4.9	19
52	Detailed Atomic Reconstruction of Extended Line Defects in Monolayer MoS2. <i>ACS Nano</i> , 2016 , 10, 541	913607	115
51	Layer-dependent modulation of tungsten disulfide photoluminescence by lateral electric fields. <i>ACS Nano</i> , 2015 , 9, 2740-8	16.7	39
50	Rotating Anisotropic Crystalline Silicon Nanoclusters in Graphene. <i>ACS Nano</i> , 2015 , 9, 9497-506	16.7	13
49	Atomic Level Distributed Strain within Graphene Divacancies from Bond Rotations. <i>ACS Nano</i> , 2015 , 9, 8599-608	16.7	20
48	Atomic Structure of Graphene Subnanometer Pores. ACS Nano, 2015, 9, 11599-607	16.7	56
47	Thermally Induced Dynamics of Dislocations in Graphene at Atomic Resolution. ACS Nano, 2015, 9, 1006	56 .8.5	27
46	Partial Dislocations in Graphene and Their Atomic Level Migration Dynamics. <i>Nano Letters</i> , 2015 , 15, 5950-5	11.5	33
45	Spatially dependent lattice deformations for dislocations at the edges of graphene. <i>ACS Nano</i> , 2015 , 9, 656-62	16.7	11
44	Atomic-scale mechanism of grain boundary motion in graphene. <i>Carbon</i> , 2015 , 84, 146-150	10.4	8
43	Hydrogen-free graphene edges. <i>Nature Communications</i> , 2014 , 5, 3040	17.4	63
42	The role of the bridging atom in stabilizing odd numbered graphene vacancies. <i>Nano Letters</i> , 2014 , 14, 3972-80	11.5	36
41	Stability and dynamics of the tetravacancy in graphene. Nano Letters, 2014, 14, 1634-42	11.5	57
40	Atomic structure and dynamics of metal dopant pairs in graphene. <i>Nano Letters</i> , 2014 , 14, 3766-72	11.5	168
39	Detailed formation processes of stable dislocations in graphene. <i>Nanoscale</i> , 2014 , 6, 14836-44	7.7	24
38	Extended Klein edges in graphene. ACS Nano, 2014 , 8, 12272-9	16.7	31
37	Atomistic processes of grain boundary motion and annihilation in graphene. <i>Journal of Physics Condensed Matter</i> , 2013 , 25, 155301	1.8	6
36	Bond length and charge density variations within extended arm chair defects in graphene. <i>ACS Nano</i> , 2013 , 7, 9860-6	16.7	32

(2006-2013)

35	Rippling graphene at the nanoscale through dislocation addition. Nano Letters, 2013, 13, 4937-44	11.5	54
34	Formation and development of dislocation in graphene. <i>Applied Physics Letters</i> , 2013 , 102, 021603	3.4	28
33	Less strained and more efficient GaN light-emitting diodes with embedded silica hollow nanospheres. <i>Scientific Reports</i> , 2013 , 3, 3201	4.9	33
32	Reduction of graphene damages during the fabrication of InGaN/GaN light emitting diodes with graphene electrodes. <i>Nanotechnology</i> , 2012 , 23, 425302	3.4	8
31	Improved emission efficiency of a-plane GaN light emitting diodes with silica nano-spheres integrated into a-plane GaN buffer layer. <i>Applied Physics Letters</i> , 2012 , 100, 191116	3.4	6
30	Growth mechanism of highly uniform InAs/GaAs quantum dot with periodic arsine interruption by metalorganic chemical vapor deposition. <i>Journal of Applied Physics</i> , 2011 , 110, 044302	2.5	3
29	Dynamics and stability of divacancy defects in graphene. <i>Physical Review B</i> , 2011 , 84,	3.3	80
28	The role of pentagonBeptagon pair defect in carbon nanotube: The center of vacancy reconstruction. <i>Applied Physics Letters</i> , 2010 , 97, 093106	3.4	14
27	Reconstruction and evaporation at graphene nanoribbon edges. <i>Physical Review B</i> , 2010 , 81,	3.3	53
26	High quality Ge epitaxial layers on Si by ultrahigh vacuum chemical vapor deposition. <i>Thin Solid Films</i> , 2009 , 517, 3990-3994	2.2	12
25	Stability of dislocation defect with two pentagon-heptagon pairs in graphene. <i>Physical Review B</i> , 2008 , 78,	3.3	93
24	Effect of C incorporation on relaxation of SiGe/Si. <i>Applied Physics Letters</i> , 2008 , 93, 221902	3.4	5
23	The formation of pentagon-heptagon pair defect by the reconstruction of vacancy defects in carbon nanotube. <i>Applied Physics Letters</i> , 2008 , 92, 043104	3.4	33
22	Effect of interstitial C incorporation on the Raman scattering of Si1 III GexCy epitaxial layer. <i>Applied Physics Letters</i> , 2008 , 92, 061906	3.4	1
21	Atomistic simulation studies of complex carbon and silicon systems using environment-dependent tight-binding potentials. <i>Scientific Modeling and Simulation SMNS</i> , 2008 , 15, 97-121		4
20	Atomistic simulation studies of complex carbon and silicon systems using environment-dependent tight-binding potentials. <i>Lecture Notes in Computational Science and Engineering</i> , 2008 , 97-121	0.3	
19	Formation of carbon nanotube semiconductor-metal intramolecular junctions by self-assembly of vacancy defects. <i>Physical Review B</i> , 2007 , 76,	3.3	30
18	Formation of flat, relaxed Si1\(\text{IGex alloys on Si(001)}\) without buffer layers. <i>Applied Physics Letters</i> , 2006 , 88, 122103	3.4	6

17	Diffusion of adatom in the selective epitaxial growth of Si(100): A molecular dynamics study. <i>Applied Physics Letters</i> , 2006 , 88, 231909	3.4	О
16	Vacancy defects and the formation of local haeckelite structures in graphene from tight-binding molecular dynamics. <i>Physical Review B</i> , 2006 , 74,	3.3	77
15	Diffusion, coalescence, and reconstruction of vacancy defects in graphene layers. <i>Physical Review Letters</i> , 2005 , 95, 205501	7.4	416
14	The emission wavelength tuning of InAs/InP quantum dots with thin GaAs, InGaAs, InP capping layers by MOCVD. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2005 , 26, 169-173	3	2
13	Microscopic study on the behavior of the {311} facet in the selective epitaxial growth of Si(100). <i>Applied Physics Letters</i> , 2004 , 85, 4624-4626	3.4	2
12	Theoretical study on the temperature-induced structural transition of the Si(1 1 3) surface. <i>Science</i> , 2004 , 559, 63-69	1.8	8
11	Isotropic/Anisotropic Selective Epitaxial Growth of Si on Local Oxidation of Silicon (LOCOS) Patterned Si (100) Substrate by Cold Wall Ultrahigh Vacuum Chemical Vapor Deposition (UHV-CVD). <i>Japanese Journal of Applied Physics</i> , 2003 , 42, 3966-3970	1.4	3
10	Structural evolution of the Si(113) surface: Ab initio and tight-binding molecular dynamics calculations. <i>Physical Review B</i> , 2003 , 68,	3.3	17
9	Heat-induced transformation of nanodiamond into a tube-shaped fullerene: a molecular dynamics simulation. <i>Physical Review Letters</i> , 2003 , 91, 265701	7.4	42
8	Catalytic decomposition of acetylene on Fe(001): A first-principles study. <i>Physical Review B</i> , 2002 , 66,	3.3	22
7	ADDIMER DIFFUSIION ON THE Si(100) SURFACE. Surface Review and Letters, 1999 , 06, 1015-1023	1.1	1
6	Addimer diffusion along the trough between dimer rows on Si(001). Surface Science, 1999 , 426, L427-L4	1 3;2 8	13
5	Ad-Dimer Diffusion between Trough and Dimer Row on Si(100). <i>Physical Review Letters</i> , 1998 , 81, 5872-	5,87,5	31
4	Ab initiopseudopotential study of the structural and electronic properties of ZnTe under high pressure. <i>Journal of Physics Condensed Matter</i> , 1997 , 9, 6619-6631	1.8	10
3	Microscopic study of the pressure-induced structural phase transition of ZnTe. <i>Physical Review B</i> , 1996 , 53, R7622-R7625	3.3	38
2	Environment-dependent tight-binding potential model. <i>Physical Review B</i> , 1996 , 53, 979-982	3.3	214
1	Role of d electrons in the zinc-blende semiconductors ZnS, ZnSe, and ZnTe. <i>Physical Review B</i> , 1995 , 52, 1459-1462	3.3	76