

Chuong V Nguyen

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.

154
papers

3,086
citations

31
h-index

47
g-index

160
ext. papers

4,013
ext. citations

3.3
avg, IF

6.1
L-index

#	Paper	IF	Citations
154	Tunable type-II band alignment and electronic structure of C ₃ N ₄ /MoSi ₂ N ₄ heterostructure: Interlayer coupling and electric field. <i>Physical Review B</i> , 2022 , 105,	3.3	9
153	Two-Dimensional Metal/Semiconductor Contact in a Janus MoSH/MoSiN van der Waals Heterostructure.. <i>Journal of Physical Chemistry Letters</i> , 2022 , 2576-2582	6.4	5
152	Two-dimensional XY monolayers (X = Al, Ga, In; Y = N, P, As) with a double layer hexagonal structure: A first-principles perspective. <i>Applied Surface Science</i> , 2022 , 590, 152998	6.7	3
151	Monoelemental two-dimensional iodine nanosheets: a first-principles study of the electronic and optical properties. <i>Journal Physics D: Applied Physics</i> , 2022 , 55, 135104	3	
150	Intriguing interfacial characteristics of the CS contact with MX (M = Mo, W; X = S, Se, Te) and MXY ((X \ Y) = S, Se, Te) monolayers.. <i>RSC Advances</i> , 2022 , 12, 12292-12302	3.7	0
149	Theoretical prediction of Janus PdXO (X = S, Se, Te) monolayers: structural, electronic, and transport properties.. <i>RSC Advances</i> , 2022 , 12, 12971-12977	3.7	0
148	First-principles insights onto structural, electronic and optical properties of Janus monolayers CrXO (X = S, Se, Te).. <i>RSC Advances</i> , 2021 , 11, 39672-39679	3.7	0
147	Electric gating and interlayer coupling controllable electronic structure and Schottky contact of graphene/Bil ₃ van der Waals heterostructure. <i>Physical Review B</i> , 2021 , 103,	3.3	13
146	Effects of La and Ce doping on electronic structure and optical properties of janus MoS ₂ monolayer. <i>Superlattices and Microstructures</i> , 2021 , 151, 106841	2.8	1
145	Nonlinear magneto-optical absorption in a finite semi-parabolic quantum well. <i>Optical and Quantum Electronics</i> , 2021 , 53, 1	2.4	5
144	Interfacial Electronic Properties and Tunable Contact Types in Graphene/Janus MoGeSiN Heterostructures. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 3934-3940	6.4	17
143	Electronic, optical, and thermoelectric properties of Janus In-based monochalcogenides. <i>Journal of Physics Condensed Matter</i> , 2021 , 33,	1.8	9
142	Ab initio prediction of semiconductivity in a novel two-dimensional SbX (X= S, Se, Te) monolayers with orthorhombic structure. <i>Scientific Reports</i> , 2021 , 11, 10366	4.9	20
141	Two-Dimensional Boron Phosphide/MoGeN van der Waals Heterostructure: A Promising Tunable Optoelectronic Material. <i>Journal of Physical Chemistry Letters</i> , 2021 , 12, 5076-5084	6.4	21
140	Point Defects in a Two-Dimensional ZnSnN ₂ Nanosheet: A First-Principles Study on the Electronic and Magnetic Properties. <i>Journal of Physical Chemistry C</i> , 2021 , 125, 13067-13075	3.8	15
139	Study of the Elastic Properties of the Energetic Molecular Crystals Using Density Functionals with van der Waals Corrections. <i>ACS Omega</i> , 2021 , 6, 642-648	3.9	4
138	Structural, elastic, and electronic properties of chemically functionalized boron phosphide monolayer.. <i>RSC Advances</i> , 2021 , 11, 8552-8558	3.7	3

137	Stacking effects in van der Waals heterostructures of blueP and Janus XYO (X = Ti, Zr, Hf; Y = S, Se) monolayers.. <i>RSC Advances</i> , 2021 , 11, 12189-12199	3.7	4
136	Outstanding elastic, electronic, transport and optical properties of a novel layered material CF: first-principles study.. <i>RSC Advances</i> , 2021 , 11, 23280-23287	3.7	3
135	First-principles study of the electronic structures and optical and photocatalytic performances of van der Waals heterostructures of SiS, P and SiC monolayers.. <i>RSC Advances</i> , 2021 , 11, 14263-14268	3.7	6
134	A van der Waals heterostructure of MoS ₂ /MoSi ₂ N ₄ : a first-principles study. <i>New Journal of Chemistry</i> , 2021 , 45, 8291-8296	3.6	20
133	Two-dimensional van der Waals graphene/transition metal nitride heterostructures as promising high-performance nanodevices. <i>New Journal of Chemistry</i> , 2021 , 45, 5509-5516	3.6	12
132	Theoretical prediction of electronic, transport, optical, and thermoelectric properties of Janus monolayers In ₂ XO (X=S,Se,Te). <i>Physical Review B</i> , 2021 , 103,	3.3	39
131	Oxygenation of Janus group III monochalcogenides: First-principles insights into GaInXO (X=S, Se, Te) monolayers. <i>Physical Review B</i> , 2021 , 104,	3.3	12
130	A theoretical study on elastic, electronic, transport, optical and thermoelectric properties of Janus SnSO monolayer. <i>Journal Physics D: Applied Physics</i> , 2021 , 54, 475306	3	3
129	Anisotropy of effective masses induced by strain in Janus MoS ₂ Se and WSe ₂ monolayers. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2021 , 134, 114826	3	2
128	Electronic and optical properties of a Janus SnSSe monolayer: effects of strain and electric field. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 11637-11643	3.6	33
127	Magneto-optical absorption in silicene and germanene induced by electric and Zeeman fields. <i>Physical Review B</i> , 2020 , 101,	3.3	17
126	Graphene hetero-multilayer on layered platinum mineral jacutingaite (Pt ₂ HgSe ₃): van der Waals heterostructures with novel optoelectronic and thermoelectric performances. <i>Journal of Materials Chemistry A</i> , 2020 , 8, 13248-13260	13	44
125	First-principles investigation of nonmetal doped single-layer BiOBr as a potential photocatalyst with a low recombination rate. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 15354-15364	3.6	45
124	Pyramidal core-shell quantum dot under applied electric and magnetic fields. <i>Scientific Reports</i> , 2020 , 10, 8961	4.9	14
123	Interlayer coupling and electric field controllable Schottky barriers and contact types in graphene/PbI ₂ heterostructures. <i>Physical Review B</i> , 2020 , 101,	3.3	45
122	Low-energy bands and optical properties of monolayer WS ₂ . <i>Optik</i> , 2020 , 209, 164581	2.5	2
121	First-principles prediction of chemically functionalized InN monolayers: electronic and optical properties.. <i>RSC Advances</i> , 2020 , 10, 10731-10739	3.7	10
120	Effects of electric field and strain engineering on the electronic properties, band alignment and enhanced optical properties of ZnO/Janus ZrSSe heterostructures.. <i>RSC Advances</i> , 2020 , 10, 9824-9832	3.7	9

119	Effects of different surface functionalization on the electronic properties and contact types of graphene/functionalized-GeC van der Waals heterostructures. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 7952-7961	3.6	15
118	A first-principles study of electronic structure and photocatalytic performance of GaN-MX (M = Mo, W; X= S, Se) van der Waals heterostructures.. <i>RSC Advances</i> , 2020 , 10, 24683-24690	3.7	9
117	Electronic structure, optoelectronic properties and enhanced photocatalytic response of GaN-GeC van der Waals heterostructures: a first principles study.. <i>RSC Advances</i> , 2020 , 10, 24127-24133	3.7	11
116	Electronic and optoelectronic properties of van der Waals heterostructure based on graphene-like GaN, blue phosphorene, SiC, and ZnO: A first principles study. <i>Journal of Applied Physics</i> , 2020 , 127, 245302	3.5	7
115	Van der Waals heterostructures of SiC and Janus MSSe (M = Mo, W) monolayers: a first principles study.. <i>RSC Advances</i> , 2020 , 10, 25801-25807	3.7	8
114	Computational prediction of electronic and optical properties of Janus Ga ₂ SeTe monolayer. <i>Journal Physics D: Applied Physics</i> , 2020 , 53, 455302	3	13
113	Stacking and electric field effects on the band alignment and electronic properties of the GeC/GaSe heterostructure. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020 , 120, 114050	3	7
112	The characteristics of defective ZrS ₂ monolayers adsorbed various gases on S-vacancies: A first-principles study. <i>Superlattices and Microstructures</i> , 2020 , 140, 106454	2.8	12
111	Electronic structure and optical performance of PbI ₂ /SnSe ₂ heterostructure. <i>Chemical Physics</i> , 2020 , 533, 110736	2.3	2
110	Computational insights into structural, electronic and optical characteristics of GeC/CN van der Waals heterostructures: effects of strain engineering and electric field.. <i>RSC Advances</i> , 2020 , 10, 2967-2974	3.7	7
109	Magneto-optical transport properties of monolayer transition metal dichalcogenides. <i>Physical Review B</i> , 2020 , 101,	3.3	44
108	Physicochemical properties of L- and DL-valine: first-principles calculations. <i>Amino Acids</i> , 2020 , 52, 425-433	3.5	2
107	Stark and Zeeman effects on the topological phase and transport properties of topological crystalline insulator thin films. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 12129-12139	3.6	
106	Electronic and photocatalytic performance of boron phosphide-blue phosphorene vdW heterostructures. <i>Applied Surface Science</i> , 2020 , 523, 146483	6.7	47
105	Magneto-optical absorption in Pöschl-Teller-like quantum well. <i>Physica B: Condensed Matter</i> , 2020 , 592, 412279	2.8	3
104	Tuning the electronic, photocatalytic and optical properties of hydrogenated InN monolayer by biaxial strain and electric field. <i>Chemical Physics</i> , 2020 , 532, 110677	2.3	8
103	Electronic, optical and photocatalytic properties of fully hydrogenated GeC monolayer. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020 , 117, 113857	3	6
102	A type-II GaSe/HfS ₂ van der Waals heterostructure as promising photocatalyst with high carrier mobility. <i>Applied Surface Science</i> , 2020 , 534, 147607	6.7	40

101	Tunable electronic properties of the dynamically stable layered mineral PtHgSe (Jacutingaite). <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 24471-24479	3.6	14
100	Janus monolayer PtSSe under external electric field and strain: A first principles study on electronic structure and optical properties. <i>Superlattices and Microstructures</i> , 2020 , 147, 106683	2.8	39
99	Strain engineering of the electro-optical and photocatalytic properties of single-layered Janus MoSSe: First principles calculations. <i>Optik</i> , 2020 , 224, 165503	2.5	4
98	Vertical two-dimensional layered conjugated porous organic network structures of poly-benzimidazobenzophenanthroline (BBL): A first-principles study. <i>Applied Physics Letters</i> , 2020 , 117, 233101	3.4	14
97	Interfacial characteristics, Schottky contact, and optical performance of a graphene/Ga2SSe van der Waals heterostructure: Strain engineering and electric field tunability. <i>Physical Review B</i> , 2020 , 102,	3.3	55
96	Electronic structures, and optical and photocatalytic properties of the BPBSe van der Waals heterostructures. <i>New Journal of Chemistry</i> , 2020 , 44, 14964-14969	3.6	5
95	First principles study of structural, optoelectronic and photocatalytic properties of SnS, SnSe monolayers and their van der Waals heterostructure. <i>Chemical Physics</i> , 2020 , 539, 110939	2.3	9
94	Investigation of strain and doping on the electronic properties of single layers of CN and CN: a first principles study.. <i>RSC Advances</i> , 2020 , 10, 27743-27751	3.7	23
93	The mechanical, electronic, optical and thermoelectric properties of two-dimensional honeycomb-like of XSb (X = Si, Ge, Sn) monolayers: a first-principles calculations.. <i>RSC Advances</i> , 2020 , 10, 30398-30405	3.7	20
92	Oxygen Vacancies in the Single Layer of Ti2CO2 MXene: Effects of Gating Voltage, Mechanical Strain, and Atomic Impurities. <i>Physica Status Solidi (B): Basic Research</i> , 2020 , 257, 2000343	1.3	14
91	Electronic and photocatalytic properties of two-dimensional boron phosphide/SiC van der Waals heterostructure with direct type-II band alignment: a first principles study.. <i>RSC Advances</i> , 2020 , 10, 32027-32033	3.7	33
90	Two-dimensional silicon bismotide (SiBi) monolayer with a honeycomb-like lattice: first-principles study of tuning the electronic properties.. <i>RSC Advances</i> , 2020 , 10, 31894-31900	3.7	21
89	Modulating the electro-optical properties of doped C3N monolayers and graphene bilayers via mechanical strain and pressure. <i>New Journal of Chemistry</i> , 2020 , 44, 15785-15792	3.6	24
88	van der Waals heterostructures based on MSSe (M = Mo, W) and graphene-like GaN: enhanced optoelectronic and photocatalytic properties for water splitting. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 20704-20711	3.6	11
87	Type-I band alignment of BX-ZnO (X = As, P) van der Waals heterostructures as high-efficiency water splitting photocatalysts: a first-principles study.. <i>RSC Advances</i> , 2020 , 10, 44545-44550	3.7	7
86	Strain effects on the electronic and optical properties of Van der Waals heterostructure MoS2/WS2: A first-principles study. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2020 , 116, 113799	3	14
85	Electronic properties and enhanced photocatalytic performance of van der Waals heterostructures of ZnO and Janus transition metal dichalcogenides. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 10351-10359 ²⁶	3.6	26
84	Understanding the electronic properties, contact types and optical performances in graphene/InN heterostructure: Role of electric gating. <i>Diamond and Related Materials</i> , 2020 , 106, 107851	3.5	7

83	Theoretical prediction of electronic and optical properties of haft-hydrogenated InN monolayers. <i>Superlattices and Microstructures</i> , 2020 , 142, 106519	2.8	4
82	Phonon-assisted cyclotron resonance in Pöschl-Teller quantum well. <i>Journal of Applied Physics</i> , 2019 , 126, 124301	2.5	12
81	One- and two-photon-induced cyclotron-phonon resonance in modified-Pöschl-Teller quantum well. <i>Applied Physics A: Materials Science and Processing</i> , 2019 , 125, 1	2.6	11
80	Tri-layered van der Waals heterostructures based on graphene, gallium selenide and molybdenum selenide. <i>Journal of Applied Physics</i> , 2019 , 125, 225304	2.5	10
79	Cyclotron-phonon resonance line-width in monolayer silicene. <i>Superlattices and Microstructures</i> , 2019 , 131, 117-123	2.8	2
78	Two-photon induced magneto-optical absorption in finite semi-parabolic quantum wells. <i>Superlattices and Microstructures</i> , 2019 , 130, 446-453	2.8	
77	Tuning the electronic properties of GaS monolayer by strain engineering and electric field. <i>Chemical Physics</i> , 2019 , 524, 101-105	2.3	6
76	Magneto-optical effect in GaAs/GaAlAs semi-parabolic quantum well. <i>Thin Solid Films</i> , 2019 , 682, 10-17	2.2	26
75	Tailoring electronic properties and Schottky barrier in sandwich heterostructure based on graphene and tungsten diselenide. <i>Diamond and Related Materials</i> , 2019 , 94, 129-136	3.5	13
74	Strain engineering and electric field tunable electronic properties of Ti ₂ CO ₂ MXene monolayer. <i>Materials Research Express</i> , 2019 , 6, 065910	1.7	4
73	Van der Waals heterostructures of P, BSe, and SiC monolayers. <i>Journal of Applied Physics</i> , 2019 , 125, 094301	2.5	45
72	Electronic and optical properties of layered van der Waals heterostructure based on MS ₂ (M = Mo, W) monolayers. <i>Materials Research Express</i> , 2019 , 6, 065060	1.7	7
71	Electric field tunable electronic properties of P-ZnO and SiC-ZnO van der Waals heterostructures. <i>Computational Materials Science</i> , 2019 , 164, 166-170	3.2	20
70	Controlling electronic properties of PtS ₂ /InSe van der Waals heterostructure via external electric field and vertical strain. <i>Chemical Physics Letters</i> , 2019 , 724, 1-7	2.5	18
69	First-principles study of metal-semiconductor contact between MX ₂ (M = Nb, Pt; X = S, Se) monolayers. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2019 , 383, 125867	2.3	2
68	Optoelectronic and solar cell applications of Janus monolayers and their van der Waals heterostructures. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 18612-18621	3.6	77
67	Strain and electric field engineering of electronic structures and Schottky contact of layered graphene/Ca(OH) ₂ heterostructure. <i>Superlattices and Microstructures</i> , 2019 , 133, 106185	2.8	3
66	Electric field tuning of dynamical dielectric function in phosphorene. <i>Chemical Physics Letters</i> , 2019 , 731, 136606	2.5	1

65	Computational understanding of electronic properties of graphene/ $\{\mathrm{PtS}\}_2$ heterostructure under electric field. <i>Applied Physics A: Materials Science and Processing</i> , 2019 , 125, 1	2.6	3
64	Tunable electronic properties of InSe by biaxial strain: from bulk to single-layer. <i>Materials Research Express</i> , 2019 , 6, 115002	1.7	2
63	Rashba spin splitting and photocatalytic properties of Ge $\{\mathrm{MS}\}$ Se (M=Mo, W) van der Waals heterostructures. <i>Physical Review B</i> , 2019 , 100,	3.3	92
62	Strain-Tunable Electronic and Optical Properties of Monolayer Germanium Monosulfide: Ab-Initio Study. <i>Journal of Electronic Materials</i> , 2019 , 48, 2902-2909	1.9	8
61	Strain and electric field tunable electronic properties of type-II band alignment in van der Waals GaSe/MoSe ₂ heterostructure. <i>Chemical Physics</i> , 2019 , 521, 92-99	2.3	15
60	Band alignment and optical features in Janus-MoSeTe/X(OH) (X = Ca, Mg) van der Waals heterostructures. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 25849-25858	3.6	36
59	Tailoring the structural and electronic properties of an SnSe/MoS van der Waals heterostructure with an electric field and the insertion of a graphene sheet. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 22140-22148	3.6	32
58	Electronic and optical properties of Janus ZrSSe by density functional theory.. <i>RSC Advances</i> , 2019 , 9, 41058-41065	3.7	45
57	Modulation of electronic properties of monolayer InSe through strain and external electric field. <i>Chemical Physics</i> , 2019 , 516, 213-217	2.3	14
56	Nonlinear optical absorption and cyclotron impurity resonance in monolayer silicene. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2019 , 105, 168-173	3	
55	Vertical strain and electric field tunable electronic properties of type-II band alignment C ₂ N/InSe van der Waals heterostructure. <i>Chemical Physics Letters</i> , 2019 , 716, 155-161	2.5	30
54	Opening a band gap in graphene by C-C bond alternation: a tight binding approach. <i>Materials Research Express</i> , 2019 , 6, 045605	1.7	4
53	Modulation of electronic properties and Schottky barrier in the graphene/GaS heterostructure by electric gating. <i>Physica B: Condensed Matter</i> , 2019 , 555, 69-73	2.8	0
52	Electronic properties and optical behaviors of bulk and monolayer ZrS ₂ : A theoretical investigation. <i>Superlattices and Microstructures</i> , 2019 , 125, 205-213	2.8	15
51	Magneto-electronic perturbation effects on the electronic phase of phosphorene. <i>Materials Research Express</i> , 2019 , 6, 026102	1.7	1
50	Theoretical investigation of electronic structure and thermoelectric properties of MX ₂ (M=Zr, Hf; X=S, Se) van der Waals heterostructures. <i>Journal of Physics and Chemistry of Solids</i> , 2019 , 126, 304-309	3.9	19
49	Investigation of cyclotron-phonon resonance in monolayer molybdenum disulfide. <i>Journal of Physics and Chemistry of Solids</i> , 2019 , 125, 74-79	3.9	14
48	Adsorption and magnetism of bilayer graphene on the MnO polar surface with oxygen vacancies in the interface: First principles study. <i>Superlattices and Microstructures</i> , 2018 , 117, 72-81	2.8	3

47	Van der Waals graphene/g-GaSe heterostructure: Tuning the electronic properties and Schottky barrier by interlayer coupling, biaxial strain, and electric gating. <i>Journal of Alloys and Compounds</i> , 2018 , 750, 765-773	5.7	45
46	Magneto-optical properties of semi-parabolic plus semi-inverse squared quantum wells. <i>Physica B: Condensed Matter</i> , 2018 , 539, 117-122	2.8	26
45	First principles study of the electronic properties and band gap modulation of two-dimensional phosphorene monolayer: Effect of strain engineering. <i>Superlattices and Microstructures</i> , 2018 , 118, 289-297	2.8	15
44	Tuning the electronic properties and Schottky barrier height of the vertical graphene/MoS2 heterostructure by an electric gating. <i>Superlattices and Microstructures</i> , 2018 , 116, 79-87	2.8	32
43	Linear and nonlinear magneto-optical properties of monolayer MoS2. <i>Journal of Applied Physics</i> , 2018 , 123, 034301	2.5	16
42	First principles study of optical properties of molybdenum disulfide: From bulk to monolayer. <i>Superlattices and Microstructures</i> , 2018 , 115, 10-18	2.8	23
41	Electric-field tunable electronic properties and Schottky contact of graphene/phosphorene heterostructure. <i>Vacuum</i> , 2018 , 149, 231-237	3.7	31
40	Tuning the Electronic and Optical Properties of Two-Dimensional Graphene-like (hbox {C} _2hbox {N}) Nanosheet by Strain Engineering. <i>Journal of Electronic Materials</i> , 2018 , 47, 4594-4603	1.9	11
39	First-principles study of electronic properties of AB-stacked bilayer armchair graphene nanoribbons under out-plane strain. <i>Indian Journal of Physics</i> , 2018 , 92, 447-452	1.4	4
38	First principle study on the electronic properties and Schottky contact of graphene adsorbed on MoS2 monolayer under applied out-plane strain. <i>Surface Science</i> , 2018 , 668, 23-28	1.8	31
37	Tuning the Electronic Properties, Effective Mass and Carrier Mobility of MoS2 Monolayer by Strain Engineering: First-Principle Calculations. <i>Journal of Electronic Materials</i> , 2018 , 47, 730-736	1.9	42
36	Electronic properties of GaSe/MoS2 and GaS/MoSe2 heterojunctions from first principles calculations. <i>AIP Advances</i> , 2018 , 8, 075207	1.5	10
35	First principles calculations of the geometric structures and electronic properties of van der Waals heterostructure based on graphene, hexagonal boron nitride and molybdenum diselenide. <i>Diamond and Related Materials</i> , 2018 , 88, 151-157	3.5	12
34	LO-phonon-assisted cyclotron resonance in a special asymmetric hyperbolic-type quantum well. <i>Superlattices and Microstructures</i> , 2018 , 120, 738-746	2.8	20
33	Magneto-optical absorption in quantum dot via two-photon absorption process. <i>Optik</i> , 2018 , 173, 263-270	2.5	3
32	Ab-initio study of electronic and optical properties of biaxially deformed single-layer GeS. <i>Superlattices and Microstructures</i> , 2018 , 120, 501-507	2.8	17
31	Interlayer coupling and electric field tunable electronic properties and Schottky barrier in a graphene/bilayer-GaSe van der Waals heterostructure. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 17899-17908	3.6	76
30	Structural and electronic properties of a van der Waals heterostructure based on silicene and gallium selenide: effect of strain and electric field. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 27856-27864	2.6	54

29	First-principles study of W, N, and O adsorption on TiB ₂ (0001) surface with disordered vacancies. <i>Superlattices and Microstructures</i> , 2018 , 123, 414-426	2.8	6
28	Fundamental exciton transitions in SiO ₂ /Si/SiO ₂ cylindrical core/shell quantum dot. <i>Journal of Applied Physics</i> , 2018 , 124, 144303	2.5	8
27	Effective Photocatalytic Activity of Mixed Ni/Fe-Base Metal-Organic Framework under a Compact Fluorescent Daylight Lamp. <i>Catalysts</i> , 2018 , 8, 487	4	39
26	Layered graphene/GaS van der Waals heterostructure: Controlling the electronic properties and Schottky barrier by vertical strain. <i>Applied Physics Letters</i> , 2018 , 113, 171605	3.4	141
25	Electronic structure, optical and photocatalytic performance of SiC-MX (M = Mo, W and X = S, Se) van der Waals heterostructures. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 24168-24175	3.6	60
24	Phonon-assisted cyclotron resonance in special symmetric quantum wells. <i>Applied Physics A: Materials Science and Processing</i> , 2018 , 124, 1	2.6	9
23	First principles study on the electronic properties and Schottky barrier of Graphene/InSe heterostructure. <i>Superlattices and Microstructures</i> , 2018 , 122, 570-576	2.8	21
22	Effect of strains on electronic and optical properties of monolayer SnS: Ab-initio study. <i>Physica B: Condensed Matter</i> , 2018 , 545, 255-261	2.8	16
21	Linear and nonlinear magneto-optical properties of monolayer phosphorene. <i>Journal of Applied Physics</i> , 2017 , 121, 045107	2.5	33
20	Phase Transition in Armchair Graphene Nanoribbon Due to Peierls Distortion. <i>Journal of Electronic Materials</i> , 2017 , 46, 3815-3819	1.9	2
19	First principles study of structural, electronic and magnetic properties of graphene adsorbed on the O-terminated MnO(111) surface. <i>Diamond and Related Materials</i> , 2017 , 74, 31-40	3.5	10
18	Simple single-emitting layer hybrid white organic light emitting with high color stability. <i>Journal of Applied Physics</i> , 2017 , 122, 134503	2.5	3
17	Out-of-plane strain and electric field tunable electronic properties and Schottky contact of graphene/antimonene heterostructure. <i>Superlattices and Microstructures</i> , 2017 , 112, 554-560	2.8	22
16	First-principles study of the structural and electronic properties of graphene/MoS ₂ interfaces. <i>Journal of Applied Physics</i> , 2017 , 122, 104301	2.5	43
15	Magneto-optical transport properties of monolayer MoS ₂ on polar substrates. <i>Physical Review B</i> , 2017 , 96,	3.3	61
14	Band gap and electronic properties of molybdenum disulphide under strain engineering: density functional theory calculations. <i>Molecular Simulation</i> , 2017 , 43, 86-91	2	10
13	First-principles study of the structural and electronic properties of graphene adsorbed on MnO(111) surfaces. <i>Computational and Theoretical Chemistry</i> , 2016 , 1098, 22-30	2	9
12	Effect of biaxial strain and external electric field on electronic properties of MoS ₂ monolayer: A first-principle study. <i>Chemical Physics</i> , 2016 , 468, 9-14	2.3	33

11	Transport properties of armchair graphene nanoribbons under uniaxial strain: A first principles study. <i>Solid State Communications</i> , 2016 , 237-238, 10-13	1.6	2
10	Band Gap Modulation of Bilayer MoS ₂ Under Strain Engineering and Electric Field: A Density Functional Theory. <i>Journal of Electronic Materials</i> , 2016 , 45, 4038-4043	1.9	13
9	Modulation of the band structure in bilayer zigzag graphene nanoribbons on hexagonal boron nitride using the force and electric fields. <i>Materials Chemistry and Physics</i> , 2015 , 154, 78-83	4.4	12
8	Effect of electric field on the electronic and magnetic properties of a graphene nanoribbon/aluminium nitride bilayer system. <i>RSC Advances</i> , 2015 , 5, 49308-49316	3.7	25
7	Tuning the electronic properties of armchair graphene nanoribbons by strain engineering. <i>Physica Scripta</i> , 2015 , 90, 015802	2.6	10
6	Substrate-induced band structure and electronic properties in graphene/Al ₂ O ₃ (0001) interface. <i>Surface Science</i> , 2015 , 632, 111-117	1.8	4
5	Dispersion-Corrected Density Functional Theory Investigations of Structural and Electronic Properties of Bulk MoS ₂ : Effect of Uniaxial Strain. <i>Nanoscale Research Letters</i> , 2015 , 10, 433	5	18
4	Electric field and substrate-induced modulation of spin-polarized transport in graphene nanoribbons on AlN semiconductors. <i>Journal of Applied Physics</i> , 2015 , 117, 174309	2.5	28
3	Template-Based Growth of Various Oxide Nanorods by Sol-Gel Electrophoresis. <i>Advanced Functional Materials</i> , 2002 , 12, 59	15.6	199
2	Organic-inorganic sol-gel coating for corrosion protection of stainless steel. <i>Journal of Materials Science Letters</i> , 2002 , 21, 251-255		39
1	Electrophoretic Growth of Lead Zirconate Titanate Nanorods. <i>Advanced Materials</i> , 2001 , 13, 1269	24	142