## Congxin Xia

## List of Publications by Year in descending order

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Version: 2024-02-01

		101496	118793
196	5,254	36	62
papers	citations	h-index	g-index
196	196	196	5319
all docs	docs citations	times ranked	citing authors
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#	Article	IF	CITATIONS
1	Highly Polarization-Sensitive, Broadband, Self-Powered Photodetector Based on Graphene/PdSe <sub>2</sub> /Germanium Heterojunction. ACS Nano, 2019, 13, 9907-9917.	7.3	420
2	Universality of electronic characteristics and photocatalyst applications in the two-dimensional Janus transition metal dichalcogenides. Physical Review B, 2018, 98, .	1.1	229
3	Tunable Room-Temperature Ferromagnetism in Two-Dimensional Cr <sub>2</sub> Te <sub>3</sub> . Nano Letters, 2020, 20, 3130-3139.	4.5	175
4	A type-II GeSe/SnS heterobilayer with a suitable direct gap, superior optical absorption and broad spectrum for photovoltaic applications. Journal of Materials Chemistry A, 2017, 5, 13400-13410.	5 <b>.</b> 2	138
5	Enhanced rectification, transport property and photocurrent generation of multilayer ReSe2/MoS2 p–n heterojunctions. Nano Research, 2016, 9, 507-516.	5 <b>.</b> 8	132
6	Interlayer coupling effects on Schottky barrier in the arsenene-graphene van der Waals heterostructures. Applied Physics Letters, 2015, 107, .	1.5	128
7	Various Structures of 2D Transitionâ€Metal Dichalcogenides and Their Applications. Small Methods, 2018, 2, 1800094.	4.6	107
8	Effects of strain and electric field on electronic structures and Schottky barrier in graphene and SnS hybrid heterostructures. Carbon, 2016, 109, 737-746.	5.4	106
9	Tuning the band gap of hematite α-Fe2O3 by sulfur doping. Physics Letters, Section A: General, Atomic and Solid State Physics, 2013, 377, 1943-1947.	0.9	98
10	The characteristics of n- and p-type dopants in SnS2monolayer nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 19674.	1.3	98
11	Enhancement of hole mobility in InSe monolayer via an InSe and black phosphorus heterostructure. Nanoscale, 2017, 9, 14682-14689.	2.8	92
12	Elastic, electronic and optical properties of the two-dimensional PtX2 (X = S, Se, and Te) monolayer. Applied Surface Science, 2018, 435, 476-482.	3.1	89
13	The feasibility of tunable p-type Mg doping in a GaN monolayer nanosheet. Acta Materialia, 2013, 61, 7720-7725.	3.8	81
14	Composition-tunable 2D SnSe <sub>2(1â°x)</sub> S <sub>2x</sub> alloys towards efficient bandgap engineering and high performance (opto)electronics. Journal of Materials Chemistry C, 2017, 5, 84-90.	2.7	81
15	Recent advances in optoelectronic properties and applications of two-dimensional metal chalcogenides. Journal of Semiconductors, 2016, 37, 051001.	2.0	75
16	Direct Wide Bandgap 2D GeSe <sub>2</sub> Monolayer toward Anisotropic UV Photodetection.  Advanced Optical Materials, 2019, 7, 1900622.	3 <b>.</b> 6	70
17	xmins:mmi="http://www.w3.org/1998/Math/MathMiL" display="inline" overflow="scroll"> <mml:mi>Sn</mml:mi> <mml:msub><mml:mi>X</mml:mi><mml:mn>2</mml:mn><td>b&gt;<td>math&gt;</td></td></mml:msub>	b> <td>math&gt;</td>	math>

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Let-tellurene/MoS (sub ) 2 (slub ) heterostructure. Journal of Materials Chemistry C, 2018, 6, 10256-10262. 2.7 50  Tuning the electronic structures and magnetism of two-dimensional porous C (sub ) 2 (slub ) N via transition metal embedding. Physical Chemistry Chemical Physics, 2016, 18, 22678-22686. 1.3 54  Electronic and magnetic properties of n-type and p-doped MoS (sub ) 2 (slub ) monolayers. RSC Advances, 2016, 6, 6772-16778. 1.7 54  Two-dimensional transition-metal dichalcogenides-based ferromagnetic van der Waals heterostructures. Nanoscale, 2017, 9, 17585-17592. 2.8 50  A theoretical simulation of small-molecules sensing on an S-vacancy SnS (sub ) 2 (slub ) monolayer. 2.8 50  hard p type dopants in the InSe monolayer via substitutional doping. Journal of Materials Science, 2017, 9, 19470-10480. 1.7 48  Tunable electronic structures and magnetism in arsenene nanosheets via transition metal doping. Journal of Materials Science, 2016, 51, 504-9513. 1.7 46  Electronic and magnetic properties of X-doped (XA-ATI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579. 2.7 49  Electronic and magnetic properties of X-doped (XA-ATI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579. 2.7 49  Electronic and magnetic properties of X-doped (XA-ATI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579. 2.7 49  Electronic and magnetic properties of X-doped (XA-ATI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579. 2.7 49  Electronic and magnetic properties of Shades and GaSe nanosheets. Physical Chemistry Chemical Physics, 2016, 18, 654-6540. 3.1 49  Tunable donor and acceptor impurity states in a WSe (sub ) 2 (sub ) monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800. 1.7 43  Electronic and magnetic properties of ShSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 81	19		2.7	59
transition metal embedding. Physical Chemistry Chemical Physics, 2016, 18, 22678-22686.  Electronic and magnetic properties of n-type and p-doped MoS sub> 2 (sub> monolayers. RSC Advances, 2016, 6, 16772-16778.  Two-dimensional transition-metal dichałcogenides-based ferromagnetic van der Waals heterostructures. Nanocade, 2017, 9, 17365-17592.  A theoretical simulation of small-molecules sensing on an S-vaccancy SnS (sub> 2 (sub> monolayer. Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 52  Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 52  1.7 48  2.8 13  2.8 2017, 52, 7207-7214.  1.8 19 10470-10480.  1.9 10470-10480.  1.10 48  2.11 48  2.12 104 105 105 105 105 105 105 105 105 105 105	20	Interlayer coupling and external electric field tunable electronic properties of a 2D type-I α-tellurene/MoS <sub>2</sub> heterostructure. Journal of Materials Chemistry C, 2018, 6, 10256-10262.	2.7	56
Two-dimensional transition-metal dichalcogenides-based ferromagnetic van der Waals  Two-dimensional transition-metal dichalcogenides-based ferromagnetic van der Waals  A theoretical simulation of small-molecules sensing on an S-vacancy SnS (sub-) 2 (sub-) monolayer.  Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 52  A theoretical simulation of small-molecules sensing on an S-vacancy SnS (sub-) 2 (sub-) monolayer.  Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 52  1.4 Experimental Science, 2016, 51, 10470-10480.  1.5 In-and p-type dopants in the InSe monolayer via substitutional doping, Journal of Materials Science, 2016, 51, 9504-9513.  1.7 48  1.8 In-and p-type dopants in the InSe monolayer via substitutional doping, Journal of Materials Science, 2016, 51, 9504-9513.  1.7 46  1.8 In-and p-type dopants in the InSe monolayer via substitutional doping, Journal of Materials Science, 2016, 51, 9504-9513.  1.8 In-and p-type dopants in the InSe monolayer via substitutional doping, Journal of Materials Science, 2016, 51, 9504-9513.  2.8 In-and p-type dopants in the InSe monolayer via substitutional doping, Journal of Materials Science, 2016, 51, 9504-9513.  2.8 In-and p-type dopants in the InSe monolayer dopants in arsenene nanosheets via transition metal doping.  1.7 46  1.8 In-and p-type dopants in the InSe monolayer dopants in arsenene nanosheets via transition metal doping.  1.8 In-and p-type dopants in the InSe monolayer dopants in arsenene nanosheets via transition of the doping of the MoS (sub-) 2 (sub-) 3 monolayer by adsorption of the MoS (sub-) 2 (sub-) 3 monolayer by adsorption of the MoS (sub-) 2 (sub-) 3 monolayer by adsorption of the MoS (sub-) 2 (sub-) 4 monolayer by 3 disorption of 4 monolayer by 3 disorption of 4 monolayer by 3 disorption of 4 monolayer by 4 monolayer by 3 disorption of 4 monolayer by	21	Tuning the electronic structures and magnetism of two-dimensional porous C <sub>2</sub> N via transition metal embedding. Physical Chemistry Chemical Physics, 2016, 18, 22678-22686.	1.3	54
heterostructures. Nanoscale, 2017, 9, 17585-17592.  A theoretical simulation of small-molecules sensing on an S-vacancy SnS sub>2 ⟨ /sub> monolayer. Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 52  n- and p-type dopants in the InSe monolayer via substitutional doping. Journal of Materials Science, 2017, 52, 7207-7214.  1.7 48  Tunable electronic structures and magnetism in arsenene nanosheets via transition metal doping. Journal of Materials Science, 2016, 51, 9504-9513.  1.7 46  Electronic and magnetic properties of X-doped (XÂ-ÂTI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579.  Characteristics of p-type Mg-doped GaS and GaSe nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 18799.  Band engineering of the MoS (sub>2 ⟨ /sub> /stanene heterostructure: strain and electrostatic gating. Nanotechnology, 2017, 28, 195702.  Tunable donor and acceptor impurity states in a WSe (sub>2 ⟨ /sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  30 Tunable donor and acceptor impurity states in a WSe (sub>2 ⟨ /sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  31 Spinac* orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  32 Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  33 Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  41 Typeâ¢1 Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.	22		1.7	54
Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.  1.3 32  2.5 n- and p-type dopants in the InSe monolayer via substitutional doping. Journal of Materials Science, 2017, 52, 7207-7214.  2.6 Tunable electronic structures and magnetism in arsenene nanosheets via transition metal doping.  2.7 Journal of Materials Science, 2016, 51, 9504-9513.  2.8 Lectronic and magnetic properties of X-doped (XĀ=ĀTI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579.  2.8 Characteristics of p-type Mg-doped GaS and GaSe nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 18799.  2.8 Band engineering of the MoS <sub>2</sub> /sub>/stanene heterostructure: strain and electrostatic gating.  2.9 Band engineering of the MoS <sub>2</sub> /stanene heterostructure: strain and electrostatic gating.  2.9 Nanotechnology, 2017, 28, 195702.  3.0 Tunable donor and acceptor impurity states in a WSe <sub>2</sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  3.1 Spinā€"orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  3.1 Spinā€"orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 8538-8164.  3.2 Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  3.1 Type†Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  3.2 Type†Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.	23	Two-dimensional transition-metal dichalcogenides-based ferromagnetic van der Waals heterostructures. Nanoscale, 2017, 9, 17585-17592.	2.8	53
Tunable electronic structures and magnetism in arsenene nanosheets via transition metal doping.  1.7 46  Tunable electronic structures and magnetism in arsenene nanosheets via transition metal doping.  1.7 46  Electronic and magnetic properties of X-doped (XÅ=ĀTI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579.  Characteristics of p-type Mg-doped GaS and GaSe nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 18799.  Band engineering of the MoS <sub>2</sub> /stanene heterostructure: strain and electrostatic gating. Nanotechnology, 2017, 28, 195702.  Tunable donor and acceptor impurity states in a WSe <sub>2</sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  1.7 43  Spinā€"orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 nonolayer. Solid State Communications, 2015, 220, 31-35.  Typeå& Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  Crosså& Gubstitution Promoted Ultrawide Bandgap up to 4.5 & V in a 2D Semiconductor: Gallium	24	A theoretical simulation of small-molecules sensing on an S-vacancy SnS <sub>2</sub> monolayer. Physical Chemistry Chemical Physics, 2017, 19, 10470-10480.	1.3	52
Electronic and magnetic properties of X-doped (XÂ=ÂTI, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579.  28 Characteristics of p-type Mg-doped GaS and GaSe nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 18799.  29 Band engineering of the MoS <sub>2</sub> /stanene heterostructure: strain and electrostatic gating.  30 Tunable donor and acceptor impurity states in a WSe <sub>2</sub> /sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  31 SpináC*orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  32 Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  33 Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  34 Typeâ& Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  34 Crosså&bubstitution Promoted Ultrawide Bandgap up to 4.5&V in a 2D Semiconductor: Gallium	25	n- and p-type dopants in the InSe monolayer via substitutional doping. Journal of Materials Science, 2017, 52, 7207-7214.	1.7	48
of Alloys and Compounds, 2016, 654, 574-579.  Characteristics of p-type Mg-doped GaS and GaSe nanosheets. Physical Chemistry Chemical Physics, 2014, 16, 18799.  1.3 44  Band engineering of the MoS <sub>2</sub> /stanene heterostructure: strain and electrostatic gating. Nanotechnology, 2017, 28, 195702.  1.3 44  Tunable donor and acceptor impurity states in a WSe <sub>2</sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  1.7 43  Spinå6"orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  Typeå6t Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  Crosså66ubstitution Promoted Ultrawide Bandgap up to 4.5ÅeV in a 2D Semiconductor: Callium	26		1.7	46
29 Band engineering of the MoS⟨sub⟩2⟨/sub⟩/stanene heterostructure: strain and electrostatic gating.  1.3 44  29 Band engineering of the MoS⟨sub⟩2⟨/sub⟩/stanene heterostructure: strain and electrostatic gating.  30 Tunable donor and acceptor impurity states in a WSe⟨sub⟩2⟨/sub⟩ monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  1.7 43  31 Spinâ€"orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  32 Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  33 Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  34 Typeâ€l Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  34 Crossâ€Bubstitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium	27	Electronic and magnetic properties of X-doped (XÂ=ÂTi, Zr, Hf) tungsten disulphide monolayer. Journal of Alloys and Compounds, 2016, 654, 574-579.	2.8	45
Nanotechnology, 2017, 28, 195702.  Tunable donor and acceptor impurity states in a WSe <sub>2</sub> monolayer by adsorption of common gas molecules. RSC Advances, 2016, 6, 82793-82800.  1.7 43  Spinå& orbit coupling effects on electronic structures in stanene nanoribbons. Physical Chemistry Chemical Physics, 2016, 18, 6534-6540.  Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  Typeå& Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.	28		1.3	44
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Chemical Physics, 2016, 18, 6534-6540.  Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.  Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.  Typeâ€I Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  Crossâ€Substitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium	30		1.7	43
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monolayer. Solid State Communications, 2015, 220, 31-35.  Type†Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.  5.2 41  Crossâ€Substitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium	32	Electronic and magnetic properties of SnSe monolayers doped by Ga, In, As, and Sb: a first-principles study. Physical Chemistry Chemical Physics, 2016, 18, 8158-8164.	1.3	42
Field Effects. Small, 2018, 14, e1800365.  5.2 41  Crossâ€Substitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium	33	Effect of structural defects on electronic and magnetic properties of pristine and Mn-doped MoS2 monolayer. Solid State Communications, 2015, 220, 31-35.	0.9	41
Crossâ€Substitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium  Thiophosphate. Advanced Materials, 2021, 33, e2008761.  11.1 41	34	Type†Transition Metal Dichalcogenides Lateral Homojunctions: Layer Thickness and External Electric Field Effects. Small, 2018, 14, e1800365.	5.2	41
	35	Crossâ€Substitution Promoted Ultrawide Bandgap up to 4.5ÂeV in a 2D Semiconductor: Gallium Thiophosphate. Advanced Materials, 2021, 33, e2008761.	11.1	41

Electrostatic gating dependent multiple-band alignments in a high-temperature ferromagnetic Mg(OH)

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#	Article	IF	CITATIONS
37	Tuning anisotropic electronic transport properties of phosphorene via substitutional doping. Physical Chemistry Chemical Physics, 2016, 18, 25869-25878.	1.3	38
38	Modulation of electronic transport properties in armchair phosphorene nanoribbons by doping and edge passivation. Scientific Reports, 2017, 7, 12799.	1.6	38
39	Reversible Half Wave Rectifier Based on 2D InSe/GeSe Heterostructure with Nearâ€Broken Band Alignment. Advanced Science, 2021, 8, 1903252.	5.6	38
40	Photovoltaic Field-Effect Photodiodes Based on Double van der Waals Heterojunctions. ACS Nano, 2021, 15, 14295-14304.	7.3	37
41	The rectifying and negative differential resistance effects in graphene/h-BN nanoribbon heterojunctions. Physical Chemistry Chemical Physics, 2016, 18, 27976-27980.	1.3	36
42	Electronic and magnetic properties of 1T-HfS2 by doping transition-metal atoms. Applied Surface Science, 2016, 383, 151-158.	3.1	36
43	First-principles studies on substitutional doping by group IV and VI atoms in the two-dimensional arsenene. Applied Surface Science, 2016, 378, 350-356.	3.1	36
44	Effects of applied strain and electric field on small-molecule sensing by stanene monolayers. Journal of Materials Science, 2017, 52, 5083-5096.	1.7	36
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