

Ángel Rubio Secades

List of Publications by Year in descending order

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Version: 2024-02-01

709
papers

55,619
citations

1070

116
h-index

2072

211
g-index

724
all docs

724
docs citations

724
times ranked

38941
citing authors

#	ARTICLE	IF	CITATIONS
1	Free electron gas in cavity quantum electrodynamics. <i>Physical Review Research</i> , 2022, 4, .	1.3	33
2	Detecting multiple chiral centers in chiral molecules with high harmonic generation. <i>Optics Express</i> , 2022, 30, 3729.	1.7	16
3	Moiré engineering of spin-orbit coupling in twisted platinum diselenide. <i>Electronic Structure</i> , 2022, 4, 014004.	1.0	8
4	Spin-orbit induced equilibrium spin currents in materials. <i>Physical Review B</i> , 2022, 105, .	1.1	3
5	First-principles modelling for time-resolved ARPES under different pump-probe conditions. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 2022, 254, 147152.	0.8	9
6	Excited-state band structure mapping. <i>Physical Review B</i> , 2022, 105, .	1.1	8
7	Nanometer-Scale Lateral p-n Junctions in Graphene/RuCl ₃ Heterostructures. <i>Nano Letters</i> , 2022, 22, 1946-1953.	4.5	25
8	Cavity-Modulated Proton Transfer Reactions. <i>Journal of the American Chemical Society</i> , 2022, 144, 4995-5002.	6.6	32
9	TMDs as a platform for spin liquid physics: A strong coupling study of twisted bilayer WSe ₂ . <i>APL Materials</i> , 2022, 10, .	2.2	19
10	A new Hall for quantum protection. <i>Science</i> , 2022, 375, 976-977.	6.0	4
11	Microscopic theory of light-induced ultrafast skyrmion excitation in transition metal films. <i>Npj Computational Materials</i> , 2022, 8, .	3.5	8
12	Comment on "Origin of symmetry-forbidden high-order harmonic generation in the time-dependent Kohn-Sham formulation". <i>Physical Review A</i> , 2022, 105, .	1.0	1
13	Moiré nematic phase in twisted double bilayer graphene. <i>Nature Physics</i> , 2022, 18, 196-202.	6.5	51
14	Polaritonic Hofstadter butterfly and cavity control of the quantized Hall conductance. <i>Physical Review B</i> , 2022, 105, .	1.1	20
15	A perspective on <i>ab initio</i> modeling of polaritonic chemistry: The role of non-equilibrium effects and quantum collectivity. <i>Journal of Chemical Physics</i> , 2022, 156, .	1.2	39
16	Strongly correlated electron-photon systems. <i>Nature</i> , 2022, 606, 41-48.	13.7	66
17	Few-Femtosecond Dynamics of Free-Free Opacity in Optically Heated Metals. <i>Physical Review X</i> , 2022, 12, .	2.8	6
18	Unconventional excitonic states with phonon sidebands in layered silicon diphosphide. <i>Nature Materials</i> , 2022, 21, 773-778.	13.3	20

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19	Probing phonon dynamics with multidimensional high harmonic carrier-envelope-phase spectroscopy. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119, .	3.3	13
20	Frequency-Dependent Sternheimer Linear-Response Formalism for Strongly Coupled Light-Matter Systems. Journal of Chemical Theory and Computation, 2022, 18, 4354-4365.	2.3	9
21	Wavefunction embedding for molecular polaritons. Journal of Chemical Physics, 2022, 157, .	1.2	18
22	Effect of spin-orbit coupling on the high harmonics from the topological Dirac semimetal Na ₃ Bi. Npj Computational Materials, 2022, 8, .	3.5	13
23	Engineering quantum materials with chiral optical cavities. Nature Materials, 2021, 20, 438-442.	13.3	120
24	Quantitative sampling of atomic-scale electromagnetic waveforms. Nature Photonics, 2021, 15, 143-147.	15.6	44
25	Toward Confined Carbyne with Tailored Properties. Nano Letters, 2021, 21, 1096-1101.	4.5	27
26	Moiré metrology of energy landscapes in van der Waals heterostructures. Nature Communications, 2021, 12, 242.	5.8	60
27	High-order harmonic generation in graphene: Nonlinear coupling of intraband and interband transitions. Physical Review B, 2021, 103, .	1.1	31
28	Vibrational coherent control of localized d electronic excitation. Nature Physics, 2021, 17, 368-373.	6.5	10
29	Quantitative Waveform Sampling on Atomic Scales. , 2021, , .		0
30	High Harmonics and Isolated Attosecond Pulses from MgO . Physical Review Applied, 2021, 15, .	1.5	26
31	Unravelling the intertwined atomic and bulk nature of localised excitons by attosecond spectroscopy. Nature Communications, 2021, 12, 1021.	5.8	32
32	Self-Consistent Potential Correction for Charged Periodic Systems. Physical Review Letters, 2021, 126, 076401.	2.9	44
33	Higher-Order Band Topology in Twisted Moiré Superlattice. Physical Review Letters, 2021, 126, 066401.	2.9	56
34	Hydrated Alkali Atoms on Copper(111): A Density Functional Theory Study. Journal of Physical Chemistry C, 2021, 125, 3868-3879.	1.5	3
35	Moiré heterostructures as a condensed-matter quantum simulator. Nature Physics, 2021, 17, 155-163.	6.5	317
36	Programmable hyperbolic polaritons in van der Waals semiconductors. Science, 2021, 371, 617-620.	6.0	58

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37	Enhanced tunable second harmonic generation from twistable interfaces and vertical superlattices in boron nitride homostructures. <i>Science Advances</i> , 2021, 7, .	4.7	73
38	Intermolecular interactions in optical cavities: An <i>ab initio</i> QED study. <i>Journal of Chemical Physics</i> , 2021, 154, 094113.	1.2	81
39	Single and double charge transfer in the Ne^+H_2 collision within time-dependent density-functional theory. <i>Physical Review A</i> , 2021, 103, .		
40	Simulating Vibronic Spectra without Born-Oppenheimer Surfaces. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 3074-3081.	2.1	8
41	Ultrafast dynamical Lifshitz transition. <i>Science Advances</i> , 2021, 7, .	4.7	38
42	Real-time observation of a correlation-driven sub-fs charge migration in ionised adenine. <i>Communications Chemistry</i> , 2021, 4, .	2.0	38
43	Identification of the Mott-insulating Charge Density Wave State in TaTe_2 . <i>Physical Review Letters</i> , 2021, 126, 196406.	2.9	27
44	Nonlinear electric conductivity and THz-induced charge transport in graphene. <i>New Journal of Physics</i> , 2021, 23, 063047.	1.2	5
45	Light-Induced Charge Transfer from Transition-Metal-Doped Aluminum Clusters to Carbon Dioxide. <i>Journal of Physical Chemistry A</i> , 2021, 125, 5878-5885.	1.1	2
46	Phonoritons as Hybridized Exciton-Photon-Phonon Excitations in a Monolayer hBN Optical Cavity. <i>Physical Review Letters</i> , 2021, 126, 227401.	2.9	18
47	Direct measurement of key exciton properties: Energy, dynamics, and spatial distribution of the wave function. <i>Natural Sciences</i> , 2021, 1, e10010.	1.0	52
48	Survival of Floquet-Bloch States in the Presence of Scattering. <i>Nano Letters</i> , 2021, 21, 5028-5035.	4.5	41
49	Strong chiral dichroism and enantiopurification in above-threshold ionization with locally chiral light. <i>Physical Review Research</i> , 2021, 3, .	1.3	9
50	The ferroelectric photo ground state of SrTiO_3 : Cavity materials engineering. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	3.3	49
51	Down-conversion processes in <i>ab initio</i> nonrelativistic quantum electrodynamics. <i>Physical Review Research</i> , 2021, 3, .	1.3	8
52	Out-of-Plane Transport of 1T-TaS_2 /Graphene-Based van der Waals Heterostructures. <i>ACS Nano</i> , 2021, 15, 11898-11907.	7.3	20
53	The 2021 ultrafast spectroscopic probes of condensed matter roadmap. <i>Journal of Physics Condensed Matter</i> , 2021, 33, 353001.	0.7	55
54	Coherent coupling between vortex bound states and magnetic impurities in 2D layered superconductors. <i>Nature Communications</i> , 2021, 12, 4668.	5.8	5

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55	Approximations based on density-matrix embedding theory for density-functional theories. <i>Electronic Structure</i> , 2021, 3, 035001.	1.0	1
56	Quantum paraelectric phase of SrTiO_3 from first principles. <i>Physical Review B</i> , 2021, 104, .	2.9	25
57	Light-Driven Extremely Nonlinear Bulk Photogalvanic Currents. <i>Physical Review Letters</i> , 2021, 127, 126601.	2.9	13
58	Nematicity Arising from a Chiral Superconducting Ground State in Magic-Angle Twisted Bilayer Graphene under In-Plane Magnetic Fields. <i>Physical Review Letters</i> , 2021, 127, 127001.	1.1	10
59	All-optical generation of antiferromagnetic magnon currents via the magnon circular photogalvanic effect. <i>Physical Review B</i> , 2021, 104, .	4.5	10
60	Engineering Three-Dimensional Moiré Flat Bands. <i>Nano Letters</i> , 2021, 21, 7519-7526.	3.3	59
61	Moiré correlations in ABCA graphene. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	2.1	65
62	Polaritonic Chemistry: Collective Strong Coupling Implies Strong Local Modification of Chemical Properties. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 508-516.	5.8	57
63	Realization of nearly dispersionless bands with strong orbital anisotropy from destructive interference in twisted bilayer MoS ₂ . <i>Nature Communications</i> , 2021, 12, 5644.	3.3	42
64	Making ab initio QED functional(s): Nonperturbative and photon-free effective frameworks for strong light-matter coupling. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	2.3	1
65	Conditional Wave Function Theory: A Unified Treatment of Molecular Structure and Nonadiabatic Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2021, , .	3.5	19
66	Common microscopic origin of the phase transitions in Ta ₂ NiS ₅ and the excitonic insulator candidate Ta ₂ NiSe ₅ . <i>Npj Computational Materials</i> , 2021, 7, .	1.1	7
67	Photoionization and transient Wannier-Stark ladder in silicon: First-principles simulations versus Keldysh theory. <i>Physical Review B</i> , 2021, 104, .	2.3	1
68	Nonadiabatic Electron Dynamics in Tunneling Junctions: Lattice Exchange-Correlation Potential. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 295-301.	1.1	13
69	Entangled photon assisted multidimensional nonlinear optics of exciton-polaritons. <i>Journal of Applied Physics</i> , 2020, 128, 113102.	1.1	17
70	Role of intraband dynamics in the generation of circularly polarized high harmonics from solids. <i>Physical Review B</i> , 2020, 102, .	2.8	29
71	How Circular Dichroism in Time- and Angle-Resolved Photoemission Can Be Used to Spectroscopically Detect Transient Topological States in Graphene. <i>Physical Review X</i> , 2020, 10, .	1.1	13
72	Parameter-free hybridlike functional based on an extended Hubbard model: DFT+U. <i>Physical Review B</i> , 2020, 102, .		

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73	Femtosecond exciton dynamics in WSe ₂ optical waveguides. Nature Communications, 2020, 11, 3567.	5.8	31
74	Dynamical amplification of electric polarization through nonlinear phononics in 2D SnTe. Npj Computational Materials, 2020, 6, .	3.5	12
75	Charge-Transfer Plasmon Polaritons at Graphene/RuCl ₃ Interfaces. Nano Letters, 2020, 20, 8438-8445.	4.5	53
76	Ultrafast Real-Time Dynamics of CO Oxidation over an Oxide Photocatalyst. ACS Catalysis, 2020, 10, 13650-13658.	5.5	11
77	Light-Matter Hybrid-Orbital-Based First-Principles Methods: The Influence of Polariton Statistics. Journal of Chemical Theory and Computation, 2020, 16, 5601-5620.	2.3	19
78	Exact exchange-correlation potential of effectively interacting Kohn-Sham systems. Physical Review A, 2020, 101, .	1.0	1
79	Light-Induced Renormalization of the Dirac Quasiparticles in the Nodal-Line Semimetal ZrSiSe. Physical Review Letters, 2020, 125, 076401.	2.9	26
80	Effect of many modes on self-polarization and photochemical suppression in cavities. Journal of Chemical Physics, 2020, 153, 104103.	1.2	44
81	Direct Measurement of Electron-Phonon Coupling with Time-Resolved ARPES. Physical Review Letters, 2020, 125, 136401.	2.9	27
82	Virial Relations for Electrons Coupled to Quantum Field Modes. Journal of Chemical Theory and Computation, 2020, 16, 6236-6243.	2.3	6
83	Chemistry in Quantum Cavities: Exact Results, the Impact of Thermal Velocities, and Modified Dissociation. Journal of Physical Chemistry Letters, 2020, 11, 7525-7530.	2.1	26
84	Giant Exciton Mott Density in Anatase TiO ₂ . Physical Review Letters, 2020, 125, 116403.	2.9	73
85	Topological phase transitions induced by disorder in magnetically doped Bi ₂ Te thin films. Physical Review B, 2020, 102, .	1.1	6
86	Coupled Cluster Theory for Molecular Polaritons: Changing Ground and Excited States. Physical Review X, 2020, 10, .	2.8	102
87	Dual boson diagrammatic Monte Carlo approach applied to the extended Hubbard model. Physical Review B, 2020, 102, .	1.1	17
88	Ultrafast transient absorption spectroscopy of the charge-transfer insulator NiO: Beyond the dynamical Franz-Keldysh effect. Physical Review B, 2020, 102, .	1.1	12
89	Nature of Symmetry Breaking at the Excitonic Insulator Transition: Ta ₂ Te ₅ . Physical Review Letters, 2020, 124, 197601.	2.9	73
90	Correlated electronic phases in twisted bilayer transition metal dichalcogenides. Nature Materials, 2020, 19, 861-866.	13.3	544

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91	Room Temperature Terahertz Electroabsorption Modulation by Excitons in Monolayer Transition Metal Dichalcogenides. <i>Nano Letters</i> , 2020, 20, 5214-5220.	4.5	14
92	Polycyclic aromatic chains on metals and insulating layers by repetitive [3+2] cycloadditions. <i>Nature Communications</i> , 2020, 11, 1490.	5.8	23
93	Electron-phonon-driven three-dimensional metallicity in an insulating cuprate. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 6409-6416.	3.3	18
94	Octopus, a computational framework for exploring light-driven phenomena and quantum dynamics in extended and finite systems. <i>Journal of Chemical Physics</i> , 2020, 152, 124119.	1.2	210
95	Local Berry curvature signatures in dichroic angle-resolved photoelectron spectroscopy from two-dimensional materials. <i>Science Advances</i> , 2020, 6, eaay2730.	4.7	57
96	Setting the photoelectron clock through molecular alignment. <i>Nature Communications</i> , 2020, 11, 2546.	5.8	26
97	Attosecond timing of the dynamical Franz-Keldysh effect. <i>JPhys Photonics</i> , 2020, 2, 025001.	2.2	13
98	Universal slow plasmons and giant field enhancement in atomically thin quasi-two-dimensional metals. <i>Nature Communications</i> , 2020, 11, 1013.	5.8	53
99	One-dimensional flat bands in twisted bilayer germanium selenide. <i>Nature Communications</i> , 2020, 11, 1124.	5.8	80
100	Relevance of the Quadratic Diamagnetic and Self-Polarization Terms in Cavity Quantum Electrodynamics. <i>ACS Photonics</i> , 2020, 7, 975-990.	3.2	105
101	Covalent C-N Bond Formation through a Surface Catalyzed Thermal Cyclodehydrogenation. <i>Journal of the American Chemical Society</i> , 2020, 142, 3696-3700.	6.6	27
102	Time-Dependent Magnons from First Principles. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 1007-1017.	2.3	21
103	High-harmonic generation from spin-polarised defects in solids. <i>Npj Computational Materials</i> , 2020, 6, .	3.5	48
104	Contrasting Efficiency of Electron-Induced Reaction at Cu(110) in Aliphatic and Aromatic Bromides. <i>Journal of the American Chemical Society</i> , 2020, 142, 9453-9459.	6.6	4
105	Floquet states in dissipative open quantum systems. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2020, 53, 225601.	0.6	21
106	Berry curvature engineering by gating two-dimensional antiferromagnets. <i>Physical Review Research</i> , 2020, 2, .	1.3	22
107	Polaritonic coupled-cluster theory. <i>Physical Review Research</i> , 2020, 2, .	1.3	57
108	Floquet dynamics in light-driven solids. <i>Physical Review Research</i> , 2020, 2, .	1.3	33

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109	Polariton panorama. <i>Nanophotonics</i> , 2020, 10, 549-577.	2.9	155
110	Light-induced topological magnons in two-dimensional van der Waals magnets. <i>SciPost Physics</i> , 2020, 9, .	1.5	18
111	Role of electron scattering on the high-order harmonic generation from solids. <i>Physical Review Research</i> , 2020, 2, .	1.3	3
112	Role of intraband dynamics on circularly polarized high-harmonic generation from solids. , 2020, , .		0
113	Electronic Structure of Materials by Ab Initio Methods: Overview. , 2020, , 207-212.		0
114	Attosecond screening dynamics mediated by electron localization in transition metals. <i>Nature Physics</i> , 2019, 15, 1145-1149.	6.5	59
115	Quantum Electrodynamical Bloch Theory with Homogeneous Magnetic Fields. <i>Physical Review Letters</i> , 2019, 123, 047202.	2.9	30
116	Maximized electron interactions at the magic angle in twisted bilayer graphene. <i>Nature</i> , 2019, 572, 95-100.	13.7	644
117	Multiflat Bands and Strong Correlations in Twisted Bilayer Boron Nitride: Doping-Induced Correlated Insulator and Superconductor. <i>Nano Letters</i> , 2019, 19, 4934-4940.	4.5	123
118	Force balance approach for advanced approximations in density functional theories. <i>Journal of Chemical Physics</i> , 2019, 151, 154107.	1.2	17
119	Light-induced anomalous Hall effect in massless Dirac fermion systems and topological insulators with dissipation. <i>New Journal of Physics</i> , 2019, 21, 093005.	1.2	34
120	Self-Consistent Density-Functional Embedding: A Novel Approach for Density-Functional Approximations. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 5209-5220.	2.3	14
121	Reduced Density-Matrix Approach to Strong Matter-Photon Interaction. <i>ACS Photonics</i> , 2019, 6, 2694-2711.	3.2	31
122	Strong-field physics in the molecular frame. <i>EPJ Web of Conferences</i> , 2019, 205, 07002.	0.1	0
123	Light-Matter Response in Nonrelativistic Quantum Electrodynamics. <i>ACS Photonics</i> , 2019, 6, 2757-2778.	3.2	79
124	Universal optical control of chiral superconductors and Majorana modes. <i>Nature Physics</i> , 2019, 15, 766-770.	6.5	48
125	Capturing vacuum fluctuations and photon correlations in cavity quantum electrodynamics with multitrajjectory Ehrenfest dynamics. <i>Physical Review A</i> , 2019, 99, .	1.0	33
126	Polarization states of high-harmonics generated in silicon from elliptical drivers. <i>EPJ Web of Conferences</i> , 2019, 205, 02022.	0.1	0

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127	Microscopic theory for the light-induced anomalous Hall effect in graphene. <i>Physical Review B</i> , 2019, 99, .	1.1	117
128	Ab initio simulation of laser-induced water decomposition close to carbon nanotubes. <i>Physical Review B</i> , 2019, 99, .	1.1	10
129	Cavity Control of Excitons in Two-Dimensional Materials. <i>Nano Letters</i> , 2019, 19, 3473-3479.	4.5	65
130	Polarization-state-resolved high-harmonic spectroscopy of solids. <i>Nature Communications</i> , 2019, 10, 1319.	5.8	60
131	Local adsorption structure and bonding of porphine on Cu(111) before and after self-metalation. <i>Journal of Chemical Physics</i> , 2019, 150, 094702.	1.2	11
132	Density-Matrix Embedding Theory Study of the One-Dimensional Hubbard–Holstein Model. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 2221-2232.	2.3	22
133	Orbital magneto-optical response of periodic insulators from first principles. <i>Npj Computational Materials</i> , 2019, 5, .	3.5	5
134	Modification of excitation and charge transfer in cavity quantum-electrodynamical chemistry. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 4883-4892.	3.3	138
135	Unraveling materials Berry curvature and Chern numbers from real-time evolution of Bloch states. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 4135-4140.	3.3	20
136	Light-matter interactions within the Ehrenfest–Maxwell–Pauli–Kohn–Sham framework: fundamentals, implementation, and nano-optical applications. <i>Advances in Physics</i> , 2019, 68, 225-333.	35.9	54
137	Attosecond Electron Localization and Screening Dynamics in Metals. , 2019, , .		0
138	Optically Driven Attosecond Electron Dynamics in III-V Semiconductors. , 2019, , .		0
139	Exciton control in a room temperature bulk semiconductor with coherent strain pulses. <i>Science Advances</i> , 2019, 5, eaax2937.	4.7	28
140	Benchmarking semiclassical and perturbative methods for real-time simulations of cavity-bound emission and interference. <i>Journal of Chemical Physics</i> , 2019, 151, 244113.	1.2	37
141	Optimal control theory for quantum electrodynamics: an initial state problem. <i>European Physical Journal B</i> , 2019, 92, 1.	0.6	6
142	Nonadiabatic quantum dynamics without potential energy surfaces. <i>Physical Review Materials</i> , 2019, 3, .	0.9	12
143	Topological Floquet engineering of twisted bilayer graphene. <i>Physical Review Research</i> , 2019, 1, .	1.3	56
144	Electronic Structure of Materials by Ab Initio Methods: Overview. , 2019, , 1-6.		0

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145	From a quantum-electrodynamical light-matter description to novel spectroscopies. <i>Nature Reviews Chemistry</i> , 2018, 2, .	13.8	182
146	Ab Initio Optimized Effective Potentials for Real Molecules in Optical Cavities: Photon Contributions to the Molecular Ground State. <i>ACS Photonics</i> , 2018, 5, 992-1005.	3.2	96
147	Ultrasensitive H2S gas sensors based on p-type WS2 hybrid materials. <i>Nano Research</i> , 2018, 11, 4215-4224.	5.8	76
148	Layered Insulator/Molecule/Metal Heterostructures with Molecular Functionality through Porphyrin Intercalation. <i>ACS Nano</i> , 2018, 12, 2677-2684.	7.3	14
149	Transient Charge and Energy Flow in the Wide-Band Limit. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 2495-2504.	2.3	34
150	Propagators for the Time-Dependent Kohn-Sham Equations: Multistep, Runge-Kutta, Exponential Runge-Kutta, and Commutator Free Magnus Methods. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 3040-3052.	2.3	69
151	Atomic-like high-harmonic generation from two-dimensional materials. <i>Science Advances</i> , 2018, 4, eaao5207.	4.7	98
152	Phonon-driven spin-Floquet magneto-valleytronics in MoS2. <i>Nature Communications</i> , 2018, 9, 638.	5.8	86
153	Application of the Real-Time Time-Dependent Density Functional Theory to Excited-State Dynamics of Molecules and 2D Materials. <i>Journal of the Physical Society of Japan</i> , 2018, 87, 041016.	0.7	10
154	Adsorption Conformation and Lateral Registry of Cobalt Porphine on Cu(111). <i>Journal of Physical Chemistry C</i> , 2018, 122, 5452-5461.	1.5	14
155	Phonon Driven Floquet Matter. <i>Nano Letters</i> , 2018, 18, 1535-1542.	4.5	63
156	Attosecond optical-field-enhanced carrier injection into the GaAs conduction band. <i>Nature Physics</i> , 2018, 14, 560-564.	6.5	123
157	Light-matter interaction in the long-wavelength limit: no ground-state without dipole self-energy. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2018, 51, 034005.	0.6	132
158	Large area planar stanene epitaxially grown on Ag(111). <i>2D Materials</i> , 2018, 5, 025002.	2.0	164
159	Ab Initio Simulation of Attosecond Transient Absorption Spectroscopy in Two-Dimensional Materials. <i>Applied Sciences (Switzerland)</i> , 2018, 8, 1777.	1.3	19
160	All-optical nonequilibrium pathway to stabilising magnetic Weyl semimetals in pyrochlore iridates. <i>Nature Communications</i> , 2018, 9, 4452.	5.8	38
161	Electronic Structure of Materials by Ab Initio Methods: Overview. , 2018, , 1-6.		1
162	Atomic and Electronic Structure of Silicene on Ag: A Theoretical Perspective. <i>Nanoscience and Technology</i> , 2018, , 159-178.	1.5	1

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163	Visualizing topological edge states of single and double bilayer Bi supported on multilayer Bi(111) films. <i>Physical Review B</i> , 2018, 98, .	1.1	40
164	Special issue in honor of Eberhard K.U. Gross for his 65th birthday. <i>European Physical Journal B</i> , 2018, 91, 1.	0.6	0
165	Cavity quantum-electrodynamical polaritonically enhanced electron-phonon coupling and its influence on superconductivity. <i>Science Advances</i> , 2018, 4, eaau6969.	4.7	140
166	Benchmarking nonequilibrium Green's functions against configuration interaction for time-dependent Auger decay processes. <i>European Physical Journal B</i> , 2018, 91, 1.	0.6	7
167	Ab initio nonrelativistic quantum electrodynamics: Bridging quantum chemistry and quantum optics from weak to strong coupling. <i>Physical Review A</i> , 2018, 98, .	1.0	126
168	Multiple-orbital effects in laser-induced electron diffraction of aligned molecules. <i>Physical Review A</i> , 2018, 98, .	1.0	14
169	High-harmonic generation from few-layer hexagonal boron nitride: Evolution from monolayer to bulk response. <i>Physical Review B</i> , 2018, 98, .	1.1	54
170	Ultrafast Modification of Hubbard U in a Strongly Correlated Material: Ab Initio High-Harmonic Generation in NiO. <i>Physical Review Letters</i> , 2018, 121, 097402.	2.9	118
171	Coupled forward-backward trajectory approach for nonequilibrium electron-ion dynamics. <i>Physical Review B</i> , 2018, 97, .	1.1	12
172	Role of intraband transitions in photocarrier generation. <i>Physical Review B</i> , 2018, 98, .	1.1	26
173	Phonon-Driven Selective Modulation of Exciton Oscillator Strengths in Anatase TiO ₂ Nanoparticles. <i>Nano Letters</i> , 2018, 18, 5007-5014.	4.5	29
174	Kinetic-Energy Density-Functional Theory on a Lattice. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 4072-4087.	2.3	7
175	First-principles simulations for attosecond photoelectron spectroscopy based on time-dependent density functional theory. <i>European Physical Journal B</i> , 2018, 91, 1.	0.6	15
176	Electronic Structure of Materials by Ab Initio Methods: Overview. , 2018, , 1-6.		0
177	Light-matter interactions via the exact factorization approach. <i>European Physical Journal B</i> , 2018, 91, 1.	0.6	44
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