

Lorenz S Cederbaum

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4596775/publications.pdf>

Version: 2024-02-01

315
papers

12,963
citations

20759

60
h-index

37111

96
g-index

324
all docs

324
docs citations

324
times ranked

5189
citing authors

#	ARTICLE	IF	CITATIONS
1	BEYOND BORN-OPPENHEIMER: Molecular Dynamics Through a Conical Intersection. Annual Review of Physical Chemistry, 2004, 55, 127-158.	4.8	694
2	Direct calculation of ionization potentials of closed-shell atoms and molecules. Theoretica Chimica Acta, 1973, 31, 239-260.	0.9	312
3	Multiconfigurational time-dependent Hartree method for bosons: Many-body dynamics of bosonic systems. Physical Review A, 2008, 77, .	1.0	280
4	Non-Hermitian electronic theory and applications to clusters. Physics Reports, 2002, 368, 1-117.	10.3	261
5	Multiply Charged Anions in the Gas Phase. Chemical Reviews, 2002, 102, 181-200.	23.0	251
6	What will it take to observe processes in 'real time'?. Nature Photonics, 2014, 8, 162-166.	15.6	220
7	Role of Excited States in the Splitting of a Trapped Interacting Bose-Einstein Condensate by a Time-Dependent Barrier. Physical Review Letters, 2007, 99, 030402.	2.9	175
8	Mechanism of Interatomic Coulombic Decay in Clusters. Physical Review Letters, 2004, 93, 263002.	2.9	166
9	Multielectron wave-packet propagation: General theory and application. Journal of Chemical Physics, 2005, 123, 044111.	1.2	165
10	Exact Quantum Dynamics of a Bosonic Josephson Junction. Physical Review Letters, 2009, 103, 220601.	2.9	163
11	Time-dependent photodissociation of methyl iodide with five active modes. Journal of Chemical Physics, 1994, 101, 5623-5646.	1.2	162
12	Interatomic Coulombic Decay in van der Waals Clusters and Impact of Nuclear Motion. Physical Review Letters, 2000, 85, 4490-4493.	2.9	156
13	Electronic decay of valence holes in clusters and condensed matter. Physical Review B, 2001, 64, .	1.1	148
14	Electronic decay in weakly bound heteroclusters: Energy transfer versus electron transfer. Journal of Chemical Physics, 2001, 115, 5076-5088.	1.2	148
15	Ultrafast correlation-driven electron dynamics. Journal of Physics B: Atomic, Molecular and Optical Physics, 2014, 47, 124002.	0.6	145
16	Short-Time Dynamics Through Conical Intersections in Macrosystems. Physical Review Letters, 2005, 94, 113003.	2.9	140
17	Ab initio calculation of interatomic decay rates by a combination of the Fano ansatz, Green's-function methods, and the Stieltjes imaging technique. Journal of Chemical Physics, 2005, 123, 204107.	1.2	135
18	Ultralong-range energy transfer by interatomic Coulombic decay in an extreme quantum system. Nature Physics, 2010, 6, 508-511.	6.5	133

#	ARTICLE	IF	CITATIONS
19	Site- and energy-selective slow-electron production through intermolecular Coulombic decay. <i>Nature</i> , 2014, 505, 661-663.	13.7	131
20	Unified view on multiconfigurational time propagation for systems consisting of identical particles. <i>Journal of Chemical Physics</i> , 2007, 127, 154103.	1.2	124
21	Reduced density matrices and coherence of trapped interacting bosons. <i>Physical Review A</i> , 2008, 78, .	1.0	124
22	General variational many-body theory with complete self-consistency for trapped bosonic systems. <i>Physical Review A</i> , 2006, 73, .	1.0	119
23	Molecular double core hole electron spectroscopy for chemical analysis. <i>Journal of Chemical Physics</i> , 2010, 132, .	1.2	111
24	Interatomic and Intermolecular Coulombic Decay. <i>Chemical Reviews</i> , 2020, 120, 11295-11369.	23.0	106
25	Valence one-electron and shake-up ionization bands of polycyclic aromatic hydrocarbons. I. Benzene, naphthalene, anthracene, naphthacene, and pentacene. <i>Journal of Chemical Physics</i> , 2001, 115, 5859-5882.	1.2	103
26	Ultrafast charge migration in 2-phenylethyl-N,N-dimethylamine. <i>Chemical Physics Letters</i> , 2008, 450, 232-235.	1.2	103
27	Coulombic Energy Transfer and Triple Ionization in Clusters. <i>Physical Review Letters</i> , 2003, 90, 153401.	2.9	98
28	Charge migration in different conformers of glycine: The role of nuclear geometry. <i>Chemical Physics</i> , 2007, 338, 320-328.	0.9	98
29	Complex absorbing potentials in the framework of electron propagator theory. I. General formalism. <i>Journal of Chemical Physics</i> , 2002, 117, 5511-5521.	1.2	96
30	Laser-induced conical intersections in molecular optical lattices. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2008, 41, 221001.	0.6	95
31	Complex absorbing potentials in the framework of electron propagator theory. II. Application to temporary anions. <i>Journal of Chemical Physics</i> , 2003, 118, 6188-6199.	1.2	92
32	Dynamic Interference of Photoelectrons Produced by High-Frequency Laser Pulses. <i>Physical Review Letters</i> , 2012, 108, 253001.	2.9	92
33	Numerically exact quantum dynamics of bosons with time-dependent interactions of harmonic type. <i>Physical Review A</i> , 2012, 86, .	1.0	92
34	Electron Correlation as the Driving Force for Charge Transfer: Charge Migration Following Ionization in N-Methyl Acetamide. <i>Journal of Physical Chemistry A</i> , 2005, 109, 409-414.	1.1	91
35	Tracing Ultrafast Interatomic Electronic Decay Processes in Real Time and Space. <i>Physical Review Letters</i> , 2007, 98, 083201.	2.9	89
36	THE MULTI-MODE VIBRONIC-COUPPLING APPROACH. <i>Advanced Series in Physical Chemistry</i> , 2004, , 323-367.	1.5	86

#	ARTICLE	IF	CITATIONS
37	Pathway from Condensation via Fragmentation to Fermionization of Cold Bosonic Systems. <i>Physical Review Letters</i> , 2005, 95, 140402.	2.9	86
38	Interatomic electronic decay processes in singly and multiply ionized clusters. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 2011, 183, 36-47.	0.8	86
39	Environmental effects on a conical intersection: A model study. <i>Faraday Discussions</i> , 2004, 127, 395.	1.6	85
40	Interatomic Electronic Decay in Endohedral Fullerenes. <i>Physical Review Letters</i> , 2006, 96, 053401.	2.9	85
41	Ionization and double ionization of small water clusters. <i>Journal of Chemical Physics</i> , 2006, 125, 204305.	1.2	84
42	Born–Oppenheimer approximation and beyond for time-dependent electronic processes. <i>Journal of Chemical Physics</i> , 2008, 128, 124101.	1.2	82
43	Strong impact of light-induced conical intersections on the spectrum of diatomic molecules. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2011, 44, 045603.	0.6	82
44	Zoo of Quantum Phases and Excitations of Cold Bosonic Atoms in Optical Lattices. <i>Physical Review Letters</i> , 2005, 95, 030405.	2.9	80
45	An efficient combination of computational techniques for investigating electronic resonance states in molecules. <i>Journal of Chemical Physics</i> , 2001, 115, 6853-6861.	1.2	79
46	Short-time dynamics through conical intersections in macrosystems. I. Theory: Effective-mode formulation. <i>Journal of Chemical Physics</i> , 2006, 124, 144103.	1.2	79
47	Potential energy surface of the CO ₂ ⁻ anion. <i>Physical Chemistry Chemical Physics</i> , 2004, 6, 42.	1.3	74
48	Light-Induced Conical Intersections: Topological Phase, Wave Packet Dynamics, and Molecular Alignment. <i>Journal of Physical Chemistry A</i> , 2012, 116, 2636-2643.	1.1	74
49	Electronic Structure of the PYP Chromophore in Its Native Protein Environment. <i>Journal of the American Chemical Society</i> , 2007, 129, 6798-6806.	6.6	72
50	Charge migration following ionization in systems with chromophore-donor and amine-acceptor sites. <i>Journal of Chemical Physics</i> , 2008, 129, 104305.	1.2	72
51	Core Ionization Initiates Subfemtosecond Charge Migration in the Valence Shell of Molecules. <i>Physical Review Letters</i> , 2016, 117, 093002.	2.9	72
52	Reactive Scattering Dynamics on Conically Intersecting Potential Energy Surfaces: The H + H ₂ Exchange Reaction. <i>Journal of Physical Chemistry A</i> , 2001, 105, 2321-2329.	1.1	71
53	The exact molecular wavefunction as a product of an electronic and a nuclear wavefunction. <i>Journal of Chemical Physics</i> , 2013, 138, 224110.	1.2	71
54	Existence of a Correlation Bound <i>i</i> -Type Anion State of C ₆₀ . <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 849-853.	2.1	71

#	ARTICLE	IF	CITATIONS
55	Electronic structure of the photoactive yellow protein chromophore: Ab initio study of the low-lying excited singlet states. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2007, 190, 241-257.	2.0	67
56	Ultrafast Interatomic Electronic Decay in Multiply Excited Clusters. <i>Physical Review Letters</i> , 2010, 105, 043004.	2.9	67
57	Fingerprints of the nodal structure of autoionizing vibrational wave functions in clusters: Interatomic Coulombic decay in Ne dimer. <i>Journal of Chemical Physics</i> , 2001, 114, 7351-7360.	1.2	64
58	Exact ground state of finite Bose-Einstein condensates on a ring. <i>Physical Review A</i> , 2005, 72, .	1.0	64
59	Scattering of an attractive Bose-Einstein condensate from a barrier: Formation of quantum superposition states. <i>Physical Review A</i> , 2009, 80, .	1.0	64
60	Resonant Auger Decay of Molecules in Intense X-Ray Laser Fields: Light-Induced Strong Nonadiabatic Effects. <i>Physical Review Letters</i> , 2011, 106, 123001.	2.9	63
61	Electron Impact Catalytic Dissociation: Two-Bond Breaking by a Low-Energy Catalytic Electron. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 4119-4122.	7.2	62
62	Light-induced conical intersections in polyatomic molecules: General theory, strategies of exploitation, and application. <i>Journal of Chemical Physics</i> , 2013, 139, 154314.	1.2	62
63	Strong interference effects in the resonant Auger decay of atoms induced by intense x-ray fields. <i>Physical Review A</i> , 2011, 83, .	1.0	60
64	Foreign imaging in Auger spectroscopy: The Si 2p spectrum of silicon tetrafluoride. <i>Physical Review Letters</i> , 1993, 71, 649-652.	2.9	59
65	Formation and Dynamics of Many-Boson Fragmented States in One-Dimensional Attractive Ultracold Gases. <i>Physical Review Letters</i> , 2008, 100, 130401.	2.9	59
66	Electron-correlation-driven charge migration in oligopeptides. <i>Chemical Physics</i> , 2013, 414, 100-105.	0.9	59
67	Conical Intersections Induced by Quantum Light: Field-Dressed Spectra from the Weak to the Ultrastrong Coupling Regimes. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 6215-6223.	2.1	59
68	Ab initio lifetimes in the interatomic Coulombic decay of neon clusters computed with propagators. <i>Journal of Chemical Physics</i> , 2007, 126, 164110.	1.2	58
69	Conical intersections induced by light: Berry phase and wavepacket dynamics. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2011, 44, 175102.	0.6	57
70	How an interacting many-body system tunnels through a potential barrier to open space. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 13521-13525.	3.3	55
71	Intermolecular Coulombic Decay in Small Biochemically Relevant Hydrogen-Bonded Systems. <i>Journal of the American Chemical Society</i> , 2011, 133, 6817-6824.	6.6	53
72	Ground-state fragmentation of repulsive Bose-Einstein condensates in double-trap potentials. <i>Physical Review A</i> , 2004, 70, .	1.0	52

#	ARTICLE	IF	CITATIONS
73	Impact of Sulfur vs Oxygen on the Low-Lying Excited States of trans-p-Coumaric Acid and trans-p-Coumaric Thio Acid. Journal of Physical Chemistry A, 2005, 109, 4623-4631.	1.1	52
74	Quantum dynamics of attractive versus repulsive bosonic Josephson junctions: Bose-Hubbard and full-Hamiltonian results. Physical Review A, 2010, 82, .	1.0	52
75	Observation of electron-transfer-mediated decay in aqueous solution. Nature Chemistry, 2017, 9, 708-714.	6.6	51
76	Methylboron Oxide, H ₃ CBO. Angewandte Chemie International Edition in English, 1989, 28, 88-90.	4.4	50
77	Intersections of potential energy surfaces of short-lived states: The complex analogue of conical intersections. Journal of Chemical Physics, 2004, 120, 3201-3214.	1.2	50
78	Multiconfigurational time-dependent Hartree method for mixtures consisting of two types of identical particles. Physical Review A, 2007, 76, .	1.0	50
79	Ultrafast Charge Migration Following Valence Ionization of 4-Methylphenol: Jumping over the Aromatic Ring. Journal of Physical Chemistry A, 2010, 114, 8676-8679.	1.1	49
80	A One-Step Four-Bond-Breaking Reaction Catalyzed by an Electron. Angewandte Chemie - International Edition, 2012, 51, 8003-8007.	7.2	48
81	Interatomic decay of inner-valence-excited states in clusters. Journal of Chemical Physics, 2006, 124, 144315.	1.2	47
82	General mapping for bosonic and fermionic operators in Fock space. Physical Review A, 2010, 81, .	1.0	47
83	Wave chaos as signature for depletion of a Bose-Einstein condensate. Physical Review A, 2012, 86, .	1.0	46
84	Impact of interatomic electronic decay processes on Xe ⁵⁴ d hole decay in the xenon fluorides. Journal of Chemical Physics, 2003, 119, 10575-10584.	1.2	45
85	Accurate multi-boson long-time dynamics in triple-well periodic traps. Physical Review A, 2011, 83, .	1.0	45
86	Electronic decay following ionization of aqueous Li ⁺ microsolvation clusters. Journal of Chemical Physics, 2005, 122, 094305.	1.2	44
87	Universality of fragmentation in the Schrödinger dynamics of bosonic Josephson junctions. Physical Review A, 2014, 89, .	1.0	44
88	Evidence for a partial breakdown of the molecular orbital picture in the ionization spectra of large saturated hydrocarbons. Journal of Chemical Physics, 1996, 105, 7583-7596.	1.2	43
89	Coherent intense resonant laser pulses lead to interference in the time domain seen in the spectrum of the emitted particles. Physical Review A, 2012, 86, .	1.0	43
90	Nuclear-wave-packet quantum interference in the intense laser dissociation of the D ₂ molecule. Physical Review A, 2013, 88, .	1.0	43

#	ARTICLE	IF	CITATIONS
91	Resonant Auger decay of the core-excited C O molecule in intense x-ray laser fields. <i>Physical Review A</i> , 2011, 84, .	1.0	42
92	Competition between Light-Induced and Intrinsic Nonadiabatic Phenomena in Diatomics. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 1624-1630.	2.1	42
93	Calculation of interatomic decay widths of vacancy states delocalized due to inversion symmetry. <i>Journal of Chemical Physics</i> , 2006, 125, 094107.	1.2	41
94	On the interatomic electronic processes following Auger decay in neon dimer. <i>Journal of Chemical Physics</i> , 2008, 129, 074307.	1.2	41
95	On the size-dependence of the static self-energy in propagator calculations. <i>Journal of Chemical Physics</i> , 1995, 103, 3578-3588.	1.2	40
96	Short-time dynamics through conical intersections in macrosystems. II. Applications. <i>Journal of Chemical Physics</i> , 2006, 124, 144104.	1.2	40
97	Photoinduced Isomerization of the Photoactive Yellow Protein (PYP) Chromophore: Interplay of Two Torsions, a HOOP Mode and Hydrogen Bonding. <i>Journal of Physical Chemistry A</i> , 2011, 115, 9237-9248.	1.1	40
98	Dynamics of interatomic Coulombic decay in quantum dots. <i>Journal of Chemical Physics</i> , 2011, 135, 144112.	1.2	40
99	Proton-Transfer Mediated Enhancement of Nonlocal Electronic Relaxation Processes in X-ray Irradiated Liquid Water. <i>Journal of the American Chemical Society</i> , 2014, 136, 18170-18176.	6.6	40
100	Correlation effects in the valence x-ray photoionization spectra of ethylene, butadiene, and hexatriene. <i>International Journal of Quantum Chemistry</i> , 1997, 63, 465-481.	1.0	39
101	<i>Ab initio</i> calculation of interatomic decay rates of excited doubly ionized states in clusters. <i>Journal of Chemical Physics</i> , 2008, 129, 244102.	1.2	39
102	Swift Loss of Coherence of Soliton Trains in Attractive Bose-Einstein Condensates. <i>Physical Review Letters</i> , 2011, 106, 240401.	2.9	39
103	Benchmark Calculations of the Energies for Binding Excess Electrons to Water Clusters. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 893-900.	2.3	39
104	Two trapped particles interacting by a finite-range two-body potential in two spatial dimensions. <i>Physical Review A</i> , 2013, 87, .	1.0	39
105	Interatomic Coulombic decay in a heteroatomic rare gas cluster. <i>Journal of Chemical Physics</i> , 2006, 124, 154305.	1.2	38
106	Many-body tunneling dynamics of Bose-Einstein condensates and vortex states in two spatial dimensions. <i>Physical Review A</i> , 2015, 92, .	1.0	38
107	Demixing of Bosonic Mixtures in Optical Lattices from Macroscopic to Microscopic Scales. <i>Physical Review Letters</i> , 2006, 97, 230403.	2.9	37
108	Many-body theory for systems with particle conversion: Extending the multiconfigurational time-dependent Hartree method. <i>Physical Review A</i> , 2009, 79, .	1.0	37

#	ARTICLE	IF	CITATIONS
109	Environment assisted electron capture. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2009, 42, 231001.	0.6	37
110	ac Stark effect in the electronic continuum and its impact on the photoionization of atoms by coherent intense short high-frequency laser pulses. <i>Physical Review A</i> , 2013, 88, .	1.0	37
111	Exact decay and tunnelling dynamics of interacting few-boson systems. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2009, 42, 044018.	0.6	36
112	Generation of Highly Damaging H_{2}O^{+} Radicals by Inner Valence Shell Ionization of Water. <i>ChemPhysChem</i> , 2010, 11, 1006-1009.	1.0	36
113	Tracing molecular electronic excitation dynamics in real time and space. <i>Journal of Chemical Physics</i> , 2010, 132, 144302.	1.2	36
114	Extreme Correlation Effects in the Elusive Bound Spectrum of C_{60}^{+} . <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 3319-3324.	2.1	36
115	Phantom vortices: hidden angular momentum in ultracold dilute Bose-Einstein condensates. <i>Scientific Reports</i> , 2017, 7, 40122.	1.6	36
116	Light-induced conical intersections for short and long laser pulses: Floquet and rotating wave approximations versus numerical exact results. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2012, 45, 135101.	0.6	35
117	Formation of satellite bands in the ionization spectra of extended systems. <i>Physical Review B</i> , 1996, 53, 13326-13339.	1.1	34
118	Time-dependent multi-orbital mean-field for fragmented Bose-Einstein condensates. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2007, 362, 453-459.	0.9	34
119	Correlation-bound anions of NaCl clusters. <i>Journal of Chemical Physics</i> , 2010, 133, 114301.	1.2	33
120	Simulation of a complex spectrum: Interplay of five electronic states and 21 vibrational degrees of freedom in C_5H_4^+ . <i>Journal of Chemical Physics</i> , 2005, 123, 204310.	1.2	32
121	Coupled-cluster theory for systems of bosons in external traps. <i>Physical Review A</i> , 2006, 73, .	1.0	32
122	Interferences in the Density of Two Bose-Einstein Condensates Consisting of Identical or Different Atoms. <i>Physical Review Letters</i> , 2007, 98, 110405.	2.9	32
123	Ultrafast electron dynamics following outer-valence ionization: The impact of low-lying relaxation satellite states. <i>Journal of Chemical Physics</i> , 2009, 130, 154305.	1.2	32
124	Interatomic Electronic Decay Driven by Nuclear Motion. <i>Physical Review Letters</i> , 2010, 105, 173401.	2.9	32
125	Ionic-Charge Dependence of the Intermolecular Coulombic Decay Time Scale for Aqueous Ions Probed by the Core-Hole Clock. <i>Journal of the American Chemical Society</i> , 2011, 133, 13430-13436.	6.6	32
126	Excitation spectra of many-body systems by linear response: General theory and applications to trapped condensates. <i>Physical Review A</i> , 2013, 88, .	1.0	32

#	ARTICLE	IF	CITATIONS
127	Controlled energy-selected electron capture and release in double quantum dots. <i>Physical Review B</i> , 2013, 88, .	1.1	32
128	Influence of Light-Induced Conical Intersection on the Photodissociation Dynamics of $D_{2^{+}}$ Starting from Individual Vibrational Levels. <i>Journal of Physical Chemistry A</i> , 2014, 118, 11908-11915.	1.1	32
129	Generic regimes of quantum many-body dynamics of trapped bosonic systems with strong repulsive interactions. <i>Physical Review A</i> , 2014, 89, .	1.0	32
130	Auger Electron Spectroscopy as a Probe of the Solution of Aqueous Ions. <i>Journal of the American Chemical Society</i> , 2009, 131, 7264-7271.	6.6	31
131	On the Cholesky decomposition for electron propagator methods: General aspects and application on C ₆₀ . <i>Journal of Chemical Physics</i> , 2010, 132, 044110.	1.2	31
132	Breaking the resilience of a two-dimensional Bose-Einstein condensate to fragmentation. <i>Physical Review A</i> , 2014, 90, .	1.0	31
133	On the doubly ionized states of Ar ₂ and their intra- and interatomic decay to Ar ₂ ³⁺ . <i>Journal of Chemical Physics</i> , 2008, 128, 014307.	1.2	30
134	Effect of Light-Induced Conical Intersection on the Photodissociation Dynamics of the $D_{2^{+}}$ Molecule. <i>Journal of Physical Chemistry A</i> , 2013, 117, 8528-8535.	1.1	30
135	All for one and one for all: accommodating an extra electron in C ₆₀ . <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 13287.	1.3	30
136	An effective Hamiltonian for the short-time dynamics at a conical intersection. <i>Molecular Physics</i> , 2006, 104, 1081-1093.	0.8	29
137	High activity of helium droplets following ionization of systems inside those droplets. <i>Physical Review B</i> , 2007, 76, .	1.1	29
138	Dynamic interference in the photoionization of He by coherent intense high-frequency laser pulses: Direct propagation of the two-electron wave packets on large spatial grids. <i>Physical Review A</i> , 2016, 93, .	1.0	29
139	Exploring Interatomic Coulombic Decay by Free Electron Lasers. <i>Physical Review Letters</i> , 2011, 107, 273002.	2.9	28
140	Towards controlling the dissociation probability by light-induced conical intersections. <i>Faraday Discussions</i> , 2016, 194, 479-493.	1.6	28
141	Direct Signatures of Light-Induced Conical Intersections on the Field-Dressed Spectrum of Na ₂ . <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 2739-2745.	2.1	28
142	Extrapolating bound state data of anions into the metastable domain. <i>Journal of Chemical Physics</i> , 2004, 121, 6628-6633.	1.2	27
143	Using pH-Value To Control Intermolecular Electronic Decay. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 1306-1309.	7.2	27
144	Anions of Xenon Clusters Bound by Long-Range Electron Correlations. <i>Physical Review Letters</i> , 2011, 107, 133401.	2.9	27

#	ARTICLE	IF	CITATIONS
145	Bornâ€œOppenheimer approximation in optical cavities: from success to breakdown. <i>Chemical Science</i> , 2021, 12, 1251-1258.	3.7	27
146	Stable Free Dianionic Silicoï¿½Carbon Clusters. <i>Angewandte Chemie International Edition in English</i> , 1997, 36, 1889-1891.	4.4	26
147	Nonlocal Effects in the Core Ionization and Auger Spectra of Small Ammonia Clusters. <i>Journal of Physical Chemistry B</i> , 2011, 115, 5441-5447.	1.2	26
148	Damaging Intermolecular Energy and Proton Transfer Processes in Alphaâ€œParticleâ€œIrradiated Hydrogenâ€œBonded Systems. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 17023-17027.	7.2	26
149	Bound states of negatively charged ions induced by a magnetic field. <i>Physical Review A</i> , 2000, 61, .	1.0	25
150	Ultrafast Photoinitiated Long-Range Electron Transfer in Cyclophane-Bridged Zincporphyrinâ€œQuinone Complexes via Conical Intersections. <i>Journal of Physical Chemistry B</i> , 2004, 108, 19049-19055.	1.2	24
151	Impact of nuclear dynamics on interatomic Coulombic decay in a He dimer. <i>Physical Review A</i> , 2010, 82, .	1.0	24
152	The effect of light-induced conical intersections on the alignment of diatomic molecules. <i>Chemical Physics</i> , 2012, 399, 146-150.	0.9	24
153	Nuclear dynamics during the resonant Auger decay of water molecules. <i>Journal of Chemical Physics</i> , 2009, 130, 154307.	1.2	23
154	Ultrafast charge separation driven by differential particle and hole mobilities. <i>Journal of Chemical Physics</i> , 2011, 134, 024303.	1.2	23
155	Recursive formulation of the multiconfigurational time-dependent Hartree method for fermions, bosons and mixtures thereof in terms of one-body density operators. <i>Chemical Physics</i> , 2012, 401, 2-14.	0.9	23
156	Charge transfer driven by electron correlation: A non-Dyson propagator approach. <i>Journal of Chemical Physics</i> , 2005, 122, 134104.	1.2	22
157	Allene and pentatetraene cations as models for intramolecular charge transfer: Vibronic coupling Hamiltonian and conical intersections. <i>Journal of Chemical Physics</i> , 2005, 122, 144320.	1.2	22
158	Quantum dynamics in macrosystems with several coupled electronic states: Hierarchy of effective Hamiltonians. <i>Journal of Chemical Physics</i> , 2007, 127, 124107.	1.2	22
159	Overlap of exact and Gross-Pitaevskii wave functions in Bose-Einstein condensates of dilute gases. <i>Physical Review A</i> , 2016, 94, .	1.0	22
160	Suppression of electron correlation and of autoionization by strong laser fields. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 1999, 32, L279-L284.	0.6	21
161	Efficient generation and properties of mesoscopic quantum superposition states in an attractive Boseâ€œEinstein condensate threaded by a potential barrier. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2009, 42, 091004.	0.6	21
162	Optimal time-dependent lattice models for nonequilibrium dynamics. <i>New Journal of Physics</i> , 2011, 13, 043003.	1.2	21

#	ARTICLE	IF	CITATIONS
163	Photoionization of hydrogen atoms by coherent intense high-frequency short laser pulses: Direct propagation of electron wave packets on large spatial grids. <i>Physical Review A</i> , 2013, 88, .	1.0	21
164	Controlling the velocities and the number of emitted particles in the tunneling to open space dynamics. <i>Physical Review A</i> , 2014, 89, .	1.0	21
165	Discovery of a new class of stable gas-phase dianions: Mixed oxygen-carbon cluster OCn^{2-} ($n=5-19$). <i>Journal of Chemical Physics</i> , 2002, 117, 7002-7009.	1.2	20
166	Influence of Delocalization on the Stability of Dianions: A Study of a Systematic Series of Dianions with Growing Electronic Localization. <i>Journal of the American Chemical Society</i> , 2003, 125, 9531-9537.	6.6	20
167	Impact of intense laser pulses on the autoionization dynamics of the 2s2p doubly excited state of He. <i>Physical Review A</i> , 2017, 96, .	1.0	20
168	Dianionic Tetraborates Do Exist as Stable Entities. <i>Journal of the American Chemical Society</i> , 2002, 124, 10903-10910.	6.6	19
169	Microsolvation of Li^+ in Water Analyzed by Ionization and Double Ionization. <i>Journal of Physical Chemistry A</i> , 2004, 108, 5831-5844.	1.1	19
170	Fragmented Metastable States Exist in an Attractive Bose-Einstein Condensate for Atom Numbers Well above the Critical Number of the Gross-Pitaevskii Theory. <i>Physical Review Letters</i> , 2008, 100, 040402.	2.9	19
171	Interatomic Coulombic decay in a He dimer: <i>Ab initio</i> potential-energy curves and decay widths. <i>Physical Review A</i> , 2010, 82, .	1.0	19
172	An Excited Electron Avoiding a Positive Charge. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2300-2303.	2.1	19
173	Native hydrogen bonding network of the photoactive yellow protein (PYP) chromophore: Impact on the electronic structure and photoinduced isomerization. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2012, 234, 123-134.	2.0	19
174	Detecting ultrafast interatomic electronic processes in media by fluorescence. <i>New Journal of Physics</i> , 2014, 16, 102002.	1.2	19
175	The exact wavefunction of interacting N degrees of freedom as a product of N single-degree-of-freedom wavefunctions. <i>Chemical Physics</i> , 2015, 457, 129-132.	0.9	19
176	Damaging Intermolecular Energy and Proton Transfer Processes in α -Particle-Irradiated Hydrogen-Bonded Systems. <i>Angewandte Chemie</i> , 2018, 130, 17269-17273.	1.6	19
177	Real-time observation of X-ray-induced intramolecular and interatomic electronic decay in CH_2I_2 . <i>Nature Communications</i> , 2019, 10, 2186.	5.8	19
178	On the intermolecular Coulombic decay of singly and doubly ionized states of water dimer. <i>Journal of Chemical Physics</i> , 2010, 133, 154307.	1.2	18
179	Time-resolved pump-probe spectroscopy to follow valence electronic motion in molecules: Theory. <i>Physical Review A</i> , 2013, 88, .	1.0	18
180	The exact wavefunction factorization of a vibronic coupling system. <i>Journal of Chemical Physics</i> , 2014, 140, 054104.	1.2	18

#	ARTICLE	IF	CITATIONS
181	Intrinsic and light-induced nonadiabatic phenomena in the NaI molecule. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 19656-19664.	1.3	18
182	Impact of cavity on interatomic Coulombic decay. <i>Nature Communications</i> , 2021, 12, 4083.	5.8	18
183	Linewidth and lifetime of atomic levels and the time evolution of spectra and coincidence spectra. <i>Physical Review A</i> , 2010, 81, .	1.0	17
184	Exact many-body wave function and properties of trapped bosons in the infinite-particle limit. <i>Physical Review A</i> , 2017, 96, .	1.0	17
185	Variance of an anisotropic Bose-Einstein condensate. <i>Chemical Physics</i> , 2018, 509, 45-54.	0.9	17
186	Interatomic Coulombic electron capture from first principles. <i>Physical Review A</i> , 2018, 98, .	1.0	17
187	Finite-size effects and quantum phonon fluctuations in the optical absorption edge of dimerized chains. <i>Physical Review B</i> , 1997, 55, 1481-1485.	1.1	16
188	Ionization of the xenon fluorides. <i>Journal of Chemical Physics</i> , 2003, 119, 7763-7771.	1.2	16
189	Electronic structure of isolated PtX_6^{2-} ($X=\text{F,Cl,Br}$) dianions. <i>Journal of Chemical Physics</i> , 2003, 118, 1747-1755.	1.2	16
190	Jahn-Teller effect for short-lived states: Study of the complex potential energy surfaces. <i>Journal of Chemical Physics</i> , 2004, 121, 5.	1.2	16
191	Quantum hydrodynamics: Mixed states, dissipation, and a new hybrid quantum-classical approach. <i>International Journal of Quantum Chemistry</i> , 2004, 100, 1153-1162.	1.0	16
192	Accurate Quantum Chemistry in Single Precision Arithmetic: Correlation Energy. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 320-326.	2.3	16
193	Time-resolved pump-probe spectroscopy to follow valence electronic motion in molecules: Application. <i>Physical Review A</i> , 2014, 90, .	1.0	16
194	Following the Birth of a Nanoplasma Produced by an Ultrashort Hard-X-Ray Laser in Xenon Clusters. <i>Physical Review X</i> , 2018, 8, .	2.8	16
195	Diisocyanogen or Isocyanogen?. <i>Angewandte Chemie International Edition in English</i> , 1989, 28, 761-762.	4.4	15
196	Fragmented many-body states of definite angular momentum and stability of attractive three-dimensional condensates. <i>Physical Review A</i> , 2010, 82, .	1.0	15
197	Interrelation between the Distributions of Kinetic Energy Release and Emitted Electron Energy following the Decay of Electronic States. <i>Physical Review Letters</i> , 2011, 107, 173001.	2.9	15
198	Ultrafast reorganization of the hole charge created upon outer-valence ionization of porphyrins. <i>Chemical Physics</i> , 2012, 399, 245-251.	0.9	15

#	ARTICLE	IF	CITATIONS
199	Photodissociation of D2+ induced by linearly chirped laser pulses. <i>Journal of Chemical Physics</i> , 2015, 143, 014305.	1.2	15
200	Barrierless Single-Electron-Induced <i>cis</i> → <i>trans</i> Isomerization. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 10470-10473.	7.2	15
201	Dynamic interference in the resonance-enhanced multiphoton ionization of hydrogen atoms by short and intense laser pulses. <i>Chemical Physics</i> , 2018, 509, 145-150.	0.9	15
202	Core-level interatomic Coulombic decay in van der Waals clusters. <i>Physical Review Research</i> , 2020, 2, .	1.3	15
203	Ultrafast excited-state charge transfer at a conical intersection: effects of an environment. <i>Computer Physics Communications</i> , 2005, 169, 95-98.	3.0	14
204	Properties of fragmented repulsive condensates. <i>Physical Review A</i> , 2005, 71, .	1.0	14
205	Stable and Long-Lived Trianions in the Gas Phase. <i>Journal of Physical Chemistry A</i> , 2005, 109, 11401-11406.	1.1	14
206	Combined experimental-theoretical study of the lower excited singlet states of paravinyl phenol, an analog of the paracoumaric acid chromophore. <i>Journal of Chemical Physics</i> , 2006, 125, 204303.	1.2	14
207	Excitation spectra of fragmented condensates by linear response: General theory and application to a condensate in a double-well potential. <i>Physical Review A</i> , 2012, 86, .	1.0	14
208	Dynamics and symmetries of a repulsively bound atom pair in an infinite optical lattice. <i>Physical Review A</i> , 2012, 86, .	1.0	14
209	Tracking the photodissociation probability of D2+ induced by linearly chirped laser pulses. <i>Journal of Chemical Physics</i> , 2016, 144, 074309.	1.2	14
210	Ultrafast Intermolecular Energy Transfer from Vibrations to Electronic Motion. <i>Physical Review Letters</i> , 2018, 121, 223001.	2.9	14
211	Ab initio complex potential energy curves of the He*(1s2p 1P)–Li dimer. <i>Journal of Chemical Physics</i> , 2020, 152, 184303.	1.2	14
212	Orbital Picture of Ionization and Its Breakdown in Nanoarrays of Quantum Dots. <i>Physical Review Letters</i> , 2002, 89, 133003.	2.9	13
213	Interatomic relaxation effects in double core ionization of chain molecules. <i>Journal of Chemical Physics</i> , 2012, 137, 154316.	1.2	13
214	Unified view on linear response of interacting identical and distinguishable particles from multiconfigurational time-dependent Hartree methods. <i>Journal of Chemical Physics</i> , 2014, 140, 034108.	1.2	13
215	The best orbital and pair function for describing ionic and excited states on top of the exact ground state. <i>Journal of Chemical Physics</i> , 2014, 141, 194102.	1.2	13
216	Enhanced many-body effects in the excitation spectrum of a weakly interacting rotating Bose-Einstein condensate. <i>Physical Review A</i> , 2018, 98, .	1.0	13

#	ARTICLE	IF	CITATIONS
217	Valence ionization spectra of disubstituted s-tetrazines: strong correlation effects induced by substitution. <i>Journal of the American Chemical Society</i> , 1990, 112, 94-102.	6.6	12
218	Extended Two-Particle Green's Functions and Optical Potentials for Two-Particle Scattering by Many-Body Targets. <i>Annals of Physics</i> , 1996, 252, 276-299.	1.0	12
219	Transition from Rydberg to giant-dipole-moment states of hydrogen atoms in crossed fields: A suggestion for an experiment. <i>Physical Review A</i> , 1999, 59, 3695-3700.	1.0	12
220	Stability of Negatively Charged Ions Moving in a Magnetic Field. <i>Physical Review Letters</i> , 2001, 86, 5450-5453.	2.9	12
221	Cyclic Carbon Cluster Dianions and Their Aromaticity. <i>Journal of the American Chemical Society</i> , 2002, 124, 3163-3168.	6.6	12
222	Effect of relativity on the ionization spectra of the xenon fluorides XeFn (n=2, 4, 6). <i>Journal of Chemical Physics</i> , 2005, 122, 214302.	1.2	12
223	How many bound valence states does the C ₆₀ ²⁻ anion have?. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 10840-10845.	1.3	12
224	Communication: Substantial impact of the orientation of transition dipole moments on the dynamics of diatomics in laser fields. <i>Journal of Chemical Physics</i> , 2018, 149, 181101.	1.2	12
225	Charge separated states of endohedral fullerene Li@C ₂₀ . <i>Journal of Chemical Physics</i> , 2019, 151, 114306.	1.2	12
226	Long-lived Gas-phase Dianions Containing Tetrahedrally Coordinated Oxygen Atoms: O(BN) and O(C ₂). <i>Journal of Physical Chemistry A</i> , 2002, 106, 1406-1408.	1.1	11
227	Magnetically induced anions. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 4981.	1.3	11
228	Ionization spectra and electronic decay in small iodide clusters: Fully relativistic results. <i>Journal of Chemical Physics</i> , 2006, 125, 034309.	1.2	11
229	PtF ₆ ²⁻ dianion and its detachment spectrum: A fully relativistic study. <i>Journal of Chemical Physics</i> , 2007, 126, 144310.	1.2	11
230	Photodetachment spectra of the PtX ₄ ²⁻ (X=F,Cl,Br) dianions and their Jahn-Teller distortions: A fully relativistic study. <i>Journal of Chemical Physics</i> , 2008, 129, 174302.	1.2	11
231	Elastic scattering of a Bose-Einstein condensate at a potential landscape. <i>Journal of Physics: Conference Series</i> , 2014, 488, 012032.	0.3	11
232	Electron-correlation driven capture and release in double quantum dots. <i>Journal of Physics Condensed Matter</i> , 2016, 28, 075301.	0.7	11
233	Attractive Bose-Einstein condensates in anharmonic traps: Accurate numerical treatment and the intriguing physics of the variance. <i>Chemical Physics</i> , 2018, 515, 287-298.	0.9	11
234	Inner-valence ionization of molecular anions and ultrafast relaxation by electron emission. <i>Chemical Physics Letters</i> , 2000, 324, 416-422.	1.2	10

#	ARTICLE	IF	CITATIONS
235	Systematic corrections to the equivalent core model. <i>Journal of Chemical Physics</i> , 2002, 116, 8723-8730.	1.2	10
236	Competitive charge- and energy-transfer processes following core ionization in the Na ⁺ CO cluster. <i>Journal of Chemical Physics</i> , 2005, 123, 154308.	1.2	10
237	Trapping of cold atoms in optical lattices by the quadrupole force. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2007, 362, 215-220.	0.9	10
238	Strong enhancement of cage effects in water photolysis caused by interatomic Coulombic decay. <i>Journal of Chemical Physics</i> , 2016, 144, 164307.	1.2	10
239	Many-body effects in the excitation spectrum of weakly interacting Bose-Einstein condensates in one-dimensional optical lattices. <i>Physical Review A</i> , 2017, 95, .	1.0	10
240	Caged-electron states and split-electron states in the endohedral alkali C ₆₀ . <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 11837-11843.	1.3	10
241	How small can a Peierls dimerized chain be?. <i>Solid State Communications</i> , 1998, 106, 733-737.	0.9	9
242	Equivalent core model: Extended theory and applications. <i>Journal of Chemical Physics</i> , 2003, 118, 2081-2091.	1.2	9
243	Build-up of coherence between initially-independent subsystems: The case of Bose-Einstein condensates. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2009, 373, 301-304.	0.9	9
244	Influence of caged noble-gas atom on the superatomic and valence states of C ₆₀ ⁺ . <i>Molecular Physics</i> , 2015, 113, 2964-2969.	0.8	9
245	Observation of fast and slow interatomic Coulombic decay in argon dimers induced by electron-impact ionization. <i>Physical Review A</i> , 2017, 96, .	1.0	9
246	Bound electronic states of the smallest fullerene C ₂₀ ⁻ anion. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 17434-17441.	1.3	9
247	Caged-Electron States in Endohedral Li Fullerenes. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 7617-7622.	2.1	9
248	Polaritonic States of Matter in a Rotating Cavity. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 6056-6061.	2.1	9
249	Endocircular Li Carbon Rings. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 16649-16654.	7.2	9
250	Quantum light-induced nonadiabatic phenomena in the absorption spectrum of formaldehyde: Full- and reduced-dimensionality studies. <i>Journal of Chemical Physics</i> , 2020, 153, 234302.	1.2	9
251	Cooperative molecular structure in polaritonic and dark states. <i>Journal of Chemical Physics</i> , 2022, 156, 184102.	1.2	9
252	Quantum Phonon Fluctuations in Mesoscopic Dimerized Systems. <i>Journal of the Physical Society of Japan</i> , 1999, 68, 1954-1962.	0.7	8

#	ARTICLE	IF	CITATIONS
253	Charge transfer in the Cl ⁺ CO cluster induced by core ionization. <i>Journal of Chemical Physics</i> , 2005, 122, 104304.	1.2	8
254	Theory of magnetically induced anions. <i>Physical Review A</i> , 2007, 75, .	1.0	8
255	Exploring Nonadiabatic Effects by Recoil of Fast Photoelectrons. <i>Physical Review Letters</i> , 2009, 103, 133001.	2.9	8
256	Exploring Protonation and Deprotonation Effects with Auger Electron Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2733-2737.	2.1	8
257	Fractional driven-damped oscillator and its general closed form exact solution. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2018, 505, 744-762.	1.2	8
258	A Concerted Synchronous [2 + 2] Cycloreversion Repair Catalyzed by Two Electrons. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 6973-6977.	2.1	8
259	Striking Generic Impact of Light-Induced Non-Adiabaticity in Polyatomic Molecules. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 5324-5329.	2.1	8
260	Suppression of X-ray-Induced Radiation Damage to Biomolecules in Aqueous Environments by Immediate Intermolecular Decay of Inner-Shell Vacancies. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 7146-7150.	2.1	8
261	Coexistence of short- and large-scale phase variations in a charge-density wave weakly coupled to impurities. <i>Physical Review B</i> , 1995, 52, 11845-11852.	1.1	7
262	First-order static excitation potential: a Scheme for excitation energies and transition moments. <i>Physical Review A</i> , 1998, 57, 4311-4321.	1.0	7
263	Dynamical Green's function and an exact optical potential for electron-molecule scattering including nuclear dynamics. <i>Physical Review A</i> , 1999, 60, 2983-2999.	1.0	7
264	Structural and magnetic transitions in ensembles of mesoscopic Peierls rings in a magnetic flux. <i>Physical Review B</i> , 1999, 60, 6646-6654.	1.1	7
265	Vibronic Resonances Arising from Conically Intersecting Electronic States. <i>Journal of Physical Chemistry A</i> , 2002, 106, 4320-4335.	1.1	7
266	Resonances and pseudoresonances in a potential with attractive coulomb tail: A study using analytic-continuation techniques. <i>International Journal of Quantum Chemistry</i> , 2003, 94, 75-92.	1.0	7
267	Quantum dynamics through conical intersections in macrosystems: Combining effective modes and time-dependent Hartree. <i>Chemical Physics</i> , 2008, 347, 78-96.	0.9	7
268	Electron transfer mediated decay in HeLi ₂ cluster: Potential energy surfaces and decay widths. <i>Journal of Chemical Physics</i> , 2019, 150, 164309.	1.2	7
269	Tracing charge transfer in argon dimers by XUV-pump IR-probe experiments at FLASH. <i>Journal of Chemical Physics</i> , 2019, 151, 084314.	1.2	7
270	Quantum Effects Dominating the Interatomic Coulombic Decay of an Extreme System. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 6600-6605.	2.1	7

#	ARTICLE	IF	CITATIONS
271	Fano interferences in environment-enabled electron capture. <i>Physical Review A</i> , 2021, 103, .	1.0	7
272	Competition between interatomic Coulombic decay and autoionization of doubly-excited atoms. <i>Chemical Physics Letters</i> , 2020, 754, 137571.	1.2	7
273	Number fluctuations of cold, spatially split bosonic objects. <i>Physical Review A</i> , 2011, 84, .	1.0	6
274	Interatomic Coulombic electron capture in atomic, molecular, and quantum dot systems. <i>EPJ Web of Conferences</i> , 2015, 84, 07002.	0.1	6
275	The All-Seeing Eye of Resonant Auger Electron Spectroscopy: A Study on Aqueous Solution Using Tender X-rays. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 4457-4462.	2.1	6
276	Coupled-cluster theory for bosons in rings and optical lattices. <i>Computational and Theoretical Chemistry</i> , 2006, 768, 151-158.	1.5	5
277	Nonadditivity and anisotropy of the polarizability of clusters: Relativistic finite-field calculations for the Xe dimer. <i>Physical Review A</i> , 2010, 81, .	1.0	5
278	Efficient computation of adiabatic populations in multi-mode Jahn-Teller systems through the use of effective vibrational modes. <i>Journal of Chemical Physics</i> , 2011, 135, 174110.	1.2	5
279	Efficient computation of adiabatic electronic populations in multi-mode vibronic systems: Theory, implementation, and application. <i>Journal of Chemical Physics</i> , 2012, 137, 114110.	1.2	5
280	Tracing electron solvation in Li^+ in water. <i>Journal of Chemical Physics</i> , 2012, 137, 114110.	1.2	5
281	Time-resolved observation of interatomic excitation-energy transfer in argon dimers. <i>Journal of Chemical Physics</i> , 2017, 146, 104305.	1.2	5
282	Efficient non-resonant intermolecular vibrational energy transfer. <i>Molecular Physics</i> , 2019, 117, 1950-1955.	0.8	5
283	Electron spectroscopic study of nanoplasma formation triggered by intense soft x-ray pulses. <i>Journal of Chemical Physics</i> , 2019, 151, 184305.	1.2	5
284	Bound states and symmetry breaking of the ring $\text{C}_{20}^{\text{H}_8}$ anion. <i>Journal of Chemical Physics</i> , 2020, 152, 244307.	1.2	5
285	Electron attachment to a proton in water by interatomic Coulombic electron capture: An R-matrix study. <i>Physical Review A</i> , 2021, 104, .	1.0	5
286	MCTDHB Physics and Technologies: Excitations and Vorticity, Single-Shot Detection, Measurement of Fragmentation, and Optimal Control in Correlated Ultra-Cold Bosonic Many-Body Systems. , 2016, , 23-49.		5
287	High intensity x-ray interaction with a model bio-molecule system: double-core-hole states and fragmentation of formamide. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2020, 53, 244005.	0.6	5
288	Storing and releasing Mg by C12 carbon ring. <i>Chemical Physics Letters</i> , 2022, 799, 139554.	1.2	5

#	ARTICLE	IF	CITATIONS
289	Quantum States of Magnetically Induced Anions. <i>Physical Review Letters</i> , 2005, 95, 113002.	2.9	4
290	Kinetic energy release in fragmentation processes following electron emission: A time-dependent approach. <i>Journal of Chemical Physics</i> , 2012, 136, 114111.	1.2	4
291	Probing the interface of doped isotopically mixed helium droplets by the directional anisotropy of interatomic Coulombic decay. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 18167.	1.3	4
292	Polarization and site dependence of interatomic relaxation effects in double core hole states. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2013, 46, 164012.	0.6	4
293	Quantum oscillations between close-lying states mediated by the electronic continuum in intense high-frequency pulses. <i>Physical Review A</i> , 2015, 91, .	1.0	4
294	Many-Body Effects in Fragmented, Depleted, and Condensed Bosonic Systems in Traps and Optical Cavities by MCTDHB and MCTDH-X. , 2018, , 93-115.		4
295	Signature of the neighbor's quantum nuclear dynamics in the electron transfer mediated decay spectra. <i>Chemical Science</i> , 2021, 12, 9379-9385.	3.7	4
296	Theory of double ionization of a neighboring molecule by interatomic Coulombic decay. <i>Physical Review A</i> , 2021, 103, .	1.0	4
297	Numerically-Exact Schrödinger Dynamics of Closed and Open Many-Boson Systems with the MCTDHB Package. , 2013, , 81-92.		4
298	Quantum Many-Body Dynamics of Trapped Bosons with the MCTDHB Package: Towards New Horizons with Novel Physics. , 2015, , 63-86.		4
299	Extremely narrow peaks in predissociation of sodium dimer due to rovibronic coupling. <i>Journal of Chemical Physics</i> , 2004, 121, 3527-3532.	1.2	3
300	Field Operators in Real Space. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3009-3014.	1.1	3
301	Fragmentation of Molecules by Virtual Photons from Remote Neighbors. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 8964-8969.	2.1	3
302	Signatures of light-induced nonadiabaticity in the field-dressed vibronic spectrum of formaldehyde. <i>Journal of Chemical Physics</i> , 2021, 154, 124308.	1.2	3
303	Vorticity, Variance, and the Vigor of Many-Body Phenomena in Ultracold Quantum Systems: MCTDHB and MCTDH-X. , 2016, , 79-96.		3
304	Core-hole Hamiltonians and corrected equivalent core model for systems with equivalent atoms. <i>Journal of Chemical Physics</i> , 2003, 119, 12138-12152.	1.2	2
305	Weak pinning of the charge-density wave revisited. <i>Synthetic Metals</i> , 1997, 86, 2225-2226.	2.1	1
306	Strong impact of protonation and deprotonation on intermolecular Coulombic decay. <i>Journal of Physics: Conference Series</i> , 2012, 388, 022042.	0.3	1

#	ARTICLE	IF	CITATIONS
307	Endocircular Li Carbon Rings. <i>Angewandte Chemie</i> , 2021, 133, 16785-16790.	1.6	1
308	Ultrafast Charge Migration Following Ionization in Oligopeptides. <i>Springer Series in Chemical Physics</i> , 2009, , 586-588.	0.2	1
309	Effect of quantum phonon fluctuations on optical properties of finite semiconducting chains. Adiabatic and nonadiabatic results. <i>Synthetic Metals</i> , 1997, 85, 1101-1102.	2.1	0
310	Electron-phonon coupling in a one-band MX-chain model. A numerical study. <i>Synthetic Metals</i> , 1997, 86, 2221-2222.	2.1	0
311	Interplay between dia- and paramagnetism in ensembles of mesoscopic peierls rings in a magnetic flux. <i>Solid State Communications</i> , 1998, 108, 607-612.	0.9	0
312	Quantum phonon fluctuations in mesoscopic Peierls rings threaded by a magnetic flux. <i>Synthetic Metals</i> , 1999, 101, 345-346.	2.1	0
313	Structural change in mesoscopic Peierls chains. <i>Synthetic Metals</i> , 1999, 101, 394.	2.1	0
314	Electronic and lattice excitations in nonuniform one-dimensional clusters. <i>Synthetic Metals</i> , 1999, 102, 1581.	2.1	0
315	On the Endocircular Li@C16 System. <i>Frontiers in Chemistry</i> , 2022, 10, 813563.	1.8	0