

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

23879

60
h-index

42259

96
g-index

324
all docs

324
docs citations

324
times ranked

5831
citing authors

#	ARTICLE	IF	CITATIONS
1	On the Endocircular Li@C16 System. <i>Frontiers in Chemistry</i> , 2022, 10, 813563.	1.8	0
2	Storing and releasing Mg by C12 carbon ring. <i>Chemical Physics Letters</i> , 2022, 799, 139554.	1.2	5
3	Cooperative molecular structure in polaritonic and dark states. <i>Journal of Chemical Physics</i> , 2022, 156, 184102.	1.2	9
4	Bornâ€“Oppenheimer approximation in optical cavities: from success to breakdown. <i>Chemical Science</i> , 2021, 12, 1251-1258.	3.7	27
5	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
6	Caged-electron states and split-electron states in the endohedral alkali C₆₀. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 11837-11843.	1.3	10
7	Theory of double ionization of a neighboring molecule by interatomic Coulombic decay. <i>Physical Review A</i> , 2021, 103, .	1.0	4
8	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
9	Polaritonic States of Matter in a Rotating Cavity. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 6056-6061.	2.1	9
10	Endocircular Li Carbon Rings. <i>Angewandte Chemie</i> , 2021, 133, 16785-16790.	1.6	1
11	Endocircular Li Carbon Rings. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 16649-16654.	7.2	9
12	Impact of cavity on interatomic Coulombic decay. <i>Nature Communications</i> , 2021, 12, 4083.	5.8	18
13	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
14	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
15	Fano interferences in environment-enabled electron capture. <i>Physical Review A</i> , 2021, 103, .	1.0	7
16	Interatomic and Intermolecular Coulombic Decay. <i>Chemical Reviews</i> , 2020, 120, 11295-11369.	23.0	106
17	Fragmentation of Molecules by Virtual Photons from Remote Neighbors. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 8964-8969.	2.1	3
18	Bound states and symmetry breaking of the ring C20âˆ’ anion. <i>Journal of Chemical Physics</i> , 2020, 152, 244307.	1.2	5

#	ARTICLE	IF	CITATIONS
19	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
20	Ab initio complex potential energy curves of the He*(1s2p 1P) \leftrightarrow Li dimer. <i>Journal of Chemical Physics</i> , 2020, 152, 184303.	1.2	14
21	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
22	Competition between interatomic Coulombic decay and autoionization of doubly-excited atoms. <i>Chemical Physics Letters</i> , 2020, 754, 137571.	1.2	7
23	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
24	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
25	Core-level interatomic Coulombic decay in van der Waals clusters. <i>Physical Review Research</i> , 2020, 2, .	1.3	15
26	Efficient non-resonant intermolecular vibrational energy transfer. <i>Molecular Physics</i> , 2019, 117, 1950-1955.	0.8	5
27	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
28	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
29	Real-time observation of X-ray-induced intramolecular and interatomic electronic decay in CH ₂ I ₂ . <i>Nature Communications</i> , 2019, 10, 2186.	5.8	19
30	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
31	Charge separated states of endohedral fullerene Li@C ₂₀ . <i>Journal of Chemical Physics</i> , 2019, 151, 114306.	1.2	12
32	Caged-Electron States in Endohedral Li Fullerenes. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 7617-7622.	2.1	9
33	Many-Body Effects in Fragmented, Depleted, and Condensed Bosonic Systems in Traps and Optical Cavities by MCTDHB and MCTDH-X. , 2018, , 93-115.		4
34	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
35	Variance of an anisotropic Bose-Einstein condensate. <i>Chemical Physics</i> , 2018, 509, 45-54.	0.9	17
36	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

#	ARTICLE	IF	CITATIONS
37	Damaging Intermolecular Energy and Proton Transfer Processes in Alpha-Particle-Irradiated Hydrogen-Bonded Systems. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 17023-17027.	7.2	26
38	Damaging Intermolecular Energy and Proton Transfer Processes in Alpha-Particle-Irradiated Hydrogen-Bonded Systems. <i>Angewandte Chemie</i> , 2018, 130, 17269-17273.	1.6	19
39	Ultrafast Intermolecular Energy Transfer from Vibrations to Electronic Motion. <i>Physical Review Letters</i> , 2018, 121, 223001.	2.9	14
40	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
41	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
42	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
43	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
44	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
45	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
46	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
47	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
48	Interatomic Coulombic electron capture from first principles. <i>Physical Review A</i> , 2018, 98, .	1.0	17
49	Bound electronic states of the smallest fullerene C_{20}^{-} anion. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 17434-17441.	1.3	9
50	Observation of electron-transfer-mediated decay in aqueous solution. <i>Nature Chemistry</i> , 2017, 9, 708-714.	6.6	51
51	Phantom vortices: hidden angular momentum in ultracold dilute Bose-Einstein condensates. <i>Scientific Reports</i> , 2017, 7, 40122.	1.6	36
52	Intrinsic and light-induced nonadiabatic phenomena in the NaI molecule. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 19656-19664.	1.3	18
53	Competition between Light-Induced and Intrinsic Nonadiabatic Phenomena in Diatomics. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 1624-1630.	2.1	42
54	Time-resolved observation of interatomic excitation-energy transfer in argon dimers. <i>Journal of Chemical Physics</i> , 2017, 146, 104305.	1.2	5

#	ARTICLE	IF	CITATIONS
55	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
56	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
57	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
58	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
59	Electron-correlation driven capture and release in double quantum dots. <i>Journal of Physics Condensed Matter</i> , 2016, 28, 075301.	0.7	11
60	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
61	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
62	Tracking the photodissociation probability of D ₂ ⁺ induced by linearly chirped laser pulses. <i>Journal of Chemical Physics</i> , 2016, 144, 074309.	1.2	14
63	How many bound valence states does the C ₆₀ ⁻ anion have?. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 10840-10845.	1.3	12
64	Towards controlling the dissociation probability by light-induced conical intersections. <i>Faraday Discussions</i> , 2016, 194, 479-493.	1.6	28
65	Core Ionization Initiates Subfemtosecond Charge Migration in the Valence Shell of Molecules. <i>Physical Review Letters</i> , 2016, 117, 093002.	2.9	72
66	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
67	Field Operators in Real Space. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3009-3014.	1.1	3
68	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
69	Vorticity, Variance, and the Vigor of Many-Body Phenomena in Ultracold Quantum Systems: MCTDHB and MCTDH-X. , 2016, , 79-96.		3
70	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
71	Photodissociation of D ₂ ⁺ induced by linearly chirped laser pulses. <i>Journal of Chemical Physics</i> , 2015, 143, 014305.	1.2	15
72	Interatomic Coulombic electron capture in atomic, molecular, and quantum dot systems. <i>EPJ Web of Conferences</i> , 2015, 84, 07002.	0.1	6

#	ARTICLE	IF	CITATIONS
73	Barrierless Single-Electron-Induced <i>cis</i> → <i>trans</i> Isomerization. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 10470-10473.	7.2	15
74	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
75	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
76	Tracing electron solvation in Li^+ in water. <i>Journal of Chemical Physics</i> , 2014, 140, 034108.	1.2	13
77	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
78	Quantum Many-Body Dynamics of Trapped Bosons with the MCTDHB Package: Towards New Horizons with Novel Physics. , 2015, , 63-86.		4
79	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
80	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
81	Detecting ultrafast interatomic electronic processes in media by fluorescence. <i>New Journal of Physics</i> , 2014, 16, 102002.	1.2	19
82	Universality of fragmentation in the Schrödinger dynamics of bosonic Josephson junctions. <i>Physical Review A</i> , 2014, 89, .	1.0	44
83	Time-resolved pump-probe spectroscopy to follow valence electronic motion in molecules: Application. <i>Physical Review A</i> , 2014, 90, .	1.0	16
84	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
85	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
86	Breaking the resilience of a two-dimensional Bose-Einstein condensate to fragmentation. <i>Physical Review A</i> , 2014, 90, .	1.0	31
87	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
88	What will it take to observe processes in 'real time?'. <i>Nature Photonics</i> , 2014, 8, 162-166.	15.6	220
89	Site- and energy-selective slow-electron production through intermolecular Coulombic decay. <i>Nature</i> , 2014, 505, 661-663.	13.7	131
90	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

#	ARTICLE	IF	CITATIONS
91	All for one and one for all: accommodating an extra electron in C ₆₀ . Physical Chemistry Chemical Physics, 2014, 16, 13287.	1.3	30
92	Ultrafast correlation-driven electron dynamics. Journal of Physics B: Atomic, Molecular and Optical Physics, 2014, 47, 124002.	0.6	145
93	The exact wavefunction factorization of a vibronic coupling system. Journal of Chemical Physics, 2014, 140, 054104.	1.2	18
94	Elastic scattering of a Bose-Einstein condensate at a potential landscape. Journal of Physics: Conference Series, 2014, 488, 012032.	0.3	11
95	Nuclear-wave-packet quantum interference in the intense laser dissociation of the D ₂ molecule. Physical Review A, 2013, 88, .	1.0	43
96	ac Stark effect in the electronic continuum and its impact on the photoionization of atoms by coherent intense short high-frequency laser pulses. Physical Review A, 2013, 88, .	1.0	37
97	Excitation spectra of many-body systems by linear response: General theory and applications to trapped condensates. Physical Review A, 2013, 88, .	1.0	32
98	Extreme Correlation Effects in the Elusive Bound Spectrum of C ₆₀ ⁺ . Journal of Physical Chemistry Letters, 2013, 4, 3319-3324.	2.1	36
99	Photoionization of hydrogen atoms by coherent intense high-frequency short laser pulses: Direct propagation of electron wave packets on large spatial grids. Physical Review A, 2013, 88, .	1.0	21
100	Probing the interface of doped isotopically mixed helium droplets by the directional anisotropy of interatomic Coulombic decay. Physical Chemistry Chemical Physics, 2013, 15, 18167.	1.3	4
101	Electron-correlation-driven charge migration in oligopeptides. Chemical Physics, 2013, 414, 100-105.	0.9	59
102	The exact molecular wavefunction as a product of an electronic and a nuclear wavefunction. Journal of Chemical Physics, 2013, 138, 224110.	1.2	71
103	Existence of a Correlation Bound <i>s</i> -Type Anion State of C ₆₀ . Journal of Physical Chemistry Letters, 2013, 4, 849-853.	2.1	71
104	Effect of Light-Induced Conical Intersection on the Photodissociation Dynamics of the D ₂ ⁺ Molecule. Journal of Physical Chemistry A, 2013, 117, 8528-8535.	1.1	30
105	Light-induced conical intersections in polyatomic molecules: General theory, strategies of exploitation, and application. Journal of Chemical Physics, 2013, 139, 154314.	1.2	62
106	Controlled energy-selected electron capture and release in double quantum dots. Physical Review B, 2013, 88, .	1.1	32
107	Polarization and site dependence of interatomic relaxation effects in double core hole states. Journal of Physics B: Atomic, Molecular and Optical Physics, 2013, 46, 164012.	0.6	4
108	Two trapped particles interacting by a finite-range two-body potential in two spatial dimensions. Physical Review A, 2013, 87, .	1.0	39

#	ARTICLE	IF	CITATIONS
109	Time-resolved pump-probe spectroscopy to follow valence electronic motion in molecules: Theory. <i>Physical Review A</i> , 2013, 88, .	1.0	18
110	Numerically-Exact Schrödinger Dynamics of Closed and Open Many-Boson Systems with the MCTDHB Package. , 2013, , 81-92.		4
111	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
112	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
113	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
114	Interatomic relaxation effects in double core ionization of chain molecules. <i>Journal of Chemical Physics</i> , 2012, 137, 154316.	1.2	13
115	Dynamic Interference of Photoelectrons Produced by High-Frequency Laser Pulses. <i>Physical Review Letters</i> , 2012, 108, 253001.	2.9	92
116	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
117	Numerically exact quantum dynamics of bosons with time-dependent interactions of harmonic type. <i>Physical Review A</i> , 2012, 86, .	1.0	92
118	Dynamics and symmetries of a repulsively bound atom pair in an infinite optical lattice. <i>Physical Review A</i> , 2012, 86, .	1.0	14
119	Strong impact of protonation and deprotonation on intermolecular Coulombic decay. <i>Journal of Physics: Conference Series</i> , 2012, 388, 022042.	0.3	1
120	Wave chaos as signature for depletion of a Bose-Einstein condensate. <i>Physical Review A</i> , 2012, 86, .	1.0	46
121	Coherent intense resonant laser pulses lead to interference in the time domain seen in the spectrum of the emitted particles. <i>Physical Review A</i> , 2012, 86, .	1.0	43
122	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
123	Exploring Protonation and Deprotonation Effects with Auger Electron Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 2733-2737.	2.1	8
124	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
125	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
126	A One-Step Four-Bond-Breaking Reaction Catalyzed by an Electron. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 8003-8007.	7.2	48

#	ARTICLE	IF	CITATIONS
127	The effect of light-induced conical intersections on the alignment of diatomic molecules. <i>Chemical Physics</i> , 2012, 399, 146-150.	0.9	24
128	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
129	Ultrafast reorganization of the hole charge created upon outer-valence ionization of porphyrins. <i>Chemical Physics</i> , 2012, 399, 245-251.	0.9	15
130	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
131	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
132	An Excited Electron Avoiding a Positive Charge. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2300-2303.	2.1	19
133	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
134	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
135	Accurate Quantum Chemistry in Single Precision Arithmetic: Correlation Energy. <i>Journal of Chemical Theory and Computation</i> , 2011, 7, 320-326.	2.3	16
136	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
137	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
138	Dynamics of interatomic Coulombic decay in quantum dots. <i>Journal of Chemical Physics</i> , 2011, 135, 144112.	1.2	40
139	Using pH-Value To Control Intermolecular Electronic Decay. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 1306-1309.	7.2	27
140	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
141	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
142	Optimal time-dependent lattice models for nonequilibrium dynamics. <i>New Journal of Physics</i> , 2011, 13, 043003.	1.2	21
143	Ultrafast charge separation driven by differential particle and hole mobilities. <i>Journal of Chemical Physics</i> , 2011, 134, 024303.	1.2	23
144	Swift Loss of Coherence of Soliton Trains in Attractive Bose-Einstein Condensates. <i>Physical Review Letters</i> , 2011, 106, 240401.	2.9	39

#	ARTICLE	IF	CITATIONS
145	Exploring Interatomic Coulombic Decay by Free Electron Lasers. Physical Review Letters, 2011, 107, 273002.	2.9	28
146	Number fluctuations of cold, spatially split bosonic objects. Physical Review A, 2011, 84, .	1.0	6
147	Resonant Auger decay of the core-excited C $\langle\text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline">\langle\text{mml:msup}>\langle\text{mml:mrow}/>\langle\text{mml:mo}>*\langle\text{mml:mo}>\langle\text{mml:msup}>\langle\text{mml:math}>O$ molecule in intense x-ray laser fields. Physical Review A, 2011, 84, .	1.0	42
148	Accurate multi-boson long-time dynamics in triple-well periodic traps. Physical Review A, 2011, 83, .	1.0	45
149	Strong interference effects in the resonant Auger decay of atoms induced by intense x-ray fields. Physical Review A, 2011, 83, .	1.0	60
150	Interrelation between the Distributions of Kinetic Energy Release and Emitted Electron Energy following the Decay of Electronic States. Physical Review Letters, 2011, 107, 173001.	2.9	15
151	Anions of Xenon Clusters Bound by Long-Range Electron Correlations. Physical Review Letters, 2011, 107, 133401.	2.9	27
152	Resonant Auger Decay of Molecules in Intense X-Ray Laser Fields: Light-Induced Strong Nonadiabatic Effects. Physical Review Letters, 2011, 106, 123001.	2.9	63
153	Strong impact of light-induced conical intersections on the spectrum of diatomic molecules. Journal of Physics B: Atomic, Molecular and Optical Physics, 2011, 44, 045603.	0.6	82
154	Efficient computation of adiabatic populations in multi-mode Jahn-Teller systems through the use of effective vibrational modes. Journal of Chemical Physics, 2011, 135, 174110.	1.2	5
155	Interatomic Coulombic decay in a He dimer: <i>Ab initio</i> potential-energy curves and decay widths. Physical Review A, 2010, 82, .	1.0	19
156	Fragmented many-body states of definite angular momentum and stability of attractive three-dimensional condensates. Physical Review A, 2010, 82, .	1.0	15
157	Nonadditivity and anisotropy of the polarizability of clusters: Relativistic finite-field calculations for the Xe dimer. Physical Review A, 2010, 81, .	1.0	5
158	Impact of nuclear dynamics on interatomic Coulombic decay in a He dimer. Physical Review A, 2010, 82, .	1.0	24
159	Generation of Highly Damaging H ₂ O ⁺ Radicals by Inner Valence Shell Ionization of Water. ChemPhysChem, 2010, 11, 1006-1009.	1.0	36
160	Ultralong-range energy transfer by interatomic Coulombic decay in an extreme quantum system. Nature Physics, 2010, 6, 508-511.	6.5	133
161	Molecular double core hole electron spectroscopy for chemical analysis. Journal of Chemical Physics, 2010, 132, .	1.2	111
162	Quantum dynamics of attractive versus repulsive bosonic Josephson junctions: Bose-Hubbard and full-Hamiltonian results. Physical Review A, 2010, 82, .	1.0	52

#	ARTICLE	IF	CITATIONS
163	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
164	General mapping for bosonic and fermionic operators in Fock space. <i>Physical Review A</i> , 2010, 81, .	1.0	47
165	Interatomic Electronic Decay Driven by Nuclear Motion. <i>Physical Review Letters</i> , 2010, 105, 173401.	2.9	32
166	Ultrafast Interatomic Electronic Decay in Multiply Excited Clusters. <i>Physical Review Letters</i> , 2010, 105, 043004.	2.9	67
167	On the Cholesky decomposition for electron propagator methods: General aspects and application on C60. <i>Journal of Chemical Physics</i> , 2010, 132, 044110.	1.2	31
168	Correlation-bound anions of NaCl clusters. <i>Journal of Chemical Physics</i> , 2010, 133, 114301.	1.2	33
169	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
170	Ultrafast Charge Migration Following Valence Ionization of 4-Methylphenol: Jumping over the Aromatic Ring ^{>â€</sup>. <i>Journal of Physical Chemistry A</i>, 2010, 114, 8676-8679.}	1.1	49
171	Tracing molecular electronic excitation dynamics in real time and space. <i>Journal of Chemical Physics</i> , 2010, 132, 144302.	1.2	36
172	Exploring Nonadiabatic Effects by Recoil of Fast Photoelectrons. <i>Physical Review Letters</i> , 2009, 103, 133001.	2.9	8
173	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
174	Exact Quantum Dynamics of a Bosonic Josephson Junction. <i>Physical Review Letters</i> , 2009, 103, 220601.	2.9	163
175	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
176	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
177	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
178	Nuclear dynamics during the resonant Auger decay of water molecules. <i>Journal of Chemical Physics</i> , 2009, 130, 154307.	1.2	23
179	Build-up of coherence between initially-independent subsystems: The case of Boseâ€™Einstein condensates. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2009, 373, 301-304.	0.9	9
180	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

#	ARTICLE	IF	CITATIONS
181	Environment assisted electron capture. Journal of Physics B: Atomic, Molecular and Optical Physics, 2009, 42, 231001.	0.6	37
182	Efficient generation and properties of mesoscopic quantum superposition states in an attractive Bose-Einstein condensate threaded by a potential barrier. Journal of Physics B: Atomic, Molecular and Optical Physics, 2009, 42, 091004.	0.6	21
183	Ultrafast Charge Migration Following Ionization in Oligopeptides. Springer Series in Chemical Physics, 2009, , 586-588.	0.2	1
184	Ultrafast charge migration in 2-phenylethyl-N,N-dimethylamine. Chemical Physics Letters, 2008, 450, 232-235.	1.2	103
185	Quantum dynamics through conical intersections in macrosystems: Combining effective modes and time-dependent Hartree. Chemical Physics, 2008, 347, 78-96.	0.9	7
186	Reduced density matrices and coherence of trapped interacting bosons. Physical Review A, 2008, 78, .	1.0	124
187	Fragmented Metastable States Exist in an Attractive Bose-Einstein Condensate for Atom Numbers Well above the Critical Number of the Gross-Pitaevskii Theory. Physical Review Letters, 2008, 100, 040402.	2.9	19
188	Born-Oppenheimer approximation and beyond for time-dependent electronic processes. Journal of Chemical Physics, 2008, 128, 124101.	1.2	82
189	Charge migration following ionization in systems with chromophore-donor and amine-acceptor sites. Journal of Chemical Physics, 2008, 129, 104305.	1.2	72
190	Photodetachment spectra of the PtX ₄ ²⁻ (X=F,Cl,Br) dianions and their Jahn-Teller distortions: A fully relativistic study. Journal of Chemical Physics, 2008, 129, 174302.	1.2	11
191	<i>Ab initio</i> calculation of interatomic decay rates of excited doubly ionized states in clusters. Journal of Chemical Physics, 2008, 129, 244102.	1.2	39
192	On the doubly ionized states of Ar ₂ and their intra- and interatomic decay to Ar ₂ ³⁺ . Journal of Chemical Physics, 2008, 128, 014307.	1.2	30
193	On the interatomic electronic processes following Auger decay in neon dimer. Journal of Chemical Physics, 2008, 129, 074307.	1.2	41
194	Multiconfigurational time-dependent Hartree method for bosons: Many-body dynamics of bosonic systems. Physical Review A, 2008, 77, .	1.0	280
195	Formation and Dynamics of Many-Boson Fragmented States in One-Dimensional Attractive Ultracold Gases. Physical Review Letters, 2008, 100, 130401.	2.9	59
196	Laser-induced conical intersections in molecular optical lattices. Journal of Physics B: Atomic, Molecular and Optical Physics, 2008, 41, 221001.	0.6	95
197	PtF ₆ ²⁻ dianion and its detachment spectrum: A fully relativistic study. Journal of Chemical Physics, 2007, 126, 144310.	1.2	11
198	Quantum dynamics in macrosystems with several coupled electronic states: Hierarchy of effective Hamiltonians. Journal of Chemical Physics, 2007, 127, 124107.	1.2	22

#	ARTICLE	IF	CITATIONS
199	High activity of helium droplets following ionization of systems inside those droplets. <i>Physical Review B</i> , 2007, 76, .	1.1	29
200	Multiconfigurational time-dependent Hartree method for mixtures consisting of two types of identical particles. <i>Physical Review A</i> , 2007, 76, .	1.0	50
201	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
202	Role of Excited States in the Splitting of a Trapped Interacting Bose-Einstein Condensate by a Time-Dependent Barrier. <i>Physical Review Letters</i> , 2007, 99, 030402.	2.9	175
203	Tracing Ultrafast Interatomic Electronic Decay Processes in Real Time and Space. <i>Physical Review Letters</i> , 2007, 98, 083201.	2.9	89
204	Theory of magnetically induced anions. <i>Physical Review A</i> , 2007, 75, .	1.0	8
205	Unified view on multiconfigurational time propagation for systems consisting of identical particles. <i>Journal of Chemical Physics</i> , 2007, 127, 154103.	1.2	124
206	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
207	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
208	Charge migration in different conformers of glycine: The role of nuclear geometry. <i>Chemical Physics</i> , 2007, 338, 320-328.	0.9	98
209	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
210	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
211	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
212	Interatomic Electronic Decay in Endohedral Fullerenes. <i>Physical Review Letters</i> , 2006, 96, 053401.	2.9	85
213	Coupled-cluster theory for bosons in rings and optical lattices. <i>Computational and Theoretical Chemistry</i> , 2006, 768, 151-158.	1.5	5
214	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
215	Ionization spectra and electronic decay in small iodide clusters: Fully relativistic results. <i>Journal of Chemical Physics</i> , 2006, 125, 034309.	1.2	11
216	Ionization and double ionization of small water clusters. <i>Journal of Chemical Physics</i> , 2006, 125, 204305.	1.2	84

#	ARTICLE	IF	CITATIONS
217	Combined experimental-theoretical study of the lower excited singlet states of paraviny phenol, an analog of the paracoumaric acid chromophore. <i>Journal of Chemical Physics</i> , 2006, 125, 204303.	1.2	14
218	An effective Hamiltonian for the short-time dynamics at a conical intersection. <i>Molecular Physics</i> , 2006, 104, 1081-1093.	0.8	29
219	Demixing of Bosonic Mixtures in Optical Lattices from Macroscopic to Microscopic Scales. <i>Physical Review Letters</i> , 2006, 97, 230403.	2.9	37
220	Interatomic decay of inner-valence-excited states in clusters. <i>Journal of Chemical Physics</i> , 2006, 124, 144315.	1.2	47
221	Short-time dynamics through conical intersections in macrosystems. II. Applications. <i>Journal of Chemical Physics</i> , 2006, 124, 144104.	1.2	40
222	Interatomic Coulombic decay in a heteroatomic rare gas cluster. <i>Journal of Chemical Physics</i> , 2006, 124, 154305.	1.2	38
223	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
224	Coupled-cluster theory for systems of bosons in external traps. <i>Physical Review A</i> , 2006, 73, .	1.0	32
225	General variational many-body theory with complete self-consistency for trapped bosonic systems. <i>Physical Review A</i> , 2006, 73, .	1.0	119
226	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
227	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
228	Charge transfer in the Cl ⁺ CO cluster induced by core ionization. <i>Journal of Chemical Physics</i> , 2005, 122, 104304.	1.2	8
229	Electronic decay following ionization of aqueous Li ⁺ microsolvation clusters. <i>Journal of Chemical Physics</i> , 2005, 122, 094305.	1.2	44
230	Ab initio calculation of interatomic decay rates by a combination of the Fano ansatz, Green's-function methods, and the Stieltjes imaging technique. <i>Journal of Chemical Physics</i> , 2005, 123, 204107.	1.2	135
231	Simulation of a complex spectrum: Interplay of five electronic states and 21 vibrational degrees of freedom in C5H4 ⁺ . <i>Journal of Chemical Physics</i> , 2005, 123, 204310.	1.2	32
232	Exact ground state of finite Bose-Einstein condensates on a ring. <i>Physical Review A</i> , 2005, 72, .	1.0	64
233	Pathway from Condensation via Fragmentation to Fermionization of Cold Bosonic Systems. <i>Physical Review Letters</i> , 2005, 95, 140402.	2.9	86
234	Short-Time Dynamics Through Conical Intersections in Macrosystems. <i>Physical Review Letters</i> , 2005, 94, 113003.	2.9	140

#	ARTICLE	IF	CITATIONS
235	Properties of fragmented repulsive condensates. <i>Physical Review A</i> , 2005, 71, .	1.0	14
236	Charge transfer driven by electron correlation: A non-Dyson propagator approach. <i>Journal of Chemical Physics</i> , 2005, 122, 134104.	1.2	22
237	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
238	Zoo of Quantum Phases and Excitations of Cold Bosonic Atoms in Optical Lattices. <i>Physical Review Letters</i> , 2005, 95, 030405.	2.9	80
239	Quantum States of Magnetically Induced Anions. <i>Physical Review Letters</i> , 2005, 95, 113002.	2.9	4
240	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
241	Impact of Sulfur vs Oxygen on the Low-Lying Excited States of trans-p-Coumaric Acid and trans-p-Coumaric Thio Acid. <i>Journal of Physical Chemistry A</i> , 2005, 109, 4623-4631.	1.1	52
242	Stable and Long-Lived Trianions in the Gas Phase. <i>Journal of Physical Chemistry A</i> , 2005, 109, 11401-11406.	1.1	14
243	Multielectron wave-packet propagation: General theory and application. <i>Journal of Chemical Physics</i> , 2005, 123, 044111.	1.2	165
244	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
245	Extrapolating bound state data of anions into the metastable domain. <i>Journal of Chemical Physics</i> , 2004, 121, 6628-6633.	1.2	27
246	Ground-state fragmentation of repulsive Bose-Einstein condensates in double-trap potentials. <i>Physical Review A</i> , 2004, 70, .	1.0	52
247	Intersections of potential energy surfaces of short-lived states: The complex analogue of conical intersections. <i>Journal of Chemical Physics</i> , 2004, 120, 3201-3214.	1.2	50
248	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
249	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
250	Mechanism of Interatomic Coulombic Decay in Clusters. <i>Physical Review Letters</i> , 2004, 93, 263002.	2.9	166
251	THE MULTI-MODE VIBRONIC-COUPLING APPROACH. <i>Advanced Series in Physical Chemistry</i> , 2004, , 323-367.	1.5	86
252	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

#	ARTICLE	IF	CITATIONS
253	Potential energy surface of the CO ₂ ⁻ anion. <i>Physical Chemistry Chemical Physics</i> , 2004, 6, 42.	1.3	74
254	Environmental effects on a conical intersection: A model study. <i>Faraday Discussions</i> , 2004, 127, 395.	1.6	85
255	Ultrafast Photoinitiated Long-Range Electron Transfer in Cyclophane-Bridged Zincporphyrin ⁺ Quinone Complexes via Conical Intersections. <i>Journal of Physical Chemistry B</i> , 2004, 108, 19049-19055.	1.2	24
256	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
257	BEYOND BORN-OPPENHEIMER: Molecular Dynamics Through a Conical Intersection. <i>Annual Review of Physical Chemistry</i> , 2004, 55, 127-158.	4.8	694
258	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
259	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
260	Ionization of the xenon fluorides. <i>Journal of Chemical Physics</i> , 2003, 119, 7763-7771.	1.2	16
261	Electronic structure of isolated PtX ₆ ²⁻ (X=F,Cl,Br) dianions. <i>Journal of Chemical Physics</i> , 2003, 118, 1747-1755.	1.2	16
262	Coulombic Energy Transfer and Triple Ionization in Clusters. <i>Physical Review Letters</i> , 2003, 90, 153401.	2.9	98
263	Impact of interatomic electronic decay processes on Xe ⁺ 5d hole decay in the xenon fluorides. <i>Journal of Chemical Physics</i> , 2003, 119, 10575-10584.	1.2	45
264	Magnetically induced anions. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 4981.	1.3	11
265	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
266	Equivalent core model: Extended theory and applications. <i>Journal of Chemical Physics</i> , 2003, 118, 2081-2091.	1.2	9
267	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
268	Discovery of a new class of stable gas-phase dianions: Mixed oxygen-carbon cluster OC _n ²⁻ (n=5-19). <i>Journal of Chemical Physics</i> , 2002, 117, 7002-7009.	1.2	20
269	Systematic corrections to the equivalent core model. <i>Journal of Chemical Physics</i> , 2002, 116, 8723-8730.	1.2	10
270	Orbital Picture of Ionization and Its Breakdown in Nanoarrays of Quantum Dots. <i>Physical Review Letters</i> , 2002, 89, 133003.	2.9	13

#	ARTICLE	IF	CITATIONS
271	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
272	Vibronic Resonances Arising from Conically Intersecting Electronic States. <i>Journal of Physical Chemistry A</i> , 2002, 106, 4320-4335.	1.1	7
273	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
274	Dianionic Tetraborates Do Exist as Stable Entities. <i>Journal of the American Chemical Society</i> , 2002, 124, 10903-10910.	6.6	19
275	Cyclic Carbon Cluster Dianions and Their Aromaticity. <i>Journal of the American Chemical Society</i> , 2002, 124, 3163-3168.	6.6	12
276	Multiply Charged Anions in the Gas Phase. <i>Chemical Reviews</i> , 2002, 102, 181-200.	23.0	251
277	Non-Hermitian electronic theory and applications to clusters. <i>Physics Reports</i> , 2002, 368, 1-117.	10.3	261
278	Electronic decay of valence holes in clusters and condensed matter. <i>Physical Review B</i> , 2001, 64, .	1.1	148
279	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
280	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
281	Stability of Negatively Charged Ions Moving in a Magnetic Field. <i>Physical Review Letters</i> , 2001, 86, 5450-5453.	2.9	12
282	Electronic decay in weakly bound heteroclusters: Energy transfer versus electron transfer. <i>Journal of Chemical Physics</i> , 2001, 115, 5076-5088.	1.2	148
283	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
284	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
285	Inner-valence ionization of molecular anions and ultrafast relaxation by electron emission. <i>Chemical Physics Letters</i> , 2000, 324, 416-422.	1.2	10
286	Bound states of negatively charged ions induced by a magnetic field. <i>Physical Review A</i> , 2000, 61, .	1.0	25
287	Interatomic Coulombic Decay in van der Waals Clusters and Impact of Nuclear Motion. <i>Physical Review Letters</i> , 2000, 85, 4490-4493.	2.9	156
288	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
289	Suppression of electron correlation and of autoionization by strong laser fields. Journal of Physics B: Atomic, Molecular and Optical Physics, 1999, 32, L279-L284.	0.6	21
290	Transition from Rydberg to giant-dipole-moment states of hydrogen atoms in crossed fields: A suggestion for an experiment. Physical Review A, 1999, 59, 3695-3700.	1.0	12
291	Dynamical Green's function and an exact optical potential for electron-molecule scattering including nuclear dynamics. Physical Review A, 1999, 60, 2983-2999.	1.0	7
292	Structural and magnetic transitions in ensembles of mesoscopic Peierls rings in a magnetic flux. Physical Review B, 1999, 60, 6646-6654.	1.1	7
293	Quantum phonon fluctuations in mesoscopic Peierls rings threaded by a magnetic flux. Synthetic Metals, 1999, 101, 345-346.	2.1	0
294	Structural change in mesoscopic Peierls chains. Synthetic Metals, 1999, 101, 394.	2.1	0
295	Electronic and lattice excitations in nonuniform one-dimensional clusters. Synthetic Metals, 1999, 102, 1581.	2.1	0
296	How small can a Peierls dimerized chain be?. Solid State Communications, 1998, 106, 733-737.	0.9	9
297	Interplay between dia- and paramagnetism in ensembles of mesoscopic peierls rings in a magnetic flux. Solid State Communications, 1998, 108, 607-612.	0.9	0
298	First-order static excitation potential: a scheme for excitation energies and transition moments. Physical Review A, 1998, 57, 4311-4321.	1.0	7
299	Finite-size effects and quantum phonon fluctuations in the optical absorption edge of dimerized chains. Physical Review B, 1997, 55, 1481-1485.	1.1	16
300	Effect of quantum phonon fluctuations on optical properties of finite semiconducting chains. Adiabatic and nonadiabatic results. Synthetic Metals, 1997, 85, 1101-1102.	2.1	0
301	Electron-phonon coupling in a one-band MX-chain model. A numerical study. Synthetic Metals, 1997, 86, 2221-2222.	2.1	0
302	Weak pinning of the charge-density wave revisited. Synthetic Metals, 1997, 86, 2225-2226.	2.1	1
303	Stable Free Dianionic Silicon-Carbon Clusters. Angewandte Chemie International Edition in English, 1997, 36, 1889-1891.	4.4	26
304	Correlation effects in the valence x-ray photoionization spectra of ethylene, butadiene, and hexatriene. International Journal of Quantum Chemistry, 1997, 63, 465-481.	1.0	39
305	Extended Two-Particle Green's Functions and Optical Potentials for Two-Particle Scattering by Many-Body Targets. Annals of Physics, 1996, 252, 276-299.	1.0	12
306	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

#	ARTICLE	IF	CITATIONS
307	Formation of satellite bands in the ionization spectra of extended systems. <i>Physical Review B</i> , 1996, 53, 13326-13339.	1.1	34
308	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
309	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
310	Time-dependent photodissociation of methyl iodide with five active modes. <i>Journal of Chemical Physics</i> , 1994, 101, 5623-5646.	1.2	162
311	Foreign imaging in Auger spectroscopy: The Si 2p spectrum of silicon tetrafluoride. <i>Physical Review Letters</i> , 1993, 71, 649-652.	2.9	59
312	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
313	Methylboron Oxide, H ₃ C ₂ B ₂ O. <i>Angewandte Chemie International Edition in English</i> , 1989, 28, 88-90.	4.4	50
314	Diisocyanogen or Isocyanogen?. <i>Angewandte Chemie International Edition in English</i> , 1989, 28, 761-762.	4.4	15
315	Direct calculation of ionization potentials of closed-shell atoms and molecules. <i>Theoretica Chimica Acta</i> , 1973, 31, 239-260.	0.9	312