

Nimrod Moiseyev

List of Publications by Year in descending order

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131
papers

7,315
citations

117571

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133
times ranked

3841
citing authors

#	ARTICLE	IF	CITATIONS
1	Enhanced Coupling of Electron and Nuclear Spins by Quantum Tunneling Resonances. <i>Physical Review Letters</i> , 2022, 128, 013401.	2.9	3
2	Encircling exceptional points of Bloch waves: mode conversion and anomalous scattering. <i>Journal Physics D: Applied Physics</i> , 2022, 55, 235301.	1.3	6
3	Complex energies and transition dipoles for shape-type resonances of uracil anion from stabilization curves via Pad $\hat{\circ}$. <i>Journal of Chemical Physics</i> , 2022, 156, .	1.2	2
4	Transfer of information through waveguides near an exceptional point. <i>Physical Review A</i> , 2021, 103, .	1.0	10
5	Uniform vs Partial Scaling within Resonances via Pad $\hat{\circ}$ Based on the Similarities to Other Non-Hermitian Methods: Illustration for the Beryllium 1 <i>s</i> ² <i>p</i> ³ State. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 3435-3444.	2.3	4
6	Variational Solutions for Resonances by a Finite-Difference Grid Method. <i>Molecules</i> , 2021, 26, 5248.	1.7	0
7	Complex absorbing potentials for stark resonances. <i>International Journal of Quantum Chemistry</i> , 2020, 120, e26067.	1.0	2
8	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
9	Ab initio complex potential energy curves of the He*(1s2p 1P) $\hat{\circ}$ Li dimer. <i>Journal of Chemical Physics</i> , 2020, 152, 184303.	1.2	14
10	Evidence of Nonrigidity Effects in the Description of Low-Energy Anisotropic Molecular Collisions of Hydrogen Molecules with Excited Metastable Helium Atoms. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 2450-2459.	2.3	4
11	Laser Control of Resonance Tunneling via an Exceptional Point. <i>Physical Review Letters</i> , 2020, 124, 253202.	2.9	13
12	Ab Initio Complex Transition Dipoles between Autoionizing Resonance States from Real Stabilization Graphs. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 5601-5609.	2.1	7
13	Linking Scalar Elastodynamics and Non-Hermitian Quantum Mechanics. <i>Physical Review Applied</i> , 2020, 13, .	1.5	14
14	Ab-initio theory of photoionization via resonances. <i>Journal of Chemical Physics</i> , 2019, 150, 204111.	1.2	11
15	Quantum Effects in Cold Molecular Collisions from Spatial Polarization of Electronic Wave Function. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 855-863.	2.1	13
16	Quantum uncertainties and Heisenberg-like uncertainty relations for a weak measurement scheme involving two arbitrary noncommuting observables. <i>Physical Review A</i> , 2018, 97, .	1.0	2
17	Light Stops at Exceptional Points. <i>Physical Review Letters</i> , 2018, 120, 013901.	2.9	138
18	Polarization dependence of the propagation constant of leaky guided modes. <i>Physical Review A</i> , 2018, 97, .	1.0	7

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19	Simple Closed-Form Expression for Penning Reaction Rate Coefficients for Cold Molecular Collisions by Non-Hermitian Time-Independent Adiabatic Scattering Theory. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 236-241.	2.3	8
20	Exceptional points in the Riesz-Feller Hamiltonian with an impenetrable rectangular potential. <i>Physical Review A</i> , 2018, 98, .	1.0	3
21	The effect of large autoionization decay rates (resonance widths) on cold molecular cross-sections and the reflection phenomenon. <i>Chemical Physics</i> , 2018, 515, 88-93.	0.9	1
22	Forces on nuclei moving on autoionizing molecular potential energy surfaces. <i>Journal of Chemical Physics</i> , 2017, 146, 024101.	1.2	4
23	Ab Initio Complex Potential Energy Surfaces From Standard Quantum Chemistry Packages. <i>Advances in Quantum Chemistry</i> , 2017, 74, 321-346.	0.4	11
24	Adiabatic Variational Theory for Cold Atom-Molecule Collisions: Application to a Metastable Helium Atom Colliding with <i>ortho</i> - and <i>para</i> -Hydrogen Molecules. <i>Journal of Physical Chemistry A</i> , 2017, 121, 2194-2198.	1.1	15
25	Polyatomic <i>ab Initio</i> Complex Potential Energy Surfaces: Illustration of Ultracold Collisions. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 1682-1690.	2.3	25
26	The boomerang effect in electron-hydrogen molecule scattering as determined by time-dependent calculations. <i>Journal of Chemical Physics</i> , 2017, 146, 204303.	1.2	6
27	On the calculation of resonances by analytic continuation of eigenvalues from the stabilization graph. <i>Journal of Chemical Physics</i> , 2017, 147, 014101.	1.2	19
28	Adiabatic perturbation theory for atoms and molecules in the low-frequency regime. <i>Journal of Chemical Physics</i> , 2017, 147, 224101.	1.2	4
29	Directly probing anisotropy in atom-molecule collisions through quantum scattering resonances. <i>Nature Physics</i> , 2017, 13, 35-38.	6.5	99
30	Molecular resonances by removing complex absorbing potentials via Pad \hat{C} ; Application to CO \hat{v} and N $\hat{2}\hat{v}$. <i>Journal of Chemical Physics</i> , 2016, 145, 164111.	1.2	19
31	On the Equivalence of Different Methods for Calculating Resonances: From Complex Gaussian Basis Set to Reflection-Free Complex Absorbing Potentials via the Smooth Exterior Scaling Transformation. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 2542-2552.	2.3	13
32	Dynamically encircling an exceptional point for asymmetric mode switching. <i>Nature</i> , 2016, 537, 76-79.	13.7	684
33	Atomic and Molecular Complex Resonances from Real Eigenvalues Using Standard (Hermitian) Electronic Structure Calculations. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3098-3108.	1.1	37
34	Light-induced conical intersection effect enhancing the localization of molecules in optical lattices. <i>Physical Review A</i> , 2015, 92, .	1.0	15
35	Characteristic footprints of an exceptional point in the dynamics of Li dimer under a laser field. <i>Journal of Chemical Physics</i> , 2015, 143, 154308.	1.2	5
36	Perturbation theory for quasienergy Floquet solutions in the low-frequency regime of the oscillating electric field. <i>Physical Review A</i> , 2015, 91, .	1.0	13

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37	Advantages of complex scaling only the most diffuse basis functions in simultaneous description of both resonances and bound states. <i>Molecular Physics</i> , 2015, 113, 3141-3146.	0.8	10
38	Adiabatic theory for anisotropic cold molecule collisions. <i>Journal of Chemical Physics</i> , 2015, 143, 074114.	1.2	17
39	Helium in chirped laser fields as a time-asymmetric atomic switch. <i>Journal of Chemical Physics</i> , 2014, 141, 014307.	1.2	19
40	Entanglement and Spin Squeezing in Non-Hermitian Phase Transitions. <i>Physical Review Letters</i> , 2014, 113, 250401.	2.9	116
41	Conditions for the applicability of the Kramers-Henneberger approximation for atoms in high-frequency strong laser fields. <i>Physical Review A</i> , 2014, 90, .	1.0	13
42	Time-asymmetric quantum-state-exchange mechanism. <i>Physical Review A</i> , 2013, 88, .	1.0	93
43	Chemistry in high-frequency strong laser fields: the story of HeS molecule. <i>Molecular Physics</i> , 2013, 111, 1814-1822.	0.8	5
44	Breakdown of adiabatic transfer of light in waveguides in the presence of absorption. <i>Physical Review A</i> , 2013, 88, .	1.0	52
45	Asymmetric effect of slowly varying chirped laser pulses on the adiabatic state exchange of a molecule. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2012, 45, 051002.	0.6	34
46	Scattering from a waveguide by cycling a non-Hermitian degeneracy. <i>Physical Review A</i> , 2012, 85, .	1.0	41
47	Distinguishing between aligned and randomly oriented polar molecules by using a combination of strong laser field with a weak static field. <i>Molecular Physics</i> , 2012, 110, 1721-1728.	0.8	2
48	Resonance energies, lifetimes and complex energy potential curves from standard wave-packet calculations. <i>Molecular Physics</i> , 2012, 110, 537-546.	0.8	4
49	Ab-initio complex molecular potential energy surfaces by the back-rotation transformation method. <i>Chemical Physics Letters</i> , 2012, 524, 84-89.	1.2	12
50	On the observability and asymmetry of adiabatic state flips generated by exceptional points. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2011, 44, 435302.	0.7	170
51	Fingerprints of exceptional points in the survival probability of resonances in atomic spectra. <i>Physical Review A</i> , 2011, 84, .	1.0	30
52	Diverging Rabi Oscillations in Subwavelength Photonic Lattices. <i>Physical Review Letters</i> , 2011, 106, 073901.	2.9	21
53	Electron relaxation in quantum dots by the interatomic Coulombic decay mechanism. <i>Physical Review B</i> , 2011, 83, .	1.1	35
54	Evaluation of partial widths and branching ratios from resonance wave functions. <i>Physical Review A</i> , 2010, 82, .	1.0	13

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55	The absolute position of a resonance peak. Journal of Physics B: Atomic, Molecular and Optical Physics, 2010, 43, 185205.	0.6	25
56	Spanning the Hilbert space with an even tempered Gaussian basis set. International Journal of Quantum Chemistry, 2009, 109, 2996-3002.	1.0	24
57	Feshbach Resonances: The Branching of Quantum Mechanics into Hermitian and Non-Hermitian Formalisms. Journal of Physical Chemistry A, 2009, 113, 7660-7666.	1.1	4
58	Suppression of Feshbach Resonance Widths in Two-Dimensional Waveguides and Quantum Dots: A Lower Bound for the Number of Bound States in the Continuum. Physical Review Letters, 2009, 102, 167404.	2.9	64
59	Visualization of Branch Points in $\langle \text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"} \langle \text{mml:mi mathvariant="script"} \rangle P \langle \text{mml:mi mathvariant="script"} \rangle T \langle \text{mml:mi} \rangle \langle \text{mml:math} \rangle$ -Symmetric Waveguides. Physical Review Letters, 2008, 101, 080402.	2.9	793
60	Reflection-free complex absorbing potential for electronic structure calculations: Feshbach-type autoionization resonances of molecules. Journal of Chemical Physics, 2007, 127, 034105.	1.2	36
61	Attosecond laser pulse synthesis using bichromatic high-order harmonic generation. Physical Review A, 2006, 74, .	1.0	33
62	Dipole and quadrupole forces exerted on atoms in laser fields: The nonperturbative approach. Physical Review A, 2006, 74, .	1.0	11
63	Ab initio calculation of harmonic generation spectra of helium using a time-dependent non-Hermitian formalism. Physical Review A, 2006, 74, .	1.0	15
64	Resonance solutions of the nonlinear Schrödinger equation: Tunneling lifetime and fragmentation of trapped condensates. Physical Review A, 2005, 72, .	1.0	40
65	Dynamical symmetry analysis of ionization and harmonic generation of atoms in bichromatic laser pulses. International Journal of Quantum Chemistry, 2005, 103, 824-840.	1.0	4
66	Adiabatic theorem for non-Hermitian time-dependent open systems. Physical Review A, 2005, 72, .	1.0	64
67	Resonance positions and lifetimes for flexible complex absorbing potentials. Physical Review A, 2005, 72, .	1.0	35
68	Non-Hermitian quantum mechanics versus the conventional quantum mechanics: Effect of the relative phasing of bichromatic fields on high-order harmonic generation. Physical Review A, 2004, 69, .	1.0	6
69	The Resonance Phenomena Associated with the Time Asymmetry in Non-Hermitian Quantum Mechanics. International Journal of Theoretical Physics, 2003, 42, 2131-2143.	0.5	0
70	On the collapse and restoration of condensates in dimensions in the mean-field approximation. Israel Journal of Chemistry, 2003, 43, 267-277.	1.0	0
71	High harmonic generation spectra of aligned benzene in circular polarized laser field. Journal of Chemical Physics, 2003, 118, 8726-8738.	1.2	31
72	Ionization and high-order harmonic generation in aligned benzene by a short intense circularly polarized laser pulse. Physical Review A, 2003, 68, .	1.0	65

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73	Stability and instability of dipole selection rules for atomic high-order-harmonic-generation spectra in two-beam setups. <i>Physical Review A</i> , 2002, 65, .	1.0	11
74	High-order harmonic generation by molecules of discrete rotational symmetry interacting with circularly polarized laser field. <i>Physical Review A</i> , 2001, 64, .	1.0	64
75	Phases and amplitudes of recurrences in autocorrelation function by a simple classical trajectory method. <i>Journal of Chemical Physics</i> , 2001, 115, 10608-10620.	1.2	7
76	Crossover phenomena and resonances in quantum systems. <i>Physical Review A</i> , 2001, 64, .	1.0	22
77	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
78	Non-Hermitian formulation of interference effect in scattering experiments. <i>Journal of Chemical Physics</i> , 2000, 113, 6088-6095.	1.2	23
79	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
80	Trapping of an Electron due to Molecular Vibrations. <i>Physical Review Letters</i> , 2000, 84, 1681-1684.	2.9	70
81	High Harmonic Generation of Soft X-Rays by Carbon Nanotubes. <i>Physical Review Letters</i> , 2000, 85, 5218-5221.	2.9	75
82	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
83	Scattering matrix determination by asymptotic analysis of complex scaled resonance wave functions: Model Cl+H2 nonadiabatic dynamics. <i>Journal of Chemical Physics</i> , 1999, 111, 7187-7196.	1.2	7
84	Crossed-beam experiment: High-order harmonic generation and dynamical symmetry. <i>Physical Review A</i> , 1999, 60, 2585-2586.	1.0	18
85	Quantum theory of resonances: calculating energies, widths and cross-sections by complex scaling. <i>Physics Reports</i> , 1998, 302, 212-293.	10.3	882
86	Derivations of universal exact complex absorption potentials by the generalized complex coordinate method. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 1998, 31, 1431-1441.	0.6	132
87	Selection Rules for the High Harmonic Generation Spectra. <i>Physical Review Letters</i> , 1998, 80, 3743-3746.	2.9	233
88	Classical versus quantum harmonic-generation spectrum of a driven anharmonic oscillator in the high-frequency regime. <i>Physical Review A</i> , 1998, 57, 1345-1354.	1.0	9
89	Fingerprints of Broad Overlapping Resonances in the H ₂ Cross Section. <i>Physical Review Letters</i> , 1998, 81, 2221-2224.	2.9	38
90	Selective quasienergies from short time cross-correlation probability amplitudes by the filter-diagonalization method. <i>Physical Review E</i> , 1998, 58, 376-381.	0.8	6

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91	Photoabsorption probability for a system governed by a time-dependent Hamiltonian through the $(t, t\hat{\epsilon}^2)$ formalism. <i>Journal of Chemical Physics</i> , 1997, 106, 6839-6847.	1.2	7
92	Localization of multiphoton ionization/dissociation resonance wave functions in AC fields. <i>International Journal of Quantum Chemistry</i> , 1997, 63, 279-285.	1.0	4
93	Complex scaling of abinitiomolecular potential surfaces. <i>Journal of Chemical Physics</i> , 1996, 104, 6192-6195.	1.2	8
94	Transition state resonances by complex scaling: A three-dimensional study of ClHCl. <i>Journal of Chemical Physics</i> , 1995, 103, 8468-8476.	1.2	39
95	The solution of the time-dependent Schrödinger equation by the $(t, t\hat{\epsilon}^M)$ method: Complex scaled multiphoton ionization/dissociation resonance wave functions are square integrable. <i>Journal of Chemical Physics</i> , 1994, 101, 9716-9718.	1.2	17
96	The solution of the time-dependent Schrödinger equation by the $(t, t\hat{\epsilon}^M)$ method: Multiphoton ionization/dissociation probabilities in different gauges of the electromagnetic potentials. <i>Journal of Chemical Physics</i> , 1994, 100, 7310-7318.	1.2	36
97	The complex coordinate scattering theory and its application to the study of the surface asymmetry effect in helium diffraction from copper. <i>International Journal of Quantum Chemistry</i> , 1993, 46, 343-363.	1.0	5
98	Cumulative reaction probability from Siegert eigenvalues: Model studies. <i>Journal of Chemical Physics</i> , 1993, 98, 9618-9623.	1.2	25
99	The solution of the time-dependent Schrödinger equation by the $(t, t\hat{\epsilon}^M)$ method: Theory, computational algorithm and applications. <i>Journal of Chemical Physics</i> , 1993, 99, 4590-4596.	1.2	220
100	Absorbing boundary conditions by the partial integration exterior scaling method. <i>Journal of Chemical Physics</i> , 1993, 99, 7703-7708.	1.2	9
101	Cumulative reaction probability by the complex coordinate scattering theory. <i>Journal of Chemical Physics</i> , 1993, 98, 6327-6331.	1.2	7
102	Cumulative reaction probabilities using Padé analytical continuation procedures. <i>Journal of Chemical Physics</i> , 1993, 99, 3509-3515.	1.2	7
103	Quantum mechanical thermal rate constants using flux-flux correlation functions and Padé analytical continuation procedures. <i>Journal of Chemical Physics</i> , 1993, 98, 8601-8605.	1.2	10
104	Gas/surface complex coordinate scattering theory: HD/Ag(111), HD/Pt(111) rotationally inelastic transition intensities. <i>Journal of Chemical Physics</i> , 1992, 96, 2347-2355.	1.2	20
105	The complex coordinate scattering theory and the Kohn variational method: A general formulation and application to long range potentials. <i>Journal of Chemical Physics</i> , 1992, 97, 6443-6450.	1.2	18
106	The complex coordinate scattering theory: Broken inversion symmetry of corrugated surfaces in helium diffraction from Cu(115). <i>Journal of Chemical Physics</i> , 1992, 97, 2804-2808.	1.2	11
107	Foreword by the Guest Editor of this Issue. <i>Israel Journal of Chemistry</i> , 1991, 31, 273-273.	1.0	0
108	Resonances, Cross Sections, and Partial Widths by the Complex Coordinate Method. <i>Israel Journal of Chemistry</i> , 1991, 31, 311-322.	1.0	38

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109	A theory of He diffraction and resonance scattering from Cu(115) by the complex coordinate method. Journal of Chemical Physics, 1991, 94, 1636-1642.	1.2	21
110	Gas surface scattering cross section by the complex coordinate method. Journal of Chemical Physics, 1991, 94, 6330-6333.	1.2	18
111	Resonance Positions and Widths for Time-Periodic Hamiltonians by the Complex Coordinate Method. Israel Journal of Chemistry, 1990, 30, 107-114.	1.0	14
112	Partial widths obtained by the complex resonance-scattering theory. Physical Review A, 1990, 42, 255-260.	1.0	45
113	Tunneling rates in bound systems using smooth exterior complex scaling within the framework of the finite basis set approximation. Journal of Chemical Physics, 1990, 93, 3413-3419.	1.2	60
114	Perturbation analysis of gas-surface diffractive selective adsorption resonance states. Molecular Physics, 1989, 66, 465-478.	0.8	5
115	Resonance transition probabilities by the complex Lanczos recursion method. Journal of Chemical Physics, 1988, 89, 6836-6840.	1.2	16
116	Complex quasiprobability for atoms trapped on surfaces: A novel application of the complex coordinate method. Journal of Chemical Physics, 1988, 88, 5864-5870.	1.2	10
117	Highly excited vibrational states by adiabatic vs self-consistent field methods. Journal of Chemical Physics, 1987, 86, 2146-2151.	1.2	14
118	Application of the complex rotation method to the study of resonance states of atoms at a corrugated surface. Journal of Chemical Physics, 1987, 86, 1048-1054.	1.2	23
119	Resonances from the complex dilated Hamiltonians in a dilation-adapted basis set with a new stabilization parameter. Journal of Chemical Physics, 1986, 84, 3931-3936.	1.2	13
120	On the "New possibility of chemical bonding": Anti-resonance phenomena. Chemical Physics Letters, 1984, 106, 354-355.	1.2	3
121	Motion of wave packets in regular and chaotic systems. Journal of Chemical Physics, 1983, 79, 5945-5950.	1.2	37
122	Resonance states by the generalized complex variational method. Molecular Physics, 1982, 47, 585-598.	0.8	50
123	Studies of multi-channel resonances by the complex scaling method. Molecular Physics, 1981, 42, 129-139.	0.8	18
124	Study of predissociation resonances by the complex coordinate method. International Journal of Quantum Chemistry, 1981, 20, 835-842.	1.0	18
125	Cusps, \hat{I} trajectories, and the complex virial theorem. Journal of Chemical Physics, 1981, 74, 4739-4740.	1.2	99
126	Criteria of accuracy of resonance eigenvalues. International Journal of Quantum Chemistry, 1980, 17, 1201-1211.	1.0	26

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127	Association of resonance states with the incomplete spectrum of finite complex-scaled Hamiltonian matrices. <i>Physical Review A</i> , 1980, 22, 618-624.	1.0	101
128	Autoionizing states of H_2 and H_2^+ using the complex-scaling method. <i>Physical Review A</i> , 1979, 20, 814-817.	1.0	174
129	Fermi and Coulomb correlations in the 21 S state of the helium isoelectronic sequence. <i>Theoretica Chimica Acta</i> , 1977, 45, 61-67.	0.9	17
130	The Gaussian potential: Bound states in the continuum?. <i>Theoretica Chimica Acta</i> , 1976, 41, 321-328.	0.9	6
131	The RVP Method – From Real Ab-Initio Calculations to Complex Energies and Transition Dipoles. <i>Frontiers in Physics</i> , 0, 10, .	1.0	2