

# Karel Houfek

## List of Publications by Year in descending order

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34  
papers

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citations

840776

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839539

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34  
all docs

34  
docs citations

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times ranked

295  
citing authors

#	ARTICLE	IF	CITATIONS
1	Vibronal excitation in the $\text{HCO}^+$ system: Nonlocal model of $\text{HCO}^+$ vibrational excitation in the $\text{HCO}^+$ system. Physical Review Letters, 2022, 105, .	2.5	7
2	Vibronic Coupling through the Continuum in the $\text{HCO}^+$ System. Physical Review Letters, 2022, 129, .	7.8	11
3	Electron-impact vibrational excitation of isocyanic acid HNCO. Physical Review A, 2020, 102, .	2.5	1
4	Electron-impact vibrational excitation of isocyanic acid HNCO. Physical Review A, 2020, 102, .	2.5	7
5	Associative detachment in $\text{Li}+\text{H}\hat{\text{a}}^+$ collisions. European Physical Journal D, 2018, 72, 1.	1.3	3
6	Dissociative recombination by frame transformation to Siegert pseudostates: A comparison with a numerically solvable model. Physical Review A, 2018, 97, .	2.5	5
7	Converged and consistent high-resolution low-energy electron- $\text{H}^+$ hydrogen scattering. I. Data below $10^{-4}$ eV. Atomic Data and Nuclear Data Tables, 2018, 119, 303-313.	2.4	4
8	Resonances and Dissociative Electron Attachment in HNCO. Physical Review Letters, 2018, 121, 143402.	7.8	25
9	Time-dependent formulation of the two-dimensional model of resonant electron collisions with diatomic molecules and interpretation of the vibrational excitation cross sections. Physical Review A, 2017, 95, .	2.5	1
10	Role of electronic correlations in photoionization of $\text{NO}^+$ in the vicinity of the $\text{A}^1\Sigma^+$ conical intersection. Physical Chemistry Chemical Physics, 2017, 19, 19673-19682.	2.8	9
11	Reducing the dimensionality of grid based methods for electron-atom scattering calculations below ionization threshold. Computer Physics Communications, 2017, 213, 46-51.	7.5	1
12	Isotopic effects in the interaction of $\text{H}^+$ with $\text{D}_2$ and $\text{H}_2$ .	2.5	6
13	New version of hex-ecs, the B-spline implementation of exterior complex scaling method for solution of electron- $\text{H}^+$ hydrogen scattering. Computer Physics Communications, 2016, 204, 216-217.	7.5	2
14	Lowest autodetachment state of the water anion. European Physical Journal D, 2016, 70, 1.	1.3	4
15	Recent developments in R-matrix applications to molecular processes. Journal of Physics: Conference Series, 2015, 635, 072058.	0.4	0
16	Interaction of $\text{O}\hat{\text{a}}^+$ and $\text{H}_2$ at low temperatures. Journal of Chemical Physics, 2015, 142, 014304.	3.0	12
17	Collisions of electrons with hydrogen atoms II. Low-energy program using the method of the exterior complex scaling. Computer Physics Communications, 2014, 185, 2903-2912.	7.5	6
18	Collisions of electrons with hydrogen atoms I. Package outline and high energy code. Computer Physics Communications, 2014, 185, 2893-2902.	7.5	5

#	ARTICLE	IF	CITATIONS
19	Resonant inelastic collisions of electrons with diatomic molecules. Nuclear Instruments & Methods in Physics Research B, 2012, 279, 71-75.	1.4	1
20	Dissociative electron attachment and vibrational excitation of CF <sub>3</sub> Cl: Effect of two vibrational modes revisited. Physical Review A, 2011, 84, .	2.5	14
21	Electron scattering in HCl: An improved nonlocal resonance model. Physical Review A, 2010, 81, .	2.5	16
22	Comparison of the Chebyshev Method and the Generalized Crank-Nicholson Method for time Propagation in Quantum Mechanics. , 2010, , .		2
23	Efficient Numerical Solution of Time-Dependent Multichannel One-Dimensional or Radial Problems In Quantum Mechanics. , 2009, , .		0
24	On irregular oscillatory structures in resonant vibrational excitation cross-sections in diatomic molecules. Chemical Physics, 2008, 347, 250-256.	1.9	11
25	Probing the nonlocal approximation to resonant collisions of electrons with diatomic molecules. Physical Review A, 2008, 77, .	2.5	19
26	Efficient Numerical Solution of Coupled Radial Differential Equations in Multichannel Scattering Problems. , 2008, , .		0
27	Giant structures in low-energy electron deuterium-iodide elastic scattering cross section. Physical Review A, 2007, 75, .	2.5	1
28	Long-lived states of molecular hydrogen anion. AIP Conference Proceedings, 2007, , .	0.4	0
29	Numerically solvable model for resonant collisions of electrons with diatomic molecules. Physical Review A, 2006, 73, .	2.5	17
30	Dissociative electron attachment and vibrational excitation of H <sub>2</sub> by low-energy electrons: Calculations based on an improved nonlocal resonance model. II. Vibrational excitation. Physical Review A, 2006, 73, .	2.5	37
31	Nonlocal model of dissociative electron attachment and vibrational excitation of NO. Physical Review A, 2005, 71, .	2.5	43
32	Dissociative electron attachment and vibrational excitation of H <sub>2</sub> by low-energy electrons: Calculations based on an improved nonlocal resonance model. Physical Review A, 2004, 70, .	2.5	55
33	Calculation of rate constants for dissociative attachment of low-energy electrons to hydrogen halides HCl, HBr, and HI and their deuterated analogs. Physical Review A, 2002, 66, .	2.5	14
34	Dissociative attachment of low-energy electrons to vibrationally excited hydrogen molecules. European Physical Journal D, 2002, 52, 29-40.	0.4	4