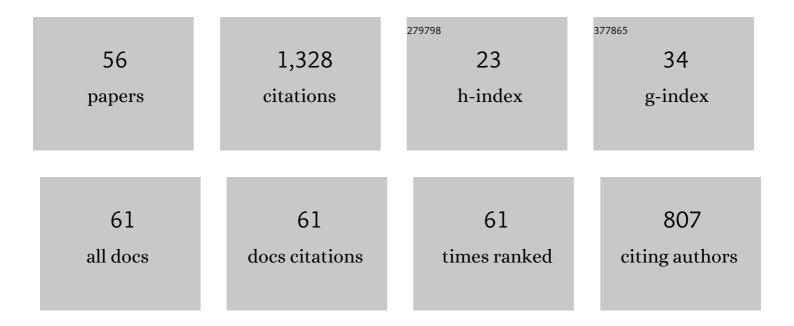
Raymond Bemish

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	The Ar–C2H2 intermolecular potential from high resolution spectroscopy and ab initio theory: A case for multicenter interactions. Journal of Chemical Physics, 1993, 99, 8585-8598.	3.0	92
2	Photodissociation of molecules oriented by dc electric fields: Determining photofragment angular distributions. Journal of Chemical Physics, 1994, 101, 9447-9456.	3.0	71
3	Electron Impact Ionization in Helium Nanodroplets:Â Controlling Fragmentation by Active Cooling of Molecular Ions. Journal of the American Chemical Society, 2004, 126, 11283-11292.	13.7	65
4	Probing Charge-Transfer Processes in Helium Nanodroplets by Optically Selected Mass Spectrometry (OSMS):Â Charge Steering by Long-Range Interactions. Journal of the American Chemical Society, 2005, 127, 7235-7242.	13.7	65
5	Photofragment vibrational, rotational, and translational distributions for N2–HF (v=1). Journal of Chemical Physics, 1994, 101, 9457-9468.	3.0	61
6	The C(3P) + NO(X2Î) → O(3P) + CN(X2Σ+), N(2D)/N(4S) + CO(X1Σ+) reaction: Rates, branching ratios, and final states from 15 K to 20 000 K. Journal of Chemical Physics, 2018, 149, 094305.	3.0	43
7	Infrared spectroscopy and ab initio potential energy surface for Ne–C2H2 and Ne–C2HD complexes. Journal of Chemical Physics, 1998, 109, 8968-8979.	3.0	41
8	Aqueous phosphoric acid as a mild reagent for deprotection of the t-butoxycarbonyl group. Tetrahedron Letters, 2003, 44, 8113-8115.	1.4	41
9	Adsorption of Acridine Orange at a C8,18/Water/Acetonitrile Interface. Journal of Physical Chemistry B, 2007, 111, 10208-10216.	2.6	41
10	Solvent structures of mixed water/acetonitrile mixtures at chromatographic interfaces from computer simulations. Physical Chemistry Chemical Physics, 2008, 10, 4765.	2.8	40
11	Molecular control using dc electric fields: quenching of the tunneling in HF dimer. Chemical Physics Letters, 1996, 251, 182-188.	2.6	38
12	The N2H+–He intermolecular potential energy surface: A vibrational adiabatic correction. Journal of Chemical Physics, 1997, 106, 8672-8680.	3.0	36
13	Exhaustive state-to-state cross sections for reactive molecular collisions from importance sampling simulation and a neural network representation. Journal of Chemical Physics, 2019, 150, 211101.	3.0	35
14	Computational study of collisions between O(3P) and NO(2Î) at temperatures relevant to the hypersonic flight regime. Journal of Chemical Physics, 2014, 141, 164319.	3.0	34
15	Abinitioand scaled potential energy surfaces for Ar–C2H2: Comparison with scattering and spectroscopic experiments. Journal of Chemical Physics, 1996, 105, 10462-10471.	3.0	32
16	Molecular Dynamics Simulations of 1-Ethyl-3-methylimidazolium Bis[(trifluoromethyl)sulfonyl]imide Clusters and Nanodrops. Journal of Physical Chemistry A, 2015, 119, 352-368.	2.5	32
17	The N(⁴ S) + O ₂ (X ³ Σâ^'g) ↔ O(³ P) + NO(X ² Î) reaction: thermal and vibrational relaxation rates for the ² A′, ⁴ A′ and ² A′′ states. Physical Chemistry Chemical Physics, 2020, 22, 3927-3939.	2.8	30
18	Probing the dynamics of weakly bound complexes using high-resolution laser spectroscopy. Faraday Discussions, 1994, 97, 57.	3.2	29

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19	Reactive collisions for NO(² Î) + N(⁴ S) at temperatures relevant to the hypersonic flight regime. Physical Chemistry Chemical Physics, 2017, 19, 2392-2401.	2.8	29
20	Collision-induced dissociation of positive and negative copper oxide cluster ions generated by direct laser desorption/ionization of copper oxide. International Journal of Mass Spectrometry and Ion Processes, 1990, 102, 115-132.	1.8	27
21	Near-infrared laser spectroscopy of the Ar–C2HD complex:. Chemical Physics Letters, 1997, 281, 272-280.	2.6	26
22	Fragmentation of HCN in optically selected mass spectrometry: Nonthermal ion cooling in helium nanodroplets. Journal of Chemical Physics, 2005, 123, 141103.	3.0	26
23	Electrospray of an Energetic Ionic Liquid Monopropellant for Multi-Mode Micropropulsion Applications. , 2015, , .		26
24	Molecular Mechanisms Underlying Solute Retention at Heterogeneous Interfaces. Journal of Physical Chemistry Letters, 2017, 8, 4600-4607.	4.6	26
25	Krypton charge exchange cross sections for Hall effect thruster models. Journal of Applied Physics, 2013, 113, 163301.	2.5	25
26	Electrospray of 1-Butyl-3-Methylimidazolium Dicyanamide Under Variable Flow Rate Operations. Journal of Propulsion and Power, 2014, 30, 1701-1710.	2.2	25
27	Accurate reproducing kernel-based potential energy surfaces for the triplet ground states of N ₂ O and dynamics for the N + NO ↔ O + N ₂ and N ₂ + O → 2N + O reactions. Physical Chemistry Chemical Physics, 2020, 22, 18488-18498.	2.8	24
28	High resolution infrared molecular beam spectroscopy of cyanoacetylene clusters. Journal of Chemical Physics, 1995, 103, 8828-8839.	3.0	23
29	Communication: Equilibrium rate coefficients from atomistic simulations: The O(3P) + NO(2Î) → O2(<i>X</i> 3Σ <i>g</i> â^') + N(4S) reaction at temperatures relevant to the hypersonic flight regime. Journal of Chemical Physics, 2015, 142, 091104.	3.0	22
30	The ethylene–carbon dioxide complex: A double internal rotor. Journal of Chemical Physics, 1995, 103, 7788-7795.	3.0	18
31	Electron-Catalyzed Mutual Neutralization of Various Anions with <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msup><mml:mi>Ar</mml:mi><mml:mo>+</mml:mo></mml:msup>: Evidence of a New Plasma Process. Physical Review Letters. 2011. 106. 018302.</mml:math 	7.8	18
32	Collision-induced rotational excitation in N2+(2Σg+,v=)–Ar: Comparison of computations and experiment. Journal of Chemical Physics, 2016, 144, 224307.	3.0	16
33	From in silica to in silico: retention thermodynamics at solid–liquid interfaces. Physical Chemistry Chemical Physics, 2018, 20, 18610-18622.	2.8	15
34	The structure and intermolecular dynamics of the nitrous oxide–ethylene complex: Experiment and ab initio theory. Journal of Chemical Physics, 1996, 104, 4411-4418.	3.0	13
35	Dynamics on Multiple Potential Energy Surfaces: Quantitative Studies of Elementary Processes Relevant to Hypersonics. Journal of Physical Chemistry A, 2020, 124, 6255-6269.	2.5	13
36	Machine Learning for Observables: Reactant to Product State Distributions for Atom–Diatom Collisions. Journal of Physical Chemistry A, 2020, 124, 7177-7190.	2.5	12

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37	Emission-Excitation Cross Sections Relevant to Krypton-Propelled Electric Thrusters. Journal of Propulsion and Power, 2015, 31, 725-736.	2.2	11
38	Capillary ionic liquid electrospray: beam compositional analysis by orthogonal time-of-flight mass spectrometry. Journal of Fluid Mechanics, 2021, 928, .	3.4	11
39	The argon–diacetylene complex: An example of distributed interactions and transferable potentials. Journal of Chemical Physics, 1996, 105, 10171-10177.	3.0	10
40	Mass Spectrometry of Selected Ionic Liquids in Capillary Electrospray at Nanoliter Volumetric Flow Rates. , 2016, , .		10
41	The C(³ P) + O ₂ (³ Σ _g ^{â^{-*}}) â ⁺ CO ₂ â CO(¹ Σ ⁺) + O(¹ D)/O(³ P) reaction: thermal and vibrational relaxation rates from 15 K to 20 000 K. Physical Chemistry Chemical Physics, 2021, 23, 11251-11263.	†" 2.8	10
42	Machine learning product state distributions from initial reactant states for a reactive atom–diatom collision system. Journal of Chemical Physics, 2022, 156, 034301.	3.0	10
43	Oxidation of carbamate-protected alkylhydrazines to the corresponding hydrazones under Swern conditions. Tetrahedron Letters, 2001, 42, 1453-1454.	1.4	8
44	Formation of cold ion-neutral clusters using superfluid helium nanodroplets. Review of Scientific Instruments, 2010, 81, 054101.	1.3	8
45	Orthogonal time-of-flight mass spectrometry of an ion beam with a broad kinetic energy profile. Review of Scientific Instruments, 2017, 88, 105111.	1.3	7
46	Communication: Vibrational relaxation of CO($ll \hat{E}$) in collision with Ar(1S) at temperatures relevant to the hypersonic flight regime. Journal of Chemical Physics, 2017, 146, 111102.	3.0	6
47	A Combined Experimental and Theoretical Treatment of Ionic Liquid Thermal Dissociation. , 2017, , .		6
48	Mesospheric implications for the reaction of Si ⁺ with O ₂ (<i>a</i> ¹ î" _g). Geophysical Research Letters, 2010, 37, .	4.0	5
49	A new technique to study plasma chemistry kinetics. Journal of Physics: Conference Series, 2011, 300, 012007.	0.4	4
50	Application of a First Generation Collisional Radiative Model for Iodine to Optical Emissions from the Plume of an Iodine Hall Effect Thruster. , 2017, , .		4
51	A guided-ion beam study of the collisions and reactions of I+ and I2+ with I2. Journal of Chemical Physics, 2015, 142, 074301.	3.0	3
52	Integral cross section measurements and product recoil velocity distributions of Xe2+ + N2 hyperthermal charge-transfer collisions. Journal of Chemical Physics, 2016, 145, 044309.	3.0	2
53	Quantum and quasiclassical trajectory studies of rotational relaxation in Ar–N ₂ ⁺ collisions. Physical Chemistry Chemical Physics, 2017, 19, 27945-27951.	2.8	2
54	Solvated Ion Cluster Dissociation Rates for Ionic Liquid Electrospray Propellants. , 2019, , .		1

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#	Article	IF	CITATIONS
55	Aqueous Phosphoric Acid as a Mild Reagent for Deprotection of the t-Butoxycarbonyl Group ChemInform, 2004, 35, no.	0.0	Ο
56	Photoemission resulting from collisions of I ₂ with 5–100 eV electrons. Journal of Physics B: Atomic, Molecular and Optical Physics, 2020, 53, 215201.	1.5	0