

James M Farrar

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Experiment and theory elucidate the pathways for H ₃ ⁺ formation in the ultrafast double ionization in methanol. <i>Natural Sciences</i> , 2022, 2, .	2.1	0
2	Chemi-ionization Reactions and Basic Stereodynamical Effects in Collisions of Atom-Molecule Reagents. <i>Journal of Physical Chemistry A</i> , 2021, 125, 3307-3315.	2.5	8
3	Stereo-dynamical effects in chemi-ionization reactions of atmospheric O ₂ and N ₂ molecules promoted by collisions with Ne*(3P ₂ ,0) atoms. <i>Chemical Physics Letters</i> , 2021, 778, 138813.	2.6	5
4	Quantum-State Controlled Reaction Channels in Chemi-ionization Processes: Radiative (Optical-Physical) and Exchange (Oxidative-Chemical) Mechanisms. <i>Accounts of Chemical Research</i> , 2020, 53, 2248-2260.	15.6	27
5	A New Insight on Stereo-Dynamics of Penning Ionization Reactions. <i>Frontiers in Chemistry</i> , 2019, 7, 445.	3.6	13
6	A Velocity Map Imaging Study of the Reactions of O ⁺ (4S) With CH ₄ . <i>Frontiers in Chemistry</i> , 2019, 7, 227.	3.6	1
7	Angular and energy distributions of fragment ions in dissociative double photoionization of acetylene molecules in the 31.9-50.0 eV photon energy range. <i>Journal of Chemical Physics</i> , 2016, 145, 114308.	3.0	13
8	Velocity Map Imaging Study of Ion-Radical Chemistry: Charge Transfer and Carbon-Carbon Bond Formation in the Reactions of Allyl Radicals with C ⁺ . <i>Journal of Physical Chemistry A</i> , 2016, 120, 6122-6128.	2.5	9
9	Imaging ion-molecule reactions: Charge transfer and halide transfer reactions of O ⁺ with CH ₃ Cl, CH ₃ Br, and CH ₃ I. <i>International Journal of Mass Spectrometry</i> , 2015, 377, 93-100.	1.5	11
10	Velocity Map Imaging Study of Charge-Transfer and Proton-Transfer Reactions of CH ₃ Radicals with H ₃ ⁺ . <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 1684-1689.	4.6	51
11	Ion-molecule reaction dynamics: Velocity map imaging studies of N ⁺ and O ⁺ with CD ₃ OD. <i>Journal of Chemical Physics</i> , 2015, 143, 084304.	3.0	6
12	Kinetic Energy Release in molecular dications fragmentation after VUV and EUV ionization and escape from planetary atmospheres. <i>Planetary and Space Science</i> , 2014, 99, 149-157.	1.7	49
13	Ion imaging study of dissociative charge transfer in the N ₂ ⁺ + CH ₄ system. <i>Journal of Chemical Physics</i> , 2013, 138, 124304.	3.0	9
14	Imaging ion-molecule reactions: Charge transfer and C-N bond formation in the C ⁺ + NH ₃ system. <i>Journal of Chemical Physics</i> , 2012, 136, 204305.	3.0	24
15	Ion imaging study of reaction dynamics in the N ⁺ + CH ₄ system. <i>Journal of Chemical Physics</i> , 2012, 137, 154312.	3.0	15
16	Singlet and triplet state dynamics of charge and hydride transfer reactions of OD ⁺ (X ³ Σ ⁺) with propyne. <i>International Journal of Mass Spectrometry</i> , 2009, 280, 154-161.	1.5	2
17	Vibrational-Rotational Energy Distributions in the Reaction O ⁺ + D ₂ → OD + D ⁺ . <i>Journal of Physical Chemistry A</i> , 2009, 113, 15233-15239.	2.5	2
18	Hydride transfer reaction dynamics of OD ⁺ +C ₃ H ₆ . <i>Journal of Chemical Physics</i> , 2007, 126, 244315.	3.0	2

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19	Low energy crossed beam studies of OD ⁺ and D ₂ O ⁺ with C ₂ H ₄ : covalent and electrostatic complexes. <i>Physica Scripta</i> , 2007, 76, C48-C55.	2.5	0
20	Reaction dynamics of OH ⁺ ($\hat{1}\hat{1}\hat{3}$) + C ₂ H ₂ studied with crossed beams and density functional theory calculations. <i>Journal of Chemical Physics</i> , 2006, 125, 133117.	3.0	3
21	Dynamics study of the reaction OH ⁺ + C ₂ H ₂ → C ₂ H ⁺ + H ₂ O with crossed beams and density-functional theory calculations. <i>Journal of Chemical Physics</i> , 2006, 124, 124317.	3.0	7
22	Experimental and theoretical studies of charge transfer and hydride transfer in the reactions of OD ⁺ and C ₂ H ₄ . <i>International Journal of Mass Spectrometry</i> , 2005, 241, 271-282.	1.5	8
23	Quantum State-Resolved Study of the Four-Atom Reaction OH ⁻ ($\hat{X}\hat{1}\hat{g}^+$) + D ₂ ($\hat{X}\hat{1}\hat{g}^+$, $\nu = 0$) → HOD($\hat{X}\hat{1}\hat{A}\hat{g}^+$, $\nu\hat{a}\hat{e}^-$) _{2.5-D} (1S). <i>Journal of Physical Chemistry A</i> , 2005, 109, 6392-6396.	2.5	5
24	Proton transfer dynamics of the reaction H ₃ O ⁺ (NH ₃ , H ₂ O)NH ₄ ⁺ studied using the crossed molecular beam technique. <i>Journal of Chemical Physics</i> , 2004, 120, 199-205.	3.0	12
25	Experimental and theoretical studies of charge transfer and deuterium ion transfer between D ₂ O ⁺ and C ₂ H ₄ . <i>Journal of Chemical Physics</i> , 2004, 121, 3495-3506.	3.0	10
26	Spectroscopy and reactivity of size-selected Mg ^[sup +] -ammonia clusters. <i>Journal of Chemical Physics</i> , 2004, 121, 8375.	3.0	20
27	Reaction Dynamics of H ₂ O ⁺ (D ₂ O ⁺) + NH ₃ Studied with Crossed Molecular Beams and Density Functional Theory Calculations. <i>Journal of Physical Chemistry A</i> , 2004, 108, 9876-9886.	2.5	11
28	Size-dependent reactivity in open shell metal-ion polar solvent clusters: spectroscopic probes of electronic-vibration coupling, oxidation and ionization. <i>International Reviews in Physical Chemistry</i> , 2003, 22, 593-640.	2.3	50
29	CHEMICAL REACTIONS: Steric and Solvent Effects in Ionic Reactions. <i>Science</i> , 2002, 295, 2222-2223.	12.6	1
30	Photodissociation Spectra for Size-Selected Sr ⁺ (CH ₃ OH) _n and Sr ⁺ (CH ₃ OD) _n Clusters. <i>Journal of Physical Chemistry A</i> , 2002, 106, 9993-9998.	2.5	7
31	Hot and Cold Clusters: Photodissociation of Sr ⁺ (CH ₃ OD) _n through Vibrationally Excited Intermediates. <i>Journal of Physical Chemistry A</i> , 2002, 106, 11882-11890.	2.5	6
32	Spectroscopy and reactivity of size-selected Mg ⁺ methanol clusters. <i>Journal of Chemical Physics</i> , 2001, 114, 6180-6189.	3.0	36
33	Dynamics of the OH ⁺ + D ₂ isotope exchange reaction: Reactive and nonreactive decay of the collision complex. <i>Journal of Chemical Physics</i> , 2000, 113, 581-595.	3.0	9
34	Dynamics of the reaction of O ⁻ with D ₂ at low collision energies: reagent rotational energy effects. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 679-685.	2.8	1
35	Vibrational state-resolved study of the O ⁺ + H ₂ reaction: Isotope effects on the product energy partitioning. <i>Journal of Chemical Physics</i> , 1999, 111, 7348-7358.	3.0	7
36	Cluster size specific chemistry: deuterium atom pickup in Sr ⁺ solvated by ammonia. <i>Chemical Physics Letters</i> , 1999, 304, 350-356.	2.6	12

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37	Spectroscopic studies of mass selected clusters of Sr ⁺ solvated by H ₂ O and D ₂ O. Journal of Chemical Physics, 1999, 111, 8469-8480.	3.0	49
38	Ion Reaction Dynamics. Annual Review of Physical Chemistry, 1995, 46, 525-554.	10.8	33
39	Frequency- and time-resolved cluster photodissociation dynamics in Sr ⁺ (H ₂ O) _n , Sr ⁺ (NH ₃) _n and Sr ⁺ (CH ₃ OH) _n . Journal of the Chemical Society, Faraday Transactions, 1993, 89, 1457.	1.7	27
40	Highlights: Ventures in freshman chemistry. Journal of Chemical Education, 1993, 70, 847.	2.3	0
41	Dynamics of the reaction of C ⁽⁺⁾ with HCl. Astrophysical Journal, 1988, 335, 491.	4.5	10
42	Reactive scattering from double-minimum potentials. Faraday Discussions of the Chemical Society, 1987, 84, 281.	2.2	6
43	Dynamics of the reaction of N ⁺ with H ₂ . V. Reactive and nonreactive scattering of N ⁺ (3P) at relative energies below 3.6 eV. Journal of Chemical Physics, 1980, 73, 3750-3762.	3.0	27
44	Ethyl trichloroacetate hydrolysis. I. Kinetic evidence for a common tetrahedral intermediate in the acid-catalyzed and water-catalyzed hydrolyses. Journal of the American Chemical Society, 1975, 97, 2250-2254.	13.7	7
45	Entropies of dissociation of some moderately strong acids. Journal of the American Chemical Society, 1969, 91, 6057-6062.	13.7	57