# Millard H Alexander

List of Publications by Year
in descending order



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|---|---|---|---|
| 175 papers | 7,428 citations | 44042 48 h-index | 71651 76 g-index |
| 179 all docs | 179 docs citations | 179 times ranked | 1660 citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | Enhanced reactivity of fluorine with para-hydrogen in cold interstellar clouds by resonance-induced quantum tunnelling. Nature Chemistry, 2019, 11, 744-749. | 6.6 | 34 |
| 2 | Accurate characterization of the lowest triplet potential energy surface of SO2 with a coupled cluster method. Journal of Chemical Physics, 2019, 150, 144303. | 1.2 | 2 |
| 3 | Experimental and theoretical investigation of the temperature dependent electronic quenching of O(1D) atoms in collisions with Kr. Journal of Chemical Physics, 2018, 148, 124311. | 1.2 | 12 |
| 4 | Photoabsorption Assignments for the Cl̃ƒ1B2 â†•Xl̃ƒ1A1 Vibronic Transitions of SO2, Using New Ab Initio Potential Energy and Transition Dipole Surfaces. Journal of Physical Chemistry A, 2017, 121, 1012-1021. | 1.1 | 18 |
| 5 | Final State Resolved Quantum Predissociation Dynamics of SO<sub>2</sub>(Cl̃ƒ<sup>1</sup>B̃<sub>2</sub>) and Its Isotopomers via a Crossing with a Singlet Repulsive State. Journal of Physical Chemistry A, 2017, 121, 4930-4938. | 1.1 | 9 |
| 6 | First-principles C band absorption spectra of SO<sub>2</sub> and its isotopologues. Journal of Chemical Physics, 2017, 146, 154305. | 1.2 | 7 |
| 7 | The interaction of NO(X2Î) with H2: Ab initio potential energy surfaces and bound states. Journal of Chemical Physics, 2017, 146, 114301. | 1.2 | 17 |
| 8 | Accurate transport properties for O(3P)â€“H and O(3P)â€“H2. Journal of Chemical Physics, 2016, 145, 164309. | 1.2 | 11 |
| 9 | New ab initio adiabatic potential energy surfaces and bound state calculations for the singlet ground Xl̃ƒ1A1 and excited Cl̃ƒ1B2(21Aâ€²) states of SO2. Journal of Chemical Physics, 2016, 144, 174301. | 1.2 | 17 |
| 10 | Chemical Control and Spectral Fingerprints of Electronic Coupling in Carbon Nanostructures. Journal of Physical Chemistry C, 2016, 120, 29476-29483. | 1.5 | 2 |
| 11 | Publisherâ€™s Note: â€œNew ab initio adiabatic potential energy surfaces and bound state calculations for the singlet ground Xl̃ƒ1A1 and excited Cl̃ƒ1B2(21Aâ€²) states of SO2â€•[J. Chem. Phys. 144, 174301 (2016)]. Journal of Chemical Physics, 2016, 144, 209901. | 1.2 | 0 |
| 12 | Resonances in rotationally inelastic scattering of NH3 and ND3 with H2. Journal of Chemical Physics, 2015, 143, 044312. | 1.2 | 34 |
| 13 | Theoretical investigation of the dynamics of O(1Dâ†'3P) electronic quenching by collision with Xe. Journal of Chemical Physics, 2015, 143, 054306. | 1.2 | 7 |
| 14 | Electronic quenching of O(1D) by Xe: Oscillations in the product angular distribution and their dependence on collision energy. Journal of Chemical Physics, 2015, 143, 054307. | 1.2 | 4 |
| 15 | State-Specific Collision Dynamics of Molecular Super Rotors with Oriented Angular Momentum. Journal of Physical Chemistry A, 2015, 119, 12471-12479. | 1.1 | 14 |
| 16 | A finite-element visualization of quantum reactive scattering. II. Nonadiabaticity on coupled potential energy surfaces. Journal of Chemical Physics, 2015, 142, 034108. | 1.2 | 1 |
| 17 | Spectroscopic observation of resonances in the F + H <sub>2</sub> reaction. Science, 2015, 349, 510-513. | 6.0 | 98 |
| 18 | Rotationally inelastic scattering of OH by molecular hydrogen: Theory and experiment. Journal of Chemical Physics, 2015, 142, 204310. | 1.2 | 34 |

| # | Article | IF | Citations |
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| 19 | Theoretical investigation of the relaxation of the bending mode of ${m CH}_2(ilde{X})$ CH 2(XÌƒ) by collisions with helium. Journal of Chemical Physics, 2014, 141, 214305. | 1.2 | 2 |
| 20 | Collision dynamics of symmetric top molecules: A comparison of the rotationally inelastic scattering of CD3 and ND3 with He. Journal of Chemical Physics, 2014, 140, 134308. | 1.2 | 10 |
| 21 | A MATLAB-based finite-element visualization of quantum reactive scattering. I. Collinear atom-diatom reactions. Journal of Chemical Physics, 2014, 141, 024118. | 1.2 | 4 |
| 22 | The interaction of OH(X2Î) with H2: Ab initio potential energy surfaces and bound states. Journal of Chemical Physics, 2014, 141, 174309. | 1.2 | 26 |
| 23 | Transport Properties for Systems with Deep Potential Wells: H + O<sub>2</sub>. Journal of Physical Chemistry A, 2014, 118, 11935-11942. | 1.1 | 13 |
| 24 | Theoretical investigation of intersystem crossing between the aÌƒA11 and XÌƒB13 states of CH2 induced by collisions with helium. Journal of Chemical Physics, 2014, 141, 064312. | 1.2 | 4 |
| 25 | The rate of the FÂ+ÂH2 reaction at very low temperatures. Nature Chemistry, 2014, 6, 141-145. | 6.6 | 105 |
| 26 | State-to-state quantum dynamics of the F + HCl (vi = 0, ji = 0) â†' HF(vf, jf) + Cl reaction on the ground state potential energy surface. Physical Chemistry Chemical Physics, 2013, 15, 15347. | 1.3 | 15 |
| 27 | Rotationally inelastic scattering of CD3 and CH3 with He: comparison of velocity map-imaging data with quantum scattering calculations. Chemical Science, 2013, 4, 4199. | 3.7 | 16 |
| 28 | Exact quantum scattering calculations of transport properties: CH2($ilde{X}^3$XÌƒ3B1,) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 3 | 1.2 | 14 |
| 29 | Exact quantum scattering calculations of transport properties for the H2Oâ€“H system. Journal of Chemical Physics, 2013, 139, 194309. | 1.2 | 16 |
| 30 | Uncloaking the Quantum Nature of Inelastic Molecular Collisions. Science, 2013, 341, 1076-1077. | 6.0 | 11 |
| 31 | Theoretical study of the vibrational relaxation of the methyl radical in collisions with helium. Journal of Chemical Physics, 2013, 138, 104317. | 1.2 | 10 |
| 32 | Theoretical investigation of rotationally inelastic collisions of CH2(XÌƒ) with helium. Journal of Chemical Physics, 2012, 136, 224306. | 1.2 | 20 |
| 33 | Resonances in rotationally inelastic scattering of OH(X2Î) with helium and neon. Journal of Chemical Physics, 2012, 136, 144308. | 1.2 | 24 |
| 34 | Exact quantum scattering calculation of transport properties for free radicals: OH(X2Î)â€“helium. Journal of Chemical Physics, 2012, 137, 094306. | 1.2 | 11 |
| 35 | Spin-orbit quenching of Cl(2P1/2) by H2. Journal of Chemical Physics, 2012, 136, 124312. | 1.2 | 4 |
| 36 | Theoretical investigation of rotationally inelastic collisions of the methyl radical with helium. Journal of Chemical Physics, 2011, 135, 064306. | 1.2 | 20 |

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| 37 | Communication: Non-adiabatic coupling and resonances in the F + H2 reaction at low energies. Journal of Chemical Physics, 2011, 134, 231101. | 1.2 | 45 |
| 38 | Chemical Kinetics Under Test. Science, 2011, 331, 411-412. | 6.0 | 9 |
| 39 | Theoretical investigation of rotationally inelastic collisions of CH2(Ã£) with helium. Journal of Chemical Physics, 2011, 134, 154307. | 1.2 | 27 |
| 40 | Depolarization in H<sub>2</sub>Oâ€“He collisions. Molecular Physics, 2010, 108, 1159-1169. | 0.8 | 13 |
| 41 | Low-energy inelastic collisions of OH radicals with He atoms and<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mi mathvariant="normal">D</mml:mi><mml:mrow><mml:mn>2</mml:mn></mml:mrow></mml:msub></mml:mrow></mml:math>mole Physical Review A, 2010, 82, . | 1.0 | 44 |
| 42 | Time-dependent wavepacket investigation of state-to-state reactive scattering of Cl with para-H2 including the open-shell character of the Cl atom. Journal of Chemical Physics, 2010, 132, 034308. | 1.2 | 38 |
| 43 | State-to-state inelastic scattering of Stark-decelerated OH radicals with Ar atoms. Physical Chemistry Chemical Physics, 2010, 12, 10660. | 1.3 | 57 |
| 44 | Joint Experimentalâ^'Theoretical Investigation of the Lower Bound States of the NO(X<sup>2</sup>Î)-Kr Complex. Journal of Physical Chemistry A, 2009, 113, 7366-7375. | 1.1 | 22 |
| 45 | Tensor cross sections and collisional depolarization of OH(Xâ€‰2Î) in collisions with helium. Journal of Chemical Physics, 2009, 130, 164315. | 1.2 | 40 |
| 46 | Nonadiabatic Interactions in the Cl + H <sub>2</sub> Reaction Probed by ClH <sub>2</sub> <sup>-</sup> and ClD <sub>2</sub> <sup>-</sup> Photoelectron Imaging. Science, 2008, 319, 72-75. | 6.0 | 74 |
| 47 | The Extent of Nonâ€“Born-Oppenheimer Coupling in the Reaction of Cl( <sup>2</sup> P ) with para- H <sub>2</sub>. Science, 2008, 322, 573-576. | 6.0 | 95 |
| 48 | Nonadiabatic effects in the photodetachment of ClH2â^'. Journal of Chemical Physics, 2008, 128, 084312. | 1.2 | 10 |
| 49 | Role of van der Waals resonances in the vibrational relaxation of HF by collisions with H atoms. Journal of Chemical Physics, 2007, 127, 114301. | 1.2 | 21 |
| 50 | Breakdown of the Born-Oppenheimer Approximation in the F+ o -D <sub>2</sub> â†' DF + D Reaction. Science, 2007, 317, 1061-1064. | 6.0 | 149 |
| 51 | An ab initio investigation of the O(3P)â€“H2(1Î£+g) van der Waals well. Physical Chemistry Chemical Physics, 2006, 8, 4420-4426. | 1.3 | 17 |
| 52 | Product multiplet branching in the O(1D)+H2â†'OH(2Î)+H reaction. Journal of Chemical Physics, 2004, 121, 5221-5235. | 1.2 | 69 |
| 53 | Angular distributions for the F+H2â†'HF+H reaction: The role of the F spin-orbit excited state and comparison with molecular beam experiments. Journal of Chemical Physics, 2004, 121, 5812-5820. | 1.2 | 24 |
| 54 | Reactivity of the F spinâ€“orbit excited state in the F + HD reaction: Product translational and rotational energy distributions. Physical Chemistry Chemical Physics, 2004, 6, 5018-5025. | 1.3 | 12 |

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| 55 | Details and consequences of the nonadiabatic coupling in the Cl(2P)â€‰+â€‰H2reaction. Faraday Discussions, 2004, 127, 59-72. | 1.6 | 48 |
| 56 | A simple theoretical study of the ClH2? photoelectron spectrum. Physical Chemistry Chemical Physics, 2004, 6, 4984. | 1.3 | 15 |
| 57 | The dynamics of the prototype abstraction reaction Cl(2P3/2,1/2)+ H2: A comparison of crossed molecular beam experiments with exact quantum scattering calculations on coupled ab initio potential energy surfaces. Physical Chemistry Chemical Physics, 2004, 6, 5007. | 1.3 | 36 |
| 58 | Experimental and theoretical investigation of the AlH bâ€Š3Î£âˆ’â€“aâ€‰3Î electronic transition. Journal of Chemical Physics, 2003, 118, 10477-10484. | 1.2 | 9 |
| 59 | Theoretical investigation of the lower bend-stretch states of the Clâˆ’H2 anion complex and its isotopomers. Journal of Chemical Physics, 2003, 118, 9637-9642. | 1.2 | 30 |
| 60 | An ab initio based model for the simulation of multiple 2P atoms embedded in a cluster of spherical ligands, with application to Al in solid para-hydrogen. Journal of Chemical Physics, 2002, 117, 5311-5318. | 1.2 | 2 |
| 61 | Laser spectroscopic study of the SiAr van der Waals complex. Journal of Chemical Physics, 2002, 116, 9239-9248. | 1.2 | 9 |
| 62 | Theoretical Study of the Validity of the Born-Oppenheimer Approximation in the Cl + H2rightarrow HCl + H Reaction. Science, 2002, 296, 715-718. | 6.0 | 138 |
| 63 | Experimental and Theoretical Study of State-Resolved Electronically Inelastic Collisions of Highly Rotationally Excited CN(A2Î) with Argon and Helium:â€‰ The Role of Gateway Levels. Journal of Physical Chemistry A, 2002, 106, 8345-8354. | 1.1 | 12 |
| 64 | Electronic spectroscopy and excited state dynamics of the Alâ€“H2/D2 complex. Faraday Discussions, 2001, 118, 387-404. | 1.6 | 13 |
| 65 | Fully State-Resolved Differential Cross Sections for the Inelastic Scattering of the Open-Shell NO Molecule by Ar. Science, 2001, 294, 832-834. | 6.0 | 108 |
| 66 | Experimental and theoretical investigation of the rotational structure of the Alâ€“H2/D2 complex. Journal of Chemical Physics, 2001, 114, 8938-8947. | 1.2 | 15 |
| 67 | Experimental and theoretical study of Î›-doublet resolved rotationally inelastic collisions of highly rotationally excited CH(Aâ€Š2Î”,v=0) with Ar. Journal of Chemical Physics, 2001, 115, 800-809. | 1.2 | 12 |
| 68 | State-resolved rotationally inelastic collisions of highly rotationally excited CN(Aâ€Š2Î) with helium: Influence of the interaction potential. Journal of Chemical Physics, 2001, 115, 8393-8402. | 1.2 | 28 |
| 69 | State-to-state rate coefficients for transfer from the rotational levels J = 7.5, 20.5, 31.5 and 40.5 in NO(X 2Î1/2, v = 2) in collisions with He, Ar and N2 and for J=7.5, 20.5 and 31.5 in collisions with NO: comparisons between experiment and theory. Physical Chemistry Chemical Physics, 2000, 2, 473-479. | 1.3 | 18 |
| 70 | An investigation of the F+H2 reaction based on a full ab initio description of the open-shell character of the F(2P) atom. Journal of Chemical Physics, 2000, 113, 11084-11100. | 1.2 | 238 |
| 71 | A joint theoreticalâ€“experimental investigation of the lower bound states of the NO(Xâ€Š2Î)â€“Ar complex. Journal of Chemical Physics, 2000, 113, 73-85. | 1.2 | 31 |
| 72 | Experimental and theoretical study of the electronic spectrum of the BAr2 complex: Transition to the excited valence B(2s2p2â€Š2D) state. Journal of Chemical Physics, 2000, 112, 5037-5043. | 1.2 | 14 |

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| 73 | Potential energy surfaces for and energetics of the weakly-bound Alâ€“H2 and Bâ€“H2 complexes. Journal of Chemical Physics, 2000, 112, 5722-5730. | 1.2 | 31 |
| 74 | Experimental and theoretical study of rotationally inelastic collisions of highly rotationally excited CN(Aâ€Š2Î) with Ar. Journal of Chemical Physics, 2000, 112, 4474-4484. | 1.2 | 36 |
| 75 | A new, fully ab initio investigation of the NO(Xâ€‰2Î)Ar system. I. Potential energy surfaces and inelastic scattering. Journal of Chemical Physics, 1999, 111, 7426-7434. | 1.2 | 109 |
| 76 | A new, fully ab initio investigation of the ArNO(Xâ€‰2Î) system. II. Bound states of the Arâ€“NO complex. Journal of Chemical Physics, 1999, 111, 7435-7439. | 1.2 | 36 |
| 77 | A combined experimental and theoretical study of rotational energy transfer in collisions between NO(Xâ€‰2Î1/2, v=3,J) and He, Ar and N2 at temperatures down to 7 K. Journal of Chemical Physics, 1998, 109, 3882-3897. | 1.2 | 74 |
| 78 | Spinâ€“orbit effects in the reaction of F(2P) with H2. Journal of Chemical Physics, 1998, 109, 5710-5713. | 1.2 | 131 |
| 79 | Theoretical investigation of weakly-bound complexes of O(3P) with H2. Journal of Chemical Physics, 1998, 108, 4467-4477. | 1.2 | 35 |
| 80 | Experimental and theoretical study of the AlNe complex. Journal of Chemical Physics, 1998, 108, 3522-3530. | 1.2 | 26 |
| 81 | Inelastic collisions of fine structure and Î›-doublet resolved rotational states of PH(Aâ€‰3Î,â€‰v=0) with helium. Journal of Chemical Physics, 1997, 106, 7642-7653. | 1.2 | 18 |
| 82 | A collaborative theoretical and experimental study of the structure and electronic excitation spectrum of the Bar and Bar2 complexes. Journal of Chemical Physics, 1997, 106, 6320-6331. | 1.2 | 33 |
| 83 | Adiabatic and diabatic potential-energy surfaces of the CN(Xâ€‰2Î£+,Aâ€‰2Î)Ne complex and nonadiabatic predissociation dynamics. Journal of Chemical Physics, 1997, 107, 7148-7162. | 1.2 | 27 |
| 84 | Stateâ€resolved inelastic collisions of single rotational, fineâ€structure, and Î› doublet levels of NH(Aâ€‰3Î) with helium: A combined experimental and theoretical study. Journal of Chemical Physics, 1996, 104, 1325-1337. | 1.2 | 28 |
| 85 | The rotational relaxation of NH(câ€‰1Î) in collisions with Ar: A combined theoretical and experimental investigation. Journal of Chemical Physics, 1995, 102, 4069-4083. | 1.2 | 28 |
| 86 | Theoretical study of the interaction of AlH(Xâ€‰1Î£+,Aâ€‰1Î) with Ar: Potential energy surfaces and bendâ€“stretch levels of the ArAlH(X,A) van der Waals complex. Journal of Chemical Physics, 1995, 102, 2413-2425. | 1.2 | 23 |
| 87 | Fullyab initioinvestigation of bound and predissociating states of the NeOH(X) complex. Journal of Chemical Physics, 1995, 103, 3400-3417. | 1.2 | 27 |
| 88 | Ab initio potential energy surfaces and quantum scattering studies of NO(Xâ€‰2Î) with He: Î›â€doublet resolved rotational and electronic fineâ€structure transitions. Journal of Chemical Physics, 1995, 103, 6973-6983. | 1.2 | 53 |
| 89 | Experimental investigation of weakly bound B(2p,3s)â€“H2/D2complexes through laser fluorescence excitation spectroscopy. Journal of Chemical Physics, 1995, 103, 7966-7974. | 1.2 | 28 |
| 90 | Experimental and theoretical study of the Bâ€“Ne nonbonding interaction: The freeâ€boundBâ€‰2Î£+â€“Xâ€‰2Î electronic transition. Journal of Chemical Physics, 1995, 103, 2779-2786. | 1.2 | 36 |

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| 91 | Theoretical investigation of weaklyâ€bound complexes of B with H2. Journal of Chemical Physics, 1995, 103, 7956-7965. | 1.2 | 44 |
| 92 | Theoretical Study of Bound States of Ar-NO. The Journal of Physical Chemistry, 1994, 98, 1073-1079. | 2.9 | 28 |
| 93 | Theoretical studies of He(1S)+CH(Xâ€‰2Î). II. Fullyabinitiocross sections for the inelastic scattering and comparison with experiment. Journal of Chemical Physics, 1994, 100, 1338-1349. | 1.2 | 23 |
| 94 | Potential energy surfaces for the interaction of BH(Xâ€‰1Î£+,Aâ€‰1Î) with Ar and a theoretical investigation of the stretchâ€bend levels of the ArBH(A) van der Waals molecule. Journal of Chemical Physics, 1994, 101, 2887-2902. | 1.2 | 67 |
| 95 | Flux redistribution during the photodissociation of ClNO in theT1state. Journal of Chemical Physics, 1994, 101, 4722-4734. | 1.2 | 17 |
| 96 | Potential energy surfaces for the interaction of CH(Xâ€‰2Î,Bâ€‰2Î£â^') with Ar and an assignment of the stretchâ€bend levels of the ArCH(B) van der Waals molecule. Journal of Chemical Physics, 1994, 101, 4547-4560. | 1.2 | 55 |
| 97 | The use of the current density in the analysis of molecular photodissociation. Journal of Chemical Physics, 1994, 101, 8663-8673. | 1.2 | 8 |
| 98 | On the generation of preferential Î›â€doublet populations in the collisional relaxation of highly rotationally excited CH(Xâ€‰2Î). Journal of Chemical Physics, 1994, 101, 7468-7479. | 1.2 | 36 |
| 99 | Scattering of NH3 by orthoâ€•and paraâ€H2: Expansion of the potential and collisional propensity rules. Journal of Chemical Physics, 1993, 98, 4662-4671. | 1.2 | 38 |
| 100 | A joint experimental and theoretical study of Aâ€‰2Îâ†'Xâ€‰2Î£+ electronic energy transfer in the CN molecule induced by collisions with helium. Journal of Chemical Physics, 1993, 98, 8580-8592. | 1.2 | 52 |
| 101 | Differential and integral cross sections for the inelastic scattering of NO (Xâ€‰2Î) by Ar based on a new ab initio potential energy surface. Journal of Chemical Physics, 1993, 99, 7725-7738. | 1.2 | 99 |
| 102 | Spinâ€“orbit branching in the photofragmentation of HCl. Journal of Chemical Physics, 1993, 99, 1752-1764. | 1.2 | 114 |
| 103 | Adiabatic and approximate diabatic potential energy surfaces for the B...H2van der Waals molecule. Journal of Chemical Physics, 1993, 99, 6014-6026. | 1.2 | 73 |
| 104 | Adiabatic representations for the study of flux redistribution during photodissociation involving coupled electronic states: The effect of vibrational excitation on the photofragmentation of CH3I. Journal of Chemical Physics, 1993, 98, 6196-6207. | 1.2 | 21 |
| 105 | Quantum flux redistribution during molecular photodissociation. Journal of Chemical Physics, 1992, 97, 2527-2535. | 1.2 | 47 |
| 106 | The study of flux redistribution during molecular photodissociation: Adiabatic and diabatic analyses and application to the dissociation of CH3I. Journal of Chemical Physics, 1992, 97, 4836-4845. | 1.2 | 31 |
| 107 | Quantum flux studies of the mechanism of Ca(4s5pâ€‰1P)â†'Ca(4s5pâ€‰3P) collisions. Journal of Chemical Physics, 1992, 96, 6672-6680. | 1.2 | 18 |
| 108 | Potential energy hypersurfaces for the interaction of NO with the Ag(111) surface. Journal of Chemical Physics, 1991, 94, 8454-8467. | 1.2 | 17 |

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| 109 | Collisions of NO(Xâ€‰2Î) with a Ag(111) surface: New quantum scattering studies based on a semiempirical potential energy surface. Journal of Chemical Physics, 1991, 94, 8468-8478. | 1.2 | 15 |
| 110 | Quantum study of the redistribution of flux during inelastic collisions. Journal of Chemical Physics, 1991, 95, 8931-8940. | 1.2 | 30 |
| 111 | Energetics and spin selectivity in the infrared multiphoton dissociation HN3(XÌƒ1Aâ€™)â†'N2(X1Î£g+)+NH(X3Î£â¯,a1Î”). AIP Conference Proceedings, 1989, , . | 0.3 | 0 |
| 112 | Rotational energy transfer in HF: A computational study. Journal of Chemical Physics, 1989, 91, 7563-7589. | 1.2 | 9 |
| 113 | The inelastic scattering of 2Î [case (b)] molecules and an understanding of the differing Î› doublet propensities for molecules of Ï€ vs Ï€3 orbital occupancy. Journal of Chemical Physics, 1989, 91, 839-848. | 1.2 | 130 |
| 114 | A logâ€derivative propagation scheme for the exact solution of twoâ€state curve crossing problems. Journal of Chemical Physics, 1989, 91, 2388-2395. | 1.2 | 16 |
| 115 | Quantum scattering studies of electronically inelastic collisions of CN (Xâ€‰2Î£+,â€‰Aâ€‰2Î) with He. Journal of Chemical Physics, 1989, 91, 5425-5439. | 1.2 | 113 |
| 116 | A semiclassical treatment of rotationally electronically inelastic scattering of NO from Ag(111). Journal of Chemical Physics, 1989, 90, 575-586. | 1.2 | 20 |
| 117 | Theoretical study of Ca(4s5pâ€‰1P)â‡„Ca(4s5pâ€‰3P) transitions in collision with noble gases: Integral cross sections and alignment effects. Journal of Chemical Physics, 1989, 91, 1658-1667. | 1.2 | 30 |
| 118 | Mechanism of and alignment effects in spinâ€“changing collisions involving atoms in 1P electronic states: Ca(4s5pâ€‰1P)+noble gases. Journal of Chemical Physics, 1989, 90, 5373-5385. | 1.2 | 35 |
| 119 | Inelastic collisions of OH(Xâ€‰2Î) with paraâ€H2: Î›â€doublet and hyperfineâ€structure transitions. Journal of Chemical Physics, 1988, 88, 6931-6937. | 1.2 | 43 |
| 120 | Adiabatic and diabatic potential energy surfaces for collisions of CN(Xâ€‰2Î£+,â€‰Aâ€‰2Î) with He. Journal of Chemical Physics, 1988, 89, 3139-3151. | 1.2 | 208 |
| 121 | Inelastic collisions of CaF(Aâ€‰2Î) with He and Ar: Quantum calculations and adiabatic analysis. Journal of Chemical Physics, 1988, 88, 3581-3589. | 1.2 | 13 |
| 122 | Energetics and spinâ€and Î›â€doublet selectivity in the infrared multiphoton dissociation HN3(XÌƒâ€‰1Aâ€™)â†'N2(Xâ€‰1Î£+g)+NH(X3Î£â¯,aâ€‰1Î”): Theory. Journal of Chemical Physics, 1988, 89, 1388-1400. | 1.2 | 78 |
| 123 | Quantum theory and collisional propensity rules for rotationally inelastic collisions between polyatomic molecules (NH3 and CO2) and an uncorrugated surface. Journal of Chemical Physics, 1988, 89, 790-800. | 1.2 | 5 |
| 124 | Clarification of the electronic asymmetry of Î› doublets in 3Î electronic states of diatomic molecules. Journal of Chemical Physics, 1987, 87, 7118-7124. | 1.2 | 22 |
| 125 | A theoretical study of alignment effects in collisions of N2 with a Ag surface. Journal of Chemical Physics, 1987, 87, 4937-4947. | 1.2 | 22 |
| 126 | Quantum closeâ€coupled studies of collisions of NO(Xâ€‰2Î) with a Ag(111) surface. Journal of Chemical Physics, 1987, 87, 3218-3231. | 1.2 | 49 |

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| 127 | Theoretical study of Ca(4s5p 1P)â†'Ca(4s5p 3P) transitions in collisions with He: Integral cross sections and alignment effects. Journal of Chemical Physics, 1987, 86, 4790-4800. | 1.2 | 67 |
| 128 | The infiniteâ€order sudden approximation for collisions involving molecules in Î electronic states: A new derivation and calculations of rotationally inelastic cross sections for NO(Xâ€‰2Î)+He and Ar. Journal of Chemical Physics, 1986, 85, 5652-5659. | 1.2 | 77 |
| 129 | Theory of Stark spectroscopy of molecules in1Î electronic states: Coherence effects and quantum beats. Journal of Chemical Physics, 1986, 85, 134-145. | 1.2 | 17 |
| 130 | Rotationally inelastic collisions between a molecule in a 2S+1Î£ electronic state and an openâ€shell target: General quantum analysis and experimental measurement of stateâ€resolved cross sections for CaCl(Xâ€‰2Î£+)+NO(Xâ€‰2Î£). Journal of Chemical Physics, 1986, 84, 1547-1553. | 1.2 | 19 |
| 131 | Rotationally inelastic collisions between a molecule in a 2Î£ electronic state and a 2S atom: Sudden factorization, scaling, and symmetry relations. Journal of Chemical Physics, 1986, 85, 1859-1865. | 1.2 | 7 |
| 132 | Collision induced transitions between2Î and2Î£ states of diatomic molecules: Quantum theory and collisional propensity rules. Journal of Chemical Physics, 1986, 84, 100-113. | 1.2 | 94 |
| 133 | On the physical origin of propensity rules in collisions involving molecules in2Î£ electronic states. Journal of Chemical Physics, 1986, 84, 3049-3058. | 1.2 | 51 |
| 134 | Quantum treatment of rotationally inelastic collisions involving molecules in II electronic states: New derivation of the coupling potential. Chemical Physics, 1985, 92, 337-344. | 0.9 | 256 |
| 135 | Collisional effects in Stark spectroscopy of molecules in1Î electronic states. Journal of Chemical Physics, 1985, 83, 3340-3348. | 1.2 | 15 |
| 136 | Rotationally inelastic collisions between a diatomic molecule in a 2S+1Î£ electronic state and a 2S atom: The optimal choice for the totalâ€j representation. Journal of Chemical Physics, 1985, 83, 5060-5067. | 1.2 | 15 |
| 137 | Inelastic collisions of CaCl(Xâ€‰2Î£+) with Ar: A collaborative theoretical and experimental study. Journal of Chemical Physics, 1985, 83, 556-566. | 1.2 | 29 |
| 138 | Collisionâ€induced transitions between molecular hyperfine levels: Quantum formalism, propensity rules, and experimental study of CaBr(Xâ€‰2Î£+)+Ar. Journal of Chemical Physics, 1985, 83, 2191-2200. | 1.2 | 86 |
| 139 | Clarification of the electronic asymmetry in Îâ€state Î› doublets with some implications for molecular collisions. Journal of Chemical Physics, 1984, 80, 4325-4332. | 1.2 | 136 |
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