Robert W. Field

List of Publications by Year in Descending Order

243 papers	7,962 citations	46 h-index	76 g-index
249 ext. papers	8,503 ext. citations	4.2 avg, IF	5.69 L-index

#	Paper	IF	Citations
243	Enantioselective orientation of chiral molecules induced by terahertz pulses with twisted polarization. *Physical Review Research*, **2021**, 3,	3.9	7
242	Substitution Reactions in the Pyrolysis of Acetone Revealed through a Modeling, Experiment, Theory Paradigm. *Journal of the American Chemical Society*, **2021**, 143, 3124-3142	16.4	9
241	Nonlinear rotational spectroscopy reveals many-body interactions in water molecules. *Proceedings of the National Academy of Sciences of the United States of America*, **2021**, 118,	11.5	2
240	Photodissociation of dicarbon: How nature breaks an unusual multiple bond.. *Proceedings of the National Academy of Sciences of the United States of America*, **2021**, 118,	11.5	3
239	Long-range model of vibrational autoionization in core-nonpenetrating Rydberg states of NO.. *Journal of Chemical Physics*, **2021**, 155, 244303	3.9	0
238	One-colour (~220 nm) resonance-enhanced (S1 Σ+1 ← X1 Σ+g) multi-photon dissociation of acetylene: probe of the C2 A1 Σ+u ← X1 Σ+g band by frequency-modulation spectroscopy. *Molecular Physics*, **2020**, 118, e1724340	1.7	0
237	Photodissociation transition states characterized by chirped pulse millimeter wave spectroscopy. *Proceedings of the National Academy of Sciences of the United States of America*, **2020**, 117, 146-151	11.5	5
236	Preparation of high orbital angular momentum Rydberg states by optical-millimeter-wave STIRAP. *Journal of Chemical Physics*, **2020**, 153, 084301	3.9	2
235	Determination of the sign of the population difference in a two-level system by frequency-modulation spectroscopy. *Molecular Physics*, **2020**, 118, e1660007	1.7	0
234	Analysis of vibrational autoionization of CaF Rydberg states. *Journal of Chemical Physics*, **2019**, 150, 154305	3.9	4
233	Call for Papers for the Journal of Molecular Spectroscopy Special Issue on Electronic Spectra and Structure of Small Molecules in Honor of Anthony J. Merer. *Journal of Molecular Spectroscopy*, **2019**, 356, 46	1.3	
232	Visible and ultraviolet laser spectroscopy of ThF. *Journal of Molecular Spectroscopy*, **2019**, 358, 1-16	1.3	6
231	Anomalous Intensities in the 2+1 REMPI Spectrum of the E Σ+ ← X Σ+ Transition of CO. *Journal of Physical Chemistry A*, **2019**, 123, 2780-2788	2.8	
230	Precision spectroscopy and comprehensive analysis of perturbations in the A1Σ+(v = 0) state of 13C18O. *Molecular Physics*, **2019**, 117, 79-96	1.7	6
229	Roadmap on STIRAP applications. *Journal of Physics B: Atomic, Molecular and Optical Physics*, **2019**, 52, 202001	1.3	54
228	One-color (212–220 nm) resonantly-enhanced (S1 ← X1 Σ+g) multi-photon dissociation of acetylene. *Journal of Molecular Spectroscopy*, **2019**, 361, 24-33	1.3	2
227	The frequency-domain infrared spectrum of ammonia encodes changes in molecular dynamics caused by a DC electric field. *Proceedings of the National Academy of Sciences of the United States of America*, **2019**, 116, 23444-23447	11.5	10

226	The dicarbon bonding puzzle viewed with photoelectron imaging. *Nature Communications*, **2019**, 10, 5199	17.4	7
225	Tribute to Hai-Lung Dai. *Journal of Physical Chemistry A*, **2019**, 123, 10463-10464	2.8	
224	Probing the predissociated levels of the S state of acetylene via H-atom fluorescence and photofragment fluorescence action spectroscopy. *Journal of Chemical Physics*, **2018**, 149, 174309	3.9	4
223	Experimental studies of the NaCs 12(0) [7]state: Spin-orbit and non-adiabatic interactions and quantum interference in the 12(0) [7]and 11(0) [5]emission spectra. *Journal of Chemical Physics*, **2017**, 146, 104302	3.9	3
222	Encoding of vinylidene isomerization in its anion photoelectron spectrum. *Science*, **2017**, 358, 336-339	33.3	36
221	Perturbations in the A1v = 0 state of 12C18O investigated via complementary spectroscopic techniques. *Molecular Physics*, **2017**, 115, 3178-3191	1.7	6
220	Direct single-shot observation of millimeter-wave superradiance in Rydberg-Rydberg transitions. *Physical Review A*, **2017**, 95,	2.6	7
219	Fourier transform emission spectra and deperturbation analysis of the A 2 X 2 and B 2 X 2 electronic transitions of ZnH. *Journal of Molecular Spectroscopy*, **2017**, 340, 21-28	1.3	
218	Coherent laser-millimeter-wave interactions en route to coherent population transfer. *Journal of Chemical Physics*, **2017**, 147, 144201	3.9	3
217	Stark Interference of Electric and Magnetic Dipole Transitions in the A-X Band of OH. *Physical Review Letters*, **2016**, 116, 153001	7.4	2
216	Saddle point localization of molecular wavefunctions. *Scientific Reports*, **2016**, 6, 33068	4.9	5
215	The origin of unequal bond lengths in the C (1)B2 state of SO2: Signatures of high-lying potential energy surface crossings in the low-lying vibrational structure. *Journal of Chemical Physics*, **2016**, 144, 144313	3.9	9
214	Observation of b2 symmetry vibrational levels of the SO2 C (1)B2 state: Vibrational level staggering, Coriolis interactions, and rotation-vibration constants. *Journal of Chemical Physics*, **2016** , 144, 144311	3.9	12
213	The rotation-vibration structure of the SO2 C (1)B2 state explained by a new internal coordinate force field. *Journal of Chemical Physics*, **2016**, 144, 144312	3.9	13
212	VIS and VUV spectroscopy of 12C17O and deperturbation analysis of the A1 1 levels. *RSC Advances*, **2016**, 6, 31588-31606	3.7	8
211	Perspective: The first ten years of broadband chirped pulse Fourier transform microwave spectroscopy. *Journal of Chemical Physics*, **2016**, 144, 200901	3.9	83
210	Electric potential invariants and ions-in-molecules effective potentials for molecular Rydberg states. *Journal of Chemical Physics*, **2016**, 145, 234301	3.9	3
209	Spectroscopy and perturbation analysis of the CO A1X1 (2,0), (3,0) and (4,0) bands. *Molecular Physics*, **2016**, 114, 627-636	1.7	14

208	A Molecular Precursor to Phosphaethyne and Its Application in Synthesis of the Aromatic 1,2,3,4-Phosphatriazolate Anion. *Journal of the American Chemical Society*, **2016**, 138, 6731-4	16.4	26
207	Product Branching in the Low Temperature Reaction of CN with Propyne by Chirped-Pulse Microwave Spectroscopy in a Uniform Supersonic Flow. *Journal of Physical Chemistry Letters*, **2015**, 6, 1599-604	6.4	37
206	Does Infrared Multiphoton Dissociation of Vinyl Chloride Yield Cold Vinylidene?. *Journal of Physical Chemistry Letters*, **2015**, 6, 2457-62	6.4	4
205	Edge effects in chirped-pulse Fourier transform microwave spectra. *Journal of Molecular Spectroscopy*, **2015**, 312, 54-57	1.3	7
204	Millimeter-wave optical double resonance schemes for rapid assignment of perturbed spectra, with applications to the C (1)B(2) state of SO2. *Journal of Chemical Physics*, **2015**, 142, 144201	3.9	17
203	Ultrafast isomerization initiated by X-ray core ionization. *Nature Communications*, **2015**, 6, 8199	17.4	74
202	Probing cis-trans isomerization in the S1 state of C2H2 via H-atom action and hot band-pumped IR-UV double resonance spectroscopies. *Journal of Chemical Physics*, **2015**, 143, 084310	3.9	9
201	Communication: Observation of local-bender eigenstates in acetylene. *Journal of Chemical Physics*, **2015**, 143, 071101	3.9	3
200	Direct detection of RydbergRydberg millimeter-wave transitions in a buffer gas cooled molecular beam. *Chemical Physics Letters*, **2015**, 640, 124-136	2.5	18
199	Observation of the simplest Criegee intermediate CH2OO in the gas-phase ozonolysis of ethylene. *Science Advances*, **2015**, 1, e1400105	14.3	66
198	Spectroscopic characterization of isomerization transition states. *Science*, **2015**, 350, 1338-42	33.3	35
197	Simplified Cartesian basis model for intrapolyad emission intensities in the bent-to-linear electronic transition of acetylene. *Journal of Physical Chemistry A*, **2015**, 119, 857-65	2.8	3
196	A chirped-pulse Fourier-transform microwave/pulsed uniform flow spectrometer. I. The low-temperature flow system. *Journal of Chemical Physics*, **2014**, 141, 154202	3.9	34
195	A Signature of Roaming Dynamics in the Thermal Decomposition of Ethyl Nitrite: Chirped-Pulse Rotational Spectroscopy and Kinetic Modeling. *Journal of Physical Chemistry Letters*, **2014**, 5, 3641-8	6.4	26
194	Gas-Phase Structure Determination of Dihydroxycarbene, One of the Smallest Stable Singlet Carbenes. *Angewandte Chemie*, **2014**, 126, 4173-4176	3.6	3
193	Chirped-Pulse millimeter-Wave spectroscopy for dynamics and kinetics studies of pyrolysis reactions. *Physical Chemistry Chemical Physics*, **2014**, 16, 15739-15751	3.6	43
192	Reduced dimension rovibrational variational calculations of the S(1) state of C2H2. II. The S(1) rovibrational manifold and the effects of isomerization. *Journal of Chemical Physics*, **2014**, 140, 024313	3.9	9
191	A chirped-pulse Fourier-transform microwave/pulsed uniform flow spectrometer. II. Performance and applications for reaction dynamics. *Journal of Chemical Physics*, **2014**, 141, 214203	3.9	46

190	Full dimensional Franck-Condon factors for the acetylene □□(1)Au-X (1)□(g)(+) transition. II. Vibrational overlap factors for levels involving excitation in ungerade modes. *Journal of Chemical Physics*, **2014**, 141, 134305	3.9	8
189	A new approach toward transition state spectroscopy. *Faraday Discussions*, **2013**, 163, 33-57; discussion 117-38	3.6	34
188	High resolution spectroscopy and perturbation analysis of the CO A1□X1□ (0,0) and (1,0) bands. *Molecular Physics*, **2013**, 111, 2163-2174	1.7	28
187	Laser-induced fluorescence study of the S1 state of doubly-substituted 13C acetylene and harmonic force field determination. *Journal of Physical Chemistry A*, **2013**, 117, 13696-703	2.8	9
186	High-accuracy estimates for the vinylidene-acetylene isomerization energy and the ground state rotational constants of :C?CH2. *Journal of Physical Chemistry A*, **2013**, 117, 11679-83	2.8	32
185	Chirped-pulse millimeter-wave spectroscopy: spectrum, dynamics, and manipulation of Rydberg-Rydberg transitions. *Journal of Chemical Physics*, **2013**, 138, 014301	3.9	17
184	Broadband velocity modulation spectroscopy of HfF+: Towards a measurement of the electron electric dipole moment. *Chemical Physics Letters*, **2012**, 546, 1-11	2.5	41
183	The □1Au state of acetylene: ungerade vibrational levels in the region 45,800□46,550 cm□1. *Molecular Physics*, **2012**, 110, 2707-2723	1.7	19
182	Commensurate two-quantum coherences induced by time-delayed THz fields. *Physical Review Letters*, **2012**, 109, 123603	7.4	57
181	Production of a beam of highly vibrationally excited CO using perturbations. *Journal of Chemical Physics*, **2012**, 136, 214201	3.9	11
180	Laser-induced fluorescence studies of HfF+ produced by autoionization. *Journal of Chemical Physics* , **2011**, 135, 154308	3.9	11
179	Molecular orientation and alignment by intense single-cycle THz pulses. *Physical Review Letters*, **2011**, 107, 163603	7.4	218
178	Design and evaluation of a pulsed-jet chirped-pulse millimeter-wave spectrometer for the 70-102 GHz region. *Journal of Chemical Physics*, **2011**, 135, 024202	3.9	61
177	Cis-trans isomerization in the S1 state of acetylene: identification of cis-well vibrational levels. *Journal of Chemical Physics*, **2011**, 134, 244310	3.9	20
176	Reduced dimension discrete variable representation study of cis-trans isomerization in the S1 state of C2H2. *Journal of Chemical Physics*, **2011**, 134, 244311	3.9	23
175	Chirped-pulse millimeter-wave spectroscopy of Rydberg-Rydberg transitions. *Physical Review Letters*, **2011**, 107, 143001	7.4	21
174	Determination of the ground electronic state in transition metal halides: ZrF. *Journal of Physical Chemistry A*, **2011**, 115, 9620-32	2.8	5
173	Spectral signatures of inter-system crossing mediated by energetically distant doorway levels: examples from the acetylene S1 state. *Journal of Physical Chemistry A*, **2011**, 115, 11921-43	2.8	5

172	A quantum defect model for the s, p, d, and f Rydberg series of CaF. *Journal of Chemical Physics*, **2011**, 134, 114313	3.9	23
171	Rotational analysis and deperturbation of the A2□-X2□ and B'2□ -X2□ emission spectra of MgH. *Journal of Chemical Physics*, **2011**, 135, 094308	3.9	24
170	Deconvolution of spectral data using a doorway-coupling model Hamiltonian. *Journal of Chemical Physics*, **2010**, 132, 134302	3.9	4
169	Time-Dependent Center-of-Gravity Metric Determines Key Dynamical Features of Doorway-Mediated Intersystem Crossing. *Journal of Physical Chemistry Letters*, **2010**, 1, 2144-2148	6.4	1
168	The Stark effect in Rydberg states of a highly polar diatomic molecule: CaF. *Journal of Chemical Physics*, **2009**, 131, 064301	3.9	6
167	Interference in acetylene intersystem crossing acts as the molecular analog of Young's double-slit experiment. *Proceedings of the National Academy of Sciences of the United States of America*, **2009**, 106, 2510-4	11.5	6
166	Stretch-bend combination polyads in the □□1Au state of acetylene, C2H2. *Journal of Molecular Spectroscopy*, **2009**, 256, 256-278	1.3	22
165	Ab initio investigation of high multiplicity □□ optical transitions in the spectra of CN and isoelectronic species. *Journal of Molecular Spectroscopy*, **2009**, 258, 6-12	1.3	8
164	Room-temperature metal-hydride discharge source, with observations on NiH and FeH. *Journal of Physical Chemistry A*, **2009**, 113, 13159-66	2.8	16
163	Perturbations in the 4□ level of the {tf="PS7CD9"{raise7pttilde lower7ptAclap}^lower4pt{hskip-.5pt1}hskip-1ptA_u} state of acetylene, C2H2This article is part of a Special Issue on Spectroscopy at the University of New Brunswick in honour of Colan Linton and Ron Lees.. *Canadian Journal of Physics*, **2009**, 87, 437-441	1.1	6
162	Separation of long-range and short-range interactions in Rydberg states of diatomic molecules. *Journal of Chemical Physics*, **2008**, 128, 194301	3.9	15
161	Direct observation of the symmetric stretching modes of □□1 A u acetylene by pulsed supersonic jet laser induced fluorescence. *Molecular Physics*, **2008**, 106, 1867-1877	1.7	16
160	Darling-Dennison resonance and Coriolis coupling in the bending overtones of the A 1A(u) state of acetylene, C2H2. *Journal of Chemical Physics*, **2008**, 129, 054304	3.9	25
159	Polarization dependence of transition intensities in double resonance experiments: unresolved spin doublets. *Journal of Chemical Physics*, **2008**, 128, 014301	3.9	11
158	Experimental mapping of the absolute magnitude of the transition dipole moment function □(R) of the Na2 A □u+-X □g+ transition. *Physical Review A*, **2008**, 77,	2.6	16
157	Evolution of chemical bonding during HCN HNC isomerization as revealed through nuclear quadrupole hyperfine structure. *Angewandte Chemie - International Edition*, **2008**, 47, 2969-72	16.4	18
156	Theoretical emission spectra of HNC(. *Chemical Physics Letters*, **2008**, 455, 145-150	2.5	0
155	Contrasting singlet-triplet dynamical behavior of two vibrational levels of the acetylene S1 2(1)3(1)B2 polyad. *Journal of Physical Chemistry A*, **2007**, 111, 12534-7	2.8	6

154	The pure rotational spectrum of CrCN (X 6 Σ+): an unexpected geometry and unusual spin interactions. *Molecular Physics*, **2007**, 105, 585-597	1.7	29
153	Intramolecular Dynamics in the Frequency Domain. *Advances in Chemical Physics*, **2007**, 463-490		7
152	Studies of intersystem crossing dynamics in acetylene. *Journal of Chemical Physics*, **2007**, 126, 184307	3.9	6
151	Observation of the A1A" state of isocyanogen. *Journal of Chemical Physics*, **2007**, 126, 244307	3.9	2
150	New spectroscopic data, spin-orbit functions, and global analysis of data on the A 1Sigmau+ and b 3Piu states of Na2. *Journal of Chemical Physics*, **2007**, 127, 044301	3.9	37
149	Resonance between electronic and rotational motions in Rydberg states of CaF. *Molecular Physics*, **2007**, 105, 1661-1673	1.7	10
148	The spin-orbit and rotational constants for the N2 C" 5Pi(ui) (v = 3) state. *Journal of Chemical Physics*, **2006**, 124, 081103	3.9	11
147	Broad shape resonance effects in CaF Rydberg states. *Journal of Chemical Physics*, **2006**, 124, 194302	3.9	3
146	Accurate inertias for large-amplitude motions: improvements on prevailing approximations. *Journal of Physical Chemistry A*, **2006**, 110, 7406-13	2.8	26
145	Electronic signatures of large amplitude motions: dipole moments of vibrationally excited local-bend and local-stretch states of S0 acetylene. *Journal of Physical Chemistry B*, **2006**, 110, 18912-20	3.4	17
144	Laboratory Measurements of the Hyperfine Structure of H 14 N 12 C and D 14 N 12 C. *Astrophysical Journal*, **2006**, 649, L53-L56	4.7	23
143	Differential temperature laser induced fluorescence spectroscopy. *Chemical Physics*, **2006**, 324, 709-720	2.3	2
142	Infrared–infrared double resonance spectroscopy of 39K2: The 1 3Σg state. *Chemical Physics Letters*, **2006**, 431, 267-271	2.5	17
141	Multi-isotopologue analyses of new vibration–rotation and pure rotation spectra of ZnH and CdH. *Journal of Molecular Spectroscopy*, **2006**, 237, 87-96	1.3	19
140	Millimeter-wave-detected, millimeter-wave optical polarization spectroscopy. *Journal of Chemical Physics*, **2005**, 123, 141102	3.9	6
139	Quantum defect theory of dipole and vibronic mixing in Rydberg states of CaF. *Journal of Chemical Physics*, **2005**, 122, 184314	3.9	22
138	Properties of nearly one-electron molecules. II. Application to the Rydberg spectrum of CaF. *Journal of Chemical Physics*, **2005**, 123, 084319	3.9	12
137	Properties of nearly one-electron molecules. I. An iterative Green function approach to calculating the reaction matrix. *Journal of Chemical Physics*, **2005**, 123, 084318	3.9	10

136	New S1 state vibrational and T3, 2, 1 spin□otational assignments in the vicinity of the acetylene A□1Au□X□1□g+V03K01 band. *Journal of Molecular Spectroscopy*, **2004**, 228, 565-579	1.3	11
135	"Spectrum-only" assignment of core-penetrating and core-nonpenetrating Rydberg states of calcium monofluoride. *Canadian Journal of Chemistry*, **2004**, 82, 791-803	0.9	14
134	Contrasting origins of the isomerization barriers for vinylidene, fluorovinylidene, and difluorovinylidene. *Journal of Chemical Physics*, **2003**, 118, 4037-4044	3.9	21
133	New vibrational assignments in the □ Au-[Xtilde] 1□ g electronic transition of acetylene, C2H2: the v?1 frequency. *Molecular Physics*, **2003**, 101, 663-673	1.7	21
132	Unexpected simplicity in the S1□0 dispersed fluorescence spectra of 13C2H2. *Journal of Chemical Physics*, **2002**, 116, 7939-7947	3.9	4
131	Semiclassical modeling of Rydberg wave-packet dynamics in diatomic molecules: Average decoupling theory. *Physical Review A*, **2002**, 65,	2.6	4
130	The Mechanism of Surface Electron Ejection by Laser Excited Metastable Molecules. *Journal of Physical Chemistry A*, **2002**, 106, 1122-1130	2.8	8
129	Comparison of CaF, ZnF, CaO, and ZnO: Their Anions and Cations in Their Ground and Low-Lying Excited States. *ACS Symposium Series*, **2002**, 238-259	0.4	14
128	Baseline subtraction using robust local regression estimation. *Journal of Quantitative Spectroscopy and Radiative Transfer*, **2001**, 68, 179-193	2.1	96
127	Core-Penetrating Rydberg Series of BaF: Single-State and Two-State Fits of New Electronic States in the 4.4 . *Journal of Molecular Spectroscopy*, **2001**, 205, 197-220	1.3	7
126	An assumption-violating application of the Lawrance□Knight deconvolution procedure: A retrieval of electronic coupling mechanisms underlying complex spectra. *Journal of Chemical Physics*, **2001**, 114, 6557-6561	3.9	9
125	The infrared-ultraviolet dispersed fluorescence spectrum of acetylene: New classes of bright states. *Journal of Chemical Physics*, **2001**, 114, 7424-7442	3.9	35
124	The dynamics of "stretched molecules": experimental studies of highly vibrationally excited molecules with stimulated emission pumping. *Annual Review of Physical Chemistry*, **2001**, 52, 811-52	15.7	99
123	Relabeling and classification of the Rydberg states. *Journal of Chemical Physics*, **2001**, 114, 7859-7865	3.9	15
122	Detection of OH(X,v□□□via the B 2□X 2□Transition and Properties of the B 2□ State. *Journal of Physical Chemistry A*, **2001**, 105, 6030-6037	2.8	5
121	Observation of Coriolis Coupling between nu(2) + 4nu(4) and 7nu(4) in Acetylene &Xtilde;(1)Sigma(+)(g) by Stimulated Emission Pumping Spectroscopy. *Journal of Molecular Spectroscopy*, **2000**, 199, 265-274	1.3	24
120	Remarks on the signs of g factors in atomic and molecular Zeeman spectroscopy. *Molecular Physics*, **2000**, 98, 1597-1601	1.7	20
119	A statistical approach for the study of singlet□riplet interactions in small polyatomic molecules. *Journal of Chemical Physics*, **2000**, 113, 6640-6651	3.9	14

118	Rovibrational spectroscopy of the v=6 manifold in 12C2H2 and 13C2H2. *Journal of Chemical Physics*, **2000**, 113, 7376-7383	3.9	21
117	Acetylene at the Threshold of Isomerization. *Journal of Physical Chemistry A*, **2000**, 104, 3073-3086	2.8	105
116	State-by-state assignment of the bending spectrum of acetylene at 15 000 cm□1: A case study of quantum-classical correspondence. *Journal of Chemical Physics*, **1999**, 111, 600-618	3.9	94
115	Toward a global and causal understanding of the unusual Rydberg state potential energy curves of the heteronuclear rare gas dimers. *Journal of Chemical Physics*, **1999**, 110, 10653-10656	3.9	25
114	Local mode behavior in the acetylene bending system. *Journal of Chemical Physics*, **1999**, 110, 845-859	3.9	67
113	The Predissociation Mechanism for 2Sigma+ Rydberg States of CaCl. *Journal of Molecular Spectroscopy*, **1999**, 193, 412-417	1.3	12
112	Diabatic analysis of the electronic states of hydrogen chloride. *Journal of Chemical Physics*, **1998**, 109, 8374-8387	3.9	27
111	A unified model of the dynamics and spectroscopy of the g 3□□and E 1□ states of hydrogen chloride. *Journal of Chemical Physics*, **1998**, 108, 984-989	3.9	3
110	Pure bending dynamics in the acetylene X 1□+ state up to 15 000 cm□1 of internal energy. *Journal of Chemical Physics*, **1998**, 109, 121-133	3.9	100
109	Numerical pattern recognition analysis of acetylene dispersed fluorescence spectra. *Journal of Chemical Physics*, **1998**, 108, 7100-7113	3.9	54
108	Anomalously slow intramolecular vibrational redistribution in the acetylene X 1□+ state above 10 000 cm□1 of internal energy. *Journal of Chemical Physics*, **1998**, 109, 3831-3840	3.9	43
107	Spectroscopic investigation of the generation of □somerization□states: Eigenvector analysis of the bend-CP stretch polyad. *Journal of Chemical Physics*, **1998**, 109, 492-503	3.9	33
106	The effects of triplet perturbers on photophysical processes in the □□1Au state of acetylene. *Journal of Chemical Physics*, **1997**, 106, 3423-3426	3.9	11
105	Observation of the □somerization states□of HCP by stimulated emission pumping spectroscopy: Comparison between theory and experiment. *Journal of Chemical Physics*, **1997**, 106, 2980-2983	3.9	31
104	Laser excited metastable states of acetylene in the 5.5□5.7 eV region. *Journal of Chemical Physics*, **1997**, 107, 49-53	3.9	24
103	Sideband optical□optical double resonance Zeeman spectroscopy. III. Analysis of composite lines and selective detection. *Journal of Chemical Physics*, **1997**, 107, 4189-4193	3.9	4
102	Sideband optical□optical double resonance Zeeman spectroscopy. II. Studies of NiH, PdD, and PtH. *Journal of Chemical Physics*, **1997**, 107, 4179-4188	3.9	12
101	Extended cross correlation: A technique for spectroscopic pattern recognition. *Journal of Chemical Physics*, **1997**, 107, 8349-8356	3.9	30

100	Realistic representation of the induced electric dipole moment of a polarizable ligand: The missing factor in the Rittner polarization model. *Journal of Chemical Physics*, **1997**, 106, 10379-10382	3.9	7
99	Rydberg series of BaF: peturbation facilitated studies of core ion penetrating states. *Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences*, **1997**, 355, 1507-1526	3	16
98	Vibrational Energy Flow in Highly Excited Molecules: Role of Intramolecular Vibrational Redistribution. *The Journal of Physical Chemistry*, **1996**, 100, 12735-12756		624
97	Detection of DCl by multiphoton ionization and determination of DCl and HCl internal state distributions. *Journal of Chemical Physics*, **1996**, 105, 10251-10262	3.9	14
96	Highly vibrationally excited 12C2H2 in the X 1 g state: Complementarity of absorption and dispersed fluorescence spectra. *Journal of Chemical Physics*, **1996**, 105, 11357-11359	3.9	58
95	Dispersed Fluorescence Spectrum of Acetylene from the 1 Au Origin: Recognition of Polyads and Test of Multiresonant Effective Hamiltonian Model for the X State. *The Journal of Physical Chemistry*, **1996**, 100, 7797-7809		85
94	Core-Penetrating Rydberg Series of BaF: New Electronic States in then* 4 Region. *Journal of Molecular Spectroscopy*, **1996**, 179, 99-124	1.3	14
93	Stimulated emission pumping spectroscopy of HCP near the isomerization barrier: EVIB 5 315 cm 1. *Journal of Chemical Physics*, **1996**, 105, 7383-7401	3.9	37
92	A Trans-Pacific Collaboration: Dynamical Experiments and Experiences. *Laser Chemistry*, **1995**, 15, 79-80		0
91	A multichannel quantum defect fit to the n*=6 8 core-penetrating s~p~d supercomplexes of CaF. *Journal of Chemical Physics*, **1995**, 103, 6459-6466	3.9	20
90	Electronic control of the spin orbit branching ratio in the photodissociation and predissociation of HCl. *Journal of Chemical Physics*, **1995**, 103, 6811-6814	3.9	92
89	The electronic structure of LaO: Ligand field versus ab initio calculations. *Journal of Chemical Physics*, **1995**, 103, 8004-8013	3.9	36
88	Sideband optical optical double resonance Zeeman spectroscopy. I. Theory of saturation and line shape behavior. *Journal of Chemical Physics*, **1995**, 102, 8295-8307	3.9	4
87	How robust are molecular properties? A stability criterion for eigenstates. *Journal of Chemical Physics*, **1995**, 102, 337-345	3.9	1
86	The Acetylene S0 Surface: From Dispersed Fluorescence Spectra to Polyads to Dynamics. *Zeitschrift Fur Elektrotechnik Und Elektrochemie*, **1995**, 99, 555-560		45
85	CONTINUOUS WAVE PERTURBATION-FACILITATED OPTICAL OPTICAL DOUBLE RESONANCE SPECTROSCOPY OF Na2 and Li2. *Advanced Series in Physical Chemistry*, **1995**, 251-277		1
84	New Scheme for Extracting Molecular Dynamics from Spectra: Case Study on Vibrationally Highly Excited Acetylene. *Laser Chemistry*, **1994**, 14, 183-190		10
83	Laser spectroscopic studies of terbium monofluoride: ligand field assignments of some f8[7F]sp-f8[7F]s2 transitions. *Molecular Physics*, **1994**, 83, 881-905	1.7	15

82	Frequency-modulation enhanced magnetic rotation spectroscopy of PdH, PdD, NiH, and CuH. *Journal of Chemical Physics*, **1994**, 100, 6347-6358	3.9	22
81	The 1 van der Waals and 0+ double well potentials of Xe 6s[3/2]01+Kr 1S0 determined from tunable vacuum ultraviolet laser spectroscopy. *Journal of Chemical Physics*, **1994**, 101, 10242-10251	3.9	16
80	Frequency-modulation enhanced magnetic rotation spectroscopy: A sensitive and selective absorption scheme for paramagnetic molecules. *Journal of Chemical Physics*, **1994**, 100, 6331-6346	3.9	29
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