

# Martin Quack

## List of Publications by Citations

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L-index

#	Paper	IF	Citations
299	Specific Rate Constants of Unimolecular Processes II. Adiabatic Channel Model. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1974</b> , 78, 240-252		483
298	Spectra and Dynamics of Coupled Vibrations in Polyatomic Molecules. <i>Annual Review of Physical Chemistry</i> , <b>1990</b> , 41, 839-874	15.7	405
297	How important is parity violation for molecular and biomolecular chirality?. <i>Angewandte Chemie - International Edition</i> , <b>2002</b> , 41, 4618-30	16.4	344
296	Detailed symmetry selection rules for reactive collisions. <i>Molecular Physics</i> , <b>1977</b> , 34, 477-504	1.7	282
295	Theory of unimolecular reactions induced by monochromatic infrared radiation. <i>Journal of Chemical Physics</i> , <b>1978</b> , 69, 1282	3.9	282
294	Potential energy surfaces, quasiadiabatic channels, rovibrational spectra, and intramolecular dynamics of (HF) <sub>2</sub> and its isotopomers from quantum Monte Carlo calculations. <i>Journal of Chemical Physics</i> , <b>1991</b> , 95, 28-59	3.9	260
293	Structure and Dynamics of Chiral Molecules. <i>Angewandte Chemie International Edition in English</i> , <b>1989</b> , 28, 571-586		254
292	Tridiagonal Fermi resonance structure in the IR spectrum of the excited CH chromophore in CF <sub>3</sub> H. <i>Journal of Chemical Physics</i> , <b>1984</b> , 81, 3779-3791	3.9	247
291	Anchoring the water dimer potential energy surface with explicitly correlated computations and focal point analyses. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 690-701	3.9	239
290	IR-laser photochemistry. <i>Chemical Reviews</i> , <b>1987</b> , 87, 181-216	68.1	233
289	Complex Formation in Reactive and Inelastic Scattering: Statistical Adiabatic Channel Model of Unimolecular Processes III. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1975</b> , 79, 170-183		232
288	High-resolution spectroscopic studies and theory of parity violation in chiral molecules. <i>Annual Review of Physical Chemistry</i> , <b>2008</b> , 59, 741-69	15.7	193
287	On the measurement of the parity violating energy difference between enantiomers. <i>Chemical Physics Letters</i> , <b>1986</b> , 132, 147-153	2.5	186
286	Vibrational spectrum and potential energy surface of the CH chromophore in CHD <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>1988</b> , 88, 5408-5432	3.9	183
285	A new six-dimensional analytical potential up to chemically significant energies for the electronic ground state of hydrogen peroxide. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 2565-2587	3.9	165
284	Unimolecular Processes V: Maximum Free Energy Criterion for the High Pressure Limit of Dissociation Reactions. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1977</b> , 81, 329-337		159
283	Resonance fluorescence of aniline vapour. <i>Journal of Molecular Spectroscopy</i> , <b>1972</b> , 43, 87-116	1.3	149

282	Global analysis of the high resolution infrared spectrum of methane $^{12}\text{CH}_4$ in the region from 0 to $4800\text{cm}^{-1}$ . <i>Chemical Physics</i> , <b>2009</b> , 356, 131-146	2.3	146
281	HF dimer: Empirically refined analytical potential energy and dipole hypersurfaces from ab initio calculations. <i>Journal of Chemical Physics</i> , <b>1998</b> , 108, 10096-10115	3.9	146
280	Ab initio calculation of molecular energies including parity violating interactions. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 7263-7285	3.9	140
279	Structure and dynamics of the excited $\text{CH}^+$ chromophore in $(\text{CF}_3)_3\text{CH}$ . <i>Journal of Chemical Physics</i> , <b>1985</b> , 82, 1186-1194	3.9	136
278	Tridiagonal Fermi resonance structure in the vibrational spectrum of the CH chromophore in $\text{CHF}_3$ . II. Visible spectra. <i>Journal of Chemical Physics</i> , <b>1987</b> , 86, 634-646	3.9	128
277	High-resolution interferometric Fourier transform infrared absorption spectroscopy in supersonic free jet expansions: carbon monoxide, nitric oxide, methane, ethyne, propyne, and trifluoromethane. <i>The Journal of Physical Chemistry</i> , <b>1988</b> , 92, 5455-5466		124
276	Vibrational spectra of $(\text{HF})_2$ , $(\text{HF})_n$ and their D-isotopomers: Mode selective rearrangements and nonstatistical unimolecular decay. <i>Chemical Physics</i> , <b>1989</b> , 139, 31-53	2.3	112
275	Methane line parameters in the HITRAN2012 database. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , <b>2013</b> , 130, 201-219	2.1	110
274	Influence of parity violating weak nuclear potentials on vibrational and rotational frequencies in chiral molecules. <i>Physical Review Letters</i> , <b>2000</b> , 84, 3807-10	7.4	103
273	Spectroscopy and dynamics of the isolated CH chromophore in $\text{CD}_3\text{H}$ : Experiment and theory. <i>Chemical Physics Letters</i> , <b>1984</b> , 109, 563-569	2.5	102
272	Multiconfiguration linear response approach to the calculation of parity violating potentials in polyatomic molecules. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 3148-3158	3.9	100
271	Dipole moment function and equilibrium structure of methane in an analytical, anharmonic nine-dimensional potential surface related to experimental rotational constants and transition moments by quantum Monte Carlo calculations. <i>Journal of Chemical Physics</i> , <b>1994</b> , 101, 3588-3602	3.9	100
270	Electroweak quantum chemistry of alanine: parity violation in gas and condensed phases. <i>ChemPhysChem</i> , <b>2000</b> , 1, 57-60	3.2	95
269	Mode selective stereomutation tunnelling in hydrogen peroxide isotopomers. <i>Chemical Physics Letters</i> , <b>1999</b> , 300, 312-320	2.5	94
268	The wave packet motion and intramolecular vibrational redistribution in $\text{CH}_3$ molecules under infrared multiphoton excitation. <i>Journal of Chemical Physics</i> , <b>1991</b> , 95, 4854-4876	3.9	94
267	High resolution interferometric FTIR spectroscopy of $(\text{HF})_2$ : analysis of a low frequency fundamental near $400\text{ cm}^{-1}$ . <i>Molecular Physics</i> , <b>1987</b> , 62, 1047-1064	1.7	94
266	The far infrared pure rotational spectrum and the Coriolis coupling between $\nu_3$ and $\nu_8$ in $\text{CH}_3\text{ClF}_2$ . <i>Molecular Physics</i> , <b>1989</b> , 68, 745-758	1.7	93
265	Vibrational overtone intensities of the isolated CH and CD chromophores in fluoroform and chloroform. <i>Chemical Physics Letters</i> , <b>1986</b> , 123, 197-202	2.5	90

- 264 Infrared laser chemistry and the dynamics of molecular multiphoton excitation. *Infrared Physics*, **1989**, 29, 441-466 87
- 263 Combined high resolution infrared and microwave study of bromochlorofluoromethane. *Journal of Chemical Physics*, **1997**, 106, 7558-7570 3.9 86
- 262 Vibrational spectrum, dipole moment function, and potential energy surface of the CH chromophore in CHX3 molecules. *Journal of Chemical Physics*, **1989**, 91, 6698-6713 3.9 84
- 261 Is the stereomutation of methane possible?. *Journal of Computational Chemistry*, **1995**, 16, 207-225 3.5 83
- 260 Unimolecular Processes IV: Product State Distributions after Dissociation. *Zeitschrift Fur Elektrochemie Und Elektrochemie*, **1975**, 79, 469-475 82
- 259 A critical analysis of electronic density functionals for structural, energetic, dynamic, and magnetic properties of hydrogen fluoride clusters. *Journal of Computational Chemistry*, **1997**, 18, 1695-1719 3.5 81
- 258 Molecular quantum dynamics from high resolution spectroscopy and laser chemistry. *Journal of Molecular Structure*, **1993**, 292, 171-195 3.4 81
- 257 Time-dependent processes in polyatomic molecules during and after intense infrared irradiation. *Faraday Discussions of the Chemical Society*, **1983**, 75, 197-210 80
- 256 Global analytical potential hypersurfaces for large amplitude nuclear motion and reactions in methane. I. Formulation of the potentials and adjustment of parameters to ab initio data and experimental constraints. *Journal of Chemical Physics*, **1998**, 109, 10628-10643 3.9 79
- 255 A new ab initio based six-dimensional semi-empirical pair interaction potential for HF. *Chemical Physics Letters*, **1996**, 261, 35-44 2.5 78
- 254 Reaction Dynamics and Statistical Mechanics of the Preparation of Highly Excited States by Intense Infrared Radiation. *Advances in Chemical Physics*, **2007**, 395-473 77
- 253 Global Analytical Potential Hypersurface for Large Amplitude Nuclear Motion and Reactions in Methane II. Characteristic Properties of the Potential and Comparison to Other Potentials and Experimental Information. *Journal of Physical Chemistry A*, **2004**, 108, 3166-3181 2.8 77
- 252 Molecular Spectroscopy and Molecular Dynamics: Theory and Experiment. *Zeitschrift Fur Elektrochemie Und Elektrochemie*, **1995**, 99, 231-245 75
- 251 Multiple anharmonic resonances in the vibrational overtone spectra of CHClF2. *Molecular Physics*, **1985**, 56, 727-735 1.7 72
- 250 High-resolution Fourier transform infrared and cw-diode laser cavity ringdown spectroscopy of the  $\nu_2 + 2\nu_3$  band of methane near 7510  $\text{cm}^{-1}$  in slit jet expansions and at room temperature. *Journal of Chemical Physics*, **2002**, 116, 6045-6055 3.9 70
- 249 Spectral bandshape and intensity of the C-H chromophore in the infrared spectra of CF3H and C4F9H. *Chemical Physics Letters*, **1980**, 72, 342-347 2.5 69
- 248 Current Aspects of Unimolecular Reactions. *International Reviews in Physical Chemistry*, **1981**, 1, 97-147 7 69
- 247 High-resolution cavity ring-down absorption spectroscopy of nitrous oxide and chloroform using a near-infrared cw diode laser. *Chemical Physics Letters*, **1998**, 289, 527-534 2.5 68

246	Observation and assignment of the hydrogen bond exchange disrotatory in-plane bending vibration $\nu_5$ in (HF) <sub>2</sub> . <i>Chemical Physics Letters</i> , <b>1990</b> , 171, 517-524	2.5	68
245	High resolution spectroscopy and the first global analysis of the Tetradecad region of methane 12CH <sub>4</sub> . <i>Physical Chemistry Chemical Physics</i> , <b>2013</b> , 15, 10071-93	3.6	66
244	Combined multidimensional anharmonic and parity violating effects in CDBrClF. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 11228-11240	3.9	66
243	Vibrational overtone spectra and vibrational dynamics of CFHCl <sub>2</sub> and (CH <sub>3</sub> ) <sub>2</sub> CFH. <i>Molecular Physics</i> , <b>1984</b> , 53, 257-264	1.7	64
242	Unimolecular reactions and energy transfer of highly excited molecules 175-238		64
241	Spectrum and dynamics of the CH chromophore in CD <sub>2</sub> HF. I. Vibrational Hamiltonian and analysis of rovibrational spectra. <i>Chemical Physics Letters</i> , <b>1992</b> , 190, 581-589	2.5	62
240	Accurate quantum Monte Carlo calculations of the tunneling splitting in (HF) <sub>2</sub> on a six-dimensional potential hypersurface. <i>Chemical Physics Letters</i> , <b>1995</b> , 234, 71-76	2.5	60
239	Potential energy surface and energy levels of (HF) <sub>2</sub> and its D isotopomers. <i>Molecular Physics</i> , <b>1990</b> , 69, 791-801	1.7	60
238	Observation and assignment of tunnelling-rotational transitions in the far infrared spectrum of (HF) <sub>2</sub> . <i>Molecular Physics</i> , <b>1988</b> , 65, 1025-1045	1.7	60
237	Master Equations for Photochemistry with Intense Infrared Light. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1979</b> , 83, 757-775		59
236	A spectroscopic determination of the methyl radical recombination rate constant in shock waves. <i>Chemical Physics Letters</i> , <b>1976</b> , 39, 304-309	2.5	59
235	Cw cavity ring-down infrared absorption spectroscopy in pulsed supersonic jets: nitrous oxide and methane. <i>Chemical Physics Letters</i> , <b>1999</b> , 314, 273-281	2.5	57
234	Intramolecular energy transfer and vibrational redistribution in chiral molecules: experiment and theory. <i>Faraday Discussions</i> , <b>1994</b> , 99, 49	3.6	57
233	On the densities and numbers of rovibronic states of a given symmetry species: Rigid and nonrigid molecules, transition states, and scattering channels. <i>Journal of Chemical Physics</i> , <b>1985</b> , 82, 3277-3283	3.9	57
232	Molecular chirality and the fundamental symmetries of physics: influence of parity violation on rovibrational frequencies and thermodynamic properties. <i>Chirality</i> , <b>2001</b> , 13, 745-53	2.1	56
231	Parity Violation Dominates the Dynamics of Chirality in Dichlorodisulfane. <i>Angewandte Chemie - International Edition</i> , <b>2001</b> , 40, 4195-4198	16.4	56
230	Inversion Tunneling in Aniline from High Resolution Infrared Spectroscopy and an Adiabatic Reaction Path Hamiltonian Approach. <i>Zeitschrift Fur Physikalische Chemie</i> , <b>1999</b> , 209, 1-19	3.1	56
229	Evidence for the (HF) <sub>5</sub> complex in the HF stretching FTIR absorption spectra of pulsed and continuous supersonic jet expansions of hydrogen fluoride. <i>Chemical Physics Letters</i> , <b>1993</b> , 208, 446-452	2.5	56

228	Intramolecular vibrational redistribution and unimolecular reaction: Concepts and new results on the femtosecond dynamics and statistics in CHBrClF. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1997</b> , 101, 311-328		55
227	Roadmap on STIRAP applications. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , <b>2019</b> , 52, 202001	1.3	54
226	Synchrotron-based highest resolution Fourier Transform infrared spectroscopy of naphthalene (C <sub>10</sub> H <sub>8</sub> ) and indole (C <sub>8</sub> H <sub>7</sub> N) and its application to astrophysical problems. <i>Faraday Discussions</i> , <b>2011</b> , 150, 71-99; discussion 113-60	3.6	54
225	High resolution infrared spectroscopy and global vibrational analysis for the CH <sub>3</sub> D and CHD <sub>3</sub> isotopomers of methane. <i>Molecular Physics</i> , <b>2010</b> , 108, 1209-1240	1.7	54
224	Representation of parity violating potentials in molecular main chiral axes. <i>Chemical Physics Letters</i> , <b>1999</b> , 303, 547-557	2.5	54
223	Femtosecond quantum dynamics of functional groups under coherent infrared multiphoton excitation as derived from the analysis of high-resolution spectra. <i>The Journal of Physical Chemistry</i> , <b>1993</b> , 97, 12574-12590		54
222	Vibrational dynamics of (HF) <sub>n</sub> aggregates from an ab initio based analytical (1+2+3)-body potential. <i>Journal of Molecular Structure</i> , <b>1993</b> , 294, 33-36	3.4	53
221	Overtone intensities and dipole moment surfaces for the isolated CH chromophore in CHD <sub>3</sub> and CHF <sub>3</sub> : Experiment and ab initio theory. <i>Journal of Chemical Physics</i> , <b>1990</b> , 93, 7097-7109	3.9	53
220	On FTIR Spectroscopy in Asynchronously Pulsed Supersonic Free Jet Expansions and on the Interpretation of Stretching Spectra of HF Clusters. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1995</b> , 99, 457-468		52
219	Fermi resonance structure and femtosecond quantum dynamics of a chiral molecule from the analysis of vibrational overtone spectra of CHBrClF. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1996</b> , 100, 1853-1875		52
218	Struktur und Dynamik chiraler Molek�le. <i>Angewandte Chemie</i> , <b>1989</b> , 101, 588-604	3.6	52
217	Statistical mechanics and dynamics of molecular fragmentation. <i>Societa Italiana Di Fisica Nuovo Cimento B-General Physics, Relativity Astronomy and Mathematical Physics and Methods</i> , <b>1981</b> , 63, 358-377		52
216	Survey of the high resolution infrared spectrum of methane (( <sup>12</sup> )CH <sub>4</sub> and ( <sup>13</sup> )CH <sub>4</sub> ): partial vibrational assignment extended towards 12,000 cm <sup>-1</sup> . <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 234302	3.9	51
215	Quantitative comparison between detailed (state selected) relative rate data and averaged (thermal) absolute rate data for complex forming reactions. <i>The Journal of Physical Chemistry</i> , <b>1979</b> , 83, 150-158		50
214	High-resolution IR spectrum of fluoroform: a close resonance. <i>Chemical Physics Letters</i> , <b>1981</b> , 80, 439-444	1.5	50
213	High-resolution rovibrational analysis of vibrational states of A <sub>2</sub> symmetry of the dideuterated methane CH <sub>2</sub> D <sub>2</sub> : the levels $\Gamma$ and $\Gamma + \Theta$ . <i>Molecular Physics</i> , <b>2006</b> , 104, 3371-3386	1.7	49
212	Wie wichtig ist Parit�tsverletzung f�r die molekulare und biomolekulare Chiralit�t?. <i>Angewandte Chemie</i> , <b>2002</b> , 114, 4812-4825	3.6	49
211	Group additivity and overtone intensities for the isolated CH chromophore. <i>Chemical Physics Letters</i> , <b>1984</b> , 112, 387-392	2.5	49



210	High resolution FTIR spectra of $^{12}\text{CF}_3\text{I}$ , $^{13}\text{CF}_3\text{I}$ and $^{12}\text{CF}_3\text{ } ^{79}\text{Br}$ near 1050 $\text{cm}^{-1}$ and 550 $\text{cm}^{-1}$ . <i>Molecular Physics</i> , <b>1985</b> , 55, 255-275	1.7	49
209	High-resolution near infrared spectroscopy and vibrational dynamics of dideuteromethane ( $\text{CH}_2\text{D}_2$ ). <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 2218-31	2.8	48
208	Ab initio calculations of mode selective tunneling dynamics in $^{12}\text{CH}_3\text{OH}$ and $^{13}\text{CH}_3\text{OH}$ . <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 5534-5544	3.9	48
207	The dependence of rate coefficients and product yields upon fluence, intensity, and time in unimolecular reactions induced by monochromatic infrared radiation. <i>Journal of Chemical Physics</i> , <b>1980</b> , 73, 247-255	3.9	48
206	High resolution Fourier transform spectroscopy of $\text{CH}_2\text{D}_2$ in the region 2350-2650 $\text{cm}^{-1}$ : the bands $\nu_5 + \nu_7$ , $2\nu_9$ , $\nu_3 + \nu_4$ , $\nu_3 + \nu_7$ and $\nu_5 + \nu_9$ . <i>Physical Chemistry Chemical Physics</i> , <b>2005</b> , 7, 1142-50	3.6	47
205	Molecular infrared spectra and molecular motion. <i>Journal of Molecular Structure</i> , <b>1995</b> , 347, 245-266	3.4	47
204	Stereomutation dynamics in hydrogen peroxide. <i>Chemical Physics</i> , <b>2007</b> , 338, 90-105	2.3	46
203	Mode Selective Stereomutation and Parity Violation in Disulfane Isotopomers $\text{H}_2\text{S}_2$ , $\text{D}_2\text{S}_2$ , $\text{T}_2\text{S}_2$ . <i>Helvetica Chimica Acta</i> , <b>2001</b> , 84, 1846-1861	2	46
202	High-resolution infrared spectrum and analysis of the $\nu_{11}$ , $A_{2u}(B_2)$ fundamental band of $^{12}\text{C}_6\text{H}_6$ and $^{13}\text{C}_6\text{H}_6$ . <i>Molecular Physics</i> , <b>1990</b> , 71, 759-768	1.7	46
201	Statistical Methods in Scattering <b>1981</b> , 199-276		46
200	Analysis of the rovibrational spectrum of $^{13}\text{CH}_4$ in the Octad range. <i>Journal of Molecular Spectroscopy</i> , <b>2013</b> , 291, 33-47	1.3	45
199	Absolute rate parameters for infrared photochemistry: $\text{CF}_3\text{I} + \text{I}$ . <i>Journal of Chemical Physics</i> , <b>1982</b> , 76, 955-965	3.9	45
198	Parity Violation in Chiral Molecules. <i>Chimia</i> , <b>2005</b> , 59, 530-538	1.3	44
197	Analytical three-body interaction potentials and hydrogen bond dynamics of hydrogen fluoride aggregates, $(\text{HF})_n$ , $n \leq 8$ . <i>Journal of Molecular Structure</i> , <b>2001</b> , 599, 381-425	3.4	44
196	Investigation of fermi resonances in $\text{CH}_3\text{X}$ molecules with an internal-coordinate hamiltonian. <i>Journal of the Chemical Society, Faraday Transactions 2</i> , <b>1988</b> , 84, 1371		44
195	Quantum-mechanical wavepacket dynamics of the CH group in symmetric top $\text{XCH}_3$ compounds using effective Hamiltonians from high-resolution spectroscopy. <i>Journal of the Chemical Society, Faraday Transactions 2</i> , <b>1986</b> , 82, 1173		44
194	FTIR spectroscopy of hydrogen fluoride clusters in synchronously pulsed supersonic jets. Isotopic isolation, substitution and 3-d condensation. <i>Chemical Physics Letters</i> , <b>1997</b> , 269, 29-38	2.5	43
193	Quasiadiabatic channels and effective transition-state barriers for the disrotatory in-plane hydrogen-bond exchange motion in $(\text{HF})_2$ . <i>Chemical Physics Letters</i> , <b>1991</b> , 183, 187-194	2.5	42

192	Molecular Spectra, Reaction Dynamics, Symmetries and Life. <i>Chimia</i> , <b>2003</b> , 57, 147-160	1.3	41
191	The NH and ND stretching fundamentals of 14ND2H. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 7893-7902	3.9	41
190	High-resolution FTIR spectroscopy of CHClF2 in a supersonic free jet expansion. <i>Chemical Physics Letters</i> , <b>1988</b> , 152, 275-280	2.5	41
189	Sequence structure in the high-resolution infrared spectrum of trifluoropropyne. <i>Chemical Physics Letters</i> , <b>1982</b> , 90, 370-374	2.5	41
188	Fermi resonance in CHX3: a hamiltonian in symmetrized curvilinear internal coordinates. <i>Chemical Physics Letters</i> , <b>1987</b> , 140, 512-519	2.5	40
187	On the validity of the quasidecoupled approximation for molecular infrared-multiphoton excitation. <i>Journal of Chemical Physics</i> , <b>1985</b> , 83, 3805-3812	3.9	40
186	Master equations for photochemistry with intense infrared light (III) the influence of molecular parameters in URIMIR. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , <b>1979</b> , 83, 1287-1293		40
185	Isotopic chirality and molecular parity violation. <i>Angewandte Chemie - International Edition</i> , <b>2005</b> , 44, 3623-6	16.4	39
184	Isotope selective overtone spectroscopy of CHCl3 by vibrationally assisted dissociation and photofragment ionization. <i>Journal of Chemical Physics</i> , <b>1996</b> , 104, 7426-7430	3.9	39
183	On the measurement of CP-violating energy differences in matter-antimatter enantiomers. <i>Chemical Physics Letters</i> , <b>1994</b> , 231, 421-428	2.5	39
182	Global analytical potential energy surface for large amplitude nuclear motions in ammonia. <i>Journal of Physical Chemistry B</i> , <b>2005</b> , 109, 8439-51	3.4	38
181	The permanent electric dipole moment of CH2D2: FIR el spectroscopy. <i>Molecular Physics</i> , <b>1996</b> , 89, 297-313	1.7	38
180	Synthesis, Structure, High-Resolution Spectroscopy, and Laser Chemistry of Fluorooxirane and 2,2-[2H2]-Fluorooxirane. <i>Angewandte Chemie International Edition in English</i> , <b>1997</b> , 36, 140-143		37
179	Global Analysis of the Infrared Spectrum of 13CH4: Lines in the Region 0 to 3200 cm <sup>-1</sup> . <i>Chimia</i> , <b>2008</b> , 62, 273-276	1.3	37
178	High resolution rovibrational spectroscopy of chiral and aromatic compounds. <i>ChemPhysChem</i> , <b>2007</b> , 8, 1271-81	3.2	37
177	Tunneling dynamics of the NH chromophore in NHD2 during and after coherent infrared excitation. <i>Journal of Chemical Physics</i> , <b>2003</b> , 118, 643-658	3.9	37
176	Infrared spectrum and dynamics of the hydrogen bonded dimer (HF)2. <i>Infrared Physics</i> , <b>1989</b> , 29, 535-539		37
175	High resolution interferometric Fourier transform infrared absorption spectroscopy in a supersonic free jet expansion. <i>Molecular Physics</i> , <b>1987</b> , 60, 237-248	1.7	37



174	Isotopomer-selective overtone spectroscopy by ionization detected IR+UV double resonance of jet-cooled aniline. <i>Chemical Physics Letters</i> , <b>1998</b> , 298, 320-328	2.5	36
173	Rovibrational analysis of the $2\bar{B}$ , $3\bar{B}$ and $\bar{A}$ bands of CHCl <sub>2</sub> F measured at 170 and 298 K by high-resolution FTIR spectroscopy. <i>Molecular Physics</i> , <b>2007</b> , 105, 541-558	1.7	36
172	How do Parity Violating Weak Nuclear Interactions Influence Rovibrational Frequencies in Chiral Molecules?. <i>Zeitschrift Fur Physikalische Chemie</i> , <b>2000</b> , 214,	3.1	36
171	IR laser chemistry. <i>Infrared Physics and Technology</i> , <b>1995</b> , 36, 365-380	2.7	36
170	Fermi resonance in the overtone spectra of the CH chromophore in tribromomethane. 2. Visible spectra. <i>The Journal of Physical Chemistry</i> , <b>1991</b> , 95, 1201-1209		36
169	Infrared-multiphoton excitation and wave packet motion of the harmonic and anharmonic oscillators: Exact solutions and quasiresonant approximation. <i>Journal of Chemical Physics</i> , <b>1989</b> , 90, 6320-6327	2.9	36
168	On the possibility of mode-selective IR-multiphoton excitation of ozone. <i>Chemical Physics Letters</i> , <b>1984</b> , 105, 147-152	2.5	36
167	Global analytical potential energy surface for the electronic ground state of NH <sub>3</sub> from high level ab initio calculations. <i>Journal of Physical Chemistry A</i> , <b>2013</b> , 117, 7502-22	2.8	35
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9	Fixierte Konstanten. <i>Nachrichten Aus Der Chemie</i> , <b>2015</b> , 63, 515-521	0.1	1
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