Cho Minhaeng

List of Publications by Citations

310 papers	13,345 citations	56 h-index	103 g-index
335 ext. papers	14,496 ext. citations	6.1 avg, IF	6.8 L-index

#	Paper	IF	Citations
310	Two-dimensional spectroscopy of electronic couplings in photosynthesis. *Nature*, **2005**, 434, 625-8	50.4	992
309	Coherent two-dimensional optical spectroscopy. *Chemical Reviews*, **2008**, 108, 1331-418	68.1	621
308	CHROMOPHORE-SOLVENT DYNAMICS. *Annual Review of Physical Chemistry*, **1996**, 47, 109-134	15.7	579
307	Exciton analysis in 2D electronic spectroscopy. *Journal of Physical Chemistry B*, **2005**, 109, 10542-56	3.4	360
306	Instantaneous normal mode analysis of liquid water. *Journal of Chemical Physics*, **1994**, 100, 6672-6683	3.9	315
305	Two-Dimensional Optical Spectroscopy		314
304	Two photon absorption properties of 1,3,5-tricyano-2,4,6-tris(styryl)benzene derivatives. *Journal of the American Chemical Society*, **2001**, 123, 10039-45	16.4	289
303	The Integrated Photon Echo and Solvation Dynamics. *The Journal of Physical Chemistry*, **1996**, 100, 11944-11953		248
302	Correlation between electronic and molecular structure distortions and vibrational properties. II. Amide I modes of NMA–D2O complexes. *Journal of Chemical Physics*, **2003**, 118, 3491-3498	3.9	231
301	Off-resonant transient birefringence in liquids. *Journal of Chemical Physics*, **1993**, 99, 2410-2428	3.9	231
300	Amide I vibrational dynamics of N-methylacetamide in polar solvents: the role of electrostatic interactions. *Journal of Physical Chemistry B*, **2005**, 109, 11016-26	3.4	201
299	Molecular dynamics simulation study of N-methylacetamide in water. I. Amide I mode frequency fluctuation. *Journal of Chemical Physics*, **2003**, 119, 2247-2255	3.9	193
298	Ultrafast solvent dynamics: Connection between time resolved fluorescence and optical Kerr measurements. *Journal of Chemical Physics*, **1992**, 96, 5033-5038	3.9	191
297	Two-photon absorption and nonlinear optical properties of octupolar molecules. *Journal of the American Chemical Society*, **2001**, 123, 10658-67	16.4	177
296	Infrared probes for studying the structure and dynamics of biomolecules. *Chemical Reviews*, **2013**, 113, 5817-47	68.1	171
295	Molecular dynamics simulation study of N-methylacetamide in water. II. Two-dimensional infrared pump–probe spectra. *Journal of Chemical Physics*, **2003**, 119, 2256-2263	3.9	170
294	Nitrile and thiocyanate IR probes: quantum chemistry calculation studies and multivariate least-square fitting analysis. *Journal of Chemical Physics*, **2008**, 128, 134506	3.9	154

293	The short-time dynamics of solvation. *Journal of Chemical Physics*, **1994**, 100, 6700-6708	3.9	141
292	Femtosecond characterization of vibrational optical activity of chiral molecules. *Nature*, **2009**, 458, 310-3	50.4	134
291	Photon echoes and related four-wave-mixing spectroscopies using phase-locked pulses. *Journal of Chemical Physics*, **1992**, 96, 5618-5629	3.9	134
290	Local Amide I Mode Frequencies and Coupling Constants in Polypeptides. *Journal of Physical Chemistry B*, **2003**, 107, 9132-9138	3.4	131
289	Non-Gaussian statistics of amide I mode frequency fluctuation of N-methylacetamide in methanol solution: linear and nonlinear vibrational spectra. *Journal of Chemical Physics*, **2004**, 120, 1477-90	3.9	124
288	Amide I modes in the N-methylacetamide dimer and glycine dipeptide analog: Diagonal force constants. *Journal of Chemical Physics*, **2003**, 118, 6915-6922	3.9	124
287	Amide I modes of tripeptides: Hessian matrix reconstruction and isotope effects. *Journal of Chemical Physics*, **2003**, 119, 1451-1461	3.9	116
286	Nitrile and thiocyanate IR probes: molecular dynamics simulation studies. *Journal of Chemical Physics*, **2008**, 128, 154504	3.9	113
285	Nonequilibrium photoinduced electron transfer. *Journal of Chemical Physics*, **1995**, 103, 595-606	3.9	112
284	Beta-azidoalanine as an IR probe: application to amyloid Abeta(16-22) aggregation. *Journal of Physical Chemistry B*, **2008**, 112, 10352-7	3.4	97
283	Nonlinear optical and two-photon absorbtion properties of 1,3,5-tricyano-2,4,6-tris(styryl)benzene-containing octupolar oligomers. *Chemistry - A European Journal*, **2002**, 8, 3907-16	4.8	93
282	Simulation studies of amide I IR absorption and two-dimensional IR spectra of beta hairpins in liquid water. *Journal of Physical Chemistry B*, **2005**, 109, 11789-801	3.4	93
281	Nanometric Water Channels in Water-in-Salt Lithium Ion Battery Electrolyte. *Journal of the American Chemical Society*, **2018**, 140, 15661-15667	16.4	88
280	Vibrational Spectroscopic Map, Vibrational Spectroscopy, and Intermolecular Interaction. *Chemical Reviews*, **2020**, 120, 7152-7218	68.1	87
279	Vibrational Probes: From Small Molecule Solvatochromism Theory and Experiments to Applications in Complex Systems. *Accounts of Chemical Research*, **2017**, 50, 968-976	24.3	81
278	Local Amide I Mode Frequencies and Coupling Constants in Multiple-Stranded Antiparallel β-Sheet Polypeptides. *Journal of Physical Chemistry B*, **2004**, 108, 20397-20407	3.4	79
277	Computational vibrational spectroscopy of peptides and proteins in one and two dimensions. *Accounts of Chemical Research*, **2009**, 42, 1280-9	24.3	76
276	Classical and quantum mechanical/molecular mechanical molecular dynamics simulations of alanine dipeptide in water: comparisons with IR and vibrational circular dichroism spectra. *Journal of Chemical Physics*, **2008**, 128, 105106	3.9	75

275	Amide I Modes of Helical Polypeptide in Liquid Water: Conformational Fluctuation, Phase Correlation, and Linear and Nonlinear Vibrational Spectra. *Journal of Physical Chemistry B*, **2004**, 108, 9333-9345	3.4	74
274	Vibrational solvatochromism and electrochromism of infrared probe molecules containing C?O, C?N, C=O, or C-F vibrational chromophore. *Journal of Chemical Physics*, **2011**, 134, 154513	3.9	73
273	Two-Color Pump Probe Spectroscopies of Two- and Three-Level Systems: 2-Dimensional Line Shapes and Solvation Dynamics. *Journal of Physical Chemistry A*, **2003**, 107, 5903-5912	2.8	72
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271	Computational spectroscopy of ubiquitin: comparison between theory and experiments. *Journal of Chemical Physics*, **2007**, 126, 045102	3.9	70
270	Azido-derivatized compounds as IR probes of local electrostatic environment: Theoretical studies. *Journal of Chemical Physics*, **2008**, 129, 174512	3.9	69
269	Amide I vibrational modes in glycine dipeptide analog: Ab initio calculation studies. *Journal of Chemical Physics*, **2002**, 117, 740-750	3.9	69
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267	Molecular Polarizability and First Hyperpolarizability of Octupolar Molecules: Donor-Substituted Triphenylmethane Dyes. *Journal of the American Chemical Society*, **1998**, 120, 10921-10927	16.4	67
266	Vibrational solvatochromism and electrochromism: coarse-grained models and their relationships. *Journal of Chemical Physics*, **2009**, 130, 094505	3.9	66
265	Octupolar Crystals for Nonlinear Optics: 1,3,5-Trinitro-2,4,6-tris(styryl)benzene Derivatives. *Chemistry of Materials*, **2001**, 13, 1438-1440	9.6	66
264	Coherent two-dimensional Raman scattering: Frequency-domain measurement of the intra- and intermolecular vibrational interactions. *Journal of Chemical Physics*, **1998**, 108, 1326-1334	3.9	66
263	Amide I IR, VCD, and 2d IR spectra of isotope-labeled helix in liquid water: Numerical simulation studies. *International Journal of Quantum Chemistry*, **2005**, 104, 616-634	2.1	65
262	Inter-peptide interaction and delocalization of amide I vibrational excitons in myoglobin and flavodoxin. *Journal of Chemical Physics*, **2002**, 117, 6821-6832	3.9	63
261	Correlation between electronic and molecular structure distortions and vibrational properties. I. Adiabatic approximations. *Journal of Chemical Physics*, **2003**, 118, 3480-3490	3.9	62
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259	Amide I vibrational circular dichroism of dipeptide: Conformation dependence and fragment analysis. *Journal of Chemical Physics*, **2004**, 120, 4383-92	3.9	60
258	Vibrational solvatochromism of nitrile infrared probes: beyond the vibrational Stark dipole approach. *Physical Chemistry Chemical Physics*, **2016**, 18, 18094-111	3.6	59

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252	Two-Photon Absorption and Second Hyperpolarizability of the Linear Quadrupolar Molecule. *Journal of Physical Chemistry A*, **2000**, 104, 11033-11040	2.8	55
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