

Carlos R Baiz

List of Publications by Year in descending order

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Version: 2024-02-01

66
papers

1,872
citations

257101

24
h-index

276539

41
g-index

69
all docs

69
docs citations

69
times ranked

1696
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Interfacial Dynamics in Inverted-Headgroup Lipid Membranes. <i>Journal of Chemical Physics</i> , 2022, 156, 075102. | 1.2 | 3 |
| 2 | Dynamic effect of polymers at the surfactant-water interface: an ultrafast study. <i>Soft Matter</i> , 2022, 18, 1793-1800. | 1.2 | 4 |
| 3 | Origin of thiocyanate spectral shifts in water and organic solvents. <i>Journal of Chemical Physics</i> , 2022, 156, 104106. | 1.2 | 6 |
| 4 | Rapid and Sequential Dual Oxime Ligation Enables De Novo Formation of Functional Synthetic Membranes from Water-Soluble Precursors. <i>Angewandte Chemie - International Edition</i> , 2022, 61, . | 7.2 | 4 |
| 5 | Generative Adversarial Neural Networks for Denoising Coherent Multidimensional Spectra. <i>Journal of Physical Chemistry A</i> , 2022, 126, 3816-3825. | 1.1 | 8 |
| 6 | Lanthanide-dependent coordination interactions in lanmodulin: a 2D IR and molecular dynamics simulations study. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 21690-21700. | 1.3 | 8 |
| 7 | Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry C</i> , 2021, 125, 4342-4342. | 1.5 | 0 |
| 8 | Interfacial Dynamics in Lipid Membranes: The Effects of Headgroup Structures. <i>Journal of Physical Chemistry B</i> , 2021, 125, 1343-1350. | 1.2 | 23 |
| 9 | Bursting the bubble: A molecular understanding of surfactant-water interfaces. <i>Journal of Chemical Physics</i> , 2021, 154, 170901. | 1.2 | 7 |
| 10 | Short- and long-range crowding effects on water's hydrogen bond networks. <i>Cell Reports Physical Science</i> , 2021, 2, 100419. | 2.8 | 15 |
| 11 | Pump Slice Amplitudes: A Simple and Robust Method for Connecting Two-Dimensional Infrared and Fourier Transform Infrared Spectra. <i>Journal of Physical Chemistry A</i> , 2021, 125, 6498-6504. | 1.1 | 15 |
| 12 | Molecular Mechanism of Cell Membrane Protection by Sugars: A Study of Interfacial H-Bond Networks. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 9602-9607. | 2.1 | 17 |
| 13 | Proton-modulated interactions of ions with transport sites of prokaryotic and eukaryotic NCX prototypes. <i>Cell Calcium</i> , 2021, 99, 102476. | 1.1 | 2 |
| 14 | Infrared spectroscopy probes ion binding geometries. <i>Methods in Enzymology</i> , 2021, 651, 157-191. | 0.4 | 5 |
| 15 | Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry A</i> , 2021, 125, 1680-1680. | 1.1 | 0 |
| 16 | Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry B</i> , 2021, 125, 1973-1973. | 1.2 | 0 |
| 17 | Fast Dynamics of Lipid Mixtures Investigated with Vibrational Spectroscopy. <i>Biophysical Journal</i> , 2020, 118, 85a. | 0.2 | 0 |
| 18 | Ultrafast Dynamics at Lipid-Water Interfaces. <i>Accounts of Chemical Research</i> , 2020, 53, 1860-1868. | 7.6 | 30 |

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|----|--|------|-----------|
| 19 | Ions Slow Water Dynamics at Nonionic Surfactant Interfaces. <i>Journal of Physical Chemistry B</i> , 2020, 124, 11895-11900. | 1.2 | 8 |
| 20 | Ultrafast Spectroscopy of Lipid-Water Interfaces: Transmembrane Crowding Drives H-Bond Dynamics. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 4093-4098. | 2.1 | 22 |
| 21 | Ultrafast Dynamics at the Lipid-Water Interface: DMSO Modulates H-Bond Lifetimes. <i>Langmuir</i> , 2020, 36, 6502-6511. | 1.6 | 27 |
| 22 | Molecular heterogeneity in aqueous cosolvent systems. <i>Journal of Chemical Physics</i> , 2020, 152, 190901. | 1.2 | 17 |
| 23 | Phase Transition in a Heterogeneous Membrane: Atomically Detailed Picture. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 5263-5267. | 2.1 | 5 |
| 24 | Vibrational Spectroscopic Map, Vibrational Spectroscopy, and Intermolecular Interaction. <i>Chemical Reviews</i> , 2020, 120, 7152-7218. | 23.0 | 205 |
| 25 | Calcium-Lipid Interactions Observed with Isotope-Edited Infrared Spectroscopy. <i>Biophysical Journal</i> , 2020, 118, 2694-2702. | 0.2 | 9 |
| 26 | Liquid-Liquid Phase Separation Produces Fast H-Bond Dynamics in DMSO-Water Mixtures. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 1903-1908. | 2.1 | 28 |
| 27 | Spatial Control of the Self-assembled Block Copolymer Domain Orientation and Alignment on Photopatterned Surfaces. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 23399-23409. | 4.0 | 7 |
| 28 | Slow Oil, Slow Water: Long-Range Dynamic Coupling across a Liquid-Liquid Interface. <i>Journal of the American Chemical Society</i> , 2020, 142, 8063-8067. | 6.6 | 18 |
| 29 | Interactive Tools for Teaching Fourier Transforms. <i>The Biophysicist</i> , 2020, 1, . | 0.1 | 2 |
| 30 | Interfacial H-Bond Dynamics in Reverse Micelles: The Role of Surfactant Heterogeneity. <i>Langmuir</i> , 2019, 35, 11463-11470. | 1.6 | 23 |
| 31 | Non-Additive Effects of Binding Site Mutations in Calmodulin. <i>Biochemistry</i> , 2019, 58, 2730-2739. | 1.2 | 10 |
| 32 | Site-Specific Peptide Probes Detect Buried Water in a Lipid Membrane. <i>Biophysical Journal</i> , 2019, 116, 1692-1700. | 0.2 | 13 |
| 33 | Empirical S=O stretch vibrational frequency map. <i>Journal of Chemical Physics</i> , 2019, 151, 234107. | 1.2 | 16 |
| 34 | Ultrafast pH-jump two-dimensional infrared spectroscopy. <i>Optics Letters</i> , 2019, 44, 4937. | 1.7 | 5 |
| 35 | Coordination to lanthanide ions distorts binding site conformation in calmodulin. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, E3126-E3134. | 3.3 | 90 |
| 36 | Physiological Calcium Concentrations Slow Dynamics at the Lipid-Water Interface. <i>Biophysical Journal</i> , 2018, 115, 1541-1551. | 0.2 | 30 |

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|----|--|-----|-----------|
| 37 | Vibrational Relaxation in EDTA Is Ion-Dependent. <i>Journal of Physical Chemistry A</i> , 2018, 122, 6585-6592. | 1.1 | 11 |
| 38 | Crowding Stabilizes DMSOâ€“Water Hydrogen-Bonding Interactions. <i>Journal of Physical Chemistry B</i> , 2018, 122, 5984-5990. | 1.2 | 37 |
| 39 | Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 11375-11379. | 7.2 | 94 |
| 40 | Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures. <i>Angewandte Chemie</i> , 2017, 129, 11533-11537. | 1.6 | 25 |
| 41 | Titelbild: Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures (Angew.) Tj ETQq1 1 0.784314 rgBT /Overlo | 1.6 | 0 |
| 42 | An Empirical IR Frequency Map for Ester Câ•O Stretching Vibrations. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3888-3896. | 1.1 | 54 |
| 43 | Studying Proteinâ€“Protein Binding through T-Jump Induced Dissociation: Transient 2D IR Spectroscopy of Insulin Dimer. <i>Journal of Physical Chemistry B</i> , 2016, 120, 5134-5145. | 1.2 | 42 |
| 44 | Structural Disorder of Folded Proteins: Isotope-Edited 2D IR Spectroscopy and Markov State Modeling. <i>Biophysical Journal</i> , 2015, 108, 1747-1757. | 0.2 | 23 |
| 45 | Visualizing KcsA Conformational Changes upon Ion Binding by Infrared Spectroscopy and Atomistic Modeling. <i>Journal of Physical Chemistry B</i> , 2015, 119, 5824-5831. | 1.2 | 25 |
| 46 | Ultrafast 2D IR microscopy. <i>Optics Express</i> , 2014, 22, 18724. | 1.7 | 69 |
| 47 | A Molecular Interpretation of 2D IR Protein Folding Experiments with Markov State Models. <i>Biophysical Journal</i> , 2014, 106, 1359-1370. | 0.2 | 48 |
| 48 | Amide I Two-Dimensional Infrared Spectroscopy: Methods for Visualizing the Vibrational Structure of Large Proteins. <i>Journal of Physical Chemistry A</i> , 2013, 117, 5955-5961. | 1.1 | 29 |
| 49 | Direct observation of ground-state lactamâ€“lactim tautomerization using temperature-jump transient 2D IR spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013, 110, 9243-9248. | 3.3 | 50 |
| 50 | Ultrafast equilibrium and non-equilibrium chemical reaction dynamics probed with multidimensional infrared spectroscopy. <i>International Reviews in Physical Chemistry</i> , 2012, 31, 367-419. | 0.9 | 34 |
| 51 | Coherent two-dimensional infrared spectroscopy: Quantitative analysis of protein secondary structure in solution. <i>Analyst, The</i> , 2012, 137, 1793. | 1.7 | 65 |
| 52 | Local-Mode Approach to Modeling Multidimensional Infrared Spectra of Metal Carbonyls. <i>Journal of Physical Chemistry A</i> , 2011, 115, 5354-5363. | 1.1 | 24 |
| 53 | Molecular Theory and Simulation of Coherence Transfer in Metal Carbonyls and Its Signature on Multidimensional Infrared Spectra. <i>Journal of Physical Chemistry B</i> , 2011, 115, 5322-5339. | 1.2 | 38 |
| 54 | Ultrabroadband detection of a mid-IR continuum by chirped-pulse upconversion. <i>Optics Letters</i> , 2011, 36, 187. | 1.7 | 99 |

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|----|---|-----|-----------|
| 55 | Ultrafast Vibrational Stark-Effect Spectroscopy: Exploring Charge-Transfer Reactions by Directly Monitoring the Solvation Shell Response. <i>Journal of the American Chemical Society</i> , 2010, 132, 12784-12785. | 6.6 | 27 |
| 56 | Solvent-Dependent Spectral Diffusion in a Hydrogen Bonded "Vibrational Aggregate". <i>Journal of Physical Chemistry A</i> , 2010, 114, 10590-10604. | 1.1 | 67 |
| 57 | Transient Vibrational Echo versus Transient Absorption Spectroscopy: A Direct Experimental and Theoretical Comparison. <i>Applied Spectroscopy</i> , 2010, 64, 1037-1044. | 1.2 | 5 |
| 58 | Two-Dimensional Infrared Spectroscopy of Dimanganese Decacarbonyl and Its Photoproducts: An Ab Initio Study. <i>Journal of Physical Chemistry A</i> , 2009, 113, 9617-9623. | 1.1 | 21 |
| 59 | Orientalional Dynamics of Transient Molecules Measured by Nonequilibrium Two-Dimensional Infrared Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2009, 113, 8907-8916. | 1.1 | 29 |
| 60 | Beyond 7-Azaindole: Conjugation Effects on Intermolecular Double Hydrogen-Atom Transfer Reactions. <i>Journal of Physical Chemistry A</i> , 2009, 113, 4862-4867. | 1.1 | 16 |
| 61 | Structurally Selective Geminate Rebinding Dynamics of Solvent-Caged Radicals Studied with Nonequilibrium Infrared Echo Spectroscopy. <i>Journal of the American Chemical Society</i> , 2009, 131, 13590-13591. | 6.6 | 32 |
| 62 | Two-Dimensional Infrared Spectroscopy of Metal Carbonyls. <i>Accounts of Chemical Research</i> , 2009, 42, 1395-1404. | 7.6 | 98 |
| 63 | Ultrafast nonequilibrium Fourier-transform two-dimensional infrared spectroscopy. <i>Optics Letters</i> , 2008, 33, 2533. | 1.7 | 50 |
| 64 | Multilevel vibrational coherence transfer and wavepacket dynamics probed with multidimensional IR spectroscopy. <i>Journal of Chemical Physics</i> , 2008, 129, 084503. | 1.2 | 67 |
| 65 | Theoretical Studies of Conjugation Effects on Excited State Intramolecular Hydrogen-Atom Transfer Reactions in Model Systems. <i>Journal of Physical Chemistry A</i> , 2007, 111, 10139-10143. | 1.1 | 11 |
| 66 | Rapid and Sequential Dual Oxime Ligation Enables De Novo Formation of Functional Synthetic Membranes from Water-Soluble Precursors. <i>Angewandte Chemie</i> , 0, , . | 1.6 | 0 |