

Carlos R Baiz

List of Publications by Year in descending order

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66
papers

1,872
citations

236925

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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.	3.0	3
2	Dynamic effect of polymers at the surfactant-water interface: an ultrafast study. <i>Soft Matter</i> , 2022, 18, 1793-1800.	2.7	4
3	Origin of thiocyanate spectral shifts in water and organic solvents. <i>Journal of Chemical Physics</i> , 2022, 156, 104106.	3.0	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, .	13.8	4
5	Generative Adversarial Neural Networks for Denoising Coherent Multidimensional Spectra. <i>Journal of Physical Chemistry A</i> , 2022, 126, 3816-3825.	2.5	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.	2.8	8
7	Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry C</i> , 2021, 125, 4342-4342.	3.1	0
8	Interfacial Dynamics in Lipid Membranes: The Effects of Headgroup Structures. <i>Journal of Physical Chemistry B</i> , 2021, 125, 1343-1350.	2.6	23
9	Bursting the bubble: A molecular understanding of surfactant-water interfaces. <i>Journal of Chemical Physics</i> , 2021, 154, 170901.	3.0	7
10	Short- and long-range crowding effects on water's hydrogen bond networks. <i>Cell Reports Physical Science</i> , 2021, 2, 100419.	5.6	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.	2.5	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.	4.6	17
13	Proton-modulated interactions of ions with transport sites of prokaryotic and eukaryotic NCX prototypes. <i>Cell Calcium</i> , 2021, 99, 102476.	2.4	2
14	Infrared spectroscopy probes ion binding geometries. <i>Methods in Enzymology</i> , 2021, 651, 157-191.	1.0	5
15	Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry A</i> , 2021, 125, 1680-1680.	2.5	0
16	Viewpoint on ACS PHYS Division Sponsored Virtual Seminars. <i>Journal of Physical Chemistry B</i> , 2021, 125, 1973-1973.	2.6	0
17	Fast Dynamics of Lipid Mixtures Investigated with Vibrational Spectroscopy. <i>Biophysical Journal</i> , 2020, 118, 85a.	0.5	0
18	Ultrafast Dynamics at Lipid-Water Interfaces. <i>Accounts of Chemical Research</i> , 2020, 53, 1860-1868.	15.6	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.	2.6	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.	4.6	22
21	Ultrafast Dynamics at the Lipid-Water Interface: DMSO Modulates H-Bond Lifetimes. <i>Langmuir</i> , 2020, 36, 6502-6511.	3.5	27
22	Molecular heterogeneity in aqueous cosolvent systems. <i>Journal of Chemical Physics</i> , 2020, 152, 190901.	3.0	17
23	Phase Transition in a Heterogeneous Membrane: Atomically Detailed Picture. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 5263-5267.	4.6	5
24	Vibrational Spectroscopic Map, Vibrational Spectroscopy, and Intermolecular Interaction. <i>Chemical Reviews</i> , 2020, 120, 7152-7218.	47.7	205
25	Calcium-Lipid Interactions Observed with Isotope-Edited Infrared Spectroscopy. <i>Biophysical Journal</i> , 2020, 118, 2694-2702.	0.5	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.	4.6	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.	8.0	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.	13.7	18
29	Interactive Tools for Teaching Fourier Transforms. <i>The Biophysicist</i> , 2020, 1, .	0.3	2
30	Interfacial H-Bond Dynamics in Reverse Micelles: The Role of Surfactant Heterogeneity. <i>Langmuir</i> , 2019, 35, 11463-11470.	3.5	23
31	Non-Additive Effects of Binding Site Mutations in Calmodulin. <i>Biochemistry</i> , 2019, 58, 2730-2739.	2.5	10
32	Site-Specific Peptide Probes Detect Buried Water in a Lipid Membrane. <i>Biophysical Journal</i> , 2019, 116, 1692-1700.	0.5	13
33	Empirical S=O stretch vibrational frequency map. <i>Journal of Chemical Physics</i> , 2019, 151, 234107.	3.0	16
34	Ultrafast pH-jump two-dimensional infrared spectroscopy. <i>Optics Letters</i> , 2019, 44, 4937.	3.3	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.	7.1	90
36	Physiological Calcium Concentrations Slow Dynamics at the Lipid-Water Interface. <i>Biophysical Journal</i> , 2018, 115, 1541-1551.	0.5	30

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37	Vibrational Relaxation in EDTA Is Ion-Dependent. <i>Journal of Physical Chemistry A</i> , 2018, 122, 6585-6592.	2.5	11
38	Crowding Stabilizes DMSOâ€“Water Hydrogen-Bonding Interactions. <i>Journal of Physical Chemistry B</i> , 2018, 122, 5984-5990.	2.6	37
39	Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 11375-11379.	13.8	94
40	Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures. <i>Angewandte Chemie</i> , 2017, 129, 11533-11537.	2.0	25
41	Titelbild: Quantifying Hydrogenâ€“Bond Populations in Dimethyl Sulfoxide/Water Mixtures (Angew.) Tj ETQq1 1 0.784314 rgBT /Overlo	2.0	0
42	An Empirical IR Frequency Map for Ester Câ•O Stretching Vibrations. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3888-3896.	2.5	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.	2.6	42
44	Structural Disorder of Folded Proteins: Isotope-Edited 2D IR Spectroscopy and Markov State Modeling. <i>Biophysical Journal</i> , 2015, 108, 1747-1757.	0.5	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.	2.6	25
46	Ultrafast 2D IR microscopy. <i>Optics Express</i> , 2014, 22, 18724.	3.4	69
47	A Molecular Interpretation of 2D IR Protein Folding Experiments with Markov State Models. <i>Biophysical Journal</i> , 2014, 106, 1359-1370.	0.5	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.	2.5	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.	7.1	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.	2.3	34
51	Coherent two-dimensional infrared spectroscopy: Quantitative analysis of protein secondary structure in solution. <i>Analyst, The</i> , 2012, 137, 1793.	3.5	65
52	Local-Mode Approach to Modeling Multidimensional Infrared Spectra of Metal Carbonyls. <i>Journal of Physical Chemistry A</i> , 2011, 115, 5354-5363.	2.5	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.	2.6	38
54	Ultrabroadband detection of a mid-IR continuum by chirped-pulse upconversion. <i>Optics Letters</i> , 2011, 36, 187.	3.3	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.	13.7	27
56	Solvent-Dependent Spectral Diffusion in a Hydrogen Bonded α -Vibrational Aggregate. <i>Journal of Physical Chemistry A</i> , 2010, 114, 10590-10604.	2.5	67
57	Transient Vibrational Echo versus Transient Absorption Spectroscopy: A Direct Experimental and Theoretical Comparison. <i>Applied Spectroscopy</i> , 2010, 64, 1037-1044.	2.2	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.	2.5	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.	2.5	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.	2.5	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.	13.7	32
62	Two-Dimensional Infrared Spectroscopy of Metal Carbonyls. <i>Accounts of Chemical Research</i> , 2009, 42, 1395-1404.	15.6	98
63	Ultrafast nonequilibrium Fourier-transform two-dimensional infrared spectroscopy. <i>Optics Letters</i> , 2008, 33, 2533.	3.3	50
64	Multilevel vibrational coherence transfer and wavepacket dynamics probed with multidimensional IR spectroscopy. <i>Journal of Chemical Physics</i> , 2008, 129, 084503.	3.0	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.	2.5	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, , .	2.0	0