

Michael D Fayer

List of Publications by Year in descending order

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141
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6,897
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50170

46
h-index

69108

77
g-index

144
all docs

144
docs citations

144
times ranked

4905
citing authors

#	ARTICLE	IF	CITATIONS
1	Water Dynamics and Structure of Highly Concentrated LiCl Solutions Investigated Using Ultrafast Infrared Spectroscopy. <i>Journal of the American Chemical Society</i> , 2022, 144, 4233-4243.	6.6	23
2	Probing Lattice Dynamics in Two-Dimensional Inorganic Pseudohalide Perovskites with Ultrafast Infrared Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2022, 126, 10145-10158.	1.5	4
3	Free Volume Element Sizes and Dynamics in Polystyrene and Poly(methyl methacrylate) Measured with Ultrafast Infrared Spectroscopy. <i>Journal of the American Chemical Society</i> , 2021, 143, 3583-3594.	6.6	20
4	Orientational Pair Correlations and Local Structure of Benzonitrile from Molecular Dynamics Simulations with Comparisons to Experiments. <i>Journal of Physical Chemistry B</i> , 2021, 125, 3163-3177.	1.2	3
5	Complex Formation and Dissociation Dynamics on Amorphous Silica Surfaces. <i>Journal of Physical Chemistry B</i> , 2021, 125, 4566-4581.	1.2	7
6	Distinguishing steric and electrostatic molecular probe orientational ordering via their effects on reorientation-induced spectral diffusion. <i>Journal of Chemical Physics</i> , 2021, 154, 244104.	1.2	0
7	Long Vibrational Lifetime R-Selenocyanate Probes for Ultrafast Infrared Spectroscopy: Properties and Synthesis. <i>Journal of Physical Chemistry B</i> , 2021, 125, 8907-8918.	1.2	7
8	Ultrafast Dynamics and Liquid Structure in Mesoporous Silica: Propagation of Surface Effects in a Polar Aprotic Solvent. <i>Journal of Physical Chemistry B</i> , 2021, 125, 10018-10034.	1.2	4
9	Identical Water Dynamics in Acrylamide Hydrogels, Polymers, and Monomers in Solution: Ultrafast IR Spectroscopy and Molecular Dynamics Simulations. <i>Journal of the American Chemical Society</i> , 2021, 143, 14855-14868.	6.6	13
10	Proton Transfer from a Photoacid to a Water Wire: First Principles Simulations and Fast Fluorescence Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2021, 125, 12539-12551.	1.2	11
11	Proton Transfer Dynamics in the Aprotic Proton Accepting Solvent 1-Methylimidazole. <i>Journal of Physical Chemistry B</i> , 2020, 124, 7897-7908.	1.2	6
12	Single Ensemble Non-exponential Photoluminescent Population Decays from a Broadband White-Light-Emitting Perovskite. <i>Journal of the American Chemical Society</i> , 2020, 142, 16622-16631.	6.6	44
13	CLS Next Gen: Accurate Frequency-Frequency Correlation Functions from Center Line Slope Analysis of 2D Correlation Spectra Using Artificial Neural Networks. <i>Journal of Physical Chemistry A</i> , 2020, 124, 5979-5992.	1.1	18
14	Amorphous polymer dynamics and free volume element size distributions from ultrafast IR spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 13949-13958.	3.3	18
15	Structural Dynamics in Ionic Liquid Thin Films: The Effect of Cation Chain Length. <i>Journal of Physical Chemistry C</i> , 2020, 124, 4179-4189.	1.5	20
16	Enhanced Menshutkin S_{N2} Reactivity in Mesoporous Silica: The Influence of Surface Catalysis and Confinement. <i>Journal of the American Chemical Society</i> , 2020, 142, 5636-5648.	6.6	25
17	Controlling the Dynamics of Ionic Liquid Thin Films via Multilayer Surface Functionalization. <i>Journal of the American Chemical Society</i> , 2020, 142, 9482-9492.	6.6	25
18	Microstructural and Dynamical Heterogeneities in Ionic Liquids. <i>Chemical Reviews</i> , 2020, 120, 5798-5877.	23.0	277

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19	Effects of pore size on water dynamics in mesoporous silica. <i>Journal of Chemical Physics</i> , 2020, 152, 154704.	1.2	23
20	Pulse-shaped chopping: Eliminating and characterizing heat effects in ultrafast infrared spectroscopy. <i>Journal of Chemical Physics</i> , 2020, 153, 204201.	1.2	6
21	Reorientation-induced Stokes shifts caused by directional interactions in electronic spectroscopy: Fast dynamics of poly(methyl methacrylate). <i>Journal of Chemical Physics</i> , 2019, 150, 194201.	1.2	2
22	Dynamics of Water Molecules and Ions in Concentrated Lithium Chloride Solutions Probed with Ultrafast 2D IR Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2019, 123, 7628-7639.	1.2	33
23	Dynamical properties of a room temperature ionic liquid: Using molecular dynamics simulations to implement a dynamic ion cage model. <i>Journal of Chemical Physics</i> , 2019, 151, 154502.	1.2	16
24	Bulk-like and Interfacial Water Dynamics in Nafion Fuel Cell Membranes Investigated with Ultrafast Nonlinear IR Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2019, 123, 9408-9417.	1.2	22
25	Water Dynamics in Nanoporous Silica: Ultrafast Vibrational Spectroscopy and Molecular Dynamics Simulations. <i>Journal of Physical Chemistry C</i> , 2019, 123, 5790-5803.	1.5	32
26	Tracking Aqueous Proton Transfer by Two-Dimensional Infrared Spectroscopy and ab Initio Molecular Dynamics Simulations. <i>ACS Central Science</i> , 2019, 5, 1269-1277.	5.3	47
27	Fast dynamics of a hydrogen-bonding glass forming liquid: Chemical exchange-induced spectral diffusion in 2D IR spectroscopy. <i>Journal of Chemical Physics</i> , 2019, 150, 124507.	1.2	6
28	Imidazole and 1-Methylimidazole Hydrogen Bonding and Nonhydrogen Bonding Liquid Dynamics: Ultrafast IR Experiments. <i>Journal of Physical Chemistry B</i> , 2019, 123, 2094-2105.	1.2	13
29	Influence of Water on Carbon Dioxide and Room Temperature Ionic Liquid Dynamics: Supported Ionic Liquid Membrane vs the Bulk Liquid. <i>Journal of Physical Chemistry B</i> , 2018, 122, 2389-2395.	1.2	8
30	Orientalional Pair Correlations in a Dipolar Molecular Liquid: Time-Resolved Resonant and Nonresonant Pump-Probe Spectroscopies. <i>Journal of Physical Chemistry B</i> , 2018, 122, 12147-12153.	1.2	6
31	Ion-Molecule Complex Dissociation and Formation Dynamics in LiCl Aqueous Solutions from 2D IR Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2018, 122, 10582-10592.	1.2	35
32	Dynamics and Microstructures of Nicotine/Water Binary Mixtures near the Lower Critical Solution Temperature. <i>Journal of Physical Chemistry B</i> , 2018, 122, 9538-9548.	1.2	7
33	Preface: Special Topic on Chemical Physics of Ionic Liquids. <i>Journal of Chemical Physics</i> , 2018, 148, 193501.	1.2	12
34	Extraordinary Slowing of Structural Dynamics in Thin Films of a Room Temperature Ionic Liquid. <i>ACS Central Science</i> , 2018, 4, 1065-1073.	5.3	33
35	The influence of hydrophilicity on the orientational dynamics and structures of imidazolium-based ionic liquid/water binary mixtures. <i>Journal of Chemical Physics</i> , 2018, 149, 044501.	1.2	21
36	Water Dynamics in Polyacrylamide Hydrogels. <i>Journal of the American Chemical Society</i> , 2018, 140, 9466-9477.	6.6	53

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37	Dynamically Disordered Lattice in a Layered Pb-I-SCN Perovskite Thin Film Probed by Two-Dimensional Infrared Spectroscopy. <i>Journal of the American Chemical Society</i> , 2018, 140, 9882-9890.	6.6	49
38	Two dimensional infrared spectroscopic investigations of protein and enzyme dynamics. <i>FASEB Journal</i> , 2018, 32, .	0.2	0
39	Dynamics in a Room-Temperature Ionic Liquid from the Cation Perspective: 2D IR Vibrational Echo Spectroscopy. <i>Journal of the American Chemical Society</i> , 2017, 139, 2408-2420.	6.6	41
40	Direct observation of dynamic crossover in fragile molecular glass formers with 2D IR vibrational echo spectroscopy. <i>Journal of Chemical Physics</i> , 2017, 146, 124505.	1.2	15
41	Proton Transfer in Perfluorosulfonic Acid Fuel Cell Membranes with Differing Pendant Chains and Equivalent Weights. <i>Journal of Physical Chemistry B</i> , 2017, 121, 4544-4553.	1.2	9
42	Dynamics in a Water Interfacial Boundary Layer Investigated with IR Polarization-Selective Pump-Probe Experiments. <i>Journal of Physical Chemistry B</i> , 2017, 121, 4530-4537.	1.2	24
43	Guest Hydrogen Bond Dynamics and Interactions in the Metal-Organic Framework MIL-53(Al) Measured with Ultrafast Infrared Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2017, 121, 11880-11890.	1.5	21
44	Enhanced nonlinear spectroscopy for monolayers and thin films in near-Brewster's angle reflection pump-probe geometry. <i>Journal of Chemical Physics</i> , 2017, 146, .	1.2	28
45	Dynamics of a Room Temperature Ionic Liquid in Supported Ionic Liquid Membranes vs the Bulk Liquid: 2D IR and Polarized IR Pump-Probe Experiments. <i>Journal of the American Chemical Society</i> , 2017, 139, 311-323.	6.6	55
46	Ultrafast to Ultraslow Dynamics of a Langmuir Monolayer at the Air/Water Interface Observed with Reflection Enhanced 2D IR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2017, 139, 16518-16527.	6.6	40
47	Discontinuity in Fast Dynamics at the Glass Transition of <i>ortho</i> -Terphenyl. <i>Journal of Physical Chemistry B</i> , 2017, 121, 10417-10428.	1.2	3
48	Carbon Dioxide in a Supported Ionic Liquid Membrane: Structural and Rotational Dynamics Measured with 2D IR and Pump-Probe Experiments. <i>Journal of the American Chemical Society</i> , 2017, 139, 11222-11232.	6.6	48
49	Hydrogen Bonding versus π - π Stacking Interactions in Imidazolium-Oxalato borate Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2017, 121, 7173-7179.	1.2	47
50	Impact of Hydrogen Bonding on the Dynamics and Structure of Protic Ionic Liquid/Water Binary Mixtures. <i>Journal of Physical Chemistry B</i> , 2017, 121, 8564-8576.	1.2	46
51	The influence of mesoscopic confinement on the dynamics of imidazolium-based room temperature ionic liquids in polyether sulfone membranes. <i>Journal of Chemical Physics</i> , 2017, 147, 194502.	1.2	15
52	Water-anion hydrogen bonding dynamics: Ultrafast IR experiments and simulations. <i>Journal of Chemical Physics</i> , 2017, 146, 234501.	1.2	38
53	Alkyl Chain Length Dependence of the Dynamics and Structure in the Ionic Regions of Room-Temperature Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2016, 120, 7488-7501.	1.2	36
54	Carbon dioxide in an ionic liquid: Structural and rotational dynamics. <i>Journal of Chemical Physics</i> , 2016, 144, 104506.	1.2	49

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55	Quasi-rotating frame: accurate line shape determination with increased efficiency in noncollinear 2D optical spectroscopy. <i>Journal of the Optical Society of America B: Optical Physics</i> , 2016, 33, 1143.	0.9	11
56	Unraveling the dynamics and structure of functionalized self-assembled monolayers on gold using 2D IR spectroscopy and MD simulations. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, 4929-4934.	3.3	48
57	Ionic Liquid Dynamics Measured with 2D IR and IR Pump-Probe Experiments on a Linear Anion and the Influence of Potassium Cations. <i>Journal of Physical Chemistry B</i> , 2016, 120, 5842-5854.	1.2	39
58	The Influence of Water on the Alkyl Region Structure in Variable Chain Length Imidazolium-Based Ionic Liquid/Water Mixtures. <i>Journal of Physical Chemistry B</i> , 2016, 120, 10350-10357.	1.2	13
59	Orientalional Dynamics of a Functionalized Alkyl Planar Monolayer Probed by Polarization-Selective Angle-Resolved Infrared Pump-Probe Spectroscopy. <i>Journal of the American Chemical Society</i> , 2016, 138, 14057-14065.	6.6	5
60	Critical Slowing of Density Fluctuations Approaching the Isotropic-Nematic Transition in Liquid Crystals: 2D IR Measurements and Mode Coupling Theory. <i>Journal of Physical Chemistry B</i> , 2016, 120, 7003-7015.	1.2	10
61	Ionic Liquid versus Li ⁺ Aqueous Solutions: Water Dynamics near Bistriflimide Anions. <i>Journal of Physical Chemistry B</i> , 2016, 120, 9997-10009.	1.2	24
62	Water of Hydration Dynamics in Minerals Gypsum and Bassanite: Ultrafast 2D IR Spectroscopy of Rocks. <i>Journal of the American Chemical Society</i> , 2016, 138, 9694-9703.	6.6	39
63	Water Dynamics in 1-Alkyl-3-methylimidazolium Tetrafluoroborate Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2016, 120, 11523-11538.	1.2	27
64	Coupling of Carbon Dioxide Stretch and Bend Vibrations Reveals Thermal Population Dynamics in an Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2016, 120, 549-556.	1.2	29
65	Structural and Rotational Dynamics of Carbon Dioxide in 1-Alkyl-3-methylimidazolium Bis(trifluoromethylsulfonyl)imide Ionic Liquids: The Effect of Chain Length. <i>Journal of Physical Chemistry B</i> , 2016, 120, 6698-6711.	1.2	27
66	Separation of experimental 2D IR frequency-frequency correlation functions into structural and reorientation-induced contributions. <i>Journal of Chemical Physics</i> , 2015, 143, 124505.	1.2	42
67	Dynamics of water, methanol, and ethanol in a room temperature ionic liquid. <i>Journal of Chemical Physics</i> , 2015, 142, 212408.	1.2	75
68	Dynamics of Dihydrogen Bonding in Aqueous Solutions of Sodium Borohydride. <i>Journal of Physical Chemistry B</i> , 2015, 119, 3546-3559.	1.2	16
69	The Influence of Cholesterol on Fast Dynamics Inside of Vesicle and Planar Phospholipid Bilayers Measured with 2D IR Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2015, 119, 8852-8862.	1.2	15
70	Structural Influences on the Fast Dynamics of Alkylsiloxane Monolayers on SiO ₂ Surfaces Measured with 2D IR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2015, 119, 16811-16823.	1.5	19
71	Proton Transfer in Ionic and Neutral Reverse Micelles. <i>Journal of Physical Chemistry B</i> , 2015, 119, 6024-6034.	1.2	52
72	Observation and theory of reorientation-induced spectral diffusion in polarization-selective 2D IR spectroscopy. <i>Journal of Chemical Physics</i> , 2015, 142, 184505.	1.2	46

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73	Molecular Anion Hydrogen Bonding Dynamics in Aqueous Solution. <i>Journal of Physical Chemistry B</i> , 2015, 119, 13407-13415.	1.2	36
74	New divergent dynamics in the isotropic to nematic phase transition of liquid crystals measured with 2D IR vibrational echo spectroscopy. <i>Journal of Chemical Physics</i> , 2014, 141, 194502.	1.2	19
75	Structural dynamics inside a functionalized metal-organic framework probed by ultrafast 2D IR spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 18442-18447.	3.3	76
76	Theory of third-order spectroscopic methods to extract detailed molecular orientational dynamics for planar surfaces and other uniaxial systems. <i>Journal of Chemical Physics</i> , 2014, 140, 144702.	1.2	13
77	Size-dependent ultrafast structural dynamics inside phospholipid vesicle bilayers measured with 2D IR vibrational echoes. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 918-923.	3.3	28
78	Length Scales and Structural Dynamics in Nematogen Pseudonematic Domains Measured with 2D IR Vibrational Echoes and Optical Kerr Effect Experiments. <i>Journal of Physical Chemistry B</i> , 2014, 118, 7856-7868.	1.2	12
79	Dynamics of Molecular Monolayers with Different Chain Lengths in Air and Solvents Probed by Ultrafast 2D IR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2014, 118, 523-532.	1.5	18
80	Dynamics and structure of room temperature ionic liquids. <i>Chemical Physics Letters</i> , 2014, 616-617, 259-274.	1.2	70
81	Ultrafast Structural Dynamics Inside Planar Phospholipid Multibilayer Model Cell Membranes Measured with 2D IR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2013, 135, 11063-11074.	6.6	41
82	Dynamics in the Isotropic Phase of Nematogens Using 2D IR Vibrational Echo Measurements on Natural-Abundance ^{13}C and Extended Lifetime Probes. <i>Journal of Physical Chemistry B</i> , 2013, 117, 15060-15071.	1.2	35
83	Orientational Dynamics in a Lyotropic Room Temperature Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2013, 117, 14775-14784.	1.2	34
84	Structural Dynamics at Monolayer-Liquid Interfaces Probed by 2D IR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2013, 117, 1409-1420.	1.5	23
85	Theoretical examination of picosecond phenol migration dynamics in phenylacetylene solution. <i>Chemical Physics</i> , 2013, 422, 175-183.	0.9	2
86	Dynamics of Isolated Water Molecules in a Sea of Ions in a Room Temperature Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2013, 117, 623-635.	1.2	58
87	Dynamics in the Interior of AOT Lamellae Investigated with Two-Dimensional Infrared Spectroscopy. <i>Journal of the American Chemical Society</i> , 2013, 135, 5118-5126.	6.6	26
88	The Influence of Lithium Cations on Dynamics and Structure of Room Temperature Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2013, 117, 9768-9774.	1.2	44
89	Excitation transfer induced spectral diffusion and the influence of structural spectral diffusion. <i>Journal of Chemical Physics</i> , 2012, 137, 064109.	1.2	12
90	Water Dynamics in Water/DMSO Binary Mixtures. <i>Journal of Physical Chemistry B</i> , 2012, 116, 5479-5490.	1.2	155

#	ARTICLE	IF	CITATIONS
91	Water Dynamics in Divalent and Monovalent Concentrated Salt Solutions. <i>Journal of Physical Chemistry B</i> , 2012, 116, 13781-13792.	1.2	65
92	Fast Dynamics of HP35 for Folded and Urea-Unfolded Conditions. <i>Journal of Physical Chemistry B</i> , 2012, 116, 11024-11031.	1.2	22
93	Orientalional Dynamics of Room Temperature Ionic Liquid/Water Mixtures: Water-Induced Structure. <i>Journal of Physical Chemistry B</i> , 2012, 116, 1777-1787.	1.2	97
94	Dynamics of Functionalized Surface Molecular Monolayers Studied with Ultrafast Infrared Vibrational Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2012, 116, 23428-23440.	1.5	30
95	Ribonuclease S Dynamics Measured Using a Nitrile Label with 2D IR Vibrational Echo Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2012, 116, 4034-4042.	1.2	100
96	Conformational Dynamics and Stability of HP35 Studied with 2D IR Vibrational Echoes. <i>Journal of the American Chemical Society</i> , 2012, 134, 12118-12124.	6.6	45
97	Dynamics of Water Interacting with Interfaces, Molecules, and Ions.. <i>Accounts of Chemical Research</i> , 2012, 45, 3-14.	7.6	242
98	Protein Dynamics Studied with Ultrafast Two-Dimensional Infrared Vibrational Echo Spectroscopy. <i>Accounts of Chemical Research</i> , 2012, 45, 1866-1874.	7.6	66
99	Protein Dynamics in Cytochrome P450 Molecular Recognition and Substrate Specificity Using 2D IR Vibrational Echo Spectroscopy. <i>Journal of the American Chemical Society</i> , 2011, 133, 3995-4004.	6.6	60
100	Time-Dependent Fifth-Order Bands in Nominally Third-Order 2D IR Vibrational Echo Spectra. <i>Journal of Physical Chemistry A</i> , 2011, 115, 9714-9723.	1.1	9
101	Two-Dimensional IR Spectroscopy of Protein Dynamics Using Two Vibrational Labels: A Site-Specific Genetically Encoded Unnatural Amino Acid and an Active Site Ligand. <i>Journal of Physical Chemistry B</i> , 2011, 115, 11294-11304.	1.2	92
102	Water in a Crowd. <i>Physiology</i> , 2011, 26, 381-392.	1.6	27
103	Structural Dynamics of a Catalytic Monolayer Probed by Ultrafast 2D IR Vibrational Echoes. <i>Science</i> , 2011, 334, 634-639.	6.0	117
104	Dynamics of the folded and unfolded villin headpiece (HP35) measured with ultrafast 2D IR vibrational echo spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011, 108, 3578-3583.	3.3	105
105	Analysis of Water in Confined Geometries and at Interfaces. <i>Annual Review of Analytical Chemistry</i> , 2010, 3, 89-107.	2.8	240
106	Orientalional and Translational Dynamics of Polyether/Water Solutions. <i>Journal of Physical Chemistry B</i> , 2010, 114, 5350-5358.	1.2	37
107	Water dynamics in large and small reverse micelles: From two ensembles to collective behavior. <i>Journal of Chemical Physics</i> , 2009, 131, 014704.	1.2	170
108	Water Dynamics and Interactions in Water~Polyether Binary Mixtures. <i>Journal of the American Chemical Society</i> , 2009, 131, 5530-5539.	6.6	68

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109	Water Dynamics in Salt Solutions Studied with Ultrafast Two-Dimensional Infrared (2D IR) Vibrational Echo Spectroscopy. <i>Accounts of Chemical Research</i> , 2009, 42, 1210-1219.	7.6	123
110	Dynamics of Liquids, Molecules, and Proteins Measured with Ultrafast 2D IR Vibrational Echo Chemical Exchange Spectroscopy. <i>Annual Review of Physical Chemistry</i> , 2009, 60, 21-38.	4.8	188
111	Water inertial reorientation: Hydrogen bond strength and the angular potential. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 5295-5300.	3.3	181
112	Water at the Surfaces of Aligned Phospholipid Multibilayer Model Membranes Probed with Ultrafast Vibrational Spectroscopy. <i>Journal of the American Chemical Society</i> , 2008, 130, 13927-13937.	6.6	135
113	Solute-Solvent Complex Switching Dynamics of Chloroform between Acetone and Dimethylsulfoxide Two-Dimensional IR Chemical Exchange Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2008, 112, 13906-13915.	1.2	44
114	Direct observation of fast protein conformational switching. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 8619-8624.	3.3	125
115	1P-102 Fast protein conformational switching observed with 2D-IR vibrational echo (Invited Talk for) TJ ETQq1 1 0.784314 rgBT /Overlap	0.0	0
116	Disulfide bond influence on protein structural dynamics probed with 2D-IR vibrational echo spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 19309-19314.	3.3	61
117	Neuroglobin dynamics observed with ultrafast 2D-IR vibrational echo spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 16116-16121.	3.3	71
118	Water Dynamics in Nafion Fuel Cell Membranes: The Effects of Confinement and Structural Changes on the Hydrogen Bond Network. <i>Journal of Physical Chemistry C</i> , 2007, 111, 8884-8891.	1.5	121
119	Probing dynamics of complex molecular systems with ultrafast 2D IR vibrational echo spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 1533.	1.3	93
120	Hydrogen Bond Lifetimes and Energetics for Solute/Solvent Complexes Studied with 2D-IR Vibrational Echo Spectroscopy. <i>Journal of the American Chemical Society</i> , 2007, 129, 4328-4335.	6.6	82
121	Dynamics in Supercooled Ionic Organic Liquids and Mode Coupling Theory Analysis. <i>Journal of Physical Chemistry A</i> , 2006, 110, 10384-10391.	1.1	83
122	Cytochrome c552 Mutants: Structure and Dynamics at the Active Site Probed by Multidimensional NMR and Vibration Echo Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2006, 110, 18803-18810.	1.2	18
123	Dynamics of Proteins Encapsulated in Silica Sol-Gel Glasses Studied with IR Vibrational Echo Spectroscopy. <i>Journal of the American Chemical Society</i> , 2006, 128, 3990-3997.	6.6	65
124	Vibrational Echo and Pump-Probe Spectroscopic Studies of the Dynamics of Water Molecules Confined to Nanoscopic Dimensions. , 2006, , 195-203.		5
125	The Influence of Aqueous versus Glassy Solvents on Protein Dynamics: Vibrational Echo Experiments and Molecular Dynamics Simulations. <i>Journal of the American Chemical Society</i> , 2005, 127, 14279-14289.	6.6	96
126	Orientational dynamics of water confined on a nanometer length scale in reverse micelles. <i>Journal of Chemical Physics</i> , 2005, 122, 174501.	1.2	219

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127	Water dynamics: dependence on local structure probed with vibrational echo correlation spectroscopy. <i>Chemical Physics Letters</i> , 2004, 386, 295-300.	1.2	131
128	Vibrational echo experiments on red blood cells: Comparison of the dynamics of cytoplasmic and aqueous hemoglobin. <i>Chemical Physics Letters</i> , 2004, 392, 324-329.	1.2	21
129	Using ultrafast infrared multidimensional correlation spectroscopy to aid in vibrational spectral peak assignments. <i>Chemical Physics Letters</i> , 2003, 381, 139-146.	1.2	36
130	Ultrafast heterodyne detected infrared multidimensional vibrational stimulated echo studies of hydrogen bond dynamics. <i>Chemical Physics Letters</i> , 2003, 374, 362-371.	1.2	75
131	Myoglobin-CO Conformational Substate Dynamics: 2D Vibrational Echoes and MD Simulations. <i>Biophysical Journal</i> , 2002, 82, 3277-3288.	0.2	52
132	Short time dynamics in the isotropic phase of liquid crystals: the aspect ratio and the power law decay. <i>Chemical Physics Letters</i> , 2002, 366, 82-87.	1.2	50
133	FASTPROTEINDYNAMICSPROBED WITHINFRAREDVIBRATIONALECHOEXPERIMENTS. <i>Annual Review of Physical Chemistry</i> , 2001, 52, 315-356.	4.8	139
134	Effect of chromophore diffusion on electronic excitation transfer in micellar systems. <i>Chemical Physics Letters</i> , 1997, 276, 274-281.	1.2	12
135	Theory of Electronic Excitation Transfer in Polymer Micelles and Lamellae. <i>Molecular Crystals and Liquid Crystals</i> , 1996, 283, 173-177.	0.3	2
136	Quantitative measurement of chain contraction in a solid blend of two incompatible polymers: poly(methyl methacrylate)/poly(vinyl acetate). <i>Macromolecules</i> , 1990, 23, 111-120.	2.2	50
137	Hole burning line shapes in a two-dimensional glass: A model for hole burning line shapes of molecules on surfaces. <i>Chemical Physics Letters</i> , 1990, 168, 371-378.	1.2	9
138	Shocked molecular solids: Vibrational up pumping, defect hot spot formation, and the onset of chemistry. <i>Journal of Chemical Physics</i> , 1990, 92, 3798-3812.	1.2	269
139	Fluorescence depolarization of chromophores in polymeric solids. <i>Macromolecules</i> , 1989, 22, 874-879.	2.2	11
140	Short polymer chain statistics and the relationship to end to end electronic excitation transport: random walks with variable-step lengths. <i>Macromolecules</i> , 1988, 21, 1145-1154.	2.2	9
141	Probing intermolecular interactions with picosecond photon echo experiments. <i>Accounts of Chemical Research</i> , 1987, 20, 120-126.	7.6	26