

Ilan Benjamin

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/7207933/publications.pdf>

Version: 2024-02-01

48
papers

2,235
citations

257450

24
h-index

214800

47
g-index

49
all docs

49
docs citations

49
times ranked

1377
citing authors

#	ARTICLE	IF	CITATIONS
1	Deconstructing the Local Intermolecular Ordering and Dynamics of Liquid Chloroform and Bromoform. <i>Journal of Physical Chemistry B</i> , 2021, 125, 3629-3637.	2.6	1
2	Antagonistic Role of Aqueous Complexation in the Solvent Extraction and Separation of Rare Earth Ions. <i>ACS Central Science</i> , 2021, 7, 1908-1918.	11.3	18
3	Molecular Dynamics Studies on the Effect of Surface Roughness and Surface Tension on the Thermodynamics and Dynamics of Hydronium Ion Transfer Across the Liquid/Liquid Interface. <i>Journal of Physical Chemistry B</i> , 2020, 124, 8711-8718.	2.6	4
4	Transfer of an erbium ion across the water/dodecane interface: Structure and thermodynamics via molecular dynamics simulations. <i>Chemical Physics Letters</i> , 2019, 737, 136825.	2.6	11
5	Hydronium ion at the water/1,2-dichloroethane interface: Structure, thermodynamics, and dynamics of ion transfer. <i>Journal of Chemical Physics</i> , 2019, 151, 094701.	3.0	9
6	Nanoscale view of assisted ion transport across the liquid-liquid interface. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 18227-18232.	7.1	68
7	Miscibility at the immiscible liquid/liquid interface: A molecular dynamics study of thermodynamics and mechanism. <i>Journal of Chemical Physics</i> , 2018, 148, 034707.	3.0	8
8	Structure and Dynamics of Host/Guest Complexation at the Liquid/Liquid Interface: Implications for Inverse Phase Transfer Catalysis. <i>Journal of Physical Chemistry C</i> , 2017, 121, 4999-5011.	3.1	4
9	On the local intermolecular ordering and dynamics of liquid chloroform. <i>Journal of Molecular Liquids</i> , 2017, 248, 121-126.	4.9	7
10	SN2 Reaction Rate Enhancement by β -Cyclodextrin at the Liquid/Liquid Interface. <i>Journal of Physical Chemistry C</i> , 2017, 121, 19209-19217.	3.1	6
11	Geometric and energetic considerations of surface fluctuations during ion transfer across the water-immiscible organic liquid interface. <i>Journal of Chemical Physics</i> , 2016, 145, 014701.	3.0	28
12	Unusual Structure and Dynamics at Silica/Methanol and Silica/Ethanol Interfaces—A Molecular Dynamics and Nonlinear Optical Study. <i>Journal of Physical Chemistry B</i> , 2016, 120, 1569-1578.	2.6	36
13	Reaction Dynamics at Liquid Interfaces. <i>Annual Review of Physical Chemistry</i> , 2015, 66, 165-188.	10.8	53
14	Mechanism and Dynamics of Molecular Exchange at the Silica/Binary Solvent Mixtures Interface. <i>Journal of Physical Chemistry A</i> , 2015, 119, 12073-12081.	2.5	5
15	β -Cyclodextrin at the Water/1-Bromobutane Interface: Molecular Insight into Reverse Phase Transfer Catalysis. <i>Langmuir</i> , 2015, 31, 5086-5092.	3.5	13
16	Photoinduced Excited State Electron Transfer at Liquid/Liquid Interfaces. <i>Journal of Physical Chemistry B</i> , 2014, 118, 7703-7714.	2.6	9
17	Recombination, Dissociation, and Transport of Ion Pairs across the Liquid/Liquid Interface. Implications for Phase Transfer Catalysis. <i>Journal of Physical Chemistry B</i> , 2013, 117, 4325-4331.	2.6	36
18	Effect of a Phase Transfer Catalyst on the Dynamics of an S_N2 Reaction. A Molecular Dynamics Study. <i>Journal of Physical Chemistry C</i> , 2011, 115, 2290-2296.	3.1	20

#	ARTICLE	IF	CITATIONS
19	A model SN2 reaction "on water"™ does not show rate enhancement. <i>Chemical Physics Letters</i> , 2011, 508, 59-62.	2.6	8
20	A molecular dynamics/EVB study of an SN2 reaction in water clusters. <i>Chemical Physics Letters</i> , 2010, 492, 220-225.	2.6	9
21	Molecular dynamics study of hydrated alkali and halide ions in liquid nitrobenzene. <i>Journal of Electroanalytical Chemistry</i> , 2010, 650, 41-46.	3.8	14
22	A Molecular Dynamics~Empirical Valence Bond Study of an S_N2 Reaction at the Water/Chloroform Interface. <i>Journal of Physical Chemistry C</i> , 2010, 114, 1154-1163.	3.1	19
23	Microhydration effects on a model SN2 reaction in a nonpolar solvent. <i>Journal of Chemical Physics</i> , 2009, 130, 194502.	3.0	11
24	Solute dynamics at aqueous interfaces. <i>Chemical Physics Letters</i> , 2009, 469, 229-241.	2.6	22
25	Free Energy of Transfer of Hydrated Ion Clusters from Water to an Immiscible Organic Solvent. <i>Journal of Physical Chemistry B</i> , 2009, 113, 9296-9303.	2.6	44
26	Structure and Dynamics of Hydrated Ions in a Water-Immiscible Organic Solvent. <i>Journal of Physical Chemistry B</i> , 2008, 112, 15801-15806.	2.6	32
27	Empirical valence bond model of an SN2 reaction in polar and nonpolar solvents. <i>Journal of Chemical Physics</i> , 2008, 129, 074508.	3.0	19
28	Static and Dynamic Electronic Spectroscopy at Liquid Interfaces. <i>Chemical Reviews</i> , 2006, 106, 1212-1233.	47.7	87
29	Ion Distributions near a Liquid-Liquid Interface. <i>Science</i> , 2006, 311, 216-218.	12.6	229
30	Ion distributions at the nitrobenzene~water interface electrified by a common ion. <i>Journal of Electroanalytical Chemistry</i> , 2006, 593, 142-158.	3.8	42
31	Hydration Shell Exchange Dynamics during Ion Transfer Across the Liquid/Liquid Interface. <i>Journal of Physical Chemistry B</i> , 2005, 109, 16455-16462.	2.6	36
32	Molecular dynamics study of the vibrational relaxation of OCl and OCl~ in the bulk and the surface of water and acetonitrile. <i>Journal of Molecular Liquids</i> , 2004, 110, 133-139.	4.9	15
33	Electron transfer at the interface between water and self-assembled monolayers. <i>Chemical Physics Letters</i> , 2004, 385, 79-84.	2.6	6
34	Influence of Surface Tension on Adsorbate Molecular Rotation at Liquid/Liquid Interfaces. <i>Journal of Physical Chemistry B</i> , 2004, 108, 15443-15445.	2.6	12
35	Vibrational relaxation at water surfaces. <i>Journal of Chemical Physics</i> , 2002, 117, 4532-4541.	3.0	26
36	Transfer of a Tetramethylammonium Ion across the Water~Nitrobenzene Interface:~Potential of Mean Force and Nonequilibrium Dynamics. <i>Journal of Physical Chemistry A</i> , 1999, 103, 10274-10279.	2.5	69

#	ARTICLE	IF	CITATIONS
37	Reorganization free energy for electron transfer reactions at liquid/liquid interfaces. <i>Electrochimica Acta</i> , 1998, 44, 133-138.	5.2	27
38	Molecular dynamics simulation of the water nitrobenzene interface. <i>Journal of Electroanalytical Chemistry</i> , 1998, 450, 335-345.	3.8	70
39	Solvent Effects on Electronic Spectra at Liquid Interfaces. A Continuum Electrostatic Model. <i>Journal of Physical Chemistry A</i> , 1998, 102, 9500-9506.	2.5	56
40	MOLECULAR STRUCTURE AND DYNAMICS AT LIQUID-LIQUID INTERFACES. <i>Annual Review of Physical Chemistry</i> , 1997, 48, 407-451.	10.8	271
41	Electric field effects on the structure and dynamics at a liquid liquid interface. <i>Journal of Electroanalytical Chemistry</i> , 1995, 391, 1-10.	3.8	57
42	Transfer of Small Ions across the Water/1,2-Dichloroethane Interface. <i>The Journal of Physical Chemistry</i> , 1995, 99, 9974-9985.	2.9	90
43	Theory and Computer Simulations of Solvation and Chemical Reactions at Liquid Interfaces. <i>Accounts of Chemical Research</i> , 1995, 28, 233-239.	15.6	80
44	Dynamics of ion desorption from the liquidâ€”vapor interface of water. <i>Chemical Physics Letters</i> , 1993, 202, 379-383.	2.6	18
45	Dynamics of ion transfer across a liquidâ€”liquid interface: A comparison between molecular dynamics and a diffusion model. <i>Journal of Chemical Physics</i> , 1992, 96, 577-585.	3.0	64
46	Theoretical study of the water/1,2-dichloroethane interface: Structure, dynamics, and conformational equilibria at the liquidâ€”liquid interface. <i>Journal of Chemical Physics</i> , 1992, 97, 1432-1445.	3.0	281
47	Theoretical study of ion solvation at the water liquidâ€”vapor interface. <i>Journal of Chemical Physics</i> , 1991, 95, 3698-3709.	3.0	134
48	Molecular dynamics study of the free energy functions for electron-transfer reactions at the liquid-liquid interface. <i>The Journal of Physical Chemistry</i> , 1991, 95, 6675-6683.	2.9	47