

Vance W Jaeger

List of Publications by Year in descending order

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13
papers

671
citations

933447

10
h-index

1125743

13
g-index

13
all docs

13
docs citations

13
times ranked

890
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | The General AMBER Force Field (GAFF) Can Accurately Predict Thermodynamic and Transport Properties of Many Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2015, 119, 5882-5895. | 2.6 | 319 |
| 2 | Structure, Dynamics, and Activity of Xylanase Solvated in Binary Mixtures of Ionic Liquid and Water. <i>ACS Chemical Biology</i> , 2013, 8, 1179-1186. | 3.4 | 84 |
| 3 | Comparison of Three Ionic Liquid-Tolerant Cellulases by Molecular Dynamics. <i>Biophysical Journal</i> , 2015, 108, 880-892. | 0.5 | 62 |
| 4 | Diatom Mimics: Directing the Formation of Biosilica Nanoparticles by Controlled Folding of Lysine-Leucine Peptides. <i>Journal of the American Chemical Society</i> , 2014, 136, 15134-15137. | 13.7 | 54 |
| 5 | The Structure of the Diatom Silaffin Peptide R5 within Freestanding Two-Dimensional Biosilica Sheets. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 8277-8280. | 13.8 | 34 |
| 6 | Biomimetic Growth of Ultrathin Silica Sheets Using Artificial Amphiphilic Peptides. <i>Advanced Materials Interfaces</i> , 2015, 2, 1500282. | 3.7 | 33 |
| 7 | Photodegradation of methyl orange and 2,3-butanedione on titanium-dioxide nanotube arrays efficiently synthesized on titanium coils. <i>Applied Catalysis B: Environmental</i> , 2011, 110, 6-13. | 20.2 | 31 |
| 8 | Destabilization of Human Serum Albumin by Ionic Liquids Studied Using Enhanced Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2016, 120, 12079-12087. | 2.6 | 26 |
| 9 | Acetylation dictates the morphology of nanophase biosilica precipitated by a 14-amino acid leucine-lysine peptide. <i>Journal of Peptide Science</i> , 2017, 23, 141-147. | 1.4 | 11 |
| 10 | Structure Characterization and Properties of Metal-Surfactant Complexes Dispersed in Organic Solvents. <i>Langmuir</i> , 2015, 31, 9006-9016. | 3.5 | 10 |
| 11 | Interpretation of Interfacial Protein Spectra with Enhanced Molecular Simulation Ensembles. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 698-707. | 5.3 | 4 |
| 12 | Die Struktur des Silaffin-Peptides R5 aus Diatomeen in freistehenden zweidimensionalen Biosilikatwänden. <i>Angewandte Chemie</i> , 2017, 129, 8390-8394. | 2.0 | 2 |
| 13 | Structural Insights into Self-Assembled Aerosol-OT Aggregates in Aqueous Media Using Atomistic Molecular Dynamics. <i>Journal of Physical Chemistry B</i> , 2021, 125, 13789-13803. | 2.6 | 1 |