Journal of Computer-Aided Molecular Design
Incorporating Perspectives in Drug Discovery and Design
Editors-in-Chief: F. Gago; T.R. Stouch; R. Lewis

► 100% of authors who answered a survey reported that they would definitely publish or probably publish in the journal again

The *Journal of Computer-Aided Molecular Design* provides a form for disseminating information on both the theory and the application of computer-based methods in the analysis and design of molecules. The scope of the journal encompasses papers which report new and original research and applications in the following areas: theoretical chemistry;

computational chemistry;

computer and molecular graphics;

molecular modeling;

protein engineering;

drug design;

expert systems;

general structure-property relationships;

molecular dynamics;

chemical database development and usage.

Contributions on computer-aided molecular modeling studies in pharmaceutical, polymer, materials and surface sciences, as well as other molecular-based disciplines, are particularly welcome.

As of Volume 15, a number of issues per volume will be dedicated to *Perspectives in Drug Discovery and Design* and from time to time overviews will be included in `regular’ JCAMD issues.

Impact Factor: 2.356 (2017), Journal Citation Reports®

On the homepage of *Journal of Computer-Aided Molecular Design* at [springer.com](http://springer.com) you can

► Sign up for our Table of Contents Alerts

► Get to know the complete Editorial Board

► Find submission information