

ISSN: 2332-5968 Volume 14, Number 1, January 2026



Computational Chemistry



ISSN: 2332-5968



<https://www.scirp.org/journal/cc>

Journal Editorial Board

ISSN 2332-5968 (Print) ISSN 2332-5984 (Online)

<https://www.scirp.org/journal/cc>

Editor-in-Chief

Prof. Anwar El-Shahawy

Assiut University, Egypt

Editorial Board

Dr. Ali Khalaf Al-Matar

University of Jordan, Jordan

Dr. Enrique Arribas

University of Castilla-La Mancha, Spain

Prof. Nagarajan Arumugam

PSG College of Pharmacy, India

Dr. Tamafo Fouegue Aymard Didier

The University of Bertoua, Cameroon

Prof. Todor Dudev

Sofia University, Bulgaria

Dr. Natalja Fjodorova

National Institute of Chemistry, Slovenia

Dr. Soumadwip Ghosh

La Jolla, USA

Dr. Selcuk Gümüs

Yuzuncu Yil University, Türkiye

Dr. Weiwei Han

Jilin University, China

Prof. Alireza Heidari

California South University, USA

Prof. Parasuraman Jaisankar

Indian Institute of Chemical Biology, India

Dr. C. Gopi Mohan

Amrita Institute of Medical Sciences and Research Center, India

Dr. Floriano Paes Silva Junior

Instituto Oswaldo Cruz, Brazil

Dr. Carlos P. Sosa

Cray Inc and University of Minnesota, USA

Prof. Hilal S. Wahab

Al-Nahrain University, Iraq

Dr. Carline Josette Weinberg

Roumanian Academy, Romania

Dr. Julliane Yoneda

Universidade Federal Fluminense, Brazil

Prof. Abdelkader Zarrouk

Mohammed First University, Morocco

Table of Contents

Volume 14 Number 1

January 2026

**Quantum Dynamical Effects in 2-(2-Furyl)-3-Hydroxychromone Using
Path Integral Molecular Dynamics**

A. B. E. Ali, D. P. Foulla, Y. Kanabet, P. Wang-Yang, D. Lissouck, S.-J. Koyambo-Konzapa 1

Computational Chemistry (CC)

Journal Information

SUBSCRIPTIONS

The *Computational Chemistry* (Online at Scientific Research Publishing, <https://www.scirp.org/>) is published quarterly by Scientific Research Publishing, Inc., USA.

Subscription rates:

Print: \$39 per issue.

To subscribe, please contact Journals Subscriptions Department, E-mail: sub@scirp.org

SERVICES

Advertisements

Advertisement Sales Department, E-mail: service@scirp.org

Reprints (minimum quantity 100 copies)

Reprints Co-ordinator, Scientific Research Publishing, Inc., USA.

E-mail: sub@scirp.org

COPYRIGHT

Copyright and reuse rights for the front matter of the journal:

Copyright © 2026 by Scientific Research Publishing Inc.

This work is licensed under the Creative Commons Attribution International License (CC BY).

<http://creativecommons.org/licenses/by/4.0/>

Copyright for individual papers of the journal:

Copyright © 2026 by author(s) and Scientific Research Publishing Inc.

Reuse rights for individual papers:

Note: At SCIRP authors can choose between CC BY and CC BY-NC. Please consult each paper for its reuse rights.

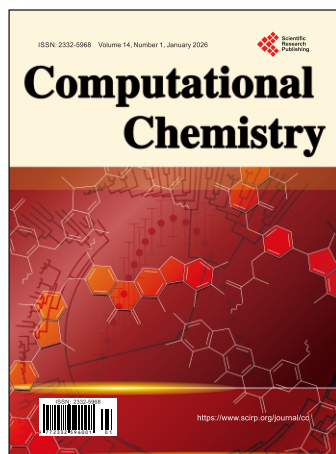
Disclaimer of liability

Statements and opinions expressed in the articles and communications are those of the individual contributors and not the statements and opinion of Scientific Research Publishing, Inc. We assume no responsibility or liability for any damage or injury to persons or property arising out of the use of any materials, instructions, methods or ideas contained herein. We expressly disclaim any implied warranties of merchantability or fitness for a particular purpose. If expert assistance is required, the services of a competent professional person should be sought.

PRODUCTION INFORMATION

For manuscripts that have been accepted for publication, please contact:

E-mail: cc@scirp.org



Computational Chemistry

ISSN: 2332-5968 (Print), 2332-5984 (Online)

<https://www.scirp.org/journal/cc>

Editorial Board

Dr. Ali Khalaf Al-Matar

Dr. Enrique Arribas

Prof. Nagarajan Arumugam

Dr. Tamafo Fouegue Aymard Didier

Prof. Todor Dudev

Dr. Natalja Fjodorova

Dr. Soumadwip Ghosh

Dr. Selcuk Gümüs

Dr. Weiwei Han

Prof. Alireza Heidari

Prof. Parasuraman Jaisankar

Dr. C. Gopi Mohan

Dr. Floriano Paes Silva Junior

Dr. Carlos P. Sosa

Prof. Hilal S. Wahab

Dr. Carline Josette Weinberg

Dr. Julliane Yoneda

Prof. Abdelkader Zarrouk

Subject Coverage

Computational Chemistry (CC) is a unique platform for publishing novel ideas, research outcomes and fundamental advances in all aspects of computational chemistry. All manuscripts must be prepared in English, and are subject to a rigorous and fair peer-review process. The journal publishes original papers including but not limited to the following fields:

- *Ab Initio* Quantum Mechanics
- Chemical Artificial Intelligence
- Chemical Computer-Assisted Instruction
- Cheminformatics
- Chemistry Database
- Code Benchmarking
- Computational Analytical Chemistry
- Computational Biochemistry
- Computational Biophysics
- Computational Colloidal Chemistry
- Computational Electro-Chemistry
- Computational Nanochemistry
- Computational Surface Chemistry
- Computer Molecular Modeling
- Computer-Aided Drug Design
- Data Acquisition, Statistic Analysis and Other Applications
- Density Functional Theory
- Force Field Development and Testing
- Molecular Design
- Molecular Dynamics
- Molecular Mechanics
- Molecular Modeling
- Molecular Simulation
- Molecular Structure Modeling and Image Display
- Quantum Chemistry Calculation
- Semi-Empirical Quantum Mechanics
- Statistical Mechanics
- Theoretical Inorganic Chemistry
- Theoretical Organic Chemistry
- Theoretical Thermo Chemistry

We are also interested in: 1) Short reports—2-5 page papers where an author can present an idea with theoretical background, but has not yet completed the research needed for a complete paper or an author presents preliminary data; 2) Short communications—2-5 page papers; 3) Technical notes—2-5 page papers; 4) Letters to the Editor (the number of pages is not restricted); 5) Reviews or book reviews—comments and critiques (the number of pages is not restricted); 6) Advertisement – 1-2 page papers; 7) News Letters – 1-5 page papers.

Website and E-Mail

<https://www.scirp.org/journal/cc>

E-mail: cc@scirp.org