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Education Information

Yildiz Technical University, Fen Edebiyat Fakültesi, Fizik, Turkey Continues

Research Areas

Physics, General Physics, Statistical physics, thermodynamic and nonlinear dynamic systems, Condensed Matter 1: Structural, Mechanical and Thermal Properties, Thermal Properties of Condensed Matter, Natural Sciences

Academic Titles / Tasks

Research Assistant, Yildiz Technical University, Faculty Of Arts & Science, Department Of Physics, 1998 - Continues

Published journal articles indexed by SCI, SSCI, and AHCI

- I. **Insight into the structural, thermal and ion transport properties of solid and liquid Mg₃N₂: a model potential and NPT molecular dynamics simulation**
Aydin Y., Gunay S. D., AKDERE Ü., TAŞSEVEN Ç.
MOLECULAR SIMULATION, vol.48, no.7, pp.602-609, 2022 (SCI-Expanded)
- II. **Influence of repeating sequence on structural and thermal stability of crystalline domain of bombyx mori silk fibroin**
Aksakal B., Akdere Ü., Günay S. D., Çağın T., Taşseven Ç.
MATERIALS RESEARCH EXPRESS, vol.6, no.12, 2019 (SCI-Expanded)
- III. **Ordering and diffusion in liquid magnesium antimonide (Mg₃Sb₂) from hypernetted-chain theory and molecular dynamics simulation**
AKDERE Ü., GÜNAY S. D., TAŞSEVEN Ç.
Ionics, vol.25, no.6, pp.2711-2717, 2019 (SCI-Expanded)
- IV. **Thermal expansion and heat capacities of AgBr and AgCl at solid and liquid phases from molecular dynamics simulation**
Akdere U.
International Journal of Modern Physics B, vol.29, no.14, 2015 (SCI-Expanded)
- V. **Hypernetted chain calculations of molten uranium dioxide: Comparison of rigid ion potentials**
GÜNAY S. D., AKDERE Ü., TAŞSEVEN Ç.
JOURNAL OF MOLECULAR LIQUIDS, vol.173, pp.124-129, 2012 (SCI-Expanded)
- VI. **THERMOPHYSICAL PROPERTIES OF URANIUM DIOXIDE: A MOLECULAR DYNAMICS STUDY OF SOLID**

AND LIQUID PHASES OF STOICHIOMETRIC UO₂

GÜNAY S. D., KAVANOZ H. B., AKDERE Ü., TAŞSEVEN Ç.

INTERNATIONAL JOURNAL OF MODERN PHYSICS B, vol.25, pp.3211-3223, 2011 (SCI-Expanded)

VII. MOLECULAR DYNAMICS SIMULATION OF UO₂: AN ALTERNATIVE RIGID ION MODEL POTENTIAL

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INTERNATIONAL JOURNAL OF MODERN PHYSICS B, vol.25, no.9, pp.1201-1210, 2011 (SCI-Expanded)

VIII. Freezing in halide salts

AKDERE Ü., YILMAZ M., KAVANOZ H. B., TAŞSEVEN Ç.

ACTA PHYSICA POLONICA A, vol.113, no.6, pp.1659-1670, 2008 (SCI-Expanded)

Articles Published in Other Journals

I. Thermophysical Properties of Anti-Parallel β -Sheets with Bombyx mori Silk Nanostructures [Gly-Ser-Gly-Ala-Gly-Ala]_n and [Gly-Ala]_n

Aksakal B., Günay S. D., Akdere Ü., Çağın T., Taşseven Ç.

materials science forum, vol.856, pp.70-73, 2016 (Scopus)

II. Thermomechanical properties of anti-parallel β -sheets with Bombyx mori silk nano structures [Gly-Ser-Gly-Ala-Gly-Ala]_n and [Gly-Ala]_n

Akdere Ü., Aksakal B., Günay S. D., Taşseven Ç., Çağın T.

Materials Science Forum, vol.851, pp.74-77, 2016 (Scopus)

III. Classical Molecular Dynamics Simulation of Mg₃Bi₂ at Solid and Liquid Phase

GÜNAY S. D., KAYADIBI F., AKDERE Ü., TAŞSEVEN Ç.

Balkan Physics Letters, vol.15, pp.151013, 2009 (Peer-Reviewed Journal)

IV. The Structure and Ionic Transport of Liquid Semiconductor NiTe

Akdere Ü., Günay S. D., Kavanoz H. B., Taşseven Ç.

Balkan Physics Letters, vol.15, pp.1-7, 2009 (Peer-Reviewed Journal)

Refereed Congress / Symposium Publications in Proceedings

I. THERMOPHYSICAL PROPERTIES OF NANO-CRYSTALLITE MAGNESIUM NITRIDE VIA MOLECULAR DYNAMICS SIMULATION AT CONSTANT PRESSURE

Aydın Y., Akdere Ü., Günay S. D., Taşseven Ç.

TURKISH PHYSICAL SOCIETY 35 INTERNATIONAL PHYSICS CONGRESS, Muğla, Turkey, 4 - 08 September 2019, pp.351

II. NVE SIMULATION AND HYPERNETTED-CHAIN CALCULATION OF THE SUPERIONIC CONDUCTOR LITHIUM NITRIDE

Akdere Ü., Günay S. D.

TURKISH PHYSICAL SOCIETY 35 INTERNATIONAL PHYSICS CONGRESS, Muğla, Turkey, 4 - 08 September 2019, pp.207

III. INTERACTION POTENTIAL AND LOCAL STRUCTURE OF LIQUID MAGNESIUM NITRIDE VIA HYPERNETTED-CHAIN THEORY

Aydın Y., Akdere Ü., Günay S. D., Taşseven Ç.

TURKISH PHYSICAL SOCIETY 35 INTERNATIONAL PHYSICS CONGRESS, Muğla, Turkey, 4 - 08 September 2019, pp.342

IV. Static Structure and Ionic Diffusion of Liquid Magnesium Nitrate

AKDERE Ü., AYDIN Y., GÜNAY S. D., TAŞSEVEN Ç.

ICSEEC: SUSTAINABLE ENERGY AND ENERGY CALCULATIONS, Muğla, Turkey, 12 April 2019, pp.49

V. Simulation of Pullout Test on Crystallite Segment of Antiparallel beta-Sheets of Bombyx Mori Silk Fibroin

Uğuz C., AKDERE Ü., GÜNAY S. D., TAŞSEVEN Ç.

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- VI. **Hydrogen Bond Analysis of Bombyx Mori Silk Fibroin by MD Simulation at Room Temperature**
Akdere Ü., Günay S. D., Taşseven Ç.
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- VII. **Simulations of Mechanical Properties of Crystalline Units of Bombyx Mori Silkworm Silk**
Akdere Ü., Uğuz C.
Turkish Physical Society 34 International Physics Congress, Muğla, Turkey, 5 - 09 September 2018, pp.297
- VIII. **Hydrogen Bond Analysis of Bombyx Mori Silk Fibroin by MD Simulation at Room Temperature**
AKDERE Ü., GÜNAY S. D., TAŞSEVEN Ç.
Turkish Physical Society 34rd International Physics Congress, 5 - 09 September 2018
- IX. **Structural phase transition of Mg₃As₂**
AKDERE Ü., GÜNAY S. D., TAŞSEVEN Ç.
Turkish Physical Society 33rd International Physics Congress, Muğla, Turkey, 6 - 10 September 2017
- X. **Temperature dependence of mechanical properties of anti-parallel beta sheets with bombyx-mori silk nano structures along chain direction**
AKDERE Ü., GÜNAY S. D., AKSAKAL B., ÇAĞIN T., TAŞSEVEN Ç.
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- XI. **Polar/Antipolar - Antiparallel beta sheets of bombyx mori silk nano crystallites solvated in different type of water**
Uğuz C., TAŞSEVEN Ç., GÜNAY S. D., AKDERE Ü.
JAPMED 10, İzmir, Turkey, 4 - 08 July 2017
- XII. **Low- and high- temperature structural analysis of magnesium antimonide**
GÜNAY S. D., AKDERE Ü., TAŞSEVEN Ç.
1st International Conference on Energy and Thermal Engineering, ICTE 2017, İstanbul, Turkey, 25 - 28 April 2017
- XIII. **Analysis of Mechanical Behavior of Anti Parallel Beta Sheets with Bombyx Mori Silk Nano Structures Along Inter Sheet Direction**
Akdere Ü., Günay S. D., Aksakal B., Çağın T., Taşseven Ç.
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- XIV. **Thermomechanical properties of anti parallel sheets with Bombyx mori silk nano structures Gly Ser Gly Ala Gly Ala n and Gly Ala n**
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- XV. **Thermophysical properties of anti parallel sheets with Bombyx mori silk nano structures Gly Ser Gly Ala Gly Ala n and Gly Ala n**
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- XVI. **Thermophysical Properties of alpha-Pu₂O₃: A New Potential Model**
GÜNAY Ş., AKGENC B., AKDERE Ü., TAŞSEVEN Ç.
3rd International Congress on Advances in Applied Physics and Materials Science, Antalya, Turkey, 24 - 28 April 2013, vol.1569, pp.208-211
- XVII. **NPT simulation and hypernetted-chain calculations of SrCl₂**
Akgenc B., AKDERE Ü., GÜNAY Ş., TAŞSEVEN Ç.
3rd International Advances in Applied Physics and Materials Science Congress, APMAS 2013, Antalya, Turkey, 24 - 28 April 2013, vol.1569, pp.15-18
- XVIII. **Thermophysical Properties of AgBr and AgCl from Molecular Dynamics Simulation**
Akdere Ü., Günay S. D., Taşseven Ç.

- Turkish Physical Society 29 International Physics Congress, Muğla, Turkey, 5 - 08 September 2012, pp.524
- XIX. Liquid State Theory of Plutonium Dioxide: Hypernetted Chain Approximation**
Akdere Ü., Akgeç B., Taşseven Ç., Günay S. D.
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- XX. A Study of Low Temperature Thermal Expansion and Structural Behaviour of Stromeyerite AgCuS**
Akdere Ü., Öztekin H. Ö., Yılmaz M., Kavanoz H. B.
Turkish Physical Society 29 International Physics Congress, Muğla, Turkey, 5 - 08 September 2012, pp.655
- XXI. Temperature Dependence of Ionic Diffusion of Thorium Dioxide**
Akdere Ü., Akgeç B., Günay S. D., Taşseven Ç.
Turkish Physical Society 29 International Physics Congress, Muğla, Turkey, 5 - 08 September 2012, pp.156
- XXII. Studying Superionic Phase Transition of Plutonium Dioxide From Molecular Dynamics Simulation**
Akdere Ü., Günay S. D., Akgeç B., Taşseven Ç.
Turkish Physical Society 29 International Physics Congress, Muğla, Turkey, 5 - 08 September 2012, pp.578
- XXIII. Thermophysical Properties of Liquid Thallium Halides (TlBr, TlCl, TlI) from Hypernetted Chain Theory**
Akdere Ü., Yılmaz M., Kavanoz H. B., Öztekin H. Ö.
Turkish Physical Society 28 International Physics Congress, Muğla, Turkey, 6 - 09 September 2011, pp.825
- XXIV. Hypernetted Chain Theory of Liquid GeO₂: Preliminary Results**
Akdere Ü., Günay S. D., Taşseven Ç.
Turkish Physical Society 28 International Physics Congress, Muğla, Turkey, 6 - 09 September 2011, pp.648
- XXV. HNC Calculations of Thermodynamic Properties of Molten AgI**
Günay S. D., Çalışkan M., Akdere Ü., Taşseven Ç.
Turkish Physical Society 28th International Physics Conference, 6 - 09 September 2011
- XXVI. Molecular dynamics simulation of thermal properties of Ag₃SI**
Öztekin H., KAVANOZ H. B., AKDERE Ü., YILMAZ M., Taşseven Ç.
7th International Conference of the Balkan Physical Union, Alexandroupoli, Greece, 9 - 13 September 2009, vol.1203, pp.252-254
- XXVII. Classical Molecular Dynamics Simulation of Mg₃Bi₂ at Solid and Liquid Phase**
Akdere Ü., Kayadibi F., Günay S. D., Taşseven Ç.
Turkish Physical Society 25th International Physics Congress, Muğla, Turkey, 25 - 29 August 2008, pp.176
- XXVIII. Static Structure and Ionic Transport in Molten NiTe and NiTe₂**
GÜNAY S. D., Akdere Ü., Taşseven Ç.
Turkish Physical Society 25th International Physics Conference, 25 - 29 August 2008
- XXIX. The structure and transport properties of β-Mg₃Bi₂ in superionic conduction and molten phase**
AKDERE Ü., GÜNAY Ş.
6TH International Conference of the Balkan Physical Union, İstanbul, Turkey, 22 August 2006 - 26 August 2007, vol.899, pp.577
- XXX. Transport properties of uranium dioxide by molecular dynamics simulation**
Gunay S. D., Akdere U., Kavanoz B., Taşseven Ç.
International Conference on Computational Methods in Science and Engineering, Corfu, Greece, 25 - 30 September 2007, vol.2, pp.1212-1215
- XXXI. The Static Structure of Liquid Semiconductors NiTe and NiTe₂ Preliminary Results**
GÜNAY S. D., AKDERE Ü., KOSAVALI F., TAŞSEVEN Ç.
AIP Conference Proceedings, 22 - 26 August 2006, vol.899, pp.616
- XXXII. Equilibrium Molecular Dynamics Calculations of The Transport Properties of Molten Thallium Halides**
GÜNAY S. D., AKDERE Ü., KOSAVALI F., TAŞSEVEN Ç.
AIP Conference Proceedings, 22 - 26 August 2006, vol.899, pp.605
- XXXIII. Hypernetted Chain Theory and Molecular Dynamic Simulation of Static Structure and Ionic Transport in Molten UO₂ and UCl₃**
Günay S. D., Akdere Ü., Kavanoz H. B., Yılmaz M., Taşseven Ç.

Turkish Physical Society 23th Conference, Muğla, Turkey, 13 - 16 September 2005

XXXIV. Static Criteria for Freezing of Liquid Halide Salts

Akdere Ü., Yılmaz M., Kavanoz H. B., Taşseven Ç.

Turkish Physical Society 23 International Physics Congress, Muğla, Turkey, 13 - 16 September 2005, pp.641

XXXV. The Static and Thermodynamic Properties of Molten UO₂ and UCl₃ (Preliminary Results)

Akdere Ü., Taşseven Ç., Kavanoz H. B.

The Fifth International Conference on Chemical Physics, İstanbul, Turkey, 31 October - 01 November 2002, pp.1-2

Supported Projects

GÜNAY S. D., Aydın Y., AKDERE Ü., TAŞSEVEN Ç., Project Supported by Higher Education Institutions, Magnezyum-Azot Grubu Bileşiklerinin Bilgisayar Simülasyonu Yöntemleriyle Fiziksel Özelliklerinin İncelenmesi, 2018 - 2020

Akdere Ü., TUBITAK Project, Thermo-mechanical Properties of Silk Fibroin Fibers from Molecular Dynamics Simulation for High Performance Applications, 2016 - 2018

Akdere Ü., Taşseven Ç., GÜNAY S. D., Project Supported by Higher Education Institutions, Süperiyonik İletkenlerin Yüksek Sıcaklık Davranışları, 2003 - 2007

Metrics

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