

## Prof.Dr. MURAT DURANDURDU

### Kişisel Bilgiler

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### Eğitim Bilgileri

Doktora, Ohio University, College Of Science, Amerika Birleşik Devletleri 1999 - 2002

Yüksek Lisans, Virginia Polytechnic Institute and State University, College Of Engineering, Materials Science And Engineering, Amerika Birleşik Devletleri 1997 - 1999

Yüksek Lisans-Tezsiz, Rutgers, The State University of New Jersey, College Of Science, Physics, Amerika Birleşik Devletleri 1995 - 1997

Lisans, Karadeniz Teknik Üniversitesi, Fen Edebiyat Fakültesi, Fizik, Türkiye 1988 - 1992

### Yaptığı Tezler

Doktora, Polyamorphism in Semiconductors, Ohio University, College Of Science, Physics, 2002

Yüksek Lisans, Molecular Statics Simulation in Aluminum ?, Virginia Polytechnic Institute And State University, College Of Engineering , Materials Science And Engineering , 1999

### Araştırma Alanları

Fizik, Yoğun Madde 1:Yapısal, Mekanik ve Termal Özellikler , Durum yoğunluğu, faz dengesi ve faz geçişleri, Temel Bilimler

### Akademik Unvanlar / Görevler

Prof.Dr., Abdullah Gül Üniversitesi, Mühendislik Fakültesi, Nanoteknoloji Mühendisliği, 2017 - Devam Ediyor

Prof.Dr., Charles Üniversitesi, Faculty of Science , Department of Physical and Macromolecular Chemistry, 2020 - 2021

Doç.Dr., Abdullah Gül Üniversitesi, Mühendislik Fakültesi, Nanoteknoloji Mühendisliği, 2014 - 2017

Doç.Dr., Texas Tech University, College Of Science, Physics, 2011 - 2014

Yrd.Doç.Dr., University of Texas at El Paso, College Of Science, Physics, 2004 - 2011

Araştırma Görevlisi Dr., The University of Michigan, College Of Engineering, Materials Science And Engineering, 2002 - 2004

Araştırma Görevlisi, Ohio University, College Of Science, Physics, 1999 - 2002

Araştırma Görevlisi, Virginia Polytechnic Institute and State University, College Of Engineering, Materials Science And Engineering, 1997 - 1999

Araştırma Görevlisi, Karadeniz Teknik Üniversitesi, Fen Edebiyat Fakültesi, Fizik, 1993 - 1994

## Akademik İdari Deneyim

Anabilim/Bilim Dalı Başkanı, Abdullah Gül Üniversitesi, 2015 - 2022

Anabilim/Bilim Dalı Başkanı, Abdullah Gül Üniversitesi, İleri Malzemeler Ve Nanoteknoloji, 2015 - 2022

Bölüm Başkanı, Abdullah Gül Üniversitesi, 2015 - 2022

## SCI, SSCI ve AHCI İndekslerine Giren Dergilerde Yayınlanan Makaleler

- I. **Amorphous BC5 from first principles calculations**  
DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.592, 2022 (SCI-Expanded)
- II. **Theoretical investigation of substituent effects on the relative stabilities and electronic structure of [BnXn](2-) clusters**  
TAHAOĞLU D., ALKAN F., DURANDURDU M.  
JOURNAL OF MOLECULAR MODELING, cilt.27, sa.12, 2021 (SCI-Expanded)
- III. **Formation of a very high-density amorphous phase of carbon and its crystallization into a simple cubic structure at high pressure**  
DURANDURDU M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.200, 2021 (SCI-Expanded)
- IV. **Amorphous boron phosphide: An ab initio investigation**  
BOLAT S., Durandurdu M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.570, 2021 (SCI-Expanded)
- V. **Amorphous zircon at high pressure**  
BOLAT S., DURANDURDU M.  
JOURNAL OF PHYSICS AND CHEMISTRY OF SOLIDS, cilt.153, 2021 (SCI-Expanded)
- VI. **A first principles study of amorphous and crystalline silicon tetraboride**  
Karacaoglu A. O., DURANDURDU M.  
MATERIALS CHEMISTRY AND PHYSICS, cilt.258, 2021 (SCI-Expanded)
- VII. **Stoichiometric amorphous boron carbide (BC)**  
Yildiz T. A., DURANDURDU M.  
JOURNAL OF MATERIALS SCIENCE, cilt.55, sa.30, ss.14709-14716, 2020 (SCI-Expanded)
- VIII. **Amorphous silicon hexaboride at high pressure**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.100, sa.14, ss.1818-1833, 2020 (SCI-Expanded)
- IX. **Ab initio simulation of amorphous BC3**  
DURANDURDU M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.178, 2020 (SCI-Expanded)
- X. **Amorphous silicon triboride: A first principles study**  
Ozlem A., Karacaoglan C., DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.536, 2020 (SCI-Expanded)
- XI. **High pressure modifications in amorphous boron suboxide: An ab initio study**  
DURANDURDU M.  
CERAMICS INTERNATIONAL, cilt.46, sa.5, ss.5968-5975, 2020 (SCI-Expanded)
- XII. **Tetrahedral amorphous boron nitride: A hard material**  
DURANDURDU M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.103, sa.2, ss.973-978, 2020 (SCI-Expanded)
- XIII. **Amorphous boron carbide from ab initio simulations**  
Yildiz T. A., DURANDURDU M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.173, 2020 (SCI-Expanded)
- XIV. **Pressure-induced amorphization, mechanical and electronic properties of zeolitic imidazolate**

**framework (ZIF-8)**

ERKARTAL M., DURANDURDU M.

MATERIALS CHEMISTRY AND PHYSICS, cilt.240, 2020 (SCI-Expanded)

- XV. **Phase transition of ZrN under pressure**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.99, sa.8, ss.942-955, 2019 (SCI-Expanded)
- XVI. **Hydrogenated amorphous boron nitride: A first principles study**  
Uchoyuk T. A., DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.502, ss.159-163, 2018 (SCI-Expanded)
- XVII. **Amorphous zirconia at high pressure**  
DURANDURDU M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.101, sa.12, ss.5411-5418, 2018 (SCI-Expanded)
- XVIII. **Solute aggregation in Ca<sub>72</sub>Zn<sub>28</sub> metallic glass**  
TAHAOĞLU D., DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.500, ss.410-416, 2018 (SCI-Expanded)
- XIX. **Amorphous magnesium silicide**  
DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.498, ss.118-124, 2018 (SCI-Expanded)
- XX. **An in-depth investigation of Mg-Zn-Ca metallic glasses: A first principles study**  
ERKARTAL M., DURANDURDU M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.153, ss.326-337, 2018 (SCI-Expanded)
- XXI. **Pressure-Induced Amorphization of MOF-5: A First Principles Study**  
ERKARTAL M., DURANDURDU M.  
CHEMISTRYSELECT, cilt.3, sa.28, ss.8056-8063, 2018 (SCI-Expanded)
- XXII. **Hard boron rich boron nitride nanoglasses**  
Cetin A. O., DURANDURDU M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.101, sa.5, ss.1929-1939, 2018 (SCI-Expanded)
- XXIII. **Permanent densification of amorphous zinc oxide under pressure: A first principles study**  
TAHAOĞLU D., DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.481, ss.27-32, 2018 (SCI-Expanded)
- XXIV. **Amorphous silicon hexaboride: a first-principles study**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.98, sa.30, ss.2723-2733, 2018 (SCI-Expanded)
- XXV. **MgCu metallic glass**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.98, sa.8, ss.633-645, 2018 (SCI-Expanded)
- XXVI. **Two successive amorphous-to-amorphous phase transformations in TiO<sub>2</sub>**  
DURANDURDU M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.100, sa.9, ss.3903-3911, 2017 (SCI-Expanded)
- XXVII. **Densification of amorphous boron under pressure**  
DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.471, ss.274-279, 2017 (SCI-Expanded)
- XXVIII. **Amorphous zirconia: ab initio molecular dynamics simulations**  
DURANDURDU M.  
Philosophical Magazine, cilt.97, sa.16, ss.1334-1345, 2017 (SCI-Expanded)
- XXIX. **High-density amorphous phase of CdO**  
DURANDURDU M.  
Journal of Non-Crystalline Solids, cilt.463, ss.64-67, 2017 (SCI-Expanded)
- XXX. **Ferromagnetism in amorphous MgO**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.97, sa.24, ss.2129-2141, 2017 (SCI-Expanded)

- XXXI. **Nanosegregated amorphous AlBN2 alloy**  
DURANDURDU M.  
Philosophical Magazine, cilt.96, sa.30, ss.3200-3210, 2016 (SCI-Expanded)
- XXXII. **Polyamorphism in Aluminum Nitride: A First Principles Molecular Dynamics Study**  
DURANDURDU M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.99, sa.5, ss.1594-1600, 2016 (SCI-Expanded)
- XXXIII. **n-type conductivity in Si-doped amorphous AlN: an ab initio investigation**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.96, sa.11, ss.1110-1121, 2016 (SCI-Expanded)
- XXXIV. **Pressure-induced phase transformations in amorphous arsenic**  
DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.437, ss.6-9, 2016 (SCI-Expanded)
- XXXV. **Local structure of As2O3 glass from first principles simulations**  
DURANDURDU M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.436, ss.18-21, 2016 (SCI-Expanded)
- XXXVI. **Amorphous boron nitride at high pressure**  
DURANDURDU M.  
PHILOSOPHICAL MAGAZINE, cilt.96, sa.18, ss.1950-1964, 2016 (SCI-Expanded)
- XXXVII. **Hexagonal nanosheets in amorphous BN: A first principles study**  
Durandurdu M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.427, ss.41-45, 2015 (SCI-Expanded)
- XXXVIII. **Novel high-pressure phase of ZrO2: An ab initio prediction**  
Durandurdu M.  
JOURNAL OF SOLID STATE CHEMISTRY, cilt.230, ss.233-236, 2015 (SCI-Expanded)
- XXXIX. **High-pressure phase transitions of TiN: an ab initio constant pressure study**  
Durandurdu M.  
PHILOSOPHICAL MAGAZINE, cilt.95, sa.22, ss.2376-2384, 2015 (SCI-Expanded)
- XL. **Liquid boron and amorphous boron: An ab initio molecular dynamics study**  
Durandurdu M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.417, ss.10-14, 2015 (SCI-Expanded)
- XLI. **Uncovering Nanoclusters in Amorphous AlN: An Ab Initio Study**  
Durandurdu M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.98, sa.4, ss.1095-1098, 2015 (SCI-Expanded)
- XLII. **Atomic structure of amorphous CdO from first principles simulations**  
Durandurdu M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.412, ss.11-15, 2015 (SCI-Expanded)
- XLIII. **New high-pressure phase of MgH2: An ab initio constant-pressure study**  
Durandurdu M.  
EPL, cilt.105, sa.4, 2014 (SCI-Expanded)
- XLIV. **Atomic structure of amorphous Mg40Cu35Ti25 alloy: An ab initio molecular dynamics study**  
Durandurdu M.  
SOLID STATE COMMUNICATIONS, cilt.154, ss.30-33, 2013 (SCI-Expanded)
- XLV. **Ab initio modeling of metallic Pd80Si20 glass**  
Durandurdu M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.65, ss.44-47, 2012 (SCI-Expanded)
- XLVI. **Nanoscale icosahedral packing in amorphous Mg50Ni50: An ab initio study**  
Tetik E., Durandurdu M., KARADAĞ F.  
EPL, cilt.100, sa.2, 2012 (SCI-Expanded)
- XLVII. **Nanosize icosahedral quasicrystal in Mg90Ca10 glass: An ab initio molecular dynamics study**  
Durandurdu M.  
JOURNAL OF CHEMICAL PHYSICS, cilt.137, sa.3, 2012 (SCI-Expanded)

- XLVIII. Formation of Cotunnite Phase in ZrO<sub>2</sub> under Uniaxial Stress: A First Principles Study**  
Ozturkz H., Durandurdu M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.94, sa.3, ss.932-937, 2011 (SCI-Expanded)
- XLIX. Formation of a two-dimensional layered structure in silica under shear stresses: An ab initio study**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.81, sa.17, 2010 (SCI-Expanded)
- L. Formation of a C<sub>2</sub>cm phase in SnS at high pressure; an ab initio constant pressure study**  
Alptekin S., Durandurdu M.  
SOLID STATE COMMUNICATIONS, cilt.150, ss.870-874, 2010 (SCI-Expanded)
- LI. Vibrational properties of amorphous germanium under pressure and its thermal expansion and Gruneisen parameters**  
Durandurdu M.  
JOURNAL OF NON-CRYSTALLINE SOLIDS, cilt.356, ss.977-981, 2010 (SCI-Expanded)
- LII. Ab initio molecular dynamics study of pressure-induced phase transformation in KCl**  
Durandurdu M.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.48, sa.3, ss.672-676, 2010 (SCI-Expanded)
- LIII. Formation of Anatase Phase in HfO<sub>2</sub> in Tensile Stress: An Ab Initio Study**  
Durandurdu M.  
JOURNAL OF THE AMERICAN CERAMIC SOCIETY, cilt.93, sa.5, ss.1467-1469, 2010 (SCI-Expanded)
- LIV. Orthorhombic intermediate phases for the wurtzite-to-rocksalt phase transformation of CdSe: An ab initio constant pressure study**  
Durandurdu M.  
CHEMICAL PHYSICS, cilt.369, ss.55-58, 2010 (SCI-Expanded)
- LV. Pressure-induced phase transformation of BaS: An ab initio constant pressure study**  
Durandurdu M.  
CHEMICAL PHYSICS, cilt.367, ss.80-82, 2010 (SCI-Expanded)
- LVI. First principles study of structural phase stability of wide-gap semiconductors MgTe, MgS and MgSe**  
Gokoglu G., Durandurdu M., Gulseren O.  
COMPUTATIONAL MATERIALS SCIENCE, cilt.47, sa.2, ss.593-598, 2009 (SCI-Expanded)
- LVII. Expanded phase of ZrO<sub>2</sub>: An ab initio constant-pressure study**  
Durandurdu M.  
EPL, cilt.88, sa.6, 2009 (SCI-Expanded)
- LVIII. New transformation mechanism for a zinc-blende to rocksalt phase transformation in MgS**  
Durandurdu M.  
JOURNAL OF PHYSICS-CONDENSED MATTER, cilt.21, sa.45, 2009 (SCI-Expanded)
- LIX. Fcc-to-bct phase transformation of aluminum under triaxial stresses: an ab initio constant pressure study**  
Durandurdu M.  
EUROPEAN PHYSICAL JOURNAL B, cilt.72, sa.2, ss.241-245, 2009 (SCI-Expanded)
- LX. Pressure-induced phase transition in AlN: An ab initio molecular dynamics study**  
Durandurdu M.  
JOURNAL OF ALLOYS AND COMPOUNDS, cilt.480, sa.2, ss.917-921, 2009 (SCI-Expanded)
- LXI. Formation of an anataselike phase in silica under anisotropic stress: An ab initio constant-pressure study**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.80, sa.2, 2009 (SCI-Expanded)
- LXII. High-density amorphous phase of GeS<sub>2</sub> glass under pressure**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.79, sa.20, 2009 (SCI-Expanded)
- LXIII. High-pressure phases of ZrO<sub>2</sub>: An ab initio constant-pressure study**  
Ozturk H., Durandurdu M.

- PHYSICAL REVIEW B, cilt.79, sa.13, 2009 (SCI-Expanded)
- LXIV. **The structural phase transition of ZnSe under hydrostatic and nonhydrostatic compressions: an ab initio molecular dynamics study**  
Durandurdu M.  
JOURNAL OF PHYSICS-CONDENSED MATTER, cilt.21, sa.12, 2009 (SCI-Expanded)
- LXV. **Pressure-induced phase transition of BeO**  
ALPTEKIN S., Durandurdu M.  
SOLID STATE COMMUNICATIONS, cilt.149, ss.345-348, 2009 (SCI-Expanded)
- LXVI. **Pressure-induced phase transition in wurtzite ZnS: An ab initio constant pressure study**  
Durandurdu M.  
JOURNAL OF PHYSICS AND CHEMISTRY OF SOLIDS, cilt.70, ss.645-649, 2009 (SCI-Expanded)
- LXVII. **An ab initio constant-pressure study of pressure-induced phase transition of MgSe**  
Oezduran M., Durandurdu M.  
EPL, cilt.84, sa.5, 2008 (SCI-Expanded)
- LXVIII. **Pressure-induced phase transition of zinc-blende AlN: An ab initio molecular dynamics study**  
Durandurdu M.  
JOURNAL OF PHYSICS AND CHEMISTRY OF SOLIDS, cilt.69, sa.11, ss.2894-2897, 2008 (SCI-Expanded)
- LXIX. **Phase transformation of 6H-SiC at high pressure: An ab initio constant-pressure study**  
Eker S., Durandurdu M.  
EPL, cilt.84, sa.2, 2008 (SCI-Expanded)
- LXX. **Diamond to beta-Sn phase transition of silicon under hydrostatic and nonhydrostatic compressions**  
Durandurdu M.  
JOURNAL OF PHYSICS-CONDENSED MATTER, cilt.20, sa.32, 2008 (SCI-Expanded)
- LXXI. **New B2O3 crystals predicted from concurrent molecular dynamics simulations and first-principles calculations**  
HUANG L., Durandurdu M., KIEFFER J.  
JOURNAL OF PHYSICAL CHEMISTRY C, cilt.111, sa.37, ss.13712-13720, 2007 (SCI-Expanded)
- LXXII. **Structural phase transition of gold under uniaxial, tensile, and triaxial stresses: An ab initio study**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.76, sa.2, 2007 (SCI-Expanded)
- LXXIII. **Ab initio simulations of the structural phase transformation of 2H-SiC at high pressure**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.75, sa.23, 2007 (SCI-Expanded)
- LXXIV. **Transformation pathways of silica under high pressure**  
HUANG L., Durandurdu M., KIEFFER J.  
NATURE MATERIALS, cilt.5, sa.12, ss.977-981, 2006 (SCI-Expanded)
- LXXV. **Ab initio molecular dynamics study of pressure-induced phase transition in ZnS**  
MARTINEZ I., Durandurdu M.  
JOURNAL OF PHYSICS-CONDENSED MATTER, cilt.18, sa.41, ss.9483-9491, 2006 (SCI-Expanded)
- LXXVI. **Transition pathway in GaAs under uniaxial stress: an ab initio study**  
Durandurdu M.  
JOURNAL OF PHYSICS-CONDENSED MATTER, cilt.18, sa.20, ss.4887-4894, 2006 (SCI-Expanded)
- LXXVII. **Electronic and mechanical properties of wurtzite type SiC nanowires**  
Durandurdu M.  
PHYSICA STATUS SOLIDI B-BASIC SOLID STATE PHYSICS, cilt.243, sa.5, 2006 (SCI-Expanded)
- LXXVIII. **Ab initio modeling of small diameter silicon nanowires**  
Durandurdu M.  
PHYSICA STATUS SOLIDI B-BASIC SOLID STATE PHYSICS, cilt.243, sa.2, 2006 (SCI-Expanded)
- LXXIX. **Ab initio simulation of polyamorphic phase transition in hydrogenated silicon**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.73, sa.3, 2006 (SCI-Expanded)

- LXXX. **Phase transition of GeSe<sub>2</sub> at high pressure**  
Durandurdu M.  
PHYSICA STATUS SOLIDI B-BASIC SOLID STATE PHYSICS, cilt.242, sa.15, ss.3085-3090, 2005 (SCI-Expanded)
- LXXXI. **Cmcm phase of GeS at high pressure**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.72, sa.14, 2005 (SCI-Expanded)
- LXXXII. **Ab initio simulation of the rhombohedral-to-simple-cubic transition in arsenic**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.72, sa.7, 2005 (SCI-Expanded)
- LXXXIII. **Structural phase transition of germanium under uniaxial stress: An ab initio study**  
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PHYSICAL REVIEW B, cilt.71, sa.5, 2005 (SCI-Expanded)
- LXXXIV. **Mechanically controlled, seeded formation of a nanoscale metastable phase in ionic compounds**  
Palko J., Durandurdu M., Kieffer J.  
NANO LETTERS, cilt.4, sa.9, ss.1769-1773, 2004 (SCI-Expanded)
- LXXXV. **Pressure-induced amorphous-to-amorphous phase transition in GaAs**  
Durandurdu M.  
PHYSICAL REVIEW B, cilt.70, sa.8, 2004 (SCI-Expanded)
- LXXXVI. **High-pressure phases of amorphous and crystalline silicon**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.67, sa.21, 2003 (SCI-Expanded)
- LXXXVII. **Pressure-induced structural phase transition of paracrystalline silicon**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.66, sa.20, 2002 (SCI-Expanded)
- LXXXVIII. **Ab initio simulation of pressure-induced low-energy excitations in amorphous silicon**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.66, sa.15, 2002 (SCI-Expanded)
- LXXXIX. **Ab initio simulation of high-pressure phases of GaAs**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.66, sa.4, 2002 (SCI-Expanded)
- XC. **First-order pressure-induced polyamorphism in germanium**  
Durandurdu M., DRABOLD D.  
Physical Review B - Condensed Matter and Materials Physics, cilt.66, sa.4, ss.412011-412014, 2002 (SCI-Expanded)
- XCI. **Simulation of pressure-induced polyamorphism in a chalcogenide glass GeSe<sub>2</sub>**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.65, sa.10, 2002 (SCI-Expanded)
- XCII. **Ab initio simulation of first-order amorphous-to-amorphous phase transition of silicon**  
Durandurdu M., DRABOLD D.  
PHYSICAL REVIEW B, cilt.64, sa.1, 2001 (SCI-Expanded)
- XCIII. **Approximate ab initio calculations of electronic structure of amorphous silicon**  
Durandurdu M., DRABOLD D., MOUSSEAU N.  
PHYSICAL REVIEW B, cilt.62, sa.23, ss.15307-15310, 2000 (SCI-Expanded)

## **Hakemli Kongre / Sempozyum Bildiri Kitaplarında Yer Alan Yayınlar**

- I. **Derivative Structures of Closo Boron Hydride Cluster, B<sub>16</sub>X<sub>16</sub>:A DFT Study**  
TAHAOĞLU D., DURANDURDU M.  
CSC2018: International Computational Science Congress, 26 - 28 Ekim 2018
- II. **CMCM INTERMEDIATE PHASE FOR THE HCP-TO-BCC PHASE TRANSITION IN MAGNESIUM**  
DURANDURDU M., Çetin A. Ö., Üçhöyük T. A.

3.ULUSLARARASI MALZEME BILIMI VE TEKNOLOJISI KONFERANSI (KAPADOKYA), 17 - 19 Eylül 2018

III. **Stability Investigation of B16H162- Derivatives by DFT Calculations**

TAHAOĞLU D., DURANDURDU M.

International Conference on Materilas Science and Technology in Cappadocia, 17 - 19 Eylül 2018

IV. **Effects of Production Methods on Structural and Mechanical Properties for Metallic Ca72Zn28 Glass- A DFT Study**

TAHAOĞLU D., DURANDURDU M.

The Internatnional Conference On Materials Science Mechanical And Automation Engineerings And Technology (IMSMATEC'18), 10 - 12 Nisan 2018

V. **Polyamorphism in TiO2**

DURANDURDU M.

ncon-International Congress on Chemistry and Materials Science, 5 - 07 Ekim 2017

VI. **Metallic Ca72Zn28 Glass from First Principle Molecular Dynamics Simulations**

TAHAOĞLU D., DURANDURDU M.

International Congress on Chemistry and Materials Science, Ankara, Türkiye, 5 - 07 Ekim 2017

VII. **Investigation of Phase Transformation of Amorphous Zinc Oxide under Pressure: a DFT Study**

TAHAOĞLU D., DURANDURDU M.

Türk Fizik Derneği 33. Uluslararası Fizik Kongresi, Bodrum, Türkiye, 6 - 10 Eylül 2017

VIII. **POLYAMORPHIC PHASE TRANSITION IN ALUMINUM NITRIDE**

DURANDURDU M.

Uluslararası Malzeme Bilimi ve Teknolojisi Konferansı Kapadokya (IMSTEC'16), 6 - 08 Nisan 2016

IX. **Hidden Nanoclusters in Amorphous AlN A First Principles Study**

DURANDURDU M.

International Semiconductor Science and Technology Conference 2015 (ISSTC2015), 11 - 13 Mayıs 2015

## Desteklenen Projeler

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