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#### Education Information

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- I. Doctorate, Gazi University, Fen Bilimleri Enstitüsü, Fizik (Dr), Turkey 1996 - 2000
- II. Postgraduate, Gazi University, Fen Bilimleri Enstitüsü, Fizik (YI) (Tezli), Turkey 1993 - 1996
- III. Undergraduate, Gazi University, Fen-Edebiyat Fakültesi, Fizik Pr., Turkey 1989 - 1993

#### Foreign Languages

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- I. English, B2 Upper Intermediate

#### Research Areas

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Physics

#### Academic Titles / Tasks

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- I. Professor, Gazi University, Fen Fakültesi, Fizik, 2014 - Continues
- II. Assistant Professor, Gazi University, Fen Fakültesi, Fizik, 2003 - 2008
- III. Research Assistant, Gazi University, Fen Fakültesi, Fizik, 1994 - 2003

#### Advising Theses

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- I. ÇİFTÇİ Y., The Investigation Of Hydrogen Storage On Some Light Elements Decorated Two Dimensional Carbon Allotropes Using Ab-Initio Methods, Doctorate, İ.ALP(Student), 2019
- II. ÇİFTÇİ Y., LuXO<sub>3</sub> (X=Al, Ga, In) perovskit bileşiklerinin temel fiziksel özelliklerinin ab initio yöntemlerle incelenmesi, Postgraduate, G.SÜRÜCÜ(Student), 2018
- III. ÇİFTÇİ Y., Bazı AIBIVC<sub>2</sub>V tipinde yarıiletken kalkopiritlerin temel fiziksel özelliklerinin ab-initio metot ile incelenmesi, Doctorate, B.KOÇAK(Student), 2016
- IV. ÇİFTÇİ Y., Bazı ikili ve/veya üçlü alaşımların temel fiziksel özelliklerinin yoğunluk fonksiyoneli teorisi ile incelenmesi, Doctorate, Y.MOĞULKOÇ(Student), 2013
- V. ÇİFTÇİ Y., THE INVESTIGATION SOME THE STRUCTURAL, ELECTRONIC AND ELASTIC PROPERTIES OF ThX (X =N, As, P, Sb) BY AB INITIO METHOD, Postgraduate, S.ARSLAN(Student), 2008
- VI. ÇİFTÇİ Y., STRUCTURAL PHASE TRANSITIONS AND ELASTIC STABILITY WITH EMBEDDED ATOM METHOD, Postgraduate, R.YİĞİTER(Student), 2007

- VII. ÇİFTÇİ Y., Embedded-atom modeli ile yapısal faz dönüşümleri ve elastik kararlılık, Postgraduate, R.YİĞİTER(Student), 2007

Published journal articles indexed by SCI, SSCI, and AHCI

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- I. **Thermodynamics evaluation of half-metallic ferromagnetism in N-doped  $\text{XTiO}_3$  ( $X = \text{Ca, Sr, and Ba}$ ) systems: DFT calculations**  
Elahi I., ÇİFTÇİ Y., Akbar W.  
Materials Today Communications, vol.34, 2023 (SCI-Expanded)
- II. **The mechanical, dynamical, thermodynamical properties and elastic anisotropies of cubic  $\text{YbAu}$  compound under pressure**  
KARS DURUKAN İ., EVECEN M., ÇİFTÇİ Y.  
MATERIALS TODAY COMMUNICATIONS, vol.33, 2022 (SCI-Expanded)
- III. **A first-principles prediction on the structural, electronic, elastic, phonon, and transport properties of  $\text{BaSiN}_2$**   
ÇİFTÇİ Y., ALP İ.  
INDIAN JOURNAL OF PHYSICS, vol.96, no.14, pp.4131-4141, 2022 (SCI-Expanded)
- IV. **Systematic study of optoelectronic and thermoelectric properties of new lead-free halide double perovskites  $\text{A}(2)\text{KGe}(6)$  ( $A = \text{Cs, Rb}$ ) for solar cell applications via ab-initio calculations**  
Mukhtar M. W., Ramzan M., Rashid M., Hussain A., Naz G., ÇİFTÇİ Y., Dahshan A., Znaidia S.  
MATERIALS SCIENCE AND ENGINEERING B-ADVANCED FUNCTIONAL SOLID-STATE MATERIALS, vol.285, 2022 (SCI-Expanded)
- V. **Pressure effects on structural, electronic and anisotropic elastic properties of Si doped  $\text{RuGe}$  compound with different concentrations by first-principles calculations**  
ÇİFTÇİ Y., ÇOBAN C., EVECEN M., Durukan I. K.  
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- VI. **First-principles study on B2 based  $\text{XAl}$  ( $X = \text{Rh, Ru}$ ) compounds**  
KARS DURUKAN İ., ÇİFTÇİ Y.  
PHYSICA SCRIPTA, vol.96, no.12, 2021 (SCI-Expanded)
- VII. **First-principles calculations of vibrational and optical properties of half-Heusler  $\text{NaScSi}$**   
Kars Durukan İ., Oztekin Ciftci Y.  
INDIAN JOURNAL OF PHYSICS, vol.95, no.11, pp.2303-2312, 2021 (SCI-Expanded)
- VIII. **Ab-initio study on physical properties of intermetallic  $\text{LiPb}$  compound**  
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- IX. **Band Alignment in Monolayer Boron Phosphide with Janus  $\text{MoSSe}$  Heterobilayers under Strain and Electric Field**  
Mogulkoc Y., Caglayan R., Çiftci Y.  
PHYSICAL REVIEW APPLIED, vol.16, no.2, 2021 (SCI-Expanded)
- X. **Mechanical and dynamic properties of stable two-dimensional boron-substituted  $\text{ThMoB}_4$ -type graphene: First-Principles Study**  
Aydın S., Çiftci Y.  
MATERIALS TODAY COMMUNICATIONS, vol.27, 2021 (SCI-Expanded)
- XI. **Ligand-free fabrication of  $\text{Au/TiO}_2$  nanostructures for plasmonic hot-electron-driven photocatalysis: Photoelectrochemical water splitting and organic-dye degradation**  
Celebi N., Aydın M. Y., Soysal F., Çiftci Y., Salimi K.  
JOURNAL OF ALLOYS AND COMPOUNDS, vol.860, 2021 (SCI-Expanded)
- XII. **Pressure effects on electronic, elastic, and vibration properties of metallic antiperovskite  $\text{PbNCa}_3$  by ab initio calculations**

- Ciftci Y., Evecen M., Alp İ.  
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- XIII. **Theoretical investigation of the electronic structure, elastic, dynamic properties of intermetallic compound NiBe under pressure**  
Evecen M., ÇİFTÇİ Y.  
EUROPEAN PHYSICAL JOURNAL B, vol.94, no.1, 2021 (SCI-Expanded)
- XIV. **The pressure effect on optoelectronic and mechanical properties of chalcopyrite BeSiN<sub>2</sub>**  
Ciftci Y., Alp İ.  
MATERIALS TODAY COMMUNICATIONS, vol.24, 2020 (SCI-Expanded)
- XV. **Anisotropic Elastic, Electronic and Vibrational Properties of the Semiconductor AgScX (X = Ge, C) Compounds**  
Kars Durukan İ., Çiftci Y.  
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- XVI. **A Theoretical Study of Pressure-Induced Effects on Phase Transition and Elastic Properties of AsTh Compound**  
ÇİFTÇİ Y., ATEŞER E.  
JOURNAL OF ELECTRONIC MATERIALS, vol.49, no.3, pp.2086-2094, 2020 (SCI-Expanded)
- XVII. **Green-Emitting Lead-Free Cs<sub>4</sub>SnBr<sub>6</sub> Zero-Dimensional Perovskite Nanocrystals with Improved Air Stability**  
Chiara R., ÇİFTÇİ Y., Queloz V. I. E., Nazeeruddin M. K., Grancini G., Malavasi L.  
JOURNAL OF PHYSICAL CHEMISTRY LETTERS, vol.11, no.3, pp.618-623, 2020 (SCI-Expanded)
- XVIII. **Exploring the role of halide mixing in lead-free BZA(2)SnX(4) two dimensional hybrid perovskites**  
Pisanu A., Coduri M., Morana M., ÇİFTÇİ Y., Rizzo A., Listorti A., Gaboardi M., Bindi L., Queloz V. I. E., Milanese C., et al.  
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- XIX. **Physical properties of ternary thallium chalcogenes Tl<sub>2</sub>(M)MQ<sub>3</sub> (M = Zr, Hf; Q = S, Se, Te) via ab-initio calculations**  
ATEŞER E., Okvuran O., ÇİFTÇİ Y., ÖZİŞİK H., DELİGÖZ E.  
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- XX. **The Effect of Pressure on Elastic Anisotropy, Vibration and Optical Properties of a AgScSi Compound**  
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JOURNAL OF ELECTRONIC MATERIALS, vol.48, no.6, pp.4050-4056, 2019 (SCI-Expanded)
- XXI. **Analysis of the structural, electronic, elastic and thermodynamic properties of CuAl<sub>2</sub>X<sub>4</sub> (X = O, S) spinel structure**  
Obeid M. M., Mogulkoc Y., Edrees S. J., Ciftci Y., Shukur M. M., Al-Marzooqee M. M. H.  
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- XXII. **First principle study of structural, electronic, mechanical, dynamic and optical properties of half-Heusler compound LiScSi under pressure**  
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- XXIII. **First-principles hydrogen adsorption properties of Li-decorated ThMoB<sub>4</sub>-type graphene**  
Alp İ., AYDIN S., ÇİFTÇİ Y.  
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- XXIV. **Physical Properties of Superhard Diamond-Like BC<sub>5</sub> from a First-Principles Study**  
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- XXV. **DFT-based ab-initio study of half-Heusler KCaP compound**  
Mogulkoc Y., Ciftci Y., Surucu G.  
JOURNAL OF OPTOELECTRONICS AND ADVANCED MATERIALS, vol.20, no.1-2, pp.61-68, 2018 (SCI-Expanded)
- XXVI. **First principle and tight-binding study of strained SnC**  
Mogulkoc Y., Modarresi M., Mogulkoc A., Ciftci Y., Alkan B.  
JOURNAL OF PHYSICS AND CHEMISTRY OF SOLIDS, vol.111, pp.458-463, 2017 (SCI-Expanded)

- XXVII. **A density functional study of chalcopyrite MgGeSb<sub>2</sub>**  
Kocak B., ÇİFTÇİ Y.  
INDIAN JOURNAL OF PHYSICS, vol.91, no.12, pp.1487-1492, 2017 (SCI-Expanded)
- XXVIII. **Structural phase transition, electronic, elastic, and vibrational properties of LiAl intermetallic compound: insights from first-principles calculations**  
Mogulkoc Y., Ciftci Y., Surucu G.  
CANADIAN JOURNAL OF PHYSICS, vol.95, no.8, pp.691-698, 2017 (SCI-Expanded)
- XXIX. **First-principles calculations of Mg<sub>1-x</sub>Cu<sub>x</sub>SiP<sub>2</sub> alloys with x=0.0, 0.25, 0.5, 0.75 and 1.0**  
Kocak B., Ciftci Y.  
JOURNAL OF ALLOYS AND COMPOUNDS, vol.705, pp.211-217, 2017 (SCI-Expanded)
- XXX. **First-Principles Study on the Structural, Elastic, Electronic and Vibrational Properties of Scandium Based Intermetallic Compounds (ScX, X = Co, Rh and Ir) Under Pressure**  
EVECEN M., ÇİFTÇİ Y.  
JOURNAL OF NANOELECTRONICS AND OPTOELECTRONICS, vol.12, no.2, pp.100-108, 2017 (SCI-Expanded)
- XXXI. **The structural, elastic, electronic and dynamical properties of chalcopyrite semiconductor BeGeAs<sub>2</sub> from first-principles calculations**  
ÇİFTÇİ Y., EVECEN M., ALDIRMAZ E.  
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- XXXII. **Structural and Thermoelectronic Properties of Chalcopyrite MgSiX<sub>2</sub> (X = P, As, Sb)**  
Kocak B., Ciftci Y., Surucu G.  
JOURNAL OF ELECTRONIC MATERIALS, vol.46, no.1, pp.247-264, 2017 (SCI-Expanded)
- XXXIII. **Structural, electronic, elastic, optical, and vibrational properties of HfXSb (X = Co, Rh, Ru) half-Heusler compounds: an ab initio study**  
Coban C., Ciftci Y., Colakoglu K.  
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- XXXIV. **Electronic and optical properties of bilayer blue phosphorus**  
Mogulkoc Y., Modarresi M., Mogulkoc A., Ciftci Y.  
COMPUTATIONAL MATERIALS SCIENCE, vol.124, pp.23-29, 2016 (SCI-Expanded)
- XXXV. **The Structural, Elastic, Electronic, Thermodynamic and Vibrational Properties of Protactinium Monocarbide (PaC) from First-Principles Calculations**  
Ciftci Y., Coban C., AYDIN S., Tatar A., Evecen M.  
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- XXXVI. **Effect of pressure on structural, electronic, mechanical and optical properties of ruthenium diboride with oP(12)-type structure**  
AYDIN S., Ciftci Y., Mogulkoc Y., Tatar A.  
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- XXXVII. **Unraveling Thermal and Dynamical Properties of the Cubic BaVO<sub>3</sub> Perovskite from First-Principles Calculation**  
Mebrouki M., Ouahrani T., Ciftci Y.  
INTERNATIONAL JOURNAL OF THERMOPHYSICS, vol.37, no.7, 2016 (SCI-Expanded)
- XXXVIII. **Ab-initio calculations of semiconductor MgGeP<sub>2</sub> and MgGeAs<sub>2</sub>**  
Kocak B., Ciftci Y.  
Materials Research Bulletin, vol.77, pp.300-306, 2016 (SCI-Expanded)
- XXXIX. **Determination of the basic physical properties of semiconductor chalcopyrite type MgSnT<sub>2</sub> (T = P, As, Sb) from first-principles calculations**  
Kocak B., ÇİFTÇİ Y.  
JOURNAL OF MATERIALS RESEARCH, vol.31, no.10, pp.1518-1531, 2016 (SCI-Expanded)
- XL. **Electronic structure and thermoelectric properties of half-Heusler compounds with eight electron valence count-KScX (X = C and Ge)**  
ÇİFTÇİ Y., Mahanti S. D.  
JOURNAL OF APPLIED PHYSICS, vol.119, no.14, 2016 (SCI-Expanded)

- XLI. **Some new members of MAX family including light-elements: Nanolayered Hf<sub>2</sub>XY (X= Al, Si, P and Y=B, C, N)**  
 AYDIN S., Tatar A., ÇİFTÇİ Y.  
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- XLII. **Electronic structure and elastic properties of AgZn under pressure from first-principles calculations**  
 ÇİFTÇİ Y.  
 CANADIAN JOURNAL OF PHYSICS, vol.94, no.3, pp.328-333, 2016 (SCI-Expanded)
- XLIII. **Ab Initio Calculations on the Structural, Mechanical, Electronic, Dynamic, and Optical Properties of Semiconductor Half-Heusler Compound ZrPdSn**  
 ÇİFTÇİ Y., ÇOBAN C.  
 ZEITSCHRIFT FÜR NATURFORSCHUNG SECTION A-A JOURNAL OF PHYSICAL SCIENCES, vol.71, no.2, pp.135-143, 2016 (SCI-Expanded)
- XLIV. **Thermo-Elastic and Lattice Dynamical Properties of Pd<sub>3</sub>X (X = Ti, Zr, Hf) Alloys: An Ab Initio Study**  
 Surucu G., Colakoglu K., Ciftci Y., Ozisik H. B., Deligoz E.  
 BRAZILIAN JOURNAL OF PHYSICS, vol.45, no.6, pp.604-614, 2015 (SCI-Expanded)
- XLV. **Lattice dynamical and thermodynamical properties of ReB<sub>2</sub>, RuB<sub>2</sub>, and OsB<sub>2</sub> compounds in the ReB<sub>2</sub> structure**  
 Deligoz E., Colakoglu K., Ciftci Y.  
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- XLVI. **A theoretical study for thorium monocarbide (ThC)**  
 AYDIN S., Tatar A., Ciftci Y.  
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- XLVII. **Lattice dynamical properties of TcB<sub>2</sub> compound**  
 Deligoz E., Colakoglu K., Ozisik H. B., Ciftci Y.  
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- XLVIII. **First principles studies of elastic, electronic and optical properties of chalcopyrite semiconductor ZnSnP<sub>2</sub>**  
 Sahin S., Ciftci Y., Colakoglu K., Korozlu N.  
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- XLIX. **Superhard transition metal tetranitrides: XN<sub>4</sub> (X=Re, Os, W)**  
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- L. **Ab initio calculations on the structural and lattice dynamical properties of TmX (X=As, P) compounds**  
 Coban C., Colakoglu K., Ciftci Y.  
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- LI. **Structural, electronic, elastic, thermodynamic and vibration properties of TbN compound from first principles calculations**  
 Ciftci Y., Ozayman M., Surucu G., Colakoglu K., Deligoz E.  
 SOLID STATE SCIENCES, vol.14, no.3, pp.401-408, 2012 (SCI-Expanded)
- LII. **A first-principle study of the structural, elastic, lattice dynamical and thermodynamic properties of PrX (X=P, As)**  
 Kocak B., Ciftci Y., Colakoglu K., Deligoz E.  
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- LIII. **The structural, elastic and thermodynamic properties of intermetallic compound CeGa<sub>2</sub>**  
 ÇİFTÇİ Y., Colakoglu K., ÇOBAN C., DELİGÖZ E.  
 CENTRAL EUROPEAN JOURNAL OF PHYSICS, vol.10, no.1, pp.197-205, 2012 (SCI-Expanded)
- LIV. **Lattice vibrational properties of Al<sub>2</sub>X (X = Sc, Y) from density functional theory calculations**  
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- LV. **A Molecular Dynamics Study on Au**

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- LVI. **Vibrational properties of Re<sub>2</sub>N and Re<sub>3</sub>N compounds**  
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SOLID STATE COMMUNICATIONS, vol.151, no.17, pp.1122-1127, 2011 (SCI-Expanded)
- LVII. **Structural, elastic, electronic and thermodynamic properties of Nd<sub>2</sub>Te via first principle calculations**  
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- LVIII. **The first principles study on the TmSb compound**  
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- LIX. **Ab initio study of structural, elastic and vibrational properties of praseodymium chalcogenides**  
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- LX. **First-principles studies of CaX (X = In, Tl) intermetallic compounds**  
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- LXI. **The structural, electronic and optical properties of Cd<sub>x</sub>Zn<sub>1-x</sub>Se ternary alloys**  
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- LXII. **The structural, elastic and vibrational properties of the DyX (X = P, As) compounds**  
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- LXIII. **Ab initio study of PrAg intermetallic compound**  
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- LXIV. **Mechanical and phonon properties of the superhard LuB<sub>2</sub>, LuB<sub>4</sub>, and LuB<sub>12</sub> compounds**  
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- LXV. **First principles predictions on mechanical and physical properties of HoX (X = As, P)**  
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- LXVI. **Structural, elastic, and lattice dynamical properties of YB<sub>2</sub> compound**  
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- LXVII. **Structural, elastic, electronic, and thermodynamic properties of PrN from first principles calculations**  
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- LXVIII. **Electronic and mechanical properties of the PdN: A first-principles study**  
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- LXIX. **The structural, electronic, elastic, vibrational, and thermodynamic properties of HoX (X=Sb, Bi)**  
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- LXX. **The electronic and optical properties of Zn<sub>1-x</sub>CaxSe mixed alloys**  
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- LXXI. **Elastic, electronic, and vibrational properties of RhN compound**  
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- JOURNAL OF MATERIALS SCIENCE, vol.45, no.14, pp.3720-3726, 2010 (SCI-Expanded)
- LXXII. **Phonon dispersion and thermodynamical properties in ZrB<sub>2</sub>, NbB<sub>2</sub>, and MoB<sub>2</sub>**  
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- LXXIII. **Lattice dynamical and thermodynamical properties of HfB<sub>2</sub> and TaB<sub>2</sub> compounds**  
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- LXXIV. **The structural, electronic, elastic, phonon, and thermodynamical properties of the SmX (X = P, Sb, Bi) compounds**  
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- LXXV. **Lattice dynamical properties of ScB<sub>2</sub>, TiB<sub>2</sub>, and VB<sub>2</sub> compounds**  
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- LXXVI. **The structural, thermodynamical and elastic properties of TiO**  
ÇİFTÇİ Y., Unlu Y., Colakoglu K., DELİGÖZ E.  
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- XXX. **The effect of Rh, Ir doping on the structural and elastic properties of CoAl: An ab-initio study**  
KARS DURUKAN İ., ÇİFTÇİ Y.

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- XXXI. **A first-principles study of B2 TcSi alloyed with Mn,Re**  
KARS DURUKAN İ., ÇİFTÇİ Y.

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- XXXII. **Structural, elastic, electronic and vibrational properties of CdSr compound : First-principles study pressure**  
KARS DURUKAN İ., ÇİFTÇİ Y.

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- XXXIII. **Ab-INITIO STUDY OF INTERMETALLIC BaZnIN B2 STRUCTURE FOR ELASTIC AND THERMODYNAMIC PROPERTIES UNDER PRESSURE**  
KARS DURUKAN İ., ÇİFTÇİ Y.

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- XXXIV. **EFFECTS OF COMPOSITIONS ON THE PROPERTIES OF Cu DOPED AuCdSHAPE MEMORY ALLOY : FIRST-PRINCIPLES STUDY**  
ÇİFTÇİ Y., Durukan İ. K.

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- XXXV. **BOR INCORPORATION INTO YRh in B2 STRUCTURE: An AB-INITIO STUDY**  
ÇİFTÇİ Y., ALP İ.

1. Uluslararası 23 Nisan Ulusal Egemenlik ve çocuk Kongresi, Ankara, Turkey, 20 - 23 April 2019, pp.170-174

- XXXVI. **THEORETICAL STUDY OF STRUCTURAL, ELASTIC ANISOTROPY, ELECTRONIC, AND VIBRATIONAL PROPERTIES OF CuBe COMPOUND**  
KARS DURUKAN İ., ÇİFTÇİ Y.

1. ULUSLARARASI 23 NİSAN MULTİDİSİPLİNER ÇALIŞMALAR KONGRESİ, 20 - 23 April 2019

- XXXVII. **FIRST-PRINCIPLES STUDY OF STRUCTURAL, ELASTIC, ELECTRONIC, THERMODYNAMIC PROPERTIES OF RuTi COMPOUND in B2 STRUCTURE**  
KARS DURUKAN İ., ÇİFTÇİ Y.

1. ULUSLARARASI 23 NİSAN MULTİDİSİPLİNER ÇALIŞMALAR KONGRESİ, 20 - 23 April 2019

- XXXVIII. **The Influence of Phosphorus Dopant on the Structural and Mechanical Properties of Silicon**  
Ikhmayies S., ÇİFTÇİ Y.

Energy Technologies Symposium, San-Antonio, Northern Mariana Islands, 01 January 2019, pp.201-211

- XXXIX. **The Influence of Boron Dopant on the Structural and Mechanical Properties of Silicon: First Principles Study**  
Ikhmayies S., ÇİFTÇİ Y.

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- XL. **The Pressure Effects on Physical Properties of CeZn Compound in B2 Structure**  
KARS DURUKAN İ., ÇİFTÇİ Y.

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- XLI. **A First-Principles Study of Pressure-induced Effects on Phase Transition, Elastic and Electronic Properties of CeTe Compound**  
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- XLII. **Pressure Effects on Electronic and Elastic Properties of Half-Heusler Semiconductor AgScSi for Optoelectronic Applications**  
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- XLIII. **Elastic and Electronic Properties of Zr Doped TiB ( $Zr_xTi_{1-x}B$ )Alloys By: First-Principles Study**  
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- XLIV. **CeP : Structural, Electronic, Thermodynamic And Vibrational Properties By First Principles Calculations**  
KARS DURUKAN İ., ÇİFTÇİ Y.  
Computational Sciences Congress, 26 - 28 October 2018
- XLV. **Structural, Optic and Vibrational Properties of CeTi Using First-Principles Study**  
KARS DURUKAN İ., ÇİFTÇİ Y.  
Computational Sciences Congress, 26 - 28 October 2018
- XLVI. **A First-Principles Study of Pressure-induced Effects on Phase Transition, Elastic, Electronic and Vibrational Properties of CeSb Compound**  
KARS DURUKAN İ., ÇİFTÇİ Y.  
Computational Sciences Congress, 26 - 28 October 2018
- XLVII. **Hydrogen Storage on Ca-decorated Rectangular Carbon Pentaheptite**  
Alp İ., Çiftci Y.  
II. International Congress on New Trends in Science, Engineering and Technology, Italy, 4 - 06 September 2018, pp.171
- XLVIII. **Structural, Electronic And Elastic Properties of NaYC Compound Under Pressure**  
ÇİFTÇİ Y., ALP İ.  
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- L. **The Physical Properties of Ternary NaYSn Compound: A First-Principles Study**  
ALP İ., ÇİFTÇİ Y.  
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- LI. **Electronic and Optical Properties of NaYPb Compound**  
ALP İ., ÇİFTÇİ Y.  
5th International conference on materials science and nanotechnology for next generation (MSNG2018), Turkey, 04 October 2018
- LII. **Some Physical Properties of Optoelectronic Functional Material NaYSi : First-Principles Study**  
ÇİFTÇİ Y., ALP İ.  
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- LIII. **Hydrogen Adsorption Ability of Mg-decorated Graphene Allotrope**  
Alp İ., Çiftci Y.  
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- LIV. **LiPb : Structural, Elastic, Electronic, Thermodynamic And Vibrational Properties By First Principles Calculations**  
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- LV. **The Thermodynamical Properties of AsPr via First Principles**  
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- LVI. **High Pressure Physical Properties of AsPu Compound**  
ALP İ., ÇİFTÇİ Y.
6. European Conference on Renewable Energy Systems (ECRES 2018), İstanbul, Turkey, 25 June 2018
- LVII. **Exploration of Thermoelectricity in NaScGe Half Heusler Compound: A First Principle Study**  
ALP İ., ÇİFTÇİ Y.
6. European Conference on Renewable Energy Systems (ECRES 2018), İstanbul, Turkey, 25 June 2018
- LVIII. **A Theoretical Study of Pressure-Induced Effects on Phase Transition and Elastic Properties of AsTh Compound**  
ÇİFTÇİ Y., ALP İ.
6. European Conference on Renewable Energy Systems (ECRES 2018), İstanbul, Turkey, 25 June 2018
- LIX. **Hydrogen Storage on Li-decorated YCrB4-type Graphene Allotrope**  
ALP İ., ÇİFTÇİ Y.
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- LX. **The Pressure Effects on Physical Properties of AlNi Compound in B2 Structure**  
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ÇİFTÇİ Y., EVECEN M., ALP İ.  
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ALP İ., ÇİFTÇİ Y.  
Turkish Physical Society 33rd International Physics Congress, 6 - 10 September 2017, pp.655
- LXVI. **Pressure effects on electronic, elastic and vibration properties of metallic antiprevskite PbNCa3 by ab initio calculations**  
ÇİFTÇİ Y., EVECEN M., ALP İ.  
Turkish Physical Soociety 33rd International Physics Congress, 6 - 10 September 2017
- LXVII. **ELECTRONIC AND MECHANICAL PROPERTIES OF HALF HEUSLER SEMICONDUCTOR LiScSn VIA FIRST-PRINCIPLES CALCULATIONS**  
MOĞULKOÇ Y., ÇİFTÇİ Y.  
TURKISH PHYSICAL SOCIETY 33rd INTERNATIONAL PHYSICS CONGRESS, 6 - 10 September 2017
- LXVIII. **ELECTRONIC AND OPTICAL PROPERTIES OF SINGLE-LAYER STRAINED SNC**  
MOĞULKOÇ Y., MODARRESİ M., MOĞULKOÇ A., ÇİFTÇİ Y., ALKAN Ş. B.  
TURKISH PHYSICAL SOCIETY 33rd INTERNATIONAL PHYSICS CONGRESS, Muğla, Turkey, 6 - 10 September 2017
- LXIX. **The pressure effect on physical properties of CuTiS2**  
ALP İ., ÇİFTÇİ Y.  
Turkish Physical Society 33.rd International Physics Congress, 6 - 10 September 2017
- LXX. **RhAl: Structural, elastic, electronic, thermodynamic and vibrational properties by first principles calculations**

ÖZCAN A., ÇİFTÇİ Y., ALP İ.

Turkish Physical Society 33rd International Physical Congress, 6 - 10 September 2017, pp.636

- LXXI. **An ab-initio study of structural, elastic, electronic and optical properties for chalcopyrite semiconductor CaSiN<sub>2</sub> compound**  
ÇİFTÇİ Y., ALP İ., MOĞULKOÇ Y.  
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- LXXII. **HIGH-PRESSURE EFFECT ON ZnSiN<sub>2</sub>: A FIRST-PRINCIPLES STUDY**  
ALP İ., ÇİFTÇİ Y.  
5. European Conference on Renewable Energy Systems, Saraybosna, Bosnia And Herzegovina, 27 - 30 August 2017, pp.75
- LXXIII. **THE CORROSION BEHAVIOUR OF ALUMINIUM MATRIXCOMPOSITES IN AQUEOUS 3.5 M NaCl SOLUTION**  
AYTAÇ A., KEZBAN S., ÇİFTÇİ Y.  
1th INTERNATIONAL CONFERENCE ON ENERGY AND THERMAL ENGINEERING, İstanbul, Turkey, 25 - 28 April 2017, pp.920
- LXXIV. **Ag Dekore Edilmiş ThMoB<sub>4</sub> Tipi Grafen Allotropunun Hidrojen Depolama Özelliklerinin İlk İlkeler Yöntemi İle İncelenmesi**  
ALP İ., ÇİFTÇİ Y., AYDIN S.  
22. Yoğun Madde Fiziği Ankara Toplantısı, Ankara, Turkey, 16 December 2016
- LXXV. **A FIRST PRINCIPLES STUDIES ON STRUCTURAL AND ELASTIC PROPERTIES OF Sr<sub>2</sub>Ge COMPOUND**  
ALP İ., ÇİFTÇİ Y.  
1st International Mediterranean Science and Engineering Congress (IMSEC 2016), Adana, Turkey, 26 - 28 October 2016
- LXXVI. **THE PRESSURE DEPENDENCE OF THE SIGNIFICANT PHYSICAL PROPERTIES OF SEMICONDUCTOR Ba<sub>2</sub>Si**  
ALP İ., ÇİFTÇİ Y.  
1st International Mediterranean Science and Engineering Congress (IMSEC 2016), Adana, Turkey, 26 - 28 October 2016
- LXXVII. **A THEORETICAL STUDY ON THE STRUCTURAL ELASTIC AND VIBRATIONAL PROPERTIES OF Sr<sub>2</sub>Si COMPOUND**  
ÇİFTÇİ Y., ALP İ.  
1st International Mediterranean Science and Engineering Congress (IMSEC 2016), Adana, Turkey, 26 - 28 October 2016
- LXXVIII. **AB INITIO CALCULATIONS FOR THE OPTICAL AND PHONON PROPERTIES OF ANTI FLUORITE SEMICONDUCTOR Sr<sub>2</sub>Sn**  
ÇİFTÇİ Y., ALP İ.  
1st International Mediterranean Science and Engineering Congress (IMSEC 2016), Adana, Turkey, 26 - 28 October 2016
- LXXIX. **A theoretical Study of Structural Elastic Vibration and Optic Properties for Half Heusler Semiconductor NaAlC Compound Using ab initio Method**  
ÇİFTÇİ Y.  
2nd International Congress on The World of Technology and Advanced Materials, Kırşehir, Turkey, 28 September - 02 October 2016
- LXXX. **Theoretical Investigations of Half Heusler Semiconductor NaMgP Compound with MgAgAs Structure**  
Alp İ., Çiftci Y.  
2nd International Congress on The World of Technology and Advanced Materials, Kırşehir, Turkey, 28 September - 02 October 2016
- LXXXI. **An ab initio Study of Structural Elastic Vibration and Optic Properties for Semiconductor NaAlGe Compound**  
ÇİFTÇİ Y., ALP İ.  
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02 October 2016

- LXXXII. **Structural Elastic Electronic Vibrational of NaMgN Compound with MgAgAs Structure Under Pressure**  
ALP İ., ÇİFTÇİ Y.  
2nd International Congress on The World of Technology and Advanced Materials, Kırşehir, Turkey, 28 September - 02 October 2016
- LXXXIII. **A DFT Study on Structural Phase Transition and Mechanical Properties of LiAl**  
MOĞULKOÇ Y., ÇİFTÇİ Y., SÜRÜCÜ G.  
2nd International Congress on The World of Technology and Advanced Material, Kırşehir, Turkey, 28 September - 02 October 2016
- LXXXIV. **First Principles Study on Electronic and Thermoelectric Properties for Half Heusler Semiconductor NaAlC Compound**  
ÇİFTÇİ Y.  
2nd International Congress on The World of Technology and Advanced Materials, 28 September - 02 October 2016
- LXXXV. **First Principles Study on Half Heusler NaMgAs Compound**  
ALP İ., ÇİFTÇİ Y.  
2nd International Congress on The World of Technology and Advanced Materials, Kırşehir, Turkey, 28 September - 02 October 2016
- LXXXVI. **A DFT Study on Structural Phase Transition and Mechanical Properties of LiAl**  
MOĞULKOÇ Y., ÇİFTÇİ Y., SÜRÜCÜ G.  
2nd International Congress On The World Of Technology And Advanced Materials, 28 September - 02 October 2016
- LXXXVII. **Study of Structural Elastic Electronic Bonding Behaviour and Vibrational Properties in ScCd Compound By First Principles**  
Çiftci Y., Alp İ.  
Turkish Physical Society 32nd International Physics Congress, Muğla, Turkey, 6 - 09 September 2016
- LXXXVIII. **FIRST PRINCIPLES INVESTGATIONS OF STRUCTURAL ELASTIC ELECTRONIC THERMODYNAMIC AND VIBRATIONAL PROPERTIES OF LACD COMPOUND**  
ÇİFTÇİ Y., ALP İ., MOĞULKOÇ Y., ÖZCAN A.  
TURKISH PHYSICAL SOCIETY 32. INTERNATIONAL PHYSICS CONGRESS, 6 - 09 September 2016
- LXXXIX. **Electronic and optical properties of stack depended blue phosphorus**  
MOĞULKOÇ Y., Modarresi M., Mogulkoç A., ÇİFTÇİ Y.  
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- XC. **Pressure dependence of physical properties for YbHg compound**  
Alp İ., ÇİFTÇİ Y.  
Turkish Physical Society 32 nd Internatioanl Physics congress, 6 - 09 September 2016
- XCI. **The first principles investigation of LuXO<sub>3</sub> X Al Ga In cubic Preovskite Oxides**  
Sürücü G., ÇİFTÇİ Y., SÜRÜCÜ G.  
Turkish Physical Society 32 nd International Physics Congress, 6 - 09 September 2016
- XCII. **Ab initio Calculations of YbCd Intermetallic Compounds in B2 Structure**  
ÇİFTÇİ Y., Alp İ., ÖZCAN A.  
Turkish Physical Society 32 nd International Physical Congress, Turkey, 6 - 09 September 2016
- XCIII. **PRESSURE DEPENDENCE OF THE PHYSICAL PROPERTIES FOR YbHg COMPOUND**  
Alp İ., Çiftci Y.  
Turkish Physical Society 32nd International Physics Congress, Muğla, Turkey, 6 - 09 September 2016
- XCIV. **THE FIRST PRINCIPLES CALCULATIONS FOR THE STRUCTURAL ELASTIC ELECTRONIC AND VIBRATIONAL PROPERTIES OF ScHg**  
ALP İ., ÇİFTÇİ Y.  
Turkish Physical Society 32nd International Physics Congress, Muğla, Turkey, 6 - 09 September 2016
- XCV. **The First Principles Calculations for the Structural Elastic Electronic and Vibrational Properties of**



## ScHg

Alp İ., Çiftci Y.

Turkish Physical Society 32 nd International Physics Congress, Muğla, Turkey, 6 - 09 September 2016, pp.334

- XCVI. **AB INITIO CALCULATIONS OF YBCD INTERMETALLIC COMPOUND IN B2 STRUCTURE**  
ÇİFTÇİ Y., ALP İ., ÖZCAN A.  
TURKISH PHYSICAL SOCIETY 32. INTERNATIONAL PHYSICS CONGRESS, 6 - 09 September 2016
- XCVII. **First Principles Investigations of Structural Elastic Electronic Thermodynamic and Vibrational Properties of LaCd Compound**  
Çiftci Y., Alp İ., Moğulkoç Y., Özcan A.  
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- XCVIII. **The First Principles Investigation Of LuXO3 Al Ga and In Cubic Perovskite Oxides**  
sürücü G., ÇİFTÇİ Y., SÜRÜCÜ G.  
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- XCIX. **Hydrogen adsorption on lithium and beryllium decorated planar ThMoB4 type graphene allotrope**  
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- CI. **Hydrogen storage on nano level optimized pentaheptite graphene allotrope with sodium decoration**  
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- CII. **Hydrogen adsorption on Ca decorated hexagonal planar carbon pentaheptite**  
ALP İ., ÇİFTÇİ Y., AYDIN S.  
12th Nanoscience and Nanotechnology Conference, Kocaeli, Turkey, 3 - 05 June 2016
- CIII. **Theoretical investigation of the electronic structure elastic dynamic properties of intermetallic compound NiBe under pressure**  
EVECEN M., ÇİFTÇİ Y., ALDIRMAZ E.  
2 nd International Conference on Organic Electronic Material Technologies (OEMT2016), 17 - 19 May 2016
- CIV. **Ab initio calculations of semiconductor LiScGe compound**  
Çiftci Y., Alp İ.  
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Alp İ., Çiftci Y.  
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- CVI. **Structural mechanical electronic dynamic and optical properties of semiconductor half Heusler compound LiScSi under pressure from first principle study**  
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- CVII. **First Principles Study on UB6 for Nuclear Applications**  
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- CIX. **THE ELECTRONIC STRUCTURE ELASTIC and VIBRATIONAL PROPERTIES of InPd UNDER PRESSURE**

via **FIRST PRINCIPLES CALCULATIONS**

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- CX. **Analysis of the Structural, Electronic and Optic Properties of Ni Doped MgSiP<sub>2</sub> Semiconductor Chalcopyrite Compound**  
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- CXV. **B<sub>2</sub> Fazdaki Şekil Hafızalı Alaşım TaRu'nun Basınca Bağlı Fiziksel Özelliklerinin incelenmesi**  
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ÇİFTÇİ Y., ALP İ.  
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- CXIX. **V<sub>2</sub>PC Bileşiğinin Yapısal, Mekanik, Elektronik, Titreşimsel ve Termodinamik Özelliklerinin Ab-initio Yöntemle İncelenmesi**  
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- II. ÇİFTÇİ Y., Project Supported by Higher Education Institutions, REc2 (RE= La, Ce, Pr, Pa, U) nadir toprak karbür bileşiklerinin temel fiziksel özelliklerinin ab-initio yöntemlerle incelenmesi, 2012 - 2015
- III. ÇİFTÇİ Y., Project Supported by Higher Education Institutions, Tb pniktidlerinin yapısal, mekaniksel, elektronik ve titreşimsel özelliklerinin ab-initio yöntemlerle incelenmesi, 2010 - 2012
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## Metrics

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Publication: 280

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H-Index (WoS): 19

H-Index (Scopus): 19

## Congress and Symposium Activities

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- I. Nanotechnology Research and Applications, Attendee, Praha, Czech Republic, 2021
- II. International Conference on Advanced Materials Science and Engineering and High Tech Device Applications (ICMATSE 20), Invited Speaker, Ankara, Turkey, 2020

## Coaching Duties

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## Refereeing Duties

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