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Eğitim Bilgileri

- I. Doktora, Gazi Üniversitesi, Fen Bilimleri Enstitüsü, Fizik (Dr), Türkiye 1996 - 2000
- II. Yüksek Lisans, Gazi Üniversitesi, Fen Bilimleri Enstitüsü, Fizik (Yl) (Tezli), Türkiye 1993 - 1996
- III. Lisans, Gazi Üniversitesi, Fen-Edebiyat Fakültesi, Fizik Pr., Türkiye 1989 - 1993

Yabancı Diller

- I. İngilizce, B2 Orta Üstü

Araştırma Alanları

Fizik, Temel Bilimler, Mühendislik ve Teknoloji

Akademik Unvanlar / Görevler

- I. Prof. Dr., Gazi Üniversitesi, Fen Fakültesi, Fizik, 2014 - Devam Ediyor
- II. Yrd. Doç. Dr., Gazi Üniversitesi, Fen Fakültesi, Fizik, 2003 - 2008
- III. Araştırma Görevlisi, Gazi Üniversitesi, Fen Fakültesi, Fizik, 1994 - 2003

Yönetilen Tezler

- I. ÇİFTÇİ Y., Bazı hafif elementlerle dekore edilmiş iki boyutlu karbon allotroplarının üzerine hidrojen depolama özelliklerinin ab-initio yöntemlerle incelenmesi, Doktora, İALP(Öğrenci), 2019
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- I. Akademik Kadroya Atama-Yardımcı Doçentlik, Akademik Kadroya Atama-Yardımcı Doçentlik, University of Engineering and Technology, Eylül, 2022

SCI, SSCI ve AHCI İndekslerine Giren Dergilerde Yayınlanan Makaleler

- I. **Computational analysis of the physical properties of AlNi intermetallic compound: pressure effect**
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- VII. **A first-principles prediction on the structural, electronic, elastic, phonon, and transport properties of BaSiN₂**
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- VIII. **Pressure effects on structural, electronic and anisotropic elastic properties of Si doped RuGe compound with different concentrations by first-principles calculations**
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- IX. **Systematic study of optoelectronic and thermoelectric properties of new lead-free halide double perovskites A(2)KGaI(6)(A = Cs, Rb) for solar cell applications via ab-initio calculations**
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- X. **Equiatomic Quaternary CoXCrAl (X = V, Nb, and Ta) Heusler Compounds: Insights from DFT Calculations**

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- XI. **First-principles study on B2 based XAl(X = Rh, Ru)compounds**
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- XII. **First-principles calculations of vibrational and optical properties of half-Heusler NaScSi**
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- XIII. **Ab-initio study on physical properties of intermetallic LiPb compound**
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- XIV. **Band Alignment in Monolayer Boron Phosphide with Janus MoSSe Heterobilayers under Strain and Electric Field**
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- XV. **Mechanical and dynamic properties of stable two-dimensional boron-substituted ThMoB4-type graphene: First-Principles Study**
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- XVI. **Ligand-free fabrication of Au/TiO₂ nanostructures for plasmonic hot-electron-driven photocatalysis: Photoelectrochemical water splitting and organic-dye degradation**
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- XIX. **The pressure effect on optoelectronic and mechanical properties of chalcopyrite BeSiN₂**
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- XX. **A Theoretical Study of Pressure-Induced Effects on Phase Transition and Elastic Properties of AsTh Compound**
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- XXI. **Anisotropic Elastic, Electronic and Vibrational Properties of the Semiconductor AgScX (X = Ge, C) Compounds**
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- XXII. **Green-Emitting Lead-Free Cs₄SnBr₆ Zero-Dimensional Perovskite Nanocrystals with Improved Air Stability**
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- XXV. **The Effect of Pressure on Elastic Anisotropy, Vibration and Optical Properties of a AgScSi Compound**
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- XXVII. **Analysis of the structural, electronic, elastic and thermodynamic properties of CuAl₂X₄ (X = O, S) spinel structure**
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- XXXII. **A density functional study of chalcopyrite MgGeSb₂**
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- XL.** **The Structural, Elastic, Electronic, Thermodynamic and Vibrational Properties of Protactinium Monocarbide (PaC) from First-Principles Calculations**
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- LV. **First principles LDA+U and GGA+U study of HfO₂: Dependence on the effective U parameter**
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- LVI. **Ab initio study of the structural, elastic, thermodynamic, electronic and vibration properties of TbMg intermetallic compound**
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- LXI. **A theoretical study for thorium monocarbide (ThC)**
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- LXIII. **Lattice dynamical properties of TcB₂ compound**
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- LXIV. **Ab initio calculations on the structural and lattice dynamical properties of TmX (X=As, P) compounds**
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- LXV. **First principles studies of elastic, electronic and optical properties of chalcopyrite semiconductor ZnSnP₂**
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- LXX. **A first-principle study of the structural, elastic, lattice dynamical and thermodynamic properties of PrX (X=P, As)**
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Metrikler

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Hakemlik Görevleri
