

Prof. Dr. HASAN HAMİT YURTSEVEN

Kişisel Bilgiler

İş Telefonu: [+90 210 505 6](tel:+902105056)

E-posta: hamit@metu.edu.tr

Web: <https://avesis.metu.edu.tr/hamit>

Uluslararası Araştırmacı ID'leri

ScholarID: [CvmKz5IAAAAJ](https://scholar.google.com/citations?user=CvmKz5IAAAAJ)

ORCID: [0000-0002-7745-6490](https://orcid.org/0000-0002-7745-6490)

Publons / Web Of Science ResearcherID: [ABA-9994-2020](https://publons.com/author/99942020/ABA-9994-2020)

ScopusID: [9334413200](https://scopus.com/authid/detail.uri?authorId=9334413200)

Yoksis Araştırmacı ID: [11797](https://yoksis.metu.edu.tr/11797)

Eğitim Bilgileri

Doktora, University of London-Kings College London, Birleşik Krallık 1980 - 1984

Lisans, Hacettepe Üniversitesi, Mühendislik Fakültesi, Türkiye 1972 - 1977

Yaptığı Tezler

Doktora, Analysis of the Critical Phenomena using Raman data, University Of London-Kings College London, 1984

Araştırma Alanları

Temel Bilimler

Akademik Unvanlar / Görevler

Prof. Dr., Orta Doğu Teknik Üniversitesi, Fen Edebiyat Fakültesi, Fizik Bölümü, 2003 - Devam Ediyor

Prof. Dr., İstanbul Teknik Üniversitesi, Fen-Edebiyat Fakültesi, 1994 - 2003

Doç. Dr., İstanbul Teknik Üniversitesi, Fen-Edebiyat Fakültesi, 1988 - 1994

Yrd. Doç. Dr., İstanbul Teknik Üniversitesi, Fen-Edebiyat Fakültesi, 1987 - 1988

Araştırma Görevlisi, Ankara Üniversitesi, Fen Fakültesi, 1985 - 1987

Verdiği Dersler

Thermodynamics and Statistical Mechanics, Lisans, 2020 - 2021

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- V. **Calculation of the phase diagram of n-alkanes (C_nH_{2n+2}) by the Landau mean field theory**
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- VII. **Magnetic transitions in two novel mixed-valence iron(II)-iron(III) metal formate frameworks: Two**

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- VIII. **Calculation of the magnetization and the heat capacity for $[\text{NH}_2 \text{NH}_3][\text{M}(\text{HCOO})_3]$ ($\text{M}=\text{Fe}, \text{Mn}$) near the phase transitions**
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- IX. **Calculation of the spin-lattice relaxation time and the activation energy near the IV-III phase transition in pyridinium fluorosulfonate (C_5NH_6)FSO₃**
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- X. **Magnetic, thermal and ferroelectric properties of MOFs (MHyM , $\text{M} = \text{Fe}, \text{Mn}$) close to phase transitions**
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- XII. **Landau phenomenological model for the alpha-INC (incommensurate)-beta transition in quartz**
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- XIII. **Investigation of vibrational, elastic and dielectric properties of cubic gauche nitrogen (cg-N)**
Akay Ö., YURTSEVEN H. H.
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- XIV. **Calculation of the damping constant and the relaxation time of the LA mode in the incommensurate phase of quartz**
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- XV. **Phenomenological approaches on the Nd³⁺ doped ferroelectric LaBGeO₅**
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- XXIII. **Raman bandwidths calculated for the librational (α -phase) and internal (δ , δ_{loc} and δ phases) modes in solid N_2 using pseudospin-phonon coupling (PS) and energy-fluctuation (EF) models**
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- XXVI. **Calculation of the thermodynamic functions using a mean field model for the fluid-solid transition in nitrogen**
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- XXVII. **Liquid-(Solid plus Liquid) Transitions in a Two-Component System of $(\text{CH}_3)\text{CCl}_3 + \text{CCl}_4$**
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- XXXII. **Analysis and calculation of the magnetization, magnetic susceptibility and the specific heat close to phase transitions in heterometallics**
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- XXXIII. **Damping constant and the inverse relaxation time calculated as a function of pressure using the X-ray diffraction data close to the cubic-tetragonal phase transition in SrTiO_3**
YURTSEVEN H. H., Kiraci A.
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- XXXV. **Calculation of the thermodynamic quantities of perovskite metal organics DMAKCr and perovskite HyFe close to the weakly first-order relaxor-like structural transformation using the mean field theory**
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- XXXIX. **CALCULATION OF THE PHASE DIAGRAMS FOR THE FLUID-SOLID AND SOLID-SOLID ($\delta\delta(\text{loc})$ - ϵ) TRANSITIONS IN MOLECULAR NITROGEN BY USING MEAN FIELD MODEL**
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- XL. **Calculation of the heat capacity Cp from the temperature-induced and pressure-induced Raman frequency shifts for solid benzene, naphthalene and anthracene**
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- XLI. **Calculation of the T-P Phase Diagrams for the Halogenomethane Compounds (CCl₄-nBrn, n=0, 1, 2, 4) Using the Mean Field Theory**
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- XLII. **Temperature and the field dependence of the magnetization close to order-disorder phase transitions in DMMn and the chromium-doped DMMn**
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- XLIV. **Thermodynamic Parameters of Cholesteric/Smectic A Transition in Cholesteric Myristate and Its Binary Mixture CM/PCPB**
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- XLV. **Correlations between the resonant frequency shifts and the thermodynamic quantities for the alpha-beta transition in quartz**
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- XLVI. **T-P Phase Diagram of Nitrogen at High Pressures**
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- XLVII. **Analysis of the Specific Heat of Ru Doped LiKS₀₄ Close to Phase Transitions**
YURTSEVEN H. H., Tirpanci D. V., KARACALI H.
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- XLVIII. **Pressure-dependent Raman modes near the cubic-tetragonal transition in strontium titanate**

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- XLIX. **Calculation of the Raman frequency and the damping constant (linewidth) of the stretching modes for the metal-organic compound DMMg close to the paraelectric-ferroelectric transitions**
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- L. **Calculation of the C-P-C-V as a function of temperature close to the melting point in benzene**
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- LI. **Calculation of the Raman and IR frequencies from the volume data at high pressures in N-2**
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- LIII. **ANALYSIS OF THE INTEGRATED INTENSITY OF THE CENTRAL PEAKS CALCULATED AS A FUNCTION OF TEMPERATURE IN THE FERROELECTRIC PHASE OF LITHIUM TANTALATE**
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- LIV. **Calculation of the P-T phase diagram of nitrogen using a mean field model**
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- LVII. **Pressure dependence of the heat capacity near the melting point in benzene**
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- LVIII. **Calculation of the tilt angle and susceptibility for the alpha-beta transition in quartz using a mean field model**
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- LIX. **Analysis of the susceptibility of condensed oxygen under high pressures and in strong magnetic fields**
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- LX. **Analysis of the Magnetic Susceptibility as a Function of Temperature for the alpha, beta and gamma Phases of Oxygen**
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- LXI. **Temperature dependence of the damping constant and the relaxation time close to the tetragonal-cubic phase transition in SrZrO₃**
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- LXII. **Raman frequencies calculated as functions of temperature and pressure using volume data for solid phase I of benzene**

- YURTSEVEN H. H., Ozdemir H.
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- LXIII. **Calculation of the Raman frequency shifts for the alpha phase of solid oxygen**
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- LXIV. **Analysis of vibrational spectra in relation to the thermodynamic quantities close to the phase transitions in ammonium halides (NH₄F and NH₄I)**
Tari O., YURTSEVEN H. H.
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- LXV. **Pressure dependence of the Raman frequency calculated from the volume data close to the ferroelectric-paraelectric transition in PbTiO₃**
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- LXVI. **Temperature dependence of the polarization, dielectric constant, damping constant and the relaxation time close to the ferroelectric-paraelectric phase transition in LiNbO₃**
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- LXIX. **Variation of the molar volume with the temperature and the Pippard relations near the melting point in benzene**
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- LXX. **Temperature and pressure dependence of the Raman frequency shifts in anthracene**
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- LXXI. **Analysis of the Heat Capacity for Pure CH₄ and CH₄/CCl₄ on Graphite Near the Melting Point and Calculation of the T-X Phase Diagram for(CH₃)CCl₃+CCl₄**
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- LXXII. **Analysis of the orientational order parameter and the magnetic field as functions of temperature for p-anisaldazine**
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- LXXVI. **Analysis of the dielectric constant, spontaneous polarization and the coersive field close to the phase transitions in ferroelectric K_2CoCl_4**
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- LXXVII. **Calculation of the soft-mode frequency for the alpha - beta transition in quartz**
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