### **Prof. HASAN HAMİT YURTSEVEN**

### **Personal Information**

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### **Education Information**

Doctorate, University of London-Kings College London, United Kingdom 1980 - 1984 Undergraduate, Hacettepe University, Mühendislik Fakültesi, Turkey 1972 - 1977

### Dissertations

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

#### **Research Areas**

**Natural Sciences** 

### Academic Titles / Tasks

Professor, Middle East Technical University, Faculty of Arts and Sciences, Department of Physics, 2003 - Continues Professor, Istanbul Technical University, Fen-Edebiyat Fakültesi, 1994 - 2003 Associate Professor, Istanbul Technical University, Fen-Edebiyat Fakültesi, 1988 - 1994 Assistant Professor, Istanbul Technical University, Fen-Edebiyat Fakültesi, 1987 - 1988 Research Assistant, Ankara University, Fen Fakültesi, 1985 - 1987

#### Courses

Thermodynamics and Statistical Mechanics, Undergraduate, 2020 - 2021

### **Advising Theses**

YURTSEVEN H. H., Thermodynamic properties of the high-pressure phases in solid nitrogen close to phase transitions, Doctorate, Ö.Akay(Student), 2020

YURTSEVEN H. H., Critical behaviour of the thermodynamic quantities close to phase transitions in molecular crystals using raman data, Doctorate, H.ÖZDEMİR(Student), 2017

YURTSEVEN H. H., First order and second order phase transitions in quartz and related materials, Doctorate, M.CEM(Student), 2017

YURTSEVEN H. H., Investigation of the dynamic properties of ferroelectric crystals close to phase transitions, Doctorate, A.KİRACI(Student), 2015

YURTSEVEN H. H., Calculation of the thermodynamic and spectroscopic quantities in molecular crystals close to the phase transitons, Doctorate, D.KAVRUK(Student), 2011

YURTSEVEN H. H., Critical behaviour of the thermodynamic quantities for the thermotropic and ferroelectric liquid crystals close to the phase transitions, Doctorate, E.KİLİT(Student), 2011

YURTSEVEN H. H., Calculation of the Raman frequencies using volume data in various phases of solid nitrogen and benzene, Postgraduate, E.Çetinbaş(Student), 2011

YURTSEVEN H. H., Alpha - beta transition in quartz, Postgraduate, M.Cem(Student), 2011

YURTSEVEN H. H., Calculation of phase diagrams and the thermodynamic quantities from the mean field models close to phase transitions in molecular and liquid crystals, Doctorate, S.ŞEN(Student), 2009

YURTSEVEN H. H., Correlations between the spectroscopic parameters and the thermodynamic quantities for systems exhibiting phase transitions, Doctorate, H.Karaçalı(Student), 2005

YURTSEVEN H. H., Correlations between the spectroscopic parameters and the thermodynamic quantites for systems exhibiting phase transitions, Doctorate, H.KARAÇALI(Student), 2005

YURTSEVEN H. H., The Analysis of various systems by using landau mean field theory, Doctorate, Y.ENGINER(Student), 2001

YURTSEVEN H. H., Amonyum halojenler ve sıvı kristallerde faz geçişlerinin incelenmesi, Doctorate, A.TÜBLEK(Student), 2001

### Published journal articles indexed by SCI, SSCI, and AHCI

 I. Tricritical behavior of the smectic-hexatic phase transitions in binary mixtures using the landau mean field theory YURTSEVEN H. H., Kilit Doğan E.

PHASE TRANSITIONS, vol.96, no.6, pp.383-399, 2023 (SCI-Expanded)

II. Calculation of the T - X phase diagram and the thermodynamic quantities for the binary mixtures of Tetradecane + Hexadecane using the Landau mean field model YURTSEVEN H. H., Tari O.

Physics and Chemistry of Liquids, vol.61, no.5, pp.340-364, 2023 (SCI-Expanded)

- III. Investigation of the complex magnetic behavior of Ni46.86Co2.91Mn38.17Sn12.06(at%) magnetic shape memory alloy at low temperatures
  Ylldlrlm O., Yuce S., Bruno N. M., Doğan E. K., YURTSEVEN H. H., DUMAN E., EMRE Ş. B.
  Physica Scripta, vol.97, no.8, 2022 (SCI-Expanded)
- IV. Calculation of the T X phase diagram of tetradecane + hexadecane and tetradecane + pentadecane under high pressure by the landau mean field theory
  Tari O., YURTSEVEN H. H.
  Fluid Phase Equilibria, vol.559, 2022 (SCI-Expanded)

V. Calculation of the phase diagram of n-alkanes (CnH2n+2) by the Landau mean field theory YURTSEVEN H. H., Kilit Dogan E.

Fluid Phase Equilibria, vol.556, 2022 (SCI-Expanded)

 VI. Calculation of the T-X phase diagram for the first-order smectic-hexatic transitions in binary mixtures
 Kilit Dogan E., YURTSEVEN H. H.

Physics of Fluids, vol.34, no.4, 2022 (SCI-Expanded)

VII. Magnetic transitions in two novel mixed-valence iron(II)-iron(III) metal formate frameworks: Two

sublattice model YURTSEVEN H. H., Tari O. Journal of Magnetism and Magnetic Materials, vol.546, 2022 (SCI-Expanded) VIII. Calculation of the magnetization and the heat capacity for [NH2 NH3][M(HCOO)3] (M=Fe, Mn) near the phase transitions YURTSEVEN H. H., Doğan E. Ferroelectrics, vol.600, no.1, pp.73-87, 2022 (SCI-Expanded) IX. Calculation of the spin-lattice relaxation time and the activation energy near the IV-III phase transition in pyridinium fluorosulfonate (C5NH6)FSO3 Kara N., Kiraci A., YURTSEVEN H. H. Ferroelectrics, vol.589, no.1, pp.45-54, 2022 (SCI-Expanded) X. Magnetic, thermal and ferroelectric properties of MOFs (MHyM, M = Fe, Mn) close to phase transitions YURTSEVEN H. H., Kilit Dogan E. Journal of Magnetism and Magnetic Materials, vol.540, 2021 (SCI-Expanded) XI. Calculation of the Liquid-Solid Phase Diagram and the Thermodynamic Quantities of the Binary System of Tetradecane and Hexadecane Using the Mean Field Theory Yurtseven H., Emirosmanoglu T., Tari O. JOURNAL OF SOLUTION CHEMISTRY, vol.50, no.11-12, pp.1335-1362, 2021 (SCI-Expanded) XII. Landau phenomenological model for the alpha-INC (incommensurate)-beta transition in quartz Yurtseven H. H., Ates S. PHILOSOPHICAL MAGAZINE, vol.101, no.21, pp.2331-2353, 2021 (SCI-Expanded) XIII. Investigation of vibrational, elastic and dielectric properties of cubic gauche nitrogen (cg-N) Akay Ö., YURTSEVEN H. H. Optik, vol.236, 2021 (SCI-Expanded) XIV. Calculation of the damping constant and the relaxation time of the LA mode in the incommensurate phase of quartz Ates S., Yurtseven H. FERROELECTRICS, vol.573, no.1, pp.9-22, 2021 (SCI-Expanded) XV. Phenomenological approaches on the Nd3+ doped ferroelectric LaBGe05 Kara N., Kiraci A., YURTSEVEN H. H. FERROELECTRICS, vol.572, no.1, pp.13-26, 2021 (SCI-Expanded) XVI. Calculation of the inverse relaxation time and the activation energy as a function of temperature for the Raman modes close to the phase transitions in solid nitrogen Yurtseven H., Akay O. JOURNAL OF MOLECULAR STRUCTURE, vol.1226, 2021 (SCI-Expanded) XVII. Calculation of Magnetization and Magnetic Susceptibility Close to Magnetic Phase Transitions in (CH3)2NH2FeIIINiII(HCOO)6 and (CH3)2NH2FeIIICuII(HCOO)6 YURTSEVEN H. H., Tari O. Journal of Low Temperature Physics, vol.202, pp.145-159, 2021 (SCI-Expanded) XVIII. Analysis of the IR-Raman Modes and the Heat Capacity Near the  $\alpha$ -Inc- $\beta$  Transitions in Quartz YURTSEVEN H. H., Günay E., Karacali H., Ateş S. Ferroelectrics, vol.577, no.1, pp.125-142, 2021 (SCI-Expanded) Temperature dependence of the piezoelectric resonance frequency in relation to the anomalous XIX. strain near the incommensurate phase of quartz Ates S., Yurtseven H. FERROELECTRICS, vol.571, no.1, pp.39-49, 2021 (SCI-Expanded) XX. Calculation of the thermodynamic functions from the Raman frequency shifts close to the epsilon delta(loc) - delta transitions and Pippard relations in nitrogen Akay O., Yurtseven H. MODERN PHYSICS LETTERS B, vol.34, no.33, 2020 (SCI-Expanded)

XXI.	Thermodynamic study on the magnetic transition and structural phase transition in [(CH3)(2)NH2]
	[Na0.5Fe0.5(HCOO)(3)] by using the Landau phenomenological model
	Yurtseven H., Tari O.
	JOURNAL OF APPLIED PHYSICS, vol.128, no.20, 2020 (SCI-Expanded)
XXII.	Magnetization Studied as a Function of Temperature and Magnetic Field for Ferromagnetic
	Transition in DMNaFe
	Kilit Dogan E., Yurtseven H.
	JOURNAL OF ELECTRONIC MATERIALS, vol.49, no.11, pp.6388-6393, 2020 (SCI-Expanded)
XXIII.	Raman bandwidths calculated for the librational ( a -phase) and internal (8, d loc and d phases)
	modes in solid N 2 using pseudospin-phonon coupling (PS) and energy-fluctuation (EF) models
	Yurtseven H., Akay O.
	JOURNAL OF MOLECULAR STRUCTURE, vol.1217, 2020 (SCI-Expanded)
XXIV.	Calculation of the Raman frequency and linewidth of vibrons using anharmonic self energy model
	for the ε, δloc and δ phases in solid nitrogen
	Akay Ö., YURTSEVEN H. H.
	Optik, vol.219, 2020 (SCI-Expanded)
XXV.	Landau mean-field model with the cubic term for the alpha-beta transition in quartz
	Ates S., Yurtseven H.
	BULLETIN OF MATERIALS SCIENCE, vol.43, no.1, 2020 (SCI-Expanded)
XXVI.	Calculation of the thermodynamic functions using a mean field model for the fluid-solid transition in
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	AKAY Ö., YURTSEVEN H. H.
	Revue Roumaine de Chimie, vol.65, no.5, pp.425-432, 2020 (SCI-Expanded)
XXVII.	Liquid-(Solid plus Liquid) Transitions in a Two-Component System of (CH3)CCl3 + CCl4
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XXVIII.	Sm A - ferroelectric-Sm C transition in the mixture of C7+10.0.4 and in the ferroelectric C7
777 III.	Aksoy S., Kurt M., Yurtseven H.
	FERROELECTRICS, vol.554, no.1, pp.21-29, 2020 (SCI-Expanded)
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	formate frameworks
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XXX.	Temperature dependence of the IR mode frequency calculated from the volume data for DMNaFE
	Alguel G., Enginer Y., Yurtseven H.
	FERROELECTRICS, vol.551, no.1, pp.229-234, 2019 (SCI-Expanded)
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	KİRACI A., Yurtseven H.
	FERROELECTRICS, vol.551, no.1, pp.235-244, 2019 (SCI-Expanded)
XXXII.	Analysis and calculation of the magnetization, magnetic susceptibility and the specific heat close to
	phase transitions in heterometallics
	Dugan E. K., Yurtseven H.
	CURRENT APPLIED PHYSICS, vol.19, no.10, pp.1096-1102, 2019 (SCI-Expanded)
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 SrTiO3
	YURTSEVEN H. H., Kiraci A.
	FERROELECTRICS, vol.551, no.1, pp.143-151, 2019 (SCI-Expanded)
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	metal organic frameworks
	Kurt M., Yurtseven H., Kurt A., Aksoy S.
	CHINESE PHYSICS B, vol.28, no.6, 2019 (SCI-Expanded)

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
	Yurtseven H., Tari O.
	INTERNATIONAL JOURNAL OF MODERN PHYSICS B, vol.33, no.11, 2019 (SCI-Expanded)
XXXVI.	Temperature dependence of magnetic and thermal properties of chiral HyFe and HyMn close to
	phase transitions by using the Landau mean field model
	Tari O., Yurtseven H.
	MATERIALS CHEMISTRY AND PHYSICS, vol.228, pp.118-123, 2019 (SCI-Expanded)
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XXXVII.	Calculation of the Raman and IR frequencies as order parameters and the damping constant
	(FWHM) close to phase transitions in methylhydrazinium structures
	KURT M., YURTSEVEN H. H., Kurt A.
	JOURNAL OF MOLECULAR STRUCTURE, vol.1181, pp.488-492, 2019 (SCI-Expanded)
XXXVIII.	Resonant frequency shifts related to the elastic constants near the alpha-beta transition in quartz
	YURTSEVEN H. H., Ates S.
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	epsilon) TRANSITIONS IN MOLECULAR NITROGEN BY USING MEAN FIELD MODEL
	Yurtseven H., Akay O.
	REVUE ROUMAINE DE CHIMIE, vol.64, no.3, pp.249-258, 2019 (SCI-Expanded)
XL.	Calculation of the heat capacity Cp from the temperature-induced and pressure-induced Raman
	frequency shifts for solid benzene, naphthalene and anthracene
	Ozdemir H., Yurtseven H. H.
	Optik, vol.178, pp.774-784, 2019 (SCI-Expanded)
XLI.	Calculation of the T-P Phase Diagrams for the Halogenomethane Compounds (CCl4-nBrn, n=0, 1, 2,
	4) Using the Mean Field Theory
	Yurtseven H., Isik S. B., Dogan E. K.
	PHYSICS OF THE SOLID STATE, vol.61, no.2, pp.201-206, 2019 (SCI-Expanded)
XLII.	Temperature and the field dependence of the magnetization close to order-disorder phase
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	transitions in DMMn and the chromium-doped DMMn
	YURTSEVEN H. H., Doğan E.
	POLYHEDRON, vol.154, pp.132-137, 2018 (SCI-Expanded)
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	to the Sm A - Sm C* transition in C7
	AKSOY S., YURTSEVEN H. H., KURT M.
	CANADIAN JOURNAL OF PHYSICS, vol.96, no.7, pp.688-692, 2018 (SCI-Expanded)
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	Binary Mixture CM/PCPB
	YURTSEVEN H. H., Doğan E.
	RUSSIAN JOURNAL OF PHYSICAL CHEMISTRY A, vol.92, no.6, pp.1208-1212, 2018 (SCI-Expanded)
XLV.	Correlations between the resonant frequency shifts and the thermodynamic quantities for the alpha-
	beta transition in quartz
	Lider M. C., YURTSEVEN H. H.
	JOURNAL OF MOLECULAR STRUCTURE, vol.1159, pp.1-4, 2018 (SCI-Expanded)
XLVI.	T-P Phase Diagram of Nitrogen at High Pressures
	Algul G., Enginer Y., YURTSEVEN H. H.
	INTERNATIONAL JOURNAL OF THERMOPHYSICS, vol.39, no.5, 2018 (SCI-Expanded)
XLVII.	Analysis of the Specific Heat of Ru Doped LiKSO4 Close to Phase Transitions
	YURTSEVEN H. H., Tirpanci D. V., KARACALİ H.
	HIGH TEMPERATURE, vol.56, no.3, pp.462-465, 2018 (SCI-Expanded)
XLVIII.	Pressure-dependent Raman modes near the cubic-tetragonal transition in strontium titanate
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KİRACI A., YURTSEVEN H. H.

JOURNAL OF THE AMERICAN CERAMIC SOCIETY, vol.101, no.3, pp.1344-1355, 2018 (SCI-Expanded)

XLIX.Calculation of the Raman frequency and the damping constant (linewidth) of the stretching modes<br/>for the metal-organic compound DMMg close to the paraelectric-ferroelectric transitions<br/>YURTSEVEN H. H., Aslan A.

FERROELECTRICS, vol.526, no.1, pp.9-15, 2018 (SCI-Expanded)

L. Calculation of the C-P-C-V as a function of temperature close to the melting point in benzene Tari O., YURTSEVEN H. H.

PHYSICS AND CHEMISTRY OF LIQUIDS, vol.56, no.5, pp.660-673, 2018 (SCI-Expanded)

LI. Calculation of the Raman and IR frequencies from the volume data at high pressures in N-2 AKAY Ö., YURTSEVEN H. H.

OPTIK, vol.160, pp.227-233, 2018 (SCI-Expanded)

LII. Calculation of the thermodynamic quantities from the Raman frequency shifts and the Pippard relations for the alpha, -alpha and alpha ' phases of oxygen Tari O., YURTSEVEN H. H.

OPTIK, vol.158, pp.699-704, 2018 (SCI-Expanded)

- LIII. ANALYSIS OF THE INTEGRATED INTENSITY OF THE CENTRAL PEAKS CALCULATED AS A FUNCTION OF TEMPERATURE IN THE FERROELECTRIC PHASE OF LITHIUM TANTALATE Kiraci A., Yurtseven H. H. THERMAL SCIENCE, vol.22, 2018 (SCI-Expanded)
- LIV. Calculation of the P-T phase diagram of nitrogen using a mean field model Enginer Y., Algul G., YURTSEVEN H. H. INTERNATIONAL JOURNAL OF MODERN PHYSICS B, vol.31, no.31, 2017 (SCI-Expanded)
- LV. Analysis of Magnetization as a Function of Temperature for CoMn1-x Fe (x) Ge
  YÜCE EMRE S., Doğan E., Emre B., Bruno N. M., Karaman I., YURTSEVEN H. H.
  JOURNAL OF SUPERCONDUCTIVITY AND NOVEL MAGNETISM, vol.30, no.12, pp.3587-3594, 2017 (SCI-Expanded)
- LVI. Calculation of the Spontaneous Polarization and the Dielectric Constant for the Ferroelectric N(CH3)
  (4)HSO4 Using the Mean Field Model
  YURTSEVEN H. H., ÇELİK M., KARACALİ H.
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- LVII. **Pressure dependence of the heat capacity near the melting point in benzene** YURTSEVEN H. H., Tari O. JOURNAL OF MOLECULAR LIQUIDS, vol.241, pp.59-63, 2017 (SCI-Expanded)

# LVIII. Calculation of the tilt angle and susceptibility for the alpha-beta transition in quartz using a mean field model

YURTSEVEN H. H., Ipekoglu U., Ates S.

MODERN PHYSICS LETTERS B, vol.31, no.9, 2017 (SCI-Expanded)

# LIX. Analysis of the susceptibility of condensed oxygen under high pressures and in strong magnetic fields

Doğan E., YURTSEVEN H. H.

JOURNAL OF MOLECULAR STRUCTURE, vol.1131, pp.236-241, 2017 (SCI-Expanded)

## LX. Analysis of the Magnetic Susceptibility as a Function of Temperature for the alpha, beta and gamma Phases of Oxygen

YURTSEVEN H. H., Avci C.

JOURNAL OF SUPERCONDUCTIVITY AND NOVEL MAGNETISM, vol.30, no.3, pp.831-838, 2017 (SCI-Expanded)

### LXI. Temperature dependence of the damping constant and the relaxation time close to the tetragonalcubic phase transition in SrZrO3 YURTSEVEN H. H., Kiraci A.

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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. OPTIK, vol.144, pp.224-231, 2017 (SCI-Expanded) LXIII. Calculation of the Raman frequency shifts for the alpha phase of solid oxygen YURTSEVEN H. H., Tari O. OPTIK, vol.128, pp.113-120, 2017 (SCI-Expanded) 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. OPTIK, vol.135, pp.244-251, 2017 (SCI-Expanded) LXV. Pressure dependence of the Raman frequency calculated from the volume data close to the ferroelectric-paraelectric transition in PbTiO3 YURTSEVEN H. H., KİRACI A. FERROELECTRICS, vol.520, no.1, pp.245-255, 2017 (SCI-Expanded) LXVI. Temperature dependence of the polarization, dielectric constant, damping constant and the relaxation time close to the ferroelectric-paraelectric phase transition in LiNbO3 Kiraci A., YURTSEVEN H. H. OPTIK, vol.132, pp.183-191, 2017 (SCI-Expanded) LXVII. Calculation of the raman frequency, damping constant (Linewidth) and the relaxation time near the tetragonal-cubic transition in PbTiO3 KİRACI A., YURTSEVEN H. H. OPTIK, vol.142, pp.311-319, 2017 (SCI-Expanded) LXVIII. Damping Constant (Linewidth) and the Relaxation Time of the Brillouin LA Mode for the Ferroelectric-Paraelectric Transition in PbZr1-xTix03 YURTSEVEN H. H., Kiraci A. IEEE TRANSACTIONS ON ULTRASONICS FERROELECTRICS AND FREQUENCY CONTROL, vol.63, no.10, pp.1647-1655, 2016 (SCI-Expanded) LXIX. Variation of the molar volume with the temperature and the Pippard relations near the melting point in benzene Tari O., YURTSEVEN H. H. JOURNAL OF MOLECULAR LIQUIDS, vol.220, pp.883-887, 2016 (SCI-Expanded) LXX. Temperature and pressure dependence of the Raman frequency shifts in anthracene Ozdemir H., YURTSEVEN H. H. INDIAN JOURNAL OF PURE & APPLIED PHYSICS, vol.54, no.8, pp.489-494, 2016 (SCI-Expanded) 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 YURTSEVEN H. H., Yilmaz A. FRONTIERS IN PHYSICS, vol.4, 2016 (SCI-Expanded) LXXII. Analysis of the orientational order parameter and the magnetic field as functions of temperature for p-anisaldazine YURTSEVEN H. H., Kaya E. JOURNAL OF THE KOREAN PHYSICAL SOCIETY, vol.68, no.11, pp.1317-1321, 2016 (SCI-Expanded) LXXIII. Temperature effect on the structural ferroelectric - paraelectric transition in Li2Ge7015 YURTSEVEN H. H., Tari O. FERROELECTRICS, vol.494, no.1, pp.170-181, 2016 (SCI-Expanded) LXXIV. Molar volume calculated at various pressures and the Pippard relations close to the melting point in benzene YURTSEVEN H. H., Tari O. PHYSICS AND CHEMISTRY OF LIQUIDS, vol.54, no.2, pp.182-192, 2016 (SCI-Expanded) LXXV. Analysis of the peak position and linewidth as a function of temperature for the phase transitions in LiKSO4 YURTSEVEN H. H., KURT M.

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