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### Advising Theses

TOMAK M., Growth and morphological characterization of intrinsic hydrogenated amorphous silicon thin film for a-Si:H/c-Si heterojunction solar cells, Doctorate, Ö.PEHLİVAN(Student), 2013

TOMAK M., First-principles study of GaAs/AlAs nanowire heterostructures, Doctorate, S.ŞENOZAN(Student), 2012

TOMAK M., Density functional investigation of nano-structures, Doctorate, O.ÜZENGİ(Student), 2010

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TOMAK M., Ensemble monte carlo simulation of quantum well infrared photodetectors, Doctorate, S.MEMİŞ(Student), 2006

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TOMAK M., Electron-electron interactions in a two dimensional quantum dot, Doctorate, N.AKMAN(Student), 1999

TOMAK M., Molecular dynamics computer simulation of solid and liquid metal alloys, Postgraduate, D.GÜLTOPRAK(Student), 1999

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- I. The effect of Coulomb interaction on optical absorption of a quantum well wire  
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- II. **Exciton related nonlinear optical properties of a quantum wire**  
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- V. **Adsorption of RuSex (x=1-5) cluster on Se-doped graphene: First principle calculations**  
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- VI. **Study of the a-Si:H/c-Si Heterointerface by Ex-Situ Spectroscopic Ellipsometry, Infrared Spectroscopy, and Solar Cell Modeling**  
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- IX. **Electronic and geometric structure of AuxCuy clusters studied by density functional theory**  
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- XI. **Structural characterization of intrinsic a-Si:H thin films for silicon heterojunction solar cells**  
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- XV. **Thermodynamics of a two-dimensional interacting Bose gas trapped in a quartic potential**  
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- XVI. **NH<sub>3</sub> and H<sub>2</sub>S adsorption on Au<sub>3</sub>Pt<sub>3</sub> cluster studied by a first principles calculation**  
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- XIX. **Bismuth doping of graphene**  
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#### **Non Academic Experience**

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