About IIT Kharagpur







Kharagpur - a dusty town tucked away in the eastern corner of India, famous until 1950 as home to the longest railway platform in the world - became the nursery where the seed of the IIT system was planted in 1951. IIT Kharagpur started its journey in the old Hijli Detention Camp in Eastern India, where some of the country's great freedom fighters toiled and sacrificed their lives for India's independence. Spurred by the success of IIT Kharagpur, four younger IITs sprouted around the country in the two following decades, and from these five came thousands of IITians, the brand ambassadors of modern India. It was the success of this one institution at Kharagpur that wrote India's technological odyssey.

The Institute takes pride in its relentless effort to provide the best platform for both education as well as research in the areas of science and technology, infrastructure designs, entrepreneurship, law, management, and medical science and technology. IITKGP is not just the place to study technology, it is the place where students are taught to dream about the future of technology and beam across disciplines, making differences enough to change the world.

Program Features/ Structure

Classroom lectures - 50%

Numerical / Problem solving, Case study and Activity - **25%**

Hands-on work with molecular dynamics coding and software - **25%**

Program Fee

Nil for AICTE-QIP sponsored participants

For students - INR 5,000 + GST@18% per participant

For others - INR 15,000 + GST @ 18% per participant

Last day of Registration



May 2020

How to Apply

Use the link: https://erp.iitkgp.ac.in/CEP/courses.htm to apply ONLINE.



Payment if applicable is to be done **ONLINE** after getting short listed for the program.

Contact Us

Dr. Puneet Kumar Patra, Co-ordinator, Department of Civil Engineering Indian Institute of Technology Kharagpur Phone: +91-3222-283414 Email: <u>puneet.patra@civil.iitkgp.ac.in</u>

Program Schedule

2020 (9:00 AM - 5:00 PM)

1 week. 8 - 12 June

IIT Kharagpur - Civil

Who will benefit (Eligibility)

1. For faculty/students of Civil, Mechanical, Aerospace,

Materials / Nanotechnology,

Physics, Chemistry, and any

2. Industrial researchers

AC accommodation will be provided to the AICTE-QIP sponsored participants at

the campus Guesthouse. Limited AC accommodation is available for others on

first-come-first-serve basis at

Apply Now

Accommodation

Metallurgy, Chemical,

other allied.

no extra cost.

Engg. Department.

and Metallurgical

and Materials

Science Engg. Department

and Venue





QUALITY IMPROVEMENT PROGRAMME

Indian Institute of Technology Kharagpur 2020

Molecular Dynamics Simulations in Materials Science & Engineering

1 Week

8 – 12 June 2020

Introduction /Overview

Atomistic modeling of materials has become a popular approach to explore the properties of materials from a modern perspective. Here, a material is simulated by directly computing the dynamics of its constituent atoms from first-principle kinematics. Over the last decade, this technique has witnessed tremendous interest with the advent of enhanced capabilities of modern hardware and algorithms. The practical utility of this method spans several traditional engineering and science disciplines, and the potential applications are limitless. Using the same method, one can find the mechanical properties of a material as well as engineer new materials to have better properties.

Program Objectives

This course aims at introducing the fundamental and practical aspects of atomistic simulations. Both theoretical and hands-on practical aspects shall be covered comprehensively. The theory lectures will commence with a gentle introduction to the basic philosophy of molecular dynamics simulations and gradually progress towards the finer details like structural optimization, time-integrators, thermostats and barostats, etc. Finally, the lectures will demonstrate applications of these tools from the standpoint of fundamental research and domains of industrial applications. Apart from the theory classes, there will be a laboratory component, where the participants will implement the algorithms taught in the theory classes.

This course is designed such that it is beneficial to people from various backgrounds including physics, metallurgy and materials engineering, chemical engineering, civil engineering, mechanical engineering, etc. The topics covered span a broad spectrum of interests ranging from purely fundamental research to industrial R&D.

What you will learn

Program Content

Introduction to molecular statics and dynamics; length and time scales associated with MD simulations

Difference between continuum and atomistic mechanics

Bonding and interatomic forces; models of various interatomic potentials

Deriving force fields from potentials; creating new potentials

Applications of optimization algorithms in molecular statics simulations

Time integration algorithms - Leapfrog, Velocity Verlet, Runge-Kutta; symplectic and non-symplectic algorithms

Thermostats and Barostats - Nose-Hoover, Nose-Hoover chains, Nose-Poincare, Gaussian isokinetic

Relationship with classical thermodynamics different laws of thermodynamics, ergodicity, canonical ensemble, Green-Kubo relations

Filtering and visualization in atomistic simulations; exploring structural and mechanical behavior

Computing transport properties - thermal conductivity, viscosity - using molecular dynamics

Challenges in molecular dynamics simulations

Applications in fundamental and industrial research

Computing elastic constants of crystals with molecular statics simulations

Computing thermal expansion coefficient with molecular dynamics

Extracting thermal conductivity with non-equilibrium molecular dynamics

About the Faculty

Dr. Puneet Kumar Patra



Dr. Puneet Kumar Patra is an Assistant Professor in the Department of Civil Engineering, IIT Kharagpur, Prior to this, he worked as a postdoctoral research scholar in the Department of Biomedical Engineering and Mechanics, Virginia Polytechnic Institute and State University. His research interests include nonequilibrium statistical mechanics. thermodynamics, molecular dynamics simulations and thermal transport characteristics in low dimensional systems. He holds a PhD degree in Mechanics from IIT Kharagpur and a B.Tech. (H) degree in Civil Engineering also from IIT Kharagpur.

Dr. Amlan Dutta



Dr. Amlan Dutta is an Assistant Professor in the Department of Metallurgical and Materials Engineering, IIT Kharagpur. He has also worked as a scientist at the Dept. of condensed matter physics and material science of the S.N. Bose National Centre for Basic Sciences, Kolkata. His research interests involve atomistic modeling, general scientific computing, data science and materials informatics, theoretical micromechanics of solids, multiscale modeling in materials science, etc. He holds a PhD from the Variable Energy Cyclotron Centre, Dept. of atomic energy, Govt. of India.