# PX923-10 Biomolecular simulation

#### 24/25

#### **Department**

**Physics** 

Level

**Taught Postgraduate Level** 

Module leader

Nicholas Hine

Credit value

10

**Module duration** 

5 weeks

**Assessment** 

50% coursework, 50% exam

**Study location** 

University of Warwick main campus, Coventry

### **Description**

### Introductory description

N/A.

Module web page

### **Module aims**

The aim of the module is to describe computational methods used in physical chemistry and their application to study the structure and dynamics of biological molecules.

A number of methods in molecular simulation, and their respective theoretical foundations, will be presented. There will be an emphasis on approaches used to probe biological phenomena and calculations of properties pertinent to biological systems.

### **Outline syllabus**

This is an indicative module outline only to give an indication of the sort of topics that may be covered. Actual sessions held may differ.

- 1. Force fields for biomolecular simulation.
- 2. Free energy calculations. Thermodynamic integration and perturbation; potentials of mean

force.

- 3. Advanced sampling. Biased sampling methods, replica exchange molecular dynamics. Coarse-graining for biomolecular simulation.
- 4. Calculated properties. Energetics, structure and dynamics; relationship to experimentally determined quantities.
- 5. Example applications. Drug-enzyme binding free energies and computer-aided drug design, protein structure prediction, delivery of pharmaceuticals across lipid membranes, the interface between nanomaterials and biological systems.

### **Learning outcomes**

By the end of the module, students should be able to:

- Students will be able to describe the theoretical background and application of a selection of computational methods (including molecular dynamics, free energy calculations and techniques to study rare events) in biophysical chemistry.
- Students will have an appreciation of the advantages and disadvantages of different computational methods in the context of solving particular biological problems (e.g. protein folding, membrane transport, drug binding).
- Students will be able to evaluate the strengths and weaknesses of biomolecular simulation studies found in the literature.
- Students will be able to perform and analyse computational calculations involving biological molecules and relate them to experimental data.

### Indicative reading list

- A.R. Leach, Molecular Modelling: principles and applications, Longman (1996).
- M. Tuckerman, Statistical Mechanics, Oxford Graduate Texts (2010).
- D. Frenkel & B. Smit, Understanding Molecular Simulation, From Algorithms to Applications (2001).
- A selection of papers from the recent research literature, that may change from year to year, will also be provided.

### Subject specific skills

Students will be able to describe the theoretical background and application of a selection of computational methods (including molecular dynamics, free energy calculations and techniques to study rare events) in biophysical chemistry.

Students will have an appreciation of the advantages and disadvantages of different computational methods in the context of solving particular biological problems (e.g. protein folding, membrane transport, drug binding).

Students will be able to evaluate the strengths and weaknesses of biomolecular simulation studies found in the literature.

Students will be able to perform and analyse computational calculations involving biological molecules and relate them to experimental data.

#### Transferable skills

Biochemistry knowledge, programming, data analysis

# **Study**

### Study time

Required **Type** 

5 sessions of 2 hours (10%) Lectures Practical classes 2 sessions of 2 hours (4%)

61 hours (61%) Private study 25 hours (25%) Assessment

Total 100 hours

## **Private study description**

Reading etc

#### Costs

No further costs have been identified for this module.

#### Assessment

You do not need to pass all assessment components to pass the module.

#### **Assessment group C1**

Weighting Study time Eligible for self-certification

Assessment component

Computational Workshop Projects 1 and 50% 2

20 hours No

One piece of assessed work based on the workshops. Write up of the 2 computational workshops.

#### Weighting Study time Eligible for self-certification

Reassessment component is the same

**Assessment component** 

Viva Voce Exam 50% 5 hours No

Duration 15 minutes. On the core material.

Reassessment component is the same

#### Feedback on assessment

-Written feedback on assessed work. Verbal discussion during viva voce exam. Written summary of viva performance.

Past exam papers for PX923

# **Availability**

There is currently no information about the courses for which this module is core or optional.