

## سمینار هفتگی ماده چگال نرم

How to parametrize forcefields in coarse-graining approaches

## **Abstract**

In this lecture, I try to introduce some coarse-graining approaches, their applications, and their limitations. Then I try to introduce the parametrization method of the Shape-base Coarse-graining (SBCG) approach. It should be noted that deriving potentials for coarse-grained Molecular Dynamics (MD) simulations are frequently done by solving an inverse problem. Methods like Iterative-Boltzmann-Inversion (IBI) or Inverse Monte Carlo (IMC) have been widely used to solve this problem. The solution obtained by the application of these methods guarantees a match in the radial distribution function (RDF) between the underlying fine-grained system and the derived coarse-grained system. In the SBCG approach, the IBI method is used for parameterization. We are going to look deeper into these methods in the context of the SBCG approach.

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