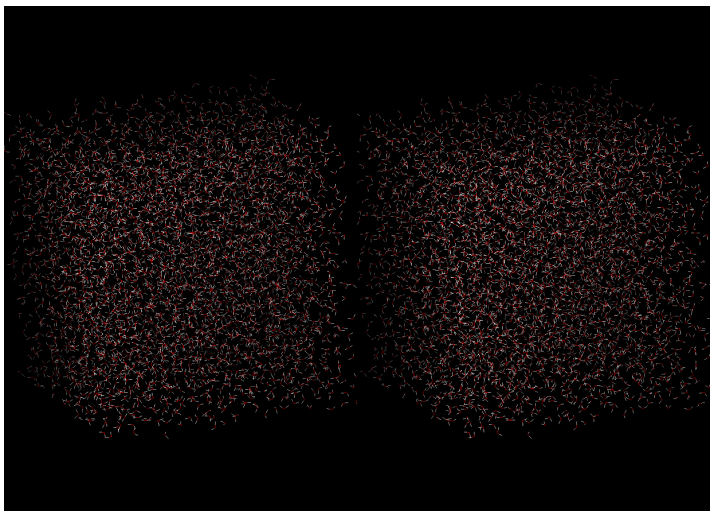
Molecular Dynamics
(Tinker)
An EMBnet introductory course



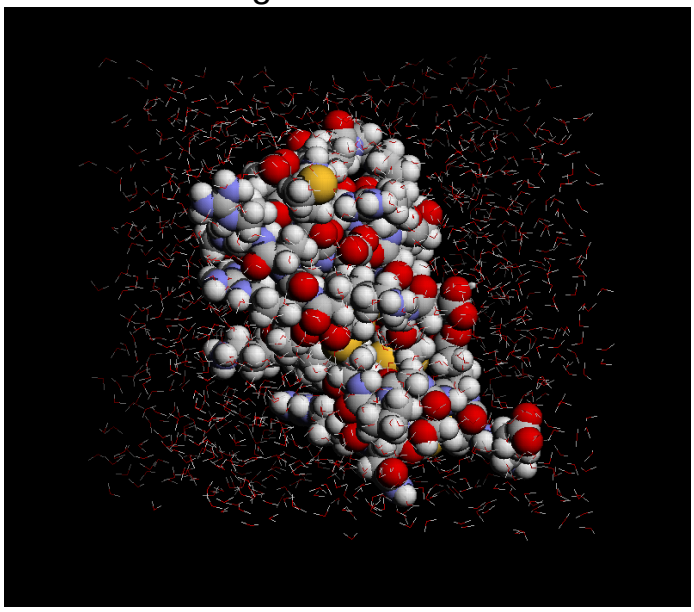
A practical introduction to **Molecular Dynamics simulations using TINKER**.

EMBnet/CNB
Salón de Actos, CNB, Madrid
10-11 noviembre 2008
09:30-16:30

Topics covered include:

- Overview of methods used in Molecular Dynamics
- Practical sessions using TINKER
  - Protein simulation *in vacuo* and implicit solvent models
  - Molecular Dynamics with periodic boundary conditions
  - Protein simulation in fully atomistic solvent models

The course will rely on e-learning tutoring for extended training.



Assistance limited to 20 places
For further information:
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with
Subject: EMBNET COURSE 2008

Web Site:
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EMBnet provides this course for free, but places are limited to 20.