

**数学与系统科学研究院**

**计算数学所学术报告**

**报告人: Dr. Zhongyuan Zhou**

**(Senior Research Associate Department  
of Chemistry University of Kansas USA)**

**报告题目:**

**A moving-grid approach in phase  
space: A promising method for the  
study of dynamics of many-electron  
atomic and molecular systems**

**邀请人: 洪佳林研究员**

**报告时间: 2009年6月30日(周二)**

**上午 10:00—11:00**

**报告地点: 科技综合楼三层 301**

**计算数学所报告厅**

## Abstract:

A moving-grid approach based on coupled coherent states and electron trajectories is developed for the study of dynamics of many-electron atomic and polyatomic molecular systems. In this approach, the initial state and trajectories are determined by using Monte Carlo method. And then the initial state propagates on the moving grids guided by the electron trajectories. As a demonstration, this approach is applied to the study of hydrogen molecule (H<sub>2</sub>).

**欢迎大家参加！**