

请各位老师在下方空白处插入一张一寸免冠照片，设置高度为 4.80 厘米，宽度为 3.65 厘米，居中对齐



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	主讲课程	-		
	科研方向	钢铁冶金、氢冶金、分子动力学		
教育及工作经历	2013.09—2017.06 北京科技大学高等工程师学院冶金工程（卓越计划）专业本科学习 2017.09—2023.06 北京科技大学冶金与生态工程学院冶金工程专业博士学习 2021.11—2022.12 多伦多大学冶金工程专博士业联合培养学习 2023.08 至今 北京科技大学金属冶炼重大事故防控技术支撑基地博士后工作			

代表性成果（包含论文、著作、获奖、专利、项目等）

获奖：

北京科技大学校长奖章、博士国家奖学金、硕士国家奖学金

代表论文：

- [1] **Chunhe Jiang**, Wang Liang, Kejiang Li, et al. A reactive molecular dynamics study of thermal pyrolysis behavior and mechanisms of lignin during the hydrothermal process: The function of the water molecules. *Bioresource Technology*, 2023, 368: 128338.
- [2] **Chunhe Jiang**, Kejiang Li, Mansoor Barati, et al. The interaction mechanism between molten SiO₂-Al₂O₃-CaO slag and graphite with different crystal orientations: Experiment and Ab initio molecular dynamics simulation. *Ceramics International*, 2023, 49(5), 8295-8301.
- [3] **Chunhe Jiang**, Mansoor Barati, Kejiang Li, et al. The dissolutive wetting behavior between liquid iron and carbonaceous materials: Experiment and ReaxFF Molecular Dynamics Simulation. *Journal of Molecular Liquids*, 2022, 360: 119435.
- [4] **Chunhe Jiang**, Kejiang Li, Jianliang Zhang, et al. Structural characteristics of liquid iron with various carbon contents based on atomic simulation. *Journal of Molecular Liquids*, 2021, 342: 116957.
- [5] **Chunhe Jiang**, Jianliang Zhang, Kejiang Li, et al. Influence of graphite crystalline orientation on the carbon dissolution reaction in liquid iron: A ReaxFF molecular dynamics simulation study. *Journal of Molecular Liquids*, 2021, 335: 115688.
- [6] **Chunhe Jiang**, Zixin Xiong, Yushan Bu et al. Study on the Structure and Properties of High-Calcium Coal Ash in the High-Temperature Zone of a Blast Furnace: A Molecular Dynamics Simulation Investigation. *JOM*, 2020, 72(7): 2713-2720.
- [7] **Chunhe Jiang**, Zixin Xiong, Kejiang Li et al. Molecular dynamics simulation study on the wetting behavior of liquid iron and graphite. *Journal of Molecular Liquids*, 2020, 311; 113550.
- [8] **Chunhe Jiang**, Haoxiang Zhang, Zixin Xiong et al. Molecular dynamics investigations on the effect of Na₂O on the structure and properties of blast furnace slag under different basicity conditions. *Journal of Molecular Liquids*, 2020, 299: 112195.

- [9] **Chunhe Jiang**, Kejiang Li, Jianliang Zhang et al. The effect of CaO and MgO on the structure and properties of coal ash in the blast furnace: A molecular dynamics simulation and thermodynamic calculation. *Chemical Engineering Science*, 2019, 210: 115226.
- [10] **Chunhe Jiang**, Kejiang Li, Jianliang Zhang et al. Effect of MgO/Al₂O₃ ratio on the structure and properties of blast furnace slags: A molecular dynamics simulation. *Journal of Non-Crystalline Solids*, 2018, 502: 76-82.
- [11] **Chunhe Jiang**, Kejiang Li, Jianliang Zhang et al. The effect of CaO(MgO) on the structure and properties of aluminosilicate system by molecular dynamics simulation. *Journal of Molecular Liquids*, 2018, 268: 762-769.
- [12] **Chunhe Jiang**, Kejiang Li, Jianliang Zhang et al. Molecular dynamics simulation on the effect of MgO/Al₂O₃ ratio on structure and properties of blast furnace slag under different basicity conditions[J]. *Metallurgical and Materials Transactions B*, 2019, 50(1): 367-375.
- [13] 张建良, **姜春鹤**, 李克江, 等. 高炉内渣铁焦界面润湿行为研究现状及展望. *钢铁*, 2021, 56(11): 10-18.