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您现在的位置: [首页](#) > [教师简介](#) > Content

教师简介

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【个人简介】

王怀谦，博士，副教授，硕士生导师。

2011年6月毕业于四川大学原子与分子物理专业，获工学博士学位，获宝钢教育奖，入选2013年福建省高校杰出青年科研人才计划、2015年泉州市高层次创新创业人才项目。主要从事纳米团簇结构与物性研究，先后主持承担国家自然科学基金项目。已在国际SCI杂志《J. Chem. Phys.》、《J. Phys. Chem.》、《Sci. China-Phys. Chem.》、《Adv. Mater.》等公开发表科研论文40多篇，其中有21篇

的国际著名刊物上，第一或通讯作者发表JCR Top
累积影响因子大于90，研究论文受到广泛关注，
内的国际权威SCI期刊累计引用393篇次。现为
Chem. Phys.》、《J. Chem. Phys.》、《J. Mater
Mater》、《Nanoscale》、《New J. Chem.》、《
家，至今已审稿60余篇次。

【受教育经历】

(1)2011.06 四川大学，原子与分子物理研究所，博士

(2)2006.06 鲁东大学，物理与电子工程学院，学士

【研究工作经历】

(1) 2013年7月~至今，华侨大学，工学院，副教授

(2) 2015年2月~2015年8月，工程物理研究院北京

(3) 2011年7月~2013年6月，华侨大学，工学院，讲师

【获奖情况】

(1) 教育部“宝钢教育奖”

- (2) 福建省新世纪优秀人才计划
- (3) 福建省高校杰出青年科研人才培育计划
- (4) 泉州市引进高层次创业创新人才
- (5) 泉州市科学技术奖(自然科学奖)二等奖(排名)
- (6) 第十二届福建省自然科学优秀学术论文三等
- (7) 第八届泉州市自然科学优秀学术论文二等奖
- (8) 华侨大学百门优质课程《大学物理》
- (9) 华侨大学就业创业导师
- (10) 华侨大学2012~2014学年优秀班主任
- (11) 华侨大学2012年首届青年教师“精彩一堂课”竞

【主持项目】

- (1) 国家自然科学基金(11247257)
- (2) 福建省新世纪优秀人才项目(2014FJ-NCET-ZR)
- (3) 福建省高校杰出青年科研人才计划项目(JA13)
- (4) 福建省自然科学基金面上项目(2017J01001)
- (5) 福建省自然科学基金青年项目(2012J05005)
- (6) 泉州市高层次人才项目(2018C077R)
- (7) 中央高校基本科研项目(JB-ZR1201)

(8) 华侨大学引进人才科研启动项目(11BS421)

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- (2) Yi-Wei Fan, **Huai-Qian Wang*** and Hui-Fang Li, Density functional theory calculations coupled with stochastic resonance. **Part A: Molecular and Biomolecular Spectroscopy**, JCR 二区, IF:3.232
- (3) **Huai-Qian Wang*** and Hui-Fang Li, Probing the structure of vanadium dioxide clusters by density functional theory photoelectron spectroscopy. **Journal of Chemical Physics**, JCR 一区, IF:3.333, Times Cited:16
- (4) Hui-Fang Li and **Huai-Qian Wang***, Probing the structure of metal-doped golden cage nanoclusters: $M@Au_n$. **Chemical Physics** 16 (2014) 244-254 (SCI, top期刊)
- (5) **Huai-Qian Wang***, Hui-Fang Li, and Xiao-Yu Kuang, Properties of small vanadium monoxide clusters. **Journal of Chemical Physics** (2012) 5272-5283(SCI, top期刊, JCR 一区, IF:4.493)
- (6) **Huai-Qian Wang***, Xiao-Yu Kuang and Hui-Fang Li, Structure and electronic properties of bimetallic copper-gold clusters. **Physical Chemistry Chemical Physics**, JCR 一区, IF: 4.493, Times Cited: 79
- (7) Hui-Fang Li, **Huai-Qian Wang***, and Xiao-Yu Kuang, Structure and electronic properties of transition-metal ions in tetrahedral and octahedral clusters. **Journal of Chemical Physics** (2013) 124701-124708(SCI, top期刊, JCR 一区, IF:4.493)

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- (8) **Huai-Qian Wang**, Xiao-Yu Kuang, Hui-Fang Li, et al. The difference of local structure between EPR and XRD of NiSiF₆·6H₂O crystal. **Journal of Alloys and Compounds** 500 (2011) 1-6, IF:4.175, Times Cited: 3)
 - (9) Yi-Wei Fan, **Huai-Qian Wang*** and Hui-Fang Li, Properties of anionic europium-doped silicon clusters: comparison of experimental photoelectron spectroscopy and theoretical calculations. **Journal of Solid State Chemistry** (SCI, JCR三区, IF:1.822)
 - (10) Yi-Wei Fan, **Huai-Qian Wang*** and Hui-Fang Li, Properties of rare-earth doped silicon-based clusters: comparison of experimental photoelectron spectroscopy and theoretical calculations. **Journal of Solid State Chemistry** (2019) 221 (SCI, JCR二区, IF:1.871)
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 - (12) **Huai-Qian Wang*** and Hui-Fang Li, Structural characterization of rare-earth doped silicon-based nanoclusters. **RSC Advances** 5 (2015) 94685-94694
 - (13) **Huai-Qian Wang**, Xiao-Yu Kuang, and Hui-Fang Li, Properties of gold cluster anions doped with zinc: comparison of experimental photoelectron spectroscopy and theoretical calculations. **Chemistry A** 113 (2009) 14022-14028 (SCI, JCR三区)
 - (14) **Huai-Qian Wang***, Hui-Fang Li and Li-Xin Zhang, Synthesis and properties of rare-earth doped silicon-based clusters (M=Cr, Mn; q=0, -1) with adjustable magnetic properties. **Magnetic Materials** 344 (2013) 79-84 (SCI, JCR三区)
 - (15) **Huai-Qian Wang***, Hui-Fang Li, Jia-Xian Wan and Hui-Fang Li, Synthesis and properties of Al-doped niobium clusters. **Journal of Molecular Modeling** 18 (2012) 2993-3001 (SCI, JCR三区)

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- (17) **Huai-Qian Wang**, Xiao-Yu Kuang, and Hui-Fang Li, The structure of silver chloride. **Molecular Physics** 107 (2009) 621-626 (IF:2.291, Times Cited: 1)
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- (19) **Huai-Qian Wang*** and Hui-Fang Li, Density functional theory study of the electronic and magnetic properties of Nb^nO_n clusters. **Theoretical Chemistry Accounts** 114 (2013) 1006-1012 (SCI, IF:2.291, Times Cited: 7)
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