



卢胜杰副教授个人简介

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卢胜杰, 男, 博士, 副教授, 山东菏泽人, 出生于1988年2月, 2012年毕业于烟台大学应用化学系, 2017年获得中国科学院化学研究所物理化学专业博士学位, 攻读博士期间获得“中国科学院大学三好学生”、“中国科学院化学研究所所长优秀奖奖学金”、“中国科学院化学研究所青年科学优秀奖”、“中国科学院化学研究所优秀毕业生”等荣誉称号, 2017年9月通过“优秀博士引进计划”加入菏泽学院化学化工学院。菏泽学院多孔纳米功能材料研究所青年骨干。目前主持山东省自然科学基金项目(博士基金)一项(第八副族金属掺杂硅半导体团簇的光电子能谱和理论计算研究, ZR2018BB040, 9万)。曾参与国家自然科学基金面上项目一项(碱金属卤盐、亚硝酸盐、醋酸盐在水中溶解的微观机理研究, 21543007)和973项目一项。主要研究方向为金属掺杂半导体团簇的结构和性质、贵金属团簇催化C-H键及C-C键偶合反应机理。在Nanoscale, J. Phys. Chem.A, J. Phys. Chem. C, Phys. Chem. Chem. Phys., J. Chem. Phys., Sci. Rep., RSC Adv., Chem. Phys. Lett.等国际期刊累计发表SCI论文10余篇, 其中中科院SCI一区2篇, 中科院SCI二区10篇, 中科院SCI三区4篇。

代表性论文:

- (1) **Sheng-Jie Lu**, Xi-Ling Xu*, Hong-Guang Xu*, and Wei-Jun Zheng*, Structural evolution and bonding properties of Au₂Siⁿ⁻/₀ (n = 1-7) clusters: Anion photoelectron spectroscopy and theoretical calculations, J. Chem. Phys., 2018, 148:244306
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- (3) **Sheng-Jie Lu***, Probing the structural evolution and bonding properties of Pt_nC₂⁻/₀ (n = 1-7) clusters by density functional calculations, Chem. Phys. Lett., 2018, 699: 218-222
- (4) **Sheng-Jie Lu***, Probing the geometric structures and electronic properties of anionic and neutral Pt₃C₂ clusters by density functional calculations, Chem. Phys. Lett., 2018, 694:70-74
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- (10) Quoc Tri Tran, **Sheng-Jie Lu**, Li-Juan Zhao, Xi-Ling Xu, Hong-Guang Xu, Van Tan Tran*, Jun Li* and Wei-Jun Zheng*, [Spin-orbit splittings and low-lying electronic states of AuSi and AuGe: anion photoelectron spectroscopy and ab initio calculations](#), J. Phys. Chem. A, 2018, 122: 3374-3382
- (11) Xue Wu, **Sheng-jie Lu**, Xiaoqing Liang, Xiaoming Huang, Ying Qin, Maodu Chen, Jijun Zhao*, Hong-Guang Xu, R. Bruce King, and Weijun Zheng*, Structures and electronic properties of B₃Siⁿ⁻ (n = 4-10) clusters: a combined ab initio and experimental study, J. Chem. Phys., 2017, 146: 044306
- (12) Yuan-Yuan Jin, **Sheng-Jie Lu**, Andreas Hermann, Xiao-Yu Kuang*, Chuan-Zhao Zhang, Cheng Lu*, Hong-Guang Xu*, and Wei-Jun Zheng, Probing the structural evolution of ruthenium doped germanium clusters: photoelectron spectroscopy and density functional theory calculations, Sci. Rep., 2016, 6: 30116
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