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用Hammerling交换势计算碱金属原子高激发态能级

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Calculation of energy levels of alkali-metal-atoms highly excited states by using Hammerling exchange potential

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摘要 考虑了碱金属原子高激发态价电子与原子实电子coulomb相互作用、交换相互作用以及极化相互作用(长程关联效应),计算了碱金属原子Li,Na,K高激发态($6 \leq N \leq 15, l \leq 2$)波函数及能级.计算结果表明:交换势取为GKS势时,能级计算值比实验值偏低;取为hammerling势时,计算值与实验值符合很好.

关键词: [高激发态](#) [X_a法](#) [hammerling交换势](#)

Abstract: Considering the electron coulomb interaction and exchange interaction and polarization interaction (electron correlation effect)between the highly excited state electron of the alkali metal atom and the ion core electrons,we calculated the wave function and energy levels of alkali metal atom Li,Na,K highly excited states($6 \leq N \leq 15, l \leq 2$).The calculated results indicate that our energy levels values are lower than experimental data when the highly excited states exchange potential is replaced by GKS potential; If the highly excited states exchange potential is replaced by the hammerling potential our theoretical results are in good agreement with the experiment.

Key words: [highly excited state](#) [X_a method](#) [hammerling exchange potential](#)

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