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Geometric phase driven predissociation: Lifetimes of $2^2 A'$ levels of \mathbf{H}_3^{1} JUAN BLANDON, VIATCHESLAV KOKOOULINE, University of Central Florida — We discuss the role of the geometric phase in predissociation dynamics of vibrational states near a conical intersection of two electronic potential surfaces of a D_{3h} molecule. For quantitative description of the predissociation driven by the coupling near a conical intersection, we developed a method for calculating lifetimes and positions of vibrational predissociated states (Feshbach resonances) for X_3 molecule. The method takes into account the two coupled three-body potential energy surfaces, which are degenerate at the intersection. As an example, we apply the method to obtain lifetimes and positions of resonances of predissociated vibrational levels of the $2^2 A'$ electronic state of the \mathbf{H}_3 molecule. The three-body recombination rate coefficient for the $\mathbf{H}+\mathbf{H}+\mathbf{H} \rightarrow \mathbf{H}_2+\mathbf{H}$ process is estimated.

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