## UNDERSTANDING $\mathrm{CH}_{5}^{+}$SPECTROSCOPY FROM A PARTICLE-ON-A-SPHERE MODEL

## MICHAEL P. DESKEVICH, CHANDRA SAVAGE, DAVID J. NESBITT, JILA, Department of Chemistry and Biochemistry, University of Colorado, Boulder, Colorado 80309.

Due to the low barrier to H rearrangement in $\mathrm{CH}_{5}^{+}$, a good approximation is to separate the angular and radial H coordinates. By fixing $\mathrm{R}_{C H}$ at a constant value, the 15 degree-of-freedom problem becomes a more computationally feasible 10 degree-of-freedom problem. The reduced dimensional problem is well suited for capturing the essential low energy, large amplitude bending/rotation dynamics. The "particle-on-a-sphere" (POS) model, which has been shown to provide good experimental agreement in $\mathrm{XH}_{n}(\mathrm{n}=2-4)$ systems is extended to accommodate a 5 hydrogen system. Building on past success with the $\mathrm{XH}_{n}$ systems, we use the $\mathrm{XH}_{5}$ POS model to calculate the patterns of the low J rovibrational spectrum, facilitating the understanding of the jet-cooled $\mathrm{CH}_{5}^{+}$spectrum.

