

# Detectability of Prolate Symmetric-Top Molecules in Diffuse Clouds by a “Hot-Axis Effect”

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The aim of this study is to detect prolate symmetric-top molecules in diffuse clouds by observations of absorption lines by a “hot-axis effect.” Molecules in low excitation temperatures in diffuse clouds can be detected by only absorption. A prolate symmetric-top molecule, such as CH<sub>3</sub>CN and CH<sub>3</sub>CCH, has a special advantage to detect absorption lines. The  $K$  rotation (rotation around the red dotted lines in Figure 1) is excited by collisions in a high kinetic (collisional) temperature and is not cooled by radiation. On the other hand, the  $J$  rotation is cooled well by radiation due to a permanent dipole moment and the low radiative temperature as shown in Figure 1. As a result, the special advantage is that the  $J = K$  rotational levels have population concentrations as shown in Figure 2. We call the population concentrations a “hot-axis effect.” Absorption lines of the  $J + 1 \leftarrow J (= K)$  and  $\Delta K = 0$  transitions indicated by the red arrows in Figure 2 can be strong. To calculate the population concentrations, both of radiation and collisions should be considered simultaneously. In the case of a linear molecule the equation for rotational distribution by radiation and collisions was derived by Oka *et al.* [1]. Recently Araki *et al.* derived the equation in the case of a symmetric-top molecule having the hot-axis effect [2]. Using the equation the population concentrations of CH<sub>3</sub>CN and CH<sub>3</sub>CCH were calculated and the intensities of absorption lines were estimated. Therefore it is suggested that the absorption lines enhanced by the hot-axis effect can be the powerful probe of the molecules in diffuse clouds.

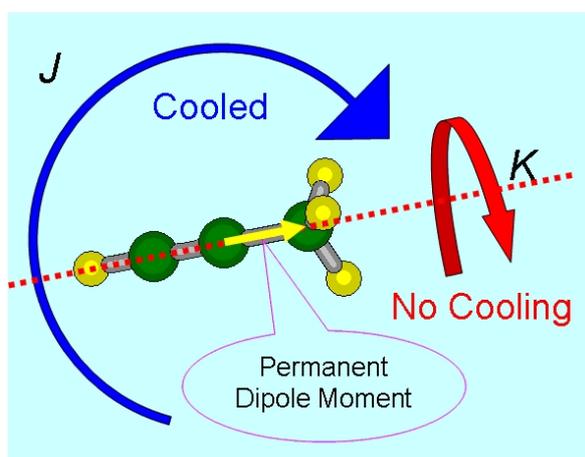


Figure 1: Hot-axis effect for a prolate symmetric top molecule. The red dotted line is the hot axis.

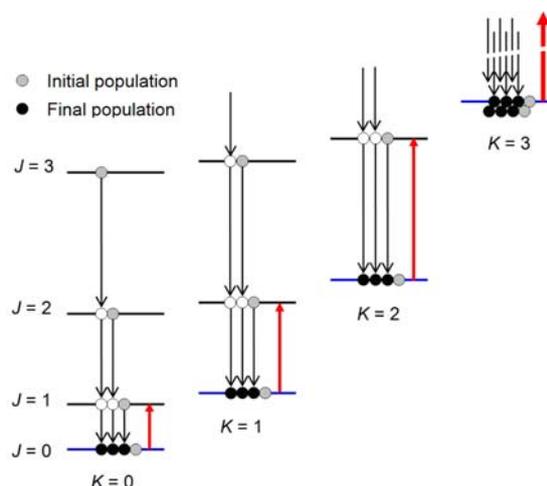


Figure 2: Population concentrations at the  $J = K$  levels for a prolate symmetric top molecule. The emissions are drawn by the black arrows. The  $J = K$  levels are described by the blue bars and the strong absorptions are by the red arrows.

## References

- [1] Oka *et al.*, 2013, ApJ, **773**, 42.
- [2] Araki *et al.*, 2014, AJ, in press.