DO WE NEED TO KNOW THE TEMPERATURE IN PRESTELLAR CORES?

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Received 2007 August 14; accepted 2007 September 26; published 2007 October 16

ABSTRACT

Molecular line observations of starless (prestellar) cores combined with a chemical evolution modeling and radiative transfer calculations are a powerful tool to study the earliest stages of star formation. However, conclusions drawn from such a modeling may noticeably depend on the assumed thermal structure of the cores. The assumption of isothermality, which may work well in chemo-dynamical studies, becomes a critical factor in molecular line formation simulations. We argue that even small temperature variations, which are likely to exist in starless cores, can have a nonnegligible effect on the interpretation of molecular line data and derived core properties. In particular, “chemically pristine” isothermal cores (low depletion) can have centrally peaked C^{18}O and C^{14}S radial intensity profiles, while having ringlike intensity distributions in models with a colder center and/or warmer envelope assuming the same underlying chemical structure. Therefore, derived molecular abundances based on oversimplified thermal models may lead to a misinterpretation of the line data.

Subject headings: astrochemistry — ISM: molecules — line: profiles — stars: formation

1. INTRODUCTION

The cooling in starless (or prestellar) cores is believed to be very effective, so it is customary to model these cores as isothermal at the temperature of \(~9–10\) K. In general, this assumption is confirmed both observationally and theoretically. However, different authors disagree on possible minor deviations from the uniform temperature, which may exist in these objects.

More precisely, various attempts to model the gas thermal balance in starless cores basically reach the same conclusion that the gas and dust temperatures are well coupled in the core center and have values of \(5–7\) K (e.g., Goldsmith 2001; Zucconi et al. 2001; Galli et al. 2002; Lesaffre et al. 2005). At a low-density periphery, heating processes due to cosmic rays and some UV irradiation take over, and the gas temperature rises to \(15\) K and higher (depending on the UV field strength), at the same time decoupling from the dust temperature.

On the other hand, recent observational evidence for the temperature structure in starless cores is somewhat controversial. Tafalla et al. (2002, 2004) have found an isothermal structure at \(\sim 9\) K for a sample of starless cores, including the L1544 and L1498 cores. Similarly, van der Tak et al. (2005) have found from \(^2\)H/D\(^2\) observations that the gas temperature is nearly constant or at least does not drop inward significantly in L1544. High-resolution observations have led Crapsi et al. (2007) to the conclusion that the gas temperature does go down in the very center of the L1544 core. A similar thermal structure (a few K in the center and more than \(10\) K in the envelope) is inferred for the L183 starless core by Pagani et al. (2007) and for the B68 globule by Pineda & Bensch (2007). Quite surprisingly, a reversed temperature radial profile for B68 is found by Bergin et al. (2006). These authors argue that the envelope of this globule is actually colder than its central region.

These deviations from isothermality would have only a subtle effect on results of dynamical and chemical modeling. However, there is an aspect that is more sensitive to temperature variations. In order to interpret observational data in terms of abundances and velocities, radiation transfer in molecular lines is commonly utilized. The synthetic line profiles depend critically on the radial distribution of the excitation temperature, which is a nonlinear function of the kinetic temperature \(T\). The importance of adequate temperature understanding has already been recognized, for example, by Di Francesco et al. (2002). These authors found that utilization of the nonuniform temperature profile with decreasing \(T\) at smaller radii increases abundances of C^{18}O, CS, and HCO\(^+\), inferred for B68, by factors of \(2–3\) in comparison with the isothermal case.

In this Letter, we study the influence of a nonuniform \(T\) distribution on synthetic integrated intensity \(J_{\text{int}}\) radial profiles for various CO and CS (and their isotopomers) transitions, using the L1544 and L1521E starless cores as templates. Observational data are taken from Tafalla et al. (2002) and Tafalla & Santiago (2004). The radiation transfer model as well as the assumed prestellar core chemical and thermal structure are briefly outlined in \S\ 2. Results for the detailed chemical modeling are presented in \S\ 3. Influence of the thermal structure on the inferred analytical abundance representations is discussed in \S\ 4. Results are summarized in \S\ 5.

2. MODEL

Synthetic radial profiles of the integrated intensities in a prestellar core are computed with the Monte Carlo code URAN(IA), described by Pavlyuchenkov & Shustov (2004). We utilize density distributions for the L1544 and L1521E cores, represented in Tafalla et al. (2002) and Tafalla & Santiago (2004) as \(n(r) = n_0/[1 + (r/r_0)^p]\), with \(p = 2.4\), \(r_0 = 2800\) AU, \(n_0 = 1.35 \times 10^6\) cm\(^{-3}\) for L1544 and \(p = 2\), \(r_0 = 4200\) AU, \(n_0 = 2.7 \times 10^5\) cm\(^{-3}\) for L1521E. A model core is assumed to be static, with a uniform radial microturbulent velocity \(v_t\) of 140 m s\(^{-1}\). We tried other \(v_t\) values as well and found that our conclusions do not depend on particular assumptions about this quantity.

Three different temperature profiles are considered (Fig. 1). In models designated “uni,” \(T\) is assumed to be 8.75 K in L1544-
based runs and 10 K in L1521E-based runs at all radii. In “warm”
models $T$ decreases from 15 K at the core surface to $\sim$5 K at
$r = 0$ (the core center). The “cold” temperature profile is flat
with $T = 8.75$ K (L1544) or 10 K (L1521E) at $r/r_{\text{core}} > 0.2$ and
linearly decreases at smaller radii to 4 K at the center.

Two different approaches for estimating the chemical struc-
ture of a starless core are employed. In §3 the CO and CS
abundances are obtained through a detailed chemical modeling
of a core. We use the same chemical model as in Pavlyuchenkov
et al. (2006) and the density profile for the L1544 core. The
core is assumed to be isothermal ($T = 10$ K). We allow the
chemistry to evolve for 2 Myr and then use the resulting mol-
ecular abundances at various times to model the emergent line
profiles. Because of the assumption of uniform $T = 10$ K in
the chemical models, they are formally inconsistent with the
radiation transfer models. However, we calculated the same
chemical models with the above nonuniform temperature pro-
files and confirmed that the chemistry is not affected by these
relatively small $T$ variations.

In §4 we use the density profiles for L1544 and L1521E and
approximate the radial abundance profiles with predefined an-
alytic expressions. Only C^{16}O (1–0) (L1544) and C^{18}O (2–1)
(L1521E) transitions are considered in that section.

3. LINE INTENSITIES WITH COMPLETE CHEMISTRY

In the chemical modeling, we use two parameter sets. In the
first set we assume a low sticking probability of $S = 0.3$ (low-
depletion case). The model core is illuminated by interstellar
UV field with $G = 0.2$ and penetrated by cosmic rays with an
ionization rate of $\dot{\chi} = 3 \times 10^{-17}$ s$^{-1}$. This case is intended to
represent “chemically pristine” starless cores like CB17 (Pav-
luyuchenkov et al. 2006) or L1521E (Hirota et al. 2002). The
other parameter set is more appropriate for “chemically evolved”
cores (like L1544 or L1498). In this high-depletion case, we adopt $S = 1$, $\dot{\chi} = 3 \times 10^{-18}$ s$^{-1}$, and the same UV
field. Integrated intensity profiles for the considered transitions,
temperatures, and chemical parameters, convolved with the an-
tenna beams, corresponding to observations in Tafalla et al.
(2002) are shown in Figure 2. The CO abundance in the low-
depletion case is not significantly time-dependent, so results
after 2 Myr of evolution are shown. In the other plots, time
steps are chosen, which in the uni model produce results close
to observational data for L1544 (Tafalla et al. 2002).

In the high-depletion case, all molecules are heavily depleted
[$X(\text{max})/X(r = 0) \approx 1000$ for CO], and it is of no surprise that
the uni and cold models do not differ from each other. Both
optically thin (C^{16}O, C^{18}S) and optically thick lines are gen-
erated at intermediate radii and are not affected by the tem-
perature in the center. However, a warmer envelope (warm
model) results in a brighter emission, with the effect being
strongest for the optically thick CS (2–1) transition. This may
have a direct impact on the analysis of observational data. If
we use the adopted chemical model to interpret this emission
erroneously as arising from an isothermal object, we would
overestimate the core age by a factor of 1.5.

In the low-depletion case with uniform $T$ (Fig. 2, bottom left),
the central hole in the C^{16}O integrated intensity distribution
does not show up at all. Even though the depletion is still quite
significant [$X(\text{max})/X(r = 0) \approx 60$], the absolute molecule number
density peaks toward the center. Also, the uniform $T$ results in
a relatively high excitation temperature. However, the central
depression in $J_{\text{uni}}$ readily appears if we adopt a nonuniform tem-
perature profile, with $T$ decreasing toward the center or in-
creasing in the envelope. In the former case, the ringlike intensity profile
arises because of the weaker central emission; in the latter case
it is caused by the brighter envelope.

The temperature dependence of the C^{18}S and CS profiles is
much weaker, because these profiles are mainly formed in the
core envelope. As in the high-depletion case, the isothermal
model and the model with a cold center do not differ from
each other, while the emission in the warmer envelope model
is brighter than in other models.

Thus, whether or not the moderate depletion is visible as a
central hole in the C^{16}O (1–0) $J_{\text{uni}}$ distribution depends on the
adopted thermal structure. A chemically evolved core always
looks like a chemically evolved core, while a less evolved
core may appear to be chemically mature just because its tem-
perature structure is not taken into account properly.

One may suggest that the temperature structure can be inferred
from transition ratios, but in this case the interpretation is also
not straightforward. In Figure 3 (top panel) we show the $T$
profiles that would have been derived from the ratios of inte-
grated intensities C^{18}O (2–1)/C^{16}O (1–0), assuming that these

![Figure 1: Model temperature profiles.](image)

![Figure 2: Integrated intensity profiles for C^{16}O (1–0), C^{18}S (2–1), and CS (2–1) transitions, calculated with different assumptions on the core thermal structure.](image)
transitions are thermalized and optically thin. Cores with warmer envelopes do stand out on this plot, but the $^{13}\text{CO} (2–1)/(1–0)$ ratios do not allow one to distinguish between cold and uni models. In particular, the central temperature drop in cold models is only formally present.

Some hints can possibly be extracted from the analysis of higher transitions. On the bottom panel of Figure 3 we show the computed $^{13}\text{CO} (3–2)/^{13}\text{CO} (2–1)$ integrated intensity ratios. In the pristine models the $(2–1)$ line is saturated (i.e., its optical depth is greater than 1), while the $(3–2)$ transition is (moderately) optically thin. Because of that, higher density at the core center enhances the $(3–2)$ transition relative to the $(2–1)$ transition. Thus, the $^{13}\text{CO} (3–2)/^{13}\text{CO} (2–1)$ ratio is higher at small $r$ in uni and cold pristine models. In warm models the computed $^{13}\text{CO} (3–2)/^{13}\text{CO} (2–1)$ ratio is large because in the warm envelope collisions more effectively populate higher levels.

In uni and cold evolved models, both $^{13}\text{CO} (2–1)$ and $^{13}\text{CO} (3–2)$ lines are optically thin, and the $(2–1)$ transition does not show the saturation effects. At the same time, the $(3–2)$ line is weak due to low density at the periphery and significant depletion at the core center. Thus, the $^{13}\text{CO} (3–2)/^{13}\text{CO} (2–1)$ ratio is low.

Note the significant difference in $^{13}\text{CO} (3–2)/^{13}\text{CO} (2–1)$ ratios between the cold and uni pristine models in the low-depletion case (Fig. 3, bottom panel, solid and long-dashed curves). This implies that in surveys of “chemically young” starless cores, which are currently under way, it is still desirable to pay specific attention to higher transitions in order to assess the thermal structure of these objects.

Another issue, which has been brought to our attention by the referee, is related to isotope selective photodissociation, which preferentially destroys $^{13}\text{CO}$. This process is not taken into account in our model, but it may affect the ability of $^{13}\text{CO}$ to trace warm envelopes of starless cores. From this point of view, $^{13}\text{CO}$ transitions may be better envelope probes.

4. ANALYTICAL FITS

To study the sensitivity of derived molecular abundances on the temperature structure, we applied the technique, utilized by Tafalla and co-authors for L1544 (Tafalla et al. 2002) and L1521E (Tafalla & Santiago 2004); i.e., we approximate the molecular abundance profiles with $X(r) = X_0 \exp[-m(r)/n_d]$ and combine them with the three $T$ profiles described above.

![Fig. 4.—Results of the analytical fits for uni (left column), cold (middle column), and warm (right column) models. Top: Observed (connected filled circles) and computed $J_{15}$ profiles for the best-fit $X_r-n_r$ combination (triangles), as well as for bracketing combinations marked with crosses and boxes on bottom panels. Bottom: $X_r-n_r$ diagrams indicating the quality of agreement with observations. The darkest color corresponds to the best agreement. The best-fit positions for each $T$ profile are also marked with triangles.](image-url)
\[ L1544. – This is a well-known example of a starless core with some species, like CO, CS, C_2S, being strongly depleted. Thus, according to the results of § 3, we do not expect details of the thermal structure to have any significant effect on the interpretation of observational data.

With each of the three temperature profiles, we computed a series of RT models for different values of \( X_0 \) and \( n_d \) and compared the resulting distributions of the integrated intensity with the observed distributions. In Figure 4 we present the results of the comparison as a function of \( X_0 \) and \( n_d \). The best-fit positions in each plot are indicated by triangles. Combinations that bracket the scatter in observational points for the C^{18}O (1–0) transition are shown with crosses and boxes. The error bars in Figure 4 represent the extent of scatter in the observed integrated emission as a function of core radius (e.g., Tafalla et al. 2002, Fig. 3).

All the considered models do formally agree on the inferred values of \( X_0 \approx 10^{-7} \) and \( n_d = 8 \times 10^{4} \text{ cm}^{-3} \). However, the overall range of these parameters allowed by the scatter in \( J_{\text{int}} \) differs for various models. It appears that observations are marginally consistent with abundance profiles in which CO molecules are either totally depleted (crosses) or very moderately depleted (boxes). As the scatter mostly reflects the complex internal structure of the core, a difference between the bracketing \( X_0, n_d \) combinations may, to a certain degree, serve as a quantitative manifestation of a chemical inhomogeneity of the core.

\[ L1521E. – Unlike L1544, the starless core L1521E seems to be richer in gas phase CO, which is interpreted as a sign of its chemical youth (Hirota et al. 2002; Tafalla & Santiago 2004). Indeed, if we use the same exponential \( X(r) \) profile as before and a uniform \( T \) of 10 K, we find the best agreement with observations of the C^{18}O (2–1) transition for \( X_0 = 10^{-7} \) and \( n_d = 10^{4} \text{ cm}^{-3} \). This \( n_d \) value is greater than the central density \( n_0 \) of L1521E; i.e., the abundance distribution is nearly uniform. This is why in the uni model our conclusions are identical to those obtained by Tafalla & Santiago (2004).

However, the fit is different if we assume cold or warm \( T \) profiles. Neither centrally depleted nor constant abundances are able to reproduce the observed \( J_{\text{int}} \) in L1521E. In order to check other possible inferences for the L1521E chemical composition, we employed a different \( X(r) \) profile as well, assuming that \( X(r) = X_0 \exp (-r/r_{\text{CO}}) \). In Figure 5 we compare the fitting results for the C^{18}O (2–1) transition in the L1521E core to the observational data, given in Tafalla & Santiago (2004). A combination of uniform \( T \) and nearly constant abundance is able to fit the observations almost perfectly (solid line). To reproduce the observed \( J_{\text{int}}(r) \) in a core with a nonuniform \( T \) distribution, one need a centrally peaked abundance profile (Fig. 5). Again, as in a nonevolved chemical model, in models with lower central \( T \) we may get either flat or even depressed \( J_{\text{int}}(r) \) despite the fact that the underlying abundance is growing toward the core center. The fit quality is not high, especially, in the warm model, but this problem can be solved by using an abundance profile with central flattening (like \( \exp \left[-r^2/r_{\text{CO}}^2\right] \)) or by constructing an adequate chemical model.

5. CONCLUSIONS

The purpose of this study is to draw attention to the problems related to the thermal structure of prestellar cores. We demonstrate that realistic radial temperature profiles can produce quite different line intensities, even when the underlying chemical structure is the same. Effects of a nonuniform \( T \) distribution can mimic effects of chemical evolution and thus may hamper the interpretation of observations. This is especially true for “pristine” starless cores and is less important for chemically evolved cores. Finally, we would like to note that in this study we only consider transitions of CO and CS molecules. While ammonia observations represent a more conventional tool to measure the temperature in the interstellar medium, they are probably prone to the same kind of uncertainty.

This study is supported by the DFG grant “Molecular Cloud Chemistry” HE 1935/21-1. D. W. acknowledges support from RFPI grant 07-02-01031 and the Russian Science Support Foundation grant.

REFERENCES

Bergin, E. A., Maret, S., van der Tak, F. F. S., Alves, J., Carmody, S. M., & Lada, Ch. J. 2006, Apj, 645, 369
Crapsi, A., Caselli, P., Walmsley, M. C., & Tafalla, M. 2007, A&A, 470, 221
Di Francesco, J., Hogerheijde, M. R., Welch, W. J., & Bergin, E. A. 2002, AJ, 124, 2749
Galili, D., Walmsley, M., & Gonçalves, J. 2002, A&A, 394, 275
Goldsmith, P. F. 2001, Apj, 557, 736
Hirota, T., Ito, T., & Yamamoto, S. 2002, Apj, 565, 359
Lesaffre, P., Belloche, A., Chièze, J.-P., André, P. 2005, A&A, 443, 961
Pagani, L., Bacmann, A., Cabrit, S., & Vastel, C. 2007, A&A, 467, 179
Pavlyuchenkov, Ya., Wiebe, D., Launhardt, R., & Henning, Th. 2006, Apj, 645, 1212
Pavlyuchenkov, Ya. N., & Shustov, B. M. 2004, Astron. Rep., 48, 315
Pineda, J. L., & Bensch, F. 2007, A&A, 470, 615
Tafalla, M., Myers, P. C., Caselli, P., & Walmsley, C. M. 2004, A&A, 416, 191
Tafalla, M., Myers, P. C., Caselli, P., Walmsley, C. M., & Comito, C. 2002, Apj, 569, 815
Tafalla, M., & Santiago, J. 2004, A&A, 414, L53
van der Tak, F. F. S., Caselli, P., & Ceccarelli, C. 2005, A&A, 439, 195
Zucconi, A., Walmsley, C. M., & Galli, D. 2001, A&A, 376, 650