Chemical sensitivity to the ratio of the cosmic-ray ionization rates of He and H$_2$ in dense clouds

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ABSTRACT

Aims. To determine whether or not gas-phase chemical models with homogeneous and time-independent physical conditions explain the many observed molecular abundances in astrophysical sources, it is crucial to estimate the uncertainties in the calculated abundances and compare them with the observed abundances and their uncertainties. Non linear amplification of the error and bifurcation may limit the applicability of chemical models. Here we study such effects on dense cloud chemistry.

Methods. Using a previously studied approach to uncertainties based on the representation of rate coefficient errors as log normal distributions, we attempted to apply our approach using as input a variety of different elemental abundances from those studied previously. In this approach, all rate coefficients are varied randomly within their log normal (Gaussian) distribution, and the time-dependent chemistry calculated anew many times so as to obtain good statistics for the uncertainties in the calculated abundances.

Results. Starting with so-called “high-metal” elemental abundances, we found bimodal rather than Gaussian like distributions for the abundances of many species and traced these strange distributions to an extreme sensitivity of the system to changes in the ratio of the cosmic ray ionization rate $\zeta_{\text{He}}$ for He and that for molecular hydrogen $\zeta_{\text{H}_2}$. The sensitivity can be so extreme as to cause a region of bistability, which was subsequently found to be more extensive for another choice of elemental abundances. To the best of our knowledge, the bistable solutions found in this way are the same as found previously by other authors, but it is best to think of the ratio $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ as a control parameter perpendicular to the “standard” control parameter $\zeta/\zeta_\text{H}_2$.

Key words. astrochemistry – ISM: abundances – ISM: clouds – ISM: molecules

1. Introduction

The equations describing the chemical evolution of quiescent cores (also known as molecular clouds) are highly non-linear and can result in extreme sensitivity to the initial conditions or to some of the parameters. For example, the elemental ratio of carbon to oxygen is well known to be an important parameter for dense cloud chemistry (Terzivela & Herbst 1998). Another well-known consequence of this non-linearity is the presence of bistable regions for some ranges of model parameters (temperature, density, cosmic-ray ionization rate, rate coefficients and elemental abundances). The phenomenon of bistability is characterized by two stable solutions for chemical abundances at steady-state for the same set of parameters. Bistability in dark cloud conditions, i.e. at low temperature, was first discovered by Le Bourlot et al. (1993) and subsequently studied by a large number of authors (Le Bourlot et al. 1995; Shalabiea & Greenberg 1995; Lee et al. 1998; Pineau des Forêts & Roueff 2000; Charnley & Markwick 2003; Boger & Sternberg 2006). The two solutions of this bistability are characterized by a large difference in the C/O and O$_2$/O ratios and in the ionization fraction. For this last reason, they were named the High and Low Ionization Phases (HIP and LIP). Le Bourlot et al. (1995) showed that the region of bistability can be mapped out by variations in density, temperature, cosmic-ray ionization rate, and elemental depletions within specific ranges. Pineau des Forêts & Roueff (2000) later showed that even variations in rate coefficients can lead to the two solutions. Recently, Boger & Sternberg (2006) identified the chemical mechanisms of this phenomenon as a cycle involving the H$_3^+$, O$_2$, and S$^+$ species.

In two previous papers, we presented a Monte Carlo method to compute the theoretical error of abundances in chemical models due to uncertainties in rate coefficients and physical conditions (Wakelam et al. 2005, 2006). The errors in the rate coefficients are defined by a log-normal distribution. As a consequence, the resulting abundances at steady-state usually follow a Gaussian distribution (Wakelam et al. 2005). When applying this method, we found that the chemical abundances are very sensitive to the ratio between the ionization rates of helium $\zeta_{\text{He}}$ and molecular hydrogen $\zeta_{\text{H}_2}$ caused by cosmic-ray particles. Characterized by a variation of several orders of magnitude in some abundances for small variations of $\zeta_{\text{He}}/\zeta_{\text{H}_2}$, this sensitivity can lead to bimodal rather than Gaussian distributions for the abundances. For certain ranges of the parameters, it can even lead to bistability, with the ratio $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ appearing to be a control parameter.

This paper, which describes the sensitivity and some of its consequences, is organized as follows. In Sect. 2, we briefly present the chemical model and the uncertainty method used for
Table 1. High- and low-metal elemental abundances with respect to H$_2$.

<table>
<thead>
<tr>
<th>Element</th>
<th>High-Metal</th>
<th>Low-Metal (Intermediate)</th>
</tr>
</thead>
<tbody>
<tr>
<td>He</td>
<td>2.8 × 10$^{-4}$</td>
<td>2.8 × 10$^{-5}$</td>
</tr>
<tr>
<td>N</td>
<td>4.28 × 10$^{-3}$</td>
<td>4.28 × 10$^{-5}$</td>
</tr>
<tr>
<td>O</td>
<td>3.52 × 10$^{-4}$</td>
<td>3.52 × 10$^{-4}$</td>
</tr>
<tr>
<td>C$^+$</td>
<td>1.46 × 10$^{-3}$</td>
<td>1.46 × 10$^{-4}$</td>
</tr>
<tr>
<td>S$^+$</td>
<td>1.6 × 10$^{-5}$</td>
<td>1.6 × 10$^{-7}$</td>
</tr>
<tr>
<td>Si$^+$</td>
<td>1.6 × 10$^{-6}$</td>
<td>1.6 × 10$^{-8}$</td>
</tr>
<tr>
<td>Fe$^+$</td>
<td>6.0 × 10$^{-7}$</td>
<td>6.0 × 10$^{-9}$</td>
</tr>
<tr>
<td>Na$^+$</td>
<td>4.0 × 10$^{-2}$</td>
<td>4.0 × 10$^{-9}$</td>
</tr>
<tr>
<td>Mg$^+$</td>
<td>1.4 × 10$^{-2}$</td>
<td>1.4 × 10$^{-8}$</td>
</tr>
<tr>
<td>P$^+$</td>
<td>6.0 × 10$^{-4}$</td>
<td>6.0 × 10$^{-9}$</td>
</tr>
<tr>
<td>Cl$^+$</td>
<td>8.0 × 10$^{-7}$</td>
<td>8.0 × 10$^{-9}$</td>
</tr>
</tbody>
</table>

Note: The intermediate-metal case differs from the low-metal case by the S$^+$ abundance only, as is indicated in brackets in the table.

this analysis. In a somewhat unorthodox way, we have decided to present, in Sect. 3, the manner in which the sensitivity was found while studying the use of so-called “high-metal” elemental abundances. Section 4 shows the influence of the elemental abundances. Finally, the last section presents our conclusions.

2. Chemical model and method of uncertainty

We used a gas-phase time-dependent chemical model with the osu.2003$^1$ network reported by Smith et al. (2004) (4233 reactions, 421 species and 12 elements). The model computes the evolution of the species for a fixed temperature and density. The initial conditions are all atomic except for molecular hydrogen, while the chemical model and database are the same as used in Wakelam et al. (2006). For this work, we will present the results using three different sets of elemental abundances: the high-, low- and intermediate-metal cases. The high- and low-metal elemental abundances have been defined by Graedel et al. (1982) and are listed in Table 1. The intermediate case has the same abundances as the low-metal one except that the amount of the element sulfur is raised to 2 × 10$^{-6}$ compared with H$_2$ (also given in Table 1). The other parameters are the typical ones for quiescent cores: a kinetic temperature of 10 K, an H$_2$ density of 10$^4$ cm$^{-3}$, a visual extinction of 10 so that the photochemistry driven by the external UV photons does not occur, and a fixed cosmic-ray ionization rate $\zeta$ of 1.3 × 10$^{-17}$ s$^{-1}$.

The method used in this analysis to derive uncertainties in abundance is described in detail in Wakelam et al. (2005). The uncertainties in rate coefficients derive from the UMIST database when available, but for most of the reactions, including the ionization rates of H$_2$ and He, we have assumed a factor of 2 uncertainty. Note that the uncertainty in ionization rates is not an uncertainty in the parameter $\zeta$, but an uncertainty in the multiplicative factor that distinguishes the individual rates. Although ionization rates for H$_2$ and He have been used in a large number of models, the current rates used in terms of $\zeta$ derive back to the Ph. D. thesis of Black (1975), where the factors are 0.93 for the production of H$_2^+$ ($\zeta_{H_2}^{H_2^+}$) and 0.5 for the production of He$^+$ ($\zeta_{H_2}^{He^+}$), leading to a ratio $\zeta_{H_2}^{H_2^+}/\zeta_{H_2}^{He^+}$ of 0.54, if one neglects the minor channels for H$_2$. An older ratio of unity was used by Herbst & Klemperer (1973). New estimates, involving both direct ionization and secondary ionization by energetic electrons, are sorely needed (Black, private communication) based on the treatment of Dalgarno et al. (1999) for a mixture of atomic and molecular hydrogen and helium.

The uncertainty method consists of generating $N$ new sets of rate coefficients by replacing each coefficient $k_i$ by a random value consistent with its uncertainty factor $F_i$. We assume a normal distribution of log $k_i$ with a standard deviation $\sigma_i = \log F_i$. We run the model for each set $j$, which produces, for each species, $N$ values of the fractional abundances $X_j(t)$ at a time $t$. For this work, we ignore the uncertainty in physical conditions and consider the uncertainty in the rate coefficients only. A total of 2000 different runs was made for each set of parameters studied; this number is large enough to achieve statistical significance (see Wakelam et al. 2006).

3. The sensitivity to $\zeta_{H_2}/\zeta_{He}$ in the high-metal case

The composition of the theoretical error due to the uncertainties in rate coefficients was first applied to the case of high-metal elemental abundances. Figure 1 shows the results of the random variation of all the rate coefficients for the O$_2$ abundance as a function of time. As can be seen from this figure, the distribution of the abundance is spread into two peaks after 10$^7$ yr. The elemental abundances are for the high-metal case.

Fig. 1. Density of probability of the O$_2$ abundance as a function of time (left panel). The right panel shows the histogram of the abundance at 10$^7$ yr. The elemental abundances are for the high-metal case.

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$^1$ http://www.physics.ohio-state.edu/~eric/research_files/
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H₂, and the dissociative recombination of H₂⁺ to form H₂ and H.
The thick lines (hist1) represent histograms for all runs while the thin lines (hist2) represent those ending up with the higher O₂ abundance. For reactions not responsible for the bimodal distributions, there should be no difference between the pairs of histograms except that the total number of runs should be lower for the second one. For most of the reactions, we indeed found no difference, as for the reaction H₂⁺ + e– → H₂ + H in Fig. 3. In fact, there were only two exceptions: the direct ionization rate coefficients of He and H₂ by cosmic rays (ζ₃He and ζ₃H₂), and their marked effect can be seen in Fig. 3. The high abundance of O₂ tends to be formed with lower values of ζ₃He and higher values of ζ₃H₂, as can be seen most clearly in the lower panels of Fig. 3, in which the ratios of the two histograms are plotted. It is then convenient to consider the sensitivity of the abundance of O₂ and other species to the ratio ζ₃He/ζ₃H₂ rather than to their absolute values. To confirm their importance, we randomly varied only ζ₃He and ζ₃H₂, keeping the other rate coefficients constant, and we obtained the same bimodal distribution as with the fully random method at steady state. The dispersion of the abundances in this case is then almost as large as when we varied all the rate coefficients, as can be seen in the dotted lines in Fig. 2.

The sensitivity of the chemical abundances to the ratio ζ₃He/ζ₃H₂ can be explored in another way. In Fig. 4 (left panel), we show the steady-state abundance of O₂, computed by randomly varying all the rate coefficients as a function of ζ₃He/ζ₃H₂. For comparison, the abundance computed with the random variation of ζ₃He and ζ₃H₂ only is shown in the middle plot of the figure. The smaller vertical dispersion is caused by not varying the other rate coefficients. Finally and for more clarity, we depict, in the lower panel, the calculated O₂ abundance as a function of ζ₃He/ζ₃H₂ obtained by calculations in which the ionization rates are not varied randomly. Rather, the solid lines show values...
of the steady-state O$_2$ abundance computed for 5 values of $\xi_{He}/\xi_{H_2}$ (between 2 and 2) with $\xi_{H_2}$ varied to account for the range of the abscissa between 0.05 and 5. In the same panel, dotted lines show the O$_2$ abundances computed for 5 values of $\xi_{H_2}$ with $\xi_{He}$ varied to account for the range of the abscissa. The abundances computed by each method are similar especially at the inflection point, confirming that the choice of the ionization ratio as a parameter of the sensitivity is a reasonable one.

The abundance of O$_2$ decreases by two orders of magnitude (from $10^{-5}$ to $10^{-7}$) when $\xi_{He}/\xi_{H_2}$ goes from 0.3 to 0.8, and the inflection point of the curve occurs at the standard value of $\xi_{He}/\xi_{H_2}$ used in osu.2003 (0.5,0.93). In Fig. 5 we show some other examples of this sensitivity by using five different values of $\xi_{He}$, as in the solid lines of the right panel of Fig. 4. The abundance of atomic carbon has an almost equal sensitivity albeit opposite to that of O$_2$, while the H$_2$O abundance shows a similar profile to O$_2$ except for a local maximum at $\xi_{He}/\xi_{H_2}$ around 0.2. Although the electronic abundance increases towards higher $\xi_{He}/\xi_{H_2}$, it varies only a factor of 2.5 between $\xi_{He}/\xi_{H_2}$ = 0.1 and $\xi_{He}/\xi_{H_2}$ = 3. The abundance of the H$^+$ ion, which does not show any bimodal distribution, is linear with $\xi_{He}/\xi_{H_2}$. The dispersion of the abundances, for a given ratio $\xi_{He}/\xi_{H_2}$, due to the variation of $\xi_{He}$ by a factor of 2 around the standard value $\xi_{He}$, depends on the species: it is smaller for H$_2$O, O$_2$ and e$^-$ than for C and H$^+$, for instance, but remains anyway below 0.5 dex.

4. Sensitivity for different model parameters

4.1. Other choices of elemental abundances

The calculated sensitivity depends on the elemental abundances used for the modeling. In Fig. 6, we show the C, H$_2$O, e$^-$ and H$^+$ fractional abundances as a function of $\xi_{He}/\xi_{H_2}$ for the low-, intermediate-, and high-metal elemental abundances (see Sect. 2). The sensitivity plots, as well as most others later in the paper, are obtained by fixing the helium ionization rate at its standard value so that no dispersion is seen. The H$_2$ density remains at $10^4$ cm$^{-3}$ and the temperature at 10 K. Although the sensitivity exists for the three cases, the inflection point is shifted to higher values of $\xi_{He}/\xi_{H_2}$ for the low-metal case. For the intermediate case, on the other hand, the sensitivity seems to be stronger, leading to an abrupt jump in the abundance of C for $\xi_{He}/\xi_{H_2}$ = 1.1. A similar abrupt change can be seen for the electronic abundance. This jump is actually a manifestation of bistability (see Sect. 5). Note that for $\xi_{He}/\xi_{H_2}$ < 1.1, the abundances for the intermediate-metal case tend to be close to the low-metal ones whereas for $\xi_{He}/\xi_{H_2}$ > 1.1 they are similar to the high-metal ones.

When one runs the uncertainty method with a factor of two uncertainty in $\xi_{He}$ and $\xi_{H_2}$, the values of $\xi_{He}/\xi_{H_2}$ are spread approximatively between 0.13 and 2.2. The distribution of the abundances computed with this method thus depends on the sensitivity within this range of values. The distribution of the steady-state abundances of C computed by variation of all the rate coefficients are shown in Fig. 7 for the three sets of abundances. Because the inflection point for the low-metal case is shifted to $\xi_{He}/\xi_{H_2}$ = 2, the uncertainty method gives Gaussian-like distributions although slightly asymmetrical towards higher C abundances. This explains why we did not see the sensitivity effect in our previous work on rate coefficient uncertainties in dark clouds (see Wakelam et al. 2006). For high-metal elemental abundances, the point of inflection occurs close to the standard value of $\xi_{He}/\xi_{H_2}$ = 0.54 and the distribution for the C abundance is bimodal. When using the intermediate-metal abundances, the
uncertainty method gives two distinct peaks with no solutions between them. The peak at higher abundance is the smaller since the point of inflection occurs at values of $\xi_{\text{He}}/\xi_{\text{H}} = 1.1$ somewhat larger than the standard. The lack of any model runs with steady-state abundances in between the two distributions stems from the sharpness of the curve around the inflection point and is a consequence of the bistability.

In addition to its dependence on metallicity, the sensitivity to $\xi_{\text{He}}/\xi_{\text{H}}$ depends strongly on the C/O elemental abundance ratio. In Fig. 8, we show the steady-state abundance of atomic carbon as a function of $\xi_{\text{He}}/\xi_{\text{H}}$, computed for the three sets of elemental abundances, each with three different values of C/O – 0.2, 0.4 (the standard one), and 0.6 – obtained by maintaining a constant carbon abundance while varying the oxygen. The lower the C/O ratio is, the more the inflection point is shifted towards higher values of $\xi_{\text{He}}/\xi_{\text{H}}$. As an example, the inflection point for the high-metal case occurs around 0.2 for C/O = 0.6 whereas it rises to 1.5 for C/O = 0.2. In both cases, the inflection point is too far from the standard value of $\xi_{\text{He}}/\xi_{\text{H}}$ to obtain a bimodal distribution. For the intermediate-metal case, the C/O ratio of 0.6 results in shifting the inflection point to the standard value of $\xi_{\text{He}}/\xi_{\text{H}}$, so that the analogous panel to the one in Fig. 7 would show two equal areas for the two peaks. The dispersion of the carbon abundance caused by the variation of C/O is larger for values of $\xi_{\text{He}}/\xi_{\text{H}}$ smaller than the inflection point while the abundances eventually form a single plateau after the jump.

As discussed in Wakelam & Herbst (in preparation), the helium abundance used in the classic low- and high-metal elemental abundances (see Sect. 2) is quite high (0.28 compared with H$_2$) compared with the more modern value of 0.18 measured in the interstellar medium (see, for example, Baldwin et al. 1991).

Naturally, the sensitivity will change if a lower value of He is used. Figure 9 shows the distribution of the steady-state C abundance computed with a random variation of $\xi_{\text{He}}$ and $\xi_{\text{H}}$ and the high-metal case with the lower helium abundance. In this case, we do not obtain a clear bimodal distribution but merely an asymmetrical profile, because the inflection point is shifted towards higher values of $\xi_{\text{He}}/\xi_{\text{H}}$, as shown in Fig. 9, and the random method barely reaches the sensitive portion of the curve.

4.2. Chemical database

To test the influence of the chemical network utilized, we constructed the sensitivity curves at steady-state with two other databases – Srates and rate99 – using both low-metal and high-metal abundances. The rate99 network is the UMIST list of reactions (Le Teuff et al. 2000) prior to its current updating (see http://www.udfa.net) whereas Srates is the list of reactions used by Wakelam et al. (2004) to study sulphur chemistry in hot cores. This latter network contains only five elements (H, He, C, O and S) but, as discussed in Wakelam et al. (2004), the steady-state abundances computed with it are very close to the ones computed with the larger osu.2003 network at low temperature. The ionization rates for H$_2$ and He are the same in all three databases.

Figure 10 is a sensitivity plot of the C abundance for five different values of $\xi_{\text{He}}$ between 2$\xi_{\text{He}}$ and $\xi_{\text{He}}$/2. The sensitivity obtained with Srates is similar to that for osu.2003, which indicates that the absence of the other elements (Si, Fe, Na etc.) does not change the sensitivity drastically. On the contrary, the results with rate99 show a much different relation between $\xi_{\text{He}}/\xi_{\text{H}}$ and the atomic carbon abundance, especially for the high-metal case, where the rate99 curves show (a) a significantly lower value for the point of inflection of the sensitivity, (b) a wide dispersion due to the value of $\xi_{\text{He}}$ at low values of $\xi_{\text{He}}/\xi_{\text{H}}$, and (c) a much smaller jump in the abundance. We have shown in Wakelam et al. (2006) that rate99 and osu.2003 often have differences in rate coefficients larger than the estimated random errors for unstudied reactions of a factor of 2, whereas most of the rate coefficients in Srates are taken from the osu database. Since the three networks have the same $\xi_{\text{He}}$ and $\xi_{\text{H}}$, the shape of the sensitivity can depend only on the values of other rate coefficients.

5. Bistability

In Sect. 4, we saw that the variation of $\xi_{\text{He}}/\xi_{\text{H}}$ leads to a sharp change in the abundances of some species for the intermediate-metal case (see Fig. 6). This jump is indeed a manifestation of the same bistability discovered by Le Bourlot et al. (1993), as
can be determined by comparing the region of the bistability in parameter space. The parameter usually used to study this bistability is the ratio between the cosmic-ray ionization rate $\zeta$ and the density $n_H$ but other model parameters such as the temperature and rate coefficients have been used (Le Bourlot et al. 1995; Pineau des Forêts & Roueff 2000). What we have found here is that $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ is also a control parameter which explore another dimension of the bistability. If we imagine $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ to be an orthogonal axis to $\zeta_{\text{H}}$, the bistability of Le Bourlot et al. (1993) is the same as ours when $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ is at its standard value. In Fig. 11, we show an illustrative scheme of this bistability as controlled by the two parameters $n_{\text{H}_2}$ and $\zeta_{\text{H}_2}/\zeta_{\text{H}}$. To make the bistability appear, we constructed hysteresis curves for a function of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ for different $\text{H}_2$ densities. These curves, shown in Fig. 12 for atomic carbon, were produced in two steps. First, we ran the model to compute the steady-state abundance using one value of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$, then we used this solution as the initial condition and increased $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ to compute the next point, shown as the crosses in Figs. 12 and 13. In addition, the same principle but with decreasing $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ was utilized, with results shown as diamonds. In order to vary $\zeta_{\text{H}_2}/\zeta_{\text{H}}$, we fixed the helium ionization rate and changed that of hydrogen. When two solutions exist for a range of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ values, a region of bistability exists.

To explore the domains of bistability using the new control parameter, we constructed hysteresis curves using the three different elemental abundances and five densities between $10^6$ and $10^8$ cm$^{-3}$. All the results are shown in Fig. 12 for atomic carbon. Bistability exists for all the elemental abundances but the domain of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ where the two solutions coexist are small. To show the bistability more clearly, a magnification of this region is done for $n(\text{H}_2) = 10^7$ cm$^{-3}$ and the intermediate-metal case. In both the low- and high-metal cases, bistability exists only at very low density ($10^2$ cm$^{-3}$) whereas in the intermediate-metal case, the bistability exists between $10^2$ and $10^4$ cm$^{-3}$. The increase in density shifts and smooths the inflection point, making the bistability eventually disappear. Note that this is the first time that the bistability has been found with the high-metal case (Lee et al. 1998; Boger & Sternberg 2006), because it does not exist for the standard value of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$. We have confirmed that, for the high-metal case, using the standard value of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ leads to no bistability with $n_{\text{H}_2}$ as low as $10$ cm$^{-3}$. Using $\zeta_{\text{H}_2}/\zeta_{\text{H}} = 0.16$ however, bistability exists in a very small density range of 70–90 cm$^{-3}$.

A more focused look at the two control parameters is shown in Fig. 13, where we plot the hysteresis curves of the electron abundance as a function of each parameter for the case of intermediate-metal elemental abundances. The value of $\zeta$ is set at the standard value. For the variation of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ (right panel), we used five $\text{H}_2$ densities between $10^2$ and $10^6$ cm$^{-3}$, as in the previous figure, while for the hysteresis curve as a function of $n_{\text{H}_2}$ (left panel) the standard value of $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ (0.54) was used and the bistability found to exist in the density range 60–150 cm$^{-3}$. With the other control parameter and for $n_{\text{H}_2} = 10^4$ cm$^{-3}$, the bistablity exists for $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ between 0.5 and 0.7 and the domains of bistability using both parameters overlap. For higher densities, the bistable region using $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ shifts towards higher values and so does not exist anymore in the hysteresis curve versus $n_{\text{H}_2}$. These hysteresis curves can be combined into a plot of the type shown in Fig. 11.

6. Chemical reactions involved in the sensitivity

The sensitivity of the abundances to the ratio $\zeta_{\text{H}_2}/\zeta_{\text{H}}$ is related to the non-linearity of the system. A detailed mathematical analysis of this phenomenon is in progress (Massacrier et al., in preparation). It is possible, however, to understand the
difference, between the low and high ζHe/ζH2 phases qualitatively from a chemical point of view. For this purpose, we have considered the main reactions of production and destruction of O, C, CO, S and SO, which are the main reservoirs of carbon, oxygen and sulphur for the intermediate- and low-metal elemental compositions in which sulphur is relatively abundant, and/or are the species most influenced by ζHe and ζH2. Let us first consider low values of ζHe/ζH2. The ionization of H2 leads to the formation of OH through a well-known series of reactions (see, e.g. Boger & Sternberg 2006). The radical OH then reacts with atomic oxygen to form O2. The most efficient way of destroying O2 is to form SO by reaction with S. Since the elemental abundance of atomic sulphur is much less than that of oxygen, large amounts of O2 remain in the gas phase. Atomic carbon reacts with O2 to form CO, where it is stored since CO is not efficiently destroyed. For high ζHe/ζH2 phases, the chemical processes involved in destroying CO by He+ produces C* and O more efficiently. Carbon then exchanges its charge with S so large abundances of C remain in the gas phase to deplete molecular oxygen.

Since the bistability we see is an extreme manifestation of the sensitivity to ζHe/ζH2, the chemical processes involved in bistability may be related to those discussed above, although our control parameter is different from those previously studied. Pineau des Forêts et al. (1992) and Le Bourlot et al. (1993, 1995) identified the abundance ratio H+/H2 as being important for bistability, with one ion identified with the HIP and the other with the LIP solution. The abundance ratio H+/H2 depends on the abundance of electrons and as a consequence on the efficiency of the dissociative recombination of H2 with electrons, the rate of which was quite uncertain at the time. In our analysis however, the H2 ion shows little if any sensitivity to the control parameter, although the electron abundance does show a jump between the low and high ζHe/ζH2 phases, which can then be thought of as analogous in some sense to the LIP and HIP solutions. Recently, Boger & Sternberg (2006) made another detailed analysis of the chemical differences between the so-called HIP and LIP solutions, arguing that bistability is due to a loop including H2, O2 and S+. Although this explanation is related in some aspects to that given for the sensitivity reported here, one must remember that our control parameter raises some new points, given the relation of the ionization rate of He to the formation of the important ion C+, which is a precursor for many species, and can charge exchange to form abundant C. Moreover, S+, which is prominent in the explanation of Boger & Sternberg (2006), is not the dominant charge carrier in our calculations. Note that the osu.2003 database includes the recombination of the main ions with negative grains and this does not cancel the bistability, although such a cancellation was suggested by Boger & Sternberg (2006). It may well be that the H+/H2 abundance ratio, the H2 - O2 - S+ cycle, and the ζHe/ζH2 ratio comprise pieces of a larger and more complex jigsaw puzzle. Since bistability also appears to be a function of the reaction network utilized, secondary reactions may play a more important role than heretofore realized. This possibility has to be clarified through the building of a simplified reaction network that still shows the main characteristics of bistability.

7. Observational relevance

The sensitivity to ζHe/ζH2 depends highly on the model parameters and, for this reason, it is not obvious if such effect can be detected in the interstellar medium. In an ideal situation where the cloud characteristics (at least the elemental abundances and the age) are well known and correspond to a distinct chemical regime, one could use the observations to constrain the ratio ζHe/ζH2, assuming however that no local variation of this ratio is expected. In reality, the cloud characteristics are in general far from being well constrained, and an analysis of the uncertainties in the ionization rates of H2 and He is strongly required to understand the relevance of this sensitivity for the interstellar medium. Note that ζHe/ζH2 is not sensitive to the total intensity of the cosmic-ray flux.

If the intermediate-metal elemental abundances reflect the cloud characteristics and if the ionization rate of He and H2 by cosmic-rays can vary by at least a factor of 2 within a cloud, the prediction of bistability over a wide range of densities suggests that strong chemical heterogeneities in dense clouds can be caused by this effect. In the absence of bistability an extreme sensitivity is predicted at most densities for low-metal and high-metal abundances and may cause some observed heterogeneities in abundance. However, we have to consider the age of the cloud. The bifurcation occurs just after 107 yr, starting from our standard initial conditions, and is a maximum at steady-state (≈ 107 yr). As a consequence, the chemical heterogeneity can be caused by this sensitivity only in evolved clouds (≥ 107 - 108 yr).

The large gas-phase abundances of H2O and O2 molecules predicted by dense cloud chemical models is a long standing problem. Using high values of ζHe/ζH2, we obtain much smaller abundances for these species, closer to the upper limits in dense clouds (Bergin & Snell 2002; Pagani et al. 2003). Large uncertainties in the ζHe/ζH2 ratio could then solve the problem related to the modeling of water and molecular oxygen. One has, however, to consider the overall picture, and the better agreement with H2O and O2 should not worsen the agreement for the other species. A comparison between observations and modeling should be done in a systematic way, as done by Wakelam et al. (2006). This was not the purpose of the present study, but we can already note, as an example, that the observed abundances in L134N (North peak) of OH and C4H (7.5 × 10−8 and 10−9 respectively, Ohishi et al. 1992) are reproduced by the model when using the standard value of ζHe/ζH2, but are underestimated when using the higher ratios ζHe/ζH2 that reproduce H2O and O2 abundances.

Finally, the very low H2 density (100 cm−3), at which the bistability as a function of ζHe/ζH2 is restricted to the low and high-metal elemental abundances, is not relevant for clouds with the assumed visual extinction (A_v = 10) since it would require a size of 26 pc, which is much larger than typical cloud sizes (Snell 1981).

8. Conclusions

Using a Monte Carlo uncertainty method to compute the theoretical error of molecular abundances in chemical models, we found a strong sensitivity of the steady-state abundances to the ratio between the ionization rate coefficient of helium ζHe and that of molecular hydrogen ζH2 by cosmic rays. This sensitivity to the ratio has the consequence of changing the abundance of some species by several orders of magnitude for small variations of ζHe/ζH2. Lower values of ζHe/ζH2 are characterized by large abundances of smaller molecules such as O2, H2O, SO, etc. whereas high values of ζHe/ζH2 lead to large abundances of atoms. If the sensitivity is strong enough, it can result in the phenomenon of bistability, which we located at low densities for two sets of elemental abundances – the low- and high-metal cases, and over a wide range of densities for the case of the so-called intermediate-metal abundances, in which sulfur plays a particularly prominent role. The bistability has been determined to be
the same as studied earlier by a variety of authors with other control parameters; it is perhaps simplest to consider our control parameter $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ as orthogonal to one of the others – the ratio of the overall ionization rate parameter $\zeta$ to the density $n$ – so that the bistability can be imagined to exist in a region in three-dimensional space rather than in the customary two-dimensional hysteresis picture.

It is interesting to speculate that the sensitivity detected here via our uncertainty analysis is a necessary but not sufficient condition for bistability because only the most extreme sensitivity, in which the bimodal distributions detected have no solutions in between them, leads to bistability. Unfortunately, it is not guaranteed that our type of uncertainty analysis can always locate a bistable region since finding an extreme sensitivity may require large uncertainties in some parameters, such as the gas density, to take us from our standard parametric values to the bistable region.

Several questions arise when considering the dramatic effects of small $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ variations on the gas-phase chemistry. First, even if one assumes a given flux and energy distribution of the cosmic rays, the determination of $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ suffers some significant uncertainties, with the role of secondary electrons particularly important. Besides these difficulties, one might also have to consider variations of the ratio $\zeta_{\text{He}}/\zeta_{\text{H}_2}$ in space, such as between one cloud and another, or between the center and the edges of a same cloud, and in time. Changes in the ratio imply important variations in the spectral distribution of the cosmic rays (Glassgold & Langer 1973). Such variations might be produced by the attenuation of the cosmic rays inside a cloud or the proximity of supernovae in star forming clouds. Whether such variations are expected or not and how they could affect the chemistry should be considered in the future. Finally, a knowledge of the helium elemental abundance appears to be crucial since the sensitivity region depends on this parameter.

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