A model for star formation in cosmological simulations of galaxy formation

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Abstract / We present a new model to describe the star formation process in galaxies, which includes the description of the different gas phases – molecular, atomic, and ionized – together with its metal content. The model, which will be coupled to cosmological simulations of galaxy formation, will be used to investigate the relation between the star formation rate (SFR) and the formation of molecular hydrogen. The model follows the time evolution of the molecular, atomic and ionized phases in a gas cloud and estimates the amount of stellar mass formed, by solving a set of five coupled differential equations. As expected, we find a positive, strong correlation between the molecular fraction and the initial gas density, which manifests in a positive correlation between the initial gas density and the SFR of the cloud.

Keywords / galaxies: star formation — galaxies: evolution — methods: numerical

1. Introduction

The star formation rate (SFR) is one of the most fundamental properties of galaxies. In a cosmological context, the SFR is the result of a complex network of processes that act together during a galaxy’s lifetime, such as gas cooling, star formation, chemical enrichment and feedback from supernovae and galactic nuclei. Furthermore, the amount and properties of the gas from where stars form are strongly affected by mergers, interactions and mass accretion. Theoretical and observational studies show that the most important factor determining the star formation rate of a gas cloud is its density, although the details of this process are not yet fully understood.

Galaxy formation in the context of the standard cosmological paradigm, the A Cold Dark Matter model (ΛCDM) is an extremely complex, highly non-linear process that is free from any simplifying symmetries. Numerical simulations are the methods best suited to study this process on a physical basis, since they can follow the joint evolution of dark matter and baryons, naturally capturing processes such as mergers and mass accretion. However, large uncertainties still exist in the treatment of the baryonic evolution, as the physical processes affecting this component – star formation, various forms of feedback, chemical enrichment – act at unresolved scales and are only introduced via sub-grid physics. These models then need to tune a number of free input parameters that are usually not independent from each other, making the predictions of different models sometimes inconsistent (Scannapieco et al., 2012).

Because of its central role in galaxy formation, it is extremely important for simulations to properly describe the star formation process at the resolved scales, together with the associated feedback and chemical enrichment processes, and investigate in more detail the assumed sub-grid models and their dependencies with the properties of the interstellar gas. In this work, we present a new model of star formation, which describes the star formation rate of a gas cloud considering its chemical abundance and the relative abundance of the atomic, molecular and ionized gas phases. This new model is designed to be coupled to our cosmological simulation code GADGET3 (Springel et al., 2008) and the extensions of Scannapieco et al. (2005, 2006) and Poulihazan et al. (2018). This model has already been applied to chemical evolution models – simplified semi-analytic models of the formation of an isolated galaxy – to study the properties of the Milky Way and chemical gradients (Mollá et al., 2017, 2018).

This work is organized as follows. In Sec. 2 we describe our new model; in Sec. 3 we discuss the first results of the model and in Sec. 4 we present the conclusions.
2. The star formation model

The new star formation model is based on the work of Ascasibar et al. (2015) (see also Mollà et al. (2017), Murante et al. (2010), Millán-Irigoyen et al. (2020)), and is designed to follow the time evolution of various phases – molecular, atomic and ionized – in a gas cloud, also considering its metal content. In the context of our simulations, this model can be used to better describe the star formation process and investigate its efficiency as a function of the amount of gas in the molecular and atomic phases. The model is intended as a replacement for the traditional prescription for the star formation rate used in galaxy formation simulations [Katz 1992]:

$$\psi(t) = \epsilon_s \frac{\rho}{t_f},$$  \hspace{1cm} (1)

where $\epsilon_s$ is the efficiency of star formation and $t_f$ is the free fall time of the gas. In this way, $\psi(t)$ is only a function of the total gas density, $\rho$, and does not depend explicitly on the molecular or atomic fractions.

Our new model follows the coupled evolution of the different gas phases, ionized ($i$), atomic ($a$) and molecular ($m$) mass densities, together with the stars ($s$) and metals ($z$), solving a system of 5 coupled differential equations, namely:

$$\frac{di(t)}{dt} = - \frac{i(t)}{\tau_R} + (\eta_{ion} + R) \psi(t),$$ \hspace{1cm} (2)

$$\frac{da(t)}{dt} = - \frac{a(t)}{\tau_R} + \frac{i(t)}{\tau_R} \psi(t),$$ \hspace{1cm} (3)

$$\frac{dm(t)}{dt} = \frac{a(t)}{\tau_C} - (\eta_{diss} + 1) \psi(t),$$ \hspace{1cm} (4)

$$\frac{ds(t)}{dt} = (1 - R) \psi(t),$$ \hspace{1cm} (5)

$$\frac{dz(t)}{dt} = (Z_{SN} R - Z) \psi(t).$$ \hspace{1cm} (6)

These equations model the physical processes leading to a mass exchange between the different phases: photoionization of atomic hydrogen and supernovae – with assumed efficiencies per unit of star formation rate $\eta_{ion}$ and $R$, respectively; recombination of ionized hydrogen with electrons – characterized by a time $\tau_R$ –; dissociation of molecular hydrogen – with an efficiency per unit of star formation rate $\eta_{diss}$ –; and condensation of atomic hydrogen – catalyzed by dust grains – and regulated by the time parameter $\tau_C$. In our model, stars form from the molecular and metal components – proportionally to $Z = z/(i + a + m)$, the fractional amount of metals $z$, with a typical timescale $\tau_S$, and later feed metals and ionized gas to the interstellar medium (ISM) at their death as supernovae. This process is characterized by two parameters, $Z_{SN}$ for metal enrichment and $R$ for the ionized gas. The processes leading to mass exchange between the different phases are schematized in Fig. 1.

The model has several input parameters; however, most of them are constraints that are well constrained empirically or theoretically that we take from the literature. There are two parameters, $\tau_R$ and $\tau_C$, which do not depend on the properties of the gas, and this dependency is considered in our model following Osterbrock & Ferland (2006) and Millán-Irigoyen et al. (2020).

By solving our system of coupled equations, we obtain $m(t)$, the amount of gas in the molecular phase as a function of time and, based on the strong observational correlation between molecular hydrogen and star formation [Bigiel et al. 2008, Wong & Leo Blitz 2002], we assume a star formation rate of

$$\psi(t) = \frac{m(t)}{\tau_S},$$ \hspace{1cm} (7)

for the gas cloud, which replaces the one given in Eq. 1. In this way, we have a new star formation law that couples the star formation rate and the molecular fraction.

3. Results

As a first test of our model, we investigate the evolution of the different gas phases of a gas cloud in an idealized scenario. For this purpose, we integrated the system of equations for a single cloud of a given density, assuming initial parameters for the atomic, molecular and ionized gas fractions, and for the stellar and metal fractions. The system was evolved for a period of 1 Gyr and for three different values of the initial gas density, $n = 1, 10, 100$ cm$^{-3}$. We used, in particular, the following initial conditions: $i_f(t_0) = 0.001$, $a_f(t_0) = 0.998$, $m_f(t_0) = 0.001$, $s_f(t_0) = 0$ and $Z_f(t_0) = 0.0001$ where the subscript $f$ indicates the use of fractional quantities with respect to the total amount of gas, $g = i + a + m$. As mentioned above, $i$, $a$ and $m$ are respectively the mass densities of ionized, atomic and molecular gas. Note that $g \equiv n$, given that $n$ is the number density of protons, so it can be readily transformed into mass density units.

The results for the atomic, molecular and star fractions are shown in Fig. 2. We find a positive correlation between the initial density and the final fraction of molecular gas. For the highest density considered, the conversion of atomic to molecular gas is very efficient.
4. Conclusions

We presented a new model to describe the star formation rate of a gas cloud, which considers the evolution of the atomic, molecular and ionized fraction of gas, providing an estimation of the stellar mass expected to form as a function of time. The model is specifically designed to be coupled to hydrodynamical simulations of galaxy formation in a cosmological context.

We tested our model on simplified scenarios, by applying it to the evolution of clouds with different initial properties. In particular, we studied the dependency of the molecular, atomic and stellar fractions of a gas cloud, composed at the initial time almost entirely by atomic material, assuming three different values for the total initial density. Our results showed a strong dependence of the amount of stellar mass formed in the cloud and the initial density, which follows the trends found for the atomic and molecular phases.

Our model has already been applied in simulations of the formation of a Milky Way-mass galaxy – both in an idealized scenario and in cosmological context. Compared to the standard model (Eq. 1), we find that coupling the star formation efficiency to the amount of molecular gas produces galaxies with lower stellar masses and with retarded star formation rates (Lozano et al., in preparation).

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