

# Magritte: a new Multidimensional Accelerated General-purpose Radiative Transfer code

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## ABSTRACT

The In this paper we present the basis for a new multidimensional accelerated general-purpose radiative transfer code, called Magritte.

**Key words:** radiative transfer, astrochemistry, methods: numerical

## 1 INTRODUCTION

The Radiative Transfer problem is one of the oldest and

New C++ code based on 3D-PDR Bisbas et al. (2012) improved performance. separate modules for chemistry, radiative transfer and thermal balance.

The paper is organized as follows. In section 2 we present the new computational scheme and the different modules of Magritte. Section 3 describes the different benchmarks that were done to test the code and to compare its performance with 3D-PDR. In Section 4 we present the new applications that have become feasible through the nove Finally our conclusions are discussed in Section 5.

## 2 COMPUTATIONAL SCHEME

Although Magritte is based on 3D-PDR, the whole code has been rewritten from scratch.

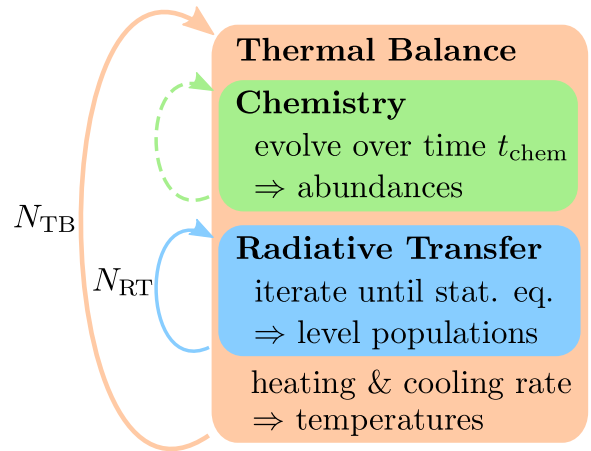
### 2.1 Data structures and memory layout

Magritte assumes an unstructured grid of cells

Data-oriented design

Much better memory scaling and improved overall performance.

Plot of memory use as function of the number of grid cells, Magritte vs. 3D-PDR.



**Figure 1.** Iteration schemes of the Magritte workflow, different boxes represent the different modules.

### 2.2 Modules

To facilitate the incorporation of Magritte in other codes, the code is structured into separate dedicated modules each handling a specific part of the calculation.

#### 2.2.1 Ray-tracing

On the length and time scales considered in Magritte simulations, light travels on straight lines. The first step in

The direction of a ray is determined by the HEALPix<sup>1</sup> discretization of the sphere Górski et al. (2005)

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<sup>1</sup> healpix.sourceforge.net

### 2.2.2 Chemistry

Magritte determines the relative abundances of a limited number of atomic and molecular species at each cell. This is done by solving the time-dependent chemistry of a self-contained network of formation and destruction reactions. The chemical network is a subset of the most recent UMIST data base of reaction rates [Woodall et al. \(2007\)](#), consisting of 320 reactions between 33 species (including electrons), and includes photoionization and photodissociation reactions in addition to the standard gas-phase chemistry.

### 2.2.3 Radiative transfer

Lines plus continuum.

The line data were taken from the Leiden Atomic and Molecular Database (LAMDA, [Schöier et al. 2005](#)).

Acceleration methods Ng-acceleration [Ng \(1974\)](#) and Rybicki-Hummer acceleration scheme [Rybicki & Hummer \(1991\)](#). The combination of both acceleration schemes yield a significant speed up in the convergence of the level populations. The approximated lambda operator for the Sobolev case is simply given by

$$\Lambda^* = 1 - \beta \quad (1)$$

Later versions will be able to treat multiple dust scattering

### 2.2.4 Thermal balance

Magritte can self-consistently determine the temperature assuming local thermal balance, i.e. equal heating and cooling rates for each cell. (Argument on time scales?)

## 2.3 Parallelization strategy

## 3 BENCHMARKS

## 4 APPLICATIONS

The modular character of Magritte allows it to be easily used in various astrophysical simulations.

## 5 CONCLUSIONS

We have presented Magritte: a new multidimensional accelerated general-purpose radiative transfer code.

Once all modules are finished and extensively tested, the source code for Magritte and its separate modules will be made freely available on [github.com/Magritte-code](https://github.com/Magritte-code).

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