

## Cosmic Ray Attenuation Module

### Introduction

In one-dimension, the code is now able to calculate the cosmic ray ionization rate in-situ from user defined input cosmic ray spectrum. The attenuation follows the prescription of Padovani+2009 using the 'continuous-slowing-down approximation (CSDA)' also known as the 'continuous energy loss regime'. This approximation will break down towards very high column densities ( $>100 \text{ g/cm}^2$ ) (see Padovani+2018 for details). The full details of the implementation are in Gaches+2019a (submitted). In one-dimension, the user defines cosmic ray spectra on either of the available surfaces. The code calculates the cosmic ray attenuation using the in-situ calculated molecular hydrogen from a user-given loss function.

### Outputs

The chemistry code includes two main cosmic-ray related output files:

- `zeta.txt`: This text file contains three columns: (x, CRIR/1.3E-17, NCOL) where CRIR is the cosmic ray ionization rate and NCOL is the column density of  $\text{H}_2$
- `OUTCRfin`: This file is the main output file for the cosmic ray spectrum. The first line is (NENE, Energy\_array), where NENE is the number of energy bins and the Energy\_array is NENE entries long. The following lines are (x, CR\_spectrum) where the CR\_spectrum is the spectrum at point x. If NENE=40 and the 1D domain has 1000 points, the file has 1001 rows and 41 columns.

### User-defined Inputs

The user can define several different inputs. The first of which is their own loss function. The public distribution contains two pre-defined loss function input files: `LE_loss_p.txt` and `LE_loss_p_2.txt`. They are the same loss function but at different energy resolutions. The user can define their own where with the format:

bins

E, LE/E<sup>16</sup>

These functions are used to interpolate over, so the number of bins can be (and should be) higher than NENE. `LE_loss_p2.txt` has 256 bins. If you define your own loss function, change the input file name in `spec_atten.F90` in the `CRSetup()` function.

The next inputs are all defined in the `params.dat` file. An example is given below for two sources:

```
-----|
Cosmic Ray Sources |
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2          |Number of CR Sources
40         |Number of CR Energy Bins
1          |Index for CR Transport - 1 = diffusive, 2 = rectilinear
SAC        |1150-tropic on Source (SAC)
Taurus_spec.dat |Input CR Spectrum as (E, F(E))
0.1         |Radial scaling in (pc)
-1         |CR Surface - 1 for +x direction, -1 for -x direction
150        |1150-tropic on Source (SAC)
IS_L_spectrum.dat |Input CR Spectrum as (E, F(E))
1          |CR Surface - 1 for +x direction, -1 for -x direction
```

In this case, `Taurus_spec.dat` is a spectrum for cosmic rays from a Taurus-like protocluster which is embedded in the center (-1) of the cloud. The other source is `IS_L_spectrum.dat` which is an external interstellar spectrum at the external surface (1). The index for CR transport is for sources only, where the physical transport scales as  $(r/\text{scale})^{\alpha}$ . The public distribution contains three example spectra:

- `IS_L_spectrum.dat` and `IS_H_spectrum.dat` are the Low and High Interstellar spectrum from Ilev+2015
- `Taurus_spec.dat` is an example protocluster Spectrum

User-defined spectrum must have the format: (e, je) where je is in units of particles/eV/s/cm<sup>2</sup>/sr. The file `IS_Spectra.pdf` shows a plot of the two different provided interstellar spectra.