

# User's Guide for Stage 2 Disk Evolution Code

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## 1 Overview

This document describes the structure, numerical scheme, and physical implementation of the Stage 2 version of the one-dimensional viscous disk evolution code. Stage 2 corresponds to the first fully time-dependent implementation of the disk energy equation coupled with irradiation heating including a finite time delay.

The code is designed for robustness, extensibility, and physical transparency, and is intended for studies of accretion and decretion disks in high-energy binary systems.

## 2 General Structure of the Code

The time integration is organized around three nested layers:

1. **Output-level time stepping** with a fixed interval  $\Delta t_{\text{out}}$ .
2. **Adaptive substepping** to reach the next output time.
3. **Local TRY/COMMIT evolution** for each substep.

The main program advances the system by repeatedly calling `evolve_try_to_target`, which ensures that the system evolves from time  $t_n$  to  $t_n + \Delta t_{\text{out}}$  using as many internal substeps as required for stability.

## 3 Main Modules

### 3.1 `mod_global`

This module stores global grid definitions and state arrays.

Key quantities include:

- Radial grid: `r(i)`, `nr`
- Surface density: `sigma_cur(i)`
- Viscosity: `nu_cur(i)`
- Thermal quantities: `Tmid_cur`, `H_cur`, `rho_cur`, `kappa_cur`, `tau_cur`
- Heating and cooling rates: `Qvis_cur`, `Qrad_cur`, `Qirr_cur`

The suffix `_cur` indicates the current committed state inside a substep.

### 3.2 evolve\_substep\_mod

This module controls the evolution over a single adaptive substep.

The central routine is:

`substep_try_and_commit()`

It performs:

1. Construction of the irradiation profile at the substep start.
2. A TRY evolution using `evolve_physics_one_substep_try`.
3. Acceptance or rejection of the step.
4. Commitment of the result to the `_cur` arrays.

No retry loop exists inside this routine; retries are handled at a higher level.

### 3.3 evolve\_try\_mod

This module contains the physics solver for a single substep.

The main routine is:

`evolve_physics_one_substep_try()`

It advances:

- Surface density  $\Sigma$
- Temperature  $T$
- Viscosity  $\nu$
- Disk structure and heating/cooling terms

for a given substep size  $\Delta t$ .

### 3.4 evolve\_to\_target (or evolve\_try\_to\_target)

This routine advances the system from  $t_n$  to  $t_n + \Delta t_{\text{out}}$  using adaptive substeps.

Algorithm:

1. Initialize `t_local = t_nd`.
2. While remaining time  $> 0$ :
  - Attempt a substep of size  $\Delta t_{\text{try}}$ .
  - On failure, reduce  $\Delta t$ .
  - On success, commit and advance `t_local`.
3. Ensure that output times are uniformly spaced.

## 4 Governing Equations

### 4.1 Surface Density Evolution

The disk surface density evolves according to the standard viscous diffusion equation:

$$\frac{\partial \Sigma}{\partial t} = \frac{3}{r} \frac{\partial}{\partial r} \left[ r^{1/2} \frac{\partial}{\partial r} \left( \nu \Sigma r^{1/2} \right) \right] + S(r, t), \quad (1)$$

where  $S(r, t)$  is an external source or sink term.

Time integration uses the  $\theta$ -method:

$$\Sigma^{n+1} = \Sigma^n + \Delta t \left[ (1 - \theta) F(\Sigma^n) + \theta F(\Sigma^{n+1}) \right]. \quad (2)$$

### 4.2 Energy Equation

At each radius, the midplane temperature evolves according to:

$$\Sigma c_V \frac{dT}{dt} = Q_{\text{vis}} + Q_{\text{irr}} - Q_{\text{rad}}. \quad (3)$$

An implicit Euler step is used:

$$\frac{T^{n+1} - T^n}{\Delta t} = \frac{Q_{\text{vis}}(T^{n+1}) + Q_{\text{irr}} - Q_{\text{rad}}(T^{n+1})}{\Sigma c_V}. \quad (4)$$

This nonlinear equation is solved locally using a damped Newton method with bracketing safeguards.

## 5 Convergence and Acceptance Criteria

Each substep is accepted if:

- No numerical failure occurs (NaN, divergence).
- $\Sigma > 0$ ,  $T > 0$ ,  $\nu > 0$  in active cells.

The Newton iteration itself enforces local convergence of the energy equation, so no global thermal balance test is required at this stage.

## 6 Irradiation Model

### 6.1 Irradiation Heating Rate

The irradiation heating rate is computed using Eq. (16) of Lee, Okazaki, and Hayasaki (2024):

$$Q_{\text{irr}}(\xi) = \frac{A_1 L_1}{2\pi r_{\text{in}}^2} \frac{1}{\xi} \left[ (1 + Q_{12}) \frac{dY}{d\xi} - \frac{\beta_1 + Q_{12}\beta_2}{\xi^2} \left( \frac{Y}{\xi} - \frac{1}{2} \frac{dY}{d\xi} \right) \right], \quad (5)$$

where  $\xi = r/r_{\text{in}}$  and  $Y = H/r$ .

At Stage 2:

- Shadowing is disabled.
- $Q_{\text{irr}}(r)$  is computed once per substep using the geometry at the substep start.

## 6.2 Time Delay (Scheme C)

Irradiation is powered by accretion luminosity with a finite delay.

Procedure:

1. Measure instantaneous accretion rate  $\dot{M}_{\text{in}}(t)$  at the inner boundary.
2. Store  $\dot{M}_{\text{in}}$  in a physical-time history buffer.
3. Define a delay time:

$$t_{\text{delay}} = f_{\text{delay}} \frac{r_{\text{in}}^2}{\nu(r_{\text{in}})}. \quad (6)$$

4. Compute delayed accretion rate  $\dot{M}_{\text{in}}(t - t_{\text{delay}})$  by linear interpolation.
5. Set irradiation luminosity:

$$L_{\text{irr}} = \eta_{\text{acc}} \dot{M}_{\text{in}}(t - t_{\text{delay}}) c^2, \quad (7)$$

optionally capped at the Eddington luminosity.

This scheme ensures causality and avoids instantaneous feedback.

## 7 Scope and Limitations of Stage 2

Stage 2 includes:

- Fully time-dependent energy equation.
- Adaptive substepping with fixed output times.
- Irradiation with time delay.

Not included at this stage:

- Shadowing effects.
- Implicit irradiation-thermal coupling.
- Stability branch tracking.
- Checkpointing of irradiation history buffers.

These are intended for later development stages.

## 8 Summary

The Stage 2 code provides a robust and physically consistent framework for studying irradiated viscous disks with delayed feedback. It forms a solid foundation for future extensions including shadowing, branch stability analysis, and global energy diagnostics.