PHYS 6260 Project Progress Report

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1 Introduction

For this project, I am attempting to model stellar structure for a variety of stars to see if I can reproduce the trends seen along the zero-age main sequence (ZAMS) from stellar data. The current setup of my model is:

- 1. Specify total mass and chemical composition (assumed to be constant throughout the star).
- 2. Calculate/approximate boundary conditions at the surface and the center based on semi-empirical scaling relations and astrophysical equations for radius r, pressure P, luminosity L, and temperature T.
- 3. Integrate the equations of stellar structure from the center outwards and from the surface inwards as a function of mass until the solutions meet.
- 4. Calculate residuals $\vec{\mathcal{R}}$ at this mass midpoint for each of the four variables above.
- 5. Determine how these residuals are affected by small perturbations to the approximated boundary conditions (total radius R, central pressure P_c , total luminosity L_s , central temperature T_c).
- 6. Use these partial results to calculate approximate perturbations to these boundary conditions that should reduce the residuals to zero.
- 7. Repeat until the residuals are small compared to the perturbed boundary conditions.

2 Project Status

The code for my project is located in a GitHub repository. Currently, all of my code is located in a single Jupyter Notebook to give me easier control over debugging; once the code can produce reasonable results, I will separate the code into python files so it is more modular.

2.1 Difficulties and Changes

I have come across a few major difficulties that led to noteworthy plan changes, each of which I will discuss in further detail below:

- 1. Correcting the ZAMS central boundary conditions
- 2. Properly implementing the perturbations calculations
- 3. Handling runtime errors during the RK4 integration

The primary issue was correcting the central boundary conditions used for ZAMS stars. Previously, I wanted to use m=r=L=0, but this led to issues with calculating the derivatives $\frac{\partial \vec{z}}{\partial m}$. I tried using small values for m and r(m) (atomic/microscopic scales), but this led to non-convergence issues because the derivatives and the integration range were so large. Instead, I'm assuming the core of mass $M_0=10^{-10}M$ is small enough to approximate constant conditions for T_c , P_c , etc., but large enough for the integration to converge by reducing the mass range. For a Sun-like star, this yields $M_0 \approx 2 \cdot 10^{20}$ kg and $r_c \approx 66$ km; though these are large on human scales, a core of this mass/radius is significantly smaller than the Moon, so I believe this approximation is reasonable. Here are my updated central boundary conditions:

•
$$M_0 = \lim_{m \to 0} m \implies \frac{M_0}{M} \approx 0$$

•
$$\rho_c = \lim_{m \to M_0} \rho(m) = 115 \frac{3}{4\pi} \frac{M}{R^3}$$

•
$$r_c = \lim_{m \to M_0} r(m) = \left(\frac{3}{4\pi} \frac{M_0}{\rho_c}\right)^{1/3}$$

•
$$P_c = \lim_{m \to M_0} P(m) = 7.701 \frac{GM^2}{R^4}$$

•
$$T_c = \lim_{m \to M_0} T(m) = \frac{2}{3} \frac{G}{k_B} \frac{Mm_u}{R}$$

•
$$L_c = \lim_{m \to M_0} L(m) = M_0 \epsilon \left(\vec{X}, \rho_c, T_c \right)$$

For the perturbations, I was applying them incorrectly. Previously, I had assumed they were multiplicative rather than additive. This has been resolved, and perturbations should now be calculated and applied correctly.

For the RK4 integration, I was getting runtime errors involving NaNs. Since RK4 integration involves calculations at the midpoints, there was a moment where the temperature gradient at the midpoint flipped from radiative to convective (or vice versa), which ended up yielding a negative temperature value in one of the \vec{k} 's This created NaNs in the calculations for ϵ . To address this, I am catching value errors for NaNs and inf during the RK4 integration; if these arise, then h is reduced and RK4 is rerun until the stepsize is small enough for the integration to yield sensible results. Since I expect this situation to arise infrequently, this seemed like a better approach than reducing the stepsize over the whole domain.

2.2 Outlook

Currently, most of the code for my project has been written and is functioning properly. I have:

- 1. Functions for the equation of state (P, ρ, T)
- 2. Functions for the remaining intermediary variables (ϵ , κ , etc.)
- 3. A function to calculate the stellar derivatives $\frac{\partial \vec{z}}{\partial m}$
- 4. A function to calculate/estimate the ZAMS boundary conditions
- 5. A function that performs one pass of integrating the equations of stellar structure
- 6. A function that calculates perturbations to the boundary conditions using the residuals $\vec{\mathcal{R}}$
- 7. A function that performs the full stellar integration (multiple integration passes and updating the boundary conditions)

There are a few minor errors I need to address, including how I am calculating the density from the equation of state. There may be additional runtime errors (likely involving the perturbations/residuals calculations, if they exist) I would need to address once this is fixed. I would also like to add H⁻ as an opacity source, as this is an important source of opacity at the surface for cooler stars that I should consider for the surface boundary conditions. I would like to have these done by the end of next week.

If I can get the ZAMS integration running properly in a reasonable amount of time, I would like to try expanding to giant stars and white dwarfs. This would require a few modifications:

- Giants and white dwarfs would need different sets of boundary conditions.
- For giant stars, I can no longer assume the chemical composition \vec{X} is constant throughout the star; rather, I would need to have a set of chemical compositions $\vec{X} = [\vec{X}_1, \vec{X}_2, \dots]$ that change based on which one produces the most energy given the conditions at that location.
- For white dwarfs, they do not produce energy via nuclear fusion, so ϵ is irrelevant. Rather, I think I would need to calculate the luminosity as $L(m) = 4\pi r^2(m)\sigma_{SB}T^4(m)$.