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COLLEGE OF ARTS & SCIENCES

SUPERCRITICAL CARBON DIOXIDE ROUND TURBULENT JETS:  
USING LARGE EDDY SIMULATIONS TO INVESTIGATE FUNDAMENTAL FLOW  
PHYSICS OF COMPLEX FLUIDS

By

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A Dissertation submitted to the  
Department of Mathematics  
in partial fulfillment of the  
requirements for the degree of  
Doctor of Philosophy

2023

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## **ACKNOWLEDGMENTS**

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## LIST OF ABBREVIATIONS

The following short list of abbreviations are used throughout this document.

AMR	Adaptive Mesh Refinement
CFD	Computational Fluid Dynamics
CO2	Carbon Dioxide
DNS	Direct Numerical Simulation
DOE	Department of Energy
ECP	Exascale Computing Project
EoS	equation of state
HMHW	half-mean half-width
HPC	high-performance computing
LES	Large Eddy Simulation
MISDC	Multi-Implicit Spectral Deferred Correction
MOL	Method of Lines
NIST	National Institute of Standards and Technology
NREL	National Renewable Energy Laboratory
ODE	Ordinary Differential Equation
PDE	Partial Differential Equation
PPM	piecewise parabolic method
PR EoS	Peng-Robinson equation of state
RANS	Reynolds-Averaged Navier-Stokes
rms	root mean square
SDC	Spectral Deferred Correction
sCO2	Supercritical Carbon Dioxide
SGS	Subgrid Scale
SMD	dynamic Smagorinsky
SRK EoS	Soave-Redlich-Kwong equation of state
TKE	Turbulent Kinetic Energy

# ABSTRACT

Supercritical carbon dioxide ( $\text{sCO}_2$ ) is of interest to a wide range of engineering problems, including carbon capture, utilization, and storage (CCUS) as well as advanced cycles for power generation. Non-ideal variations in physical properties of  $\text{sCO}_2$  impact the physics of these systems. In this study, we simulate turbulent  $\text{sCO}_2$  jets to gain a better understanding of these physics. Of particular interest is the impact of pseudo-boiling on supercritical flow dynamics. We use a second order finite volume method with adaptive mesh refinement as implemented in the ~~first-principles simulation~~ code PeleC to perform a Large Eddy Simulation (LES) of three turbulent jets of  $\text{sCO}_2$ . Additionally, we use the Soave-Redlich-Kwong equation of state to close the system and examine the impact of a cubic equation of state on the turbulent flow physics. We look at velocity and Reynolds stress profiles at different downstream locations for three cases in which the temperature of the jet and that of the ambient fluid differ. These results are then contrasted with established theory for ideal gas jets and similar numerical studies involving transcritical injection in order to capture the effects of widely varying thermal properties in the pseudocritical region.



# CHAPTER 1

## INTRODUCTION

Fluids are a vital part of everyday life - two of the most common ones being the air we breath and the water we drink. Less obvious are all the ways even these common fluids function behind the scenes in our day to day activities. Water, for example, serves an important role in various industrial settings, proving useful in areas ranging from thermal management to energy production; steam turbines alone accounted for 45% of electricity generation in the United States in 2021 [64]. Improvements to these types of systems impact technologies across a variety of fields and are thus an important area of research. One strategy toward this end, and the primary motivator of the work presented here, is the use of supercritical fluids as the working fluid of these systems.

In order to discuss supercritical fluids further, we must first go into more detail about general fluids. A *fluid* is a large collection of mutually interacting particles (e.g., molecules, atoms, etc.) in a state of constant and chaotic motion. This results in the continuous deformation of the substance under the effects of a shearing stress. The two categories of fluids that most people are familiar with are gases and liquids. *Liquids* are (mostly) incompressible and have definite volume for a set temperature and pressure. *Gases* are compressible and do not have definite volume. A simplified distinction between the two is that both gases and liquids will conform to the shape of whatever container they are in, but gases will further spread to fill all available space present.

A *supercritical fluid* is a fluid that is held above a critical temperature and pressure, at which point the distinction between a gas and liquid phase no longer exists [20, 65]. Supercritical fluids have qualities associated with both gases and liquids yet simultaneously have features that exclude them from fully being categorized as one or the other. For example, while they have viscosities akin to gases, they have solvent capabilities associated with liquids []. Similarly, while they have densities in line with liquids, they lack surface tension []. One benefit of this duality is that supercritical fluids can be fine tuned to be more gas-like or more liquid-like depending on the application at hand. This also results in ambiguity on how to actually classify them, with some sources considering them highly compressed gases [24], expanded liquids [1], or even as their own distinctly separate phase [6]. The distinction usually lies on the specifics of the regime and the application at hand.

This work is concerned with supercritical Carbon Dioxide (sCO<sub>2</sub>) in particular. sCO<sub>2</sub> has many beneficial features to a wide variety of industrial applications, as we will detail further in section 2. Many of the applications of interest to this work involve injection technologies that require a round turbulent jet configuration within the system. While much research has gone into sCO<sub>2</sub> flows, the current landscape is lacking in turbulence of this type with the context of understanding the underlying physics relating to the flow itself. For research that does involve other supercritical turbulent jets, the regimes explored in those works typically involve transcritical fluid injection within the regime of interest or explore regions outside the scope of this work.

The goal of this work is to explore the pseudo-boiling region of the pseudo-critical zone and analyze the influence of extreme thermodynamic fluctuations on turbulence statistics within the flow field. To that end, the rest of this chapter continues as follows: first, the mathematical framework for modeling compressible Newtonian fluids is provided to form the basis of the modeling done in this dissertation. Further consideration is then given to turbulence modeling and the numerical methods developed for studying turbulence to provide insight into the quantities of interest analyzed within this dissertation and the choices of numerical methods used herein. Important applications of supercritical carbon dioxide in particular are provided to motivate the problem presented in this dissertation. Existing numerical studies on supercritical fluids are reviewed to demonstrate how this dissertation fits into the current landscape of research and to emphasize the contribution the results of this work make to the field. This chapter concludes with an outline of the dissertation, the goals of the dissertation, and the main contributions made through this work.

## 1.1 Mathematics of Fluid Flow

Scale is one of the key factors to consider when developing a mathematical description of a fluid system. For example, consider modeling flow past a satellite in the exosphere vs. the flow past a turtle in the ocean; these two mediums have vastly different characteristics and would thus require different modeling techniques. Scale is also an important concept when it comes to turbulence in particular so we will begin that discussion here with our choice in perspective for the mathematical framework of our system of interest.

From a kinetics perspective, particle motion within a fluid can be broken up into two phases: particle interaction and free flight. Average time spent in free flight,  $\langle t_f \rangle$ , is typically much greater than collision time for a given interaction,  $t_c$ . The average length traveled between collisions

is known as the mean free path,  $\ell$ . Since free flight time dominates particle interaction time, this phase determines the length scale of the kinetic description of motion. In addition to this inherent physical scale, there is also a scale associated with the resolution of the problem itself,  $L$ . These two scales are important, as the mathematical description of your model depends on how these two scales compare to one another. This comparison is related through the non-dimensional Knudsen number:

$$Kn = \frac{\ell}{L} \quad (1.1)$$

Flows with large Knudsen number ( $Kn \gg 10$ ) require modeling from the kinetic or microscopic perspective as particle interactions become sparse enough compared to the scope of the problem to require a statistical mechanics framework. On the other hand, small Knudsen number flows ( $Kn \ll 0.01$ ) have a problem scale that far exceeds the particle-level interactions present, giving way to an average overall motion within the fluid. This scale dichotomy is demonstrated with the graphic in Figure 1.1. For small Knudsen flows, a continuum description of the fluid is appropriate for capturing this macroscopic behavior.

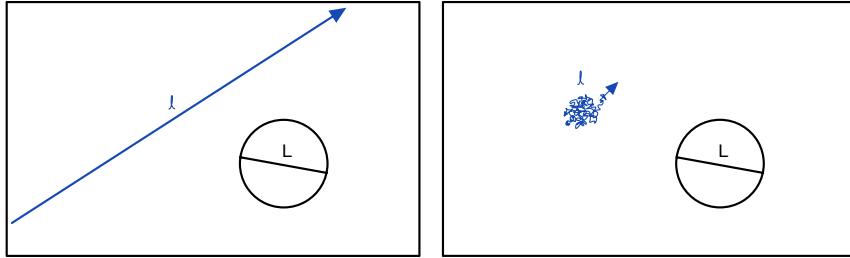


Figure 1.1: Characteristic length scale of problem,  $L$ , compared to mean free path of particles,  $\ell$ , for a flow with large Knudsen number (left) vs. small Knudsen number (right)

This work falls within the small Knudsen regime, so we will be working with the continuum description of fluids. In this section we will discuss the Navier-Stokes Equations that arise from this modeling technique and how we account for the supercritical nature of the flow through our choice of equation of state.

### 1.1.1 Continuum Description of Fluids

The continuum hypothesis assumes that the fluid has no fine structures and that it is perfectly continuous, i.e., the properties of a small subdivision are the same as other subdivisions. This allows for the approximation of physical quantities at the infinitesimal limit [28].

For example, consider a fluid with arbitrary volume  $V$  as depicted in Figure 1.2.

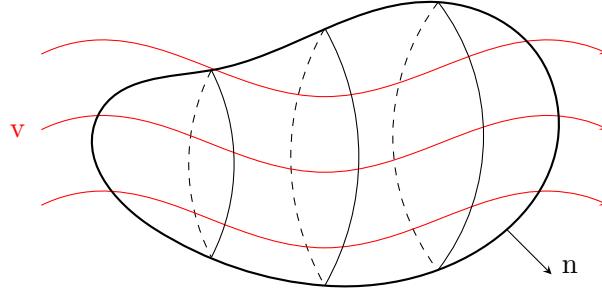


Figure 1.2: A fluid of arbitrary volume  $V$  bounded by surface  $S$  with velocity  $\mathbf{v}(\mathbf{x}, t)$ . A differential volume and surface area is given by  $dv$  and  $ds$ , respectively.  $\mathbf{n}$  is the outward-pointing unit normal vector to the surface  $S$ .

For a fluid with density  $\rho(\mathbf{x}, t)$ , mass within a small representative volume can be described with

$$\rho dv$$

Total mass in the arbitrary volume is then given by

$$\iiint_V \rho dv$$

The rate of change of mass through the volume is now

$$\begin{aligned} & \frac{d}{dt} \iiint_V \rho dv \\ &= \iiint_V \frac{\partial \rho}{\partial t} dv \end{aligned} \quad (1.2)$$

Simultaneously, overall change in mass throughout the volume can be described by the net mass flux through the surface  $S$ . Volumetric flow through a small portion of the bounding surface is given by

$$\mathbf{v} \cdot \mathbf{n} ds$$

Total mass flux through the entire surface is then

$$\iint_S \rho \mathbf{v} \cdot \mathbf{n} ds \quad (1.3)$$

Applying the divergence theorem to Eq. 1.3 yields the following volume integral

$$\iiint_V \nabla \cdot (\rho \mathbf{v}) dv \quad (1.4)$$

Assuming there is no additional source generating or leaking mass within the control volume, we can relate Eqs. 1.2 to 1.4 :

$$\begin{aligned} \iiint_V \frac{\partial \rho}{\partial t} dv &= - \iiint_V \nabla \cdot (\rho \mathbf{v}) dv \\ \iiint_V \frac{\partial \rho}{\partial t} dv + \iiint_V \nabla \cdot (\rho \mathbf{v}) dv &= 0 \\ \iiint_V \left( \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) \right) dv &= 0 \end{aligned} \quad (1.5)$$

Note the inclusion of the negative sign for the right side of the initial equality; in the surface integral formulation, the outward facing normal describes flux out of the volume, thus yielding a decrease in mass within the volume. Since Eq. 1.5 holds for any arbitrary volume  $V$ , the integrand must be identically equal to zero:

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) = 0 \quad (1.6)$$

Through the continuum hypothesis and conservation of mass, we have now arrived at the continuity equation in Eq. 1.6. This specific process demonstrates an even more fundamental relationship known as a *conservation law*. More generally, for some integrated property  $\phi$ , the rate of change of  $\phi$  within a control volume must be equal to the amount of  $\phi$  lost or gained through the boundaries of the control volume plus what is created or consumed by any sinks or sources,  $s$ , within the volume (sinks having positive orientation to match the positive orientation of the outward-facing normal  $\mathbf{n}$ ) [].

$$\frac{\partial \phi}{\partial t} + \nabla \cdot (\phi \mathbf{v}) + s = 0 \quad (1.7)$$

In addition to this concept applying to conservation of mass, as was seen in this section, the idea outlined by Eq. 1.7 applies to conservation of momentum and energy within the fluid. Together, these expressions combine to form the basis of the Navier-Stokes Equations, as will be seen in more detail in chapter 2. The important takeaway from this section is that with the continuum hypothesis and fundamental laws of physics, one can adequately capture macroscopic flow behavior for the types of flows we are interested in within this work.

### 1.1.2 Equation of State

Conservation of mass, momentum, and energy gives us five equations to describe our fluid system. For compressible flows, this is not enough information to solve for all the unknowns within the system of coupled partial differential equations. A sixth equation, known as the Equation

of State (EoS), must be chosen in order to close the system. The EoS relates three of the six unknowns: pressure, temperature, and density. Here we will briefly discuss some EoS options and their distinguishing characteristics in order to motivate the choice made for this work.

The simplest option available is the ideal gas EoS, which comes from the ideal gas law. This EoS relates density, pressure, and temperature in the following manner:

$$p = \frac{RT}{V_m} \quad (1.8)$$

where  $p$  is pressure,  $R$  is the universal gas constant,  $T$  is temperature, and  $V_m = \frac{V}{n}$  is the molar volume of the fluid (it is common to express density in terms of molar volume for sake of simplicity in writing the EoS with  $V$  being volume and  $n$  being the number of moles). The ideal gas EoS is fairly accurate for liquids and gases at moderate temperatures and low pressures. It fails at low temperatures and high pressures, especially near the transition region from gas to liquid. The inaccuracy noted in this region means this EoS would not be suitable for the area of interest within this study.

Cubic EoS generally provide more accuracy than the ideal gas EoS. The first cubic EoS was developed by van der Waal in 1873, modifying the ideal gas EoS to take into consideration the finite size of molecules and interactions between molecules (the ideal gas EoS only accounts for interactions with the container and treats molecules as point particles). Other cubic EoS can be thought of as modifications from this base form:

$$p = \frac{RT}{V_m - b} - \frac{a}{V_m^2} \quad (1.9)$$

where  $a$  and  $b$  are constants related to the pressure and temperature at the critical point,  $p_c$  and  $T_c$  respectively:

$$a = \frac{27(RT_c)^2}{64p_c}, \quad b = \frac{RT_c}{8p_c}$$

One of the main benefits of using a cubic EoS is that they can have comparable and sometimes even better accuracy compared to their higher order counterparts, thus reducing computational costs. However, it is important to take into consideration the regime of interest in addition to the fluid of interest when choosing an EoS, as each one has its own pros and cons. For example, molecule polarity and density are two factors that can have a high impact in selection between the Soave-Redlich-Kwong Equation of State (SRK EoS) and Peng-Robinson Equation of State (PR EoS) alone [23].

This work uses the SRK EoS as will be more thoroughly introduced in chapter 2. It has been shown that this equation of state is fairly accurate for the parameter regime under careful consideration in this work []. This accuracy is detailed further in chapter 3 through comparisons with data from National Institute of Standards and Technology (NIST). Overall, when adequately considered, the EoS is the key avenue to incorporating specific fluid properties into the mathematical model.

## 1.2 Turbulence

In addition to categorizing by fluid type, fluid flow can be categorized based on flow characteristics. The main two classifications of note are laminar flow and turbulent flow.

Laminar flow is denoted by fluid particles having well-defined parallel trajectories of motion, or streamlines. Streamlines do not cross, meaning adjacent layers within the fluid flow by one another with little to no mixing. From a more generalized perspective, the flow appears to be smooth. In contrast to this, turbulent flow is characterized by its unpredictable and chaotic trajectories. Streamlines do cross resulting in swirls and eddies of varying length scales which induce mixing. Turbulent flow can be qualitatively described as being rough due to this high degree of fluctuation within the velocity and pressure fields present. This generalized description is depicted in Figure 1.3.

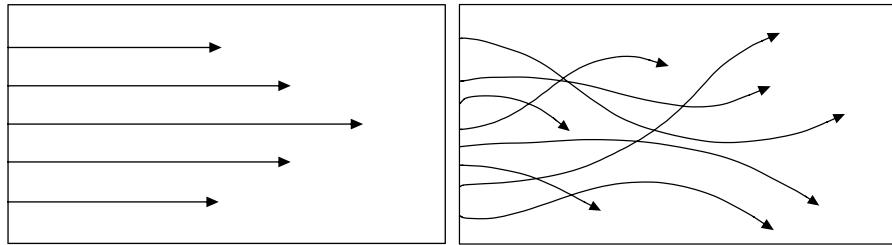


Figure 1.3: Example of streamlines in laminar (left) vs. turbulent (right) flow.

The Reynolds number is a dimensionless value that can be used to distinguish laminar flow from turbulent flow. It is defined as follows:

$$\text{Re} = \frac{\rho u L}{\mu} \quad (1.10)$$

where  $\rho$  and  $\mu$  are the density and dynamic viscosity of the fluid, respectively,  $u$  is the flow velocity, and  $L$  is a characteristic length scale associated with the given flow scenario (e.g., pipe diameter).

As is demonstrated by the ratio in Equation 1.10, the Reynolds number measures the relative effects of inertial forces compared to viscous forces within a given flow scenario. A small Reynolds number signifies the dominance of viscous forces; fluid parcels moving in tandem want to “stick together,” resulting in the sheared flow and parallel trajectories seen in laminar flow. Turbulence is then characterized by a large Reynolds number, where inertial forces take precedence. Here, deviations within the laminar flow field result in lateral mixing between shear layers. This creates eddies and random trajectories that result in the chaotic motion of turbulent flow.

This work focuses on the turbulent round jet and its associated dynamics in the context of supercritical fluids. The remainder of this section details a brief overview of important turbulence concepts and numerical methods developed for studying turbulence in order to motivate the modeling and numerical choices made within this work.

### 1.2.1 Mathematics of Turbulence

Turbulence has been recognized and investigated by scholars in some way, shape, or form for thousands of years; Lucretius being one of the earliest, describing eddy motion in his *De rerum natura* [7]. In that time, a comprehensive theory of turbulence has still not been reached. There are, however, many widely accepted hypotheses and models that at least in part explain certain aspects of turbulent flows.

Turbulent motion contains a wide range of scales. Richardson formalized this concept by describing turbulence as a compositions of eddies, each with their own characteristic size and velocity [1]. From this perspective, large eddies are said to be unstable and break up into smaller eddies, transferring energy down the line to smaller and smaller eddies. This energy cascade continues down to the molecular level where it is dissipated through molecular viscosity. Kolmogorov further developed this theory, noting that for sufficiently high Reynolds flow, the statistics of these smallest-scale motions are uniquely determined by the kinematic viscosity of the fluid,  $\nu$ , and the dissipation rate,  $\varepsilon$  [1]. These smallest eddy length, velocity, and time scales at which dissipation occurs, known as the Kolmogorov scales, are defined as follows:

$$\begin{aligned}\eta &\equiv (\nu^3/\varepsilon)^{1/4}, \\ u_\eta &\equiv (\nu\varepsilon)^{1/4}, \\ \tau_\eta &\equiv (\nu/\varepsilon)^{1/2}\end{aligned}\tag{1.11}$$

While the continuum description introduced in the previous section is still valid for turbulent flows, this vast scale separation and the random nature of fluctuations across these scales, including

a dependence on problem-specific flow configurations and boundary conditions, make universal solutions to these types of problems unattainable. To that end, numerical simulations are invaluable in studying turbulence phenomena.

### 1.2.2 Numerical Approaches to Turbulence Modeling

A vast array of numerical approaches have been developed to tackle the modeling challenges inherent to turbulence, with further development of these and new techniques remaining an active field of research today. We will briefly look at three families of numerical simulation methods in order to highlight the challenges associated with the large scale disparity included in turbulent flows. Methods of note include use of the Reynolds-Averaged Navier-Stokes (RANS) equations, direct numerical simulation (DNS), and large eddy simulation (LES), with scale coverage of each outlined in Figure 1.4.

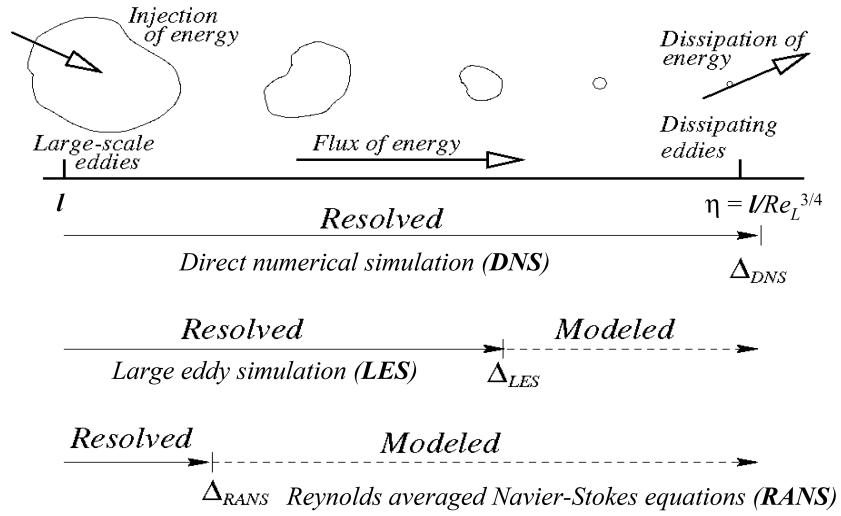


Figure 1.4: Depiction of the energy cascade present in turbulence along associated scales. Resolved vs. Modeled scales are mapped for DNS, LES, and RANS [].

At one extreme, we have DNS. The aim of DNS to resolve all scales within the turbulent flow, from the largest eddies down to the Kolmogorov scale. To do this, all components of the Navier-Stokes equations are discretized and numerically advanced. Numerical methods may include high-order finite differencing or pseudo-spectral methods depending on the type of turbulence, flow configuration, and boundary conditions at hand. While DNS can capture physics across all scales of the flow with high accuracy, this comes at a steep computational cost. Grid spacing and time

stepping are highly dependent upon the Reynolds number (approximately as  $Re^3$  [51]). Thus DNS very quickly becomes limited to lower Reynolds flows, even with high computing resources.

On the other end of the spectrum, we have RANS modeling. The RANS equations make use of the Reynolds decomposition of the flow, in which a random velocity field  $\mathbf{u}$  can be split into an ensemble average,  $\langle \mathbf{u} \rangle$ , and a fluctuating component,  $\mathbf{u}'$ , such that:

$$\begin{aligned}\mathbf{u}'(\mathbf{x}, t) &\equiv \mathbf{u}(\mathbf{x}, t) - \langle \mathbf{u}(\mathbf{x}, t) \rangle, \\ \langle \mathbf{u}(\mathbf{x}, t) \rangle &= \lim_{N \rightarrow \infty} \frac{1}{N} \sum_{i=1}^N \mathbf{u}_i(\mathbf{x}, t)\end{aligned}\tag{1.12}$$

where  $i$  is the  $i$ -th realization of a flow of identical conditions with the number of realizations  $N$  tending toward infinity. In practice, for numerical models with the RANS equations, this is implemented via time averaging:

$$\langle \mathbf{u}(\mathbf{x}, t) \rangle \approx \overline{\mathbf{u}(\mathbf{x})} \equiv \frac{1}{T} \int_{t_0}^{t_0+T} \mathbf{u}(\mathbf{x}, s) ds$$

where  $T$  is taken to be much larger than the time scale of the fluctuating components. By substituting  $\mathbf{u}(\mathbf{x}, t) = \overline{\mathbf{u}(\mathbf{x})} + \mathbf{u}'(\mathbf{x}, t)$  particular substitution into the Navier-Stokes equations, one can get a statistical description of the average flow field for the turbulent system (note - the more general decomposition given in equation 1.12 can be used in the substitution as well to take into consideration a moving average within the flow field. For many of the applications in which RANS is used, it is enough to look at the steady state given by the time averaged system). Further equations are then needed to model the additional stresses that arise from the nonlinear interactions between components of the mean flow field. Overall, simulations of the RANS equations resolve the larger scale motions associated with the average motion of the field in a statistical description of the system while the influence of fluctuating fields are incorporated via modeling. While this reduces the computational cost of the numerics, more closure issues arise from the modeling requirements introduced, which present a whole new set of choices and assumptions that are highly problem specific.

Finally, we have LES, which falls between the two extremes of the methodology spectrum presented thus far. The strategy with LES is to only resolve the energy containing scales in the system and model the influence of the smaller scales, as demonstrated in Figure 1.5. This allows for some relaxation in computational cost compared to DNS without introducing the same degree of modeling assumptions required for RANS. While still fairly computationally expensive overall, LES

can resolve more of the instantaneous flow field and capture certain physics that inherently depend on the small scale dynamics; quantities of this nature are lost or become highly model dependent in RANS which introduces more error to the simulation.

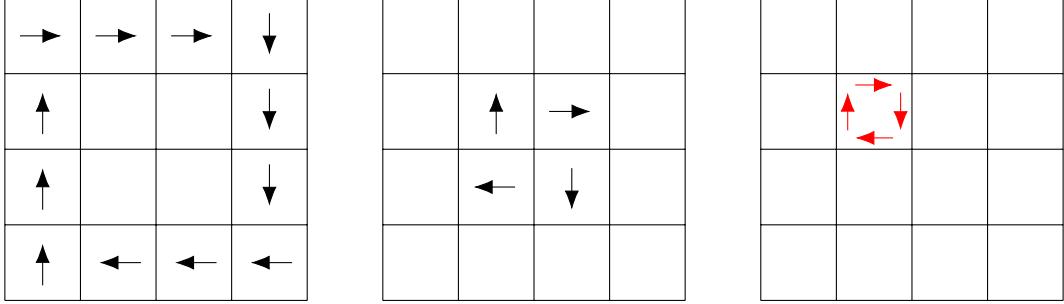


Figure 1.5: LES resolves large scale eddies (black) and models the effects of fine scale eddies (red) that are unresolved due to mesh size.

The procedure for LES involves decomposing random velocity field  $\mathbf{u}$  into filtered components,  $\bar{\mathbf{u}}$ , and residual components,  $\mathbf{u}''$ , such that:

$$\begin{aligned}\mathbf{u}''(\mathbf{x}, t) &\equiv \mathbf{u}(\mathbf{x}, t) - \overline{\mathbf{u}(\mathbf{x}, t)}, \\ \overline{\mathbf{u}(\mathbf{x}, t)} &= \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \mathbf{u}(\mathbf{r}, s) G(\mathbf{x} - \mathbf{r}, t - s) d\mathbf{r} ds\end{aligned}\quad (1.13)$$

$G$  is the convolution kernel of associated filter spatial and/or temporal cutoff  $\Delta$  and  $\tau_c$ , respectively, satisfying  $\int_{-\infty}^{\infty} \int_{-\infty}^{\infty} G(\mathbf{r}, s) d\mathbf{r} ds = 1$ . In practice, common kernels are only spatially dependent. Applying this filtering procedure to the Navier-Stokes equations yields an additional term known as the residual stress tensor that must be modeled. The filtered Navier-Stokes equations can then be discretized to numerically simulate the large eddy behavior of the system.

This general LES procedure may seem similar to what was described for the RANS model, especially in terms of the decomposition of the flow field, but there are key differences to note. Conceptually, the result of a numerical simulation using RANS describes an average behavior of turbulent flow over many instances or experiments. The output of these simulations is not a snapshot of one instance of the given flow field but rather an approximation to how the flow behaves in a statistical sense. LES does approximate a single realization of the flow field, just only including the low frequency modes of the system which contain the majority of the energy. Additionally, the components of the Reynolds decomposition obey the following rules:

$$\bar{\mathbf{u}} = \bar{\mathbf{u}}, \quad \bar{\mathbf{u}}' = 0$$

These rules do not hold for  $\bar{\mathbf{u}}$  and  $\mathbf{u}''$ :

$$\bar{\bar{\mathbf{u}}} \neq \bar{\mathbf{u}}, \quad \bar{\mathbf{u}}'' \neq 0$$

We use LES in this work to balance the computational cost of the simulation with the degree of resolution for physical phenomena within the flow. Many studies involving turbulent supercritical fluids utilize LES, as will be seen in the next section, with additional research having been conducted regarding appropriate choices in modeling for the residual stress tensor. Model selection and further details on filtering are discussed in Chapter 2.

### 1.3 Supercritical Carbon Dioxide

As mentioned earlier, supercritical fluids have many qualities that make them desirable as working fluids in a variety of systems. We now shift our focus to one particular fluid of interest: Carbon Dioxide (CO<sub>2</sub>). As seen in the phase diagram of Figure 1.6, the critical temperature,  $T_c$ , and critical pressure,  $p_c$ , of CO<sub>2</sub> are 304.128 K and 73.773 bar. The critical temperature and pressure of CO<sub>2</sub> is fairly easy to attain, making it a strong candidate for systems with high thermal outputs. Additionally, sCO<sub>2</sub> has a relatively low toxicity and environmental impact [], and is chemically stable, non-flammable, and readily available []. For these reasons, sCO<sub>2</sub> is a highly coveted alternative working fluid in many different applications, and is one of the most widely used supercritical fluids along with water [20].

In the next part of this section, we will explore some applications of sCO<sub>2</sub> in which the turbulent round jet configuration is used. Then we will go through a brief review of recent studies involving sCO<sub>2</sub>, and more generally supercritical turbulent jets, in order to demonstrate where this work fits in among current research. Finally, we outline the structure of the remainder of the dissertation.

#### 1.3.1 Applications of Interest

One of the key applications of interest that motivates this work is the use of sCO<sub>2</sub> as the working fluid in advanced cycles for power generation. sCO<sub>2</sub> has shown promise as a working fluid for both indirect cycle and direct-firing cycles [71, 72]. One example of indirect cycle improvement uses sCO<sub>2</sub> in place of water for the conventional steam-Rankine cycle. An example of where these types of configurations may prove useful is in managing thermal runoff from existing coal and natural gas combustion processes [71]. Compared to steam, sCO<sub>2</sub> is less corrosive, more thermally stable, and has increased power density. The critical point of CO<sub>2</sub> is easily accessible, and once achieved,

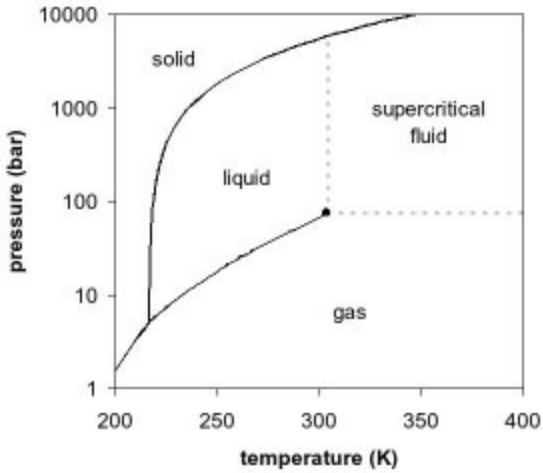


Figure 1.6: Phase diagram for Carbon Dioxide (CO<sub>2</sub>). Critical pressure,  $p_c$ , and temperature,  $T_c$ , are 73.773 bar and 304.128 K, respectively.

allows for the use of a single phase fluid design, leading to a simplified and more compact turbine (see Figure 1.7). Ultimately, this also allows for lower operation and maintenance costs [15]. The benefits of using sCO<sub>2</sub> turbines over the traditional steam design has been highly researched and has only seen an increase in momentum for implementation [14, 17, 15, 49].

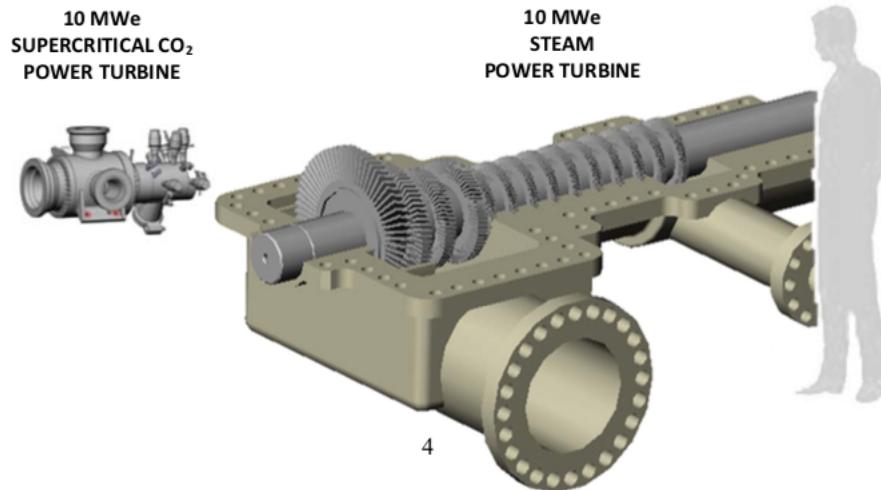


Figure 1.7: Size comparison for steam vs. sCO<sub>2</sub> turbine via Echogen Power Systems LLC [49].

An example of direct-firing cycles that use sCO<sub>2</sub> include the Allam cycle [4]. When compared to

the conventional Brayton cycle, studies show that the Allam cycle has much higher efficiency [16, 5]. Additionally, the carbon footprint for the Allam cycle is virtually zero, allowing for CO<sub>2</sub> produced from the system to be stored underground or used elsewhere, aiding in carbon sequestration efforts [21]. This two-for-one benefit of using sCO<sub>2</sub>-based cycles such as the Allam cycle has spurred much research [69, 9] and development [56] into related technologies.

Of particular importance to these applications is the round turbulent jet, as this is a major component of many injection technologies. The high densities associated with the liquid-like aspect of supercritical fluids coupled with the relatively low gas-like viscosity associated with them typically results in a high Reynolds flow, often resulting in a turbulent system. The turbulence physics of these jets is crucial in developing machinery for these systems. Current experimental research is mainly application oriented. Research into the sCO<sub>2</sub> jet's rock breaking ability has been of primary importance to Enhanced Geothermal Systems (EGS) applications [52, 53, 36, 61, 34], with additional focus being given to pipeline leakage and flow dynamics upon wall impact [68, 67], which unfortunately does not explore the underlying turbulence statistics of the flow. Enhanced recovery of unconventional reservoirs for oil and gas development is an additional area where rock breaking research is often applied [27, 26]. Chemical engineering design aspects of sCO<sub>2</sub> injection are more focused on solubility dynamics as opposed to turbulence [30, 32]. Other experiments focus on similar application specific quantities of interest, such as heat transfer and mixing, which is related in part to the turbulence dynamics [74, 37], but they also note the difficulty in experimental design for investigating these aspects of the flow under the conditions needed to replicate those in real applications [30]. Thus, numerical simulations are necessary to further explore the turbulence statistics of these flows.

### 1.3.2 Overview of Current Numerical Landscape

Numerical simulations are a necessary tool for investigating many aspects of supercritical flow fields due to the often challenging nature of experiment design for the extreme conditions inherent to these systems. Studies using DNS have been implemented to help establish benchmark test cases for other types of numerical schemes [48, 60]. Ruiz et al. use 2D DNS to simulate a mixing layer created by two streams of supercritical Oxygen and gaseous Nitrogen, using two different computational fluid dynamics (CFD) solvers to add confidence to their results [57]. A 3D DNS is used by Ries et al. to simulate a round Nitrogen jet for comparison with experimental data produced by Mayer et al. [55]. However, this study requires a reduction in Reynolds number from  $1.62 \times 10^5$ ,

based on the injection diameter, to 5300 in order to feasibly execute the computations. Li also utilizes a low Reynolds number of 1750 to study a round turbulent sCO<sub>2</sub> jet with a preconditioning scheme [33]. The RANS approach has also been implemented utilizing theory from the ideal gas case [39], but with the goal of ascertaining a more general understanding of why specifically sCO<sub>2</sub>'s rock-breaking ability is better than that of water.

Much research has gone into the development of appropriate numerical methods for investigations regarding turbulence in supercritical fluids. In order to maintain a high Reynolds flow and better capture the effects of the supercritical nature of the fluid on the turbulence dynamics, the use of LES has been explored. The impact of subgrid-scale (SGS) models in capturing transcritical and supercritical dynamics of cryogenic Nitrogen have been analyzed through comparison with the Mayer et al. experiment and highly accurate NIST data [50, 79, 80, 45, 70]. Schmitt et al. does a similar investigation using LES, then extending their investigation to include sCO<sub>2</sub> after validation with the Mayer et al. data [59]. However, this investigation uses low-pressure jets and does note the SGS models might need additional contributions to handle non-linearities and the pressure regime. While many of these investigations note that SGS models may need modification to deal with supercritical flows [59, 50, 45, 79], it is noted by Muller et al. that given a sufficiently fine grid, the influence of SGS modeling and numerical flux discretization is essentially limited to second-order moments [45]. Thus, we will be using the compressible version of the dynamic Smagorinsky SGS closures for our investigation, with further consideration of any influence of SGS model on our quantities of interest being noted later on.

Many of the numerical investigations sited thus far consider cryogenic nitrogen in order to compare with the Mayer et al. experiment on supercritical jet turbulence [41]. A wide variety of numerical investigations into jet turbulence using sCO<sub>2</sub> exist but typically explore either other parameter regimes of interest or application-specific quantities of interest. Examples of turbulent adjacent quantities of interest include fluctuation characteristics based on inlet conditions [76], effects of nozzle and aperture differences on pressure and velocity decay [8] and wave features [38], mixing between sCO<sub>2</sub> and other fluid phases [54], and energy dissipation [35]. These studies all involve high pressure jets and are commonly found in applications involving rock fracturing. Related configurations are also studied, such as the swirling-round sCO<sub>2</sub> jet [73], turbulent jet-in-crossflows [75], slot jet impingement [3], and channel flow [2].

While much of the literature thus far has explored the impact of different numerical methods on modeling supercritical fluid flows and has aimed to strengthen the validity of these simulations

in spite of the lack of experimental data available in the current landscape, a general consensus has still not been reached on how the supercritical nature of these fluids impacts the turbulence physics of these models. Thus, there remain open questions for understanding the fundamental flow behavior of turbulent jets in a supercritical environment, especially near the supercritical point, where both experimental and numerical investigations are still a challenge.

Our objective is to use LES to investigate the turbulence physics of sCO<sub>2</sub> near the critical point in order to capture the effects of widely varying thermal properties of supercritical fluids. Using the compressible Navier-Stokes equation solver, *PeleC* [25, 62], closed with the SRK EoS, we consider three cases in order to examine various quantities of interest associated with classical turbulence mechanics. These three cases are chosen to capture different areas around a peak in specific heat that is associated with the pseudo-critical region. The rest of this dissertation is outlined as follows. Chapter 2 details the model used for this study. Chapter 3 gives an overview of the numerical methods implemented through *PeleC*. Chapter 4 outlines the simulation setup with parameter choices and validation. Results for cross-case comparisons and comparisons with external work are then presented in Chapter 5, with a summary and future work detailed in Chapter 6.

# CHAPTER 2

## MODEL OVERVIEW

### 2.1 Introduction

This section details the governing equations used to simulate the sCO<sub>2</sub> jet. All models and equations described here are from *PeleC* [25, 62], a compressible hydrodynamics code for reacting flows that leverages *AMReX* [78, 77, 29] for AMR. It also leverages the *PelePhysics* library for complex physics, including chemical reactions, non-ideal EoS, and high fidelity transport models. As introduced previously, the compressible Navier-Stokes equations form the basis of the model along with the SRK EoS. Here we also detail the filtering used in the LES and the dynamic Smagorinsky (SMD) used to model the SGS dynamics [40]. We also describe the transport and thermodynamic models used for incorporating into the system real fluid dynamics involving supercritical conditions.

### 2.2 Governing Equations

We consider the three-dimensional compressible Navier-Stokes equations, presented here with Einstein notation:

$$\frac{\partial \rho}{\partial t} + \frac{\partial}{\partial x_j} (\rho u_j) = 0, \quad (2.1a)$$

$$\frac{\partial}{\partial t} (\rho u_i) + \frac{\partial}{\partial x_j} (\rho u_i u_j + p \delta_{ij} - \sigma_{ij}) = 0, \quad (2.1b)$$

$$\frac{\partial}{\partial t} (\rho E) + \frac{\partial}{\partial x_j} ((\rho E + p) u_j + q_j - \sigma_{ij} u_i) = 0 \quad (2.1c)$$

where  $\rho$  is the density,  $u_j$  is the velocity for the  $x_j$  direction,  $p$  is the pressure,  $E = e + \frac{u_i u_i}{2}$  is the total energy,  $e$  is the internal energy, and  $T$  is the temperature. Following the assumptions made for Newtonian fluids [], the diffusive fluxes are

$$\sigma_{ij} = 2\mu S_{ij} - \frac{2}{3}\mu \delta_{ij} S_{kk}, \quad q_j = -\lambda \frac{\partial T}{\partial x_j} \quad (2.2)$$

where  $S_{ij} = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right)$  is the strain-rate tensor,  $\mu$  is the dynamic viscosity, and  $\lambda$  is the thermal conductivity. Models regarding these two components are given in more detail in the next section.

$\delta_{ij}$  here is the Kronecker delta. External forces such as gravity are not included in this study. The system is closed using the SRK EoS [63] to relate pressure, density, and temperature as follows:

$$\begin{aligned} p &= \frac{RT}{V_m - b} - \frac{a\alpha}{V_m(V_m + b)}, \\ a &= \frac{0.42747R^2T_c^2}{P_c}, \\ b &= \frac{0.08664RT_c}{P_c}, \\ \alpha &= (1 + (0.48508 + 1.55171\omega - 0.15613\omega^2)(1 - T_r^{0.5}))^2 \end{aligned} \quad (2.3)$$

where  $R$  is the ideal gas constant,  $T_c$  and  $P_c$  are the critical temperature and pressure of the species, respectively,  $T_r = T/T_c$  is the reduced temperature given by the ratio of the absolute temperature to the critical temperature,  $V_m$  is the molar volume of the species, and  $\omega$  is the acentric factor of the species. All cases are run with a single species, that being CO<sub>2</sub>.

### 2.2.1 Filtered Navier-Stokes Equations

To perform the LES, we consider the filtered compressible Navier-Stokes equations as implemented by Martín, Piomelli, and Graham [40]. Here we go through a brief derivation along with the main assumptions needed. First, applying the filtering operation from Equation 1.13 to Equations 2.1, we get:

$$\frac{\partial \bar{\rho}}{\partial t} + \frac{\partial}{\partial x_j} (\bar{\rho} \bar{u}_j) = 0, \quad (2.4a)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \bar{u}_i) + \frac{\partial}{\partial x_j} (\bar{\rho} \bar{u}_i \bar{u}_j + \bar{p} \delta_{ij} - \bar{\sigma}_{ij}) = 0, \quad (2.4b)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \bar{E}) + \frac{\partial}{\partial x_j} ((\bar{\rho} \bar{E} \bar{u}_j + \bar{p} \bar{u}_j) + \bar{q}_j - \bar{\sigma}_{ij} \bar{u}_i) = 0 \quad (2.4c)$$

In order to avoid having to model  $\bar{\rho} \bar{u}_j$  in the conservation of mass equation in 2.4, Favre-filtering,  $\tilde{\cdot} = \bar{\rho} \cdot / \bar{\rho}$ , is also applied [ ]. Additionally, the following two sets of assumptions are made regarding transport terms [ ]:

$$\overline{\mu(T) S_{ij}} \simeq \mu(\tilde{T}) \widetilde{S_{ij}}, \quad \overline{\lambda(T) \frac{\partial T}{\partial x_j}} \simeq \lambda(\tilde{T}) \frac{\partial \tilde{T}}{\partial x_j} \quad (2.5a)$$

$$\tilde{\mu} = \mu(\tilde{T}), \quad \tilde{\lambda} = \lambda(\tilde{T}) \quad (2.5b)$$

The assumptions in Equations 2.5 are common in the literature regarding LES, even though transport terms are nonlinearly dependent on temperature [ ]. Justification for this type of assumption can be made where if performing a LES with course enough grid, molecular transport coefficients

remain small compared to turbulent transport coefficients []. Applying Equations 2.5 along with the Favre filter to Equations 2.4 yields the following:

$$\begin{aligned}\frac{\partial \bar{\rho}}{\partial t} + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_j) &= 0, \\ \frac{\partial}{\partial t} (\bar{\rho} \tilde{u}_i) + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_i \tilde{u}_j + \bar{p} \delta_{ij} - \tilde{\sigma}_{ij}) &= 0, \\ \frac{\partial}{\partial t} (\bar{\rho} \tilde{E}) + \frac{\partial}{\partial x_j} \left( (\bar{\rho} \tilde{E} \tilde{u}_j + \bar{p} \tilde{u}_j) + \tilde{q}_j - \bar{\sigma}_{ij} \tilde{u}_i \right) &= 0\end{aligned}\quad (2.6)$$

with filtered Equations 2.2 now given by:

$$\tilde{\sigma}_{ij} = 2\tilde{\mu} \widetilde{S_{ij}} - \frac{2}{3} \tilde{\mu} \delta_{ij} \widetilde{S_{kk}}, \quad \tilde{q}_j = -\tilde{\lambda} \frac{\partial \tilde{T}}{\partial x_j} \quad (2.7)$$

Final simplifications to Equations 2.6 come from the SGS terms. The SGS stress  $\tau_{ij}$ , SGS heat flux  $\mathcal{Q}_j$ , SGS turbulent diffusion  $\partial \mathcal{J}_j / \partial x_j$ , and SGS turbulent viscous diffusion  $\partial \mathcal{D}_j / \partial x_j$  set through the following definitions []:

$$\begin{aligned}\tau_{ij} &= \bar{\rho} (\tilde{u}_i \tilde{u}_j - \tilde{u}_i \tilde{u}_j), \\ \mathcal{Q}_j &= \bar{\rho} (\tilde{u}_j \tilde{T} - \tilde{u}_j \tilde{T}), \\ \mathcal{J}_j &= \bar{\rho} (\tilde{u}_j \tilde{u}_k \tilde{u}_k - \tilde{u}_j \tilde{u}_k \tilde{u}_k), \\ \mathcal{D}_j &= \bar{\rho} (\tilde{\sigma}_{ij} \tilde{u}_i - \tilde{\sigma}_{ij} \tilde{u}_i)\end{aligned}\quad (2.8)$$

After substituting the appropriate pieces of Equations 2.8 into the momentum and energy components of 2.6 and applying a few further assumptions, we can eventually get the following:

$$\frac{\partial \bar{\rho}}{\partial t} + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_j) = 0, \quad (2.9a)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \tilde{u}_i) + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_i \tilde{u}_j + \bar{p} \delta_{ij} - \tilde{\sigma}_{ij}) = -\frac{\partial \tau_{ij}}{\partial x_j}, \quad (2.9b)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \tilde{E}) + \frac{\partial}{\partial x_j} \left( (\bar{\rho} \tilde{E} + \bar{p}) \tilde{u}_j + \tilde{q}_j - \tilde{\sigma}_{ij} \tilde{u}_i \right) = -\frac{\partial}{\partial x_j} \left( \gamma c_v \mathcal{Q}_j + \frac{1}{2} \mathcal{J}_j - \mathcal{D}_j \right) \quad (2.9c)$$

with  $\gamma = c_p/c_v$ . The simplification above hinges upon relating enthalpy to internal energy and pressure as in the ideal gas case with  $h = e + p/\rho = c_p T$  with  $c_p$  assumed to be constant. Neither of these assumptions is consistent with supercritical fluids but is chosen here for ease of splitting Equations 2.9 into simulated vs. modeled quantities. This is also standard practice in many supercritical LES studies [], though these assumptions leave room for future work on incorporating higher order EoS into the LES derivation.

Finally, it is also noted that the divergence of the SGS heat flux and the SGS turbulent diffusion are of comparable order of magnitude while the SGS viscous diffusion is an order of magnitude

smaller [40]. Therefore, the SGS turbulent viscous diffusion is omitted from the system. The final set of equations to be discretized and advanced is given by:

$$\frac{\partial \bar{\rho}}{\partial t} + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_j) = 0, \quad (2.10a)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \tilde{u}_i) + \frac{\partial}{\partial x_j} (\bar{\rho} \tilde{u}_i \tilde{u}_j + \bar{p} \delta_{ij} - \widetilde{\sigma_{ij}}) = -\frac{\partial \tau_{ij}}{\partial x_j}, \quad (2.10b)$$

$$\frac{\partial}{\partial t} (\bar{\rho} \tilde{E}) + \frac{\partial}{\partial x_j} \left( (\bar{\rho} \tilde{E} + \bar{p}) \tilde{u}_j + \tilde{q}_j - \widetilde{\sigma_{ij}} \tilde{u}_i \right) = -\frac{\partial}{\partial x_j} \left( \gamma c_v \mathcal{Q}_j + \frac{1}{2} \mathcal{J}_j \right) \quad (2.10c)$$

with Equations 2.7 describing the filtered transport coefficients. The SGS stress  $\tau_{ij}$ , SGS heat flux  $\mathcal{Q}_j$ , and SGS turbulent diffusion  $\mathcal{J}_j$  need to be modeled in order to close the system.

## 2.2.2 Subgrid-Scale Modeling for Large Eddy Simulation

We use the SMD LES model for compressible flow as described by Martín, Piomelli, and Graham [40]. In this work, the grid provides the implicit filtering of the equations. The SGS stress tensor,  $\tau_{ij}$ , is included in the diffusive fluxes and is calculated as follows:

$$\begin{aligned} \tau_{ij} - \frac{\delta_{ij}}{3} \tau_{kk} &= -C_s^2 2 \bar{\Delta}^2 \bar{\rho} |\tilde{S}| \left( \tilde{S}_{ij} - \frac{\delta_{ij}}{3} \tilde{S}_{kk} \right) = C_s^2 \alpha_{ij}, \\ \tau_{kk} &= C_I 2 \bar{\rho} \bar{\Delta}^2 |\tilde{S}|^2 = C_I \alpha \end{aligned} \quad (2.11)$$

with  $\bar{\Delta}$  being the filter width associated with the smallest scale retained by the filtering operation ( $\bar{\Delta}$  is the grid spacing for our cases). Additionally,  $|\tilde{S}| = (2 \tilde{S}_{ij} \tilde{S}_{ij})^{1/2}$ . The two model coefficients are calculated as follows:

$$C = C_s^2 = \frac{\langle \mathcal{L}_{ij} M_{ij} \rangle}{\langle M_{kl} M_{kl} \rangle}, \quad C_I = \frac{\langle \mathcal{L}_{kk} \rangle}{\langle \beta - \hat{\alpha} \rangle} \quad (2.12)$$

where the Germano identity,  $\mathcal{L}_{ij} = T_{ij} - \hat{\tau}_{ij}$ , is used to relate the SGS stress tensor to the “resolved turbulent stresses”,  $\mathcal{L}_{ij} = (\widehat{\rho u_i \rho u_j} / \bar{\rho}) - \widehat{\rho u_i \rho u_j} / \bar{\rho}$ , and the subtest stresses,  $T_{ij} = \widehat{\rho u_i \check{u}_j} - \widehat{\rho \check{u}_i \check{u}_j}$  [22]. In this relationship, a hat denotes quantities associated with a test filter  $\widehat{G}$  which has a characteristic length of  $\widehat{\Delta}$ . The breve denotes Favre-filtered quantities using  $\widehat{G}$  (i.e.,  $\check{f} = \widehat{\rho f} / \bar{\rho}$ ). Additionally,  $M_{ij} = \beta_{ij} - \widehat{\alpha_{ij}}$  with  $\beta_{ij} = -2 \widehat{\Delta}^2 \widehat{\rho} |\tilde{S}| \left( \check{S}_{ij} - \delta_{ij} \check{S}_{kk} / 3 \right)$  and  $\beta = 2 \widehat{\Delta}^2 \widehat{\rho} |\tilde{S}|^2$ .

The SGS heat flux  $\mathcal{Q}_j$  is also modeled dynamically as in [40]:

$$\mathcal{Q}_j = -\frac{\bar{\rho} \nu_T}{Pr_T} \frac{\partial \tilde{T}}{\partial x_j} = -C \frac{\bar{\Delta}^2 \bar{\rho} |\tilde{S}|}{Pr_T} \frac{\partial \tilde{T}}{\partial x_j} \quad (2.13)$$

where  $C$  is modeled as in Equation 2.12 and the turbulent Prandtl number,  $Pr_T$ , is calculated dynamically as:

$$Pr_T = \frac{C \langle T_k T_k \rangle}{\langle \mathcal{K}_j T_j \rangle} \quad (2.14)$$

where

$$T_j = -\widehat{\Delta}^2 \widehat{\rho} |\check{S}| \frac{\partial \check{\tilde{T}}}{\partial x_j} + \overline{\Delta}^2 \bar{\rho} |\tilde{S}| \widehat{\frac{\partial \tilde{T}}{\partial x_j}}, \quad \mathcal{K}_j = \left( \frac{\widehat{\rho u_j \rho T}}{\bar{\rho}} \right) - \frac{\widehat{\rho u_j} \widehat{\rho T}}{\widehat{\bar{\rho}}}. \quad (2.15)$$

Finally, the SGS turbulent diffusion  $\mathcal{J}_j$  is modeled following the strategy proposed by Knight et al. []:

$$\mathcal{J}_j = \tilde{u}_k \tau_{jk} \quad (2.16)$$

For our simulations, we implement the three point box filter as described in [58] with a filter-grid ratio of 2, i.e.  $\widehat{\Delta} = 2\overline{\Delta}$ . With this filter, the convolution kernel from Equation 1.13 is defined as follows:

$$G(\mathbf{x} - \mathbf{r}) = \begin{cases} \frac{1}{\widehat{\Delta}} & |\mathbf{x} - \mathbf{r}| \leq \frac{\widehat{\Delta}}{2}, \\ 0 & \text{otherwise} \end{cases} \quad (2.17)$$

The discrete implementation of this filtering function is described in the numerical methods section.

As noted previously, the choice of SGS modeling is important in accurately capturing the turbulence statistics of the system. Müller et al. found that while the choice in thermodynamic modeling is crucial in capturing first-order moments, the effects SGS modeling is limited to second-order moments [45]. Therefore, our conclusions relating to the turbulence dynamics will be unaffected. That being said, [45] also notes that the choice in SGS model and numerical flux discretization had a larger than expected effect on resolved Reynolds stress profiles. Specifically, the constant Smagorinsky model yielded decaying fluctuation magnitudes during early evolution, resulting in the transition to a fully turbulent mixing zone to start from lower turbulence levels. However, this did agree with the jet break-up location inferred from mean density profiles, which were shifted slightly downstream by comparison to other SGS models. These relationships will be taken into consideration for this study as well and noted in the discussion of Reynolds stress profiles.

## 2.3 Transport Models

Modeling of transport coefficients is done through *PelePhysics* []. There are three modeling options available in *PelePhysics*; we use the *Simple* model. The *Simple* model approximates ideal gas transport coefficients using EGLib functions [19], which have the following form:

$$\ln(q_0) = \sum_{n=1}^4 a_{q,n} (\ln(T))^{n-1} \quad (2.18)$$

where  $q_0$  is the transport quantity of interest (either  $\lambda$  or  $\mu$ ) and  $a$  is the appropriate pre-calculated polynomial fit coefficient (see Table A.1). Chung's high pressure correction for viscosity and thermal conductivity are included to account for real gas dynamics [10]:

$$q = q_k + q_p \quad (2.19)$$

where  $q_k$  is the low-pressure gas transport quantity related to the ideal gas quantity  $q_0$  and  $q_p$  is the high-pressure deviation.

For viscosity, these quantities are:

$$\begin{aligned} \mu_k &= \mu_0 \left( \frac{1}{G_2} + A_6 Y \right), \\ \mu_p &= \left( \frac{36.344 \times 10^{-6} (MT_c)^{1/2}}{V_c^{2/3}} \right) A_7 Y^2 G_2 \exp \left( A_8 + \frac{A_9}{T^*} + \frac{A_{10}}{T^{*2}} \right) \end{aligned} \quad (2.20)$$

where  $M$  is the molecular weight,  $V_c$  is the critical molar volume,  $T^* = T/\epsilon_k$  is a dimensionless temperature scaling using the Lennard-Jones potential well depth [],  $Y = (\rho V_c)/6$ , and  $G_1 = (1 - 0.5Y)/(1 - Y)^3$ , and  $G_2 = \{A_1 [1 - \exp(-A_4 Y)] / Y + A_2 G_1 \exp(A_5 Y) + A_3 G_1\} / (A_1 A_4 + A_2 + A_3)$ . The constants  $A_{1-10}$  are linear functions calculated as follows:

$$A_i = a_{i0} + a_{i1}\omega + a_{i2}\mu_r^4 + a_{i3}\kappa \quad i = 1, \dots, 10 \quad (2.21)$$

where  $\mu_r$  is the reduced dipole moment of the species,  $\kappa$  is the association factor of the species, and  $a_{ij}$  are constants (see Table A.3).

Similarly, thermal conductivity components are given by:

$$\begin{aligned} \lambda_k &= \lambda_0 \left( \frac{1}{H_2} + B_6 Y \right), \\ \lambda_p &= \left( \frac{3.039 \times 10^{-4} (T_c/M)^{1/2}}{V_c^{2/3}} \right) B_7 Y^2 H_2 T_r^{1/2} \end{aligned} \quad (2.22)$$

where  $H_2 = \{B_1 [1 - \exp(-B_4 Y)] / Y + B_2 G_1 \exp(B_5 Y) + B_3 G_1\} / (B_1 B_4 + B_2 + B_3)$  and  $B_{1-7}$  are defined as:

$$B_i = b_{i0} + b_{i1}\omega + b_{i2}\mu_r^4 + b_{i3}\kappa \quad i = 1, \dots, 7 \quad (2.23)$$

where  $b_{ij}$  are constants (see Table A.4). All species-related constants mentioned in this section can also be found in Table A.2.

These high pressure corrections as derived and analyzed by Chung et al. show improved accuracy for approximating transport properties over a wide range of temperatures and pressures when compared to experimental data for a variety of species [10].

Make plots to compare with NIST data.

## 2.4 Thermodynamics and Related Quantities

Modeling of thermodynamic quantities is also done through *PelePhysics* []. In a similar fashion to equation 2.19, thermodynamic properties can be broken up into an ideal gas component and a departure from ideal gas behavior:

$$q(\phi) = q_I(\phi) + q_D(\phi) \quad (2.24)$$

where  $\phi$  is the appropriate combination of state variables involving temperature, specific volume, and pressure,  $q_I$  is the ideal state component and  $q_D$  is the departure component. These decompositions can be derived from fundamental thermodynamic relations []. For example, change in internal energy can be written as follows:

$$de = \left( \frac{\partial e}{\partial T} \right)_V dT + \left[ T \left( \frac{\partial p}{\partial T} \right)_V - p \right] dV \quad (2.25)$$

substituting in the appropriate partial derivative using the SRK EoS and integrating from an ideal state reference point  $(v, t)$  to general point far from ideal conditions  $(V, T)$  yields:

$$e(V, T) - e(v, t) = e(v, T) - e(v, t) + a \left[ \alpha - T \frac{\partial \alpha}{\partial T} \right] \int_v^V \frac{1}{V_m(V_m + b)} dV' \quad (2.26)$$

Simplifying the left- and right-hand side of Equation 2.26 and utilizing the choice of  $v \rightarrow \infty$  for the ideal state volume [], we get:

$$e(V, T) = e(\infty, T) + a \left[ \alpha - T \frac{\partial \alpha}{\partial T} \right] \int_{\infty}^V \frac{1}{V_m(V_m + b)} dV' \quad (2.27)$$

Evaluating the integral in Equation 2.27, we arrive at the final expression for the internal energy of the species described by the SRK EoS:

$$e(V, T) = e_I(T) + a \left[ T \frac{\partial \alpha}{\partial T} - \alpha \right] \frac{1}{b} \ln \left( 1 + \frac{b}{V_m} \right) \quad (2.28)$$

where  $e_I(T) = c_v T$  is the ideal internal energy which is only dependent on temperature and  $e_D(T, V) = a \left[ T \frac{\partial \alpha}{\partial T} - \alpha \right] \frac{1}{b} \ln \left( 1 + \frac{b}{V_m} \right)$  is the departure function for the internal energy. A similar procedure can be done to arrive at the decomposed enthalpy equation involving SRK EoS-specific quantities:

$$h(T, V) = h_I(T) + a \left[ T \frac{\partial \alpha}{\partial T} - \alpha \right] \frac{1}{b} \ln \left( 1 + \frac{b}{V_m} \right) + \frac{RTb}{V_m - b} - \frac{a\alpha}{V_m + b} \quad (2.29)$$

The departure function can be directly evaluated while the ideal gas portion must be approximated.

Ideal specific heat at constant volume scaled by the ideal gas constant  $c_{vI}/R$  is approximated with NASA polynomial fits [42]. These polynomial fits have the following form:

$$\psi(T) = c_{\psi,0} + c_{\psi,1}T + c_{\psi,2}T^2 + c_{\psi,3}T^3 + c_{\psi,4}T^4 + c_{\psi,5}\frac{1}{T} \quad (2.30)$$

where  $c_{\psi,0-5}$  are species-specific constants for each quantity's polynomial fit (see Table B.1). Other ideal gas quantities of interest can also be calculated in this form via integration, so that  $\psi$  here can be either ideal internal energy  $e_I/RT$ , specific heat at constant volume  $c_{vI}/R$ , or enthalpy  $h_I/RT$ . The equations for additional thermodynamic quantities, such as constant volume and constant pressure specific heat, and useful derivatives involving the SRK EoS can be found in Appendix B.2.

# CHAPTER 3

## NUMERICAL METHODS

### 3.1 Introduction

To discretize and evolve the system of partial differential equations, including the LES SGS terms, we use *PeleC* [25, 62]. *PeleC* is a highly scalable code for heterogeneous architectures that is being developed as part of the Exascale Computing Project (ECP) through the Department of Energy (DOE). For spatial discretization, *PeleC* contains a few variations of the general piecewise parabolic method (PPM) originally derived by Colella and Woodward [11]. We utilize a variation that allows for extrema preservation in the presence of steep gradients [43, 13]. *PeleC* also has two options available time-stepping: an explicit second-order Method of Lines (MOL) formulation and an iterative scheme base on a Spectral Deferred Correction (SDC) approach [46]. For this work, we use the SDC method where the time step is dynamically limited using a Courant number of 0.9.

The rest of this chapter is outlined as follows: first we introduce the general finite volume formulation for conservative systems as this forms the basis of *PeleC*, with an additional note on the foundations in *AMReX*. The remaining sections cover the numerical methods chosen from those available in *PeleC*, with motivation provided as to why they were selected for this work and a basic overview of the method. For additional details as to the specific implementations of said methods within *PeleC*, please see the sources referenced throughout, in addition to the general *PeleC* documentation page [] and Github repository []; the full code is available for download and use. See Appendix C for input parameters used in this work.

### 3.2 Finite Volume Formulation

*PeleC* utilizes a finite volume formulation [] to discretize the system given in Equation 2.10. First, we re-write a slightly modified version of this system in a condensed tensor notation, where the diffusive fluxes have been moved to the right-hand side of the original expression:

$$\frac{\partial \mathbf{U}}{\partial t} + \nabla \cdot \mathbf{F} = \nabla \cdot \mathbf{D} \quad (3.1)$$

where  $\mathbf{U}$  is the vector of conserved quantities  $(\rho, \rho\mathbf{u}, \rho E)$ ,  $\mathbf{F}$  contains the advective fluxes, and  $\mathbf{D}$  contains the diffusive terms. Note we have dropped the filter notation from the previous section

for ease and simplicity of writing but again we are working with the filtered interpretation of the compressible Navier-Stokes Equations as derived in Equation 2.10.

In the finite volume formulation, the domain is sub-divided into  $i$  cells, each with volume  $V_i$ . In *PeleC*, these cells are given by boxes in Cartesian coordinates. We then integrate the system in Equation 3.1 over each finite volume to get the following:

$$\int_{V_i} \frac{\partial \mathbf{U}}{\partial t} dV + \int_{V_i} \nabla \cdot \mathbf{F} dV = \int_{V_i} \nabla \cdot \mathbf{D} dV \quad (3.2)$$

In this formulation, a volume average of the state variables are considered through the following definition:

$$\mathbf{U}_i = \frac{1}{V_i} \int_{V_i} \mathbf{U} dV \quad (3.3)$$

Applying the definition given by Equation 3.3 to Equation 3.2 along with the divergence theorem yields the following:

$$\frac{d\mathbf{U}_i}{dt} + \frac{1}{V_i} \int_{S_i} \mathbf{F} \cdot \mathbf{n} dS = \frac{1}{V_i} \int_{S_i} \mathbf{D} \cdot \mathbf{n} dS \quad (3.4)$$

where  $S_i$  is the surface of cell  $V_i$ . To discretize further, the convective and diffusive fluxes must be approximated at the faces of each surface  $S_i$  in order to get the total flux contributions to the system:

$$\frac{d\mathbf{U}_i}{dt} + \frac{1}{V_i} \sum_j S_{i,j} \bar{\mathbf{F}}_{i,j} = \frac{1}{V_i} \sum_j S_{i,j} \bar{\mathbf{D}}_{i,j} \quad (3.5)$$

where  $S_{i,j}$  is the  $j$ -th edge of cell  $i$ , and corresponding subscripts on fluxes imply flux through that equivalent edge. Further methods may then be applied to perform the flux approximations needed and to discretize in time the system established in Equation 3.5.

### 3.2.1 Adaptive Mesh Refinement with *AMReX*

*PeleC* is built in *AMReX* [78, 77, 29], a block-structured hierarchical adaptive mesh refinement library for meshing infrastructure, including grid refinement, distributed parallelism, geometry representation, and output [25]. Data management within *AMReX* is generally structured as follows: grids are represented by Boxes. Each Box is associated with a data container known as a FAB. A MultiFAB manages a collection of FABs, which store data at each AMR level [25]. Grids are refinement ratios must be even and equally implemented in each direction. Figure 3.1 shows an example of this type of refinement hierarchy using a refinement ratio of two for three levels of mesh refinement. Grid refinement is based on user-specified error criteria. Time scales are

also appropriately refined in order to maintain necessary CFL conditions. For further details on *AMReX*, see [78].

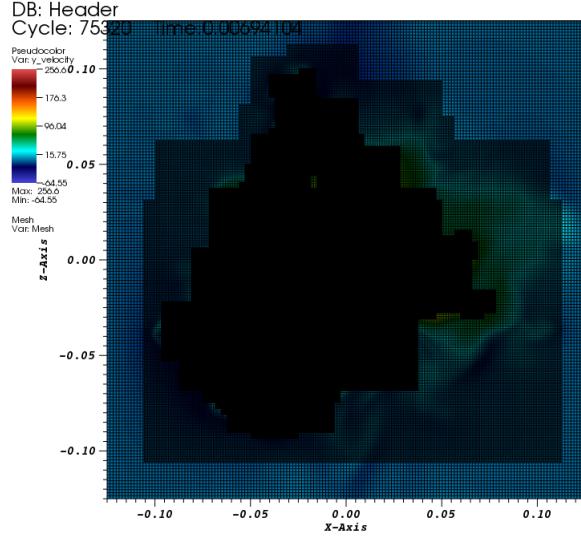


Figure 3.1: 2D slice of 3D data demonstrating mesh refinement using *AMReX*. The refinement ratio is 2 and there are 3 levels of mesh refinement.

### 3.3 Time-Stepping with Spectral Deferred Correction Methods

Here we present the general idea behind SDC methods as is detailed in [47]. Consider an ordinary differential equation (ODE) of the following form:

$$\begin{aligned} \phi_t &= F(\phi(t), t), \quad t \in [t^n, t^{n+1}], \\ \phi(t^n) &= \phi^n \end{aligned} \tag{3.6}$$

with a solution given by the following integral expression:

$$\phi(t) = \phi^n + \int_{t^n}^t F(\phi)d\tau \tag{3.7}$$

Then, given an approximation  $\phi^{(k)}(t)$  to  $\phi(t)$ , one can define a residual as follows:

$$E(t, \phi^{(k)}) = \phi^n + \int_{t^n}^t F(\phi^{(k)})d\tau - \phi^{(k)}(t) \tag{3.8}$$

Defining the error to be  $\delta^{(k)}(t) = \phi(t) - \phi^{(k)}(t)$ , one can then get the error as follows:

$$\delta^{(k)}(t) = \int_{t^n}^t \left[ F(\phi^{(k)} + \delta^{(k)}) - F(\phi^{(k)}) \right] d\tau + E(t, \phi^{(k)}) \tag{3.9}$$

Now, one can use a higher-order quadrature rule to approximate the integral in Equation 3.8. Along with using a lower-order discretization for the integral in Equation 3.9, an iterative scheme can be constructed through which order of accuracy can be improved over subsequent iterative steps, up to the order of accuracy set by the quadrature. This iteration is defined as  $\phi^{(k+1)} = \phi^{(k)} + \delta^{(k)}$ , and using Equations 3.8 and 3.9, we can write an update as follows:

$$\phi^{(k+1)}(t) = \phi^n + \int_{t^n}^t \left[ F(\phi^{(k+1)}) - F(\phi^{(k)}) \right] d\tau + \int_{t^n}^t F(\phi^{(k)}) d\tau \quad (3.10)$$

where again the first integral in Equation 3.10 is approximated with a low-order discretization and the second integral is approximated using higher-order quadrature. *AMReX* uses quadrature.

In the multi-implicit spectral deferred correction (MISDC) formulation [47],  $F$  in Equation 3.6 can be broken up into individual processes:

$$F \equiv A(\phi) + D(\phi) \quad (3.11)$$

where  $A(\phi)$  and  $D(\phi)$  are advection and diffusion processes, respectively. Note this decomposition now recovers the system formed through the finite volume formulation of Equation 3.5. This decomposition in general also includes a reaction term, but since our system is a single species with no reactions it is omitted here. These individual processes can now be treated separately. Discussion of flux treatment is given in the next section. For further details on discretizing Equation 3.10 with 3.11 and applying this to a system of ODEs as in 3.5, please see [46]. For the sake of this work, it is important to note that fluxes can be treated separately in a sort of operator splitting fashion and that accuracy can be set through quadrature choice, which for the methods used here is second-order.

The MOL and SDC time stepping formulations in *PeleC* share a considerable amount of code.

## 3.4 Flux Calculations

Flux approximations are needed at cell faces, as outlined in Section 3.2. Following the methodology outlined in for the MISDC formulation, both advective and diffusive fluxes are centered in time for part of the iterative process [47]. To get the advective fluxes, the PPM [11] is employed to approximate state values near cell faces where they are then extrapolated to  $n + 1/2$ . These values are then used in an approximate Riemann solver [12] to calculate the advective fluxes at cell faces at  $n + 1/s$ . The diffusive fluxes are discretized in space using center difference formulation, with

time-centering for the iterator given by the midpoint rule [13, 43]. Here we detail the PPM a little further and highlight it's advantages for this work.

### 3.4.1 Piecewise Parabolic Method

*PeleC* houses a few variations of the PPM, including the original formulation by Colella and Woodward [11]. The one used in this work is the “extrema preserving” variation of the PPM [13, 43] in order to aid in the handling of steep density gradients. Here we present the brief outline of the method in one dimension as is done in [44] in order to illustrate the general procedure.

First, the system in Equation 3.1 is recast in primitive form. In one dimension, this comes out be:

$$\frac{\partial \mathbf{Q}}{\partial t} + \mathbf{A} \frac{\partial \mathbf{Q}}{\partial x} = \mathbf{S}_{\mathbf{Q}} \quad (3.12)$$

where vector  $\mathbf{Q}$  houses the primitive variables  $(\rho, u, p, \rho e)$ ,  $A$  is the matrix given by  $\partial \mathbf{F} / \partial \mathbf{Q}$ , and  $\mathbf{S}_{\mathbf{Q}}$  is the diffusive terms recast in primitive form. Internal energy is now carried in  $\mathbf{Q}$  to save on the number of calls made to the EoS during the Riemann solver.

The first step in interpolating primitive values to cell edges is to take the average cross-cell difference as follows:

$$\delta q_i = \frac{1}{2} (q_{i+1} - q_{i-1}) \quad (3.13)$$

Monotonicity is enforced with van Leer limiting [66], with limited values denoted  $\delta q_i^*$ . Using these limited values, primitive variables can then be interpolated to cell edges:

$$q_{i+1/2} = q_i + \frac{1}{2} (q_{i+1} - q_i) - \frac{1}{6} (\delta q_{i+1}^* - \delta q_i^*) \quad (3.14)$$

An additional constraint is imposed to ensure  $q_{i+1/2}$  falls between adjacent cell averages.

Next, edge state values  $q_{L,i+1/2}$  and  $q_{R,i-1/2}$  can be set from  $q_{i+1/2}$  and  $q_{i-1/2}$ , respectively. An additional flattening limiter that utilizes local pressure in its construction is applied here as well. Monotonicity is also enforced again, this time across  $q_{L,i+1/2}$ ,  $q_i$ , and  $q_{R,i-1/2}$  within a single cell.

Now, limited piecewise parabolic reconstruction can be achieved by computing average values swept out by parabolic profiles across each face, with the assumption that it moves at the speed of

a characteristic wave  $\lambda_k$ . These averages have the following form:

$$\begin{aligned}\mathcal{I}_+^{(k)}(q_i) &= \frac{1}{\sigma_k \Delta x} \int_{((i+1/2)-\sigma_k)\Delta x}^{(i+1/2)\Delta x} q_i^I(x) dx, \\ \mathcal{I}_-^{(k)}(q_i) &= \frac{1}{\sigma_k \Delta x} \int_{(i-1/2)\Delta x}^{((i-1/2)+\sigma_k)\Delta x} q_i^I(x) dx\end{aligned}\quad (3.15)$$

with  $\sigma_k = |\lambda_k| \Delta t / \Delta x$ , where  $\lambda_k = \{u - c, u, u, u + c\}$  and  $\Delta t$  and  $\Delta x$  are the discretization step size in time and space, respectively. With the parabolic profile  $q_i^I(x)$  given by:

$$\begin{aligned}q_i^I(x) &= q_{R,i-1/2} + \xi(x) [q_{L,i+1/2} - q_{R,i-1/2} + q_{i,6} (1 - \xi(x))] \\ q_{i,6} &= 6q_i - 3(q_{R,i-1/2} + q_{L,i+1/2}) \\ \xi(x) &= \frac{x - x_{i-1/2}}{\Delta x}, \quad x_{i-1/2} \leq x \leq x_{i+1/2}\end{aligned}\quad (3.16)$$

the integrals in Equation 3.15 can be explicitly written as follows:

$$\begin{aligned}\mathcal{I}_+^{(k)}(q_i) &= q_{L,i+1/2} - \frac{\sigma_k}{2} \left[ q_{L,i+1/2} - q_{R,i-1/2} - \left(1 - \frac{2}{3}\sigma_k\right) q_{i,6} \right], \\ \mathcal{I}_-^{(k)}(q_i) &= q_{R,i-1/2} + \frac{\sigma_k}{2} \left[ q_{L,i+1/2} - q_{R,i-1/2} + \left(1 - \frac{2}{3}\sigma_k\right) q_{i,6} \right]\end{aligned}\quad (3.17)$$

Finally, these averages can now be used to perform a characteristic tracing to get the time-centered left and right edge states,  $q_{L,i+1/2}^{n+1/2}$  and  $q_{R,i-1/2}^{n+1/2}$ . Full expressions for these edge states can be found in [], along with the appropriate eigenvalue and eigenvectors needed for this procedure. With the edge states approximated, the advective fluxes can then be calculated via an approximate Riemann solver.

### 3.4.2 Approximate Riemann Solver

*PeleC* uses the approximate Riemann solver developed by Colella and Glaz [12]. This Riemann solver features two iterative scheme options for calculating the classical post-discontinuity state values  $\mathbf{u}^*$  and  $p^*$ : a Secant method and a Newton iteration. Full details for these implementations can be found in [12]. The key feature of this solver that is beneficial to this problem is the treatment of the EoS terms. Special consideration is given in formulation of this method to reduce the number of calls to the EoS within these iterative loops in order to make the method computationally tractable. This is of particular importance for cubic or higher order EoS.

### 3.5 Discrete LES Test Filter

A discretized form of Equation 2.17 is needed for the test filter in calculating the SMD model coefficients. *PeleC* contains a variety of filtering options. As mentioned in Chapter 2, we use the three point box filter as described in [58]. In one dimension, this discretization results in the following implementation:

$$\bar{\phi}_i = \frac{1}{24} \epsilon^2 (\phi_{i+1} + \phi_{i-1}) + \frac{1}{12} (12 - \epsilon^2) \phi_i \quad (3.18)$$

where  $\epsilon$  is the grid filter ratio, which in our case is equal to two.

# CHAPTER 4

## SIMULATIONS

### 4.1 Computational Domain

#### 4.1.1 AMR and Discretization

Our target simulation is a 3D LES with a grid size that is 100 times the smallest scale of the turbulent flow. In our setup, the domain length in each direction is:  $x = 25d$ ,  $y = 62.5d$ , and  $z = 25d$ , where the jet diameter  $d = 0.01\text{ cm}$  and the  $y$ -axis is the axial direction of the jet. This smallest scale of these turbulent flows, known as the Kolmogorov scale [31], can be approximated:

$$\eta = \left( \frac{\nu^3}{\varepsilon} \right) \quad (4.1)$$

where  $\nu = \mu/\rho$  is the kinematic viscosity and  $\varepsilon = v_j^3/d$  is used to approximate the average rate of dissipation of turbulence kinetic energy per unit mass. For these turbulent jets,  $\eta = 5.37 \times 10^{-6}\text{ cm}$ . Therefore, to keep the calculation tractable and achieve an adequate LES grid size, we implement four levels of refinement, with a refinement ratio of 2, leading to 80, 200, and 80 cells on the coarsest level in the  $x$ ,  $y$ , and  $z$  directions, respectively. This results in an initial mesh size of  $\Delta x_0 = \Delta y_0 = \Delta z_0 = 0.3125$ , leading to  $\Delta x_3 = \Delta y_3 = \Delta z_3 = 3.9062 \times 10^{-4}$ , where the subscript denotes the AMR level. These four levels of refinement occur in the region outward from the jet inlet up to a distance of  $20d$  in the  $x$  and  $z$  direction and  $60d$  in the  $y$  direction, as can be seen in Figure 4.1. The refinement criterion is given by the vorticity, specifically with  $\omega \geq 5000^{2l}$ , where  $\omega$  is the magnitude of the vorticity and  $l$  is the AMR level. For the first ten flow throughs of the simulation, mesh refinement only occurs up to one level within the refinement region to establish the flow pattern. Thereafter, the simulation proceeds with the four levels of mesh refinement until reaching steady state, whereupon statistics are collected for analysis.

#### 4.1.2 Initial and Boundary Conditions

Our inlet consists of an opening centered in the  $xz$ -plane with diameter  $d = 0.01\text{ cm}$  through which the sCO<sub>2</sub> jet is initialized. The pressure in the jet at the inlet is the same as that of the quiescent background fluid and it is given by  $p_j = p_0 = 10.1325\text{ MPa}$ . The ambient fluid remains at rest while the jet is initialized with an inflow velocity of  $v_j = 1800\text{ cm s}^{-1}$ , leading to a Reynolds

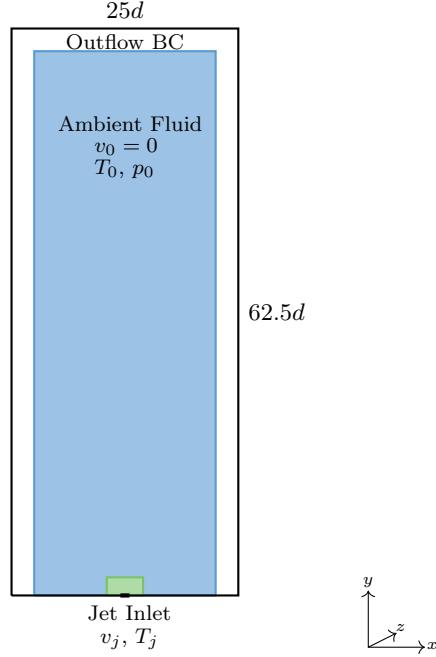


Figure 4.1: Two dimensional slice schematic of jet setup. Four levels of refinement are enforced within the green box based on proximity to jet inlet. Refinement based on vorticity criterion then occurs within the blue region. Outside the blue region, AMR is explicitly turned off to allow flow structures to be dissipated numerically and allowed to leave the domain without incurring spurious reflections.

number of the initialized jet of  $Re_j = 22910$ , with  $v_j$  and  $d$  being the reference velocity and length scale respectively. For the jet temperature and pressure conditions given, the reference density is  $\rho_j = 3.019 \times 10^{-1} \text{ g cm}^{-3}$ , as calculated via the SRK EoS in *PeleC*. To implement a turbulent inflow, we add noise to this input velocity using the mean velocity and root mean square (rms) values from a predetermined DNS velocity profile [18]. We begin by scaling the DNS values with our given jet velocity and radius. After determining which scaled DNS values we are near for the radius determined by our mesh, a linear interpolant is created to compute the mean velocity at our given radius. Noise is then added to each cylindrical component of the velocity as follows:

$$v = \langle v_{\text{DNS}} \rangle + (v'_{\text{DNS}} + \beta v'_{\text{DNS}} r_1 \sin \theta_1) \cdot r_2 \sin \theta_2 \quad (4.2a)$$

$$u = u'_{\text{DNS}} + \beta u'_{\text{DNS}} r_3 \sin \theta_3 \quad (4.2b)$$

$$w = w'_{\text{DNS}} + \beta w'_{\text{DNS}} r_4 \sin \theta_4 \quad (4.2c)$$

where  $\langle u_y \rangle$  is the mean velocity in the axial direction,  $u' = \langle u^2 \rangle^{\frac{1}{2}}$  is the rms for each velocity component, and  $\beta = 0.1$ . Each  $r_i$  and  $\theta_k$  value is randomly generated as follows:

$$r_i = \sqrt{-2.0 \log(X_i)} \quad (4.3a)$$

$$\theta_k = 2\pi X_k \quad (4.3b)$$

where  $X_n$  are random numbers between 0 and 1. The inflow parameters are finalized after being converted into Cartesian coordinates.

We implement zero gradient boundary conditions for all boundaries not involving the jet inflow. Additionally, AMR is halted at a distance of  $2.5d$  from the boundary in the  $x$  and  $z$  directions, and that of  $5d$  in the axial direction in order to ensure that all waves are dissipated, thus avoiding spurious reflections from the boundaries.

## 4.2 Case Descriptions

Three cases are investigated in this study. The first case pertains to an isothermal jet, in which the jet conditions match that of the ambient fluid for all quantities aside from velocity. The other two cases consider non-isothermal jets, in which the ambient fluid is adjusted to no longer match the temperature of the jet inflow. All three cases consider the same inflow conditions, while ambient fluid conditions are what varies between them. Temperatures for the ambient fluid in the non-isothermal cases are chosen such that the ambient fluid is supercritical but the temperature for each case falls on either side of the peak in specific heat seen near the critical point. This peak is associated with the pseudo-boiling phenomenon seen in supercritical fluids, where upon crossing the Widom line, the supercritical fluid experiences a shift from being more gas-like to more liquid-like in nature, although distinct phase separation between the two is not present. A summary of the conditions for each case can be seen in Table 4.1.

Table 4.1: Summary of Jet Parameters for each case.

Parameter	Case 1	Case 2	Case 3
$p_{in}$ MPa	10.1325	10.1325	10.1325
$p_\infty$ g/cm <sup>3</sup>	10.1325	10.1325	10.1325
$T_{in}$ K	330	330	330
$T_\infty$ K	330	350	314
$v_{in}$ (cm/s)	1800	1800	1800
$\rho_{in}$ g/cm <sup>3</sup>	10.1325	10.1325	10.1325
$R_{in}$	22911	22911	22911
$M_{in}$	0.08	0.08	0.08

The goal in choosing the ambient temperature in this way is to investigate the impact of large thermodynamic changes on the flow field of the supercritical jet. An example of this can be seen with the sharp peak in constant pressure specific heat in Figure 4.2. This choice differs from similar studies, which typically involve transcritical injection.

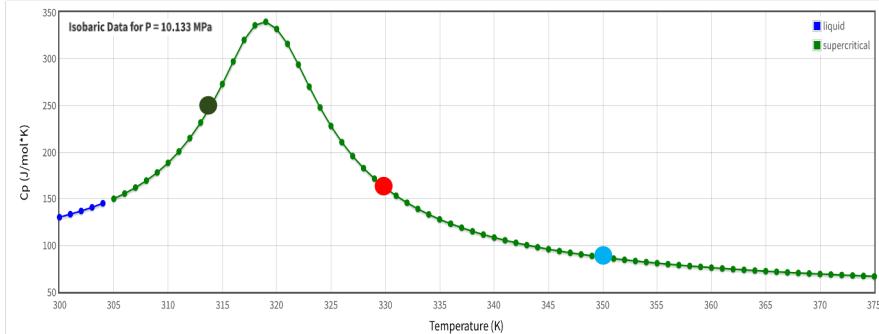


Figure 4.2: Extreme thermodynamic variation seen near critical point with ambient temperatures plotted for reference to see regime coverage of each case. Red is the jet temperature for all cases and isothermal temperature for the ambient case. Blue is the 350K ambient case. Green is the 314K ambient case.

### 4.3 Compute Time and Hardware Specifications

Simulations were run on the National Renewable Energy Laboratory (NREL) Eagle supercomputer using the Intel suite of compilers. Simulations utilized 576 MPI ranks in total, with 36 ranks spread across 16 nodes. Roughly 17,000 cells were handled per MPI rank. Dynamic load balancing of AMR levels allowed for maximum utilization of the system. In total, over 100 TB of data were

stored from the original 3D simulations run over the course of the 10 flow-throughs of the domain needed for averaging procedures. To help manage data storage issues, slicing procedures were implemented to reduce full 3D simulation data down to 2D slice data, where averaging procedures were then performed as outlined in the next section.

## 4.4 Post-Processing Procedures

Post-processing procedures were written in Python and primarily leveraged the following libraries: yt, pandas, numpy, and, matplotlib. Original 3D data at each plotfile was sliced in three different ways using yt: along the centerline to create 1D data, normal to the jet axis at various points down stream to create 2D data, and along the axial direction at  $z = 0$  to create 2D data.

For time averaging, a basic discrete averaging procedure was implemented over  $i$  saved plot files from the final 2 flow-throughs of the simulation, with 100 plots saved per flow-through:

$$\phi_{avg,slice} = \frac{1}{200} \sum_{i=1}^{200} \phi_{i,slice} \quad (4.4)$$

Radial averages were approximated with  $r = 0$  lying in the center of the  $xz$ -plane with  $r = \sqrt{x^2 + z^2}$ . The  $N$  points nearest to each fixed  $r$  distance away from the center were then averaged together:

$$\phi_{avg,r} = \frac{1}{N} \sum_{i=1}^N \phi_{i,r} \quad (4.5)$$

For quantities that were averaged in both time and radially, temporal averaging was performed first, with the radial average then taken of the time averaged slice data.

# CHAPTER 5

## RESULTS

### 5.1 Isothermal Jet

Our first case involves the isothermal jet where both jet and ambient fluid are at 330K. Quantities of interest for this case are compared against incompressible round jet statistics as outlined in Pope [51] and isothermal compressible jet turbulence features []. This case serves as a baseline for comparison with the other two cases involving non-isothermal jets.

#### 5.1.1 Flow Field Features

All images herein depict a slice of the 3D flow field at  $z = 0$ . Each figure contains an instantaneous snapshot and a time-averaged snapshot of the entire slice domain, plus an additional zoomed-in image of the time-averaged quantity of interest near the inlet. All instantaneous images are taken from the final data point of the simulation.

Figure 5.1 shows the axial velocity component of the flow field. The general spreading rate and decay of the velocity field can be seen from the instantaneous snapshot in 5.1a. The flow is mostly steady up until  $y/d = 2$  before perturbations begin. The stream mostly stays together through these initial perturbations up until  $y/d = 5$  where spreading then begins. The averaged axial velocity fields in 5.1b and 5.1c shed more light on the approximate development regions of the jet. The potential core extends up to  $y/d = 5$ , followed by the transition region where  $5 \leq y/d \leq 15$ , thereafter the jet appears fully developed. Centerline analysis of the turbulent kinetic energy later on will provide more information for these boundaries.

Figure 5.2 shows minor pressure fluctuations in the flow, scaled against the maximum pressure achieved above the ambient pressure. 5.2a shows minor pressure oscillations mirrored on each side of the jet edge in the same zone as the initial velocity fluctuations see in 5.1a. Thereafter, the oscillations become off-kilter, correlating to the beginning of the jet disintegration as seen in the velocity field. Pressure fluctuations are concentrated near the inlet and die down past the transition zone. This can be seen more clearly in 5.2b. On average, pressure fluctuations yield a minor increase within the potential core and decrease on the jet perimeter, as can be seen in 5.2c. Fluctuations also lead to a minor pressure drop on average in the transition region.

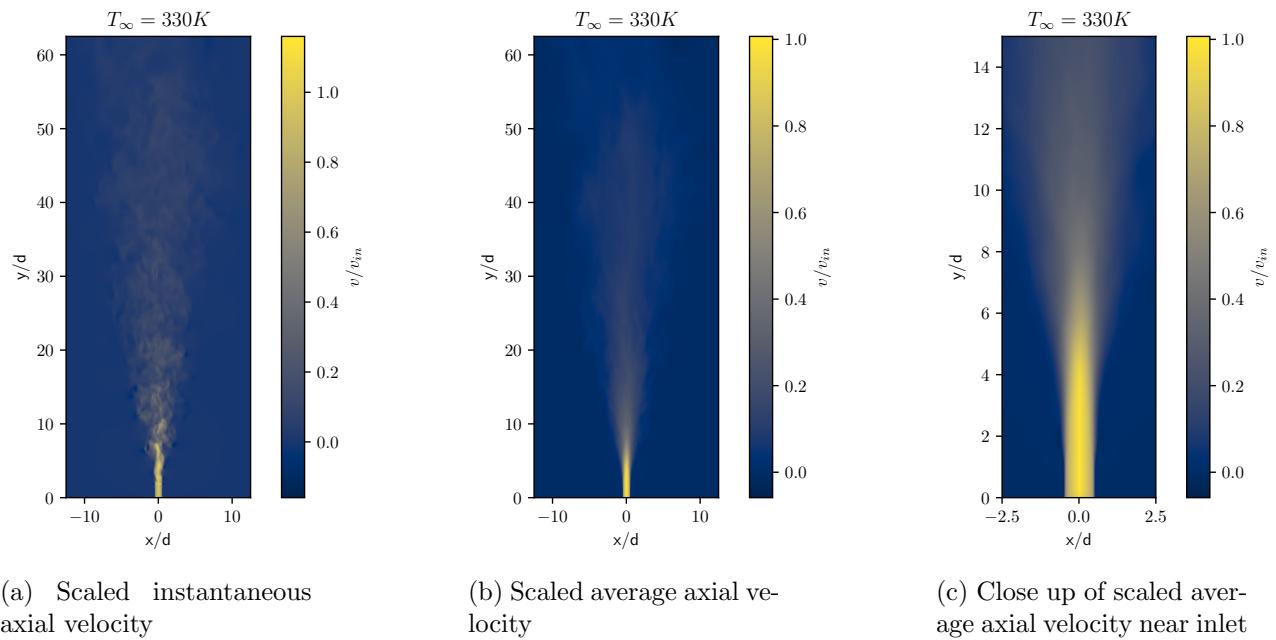


Figure 5.1: Axial velocity features of the isothermal jet

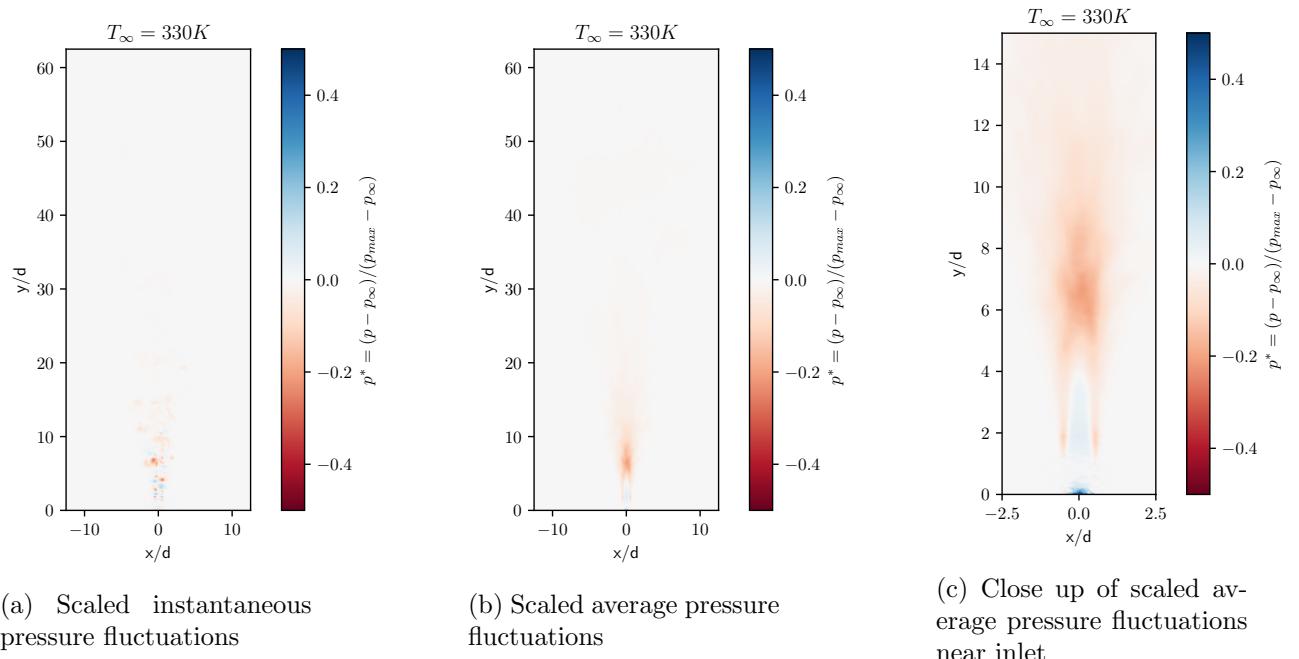


Figure 5.2: Pressure features of the isothermal jet

Figure 5.3 shows magnitude of the vorticity scaled between the maximal and minimal values. 5.3a shows the strongest vorticity occurring at the jet interface with the ambient fluid near the inlet. Past this initial stage, vortex shedding occurs and vorticity dissipates as the jet spreads. The averages in 5.3b and 5.3c both show again that the most intense vorticity occurs at the inlet along the outer edge of the jet. This high intensity remains constant until about  $y/d = 1$  before more mixing with the ambient fluid occurs as the jet spreads and the vorticity lessens in intensity. Vortices are still restricted to the jet edge until around  $y/d = 5$  where the transition to fully developed turbulence enables vortical motions to extend across the fully spread of the jet.

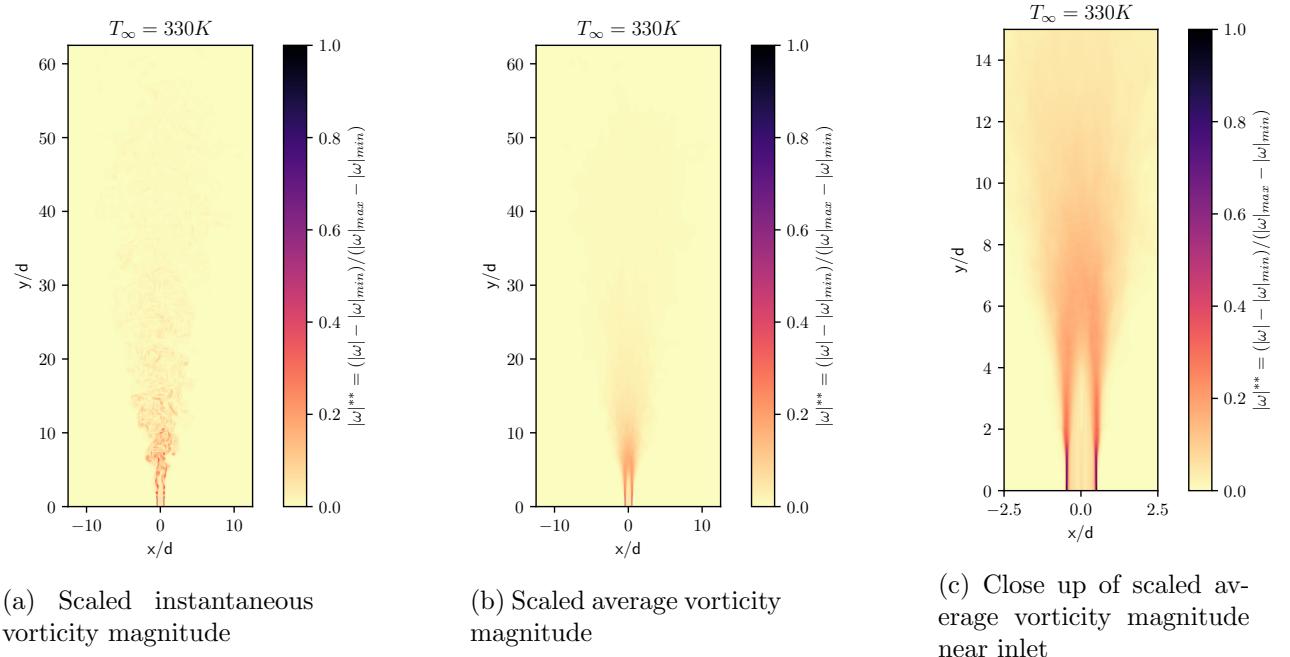


Figure 5.3: Vorticity magnitude features of the isothermal jet

### 5.1.2 Mean Flow Properties

Figure 5.4 depicts the time and radially averaged scaled axial velocity component plotted against radial distance from the centerline at multiple normal slices downstream from the inlet. The velocity is scaled by the average axially velocity value at the inflow while the radial direction is scaled by the jet diameter. These plots demonstrate the axially velocity decay as the flow progresses further downstream. As velocity value along the centerline decreases, the overall velocity profile also widens and flattens out. Typically, for the round turbulent jet, this expansion would occur in such a way

that upon specially selected scaling, these profiles would collapse into one profile after a certain point. This potential is explored further in the next figure. Here though this decay is a common feature of how the axial velocity of the jet develops as it leaves the inlet.

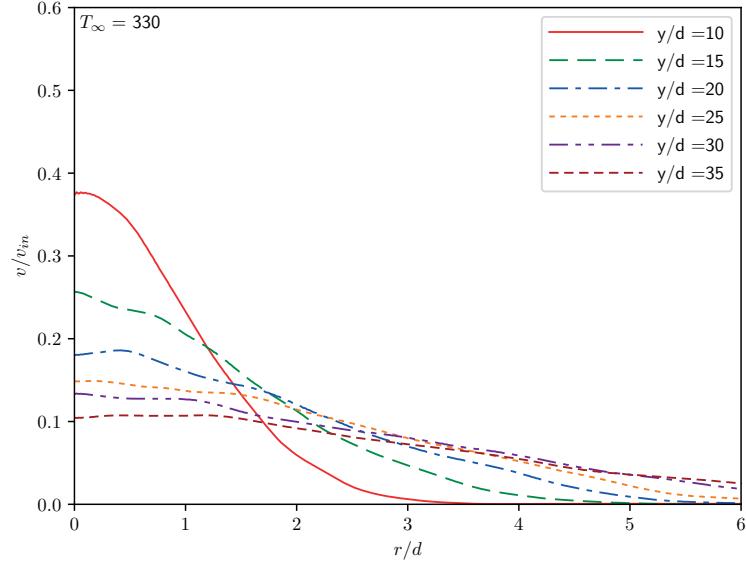


Figure 5.4: Average (both in time and radially) axial velocity scaled by inlet value plotted along radial distance from centerline. Profile decay follows similar trajectory to what is expected in incompressible round jet theory [51].

Figure 5.5 depicts time and radially averaged scaled axial velocity components plotted against the radial distance from the centerline. Each curve is made at a slice normal to the axial flow direction at different points downstream. Figure 5.5a depicts velocity curves every  $3d$  downstream from the inlet near where the transition region begins while Figure 5.5b contains plots taken every  $5d$ . The axial velocity is scaled by the centerline value  $v_c$  while the radial distance is scaled by the half width half mean (HWHM), where the velocity component is equal to half the value on the centerline  $v(r_{1/2}) = v_c/2$ .

Figure 5.5a shows the near-field axial velocity profiles. They already exhibit self-similarity collapse into one profile which is most likely the result of the inflow condition []. In the far-field slices of 5.5b, self-similarity is fairly well maintained with minor fluctuations in the center and edge of the jet. These fluctuations are most likely the result of low resolution in the time averaging of available data.

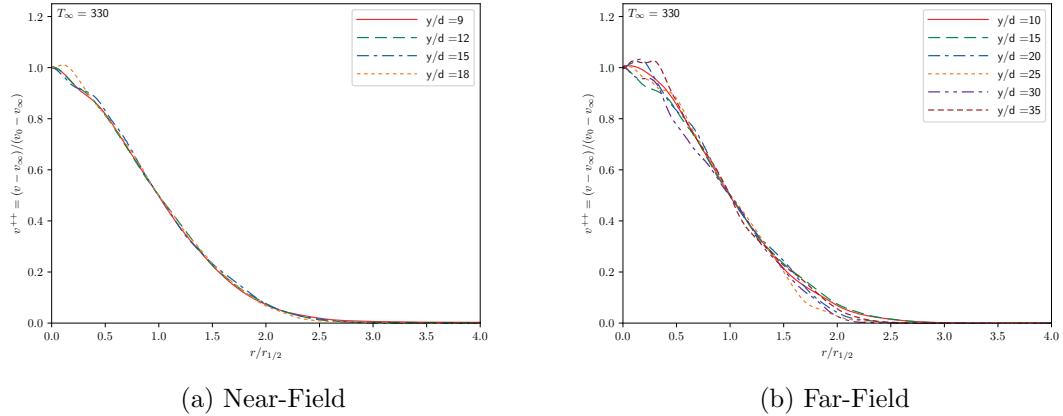


Figure 5.5: Normal slices of scaled axial velocity, averaged in both time and the radial direction. Plotted against radial direction scaled by  $r_{1/2}$ . Both near- and far-field regions demonstrate the self-similarity within the round turbulent jet.

Another common feature of round turbulent jets is the development of a linear relationship between the jet centerline value and the distance downstream. This comparison is depicted in Figure 5.6. Here, the centerline value of the axial velocity  $v_0$  is inversely linearly proportional to the distance downstream. The dip at  $y/d = 30$  may be attributed to again lack of convergence due to low sample size in time averaging.

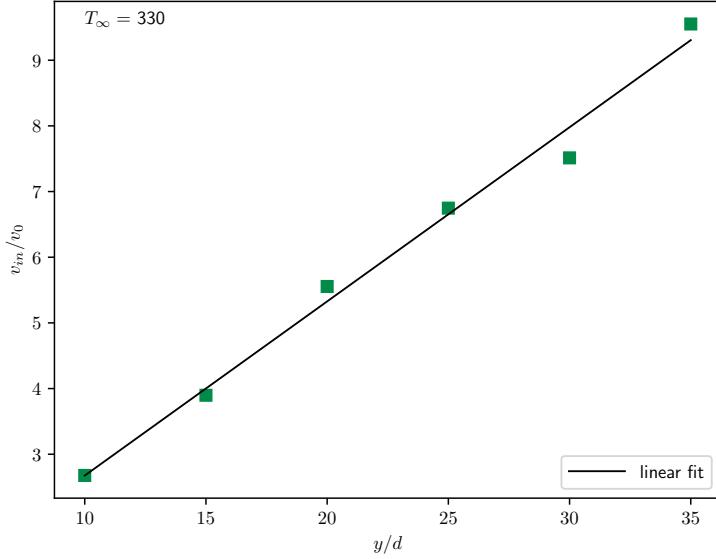


Figure 5.6: Axial inlet velocity scaled by centerline values along the axial direction. When distance downstream is scaled by jet diameter, linear decay of the centerline axial velocity is observed.

### 5.1.3 Turbulence Dynamics

Figure 5.7 shows the time and radially averaged Reynolds stresses at two points downstream from the inlet. Here, velocity components  $(u, v, w)$  correspond to the  $(r, y, \theta)$  directions, respectively. Each Reynolds stress component follows general trends associated with round turbulent jets [51], with the axial component providing the leading contribution, followed by the other two directional components closely with these two being of roughly the same magnitude, and the cross-directional component contributing the least. The cross-directional component is non-zero at the jet center; this could be due to... Note also that self-similarity is not exhibited, as each component exhibits an increase in magnitude at the center of the jet as distance downstream is increased.

Figure 5.8 shows the time average turbulent kinetic energy components along the centerline of the jet. Each component grows through the potential core region of the jet up until all components reach a peak in energy around  $y/d = 6$ , with the axial component's peak coming slightly ahead of the other two directions. Rapid decay is then observed up until around  $y/d = 15$  before a slower decay sets in up until a leveling off is achieved around  $y/d = 30$ . This rapid decay and then further progression correspond to the transition and fully developed jet regions, respectively.

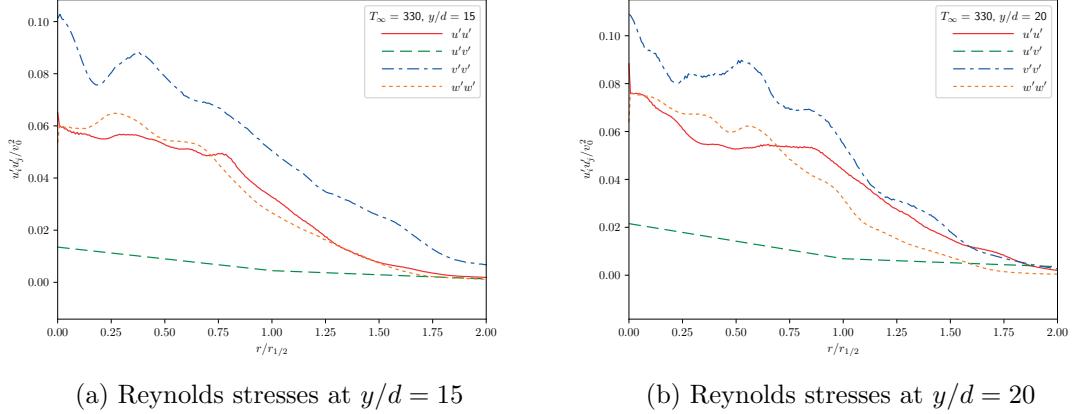


Figure 5.7: Time and radially averaged Reynolds stresses for the isothermal jet at two locations downstream. Both slices follow similar Reynolds stress relations as seen in incompressible round jet [51].

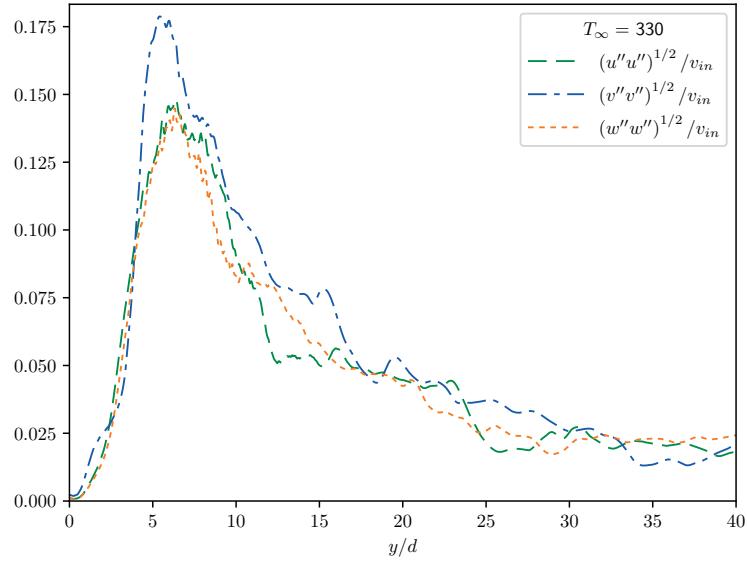


Figure 5.8: Average turbulent kinetic energy components along centerline.

## 5.2 Non-Isothermal Jets

The two non-isothermal jet cases as described in Chapter 4 are presented here. Some features are compared to the isothermal jet case while others are used for direct comparison between the two non-isothermal cases. Again, the  $350K$  ambient case depicts a jet entering a warmer ambient fluid, moving further away from the supercritical point while the  $314K$  ambient case depicts a jet entering a cooler ambient fluid with a transition over the pseudo-boiling point. It is anticipated that the  $314K$  ambient case will then vary much more prominently from the other two cases across various quantities of interest examined.

### 5.2.1 Flow Field Features

Figure 5.9 shows various axial velocity field comparisons between both the isothermal and non-isothermal jets. The instantaneous axial velocity over the whole domain is depicted in 5.9a. The  $350K$  ambient case has a laminar flow region of similar length to the  $330K$  ambient case, with the jet core staying intact up until  $y/d = 2$  before fluctuations begin to take effect. The developing region after this within  $5 \leq y/d \leq 15$  appears to be slightly more volatile in the  $350K$  ambient case compared to the  $330K$  ambient case, with the region containing finer-scale fluctuations. The spreading rate of the  $350K$  ambient case is smaller than that of the base case while overall intensity levels seem to be comparable in the downstream direction. The  $314K$  ambient case has a much shorter laminar region compared to the base case, with fluctuations beginning to become apparent at  $y/d = 1$ . Perturbations after that appear to be more symmetric in structure, with vortices along the jet edge staying in tandem instead of the oscillatory shedding apparent in the other two cases. The  $314K$  ambient case also appears to have a spreading rate comparable to the  $350K$  ambient case, but with a faster decay in intensity in the downstream direction. 5.9b helps further illuminate the difference in spread of the three jets. Here the  $350K$  ambient core intensity is more persistent downstream with a more distinct and thinner outer edge flow present than in the  $330K$  ambient case, while the  $314K$  case has a lower intensity central flow region but a comparably sized edge flow region compared to the  $330K$  ambient case. Finally, 5.9c helps depict the potential core and transition regions of the three cases more clearly. Here, the similarity between the potential cores of the  $330K$  and  $350K$  ambient cases is clear while the  $314K$  ambient case has a much shorter core length.

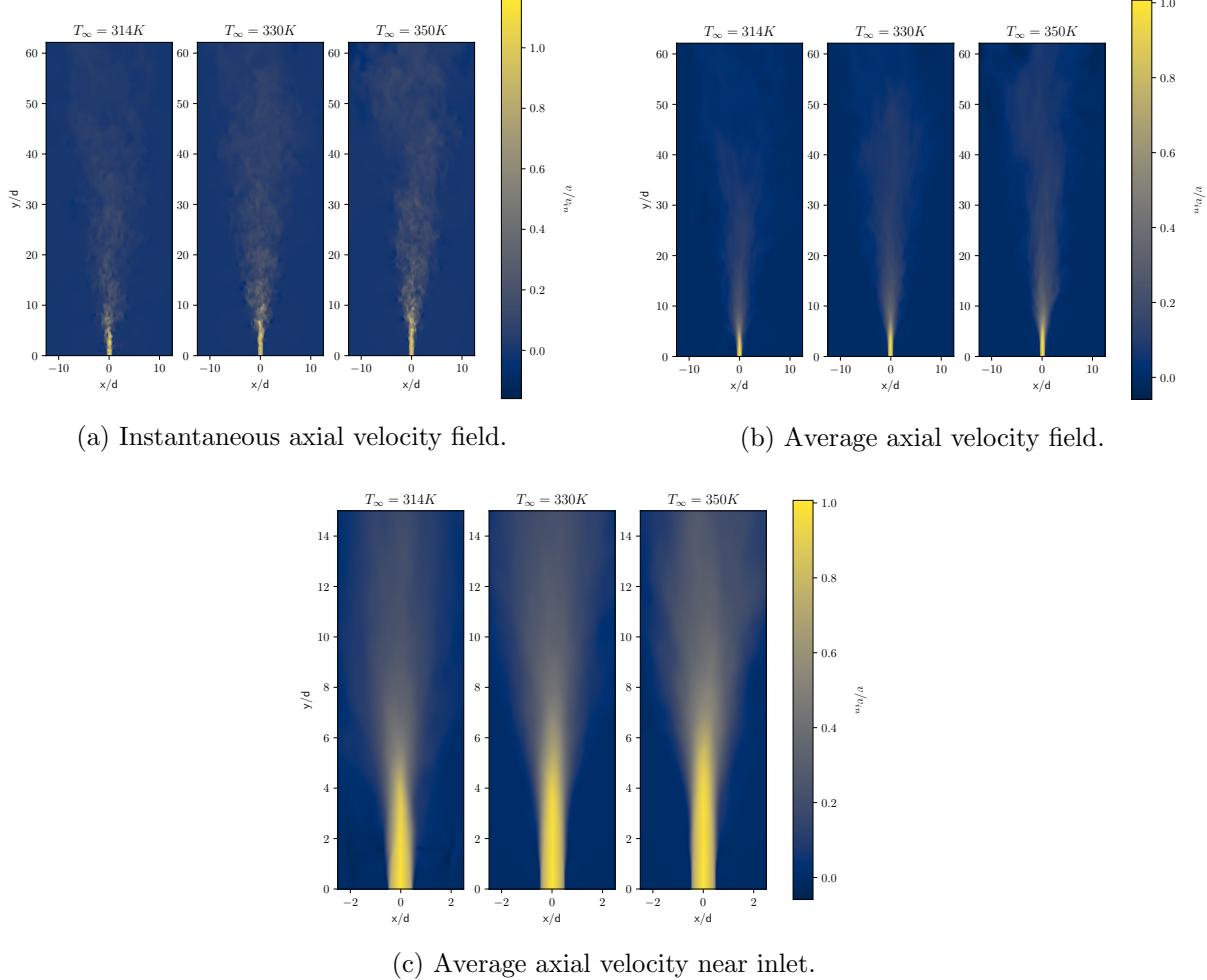


Figure 5.9: Axial velocity field comparison between 314K ambient case (left), 330K ambient case (middle), and 350K ambient case (right).

Figure 5.10 shows pressure fluctuations near the inlet for all three cases. 5.10a contains an instantaneous snapshot of these fluctuations near the inlet. The 350K ambient case is comparable to the 330K ambient case, with similar large coherent structures forming in the pressure oscillations. However, the initial pressure waves emanating from the inlet are stronger in the 350K ambient case, resulting in slightly more intense fluctuations downstream. Minor spurious perturbations are also seen downstream in the 350K ambient case. The 314K ambient case is qualitatively much different from the other two cases. Larger coherent structures are much less defined, with smaller fluctuations appearing much more prominent throughout. There is also a band of cells slightly before  $y/d = 2$  where perturbations are more concentrated. This could be due to... some sort of glsamr issue?

\*(this is right around where the grid refinement drops from the imposed high-level refinement at the inlet)\* Average pressure fluctuation can be seen in 5.10b. All cases have a slight increase in pressure isolated within the potential core surrounded by and followed with a pressure drop. The difference between these low and high pressure pockets is most prominent in the  $314K$  ambient case, followed by the  $330K$  ambient case. Numerical artifacts are extremely prominent in the  $314K$  ambient case, most likely due to.... with some similar artifacts present in the  $350K$  ambient case as well.

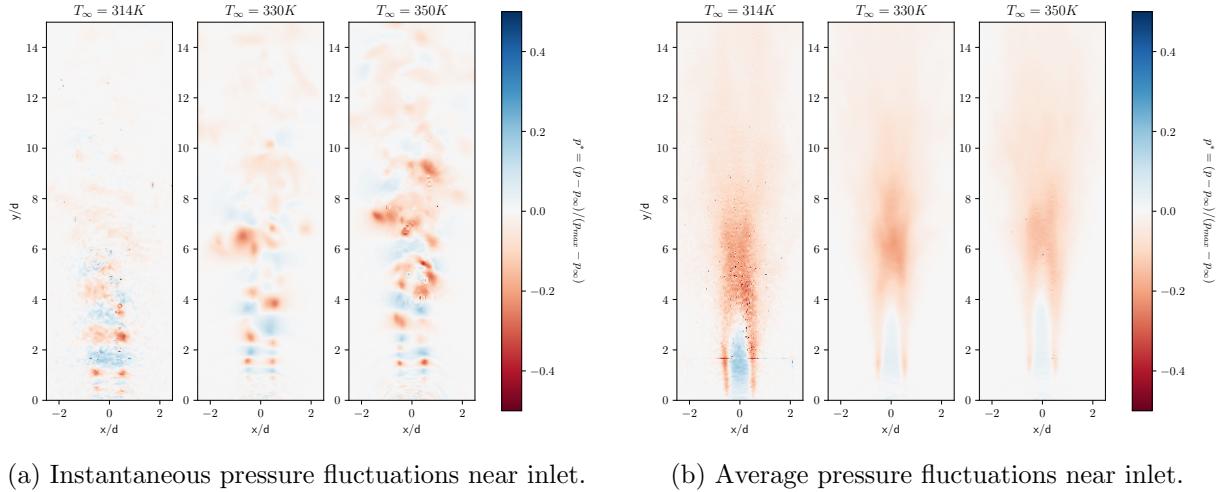


Figure 5.10: Pressure fluctuation field comparison between  $314K$  ambient case (left),  $330K$  ambient case (middle), and  $350K$  ambient case (right).

Figure 5.11 shows both instantaneous and time-averaged flow fields of the vorticity magnitude for the three cases. The vorticity magnitude features seen in 5.11a follow similar trends as were established with the axial velocity discussion in Figure 5.9a. In all three cases, the strongest vorticity occurs along the jet edge right at the inflow, leading to vortex formation and subsequent fluctuations. This steady inflow window before fluctuations lead to jet disintegration is comparable between the  $330K$  and  $350K$  ambient case. However, the  $350K$  maintains higher vorticity intensity downstream with pockets of stronger vorticity seen in the transition region of the jet compared to that of the  $330K$  ambient case. The  $314K$  ambient case has a shorter region of steady flow in and also features more prominently distinct vortex formation along the edge of the jet. These vortices roll along the edge in tandem before mixing to lead into the transition region around  $y/d = 3$ . Similar to what was seen with the  $314K$  ambient pressure field, there are some vorticity artifacts present in the mesh at the edge of the initial high-refinement zone surrounding the jet inlet. Similar

to the  $350K$  ambient case, the  $314K$  ambient case also maintains some pockets of high vorticity, but then does an overall decay downstream that is much more intense than the other two cases. The averaged vorticity magnitude fields in [5.11b](#) and [5.11c](#) help showcase the different jet regions and jet spreading. From [5.11b](#) one can see that the  $350K$  ambient case has more widespread vorticity in the transition region from around  $5 \leq y/d \leq 15$  and then maintains the intensity formed here for further along downstream compared to the  $330K$  ambient case. The  $314K$  ambient case has a much narrower vorticity range in the same region with intensity decaying much more rapidly downstream. Near the inlet, as depicted in [5.11c](#), it's easier to see that the  $314K$  ambient case begins spreading much earlier than the other two cases and has much less uniform vorticity along the jet edges. These vorticity fluctuations before the mesh refinement change may contribute to the vorticity artifacts seen? This earlier breakup also impacts the way the high-vorticity outer edges of the jet merge around the low-vorticity center at the end of the potential core, where the merging of the two jet edges across the middle meet in an almost concave point as compared to the more convex central region seen in the other two cases.

The final figure in this section, Figure [5.12](#), contains direct comparisons between the two non-isothermal cases for various quantities of interest. Figure [5.12a](#) shows temperature change between the jet and the ambient fluid. The jet more rapidly approaches the ambient fluid temperature in the  $314K$  ambient case as compared to the  $350K$  ambient case, where temperature variation from the inlet is much more gradual downstream. Figure [5.12b](#) shows the constant-pressure specific heat variation between the two cases. The added peak in specific heat can be seen in the  $314K$  ambient case as the fluid transitions through the pseudo-boiling point. This peak occurs at the jet edge in the potential core and is additionally maintained after the transition region, showing a slower transition to the ambient specific heat as compared to the  $350K$  ambient case. [5.12c](#) shows the density transition between the jet and the ambient fluid. Density change downstream in both cases occurs at similar rates, with both jets exhibiting a rapid shift during the transition region and then continuing with a much slower adjustment thereafter. The last group of images, [5.12d](#), [5.12e](#), and [5.12f](#), show the sound speed, enthalpy, and compressibility factor of the fluid, respectively. All three quantities exhibit similar trends, with the  $350K$  ambient case showing a more rapid transition to the ambient conditions as compared to the  $314K$  ambient case. These follow the same trend as seen in the temperature, but to varying degrees of intensity. For example, enthalpy appears to have an overall faster downstream decay for both cases as compared to the temperature decay, while sound speed and compressibility factor have comparable decay rates.

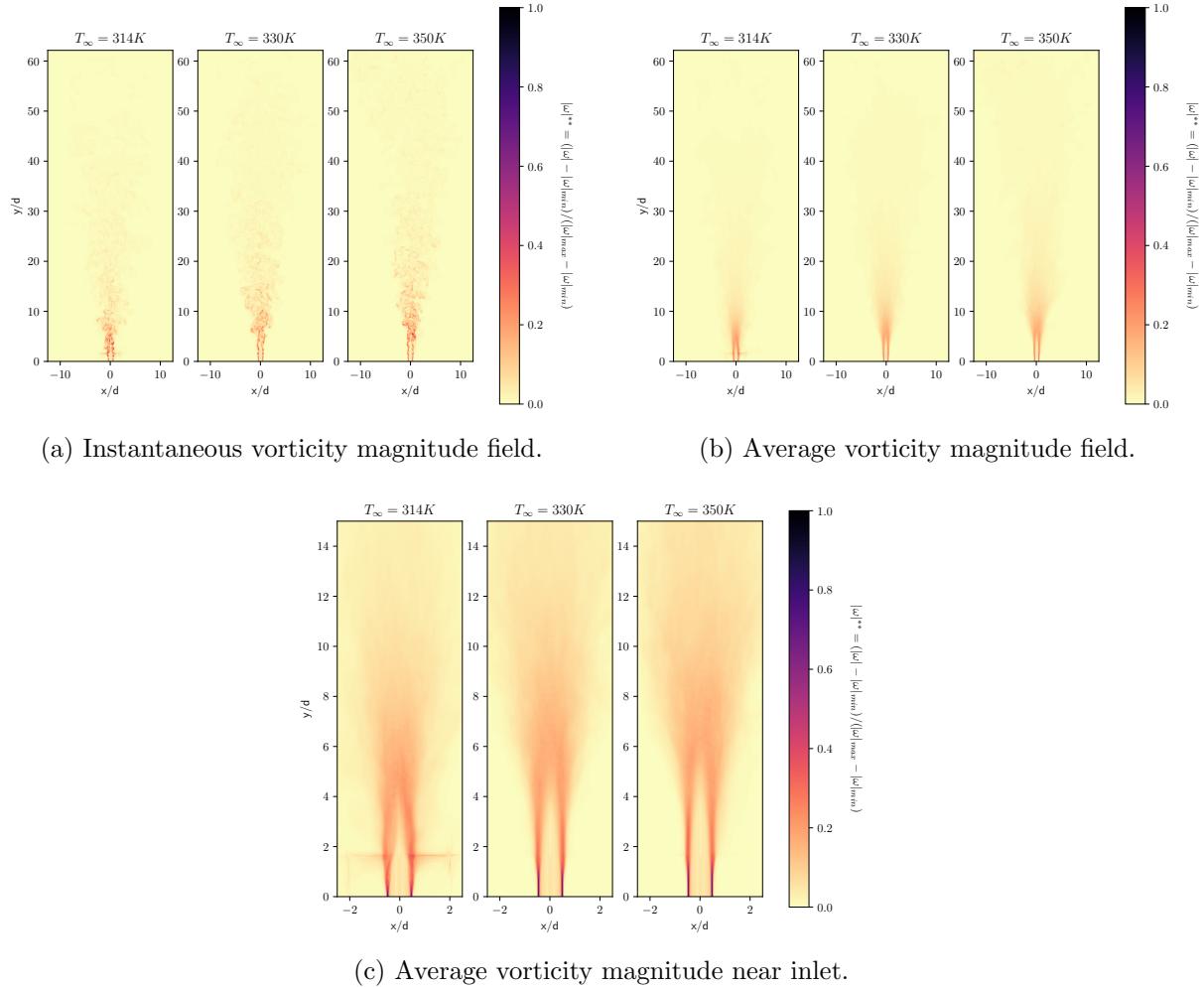


Figure 5.11: Vorticity magnitude field comparison between 314K ambient case (left), 330K ambient case (middle), and 350K ambient case (right).

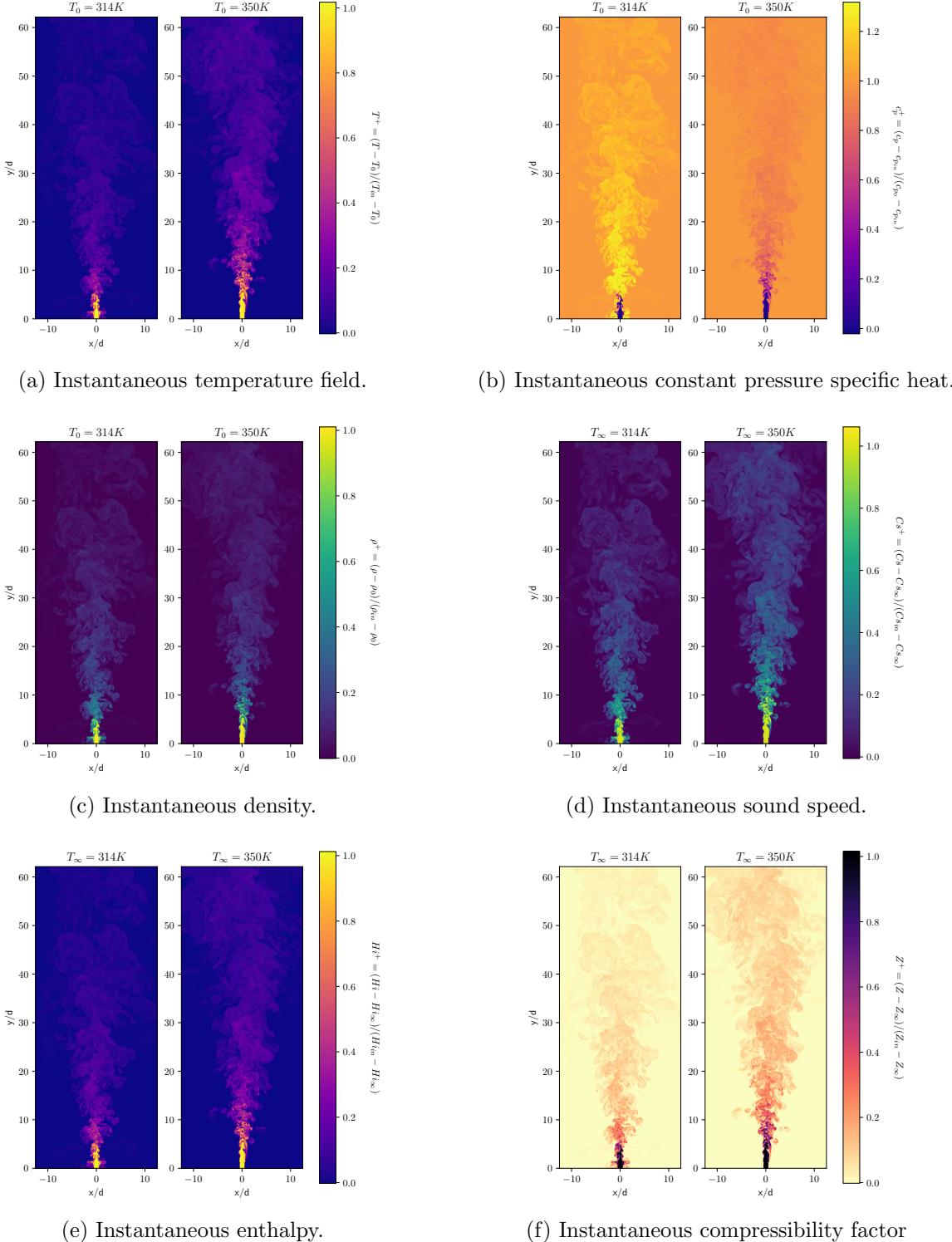


Figure 5.12: Direct comparison between 314K ambient case (left) and 350K ambient case (right) for various fluid quantities of interest.

### 5.2.2 Mean Flow Properties

Figure 5.13 depicts the time and radially averaged scaled axial velocity component plotted against radial distance from the centerline at multiple normal slices downstream from the inlet. The velocity is scaled by the average axial velocity value at the inflow while the radial direction is scaled by the jet diameter. In both cases, velocity profiles flatten as they get further downstream. Compared to the  $330K$  ambient case, the  $350K$  ambient case has a slower axial velocity decay downstream while the  $314K$  ambient case has a much faster decay. This agrees with the trends seen in the vertical slice images.

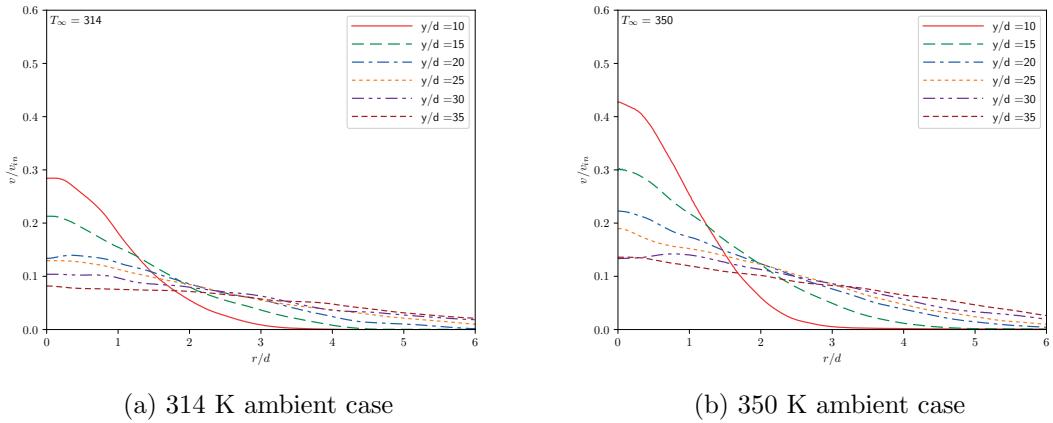


Figure 5.13: Average (both in time and radially) axial velocity scaled by inlet value plotted along radial distance from centerline.

Figures 5.14 shows a different scaling of the axial velocity profiles depicted in the previous figure. Here, the axial velocity is scaled by the centerline value at each normal slice while the radial distance is scaled by the  $r_{1/2}$ , as was done in Figure 5.5. Figures 5.14a and 5.14b show the near field profiles resulting from this scaling. As was the case with the  $330K$  ambient case, self-similarity is seen in the near field for the non-isothermal cases, where fairly agreeable profile collapse occurs at  $y/d = 15$ . The  $350K$  ambient case profiles are slightly more in-line with each other as compared to the  $314K$  ambient case, but both demonstrate a general tendency toward self-similarity still. Figures 5.14c and 5.14d however do not show as agreeable a collapse as the  $330K$  ambient case, though one still might be able to say the profiles tend toward self-similarity still. The  $314K$  ambient case has better self-similarity structure toward the centerline of the jet as compared to the  $350K$

ambient case, though the outer edge of the jet has a wider spread of profiles. The  $350K$  ambient case generally has a more uniform spread of profiles in both regions before and after the half-mean half-width (HMHW).

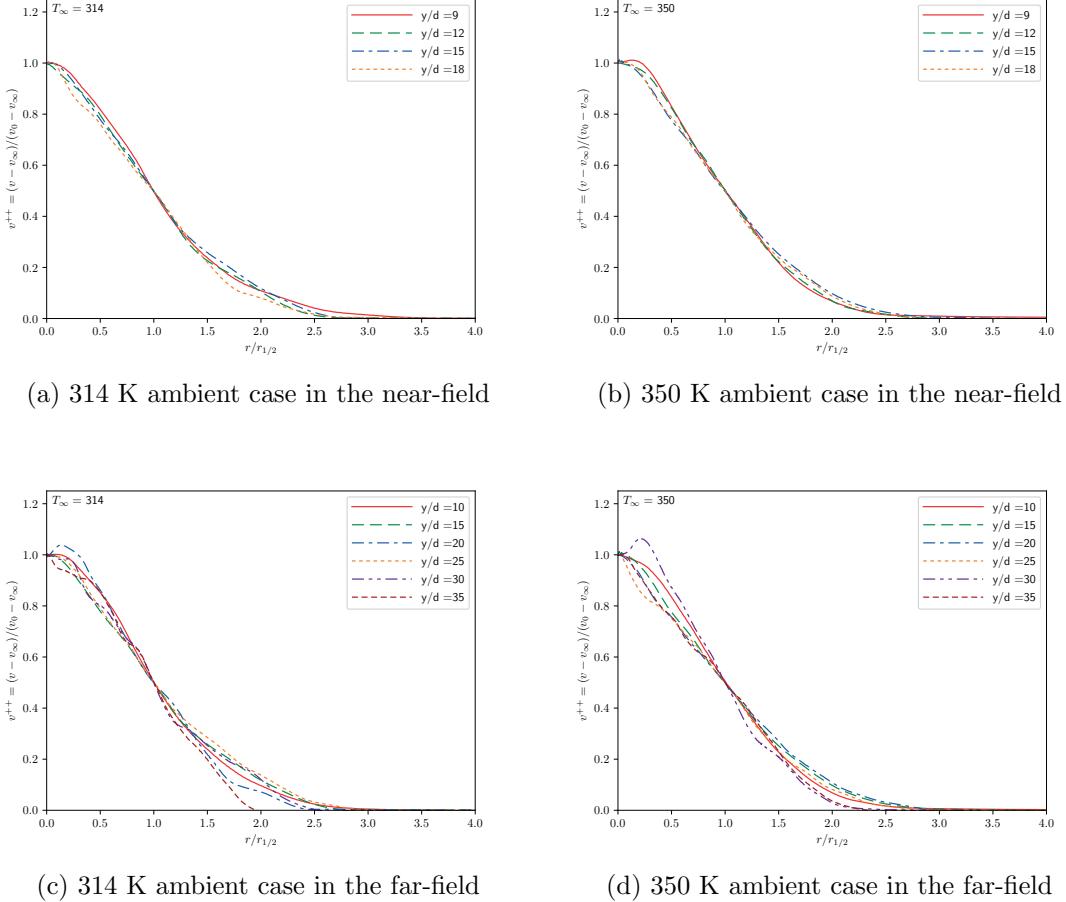


Figure 5.14: Normal slices of scaled near-field and far-field axial velocity profiles, averaged in both time and the radial direction. Plotted against radial direction scaled by  $r_{1/2}$ .

Density self-similarity is not as evident as is in axial velocity profiles, as can be seen in Figure 5.15. The  $350K$  ambient case shows more promise but the profile spread is not nearly as tight as it was with the axial velocity. The  $314K$  case is even worse, with more prominent fluctuations in the curves than any of the previous profile depictions. Far-field density in this case has the widest spread of all previous comparisons. One interesting feature of these plots can be seen on the outer edge of the jet, past the HMHW mark. For the  $314K$  ambient case, profiles exhibit an initial decay

as they dip below the  $y/d = 10$  profile before increasing to fall above this initial curve at later slices downstream. This is opposite to what is seen in the  $350K$  ambient case, where earlier slices rise above the initial curve at  $y/d = 10$  before decaying below it further downstream. This later decay is consistent with what is seen in the axial velocity self-similarity curves for both cases, while the density profiles for the  $314K$  ambient case break from this pattern.

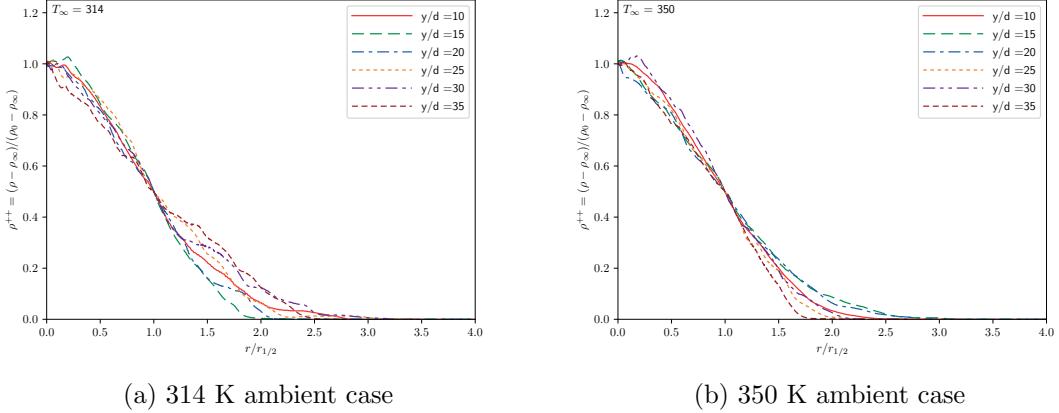


Figure 5.15: Normal slices of scaled far-field density, averaged in both time and the radial direction. Plotted against radial direction scaled by  $r_{1/2}$ .

Figure 5.16 shows the recovery of linear decay in the axial velocity along the centerline of the jet for each case. The  $350K$  ambient case has a less steep slope compared to the  $330K$  ambient case seen in Figure 5.6, corresponding to a slower decay in axial velocity along the centerline. The  $314K$  ambient case has a steeper slope compared to the  $330K$  ambient case, which corresponds to the centerline decay occurring at a faster rate than that in the base case. These findings are consistent with the analysis from Figure 5.13.

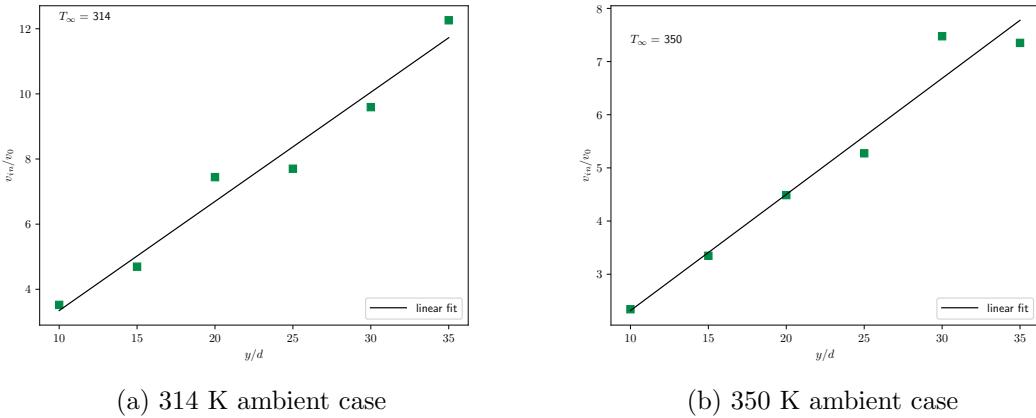


Figure 5.16: Axial inlet velocity scaled by centerline values along the axial direction. When distance downstream is scaled by jet diameter, linear decay of the centerline axial velocity is observed.

Finally, Figure 5.17 shows various quantities along the centerline for both non-isothermal cases. For all quantities presented here, centerline plots help depict the different regions of the jet. The potential core is maintained until around  $y/d = 2.5$  for the  $314K$  ambient case and until around  $y/d = 5$  for the  $350K$  ambient case. Then there is a transition region where quantities experience rapid change. This occurs up until about  $y/d = 10$  for the  $314K$  ambient case and  $y/d = 12$  for the  $350K$  ambient case. After that, a gradual leveling out occurs as the quantity of interest continues to adjust as it mixes with the ambient fluid through the remainder of the domain. Figure 5.17a shows the temperature change for each case along the centerline of the jet. The  $314K$  ambient case exhibits a more rapid decay from the inflow temperature as compared to the  $350K$  ambient case, as can be seen in the separation between the curves between  $4 \leq y/d \leq 16$ . Afterwards, they continue their respective transition toward ambient conditions at roughly the same rate. The  $350K$  ambient case never gets as close to ambient conditions as the  $314K$  ambient case due to the slower initial decline, hovering instead around  $T^+ = 0.18$  as compared to the  $T^+ = 0.14$  that the  $314K$  ambient case sits at. Enthalpy and the compressibility factor in Figures 5.17e and 5.17f respectively follow the same trend though with less separation between the two cases as compared to the temperature trajectories. Figure 5.17d shows the sound speed transition along the centerline. Compared to the previously discussed quantities, the sound speed decay appears much more gradual, with little rate difference between the two cases. The initial transition region is less steep and the region thereafter more of a continued gradual decline as opposed to a leveling out like the other quantities. Figure

[5.17b](#) shows the constant-pressure specific heat for the two non-isothermal cases. Again, neither fully transitions to the ambient condition as the jet persists through the domain. Here the jump in specific heat over the peak present near the pseudo-boiling point can be seen for the  $314K$  ambient case. This case also has a steeper initial transition than the other case does, though both have similar rates of approach toward ambient conditions after this initial shift. Figure [5.17c](#) shows the density of the two cases. This quantity follows a unique trajectory between the two cases compared to those previously discussed. Here, the initial transition rate of the two cases from the potential core to the fully developed region is nearly the same for the most part. Toward the end of this transition though, the  $314K$  ambient case begins to level out at a faster rate than the  $350K$  ambient case, allowing the curves to collapse onto each other. Then both centerline densities continue on past  $y/d = 16$  at approximately  $\rho^+ = .12$  and steadily decline toward ambient conditions thereafter.

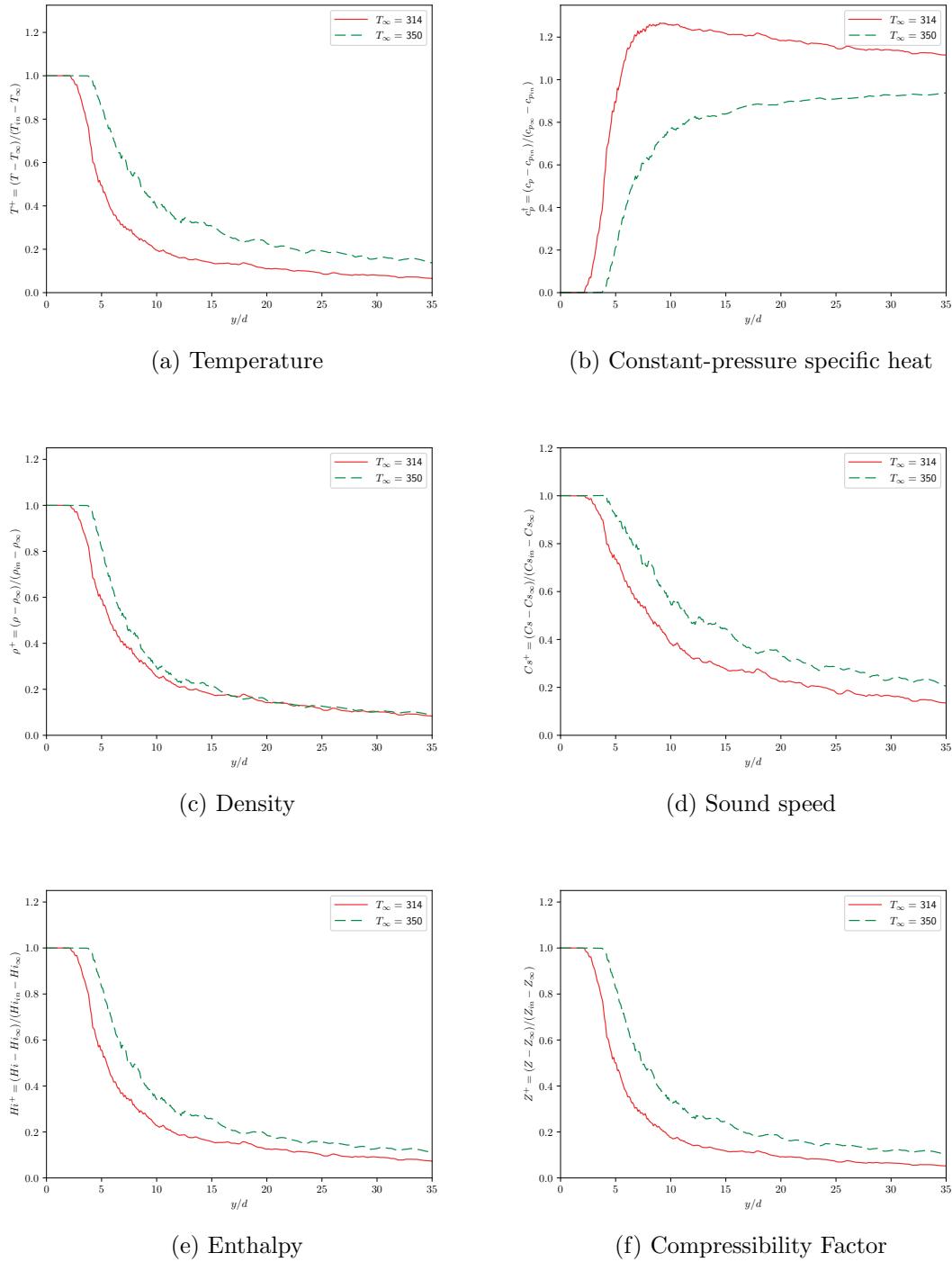


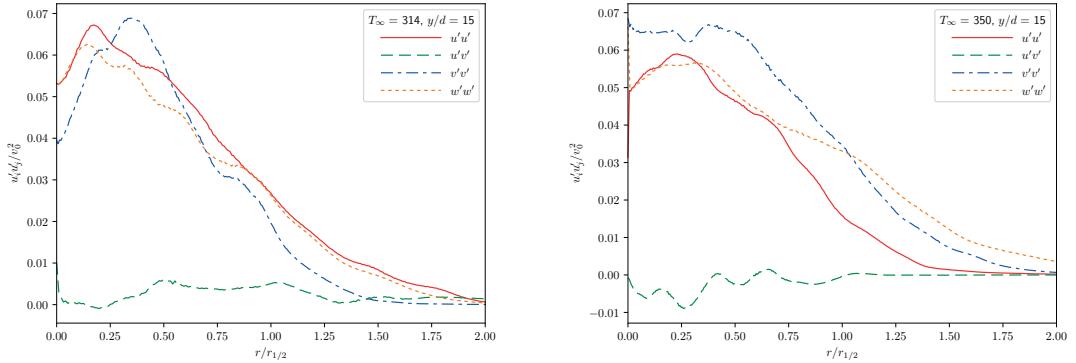
Figure 5.17: Comparison of centerline decay for various fluid quantities.

### 5.2.3 Turbulence Dynamics

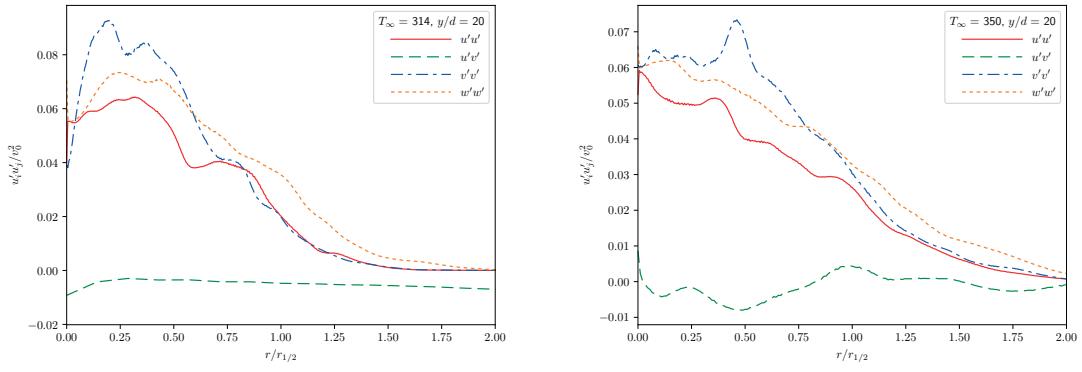
Figure 5.18 shows the resolved Reynolds stresses of each non-isothermal case at two different locations downstream. As was the case with the 330K ambient resolved Reynolds stresses in Figure 5.7, self-similarity in the axial direction is not recovered. The Reynolds stresses for both of the non-isothermal cases are generally smaller than their counterparts in the isothermal case. The stresses in Figure 5.18b generally follow the trends outlined in the isothermal case, but the remaining slices for these two cases do not follow suite, with little to no separation between the axial direction stresses and the other stresses. This could be due the effects of the SGS modeling as mentioned in [45].

Figure 5.19 shows the components of the resolved turbulent kinetic energy (TKE) for each non-isothermal case. Both cases follow the same general trends seen in the 330K ambient case, with the axial component being the strongest and the other two components being of the same magnitude. The 350K ambient case is the most similar to the 330K ambient case between the two. For this case, peak TKE occurs at approximately  $y/d = 7$ , which is slightly later than the isothermal case. Additionally, this peak is aligned with the peak seen in the other components, where for the isothermal case the axial component peak came slightly before the peak of the other two. Magnitude of the peak axial component is larger in the 350K ambient case while the magnitudes of the other directional components are similar to their counterparts in the 330K ambient case. Decay rate is also slightly higher in the 350K ambient case compared to the isothermal case. The TKE components of the 314K ambient case reach their peaks slightly earlier than the 330K ambient case at about  $y/d = 5$ . The peak of the axial component is much larger than that of the other two cases. Additionally, the other TKE components of the 314K ambient case have a smaller magnitude than their counterparts in the other cases, resulting in a larger disparity between the magnitudes of the components in this case. There is also stronger overall decay in the 314K ambient case as compared to the 350K ambient case.

Figure 5.20 further emphasizes the differences between the TKE components and overall TKE by directly comparing each quantity across all cases. For all quantities, the 314K ambient case peaks before the other two cases. For the overall resolved TKE and the axial component, the 330K ambient case is next to peak, followed by the 350K ambient case. For the cross directional components, the peaks of these two cases coincide. The 314K ambient cases experiences the strongest decay along the centerline immediately following the peak for each quantity, however, the



(a) Reynolds stresses at  $y/d = 15$  for 314 K ambient case    (b) Reynolds stresses at  $y/d = 15$  for 350 K ambient case



(c) Reynolds stresses at  $y/d = 20$  for 314 K ambient case    (d) Reynolds stresses at  $y/d = 20$  for 350 K ambient case

Figure 5.18: Time and radially averaged Reynolds stresses for the non-isothermal jets at two locations downstream.

other two cases eventually decay to the same value so that all cases exhibit a roughly equivalent leveling off as the jet progresses through the remainder of the domain.

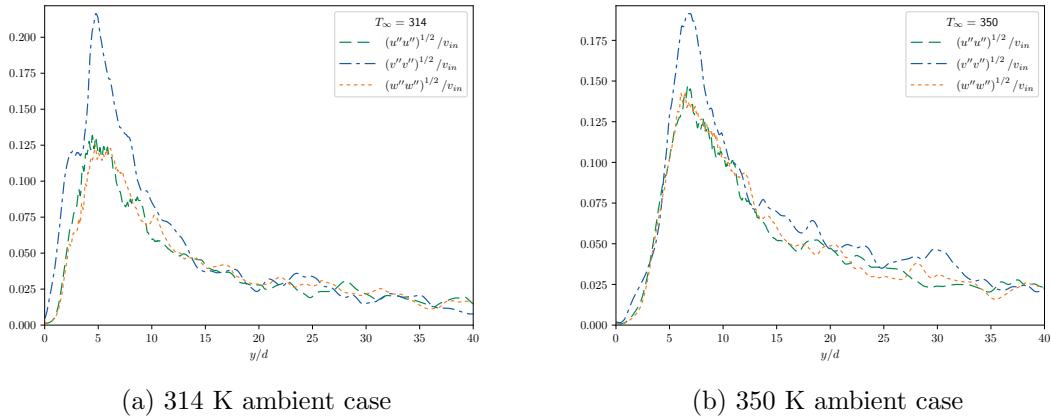


Figure 5.19: Comparison of average turbulent kinetic energy components along centerline for each non-isothermal case.

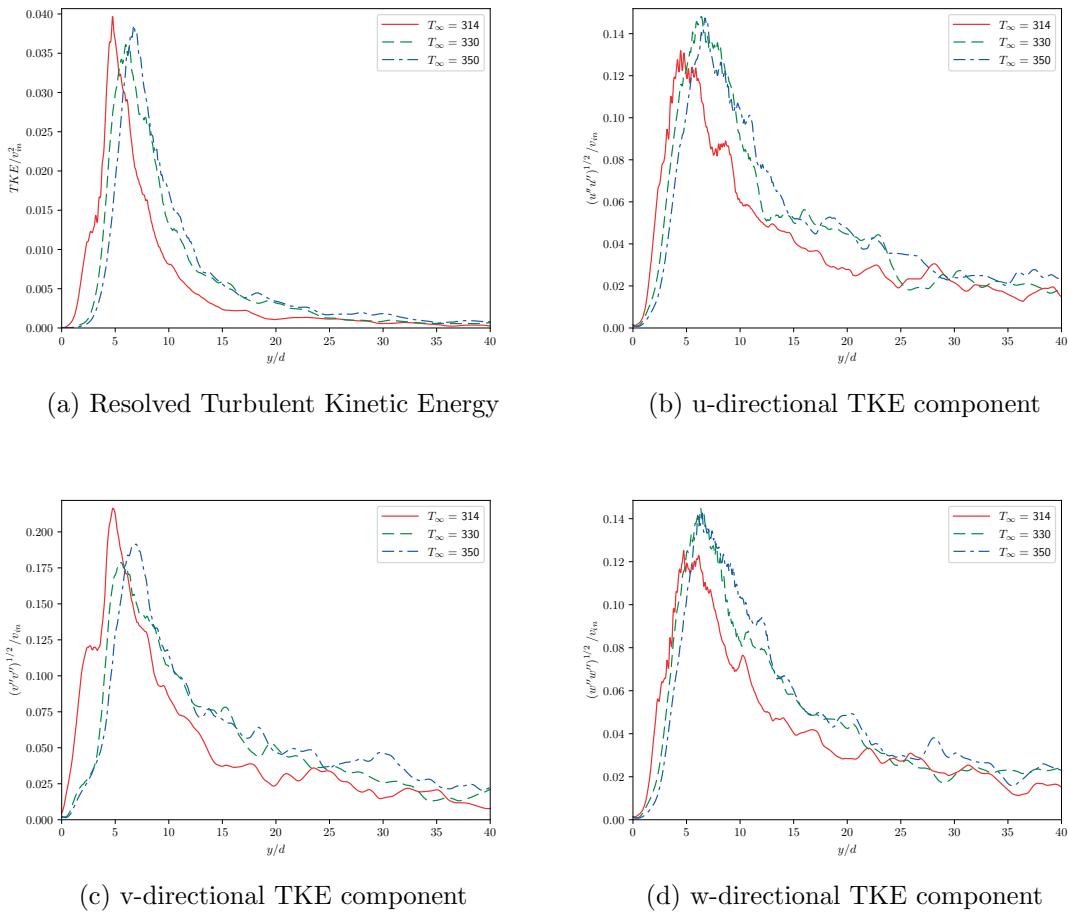


Figure 5.20: Cross-case comparisons of average turbulent kinetic energy components and total resolved TKE along centerline.

# CHAPTER 6

## CONCLUSION

In this work, we studied three cases of a supercritical round turbulent jet injected into a supercritical environment: an isothermal case, a non-isothermal case with ambient temperature further away from the critical point, and a non-isothermal case with ambient temperature closer to the critical point and specifically crossing a region of intense thermodynamic fluctuation. It was shown that the isothermal case had many similarities with ideal incompressible round jets. Many properties found in classical round turbulent jets were recovered in this case, such as self-similarity, linear decay along the centerline, and general trends associated with Reynolds stresses (although self-similarity in the Reynolds stresses was not recovered; general trends were still agreeable). The non-isothermal case further away from the critical point behaved similarly to the isothermal case with slight differences being noted in potential core length and spreading of the jet. The non-isothermal case that transited the pseudo-boiling point exhibited noticeably different behavior. These trends were compared with previous studies involving transcritical injection to see how phase lack of full phase change impacted flow characteristics.

Some limitations of this work are now noted. SGS models do not take into consideration real gas behavior, being limited to assumptions based on the ideal gas equation of state. There is room for improvement in turbulence modeling with the aim of incorporating higher order equations of state. As with any numerical simulation, discretization and resolution inherently lead to error. Ideally with more power, finer resolution can be achieved to help improve simulations. Finally, minor artifacts were noted in the non-isothermal case closest to the critical point due to the presence of steep gradients. Though the impact of these are believed to be mild, they are still present and add to cumulative error.

Future work aims to incorporate multiple species to the flow system and to investigate higher Mach flows. More specific application configurations can also be analyzed via incorporation of complex geometries.

# APPENDIX A

## TRANSPORT COEFFICIENTS

### A.1 EGLib Low Pressure Approximation Coefficients

Table A.1: Coefficients for EGLib Polynomials [19] used to model transport quantities as described in Equation 2.18

 $i$	 $a_{\mu,i}$	 $a_{\lambda,i}$
0	$-2.281\ 103\ 45 \times 10^1$	$-8.74831432$
1	$4.62954710$	$4.79275291$
2	$-5.006\ 890\ 01 \times 10^{-1}$	$-4.186\ 850\ 61 \times 10^{-1}$
3	$2.100\ 129\ 69 \times 10^{-2}$	$1.352\ 102\ 42 \times 10^{-2}$

### A.2 Chung High Pressure Corrections

Chung et al. [10] details the derivation of these general transport coefficients via their regression analysis of various fluids. The coefficients for the linear approximation are printed here for convenience as they are implemented in *PelePhysics* [].

Table A.2: Quantities used in part for Chung's High Pressure Corrections. Values gathered from EGLib approximations [19]

Transport constants	
 $M$	$4.400\ 995 \times 10^1$
 lenny-j. pot. well depth	$2.440\ 000 \times 10^2$
 lenny-j. collision diam.	3.763
 dipole moment	0.000
 polarizability	2.650
 rot. relax. coll. num.	2.100
 egtransetNLIN	1

Table A.3: Linear Coefficients Used in Calculating High Pressure Viscosity Corrections as found by Chung et al. [10]

$i$	$a_{i0}$	$a_{i1}$	$a_{i2}$	$a_{i3}$
1	6.32402	50.41190	-51.68010	1189.02000
2	$0.12102 \times 10^{-2}$	$-0.11536 \times 10^{-2}$	$-0.62571 \times 10^{-2}$	$0.37283 \times 10^{-1}$
3	5.28346	254.20900	-168.48100	3898.27000
4	6.62263	38.09570	-8.46414	31.41780
5	19.74540	7.63034	-14.35440	31.52670
6	-1.89992	-12.53670	4.98529	-18.15070
7	24.27450	3.44945	-11.29130	69.34660
8	0.79716	1.11764	$0.12348 \times 10^{-1}$	-4.11661
9	-0.23816	$0.67695 \times 10^{-1}$	-0.81630	4.02528
10	$0.68629 \times 10^{-1}$	0.34793	0.59256	-0.72663

Table A.4: Linear Coefficients Used in Calculating High Pressure Thermal Conductivity Corrections as found by Chung et al. [10]

$i$	$b_{i0}$	$b_{i1}$	$b_{i2}$	$b_{i3}$
1	2.41657	0.74824	-0.91858	121.72100
2	-0.50924	-1.50936	-49.99120	69.98340
3	6.61069	5.62073	64.75990	27.03890
4	14.54250	-8.91387	-5.63794	74.34350
5	0.79274	0.82019	-0.69369	6.31734
6	-5.86340	12.80050	9.58926	-65.52920
7	81.17100	114.15800	-60.84100	466.77500

## APPENDIX B

# ADDITIONAL THERMODYNAMIC QUANTITIES

### B.1 Ideal Gas NASA Polynomial Fits Coefficients

Table B.1: Coefficients for NASA polynomials [42] used in modeling ideal thermodynamic quantities as described in Equation 2.30

$i$	$c_{e_I,i}$	$c_{v_I,i}$	$c_{h_I,i}$
0	1.356 773 52	1.356 773 52	2.356 773 52
1	$4.492\ 298\ 39 \times 10^{-3}$	$8.984\ 596\ 77 \times 10^{-3}$	$4.492\ 298\ 39 \times 10^{-3}$
2	$-2.374\ 520\ 90 \times 10^{-6}$	$-7.123\ 562\ 69 \times 10^{-6}$	$-2.374\ 520\ 90 \times 10^{-6}$
3	$6.147\ 975\ 55 \times 10^{-10}$	$2.459\ 190\ 22 \times 10^{-9}$	$6.147\ 975\ 55 \times 10^{-10}$
4	$-2.873\ 990\ 96 \times 10^{-14}$	$-1.436\ 995\ 48 \times 10^{-13}$	$-2.873\ 990\ 96 \times 10^{-14}$

### B.2 SRK-Related Derivatives

Here are a few additional derivatives involving quantities related to the SRK EoS that are used at various points for the computations within *PelePhysics* []:

#### B.2.1 Constant-Pressure Specific Heat

$$c_p = \left( \frac{\partial h_m}{\partial T} \right)_p = \frac{\partial h_m}{\partial T} - \frac{\frac{\partial h}{\partial \tau}}{\frac{\partial p}{\partial \tau}} \frac{\partial p}{\partial T} \quad (\text{B.1})$$

where

$$\frac{\partial p}{\partial T} = \frac{R}{\tau - b_m} - \frac{\partial a_m}{\partial T} \frac{1}{\tau(\tau + b_m)} \quad (\text{B.2})$$

$$\frac{\partial p}{\partial \tau} = -\frac{RT}{(\tau - b_m)^2} + \frac{a_m(2\tau + b_m)}{[\tau(\tau + b_m)]^2} \quad (\text{B.3})$$

$$\frac{\partial h_m}{\partial \tau} = -\left( T \frac{\partial a_m}{\partial T} - a_m \right) \frac{1}{\tau(\tau + b_m)} + \frac{a_m}{(\tau + b_m)^2} - \frac{RTb_m}{(\tau - b_m)^2} \quad (\text{B.4})$$

$$\frac{\partial h_m}{\partial T} = c_p^{id} + T \frac{\partial^2 a_m}{\partial T^2} \frac{1}{b_m} \ln\left(1 + \frac{b_m}{\tau}\right) - \frac{\partial a_m}{\partial T} \frac{1}{\tau + b_m} + \frac{Rb_m}{\tau - b_m} \quad (\text{B.5})$$

### B.2.2 Sound Speed

$$a^2 = -\frac{c_p}{c_v} \tau^2 \frac{\partial p}{\partial \tau} \quad (\text{B.6})$$

### B.2.3 Primitive Derivatives Needed for Riemann Solver

$$\left. \frac{\partial p}{\partial \rho} \right|_{p,Y} = -\tau^2 \left( \frac{\partial p}{\partial \tau} - \frac{\frac{\partial e}{\partial \tau}}{\frac{\partial e}{\partial T}} \frac{\partial p}{\partial T} \right) \quad (\text{B.7})$$

$$\left. \frac{\partial p}{\partial e} \right|_{\rho,Y} = \frac{1}{c_v} \frac{\partial p}{\partial T} \quad (\text{B.8})$$

$$\frac{\partial e}{\partial \tau} = \frac{1}{\tau(\tau + b_m)} \left( a_m - T \frac{\partial a_m}{\partial T} \right) \quad (\text{B.9})$$

# APPENDIX C

## PELEC INPUT PARAMETERS FOR SIMULATIONS



Here inputs are provided to aid in reproducibility of the present simulations. Note though that the current version of *PeleC* may have different input structures compared to those displayed here.

```

File Edit Options Buffers Tools Help
# ----- INPUTS TO MAIN PROGRAM -----
#max_step = 10
max_step = 100000000
stop_time = 0.003472222 # 10 flow through times (domain height / u_jet) * 10

# PROBLEM SIZE & GEOMETRY
geometry.is_periodic = 0 0 0
geometry.coord_sys   = 0          # 0 => cart
geometry.prob_lo     = -0.125  0.0 -0.125
geometry.prob_hi     = 0.125   0.625 0.125
amr.n_cell           = 80 200 80

# >>>>>>>>> BC KEYWORDS <<<<<<<<<<<<
# Interior, UserBC, Symmetry, SlipWall, NoSlipWall
# >>>>>>>>> BC KEYWORDS <<<<<<<<<<<<
pelec.lo_bc      = "FOExtrap" "Hard" "FOExtrap"
pelec.hi_bc      = "FOExtrap" "FOExtrap" "FOExtrap"

# WHICH PHYSICS
pelec.do_hydro = 1
pelec.do_mol_AD = 0
pelec.nscbc_adv = 0
pelec.do_react = 0
pelec.ppm_type = 2
pelec.allow_negative_energy = 0
pelec.diffuse_temp = 1
pelec.diffuse_yel = 1
pelec.diffuse_spec = 0
pelec.diffuse_enth = 0
pelec.do_les = 1
pelec.les_model = 1

# DYNAMIC SMAGORINSKY
pelec.les_test_filter_type = 3
pelec.les_test_filter_fgr = 2

# TIME STEP CONTROL
pelec.dt_cutoff    = 5.e-20 # level 0 timestep below which we halt
pelec.cfl          = 0.9   # cfl number for hyperbolic system
pelec.init_shrink  = 0.1    # scale back initial timestep
pelec.change_max   = 1.05   # maximum increase in dt over successive steps

# DIAGNOSTICS & VERBOSITY
pelec.sum_interval = 1       # timesteps between computing mass
pelec.v            = 1       # verbosity in PeleC cpp files
amr.v             = 1       # verbosity in Amr.cpp
#amr.grid_log      = grdlog # name of grid logging file

# REFINEMENT / REGRIDDING
amr.max_level     = 4       # maximum level number allowed
amr.ref_ratio     = 2 2 2 2 # refinement ratio
amr.regrid_int    = 10      # how often to regrid
amr.blocking_factor = 8      # block factor in grid generation
amr.max_grid_size = 16

# CHECKPOINT FILES
amr.check_file    = chk      # root name of checkpoint file
amr.check_int     = 0.0000034722222 # chk with same frequency as plots to reduce repetition in calculations

```

Figure C.1: Input parameters for available physics options in PeleC

```
&tagging
  denerr = 1e20
  dengrad = 1e20
  max_denerr_lev = 3
  max_dengrad_lev = 3

  presserr = 1e20
  pressgrad = 1e20
  max_presserr_lev = 3
  max_pressgrad_lev = 3

  velerr = 1e20
  velgrad = 1e20
  max_velerr_lev = 3
  max_velgrad_lev = 3

  temperr = 1e20
  tempgrad = 4.0
  max_temperr_lev = 3
  max_tempgrad_lev = 3

  vorterr = 5000
  max_vorterr_lev = 3

  vfracerr = 1e20
  max_vfracerr_lev = 5
  /
&extern
!
! Specific heat ratio for air
!
! eos_gamma = 1.4
!
/
```

Figure C.2: AMR tagging criteria selected within in PeleC

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## **BIOGRAPHICAL SKETCH**

Julia Ream was born and raised in Sarasota, FL. From an early age, Julia found math to be “fun”, but didn’t come to realize how uncommon that sentiment was until much later on in life. She started her undergraduate career at Florida State University in 2013, enrolling as an Exploratory student before committing to the Humanities and later on Theatre. Upon realizing that she missed doing math (weird), she added Applied Mathematics to her roster of majors. Although it took nearly 21 years, Julia eventually realized that she actually really liked math. So, after completing a Bachelor of Arts in Humanities and a Bachelor of Science in Mathematics, she decided to pursue her Doctoral degree in Applied and Computational Mathematics, also at Florida State University.