

On the long-time transient formation of sink zones in near-critical fluids. A theoretical perspective

Yash Kulkarni¹, Roger Khayat² and Sakir Amiroudine^{1,†}

¹University of Bordeaux, CNRS, I2M Bordeaux, Talence F-33400, France

²Mechanical and Materials Engineering, Western University, Canada

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Near-critical fluids subject to simultaneous thermal quench and an imposed external acceleration field are reported to develop ‘sink zones’, where the temperature in the bulk falls below the imposed boundary value. This anomalous cooling effect persists for a long period of time corresponding to the diffusive elimination of the cold boundary layer. The sink-zone phenomenon has been captured previously in numerical simulations (Zappoli *et al.*, *J. Fluid Mech.*, vol. 316, 1996, pp. 53–72; Sharma *et al.*, *Phys. Rev E*, vol. 96, 2017, 063102; Sharma *et al.*, *Phys. Fluids*, vol. 29, 2017, 126103) and was observed experimentally (Beysens *et al.*, *Phys. Rev. E*, vol. 84, 2011, 051201). The present work provides a detailed and thorough theoretical analysis and interpretation based on matched asymptotic expansions. By examining the one-dimensional transient flow and temperature with a mean-field approach, we provide further insight into this striking phenomenon, showing that the behaviour involved is unique to highly compressible near-critical van der Waals gases. We also show that the sub-cooling phenomenon cannot be predicted in perfect gases unless very extreme conditions are applied.

Key words: general fluid mechanics, compressible boundary layers, boundary layer stability

1. Introduction

Supercritical fluids (SCFs) are a class of fluids that exist beyond the liquid–vapour critical point and may be considered as an intermediate state between a liquid and a gas, exhibiting some properties of liquids, like high density and solubility, and some properties of gases, such as low viscosity and high compressibility (Stanley 1971). On approaching the critical point, the various thermo-physical properties show a singular behaviour, such as diverging isothermal compressibility χ_T , heat conductivity λ , specific heat at constant pressure C_p

† Email address for correspondence: sakir.amiroudine@u-bordeaux.fr

and constant volume C_V and a vanishing thermal diffusivity D_T . These properties have been found to behave according to universal power laws with a different critical exponent for each property. They can be modelled on the basis of a common parameter $\mu = (T_i - T_c)/T_c$, where T_c refers to the critical temperature and T_i is the initial temperature of the fluid (Stanley 1971).

Regarding the thermal behaviour of SCFs, it was initially believed that the vanishing thermal diffusivity would have an impeding effect on the thermalization of a cell filled with a SCF when subjected to a thermal heating at the boundary. However, Nitsche & Straub (1987) observed that the thermal relaxation in the fluid cell containing supercritical SF₆ was completed in seconds as opposed to the predicted scale of days based on the thermal diffusion. This phenomenon of a critical acceleration of heating was termed the piston effect, attributed to the high compressibility near the critical point (Boukari *et al.* 1990; Onuki & Ferrell 1990; Zappoli *et al.* 1990). This phenomenon may be explained by considering a cell filled entirely with a SCF and exposed to a small temperature increase at the boundary. A very thin thermal boundary layer is formed due to the vanishing thermal diffusivity. The high-density gradients in the thermal boundary layer, in conjunction with a large isothermal compressibility in the bulk, compress the bulk fluid adiabatically. Consequently, a wave propagates throughout the bulk fluid at the speed of sound, which converts the mechanical energy of the waves into thermal energy, yielding an increase in the temperature of the bulk. The time scale on which this thermal homogenization in the bulk occurs is very small as compared to the diffusion time scale. A similar argument can be applied when the bounding wall is cooled, where the fluid in the thermal boundary layer contracts, and adiabatic cooling thermalizes the bulk fluid. This occurrence has also been termed the piston effect owing to its resemblance to a piston which compresses a gas in a closed cylinder and is the underlying phenomenon for understanding the physics governing near-critical fluids.

Carlès & Zappoli (1995) analysed the problem of a container filled with a near-critical fluid when subjected to mechanical vibrations by means of matched asymptotic expansions. Among the various characteristic regimes defined for fluid vibration, a specific low-frequency regime was highlighted wherein the bulk part of the fluid almost behaves like a solid bouncing back and forth between two highly compressible thermal boundary layers. This problem was solved with the temperatures at the walls being identical to the initial temperature (no heat or thermal quench was applied).

Due to the anomalous behaviour of the properties of near-critical fluids, these fluids have been known to exhibit various intriguing flow phenomena and instabilities when subject to thermal perturbations (Zappoli *et al.* 1996; Accary *et al.* 2005*a,b*) and simultaneous thermal quench and vibration (Amiroudine & Beysens 2008; Zappoli, Beysens & Garrabos 2016; Sharma, Erriguible & Amiroudine 2017*a,b*). A striking phenomenon, wherein the temperature of the SCF in the bulk drops below the imposed boundary temperature upon simultaneous thermal quench and vibration in zero-gravity conditions, was previously investigated numerically (Sharma, Erriguible & Amiroudine 2017*a*). This subcooling phenomenon was also observed in experiments performed under microgravity in low Earth orbit by using SF₆ at liquid-like density (Beysens, Fröhlich & Garrabos 2011).

We identify subcooled regions as sink zones, as they act like a sink for the heat within the fluid domain. Although this phenomenon was clearly observed and described numerically and in experiments, no theoretical proof has been proposed so far. In order to explain this anomalous behaviour, a transient one-dimensional matched asymptotic analysis is performed based on the initial work of Carlès & Zappoli (1995) but in the case of non-homogeneous thermal boundary conditions. It is indeed observed that the

effect of high compressibility along with the action of self-weight (due to the high acceleration) causes the temperature to change in the bulk in addition to the usual action of the piston effect. This subsequently affects the overall temperature profile thereby leading to the formation of sink zones. The explanation for the one-dimensional case can be judiciously extended to explain the sink-zone appearance in the two-dimensional numerical simulations of Sharma, Erriguible & Amiroudine (2017*a,b*). Furthermore, the appearance of sink zones is found to depend on several factors such as proximity to the critical point and acceleration.

The subject of simultaneous thermal perturbations and mechanical vibrations of SCFs have caught the attention of several researchers (Garrabos *et al.* 2007; Amiroudine & Beysens 2008; Gandikota *et al.* 2013; Lyubimova *et al.* 2019). This was primarily attributed to the intriguing flow features arising from the complex thermomechanical coupling in such fluids. Furthermore, these have also been motivated by their direct implications, for example, for g-jitters experienced by cryogenic storage reservoirs in zero gravity in space technology. Experiments and numerical simulations performed with simultaneous thermal quenches and mechanical vibrations with supercritical CO₂ and SF₆ under weightlessness suggest evidence of various types of instabilities (Wunenburger *et al.* 1999; Garrabos *et al.* 2007). The phenomenon described by Carlès & Zappoli (1995), wherein the bulk part of the fluid almost behaves like a solid, bouncing back and forth between two highly compressible thermal boundary layers, was further corroborated by the numerical studies of Jounet *et al.* (1999). Amiroudine & Beysens (2008) investigated this problem with a linearized equation of state for supercritical CO₂, and observed finger-like patterns which were attributed to thermo-vibrational instability. The stability analysis using the same model was further extended by Gandikota *et al.* (2013), wherein different types of instabilities were quantified in terms of the amplitude and frequency of the vibration. In addition, a stability diagram was reported, illustrating the critical amplitude of all instabilities (corner type, parametric, Rayleigh vibrational), providing an insight into the kind of instabilities one may expect at a given frequency and proximity to the critical point. The anomalous behaviour of the thermophysical properties in SCFs has also led to several unique observations such as the drop in temperature near the wall below the initial temperature imposed at the wall (see next).

Indeed, an unconventional cooling was observed by Beysens, Fröhlich & Garrabos (2011) in their experimental work. An interferometer cell filled with a SCF was subject to a heat pulse with boundary temperatures held constant (see figure 1*a*), and, at the end of the heat pulse, the bulk fluid was observed to cool below the initial temperature during a long thermal transient (see figure 1*b*). The heating and cooling times are comparable and of the order of the piston-effect time (≈ 1.3 s). In contrast, the time to return to the final equilibrium temperature corresponds to the diffusion time (≈ 630 s). This clearly indicates that the cooling process indeed corresponds to an isentropic piston effect and the return to equilibrium to a diffusive process.

A nearly similar observation was also reported by Zappoli *et al.* (1996) when investigating the mechanism of heat and mass transport in a side-heated square cavity in the presence of gravity. The temperature in their case was found to rise above the wall temperature near the corners. Very recently, Sharma, Erriguible & Amiroudine (2017*a,b*) analysed numerically the behaviour of a SCF under the combined effect of thermal perturbation and mechanical vibration. They considered a square cell (size $h = 7$ mm) with solid walls filled with supercritical H₂. The vertical walls were adiabatic and the top and bottom walls were thermally quenched. The initial temperature T_i was set 100 mK above the critical temperature, i.e. $T_i - T_c = 100$ mK and the amplitude of the thermal quench at the walls was $\delta T = 10$ mK. The cell was submitted to a vibration of an amplitude

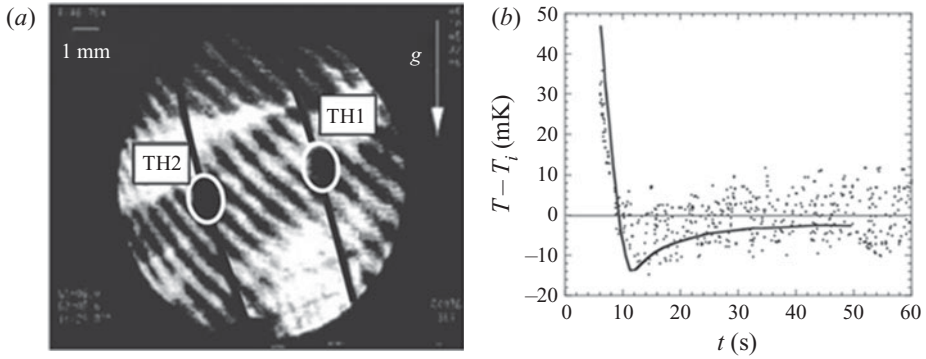


Figure 1. (a) The interferometer cell with two 0.89 mm diameter thermistors used in Earth gravity experiments in a homogeneous state. Thermistors TH1 and TH2 are identified by black dots. TH1 is used for heating. (b) Time response of thermistor TH2 (dots) during relaxation after a 100 mW, 4 s heat pulse on Earth (TH1 exhibits a similar response). A temperature undershoot below the thermostat temperature is observed. The line is the corresponding one-dimensional simulation ($T_i = T_c + 16.5$ K). Reprinted figure from: Beysens, Fröhlich & Garrabos (2011), with permission from APS (2020).

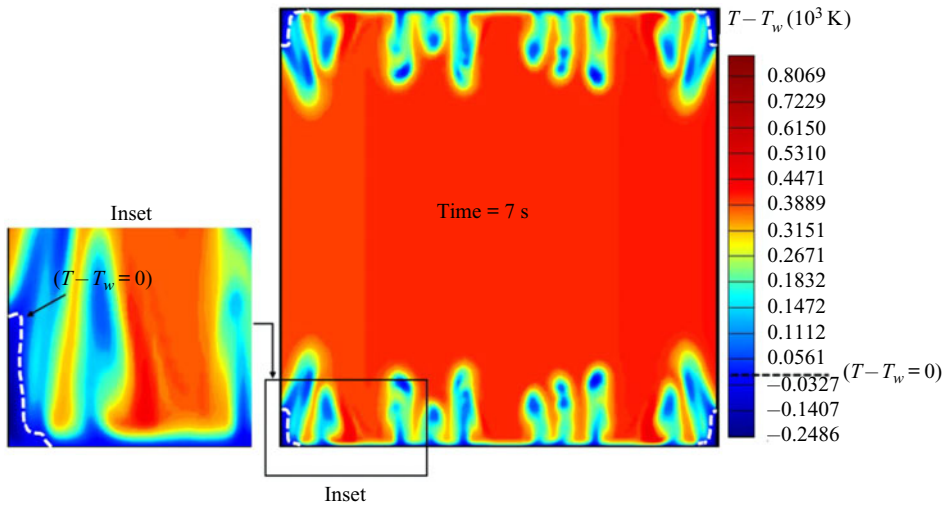


Figure 2. The $T - T_w$ contour plots for ($T_i - T_c = 100$ mK), $\delta T = 10$ mK, $f = 20$ Hz and $A = 5$ mm at time $t = 7$ s. Reprinted figure from: Sharma, Erriguible & Amiroudine (2017a,b) with permission obtained from APS (2020).

$A = 5$ mm and a frequency $f = 20$ Hz. Figure 2 shows the temperature contour plots $T - T_w$ where $T_w = T_i - \delta T$ corresponds to the wall temperature. The inset in figure 2 highlights the sink zone corresponding to $T - T_w < 0$.

2. The physical domain and problem formulation for SCFs

2.1. Statement of the problem

In this study, the existence of the observed sink zone is analysed quantitatively and demonstrated theoretically using a one-dimensional model with mean-field properties for a van der Waals gas. To our knowledge, and despite the reported experimental and numerical evidence of the sink-zone phenomenon, no theoretical proof has been proposed so far in

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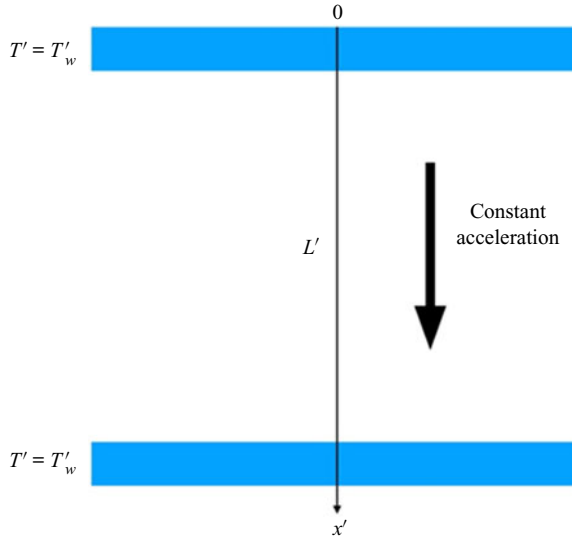


Figure 3. Sketch of the one-dimensional domain; x' varies from 0 to L' and the walls are maintained at a constant temperature T'_w such that $T'_i > T'_w > T'_c$ for all time $t' > 0$; T'_i represents the initial temperature of the bulk fluid. Both walls are thermally quenched.

the literature. The objective of the present work is to prove mathematically the appearance of the observed sink zone and elucidate the theoretical and physical mechanisms behind this phenomenon.

Consider the non-isothermal response of the one-dimensional transient flow and heat transfer of a near-critical fluid subject to a constant volume force or acceleration. The physical domain is schematically illustrated in figure 3. In what follows, a prime denotes dimensional parameters and variables, the subscript '0' refers to a perfect gas, the subscript 'i' refers to initial conditions and a subscript 'w' refers to conditions at the wall. The cell is subject to a thermal quench $\delta T'$ at both walls, i.e. $T'_w = T'_i - \delta T'$. The length of the cavity is taken as L' and the constant acceleration is imposed in the downward x' -direction.

2.2. Problem formulation

The model adopted here is an extension of the one used by Carlès & Zappoli (1995). The thermostated container is filled with a near-critical van der Waals fluid at critical density, initially at rest, in thermal equilibrium with the walls. A constant external acceleration field is imposed on the fluid cell, and the walls are thermally quenched and maintained at a value $T'_w = T'_i - \delta T'$ for all time $t' > 0$. In addition, the no-penetration condition is imposed at the walls. The boundary conditions at $x' = 0$ and $x' = L'$ are written explicitly as

$$T'(x' = 0, t') = T'(x' = L', t') = \begin{cases} T'_i, & t' = 0, \\ T'_w, & t' > 0. \end{cases} \quad (2.1a)$$

$$u'(x' = 0, t') = u'(x' = L', t') = 0. \quad (2.1b)$$

The imposed acceleration field is given by $A'H(t')$, where A' is the constant amplitude, and

$$H(t') = \begin{cases} 0, & t' < 0 \\ 1, & t' \geq 0 \end{cases} \quad (2.2)$$

is the Heaviside step function. We have purposely chosen $A' = 10\text{ g}$ (where $g = 9.81\text{ m s}^{-2}$ is the gravity) in the present study since the main idea here is to explain theoretically the emergence of the sink zone, which was reported in an earlier experiment (Beysens, Fröhlich & Garrabos 2011) and numerical simulation (Sharma, Erriguible & Amiroudine 2017a), and not necessarily to compare quantitatively against earlier results. Our present model follows closely the one adopted in the numerical simulation. Sharma, Erriguible & Amiroudine (2017a) considered a time-varying acceleration in a zero- g environment, namely $A'\omega'^2\sin(\omega't')$ (in their expression, A' represents the amplitude of displacement and ω' is the imposed frequency). But, in an effort to explain the observed sink zone, they also focused their study on a constant acceleration, $A' = 100\text{ g}$, as reported in their figure 6. The 10 g acceleration, which is different from the 100 g in Sharma, Erriguible & Amiroudine (2017a), is arbitrarily chosen in order to amplify the magnitude of the sink zone since it seems to clearly illustrate the interplay between the acceleration and the imposed thermal quench at the wall. Simultaneously, the constant 10 g acceleration could also mimic the 1 g terrestrial environment as in the experiment of Beysens, Fröhlich & Garrabos (2011).

The thermal conductivity λ' of near-critical fluids also diverges. This fact is taken into account by the introduction of the following temperature dependence:

$$\frac{\lambda'}{\lambda'_0} = 1 + \Lambda \left[\frac{T' - T'_c}{T'_c} \right]^{-1/2}, \quad (2.3)$$

where λ'_0 is the thermal conductivity of a perfect gas ($\Lambda = 0.75$ for CO_2). The real exponent, 0.57 , is slightly superior to 0.5 , which is a value based on the mean-field theory (see, for instance, Zappoli, Beysens & Garrabos 2016). However, the 0.5 exponent is convenient for our theoretical formulation and analysis, and should not lead to a different phenomenological interpretation.

We use the following non-dimensional variables for the density, temperature, pressure, velocity, time and position, respectively,

$$\rho = \frac{\rho'}{\rho'_c}, \quad T = \frac{T'}{T'_c}, \quad P = \frac{P'}{\rho'_c r' T'_c}, \quad u = \frac{u'}{\sqrt{\gamma_0 r' T'_c}}, \quad t = \frac{t'}{t'_a}, \quad x = \frac{x'}{L'}, \quad (2.4a-f)$$

where T'_c and ρ'_c are the critical values of temperature and density, respectively, and $t'_a = L'/\sqrt{\gamma_0 r' T'_c}$ is the acoustic time for a perfect gas. Here γ_0 is the specific heat ratio for a perfect gas, and r' is the universal gas constant divided by the molar mass for a perfect gas.

The equation of state used here is for a van der Waals fluid, which has been proven to model correctly the fluid flow near the critical point, in numerical simulations (Zappoli *et al.* 1996; Accary *et al.* 2005a,b) and asymptotic analyses (Peng & Robinson 1976; Carlès & Zappoli 1995). The van der Waals equation of state does not necessarily yield the exact real critical exponents. However, both numerical simulations (Zappoli *et al.* 1990, 1996) and asymptotic analyses (Ferrel & Hong 1993; Zappoli & Carlès 1995) have already demonstrated that a van der Waals equation of state can provide a wealth of phenomenological insight, while enabling a relatively simple mathematical formulation. This choice was therefore made here to facilitate the formal analysis of the process of

heat sink rather than to provide a quantitative comparison with real measurements. If a quantitative analysis is needed, one could use more sophisticated state equations such as the Peng–Robinson equation (Peng & Robinson 1976) and the NIST database or real exponents (Ferrel & Hong 1993).

We denote the kinematic viscosity and the thermal diffusivity for a perfect gas, by ν'_0 and κ'_0 , respectively. The introduction of the non-dimensional variables leads to the following equations for mass, momentum and energy conservation, as well as the van der Waal equation of state

$$\left. \begin{aligned} \rho_t + (\rho u)_x &= 0, \\ \rho(u_t + uu_x) &= -\frac{1}{\gamma_0} p_x + \frac{4}{3} \varepsilon u_{xx} + \rho A H(t), \\ \rho(T_t + uT_x) &= -(\gamma_0 - 1) \left(p + \frac{9}{8} \rho^2 \right) u_x + \frac{4}{3} \varepsilon \gamma_0 (\gamma_0 - 1) u_x^2 \\ &\quad + \varepsilon \left[\frac{\gamma_0}{Pr_0} \left(1 + \frac{\Lambda}{\sqrt{T-1}} \right) T_{xx} - \frac{\Lambda}{2} T_x^2 (T-1)^{-3/2} \right], \\ P &= \frac{\rho T}{1 - \rho/3} - \frac{9}{8} \rho^2. \end{aligned} \right\} \quad (2.5)$$

Here, $\varepsilon = (t'_a/t'_d)Pr_0$, where $t'_d = L'^2/\kappa'_0$ is the characteristic diffusion time, $Pr_0 = \nu'_0/\kappa'_0$ is the reference Prandtl number for a perfect gas and $\mu = (T'_i - T'_c/T'_c)$ is a measure of the departure of the initial temperature from the critical temperature. We note that both ε and μ are typically very small. For a 10 mm cell filled with CO₂, $\varepsilon \approx 10^{-8}$. However, μ cannot be less than 10^{-6} for the continuum hypothesis to hold. Finally, $A = A'L'/\gamma_0 r'T'_c$ is the non-dimensional imposed acceleration. For the CO₂ cell, $A = 1.22 \times 10^{-5}$ which corresponds to a dimensional acceleration of $A' = 10$ g.

Initially, the fluid is assumed to be at rest, and (2.5) are solved subject to the following initial and boundary conditions:

$$\left. \begin{aligned} \rho(x, t = 0) &= 1, \quad u(x, t = 0) = 0, \\ T(x, t = 0) &\equiv T_i = 1 + \mu, \quad P(x, t = 0) \equiv P_i = \frac{3}{2}(1 + \mu) - \frac{9}{8}, \\ u(x = 0, t) &= u(x = 1, t) = 0, \quad T(x = 0, t) = T(x = 1, t) = T_w. \end{aligned} \right\} \quad (2.6)$$

(2.7a,b)

Here, T_w represents the dimensionless wall temperature at time $t > 0$. When cast in terms of the dimensionless thermal quench $\delta T = aA$, it reads

$$T_w(t) = T_i - \delta TH(t). \quad (2.8)$$

In this study, we assume the dimensionless thermal quench and imposed acceleration to be of the same order of magnitude so that a remains of order one. For a thermal quench of 10 mK, $a = 2.7$ for the CO₂ cell described above.

We emphasize that the aim of the present one-dimensional analysis is to predict the non-isothermal transient response of a slab shaped container subject to a volume force acceleration (see figure 3). In the study of Carlès & Zappoli (1995), the walls were not thermally quenched but rather maintained at the same initial temperature of the fluid. Additionally, in the present work, a constant instead of a periodic acceleration field is imposed in order to demonstrate the appearance of sink zones in the bulk near the

boundary. It is worth noting that the analysis carried out by Carlès & Zappoli (1995) captured only the steady state and sustained oscillatory behaviours. In the present study, the entire transient phenomenon is captured in order to understand and highlight the striking sub-cooling of the bulk observed in previous numerical and experimental works. The present study constitutes the theoretical proof of the observed and numerically predicted phenomenon. The analysis is performed in the bulk first. Due to the singularity in the problem (the walls are maintained at a constant temperature), it is necessary to introduce boundary layers near the walls of the cell.

Several characteristic time scales are inherent to the fluid flow and heat transfer near the critical point. In particular, we mention the acoustic time scale reflected here by the normalized time t , the piston-effect time scale reflected by $\tau = (\varepsilon/\mu^{3/2})t$ (Zappoli *et al.* 1990) and the thermal diffusion time scale. The latter time scale is very long and tends to infinity upon approaching the critical point. Consequently, we shall focus in this study on the acoustic and piston-effect time scales as the phenomenon of sink zone is observed on the piston-effect time scale (see § 4). Although the formulation is presented in dimensionless form, the results will be reported in terms of dimensional variables and parameters, in an effort to elucidate realistically the underlying phenomena.

3. The flow and temperature fields at the acoustic time scale

Although no sink zone is expected to emerge at the acoustic time scale, it is useful to consider this scale in some detail here as it illustrates the implementation of the matched asymptotic approach, providing an additional insight into the distinct physical features in the propagation of acoustic waves in the cell domain. In addition, this section will pave the way towards understanding the mathematical and physical complexity inherent to the flow and temperature fields at the piston-effect time scale.

We begin by observing that at the acoustic time scale, as the physical time is already scaled by the acoustic time, there is no need to introduce any further change of variables.

3.1. The flow and temperature in the bulk region

The physical variables in the problem are written as a superposition of the initial state and first-order perturbations on the t scale with corresponding order coefficients $A_{1 \leq i \leq 4}$ as

$$\left. \begin{aligned} \rho(x, t) &= 1 + A_1 \bar{\rho}(x, t) + o(A_1), & u(x, t) &= A_2 \bar{u}(x, t) + o(A_2), \\ T(x, t) &= T_i + A_3 \bar{T}(x, t) + o(A_3), & P(x, t) &= P_i + A_4 \bar{P}(x, t) + o(A_4). \end{aligned} \right\} \quad (3.1)$$

Introducing expressions (3.1) into the governing equations, the set of (2.5), and examining the balance of terms, provides the order coefficients which turn out to be all equal to the imposed non-dimensional acceleration field magnitude: $A_i = A$ for any i . Thus, we formally write

$$\left. \begin{aligned} \rho(x, t) &= 1 + A \bar{\rho}(x, t), & u(x, t) &= A \bar{u}(x, t), \\ P(x, t) &= P_i + A \bar{P}(x, t), & T(x, t) &= T_i + A \bar{T}(x, t), \end{aligned} \right\} \quad (3.2)$$

which, when introduced in (2.5), lead to the following closed set of first-order equations in the bulk region for the perturbation field:

$$\bar{\rho}_t = -\bar{u}_x, \quad \bar{u}_t = -\frac{1}{\gamma_0} \bar{P}_x + H(t), \quad -\alpha \bar{u}_x = \bar{T}_t, \quad \bar{P} = \frac{3}{2} \bar{T}, \quad (3.3a-d)$$

where $\alpha = (\gamma_0 - 1)(P_i + \frac{9}{8})$.

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After successive substitutions and eliminating the density, pressure and temperature, we obtain the following non-homogeneous wave equation for the velocity and temperature fields:

$$\bar{u}_{tt}(x, t) = c^2 \bar{u}_{xx}(x, t) + \delta(t), \tag{3.4a}$$

$$\bar{T}_{tt}(x, t) = c^2 \bar{T}_{xx}(x, t), \tag{3.4b}$$

where $\delta(t)$ is the Dirac delta function, reflecting the initial suddenly imposed acceleration and thermal quench at the boundary, as reflected in the second equation in (2.5) and (2.8).

Here, $c = \sqrt{\frac{3}{2}(\alpha/\gamma_0)}$ is the speed of sound for a van der Waals fluid relative to the speed of sound of a perfect gas. Since the fluid is initially at rest, equation (3.4a) is solved subject to $\bar{u}(x, t = 0) = 0$, as well as homogeneous conditions at the walls: $\bar{u}(x = 0, t) = \bar{u}(x = 1, t) = 0$. We recall that no boundary conditions can be imposed at this stage. In this case, we first tackle the solution of (3.4a) for the velocity, which may be conveniently written in the Laplace domain as

$$\bar{u}(x, s) = -\frac{1}{s^2} \left(\frac{e^{-(s/c)x}}{e^{-s/c} + 1} + \frac{e^{(s/c)x}}{e^{s/c} + 1} - 1 \right). \tag{3.5}$$

We note that, throughout this work, we use the same notation for any variable and its Laplace transform to avoid the proliferation of notations. From the solution of the velocity field, the transformed bulk temperature profile can be obtained from (3.4b) as

$$\bar{T}(x, s) = -\frac{2\gamma_0 c}{3} \frac{1}{s^2} \left(\frac{e^{-(s/c)x}}{e^{-s/c} + 1} - \frac{e^{(s/c)x}}{e^{s/c} + 1} \right). \tag{3.6}$$

We observe that, although the velocity profile (3.5) satisfies the homogeneous (no-penetration) conditions at the walls, this is not the case for the temperature, necessitating the presence of a boundary layer at each wall. In the matched asymptotic analysis, boundary conditions for the bulk are derived from matching with the boundary layers. However, it will be shown later in this section that the order coefficient for velocity in the boundary layer is much smaller than in the bulk. Consequently, the boundary conditions for the velocity in the bulk are the actual physical no-penetration conditions introduced in (2.7a,b). Finally, and as we shall see in § 5, although the velocity takes the same form as in the case of a perfect gas, the temperature assumes a slightly different form as a result of the additional coupling with the density in that case (see (3.3a–d)).

We note that solutions (3.5) and (3.6) yield time-oscillatory velocity and temperature, as well as oscillatory density and pressure. By taking the time average of the solution over one cycle, the following temperature profile is obtained:

$$\langle \bar{T} \rangle(x) = -\frac{\gamma_0}{3} (1 - 2x), \tag{3.7}$$

which suggests that the average profile does not depend on the imposed acceleration, wall temperature and proximity to the critical point. Although the profile $\bar{T}(x, t)$ itself depends on the proximity to the critical point (c depends on it), averaging it over one cycle cancels its effect. However, $\bar{T}(x, t)$ is multiplied by A (see (3.2)), i.e. by the non-dimensional acceleration whose value for a 10 g acceleration (considered here) is of the order of $A \approx 10^{-5}$. It can be shown that this average temperature is above the wall temperature on the acoustic time scale. In this case, $\langle \bar{T} \rangle_{max}(x = 1) \simeq 0.46$ and $\langle \bar{T} \rangle_{min}(x = 0) \simeq -0.46$

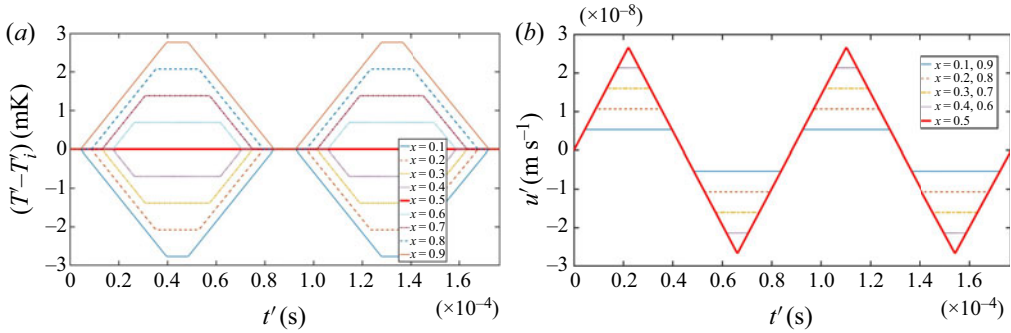


Figure 4. Transient (a) temperature profiles and (b) velocity profiles at the acoustic time scale for different (dimensionless) locations for $0 < x < 1$ in the bulk region. The walls are maintained at constant temperature T'_w such that $T'_i > T'_w > T'_c$ for $t' > 0$. The horizontal axis represents the dimensional time t' (acoustic time scale). The fluid is CO₂ with critical temperature of 304.14 K filled inside a 10 mm thick one-dimensional slab under 10 g acceleration. Initial temperature is 0.2 K above the critical point giving the value of dimensionless speed of sound of $c = 0.8$. The angular points or sharp corners are a result of choosing the acceleration as a step function. Note from figure 4(a) that $T' - T'_i \leq 0$ for $0 < x \leq 0.5$ and $T' - T'_i \geq 0$ for $0.5 \leq x < 1$.

for $\gamma_0 = 1.4$. So $\langle \bar{T} \rangle_{min}$ is above the boundary value temperature $\bar{T}_w = -\delta T' / AT'_c \sim -5.4$ with $A = 1.22 \times 10^{-5}$. We note that, in contrast to the boundary layer regions, the velocity and temperature fields do not depend on the thermal quench in the bulk region.

Solutions (3.5) and (3.6) are numerically inverted and the oscillatory profiles for the temperature and velocity fields are obtained, as shown in figure 4.

The velocity and temperature profiles are periodic with a dimensionless period of $2/c \approx 2.5$ (which corresponds approximately to 8.8×10^{-5} s and corroborates well the results in figure 4). This period corresponds to the propagation of the sound speed back and forth in the one-dimensional cell. While the velocity profiles in figure 4(a) are symmetric about $x = 0.5$ (the profiles for $x > 0.5$ are identical to the ones for $x < 0.5$), the temperature profiles in figure 4(b) are antisymmetric with the same dimensionless period of $2/c \approx 2.5$. This distinction in symmetry, although not possible to confirm from (3.4), becomes evident when replacing x by $1 - x$ in solutions (3.5) and (3.6). In spite of the space–time coupling between the velocity and the temperature in (3.3a–d) and (3.4), which results in similar waveforms (trapezoidal due to step function), solutions (3.5) and (3.6) clearly indicate the velocity and the temperature are not in phase. In fact, figure 4 indicates that the velocity reaches its maximum at a location $x < 0.5$ earlier than the temperature reaching its minimum. On the other hand, at a location $x > 0.5$, the velocity reaches its maximum later than the temperature. In general, the delay decreases with diminishing distance from the middle of the cell. Finally, we note that, although the temperature remains equal to its initial value at all time in the middle of the cell ($x = 0.5$), the velocity continues to fluctuate with time as shown in figure 4.

As stated above, only the velocity boundary conditions were used and no boundary conditions were imposed on the temperature when obtaining the bulk temperature profile at the acoustic time scale. More importantly, the solution is thus independent of the thermal quench magnitude. Consequently, this profile is solely due to the presence of the suddenly imposed acceleration field A initially. This in turn causes the pressure and consequently the temperature through the last equation in (3.3a–d), to oscillate about an average linear profile with constant gradient. Later, it will be shown that this behaviour is also exhibited

for a perfect gas. We finally note that no sink zone is developed in the bulk at this acoustic time scale.

3.2. The top and bottom boundary layers

The analysis is focused on the top boundary layer located at $x = 0$ as the sink zone is most likely to appear on the top layer due to the gravity field which is directed downwards. In order to capture this phenomenon in this confined region, which represents the thin thermal diffusion layer due to the vanishing of the thermal diffusion coefficient, a new space variable z is introduced and is defined by $z = x/\delta$. Here $\delta = \sqrt{\varepsilon\mu^{1/2}}$ is the boundary layer thickness. All the variables are once again perturbed about the initial values. However, in this case, terms are expanded for $x \rightarrow 0$ and $\delta \rightarrow 0$ with z fixed. In particular, a coupling arises between the diffusion term and work of the pressure forces in the energy equation, which represents the thermomechanical coupling. Such a coupling is characteristic of near-critical boundary layers and absent in classical perfect gases (see § 5). Similar to the case in the bulk region, the balance of terms yield the following expansions:

$$\left. \begin{aligned} \rho(z, t) &= 1 + \frac{A}{\mu} \hat{\rho}(z, t), & u(z, t) &= A \sqrt{\frac{\varepsilon}{\mu}} \hat{u}(z, t), \\ T(z, t) &= T_i + A \hat{T}(z, t), & P(z, t) &= P_i + A \hat{P}(z, t). \end{aligned} \right\} \quad (3.8)$$

It is now clear that the order coefficient of the velocity is much smaller compared to the one in the bulk, and hence the validity of the bulk boundary (no-penetration) conditions on the velocity is fully justified. Introducing expressions (3.8) in the set of equations (2.5) leads to the following system of equations:

$$\hat{\rho}_t = -\hat{u}_z, \quad \hat{P}_z = 0, \quad K \hat{u}_z = \frac{2}{3} \hat{T}_{zz}, \quad \hat{P} = \frac{3}{2} \hat{T} + \frac{9}{4} \hat{\rho}, \quad (3.9a-d)$$

where $K = \frac{2}{3}(\alpha\gamma_0\Lambda)Pr_0$.

The boundary conditions are

$$\hat{T}(z = 0, t) = -aH(t), \quad \hat{u}(z = 0, t) = 0. \quad (3.10a,b)$$

Here, we recall $a = \delta T/A = O(1)$. An additional boundary condition is needed in the limit $z \rightarrow \infty$ in order to close the problem. It corresponds to the matching condition on the temperature

$$\lim_{z \rightarrow \infty} \hat{T}(z, t) \sim \lim_{x \rightarrow 0} \bar{T}(x, t). \quad (3.11)$$

The equation for the temperature is readily derived from the set (3.9a-d) as

$$\hat{T}_t = \frac{1}{K} \hat{T}_{zz} + \frac{2}{3} \hat{P}_t = \frac{1}{K} \hat{T}_{zz} + \bar{T}_t(x = 0, t). \quad (3.12)$$

Here, we note that the pressure is known since (3.9a-d) indicates that the pressure is constant across the boundary layer, and is therefore equal to the pressure in the bulk region.

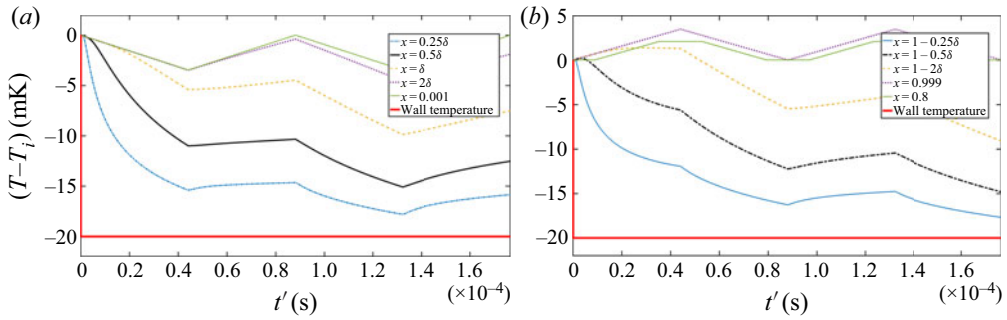


Figure 5. Transient temperature profile of $T' - T'_i$ in the top (a) and bottom (b) boundary layers at the acoustic time scale for six different positions. The thermal quench of 20 mK is imposed at both walls and all other parameters are exactly the same as in figure 4. All profiles are oscillatory with a dimensionless period of $2/c$. We observe that the angular points are the direct result of the non-sinusoidal wave patterns shown in figure 4.

In this case, the temperature deviation in the Laplace domain becomes

$$\hat{T}(z, s) = -\frac{a}{s} e^{-z\sqrt{Ks}} + \bar{T}(x = 0, s)(1 - e^{-z\sqrt{Ks}}). \tag{3.13}$$

The temperature in the bottom layer is obtained similarly by introducing $z^* = (1 - x)/\delta$, yielding the counterpart solution of (3.13)

$$\hat{T}(z^*, s) = -\frac{a}{s} e^{-z^*\sqrt{Ks}} + \bar{T}(x = 1, s)(1 - e^{-z^*\sqrt{Ks}}). \tag{3.14}$$

Noting that $\bar{T}(x = 1, s) = -\bar{T}(x = 0, s) = (1/s^2)(2\gamma_0 c/3)((e^{s/c} - 1)/(e^{s/c} + 1))$, (3.13) and (3.14) indicate that while the temperature is antisymmetric in the bulk region, the symmetry is broken by the imposition of the thermal quench at the walls. A more useful expression than the separate representations in the bulk and boundary layer regions is the composite solution. Recalling that $z = x/\delta$ and $z^* = (1 - x)/\delta$, the composite temperature can be expressed in terms of x by combining (3.6), (3.13) and (3.14)

$$T(x, t) = T_i + A[\bar{T}(x, t) + \hat{T}(z, t) - \lim_{z \rightarrow \infty} \hat{T}(z, t) + \hat{T}(z^*, t) - \lim_{z^* \rightarrow \infty} \hat{T}(z^*, t)]. \tag{3.15}$$

Once again, due to the complex expressions of the temperature field, the analytical inversion is not possible. Hence, numerical inversion is performed using the Euler algorithm described by Abate & Whitt (2006). The algorithm is an implementation of Fourier-series method, using Euler summation to accelerate the convergence of the final infinite series appearing while evaluating the Bromwich integral for the inverse Laplace transform of a complex valued function.

Figure 5 illustrates the evolution of the temperature at different positions in the boundary layer regions. Although all profiles are oscillatory with a non-dimensional period of $2/c$, periodicity is essentially lost very near the wall, with a sudden drop in the temperature. A little further from the wall, the initial drop softens, and smooth oscillations begin to appear. Further still, the oscillations begin to show a rapid change around the peak. The triangular wave pattern appears at a location close to the walls. The temperature ramps up at an increasingly slower rate when approaching the peak to diminish rapidly thereafter. The difference in the growth and diminishing rates is reflected in the dissymmetry of the oscillations. Far away from the wall, the behaviour overlaps that of the bulk (reported earlier in figure 4); the change in slope becomes discontinuous, and the temperature ramps up and down at the same rate (symmetric wave). As expected, the temperature remains at all time above the temperature of the walls: no sink zone is observed.

4. The flow and temperature fields at the piston-effect time scale

The second time scale that appears in fluids near the critical point is the piston-effect time scale. It represents the time at which the temperature is homogenized (to first order) when a cell is heated on one end. A new time variable $\tau = (\varepsilon/\mu^{3/2})t$ relative to the piston-effect time scale (Zappoli *et al.* 1990, 1996; Zappoli, Beysens & Garrabos 2016) is introduced to capture the phenomena at this time scale. Here again, we examine the response subject to the imposed acceleration and thermal wall quench.

4.1. The bulk region

In this region, the orders of the density, temperature, pressure and velocity are determined as a function of the amplitude A and proximity to the critical point μ as

$$\left. \begin{aligned} \rho(x, \tau) &= 1 + A\bar{\rho}(x, \tau), & u(x, \tau) &= A\frac{\varepsilon}{\mu^{3/2}}\bar{u}(x, \tau), \\ P(x, \tau) &= P_i + A\bar{P}(x, \tau), & T(x, \tau) &= T_i + A\bar{T}(x, \tau), \end{aligned} \right\} \quad (4.1)$$

with the corresponding linearized equations

$$\bar{\rho}_\tau = -\bar{u}_x, \quad \bar{P}_x = \gamma_0 H(\tau), \quad -\alpha\bar{u}_x = \bar{T}_\tau, \quad \bar{P} = \frac{2}{3}\bar{T}. \quad (4.2a-d)$$

Comparing systems (3.3a-d) and (4.2a-d), we note interesting similarities and differences between the acoustic and piston-effect time scales in the bulk region. At the piston-effect time scale, the pressure as well the temperature gradients are solely due to the action of the imposed acceleration. At the acoustic time scale, additional contribution results from inertia in the second (momentum) in (3.3a-d). The absence of the acceleration term at the piston-effect time scale in the momentum equation in (4.3) leads to a break in the symmetry exhibited at the acoustic time scale. At both time scales, the diffusion term in the energy equation is negligible, and a wall temperature condition is therefore not needed for the bulk solution. We finally note that the density is decoupled from the remaining variables at both scales as it is dominated by the pressure and temperature in the equation of state. This is not the case in the boundary layer regions, as we shall see and already saw in (3.9a-d).

The solution of (4.2a-d) leads to the following general expressions for the temperature and velocity in the Laplace domain:

$$\bar{T}(x, s) = \frac{2}{3} \left[\frac{\gamma_0 x}{s} + F(s) \right] = \frac{2}{3} \bar{P}(x, s), \quad (4.3)$$

$$\bar{u}(x, s) = -\frac{\gamma_0}{3\alpha} x^2 - \frac{2}{3} \frac{s}{\alpha} F(s)x + G(s), \quad (4.4)$$

where $F(s)$ and $G(s)$ are unknown functions which will be determined from matching. In comparison to the acoustic time scale, we can already see two important distinctions. We first observe that the symmetry in the velocity and temperature fields earlier predicted at the acoustic scale in the bulk region (see figure 4) is now broken. This becomes evident when replacing x by $1 - x$ in (4.3). We also anticipate that the order of magnitude of the velocity in the boundary layers is comparable to that in the bulk region. Consequently, and unlike the case at the acoustic time scale, the no-penetration condition cannot apply for the velocity (4.4) in the bulk region. Hence, at the piston-effect time scale, the boundary layers at the two walls are both hydrodynamic and thermal, with matching conditions needed to obtain the complete profiles of the velocity and the temperature. We finally anticipate the

pressure to be independent of position across the two boundary layers. Consequently, the last equation in system (4.2a–d) and (4.3) suggest, and as we shall show, that $F(s)$ actually turns out to be the pressure in the top boundary layer.

4.2. The top and bottom boundary layers

For the top layer, we again introduce the following change of variable $z = x/\delta$, where we recall $\delta = \sqrt{\varepsilon\mu^{1/2}}$, and set

$$\left. \begin{aligned} \rho(z, \tau) &= 1 + \frac{A}{\mu} \hat{\rho}(z, \tau), & u(z, \tau) &= A \frac{\varepsilon}{\mu^{3/2}} \hat{u}(z, \tau), \\ T(z, \tau) &= T_i + A \hat{T}(z, \tau), & P(z, \tau) &= P_i + A \hat{P}(z, \tau), \end{aligned} \right\} \quad (4.5)$$

leading to the following equations for the perturbation field:

$$\hat{\rho}_\tau = -\hat{u}_z, \quad \hat{P}_z = 0, \quad K\hat{u}_z = \frac{2}{3}\hat{T}_{zz}, \quad \hat{P} = \frac{3}{2}\hat{T} + \frac{9}{4}\hat{\rho}. \quad (4.6a-d)$$

Interestingly, system (4.6a–d) is the same as (3.9a–d) at the acoustic time scale when τ is replaced by t . However, the solution in each case is not the same since it is matched on to different temperatures in the bulk region. Both (3.9a–d) and (4.6a–d) indicate that, close to the walls, the diffusion term is no longer negligible, causing a thin boundary layer to form, and a coupling occurs between the density and temperature terms through the state equation which gives rise to the thermomechanical coupling.

Upon solving system (4.6a–d) subject to the top wall conditions, and matching with the bulk field (4.3) and (4.4), we obtain, in particular, the temperature and velocity perturbations near the top wall

$$\hat{T}(z, s) = -\frac{a}{s} e^{-z\sqrt{Ks}} + \frac{2}{3} \hat{P}_T(s) (1 - e^{-z\sqrt{Ks}}), \quad (4.7)$$

$$\hat{u}(z, s) = \frac{2}{3} \frac{1}{\sqrt{Ks}} \left(\frac{2}{3} s \hat{P}_T(s) + a \right) (e^{-z\sqrt{Ks}} - 1). \quad (4.8)$$

Here, $\hat{P}_T(s)$ is the pressure transform in the top layer. Matching the temperature in the bulk and top layer regions by taking the limit $x \rightarrow 0$ of (4.3) and $z \rightarrow \infty$ of (4.7), we find that $F(s) = \hat{P}_T(s)$.

Similarly, matching (4.4) and (4.7) for the velocity, we obtain $G(s) = -\frac{2}{3}(1/\sqrt{Ks})(\frac{2}{3}s\hat{P}_T(s) + a)$. We note that the pressure $\hat{P}_T(s)$ remains unknown at this stage, and will be determined next once the matching near the bottom layer is secured.

A similar treatment is implemented in the boundary layer near the bottom wall by letting $z^* = (1 - x)/\delta$. In this case, the governing equations are

$$\hat{\rho}_\tau = \hat{u}_{z^*}, \quad \hat{P}_{z^*} = 0, \quad -K\hat{u}_{z^*} = \frac{2}{3}\hat{T}_{z^*z^*}, \quad \hat{P} = \frac{3}{2}\hat{T} + \frac{9}{4}\hat{\rho}, \quad (4.9a-d)$$

yielding the following temperature and velocity perturbations in the Laplace domain:

$$\hat{T}(z^*, s) = -\frac{a}{s} e^{-z^*\sqrt{Ks}} + \frac{2}{3} \hat{P}_B(s) (1 - e^{-z^*\sqrt{Ks}}), \quad (4.10)$$

$$\hat{u}(z^*, s) = -\frac{2}{3} \frac{1}{\sqrt{Ks}} \left(\frac{2}{3} s \hat{P}_B(s) + a \right) (e^{-z^*\sqrt{Ks}} - 1). \quad (4.11)$$

Matching the temperatures from (4.3) and (4.10) by setting $\bar{T}(x \rightarrow 1, s) \sim \hat{T}(z^* \rightarrow \infty, s)$ indicates that the pressure transform in the bottom layer becomes $\hat{P}_B(s) = F(s) + \gamma_0/s$.

On the long-time transient formation of sink zones

Finally, $F(s)$ is determined from (4.4) and (4.11) by setting $\bar{u}(x \rightarrow 1, s) \sim \hat{u}(z^* \rightarrow \infty, s)$, yielding $F(s) = -\gamma_0/2s - \frac{3}{2}a(\beta/s)(\sqrt{s} + \beta)^{-1}$, where $\beta = \frac{3}{4}(\sqrt{K}/\alpha)$.

4.3. The sink-zone phenomenon

Noting that $L^{-1}[(\beta/s)(\sqrt{s} + \beta)^{-1}] = 1 - e^{\beta^2\tau} \operatorname{erfc}(\beta\sqrt{\tau})$, the pressure distribution becomes readily and explicitly available in the three regions

$$P(x, \tau) = P_i - \frac{3}{2}\delta T[1 - e^{\beta^2\tau} \operatorname{erfc}(\beta\sqrt{\tau})] + \begin{cases} -A\frac{\gamma_0}{2}, & x \sim 0^+ \\ A\gamma_0\left(x - \frac{1}{2}\right), & 0 < x < 1 \\ +A\frac{\gamma_0}{2}, & x \sim 1 \end{cases} \quad (4.12)$$

Clearly, the transient response is directly related to the thermal quench at the walls. Interestingly, and unlike the temperature and velocity, the initial pressure profile across the three regions diminishes by the same factor $\frac{3}{2}\delta T$ (uniformly across the three regions). At any time, the pressure exhibits the same gradient across the bulk region.

As to the temperature, unlike the pressure, the analytical inversion in the top and bottom boundary layers is not possible, and the inversion will be carried out numerically. Substituting $F(s)$ in (4.3), the analytical inversion leads to an explicit time and position dependence for the temperature in the bulk region

$$T(0 < x < 1, \tau) = T_w + \frac{2}{3}AH(\tau)\gamma_0(x - \frac{1}{2}) + \delta T e^{\beta^2\tau} \operatorname{erfc}(\beta\sqrt{\tau}). \quad (4.13)$$

Profile (4.13) illustrates analytically the conditions for the sink zone to emerge in the bulk region, and corresponds to the basic idea behind the present work. The composite solution for the temperature transform is obtained by combining (4.3), (4.7) and (4.10).

The evolution of the temperature profiles in the bulk and boundary layers is illustrated in figure 6, where the temperature is plotted against position at different times. Since the profiles in the boundary layers are obliterated over the range $0 < x < 1$ (middle sections of figure 6), they are shown over a small x range near the top and bottom walls in the left and right sections of the figure, respectively. Figure 6(a) shows a wider perspective over the range $0 < t' < 25$ s, whereas figure 6(b) depicts the details for $5 < t' < 25$ s.

We have also added the profiles corresponding to a zero acceleration (dashed lines). In all three regions, the temperature drops significantly with time, especially initially. The profiles in the middle section of figure 6(a,b) reflect the behaviour in the bulk region. We first observe that the temperature gradient $\frac{2}{3}A\gamma_0$ is constant and independent of time. There are two contributing terms in (4.13) to the temperature perturbation, the first term $\frac{2}{3}A\gamma_0(x - \frac{1}{2})$ represents the coupling between the imposed acceleration and the pressure gradient. The second term $e^{\beta^2\tau} \operatorname{erfc}(\beta\sqrt{\tau})$ is the signature of the piston effect in the heat transfer process. It is positive and monotonically decreasing with time. The profiles are comparable to those numerically calculated by Sharma, Erriguible & Amiroudine (2017a,b).

The transient behaviour becomes evident in the limits of small and large times. For this, we note the following useful asymptotic forms of the transient term, which will be used in

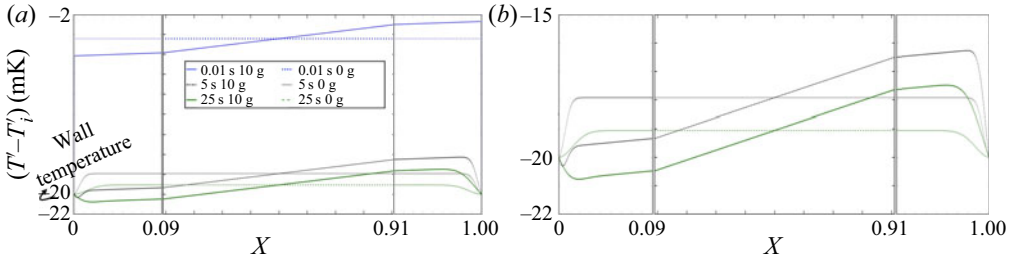


Figure 6. Temperature profiles at the piston-effect time scale against position x . Profiles shown for $0 < t' < 25$ s in (a), and for $5 < t' < 25$ s in (b). The bulk region is shown for $0.09 \leq x \leq 0.91$, along with the corresponding boundary layers. The wall temperature T'_w corresponds to the (-20 mK) horizontal line. The conditions are the same as at the acoustic time scale except that the thermal quench here is 20 mK . Dotted lines correspond to zero gravity and plain lines to an acceleration of 10 g .

subsequent discussions:

$$e^{\beta^2 \tau} \operatorname{erfc}(\beta \sqrt{\tau}) \approx \begin{cases} 1 - \frac{2}{\sqrt{\pi}} \beta \sqrt{\tau}, & (\tau \ll 1) \\ \frac{1}{\beta \sqrt{\pi \tau}}, & (\tau \gg 1) \end{cases} \quad (4.14)$$

Consequently, (4.13) and (4.14) suggest that, in the absence of an imposed acceleration ($A = 0$), the temperature will everywhere always be larger than the wall temperature $T_w = T_i - \delta T$, and this case represents the usual process of heat diffusion, with the pressure remaining independent of the position across the three regions and below the initial pressure at all time as per (4.12). Simultaneously, the temperature is independent of the position (flat dashed lines in the middle section of figure 6) with narrow boundary layers. Now, for $A \neq 0$, there is a competition between the acceleration and the piston-effect thermal quench. It is clear from (4.14) that the temperature can be below that of the wall only in the upper half-region ($x < 1/2$); $T(x > 1/2, \tau) > T_w$ at any time. The early sharp drop in the bulk temperature depicted in figure 6(a) is reflected in the first equation of (4.14). The saturation shown in the figure at the later time is reflected in the second equation of (4.14) through the $1/\sqrt{\tau}$ behaviour. In this case, the effect of the thermal quench, despite its imposition at the walls, is no longer felt across the bulk region. The monotonic decrease with time reflected in figure 6 and (4.13) is typical of the bulk region, and will not always be present near the walls. The profiles reveal that the temperature decreases below the boundary value at a longer time ($t' \simeq 25 \text{ s}$ in figure 6, the piston-effect time scale) near the upper wall of the cell. More generally, this time may be estimated from (4.14) to be: $(\pi/4\beta^2)(1 - \gamma_0/3a)^2 < \tau_{PE} < 9a^2/\pi\gamma_0^2\beta^2$.

Concerning the complex dynamics near the walls, additional insight is gained by examining (4.7) and (4.10) for small z and z^* , respectively. In this case, the analytical inversion becomes possible. Noting that $L^{-1}[(s + \beta\sqrt{s})^{-1}] = e^{\beta^2 \tau} \operatorname{erfc}(\beta \sqrt{\tau})$, we obtain

$$T(z \rightarrow 0, \tau) \approx T_w - \sqrt{K} \left[\frac{\gamma_0 A - 3\delta T}{3\sqrt{\pi \tau}} + \delta T \beta e^{\beta^2 \tau} \operatorname{erfc}(\beta \sqrt{\tau}) \right] z, \quad (4.15)$$

$$T(z^* \rightarrow 0, \tau) \approx T_w + \sqrt{K} \left[\frac{\gamma_0 A + 3\delta T}{3\sqrt{\pi \tau}} - \delta T \beta e^{\beta^2 \tau} \operatorname{erfc}(\beta \sqrt{\tau}) \right] z^*. \quad (4.16)$$

Figure 6 indicates that while the temperature drops everywhere like $1/\sqrt{\tau}$ at a significant rate initially. The early behaviour is reflected from (4.15) and (4.16) in the sharp slopes near the walls as the term involving $1/\sqrt{\tau}$ is dominant for small τ . At a later time, on recalling (4.14), the effect of the quench is annihilated in (4.15) and (4.16) as the terms proportional to δT cancel out, yielding a temperature slope essentially equal to $-A(\gamma_0/3)\sqrt{K/\pi\tau}$ near both walls, a behaviour that also explains the saturation with time shown in figure 6 near the two walls. More importantly, (4.16) confirms, as figure 6 indicates, that a sink zone cannot form near the bottom wall for short and long time. On the other hand, a sink zone can form near the top wall. The temperature can drop by an amount essentially equal to $A(\gamma_0/3)\sqrt{K/\pi\tau}$ that subsides over the piston-effect time scale, independently of the thermal quench level.

This result is similar to the experimental observations of Beysens, Fröhlich & Garrabos (2011) in which an internal heat source is applied in a cell containing a SCF with thermostated walls (see figure 1). A hot boundary layer compresses (heats) the bulk fluid, and a cold boundary layer expands (cools) the bulk fluid. When the heating stops, the pressure ceases to rise due to the cold layer. However, the pressure drop due to the cold layer still continues as the layer slowly vanishes by diffusion. This results in a temperature drop in the bulk sample that can be reached beyond the initial equilibrium temperature as can be seen from figure 1 (Beysens, Fröhlich & Garrabos 2011). The effect is more important when the cold boundary layer is thinner, making the heat exchange more efficient. This is especially true when the cold boundary layer is convected by buoyancy due to the presence of the acceleration. The cooling effect is then more pronounced when the acceleration is larger. Our theoretical results corroborate well this experimental study. Moreover, our predictions are in agreement with the numerical results (Sharma, Erriguible & Amiroudine 2017a,b).

Figure 7 shows the evolution of the temperature field as a function of time over both boundary layer and bulk regions. As can be seen from this figure, the top boundary exhibits a negative temperature (the temperature dips below the imposed wall temperature, which in this case is $T'_w = T'_i - 20$ mK). A minimum temperature, similar to the one observed by Beysens, Fröhlich & Garrabos (2011) reported above in figure 1, is also predicted and the time at which this minimum temperature occurs increases with the increase in distance from the top wall. On the other hand, the bottom boundary layer does not exhibit a minimum. The presence and absence of the minimum near the top and bottom walls, respectively, can be confirmed from (4.15) and (4.16). Approximately, for a short time and close to the top wall, (4.15) suggests, after using (4.14), that the minimum occurs at a time $\tau \simeq (3a - \gamma_0)/6a\beta^2$.

4.4. *The mechanism behind the sink-zone formation*

We are now in a position to elucidate the physical mechanism behind the temperature drop. To aid the discussion, we also include contour and surface plots in figure 8 for the temperature as function of time and position. We recall that the acceleration is acting downward in the x -direction (see figure 3). The heavier fluid in the top thermal boundary layer and the bulk exerts a force (due to its self-weight) on the fluid beneath it, which leads to a high stratification of the density field in the whole domain (due the high compressibility). Consequently, the fluid is compressed near the bottom, while near the top the fluid expands in addition to the usual expansion caused by the piston effect. Therefore, the overall temperature profile is a superimposed effect of the expansion of fluid due to the piston effect and compression or expansion due to the self-weight, depending on the location of the fluid element from the bottom. This results in the temperature near the

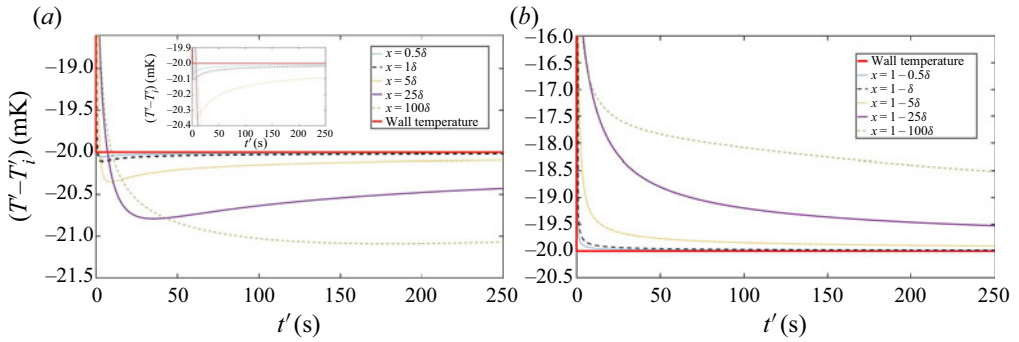


Figure 7. Temperature profile at a fixed z as a function of time (a) in the left boundary layer and (b) in the right boundary layer. The conditions are the same as in previous plots.

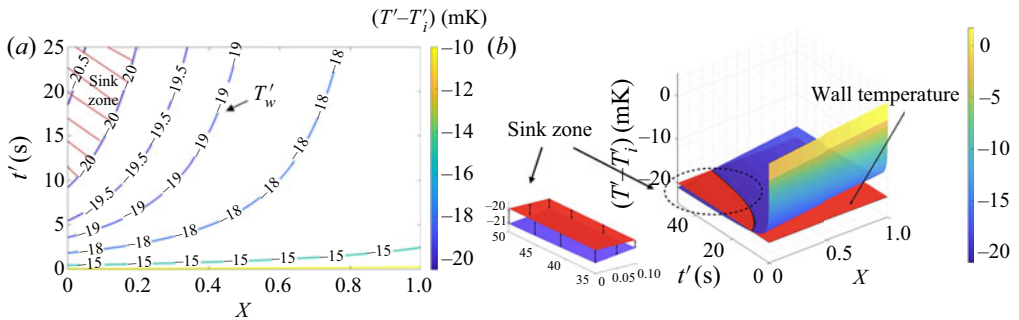


Figure 8. Temperature contour plot (a) and surface (b) as a function of the time and position at the piston-effect time scale. The thermal quench is 20 mK.

bottom being higher than in a weightlessness condition, as can be clearly seen from 6 and 8 (the temperature is higher than -20 mK). At subsequent times, as the temperature of the fluid in the thermal boundary layer drops due to cooling, the density in the thermal boundary layer increases while the thermal expansion (contraction in the present case) increases (due to its diverging behaviour when approaching the critical point). While the fluid in the bottom thermal boundary layer is already in a stable configuration, with cold and heavier fluid being under the lighter bulk fluid, the scenario is the opposite for the fluid in the top thermal boundary layer. A heavier and colder fluid (in the top thermal boundary layer) not only compresses the fluid beneath it but also moves downwards, causing the temperature to drop at a faster rate as compared to the weightlessness condition. The combined action of compression of the bulk fluid by this heavy fluid and action of self-weight causes a faster decrease in the temperature near the top as compared to the bottom. Consequently, while the temperature of the fluid near the top reaches the wall temperature, the fluid near the bottom is yet to attain the imposed wall temperature and thus continues to expand. This temperature gradient near the bottom, which corresponds to the temperature in the bulk (and it is constant, as can be seen from figure 6), thus causes the fluid to expand, thereby decreasing the temperature everywhere. As a result, the temperature near the top wall falls below the boundary value and extends into the bulk fluid with time as can be seen in figures 6 and 8). The temperature of the fluid continues to fall below the wall temperature until the fluid in the thermal boundary layer

starts to expand due to the heating from the top wall and thus exerting an opposite effect, i.e. compressing the bulk fluid.

When both these opposite effects (the expansion and compression of the bulk due to the bottom and top boundary layers, respectively) become balanced, the temperature stops falling below the boundary value. Subsequently, it tends to return to the imposed values as imposed by the boundary conditions (see figure 7).

5. The special case of a perfect gas

In an effort to contrast the behaviour of the van der Waals gas near criticality, we consider the case of a perfect gas of constant properties, with governing equations deduced from system (2.5) by setting $\Lambda = 0$ and $P = \rho T$, subject to the same initial and boundary conditions. We observe that only two time scales exist in the case of a perfect gas, namely the acoustic and diffusive time scales (Kassoy 1979).

5.1. At the acoustic time scale

Perturbing about the initial values, exactly as in (3.2) for a van der Waals gas, yields the following equations:

$$\bar{\rho}_t = -\bar{u}_x, \quad \bar{u}_t = -\frac{1}{\gamma_0} \bar{P}_x + H(t), \quad -(\gamma_0 - 1)\bar{u}_x = \bar{T}_t, \quad \bar{P} = \bar{\rho} + \bar{T}. \quad (5.1a-d)$$

Interestingly, the flow and temperature fields are coupled to the density, which is not the case for a van der Waals gas, as equation (3.3a-d) suggests. In this case, the equation and solution for the velocity in the bulk region take the same form as (3.4) and (3.5) with $c = 1$, as expected. The temperature takes a slightly different form from (3.5)

$$\bar{T}(x, s) = \frac{\gamma_0 - 1}{s^2} \left(\frac{e^{sx}}{1 + e^s} - \frac{e^{-sx}}{1 + e^{-s}} \right). \quad (5.2)$$

As in the case of a van der Waals gas, the behaviour is oscillatory in time. The time average over one cycle also suggests, as does (3.7), a linear increase with position

$$\langle \bar{T} \rangle = (\gamma_0 - 1) \left(x - \frac{1}{2} \right). \quad (5.3)$$

Setting $z = x/\delta$, and noting that the boundary layer thickness is $\delta = \sqrt{\varepsilon}$ for a perfect gas, the field perturbation takes a slightly different form than (3.8)

$$\left. \begin{aligned} \rho(z, t) &= 1 + A\hat{\rho}(z, t), & u(z, t) &= A\sqrt{\varepsilon}\hat{u}(z, t), \\ T(z, t) &= T_i + A\hat{T}(z, t), & P(z, t) &= 1 + A\hat{P}(z, t). \end{aligned} \right\} \quad (5.4)$$

We see that the density perturbation for the perfect gas in the boundary layer differs from that of a near-critical fluid. This is discussed later. The density is of order A in (5.4), which is small relative to the supercritical case for a van der Waals gas for which (3.8) indicates that the density is of the order A/μ . Recall that A and μ are of the order of 10^{-5} and 10^{-3} , respectively. The corresponding equations are

$$\hat{\rho}_t = -\hat{u}_z, \quad \hat{P}_z = 0, \quad \hat{T}_t = -(\gamma_0 - 1)\hat{u}_z + \frac{\gamma_0}{Pr_0}\hat{T}_{zz}, \quad \hat{P} = \hat{T} + \hat{\rho}. \quad (5.5a-d)$$

When compared to (3.9a-d), we see that (5.5a-d) suggests that the transient term in the energy equation is no longer negligible for a perfect gas. However, the resulting

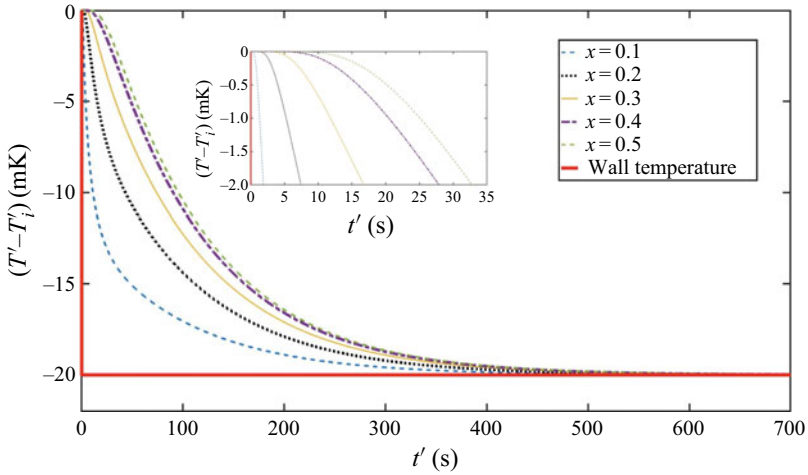


Figure 9. Temperature profiles at the diffusion time scale for the perfect gas. The lower dashed line represents the wall temperature.

temperature equation remains of the same form as (3.12), namely $\hat{T}_t = (1/Pr_0)\hat{T}_{zz} + ((\gamma_0 - 1)/\gamma_0)\hat{P}_t$.

5.2. At the diffusion time scale

The diffusion time scale in a perfect gas is defined by $\tau = \varepsilon t$ where ε is the same small parameter as defined in § 2 below (2.5). In this case, all field perturbations in the bulk region are of order A , with the corresponding equations written as

$$\bar{u}_x = 0, \quad \bar{P}_x = \gamma_0 H(\tau), \quad \bar{T}_\tau = \frac{\gamma_0}{Pr_0} \bar{T}_{xx}, \quad \bar{P} = \bar{\rho} + \bar{T}. \quad (5.6a-d)$$

In this case, and despite the presence of a pressure gradient, the velocity is negligibly small over the entire domain. In addition, the diffusive term in the energy equation ensures that the temperature perturbation satisfies: $\bar{T}(x = 0, \tau) = \bar{T}(x = 1, \tau) = -a\delta(\tau)$ at the two walls.

Consequently, and as expected, there is no formation of thermal boundary layers. In this case

$$T(x, s) = T_i - \delta T \frac{e^{x\sqrt{(Pr_0/\gamma_0)s}} + e^{(1-x)\sqrt{(Pr_0/\gamma_0)s}}}{s(1 + e^{\sqrt{(Pr_0/\gamma_0)s}})}. \quad (5.7)$$

The above equation reveals that at no time the temperature in the bulk falls below $1 - a$. This means that no quenching is observed in perfect gases.

Figure 9 shows the temperature profile as a function of time. The conditions are the same as in the previous SCF case, i.e. the fluid is at 0.2 K above the critical temperature and the applied thermal quench is 20 mK. As expected, no formation of sink zones is predicted.

6. Concluding remarks

The results of the present theoretical analysis indicate that cooling beyond the boundary value or the development of sink zones is unique to near-critical fluids, and is not predicted

for perfect gases. The novelty of the present work is a detailed quantitative theoretical analysis of the sink-zone phenomenon observed in experiments (Beysens, Fröhlich & Garrabos 2011) and numerical simulations (Sharma, Erriguible & Amiroudine 2017a). The present analysis is based on the matched asymptotic expansions, and constitutes a thorough mathematical proof of the existence of sink zones.

At the acoustic time scale in both cases, i.e. near-critical and perfect gases, the development of a weak temperature gradient is observed in the bulk region; weak in the sense that it is not possible for the temperature to drop below the boundary value unless an acceleration field is applied. The temperature at the top wall drops slightly below the initial value, while that at the bottom increases above the initial value. However, the temperature remains exactly the same as the initial temperature at the midpoint. It is further shown that the development of this gradient is independent of the wall temperature and solely depends on the acceleration field or imposed gravity. Thus, it is natural to conclude that the behaviour of the ideal gas and near-critical fluid is nearly the same at the acoustic time scale.

At the piston-effect time scale in a near-critical fluid, an unusual scenario is predicted in which ‘sink zones’ develop. From the expression of the temperature in the bulk, it is seen that, as per the initial value theorem, the initial value of the temperature matches with the final value of the acoustic time scale. This may seem to be a constraint that must be satisfied, but this is not the case since, by introducing a change of variable for a different time scale, only the phenomenon occurring at that time scale is captured and the rest is ‘filtered out’. This initial gradient again depends only on the acceleration and not on the thermal quench amplitude. The effect of the quenching is seen at time $t > 0$ at the piston-effect time scale. The thermo-mechanical coupling equilibrates the temperature in the bulk to that on the boundary, and in the process, the initial temperature profile which contains the temperature gradient, is brought down to the boundary value as expected. However, in the process, the initial gradient is not at all disturbed, and a sink zone near the top wall is observed. This is nicely interpreted in terms of density stratification. We note that the coupling of the pressure gradient with the sudden acceleration initiates the density stratification and development of the initial gradient. This gradient development is caused only by the acceleration, and happens so fast that, later, when the temperature reaches the boundary value, it remains unchanged, and sink zones are formed. From the order coefficients it is also seen that huge density fluctuations occur in the boundary layers which are 600 times larger than in perfect gases. This phenomenon may speed up any kind of chemical reaction occurring at the wall with the inner fluid and proves to have some very good or drastic applications.

The behaviour of the perfect gas at the diffusion time scale is as expected. No sink zones are observed, and a careful examination of the equations show that the acceleration term in the bulk actually disappears, with no gravity effect. The process of diffusion results in a balance between the temperatures in the bulk region and near the boundary.

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Author ORCIDs.

 Roger Khayat <https://orcid.org/0000-0001-8307-7019>;

 Sakir Amiroudine <https://orcid.org/0000-0002-2387-0012>.

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