Entropy-based viscous regularization for the multi-dimensional Euler equations in low-Mach and transonic flows

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Abstract

We present a new version of the entropy viscosity method, a viscous regularization technique for hyperbolic conservation laws, that is well-suited to low-Mach flows. By means of a low-Mach asymptotic study, new expressions for the entropy viscosity coefficients are derived. These definitions are valid for a wide range of Mach numbers, from subsonic flows (with very low Mach numbers) to supersonic flows, and no longer depend on an analytical expression for the entropy function. In addition, the entropy viscosity method is extended to Euler equations with variable area for nozzle flow problems. The effectiveness of the method is demonstrated using various 1-D and 2-D benchmark tests: flow in a converging-diverging nozzle; Leblanc shock tube; slow moving shock; strong shock for liquid phase; low-Mach flows around a cylinder and over a circular hump; and supersonic flow in a compression corner. Convergence studies are performed for both smooth solutions and solutions with shocks present.

Key words: entropy viscosity method, artificial viscosity, low-Mach regime, shock capturing, Euler equations with variable area.

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1. Introduction

Solving accurately compressible fluid equations in the low-Mach limit is an ongoing topic of research. In many engineering applications, compressibility effects require the solution of the compressible fluid equations in nearly incompressible regimes and/or for low-Mach flow problems. For example, such flows are encountered in aerodynamics in the study of airships. In the nuclear industry, flows are nearly in the incompressible regime but compressible effects cannot be neglected because of the intense heat source, and because of some postulated accident scenarios, and thus need to be accurately resolved. Hence, there is a strong interest to develop computational methods that can solve both compressible and incompressible flow problems.

When solving Euler equations for a wide range of Mach numbers, multiple

12 questions must be addressed: stability, accuracy and solution convergence in the 13 low-Mach regime. Because of the hyperbolic nature of the equations, shocks can form during transonic and supersonic flows and require the use of adequate nu-15 merical techniques to stabilize solution and correctly resolve the discontinuities. A wide range of stabilization methods are available in the literature: approxi-17 mate Riemann solvers [1], flux-limiter techniques [2, 3], pressure-based viscosity 18 methods [4], Lapidus method [5, 6, 7], and the entropy-viscosity method [8, 9], 19 among others. These numerical methods are usually developed using simple equations of state and tested for transonic and supersonic flows where the dis-21 parity between the acoustic wave speed and the fluid speed is not excessively 22 large and thus the Mach number is of order one. This approach, however, leads to a well-known accuracy problem in the low-Mach regime where the fluid velocity is smaller that the speed of sound by multiple orders of magnitude. The numerical dissipative terms become ill-scaled in the low-Mach regime and lead to the wrong numerical solution by changing the nature of the equations

solved. This behavior is well documented in the literature [10, 11, 12]. In [10], a low-Mach asymptotic study has demonstrated convergence of the compressible Euler equations to the incompressible ones. Many well-known stabilization techniques, for instance, the Roe scheme and the SUPG technique, do not yield the 31 correct solution in the low-Mach regime and suitable modifications have been 32 proposed (see [13] for the Roe scheme and [12] for the SUPG method) to en-33 sure the convergence to the correct solution while preserving the original shock stabilization properties of these schemes. Additionally, the time step size may be severely restricted when solving compressible fluid equations with an explicit time discretization because of the large disparity between the fluid velocity and 37 the speed of sound. To avoid an excessive number of explicit time steps, time preconditioners have been proposed and proved efficient [11]; however, because they modify the time derivatives in the governing equations, such acceleration techniques can only be used to obtain steady-state solutions for low-Mach flows 41 using explicit schemes. To avoid modifying the time derivatives, the temporal implicit capabilities of the MOOSE multiphysics framework [14] are used. With such a choice, low-Mach steady-state solutions can be obtained effectively while preserving the accuracy of the transient solution; however, it requires the use of nonlinear solvers. 46

In this paper, we employ the entropy viscosity method as a numerical stabilization for the inviscid Euler equation and assess its performance in the lowMach regime. The entropy viscosity method is a viscous regularization technique
introduced by Guermond et al. to solve hyperbolic systems of equations and
has successfully been applied to multi-dimensional supersonic flows with various
spatial discretization schemes [15]. It is fairly straightforward to implement, can
be used with unstructured grids, and has dissipative terms that are consistent
with the entropy minimum principle. However, it has not been evaluated in the

55 low-Mach regime.

This paper is organized as follows: in Section 2 the current definition of the entropy viscosity method is recalled and its ill-scaled nature in the low-Mach 57 regime is discussed. In Section 3, a new formulation of the viscosity residual is derived. This formulation no longer requires an analytical expression for the 59 entropy function. A low-Mach asymptotic study is carried out to adapt the 60 definition of the entropy viscosity coefficients in the incompressible limit while ensuring that the viscosity coefficients scale appropriately for all flow speeds (from low-Mach to supersonic). In Section 4, we extend the entropy viscosity method to Euler equations with variable area in order to model nozzle flows: 64 the viscous dissipative terms are adapted so that the entropy minimum principle remains satisfied. Spatial and temporal discretizations and solution tehcniques are presented in Section 5. 1-D and 2-D numerical results are provided in Section 6 for a wide range of Mach numbers: liquid and gas nozzle flow problems, 68 low-Mach flows over a cylinder and a circular bump (with Mach numbers as low as 10^{-7}), and supersonic flows in a compression corner [16]. Convergence studies are performed in 1-D in order to demonstrate the accuracy of the solution 71 technique.

2. The Entropy Viscosity Method

74 2.1. Background

Euler equations in conservative form are given by

$$\partial_t \rho + \vec{\nabla} \cdot (\rho \vec{u}) = 0 \tag{1a}$$

 $\partial_t \left(\rho \vec{u} \right) + \vec{\nabla} \cdot \left(\rho \vec{u} \otimes \vec{u} + P \mathbb{I} \right) = 0 \tag{1b}$

 $\partial_t \left(\rho E \right) + \vec{\nabla} \cdot \left[\vec{u} \left(\rho E + P \right) \right] = 0 \tag{1c}$

where ρ , $\rho \vec{u}$ and E are the density, the momentum and the total specific energy, respectively, and will be referred to as the conservative variables. \vec{u} is the fluid 79 velocity and its specific internal energy is denoted by $e = E - \frac{u^2}{2}$. An equation 80 of state, dependent upon ρ and e, is used to compute the pressure P. The tensor product $\vec{a} \otimes \vec{b}$ is such that $(\vec{a} \otimes \vec{b})_{i,j} = a_i b_j$. The identity tensor is denoted by \mathbb{I} . 82 Next, the entropy viscosity method [8, 9, 17, 18] applied to Eq. (1) is recalled. The method consists of adding dissipative terms with a viscosity coefficient modulated by the entropy production; this allows for a high-order accuracy when the solution is smooth (provided that the spatial and temporal discretizations also are high order). The derivation of the viscous regularization (or dissipa-87 tive terms) is carried out to be consistent with the entropy minimum principle; details and proofs of the derivation can be found in [15]. The viscous regularization thus obtained is valid for any equation of state as long as the physical entropy function s is concave (or -s is a convex function) with respect to the 91 internal energy e and the specific volume $1/\rho$. The Euler equations with viscous regularization become

$$\partial_t \rho + \vec{\nabla} \cdot (\rho \vec{u}) = \vec{\nabla} \cdot \left(\kappa \vec{\nabla} \rho \right) \tag{2a}$$

$$\partial_t (\rho \vec{u}) + \vec{\nabla} \cdot (\rho \vec{u} \otimes \vec{u} + P \mathbb{I}) = \vec{\nabla} \cdot \left(\mu \rho \vec{\nabla}^s \vec{u} + \kappa \vec{u} \otimes \vec{\nabla} \rho \right)$$
 (2b)

$$\partial_t (\rho E) + \vec{\nabla} \cdot [\vec{u} (\rho E + P)] = \vec{\nabla} \cdot \left(\kappa \vec{\nabla} (\rho e) + \frac{1}{2} ||\vec{u}||^2 \kappa \vec{\nabla} \rho + \rho \mu \vec{u} \vec{\nabla} \vec{u} \right)$$
(2c)

where κ and μ are positive viscosity coefficients (in units of length²/time). $\nabla^s \vec{u}$ denotes the symmetric gradient operator and guarantees the method to be rotationally invariant [15]. The viscosity coefficients are key ingredients in the viscous regularization of Eq. (2). Other stabilization approaches have been proposed in the literature, for instance, the Lapidus method [7, 5] or pressure-based viscosity methods [4]. Here, we follow the work of Guermond et al. and define

the viscosity coefficients, κ and μ , based on the local entropy production. These coefficients are numerically evaluated using the local entropy residual $R_{\rm ent}(\vec{r},t)$ defined in Eq. (3); $R_{\rm ent}(\vec{r},t)$ is known to be peaked in shocks and vanishingly small elsewhere [1].

$$R_{\rm ent}(\vec{r},t) := \partial_t s + \vec{u} \cdot \vec{\nabla} s \tag{3}$$

In the current version of the method, the ratio of κ to μ is defined through a numerical Prandlt number, $Pr = \kappa/\mu$. Pr is a user-defined parameter and 107 is usually taken in the range [0.001;1]. Since the entropy residual $R_{\rm ent}(\vec{r},t)$ may be extremely large in shocks, the definition of the viscosity coefficients 109 also includes a first-order viscosity coefficient that serves as an upper bound for 110 the entropy-based viscosity coefficients. The first-order viscosity coefficients, 111 denoted by μ_{max} and κ_{max} , are chosen so that the numerical scheme becomes 112 equivalent to an upwind scheme when the first-order coefficients are employed. 113 The upwind scheme is known to be over-dissipative but guarantees monotonicity 114 [1]. In practice, the viscosity coefficients only saturate to the first-order viscosity 115 coefficients in shocks and are much smaller elsewhere, hence avoiding the over-116 dissipation of the upwind method. The first-order viscosity coefficients $\mu_{\rm max}$ 117 and κ_{max} are equal and set proportional to the largest local eigenvalue $||\vec{u}|| + c$: 118

$$\mu_{\max}(\vec{r},t) = \kappa_{\max}(\vec{r},t) = \frac{h}{2} \left(||\vec{u}(\vec{t},\vec{r})|| + c(\vec{t},\vec{r}) \right), \tag{4}$$

where h denotes the local grid size (for higher than linear finite element representations, h is defined as the ratio of the grid size to the polynomial order of the test functions used, see Eq. 2.4 in [18]). For simplicity, the first-order viscosity coefficients will only be referred to as $\kappa_{\text{max}}(\vec{r},t)$. In practice, these

quantities are evaluated within a given cell K at quadrature points:

$$\kappa_{\text{max}}^{K}(\vec{r}_{q}, t) = \frac{h_{K}}{2} \left(||\vec{u}(t, \vec{r}_{q})|| + c(t, \vec{r}_{q}) \right),$$
(5)

where \vec{r}_q denotes the position of a quadrature point. As stated earlier, the entropy viscosity coefficients, which we denote by κ_e and μ_e , are set proportional to the entropy production evaluated by computing the local entropy residual $R_{\rm ent}$. The definitions also include the inter-element jump J[s] of the entropy flux, allowing for the detection of discontinuities other than shocks (e.g., contact). κ_e and μ_e are computed as follows

$$\mu_e^K(\vec{r}_q, t) = h_K^2 \frac{\max(|R_{\text{ent}}^K(\vec{r}_q, t)|, J^K[s](t))}{||s - \bar{s}||_{\infty}}$$
(6a)

$$\kappa_e^K(\vec{r}_q, t) = \Pr \mu_e^K(\vec{r}_q, t), \tag{6b}$$

where $||\cdot||_{\infty}$ and $\bar{\cdot}$ denote the L_{∞}-norm and the average operator over the entire computational domain, respectively. The definition of the entropy jump J[s] is spatial discretization-dependent and examples of definitions can be found in [18] for discontinuous Galerkin discretization. For continuous finite element methods (FEM), the jump of a given quantity is defined as the change of its normal derivative $(\partial_n(\cdot) = \vec{n} \cdot \vec{\nabla}(\cdot))$ across the common face separating the two elements, and will be further referred to as the inter-element jump. We take the largest value over all faces f present on the boundary ∂K of element K:

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$$J^{K}[s](t) = \max_{f \in \partial K} \max_{\vec{r}_q \in f} \left(\|\vec{u}(\vec{r}_q, t)\| \llbracket \vec{\nabla} s(\vec{r}_q, t) \cdot \vec{n}(\vec{r}_q) \rrbracket_f \right), \tag{7}$$

where $[a(\vec{r_q})]_f$ denotes the inter-element jump in $a(\vec{r})$ at quadrature point $\vec{r_q}$ on face f (the quadrature points $\vec{r_q}$ are taken on the faces f of the element K). With the definition given in Eq. (7), the jump is constant over each el-

ement K of the computational domain. The denominator $||s - \bar{s}||_{\infty}$ is used for dimensionality purposes. Currently, there are no theoretical justifications for choosing the denominator beyond a dimensionality argument. Finally, the viscosity coefficients μ and κ are as follows:

$$\mu(\vec{r},t) = \min \left(\mu_e(\vec{r},t) , \, \mu_{\max}(\vec{r},t) \right)$$
 and $\kappa(\vec{r},t) = \min \left(\kappa_e(\vec{r},t) , \, \kappa_{\max}(\vec{r},t) \right)$. (8)

Given these definitions, we have the following properties. In shock regions, the entropy viscosity coefficients will experience a peak because of entropy production and thus will saturate to the first-order viscosity. The first-order coefficients are known to be over-dissipative and will smooth out any oscillatory behavior. Elsewhere in the domain, entropy production will be small and the viscosity coefficients μ and κ will remain small. High-order accuracy for entropy-based viscous stabilization has been demonstrated using several 1-D shock tube examples and various 2-D tests [8, 9, 18].

2.2. Issues in the Low-Mach Regime

In the low-Mach Regime, a smooth flow is known to approach the isentropic 148 limit, resulting in very little entropy production. Since the entropy viscosity 149 method is directly based on the evaluation of the local entropy production, it 150 is of interest to study how the entropy viscosity coefficients μ_e and κ_e scale 151 in the low-Mach regime. In practice, the entropy residual $R_{\rm ent}$ will be very 152 small in that regime and so will be the denominator $||s-\bar{s}||_{\infty}$, thus making 153 the definition of the viscosity coefficients in Eq. (6) undetermined and likely ill-154 scaled. One possible approach would consist of expanding the numerator and 155 denominator in terms of the Mach number and deriving its limit when the Mach number goes to zero. Such derivation may not be straightforward, especially 157

for general equations of state. However, this can be avoided by noting that the entropy residual $R_{\rm ent}$ can be recast as a function of pressure, density, velocity, and speed of sound as will be shown in Eq. (9) of Section 3.1. This alternate entropy residual definition is the basis for the low-Mach analysis carried out in this paper and possesses several advantages that are detailed next.

3. An All-speed Reformulation of the Entropy Viscosity Method

In this section, the entropy residual $R_{\rm ent}$ is recast as a function of pressure,
density, velocity and speed of sound. Then, a low-Mach asymptotic study is
carried out for the Euler equations with viscous regularization in order to derive
an appropriate normalization parameter that is valid in the isentropic low-Mach
regime as well as for transonic and supersonic flows.

3.1. New Definition of the Entropy Production Residual

The first step in defining viscosity coefficients that behave well in the lowMach limit is to recast the entropy residual in terms of thermodynamic variables.

This provides physical insight on possible normalization choices that can be valid
in both low-Mach and transonic flows. The alternate definition of the entropy
residual, the derivation of which is given in Appendix A, is the following:

$$R_{\rm ent}(\vec{r},t) := \partial_t s + \vec{u} \cdot \vec{\nabla} s = \frac{\mathrm{D}s}{\mathrm{D}t} = \frac{s_e}{P_e} \left(\underbrace{\frac{\mathrm{D}P}{\mathrm{D}t} - c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t}}_{\tilde{R}_{\rm ent}(\vec{r},t)} \right), \tag{9}$$

where $\frac{D}{Dt}$ denotes the material derivative $(\frac{D}{Dt} := \frac{\partial}{\partial t} + \vec{u} \cdot \vec{\nabla})$, and x_y is the standard shorthand notation for the partial derivative of x with respect to y, e.g., $P_e := \frac{\partial P}{\partial e}$. When using the SGEOS, the coefficient of proportionality $\frac{s_e}{P_e}$ can be recast under the form $\frac{s_e}{P_e} = \frac{C_v(\gamma - 1)}{P + \gamma P_{\infty}}$ which remains positive. Note that the definition of the speed of sound can be used to show that $\widetilde{R}_{\rm ent}$ is zero

when the flow becomes isentropic. The entropy residuals $R_{\rm ent}$ and $\tilde{R}_{\rm ent}$ are proportional to one another and will experience similar variations in space and time. Thus, one may elect to employ $\tilde{R}_{\rm ent}$ instead of $R_{\rm ent}$ for the evaluation of the local entropy residual. The new expression presents several advantages which include:

- An analytical expression of the entropy function s is no longer needed: the residual $\widetilde{R}_{\rm ent}$ is evaluated using the local values of pressure, density, velocity and speed of sound. Deriving an entropy function for some complex equations of state may be difficult;
- Suitable normalizations for the residual $\widetilde{R}_{\text{ent}}$ can be devised. Examples include the pressure itself or combinations of the density, the speed of sound and the norm of the velocity, i.e., ρc^2 , $\rho c ||\vec{u}||$ or $\rho ||\vec{u}||^2$.

Denoting the normalization of $\widetilde{R}_{\rm ent}$ by norm_P, the entropy-based viscosity coefficients μ_e and κ_e can be re-defined as follows:

$$\mu_e^K(\vec{r}_q, t) = h_K^2 \frac{\max\left(|\tilde{R}_{\text{ent}}^K(\vec{r}_q, t)|, J^K(t)\right)}{\operatorname{norm}_{P}^{\mu}}, \qquad (10a)$$

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$$\kappa_e^K(\vec{r}_q, t) = h_K^2 \frac{\max\left(|\tilde{R}_{\text{ent}}^K(\vec{r}_q, t)|, J^K(t)\right)}{\text{norm}_P^{\kappa}},$$
 (10b)

195 where

$$J^{K}(t) = \max_{f \in \partial K} \max_{\vec{r}_{q} \in f} || \left(\vec{u}(\vec{r}_{q}, t) || \max \left(J^{K}[P](t), c^{2}(\vec{r}_{q}, t) || J^{K}[\rho](t) \right) \right)$$
 (10c)

Note that now the jump operator J^K acts on the variables appearing in $\widetilde{R}_{\rm ent}$, namely, pressure and density. The residual $\widetilde{R}_{\rm ent}$ and the pressure jump will only contribute to the viscosity coefficients in the shock region. On the other hand, the density jump will contribute both in contact and shock regions. The

 μ and κ coefficients are kinematic viscosities (units of m^2/s); the normalization parameters norm_P are thus in units of pressure, hence the use of the subscript P.

Note also that we are not requiring the same normalization for both μ_e and κ_e so the entropy viscosity coefficients can be different. The low-Mach asymptotic study presented next will determine the proper normalization.

205 3.2. Asymptotic Study in the Low-Mach Regime

The Euler equations with viscous stabilization, Eq. (6), bear some simi-206 larities with the Navier-Stokes equations in the sense that dissipative terms (containing second-order spatial derivatives) are present in both sets of equa-208 tions. An abundant literature exists regarding the low-Mach asymptotic of the Navier-Stokes equations [10, 11, 12, 19]. The asymptotic study presented here 210 is inspired by the work of Muller et al. [19] where an asymptotic derivation for 211 the Navier-Stokes was presented. We remind the reader that the objective is to 212 determine appropriate scaling for the entropy viscosity coefficients so that the 213 dissipative terms remain well-scaled for two limit cases: (i) the isentropic low-214 Mach limit where Euler equations degenerate to an incompressible system of 215 equations in the low-Mach limit and (ii) the non-isentropic limit with formation 216 of shocks. The isentropic limit of the Euler equations with viscous regular-217 ization should yield incompressible fluid flow solutions in the low-Mach limit, namely, that the spatial steady-state pressure variations are of the order M^2 219 and that the velocity satisfies the divergence constraint $\vec{\nabla} \cdot \vec{u}_0 = 0$ [10, 11, 12]. For non-isentropic situations, shocks may form for any value of Mach number 221 and the minimum entropy principle should still be satisfied so that numerical 222 oscillations, if any, be controlled by the entropy viscosity method independently 223 of the value of the Mach number. Our objective is to determine the appropriate 224 scaling for norm_P and norm_P in these two limit cases. 225

The first step in the study of the limit cases (i) and (ii) is to re-write Eq. (2)

in a non-dimensional manner. To do so, the following variables are introduced:

$$\rho^* = \frac{\rho}{\rho_{\infty}}, \ u^* = \frac{u}{u_{\infty}}, \ P^* = \frac{P}{\rho_{\infty} c_{\infty}^2}, \ E^* = \frac{E}{c_{\infty}^2},$$
$$x^* = \frac{x}{L_{\infty}}, \ t^* = \frac{t}{L_{\infty}/u_{\infty}}, \ \mu^* = \frac{\mu}{\mu_{\infty}}, \ \kappa^* = \frac{\kappa}{\kappa_{\infty}},$$
(11)

where the subscript ∞ denote the far-field or stagnation quantities and the superscript * stands for the non-dimensional variables. The far-field reference quantities are chosen such that the dimensionless flow quantities are of order 1.

The reference Mach number is given by

$$M_{\infty} = \frac{u_{\infty}}{c_{\infty}},\tag{12}$$

where c_{∞} is a reference value for the speed of sound. Then, the scaled Euler equations with viscous regularization are:

$$\partial_{t^*} \rho^* + \vec{\nabla}^* \cdot (\rho^* \vec{u}^*) = \frac{1}{\text{P\'e}_{\infty}} \vec{\nabla}^* \cdot (\kappa^* \vec{\nabla}^* \rho^*)$$
 (13a)

$$\partial_{t^*} \left(\rho^* \vec{u}^* \right) + \vec{\nabla}^* \cdot \left(\rho^* \vec{u}^* \otimes \vec{u}^* \right) + \frac{1}{M_{\infty}^2} \vec{\nabla}^* P^* = \frac{1}{\text{Re}_{\infty}} \vec{\nabla}^* \cdot \left(\rho^* \mu^* \vec{\nabla}^{s,*} \vec{u}^* \right) + \frac{1}{\text{P\'e}_{\infty}} \vec{\nabla}^* \cdot \left(\vec{u}^* \otimes \kappa^* \vec{\nabla}^* \rho^* \right) \quad (13b)$$

$$\partial_{t^*} \left(\rho^* E^* \right) + \vec{\nabla}^* \cdot \left[\vec{u}^* \left(\rho^* E^* + P^* \right) \right] = \frac{1}{\text{P\'e}_{\infty}} \vec{\nabla}^* \cdot \left(\kappa^* \vec{\nabla}^* (\rho^* e^*) \right)$$

$$+ \frac{M_{\infty}^2}{\text{Re}_{\infty}} \vec{\nabla}^* \cdot \left(\vec{u}^* \rho^* \mu^* \vec{\nabla}^{s,*} \vec{u}^* \right) + \frac{M_{\infty}^2}{2\text{P\'e}_{\infty}} \vec{\nabla}^* \cdot \left(\kappa^* (u^*)^2 \vec{\nabla}^* \rho^* \right) , \quad (13c)$$

where the numerical Reynolds (Re_{∞}) and Péclet $(P\acute{e}_{\infty})$ numbers are defined as:

$$\operatorname{Re}_{\infty} = \frac{u_{\infty} L_{\infty}}{\mu_{\infty}} \text{ and } \operatorname{P\acute{e}}_{\infty} = \frac{u_{\infty} L_{\infty}}{\kappa_{\infty}}.$$
 (14)

Note that the Prandlt number used in the original version of the entropy viscosity method is simply given by

$$\Pr_{\infty} = \text{P\'e}_{\infty}/\text{Re}_{\infty}$$
 (15)

The numerical Reynolds and Péclet numbers defined in Eq. (14) are related to
the entropy viscosity coefficients μ_{∞} and κ_{∞} . Thus, once a scaling (in powers of M_{∞}) is obtained for Re_{∞} and $P\acute{e}_{\infty}$, the corresponding normalization parameters
norm $_P^{\mu}$ and norm $_P^{\kappa}$ will automatically be set. For brevity, the superscripts * are
omitted in the remainder of this section.

For simplicity, we use here the ideal gas equation of state; its non-dimensionalized
expression is given by

$$P^* = (\gamma - 1) \rho^* \left(E^* - \frac{1}{2} M_{\infty}^2 (u^*)^2 \right) = (\gamma - 1) \rho^* e^*.$$
 (16)

In the low-Mach isentropic limit, shocks cannot form and the compressible
Euler equations are known to converge to the incompressible equations when the
Mach number tends to zero. When adding dissipative terms, as is the case with
the entropy viscosity method, the main properties of the low-Mach asymptotic
limit must be preserved. We begin by expanding each variable in powers of the
Mach number. As an example, the expansion for the pressure is given by:

$$P(\vec{r},t) = P_0(\vec{r},t) + P_1(\vec{r},t)M_{\infty} + P_2(\vec{r},t)M_{\infty}^2 + \dots$$
 (17)

By studying the resulting momentum equations for various powers of M_{∞} , it is

observed that the leading order and first-order pressure terms, P_0 and P_1 , are spatially constant if and only if $\text{Re}_{\infty} = \text{P\'e}_{\infty} = 1$. In this case, we have at order M_{∞}^{-2} :

$$\vec{\nabla}P_0 = 0 \tag{18a}$$

and at order M_{∞}^{-1}

$$\vec{\nabla}P_1 = 0. \tag{18b}$$

Using the scaling $\mathrm{Re}_{\infty}=\mathrm{P\acute{e}}_{\infty}=1$, the leading-order expressions for the continuity, momentum, and energy equations are:

$$\partial_t \rho_0 + \vec{\nabla} \cdot (\rho \vec{u})_0 = \vec{\nabla} \cdot (\kappa \vec{\nabla} \rho)_0 \tag{19a}$$

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$$\partial_t(\rho \vec{u})_0 + \vec{\nabla} \cdot (\rho \vec{u} \otimes \vec{u})_0 + \vec{\nabla} P_2 = \vec{\nabla} \cdot (\rho \mu \vec{\nabla}^s \vec{u} + \kappa \vec{u} \otimes \vec{\nabla} \rho)_0$$
 (19b)

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$$\partial_t(\rho E)_0 + \vec{\nabla} \cdot [\vec{u}(\rho E + P)]_0 = \vec{\nabla} \cdot (\kappa \vec{\nabla}(\rho e))_0$$
 (19c)

where the notation $(fg)_0$ means that we only keep the 0th-order terms in the product fg. The leading-order of the equation of state is given by

$$P_0 = (\gamma - 1)(\rho E)_0. {(20)}$$

Using Eq. (20), the energy equation can be recast as a function of the leadingorder pressure, P_0 , as follows:

$$\partial_t P_0 + \gamma \vec{\nabla} \cdot (\vec{u}P)_0 = \vec{\nabla} \cdot (\kappa \vec{\nabla}(P))_0. \tag{21}$$

From Eq. (18a), we infer that P_0 is spatially constant. Thus, Eq. (21) becomes

$$\frac{1}{\gamma P_0} \frac{dP_0}{dt} = -\vec{\nabla} \cdot \vec{u}_0 \tag{22}$$

262 and, at steady state, we have

$$\vec{\nabla} \cdot \vec{u}_0 = 0. \tag{23}$$

That is, the leading-order of velocity is divergence-free. The same reasoning can
be applied to the leading-order of the continuity equation (Eq. (19a)) to show
that the material derivative of the density variable is zero:

$$\frac{\mathrm{D}\rho_0}{\mathrm{D}t} := \partial_t \rho_0 + \vec{u}_0 \cdot \vec{\nabla} \cdot \rho_0 = 0.$$
 (24)

Therefore, we conclude that by setting the Reynolds and Péclet numbers to one, the incompressible fluid results are retrieved in the low-Mach limit when employing the compressible Euler equations with viscous regularization terms present. In addition, the scaling of the Prandtl number can also be obtained using Eq. (15), hence clarifying the use of the numerical Prandtl in the original entropy viscosity method [8].

272 3.3. Scaling of Re_{∞} and $P\acute{e}_{\infty}$ for non-isentropic flows

Next, we consider the non-isentropic case. Recall that even subsonic flows 273 can present shocks (for instance, a step initial condition in the pressure will trigger shock formation, independently of the Mach number). The non-dimensional 275 form of the Euler equations given in Eq. (13) provides some insight on the dom-276 inant terms as a function of the Mach number. This is particular obvious in 277 the momentum equation, Eq. (13b), where the gradient of pressure is scaled by 278 $1/M_{\infty}^2$. In the non-isentropic case, we no longer have $\frac{\vec{\nabla}P}{M_{\infty}^2} = \vec{\nabla}P_2$ and there-279 fore the pressure gradient term may need to be stabilized by some dissipative 280 terms of the same scaling so as to prevent spurious oscillations from forming. 281 By inspecting the dissipative terms presents in the the momentum equation, 282 having a dissipative term that scales as $1/M_{\infty}^2$ leads to the following three op-283 tions: (a) $\mathrm{Re}_{\infty}=M_{\infty}^2$ and $\mathrm{P\acute{e}}_{\infty}=1$, (b) $\mathrm{Re}_{\infty}=1$ and $\mathrm{P\acute{e}}_{\infty}=M_{\infty}^2$, or (c)

 $\text{Re}_{\infty} = \text{P\'e}_{\infty} = M_{\infty}^2$. Any of these choices will also affect the stabilization of the continuity and energy equations. For instance, using a Péclet number equal to M_{∞}^2 may effectively stabilize the continuity equation in the shock region but this may also add an excessive amount of dissipation for subsonic flows at the location of the contact wave. Such a behavior may not be suitable for accuracy 289 purpose, making options (b) and (c) inappropriate. The same reasoning, left to 290 the reader, can be carried out for the energy equation (Eq. (13c)) and results in 291 the same conclusion. The remaining choice, option (a), has the proper scaling: 292 in this case, only the dissipation terms involving $\vec{\nabla}^{s,*}\vec{u}^*$ scale as $1/M_{\infty}^2$ since 293 $\mathrm{Re}_{\infty} = M_{\infty}^2$, leaving the regularization of the continuity equation unaffected 294 because $P\acute{e}_{\infty} = 1$.

296 3.4. An All-speed normalization of the entropy residual

The study of the above limit cases yields two different possible scalings for the Reynolds number: $\text{Re}_{\infty}=1$ in the low-Mach limit and $\text{Re}_{\infty}=M_{\infty}^2$ for non-isentropic flows, whereas the numerical Péclet number always scales as one. In order to have a stabilization method valid for a wide range of Mach numbers, from very low-Mach to supersonic flows, these two scalings should be combined in a unique definition.

We begin with the normalization parameter norm $_P^{\kappa}$. Using the definition of the viscosity coefficients given in Eq. (10) and the scaling of Eq. (11), it can be shown that:

$$\kappa_{\infty} = \frac{\rho_{\infty} c_{\infty}^2 u_{\infty} L}{\text{norm}_{P,\infty}^{\kappa}}, \qquad (25)$$

where $\operatorname{norm}_{P,\infty}$ is the reference far-field quantity for the normalization parameter norm_P . Substituting Eq. (25) into Eq. (14) and recalling that the numerical Péclet number scales as unity, we obtain:

$$\operatorname{norm}_{P,\infty}^{\kappa} = \operatorname{P\acute{e}_{\infty}} \rho_{\infty} c_{\infty}^{2} = \rho_{\infty} c_{\infty}^{2}. \tag{26}$$

Eq. (26) provides a proper normalization factor to define the κ viscosity coefficient. The derivation for norm $_P^{\mu}$ is similar and yields

$$\operatorname{norm}_{P}^{\mu} = \operatorname{Re}_{\infty} \rho_{\infty} c_{\infty}^{2} = \begin{cases} \rho ||\vec{u}||^{2} & \text{for non-isentropic flows} \\ \rho c^{2} = \operatorname{norm}_{P}^{\kappa} & \text{for low-Mach flows} \end{cases} . (27)$$

A smooth function to transition between these two states is as follows:

$$\sigma(M) = \frac{\tanh\left(a(M - M^{\text{thresh}})\right) + |\tanh\left(a(M - M^{\text{thresh}})\right)|}{2}, \qquad (28)$$

where $M^{\rm thresh}$ is a threshold Mach number value beyond which the flow is no longer considered to be low-Mach (we use $M^{\rm thresh}=0.05$), M is the local Mach number, and the scalar a determines how rapidly the transition from norm $_P^\mu=\rho c^2$ to norm $_P^\mu=\rho \|\vec{u}\|^2$ occurs in the vicinity of $M^{\rm thresh}$ (we use a=3). It is easy to verify that

$$norm_P^{\mu} = (1 - \sigma(M))\rho c^2 + \sigma(M)\rho ||\vec{u}||^2$$
 (29)

satisfies Eq. (27). Finally, we summarize the definition of the viscosity coefficients μ and κ for completeness:

$$\kappa(\vec{r},t) = \min\left(\mu_{\max}(\vec{r},t), \kappa_e(\vec{r},t)\right), \tag{30a}$$

$$\mu(\vec{r},t) = \min\left(\mu_{\max}(\vec{r},t), \mu_e(\vec{r},t)\right), \tag{30b}$$

where the first-order viscosity is given by

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$$\kappa_{\text{max}}(\vec{r},t) = \mu_{\text{max}}(\vec{r},t) = \frac{h}{2} \left(||\vec{u}|| + c \right)$$
(30c)

and the entropy viscosity coefficients by

$$\kappa_e(\vec{r}, t) = \frac{h^2 \max(\widetilde{R}_{\text{ent}}, J)}{\rho c^2} \text{ and } \mu_e(\vec{r}, t) = \frac{h^2 \max(\widetilde{R}_{\text{ent}}, J)}{\text{norm}_D^{\mu}}$$
(30d)

with the jumps given by

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$$J = \max(||\vec{u}||[[\vec{\nabla}P \cdot \vec{n}]], ||\vec{u}||c^2[[\vec{\nabla}\rho \cdot \vec{n}]])$$
 (30e)

where norm $^{\kappa}_{P}$ is computed from Eq. (29). The jump J is a function of the jump of pressure and density gradients across the face with respect to its normal vector \vec{n} . Then, the largest value over all faces is determined and used in the definition of the viscosity coefficients. With the definition of the viscosity coefficients μ and κ proposed in Eq. (30), the dissipative terms are expected to scale appropriately for very low-Mach regimes as well for transonic and supersonic flows.

4. Extension of the entropy viscosity technique to Euler equations with variable area

Fluid flows in nozzles and in pipes of varying cross-sectional area can be modeled using the variable-area variant of the Euler equations, where the conservative variables are now multiplied by the area A. In addition, these equations differ from the standard Euler equations in that the momentum equation Eq. (31b) contains a non-conservative term proportional to the area gradient. Here, the variable area is assumed to be a smooth function of space only.

$$\partial_t (\rho A) + \vec{\nabla} \cdot (\rho \vec{u} A) = 0,$$
 (31a)

 $\partial_t (\rho \vec{u} A) + \vec{\nabla} \cdot [A (\rho \vec{u} \otimes \vec{u} + P \mathbb{I})] = P \vec{\nabla} A,$ (31b)

 $\partial_t (\rho E A) + \vec{\nabla} \cdot [\vec{u} A (\rho E + P)] = 0. \tag{31c}$

The application of the entropy viscosity method to the Euler equations with variable area is not fundamentally different to its application to the standard Euler equations. However, we need to derive the associated dissipative terms and verify that the entropy minimum principle is still satisfied. The variable-area Euler equations with viscous regularization are given below; details of the derivation are provided in Appendix B.

$$\partial_t (\rho A) + \vec{\nabla} \cdot (\rho \vec{u} A) = \vec{\nabla} \cdot \left(A \kappa \vec{\nabla} \rho \right) ,$$
 (32a)

 $\partial_t \left(\rho \vec{u} A \right) + \vec{\nabla} \cdot \left[A \left(\rho \vec{u} \otimes \vec{u} + P \mathbb{I} \right) \right] = P \vec{\nabla} A + \vec{\nabla} \cdot \left[A \left(\mu \rho \vec{\nabla}^s \vec{u} + \kappa \vec{u} \otimes \vec{\nabla} \rho \right) \right] , \tag{32b}$

$$\partial_{t} (\rho A E) + \vec{\nabla} \cdot [\vec{u} A (\rho E + P)] = \vec{\nabla} \cdot \left[A \left(\kappa \vec{\nabla} (\rho e) + \frac{1}{2} ||\vec{u}||^{2} \kappa \vec{\nabla} \rho + \rho \mu \vec{u} \vec{\nabla}^{s} \vec{u} \right) \right]. \quad (32c)$$

The dissipative terms are quite similar to the ones obtained for the standard

Euler equations: each dissipative flux is simply multiplied by the variable area A in order to ensure conservation of the dissipative flux. When assuming a constant area, Eqs. 2 are recovered.

A low-Mach asymptotic limit of the Euler equations with variable area on the same model as in Section 3.2 will lead to the divergence constraint $\nabla \cdot (\vec{u}A) = 0$ that can be recast as $\nabla \cdot \vec{u} = -\vec{u} \cdot \nabla A/A$. The gradient of the area acts as a source term and will force the fluid to accelerate or decelerate, depending on its sign.

5. Discretizations and Solution Techniques

In this section, we briefly describe the spatial and temporal discretizations and the solution techniques used to solve the system of equations Eq. (32). For conciseness, we re-write the system of equations in the following form:

$$\partial_t \mathbf{U} + \vec{\nabla} \cdot \vec{\mathbf{F}} (\mathbf{U}) = \mathbf{S} + \vec{\nabla} \cdot \mathbf{D}(\mathbf{U}) \vec{\nabla} \mathbf{U}$$
(33)

where $\mathbf{U} = [\rho A, \rho \vec{u} A, \rho E A]^T$ is the solution vector, \mathbf{F} denotes the inviscid flux

$$\vec{\mathbf{F}} \equiv \begin{bmatrix} \rho u A \\ (\rho u^2 + p) A \\ u(\rho E + P) A \end{bmatrix}$$
(34)

and **S** is a source term that contains the non-conservative term $P\vec{\nabla}A$. The term $\vec{\nabla} \cdot D(\mathbf{U})\vec{\nabla}\mathbf{U}$ stands for the artificial dissipative terms.

5.1. Spatial and Temporal Discretizations

The system of equations given in Eq. (33) is discretized using a continuous
Galerkin finite element method and temporal integrators available through the
MOOSE multiphysics framework [14].

5.1.1. Continuous Finite Elements

In order to apply the continuous finite element method, Eq. (33) is multiplied by a test function $\mathbf{W}(\vec{r})$, integrated by parts and each integral is decomposed into a sum of integrals over each element K of the discrete mesh Ω . The following weak form is obtained:

$$\sum_{K} \int_{K} \partial_{t} \mathbf{U} \mathbf{W} - \sum_{K} \int_{K} \vec{\mathbf{F}}(\mathbf{U}) \cdot \vec{\nabla} \mathbf{W} + \int_{\partial \Omega} \vec{\mathbf{F}}(\mathbf{U}) \cdot \vec{n} \mathbf{W} - \sum_{K} \int_{K} \mathbf{S} \mathbf{W} + \sum_{K} \int_{K} D(\mathbf{U}) \vec{\nabla} \mathbf{U} \cdot \vec{\nabla} \mathbf{W} - \int_{\partial \Omega} D(\mathbf{U}) \vec{\nabla} \mathbf{U} \cdot \vec{n} \mathbf{W} = 0. \quad (35)$$

The integrals over the elements K are evaluated using a numerical quadrature.

The MOOSE framework provides a wide range of test functions and quadrature

rules. Linear Lagrange polynomials are employed as test functions in the results section. Second-order spatial convergence will be demonstrated for smooth solutions.

372 5.1.2. Temporal integration

The MOOSE framework offers both first- and second-order explicit and implicit temporal integrators. In all of the numerical examples presented in Section 6, the temporal derivative will be evaluated using the second-order, backward difference temporal integrator BDF2. By considering three consecutive
solutions, \mathbf{U}^{n-1} , \mathbf{U}^n and \mathbf{U}^{n+1} , at times t^{n-1} , t^n and t^{n+1} , respectively, BDF2
can be expressed as:

$$\int_{K} \partial_{t} \mathbf{U} \mathbf{W} = \int_{K} \left(\omega_{0} \mathbf{U}^{n+1} + \omega_{1} \mathbf{U}^{n} + \omega_{2} \mathbf{U}^{n-1} \right) \mathbf{W}, \qquad (36)$$

with

$$\omega_0 = \frac{2\Delta t^{n+1} + \Delta t^n}{\Delta t^{n+1} \left(\Delta t^{n+1} + \Delta t^n\right)}, \ \omega_1 = -\frac{\Delta t^{n+1} + \Delta t^n}{\Delta t^{n+1} \Delta t^n},$$
 and
$$\omega_2 = \frac{\Delta t^{n+1}}{\Delta t^n \left(\Delta t^{n+1} + \Delta t^n\right)}$$

where $\Delta t^n = t^n - t^{n-1}$ and $\Delta t^{n+1} = t^{n+1} - t^n$.

380 5.2. Boundary conditions

Boundary conditions are implemented by performing a characteristic decomposition to compute the appropriate flux at the boundaries. Our implementation
of the subsonic boundary conditions is inspired by the method described in [20]
and was adapted for a time implicit solver. Neumann boundary conditions are
used for all of the boundary types, except for the inlet supersonic boundary that
are strongly imposed with Dirichlet boundary conditions.

For each numerical solution presented in Section 6, the type of boundary conditions used will be specified and taken among the following: supersonic inlet,
subsonic inlet (stagnation pressure boundary), subsonic outlet, and supersonic
outlet. The artificial diffusion coefficient $D(\mathbf{U})$ is set to zero at the boundary of
the computational domain so that the boundary term $\int_{\partial\Omega} D(\mathbf{U}) \vec{\nabla} \mathbf{U} \cdot \vec{n} \mathbf{W}$ stemming from the integration by parts of the artificial dissipative terms in Eq. (35)
is ignored.

5.3. Solver

A Jacobian-free-Newton-Krylov (JFNK) method is used to solve for the so-395 lution at the end of each time step. An approximate Jacobian matrix of the discretized equations was derived and implemented. Obtaining the matrix en-397 tries requires that the partial derivatives of pressure with respect to the conservative variables be known (this is relatively simple for the stiffened and ideal 399 gas equations of state but may be more complex for general equations of state). 400 The contributions of the artificial dissipative terms to the Jacobian matrix are 401 approximated by lagging the viscosity coefficients (computing them with the 402 previous solution). For instance, this is shown in Eq. (37) for the dissipative 403 terms present in the continuity equation: 404

$$\frac{\partial}{\partial \mathbf{U}} \left(\kappa \vec{\nabla} \cdot \rho \vec{\nabla} W \right) \simeq \kappa \frac{\partial}{\partial \mathbf{U}} \left(\vec{\nabla} \cdot \rho \vec{\nabla} W \right), \tag{37}$$

where **U** denotes any of the conservative variables and W denotes the component of **W** associated with the continuity equation. In the above, we have neglected $\frac{\partial \kappa}{\partial \mathbf{U}}$.

₀₈ 6. Numerical Results

1-D and 2-D numerical solutions for the Euler equations with viscous reg-409 ularization solved using the entropy viscosity method are presented here. Our 410 results show that the new definitions for the viscosity coefficients are robust in 411 the low-Mach limit as well as for for transonic and supersonic flows and that shocks are appropriately resolved. 413 The first set of 1-D simulations consist of liquid water and steam flowing in 414 a converging-diverging nozzle. This test is of interest for multiple reasons: (a) 415 a steady state can be reached (some stabilization methods are known to have 416 difficulties reaching a steady state, [2, 3]), (b) an analytical solution is available 417 and a space-time convergence study can be performed, (c) it can be performed 418 for liquid and gas phases, wherein the gas phase simulation presents a shock 419 while the liquid-phase simulation has a significantly lower Mach number. Next, 420 a 1-D shock tube test (in a straight pipe), taken from the Leblanc test-case suite 42 [21], is performed. This test is known to be more challenging than Sod shock 422 tubes and the fluid's Mach number varies spatially between 0 and 5. A convergence study is also performed to demonstrate convergence of the numerical 424 solution to the exact solution. A slow moving shock is also investigated [22]. 425 This test helps in assessing the ability of the method to damp the post-shock 426 low frequency noise (oscillations). Finally, a strong shock for a liquid phase 427 (Mach number around 0.1) is also performed [23]. 428 The initial conditions for the aforementioned 1-D test cases are given in 429 Table 1. 430 The 2-D simulations are outlined next. First, 2-D subsonic flows around a 431 cylinder [12] and over a circular hump [24] are presented for various far-field Mach numbers (as low of 10^{-7}). Numerical results of a supersonic flow over 433

a compression corner are provided to illustrate the ability of the new viscosity

$\rho_{ m left}$	$u_{ m left}$	P_{left}	$ ho_{ m right}$	$u_{\rm right}$	P_{right}					
	Leblanc shock tube (Section 6.3)									
1	0	$4 \ 10^{-2}$	10^{-3}	0	$4 \ 10^{-11}$					
St	rong sho	ck for liqu	id phase	e (Section	n 6.4)					
1000	0	10^{9}	1000	0	10^{5}					
Slow moving shock (Section 6.5)										
1	-0.81	1	3.86	-3.44	10.33					

Table 1: Initial conditions for the 1-D test cases (density in kg/m^3 , velocity in m/s, pressure in Pa).

definitions to handle supersonic flows. Convergence studies are performed when analytical solutions are available.

For each simulation, data relative to the boundary conditions, the Courant-Friedrichs-Lewy number (CFL), mesh and equation of state are provided. All of the numerical solutions presented are obtained using BDF2 as temporal integrator and linear (1-D mesh), \mathbb{P}_1 (2-D triangular mesh), and \mathbb{Q}_1 (2-D quadrangular mesh) finite elements. The spatial integrals are numerically computed using a second-order Gauss quadrature rule. Steady-state is detected in a transient simulation by monitoring the nonlinear residual before proceeding with the Newton solves for a given time step. The ideal gas [25] or stiffened gas equations of state [26] are used; a generic expression is given in Eq. (38).

$$P = (\gamma - 1)\rho(e - q) - \gamma P_{\infty} \tag{38}$$

where the parameters γ , q, and P_{∞} are fluid-dependent and are given in Table 2. The ideal gas equation of state is recovered by setting $q = P_{\infty} = 0$ in Eq. (38). The entropy function for the stiffened gas equation of state is concave and given by

$$s = C_v \ln \left(\frac{P + P_{\infty}}{\rho^{\gamma - 1}} \right),$$

where C_v is the heat capacity at constant volume.

Table 2: Stiffened Gas Equation of State parameters for steam and liquid water.

fluid	γ	$C_v (J.kg^{-1}.K^{-1})$	P_{∞} (Pa)	$q (J.kg^{-1})$
liquid water (Section 6.1)	2.35	1816	10^{9}	$-1167 \ 10^3$
steam (Section 6.2)	1.43	1040	0	$2030 \ 10^3$
liquid water (Section 6.4)	4.4	1000	$6 \ 10^8$	0

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Finally, the convergence rates are computed using the following relation

$$rate_h = \ln\left(\frac{||U_{2h} - U_{\text{exact}}||}{||U_h - U_{\text{exact}}||}\right) / \ln 2$$
(39)

where $||\cdot||$ denotes either the L₁ or L₂ norms and h is the characteristic grid size.

6.1. Liquid water in a 1-D converging-diverging nozzle

A simulation for liquid flow through a 1-D converging-diverging nozzle is 452 performed. The variable area expression is given by $A(x) = 1 + 0.5\cos(2\pi x/L)$ 453 with length L=1m. At the inlet, the stagnation pressure and temperature are 454 set to $P_0 = 1MPa$ and $T_0 = 453K$, respectively. At the outlet, only the static 455 pressure is specified: $P_s = 0.5MPa$. Initially, the liquid is at rest, the temperature is uniform and equal to the stagnation temperature and the pressure 457 linearly decreases from the stagnation pressure inlet value to the static pressure 458 outlet value. The stiffened gas equation of state is used to model the liquid 459 water with the parameters provided in Table 2. Because of the low pressure difference between the inlet and the outlet, the smooth initial conditions, and 461 the large value of P_{∞} in Eq. (38), the flow remains subsonic and thus displays no shock. A detailed derivation of the exact steady-state solution can be found 463 in [27]. A uniform mesh of 50 cells was used to obtain the numerical solution 464 and the time step size was computed using a CFL number of 750. Plots of the Mach number, density, and pressure are given at steady state in Fig. 1 for the numerical and exact solutions. The viscosity coefficients are also graphed in Fig. 1d.

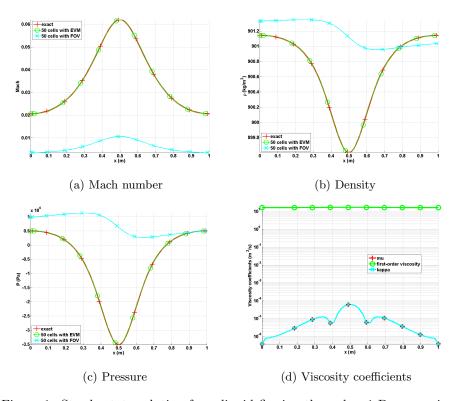


Figure 1: Steady-state solution for a liquid flowing through a 1-D converging-diverging nozzle.

In Fig. 1, the numerical solutions obtained using the first-order viscosity (FOV) and the entropy viscosity method (EVM) are plotted against the exact solution. The numerical solution obtained with the EVM and the exact solution overlap, even for a fairly coarse mesh (50 cells). On the other hand, the numerical solution obtained with the FOV does not give the correct steady state: this is an illustration of the effect of ill-scaled dissipative terms in the low-Mach limit when using the FOV. Note that the entropy viscosity coefficient is very small

compared to the first-order one (Fig. 1d): (i) the numerical solution is smooth as shown in Fig. 1 and (ii) the flow is in a isentropic low-Mach regime A convergence study was performed using the exact solution as a reference: the L₁ and L₂ norms of the error and the corresponding convergence rates are computed at steady state on various uniform meshes from 4 to 256 cells. Spatial convergence results using linear finite elements are reported in Table 3 and Table 4 for the primitive variables: density, velocity and pressure.

Table 3: L_1 norm of the error for the liquid phase in a 1-D converging-diverging nozzle at steady state.

cells	density	rate	pressure	rate	velocity	rate
4	$2.8037 \ 10^{-1}$	_	$8.4705 \ 10^5$	_	7.2737	_
8	$1.3343 \ 10^{-1}$	1.07	$4.7893 \ 10^5$	0.82	6.1493	0.24
16	$2.9373 \ 10^{-2}$	2.18	$1.0613 \ 10^5$	2.17	1.2275	2.32
32	$5.1120 \ 10^{-3}$	2.52	$1.8446 \ 10^4$	2.52	$1.8943 \ 10^{-1}$	2.69
64	$1.0558 \ 10^{-3}$	2.28	$3.7938 \ 10^3$	2.28	$3.7919 \ 10^{-2}$	2.32
128	$2.3712 \ 10^{-4}$	2.15	$8.4471 \ 10^2$	2.17	$8.5517 \ 10^{-3}$	2.15
256	$5.6058 \ 10^{-5}$	2.08	$1.9839 \ 10^2$	2.09	$2.0475 \ 10^{-3}$	2.06
512	$1.3278 \ 10^{-5}$	2.08	$4.6622 \ 10^{1}$	2.09	$4.9516 \ 10^{-4}$	2.04
1024	$3.1193 \ 10^{-6}$	2.08	$1.1755 \ 10^{1}$	1.99	$1.2379 \ 10^{-4}$	2.00

Table 4: L_2 norm of the error for the liquid phase in a 1-D converging-diverging nozzle at steady state.

cells	density	rate	pressure	rate	velocity	rate
4	$3.106397 \ 10^{-1}$	_	$5.254445 \ 10^5$	_	3.288543	_
8	$7.491623 \ 10^{-2}$	2.05	$1.636966 \ 10^5$	1.68	1.823880	0.85
16	$2.079858 \ 10^{-2}$	1.85	$4.627338 \ 10^4$	1.49	$4.990605 \ 10^{-1}$	0.87
32	$5.329627 \ 10^{-3}$	1.96	$1.180287 \ 10^4$	1.97	$1.261018 \ 10^{-1}$	1.98
64	$1.341583 \ 10^{-3}$	1.99	$2.967104 \ 10^3$	1.99	$3.160914 \ 10^{-2}$	1.99
128	$3.359766 \ 10^{-4}$	1.99	$7.428087 \ 10^2$	1.99	$7.907499 \ 10^{-3}$	1.99
256	$8.403859 \ 10^{-5}$	1.99	$1.857861 \ 10^2$	1.99	$1.977292 \ 10^{-3}$	1.99
512	$2.10075 \ 10^{-5}$	2.00	$4.7024 \ 10^{1}$	1.98	$4.9516 \ 10^{-4}$	1.99

We note that the convergence rates measured in both the L_1 and L_2 norm of the error are equal to 2; the entropy viscosity method preserves the high-order accuracy of the discretization used when the numerical solution is smooth. The new definition of the entropy viscosity coefficients behaves appropriately in the low-Mach limit.

488 6.2. Steam in a 1-D converging-diverging nozzle

We use the same nozzle geometry, initial conditions and boundary condi-489 tions as in the previously example but replace liquid water with steam and use 490 the steam parameters of the stiffened gas equation of state, Table 2. In this 491 example, compressible effects will become dominant. The pressure difference 492 between the inlet and outlet is large enough to accelerate the steam through 493 the nozzle, leading to the formation of a shock in the diverging portion of the 494 nozzle. The behavior is different from the one observed for the liquid water 495 phase in Section 6.1 because of the liquid to gas density ratio is about 1,000. 496 An exact solution at steady state is available for the gas phase [27]. The aim 497 of this section is to show that when using the new definitions of the viscosity 498 coefficients (Eq. (30)), the shock can be correctly resolved without spurious os-499 cillations. The steady-state numerical solution, obtained using a uniform mesh 500 with 500 cells, is shown in Fig. 2. The CFL was set to 80 (a high CFL value 501 can be used because the shock is stationary). 502

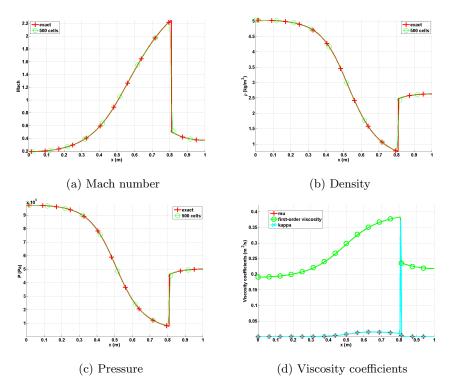


Figure 2: Steady-state solution for vapor phase flowing in a 1-D convergingdiverging nozzle.

The steady-state solution of the density, Mach number and pressure are 503 given in Fig. 2. The steady-state solution exhibits a shock around x = 0.8mand matches the exact solution. In Fig. 2d, the first-order and entropy viscosity 505 coefficients are plotted at steady state (on a log scale): the entropy viscosity coefficient is peaked in the shock region around x = 0.8m where it saturates to the first-order viscosity coefficient. Elsewhere, the entropy viscosity coefficient is small. In order to prove convergence of the numerical solution to the exact 509 solution, a convergence study is performed. Because of the presence of a shock, 510 second-order accuracy is not expected and the convergence rate of a numerical 511 solution should be 1 and 1/2 when measured in the L_1 and L_2 norms, respec-512 tively (see Theorem 9.3 in [28]). Results are reported in Table 5 and Table 6 513

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for the primitive variables: density, velocity and pressure. The convergence rates for the L₁ and L₂ norms of the error computed using Eq. (39) are in good agreement with the theoretical values.

Table 5: L_1 norm of the error for the vapor phase in a 1-D converging-diverging nozzle at steady state.

cells	density	rate	pressure	rate	velocity	rate
5	$0.72562 \ 10^{-1}$	_	$1.5657 \ 10^5$	_	173.69	_
10	$0.4165 \ 10^{-1}$	0.80	$9.6741 \ 10^4$	0.63	120.69	0.53
20	$0.20675 \ 10^{-1}$	1.01	$4.9193 \ 10^4$	0.97	72.149	0.74
40	$0.093703 \ 10^{-1}$	1.14	$2.0103 \ 10^4$	0.73	34.716	1.06
80	$0.047328 \ 10^{-1}$	0.99	$1.0208 \ 10^4$	0.98	16.082	1.11
160	$0.023965 \ 10^{-2}$	0.98	$5.1969 \ 10^3$	0.97	7.9573	1.02
320	$0.020768 \ 10^{-2}$	1.03	$2.5116 \ 10^3$	1.05	3.7812	1.07
640	$0.0059715 \ 10^{-2}$	0.98	$1.2754 \ 10^3$	0.98	1.8353	1.04

Table 6: L_2 norm of the error for the vapor phase in a 1-D converging-diverging nozzle at steady state.

cells	density	rate	pressure	rate	velocity	rate
5	$9.7144 \ 10^{-1}$	_	$2.0215 \ 10^5$	_	236.94	_
10	$5.9718 \ 10^{-1}$	0.70	$1.3024 \ 10^5$	0.63	166.56	0.51
20	$2.9503 \ 10^{-1}$	1.02	$6.6503 \ 10^4$	0.97	103.36	0.69
40	$1.8193 \ 10^{-1}$	0.69	$4.0171 \ 10^4$	0.73	66.374	0.64
80	$1.3366 \ 10^{-1}$	0.44	$2.3163 \ 10^4$	0.44	42.981	0.63
160	$9.6638 \ 10^{-2}$	0.47	$1.7263 \ 10^4$	0.42	31.717	0.44
320	$7.0896 \ 10^{-2}$	0.45	$1.2763 \ 10^4$	0.44	23.138	0.45
640	$5.2191 \ 10^{-2}$	0.44	$9.4217 \ 10^3$	0.44	16.910	0.45

517 6.3. Leblanc shock tube

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The 1-D Leblanc shock tube is a Riemann problem designed to test the robustness and the accuracy of stabilization methods. The initial conditions are given in Table 1. The ideal gas equation of state (with $\gamma = 5/3$) is used to compute the pressure. This test is computationally challenging because of the large pressure ratio at the initial interface. The computational domain consists of a 1-D straight pipe of length L = 9m with the initial interface located at

x=2m. At $t=0\,s$, the interface is removed. The numerical solution is run until $t=4\,s$ and the density, momentum and total energy profiles are given in Fig. 3, along with the exact solution. The viscosity coefficients are also plotted in Fig. 3d. These plots were run with three different uniform meshes of 800, 3200, and 6000 cells and a constant CFL=1.

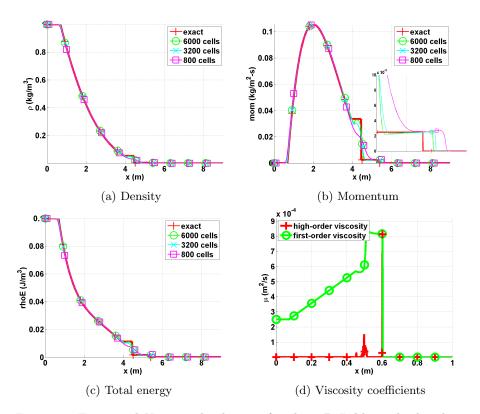


Figure 3: Exact and Numerical solutions for the 1-D Leblanc shock tube at $t=4\,s.$

The density, momentum and total energy profiles are provided in Fig. 3.

In Fig. 3b, the shock region is zoomed in for better resolution: the shock is
well resolved. We also observe that the shock position computed numerically
converges to the exact position under mesh refinement. The contact wave at x = 4.5m can be seen in Fig. 3b. The entropy viscosity coefficient profile is

shown in Fig. 3d and behaves as expected: it saturates to the first-order viscosity in the shock region, thus preventing oscillations from forming. At the location of the contact wave, a smaller peak is observed and is due to the presence of the jump terms in the definition of the entropy viscosity coefficient (Eq. (30)). The Mach number, not plotted, is of the order of 1.3 just before the shock and reaches a maximum value close to 5 in the contact region.

Once again, a convergence study is performed in order to prove convergence

Once again, a convergence study is performed in order to prove convergence
of the numerical solution to the exact solution. As in the previous example
(vapor phase in the 1-D nozzle, Section 6.2), the expected convergence rates
in the L₁ and L₂ norms are 1 and 1/2, respectively. The exact solution was
obtained by running a 1-D Riemann solver and used as the reference solution to
compute the L₁ and L₂-norms that are reported in Table 7 and Table 8 for the
conservative variables: density, momentum and total energy. The convergence
rates are again approaching their theoretical values.

Table 7: L₁ norm of the error for the 1-D Leblanc test at t = 4s.

cells	density	rate	momentum	rate	total energy	rate
100	$1.0354722 \ 10^{-2}$	_	$3.5471714 \ 10^{-3}$	_	$1.4033046 \ 10^{-3}$	_
200	$7.2680512 \ 10^{-3}$	0.51	$2.5933119 \ 10^{-3}$	0.45	$9.8611746 \ 10^{-4}$	0.51
400	$5.0825628 \ 10^{-3}$	0.52	$2.0668092 \ 10^{-3}$	0.33	$7.7844421 \ 10^{-4}$	0.34
800	$3.4025056 \ 10^{-3}$	0.58	$1.4793838 \ 10^{-3}$	0.48	$5.5702549 \ 10^{-4}$	0.48
1600	$2.1649953 \ 10^{-3}$	0.65	$9.7152832 \ 10^{-4}$	0.61	$3.5720171 \ 10^{-4}$	0.64
3200	$1.2465433 \ 10^{-3}$	0.79	$5.5937409 \ 10^{-4}$	0.79	$2.0491799 \ 10^{-4}$	0.80
6400	$6.4476928 \ 10^{-4}$	0.95	$3.0244198 \ 10^{-4}$	0.89	$1.0914891 \ 10^{-4}$	0.91
12800	$3.3950948 \ 10^{-4}$	0.93	$1.5958118 \ 10^{-4}$	0.92	$5.7909794 \ 10^{-5}$	0.91

6.4. 1-D shock tube with a liquid phase

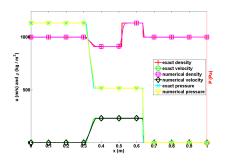
The purpose of this test is to investigate the ability of the entropy viscosity method to stabilize a strong shock with a small Mach number [23] (this reference is for a two-phase flow model but we are only interested in the initial conditions for the liquid phase): the Mach number in the shock region is of the order of 0.1.

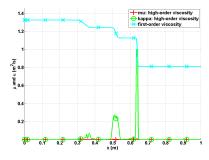
Table 8: L₂ norm of the error for the 1-D Leblanc test at t = 4s.

cells	density	rate	momentum	rate	total energy	rate
100	$5.7187851 \ 10^{-3}$	_	$1.7767236 \ 10^{-3}$	_	$7.6112265 \ 10^{-4}$	_
200	$3.8995238 \ 10^{-3}$	0.55	$1.4913161 \ 10^{-3}$	0.25	$5.5497308 \ 10^{-4}$	0.46
400	$2.8103526 \ 10^{-3}$	0.47	$1.3305301 \ 10^{-3}$	0.16	$4.6063172 \ 10^{-4}$	0.27
800	$2.1081933 \ 10^{-3}$	0.41	$1.1398931 \ 10^{-3}$	0.22	$3.7798953 \ 10^{-4}$	0.29
1600	$1.5731052 \ 10^{-3}$	0.42	$9.0394227 \ 10^{-4}$	0.33	$2.9584646 \ 10^{-4}$	0.35
3200	$1.0610667 \ 10^{-3}$	0.57	$6.2735595 \ 10^{-4}$	0.53	$2.054455 \ 10^{-4}$	0.53
6400	$7.3309974 \ 10^{-4}$	0.53	$4.4545754 \ 10^{-4}$	0.49	$1.4670834 \ 10^{-4}$	0.49
12800	$5.1020991 \ 10^{-4}$	0.52	$3.1266758 \ 10^{-4}$	0.51	$1.0299897 \ 10^{-5}$	0.51

In this case, as explained in Section 3.2, the viscosity coefficients are required to have different order of magnitude in order to ensure the correct scaling of the dissipative terms. The purpose of this test is to validate the approach presented in Section 3.2.

The stiffened gas equation of state is used to model a liquid flow with the 557 parameters given in Table 2. The computational domain of length L=1m is 558 uniformly discretized using 500 cells. The step initial conditions are given in 559 Table 1. The simulation is run with a CFL = 1 until the final time $t_{\text{final}} =$ 560 $7 ext{ } 10^{-5}s$. Results for pressure, density, velocity and the viscosity coefficients 561 are given in Fig. 4 along with the exact solution for comparison purposes. The 562 numerical solution is in good agreement with exact solution in Fig. 4a. The 563 viscosity coefficients μ and κ are not equal in the shock because the Mach number is of order 0.1. The viscosity coefficient κ saturates to the first-order 565 viscosity in the shock region around x = 0.65m and is sufficient to stabilize the numerical scheme. 567





- (a) Density, velocity and pressure profiles.
- (b) Viscosity coefficients profile.

Figure 4: Numerical solution for the 1-D liquid shock tube at at $t_{\text{final}} = 7 \cdot 10^{-5} s$.

6.5. 1-D slow moving shock

Slow moving shocks are known to produce post-shock noise of low frequency 569 that is not damped by some numerical dissipation methods [22]. The aim of 570 this simulation is to test the ability of the entropy viscosity method to dampen 57 the low frequency waves. The 1-D slow moving shock consists of a shock wave 572 moving from left to right with the initial conditions given in Table 1. The ideal 573 gas equation of state is used with a heat capacity ratio $\gamma = 1.4$. In order to make 574 the shock travel a significant distance, the final time is taken equal to t = 1.1 s. 575 A pressure boundary condition is used at the left boundary to let the rarefaction 576 and contact waves exit the domain. The numerical solution, obtained with 200 577 equally-spaced cells, is given in Fig. 5 and is compared to the exact solution 578 obtained from a Riemann solver. We use a CFL of 1. With this CFL value, 579 it takes about 50 time steps for the shock to traverse one cell. The numerical 580 results are in good agreement with the exact solution and do not display any 581 post-shock noise. The rarefaction and contact waves are not visible on Fig. 5a since they exited the computational domain through the left pressure boundary 583 condition earlier. As explained in [29], Godunov's type methods usually fail to resolve a slow moving shock because of the nature of the stabilization method: the method scales as the eigenvalue of the appropriate field. In the case of a slow moving shock, the dissipation added to the system is under-estimated and leads to post-shock noise. In the case of the entropy viscosity method, the entropy residual detects the shock position and the viscosity coefficients saturate to the first-order viscosity values in the shock region. The main difference between a Godunov's type method and the entropy viscosity method lies in the definition of the first-order viscosity coefficients that are proportional to the *local maximum* eigenvalue $||\vec{u}|| + c$ and not to the eigenvalue of the characteristic field.

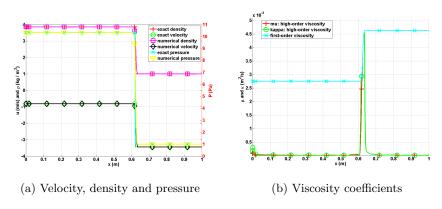


Figure 5: Slow moving shock profiles at t = 1.1s.

6.6. Subsonic flow over a 2-D cylinder

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Fluid flow over a 2-D cylinder is often used as a benchmark case to test numerical schemes in the low-Mach regime [10, 11, 12]. For this test, an analytical solution is available in the incompressible limit and is often referred to as the potential steady-state flow solution. The main features of the potential flow are the following:

- The solution is symmetric: the iso-Mach contour lines are used to assess the symmetry of the numerical solution;
- The velocity at the top of the cylinder is twice the incoming velocity set

 at the inlet;

 The spatial steady-state pressure variations are proportional to the square of inlet Mach number, i.e.,

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$$\delta P = \frac{\max(P(\vec{r})) - \min(P(\vec{r}))}{\max(P(\vec{r}))} \propto M_{\infty}^{2}$$
(40)

where δP and M_{∞} denote the spatial steady-state pressure variations and the inlet Mach number, respectively.

The computational domain consists of a 1×1 square with a circular hole of radius 0.05 in its center. A \mathbb{P}_1 triangular mesh with 4008 triangular elements is employed to discretize the geometry. The ideal gas equation of state, with $\gamma = 1.4$ is used. At the inlet, a subsonic stagnation boundary condition is used: the stagnation pressure and temperature are computed using the following relations:

$$\begin{cases}
P_0 = P \left(1 + \frac{\gamma - 1}{2} M^2 \right)^{\frac{\gamma - 1}{\gamma}} \\
T_0 = T \left(1 + \frac{\gamma - 1}{2} M^2 \right)
\end{cases}$$
(41)

A static pressure boundary condition, with static pressure $P_s = 101,325 \ Pa$, 614 is set at the outlet boundary. The implementation of the pressure boundary 615 conditions is based on [20]. A solid wall boundary condition is set for the top 616 and bottom walls of the computational domain. The simulations are run until 617 a steady state is reached (with a CFL of 40). When the residual norm (for all 618 equations) is less than 10^{-12} the steady state is considered to have been reached. 619 Several simulations are performed, with inlet Mach numbers M_{inlet} ranging 620 from 10^{-3} to 10^{-7} , and are shown in Fig. 6. The iso-Mach contour lines are 621 drawn using 30 equally-spaced intervals, from 2×10^{-10} to $M_{\rm inlet}$.

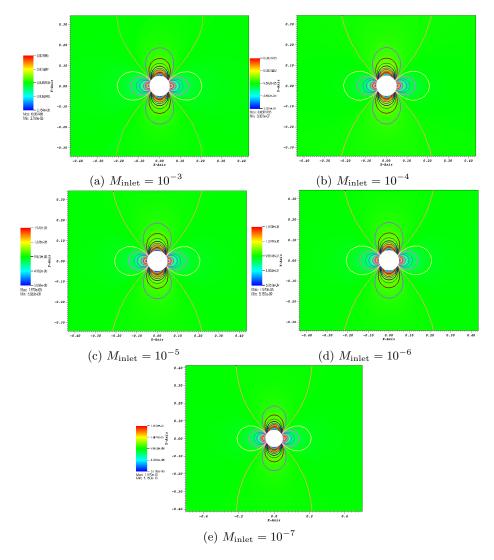


Figure 6: Iso-Mach lines for a subsonic flow over a 2-D cylinder with inlet Mach number values from ranging from 10^{-3} to 10^{-7} (steady-state solution).

The velocity at the top of the cylinder and at the inlet are given for different Mach-number values (ranging from 10^{-3} to 10^{-7}) in Table 9. The ratio of the inlet velocity to the velocity at the top of cylinder is also computed and is very close to the theoretical value of 2 that is expected in the incompressible limit.

Table 9: Velocity ratio for different Mach numbers.

Mach number	inlet velocity	velocity at the top of the cylinder	ratio
10^{-3}	$2.348 \ 10^{-3}$	$1.176 \ 10^{-3}$	1.99
10^{-4}	$2.285 \ 10^{-4}$	$1.145 \ 10^{-4}$	1.99
10^{-5}	$2.283 \ 10^{-5}$	$1.144 \ 10^{-5}$	1.99
10^{-6}	$2.283 \ 10^{-6}$	$1.144 \ 10^{-6}$	1.99
10^{-7}	$2.283 \ 10^{-7}$	$1.144 \ 10^{-7}$	1.99

In Fig. 7, the spatial steady-state variations in pressure and velocity are 627 plotted as a function of the Mach number (on a log-log scale). The spatial 628 steady-state pressure variations are expected to be of the order of M^2 in the 629 incompressible limit, which we observe. From Bernoulli's principle, this implies 630 that the velocity spatial variations should be of order M in the incompressible 631 limit, which we also observe in Fig. 7. It is known that some stabilization 632 methods, e.g., [10, 11, 12], can produce spatial steady-state pressure variations 633 with the wrong Mach-number order. Here, the entropy viscosity method yields the correct orders in the low-Mach limit. For ease of comparison, reference lines 635 with slope values of 1 and 2 are also plotted.

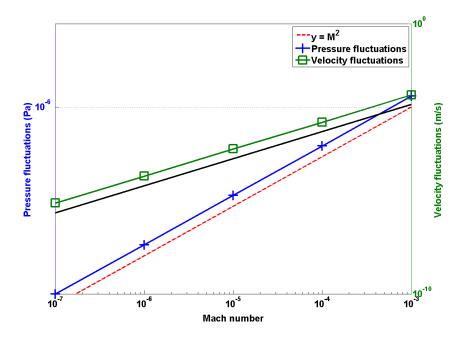


Figure 7: Log-log plot of the spatial steady-state pressure and velocity variations as a function of the far-field Mach number.

6.7. Subsonic flow over a 2-D hump

This is a another example of an internal flow configuration. It consists of 638 a channel of height L=1 m and length 3L, with a circular bump of length L639 and thickness 0.1L. The bump is located on the bottom wall at a distance L 640 from the inlet. The system is initialized with an uniform pressure P = 101,325641 Pa and temperature T = 300 K. The initial velocity is computed from the 642 inlet Mach number, the pressure, the temperature and the ideal gas equation (with $\gamma = 1.4$). Here, $C_v = 717 \ J/kg - K$. At the inlet, a subsonic stagnation 644 boundary condition is used and the stagnation pressure and temperature are computed using Eq. (41). The static pressure $P_s = 101,325 \ Pa$ is set at the 646 subsonic outlet. The results are shown in Fig. 8a, Fig. 8b, Fig. 8c and Fig. 8d for the inlet Mach numbers $M_{\infty}=0.7,\,M_{\infty}=0.01,\,M_{\infty}=10^{-4}$ and $M_{\infty}=10^{-7},$

respectively. It is expected that, for low Mach numbers, the solution does not 649 depend on the Mach number and is identical to the incompressible flow solution. 650 On the other hand, for a flow with M = 0.7, the compressible effects become 651 non negligible and a shock can form. An uniform grid of 3352 \mathbb{Q}_1 elements was 652 used to obtain the numerical solution for Mach numbers less than and equal 653 to $M_{\infty}=0.01$. A spatial mesh, once refined, was employed for the $M_{\infty}=0.7$ 654 simulation in order to better resolve the shock. A CFL of 20 was employed and 655 the simulations were run until steady state. 656

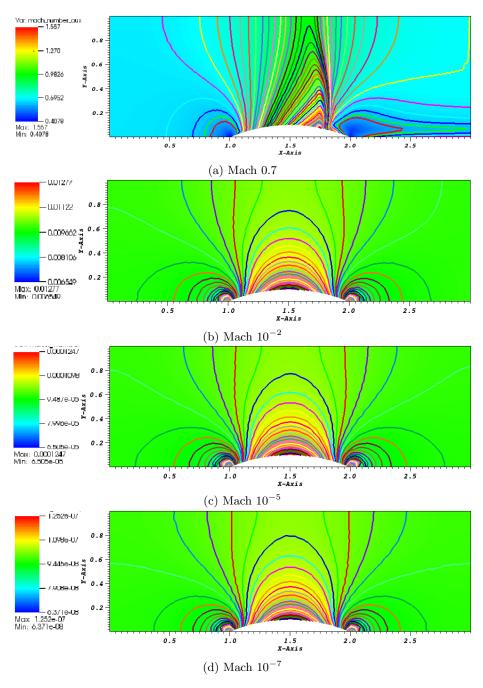


Figure 8: Iso-Mach lines for a 2-D flow over a circular bump (steady-state solution).

The results shown in Fig. 8b, Fig. 8c and Fig. 8d correspond to the lowMach regime. The iso-Mach lines are drawn ranging from the minimum and the
maximum values (provided in each legend) using 50 equally-spaced intervals.
The steady-state solution is symmetric and does not depend on the value of the
inlet Mach number, as expected in the incompressible limit.

In Fig. 8a, the steady-state numerical solution develops a shock: the compressibility effects are no longer small. The iso-Mach lines are also plotted with 50 intervals and range from 0.4 to 1.6. The shock is well resolved and does not display any instabilities or spurious oscillations.

666 6.8. Supersonic flow in a compression corner

In this last example, we consider a supersonic flow at Mach 2.5 impinging 667 on a corner with an angle of 15°. From the oblique shock theory [16], an 668 analytical solution for this supersonic flow is available and gives the downstream-669 to-upstream pressure, entropy and Mach number ratios. The initial conditions 670 are chosen to be spatially uniform: the pressure and temperature are set to P =671 $101,325 \ Pa \ and \ T = 300 \ K$, respectively. The ideal gas equation of state is used 672 with the same parameters as in Section 6.7. The initial velocity is computed from 673 the upstream Mach number. The inlet is supersonic and therefore, the pressure, 674 temperature and velocity are specified using Dirichlet boundary conditions. The outlet is also supersonic and none of the characteristics enter the domain through 676 this boundary; the values are computed by the solver.

The simulation is run with CFL=2 until steady state is reached. A 2-D mesh made of 16, 109 \mathbb{Q}_1 elements is used. The ratios for pressure, entropy and Mach number computed using the analytical (published with only two significant digits) and the numerical solutions are given in Table 10; they are in excellent agreement. The shock wave angle at steady state is also known and given by

the so-called $\theta - \beta - M$ relation:

$$\tan \theta = 2 \cot \beta \frac{M^2 \sin^2 \beta - 1}{M^2 (\gamma + \cos^2(2\beta)) + 2},$$
(42)

where θ , β and M denote the corner angle, the shock wave angle, and the upstream Mach number, respectively. For Mach 2.5 and a 15° corner angle, the analytical value for the shock wave angle is 36.94° at steady state. From Fig. 9a, the numerical value of the shock wave angle can be measured and is found to be equal to 36.9° and thus is in excellent agreement with the theory.

	analytical	numerical
Pressure	2.47	2.467
Mach number	0.74	0.741
Entropy	1.03	1.026

Table 10: Ratio of analytical and numerical downstream to upstream quantities for the compression corner problems (corner angle of 15° and inlet M=2.5 (analytical values from [16]).

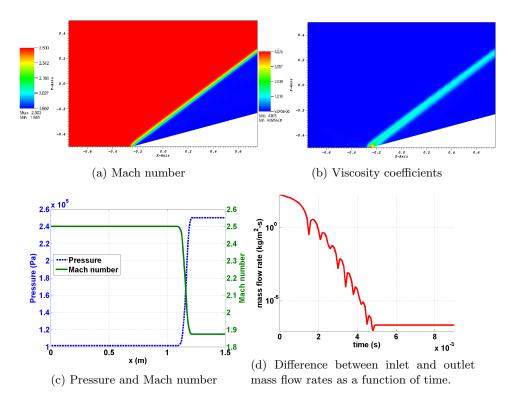


Figure 9: Steady-state solution for a flow in a 2-D compression corner.

The steady-state numerical solution is given in Fig. 9; the Mach number and the viscosity coefficients are plotted in Fig. 9a and Fig. 9b, respectively. The steady-state solution is composed of two regions of constant state separated by an oblique shock. Fig. 9b shows that the viscosity coefficient is large in the shock and small elsewhere, as expected. At the location of the corner (x = -0.25m, y = -0.5m), the viscosity coefficient is peaked because of the treatment of the wall boundary condition: at this particular node, the normal is not well defined and may cause some numerical errors. The 1-D graphs at y = 0 for the pressure and the Mach number are given in Fig. 9c: no spurious oscillations are observed and the shock is well resolved. Finally, the difference between the inlet and outlet mass flow rates is plotted in Fig. 9d and shows that a steady state has indeed been reached.

The results presented in this paper demonstrate the ability of the entropy viscosity method with the new definitions of the viscosity coefficients to correctly simulate several types of flows (from very low Mach subsonic to transonic flows) without tuning parameters.

705 7. Conclusions

A new version of the entropy viscosity method that is valid for a wide range 706 of Mach numbers has been derived and presented for the inviscid Euler equations. The definition of the viscosity coefficients is now consistent with the 708 low-Mach asymptotic limit, does not require an analytical expression for the entropy function, and is therefore applicable to a larger variety of flow regimes, 710 from very low-Mach flows to supersonic flows. The method has also been ex-711 tended to Euler equation with variable area to solve nozzle flow problems. In 712 1-D, convergence of the numerical solution to the exact solution was demon-713 strated by computing the convergence rates of the L1 and L2 norms for flows 714 in a converging-diverging nozzle and in straight pipes. For smooth solutions, 715 second-order convergence was verified; solutions with shocks converged with the 716 expected theoretical rates of 1 (L_1 -norm) and 0.5 (L_2 -norm). 717

The effectiveness of the method was also demonstrated in 2-D using a series of benchmark problems for both subsonic and supersonic flows in various geometries, with Mach numbers ranging from 10^{-7} to 2.5. For very low-Mach flows, we numerically verified that the spatial steady-state pressure variations were proportional to the square of the Mach number, as expected in the incompressible limit.

In the future, we plan to further extend the entropy viscosity method to the seven-equation two-phase flow fluid model [20]. This two-phase flow system of equations is a good candidate for two reasons: it is unconditionally hyperbolic

and degenerates to the standard Euler equations when one phase disappears.

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A. Derivation of the entropy residual as a function of density, pressure and speed of sound

The entropy residual is defined as follows:

$$R_{\rm ent}(\vec{r},t) = \partial_t s(\vec{r},t) + \vec{u} \cdot \vec{\nabla} s(\vec{r},t),$$

where all variables were defined previously. This form of the entropy residual is not suitable for the low-Mach limit as explained in Section 2.1. In this appendix, we recast the entropy residual $R_{\rm ent}(\vec{r},t)$ as a function of the primitive variables (pressure, velocity and density) and the speed of sound. The first step of this derivation is to use the chain rule, recalling that the entropy is a function of the internal energy e and the density ρ , yielding

$$R_{\rm ent}(\vec{r},t) = s_e \frac{\mathrm{D}e}{\mathrm{D}t} + s_\rho \frac{\mathrm{D}\rho}{\mathrm{D}t}$$
,.

where s_e denotes the partial derivative of s with respect to the variable e. We recall that $\frac{D}{Dt}$ denotes the material derivative. Since the internal energy e is a function of pressure P and density ρ (through the equation of state), we use again the chain rule to re-express the previous equation as a function of the material derivatives in P and ρ :

$$R_{\text{ent}}(\vec{r},t) = s_e e_P \frac{DP}{Dt} + (s_e e_\rho + s_\rho) \frac{D\rho}{Dt}$$

$$= s_e e_P \left(\frac{DP}{Dt} + \frac{1}{s_e e_P} (s_e e_\rho + s_\rho) \frac{D\rho}{Dt} \right)$$

$$= s_e e_P \left(\frac{DP}{Dt} + (\frac{e_\rho}{e_P} + \frac{s_\rho}{s_e e_P}) \frac{D\rho}{Dt} \right).$$

To prove that the term multiplying the material derivative of the density is indeed equal to the square of the speed of sound, we recall that the speed of sound is defined as the partial derivative of pressure with respect to density at constant entropy, which can be recast as a function of the entropy as follows (see Appendix A.2 of [15]):

$$c^2 := \left. \frac{\partial P}{\partial \rho} \right|_{s=cst} = P_\rho - \frac{s_\rho}{s_e} P_e \,.$$

Using the following relations (see Appendix A.1 of [15])

$$P_e = \frac{1}{e_P}$$
 and $P_\rho = -\frac{e_\rho}{e_P}$.

Substitution of these expressions into the entropy residual equation above gives Eq. (9), which is recalled below for completeness:

$$R_{\rm ent}(\vec{r},t) := \partial_t s + \vec{u} \cdot \vec{\nabla} s = \frac{\mathrm{D}s}{\mathrm{D}t} = \frac{s_e}{P_e} \left(\underbrace{\frac{\mathrm{D}P}{\mathrm{D}t} - c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t}}_{\widetilde{R}_{\rm ent}(\vec{r},t)} \right).$$

B. Derivation of the dissipative terms for the Euler equations with variable area using the entropy minimum principle

The Euler equations (without viscous regularization) with variable area are recalled here

$$\partial_t \left(\rho A \right) + \vec{\nabla} \cdot \left(\rho \vec{u} A \right) = 0 \tag{43a}$$

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$$\partial_t \left(\rho \vec{u} A \right) + \vec{\nabla} \cdot \left[A \left(\rho \vec{u} \otimes \vec{u} + P \mathbb{I} \right) \right] = P \vec{\nabla} A \tag{43b}$$

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$$\partial_t \left(\rho E A \right) + \vec{\nabla} \cdot \left[\vec{u} A \left(\rho E + P \right) \right] = 0. \tag{43c}$$

The specific entropy is a function of the density ρ and the internal energy e, i.e., $s(e, \rho)$. The above system of equations satisfies the minimum entropy principle [30],

$$A\rho\left(\partial_t s + \vec{u} \cdot \vec{\nabla} \cdot s\right) \ge 0. \tag{44}$$

The entropy function s satisfies the second law of thermodynamics, $Tds = de - \frac{P}{\rho^2}d\rho$, which implies $s_e := T^{-1}$ and $s_\rho := -PT^{-1}\rho^{-2}$. One can show that [15]

$$s_e = T^{-1} \ge 0 \text{ and } Ps_e + \rho^2 s_\rho = 0.$$
 (45)

In order to apply the entropy viscosity method to the variable-area Euler equations, dissipative terms need to be added to each equation in Eq. (43). The functional forms of these terms need to be such that the entropy residual derived with these terms present also satisfies the minimum entropy principle. To prove the minimum entropy principle, the extra terms appearing in the entropy residual are either recast as conservative terms or shown to be positive. The rest of this appendix presents this demonstration. Following [15], we first write the variable-area equations with dissipative terms:

$$\partial_t (\rho A) + \vec{\nabla} \cdot (\rho \vec{u} A) = \vec{\nabla} \cdot f \tag{46a}$$

$$\partial_t \left(\rho \vec{u} A \right) + \vec{\nabla} \cdot \left[A \left(\rho \vec{u} \otimes \vec{u} + P \mathbb{I} \right) \right] = P \vec{\nabla} A + \vec{\nabla} \cdot g \tag{46b}$$

$$\partial_t \left(\rho E A \right) + \vec{\nabla} \cdot \left[\vec{u} A \left(\rho E + P \right) \right] = \vec{\nabla} \cdot \left(h + \vec{u} \cdot g \right). \tag{46c}$$

where f, g and h are dissipative fluxes to be determined. Starting from the modified system of equations given in Eq. (46), the entropy residual is derived again. The derivation requires the following steps: express the governing laws in terms of primitive variables (ρ, \vec{u}, e) , multiply the continuity equation by ρs_{ρ} and the internal energy equation by s_{e} , and invoke multivariate chain rule, e.g., $\partial s/\partial x = s_{e}\partial e/\partial x + s_{\rho}\partial \rho/\partial x$. These steps are similar to those used for the standard Euler equations [15]. Some of the lengthy algebra is omitted here. The above steps yield:

$$A\rho\left(\partial_t s + \vec{u} \cdot \vec{\nabla} s\right) = s_e \left[\vec{\nabla} \cdot h + g : \vec{\nabla} u + \left(\frac{u^2}{2} - e\right) \vec{\nabla} \cdot f\right] + \rho s_\rho \vec{\nabla} \cdot f. \tag{47}$$

The next step consists of choosing a definition for each of the dissipative terms so that the left hand-side is positive. The right hand-side of Eq. (47) can be simplified using the relations $g = A\mu \vec{\nabla}^s \vec{u} + f \otimes \vec{u}$ and $h = \tilde{h} - 0.5||\vec{u}||^2 f$ to give

$$A\rho\left(\partial_{t}s + \vec{u}\cdot\vec{\nabla}\cdot s\right) = s_{e}\left[\vec{\nabla}\cdot\tilde{h} - e\vec{\nabla}\cdot f\right] + \rho s_{\rho}\vec{\nabla}\cdot f + As_{e}\mu\vec{\nabla}\vec{u}^{s}:\vec{\nabla}\vec{u}. \tag{48}$$

The right hand-side is now integrated by parts:

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$$A\rho \left(\partial_t s + \vec{u} \cdot \vec{\nabla} \cdot s\right) = \vec{\nabla} \cdot \left[s_e \tilde{h} - s_e e f + \rho s_\rho f\right]$$
$$-\vec{\nabla} \cdot \tilde{h} \vec{\nabla} s_e + f \cdot \vec{\nabla} (e s_e) - f \cdot \vec{\nabla} (\rho s_\rho) + A s_e \mu \vec{\nabla}^s \vec{u} : \vec{\nabla} \vec{u} \quad (49)$$

where $\vec{\nabla}^s$ is the symmetric gradient. The term $As_e\mu\vec{\nabla}^s\vec{u}:\vec{\nabla}\vec{u}$ is positive and thus, does not need any further modification. It remains to treat the other

terms of the right hand-side that we now call rhs:

$$rhs = \vec{\nabla} \cdot \left[s_e \tilde{h} - s_e e f + \rho s_\rho f \right] - \tilde{h} \cdot \vec{\nabla} s_e + f \cdot \vec{\nabla} (e s_e) - f \cdot \vec{\nabla} (\rho s_\rho) \,.$$

The first term in rhs is a conservative term. By carefully choosing a definition for \tilde{h} and f, the conservative term can be expressed as a function of the entropy s. The inclusion of the variable area in the choice of the dissipative terms is also required so that, when assuming constant area, the standard Euler equations are recovered. The following definitions for \tilde{h} and f are chosen:

$$\tilde{h} = A\kappa \vec{\nabla}(\rho e)$$
 and $f = A\kappa \vec{\nabla}\rho$,

which yields, using the chain rule,

$$rhs = \vec{\nabla} \cdot (\rho A \kappa \vec{\nabla} s) - A \kappa \underbrace{\left[\vec{\nabla} (\rho e) \vec{\nabla} s_e - \vec{\nabla} \rho \vec{\nabla} (e s_e) + \vec{\nabla} \rho \vec{\nabla} (\rho s_\rho)\right]}_{\mathbf{Q}}$$

 854 It remains to treat the term \mathbf{Q} that can be recast under a quadratic form.

Following [15], one obtain:

$$\mathbf{Q} = \rho X^t \Sigma X$$
with $X = \begin{bmatrix} \vec{\nabla} \rho \\ \vec{\nabla} e \end{bmatrix}$ and $\Sigma = \begin{bmatrix} \rho^{-2} \partial_{\rho} (\rho^2 \partial_{\rho} s) & \partial_{\rho, e} s \\ \partial_{\rho, e} s & \partial_{e, e} s \end{bmatrix}$

The matrix Σ is symmetric and identical to the matrix obtained in [15]. The sign of the quadratic form can be simply determined by studying the positiveness of the matrix Σ . In this particular case, it is required to prove that the matrix is negative definite: the quadratic form is on the right hand-side and is preceded by a negative sign. According to [15], the convexity of the opposite of the entropy

- function, i...e, -s, with respect to the internal energy e and the specific volume
- $_{\mbox{\scriptsize 862}}$ $1/\rho$ is sufficient to ensure that the matrix Σ is negative definite.
- $_{863}$ Thus, the right hand-side of the entropy residual Eq. (47) is now either recast
- 864 as conservative terms, or known to be positive. Thus, the entropy minimum
- 865 principle holds.

666 C. Entropy residual for isentropic flows

- This appendix shows that the entropy residual is zero for isentropic flows.
- 868 For convenience, we recall here the entropy residual as a function of the pressure,
- density, velocity, and speed of sound:

$$\widetilde{R}_{\text{ent}} = \frac{\mathrm{D}P}{\mathrm{D}t} - c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t} \,. \tag{50}$$

Assuming an isentropic flow, pressure is only a function of density, i.e., P= $f(\rho)$ or equivalently $\rho=f^{-1}(P)$. Using the definition of the speed of sound $c^2=\frac{\partial P}{\partial \rho}$ and the above form of the equation of state, we have

$$c^{2} = \frac{\partial P}{\partial \rho} \bigg)_{s} = \frac{dP}{d\rho} = \frac{df(\rho)}{d\rho} \,. \tag{51}$$

Using the chain rule, the entropy residual in Eq. (50) can be recast as follows and proven equal to zero:

$$\widetilde{R}_{\text{ent}} = \frac{df(\rho)}{d\rho} \frac{\mathrm{D}\rho}{\mathrm{D}t} - c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t} = c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t} - c^2 \frac{\mathrm{D}\rho}{\mathrm{D}t} = 0.$$
 (52)