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A self-contained composition space solution method for strained and curved premixed flamelets

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Abstract

Solving flames in scalar or composition space usually requires the modeling of scalar dissipation rates, or scalar gradients, which appear in the Jacobian of the transformation from the physical-space coordinates to the scalar space coordinates attached to the flame surface. Recently, Scholtissek *et al.* (2019) have discussed a self-contained solution for freely propagating premixed flames, in which the gradient distribution of the coordinate is also solved in scalar space. This approach is here extended to include curvature, strain and effects of unsteadiness. The resulting method is applied to a constant pressure and homogeneous ignition case, a stagnation flame, a tubular flame and a spherical expanding flame, capturing accurately effects of unsteadiness, strain and curvature. The results are systematically validated against physical-space solutions and experiments. In addition to the possibility of exploring these various canonical problems with a single set of equations, it is shown that another major outcome of this new formulation lies in the possibility of studying steady flames subjected to negative strain.

Keywords: premixed flamelet, progress variable space, strain, curvature, gradient equation

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1. Introduction

In most combustion systems the fraction of the flow occupied by the reaction zones stays very small. The characteristic thicknesses over which intermediate radical chemical species evolve are orders of magnitude smaller than the characteristic length of the large-scale flow motion, which carries the turbulence kinetic energy driving the mixing of the reactants. To deal with this strong multi-scale character of combustion physics [1], scalar or composition space coordinates attached to the flame surface are usually introduced, to analyse results of both experiments and simulations [2].

Such scalar space coordinates were first designed for conveniently focussing theoretical analyses on the molecular diffusion layers feeding the thin reaction zones in laminar diffusion flames [3]. Then, scalar coordinate systems have been used systematically in flame theory, up to fuel-spray flames [4]. Composition or scalar space coordinates were also introduced for turbulent combustion modeling, in diffusion flamelets [5, 6], in premixed flamelets [7–9], in conditional moment closure (CMC) [10] or again in the transport of probability density functions (PDF) [11–13]. Even though the underlying physical hypotheses required to complete the modeling can differ between these approaches, they all rely on the definition of a scalar space coordinate system centered on the thin diffusive-reactive layers.

Along these lines, strategies were developed to solve laminar and canonical combustion problems directly in scalar space. The most commonly used is the mixture fraction ($Z = 0$ in the oxidizer stream, $Z = 1$ in the fuel feeding stream), which is a passive scalar that can be related to the local equivalence ratio of the mixture in non-premixed systems and that can be defined in multicomponent flows from a set of representative species mass fractions [14, 15]. The solving of one-dimensional counter-flowing diffusion flames was for instance intensively studied in mixture fraction space, including various levels of complexity [16–19]. Sometimes scalars are combined with other coordinates. In the context of reacting mixing layers in stagnation-type flows subject to both time-varying strain and pressure, a thermal-conductivity-weighted transverse coordinate was found to be very effective [20].

The normalized progress variable ($c = 0$ in the fresh gases, $c = 1$ in the fully burnt products) is widely used for premixed flames and the solving of freely propagating laminar premixed flames in progress variable space was reported [21, 22]. Finally, the combination of both mixture fraction and progress variable allows for addressing partially premixed and non-premixed flames in a

two-dimensional scalar space [23, 24].

Up to very recently, the scalar space formalism was not self-consistent. Indeed, the jacobian for transformation between the physical and scalar coordinates contains the physical-space gradient of the scalar coordinates, which was left as a free parameter in the balance equations. To overcome this strong limitation, in [22] the balance equation for the gradient is also projected in scalar space, to come up with a fully self-consistent scalar space solution for freely propagating unstrained one-dimensional premixed flames. It is proposed in this work to extend and generalise this self-consistent solution procedure.

First, premixed flamelet equations for temperature, species mass fractions and progress variable gradient are derived incorporating effects of strain and curvature, with only flow quantities as flamelet parameters. The equation set is expressed with respect to the progress variable, which spans the scalar space. Subsequently, numerical methods for solving the flamelet equations are briefly revisited. It is shown that the asymptotic limit of the zero-dimensional and constant pressure ignition problem is well recovered by the proposed formulation. Then, three canonical configurations, namely the stagnation flame, the tubular flame and the spherical expanding flame, are simulated in progress variable space with detailed chemistry and complex transport properties. Three fuels are studied: hydrogen, methane and ethanol. By comparison to physical space reference solutions and experimental measurements of the literature, it is demonstrated that effects of unsteadiness, strain and curvature are well captured by the novel set of premixed flamelet equations, including negative stretch, opening new perspectives for turbulent combustion modeling.

2. Flame and flow parameters

2.1. Budget equations

The instationary balance equations for temperature T and the species mass fractions Y_i serve as a starting point for the flamelet model derivation. They can be expressed with respect to the 3D physical space as

$$\rho c_p \frac{\partial T}{\partial t} + \rho c_p \mathbf{u} \cdot \nabla T = \nabla \cdot (\lambda \nabla T) - \rho \sum_k^{n_s} c_{p,k} Y_k \mathbf{V}_k \cdot \nabla T + \dot{\omega}_T, \quad (1)$$

$$\rho \frac{\partial Y_i}{\partial t} + \rho \mathbf{u} \cdot \nabla Y_i = -\nabla \cdot (\rho Y_i \mathbf{V}_i) + \dot{\omega}_i, \quad (2)$$

where ρ is the density, c_p the specific heat capacity, λ the thermal conductivity, \mathbf{u} the velocity vector and \mathbf{V}_k the diffusion velocity vector of species k . The quantities $\dot{\omega}_T$ and $\dot{\omega}_i$ denote the source terms due to gas phase chemistry for temperature and species, respectively.

The progress variable is defined as a combination of species according to $Y_c = \sum_i^{n_s} \alpha_i Y_i$, where α_i is a species specific weighting factor and n_s is the total number of species. The analogous summation over Eq. (2) yields a balance equation for the progress variable:

$$\rho \frac{\partial Y_c}{\partial t} + \rho \mathbf{u} \cdot \nabla Y_c = -\nabla \cdot (\rho Y_c \mathbf{V}_c) + \dot{\omega}_c, \quad (3)$$

where the diffusive flux of the progress variable is defined as $Y_c \mathbf{V}_c = \sum_i^{n_s} \alpha_i Y_i \mathbf{V}_i$ and the source term as $\dot{\omega}_c = \sum_i^{n_s} \alpha_i \dot{\omega}_i$. A fundamental requirement for the choice of the progress variable is that it varies strictly monotonically throughout the flame front. Note that its definition depends on the fuel-oxidizer system and is not always trivial to identify [25–27].

2.2. Flame speed and displacement speed

The flame is assumed to be a thin sheet of finite thickness, which is not necessarily flat, and which can be represented by a collection of Y_c -isosurfaces (analogy to FGM [28]). The motion of each surface is described by the kinematic condition [28, 29]

$$\frac{dY_c}{dt} = \frac{\partial Y_c}{\partial t} + \mathbf{s}_f \cdot \nabla Y_c = 0, \quad (4)$$

which states that a point on the isosurface of Y_c remains on this surface for all t . For a fixed observer the flame moves with the apparent flame speed \mathbf{s}_f in the physical space, which is composed of the flow velocity vector \mathbf{u} and the flame displacement speed s_d directed opposite to the normalized gradient $\mathbf{n}_c = \nabla Y_c / |\nabla Y_c|$

$$\mathbf{s}_f = \mathbf{u} - s_d \mathbf{n}_c. \quad (5)$$

Introducing Eq. (5) in Eq. (4) one obtains

$$\rho \frac{\partial Y_c}{\partial t} + \rho \mathbf{u} \cdot \nabla Y_c = \rho s_d |\nabla Y_c|, \quad (6)$$

which further yields an expression for the local mass burning rate ρs_d of a premixed flame sheet, when further employing Eq. (3) to replace the left hand side in the above equation

$$\rho s_d = \frac{1}{|\nabla Y_c|} (-\nabla \cdot (\rho Y_c \mathbf{V}_c) + \dot{\omega}_c) . \quad (7)$$

2.3. Flame stretch, strain and curvature

Flame stretch is a key quantity influencing a flame's structure, extinction behavior, and its mass burning rate. Here, important characteristics of the flame stretch concept are briefly revisited before incorporating stretch effects into the flamelet model.

A first generic definition of flame stretch was given by Williams [30]

$$K_A = \frac{1}{A} \frac{dA}{dt} , \quad (8)$$

which describes the fractional area change of a small flame surface element moving with an infinitely thin flame sheet. In this context, extensive theory has been developed and many practical flame configurations have been studied and discussed [29, 31–33]. The formulation for flame stretch was later on extended by de Goeij and ten Thije Boonkkamp [34] considering the mass $M(t)$ in an infinitesimal volume $V(t)$ moving in the flame of finite thickness

$$K = \frac{1}{M} \frac{dM}{dt} \quad \text{with} \quad M(t) = \int_{V(t)} \rho dV . \quad (9)$$

While the formulation by Williams [30] is also referred to as weak stretch, the latter is referred to as strong stretch further accounting for the resolved diffusive layer in the vicinity of the flame [28]. The stretch rate K can be defined as [34]

$$\rho K = \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{s}_f) , \quad (10)$$

and can be rewritten using Eq. (5) according to [34]

$$\rho K = -\nabla \cdot (\rho s_d \mathbf{n}_c) , \quad (11)$$

which illustrates that the stretch rate is related to variations of the normal mass flux through

the flame surface [34]. Expanding Eq. (10), the single contributions to flame stretch can be examined [34]

$$K = \nabla \cdot \mathbf{s}_f + \frac{1}{\rho} \left(\frac{\partial \rho}{\partial t} + \mathbf{s}_f \cdot \nabla \rho \right) = \nabla \cdot \mathbf{u}_t - s_f \kappa_c + \mathbf{n}_c \cdot \nabla s_f + \frac{1}{\rho} \frac{\partial \rho}{\partial t}, \quad (12)$$

where the curvature of the Y_c -isosurface is defined as

$$\kappa_c = -\nabla \cdot \mathbf{n}_c, \quad (13)$$

the flame-tangential (flow) velocity vector is defined as $\mathbf{u}_t = \mathbf{u} - (\mathbf{n}_c \cdot \mathbf{u}) \mathbf{n}_c$ and $s_f = \mathbf{s}_f \cdot \mathbf{n}_c$ is the projected flame speed in flame-normal direction. The first and the second term on the right side of Eq. (12) can also be found from weak stretch theory [31–33] and can be interpreted as flame-tangential strain and flame stretch through instationary movement of a curved flame sheet. The third term in Eq. (12) arises from instationary flame-normal stretch, such that the flame-normal velocity s_f varies between different Y_c -isosurfaces. The fourth term in Eq. (12) originates from density variations along flame isosurfaces. Due to the fact that isosurfaces of ρ are assumed to align with Y_c -isosurfaces, the last two terms in Eq. (12) cancel, which is readily seen from the kinematic condition Eq. (4).

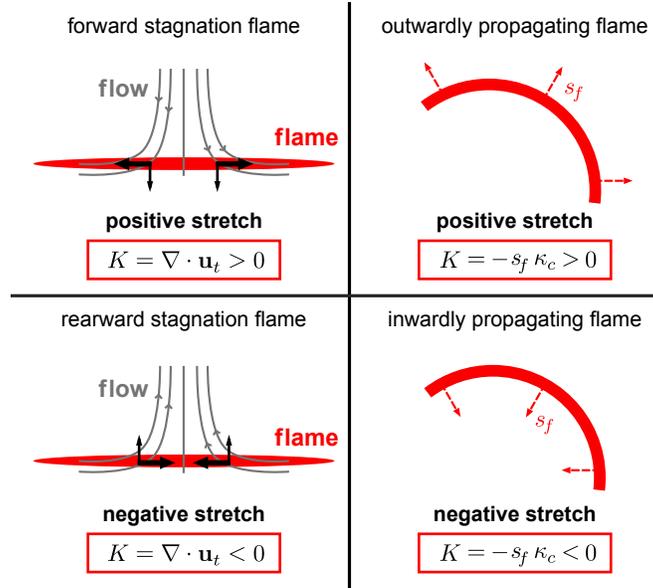


Figure 1: Schematic flame configurations which experience stretch due to strain (left column) or stretch due to curvature (right column) [35].

The relevant effects described by the first two terms in Eq. (12) are shown in Fig. 1 for flames which exhibit either stretch due to flame-tangential strain (left column) or stretch due to movement of a curved flame sheet (right column). For the latter, the flame moves in the flame-normal direction such that the flame surface increases (positive stretch, outward flame propagation) or decreases (negative stretch, inward flame propagation). Hence, stretch conditions can be readily associated with the surface-based stretch definition, see Eq. (8). For the flames in a forward or rearward stagnation flow, flame-tangential straining can be understood by considering the flow velocity components projected into the flame-tangential plane (thick arrows). Of the flames depicted in Fig. 1, only the positively stretched flames can be studied under controlled conditions in a laboratory. The rearward stagnation flame is inherently unstable and the inwardly propagating flame cannot reproducibly be ignited in perfect spherical (or cylindrical) shape before it starts to propagate. Overall, it is hardly possible to establish significant negative strain conditions with laboratory flames¹, however, large negative strain rates can be identified in turbulent flames [37, 38]. *It will be shown further below that the composition space flame model developed here allows to study also negative strain conditions, which is, with few exceptions [39, 40], not possible with conventional physical space flame solvers.*

3. Progress variable space formulation

3.1. Transformation rules

The governing equations are transformed from physical space (t, \mathbf{x}) to the progress variable space $(\tau, Y_c(t, \mathbf{x}))$, where τ represents a flamelet time defined by $\tau = t$. The corresponding temporal transformation rule reads for the exemplary scalar quantity ϕ

$$\frac{\partial \phi}{\partial t} = \frac{\partial \phi}{\partial \tau} + \frac{\partial Y_c}{\partial t} \frac{\partial \phi}{\partial Y_c}. \quad (14)$$

¹The Bunsen flame, which exhibits negative stretch, is one of the few exceptions. However, it has been realized recently that the structure of a Bunsen flame is controlled by rather complicated physics [36]. Its stretch conditions are non-uniform and have to be considered at least two-dimensional.

For any quantity $\phi = \phi(Y_c)$, the following spatial transformation rules are verified:

$$\nabla\phi = \nabla Y_c \frac{\partial\phi}{\partial Y_c}, \quad (15)$$

$$\nabla^2\phi = |\nabla Y_c|^2 \frac{\partial^2\phi}{\partial Y_c^2} + \nabla^2 Y_c \frac{\partial\phi}{\partial Y_c}. \quad (16)$$

The second order derivative $\nabla^2 Y_c$ is decomposed into its normal and curvature contribution [41]

$$\nabla^2 Y_c = \mathbf{nn} : \nabla(\nabla Y_c) + |\nabla Y_c| \nabla \cdot \mathbf{n}. \quad (17)$$

One can write $\nabla(|\nabla Y_c|) = \mathbf{n} \cdot \nabla(\nabla Y_c)$, then $\mathbf{nn} : \nabla(\nabla Y_c) = \mathbf{n} \cdot \nabla(|\nabla Y_c|) = |\nabla Y_c| (\partial|\nabla Y_c|/\partial Y_c)$, where Eq. (15) is also applied to $\nabla(|\nabla Y_c|)$. Combined with Eqs. (16) and (17), this leads to

$$\nabla^2\phi = |\nabla Y_c|^2 \frac{\partial^2\phi}{\partial Y_c^2} + |\nabla Y_c| \left(\frac{\partial|\nabla Y_c|}{\partial Y_c} + \nabla \cdot \mathbf{n} \right) \frac{\partial\phi}{\partial Y_c}. \quad (18)$$

In Eq. 18, the progress variable gradient $g_c = |\nabla Y_c|$ is introduced. This quantity plays a central role in the flamelet transformation since it relates Y_c -space to the physical space. Traditionally, g_c is used to define the scalar dissipation rate

$$\chi_c = 2D_c \nabla Y_c \cdot \nabla Y_c = 2D_c g_c^2. \quad (19)$$

While the gradient (or scalar dissipation rate) can be modeled for non-premixed flamelets [42], it has to be treated as a dependent quantity for premixed flamelets [22].

3.2. Diffusion modeling in progress variable space

Keeping in mind that the species diffusion velocity \mathbf{V}_i contains gradient information, it is transformed using the spatial transformation rule Eq. (15) [22, 43]

$$\rho Y_i \mathbf{V}_i = \rho Y_i \tilde{V}_i \nabla Y_c, \quad (20)$$

where \tilde{V}_i represents a diffusion velocity in progress variable space. Instead of gradients with respect to physical space coordinates, \tilde{V}_i analogously contains gradients with respect to the conditioning variable Y_c and can therefore be computed from conventional transport libraries on a Y_c -grid

(e.g. cantera [44] or EGLib [45, 46]). Consequently, the approach allows arbitrary diffusion models to be incorporated in progress variable space without notational overhead. In this work, a mixture averaged diffusion model [47] with a correction velocity for mass conservation [48] is applied. Additionally, thermal diffusion [45, 46] is considered for H₂-air flames.

3.3. Composition space equations

Composition space (or flamelet) equations are obtained by applying the transformation rules in Sec. 3.1 to Eqs. (1) and (2). Furthermore, noticing that $g_c = |\nabla Y_c| = \mathbf{n} \cdot \nabla Y_c$, an equation for the progress variable gradient g_c is obtained by differentiating (and transforming) the balance equation for Y_c according to

$$\mathbf{n} \cdot \nabla \left[\rho \frac{\partial Y_c}{\partial t} + \rho \mathbf{u} \cdot \nabla Y_c = -\nabla \cdot (\rho Y_c \mathbf{V}_c) + \dot{\omega}_c \right]. \quad (21)$$

Details of the derivation are available in the supplementary material. The premixed flamelet equations for species, temperature and progress variable gradient are obtained as

$$\begin{aligned} \rho \frac{\partial Y_i}{\partial \tau} &= -g_c \frac{\partial}{\partial Y_c} \left(g_c \rho Y_i \tilde{V}_i \right) + g_c \frac{\partial}{\partial Y_c} \left(g_c \rho Y_c \tilde{V}_c \right) \frac{\partial Y_i}{\partial Y_c} \\ &\quad + \rho g_c \kappa_c \left(Y_i \tilde{V}_i - Y_c \tilde{V}_c \frac{\partial Y_i}{\partial Y_c} \right) - \dot{\omega}_c \frac{\partial Y_i}{\partial Y_c} + \dot{\omega}_i, \end{aligned} \quad (22)$$

$$\begin{aligned} \rho \frac{\partial T}{\partial \tau} &= \frac{g_c}{c_p} \frac{\partial}{\partial Y_c} \left(g_c \lambda \frac{\partial T}{\partial Y_c} \right) + g_c \frac{\partial}{\partial Y_c} \left(g_c \rho Y_c \tilde{V}_c \right) \frac{\partial T}{\partial Y_c} - \rho g_c^2 \sum_k^{n_s} \frac{c_{p,k}}{c_p} Y_k \tilde{V}_k \frac{\partial T}{\partial Y_c} \\ &\quad - \rho g_c \kappa_c \left(\frac{\lambda}{\rho c_p} + Y_c \tilde{V}_c \right) \frac{\partial T}{\partial Y_c} - \dot{\omega}_c \frac{\partial T}{\partial Y_c} + \frac{\dot{\omega}_T}{c_p}, \end{aligned} \quad (23)$$

$$0 = -g_c^2 \frac{\partial^2}{\partial Y_c^2} \left(g_c \rho Y_c \tilde{V}_c \right) + g_c^2 \frac{\partial}{\partial Y_c} \left(\kappa_c \rho Y_c \tilde{V}_c \right) - \dot{\omega}_c \frac{\partial g_c}{\partial Y_c} + g_c \frac{\partial \dot{\omega}_c}{\partial Y_c} + \rho K_s g_c. \quad (24)$$

These premixed flame equations are similar to the ones presented in our earlier work [22], with additional terms to incorporate the flamelet parameters strain K_s and curvature κ_c . The quantity K_s differs from the overall stretch rate K and represents straining imposed on the flame by the flow field

$$K_s = -\frac{1}{\rho} \frac{\partial(\rho s_d)}{\partial n} = K - s_d \kappa_c \stackrel{(12)}{=} \nabla \cdot \mathbf{u}_t - (s_f + s_d) \kappa_c = \nabla \cdot \mathbf{u}_t - u_g \kappa_c, \quad (25)$$

where $u_g = \mathbf{n}_c \cdot \mathbf{u}$ is the flow velocity along the flame-normal direction. Formally, the flame stretch K can be split into flow straining K_s , and flame stretch due to the self-propagation of the (curved)

flame sheet described by $K_c = -s_d \kappa_c$. This splitting procedure has been carried out by other authors before [39, 49]. The relation (25) indicates that K_s solely depends on flow quantities. This is particularly important, since the flame displacement speed s_d is a flame response obtained by solving the premixed flamelet equations and shall therefore not appear in the definition of the strain rate, which is one of the flame control parameters.

Note that species and temperature equation are instationary in composition space, while the gradient equation is not. When prescribing a certain set of flamelet parameters the flame topology becomes implicitly fixed. Due to the fact that the gradient represents the spatial mapping from composition space to physical space it also has to be fixed for a given combination of K_s and κ_c . A similar finding was reported for FGM by van Oijen et al. [28] who showed that the equation for the progress variable in the flame-adapted coordinate system is stationary.

3.4. Boundary conditions

The conditioning variable Y_c varies on the interval $[Y_{c,\min} \ Y_{c,\max}]$ and the exact values of minimum and maximum depend on the progress variable definition. At the left boundary ($Y_{c,\min}$) Dirichlet boundary conditions are imposed

$$Y_c = Y_{c,\min} : \quad Y_i = Y_{i,0}, \quad T = T_0, \quad g_c = 0, \quad (26)$$

which are determined a-priori from the unreacted mixture of fresh gases.

For zero strain and curvature, an unstrained, planar freely-propagating flame and the constant pressure homogeneous ignition can be described by the flamelet equations (22), (23) and (24) [22]. In this case, the conditions at the right boundary ($Y_{c,\max}$) correspond to the chemical equilibrium which can be determined from the fresh gas mixture. Hence, it can also be described by Dirichlet boundary conditions

$$Y_c = Y_{c,\max} = Y_{c,\text{eq}} : \quad Y_i = Y_{i,\text{eq}}, \quad T = T_{\text{eq}}, \quad g_c = 0, \quad (27)$$

where the subscript “eq” refers to chemical equilibrium conditions.

However, for non-zero strain or curvature, the right boundary location $Y_{c,\max}$ becomes an eigenvalue of the problem since premixed flames can exhibit sub- or super-equilibrium conditions

(also called hot spots [50–52]). This implies that the Y_c -space can contract or expand and that the right boundary has to be implemented as a moving boundary. For the numerical solution, the flamelet equations, Eqs. (22)-(24), are therefore transformed to a normalized progress variable space (c -space). The procedure is described in the supplementary material.

3.5. Closure of the flamelet equations

With the flamelet gradient equation, Eq. 24, the linkage of the flamelet model to the flow is no longer inherently captured by prescribing the gradient or scalar dissipation rate as for previous premixed flamelet formulations in progress variable space [21, 23, 53]. Instead, the effects of the flow on the flamelet structure are represented directly by strain and curvature, which appear explicitly in the equations and which have been proven to be fundamental flame parameters [54]. A closure for these flame parameters is obtained by approximating K_s and κ_c by representative constant values. The approach yields an accurate and self-contained premixed flamelet model, including flamelets subject to strain and curvature. This is demonstrated in Sec. 5 by comparing flamelet solutions to well-established numerical models (formulated in the physical space) and experiments for stagnation flames, tubular flames and spherical expanding flames.

4. Numerical methods

The stationary solution of the flamelet equations is obtained with a hybrid damped Newton method [55] with an in-house C++ flame solver [56]. A straight line is prescribed as the initial solution which connects the fresh gas mixture (left boundary condition) with the corresponding equilibrium conditions (right boundary initialization). In a pseudo-transient time-stepping procedure, the initial solution is first relaxed for few time steps before a stationary solution by Newton iteration is attempted. If the stationary solver fails, the solution is advanced for a fixed number of additional time steps and the stationary solver is rerun. This procedure is repeated until convergence of the stationary solver. Starting from a coarse grid with 8-30 points, the grid is adaptively refined after each convergent solution until a user-specified target. For the evaluation of diffusion coefficients and chemical source terms, the EGlLib [45, 46] and an in-house chemistry library are used, respectively. The chemical reaction mechanism by Varga et al. [57] (12 species, 29 reactions) is applied for H_2 -air flames, the GRI 3.0 mechanism by Smith et al. [58] (53 species, 325 reactions) for CH_4 -air flames, and the mechanism by Olm et al. [59] (47 species, 250 reactions) for C_2H_5OH -air flames.

5. Self-contained scalar space flame simulations

A homogeneous reactor (ignition problem) and 1D canonical laminar flames are studied applying the previously introduced formalism. By comparison to corresponding physical space solutions the progress variable space model is validated and the numerical solutions are complemented by suitable experimental results. Table 1 contains an overview of the configurations studied. For the flame configurations, H₂-air and CH₄-air flames are considered due to the availability of extensive reference data from the literature. Fuel-lean flames are investigated in which the fuel Lewis number ($Le_{H_2} \approx 0.3$, $Le_{CH_4} \approx 1$) determines how strain and curvature affect flame structures, it is therefore denoted as the *effective Lewis number* Le [54]. In order to study flame responses for $Le > 1$, additionally C₂H₅OH-air flames are considered ($Le_{C_2H_5OH} \approx 1.6$).

The progress variable definitions for the flamelet calculations are chosen as $Y_c = Y_{H_2O} - Y_{H_2} - Y_{O_2}$ (H₂-air flames) and $Y_c = Y_{CO_2} + Y_{H_2O} + 10 Y_{H_2} - Y_{O_2}$ (CH₄- and C₂H₅OH-flames). Also other progress variable definitions would yield accurate flamelet results, often a combination of product (positive weight in definition) and educt species (negative weight) is suitable. For the hydrocarbon flames the weight for the mass fraction of H₂ was increased due to its low molecular mass. Note that it is important to include H₂ in the definition of the progress variable to correctly capture differential diffusion effects in the preheat zone of a premixed flame with a premixed flamelet model [22].

Table 1: Characteristics of several canonical combustion configurations (0D and 1D) with respect to unsteady effects and the flamelet parameters strain K_s and curvature κ_c .

case	acronym	strain	curvature	unsteadiness	reference
homogeneous reactor	HR	–	–	+	this work
freely-propagating flame	FP	–	–	–	[22]
stagnation flame	STAG	+	–	–	this work
tubular stagnation flame	TUB	+	+	–	this work
spherical expanding flame	SEF	+	+	+	this work

5.1. Constant pressure homogeneous ignition

The transient evolution of temperature and species mass fractions in the constant pressure homogeneous reactor (HR) are described neglecting convective and diffusive transport terms in

Eqs. (1) and (2), which yields the following equation set:

$$\rho \frac{\partial T}{\partial t} = \frac{\dot{\omega}_T}{c_p} \quad \text{and} \quad \rho \frac{\partial Y_i}{\partial t} = \dot{\omega}_i. \quad (28)$$

With the transient evolution of Y_i also the reaction progress variable changes. The summation over the species equation according to the definition of the progress variable yields

$$\rho \frac{\partial Y_c}{\partial t} = \dot{\omega}_c. \quad (29)$$

Defining the progress variable such that it includes species which vary during all phases of the ignition process (thus, preferably major species) allows the mapping of the thermochemical variables (here T) to the reaction progress variable in time

$$\frac{\partial T}{\partial t} = \frac{\partial Y_c}{\partial t} \frac{\partial T}{\partial Y_c}, \quad (30)$$

which defines a temporal transformation rule different from Eq. (14), since it does not consider the flamelet time τ . With this, the constant pressure homogeneous ignition is described in progress variable space by combining Eqs. (28)-(30)

$$\frac{\partial T}{\partial Y_c} = \frac{1}{c_p} \frac{\dot{\omega}_T}{\dot{\omega}_c} \quad \text{and} \quad \frac{\partial Y_i}{\partial Y_c} = \frac{\dot{\omega}_i}{\dot{\omega}_c}. \quad (31)$$

This equation set has previously been used to describe autoignition [60] and contains formally no transient term. The Eqs. (31) are one-dimensional and imply that a trajectory in composition space can be directly constructed by integrating along Y_c (assuming a monotonically varying progress variable). This model shall be denoted as the progress variable space reactor (PVR).

Figure 2 displays the constant pressure ignition of a stoichiometric CH_4 -air mixture at 1500 K. The left graph of this figure shows the evolution of the temperature and the mass fractions for CO_2 and the radical OH versus time t (*i.e.* simulated solving Eq. (28)). The middle graph shows a projection of the temperature profile into progress variable space, where $Y_c = Y_{\text{CO}_2}$. It is observed that the PVR model (Eq. (31)) accurately reproduces this progress variable space trajectory. Furthermore, the time integration (in flamelet time τ) of the premixed flamelet model (PFLT) is

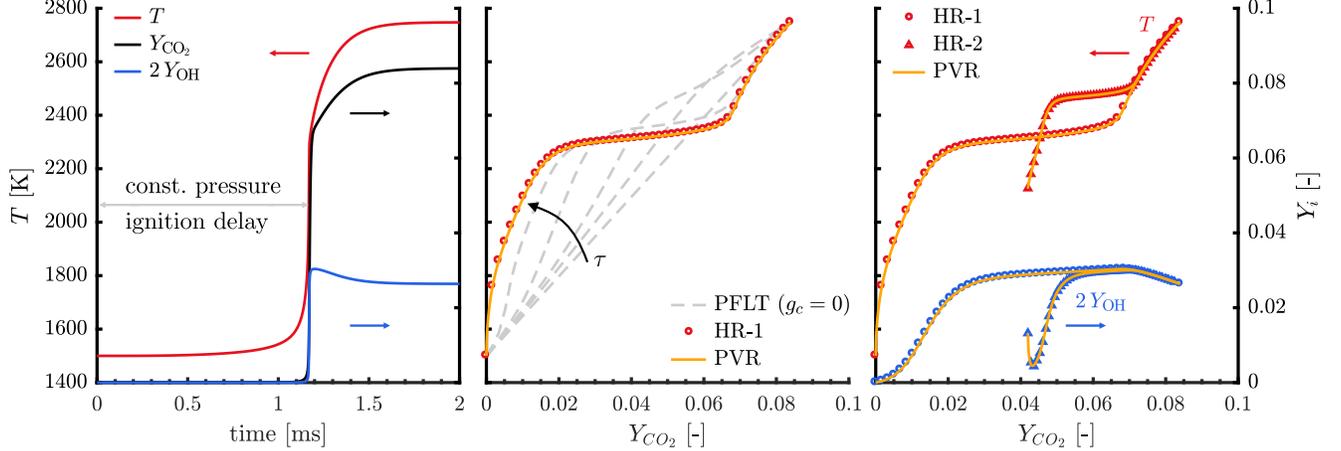


Figure 2: Left: Constant pressure ignition of a homogeneous stoichiometric CH_4 -air mixture at 1500 K versus time (case HR-1). Middle: The temperature profile for the case HR-1 mapped into progress variable space ($Y_c = Y_{\text{CO}_2}$). The temperature profile is computed with three different models: homogeneous reactor (HR), reactor in progress variable space (PVR) and the transient premixed flamelet model (PFLT, $g_c = 0$). Right: Progress variable space trajectories of two different constant pressure ignition cases recomputed with the models HR and PVR. The initial state of case HR-2 consists of a 50% fresh stoichiometric CH_4 -air mixture at 1500 K and 50% hot products with the same enthalpy level.

shown, which is first initialized with a linear profile between fresh gases and chemical equilibrium. Due to homogeneity of the mixture, the progress variable gradient $g_c = |\nabla Y_c|$ equals to zero and the premixed flamelet equations (23) and (22) reduce to

$$\rho \frac{\partial T}{\partial \tau} = -\dot{\omega}_c \frac{\partial T}{\partial Y_c} + \frac{\dot{\omega}_T}{c_p}, \quad (32)$$

$$\rho \frac{\partial Y_i}{\partial \tau} = -\dot{\omega}_c \frac{\partial Y_i}{\partial Y_c} + \dot{\omega}_i. \quad (33)$$

It is readily shown, that the stationary solution of this equation set is equal to Eqs. (31) and as seen from the mid graph in Fig. 2 the solution for the PFLT model relaxes to the solutions for the models HR and PVR for increasing τ .

These results illustrate that the behavior of the transient ignition are captured well by the stationary solution of the premixed flamelet model in progress variable space. While there are two terms in the temporal transformation rule, Eq. (14), which formally describe transient effects, $\partial T / \partial \tau$ and $\partial Y_c / \partial t$, the former can hardly be interpreted from a physical point of view. This is due to the fact that transient variations of thermodynamical quantities (i.e. species and temperature) with τ would require that the progress variable remains fixed. However, this is rarely the case for most thermochemical variables, since the conditioning variable Y_c is defined as a combination

of species and therefore dynamically linked to the full thermodynamical state through chemical reactions and thermodynamics. Also for fast reactions of radical species, the stationary solution of the premixed flamelet model is appropriate, as shown by the right graph in Fig. 2. Considering a highly reactive mixture (case HR-2), which is obtained by mixing 50% of fresh gases and 50% of hot products by mass, the PVR model (equivalently the stationary PFLT model) still accurately captures the evolution of temperature and the radical OH in progress variable space. This also applies to other radical species (not shown).

5.2. Canonical laminar premixed flames

For premixed combustion, the unstretched laminar burning velocity s_u^0 represents an important characteristic quantity which inherently contains fundamental information regarding the reactivity, diffusivity and exothermicity of a combustible mixture [61] (the superscript 0 marks the stretchless reference condition). While the reference quantity s_u^0 solely depends on the thermochemical state of the fresh gas mixture, the burning velocity s_u of a premixed flame can be influenced by multiple factors such as strain, curvature or heat-loss. It is defined at the edge of unburned gases and is thereby distinguished from the flame displacement speed s_d , which varies throughout the flame zone due to thermal expansion. The latter can be computed from a composition space flamelet solution using the transformed Eq. (7):

$$s_d = \frac{1}{\rho} \left[-\frac{\partial}{\partial Y_c} \left(g_c \rho Y_c \tilde{V}_c \right) + \frac{\dot{\omega}_c}{g_c} \right]. \quad (34)$$

The laminar burning velocity s_u is then determined with the above expression at the unburned edge of the flame, which is defined by an isotherm $T = T_u + 5\text{K}$ [49] (with T_u being the temperature of the unburnt gases). Furthermore, the absolute flame speed s_f can be defined, which marks the apparent speed of the flame for a fixed observer

$$s_f = u_g - s_u, \quad (35)$$

where u_g is the flame-normal velocity component of the flow at the unburned edge of the flame. Note that the negative sign for s_u originates from its direction opposite to the normalized gradient $\mathbf{n}_c = \nabla Y_c / |\nabla Y_c|$. For the ideal freely-propagating reference flame, s_f is constant throughout the flame

zone. However, as soon as the flame becomes stretched (due to strain and/or curvature), flame speeds are more difficult to evaluate, both experimentally and numerically [62].

Three stretched flame configurations with successively more complex conditions (see Tab. 1) are discussed in this section: the stagnation flame (STAG), the tubular stagnation flame (TUB) and the spherical expanding flame (SEF).

5.2.1. Planar stagnation flames

Planar stagnation flames have been used to study flame structures, propagation and extinction, both theoretically and experimentally, and a broad overview can be found in the review paper by Egolfopoulos et al. [36]. Among other possibilities, these types of flames can be established by directing a stream of combustible fresh gases towards a wall (single jet-wall configuration, c.f. Fig. 3), or by two counterflowing jets of the same chemical composition (symmetric twin flame configuration). The most prominent characteristic of this flame configuration is that (in the absence of wall heat losses) the only external flow effect on the flame is flame-tangential strain, which can be studied and quantified systematically [36].

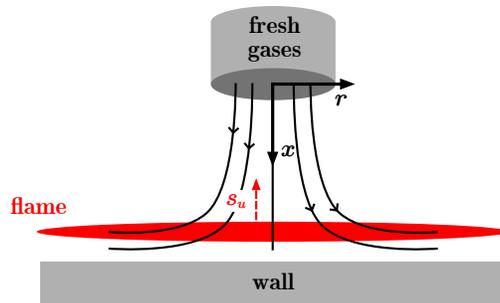


Figure 3: Schematic of a planar stagnation flame.

Physical space model. For planar flames established in a stagnation flow it has been shown that thermochemical quantities are a function of the axial coordinate x only [63]. Thus, a one-dimensional equation set can be derived from a similarity transformation which describes the flame structure along the flame centerline. Using the model by Kee et al. [63] a stagnation flame is solved imposing inflow boundary conditions from the nozzle and an adiabatic wall on the other side, with zero gradient boundary conditions for thermochemical quantities. For the flow, the wall is modeled as a slip wall, where the axial velocity becomes zero, but the radial velocity gradient is allowed

to assume finite values. The numerical model therefore corresponds to a symmetrical twin flame configuration [64].

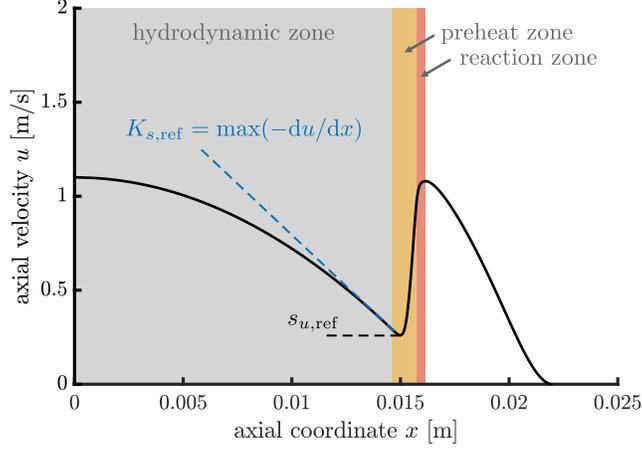


Figure 4: Flow field of the stagnation flame with the reference flame speed $s_{u,\text{ref}}$ and the reference strain rate $K_{s,\text{ref}}$, evaluated from the steepest velocity gradient upstream of the flame. Recreated according to Ref. [36].

Flow profile and strain rate. To facilitate a reasonable comparison of the numerical results with experiments, the flow field and strain profile of stagnation flames are examined. Figure 4 shows the flow field solution of the stagnation flame which can be divided into three characteristic zones: (1) the hydrodynamic zone, (2) the premixed flame’s preheat zone, where the flow velocity increases due to thermal expansion, and (3) the premixed flame’s reaction zone [36]. For this characteristic flow profile two quantities can be identified which are readily accessible with experimental methods: the minimum axial velocity upstream of the flame $s_{u,\text{ref}}$ (note that $s_u^0 \neq s_{u,\text{ref}}$) and the steepest gradient of the velocity profile upstream of the flame $K_{s,\text{ref}}$. The quantity $s_{u,\text{ref}}$ can be recorded as a function of $K_{s,\text{ref}}$ experimentally and this method is commonly used to determine s_u^0 by extrapolation to zero stretch [36, 65, 66].

Modeling strain in progress variable space. For the premixed flamelet model the strain rate is defined according to Eq. (25), which is shown for an exemplary flame configuration in Fig. 5 in physical space (coordinate x , left) and in normalized progress variable space (coordinate c , middle). Note that the strain rate $K_{s,\text{ref}}$ is proportional to the gradient of the velocity u (c.f. Fig. 4), whereas K_s , the flamelet parameter, is proportional to the gradient of the mass flux ρu . Nonetheless, both quantities agree well at the unburned edge of the flame ($c = 0$) since the effect of thermal expansion (thus, the density gradient) is small at this point. Further inspecting the profile of K_s , it is found

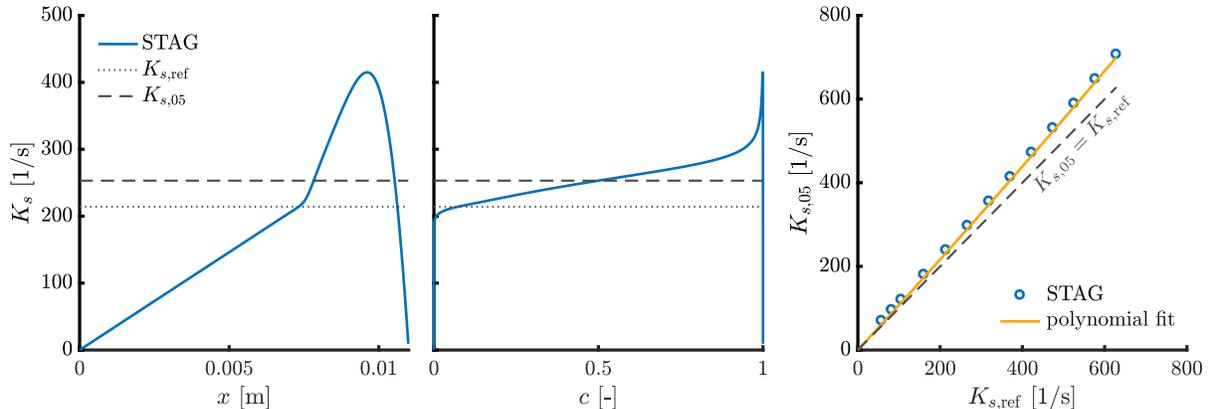


Figure 5: Strain rate K_s for a lean H_2 -air stagnation flame with $K_{s,\text{ref}} = 214\text{s}^{-1}$ versus the physical space coordinate (left) and the normalized progress variable c (middle). For the premixed flamelet model $K_{s,05}$ is prescribed which is expressed as a function of $K_{s,\text{ref}}$ to facilitate a comparison of the premixed flamelet model to experiments (right).

that K_s increases from the unburned to the burned edge of the flame. In c -space, K_s is almost linear for $0.1 < c < 0.9$, but it exhibits steep gradients towards the boundaries. In the following it will be shown that the premixed flamelet model yields accurate results for stagnation flames if the profile of K_s is approximated by a representative value. Therefore, the reference strain rate $K_{s,05}$ is introduced which is defined as the strain rate K_s evaluated at $c = 0.5$. **This choice is motivated by the fact that strain should be prescribed at the inner layer of a premixed flame to recover its flame structure accurately [39]. It should be noted that the inner structure of a premixed flame becomes stretched when conditioning on the progress variable [22] which implies that also other choices for fixing the strain rate could be reasonable.**

Comparison between experiments and numerical results. Numerical results obtained with the physical space and the progress variable space model are now compared to experiments in terms of $s_{u,\text{ref}}$ and $K_{s,\text{ref}}$. While $s_{u,\text{ref}}$ is readily evaluated from all numerical solutions, $K_{s,\text{ref}}$ is a flow control parameter of the premixed flamelet model and thus it needs to be prescribed. Similarly, $K_{s,05}$ is inaccessible from experiments. Therefore, $K_{s,\text{ref}}$ and $K_{s,05}$ are related to each other by a second order polynomial fit through the physical space computations to assign representative strain rates $K_{s,\text{ref}}$ to every premixed flamelet solution, which is shown in Fig. 5 (right).

Figure 6 (left) shows $s_{u,\text{ref}}$ versus $K_{s,\text{ref}}$ for lean H_2 -air flames ($\varphi = 0.35$) at atmospheric conditions. It contains numerical and experimental data by Vagelopoulos et al. [66] who determined laminar burning velocities with the counterflow twin flame technique. Furthermore, Fig. 6 contains the physical space solution (Eqs. (2)) for the stagnation flames (STAG), the corresponding results

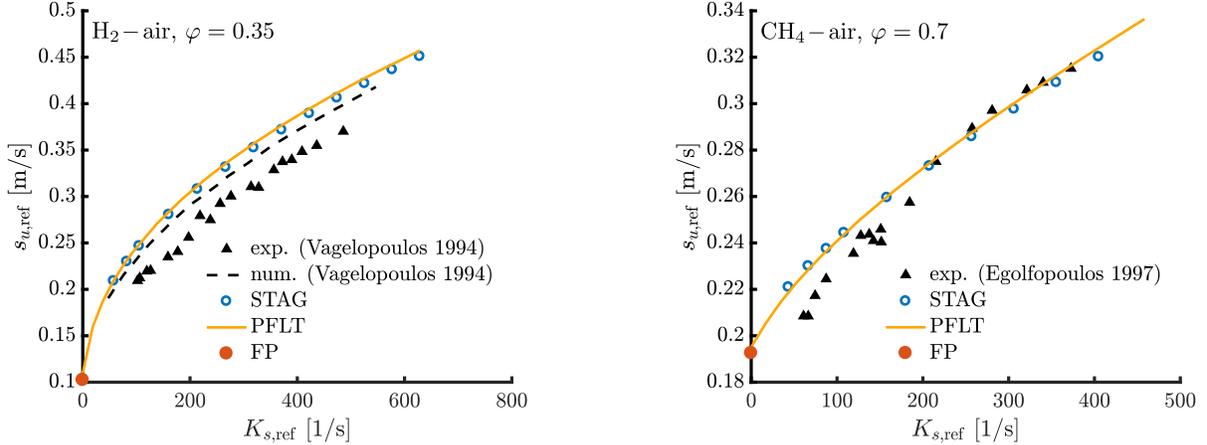


Figure 6: Reference flame speed $s_{u,\text{ref}}$ versus strain rate for atmospheric H_2 -air ($\varphi = 0.35$, left) and CH_4 -air stagnation flames ($\varphi = 0.7$, right). Experimental and numerical results by Vagelopoulos et al. [66] and Egolfopoulos et al. [64] are shown in comparison to numerical results obtained from a physical space solution method (STAG) and the premixed flamelet model in progress variable space (PFLT). The unstretched laminar burning velocity evaluated for a freely-propagating flame in physical space is shown as a reference (FP). The progress variable is defined as $Y_c = Y_{\text{H}_2\text{O}} - Y_{\text{H}_2} - Y_{\text{O}_2}$ (H_2 -air flames) and $Y_c = Y_{\text{CO}_2} + Y_{\text{H}_2\text{O}} + 10 Y_{\text{H}_2} - Y_{\text{O}_2}$ (CH_4 -flames) in the flamelet calculations.

for the premixed flamelet model (PFLT, Eqs. (22)-(24)) and the unstretched laminar burning velocity obtained for a freely-propagating flame as a reference (FP). As observed from Fig. 6, all the numerical results overpredict the experimental results (including previous simulations reported in Vagelopoulos et al. [66]), however the overall trend is well captured. Notably, the flamelet solutions agree well with the physical space reference solutions of the stagnation flames.

Similarly as for the H_2 -air flames, Fig. 6 (right) shows $s_{u,\text{ref}}$ versus $K_{s,\text{ref}}$ for lean atmospheric CH_4 -air flames ($\varphi = 0.7$). The numerical results agree well with the experimental data at strain rates $K_{s,\text{ref}} > 200 \text{ s}^{-1}$, while there are deviations for smaller strain rates. Again, the flamelet model recovers the physical space reference solutions well. Keeping in mind that s_u scales with $s_{u,\text{ref}}$, these results illustrate that the premixed flamelet model accurately captures the influence of strain on the laminar burning velocity. This is further supported by comparing the profiles of the the temperature, the progress variable gradient, and species mass fractions between the STAG and the PFLT model, which is shown in Fig. 7 for the previously discussed H_2 -air flames. It is found that the profiles obtained with the physical space reference model are very accurately recovered with the flamelet model for different strain rates. As shown in Fig. 7, this is also the case for highly diffusive species such as H_2 and radicals such as H . Similar characteristics are observed for other major and minor species (not shown for brevity).

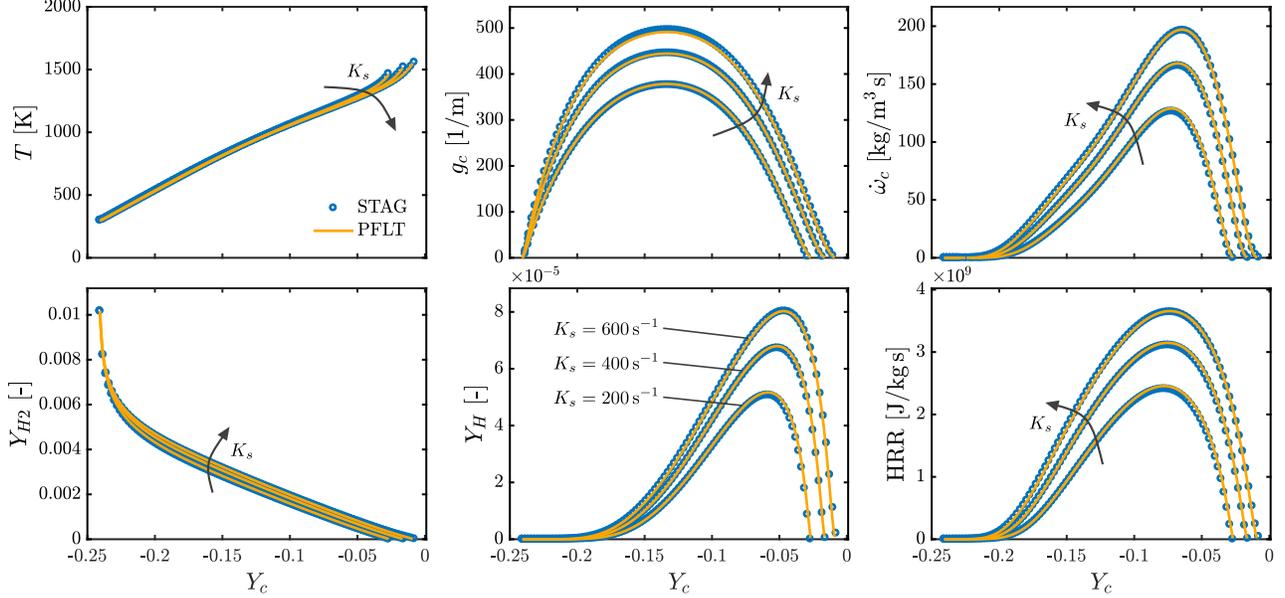


Figure 7: Profiles for the temperature, the progress variable gradient, the progress variable source term, the heat release rate (HRR), and the species mass fractions of H_2 and H for the lean H_2 -air flames studied by Vagelopoulos et al. [66] (c.f. Fig. 6, left). Results obtained with the physical space model (STAG) are compared to flamelet solutions (PFLT) for three different strain rates ($K_s = 200|400|600 \text{ s}^{-1}$). The progress variable is defined as $Y_c = Y_{\text{H}_2\text{O}} - Y_{\text{H}_2} - Y_{\text{O}_2}$.

Strain-induced effects for premixed flames with $\text{Le} < 1$. After the comparison to experimental reference data, characteristics of strained premixed flame structures are investigated in more detail with the numerical models. Positive (extensive) strain generally increases gradients in the flame [62], which is properly reflected in the flamelet model, since positive strain acts as a production term for the gradient g_c , c.f. Eq. (24). Strain thereby inherently promotes diffusive transport of reactants to the reaction zone (proportional to species gradients) but also the heat conduction away from the flame (proportional to temperature gradients). For flames with effective Lewis numbers smaller than unity, e.g. lean H_2 -air flames ($\text{Le}_{\text{H}_2} \approx 0.3$), positive strain therefore strengthens the flame [54, 62].

Figure 8 compares three lean H_2 -air flames computed with physical space solution methods (unstrained: FP, strained: STAG, $K_{s,05} = 0 / 500 / 2000 \text{ s}^{-1}$) against corresponding results obtained with the premixed flamelet model. From the right graph it is observed for an unstretched flame that the local equivalence ratio φ decreases from the fresh gas mixture (left boundary) before it returns to the original equivalence ratio in the post-oxidation zone (right boundary). The much leaner mixture for intermediate progress variables is explained by the fact that H_2 diffuses much faster than other species towards the reaction zone where it is subsequently consumed. For strained flames, it is found that the flame becomes enriched towards the right boundary which is in accordance

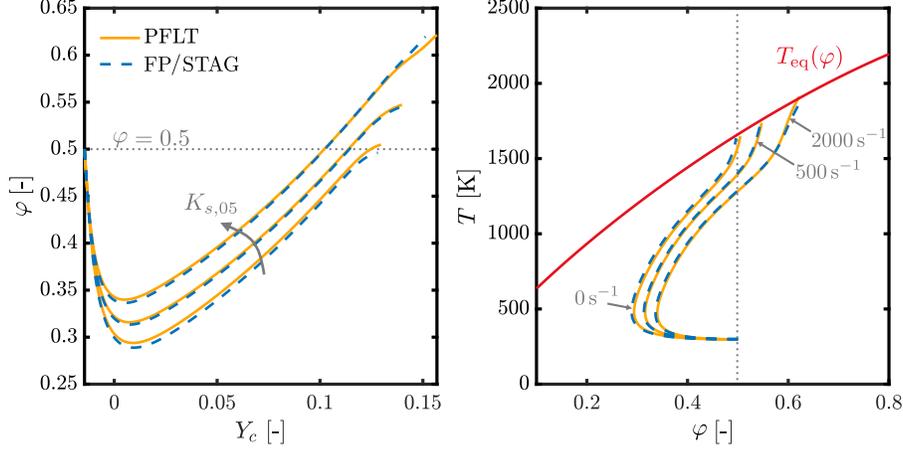


Figure 8: Left: local equivalence ratio versus progress variable for three lean H_2 -air flames at atmospheric conditions imposing different strain rates ($\varphi = 0.5$, $K_{s,05} = 0/500/2000 \text{ s}^{-1}$). The results are obtained with physical space solvers (FP/STAG) and the premixed flamelet model (PFLT). Right: local temperature versus local equivalence ratio. The solid red line marks the equilibrium temperature T_{eq} computed for H_2 -air mixtures as a function of the equivalence ratio φ . The progress variable is defined as $Y_c = Y_{\text{H}_2\text{O}} - Y_{\text{H}_2} - Y_{\text{O}_2}$ in the flamelet calculations.

with the explanation for the strain effect given above.

The right graph of Fig. 8 shows the local temperature T as a function of the local equivalence ratio φ . Furthermore, the equilibrium temperature T_{eq} is depicted, which is computed for H_2 -air mixtures as a function of φ . It is observed that unstrained and strained flames reach a local chemical equilibrium in the post oxidation zone, since the flame temperature eventually reaches a point on the profile of $T_{eq}(\varphi)$. It is found that positively strained flames with $Le < 1$ reach a higher flame temperature than the adiabatic flame temperature of their respective fresh gas mixture due to local fuel enrichment. This strain-induced effect is sometimes termed *hot spot* or *super-adiabatic* and can also be observed in turbulent combustion [50–52] (curvature can have similar effects). The comparison above further confirms that the expansion of the Y_c -space (local enrichment effect) is correctly captured by the flamelet model and the right boundary eigenvalue of the Y_c -space is well-predicted.

Extinction limits for premixed flames with $Le < 1$. Besides positive strain rates, the flamelet model in progress variable space allows prescribing negative strain rates. This is demonstrated by examining extinction limits for both, positive and negative strain rates, next. The effect of strain on the flame structure is characterized considering quantities at the inner layer, which can be defined by the position of the maximum heat release rate [36, 39]. Note that the reference quantity $s_{u,\text{ref}}$ is not suitable for this analysis since it not only depends on strain, but also on thermal expansion

which complicates a proper assessment of strain-induced effects [36].

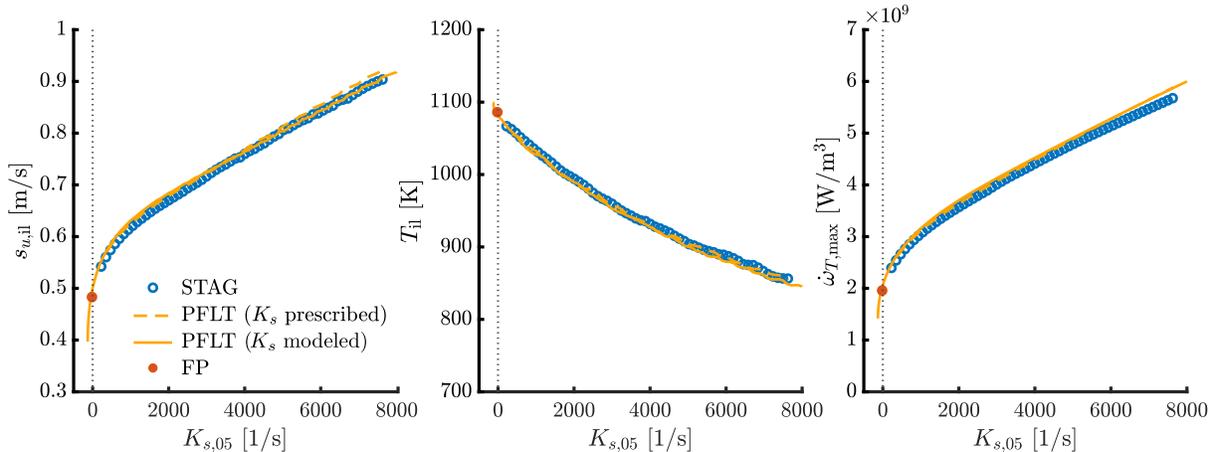


Figure 9: Inner layer quantities: burning velocity $s_{u,il}$, temperature T_{il} and maximum heat release rate $\dot{\omega}_{T,max}$ as functions of strain for lean H_2 -air flames ($\varphi = 0.5$, $T_u = 298$ K). The progress variable is defined as $Y_c = Y_{H_2O} - Y_{H_2} - Y_{O_2}$ in the flamelet calculations.

Figure 9 shows the numerical results obtained for lean H_2 -air flames ($\varphi = 0.5$) towards higher positive strain rates and also for negative strain rates. Besides results for the physical space stagnation flame model (STAG, blue circles) and the premixed flamelet model with K_s approximated by a representative value (PFLT, solid line), the latter has also been computed with the full strain rate profile extracted from corresponding STAG solutions (PFLT, dashed line). Furthermore, numerical results for unstrained flames (FP, red circle) are given for orientation.

The strain-induced strengthening of flames with $Le < 1$ is reconfirmed considering the burning velocity at the inner layer $s_{u,il}$ and the maximum heat release rate $\dot{\omega}_{T,max}$ in Fig. 9 (left and right plot, respectively) which both increase with positive strain. The inner layer temperature T_{il} decreases for $K_{s,05} > 0$ (Fig. 9, mid plot) which implies that the maximum heat release becomes shifted to lower temperatures. As long as solutions for the physical space model exist, the inner layer quantities are well reproduced by the premixed flamelet model (Eqs. (22)-(24)). The flamelet model with the prescribed strain profile virtually yields the same result compared to the model with a representative constant strain rate.

It is found that the stagnation flame extinguishes at $K_{s,05} = 7644 s^{-1}$ while the premixed flamelet model predicts solutions also for significantly higher strain rates (not explicitly shown). This is due to the fact that flames with $Le < 1$ do not become extinguished due to positive strain effects (which strengthen the flame), but due to incomplete reactions, when they are pushed against

a wall or stagnation surface. In his book, Law [54] therefore calls stagnation flames *restrained*. Approaching the stagnation surface, the thickness of the flame becomes smaller leading to reduced residence times of reactants. Increasing the strain rate, restraintment initially appears as an effect in the post-oxidation zone, where the thermochemical state does not reach equilibrium conditions anymore (lowered temperatures). Eventually increasing strain above a critical limit, residence times of reactants are so small that the flame becomes extinguished. On the contrary, the premixed flamelet model does not account for restraintment, since both boundaries are infinitely removed from the reaction zone of the premixed flame. Thus, flamelet solutions can be found for much higher strain rates and computations were aborted at $K_{s,05} = 15\,000\text{ s}^{-1}$. For modeling turbulent flames, which can exhibit very large strain rates [37], these characteristics of the premixed flamelet model could become advantageous, because highly strained canonical flames can be considered aside from artefacts induced by a given physical space configuration.

At zero strain (vertical dotted line in Fig. 9), the premixed flamelet model recovers the characteristics of the freely-propagating flame model (FP). For $K_{s,05} < 0$, it is found that the lean H_2 -air flames show a clear extinction limit since negative strain weakens the flame for $\text{Le} < 1$. The negative strain extinction limit is determined as $K_s = -130\text{ s}^{-1}$ for which the profiles of $s_{u,il}$ and $\dot{\omega}_{T,\max}$ versus $K_{s,05}$ become almost singular. To the authors' knowledge, this negative strain extinction limit is inassessible with conventional low-dimensional flame solvers.

Extinction limits for premixed flames with $\text{Le} > 1$. Opposed to the previously discussed H_2 -air flames, lean $\text{C}_2\text{H}_5\text{OH}$ -air flames exhibit an effective Lewis number larger than unity ($\text{Le}_{\text{C}_2\text{H}_5\text{OH}} \approx 1.6$) and positive strain weakens these flames [54, 62]. This strain-induced effect is displayed in Fig. 10 which shows the stagnation flame results obtained with the physical space solver (STAG) compared to flamelet solutions (PFLT, with constant K_s), and freely-propagating flame solutions for reference. Also for the $\text{C}_2\text{H}_5\text{OH}$ -air flames T_{il} (Fig. 10, mid) decreases with positive strain, indicating a shift of the maximum heat release to lower temperatures. However, it is observed that the inner layer burning velocity $s_{u,il}$ (left) and the maximum heat release rate $\dot{\omega}_{T,\max}$ (right) both decrease for $K_{s,05} > 0$. Consequently, a positive strain extinction limit can be identified for lean $\text{C}_2\text{H}_5\text{OH}$ -air flames, which is predicted as 930 s^{-1} with the physical space STAG model and 902 s^{-1} with the PFLT model (deviation: 3%). Keeping in mind that constant strain rates are prescribed for the flamelet model, the small overall differences between the two models are acceptable.

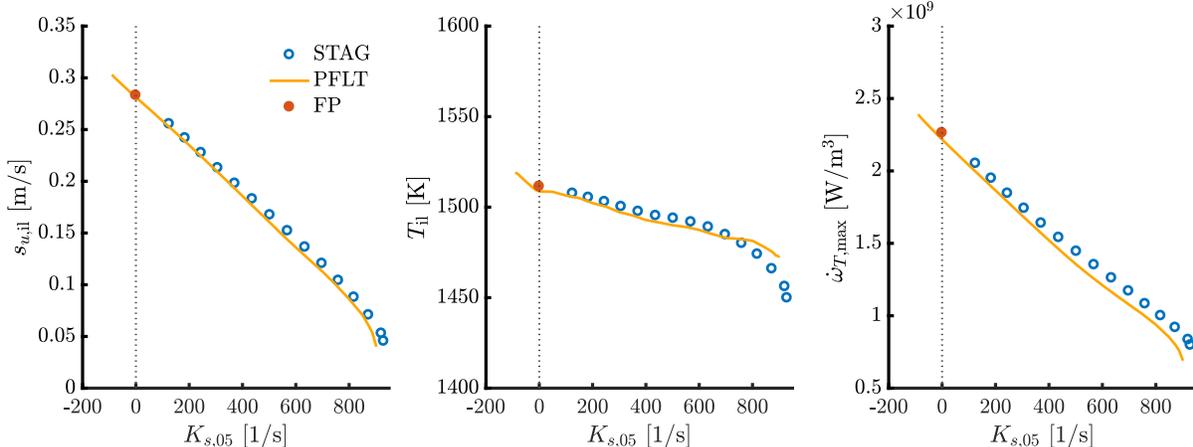


Figure 10: Inner layer quantities burning velocity $s_{u,il}$, temperature T_{il} and maximum heat release rate $\dot{\omega}_{T,max}$ as functions of strain for lean C_2H_5OH -air flames ($\varphi = 0.7$, $T_u = 363$ K). The progress variable is defined as $Y_c = Y_{CO_2} + Y_{H_2O} + 10Y_{H_2} - Y_{O_2}$ in the flamelet calculations.

Similar to H_2 -air flames, C_2H_5OH -air flames can tolerate a certain level of negative strain. It is found from Fig. 10, that the maximum flame temperature and the maximum heat release both increase with negative strain rates, which is contrary to H_2 -air flames. The negative strain extinction limit is determined as $K_{s,05} = -90 \text{ s}^{-1}$, which appears abrupt, since none of the quantities depicted in Fig. 10 visually indicates a limit. However, below this level no flamelet solution could be obtained and it remains an open question whether this is a physical effect or a numerical artifact.

The results presented in this section confirm that the premixed flamelet model captures strain effects on premixed flame structures accurately. It is further demonstrated that prescribing strain by representative values in progress variable space is a sufficiently good model, which is a valuable insight for flamelet tabulation strategies.

5.2.2. Tubular stagnation flames

Figure 11 shows the schematic of a tubular burner. A combustible fresh gas mixture is issued inwardly from a porous tube and forms a cylindrically symmetric stagnation flow. Eventually, gases exit the burner in the axial direction. With this burner setup a stationary tubular flame is established which is stretched and curved at the same time. Increasing the inlet velocity, the flame is pushed towards the stagnation line and flame strain and curvature are both increased.

Physical space model. In physical space, the tubular burner can be described by the equation set published by Wang et al. [67], which is similar to the model for stagnation flames. From a similarity transformation of the governing equations in physical space a one-dimensional equation set with

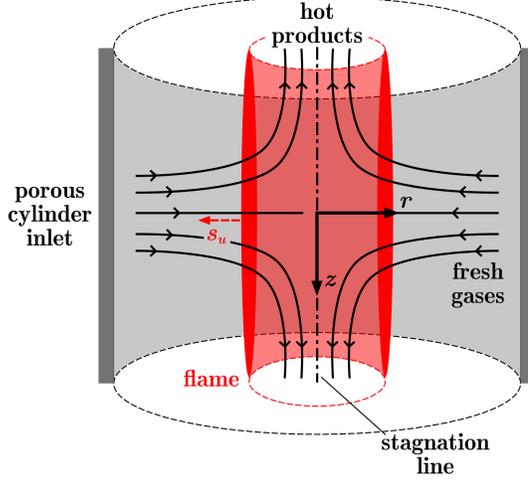


Figure 11: Schematic of a tubular burner in which a premixed tubular stagnation flame is established.

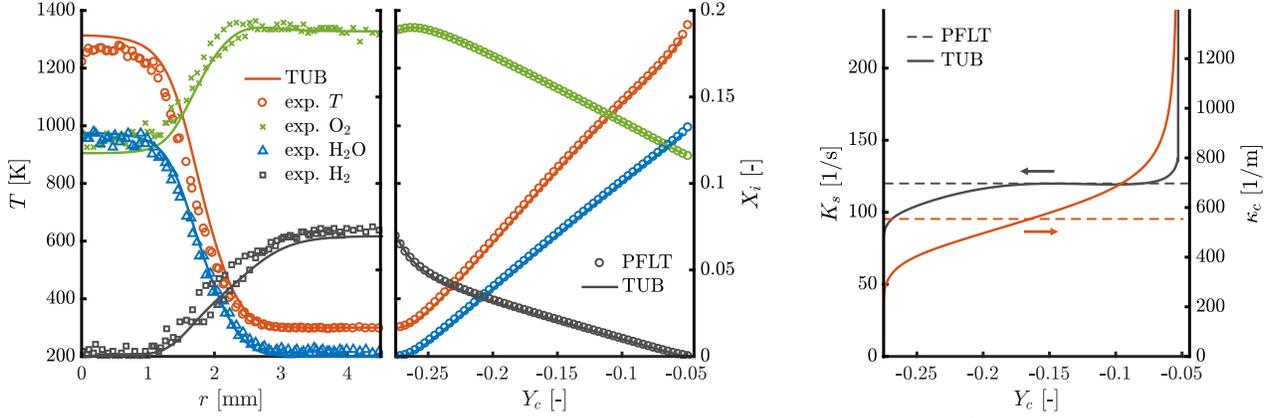


Figure 12: Flame structure of a lean H_2 -air tubular flame ($\varphi = 0.175$, $K_{s,ref} = 145\text{s}^{-1}$). Left: Comparison of the 1D flame simulation data for the tubular burner (TUB) with corresponding experimental data by Hu et al. [68]. Middle: Comparison of the physical space simulation data (TUB) to the results obtained with the premixed flamelet model (PFLT). Right: Comparison of the extracted profiles for strain K_s and curvature κ_c from the physical space numerical solution to the prescribed profiles for the premixed flamelet model. The progress variable is defined as $Y_c = Y_{\text{H}_2\text{O}} - Y_{\text{H}_2} - Y_{\text{O}_2}$ in the flamelet calculations.

respect to the radial coordinate r is obtained. This model will be used to validate the flamelet model in terms of curvature-affected premixed flames.

Comparison between experiments and numerical results. For the tubular burner, experimental data of temperature and major species concentrations are available from the work by Hu et al. [68]. Figure 12 shows the profiles of temperature, O_2 , H_2O and H_2 for an ultra lean H_2 -air tubular flame ($\varphi = 0.175$, $K_{s,ref} = 145\text{s}^{-1}$). As observed from the left graph, the position and the structure of the flame are well captured by the physical space numerical solution. While the temperature is slightly overpredicted, the numerical results recover the species profiles well. In the mid-graph the physical space numerical solution is compared to the progress variable space numerical solution

obtained with the flamelet model. It is found that the flamelet solution is essentially equivalent to the physical space solution. A small difference is found at the right boundary (post oxidation zone) which extends further for the flamelet model. Similar to the stagnation flame, this characteristic is attributed to the restraintment of the physical space flame configuration at the stagnation point. To capture both strain and curvature effects with the premixed flamelet model, representative values for K_s and κ_c were determined from the physical space solution at $c = 0.5$ and prescribed for the flamelet model.

5.2.3. Spherical expanding flames

The spherical expanding flame is established by igniting an initially quiescent gas mixture with a spark from which subsequently an outwardly propagating flame develops. For controlling the flame, experiments are carried out in closed vessels which further allow studying flame propagation under high pressures. The flame evolution is usually recorded by Schlieren or shadow photography using high speed cameras. With this technique, the burning velocity can be determined from the observation of the flame speed together with the unburned and burned gas density ratio [49] or directly from the difference of the propagation speed of the flame s_f and the fresh gas velocity u_g [49, 61, 69, 70] according to Eq. (35). Figure 13 shows a schematic of a spherical expanding flame. The flame propagates outwardly as indicated by the red arrows and it is further pushed in the same

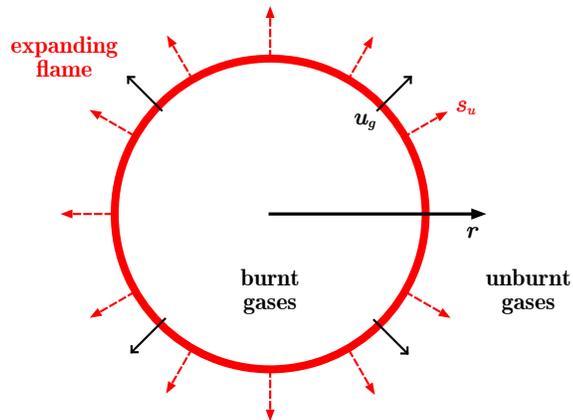


Figure 13: Schematic of a spherical expanding flame.

direction due to thermal expansion of the reaction products. This thermal expansion causes an entrainment of cold unburned gas by the flame which can be measured as the velocity of unburnt gases u_g ahead of the flame front [61, 69]. Due to symmetry, the flame can also be described by a

set of time-dependent and spatially one-dimensional differential equations [71] (with respect to the radial coordinate r).

Flame stretch and curvature. The flame's surface increases continuously with time since it is curved and it moves outwardly with the flame speed s_f . Consequently the flame front becomes stretched. The overall stretch rate of the spherical expanding flame front is determined as [35]

$$K = \frac{2}{r} \frac{dr}{dt} = \frac{2}{r} s_f = \kappa_c s_f, \quad (36)$$

where the curvature of the flame front $\kappa_c = 2/r$ evolves with time. When computing spherical expanding flames with the premixed flamelet model, strain and curvature need to be prescribed. While the definition of the curvature is readily available from the flame geometry, the premixed flamelet model further requires the strain rate *imposed by the flow* which is defined according to Eq. (25)

$$K_s = K - s_u \kappa_c \stackrel{(36)}{=} -(s_f + s_u) \kappa_c \stackrel{(35)}{=} -u_g \kappa_c. \quad (37)$$

Note that u_g is negative in this equation due to the convention $u_g = \mathbf{n}_c \cdot \mathbf{u}$. For the spherical expanding flame \mathbf{u} is directed opposite to \mathbf{n}_c , thus, the strain rate K_s is positive which is consistent with the literature. Furthermore, in the above relation the flame displacement speed s_d and the flame's burning velocity s_u are set equal which is valid only at the unburnt edge of the flame as outlined in Sec. 5.2. For spherical expanding flames usually s_f is reported as a function of K which allows to recompute the flame's curvature according to $\kappa_c = K/s_f$. However, the premixed flamelet model can only be applied if also the gas velocity u_g is reported from the experiments. Otherwise, imposed strain by the flow $K_s = -u_g \kappa_c$ and strain due to the propagation of the flame itself ($s_u \kappa_c$, which is a flame response) cannot be separated. Suitable reference data has been reported by multiple researchers for hydrogen and hydrocarbon flames [61, 69, 70, 72].

Comparison between experiments and numerical results. Figure 14 shows experimental reference data by Varea et al. [70] for lean H_2 -air flames together with results obtained with the premixed flamelet model. Similarly as for planar and tubular stagnation flames, strain and curvature are prescribed as representative constants which were determined with average values from the experiments from Eqs. (36) and (37). Overall the agreement between experiments and numerical results for the

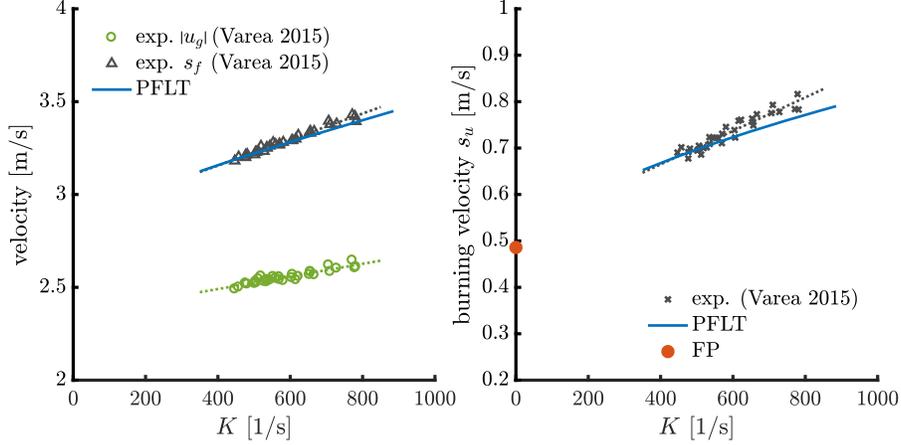


Figure 14: Left: Measurements of the apparent flame speed s_f and the velocity of the fresh gases u_g for a spherical expanding flame by Varea et al. [70] (H_2 -air, $\varphi = 0.5$). Numerical results are obtained with the premixed flamelet model (PFLT, solid line) and the dotted lines mark a linear regression of the experimental data. Right: comparison of burning velocities evaluated from the experimental data and the premixed flamelet model. The laminar burning velocity for a freely-propagating flame is shown for reference (FP). The progress variable is defined as $Y_c = Y_{\text{H}_2\text{O}} - Y_{\text{H}_2} - Y_{\text{O}_2}$ in the flamelet calculations.

stretched laminar burning velocity (Fig. 14, right graph) is good with only slight deviations towards higher stretch rates. However, these deviations are less prominent when comparing the overall flame speed s_f (Fig. 14, left graph).

Similar results are obtained for lean CH_4 -air flames for which experimental data by Varea et al. [61] and corresponding flamelet solutions are shown in Fig. 15. The overall trend and magnitude

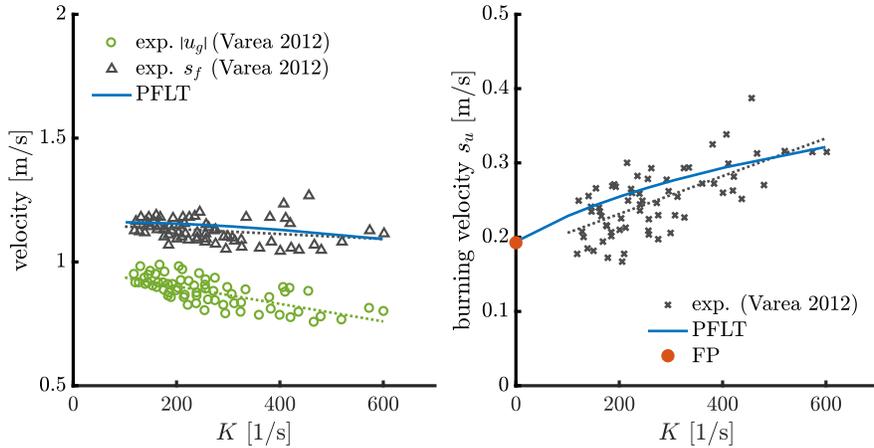


Figure 15: Left: Measurements of the apparent flame speed s_f and the velocity of the fresh gases u_g for a spherical expanding flame by Varea et al. [61] (CH_4 -air, $\varphi = 0.7$). Numerical results are obtained with the premixed flamelet model (PFLT, solid line) and the dotted lines mark a linear regression of the experimental data. Right: comparison of burning velocities evaluated from the experimental data and the premixed flamelet model. The laminar burning velocity for a freely-propagating flame is shown for reference (FP). The progress variable is defined as $Y_c = Y_{\text{CO}_2} + Y_{\text{H}_2\text{O}} + 10Y_{\text{H}_2} - Y_{\text{O}_2}$ in the flamelet calculations.

for the stretched laminar burning velocity (right graph) is well captured by the flamelet model with slightly overpredicted values for stretch rates between 150 to 400 s⁻¹. However, the comparison of the flame speed s_f (Fig. 15, left graph) shows that the deviation between numerical result and linear regression of the experimental data remains small.

Only few information is necessary to reproduce magnitude and trend of the experimentally determined burning velocities. This underlines the high flexibility of the premixed flamelet model and confirms that the model captures the interplay of strain, curvature, and unsteadiness. Given a pair of parameters, K_s and κ_c , one computation with the flamelet model yields the premixed flame response, e.g. in terms of its flame structure and burning velocity. [With few exceptions \[39\], most physical space models compute the whole transient evolution of the spherical expanding flame \(until a self-similar state\) before this information becomes available.](#)

6. Summary and conclusions

A self-contained premixed flamelet model was presented which describes premixed flames in progress variable space. The model is based on flamelet equations for temperature, species mass fractions, and the gradient of the progress variable. The latter establishes a direct link of the flamelet structure to the fundamental flame parameters strain and curvature. As already shown in an earlier work [22], treating the progress variable gradient as a dependent variable resolves a severe limitation of flamelet modeling in scalar space. Similar previous flamelet formulations [21, 23, 53] required a closure for both the shape and the magnitude of the gradient, for which a generic profile does not exist in case of premixed combustion. With the novel approach, an effective closure of the flamelet equations was obtained by prescribing the flamelet parameters strain and curvature through representative constant values.

The predictive capabilities of the new method were demonstrated for a constant pressure ignition problem and three canonical flame configurations: the stagnation flame, the tubular flame and the spherical expanding flame. Different models are necessary to describe the aforementioned canonical combustion configurations in the physical space, but all of them can be described using a single premixed flamelet model in progress variable space. A combination of physical space reference solutions and experimental measurements from the literature was used for the validation. Very accurate results were found for constant pressure ignition, flame structures, burning velocities, and

flame extinction. Furthermore, the premixed flamelet model allowed to capture strain-induced hot spots and even negative strain effects. [With few exceptions \[39, 40\], the latter remains inaccessible for conventional low-dimensional flame solvers.](#)

Multiple future applications of the premixed flamelet model are conceivable. [Its generic parametrization allows analyses of flame structures which do not have a direct canonical flame counterpart in the physical space. This facilitates parametric studies well beyond the limits of, for instance, tubular flames or stagnation flames in terms of curvature and strain. With this, the model is well suited for tabulated chemistry approaches in a CFD context since these conditions have been shown to be relevant for turbulent combustion. Further, it can be used for numerically assisted interpretation of experimental data.](#) The determination of the laminar burning velocity from experiments of stretched premixed flames usually relies on extrapolation procedures to zero stretch [36]. Extrapolation methods could be complemented using the work reported in this paper, since it allows for a continuous representation of the burning velocity over the full stretch range.

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