

DNS OF NEAR WALL DYNAMICS OF PREMIXED CH₄/AIR FLAMES

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Abstract. This work presents a numerical study on the effect of flame-wall interaction (FWI) from the viewpoint of flame dynamics. For that purpose, direct numerical simulations (DNS) employing detailed calculations of reaction rates and transport coefficients have been applied to a 2D premixed methane/air flame under atmospheric condition. Free flame (FF) and side-wall quenching (SWQ) configurations are realized by defining one lateral boundary as either a symmetry plane for the FF or a cold wall with fixed temperature at 20 °C for the SWQ case. Different components of flame stretch and Markstein number regarding tangential, normal (due to curvature) and total stretch, Ka_s , Ka_c and $Ka_{tot} = Ka_s + Ka_c$, as well as their correlations with respect to the local flame consumption speed S_L have been evaluated.

It has been shown that the FWI zone is dominated by negative flame stretch. In addition, S_L decreases with decreasing normal stretch due to curvature Ka_c while approaching the cold wall. However, S_L increases with decreasing Ka_c while approaching the symmetry boundary for the free flame case, leading to an inversion of the Markstein number Ma_{tot} based on Ka_{tot} from positive in the free flame case to negative in the SWQ case. The quenching distance evaluated based on wall-normal profiles of S_L has been found to be approximately equal to the unstretched laminar flame thickness, which compares quantitatively well with measured data from literature. The flame speed has been confirmed to scale quasi-linearly with the stretch in the FWI zone. The results reveal a distinct correlation during transition between FWI and FF regarding flame dynamics, which brings a new perspective for modeling FWI phenomena by means of flame stretch and Markstein number. To do this, the quenching effect of the wall may be reproduced by a reversed sign of the Markstein number from positive to negative in the FWI zone and by applying the general linear Markstein correlation ($S_L/S_{L,0} = 1 - Ma \cdot Ka$), leading to a decrease of the flame speed or the reaction rate in the near-wall region.

1 INTRODUCTION

In order to achieve a high power density and to improve fuel efficiency, most combustion applications take place in an enclosed environment with a relatively small volume, for instance in internal combustion or jet engines. The flame in this case propagates against a cold wall and is extinguished in the direct vicinity of the wall. As a result, the phenomenon of flame-wall interaction (FWI) significantly contributes to the flame stability and efficiency of the combustion process, but also to the pollutant emissions and lifetime of the burner [1]. The effect of FWI has been studied extensively in the last years due to its importance [2, 3, 4, 5, 6, 7, 8, 9, 10]. For instance, Dreizler and Böhm [2] and Kosaka et al. [9] performed comprehensive laser diagnostics for detailed measurements of flow velocity, temperature, species concentrations and flame structure for premixed methane/air and DME/air flames in the FWI zone, which improved the understanding of the related chemo-physical phenomena substantially. The same experimental setup has then been studied numerically by Ganter et al. [5] and Heinrich et al. [4] using different reaction-diffusion models and 2D/3D simulations, where the calculated wall-normal temperature profiles, heat fluxes and wall distances showed a good agreement with measured data. Saffman [11] measured the quenching distances for different hydrocarbon fuels and equivalence ratios. The experimental and numerical works by Häber and Suntz [6] and Strassacker et al. [7] have been focused on the effect of different wall materials, including heterogeneous wall reactions. They found that the quenching distance is only weakly dependent on the wall material and type of coating. The 3D direct numerical simulations (DNS) performed by Gruber et al. [12] for turbulent hydrogen/air flame interacting with a cold wall showed that the local flame thickness and propagation speed vary near the boundary layer, resulting in a regime change from a freely propagating flamelet to a thickened flame at the wall. In addition, the DNS results suggest that the near-wall turbulence plays an important role on the convective heat transfer by pushing the reaction zone towards the wall. Mejia et al. [8] studied the influence of wall temperatures on the flame's response to acoustic perturbations, where the flame instability can be controlled or suppressed by changing solely the temperature of the burner rim. The numerical simulation by Poinso et al. [10] showed that the wall can affect the flame evolution even if the distance between the flame and the wall is significant. Popp and Baum [13] presented a numerical study of FWI for a stoichiometric methane/air flame, where a one-step chemical reaction failed to predict heat fluxes through the wall. Zirwes et al. [14] performed 2D/3D DNS of side-wall quenching (SWQ) of premixed methane/air flames at different equivalence ratios. The calculated quenching distances however showed a consistent constant deviation from the measured data.

Despite the progresses made in the previous works, studying FWI is challenging as the relevant phenomena occur in a very thin near-wall layer with a thickness of the order of $O(0.1\text{ mm})$. Focus of the recent numerical works was also to validate the proposed modeling concepts, such as the influences of using 2D/3D grids or detailed/reduced reaction-diffusion models, on reproducing the measured data [4, 5, 7, 15, 14]. In addition, mostly under-resolved grid resolutions have been applied in these works. Moreover, flames approaching walls are dominated by unsteady effects, which further complicate the FWI phenomena and make the comparison between measurement and simulation difficult. Therefore, more fine-grained simulations of FWI with sufficiently high grid resolutions, detailed reaction and transport models, as well as well-defined boundary conditions are necessary for a thorough understanding of the underlying processes of FWI. In an attempt to achieve this, DNS have been performed in the current work for a 2D side-wall quenching (SWQ) setup operated with premixed methane/air mixtures. Detailed chemical kinetics and molecular transport models are used and the FWI zone is resolved with a

sufficiently fine grid. Special focus of the work is put on studying local flame dynamics regarding flame stretch, flame speed and the Markstein effect in the FWI zone, revealing systematically the significant role of flame dynamics for elucidating the mechanism of flame extinction at a cold wall. In comparison to our previous work shown in [16], the analysis of near wall flame dynamics has been further refined to identify different stages of flame dynamics during FWI.

2 Theoretical Basics

Flame stretch is calculated in this work by [1, 17]:

$$K = (\mathbf{I} - \vec{n}\vec{n}) : \nabla \vec{u} + S_D \nabla \cdot \vec{n} \quad (1)$$

where \mathbf{I} is the unit tensor, \vec{u} the velocity vector and \vec{n} the unit normal vector of the flame surface. The first term $K_s = (\mathbf{I} - \vec{n}\vec{n}) : \nabla \vec{u}$ on the r.h.s. in Eq. (1) is attributed to hydrodynamic stretch caused by the non-uniform flow distribution and the second term $K_c = S_d \nabla \cdot \vec{n}$ due to the motion of a curved surface with the displacement speed S_D along the flame normal direction. $\vec{n} = \frac{\nabla Y_F}{|\nabla Y_F|}$ and S_D are evaluated at the flame surface defined by the fuel mass fraction $Y_F = Y_{F,iso}$ [18]

$$S_D = -\frac{1}{|\nabla Y_F|} \frac{dY_F}{dt} = -\frac{\dot{r}_F + \nabla \cdot (\rho D_F \nabla Y_F)}{\rho |\nabla Y_F|} \quad (2)$$

\dot{r}_F is the reaction rate of fuel, ρ the density and D_F the diffusion coefficient of fuel.

The laminar flame speed is evaluated in the present work by integration of the fuel reaction rate along its local normals

$$S_L = \frac{-1}{\rho_u (Y_{F,u} - Y_{F,b})} \int \dot{r}_F d\vec{n} \quad (3)$$

where $Y_{F,u}$ and $Y_{F,b}$ are the fuel mass fractions in the unburnt and burnt mixtures. Unlike S_D , S_L represents an integral value that is not dependent on the choice of iso-surface defining the flame front.

The influence of stretch on the laminar flame speed can be derived by means of a first order Taylor approximation for moderately low stretch around $K = 0$ for $S_L = S_{L0}$ [17]:

$$S_L(K) \approx S_L(K=0) + \left. \frac{\partial S_L}{\partial K} \right|_{K=0} (K-0) \quad \text{or} \quad \frac{S_L(K)}{S_{L0}} = 1 - \frac{\mathcal{L}}{S_{L0}} K \quad (4)$$

with the Markstein length $\mathcal{L} = -\partial S_L / \partial K|_{K=0}$ and the unstretched laminar flame speed S_{L0} . Equation (4) can be rewritten as

$$\frac{S_L(K)}{S_{L0}} = 1 - \frac{\mathcal{L}}{\delta_{L0}} \frac{\delta_{L0}}{S_{L0}} K = 1 - Ma Ka \quad (5)$$

to introduce the Markstein number Ma defined as the ratio of the Markstein length and the characteristic flame thickness δ_{L0}

$$Ma = \frac{\mathcal{L}}{\delta_{L0}} = -\frac{1}{\delta_{L0}} \frac{\partial S_L}{\partial K} \quad (6)$$

and the Karlovitz number $Ka = K \delta_{L0} / S_{L0}$. In this way, Ma measures the sensitivity of flame speed to stretch. A vast variety of experimental and computational work has been conducted in the last decades regarding the effect of stretch on flame speed [19, 20, 21, 22, 23, 24, 18], which justified the asymptotic correlation proposed in Eq. (5). However, most of the previous studies have been conducted for

1D canonical flame configurations with either outwardly expanding spherical flames or counterflow flames. The spherical flame is only positively curved and the opposed counterflow flame is only positively strained in tangential direction. In addition, the influence of heat loss on flame dynamics, for instance, due to a cold wall cannot be assessed. Therefore, detailed studies on flame dynamics are needed to extend existing knowledge for the combined impact of strain and curvature considering the negative stretch range, but also the effect of heat loss, for example, due to interaction with a cold wall.

3 Numerical Setups

Figure 1 shows a sketch of the 2D computational domain, which has a rectangular shape with lengths of 32 mm in streamwise and 8 mm in spanwise direction. The boundary conditions used for the four sides of the domain are indicated in Fig.1. Premixed methane/air mixture enters the domain at the inlet from the bottom with a temperature of $T_0 = 293$ K and a pressure of $p_0 = 1$ bar. The gradients of flow velocity, temperature and all chemical scalars are set to zero at the outlet. A parabolic velocity profile is used for the inflow, with a minimum velocity of zero in the center and a maximum value of $\hat{u} = 1.6$ m/s at both lateral ends. This leads to a V-shaped flame stabilized at the center and close to the inlet plane. The right side of the domain is set as a symmetry plane and the left boundary is defined either as a cold wall with 293 K or as a symmetry plane, as illustrated in Fig.1. These two simulation setups are identified in the following with the abbreviations SWQ (side-wall quenching) and FF (free flame), meaning that the left part of flame branch propagates against a cold wall or freely towards a symmetry plane. In addition, a non-reflecting boundary condition [25] has been applied to the inlet and outlet to bypass pressure waves at these boundaries. Premixed methane/air flames at different equivalence ratios $\Phi = 0.8, 1.0$ and 1.2 are examined in this work. The laminar flame parameters shown in Tab.1 are calculated from computations of 1D unstretched freely propagating flames with Cantera. S_L^0 and δ_L^0 from these simulations are used to evaluate Ka and Ma (see Eq.(5-6)). The flame surface is identified by the mass fraction of fuel corresponding to the highest consumption rate of fuel $Y_{f_{max}}$.

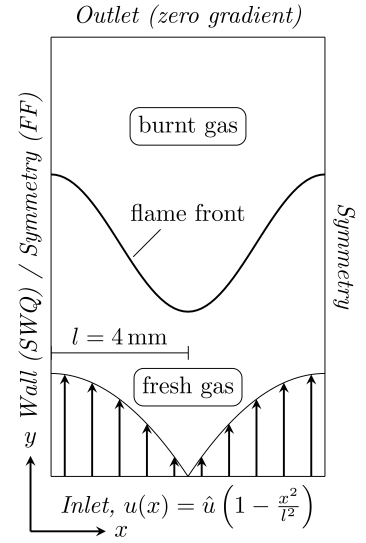


Figure 1: Computational domain and boundary conditions used for DNS.

Table 1: Laminar flame parameters from 1D flame calculations at different Φ and $p_0 = 1$ bar, $T_0 = 293$ K.

ϕ [-]	0.8	1.0	1.2
S_L^0 [cm/s]	26.17	36.41	32.19
δ_L^0 [mm]	0.536	0.445	0.472
$Y_{f_{max}}$ [-]	0.0039	0.0057	0.0068

The simulations have been performed using a DNS code developed in the framework of OpenFOAM, which includes detailed calculations of reaction rates and transport coefficients [26, 27, 28, 29, 30, 31, 32]. For the DNS the conservation equations for total mass, momentum, species masses and energy are solved in the compressible formulations. The chemical reactions of the system are described by the GRI-

3.0 mechanism [33] and molecular diffusion is considered with the mixture-averaged transport model. The simulations employ a fully implicit scheme of second order (backward) for the time derivative and a fourth order interpolation scheme for the discretization of the convective term. All diffusive terms are discretized with an unbounded scheme of fourth-order, too. The pressure-implicit split-operator (PISO) algorithm has been used for pressure correction. An equidistant grid length of $20 \mu\text{m}$ is used for the simulations, which ensures resolving the unstretched laminar flame thickness with more than 20 cells. The time step has been set to $0.25 \mu\text{s}$, corresponding to a maximum CFL number of 0.2.

4 Results and Discussions

Figure 2 from left to right shows the spatial distributions of the calculated streamwise velocity v , temperature T , mass fraction of methane Y_{CH_4} and consumption rate of methane \dot{r}_{CH_4} for the SWQ case with $\Phi = 1$. The solid lines in Fig.2a indicate streamlines for visualization of the flow direction. The V-shaped flame front can be identified by the large gradients of T and Y_{CH_4} shown in Fig.2b and c. As the temperature at the inlet is fixed at the fresh gas temperature, the flame cannot stabilize directly at the inlet plane, which leads to a slightly lifted flame above the inlet. In Fig.2a, due to thermal expansion caused by the combustion reaction, the flow is accelerated by passing through the flame front, which leads to bending of the streamlines towards the flame-normal direction. The FWI zone can be identified from the contact point of the flame surface with the wall, where the flow velocity and reaction rate decrease to 0, and the temperature decreases to the wall temperature. In this case, the flame is pushed away from the wall, which results in a temperature and velocity boundary layer downstream. The thickness of these boundary layers are smallest in the FWI zone, which means the gradients of temperature and flow velocity are largest there.

A zoomed view of the FWI zone for the same case is shown in Fig.3a, where the flame front is indicated by the white line of $Y_{CH_4} = Y_{r_{max}}$ (see Tab.1). The grid lines of the mesh are shown to validate the grid resolution, where the reaction zone with $\dot{r}_{CH_4} < 0$ is resolved by more than 10 grid points in the near-wall region. The largest wall distance has been confirmed to be $x^+ = 0.57$, which justifies the used grid resolution. The straight lines in Fig.3a indicate flame surface normals at different distances to the wall (numbered from 1 to 4). The profiles of \dot{r}_{CH_4} along these lines are depicted in Fig.3b, with the blue curve given by the solution of a 1D unstretched flame. These profiles of \dot{r}_{CH_4} are shifted along the flame normal coordinate s , so that their peak values coincide with each other. $|\dot{r}_{CH_4}|$ for the lines 3 and 4 is larger than $|\dot{r}_{CH_4}|$ obtained from the 1D solution, which is attributable to a negative flame stretch, leading to an enhanced burning rate. The reversed effect is detected for the lines 1 and 2 closer to the wall, where the calculated $|\dot{r}_{CH_4}|$ is smaller compared with the 1D solution. The flame experiences an intensive heat loss caused by the wall, which results in flame quenching or a decreased burning rate, respectively. Although not shown here, similar behavior is obtained for the other cases with $\Phi = 0.8$ and 1.2 .

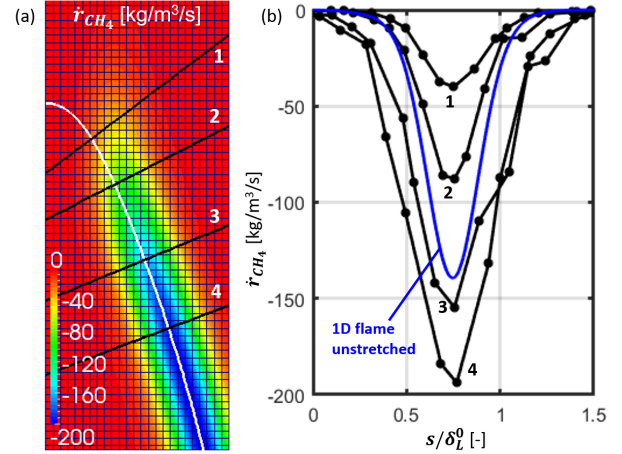


Figure 3: Contours of consumption rate of fuel \dot{r}_{CH_4} and profiles of \dot{r}_{CH_4} along flame normals for the SWQ case with $\Phi = 1$.

Figure 4 on the left shows profiles of the tangential strain component over the wall-normal distance, and in the middle for the stretch component due to curvature for the SWQ and the FF cases. $x = 0$ denotes the location of the wall in case of SWQ and the symmetry plane in case of FF. Outside the flame tip or FWI zone with $x/\delta_L^0 > 1$, Ka_s and Ka_c are similar for the SWQ and FF setups, where the flame is only weakly stretched in tangential direction with $Ka_s \approx 0.5$ and $Ka_c \approx 0$. In the near-wall region with $x/\delta_L^0 < 1$, Ka_s reverses its sign from positive in the FF case to negative in the SWQ case. Therefore, the flame is compressed lengthwise for the SWQ case, whereas it is elongated in case of FF. The mechanism is explained in Fig.4 on the right by the contours and the streamlines of the flow velocity, where the flame fronts are shown by iso-contours of the local heat release rate. The flow velocity increases towards the flame tip, which leads to a positive tangential stretch; for the SWQ case however, the flow velocity decreases while approaching the wall, which leads to a negative tangential stretch. As the flame is negatively curved at the tip of the flame, Ka_c decreases while approaching the flame tip or the wall, as shown in Fig.4 in the middle. However, the flame is curved more strongly at the tip for the FF case compared with the SWQ case, so that the normal stretch is smaller for the FF case at the flame tip.

For the SWQ flames, 3 zones with different behaviour of flame stretch can be detected, as shown in Fig.4:

- I. Ka_s decreases, whereas Ka_c or the flame curvature remains constant at $Ka_c \approx 0$. In other words, the flame is only compressed tangentially by the flow near the wall, but not curved.
- II. Ka_s continues to decrease while approaching the wall and the flame is additionally curved, leading to a negative curvature or Ka_c .
- III. Ka_s and Ka_c increase again towards zero stretch.

In the 3rd zone, the flame is so close to the wall that the reaction rate is almost zero (see Fig.5 on the right). Hence, the behavior of flame stretch during FWI is mainly attributed to the 1st and 2nd zone.

The joint effect of Ka_s and Ka_c is shown in Fig.5 on the left by the total stretch $Ka_{tot} = Ka_s + Ka_c$ over x/δ_L^0 . Although Ka_s is positive at the flame tip in the FF case (see Fig.4 on the left), Ka_{tot} is dominated by negative values near the flame tip, indicating a stronger influence of Ka_c on Ka_{tot} . While approaching the wall or flame tip at $x = 0$, Ka_{tot} starts to decrease earlier in the SWQ case compared to the FF case, which is due to the tangential compression of the flame by the wall-near flow. A reduction of Ka_s , Ka_c and Ka_{tot} with decreasing wall distance was also reported in experiments by [34] for a head-on quenching configuration with premixed methane/air flames.

Figure 5 on the right plots the normalized flame consumption speed S_L evaluated with Eq.(3) over the wall normal distance. S_L is largest at the flame tip with $x = 0$ in case of FF and smallest with $S_L = 0$ for the SWQ case, denoting a fully reversed response of S_L to flame stretch. For varying Φ , S_L from the SWQ case (blue symbol) shows a similar distribution along the wall normal distance, where S_L increases first with x for the flame branch further away from the wall and decreases again in the region close to the wall. This leads to a local maximum of S_L located at $x/\delta_L^0 \approx 1$. The result implies that the burning rate starts to decrease or the flame starts to quench towards the wall at a distance similar to the magnitude of the laminar flame thickness. In this way, the calculated quenching distances by $d_q \approx \delta_L^0$ compares well with measured data by Dreizler and Böhm [2] and Häber and Suntz [6] for a SWQ configuration of rod-stabilized premixed methane/air flames, although d_q was determined in these works based on OH*/CH* chemiluminescence and near-wall temperature profiles. Thus, the behavior of flame quenching can be reproduced or reflected by means of flame dynamics or flame speed, too.

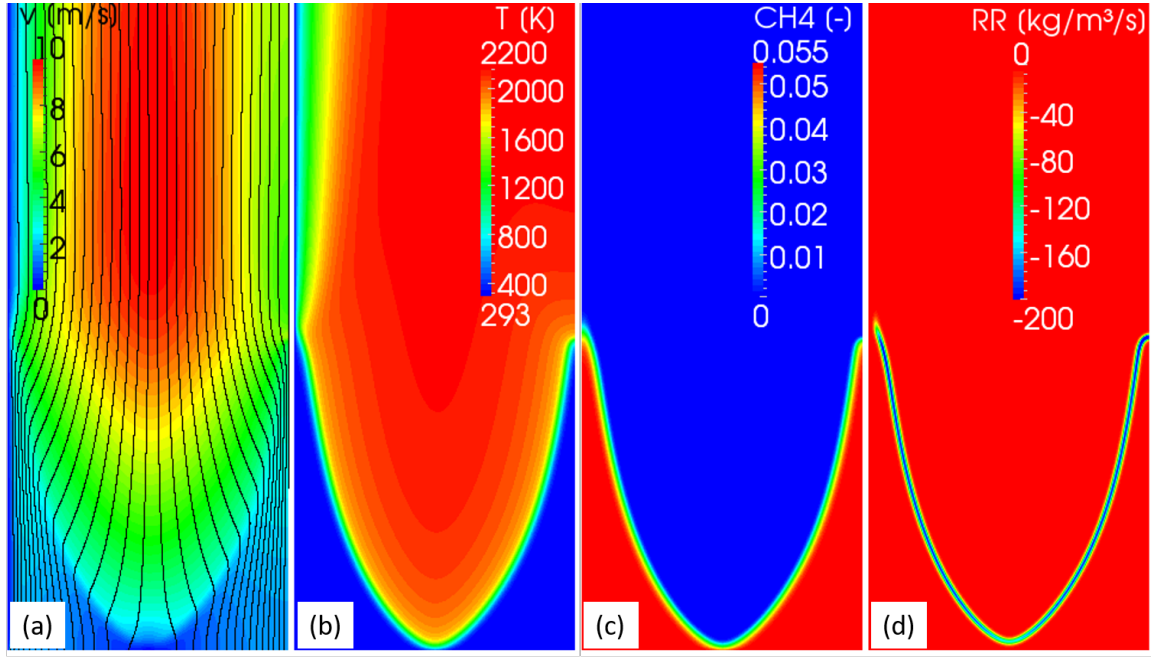


Figure 2: Calculated contours of streamwise velocity, temperature, mass fraction of CH_4 and consumption rate of CH_4 for the SWQ case with $\Phi = 1$, $p_0 = 1$ bar and $T_0 = 293$ K.

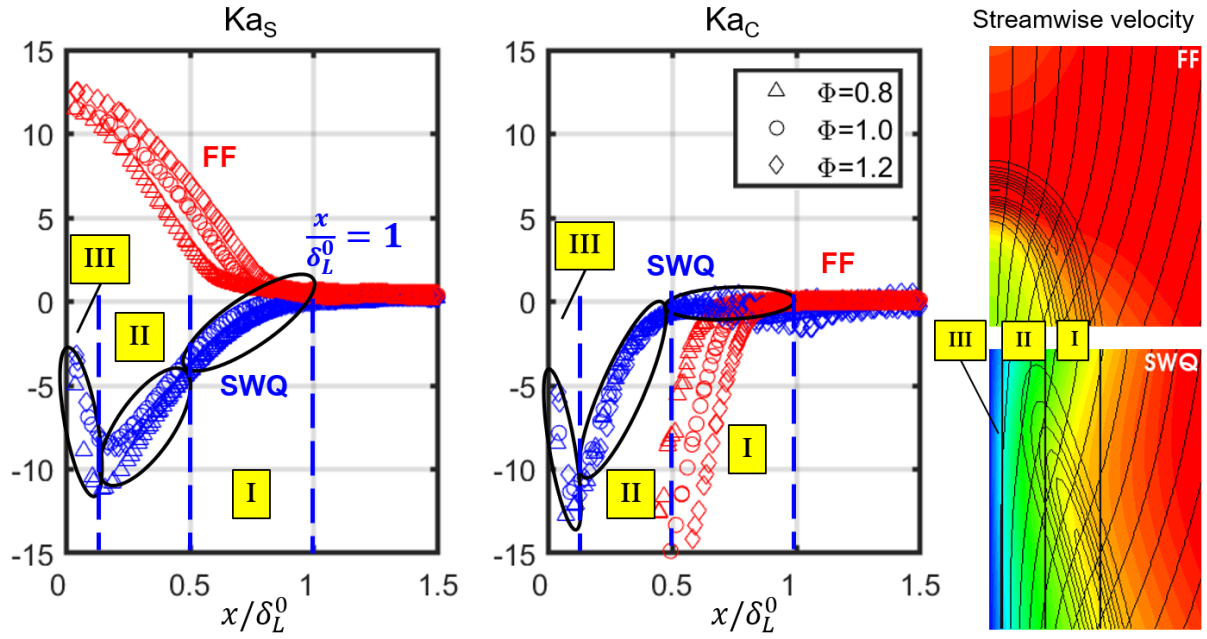


Figure 4: Comparison of tangential Ka_s (left) and normal stretch Ka_c (right) along the wall normal coordinate under SWQ (blue) and FF (red) conditions.

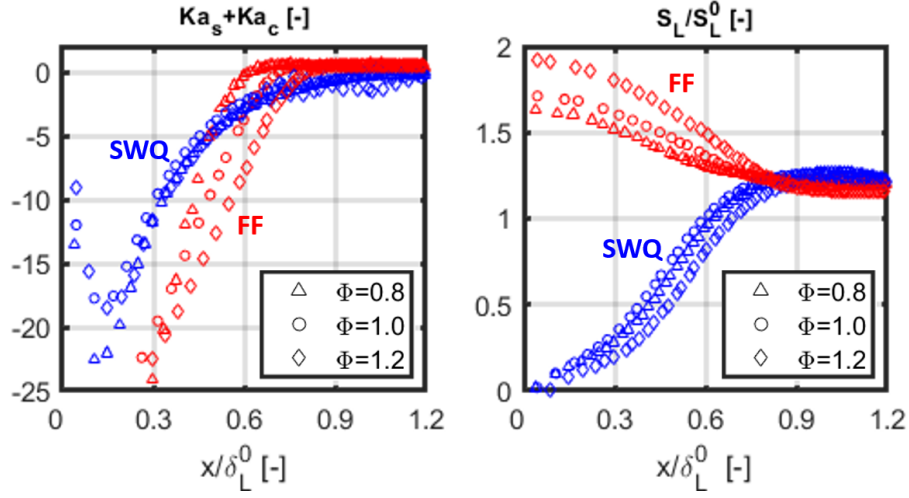


Figure 5: Comparison of Ka_{tot} (left) and S_L (right) over x at SWQ (blue) and FF (red) conditions.

In Fig.6, the flame speed is shown against the tangential stretch on the left, the normal stretch in the middle and the total stretch on the right, where the arrows indicate the direction towards the wall for SWQ or the flame tip for the FF case, respectively. A large number of data pairs of S_L vs Ka are concentrated around $Ka = 0$, which represents the ordinary flame branch without influence of the flame tip or FWI zone. As shown in Fig.4, Ka_s and Ka_c remain almost constant and close to 0 for both SWQ and FF cases in the range of $x/\delta_L^0 > 1$. The different scenarios of flame quenching close to the wall with $x/\delta_L^0 < 1$ can be detected again, similar to those for the flame stretch discussed before (see Fig.4):

- I. in the 1st stage, S_L decreases with decreasing Ka_s , while Ka_c remains almost constant.
- II. in the 2nd stage, the flame is curved and S_L decreases to 0 with both decreasing Ka_s and Ka_c

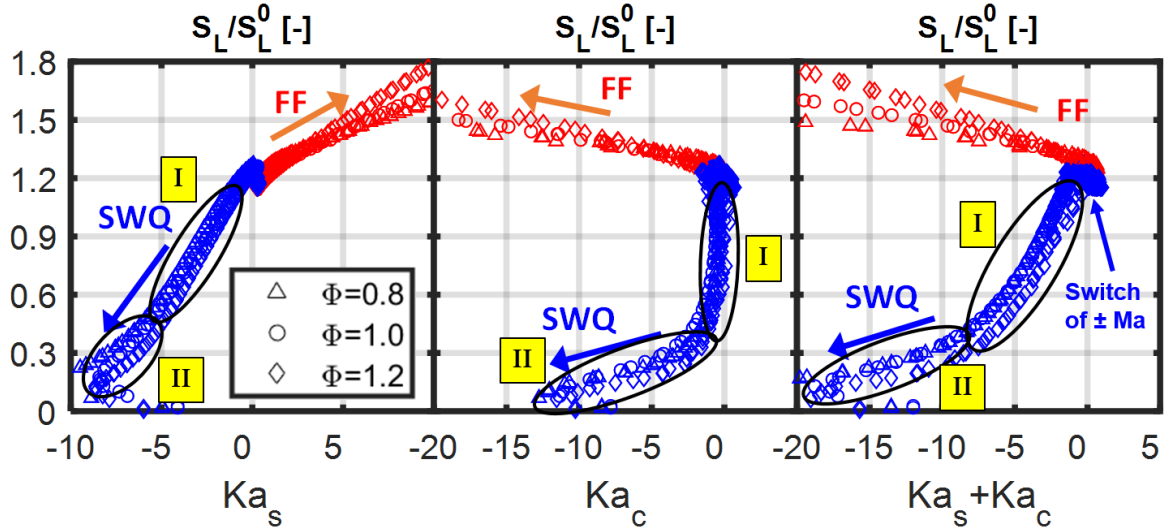


Figure 6: Correlations of flame consumption speed with flame stretch.

In Fig.6 on the left, S_L increases with Ka_s both for the FF and SWQ cases, revealing a negative Markstein number for the tangential strain $Ma_s = -\frac{dS_L/S_L^0}{dKa_s}$. The magnitude of Ma_s is however larger in the SWQ compared with the FF case, which can be detected from the slopes of the curves. Regarding the effect of flame curvature or normal stretch Ka_c on S_L shown in the middle of Fig.6, S_L decreases with decreasing Ka_c while approaching the cold wall (lower left branch) and it increases with decreasing Ka_c for the FF case (upper left branch), yielding a reversed sign of Ma_c for the SWQ and FF setups. As a consequence, Ma_{tot} considering the joint effect of Ka_s and Ka_c switches its sign from positive in the FF to negative in the SWQ case, as shown in Fig.6 on the right.

For a more quantitative analysis, the Markstein number Ma with respect to the total stretch has been evaluated from linear fitting of the relation of S_L vs. $Ka_s + Ka_c$ in Fig.7. Ma is positive for the FF case and negative for the SWQ case, indicating that Ma reverses its sign from positive to negative while approaching a cold wall, which is shown on the right of Fig.7 along with contours of the reaction rate of CH_4 . There, Ma is larger than 0 when the flame is further away from the wall, and it turns over to negative in the FWI zone. The sensitivity of S_L to Ka_s , Ka_c and Ka_{tot} are similar in magnitude for different Φ in the FWI zone, so that the FWI dynamics may be described with an almost constant $Ma \approx -0.1$ for varied stoichiometries. The result reveals a close connection between flame quenching due to interaction with a cold wall and the flame dynamics in terms of flame speed and flame stretch.

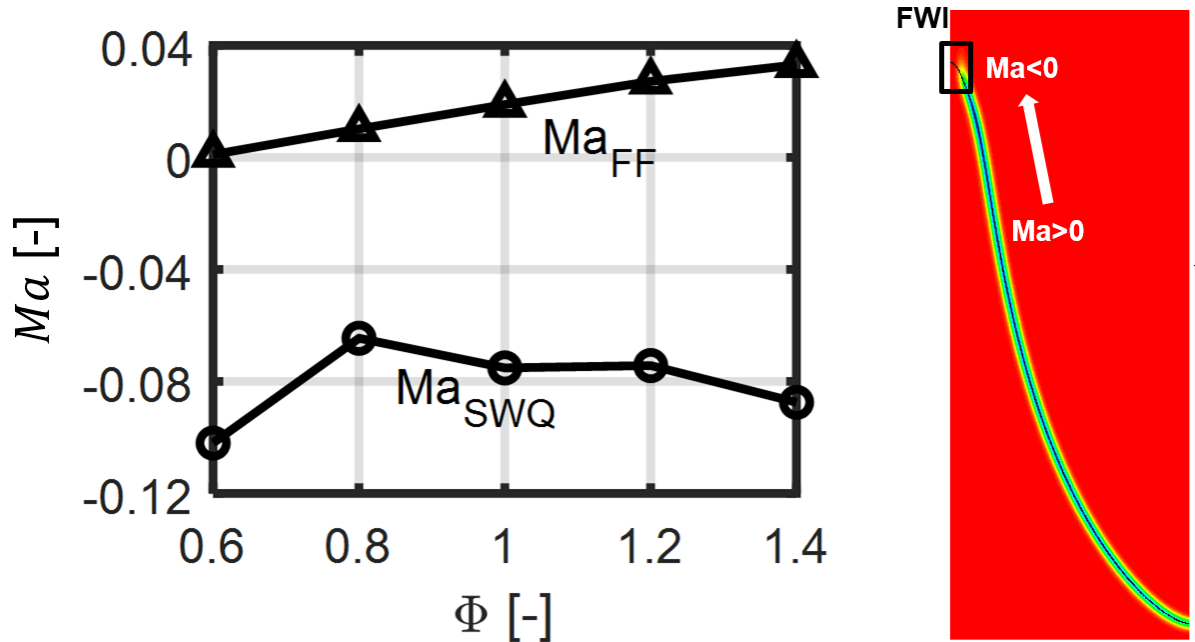


Figure 7: Calculated Markstein number for SWQ and FF flame configurations for different equivalence ratios.

5 CONCLUSIONS

Direct numerical simulations (DNS) have been performed to study flame dynamics of laminar premixed methane/air flames interacting with a cold wall, employing detailed calculations of the chemical reaction rates and transport coefficients. The flame's tangential and normal stretch Ka_s and Ka_c , as well as the

local flame consumption speed S_L have been evaluated based on the DNS results. It has been shown that the flame-wall interaction (FWI) zone is dominated by negative flame stretch due to tangential compression and negative curvature of the flame. In addition, both flame stretch and flame speed decrease while approaching the wall, which results in a reversed sign of the Markstein number Ma from positive in the free flame branch to negative in the FWI zone. The transition location of the sign of Ma has been confirmed to be located at the wall distance, that is approximately equal to the unstretched laminar flame thickness. The quenching distance d_q evaluated from this transition location of Ma with a maximum value of S_L shows a quantitatively good agreement with measured data from literature.

The obtained knowledge for flame dynamics may be used for modeling the effect of FWI. For instance, for any combustion models which use the laminar or turbulent flame speed for the description of the reaction rate, a negative Markstein number can be applied in the near-wall zone to reproduce the decrease of the burning rate. Compared with common approaches for modeling FWI with tabulated chemistry or detailed kinetics, using a Markstein analogy may be more convenient and has the potential to reduce the computational cost significantly. It should be noted that the present result with inversion of the sign of the Markstein number is valid only for flames with a positive Ma . For lean-hydrogen/air flames, Ma is naturally negative so that it will not reverse its sign, or the flame speed decreases due to the negative stretch in the FWI zone. Therefore, future work will study the influence of different fuels, wall temperatures, and turbulence on the flame dynamics during FWI.

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