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General Correlations of Iso-octane Turbulent Burning Velocities Relevant to Spark Ignition Engines

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Abstract: A better understanding of turbulent premixed flame propagation is the key for improving the efficiency of fuel consumption and reducing the emissions of spark ignition gasoline engines. In this study, we measure turbulent burning velocities (S_T) of pre-vaporized iso-octane/air mixtures over wide ranges of the equivalence ratio ($\phi = 0.9-1.25$, $Le \approx 2.94-0.93$), the root-mean-square (r.m.s.) turbulent fluctuating velocity (u' = 0-4.2 m/s), pressure p = 1-5 atm at T = 358 K and p = 0.5-3 atm at T = 373 K, where *Le* is the effective Lewis number. Results show that at any fixed *p*, *T* and *u'*, *Le* < 1 flames propagate faster than Le > 1 flames, of which the normalized iso-octane S_T/S_L data versus u'/S_L are very scattering, where S_L is the laminar burning velocity. But when the effect of *Le* is properly considered in some scaling parameters used in previous correlations, these large scattering iso-octane S_T/S_L data can be collapsed onto single curves by several modified general correlations, regardless of different ϕ , *Le*, *T*, *p*, and *u'*, showing self-similar propagation of turbulent spherical flames. The uncertainty analysis of these modified general correlations is also discussed.

Keywords: iso-octane; high-pressure turbulent burning velocity; Lewis number; general correlations; self-similar spherical flame propagation

1. Introduction

Designing a better spark ignition (SI) gasoline engine with higher thermal efficiency and lower engine emissions necessitates a thorough understanding of flame kernel initiation and its subsequent flame propagation that take place under high pressure (p), high temperature (T), high r.m.s. turbulent fluctuating velocity (u') and other complex conditions. For example, the mixture characteristics in the vicinity of an engine top dead center are frequently changed from lean to stoichiometry or even rich when the loads vary from low to high [1]. Besides, the mixture characteristics and values of u' change from cylinder to cylinder, resulting in the difference of flame initiation and subsequent flame propagation speed. Therefore, the turbulent burning velocity (S_T) has been introduced for further understanding of turbulent premixed flame propagation [2]. However, S_T data for liquid fuels (e.g., iso-octane, the major component of gasoline surrogate) under SI engine relevant conditions are rare, since most studies applied gaseous fuels (e.g., methane, propane, hydrogen at atmospheric condition [2-7] or at elevated pressure and room temperature conditions [8-16]). Whether these gaseous $S_{\rm T}$ data could be applicable to SI gasoline engines is still an open issue. To the best knowledge of the authors, the only available iso-octane S_T data which satisfied high-T, -p and -u' conditions were that reported by Lawes et al. [17] and Nguyen et al. [18] using turbulent spherical expanding flame at T = 360 K [17] and 423 K [18]. Clearly, more lean and rich iso-octane S_T data under high-T, -p, and -u'conditions are needed in order to make a detailed comparison with previous gaseous $S_{\rm T}$ data and thus address the question of similarity and difference between liquid iso-octane and gaseous data, as the first objective of this study.



Lean and rich iso-octane/air mixtures have very different Lewis number (*Le*). For instance, the lean iso-octane/air mixture at the equivalence ratio $\phi = 0.9$ has $Le \approx 2.94 >> 1$, while the rich iso-octane/air mixture at $\phi = 1.25$ has $Le \approx 0.93 < 1$. Note that *Le* is the effective Lewis number which is estimated by the ratio between thermal diffusivity and mass diffusivity with the mass diffusivity being that of the deficient reactants and the abundant inert, i.e., $Le \approx Le_{fuel}$ for lean mixture and $Le \approx Le_{oxygen}$ for the rich mixture [17]. Hence, the second objective of this paper is to seek possible general correlations of S_{T} with the consideration of the *Le* effect to represent the liquid iso-octane/air mixture at lean and rich conditions in attempt to enhance our understanding of high-pressure/temperature turbulent flame propagation that may be relevant to SI gasoline engines.

Seeking a general correlation of iso-octane S_T data is of high interest for SI engines using liquid fuels, and it is still an unresolved problem. Due to the complexity of turbulence-chemistry interactions, the different configuration of experimental facilities, and the uncertainty of measurement techniques, a general correlation to predict S_T is not always clear despite many S_T correlations available in literatures (e.g., [2–4,8–10,12–24], among others). Thus, in this work we make our efforts to analyze the scaling parameters in some selected correlations by taking the effect of *Le* into consideration based on the present measured S_T data using a single liquid fuel (iso-octane). Therefore, proper scaling parameters including the effect of *Le* and better modified general correlations using the same liquid iso-octane S_T data may be obtained. In this study, five general correlations, all including the effect of *Le*, are considered and tested by using the current measured iso-octane S_T data. Further, a mean absolute percentage error (MAPE) [23] is used to assess the accuracy of these five correlations. The first correlation is the Bradley's correlation [3]:

$$S_{\mathrm{T,}\bar{c}\,=\,0.5}/u' \sim (KLe)^{-0.3}$$
 (1)

 \bar{c} is the mean progress variable, i.e., $\bar{c} = 0$ and $\bar{c} = 1$ represent reactant and product, respectively. The stretch factor $K \approx 0.25 (u'/S_L)^2 (Re_T)^{-0.5}$ [25], where the turbulent flow Reynolds number $Re_T = u'L_I/v$, L_I is the integral length scale, and v is the kinematic viscosity of reactant. Second, by grouping the scaling parameters in correlations proposed by Kobayashi et al. [19], Chaudhuri et al. [12], and Shy et al. [16] with Le^{-1} , three modified correlations have been reported in Reference [18], respectively as shown below:

$$S_{\mathrm{T},\bar{c}} = 0.5/S_{\mathrm{L}} \sim \left[(u'/S_{\mathrm{L}})(p/p_0)Le^{-1} \right]^{0.42}$$
, where $p_0 = 1$ atm (2)

$$S_{\mathrm{T},\bar{c}} = 0.5/S_{\mathrm{L}} \sim \left(Re_{\mathrm{T},\mathrm{flame}}Le^{-1}\right)^{0.5}, \text{ where } Re_{\mathrm{T},\mathrm{flame}} = (u'\langle R \rangle)/\alpha = (u'\langle R \rangle)/(\delta_{\mathrm{L}}S_{\mathrm{L}})$$
(3)

$$S_{\mathrm{T},\bar{c}\,=\,0.5}/u' \sim \left(DaLe^{-1}\right)^{0.5}$$
, where $Da\,=\,(L_{\mathrm{I}}/u')(S_{\mathrm{L}}/\delta_{\mathrm{L}})$ (4)

 $\langle R \rangle$ is the mean turbulent spherical flame radius, δ_L is the laminar flame thickness, and $\alpha = \delta_L S_L$ is the thermal diffusivity. In Equation (3), the commonly-used integral length scale of turbulence is replaced by the average flame radius $\langle R \rangle$ and the kinematic viscosity (ν) is replaced by the thermal diffusivity ($\alpha \approx \delta_L S_L$), as proposed by Chaudhuri et al. [12,13]. Finally, the fifth correlation is that reported by Ritzinger [20]:

$$S_{\mathrm{T},\bar{c}\,=\,0.5}/S_{\mathrm{L}}\,=\,1+(0.46/Le^{0.85})(u'/S_{\mathrm{L}})^{0.5Le^{0.25}}(Re_{\mathrm{T}})^{0.25}(p/0.1\mathrm{MPa})^{0.1}$$
(5)

Equation (5) has been tested against methane/air S_T data at high-pressure gas engine conditions [26], while Equations (1)–(4) have not yet done so. To see how well these five correlations, Equations (1)–(5), could be applicable to SI engines for S_T prediction, an uncertainty comparison is performed by accessing the accuracy of these correlations based on the MAPE value which is calculated using the following equation [23].

$$MAPE(\%) = \frac{\sum_{i=1}^{n} \left| \frac{x_{exp,i} - x_{corr,i}}{x_{exp,i}} \right|}{n} \times 100$$
(6)

 $x_{\exp,i}$ is the experimental value, $x_{\text{corr},i}$ is the value predicted by the correlation, and n is the number of the data group under examination.

In short, this work measures S_T data of lean, stoichiometric, and rich iso-octane/air mixtures with Le > 1 and Le < 1 using a large dual-chamber, constant pressure/temperature, fan-stirred cruciform explosion facility, capable of generating near-isotropic turbulence. It will be shown in due course that all the scattering S_T data of iso-octane/air mixtures can be collapsed onto single curves using Equations (1)–(5). Further, the current iso-octane data are also compared with previous data proposed by Lawes et al. [17]. Finally, the uncertainty of these correlations based on the MAPE is discussed.

2. Experimental Method

Experiments are conducted in a large dual-chamber, constant-temperature, constant-pressure, fan-stirred cruciform explosion facility for S_L and S_T measurements of expanding spherical iso-octane/air flames. The reader is directed to [18,21] and references therein for detail treatment on the facility and its associated turbulence properties. For completeness, a simplified sketch of the 3D cruciform burner resided in a large pressure vessel with optical accesses is added, as shown in Figure 1, alongside the Schlieren imaging arrangement. Below are the descriptions of the experimental procedures.

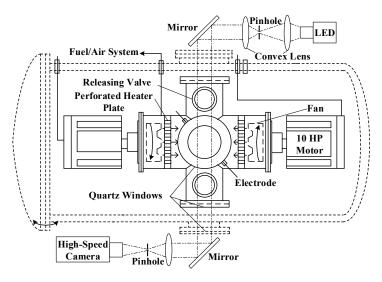


Figure 1. The high-pressure/temperature, double-chamber explosion facility for premixed turbulent combustion studies [27].

Before a run, we first vacuum the heated 3D cruciform burner (Figure 1) and then inject the appropriate mole fractions of pre-vaporized iso-octane and air into the burner to the desired initial pressure and temperature conditions. Two counter-rotating fans at a frequency of 30 Hz are turned on to well mix the iso-octane/air mixture for 4 minutes and reach a uniform temperature distribution with less than 1 °C variation in the domain of experimentation. The temperature uniformity is achieved by using effective turbulent heat convection through a pair of heated perforated plates (please see [18,21] for details). The domain of experimentation is set at $0.17 \le \langle R \rangle / R_{min} \le 0.30$ to avoid the ignition influence at the early stage of kernel development and the wall effects at the later stage of flame propagation, where the minimum wall confinement radius of the 3D cruciform bomb R_{min} is about 150 mm. For laminar flame speed measurements, we ignite the mixtures shortly (about 10 s) after the fan turn-off to allow for the decay of turbulence to quiescence, while keeping a uniform temperature distribution of the average flame radii $\langle R \rangle (t) = [A(t)/\pi]^{0.5}$, where A(t) is the area enclosed by the laminar/turbulent flame front. The flame speed d $\langle R \rangle / dt$ is obtained by the central differentiation of $\langle R \rangle$ vs. *t* data.

A typical refined gasoline consists of hundreds of different components that may vary batch by batch depending on the source of crude oil and the refinery processes, making the benchmark of S_T for gasoline difficult to measure accurately. For simplicity, iso-octane, a major surrogate component for gasoline, is used to gain insight and understanding of the underlying physical mechanism of turbulent premixed flame propagation. As such, this work measures values of S_T for iso-octane/air mixtures as a function of ϕ , u', p, T, and Le. Table 1 lists laminar properties of the iso-octane/air mixture at various conditions including T, ϕ , Le, p and corresponding laminar burning velocities (S_L). It should be noted that the selected mixtures, i.e., $\phi = 0.9$ & 1.25 at T = 358K at p = 1 atm and $\phi = 1.0$ & 1.2 at T = 373 K at p = 1 atm, have closely matched values of S_L (see Table 1) but with different Le (Le > 1 and Le < 1). As such, the effect of Le on S_T of iso-octane/air mixtures can be studied.

T (K)	ϕ	Le	<i>p</i> (atm)	$S_{\rm L}$ (cm/s)
358	0.9/1.0/1.25	2.94/1.43/0.93	1, 3, 5	41.0, 26.2, 24.8/45.0, 31.2, 27.6/40.2, 32.9, 28.2
373	1.0/1.2	1.43/ 0.93	0.5, 1, 3	49.0, 44.9, 31.0/47.0, 44.6, 36.0
423 ¹	1.0	1.43	1, 3, 5	55.0, 42.0, 36.0

Table 1. Conditions and properties of iso-octane/air mixtures.

3. High-Speed Schlieren Imaging of Turbulent Spherical Flame and Their S_T Determination

Figure 2 shows the effects of *Le* and *p* on the emergence of small scale structures and the increase of average flame propagation rate at fixed u' with increasing p for both (a) laminar and (b) turbulent cases, where all images have the same $\langle R \rangle \approx 35$ mm. At any fixed *p*, the elapsed instants after ignition show that Le < 1 flames (2nd and 4th columns) propagate much faster than Le > 1 flames (1st and 3rd columns). $S_{\rm L}$ decreases with increasing p even though the flame at 5 atm appears more wrinkling due to the emergence of cellular structures allover the flame surface especially for the case at 5 atm and $Le \approx 0.93 < 1$ (see the 4th column of the first row images in Figure 2a). For the turbulent case at constant u' = 1.4 m/s (Figure 2b), the turbulent flame propagates faster with increasing p under both Le > 1 and Le < 1 conditions, where the turbulent wrinkled flame at 5 atm are full of very small scale structures. These fine structures are mainly due to the reduction of the thickness of the laminar flamelet at high pressure that promotes the hydrodynamic instability and thus increases the wrinkled flame front surface, but they contribute only a small part for the increase of S_{T} . This is because when the flow turbulent Reynolds number ($Re_{T,flow} = u'L_I/\nu$) is kept constant for gaseous fuels (e.g., methane and syngas), $S_{\rm T}$ actually decreases with increasing pressure, similar to $S_{\rm L}$, showing a global response of burning velocities to the increase of pressure (please see Reference [11] for detail information). As such, the increase of S_T with increasing p at fixed u' is probably mainly due to the increase of $Re_{T,flow}$ at elevated pressure because $\nu \sim \rho^{-1} \sim p^{-1}$.

Figure 3a shows the averaged flame radii of iso-octane/air turbulent premixed flames versus time for *Le* > 1 (ϕ = 0.9 and 1.0) and *Le* < 1 (ϕ = 1.25) at the same *p* = 3 atm, *u'* = 1.4 m/s, and *T* = 358 K conditions. Flame speeds, $d\langle R \rangle(t)/dt$ and S_F or S_T , can be estimated from the raw data of $\langle R \rangle(t)$ within the range of 25 mm $\leq \langle R \rangle \leq 45$ mm. $d\langle R \rangle/dt$ is directly taking the time differentiation on $\langle R \rangle(t)$, while S_F is determined as the slope of the best linear-fit of $\langle R \rangle(t)$ within 25 mm $\leq \langle R \rangle \leq 45$ mm. Within this experimentation domain, S_F is just the average value of the linear increase $d\langle R \rangle/dt$ data ($S_F = \overline{dR/dt}$), as substantiated in Figure 3b. Note that these iso-octane data of $\langle R \rangle$ vs. *t* and $d\langle R \rangle/dt$ vs. *t* or $\langle R \rangle$ in Figure 3 are very similar to previous data obtained from the Leeds fan-stirred explosion bomb [17,28]. It has been shown in Reference [11] that the modified S_F using the density correction and Bradley's mean progress variable \overline{c} converting factor for Schlieren spherical flames at $\overline{c} = 0.5$ [25], i.e., $S_{T,\overline{c}=0.5} \approx (\rho_b/\rho_u)S_F(\langle R \rangle_{\overline{c}=0.1}/\langle R \rangle_{\overline{c}=0.5})^2 = (\rho_b/\rho_u)\overline{dR/dt}(\langle R \rangle_{\overline{c}=0.1}/\langle R \rangle_{\overline{c}=0.5})^2$, show good agreements between Bunsen-type and spherical flames, suggesting that S_T determined at the flame surface contour of $\overline{c} = 0.5$ may be a better representative of itself regardless of different flame geometries. The subscripts b and u indicate burned and unburned gases. For Schlieren expanding spherical turbulent premixed flames, $\langle R \rangle_{\bar{c}=0.1}/\langle R \rangle_{\bar{c}=0.5} \approx 1.4$ [22,28]. In this work, all measured $S_{\rm F}$ data are converted to turbulent burning velocities at $\bar{c} = 0.5$ ($S_{\rm T,\bar{c}} = 0.5$).

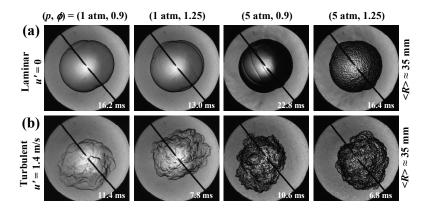


Figure 2. High-speed Schlieren images of lean and rich iso-octane/air expanding spherical premixed flames, i.e., $\phi = 0.9$ with $Le \approx 2.94 > 1$ and $\phi = 1.25$ with $Le \approx 0.93 < 1$ at 358 K, showing effects of Lewis number and pressure on the emergence of small scale structures and the increase in average flame propagation rate at fixed u' = 1.4 m/s where the flow turbulent Reynolds number ($Re_{T,flow} = u'L_I/v$) increases with pressure because $v \sim \rho^{-1} \sim p^{-1}$. All images in both (**a**) laminar and (**b**) turbulent cases have the same $\langle R \rangle \approx 35$ mm in the same view field of 110 mm × 110 mm for comparison.

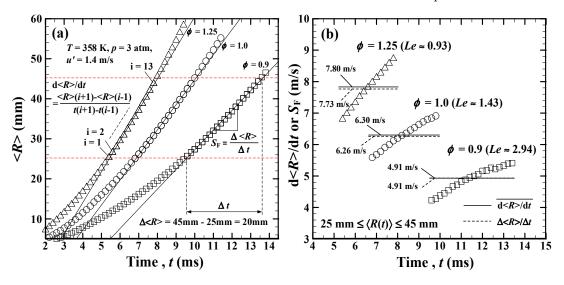


Figure 3. (a) Typical averaged flame radii of turbulent iso-octane/air premixed flames as a function of time at three different values of ϕ with Le > 1 ($\phi = 0.9$ and 1.0) and Le < 1 ($\phi = 1.25$) under the same p = 3 atm, u' = 1.4 m/s and T = 358 K conditions. (b) Two turbulent flame speeds versus time, $d\langle R \rangle/dt$ and S_F (the slope of $\langle R \rangle(t)$) as indicated in (a), which are almost equal to each other within $25 \text{ mm} \le \langle R(t) \rangle \le 45 \text{ mm}$.

4. Results and Discussion

Figure 4a presents laminar and turbulent burning velocities of iso-octane/air mixtures as a function of *p* for three different values of ϕ with different *Le* varying from 0.93 to 2.94. As increase *p*, values of S_L decrease in a minus power law manner of the form $S_L \sim p^{-n_L}$, where the exponent n_L increases with increasing *Le* from $n_L = 0.14$ at $Le \approx 0.93$ and $n_L = 0.31$ at $Le \approx 1.43$ to $n_L = 0.33$ at $Le \approx 2.94$. This suggests that Le < 1 flames are less sensitive to pressure elevation than Le > 1 flames. On the other hand, values of $S_{T,\bar{c}} = 0.5$ increase with increasing *p* in a positive power law form, i.e., $S_T \sim p^{+n_T}$, where n_T decreases with increasing *Le* from $n_T = 0.14$ at $Le \approx 0.93$ and $n_T = 0.1$ at $Le \approx 1.43$ to $n_T = 0.03$ at $Le \approx 2.94$. The latter

case for the lean iso-octane/air mixture at $\phi = 0.9$ shows that $S_T \sim p^{0.03}$ when u' = 1.4 m/s. In other words, at large $Le \approx 2.94 >> 1$ and u' = 1.4 m/s, the fine structures allover the flame front surface (Figure 2b) induced by the hydrodynamic instability at high pressure as well as the increase of $Re_{T,flow}$ due to the decrease of kinematic viscosity at elevated pressure only have a rather small influence on S_T . The reason for this tiny impact on S_T due to the increase of p when u' = 1.4 m/s is not yet clear, which deserves further studies. Figure 4b presents laminar and turbulent burning velocities as a function of temperature at four different values of u' for the stoichiometric iso-octane/air mixture at 1 atm. In general, values of S_L and $S_{T,\bar{c}} = 0.5$ increase with increasing T. $S_{T,\bar{c}} = 0.5 \sim (T/T_0)^{m_T}$, where $T_0 = 298$ K. The exponent constant $m_T = 0.07$ is small at small u' = 1.4 m/s, while the value of m_T increases to 1.07 when u' = 4.2 m/s. This suggests that $S_{T,\bar{c}} = 0.5$ is not-so-sensitive to the increase of temperature at u' = 1.4 m/s in line with previous studies by Lipatnikov et al. [29] and Fogla et al. [30] who studied the effect of density ratio on the turbulent flame speed. Note that changing the temperature is equivalent to changing the density ratio between the burned and unburned gases that has insignificant effect on S_T at low-to-moderate turbulent intensities [29,30]. However, such temperature enhancement sensitivity increases with increasing u'.

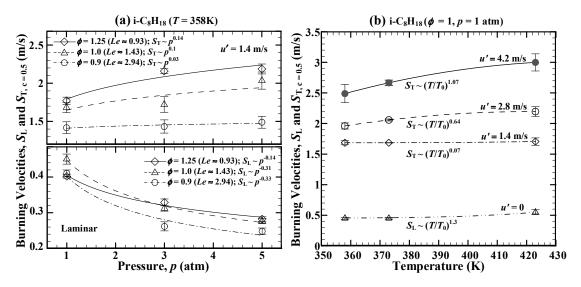


Figure 4. (a) Laminar burning velocities and turbulent burning velocities at $\bar{c} = 0.5$ of the iso-octane/air mixtures at three different ϕ with Le < 1 and Le > 1 as a function of pressure, where T = 358 K and u' = 1.4 m/s. (b) Effect of temperature on S_L and S_T at $\bar{c} = 0.5$ having four different u' varying from 0 to 4.2 m/s.

The current S_T data obtained at 373 K and 5 atm may not be sufficiently high, since much higher temperatures and pressures are typically observed in SI engines. However, the available S_T data of pre-vaporized iso-octane fuel (the major component of gasoline surrogate) even at 373 K and 5 atm are still very rare, which may be important as the first order approximation towards the practical SI engine conditions. Moreover, these experimental results are useful because they can be used to improve current models for the predication of turbulent flame speeds of pre-vaporized iso-octane fuel at high *p* and high *T*.

All present $S_{T,\bar{c}=0.5}/S_L$ of iso-octane/air mixtures at $\phi = 0.9-1.25$ at T = 358 K and/or 373 K together with previous data at $\phi = 1.0$ and at T = 423 K [18] are grouped onto two data sets having Le < 1 and Le > 1 and plotted against turbulent intensities (u'/S_L) , as shown in Figure 5. Results show that Le < 1 flames with $S_{T,\bar{c}=0.5}/S_L = 2.54(u'/S_L)^{0.53}$ having a poor goodness $R^2 = 0.58$ propagates faster than Le > 1 flames with $S_{T,\bar{c}=0.5}/S_L = 1.76(u'/S_L)^{0.64}$ and $R^2 = 0.65$, revealing large data scattering. Note that the increase of $S_{T,\bar{c}=0.5}/S_L$ with u'/S_L is not linear. There is a bending effect of $S_{T,\bar{c}=0.5}/S_L$ at higher u'/S_L for both Le < 1 and Le > 1 flames. Such bending effect has been discussed by many studies (e.g., [2,4,11,19,22,25] among others), which is mainly attributed to the turbulent flame stretch effect. At higher u'/S_L , the two fitting curves are getting closer to each other, suggesting that the effect of *Le* is getting weaker in intense turbulence. Nevertheless, it is obvious that the *Le* effect should be considered in any scaling correlations.

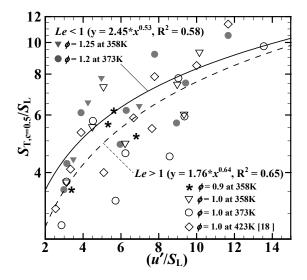


Figure 5. Normalized turbulent burning velocities at $\overline{c} = 0.5$ plotted against turbulent intensities (*u*'/SL) at different pressure p = 0.5–3 atm at 373 K and p = 1–5 atm at 358 K and/or 423 K.

First, we discuss the correlation of Equation (1) proposed by Bradley et al. [3] which is used to fit all data from Figure 5. The result is presented in Figure 6, suggesting that the Bradley's correlation can be used to fit all scattering S_T data for Le > 1 and Le < 1 flames with a better goodness of $R^2 = 0.77$ than that in Figure 5. In [3], Bradley et al. noted that Equation (1) was limited within the range of *KLe* between 0.01 and 0.63 and, if outside that range, the logarithmic plots of u_t/u'_k (or S_T/u') against *KLe* has large data scattering (please see Reference [3]). Also plotted in Figure 6 for comparison is the original version of Equation (1): $S_{T,\bar{c}=0.5}/u' = 0.88$ (*KLe*)^{-0.3} based on various gaseous mixtures [3]. The present pre-factor 0.55 is 62.5% of the previous pre-factor of 0.88 as in [3], probably due to different liquid and gaseous fuels applied.

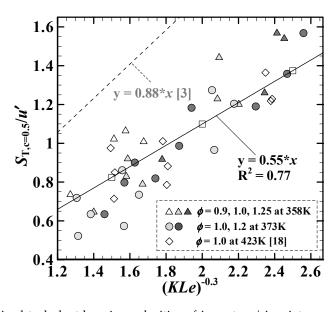


Figure 6. Normalized turbulent burning velocities of iso-octane/air mixtures using the Bradley's correlation [3]: $S_{T,\bar{c}=0.5}/u' = 0.55(KLe)^{-0.3}$. The dashed line was the original version of previous data obtained by Bradley et al. [3], where $S_{T,\bar{c}=0.5}/u' = 0.88(KLe)^{-0.3}$.

Figure 7a–c show the normalized $S_{T,\bar{c}=0.5}$ data by their corresponding values of S_L and/or u' for iso-octane/air mixtures with Le > 1 and Le < 1 using the earlier versions of Equations (2)–(4) without the consideration of *Le*. These three earlier versions of scaling parameters are respectively: (a) $(u'/S_L)(p/p_0)$ proposed by Kobayashi et al. [8,19], (b) $(Re_{T,\text{flame}})^{0.5}$ by Chaudhuri et al. [12], and (c) $(Da)^{0.5}$ by Shy et al. [10,11,16]. The results clearly show two distinct data sets with Le < 1 flames (dark grey symbols) propagating faster than Le > 1 flames (white and light grey symbols) for all three correlations as shown in Figure 7a–c. Clearly, the effect of *Le* must be taken into the consideration to obtain possible general correlations. When the aforesaid three scaling parameters are grouped with Le^{-1} , all present iso-octane $S_{T,\bar{c}=0.5}$ data can be collapsed onto single curves, as shown in Figure 7d–f, suggesting that the propagation of iso-octane turbulent premixed spherical flames is self-similar, regardless of different Le, ϕ , p, and u' applied. The goodness between these three modified correlations varies from $\mathbb{R}^2 \approx 0.89$ using Equation (2) and $\mathbb{R}^2 \approx 0.84$ using Equation (3) to $\mathbb{R}^2 \approx 0.7$ using Equation (3), all based on the same iso-octane/air $S_{T,\bar{c}=0.5}$ data.

The final $S_{T,\bar{c}=0.5}/S_L$ correlation to be discussed is Equation (5) proposed by Ritzinger [20] based on a complicated empirical relation in the form of $S_{T,\bar{c}=0.5}/S_L = 1 + (0.46/Le^{0.85})(u'/S_L)^{0.5Le^{0.25}}(Re_T)^{0.25}(p/0.1\text{MPa})^{0.1}$. Results are presented in Figure 8, where all the present iso-octane $S_{T,\bar{c}=0.5}$ data can be well collapsed onto a single curve with R² = 0.89. In general, Equations (1)–(5) can be used as general correlations for both present liquid iso-octane/air and preveous gaseous turbulent burning velocities with reasonably good R², regardless of different fuel, ϕ , *Le*, *u'*, *T* and *p* applied.

Here we apply Equation (6) to estimate MAPE and the results are shown in Figure 9. MAPE is a better statistical variable than the goodness \mathbb{R}^2 , because the latter depends strongly on the available data volume to estimate the uncertainty of these five general correlations (Equations (1)–(5)). Note that the lower the MAPE is, the more accuracy of the correlation is. All MAPE percentages for the aforesaid five general correlations are less than 13%, showing that these general correlations are reasonably good. Among them, the best is Equation (5) proposed by Ritzinger [20] with a lowest MAPE of 8.5% and the second best is Equation (2) proposed by Kobayashi et al. [8,19] with a MAPE of 9.6%. The other three general correlations (Equations (1), (3), (4)) have roughly the same MAPE (11.9%–12.9%). It should be noted that Equation (2) is probably the most convenient one to use, because it does not require any length scales of turbulence and flame.

We now discuss the iso-octane comparison between the present data and previous data obtained by Lawes et al. [17] using Equation (4). The comparison results are presented in Figure 10. The main reason that we use Equation (4) for comparison is because the Damköhler number (*Da*) to the one half power has been extensively used by Peters [31]. Besides, $Da = (L_I/u')(S_L/\delta_L)$ includes the important integral length scale of turbulence (*L*_I), but Equation (2) and Equation (3) do not include *L*_I. A good agreement between present and previous data is found. Both present and previous iso-octane $S_{T,\bar{c}=0.5}/u' = 0.082 (DaLe^{-1})^{0.5}$ with a reasonable goodness of $\mathbb{R}^2 = 0.78$.



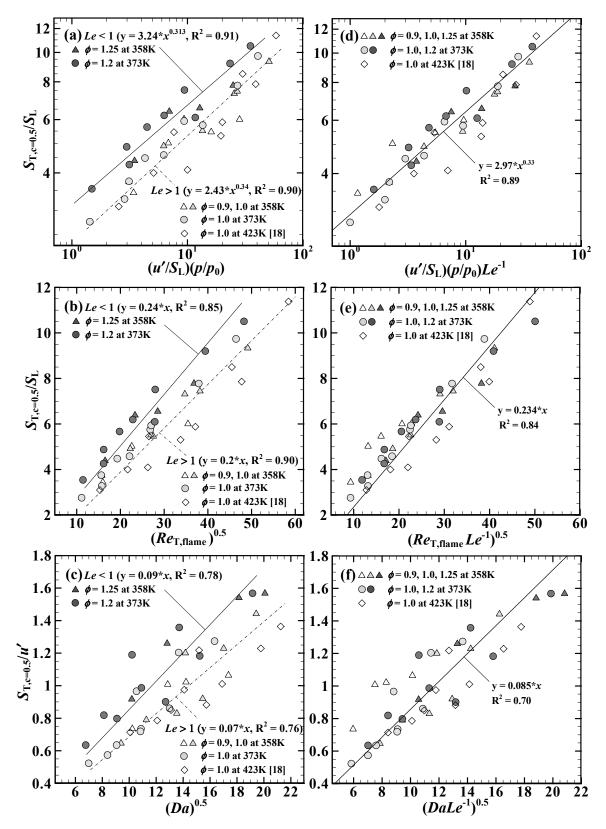


Figure 7. Left column: Normalized turbulent burning velocities for iso-octane/air mixtures plotted against three different scaling parameters: (**a**) $(u'/S_L)(p/p_0)$ [8,19]; (**b**) $(Re_{T,\text{flame}})^{0.5}$ [12]; (**c**) $(Da)^{0.5}$ [10, 11,16], all showing that Le < 1 flames propagate faster than Le > 1 flames having considerable data scattering. Right column: Same data from (**a**)–(**c**), but plotted against three modified scaling parameters grouped with Le^{-1} : (**d**) $(u'/S_L)(p/p_0)Le^{-1}$ (Equation (2)); (**e**) $(Re_{T,\text{flame}}Le^{-1})^{0.5}$ (Equation (3)); (**f**) $(DaLe^{-1})^{0.5}$ (Equation (4)), in which both Le < 1 and Le > 1 data sets are collapsed onto single curves.

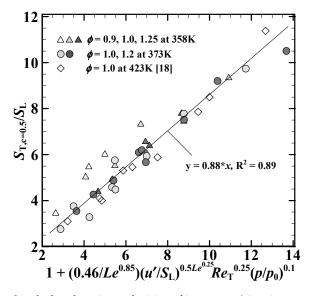


Figure 8. Normalized turbulent burning velocities of iso-octane/air mixtures using Equation (5).

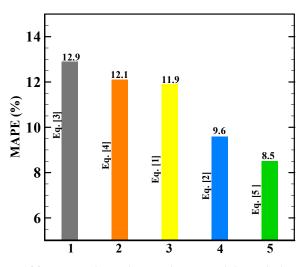


Figure 9. The accuracy of five general correlations determined through the use of mean absolute.

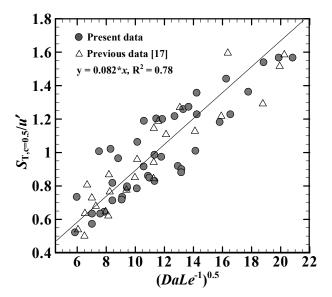


Figure 10. Comparison of present data with previous data obtained by Lawes et al. [17] using Equation (4) for the same iso-octane/air mixtures.

5. Conclusions

Turbulent burning velocities of lean, stoichiometric and rich iso-octane/air mixtures having Le < 1 and Le > 1 are measured under elevated pressures and temperatures relevant to SI gasoline engine or micro gas turbine conditions. These measurements reveal the following points:

- (1) Turbulent burning velocities increase with increasing pressure and temperature at any given r.m.s turbulent fluctuating velocities. However, the increase of turbulent burning velocities is quite small at modest u' = 1.4 m/s, which is much less sensitive with the increase of pressure and temperature as compared to that at higher values of u'.
- (2) The bending effect of $S_{T,\bar{c}=0.5}/S_L$ vs. u'/S_L curves is observed at higher u'/S_L .
- (3) *Le* < 1 flames (the rich iso-octane/air mixture) propagate faster than *Le* > 1 flames (lean and/or stoichiometric iso-octane/air mixtures) at any given conditions.
- (4) All present iso-octane $S_{T,c=0.5}$ data with Le < 1 and Le > 1 can be collapsed onto single curves regardless of different *T*, *p*, *u'*, ϕ , and *Le* used, which can be represented by five general correlations (Equations (1)–(5)), suggesting that turbulent premixed spherical flames have a self-similar propagation nature. These five general correlations have reasonable good accuracy due to their low values of MAPE (MAPE < 13%).

The fundamental understanding of flame initiation and propagation of lean iso-octane/air mixtures through the interactions of centrally-ignited outwardly-propagating premixed flame and intense turbulence at high pressure and high temperature is of importance for the future development of high-thermal-efficiency SI gasoline engines operated at fuel lean conditions. This is because premixed lean-burn technology can increase the mixture specific heat ratio and decrease the net heat loss to the engine cylinder through low temperature combustion, resulting in higher thermal efficiency and lower NO_x emissions [32]. As pointed out by Nakata et al. [32], super lean-burn combustion with a very strong tumble flow (very high u') in a long stroke cylinder is crucial to further enhance the maximum engine thermal efficiency up to 50% [33]. Hence, for the future research direction, the $S_{\rm T}$ information of iso-octane (the major component of gasoline surrogate) at extreme conditions, e.g., leaner mixtures ($\phi = 0.6$ –0.8), higher u' > 4.2 m/s and higher p > 5 atm, are needed.

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