



Analysis and optimisation of mixing in binary droplet collisions

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The collision of binary droplets plays a key role in several industrial, chemical and biological processes. In these processes, the quality of the desired outcome is strongly dependent on the mixing of the liquid droplets as they collide in mid-air. In this work, multiphase direct numerical simulations based on the volume-of-fluid method have been used to investigate the process of mixing and analyse the effects of parameters such as injection velocity, timing and collision angles. The evolution of mixing due to convection and irreversible diffusive processes has been quantified by means of the segregation parameter. To synthesise the outcome of a collision, the impact parameter has been redefined to account for the collision of non-spherical droplets. It has been found that the optimal mixing does not occur for symmetric head-on collisions, but rather at moderately asymmetrical configurations. This behaviour has been explained by analysing the velocity gradient tensor. It has been demonstrated that by breaking the symmetry, the local topology of the flow is altered and the resulting convective flows increase the contact area between the liquids, thereby augmenting the mixing process. However, it was also observed that lateral misalignment transforms the initial kinetic energy into the spinning of the merged droplets, thus preventing an enhanced mixing.

Key words: mixing enhancement, coupled diffusion and flow, drops

1. Introduction

Binary droplet collisions have been a research topic for more than a century and continue to be an active research topic due to their complexity and many applications. In its simplest

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form, the problem consists of two liquid droplets approaching each other at a constant velocity, headed for a collision. Typically, these droplets are immersed in a fluid, such as air, with a smaller viscosity than the droplets – for example, aerosol fuel droplets merging or splashing in combustion chambers (Jiang, Umemura & Law 1992; Chen 2007) in which the ignition and propagation of the flames depend on the atomisation of the fuel (Ozel-Erol *et al.* 2018). Another example concerns the drying of suspensions such as milk (Finotello *et al.* 2017), paint pigments or pharmaceutical powders (Mujumdar 2014). In these applications, large droplets may not dry out completely; on the contrary, if the droplets are too small, undesirable dust particles are produced. The main objective in the study of binary droplet collisions can be summarised as predicting the outcome of the collision for a given set of initial conditions (Ashgriz & Poo 1990). In other words, identifying the conditions under which the droplets coalesce, bounce or splash, disintegrating into several small droplets after the collision, remains important. Early studies in binary droplet collisions can be dated back to Rayleigh (1945), who reported that droplets may coalesce and break up into several smaller droplets or bounce apart after a collision. Most of the early experiments were conducted for water droplets (Adam, Lindblad & Hendricks 1968; Brazier-Smith *et al.* 1972) primarily to model atmospheric phenomena such as cloud formation. Those experiments suggest that the three main parameters that determine the outcome of the collision are the Weber number $We := \rho(d_1 + d_2)u_r^2/2\sigma$ and the impact parameter $B := 2b/(d_1 + d_2)$, where ρ is the density of the liquids, d_i ($i = 1, 2$) is the diameter of the droplets, σ is the surface tension, b is the separation distance between the centres of mass of the droplets and u_r is the magnitude of the relative velocity. The Weber number quantifies the relative kinetic energy in proportion to the interfacial energy, and the impact parameter is a measure of the off-centredness of the collision. It was found that for small We and B , the droplets undergo a slow coalescence; increasing either We or B results in the bouncing of the droplets due a layer of air trapped between the droplets. Coalescence is recovered by further increasing the relative velocity of the droplets (Orme 1997; Qian & Law 1997); however, a head-on separation or fragmentation may follow, depending on the impact parameter. This is usually summarised by a phase diagram in the We – B space, also known as the collision map, which indicates the different outcomes that are expected for a given pair of We and B (Al-Dirawi & Bayly 2019). However, the collision map is not unique and other parameters, such as the viscosity of the liquids and surrounding phase and its pressure, also determine the outcome of the collision (Chesters 1991). For example, Willis & Orme (2000) conducted experiments in a vacuum chamber, thus revealing the effect of the outer phase. Further experimental studies have focused on the interaction of non-Newtonian (Finotello *et al.* 2018) or immiscible liquids (Planchette, Lorenceau & Brenn 2010).

Theoretical and numerical studies have also addressed the main problem of the collision of binary droplets. Gopinath & Koch (2002) devised a model for a head-on collision identifying the boundaries between the regions in the collision map. Later, Roisman (2004, 2009) proposed a model giving an approximation of the flow field at early stages of a head-on collision and successfully predicted the thickness of the liquid lamella and the outer rim that is formed after coalescence. In the meantime, lattice Boltzmann (Inamuro, Tajima & Ogino 2004) and level-set (Pan & Kazuhiko 2005) simulations were being developed. Kuan, Pan & Shyy (2014) conducted a detailed study of droplet splashing using the interface-tracking method combined with adaptive mesh refinement. The applications that are based on the precise outcome of the collision of droplets are technologies that include drop-on-demand manufacturing such as bioprinting processes. The reactive jet impingement (ReJI) technology (da Conceicao Ribeiro *et al.* 2018), for example, consists

of a stream of cell-laden droplets of a gel precursor and a cross-linker that are injected and collide in mid-air to then mix and react to build the designed living tissue. This novel technique for bioprinting shows a significant improvement over other techniques in terms of cell viability and printing speed. In ReJI, as well as several adaptive manufacturing techniques, the quality of the product and throughput depends on the speed of injection and size of the droplets (Stringer & Derby 2009). This imposes a limit on the Weber number for droplet stacking (Son *et al.* 2008) and mid-air collisions (Barker *et al.* 2022). Furthermore, the complexity in the modelling of additive manufacturing processes increases from the idealised droplet collision. For example, when colliding, the morphology of the droplets does not relax to a spherical shape, requiring a redefinition of the impact parameter.

Another key aspect in the quality of printing technologies concerns the mixing of the inks as the droplets collide in mid-air. Although mixing is present in our everyday lives, quantifying mixing (Rasband 1990; Cornfeld *et al.* 2012) and controlling and optimising the mechanisms that affect it remain complex (Lin, Thiffeault & Doering 2011; Lunasin *et al.* 2012). Intuitively, one can acknowledge a state where two substances are completely mixed if the resulting material becomes homogeneous in appearance, therefore, when the gradients of the concentration of the two substances vanish (Aref 1984; Rom-Kedar, Leonard & Wiggins 1990; Walters 2000). Mixing can also be partial (Thiffeault 2012), which would correspond to an intermediate state in which, by some transformation, such as stirring, the components of the systems are transformed from a heterogeneous state into a homogeneous one (Foures, Caulfield & Schmid 2014). When the motion of the fluid is accompanied by diffusion, mixing becomes an irreversible process. In such cases, one can express the degree of mixing in terms of a passive scalar function (Mathew *et al.* 2007). The mixing process in the context of binary droplet collisions has been addressed before; this includes the work by Anilkumar, Lee & Wang (1991), who investigated experimentally the quiescent coalescence of drops of the same liquid. Liu *et al.* (2013), limiting their study to axisymmetric collisions of spherical droplets, used a front tracking algorithm combined with a set of Lagrangian particles to monitor the distribution of liquids. Later, Sun *et al.* (2015) studied a similar set-up and proposed a way of quantifying partial mixing of non-Newtonian droplets.

In this work, the mixing process resulting from a binary droplet collision of non-spherical droplets is studied. This includes identifying the conditions that result in the most efficient mixing. Most bioprinting techniques employ materials of a wide variety of properties, e.g. rheology and surface tension. However, to aid the development of such technologies, the present work is focused on understanding the mechanisms emerging from the collision of droplets with identical properties. In § 2, the underlying physical mechanisms that govern mixing in a binary droplet collision are discussed. Then, in § 3, the numerical methods and the tools for analysing the results are described, and the results are presented in § 4. Finally, conclusions are summarised in § 5.

2. Governing equations

The fluid components are modelled by three concentration fields $\alpha_i(\mathbf{x}) \in [0, 1]$, $i = 1, 2, 3$, where α_1 and α_2 correspond to liquids L1 and L2, respectively, α_3 corresponds to the surrounding gas phase and \mathbf{x} is the position vector. Here, $\alpha_i = 1$ implies the presence of component i , whereas $\alpha_i = 0$ implies its absence. Additionally, the constraint

$$\alpha_1 + \alpha_2 + \alpha_3 = 1 \tag{2.1}$$

is imposed throughout the volume of the fluid mixture Ω . The region of the liquid is denoted by $\Omega_l \subseteq \Omega$, that is, $\Omega_l = \{\mathbf{x} : \alpha_1(\mathbf{x}) + \alpha_2(\mathbf{x}) = 1\}$.

The concentration fields, α_i , follow the transport equation,

$$\partial_t \alpha_i + \nabla \cdot (\alpha_i \mathbf{u}) = \nabla \cdot (D \nabla \alpha_i), \quad \mathbf{x} \in \Omega_l, \tag{2.2}$$

for the liquid components $i = 1, 2$, and

$$\partial_t \alpha_3 + \nabla \cdot (\alpha_3 \mathbf{u}) = 0 \tag{2.3}$$

for the gas. In this way, the two liquid components are miscible, but the liquid and gas phases are immiscible. Here, D is the diffusion coefficient and \mathbf{u} is the velocity field that satisfies the Navier–Stokes equations,

$$\rho(\partial_t + \mathbf{u} \cdot \nabla) \mathbf{u} = -\nabla p + \nabla \cdot (2\rho\nu \mathbf{S}) + \sigma \kappa \hat{\mathbf{n}}_{lg} \delta_S, \tag{2.4}$$

in the incompressible limit ($\nabla \cdot \mathbf{u} = 0$). Here, $\rho := \sum_i \rho_i \alpha_i$ is the density of the fluid, and $\nu := \sum_i \nu_i \alpha_i$ corresponds to the kinematic viscosity, where the constants ρ_i and ν_i correspond to the bulk density and viscosity of the component i . Also, p is the pressure, $\mathbf{S} := (\nabla \mathbf{u} + \nabla \mathbf{u}^T)/2$ is the strain-rate tensor, σ is the surface tension of the liquid–gas interface, $\kappa := -\nabla \cdot \hat{\mathbf{n}}_{lg}$ is its mean curvature, $\hat{\mathbf{n}}_{lg}$ is the orthogonal unitary vector to the surface and δ_S represents a Dirac delta function that locates the liquid–gas interface in the three-dimensional space.

Two types of boundary conditions are imposed: these are the inlets, from which the droplets are injected, and open boundaries that allow the free movement of the flow in or out of the simulation domain specified everywhere else. As illustrated in figure 1(a), the inlets are specified at $z = 0$ as two nozzles (indicated with subscript nzl in the following) of circular cross-section of radius R_{nzl} directed into the domain at the polar and azimuth angles θ and φ , respectively. More formally, to specify the region of the inlets and the open boundaries, let us define the auxiliary function

$$G(\mathbf{x}; x_{nzl}, \theta, \varphi) := 1 - \frac{1}{R_{nzl}^2} \{[(x - x_{nzl}) \cos \varphi + y \sin \varphi]^2 \cos^2 \theta + [y \cos \varphi - (x - x_{nzl}) \sin \varphi]^2\}, \tag{2.5}$$

at the plane $z = 0$. In this way, the inlets are defined as the region where $G \geq 0$. The parameter x_{nzl} corresponds to the location of the centre of the inlet from the origin.

Each droplet is injected into the domain by a pulse of the form

$$T(t) := \begin{cases} 1 - \tanh[8 \cos(\omega t)] & \text{for } 0 \leq t \leq 2\pi/\omega = T_{pulse}, \\ 0 & \text{otherwise,} \end{cases} \tag{2.6}$$

which models the timing and duration of the pulse that injects the liquids into the simulation domain, and emulates the smooth opening and closing of a typical micro-valve used for bioprinting. The orientation of the inflow is specified by the unitary vector

$$\hat{\mathbf{J}}(\theta, \varphi) := \hat{\mathbf{e}}_x \cos \varphi \sin \theta + \hat{\mathbf{e}}_y \sin \varphi \sin \theta + \hat{\mathbf{e}}_z \cos \theta. \tag{2.7}$$

The velocity profile at the inlets is modelled by a parabolic profile, with average velocity $U_{nzl} := (dV/dt)/\pi R_{nzl}^2$. This assumption is justified by considering that the typical Reynolds number on a bioprinting micro-valve is small ($Re < 500$) and the nozzles of

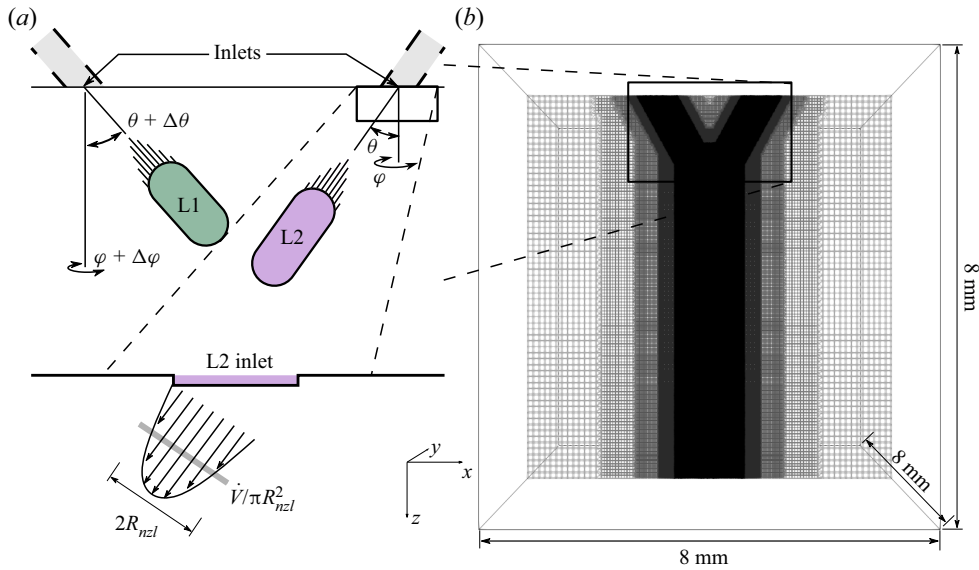


Figure 1. Simulation set-up of the binary droplet collision. (a) Schematic representation of the inlet configuration. (b) The simulation domain is a cube meshed with hexahedral cells with an increasing level of refinement closer to the path drawn by the droplets.

the micro-valves are long ($\sim 20R_{nzl}$), thus the flow field relaxes to the Poiseuille profile as opposed to a plug or a turbulent flow (da Conceicao Ribeiro *et al.* 2018). By combining (2.5)–(2.7), we prescribe the boundary condition for the flow field at the inlets, that is,

$$\mathbf{u}(\mathbf{x}, t) = \begin{cases} (U_{nzl} + \Delta U_{nzl}) T(t - \Delta t_0) G(\mathbf{x}; -x_{nzl}, \theta + \Delta\theta, \varphi + \Delta\varphi) \\ \quad \times \hat{\mathbf{J}}(\theta + \Delta\theta, \varphi + \Delta\varphi) & \text{for L1,} \\ U_{nzl} T(t) G(\mathbf{x}; +x_{nzl}, \theta, \varphi) \hat{\mathbf{J}}(\theta, \varphi) & \text{for L2,} \end{cases} \quad (2.8)$$

where, without loss of generality, the new parameters introduce asymmetry between the nozzles: ΔU_{nzl} corresponds to a velocity difference, Δt_0 represents a delay in the firing time, and $\Delta\theta$ and $\Delta\varphi$ result in a misalignment between the nozzles. Table 1 summarises the values and ranges of the parameters used in the simulations. Note that both liquids have the same material properties, that is, the density, viscosity and surface tension are equal for both droplets. As an example, figures 1 and 2 show the set-up and the average velocity resulting from (2.8).

The concentration fields α_1 and α_2 take the value 1 at the $+x_{nzl}$ and $-x_{nzl}$ inlets, and zero everywhere else on the top boundary, respectively. This is prescribed by means of the auxiliary function G :

$$\alpha_1(\mathbf{x}) = \begin{cases} 1 & \text{if } G(\mathbf{x}; -x_{nzl}, \theta + \Delta\theta, \varphi + \Delta\varphi) \geq 0, \\ 0 & \text{otherwise,} \end{cases} \quad (2.9)$$

and

$$\alpha_2(\mathbf{x}) = \begin{cases} 1 & \text{if } G(\mathbf{x}; +x_{nzl}, \theta, \varphi) \geq 0, \\ 0 & \text{otherwise.} \end{cases} \quad (2.10)$$

Parameter	Value/range	Parameter	Value/range
$\rho_1 = \rho_2$	$= 1.0 \times 10^3 \text{ kg m}^{-3}$	ρ_3	$= 1.0 \text{ kg m}^{-3}$ (for the gas)
$\nu_1 = \nu_2$	$= 1.0 \times 10^{-6} \text{ m}^2 \text{ s}^{-1}$	ν_3	$= 1.48 \times 10^{-5} \text{ m}^2 \text{ s}^{-1}$ (for the gas)
D	$= 2.0 \times 10^{-7} \text{ m}^2 \text{ s}^{-1}$	$\sigma_{13} = \sigma_{23}$	$= 72.0 \text{ mN m}^{-1}$
R_{nzt}	$= 0.25 \text{ mm}$	x_{nzt}	$= 1.125 \text{ mm}$
U_{nzt}	$= 1.0 \text{ m s}^{-1}$	$T_{pulse} = 2\pi/\omega$	$= 2.5 \text{ ms}$
θ	$= 30^\circ$	φ	$= 0^\circ$ for L1 and 180° for L2
ΔU_{nzt}	$\in [-0.8 \text{ m s}^{-1}, +0.8 \text{ m s}^{-1}]$	$\Delta t_0/T_{pulse}$	$\in [-5/16, 5/16]$
$\Delta\theta$	$\in [-10^\circ, 10^\circ]$	$\Delta\varphi$	$\in [-10^\circ, 10^\circ]$ (for L1)

Table 1. Simulation parameters.

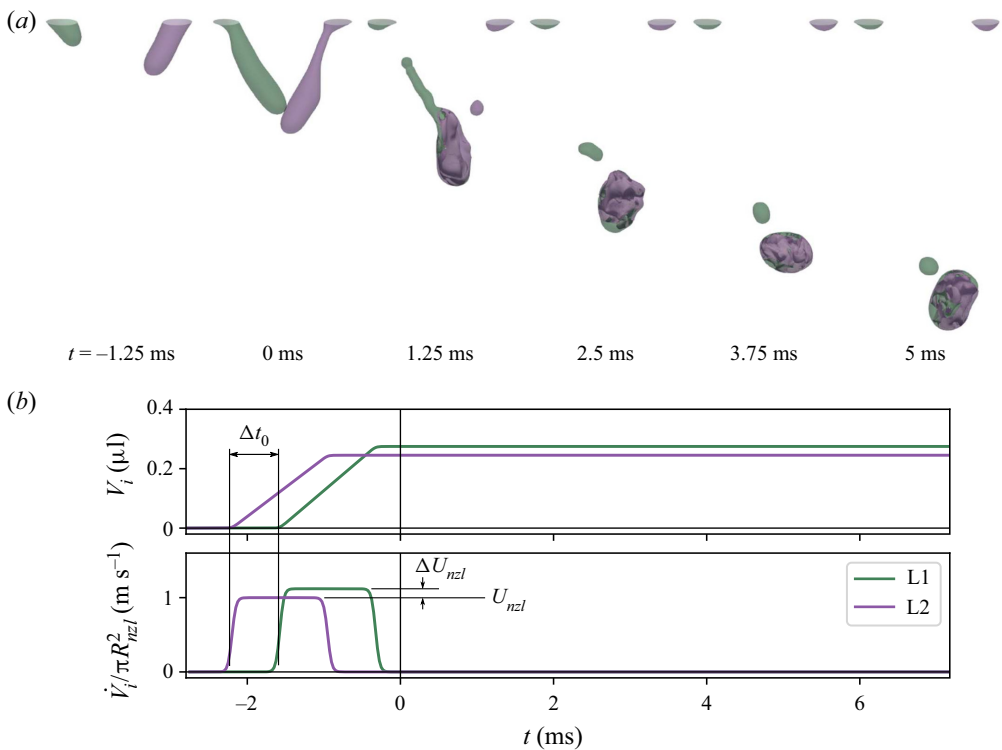


Figure 2. Evolution of the binary droplet collision. (a) Example of a time sequence of a binary droplet collision and its mixing process. The two liquids, L1 and L2, are shown by the isosurface $\alpha_i = 1/2$, $i = 1, 2$ in green and purple, respectively, and the contact surface area is highlighted in grey. (b) Evolution of the volume of the liquids (top) and their flow rate (bottom). The delay in the injection time and the velocity difference are indicated by Δt_0 and ΔU_{nzt} . The temporal coordinate has been shifted, identifying $t = 0$ as the time of contact.

3. Methods and analysis

The coupled system of advection–diffusion and momentum conservation equations in (2.2)–(2.4) is solved in an open-source, finite-volume solver with multiphase capabilities, OpenFOAM (Weller *et al.* 1998). We carry out direct numerical simulations using a second-order scheme for both temporal advancement and spatial discretisation.

The advection–diffusion equations for the concentration fields, (2.2) and (2.3), were solved using volume-of-fluid with the MULES algorithm for two sub-cycles with interface compression coefficient $c_\alpha = 1.5$. The momentum equation (2.4) is solved using the PISO-SIMPLE algorithm for the coupling of the pressure and velocity fields. The pressure is solved by an over-relaxation method with relative tolerance 10^{-7} for the pressure correction and 10^{-9} for the final pressure solution.

As can be observed in figure 1(b), the simulation domain consists of a cubic volume of 8 mm width, meshed with hexahedral cells with four levels of refinement, duplicating the density of cells approaching the path drawn by the liquids. The largest cell type is $h = 250 \mu\text{m}$ in width, which corresponds to $1/32^3$ of the domain volume, and the smallest cell type is $h = 15.625 \mu\text{m}$ in width, or $1/512^3$ of the total volume. The mesh is refined progressively such that the Kolmogorov length scale (Pope 2000), η , is maintained above the local cell size, h , that is, $\min\{\eta/h\} \approx 2.10 > 1$ for all $x \in \Omega$ and t . The increment for temporal advancement, ΔT , is adaptive in order to keep the Courant number, Co , below 0.1, which, on average, resulted in $\Delta T \sim 0.1 \mu\text{s}$.

To ensure that the parasitic currents (Harvie, Davidson & Rudman 2006) do not affect the simulation results, a different set of simulations was carried out. These consisted of a three-dimensional liquid droplet of radius R_{nzl} suspended in air with the material properties reported in table 1 and a homogeneous mesh resolution $h = 15.625 \mu\text{m}$ to match the main simulations. The kinetic energy and asphericity of the droplet are evaluated after the system arrived at a static configuration. It was found that the maximum magnitude of the parasitic currents reach at most up to 0.05 m s^{-1} , which is two orders of magnitude smaller compared to the characteristic velocity of the main simulations. Moreover, most of the analysis of the flows in the present study will be performed at the interface of the two miscible liquids where the surface tension is zero, thus this manifold does not produce parasitic currents. The asphericity, measured as the relative variance of the distance of the points in the isosurface $\alpha_1 = 1/2$ against the ideal droplet radius R_{nzl} , was found to be 2.3×10^{-4} . In conclusion, the numerical errors, arising from the discretisation of surface tension force, are negligible. We begin the simulations at $t = 0$, and they run until the droplets come out of the simulation domain, which occurs in less than 10 ms. However, in what follows, we have shifted the time variable such that contact between the droplets occurs at zero, i.e. $t \rightarrow t - t_c$, where t_c is the instant of first contact, as shown in figure 2(a).

For the parameters specified in table 1, the typical Weber number $We := 2\rho R_{nzl}u_f^2/\sigma$ found in the simulations was $We < 25$, thus we expected a slow coalescence resulting from the collision at small values of the impact parameter. The choice in the parameter ranges for the current analysis is influenced by the experimental set-up of da Conceicao Ribeiro *et al.* (2018). On the one hand, the range of values for Δt_0 and ΔU_{nzl} is unbounded for such devices. However, we observed that $\Delta t_0 \in [-5/16, 5/16]$ and $\Delta U_{nzl} \in [-0.8, 0.8] \text{ m s}^{-1}$ resulted in $B \lesssim 1$, i.e. contact between droplets. On the other hand, the angles for the alignment of the nozzle, $\Delta\theta$ and $\Delta\varphi$, are within the degrees of freedom of such types of bioprinters. These allow for adjustments that compensate for other sources of misalignment.

We will now focus on the quantification of mixing, but first, let us briefly estimate the state of turbulence of both the gas and liquid phases. Following Pope (2000), let us define the Reynolds number in the gas phase as $Re_g := U_g L_g / \nu_g$, where $U_g := \max_{gas} |\mathbf{u}|$ is the maximum velocity in the gas phase, and $L_g := 2U_g^3 / (\nu_g \max_{gas} |\nabla \times \mathbf{u}|^2)$ defines the characteristic length scale. It was found that the typical values for Re_g were approximately 10^3 , suggesting a weakly turbulent flow in the gas phase. For the liquid phase, we define

$Re_l := k^2/\nu_l\epsilon$, where $k := \overline{\mathbf{u}' \cdot \mathbf{u}'}/2$ is the turbulence kinetic energy, and $\epsilon := 2\nu_l \text{tr} \overline{\mathbf{S}^2}$ is the dissipation; the overbar corresponds to the Reynolds average, $\mathbf{u}' := \mathbf{u} - \bar{\mathbf{u}}$ and $\mathbf{S}' := (\nabla \mathbf{u}' + \nabla \mathbf{u}'^T)/2$. To estimate Re_l , we approximate $\bar{\mathbf{u}}$ by the velocity of the centre of mass of the liquid phase. This is an asymptotic approximation that, due to the large dynamic viscosity ratio between the two immiscible phases, is mostly accurate. It was found that the typical values at the instant of first contact between the droplets reached $Re_l \sim 10^3$, decaying to $Re_l \sim 200$ shortly after. Therefore, in contrast to the gas phase, this suggests a quasi-laminar flow. Although not conclusive, this analysis may give us an idea of the state of turbulence of both phases, and thus the tendency for mixing.

3.1. Quantifying mixing

We are interested in the continuous process in which mixing takes place. Therefore, to quantify partial mixing, let us define the segregation parameter between liquids L1 and L2 as

$$\chi := \frac{\langle \alpha_i^2 \rangle - \langle \alpha_i \rangle^2}{\langle \alpha_i \rangle (1 - \langle \alpha_i \rangle)} = 1 - \frac{V_1 + V_2}{V_1 V_2} \int_{\Omega_l} \alpha_1 \alpha_2 \, dV, \tag{3.1}$$

where, without loss of generality, $i = 1$ or $i = 2$ can be chosen and using (2.1) leads to the last equality in (3.1). Here, the operation $\langle \cdot \rangle$ computes the volume average in Ω_l , where α_3 is zero. Since the concentration is a conserved quantity for passive scalar mixing, any diffusive flux is assumed to be zero at the liquid–gas interface. It follows that $\langle \alpha_i \rangle = V_i/(V_1 + V_2)$, where V_i is the total volume of liquid i injected.

As the name suggests, $\chi = 0$ corresponds to an absence of variance in concentration, which implies a homogeneous or complete mixture. Conversely, $\chi = 1$ corresponds to a state in which the regions occupied by the liquids L1 and L2 do not intersect (Doering & Thiffeault 2006; Foures *et al.* 2014; Thiffeault 2021). The transport equation of the segregation parameter can be deduced by taking the material derivative and using (2.2), which results in

$$\frac{d\chi}{dt} = -2 \frac{V_1 + V_2}{V_1 V_2} \int_{\Omega_l} N \, dV, \tag{3.2}$$

where $N := D|\nabla \alpha_1|^2$, the integrand in the right-hand side of (3.2), corresponds to the scalar dissipation rate.

The scalar dissipation rate N is a positive quantity, proportional to the rate of entropy production by diffusion, which indicates the irreversibility of a mixing process (Davidson 2015). The transport equation of N is given by (Chakraborty *et al.* 2011)

$$\frac{dN}{dt} = \nabla \cdot (D \nabla N) - 2D^2 |\nabla \nabla \alpha_1|^2 - 2N \mathbf{S} : \hat{\mathbf{n}}_c \hat{\mathbf{n}}_c, \tag{3.3}$$

where $d/dt := (\partial_t + \mathbf{u} \cdot \nabla)$ is the material derivative and $\hat{\mathbf{n}}_c := \nabla \alpha_1 / |\nabla \alpha_1|$ is the unit normal vector of the contact surface. The first term in the right-hand side of (3.3) corresponds to the contribution by the fluxes driven by gradients of N . The second term is dubbed the molecular dissipation rate and represents a sink term. The last term in the right-hand side of (3.3) is called the scalar–turbulence interaction term and models the contribution by the stretching of the flow; it can be expressed in terms of the eigenvalues

$\{\lambda_i\}_{i=\alpha,\beta,\gamma}$ and eigenvectors $\{\hat{\mathbf{e}}_i\}_{i=\alpha,\beta,\gamma}$ of \mathbf{S} (Dresselhaus & Tabor 1992):

$$\mathbf{S} : \hat{\mathbf{n}}_c \hat{\mathbf{n}}_c = \sum_{i=\alpha,\beta,\gamma} (\hat{\mathbf{n}}_c \cdot \hat{\mathbf{e}}_i)^2 \lambda_i. \tag{3.4}$$

Both the segregation parameter and the scalar dissipation rate are mixing measures of the type $\langle |\nabla^m \alpha_i|^2 \rangle$, for $m = 0, 1$ (Doering & Thiffeault 2006).

To keep track of the region where the mixing takes place, one can define the contact surface area (Vervisch *et al.* 1995) between the two liquids,

$$S_c := \int_{\Omega_l} |\nabla \alpha_1| \, dV. \tag{3.5}$$

As for any diffusive process, the total mixing rate is proportional to the surface area; therefore, the higher the S_c , the quicker the segregation parameter will decrease. The evolution of the contact area is governed by the transport equation

$$\frac{dS_c}{dt} = -D \int_{\Omega_l} \frac{|\nabla \nabla \alpha_1|^2}{|\nabla \alpha_1|} \, dV - \sum_{i=\alpha,\beta,\gamma} \int_{\Omega_l} (\hat{\mathbf{n}}_c \cdot \hat{\mathbf{e}}_i)^2 \lambda_i |\nabla \alpha_1| \, dV, \tag{3.6}$$

which can be deduced using (3.3) and (3.5).

3.2. Analysis of the flow structure

To understand the effect that convective flows have in mixing, and in particular, on the terms described in (3.4), the flow field will be analysed in the region of contact between the two liquids. For that, the behaviour of the strain-rate tensor, \mathbf{S} , and, more generally, the velocity gradient tensor, $\mathbf{A} := \nabla \mathbf{u}$, will be analysed.

To illustrate the meaning of \mathbf{A} , consider a virtual swarm of particles at some position \mathbf{y} in a small region around \mathbf{x} in the three-dimensional Cartesian space that are advected by the flow field \mathbf{u} . Suppose that the motion of the particles is given by $d\mathbf{y}/dt = \mathbf{u}(\mathbf{y})$, and assume that \mathbf{u} is continuous around \mathbf{x} . Then, $d\mathbf{y}/dt \approx \mathbf{u}(\mathbf{x}) + (\mathbf{y} - \mathbf{x}) \cdot \mathbf{A}(\mathbf{x})$. This implies that $\mathbf{u}(\mathbf{x})$ corresponds to the background motion of the virtual particles, whilst \mathbf{A} describes their relative motion. Therefore, by means of the velocity gradient tensor, we can analyse the local flow field at \mathbf{x} such as the flow topology and the eigenvectors and eigenvalues of \mathbf{A} (Chong, Perry & Cantwell 1990; Meneveau 2011).

The eigenvalues of \mathbf{A} correspond to the roots of the characteristic polynomial $\det(\mathbf{A} - \lambda' \mathbf{I}) = \lambda'^3 + P\lambda'^2 + Q\lambda' + R$. The coefficients P , Q and R are called the first, second and third invariants of the transformation, and are given by

$$P = -\text{tr } \mathbf{A} = -\sum_i \lambda'_i \quad (= 0 \text{ due to incompressibility}), \tag{3.7}$$

$$Q = \frac{1}{2}[(\text{tr } \mathbf{A})^2 - \text{tr } \mathbf{A}^2] = \lambda'_1 \lambda'_2 + \lambda'_2 \lambda'_3 + \lambda'_3 \lambda'_1, \tag{3.8}$$

$$R = -\det \mathbf{A} = -\prod_i \lambda'_i. \tag{3.9}$$

We have used the primed notation to emphasise the difference between the eigenvalues of \mathbf{S} and those of \mathbf{A} . Since \mathbf{S} is a symmetric tensor, the eigenvalues are real and the eigenvectors are orthogonal, which is not the case for the eigenvalues and eigenvectors

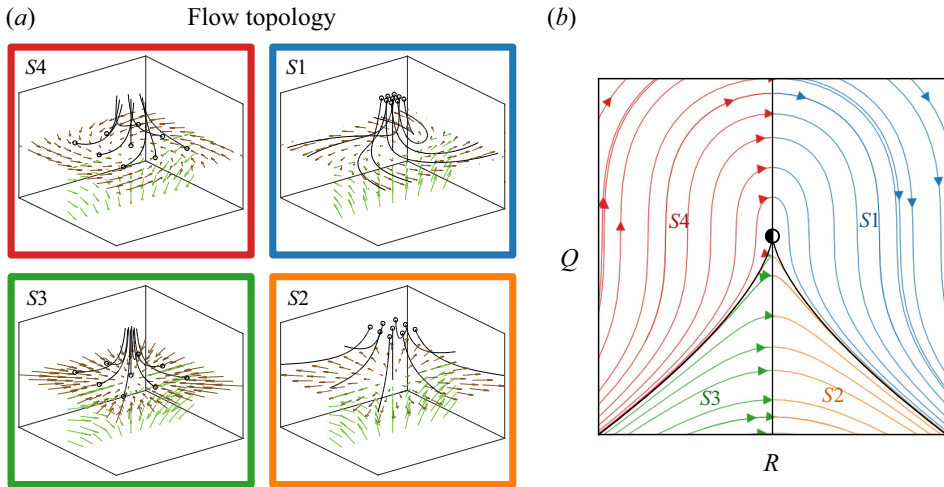


Figure 3. Flow topology and dynamics in the Q - R phase space. (a) Schematic colour-coded classification of the flow topology. The solid black lines correspond to streamlines beginning at the empty circles. The brown and green arrows represent the equal- and opposite-sign manifolds of the vector field, respectively. (b) The streamlines show the trajectories of the restricted Euler system, and the colours indicate the different flow topology regions.

of \mathbf{A} . To distinguish when the eigenvalues are real-valued or have an imaginary part, let us define the discriminant

$$\Delta := \frac{R^2}{4} + \frac{Q^3}{27}. \quad (3.10)$$

If $\Delta \leq 0$, then all three eigenvalues are real; however, if $\Delta > 0$, then \mathbf{A} has one real and two complex eigenvalues. If the eigenvalues have an imaginary part, then the nearby trajectories spiral inwards or outwards from the observation point \mathbf{x} . The real part of the eigenvalues reveals the stability at the observation point. From (3.7), it follows that the real parts of the eigenvalues cannot have the same sign, i.e. one must be positive, one must be negative, and an intermediate one can have either sign. A negative eigenvalue indicates that trajectories are attracted towards the observation point whereas positive indicates that they are repelled. Consequently, the sign of R determines the number of stable and unstable manifolds around the observation point, and when $R = 0$, the linear approximation may not be enough to determine the local stability at \mathbf{x} .

We classify the topology of the flow as follows (Chong *et al.* 1990).

- (i) $S1$, $\Delta > 0$ and $R > 0$, consists of a one-dimensional (1-D) compressive manifold combined with a two-dimensional (2-D) unstable manifold that spirals away from the focal point.
- (ii) $S2$, $\Delta \leq 0$ and $R > 0$, 1-D compressive and 2-D expansive manifolds.
- (iii) $S3$, $\Delta \leq 0$ and $R < 0$, 1-D expansive and 2-D compressive manifolds.
- (iv) $S4$, $\Delta > 0$ and $R < 0$, 1-D stretching and 2-D stable focus manifolds.

Furthermore, the topologies $S1$ and $S4$ can be subdivided depending on the sign of Q , that is, $S1^+$ and $S1^-$ ($S4^+$ and $S4^-$) for $Q \geq 0$ and $Q < 0$, respectively. The representative flows are sketched in figure 3(a) and their corresponding regions in the Q - R space in figure 3(b).

The invariants Q and R follow the equations of motion away from the interfaces (Martin *et al.* 1998):

$$\frac{dR}{dt} = \frac{2}{3} Q^2 - \text{tr}(\mathbf{A}^2 \mathbf{H}) - \nu \text{tr}(\mathbf{A}^2 \nabla^2 \mathbf{A}), \quad (3.11)$$

$$\frac{dQ}{dt} = -3R - \text{tr}(\mathbf{A} \mathbf{H}) - \nu \text{tr}(\mathbf{A} \nabla^2 \mathbf{A}), \quad (3.12)$$

where $\mathbf{H} := -(\nabla \nabla p - I \nabla^2 p/3)/\rho$ is the anisotropic pressure Hessian. The last two terms in both (3.11) and (3.12) are non-local, thus depend on the boundary conditions and the capillary forces at the liquid–gas interface. The terms proportional to ν in (3.11) and (3.12) are derived from the viscous dissipation term in the Navier–Stokes equation, thus act as sinks.

To understand the dynamics of the system in (3.11) and (3.12), let us first consider the restricted Euler system that assumes that $\mathbf{H} + \nu \nabla^2 \mathbf{A}$ is negligible. This corresponds to a state in which the pressure is isotropic and viscous damping is negligible. In this case, (3.11) and (3.12) are reduced to a closed and autonomous dynamical system whose phase space portrait is depicted in figure 3(b). At the origin ($R = 0$ and $Q = 0$), a half-stable fixed point is found, and the surrounding trajectories evolve from $R < 0$ to $R > 0$ and points in the $\Delta = 0$ manifold stay in it. This implies that the natural tendency of the reduced dynamical system is to separate the rotation-dominated from the strain-dominated regions, and the evolution takes states from $R < 0$ to $R > 0$. In other words, the $S4$ and $S3$ topologies are turning into the $S1$ and $S2$ topologies, respectively (Martin *et al.* 1998).

Furthermore, far from interfaces, the transport equations of enstrophy and strain-rate production can be expressed in terms of the invariants

$$\frac{d\zeta}{dt} = 4(R_S - R) + \nu \boldsymbol{\omega} \cdot \nabla^2 \boldsymbol{\omega}, \quad (3.13)$$

where $\zeta := |\boldsymbol{\omega}|^2/2$ is the enstrophy and $\boldsymbol{\omega} := \nabla \times \mathbf{u}$ is the vorticity, and

$$\frac{dQ_S}{dt} = -2R_S - R - \frac{1}{\rho} \text{tr} \mathbf{S} \nabla \nabla p - \nu \text{tr} \mathbf{S} \nabla^2 \mathbf{S}, \quad (3.14)$$

where $Q_S := -\text{tr} \mathbf{S}^2/2$ and $R_S := -\det \mathbf{S}$ are the second and third invariants of the strain rate, respectively. Therefore, for the restricted Euler system, (3.13) and (3.14) can be combined into

$$\frac{d}{dt} \left(\frac{\epsilon}{\nu} \right) = 12R + 2 \frac{d\zeta}{dt}, \quad (3.15)$$

where $\epsilon := -4\nu Q_S$ is the instantaneous dissipation rate of kinetic energy.

This concludes the conceptual framework that will support the analysis of our numerical simulations that will follow. However, before proceeding, these tools will be used for a similar system where an analytic approximation is available.

3.3. Axisymmetric collision of identical spherical droplets

To illustrate the concepts introduced previously, let us consider the case of an axisymmetric binary droplet collision of identical droplets. Following Roisman (2004), after the first contact and during the initial stages of the collision, the droplets undergo

a deformation in which the flow field in the vicinity of the contact surface is well approximated by

$$u_x \approx \frac{u_r x}{2l^3} (4x^2 - 3l^2), \tag{3.16}$$

$$u_\xi \approx \frac{3u_r}{4l^3} \xi (l^2 - 4x^2), \tag{3.17}$$

where x is the coordinate that runs parallel to the axis of symmetry, $\xi = (y^2 + z^2)^{1/2}$ is the distance from the axis, u_r is the magnitude of the relative velocity and l is the thickness of the liquid lamella that forms by radial expansion of the droplets at early stages. Then the velocity gradient tensor is given by

$$\mathbf{A} = \frac{3u_r}{4l^3} \begin{pmatrix} -2(l^2 - 4x^2) & 0 & 0 \\ -8xy & (l^2 - 4x^2) & 0 \\ -8xz & 0 & (l^2 - 4x^2) \end{pmatrix}. \tag{3.18}$$

At the plane of collision $x = 0$, \mathbf{A} reduces to a diagonal matrix whose second and third invariants are given by

$$R(x = 0) = \frac{27u_r^3}{32l^3} \quad \text{and} \quad Q(x = 0) = -\frac{27u_r^2}{16l^2}. \tag{3.19a,b}$$

The evaluation of the discriminant from the characteristic equation (3.10) shows that $\Delta = 0$, which corresponds to an S2 topology flow.

In this stage of the collision, the contact surface area is increasing due to the stretching motion of the flow. The scalar–turbulence interaction term evaluates to

$$-2\mathbf{S} : \hat{\mathbf{n}}_c \hat{\mathbf{n}}_c = \frac{3}{l} u_r \sim \frac{1}{N} \frac{dN}{dt}, \tag{3.20}$$

where the last relation is based on (3.3). This estimates an exponential growth rate of the scalar dissipation rate, and consequently the contact surface area.

The growth of the contact surface area cannot carry on indefinitely due to surface tension and viscous forces. Surface tension reverses the direction of the flow in a periodic fashion, leading to oscillations in the contact surface area, whilst viscous stresses dampen the motion of the flow (Roisman 2009; Roisman, Berberović & Tropea 2009).

4. Results and discussion

The simulation parameters considered for the present analysis are summarised in table 1 and are inspired by a bioprinting set-up (da Conceicao Ribeiro *et al.* 2018). To investigate the effects of differences in the configuration of the L1 inlet relative to L2, the simulations were run in four different batches: (i) varying the injection velocity, ΔU_{nzl} ; (ii) varying the relative injection time, Δt_0 ; (iii) introducing a difference in the polar angle by $\Delta\theta$; and (iv) changing the azimuth angle $\Delta\varphi$.

Additionally, to illustrate the combined effect of these configuration parameters in the dynamics of the droplet collision and mixing for an asymmetric collision presented in figures 2, 8 and 9, we set $\Delta U_{nzl} = 0.24 \text{ m s}^{-1}$, $\Delta t_0 = T_{pulse}/4$, $\Delta\theta = -4^\circ$ and $\Delta\varphi = 6^\circ$. This is to contrast against the symmetric collision presented in figures 6 and 7, in which these configuration parameters are equal to zero ($\Delta U_{nzl} = \Delta t_0 = \Delta\theta = \Delta\varphi = 0$).

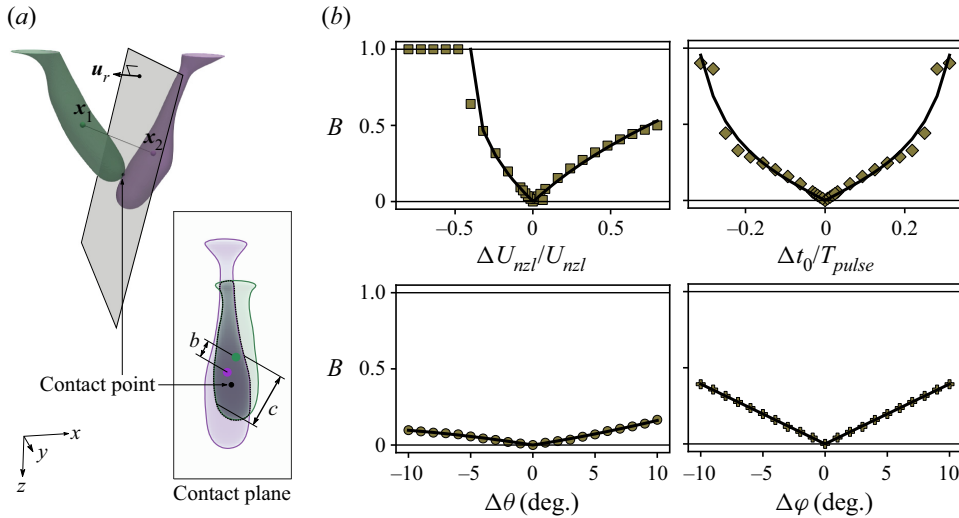


Figure 4. Calculation of the impact parameter. (a) Example of an asymmetric collision between two droplets at the moment of first contact that illustrates the construction of the dimensionless impact parameter B . (b) The impact parameter is evaluated at different set-up configurations (symbols) and the fitting of the curve in (4.9) applied to each case (solid line).

4.1. Impact of non-spherical droplets

It can be observed from figure 2(a) that the droplets do not have a spherical shape, in contrast to many studies of binary droplet collisions, but are rather elongated in the direction of motion relative to their respective nozzles. As expected, the contact point of the droplets would produce different outcomes depending on the shape of the droplet; the dimensionless impact parameter B , as defined for spherical droplets (Ashgriz & Poo 1990; Al-Dirawi & Bayly 2019), cannot be used to represent the off-centred collision in this work.

In the following, we generalise the definition of the dimensionless impact parameter for droplets that are non-spherical and continuously evolving. We will base our definition on the underlying concept of the traditional expression as stated in § 1.

Since the shape of the droplets is changing in time, the impact parameter is meaningful only at $t = 0$, which corresponds to the moment of first contact. The relative velocity between the two droplets is $\mathbf{u}_r := \mathbf{U}_1 - \mathbf{U}_2$, where \mathbf{U}_i is the velocity of the centre of mass of droplet i . The vector \mathbf{u}_r and the first contact point between the two liquids create the contact plane (see figure 4a). The shape of the droplets is then projected onto the contact plane, and if their projected images intersect, then the droplets are bound to collide; otherwise, they pass each other without contact.

The distance between the centre of mass of the droplets on the contact plane, $b := |\mathbf{b}|$, is defined such that

$$\mathbf{b} := (\mathbf{X}_1 - \mathbf{X}_2) \times \frac{\mathbf{u}_r}{|\mathbf{u}_r|}, \quad (4.1)$$

where \mathbf{X}_i , $i = 1, 2$, are the centres of mass of the liquids L1 and L2, respectively. Let P_i , $i = 1, 2$, be the projected image of droplet i onto the contact plane. Then, let $C = \bigcap_i P_i$ be the set of points at the intersection. As illustrated in figure 4(a), we measure the depth of the overlap between the projections, c , by finding the longest line in C that is parallel to

b , formally,

$$c := \max\{c' : c'b/b = (x_2 - x_1); x_1, x_2 \in C\}, \tag{4.2}$$

and $c = 0$ if $C = \emptyset$. In this way, the generalised dimensionless impact parameter can be defined as

$$B := \frac{b}{b + c}. \tag{4.3}$$

Note that $b \geq 0$ and $c \geq 0$ and if $b = 0$, then $c \neq 0$; this implies that $0 \leq B \leq 1$, where $B = 0$ corresponds to a head-on collision, and $B = 1$, which occurs if $c = 0$, implies that the droplets do not make contact during their trajectories at any time.

For spherical droplets of diameter d_i , the depth of the overlap reduces to $c = (d_1 + d_2)/2 - b$ and we recover the usual expression $B = 2b/(d_1 + d_2)$.

Let us anticipate the functional dependence of B after a small deviation from the head-on collision. This can be done by performing a small deviation around the head-on collision for the terms b and c in (4.3). Therefore, an expression can be obtained by considering

$$B \sim \frac{b_0 + \Delta b}{b_0 + \Delta b + c_0 + \Delta c} \sim \frac{\Delta b}{c_0}, \tag{4.4}$$

where b_0 and c_0 are the values for the head-on collision, and Δb and Δc represent their variation from the parameters that introduce the asymmetry (ΔU_{nzl} , Δt_0 , $\Delta\theta$ and $\Delta\varphi$) up to linear order. The last relation in (4.4) is obtained recalling $b_0 = 0$ for the head-on collision.

The expressions for Δb and c_0 can be estimated by assuming that the droplets have the morphology of a cylindrical rod of radius R_{nzl} and length $\ell_1 = T_{pulse}(U_{nzl} + \Delta U_{nzl})$ and $\ell_2 = T_{pulse}U_{nzl}$ for liquids L1 and L2. Here, c_0 corresponds to the overlapping length in the direction of \mathbf{b} . However, in the head-on collision case, c_0 is degenerate since $\mathbf{b} = 0$, thus it can take the two values depending on the source of asymmetry:

$$c_0 \approx \begin{cases} 2R_{nzl} & \text{if } \Delta\varphi \neq 0, \\ U_{nzl}T_{pulse} \cos \theta & \text{otherwise.} \end{cases} \tag{4.5}$$

To calculate Δb , we can approximate the positions of the centres of mass of the two droplets as

$$\mathbf{X}_1 \approx (t - \Delta t_0 - T_{pulse}/2)(U_{nzl} + \Delta U_{nzl})\hat{\mathbf{J}}(\theta + \Delta\theta, \Delta\varphi) - x_{nzl}\hat{\mathbf{e}}_x \tag{4.6}$$

and

$$\mathbf{X}_2 \approx (t - T_{pulse}/2)U_{nzl}\hat{\mathbf{J}}(\theta, \pi) + x_{nzl}\hat{\mathbf{e}}_x. \tag{4.7}$$

The relative velocity of the droplets is given by $\mathbf{u}_r = d(\mathbf{X}_1 - \mathbf{X}_2)/dt$. Then, by substituting (4.6) and (4.7) in (4.1), and keeping terms up to linear order, it can be shown that

$$\frac{\Delta b}{c_0} = \left[\left(\frac{\Delta t_0}{T_{pulse}} \right)^2 + \left(\frac{x_{nzl}}{U_{nzl}T_{pulse}} \right)^2 \left(\frac{\Delta\theta^2}{\cos^2 \theta} + \frac{\Delta U_{nzl}^2}{U_{nzl}^2 \sin^2 \theta} \right) + \left(\frac{x_{nzl} \Delta\varphi}{2R_{nzl}} \right)^2 \right]^{1/2}. \tag{4.8}$$

In figure 4, the value of the impact parameter calculated from the simulation results is presented. Guided by (4.8), we propose an empirical approximation to B that captures its

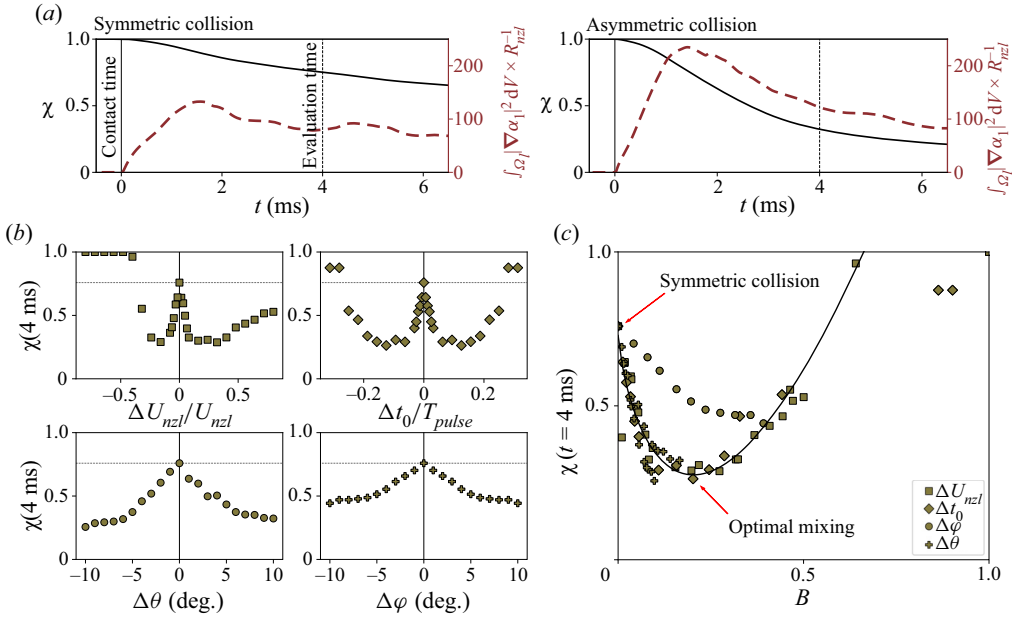


Figure 5. Quality of the mixing process after breaking the symmetry of the collision. (a) Evolution of the segregation parameter (solid line) and scalar dissipation rate (dashed line) for a symmetric and an asymmetric collision. (b) Evaluation of the segregation parameter after 4 ms after first contact for injection velocity difference, injection time difference, and polar and azimuthal angles. (c) Segregation as a function of the impact parameter. The solid line is an approximation to the common trend followed by all the points, except for those corresponding to $\Delta\varphi$.

general trend by varying each parameter:

$$B \approx \left[c_t \tan^2 \frac{\Delta t_0}{b_t T_{pulse}} + (c_\theta \Delta\theta + d_\theta \Delta\theta^2)^2 + (c_u (\Delta U_{nzi}/U_{nzi})^n - 1)^2 + (c_\varphi \Delta\varphi)^2 \right]^{1/2}, \quad (4.9)$$

where each of the constants is found by least squares optimisation. Varying ΔU_{nzi} , we find that $n = 0.4$ and $c_u = 1.43$; varying Δt_0 , gave the constants $b_t = 0.26$ and $c_t = 0.12$. Varying the angles $\Delta\theta$ and $\Delta\varphi$ resulted in $c_\theta = 0.018 \text{ rad}^{-1}$ and $d_\theta = 43.2 \text{ rad}^{-2}$, and $c_\varphi = 2.25 \text{ rad}^{-1}$, respectively.

Figure 5(a) shows the evolution of the segregation parameter χ and the scalar dissipation rate N . As the droplets collide, the mixing process begins and the segregation parameter decreases. The scalar dissipation rate shows a fast increase after the first contact between the droplets. Soon after, it arrives at a local maximum, which is reflected as an inflection point in χ at $t = 1.575 \text{ ms}$ and $t = 1.425 \text{ ms}$ for the symmetric and asymmetric collisions, respectively. From that instant onwards, χ will decrease at a lower rate. A similar behaviour was observed in all our simulations.

In figure 5(b), the segregation parameter at $t = 4 \text{ ms} = 1.6T_{pulse}$, the evaluation time, is shown for all simulation results. The evaluation time is chosen as the maximum time lapse after the collision that is common to all simulations. In other words, it is the longest interval of time for the mixing process to take place before the liquids exit the simulation domain. Since χ is a monotonically decreasing function of time, the results are similar at any other evaluation time of choice.

The four plots in figure 5(b) show the influence of the different configuration parameters in the mixing process. It can be observed from figure 5(b) that decreasing the velocity of the inlet has a strong effect on χ and reducing the velocity by $\sim 15\%$ gives a near-optimal mixing; however, decreasing by $\sim 40\%$ raises B to $B = 1$, and mixing is absent. In contrast, increasing the velocity has a smaller effect, where the optimal mixing is found at a $\sim 32\%$ increase in the inlet velocity.

The difference in injection time shows a fast improvement in the quality of the mixing, reaching an optimal mixing at $|\Delta t_0| = 0.12T_{pulse}$, which afterwards start deteriorating until $|\Delta t_0| = 0.3T_{pulse}$, when the droplets barely touch each other.

Changes in the polar and azimuthal angles, $\Delta\theta$ and $\Delta\varphi$, show an improvement in the mixing quality when breaking the symmetry of the collision. However, varying the azimuthal angle has a smaller improvement on the mixing even though the impact parameter is found to increase at a higher rate.

In summary, it can be observed in figure 5(c) that the segregation parameter is strongly dependent on the impact parameter. Over a finite time, the segregation parameter shows a high value for the symmetric collision; this is not the exception, but rather the common trend for different sources of asymmetry. The segregation parameter decreases, reaching a wide minimum at $B = 0.2 \pm 0.1$ in which optimal mixing occurs. For higher values of the impact parameter, mixing deteriorates. The overall tendency is captured by the function

$$\chi \approx a_\chi B \ln \frac{B}{b_\chi} + \chi_0, \tag{4.10}$$

where the constants $a_\chi = 2.23$ and $b_\chi = 0.551$ are found by least squares optimisation, and $\chi_0 = 0.734$ is the value of the segregation parameter for the symmetric collision. The function in (4.10) has its global minimum at $B = b_\chi/e \approx 0.203$, where e is the base of the natural logarithm. This is true except for the set of simulations where $\Delta\varphi \neq 0$. As can be observed, the segregation parameter does not decrease as much as the other cases. This does not imply that the definition of the impact parameter is inadequate. As will be discussed later, the system exhibits a deeper-level complexity, which is reflected in the infeasibility of reaching a near optimal state of mixing.

4.2. Evolution of the topology of the flow and the contact surface

To better understand why optimal mixing occurs for $B \neq 0$, we will explore in more detail the evolution of the contact surface and the flows in its vicinity.

Figure 6(a) shows the contact surface of a symmetric binary droplet collision. This corresponds to the isosurface $\alpha_1 = 1/2$ at two different times: $t = 0.5$ ms and $t = 2$ ms. It can be observed from figure 6(a) that the contact surface is flat, which is due to the reflexive symmetry of the system. The contact surface between the two liquids occurs at the mirror plane $x = 0$. This implies that $\hat{n}_c = \hat{e}_x$ for all $t \geq 0$. The x -component of the velocity field must vanish at the mirror plane. This implies that $\mathbf{u} \cdot \nabla\alpha_1 = 0$, which takes away the convective term in (2.2). Moreover, the velocity gradient tensor at the plane $x = 0$ becomes a block diagonal matrix of the form

$$\mathbf{A}^{(sym)}|_{x=0} = \begin{pmatrix} \partial_x u_x & 0 & 0 \\ 0 & \partial_y u_y & \partial_z u_y \\ 0 & \partial_y u_z & \partial_z u_z \end{pmatrix} \quad \forall t > 0. \tag{4.11}$$

This is because the component u_x of the velocity vector is an odd function and the components u_y and u_z are even functions of x , thus $\partial_x u_y = \partial_x u_z = 0$, whilst $\partial_x u_x$ may

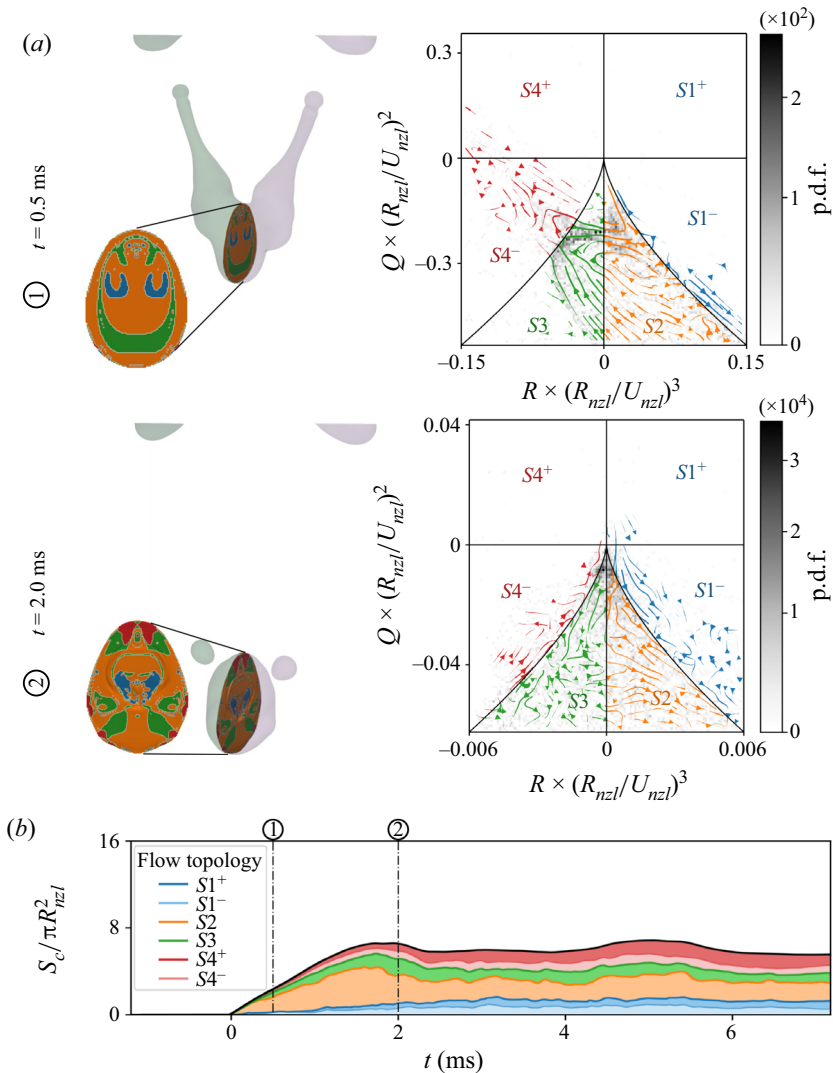


Figure 6. Topological analysis of the flow field at the contact surface. (a) Colour map showing the distribution of the different topological classifications at the isosurface $\alpha_1 = 1/2$ according to the colour code of figure 3(a) at two instants with the corresponding phase portrait of the Q - R space. The joint p.d.f. is displayed by a greyscale map in the background and by the thickness of the streamlines. (b) Evolution of the contact surface area, S_c , broken down into the different types of topologies at the surface.

not. Additionally, since $u_x = 0$ for all y and z , we find that $\partial_y u_x = \partial_z u_x = 0$ as well. In consequence, the existence of a real-valued eigenvector and an eigenvalue parallel to \hat{n}_c is guaranteed, and the vorticity can have a non-zero component only in the x direction, i.e. $\omega_y = \omega_z = 0$.

As shown in figure 6(a), the dominant flow topology at early times corresponds to $S2$, which is expected from the flow profile of the initial deformation (Roisman 2004) of the collision. Two vortical structures of type $S1$ appear near the centre and also an annular section with topology $S3$. Points with topology $S4$ cover a small area of the contact surface.

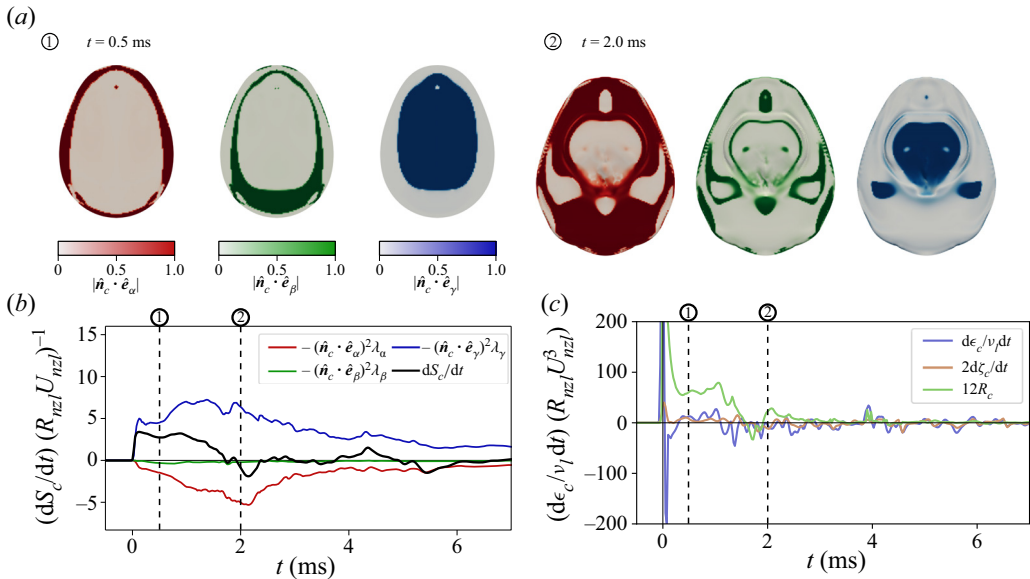


Figure 7. Eigenvector analysis of the flow field at the contact surface and growth rate of a symmetric collision. (a) Colour map showing the projections of the three eigenvectors against the orthonormal vector to the isosurface $\alpha_1 = 1/2$ at two different times. (b) Growth rate of the contact surface area and the contribution of each eigencomponent of the scalar–turbulence interaction terms. (c) Strain production at the contact surface broken down into the enstrophy production and third invariant.

In figure 6(b), the evolution of the contact surface area shows an initial increase at early times, which is slowed down and begins to oscillate. These oscillations are the product of the interplay between surface tension forces and inertia, and are dampened by viscous stresses. Additionally, the partition of the contact surface into the different topologies shows that, at early stages, the most dominant topology of the flow corresponds to S_2 . After the contact surface reaches its maximum area, the contributions from S_1 and S_4 are greater. Moreover, the decomposition for the different signs in Q reveals that, over time, an increasing fraction of the contact surface contains strong vortical structures where $Q > 0$.

Figure 7(a) complements the analysis of the flow structure. An inner region in which the direction of the most compressive eigenvalue, λ_γ , is aligned with the normal vector of the surface can be observed. This region covers most of the area of the contact surface and, according to (3.6), contributes to its growth. A thin outer region of the contact surface shows an alignment of the normal vector and the eigenvector \hat{e}_α . This implies the contraction of the surface, which can be attributed to the action of the capillary forces slowing down the radial expansion and forming a rim. An intermediate region, where the eigenvector \hat{e}_β serves as the transition between the two regions, can also be seen in figure 7(a). This is further verified by the phase portrait in the Q – R space of figure 6(a) at $t = 0.5$ ms, where the streamlines show that points of topology S_2 migrate to S_3 and back.

At $t = 2$ ms, it can be observed that although S_2 still covers most of the area, there is more presence of the other flow topologies. However, as can be seen in the Q – R space, these populations are small and stay close to the $\Delta = 0$ curve from (3.10).

Figure 7(b) shows the growth rate of the contact surface area and the respective contributions from the eigenvalue decomposition of the scalar–turbulence interaction term. The initial rise of the contact surface area can be attributed to the contribution of the most compressive eigenvalue, λ_γ . However, the contribution from the most expansive

eigenvalue, λ_α , is negative (see (3.6)) and has the effect of decreasing the surface area. It has a slower increase but eventually catches up with the λ_γ contribution. The contribution from the intermediate eigenvalue is almost negligible throughout the mixing process.

We now consider the evolution of the strain rate and enstrophy at the contact surface. For that, we will make use of (3.15) in the following form of the restricted Euler system:

$$\frac{1}{v_l} \frac{d\epsilon_C}{dt} \approx 12R_C + 2 \frac{d\zeta_C}{dt}, \quad (4.12)$$

where the subscript C in all quantities denotes integration over the contact surface, e.g. $\epsilon_C := \int \epsilon |\nabla \alpha_1| dV$. Thus ϵ_C , R_C and ζ_C are the total instantaneous dissipation of kinetic energy, third invariant and enstrophy at the contact surface.

Figure 7(c) shows the evolution of the strain-rate production at the contact surface. Beyond small fluctuations, both the strain-rate production and enstrophy production are small. In contrast, at early times ($0 \leq t \lesssim 2$ ms), the third invariant has a significant contribution. This first stage is related to the stretching of the contact surface and can be associated with the same event in figure 7(b) where dS_C/dt becomes negative.

The evolution of an asymmetric collision shows several differences, as can be seen in figures 8 and 9. From early stages, it can be observed that the isosurface $\alpha_1 = 1/2$ is no longer bound to a plane and quickly distorts to a surface with multiple protrusions and twists.

As shown in figures 8(a) and 9(a), during the early stages of the droplet collision, the lower part of the contact surface is covered by a vorticity-dominant topology that implies the twisting of the contact surface area into a pocket-like shape. The streamlines in the Q - R space show this phenomenon through the migration of points from the $S2$ topology to the $S4$. At later stages ($t > 2$ ms), a large population of vortical topologies is found in the Q - R space.

Figure 8(b) shows the fraction of the area of flow topology $S4$ that increases rapidly and soon covers most of the contact area. Followed by this, the areas of topology $S1$ also begin to dominate as expected from the reduced Euler system in the Q - R space. Compared to the symmetric collision, the fraction of area with $Q > 0$ is much larger. This is more appreciable for the topology $S4^+$, which becomes large in early stages and persistently throughout the mixing process. This implies that points in the $S3$ topology are ejected to $S4^+$, spending a short time in the $S4^-$ region of the phase space. Additionally, the topologies $S1^+$ and $S1^-$ have very similar contributions to the surface area, which is a consequence of the restricted dynamics of the Q - R phase space.

The eigenvalue analysis corroborates such trends, showing that although there is a superposition of projections of the eigenvectors, the most dominant corresponds to λ_γ (see figure 9b) and persists throughout the mixing process. Although the total contribution of the scalar-turbulence interaction term (i.e. (3.6)) is positive, dS_C/dt is negative at late times, which can be attributed to self-intersections of the contact surface and the effect of the molecular dissipation rate.

In contrast to the symmetric case, the asymmetry in the collision has a very different outcome in terms of the enstrophy and strain-rate production, as can be observed in figure 9(c). In this case, the three quantities have a similar magnitude during the first stage ($0 \leq t \lesssim 1.5$ ms). This implies that both the dissipation and vorticity are increasing at the contact surface area. Similarly, the decrease in dS_C/dt shown in figure 9(b) can also be related to the drop-off of the three quantities. Furthermore, the third invariant becomes small after $t \approx 2$ ms; however, the enstrophy and strain-rate productions are now negative.

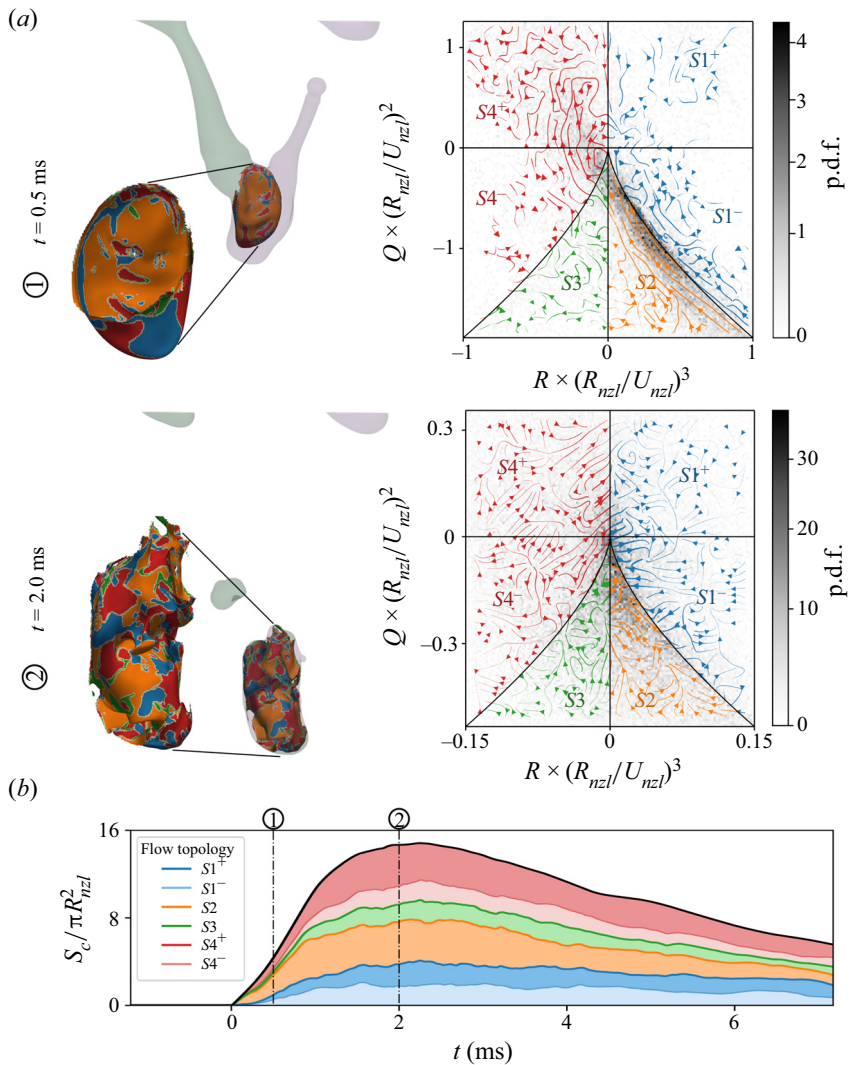


Figure 8. Topological analysis at the contact surface for an asymmetric collision ($\Delta U_{nzl} = 0.24 \text{ m s}^{-1}$, $\Delta t_0 = T_{pulse}/4$, $\Delta\theta = -4^\circ$ and $\Delta\varphi = 6^\circ$). See figure 6 for details.

Figure 10 presents a comparison of the system at different injection times Δt_0 . By changing Δt_0 exclusively, all the droplets in the simulations are guaranteed to have identical mass, initial kinetic and interfacial energy and total linear momentum, thus changing only the impact parameter. Figure 10(a) shows the evolution of the segregation parameter for comparison. The best mixing occurs at $B = 0.2$, in contrast to the worst of the four cases that corresponds to $B = 0$, the symmetric collision, and intermediate mixing for both $B = 0.033$ and $B = 0.44$, which shows a similar trend. As the symmetry is broken, the distortions to the morphology of the contact surface are more pronounced. From figure 10(b), it can be observed that the contact surface area detaches from the plane $x = 0$ and buckles against the delayed droplet (green isosurface). Figure 10(c) displays the profile of the scalar-turbulence interaction term at the contact surface, which, according to (3.6), contributes to the growth rate of its area. Among the four cases, the symmetric

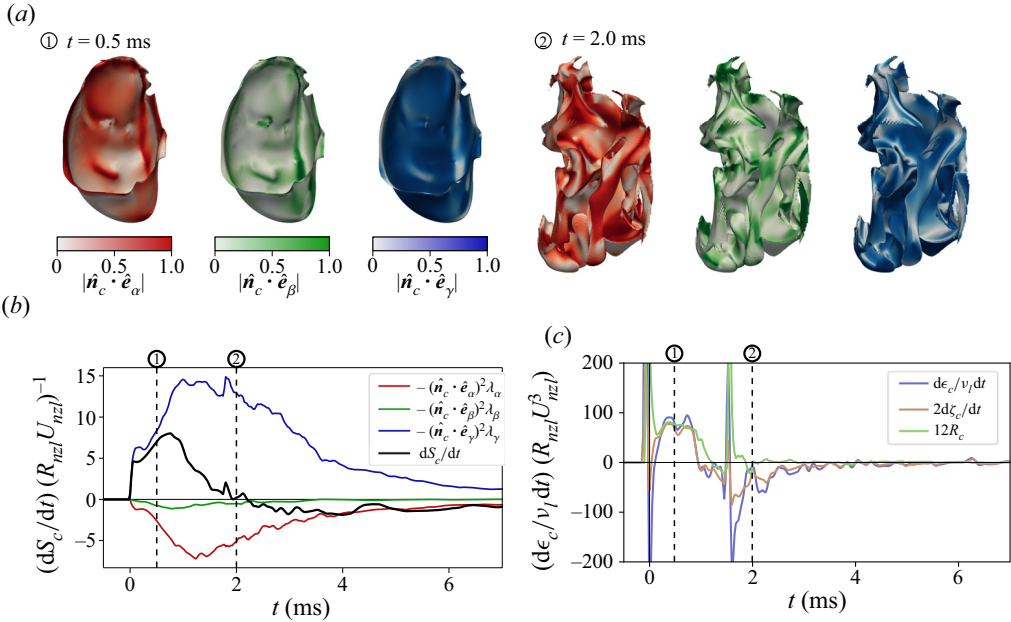


Figure 9. Eigenvector analysis at the contact surface for an asymmetric collision ($\Delta U_{nzl} = 0.24 \text{ m s}^{-1}$, $\Delta t_0 = T_{pulse}/4$, $\Delta\theta = -4^\circ$ and $\Delta\varphi = 6^\circ$). The figure layout is the same as in figure 7.

case shows the largest values of negative growth rate concentrated at the rim of the contact surface. This implies that the contact surface grows from the inside but is limited by its periphery. Therefore, as the symmetry is broken, the contact surface can now grow by buckling. However, if the impact parameter is too high, which implies that the centre of mass of the droplets is further apart, then the droplets have less interaction and their overall contact is reduced, therefore resulting in poor mixing.

After these observations, we can comment on the $\Delta\varphi \neq 0$ cases. Breaking the symmetry of the collision by a lateral misalignment transforms the initial kinetic energy of the droplets into rotational energy. As a consequence, the rotational energy is unable to produce the shear motion at the contact surface that is required to enhance mixing.

5. Conclusions

Three-dimensional direct numerical simulations based on the volume of fluid method have been conducted to analyse the mixing process for binary droplet collisions in a configuration that is representative of an industrial process such as the ReJI bioprinting technique. The quality of mixing has been quantified in terms of the segregation parameter χ , which, for a closed system, is a monotonically decreasing function of time, reflecting the irreversibility of the process.

The variation of the segregation parameter at different configurations of the set-up indicates that introducing a small asymmetry with respect to the mirror-symmetric collision can improve the quality of mixing. In order to quantify the asymmetry, a new generalised definition of the impact parameter B has been proposed for the analysis of the collision of non-spherical droplets. It was found that for most cases, the relation between χ and B follows a master curve. Therefore, the quality of mixing can be predicted by utilising

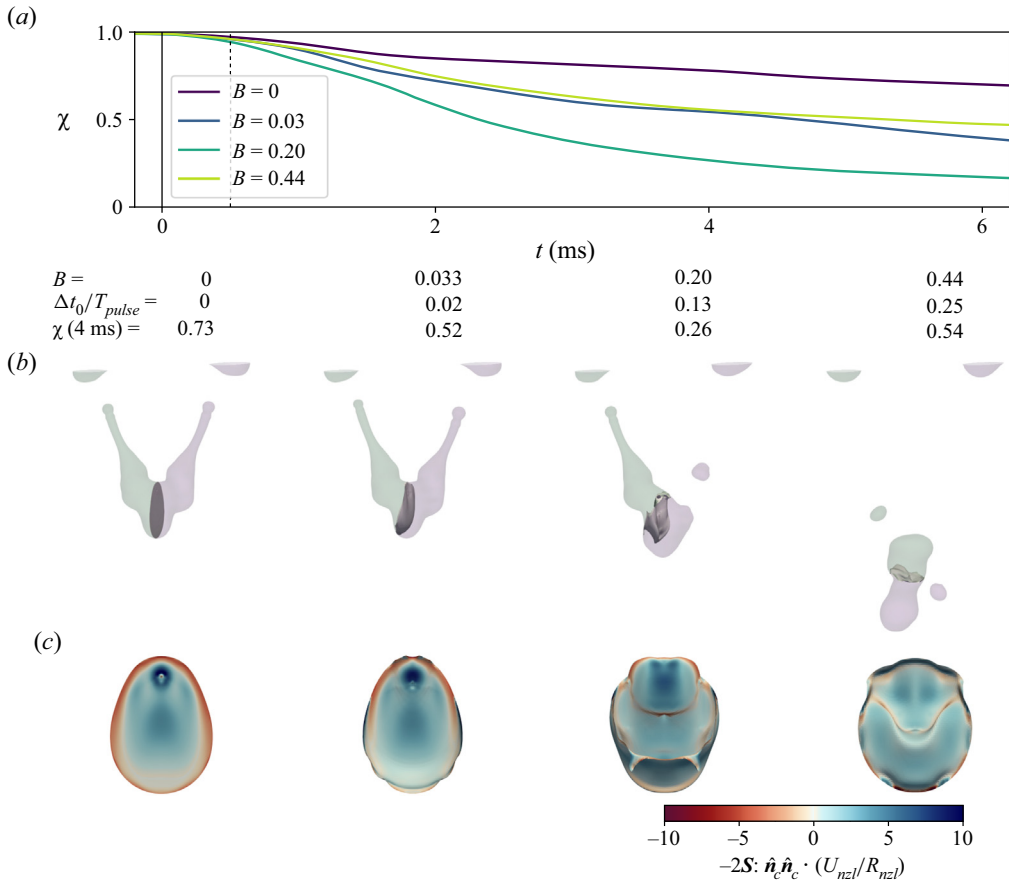


Figure 10. Comparison of a binary droplet collision at different injection times Δt_0 . (a) Segregation parameter as a function of time. The dashed line corresponds to $t = 0.5$ ms. (b) Stills of the binary droplet collision at $t = 0.5$ ms, showing the isosurface $\alpha_1 = 1/2$ in grey. (c) Profile of the scalar–turbulence interaction term at the contact surface.

such a trend. The simulation data have been used to evaluate the parameters of the master curve indicating the relation between χ and B employing least squares optimisation.

The physical explanation for the improvement of the mixing process resulting from the breaking of symmetry has been discussed. The topology analysis of the flow and an eigensystem decomposition of the velocity gradient tensor have been utilised to elucidate how the convective flows and the stretching and twisting of the contact surface between the two droplets affect the mixing process.

It was found that for a symmetric collision, the mixing process is hampered by the plane of symmetry itself, which prohibits convective flows in the region where the diffusive flux is strongest. Therefore, the mixing process is enhanced by the growth of the contact surface area. By breaking the symmetry of the collision, vortical structures that appear naturally now carry matter and stretch the area of contact between the liquids. This can be realised similarly at different combinations of the configuration parameters, therefore increasing the quality of mixing to a nearly optimal level by varying different parameters at the same time. However, if the asymmetry is too large by off-centring the collision, then mixing

becomes inefficient. It was found that the optimal mixing occurs at a moderate asymmetry in the system when the impact parameter is $B \sim 0.2$.

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