Numerical simulation of spreading drops

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**Abstract**

We consider a liquid drop that spreads on a wettable surface. Different time evolutions have been observed for the base radius $r$ depending on the relative role played by inertia, viscosity, surface tension and the wetting condition. Numerical simulations were performed to discuss the relative effect of these parameters on the spreading described by the evolution of the base radius $r(t)$ and the spreading time $t_s$. Different power law evolutions $r(t) \propto t^n$ have been observed when varying the parameters. At the early stage of the spreading, the power law $t^{1/2}$ ($n = 1/2$) is observed as long as capillarity is balanced by inertia at the contact line. When increasing the viscosity contribution, the exponent $n$ is found to increase despite the increase of the spreading time. The effect of the surface wettability is observed for liquids more viscous than water. For a small contact angle, the power law $t^{1/2}$ is then followed by the famous Tanner law $t^{1/10}$ once the drop shape has reached a spherical cap.

**Keywords:**
Drop spreading, Numerical simulation

1. Introduction

When a liquid drop contacts a wettable surface, the liquid spreads over the solid to minimize the total surface energy. The first moments of spreading tend to be rapid (some milli-seconds for millimeter sized water drops) while it takes a much longer time for viscous liquids to completely spread. From the literature [22,17,9,1,6,3,2,24,7], it appears that several time evolutions can be observed for the base radius depending if inertia, viscosity and/or gravity are involved. In addition, the effect of the substrate wettability on drop spreading is not yet clearly understood.

The evolution of the wetted surface is reported using either the base radius $r(t)$ or the wetted surface $A(t) = \pi r^2(t)$. The experimental results are usually presented in order to provide a power law evolution under the form:

$$r(t) \sim t^n, \text{ or } A(t) \sim t^{2n}. \quad (1)$$

The most popular law for drop spreading is certainly the so-called Tanner law [23,22] with $n = 1/10$. This evolution can be simply obtained by considering that the shape of the drop is a thin spherical cap and that the contact angle follows the Cox-Voinov relation. Such evolution has been found to be in agreement with several
2.1. Numerical procedure

2.1.1. Navier Stokes solver

The numerical simulations reported in this work are performed with the volume of fluid (VoF) solver developed in the JADIM code [4,12]. The one-fluid system of equations is obtained by introducing the one-fluid function $C$ used to localize one of the two phases. In this study, we define $C = 1$ in the drop (phase 1), and $C = 0$ in the surrounding air (phase 2). The one-fluid function $C$ makes possible the definition of the one fluid variables $U = C U_1 + (1 - C) U_2$ for the velocity, $P = C P_1 + (1 - C) P_2$ for the pressure, $ρ = C ρ_1 + (1 - C) ρ_2$ for the density, and $μ = C μ_1 + (1 - C) μ_2$ for the viscosity. The position of the interface is then given by the transport equation:

$$\frac{∂C}{∂t} + U ∙ ∇C = 0$$

The two fluids are assumed to be Newtonian and incompressible with no phase change. Under isothermal condition and in the absence of any surfactant, the surface tension is constant and uniform at the interface between the two fluids. In such case, the velocity field $U$ and the pressure $P$ satisfy the classical one-fluid formulation of the Navier–Stokes equations:

$$\nabla U = 0$$

$$\rho \left( \frac{∂U}{∂t} + U ∙ ∇U \right) = -∇P + μ ∙ ∇^2 U + F_σ$$

where $\Sigma$ is the viscous stress tensor, $g$ is the gravity, $F_σ$ is the capillary contribution:

$$F_σ = σ \nabla ∙ n \delta_1$$

where $σ$ is the surface tension, $n$ denotes by arbitrary choice the unit normal of the interface going out from the drop and $δ_1$ is the Dirac distribution associated to the interface.

The system of Eqs. (3)–(6) is discretized using the finite volume method. Time advancement is achieved through a third-order Runge-Kutta method for the viscous stresses. Incompressibility is satisfied at the end of each time step though a projection method. The overall algorithm is second-order accurate in both time and space. The volume fraction $C$ and the pressure $P$ are volume-centered and the velocity components are face-centered. Due to the discretization of $C$, it results a numerical thickness of the interface, the cells cut by the interface corresponding to $0 < C < 1$. The specific aspect of our approach compared to the classical VoF or Level Set methods [19–21] concerns the technique used to control the stiffness of the interface. In our approach no interface reconstruction or redistancing techniques are introduced. The interface location and stiffness are both controlled by an accurate transport algorithm based on FCT (Flux-Corrected-Transport) schemes [25]. This method leads to an interface thickness of about three grid cells due to the implementation of a specific procedure for the velocity used to transport $C$, in flow region of strong strain and shear [4]. The numerical description of the surface tension is one of the crucial points when considering systems where capillary effects control the interface shape. This interfacial force is solved using the classical CSF (continuum surface force) model developed by [5]. A classical problem of this formulation is the generation of spurious currents [13,18]. In order to decrease spurious currents intensity, a classical method introduced by [5] was employed, which consists in calculating the surface curvature from a smoothed density gradient, while the discretization of the delta function uses an un-smoothed density. The spurious currents in our code JADIM have been characterized by [12]. Their maximum magnitude evolves as 0.004σ/μ, in agreement with other codes using the Brackbill’s formulation.

2.2. Numerical modeling of the contact angle

The modeling used in this work aims to respect the physics at the contact line. It is based on two main considerations at the macroscopic scale:
The wall condition seen by the interface is a non-slip condition.

- There is a direct relation between the interface velocity and the shape of the interface characterized by the apparent contact angle.

The objective is to solve the physics at the apparent scale $L$ where the contact angle is the apparent or dynamic contact angle. At this scale the fluid obeys a non-slip condition for the resolution of the momentum equation:

$$U_W = 0$$

Hydrodynamics at a smaller scale than $L$ has an important role since it links the apparent angle to the static angle. A sub-grid modeling of such effects needs to be introduced in the simulation. This is done by using the relation derived by Cox [8] in order to link the apparent angle to the static angle:

$$g(\theta_W) = g(\theta_S) + \text{Ca} \log \left( \frac{L}{\lambda} \right)$$

where $\text{Ca} = \mu U_0 / \sigma$ is the contact line capillary number. The first volume of fluid node being located at $\Delta/2$ from the wall, it follows that $L = \Delta/2$. Considering the value of the observed slip length [14], we have used a fixed value $\lambda = 10^{-5}$ m for the slip length in (4) in the simulations reported in this work. Note that experimental observations give also a recommendation for the choice of the macroscopic length $L$ consistent with the dynamic model used in our numerical modeling. Indeed one should have $\Delta/2 = L = 10^{-4}$ to $10^{-5}$ m.

The comparison between our modeling with other variant modelings, i.e. static contact angle versus dynamic contact angle, no slip condition versus slip condition, is presented in [16]. A grid and time convergence has been performed and discussed for the different possible models. These comparisons clearly stress the strong effect of the numerical modeling used for the simulation of dynamic contact lines as illustrated in the next section, where the simulations are compared with experiments for water and more viscous liquids.

### 2.3. Comparison with experiments for water and viscous drops

We consider a spherical droplet of volume $V = 4\pi r^3/3$ deposited on an horizontal wall that enters in contact with an horizontal wettable surface at time $t = 0$. All the numerical experiments reported in this work are described by Fig. 1. The computational domain is a square domain of equal radial and vertical extension.

$$L_r = L_z = 3R_0$$

The axis of symmetry corresponds to the west boundary, the wetting conditions given by eq. (7-4) are imposed on the south boundary and classical wall conditions (no slip) are imposed on the two other boundaries (east and north). The initial conditions correspond to a drop initially located at the distance $h < R_0$ from the wall. The effect of this initial condition as well as the grid and time convergences have been performed [16] by considering different initial locations of the drop center ($h = 0.95, 0.98, 0.99, 0.995, 0.998 R_0$), different grids (corresponding to 32, 64, 128 and 167 nodes per radius) and different time steps ($\Delta t = 10^{-5}$ to $\Delta t = 10^{-8}$ s).

Squalane and water drops are considered for comparison with the experiments of Lavi & Marmur [15] Winkels et al. and [24], respectively.

We first compare the experiments of [15]. They have considered the spreading of squalane drops with $R_0 = 1$ mm. The fluid properties are $\rho = 809$ kg/m$^3$, $\mu = 0.034$ Pa s and $\sigma = 0.032$ N/m. The comparison between our simulations and the experiments is shown in Fig. 2. The normalized radius $r/R_0$ is reported versus the normalized time $\tau = \sigma t / \mu V^{1/2}$. The dynamic model is compared with the static model (a constant contact angle $\theta_S$ being imposed during all the simulation). A significant difference is observed between the dynamic model (4) and the static model ($\theta_W = \theta_S$). The dynamic models give a satisfactory agreement with the experiments of [15].

We now consider the spreading of water drops and we compare our simulations with the experiments performed by [24] for millimeter-sized water droplets in air ($R_0 = 0.5$ mm). We have selected the cases corresponding to very different contact angles $\theta_S = 115^\circ$ and $\theta_S = 0^\circ$.

Based on preliminary tests [16], the comparison with the experiments performed by [24] are performed with a grid corresponding to 128 nodes per radius $R_0$ and a time step $\Delta t = 10^{-7}$ s. The initial position is $h = 0.998 R_0$. The comparison is shown in Fig. 3 where the normalized contact line radial position $r/R_0$ is shown as function of the time normalized by the capillary-inertial time $t_0 = (\rho R_0^3 / \sigma)^{1/2}$. The agreement is very satisfactory for the two cases considered. In particular, the power law $t^{1/2}$ is clearly reproduced by the simulations. Consequently, our simulations do not show a noticeable effect of the surface wettability on the power law evolution in agreement with [24] but in opposition with [2]. Below, we consider in detail this point with additional simulations performed for a fluid ten times more viscous than water. Note also that our simulation reveals the oscillation of the contact line at the end of the
spreading. Such oscillation are not reported in the experiments of [24] and will be discussed in the following.

3. Results

We consider the effect of the main parameters controlling drop spreading: fluid properties (liquid density, liquid viscosity and surface tension), the drop size and the wetting properties of the surface. Each parameter is varied independently over a wide range of variation. All the simulations have been performed considering air as the surrounding fluid. As a consequence the density and viscosity ratio values (i.e. $\rho_2/\rho_1$ and $\mu_2/\mu_1$) are not constant but they always remain much smaller than unity so that there is no expected effect of both the viscosity and the density of the air on the spreading.

3.1. Inertial-capillary regime of spreading

We first consider low-viscous spreading. The density, the surface tension and the drop radius are varied independently while the viscosity is set fixed to the viscosity of water $\mu = 0.001$ Pa s.

**Fig. 4.** Effect of the liquid density on the drop spreading for $\sigma = 0.072$ N/m, $\mu = 0.001$ Pa s, $\theta_2 = 65^\circ$ and $R_0 = 0.5$ mm. The normalized contact line radius $r' \equiv (r - r_0)/(l_1 - l_0)$ is reported as a function of $t$ (top) and $t$ (bottom) the normalized time $t/(\rho R_0^2/\sigma)^{1/2}$.

The appearance of a capillary wave that travels from the contact line to the top of the droplet has also been reported in the simulations reported by [11] using a diffuse interface method. This figure reports the evolution of the shape of a water drop for $\theta_2 = 65^\circ$. The sequence is presented line by line for the times given in the legend. The time steps are non regular in order to show particular events observed during the spreading. At the beginning of the spreading, only the contact line is moving. The top of the drop starts to move due to the capillary wave induced by the initial impulse generated at the contact line. Surprisingly this capillary wave induces the elevation of the top of the drop. As a consequence, combined with the displacement of the contact line, the drop is elongated into two orthogonal directions resulting in the generation of a pinch event (see 2nd image ($t = 2.52$ ms), line 3). In this case, the pinch-off does not occurs resulting in the propagation of a relatively large amplitude wave. This capillary wave displaces again the top of the drop in the direction opposite to the wall. This generates a second pinch that results in the ejection of a small satellite droplet (see 2nd image ($t = 3.34$ ms), line 5). Then, the satellite droplet impacts and coalesces with the drop generating again wave oscillations, explaining the final oscillations observed for the contact line in **Fig. 4.** Such regime of droplet ejection following the deposition of drops onto
Fig. 5. Shape evolution of a water drop (radial section) for $\theta_s = 65^\circ$. The sequence is shown line by line for the times $t = 0.06, 0.36, 0.56, 1.00, 1.40, 1.86, 2.52, 2.70, 2.86, 2.98, 3.12, 3.28, 3.34, 3.52, 3.60, 4.80$ and $5.80$ ms.

A solid substrate has been identified in experiments and direct numerical simulations by [10] for smaller contact angles ($\theta_s \leq 65^\circ$). We show here that the droplet ejection can also be observed for larger contact angles.

We illustrate in Fig. 6 the effect of the density on the droplet ejection. The simulations are shown for three densities $\rho = 700$ kg/m$^3$, $\rho = 1000$ kg/m$^3$ and $\rho = 4000$ kg/m$^3$. The images corresponding to the two successive pinch-offs and to the maximum height of the satellite droplet are shown. For $\rho = 700$ kg/m$^3$ two pinch events are observed but no droplet ejection is generated by the surface oscillation, while for $\rho = 4000$ kg/m$^3$ the droplet is ejected at the first pinch event.

The effect of the surface tension is now reported in Fig. 7. The surface tension is varied from $\sigma = 0.007$ N/m to $\sigma = 0.2$ N/m. The other parameters are fixed to $\rho = 10^3$ kg/m$^3$, $\mu = 0.001$ Pa s, $R_0 = 0.5$ mm and $\theta_s = 65^\circ$. The spreading time is decreased when increasing the surface tension, since the initial impulse transmitted to the drop is increased. We also observe that the power law
evolution $t^{1/2}$ is not significantly affected by the change of the surface tension. Again the use of the normalized time $t/(ρR_0^2/σ)^{1/2}$ makes possible the collapse of the curves and clearly reveals the capillary oscillations that impact the contact line spreading.

We report in Fig. 8 the spreading time $t_{s}$ for the contact line evolution shown in Figs. 4 and 7. The spreading times deduced from simulations for water drops with radii $R_0$=0.5 mm and $R_0$=5 mm are also reported. $t_{s}$ is shown as a function of the characteristic capillary-inertial time $t_{s} = (ρR_0^2/σ)^{1/2}$. $t_{s}$ is defined as the time necessary to reach 98% of the final equilibrium radius $R_f$ of the contact line. The plot confirms that spreading is clearly controlled by the balance between inertia and capillarity. The spreading time can be estimated as:

$$t_{s} \approx 1.7 \left( \frac{ρR_0^2}{σ} \right)^{1/2}$$

The spreading time is thus increased when reducing the surface tension and increasing the drop density and the drop size.

3.2. Viscosity effect

We report in Fig. 9 the effect of the viscosity on the drop spreading. The viscosity is varied from $μ = 0.001$ Pa s to $μ = 0.07$ Pa s for $ρ = 10^3$ kg/m$^3$, $σ = 0.072$ N/m, $R_0 = 0.5$ mm and $θ_s = 65^\circ$. The spreading time is enlarged when increasing the viscosity (the curves being translated to the right) and the oscillations observed at the end of the spreading due to the propagation of capillary wave are damped. Despite the increase of the spreading time, the slope of the evolution at the early stage of the spreading increases from $t^{1/2}$ (water) to $t^{2/3}$ for the largest viscosity considered. But the duration of the power law regime is reduced because the end of the spreading is characterized by an exponential evolution of the form given by (2).

For comparison, the Tanner law $t^{1/10}$ is also shown in the figure.

There is no clear evidence that this type of evolution is observed in our simulations during the final stage of the spreading. In fact, the static contact angle $θ_s = 65^\circ$ is too large to observe such a regime. For liquids more viscous than water, we expect that the evolution is controlled by a viscous–capillary balance. The viscous stress near the contact line can be estimated as $−μ(δr/δt)/ℓ$ where $ℓ$ is the characteristic length controlling the shear stress at the contact line. Assuming an evolution of the form $ℓ = R_0^{1−p}r^p$, the balance between the viscous stress and the capillary pressure $−σR_0/r^2$ gives, after integration, the observed $t^{2/3}$ evolution:

$$r(t) \sim \left( \frac{σR_0^2}{μ} \right)^{2/23} t^{2/3}$$

for $p = 3/2$ corresponding to $ℓ \sim R_0^{1/3}t^{1/2}$. The corresponding viscous–capillary characteristic time $t_{μ} = R_0/σ$ is used in Fig. 4 (bottom) to normalize the time. For a viscosity larger than 0.01 Pa s, a satisfactory collapse of the curves is observed indicating the transition to a spreading regime controlled by the balance between capillarity and viscosity.

The evolution of the spreading time $t_s$ is reported in Fig. 10 as a function of the viscous–capillary characteristic time $t_{μ} = μR_0/σ$. Two different behaviors are clearly observed. For low viscosity, $t_s$ is found to evolve as $t_s^{4/5}$ (i.e. $μ^{4/5}$ since all the other parameters are not varying). We have no explanation of the particular evolution $t_s \propto μ^{4/5}$ revealed by the simulation for low-viscous drops. For larger viscosities an expected linear evolution $t_s \propto t_{μ}$ is observed. The linear behavior results from the linearity of the governing equations (Stokes equation) at low Reynolds number. The viscous–capillary spreading time can be estimated using the relation

$$t_s \approx \frac{55 μR_0}{σ}$$
3.3. Contact angle effect

We finally discuss the effect of the static contact angle $\theta_5$ on drop spreading. Figs. 3, 4 and 7 show that the $t^{1/2}$ power law is not affected by the surface wettability at the beginning of the spreading, when considering a fluid having the same viscosity as water for a large range of contact angles ($\theta_5 = 5^\circ$, $65^\circ$ and $115^\circ$), in agreement with the experiments reported by [24] for water drops. The wetting conditions considered by [24] ($\theta_5 = 5^\circ$, $65^\circ$ and $115^\circ$) are now simulated for a fluid with a viscosity ten times larger than that of water ($\mu = 0.01$ Pa s), the other parameters being kept the same as for water ($\sigma = 0.072$ N/m, $\rho = 10^{3}$ kg/m$^3$) and $R_0 = 0.5$ mm. Additional wetting conditions ($\theta_5 = 140^\circ$ and $\theta_5 = 160^\circ$) have also been considered.

Fig. 11 clearly reveals that the wetting condition has now a significant effect on the spreading rate. When the contact angle is increased the slope seems to decrease in agreement with the experiments of [2]. However, if the radius $r$ is normalized using the radius reached at the end of the inertial-capillary regime, all the curves collapse on a single curve showing a $t^{1/2}$ evolution independent on the contact angle $\theta_5$ in agreement with the experiments of [24].
the $t^{1/2}$ evolution, during the $t^{1/10}$ evolution and at the transition between these two regimes. The corresponding times are depicted by bullets in Fig. 11. A spherical cap is clearly identified during the $t^{1/10}$ evolution.

In the limit of a very thin spherical cap, the contact angle $\theta_W$, the drop volume $V$ and the contact line radius $r$ are linked by the geometrical relation $\theta_W \approx 4V/\pi r^2$. In the limit of a small dynamic contact angle with $\theta_S \approx 0$, the Cox relation (4) can be expressed as:

$$\theta_W^2 = g \frac{\mu}{\sigma} \frac{dr}{dt} \ln(L/\lambda)$$

Combining these two relations and integrating in time, the radius of the contact line is found to follow the famous Tanner law:

$$r(t) \sim R_0 \left( \frac{t}{t_\mu} \right)^{1/10} \quad (13)$$

For spreading drops, such evolution is thus observed for small contact angles once the drop has reached a spherical cap shape as shown in Fig. 12c. In the limit of $\theta_S \to 0$, the base radius of the spherical cap can be expressed as $r \sim R_0 \theta_S^{1/3}$ and it follows an estimation of the spreading time:

$$t_S \approx t_\mu \theta_S^{-10/3} \quad (14)$$

For the simulations reported in Fig. 11, relation (14) gives $t_S \approx 0.2$ s which is a correct order of magnitude for $\theta_S = 5^\circ$. The corresponding viscous–capillary spreading time given by relation (12) is two orders of magnitude smaller = 0.0038 s, but in agreement with the evolution reported for the larger contact angles ($\theta_S = 65^\circ, 115^\circ, 140^\circ$ and $160^\circ$).

4. Conclusion

We have considered the spreading of a drop on a horizontal wettable surface by numerical simulation. In our numerical modeling of moving contact lines, the apparent contact angle is given by the [8] relation while a zero slip length is imposed for the momentum calculation. The simulations have been satisfactorily compared with the experiments of [15] for squalane and with the experiments of [24] for water. The effects of the main parameters controlling the drop spreading have been studied separately. For viscosity of the order of that of water, the $t^{1/2}$ power law has been recovered at the early stage of the spreading for all the contact angles considered in agreement with the experiments of [24]. When increasing the viscosity, the early stage of the spreading is still characterized by a power law evolution but with an increasing exponent in agreement with the experiments of [2]. Both the viscosity and the contact angle are then shown to impact the line expansion. In particular, considering the spreading of a viscous drop on a highly wettable substance ($\theta_S = 5^\circ$), we observe two successive power law evolutions. The capillary-inertial regime $r \propto t^{1/2}$ is followed by the

![Fig. 10. Spreading time $t_S$ versus the viscous–capillary characteristic time $t_\mu = \mu R_0/\sigma$ for $\sigma = 0.072$ N/m, $\rho = 10^3$ kg/m$^3$, $R_0 = 0.5$ mm and $\theta_S = 65^\circ$. (---) $t_S \propto t_\mu^{1/3}$. (---) $t_S \propto t_\mu$.](image)

![Fig. 11. Effect of the contact angles $\theta_S$ on the time evolution of the drop base radius $r$ for $\mu = 0.01$ Pa s, $\sigma = 0.072$ N/m, $\rho = 10^3$ kg/m$^3$ and $R_0 = 0.5$ mm. The drop radial sections corresponding to a, b and c are shown in Fig. 12.](image)

For the smallest contact angle $\theta_S = 5^\circ$, we observe after the $t^{1/2}$ evolution a clear $t^{1/10}$ evolution characteristic of the Tanner Law. A similar behavior has been reported in the simulations of [11] where a regime close to the Tanner’s law is observed when increasing the viscosity while a regime close to $t^{1/2}$ is observed at the early stage of the spreading. Fig. 12 shows the shape of the drop during

![Fig. 12. Drop radial section corresponding to the times a, b and c shown in Fig. 11 ($\theta_S = 5^\circ$). (a) in the middle of the $t^{1/2}$ evolution ($t = 4 \times 10^{-4}$ s), (b) the transition between the $t^{1/2}$ and the $t^{1/10}$ evolution ($t = 3.2 \times 10^{-3}$ s) and (c) in the middle of the Tanner law evolution ($t = 10^{-2}$).](image)
Tanner $r \propto t^{1/10}$ evolution once the droplet has reached a spherical cap shape. Our simulations show that the combined effect of the viscosity and the contact angle plays an important role in the spreading of a drop. They make consistent the different experimental observations (a priori contradictory) concerning the effect of the wetting conditions.

References