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A spectral boundary integral method for simulating electrohydrodynamic flows in viscous drops

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ABSTRACT

A weakly conducting liquid droplet immersed in another leaky dielectric liquid can exhibit rich dynamical behaviors under the effect of an applied electric field. Depending on material properties and field strength, the nonlinear coupling of interfacial charge transport and fluid flow can trigger electrohydrodynamic instabilities that lead to shape deformations and complex dynamics. We present a spectral boundary integral method to simulate droplet electrohydrodynamics in a uniform electric field. All physical variables, such as drop shape and interfacial charge density, are represented using spherical harmonic expansions. In addition to its exponential accuracy, the spectral representation affords a nondissipative dealiasing method required for numerical stability. A comprehensive charge transport model, valid under a wide range of electric field strengths, accounts for charge relaxation, Ohmic conduction, and surface charge convection by the flow. A shape reparametrization technique enables the exploration of significant droplet deformation regimes. For low-viscosity drops, the convection by the flow drives steep interfacial charge gradients near the drop equator. This introduces numerical ringing artifacts that we treat via a weighted spherical harmonic expansion, resulting in solution convergence. The method and simulations are validated against experimental data and analytical predictions in the axisymmetric Taylor and Quincke electroration regimes.

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1. Introduction

A wide range of engineering applications involve liquid drops immersed in another fluid while subject to an applied electric field. Some examples include ink-jet printing [1], electrospraying [2], and microfluidic devices and pumps [3]. These systems exhibit rich dynamics due to the electric field and fluid flow coupling. When an interface separating two immiscible fluids is subject to an otherwise uniform electric field, the electric field undergoes a jump across the interface due to the mismatch in material properties. This discontinuity in the electric field induces electric stresses that can deform the interface and drive the fluid into motion.

We are interested in leaky dielectric liquids such as oils, which serve as poor conductors. Unlike electrolyte solutions where diffuse Debye layers affect the system's dynamics, leaky dielectrics are characterized by the absence of diffuse Debye layers [4]. The free charges instead concentrate on the interfaces between different phases in the system. Consequently, the electric field acting on the interfacial charge creates electric stresses along the normal and tangential directions, which

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cause deformations and fluid motion. Surface tension has a stabilizing effect in general, trying to restore the equilibrium shape. Melcher and Taylor developed a framework for studying electrohydrodynamic phenomena in leaky dielectric systems, known as the leaky dielectric model (LDM) [5]. Central to their work is a charge conservation model that describes a balance between Ohmic fluxes from the bulk, interfacial charge convection, and finite charge relaxation. It was previously shown that LDM can be derived asymptotically from electrokinetic models in the limit of strong electric fields and thin Debye layers [6,7].

This work focuses on the dynamics of a leaky dielectric drop immersed in another dielectric fluid under a uniform DC electric field. This canonical problem has been a long-standing research problem in electrohydrodynamics. In his pioneering work, Taylor [8] formulated a small-deformation theory for an isolated drop based on LDM and could predict oblate and prolate steady shapes depending on the material properties. While Taylor's theory shows good agreement with experimental data in the limit of vanishing electric capillary number Ca_E (ratio of electric to capillary forces), the discrepancy is significant at larger values of Ca_E . Therefore, other researchers attempted to extend Taylor's work by accounting for second-order effects in Ca_E [9], considering spheroidal drops [10,11], including inertial effects [12] and interfacial charge convection [13–16].

A variety of computational models have been developed to study drop dynamics under strong electric fields at finite deformations, a problem untractable using analytical theories. In the limit of negligible inertia, boundary integral equations can be used to formulate and solve the coupled electrohydrodynamic problem. Sherwood was the first to develop a boundary element method for an axisymmetric drop in an equiviscous system and applied it to capture breakup modes in prolate drops [17]. His original work was subsequently extended to study drop pair interaction [18] and to cover a wider range of fluid and electric parameters [19]. These earlier attempts used a simplified boundary condition for the electric problem, which neglected transient charge relaxation and interfacial charge convection by the flow. These two effects have recently been shown to play a significant role in drop dynamics and deformations [16]. Lanauze et al. [20] and Das and Saintillan [21] recently addressed this problem and developed axisymmetric and three-dimensional boundary element methods based on the full Melcher–Taylor LDM. The effect of charge convection was specifically addressed in [21], where it was shown to be responsible for Quincke electrorotation. These methods, however, were found to lack accuracy and stability in the regime of strong electric fields. Other numerical approaches have been used to study drop electrohydrodynamics, including immersed boundary [22], level set [23,24], and finite element methods [25–27]. More recently, finite element simulations [28,29] were also used to investigate electrohydrodynamic instabilities such as tip and equatorial streaming in drops under strong electric fields. These latter techniques all include finite fluid inertia and, with few exceptions [24], do not treat the drop surface as a sharp interface.

Improved accuracy within the boundary integral framework can be achieved using spectral methods, which rely on expansions of the shape and interfacial variables based on spherical harmonics. Such methods were recently developed to simulate electrohydrodynamics of lipid vesicles [30] and also extended to the case of individual drops and drop pairs [31–33]. These studies, however, all neglected charge relaxation and charge convection and were thus restricted to weak electric fields. Accurately capturing charge convection is especially challenging as it nonlinearly couples fluid flow and charge transport on a deformed interface. It can result in spurious aliasing errors with negative consequences for accuracy and stability. This work addresses this challenge and presents a spectral boundary integral method for the electrohydrodynamics of deformable liquid drops based on the complete Melcher–Taylor LDM. Interfacial charge convection is rigorously accounted for, and dealiasing and reparametrization techniques are implemented to improve accuracy and stability and enable long-time simulations.

The paper is organized as follows. We define the problem and discuss the governing equations and boundary conditions in Sec. 2.1, along with their non-dimensionalization in Sec. 2.2. Sec. 2.3 presents the integral form of the governing equations and boundary conditions used in developing the boundary integral method. We discuss different aspects of the numerical method in Sec. 3: the spectral representation of all variables in terms of spherical harmonics is discussed in 3.1, followed by details of the dealiasing method in Sec. 3.2. Next, in Sec. 3.3, we summarize the numerical integration methods used in this study and correction methods to ensure charge neutrality and incompressibility in Sec. 3.4. As explained in Sec. 3.5, we also use a reparametrization method to improve the numerical stability in simulations where the drop undergoes significant deformations. In Sec. 3.6, we discuss the overall convergence of our numerical method as well as the limitations on the time step size. We test and validate our computational model by applying it to a wide range of dynamical behaviors, such as the axisymmetric Taylor regime under weak electric fields in Sec. 4.1, and Quincke electrorotation under stronger electric fields in Sec. 4.2. We also investigate the dynamics of low-viscosity drops in Sec. 4.3, where charge convection plays an important role. Finally, we discuss our conclusions and possible extensions of our work in Sec. 5.

2. Problem definition

2.1. Governing equations

We consider a neutrally buoyant drop of a fluid occupying volume V^- immersed in an infinite body of another fluid V^+ while subject to a uniform electric field $\mathbf{E}_\infty = E_\infty \hat{\mathbf{e}}_z$ as depicted schematically in Fig. 1. The interface D separates the two fluid media, and the surface unit normal $\mathbf{n}(\mathbf{x})$ is pointed towards the suspending fluid. Initially, the drop is uncharged and spherical with radius r_0 . The material properties, namely the dielectric permittivities, electric conductivities, and dynamic viscosities, are denoted by $(\epsilon^\pm, \sigma^\pm, \mu^\pm)$ inside and outside the drop, respectively. Under the Taylor–Melcher leaky dielectric

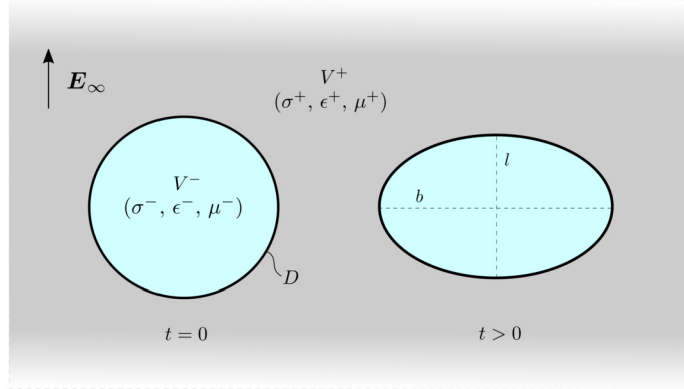


Fig. 1. Problem definition: a leaky dielectric drop with $(\sigma^-, \epsilon^-, \mu^-)$ is suspended in another leaky dielectric fluid with $(\sigma^+, \epsilon^+, \mu^+)$ and subject to an external electric field \mathbf{E}_∞ . The drop deforms and diverges from its initially spherical shape.

model [8], any net charge in the system appears on the interface D , and the bulk of the fluids remain electroneutral. Therefore, the electric potential is harmonic in the bulk:

$$\nabla^2 \varphi^\pm(\mathbf{x}) = 0, \quad \mathbf{x} \in V^\pm. \tag{1}$$

Far away from the interface, the electric field $\mathbf{E} = -\nabla\varphi$ tends to the applied electric field:

$$\mathbf{E}^+ \rightarrow \mathbf{E}_\infty = E_\infty \hat{\mathbf{e}}_z, \quad \text{as } |\mathbf{x}| \rightarrow \pm\infty. \tag{2}$$

While the tangential component of the electric field is continuous across the interface, its normal component undergoes a jump due to the mismatch in material properties:

$$\mathbf{n} \times \llbracket \mathbf{E} \rrbracket = \mathbf{0}, \quad \mathbf{x} \in D. \tag{3}$$

We define the operator $\llbracket \mathcal{F} \rrbracket := \mathcal{F}^+ - \mathcal{F}^-$ as the jump in any variable \mathcal{F} across the interface D . A surface charge density develops at the interface and follows Gauss's law,

$$q(\mathbf{x}) = \mathbf{n} \cdot \llbracket \epsilon \mathbf{E} \rrbracket, \quad \mathbf{x} \in D. \tag{4}$$

The surface charge evolves due to Ohmic currents from the bulk and convective currents on the interface. Consequently, it satisfies the conservation equation:

$$\partial_t q + \mathbf{n} \cdot \llbracket \sigma \mathbf{E} \rrbracket + \nabla_s \cdot (q\mathbf{u}) = 0, \quad \mathbf{x} \in D, \tag{5}$$

where $\nabla_s = (\mathbf{I} - \mathbf{nn}) \cdot \nabla$ is the surface gradient operator and \mathbf{u} is the fluid velocity.

Neglecting the effect of inertia and gravity, the velocity and pressure fields satisfy the Stokes and continuity equations:

$$\mu^\pm \nabla^2 \mathbf{u}^\pm - \nabla p^\pm = \mathbf{0}, \quad \nabla \cdot \mathbf{u}^\pm = 0, \quad \mathbf{x} \in V^\pm. \tag{6}$$

The velocity vector is continuous across the interface and vanishes far from it:

$$\llbracket \mathbf{u}(\mathbf{x}) \rrbracket = \mathbf{0}, \quad \mathbf{x} \in D, \tag{7}$$

$$\mathbf{u}^+(\mathbf{x}) \rightarrow \mathbf{0}, \quad \text{as } |\mathbf{x}| \rightarrow \infty. \tag{8}$$

The balance of interfacial forces requires that the jump in hydrodynamic and electric tractions across the interface balance capillary forces:

$$\llbracket \mathbf{f}^H \rrbracket + \llbracket \mathbf{f}^E \rrbracket = \gamma (\nabla_s \cdot \mathbf{n}) \mathbf{n}, \quad \mathbf{x} \in D. \tag{9}$$

We neglect Marangoni effects due to variations in surface tension, $\nabla_s \gamma = \mathbf{0}$. Hydrodynamic and electric tractions are expressed in terms of the Newtonian and Maxwell stress tensors, respectively:

$$\mathbf{f}^H = \mathbf{n} \cdot \mathbf{T}^H, \quad \mathbf{T}^H = -p\mathbf{I} + \mu(\nabla \mathbf{u} + \nabla \mathbf{u}^T), \tag{10}$$

$$\mathbf{f}^E = \mathbf{n} \cdot \mathbf{T}^E, \quad \mathbf{T}^E = \epsilon \left(\mathbf{E} \mathbf{E} - \frac{1}{2} E^2 \mathbf{I} \right). \tag{11}$$

The jump in electric tractions can be decomposed into tangential and normal components as

$$\llbracket \mathbf{f}^E \rrbracket = \llbracket \epsilon E^n \rrbracket \mathbf{E}^t + \frac{1}{2} \llbracket \epsilon (E^{n2} - E^{t2}) \rrbracket \mathbf{n} = q \mathbf{E}^t + \llbracket p^E \rrbracket \mathbf{n}, \quad (12)$$

where $p^E = \epsilon (E^{n2} - E^{t2})/2$ is the electric pressure [19]. The first term on the right-hand side represents the tangential electric stresses in leaky dielectrics, and it vanishes when both fluids are either perfect dielectrics or perfect conductors.

2.2. Non-dimensionalization

For the system described above, a dimensional analysis yields five dimensionless groups, three of which characterize the mismatch of material properties in the drop and the suspending fluid:

$$R = \frac{\sigma^+}{\sigma^-}, \quad Q = \frac{\epsilon^-}{\epsilon^+}, \quad \lambda = \frac{\mu^-}{\mu^+}. \quad (13)$$

The limits of $\lambda \rightarrow 0$ and ∞ correspond to a bubble and a rigid particle, respectively. The remaining dimensionless groups describe the system's dynamics and can be obtained by comparing the characteristic time scales in the problem. First, note that the response of each fluid phase to Ohmic conduction is characterized by the charge relaxation time:

$$\tau_c^\pm = \frac{\epsilon^\pm}{\sigma^\pm}. \quad (14)$$

The product $RQ = \tau^-/\tau^+$ is the ratio of the charge relaxation times in the two fluids and plays an important role in the dynamics of the drop [21]. The polarization time for a rigid sphere under an applied electric field is the Maxwell–Wagner relaxation time

$$\tau_{MW} = \frac{\epsilon^- + 2\epsilon^+}{\sigma^- + 2\sigma^+} = \frac{R(Q+2)}{1+2R} \tau_c^+, \quad (15)$$

which provides an approximate timescale for polarization of the drop. The accumulation of free charges on the interface creates electric forces that drive the fluid into motion on the electrohydrodynamic time scale

$$\tau_{EHD} = \frac{\mu^+}{\epsilon^+ E_\infty^2}. \quad (16)$$

Deformations away from the equilibrium spherical shape relax under the effect of surface tension on the capillary time scale

$$\tau_\gamma = \frac{\mu^+ r_0}{\gamma}. \quad (17)$$

By taking the ratios of these time scales, the two remaining dimensionless groups can be defined as

$$Ca_E = \frac{\tau_\gamma}{\tau_{EHD}} = \frac{\epsilon_\infty^2 r_0}{\gamma}, \quad Ma = \frac{\tau_{EHD}}{\tau_{MW}} = \frac{\mu^+}{\tau_{MW} \epsilon^+ E_\infty^2}. \quad (18)$$

The electric capillary number Ca_E compares electric forces versus capillary forces, while the Mason number Ma characterizes the importance of charge conduction against surface charge convection. Alternatively, we can combine Ca_E and Ma to construct a third dimensionless group Ca_{MW} to be independent of the electric field:

$$Ca_{MW} = \frac{\tau_\gamma}{\tau_{MW}} = \frac{\mu^+(1+\lambda)r_0}{\gamma \tau_{MW}} = (1+\lambda) Ca_E Ma. \quad (19)$$

For a given set of material properties, changing Ca_{MW} corresponds to varying the drop radius r_0 .

We scale the governing equations and boundary conditions using length scale r_0 , time scale τ_{MW} , pressure scale $\epsilon^+ E_\infty^2$, and the characteristic electric potential $E_\infty r_0$. In the remainder of this manuscript, all governing equations and boundary conditions are dimensionless, and results are presented in terms of the corresponding dimensionless variables.

2.3. Boundary integral formulation

The electric problem is formulated in integral form based on the solution to Laplace's equation as [17–19]

$$\varphi(\mathbf{x}_0) = -\mathbf{x}_0 \cdot \mathbf{E}_\infty - \int_D \mathbf{n} \cdot \llbracket \nabla \varphi(\mathbf{x}) \rrbracket \mathcal{G}(\mathbf{x}_0; \mathbf{x}) ds(\mathbf{x}), \quad \text{for } \mathbf{x}_0 \in V^\pm, D, \quad (20)$$

where the evaluation point \mathbf{x}_0 can be anywhere in space, and \mathbf{x} denotes the integration point on the interface. The free-space Green's function for Laplace's equation, $\mathcal{G}(\mathbf{x}_0; \mathbf{x})$, captures the electric potential due to a point charge in an unbounded domain as

$$\mathcal{G}(\mathbf{x}_0; \mathbf{x}) = \frac{1}{4\pi r}, \quad \text{where } \mathbf{r} = \mathbf{x}_0 - \mathbf{x}, r = |\mathbf{r}|. \tag{21}$$

Taking the gradient of Eq. (20) with respect to \mathbf{x}_0 and using Gauss's law (4), we derive an integral equation for the jump in the normal electric field as a function of the surface charge distribution:

$$\oint_D \llbracket E^n(\mathbf{x}) \rrbracket [\mathbf{n}(\mathbf{x}_0) \cdot \nabla_0 \mathcal{G}] ds(\mathbf{x}) - \frac{1+Q}{2(1-Q)} \llbracket E^n(\mathbf{x}_0) \rrbracket = E_\infty^n(\mathbf{x}_0) - \frac{q(\mathbf{x}_0)}{1-Q}, \quad \text{for } \mathbf{x}_0 \in D. \tag{22}$$

For a given charge distribution $q(\mathbf{x})$, Eq. (22) can determine $\llbracket E^n(\mathbf{x}) \rrbracket$, from which E^{n+} and E^{n-} follow as

$$E^{n+}(\mathbf{x}) = \frac{q(\mathbf{x}) - Q \llbracket E^n(\mathbf{x}) \rrbracket}{1-Q}, \quad E^{n-}(\mathbf{x}) = \frac{q(\mathbf{x}) - \llbracket E^n(\mathbf{x}) \rrbracket}{1-Q}. \tag{23}$$

The tangential electric field $\mathbf{E}^t(\mathbf{x}) = -\nabla_s \varphi(\mathbf{x})$ can be computed by differentiation of the electric potential (20) along the tangential direction. The interfacial jump in electric tractions $\llbracket \mathbf{f}^E \rrbracket$ follows from Eq. (12) based on the tangential and normal electric fields calculated at every point on the interface. This can be used to determine the jump in hydrodynamic tractions $\llbracket \mathbf{f}^H \rrbracket$ using the dynamic boundary condition (9) as

$$\llbracket \mathbf{f}^H \rrbracket = -\llbracket \mathbf{f}^E \rrbracket + \text{Ca}_E^{-1} (\nabla_s \cdot \mathbf{n}) \mathbf{n}, \tag{24}$$

which enters the calculation of the velocity field, as explained next.

The flow problem is also recast into a boundary integral form as [34,35]

$$\begin{aligned} \mathbf{u}(\mathbf{x}_0) = & -\frac{1}{4\pi \text{Ma} (1+\lambda)} \oint_D \llbracket \mathbf{f}^H(\mathbf{x}) \rrbracket \cdot \mathbf{G}(\mathbf{x}_0; \mathbf{x}) ds(\mathbf{x}) \\ & + \frac{1-\lambda}{4\pi (1+\lambda)} \oint_D \mathbf{u}(\mathbf{x}) \cdot \mathbf{T}(\mathbf{x}_0; \mathbf{x}) \cdot \mathbf{n}(\mathbf{x}) ds(\mathbf{x}), \quad \text{for } \mathbf{x}_0 \in D. \end{aligned} \tag{25}$$

Here, \mathbf{G} is the free-space Green's function for the Stokeslet or flow due to a unit point force in an unbounded domain, and \mathbf{T} is the corresponding stress tensor:

$$\mathbf{G}(\mathbf{x}_0; \mathbf{x}) = \frac{\mathbf{I}}{r} + \frac{\mathbf{r}\mathbf{r}}{r^3}, \quad \mathbf{T}(\mathbf{x}_0; \mathbf{x}) = 6 \frac{\mathbf{r}\mathbf{r}\mathbf{r}}{r^5}. \tag{26}$$

Note that the integral equations (20), (22) and (25) exhibit singular behaviors of different orders as \mathbf{x} approaches \mathbf{x}_0 . This is due to the singularity of the Green's function for Laplace's (21) and Stokes equations (26). Conventional quadrature schemes have poor accuracy in the presence of singular integrands and may not converge by increasing the level of discretization. Therefore, accurate numerical integration requires special treatment of the singularities, which we will discuss in Sec. 3.3.

3. Numerical methods

We solve Eqs. (22) and (25) by building upon a spectral boundary integral method introduced by Zhao et al. [36]. This method was also previously applied to study vesicle dynamics under shear and extensional flows [37,38] as well as flows of confined red blood cells [39–41]. All variables, including the interfacial shape, velocity, and charge, are represented using truncated series of spherical harmonic expansions as discussed in Sec. 3.1. Nonlinear operations, geometrical quantities (such as mean curvature), and spatial derivatives are computed accurately using a nondissipative dealiasing method discussed in Sec. 3.2. The boundary integrals are computed using a quadrature scheme for the surface collocation points, with a special treatment for the singular integrands, as summarized in Sec. 3.3. Any changes to the volume and net charge of the drop due to numerical errors are corrected as explained in Sec. 3.4 to ensure numerical stability over long simulation times. In addition, we adopt a reparameterization technique in Sec. 3.5 to minimize the high-frequency component of the interfacial shape, which improves the stability of the numerical method for cases with significant deformations. We provide a convergence analysis for our numerical method in Sec. 3.6 and discuss the time step size restrictions for our time marching scheme. The code associated with the methods and simulations in this article is publicly available at https://github.com/mfirouzn/EHD_Drop_3D [42].

At $t = 0$, the drop is uncharged with a spherical shape. The numerical algorithm used in this study follows that of Das and Saintillan [21], Firouznia and Saintillan [43], and Firouznia et al. [44]. We perform the following steps at every time iteration:

1. Given the current charge distribution $q(\mathbf{x})$ and shape of the interface, compute $\llbracket E^n(\mathbf{x}) \rrbracket$ by numerically inverting (22) using GMRES [45]. From $\llbracket E^n(\mathbf{x}) \rrbracket$, we obtain $E^{n+}(\mathbf{x})$ and $E^{n-}(\mathbf{x})$ via (23).
2. Determine the potential φ along the interface by evaluating (20).

3. Differentiate the surface potential numerically along the interface in order to obtain the tangential electric field $\mathbf{E}^t = -\nabla_s \phi$.
4. Knowing both components of the electric field, determine the jump in the electric traction $[[\mathbf{f}^E]]$ and use it to obtain $[[\mathbf{f}^H]]$ using (24).
5. Solve for the interfacial velocity using the Stokes boundary integral equation (25).
6. Compute $\partial_t q$ via (5) and update the charge distribution.
7. Update the position of the interface by advecting the grid with the normal component of the interfacial velocity: $\partial_t \mathbf{x} = (\mathbf{u} \cdot \mathbf{n})\mathbf{n}$.
8. Apply corrections to the shape and charge distribution to ensure incompressibility and charge neutrality, as discussed in Sec. 3.4.
9. Reparametrize the interfacial shape following the method discussed in Sec. 3.5 to minimize high-frequency components in the spherical harmonic expansion.

The algorithm above describes a single time-marching step. In practice, we use a second-order Runge-Kutta method as discussed in Sec. 3.6.

3.1. Surface representation

The shape of the drop is assumed to be smooth and of spherical topology. Therefore, the surface is parameterized by a truncated series of spherical harmonic expansion from a rectangular domain $\mathbb{S}^2 = \{(\theta, \phi) | \theta \in (0, \pi), \phi \in [0, 2\pi)\}$ to \mathbb{R}^3 :

$$\mathbf{x}(\theta, \phi) = \sum_{n=0}^{N-1} \sum_{m=0}^n \bar{P}_n^m(\cos \theta) (\mathbf{a}_{nm} \cos m\phi + \mathbf{b}_{nm} \sin m\phi), \tag{27}$$

where θ and ϕ are the latitude and longitude angles, and $\mathbf{s} = \{\mathbf{a}_{nm}, \mathbf{b}_{nm}\}$ are the coefficients of the expansion in a compact form [46]. The representation above yields N^2 spherical harmonic modes per each component of \mathbf{x} (total of $3N^2$ modes). The normalized associated Legendre polynomials of degree n ($n = 0, 1, 2, \dots$) and order m ($m \leq n$) are defined as

$$\bar{P}_n^m(\eta) = \frac{1}{2^n n!} \sqrt{\frac{(2n+1)(n-m)!}{2(n+m)!}} (1-\eta^2)^{m/2} \frac{d^{n+m}}{d\eta^{n+m}} (\eta^2-1)^n, \tag{28}$$

and satisfy the orthogonality condition

$$\int_{-1}^1 \bar{P}_n^m(\eta) \bar{P}_{n'}^m(\eta) d\eta = \delta_{nn'}. \tag{29}$$

Similarly, the surface charge distribution is represented as

$$q(\theta, \phi) = \sum_{n=0}^{N-1} \sum_{m=0}^n \bar{P}_n^m(\cos \theta) (\tilde{a}_{nm} \cos m\phi + \tilde{b}_{nm} \sin m\phi). \tag{30}$$

The rectangular domain \mathbb{S}^2 is discretized based on the roots of the Legendre polynomial $P_N(\cos \theta)$ along θ , and uniformly along ϕ . Forward and backward transformations are performed using the SPHEREPACK library [47,48]. Partial derivatives of a given distribution can be computed using recurrence relations for the derivatives of the associated Legendre polynomials [47,49]. Besides the spectral accuracy, the spherical harmonic representation allows for nondissipative dealiasing, which improves the numerical stability of the simulations [36].

A local coordinate system is constructed at every point $\mathbf{x}(\theta, \phi)$ on the surface of the drop, using two tangent vectors $\mathbf{a}_{1,2}$ and the unit normal \mathbf{a}_3 :

$$\mathbf{a}_1 = \partial_\theta \mathbf{x}, \quad \mathbf{a}_2 = \partial_\phi \mathbf{x}, \quad \mathbf{a}_3 = \mathbf{n} = \frac{\mathbf{a}_1 \times \mathbf{a}_2}{|\mathbf{a}_1 \times \mathbf{a}_2|}. \tag{31}$$

Consequently, the first and second fundamental forms of the drop surface have the following components:

$$L_{ij} = \mathbf{a}_i \cdot \mathbf{a}_j, \quad \text{and} \quad B_{ij} = \mathbf{a}_{i,j} \cdot \mathbf{n}, \quad (i, j = 1, 2), \tag{32}$$

which will be used in the subsequent derivations. Given a scalar function $f(\theta, \phi)$ on the surface D defined by the parameterization introduced in (27), the surface gradient $\nabla_s f$ is

$$\nabla_s f = \left(\frac{L_{22} \mathbf{a}_1 - L_{12} \mathbf{a}_2}{W^2} \right) \partial_\theta f + \left(\frac{L_{11} \mathbf{a}_2 - L_{12} \mathbf{a}_1}{W^2} \right) \partial_\phi f, \tag{33}$$

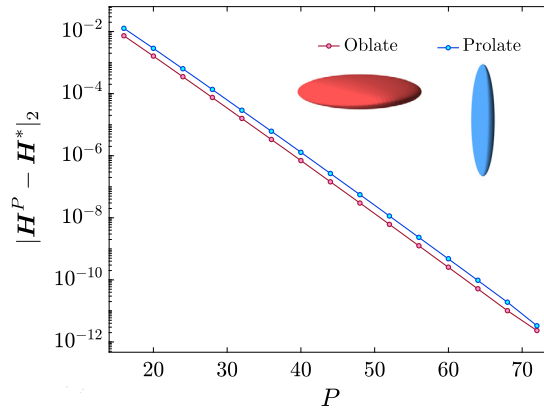


Fig. 2. Error in the mean curvature as a function of the number of modes used in the upsampled P -space. Two ellipsoids with $l/b = 5$ (prolate) and $l/b = 0.2$ (oblate) are investigated, representing the typical shapes of a drop due to EHD flows with large deformations. \mathbf{H}^* is the mean curvature computed on the finest upsampled grid and is chosen as a surrogate for the exact curvature values.

where $W = (\det \mathbf{L})^{1/2}$ is the area element. Similarly, the surface divergence of a vector field $\mathbf{v}(\theta, \phi)$ can be expressed as

$$\nabla_s \cdot \mathbf{v} = \left(\frac{L_{22} \mathbf{a}_1 - L_{12} \mathbf{a}_2}{W^2} \right) \cdot \partial_\theta \mathbf{v} + \left(\frac{L_{11} \mathbf{a}_2 - L_{12} \mathbf{a}_1}{W^2} \right) \cdot \partial_\phi \mathbf{v}. \quad (34)$$

In this study, we consider a drop of an incompressible fluid. In the absence of Marangoni effects, the capillary stress is a function of the mean curvature H :

$$H = \frac{1}{2} \text{Tr}(\mathbf{L}^{-1} \mathbf{B}) = \frac{1}{2} \left(\frac{L_{22} B_{11} - 2L_{12} B_{12} + L_{11} B_{22}}{W^2} \right). \quad (35)$$

3.2. Aliasing errors

Samples of different functions may become indistinguishable on a discrete grid by so-called aliasing [50]. This means, for instance, that a high-frequency spherical harmonic basis function may be aliased to lower frequencies and cause numerical instability in the simulations. Nonlinear operations and differentiation broaden the deformation and charge distribution spectra on a drop. As a result, the energy is moved to frequencies not resolved by the grid resolution and may alias to the resolved frequencies [50]. Therefore, it is physically consistent with the discretization level to remove the corresponding energy from the solution.

Nonlinear operations that are susceptible to aliasing in our simulations include the calculation of mean curvature, electrical stress, convective charge flux, and quadrature. We use a mesh with finer resolution $M > N$ in the simulations to prevent aliasing errors and filter the solution following any nonlinear manipulation [36]. Our numerical experiments show that $M/N = 2$ is sufficient for lower-order and polynomial nonlinearities. For higher-order and non-polynomial nonlinear manipulations such as those incurred in calculating mean curvature and charge convection, we use an adaptive algorithm that follows Rahimian et al. [49]. Given a function \mathbf{f}^M sampled over an M -grid ($2M^2$ points in physical space and M^2 spherical harmonic modes), one can interpolate the distribution on a finer P -grid ($P/M = u_f > 1$) by upsampling:

$$\mathbf{s}^P = \{\mathbf{a}_{nm}, \mathbf{b}_{nm}\}^P = \begin{cases} \{\mathbf{a}_{nm}, \mathbf{b}_{nm}\}^M, & \text{for } n \leq M \text{ and } 0 \leq m \leq n, \\ 0, & \text{for } n > M. \end{cases}$$

The upsampling factor u_f can be determined based on the desired tolerance. Fig. 2 shows how the error in the mean curvature decays as a function of the finer grid resolution P for oblate and prolate spheroids representing typical drop shapes subject to an electric field.

3.3. Numerical integration

This section summarizes the numerical integration scheme used in this study, which follows that used by Zhao et al. [36]. The boundary integrals in (20), (22), and (25) can be written in the general form

$$I(\mathbf{x}_0) = \int_D K(\mathbf{x}, \mathbf{x}_0) f(\mathbf{x}) \, ds(\mathbf{x}) = \int_{\mathbb{S}^2} K(\mathbf{x}(\theta, \phi), \mathbf{x}_0) f(\mathbf{x}(\theta, \phi)) J(\mathbf{x}(\theta, \phi)) \sin \theta \, d\theta \, d\phi, \quad (36)$$

where $f(\theta, \phi)$ is a smooth scalar function over surface D , and K is the kernel containing the Green's function for Stokes or Laplace's equations.

The approximate spacing of a mesh with $N \times 2N$ points is $h = \sqrt{A/2N^2}$, where A is the drop surface area. Due to the singular behavior of K as $\mathbf{x}_0 \rightarrow \mathbf{x}$, the drop surface D is divided into two regions where the integral I has different behaviors. Following the method of floating partition of unity [51,52,36], a local polar patch is considered centered around \mathbf{x}_0 . For any point \mathbf{x} on D , $\rho(\mathbf{x}, \mathbf{x}_0)$ is defined as the distance along the great circle that connects \mathbf{x} to \mathbf{x}_0 on \mathbb{S}^2 . In the next step, a mask function is defined based on this coordinate as

$$\eta(\rho) = \begin{cases} \exp\left(\frac{2e^{-1/t}}{t-1}\right), & \text{for } t = \rho/\rho_1 < 1, \\ 0 & \text{for } \rho \geq \rho_1, \end{cases}$$

where ρ_1 is the cutoff radius. This representation allows for accurate calculation of the integral (36) as η is a smooth function.

Next, the surface integral (36) is split into two parts:

$$I = I_1 + I_2 = \int_D K(\mathbf{x}, \mathbf{x}_0) \eta(\rho(\mathbf{x}, \mathbf{x}_0)) f(\mathbf{x}) \, ds(\mathbf{x}) + \int_D K(\mathbf{x}, \mathbf{x}_0) [1 - \eta(\rho(\mathbf{x}, \mathbf{x}_0))] f(\mathbf{x}) \, ds(\mathbf{x}). \tag{37}$$

The integrand of I_1 has support only inside the patch.

Using a transformation to the local polar coordinate system (ρ, φ) , we write:

$$I_1 = \int_0^{\rho_1} \int_D K(\mathbf{x}, \mathbf{x}_0) \eta(\rho) f(\rho, \varphi) \sin \rho \, d\rho \, d\varphi, \tag{38}$$

where the integrand is finite and periodic in φ . The integral (38) is computed using Gauss quadrature along ρ from 0 to ρ_1 , with a uniform mesh in φ from 0 to 2π . We set the patch radius on the reference sphere to $\rho_1 = \pi/\sqrt{N}$, which means the patch radius is $O(h^{1/2})$ in \mathbb{R}^3 . Inside the patch, \sqrt{N} points are considered along ρ , and $2\sqrt{N}$ points along φ . We note that the quadrature points inside the patch do not coincide with the surface mesh points. Therefore, bi-cubic spline interpolation is used to evaluate the coordinates and other functions at quadrature points of I_1 . The error of singular integration with the mentioned choice of patch size is $O(h^3)$ [52]. Higher-order accuracy can be achieved using a larger patch size at the expense of computational cost.

The second part of integral (37) is I_2 , which has a smooth integrand. Therefore, it is computed accurately as:

$$I_2 \approx \sum_{i=1}^N \sum_{j=1}^{2N} K(\mathbf{x}_{ij}, \mathbf{x}_0) \eta(\rho(\mathbf{x}_{ij}, \mathbf{x}_0)) f_{ij} J_{ij} w_{ij}, \tag{39}$$

where $\mathbf{x}_{ij} = \mathbf{x}(\theta_i, \phi_j)$ are the quadrature points, w_{ij} are the corresponding weights and J_{ij} is the Jacobian of the transformation from \mathbb{S}^2 to \mathbb{R}^3 . The quadrature (39) converges exponentially with the mesh size h , using quadrature points with Gaussian and uniform distributions along the θ and ϕ directions, respectively. Further details on the numerical integration scheme can be found in Zhao et al. [36].

3.4. Incompressibility and charge neutrality

We expect no change in the drop volume, as both fluid phases are incompressible. However, small changes in the volume occur due to numerical errors, which we correct by adjusting the shape of the drop along the normal direction [36]. In the reported simulations, the magnitude of these adjustments is smaller than $10^{-8}r_0$ at every time step, where r_0 is the initial drop radius. Similarly, the net charge (less than $10^{-17}\epsilon E_0$) is subtracted from the surface charge distribution at every time step to ensure charge neutrality throughout the simulations. The mentioned corrections are especially important for long simulation times.

3.5. Reparametrization

The fluid-fluid interface of the drop evolves and deforms. During this process, there is no physical mechanism to inhibit in-plane distortions of the grid (i.e., depletion, aggregation, and skewness) since there is no bending rigidity, in-plane shear resistance, or surface inextensibility. High-frequency components of the spherical harmonic expansions thus grow, exacerbating aliasing errors and resulting in numerical instability. In addition to correcting aliasing errors, as explained in Sec. 3.2, one must develop a reparametrization strategy that ensures stable and accurate simulations over long simulation times. Here, we use an algorithm that minimizes the high-frequency components in the spherical harmonic expansion of

the surface parametrization. This method was introduced by Veerapenani et al. [53,49] and improved by Sorgentone and Tornberg [54].

Consider an implicit representation of the interface as a smooth function $F : \mathbb{R}^3 \mapsto \mathbb{R}$ such that $F(\mathbf{x}) = 0$ for all $\mathbf{x} \in D$ where D is the drop surface. The unit normal vector can be expressed as $\mathbf{n} = \nabla F / |\nabla F|$ at every point on the surface. We define a quality metric $E : \mathcal{X} \mapsto \mathbb{R}$ where \mathcal{X} is the space of smooth functions defined over D . Now, we can view the reparameterization problem as a minimization of $E(\mathbf{x})$ subject to the constraint $F(\mathbf{x}) = 0$ as

$$\min_{\mathbf{x} \in D} \{E(\mathbf{x})\} \text{ subject to } F(\mathbf{x}) = 0. \tag{40}$$

It can be shown that the solution to the problem above is

$$(\mathbf{I} - \mathbf{n}(\mathbf{x})\mathbf{n}(\mathbf{x})) \cdot \nabla E(\mathbf{x}) = \mathbf{0}, \quad \text{and } F(\mathbf{x}) = 0. \tag{41}$$

The choice of quality metric $E(\mathbf{x})$ is not unique and could vary depending on the reparameterization strategy. [53] proposed an equivalent of the following quality metric:

$$E(\mathbf{x}) = \sum_{n=0}^{N-1} \sum_{m=0}^n \alpha_{nm} \left(|\mathbf{a}_{nm}|_2^2 + |\mathbf{b}_{nm}|_2^2 \right), \tag{42}$$

where $|\mathbf{y}|_2$ is the L^2 norm of vector \mathbf{y} , and α_{nm} is the weight for (n, m) -th spherical harmonic. E , as defined in (42), can be viewed as the spectral energy of the parametrization. We aim to minimize the high-frequency part of \mathbf{x} . Therefore, the weights α_{nm} must be small for low frequencies and larger for high frequencies. Here, we use the following perfect low-pass filter

$$\alpha_{nm} = \begin{cases} 1, & \text{for } n > N_{\text{cutoff}}, \\ 0, & \text{for } n \leq N_{\text{cutoff}}, \end{cases}$$

that was used [53,54] to simulate vesicles and surfactant-laden drops and improved long-time simulations. N_{cutoff} is the cutoff frequency and is chosen adaptively based on the energy spectrum as

$$N_{\text{cutoff}} = \min\{k \in \mathbb{N}, 1 \leq k \leq N - 1 \mid E_k/E_1 \leq P_{\text{cutoff}}\}, \tag{43}$$

where $E_k = \sum_{n=k}^{N-1} \sum_{m=0}^n \alpha_{nm} (|\mathbf{a}_{nm}|_2^2 + |\mathbf{b}_{nm}|_2^2)$ and P_{cutoff} determines the fraction of modes we penalize ($P_{\text{cutoff}} = 0.2$ in our simulations) [54].

We solve (41) by marching in pseudo-time τ along the tangential direction at every point:

$$\partial_\tau \mathbf{x} + (\mathbf{I} - \mathbf{n}(\mathbf{x})\mathbf{n}(\mathbf{x})) \cdot \nabla E(\mathbf{x}) = \mathbf{0}, \quad \text{with } \mathbf{x}(\tau = 0) = \mathbf{x}_0. \tag{44}$$

However, following (44) the volume is not necessarily conserved. In addition, the charge distribution is distorted and hence not spectrally accurate. We use the method of Sorgentone and Tornberg [54] to address these challenges. We project the linear pseudo-velocity $\partial_\tau \mathbf{x}$ along the latitudinal and longitudinal directions to obtain the angular pseudo-velocity $(\partial_\tau \theta, \partial_\tau \phi)$. This strategy comprises updates of the angular coordinates (θ, ϕ) for every point and uses the original spherical harmonic expansions for the shape and charge distribution to interpolate the updated distributions. Using the chain rule,

$$\partial_\tau \mathbf{x} = \frac{\partial \mathbf{x}}{\partial \theta} \frac{\partial \theta}{\partial \tau} + \frac{\partial \mathbf{x}}{\partial \phi} \frac{\partial \phi}{\partial \tau} = \mathbf{a}_1 \partial_\tau \theta + \mathbf{a}_2 \partial_\tau \phi. \tag{45}$$

Next, we project $\partial_\tau \mathbf{x}$ along the tangential directions by taking inner products with \mathbf{a}_1 and \mathbf{a}_2 :

$$\partial_\tau \mathbf{x} \cdot \mathbf{a}_1 = |\mathbf{a}_1|^2 \partial_\tau \theta + \mathbf{a}_1 \cdot \mathbf{a}_2 \partial_\tau \phi, \tag{46}$$

$$\partial_\tau \mathbf{x} \cdot \mathbf{a}_2 = \mathbf{a}_1 \cdot \mathbf{a}_2 \partial_\tau \theta + |\mathbf{a}_2|^2 \partial_\tau \phi. \tag{47}$$

The solution to (46) and (47) is

$$\partial_\tau \boldsymbol{\theta} = \mathbf{L}^{-1} \mathbf{w}, \tag{48}$$

where $\partial_\tau \boldsymbol{\theta} = \{\partial_\tau \theta, \partial_\tau \phi\}^T$ and $\mathbf{w} = \{\partial_\tau \mathbf{x} \cdot \mathbf{a}_1, \partial_\tau \mathbf{x} \cdot \mathbf{a}_2\}^T$, and \mathbf{L} is the surface metric defined in (32). We update the angular coordinates of every grid point by marching explicitly in pseudo-time to obtain new angles (θ^*, ϕ^*) . Using the new angular coordinates, we interpolate the position and surface charge at every point based on the spherical harmonic expansions of the grid before reparameterization (at $\tau = 0$):

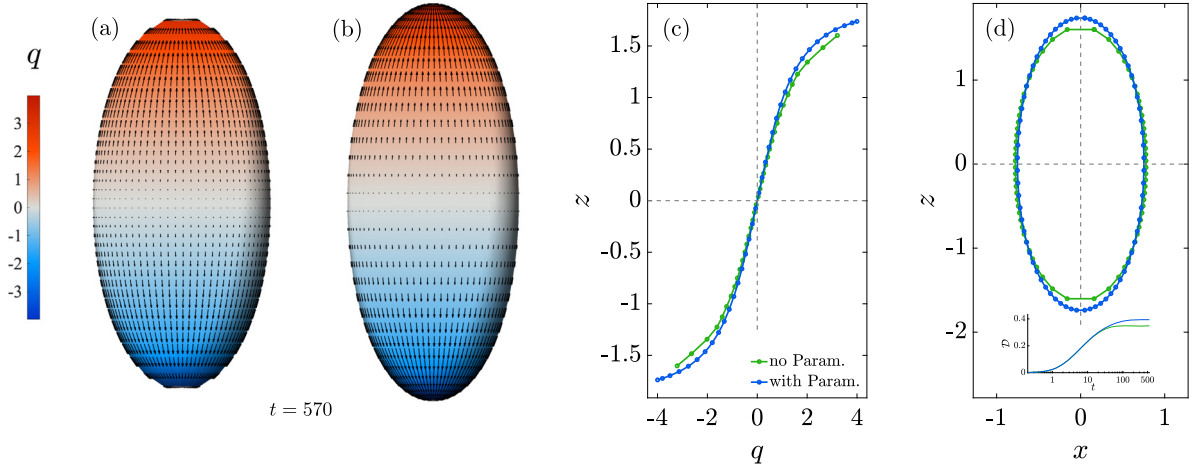


Fig. 3. (a) Large deformations and strong flow degrade the grid when a drop of system **S5** is subject to $(Ca_E, Ma) = (0.35, 5.75)$ without reparametrization. (b) Using reparametrization for the same system maintains the quality of the grid towards the steady state ($P_{\text{cutoff}} = 0.2$, $\tau = 1.0$). Both simulations are performed with $N = 18$ and $M = 2N$. The steady charge profiles and shapes are compared between the two cases in (c) and (d), respectively. Inset of (d) shows the evolution of deformation parameter \mathcal{D} .

Table 1

Systems studied using simulations and their dimensional and non-dimensional parameters. We use $\mu^+ = 0.69$ Pa s, $\gamma = 4.5$ mNm $^{-1}$, and $\epsilon^0 = 8.8542 \times 10^{-12}$ F m $^{-1}$ denotes the permittivity of vacuum.

system	Q	R	λ	Ca_E	Ma	σ^+ (S m $^{-1}$)	ϵ^+/ϵ^0	r_0 (mm)
S1	0.57	36.59	1.41	0.05 – 0.75	0.21 – 3.25	4.5×10^{-11}	5.3	0.7
S2	0.57	36.59	1.41	0.05 – 0.75	0.65 – 9.75	4.5×10^{-11}	5.3	2.1
S3	0.57	36.59	14.12	0.01 – 1.44	0.28 – 5.0	4.5×10^{-11}	5.3	0.25, 0.75, 1.25, 1.75
S4	0.60	10.56	0.07	0.15 – 1.0	0.43 – 2.86	3.8×10^{-11}	4.7	1.9
S5	1.37	0.1	1.0	0.35	5.75	4.5×10^{-11}	5.3	1.37

$$\mathbf{x}^*(\theta^*, \phi^*) = \sum_{n=0}^{N-1} \sum_{m=0}^n \bar{P}_n^m(\cos \theta^*) \left(\mathbf{a}_{nm}^0 \cos m\phi^* + \mathbf{b}_{nm}^0 \sin m\phi^* \right), \quad (49)$$

$$q^*(\theta^*, \phi^*) = \sum_{n=0}^{N-1} \sum_{m=0}^n \bar{P}_n^m(\cos \theta^*) \left(\tilde{a}_{nm}^0 \cos m\phi^* + \tilde{b}_{nm}^0 \sin m\phi^* \right). \quad (50)$$

Finally, a forward (FSHT) and backward (BSHT) spherical harmonic transformation of \mathbf{x}^* and q^* yields distributions on a standard grid that is spaced uniformly along ϕ and Gaussian along θ :

$$\mathbf{x}^* \xrightarrow{\text{FSHT}} \{\mathbf{a}_{nm}, \mathbf{b}_{nm}\} \xrightarrow{\text{BSHT}} \mathbf{x}, \quad (51)$$

$$q^* \xrightarrow{\text{FSHT}} \{\tilde{a}_{nm}, \tilde{b}_{nm}\} \xrightarrow{\text{BSHT}} q. \quad (52)$$

Fig. 3 shows the effect of reparametrization when a drop of system **S5** (see Table 1 for material properties) undergoes large prolate deformations when subject to $(Ca_E, Ma) = (0.35, 5.75)$. In the absence of reparametrization, the grid quality decreases, especially near the drop poles where the grid points are depleted. This results in an inaccurate prediction of the overall behavior of this system, especially considering the fact that the charge concentration is maximum around the poles. We note that this system was previously studied by Lac and Homsy [19] using an axisymmetric model and a simplified charge conservation equation that only accounted for Ohmic conduction. We observe that including interfacial charge convection amplifies the deformation and has a quantitative effect on the behavior of the system, consistent with previous studies on the dynamics of prolate drops [26,20].

3.6. Convergence analysis and time step restriction

We study the overall convergence of our numerical method in Fig. 4(a) for model system **S2**, which reaches a steady state. For the purpose of estimating errors, we take the numerical solution on the finest grid, $N^* = 32$, as a surrogate for the exact solution. We plot in Fig. 4(a) the relative L_2 errors in the steady-state charge density distribution, surface velocity field, and drop shape. Convergence of these fields is observed with a rate $\sim O(N^{-\alpha})$ where $4.11 \leq \alpha \leq 5.44$. The relative

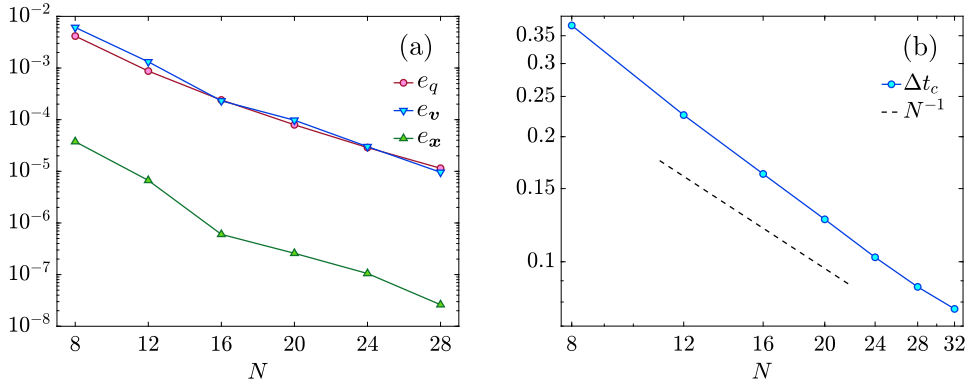


Fig. 4. (a) Numerical convergence for drop shape, charge density distribution, and the velocity field. The numerical solution with $N^* = 32$ is considered as the reference. The plot shows the relative error for each field, defined in Eq. (53), as a function of grid resolution N . (b) Critically stable time step size Δt_c as a function of N . The dashed line represents a first-order variation $O(N^{-1})$. A drop of system **S2** is considered in both cases, with $M = 2N$ and $(Ca_E, Ma) = (0.2, 2.44)$.

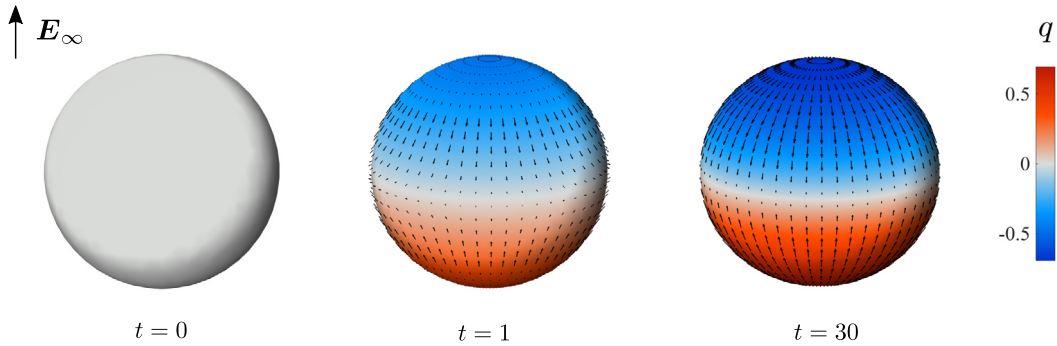


Fig. 5. Axisymmetric regime: evolution of a drop of system **S2** subject to a uniform electric field with $Ca_E = 0.2$ ($N = 8$, $M = 3N$). Colors show the interfacial charge density q , while arrows show the interfacial fluid velocity. Also see Video 1 in the Supplemental Material.

error plotted in Fig. 4(a) is defined as:

$$e_{\mathbf{g}} = \frac{\sqrt{\int_D |\mathbf{g} - \mathbf{g}^*|^2 ds}}{\sqrt{\int_D |\mathbf{g}^*|^2 ds}}, \quad \text{where } \mathbf{g} \in \{q, \mathbf{v}, \mathbf{x}\}. \quad (53)$$

An efficient time marching scheme needs to strike a balance between computational costs and mitigating numerical errors in the charge density distribution and drop shape. The most computationally expensive part of our numerical method is the solution of the Stokes and Laplace's integral equations, which we limit to twice per time step. Our numerical experiments show that an explicit second-order Runge-Kutta scheme performs sufficiently well in most cases, and all simulations shown in the paper were performed using that scheme.

Fig. 4(b) shows the critical time step size for numerical stability, Δt_c , as a function of resolution N for the same model system **S2**. We determine Δt_c using the bisection method by starting from a large time step value. A time horizon of $T = 40$ (in units of τ_{MW}) is chosen such that the system reaches steady state. If the simulation is found to be numerically unstable, we divide the time step size in half and repeat the process until we reach the maximum time step size required for stability, Δt_c . As shown in Fig. 4(b), our method shows first-order behavior with $\Delta t_c \sim O(N^{-1})$. Numerical tests further show that increasing the electric capillary number up to moderate values $Ca_E \leq 0.5$ only reduces Δt_c by a small amount (less than 3%).

4. Numerical results

In this section, we compare the results of our numerical method with existing theoretical solutions and previous computational and experimental studies on electrohydrodynamic flows in drops. In all simulations, the drop is initially uncharged ($q = 0$) and spherical. Table 1 lists all physical systems considered in this study, including their dimensional and non-dimensional parameters.

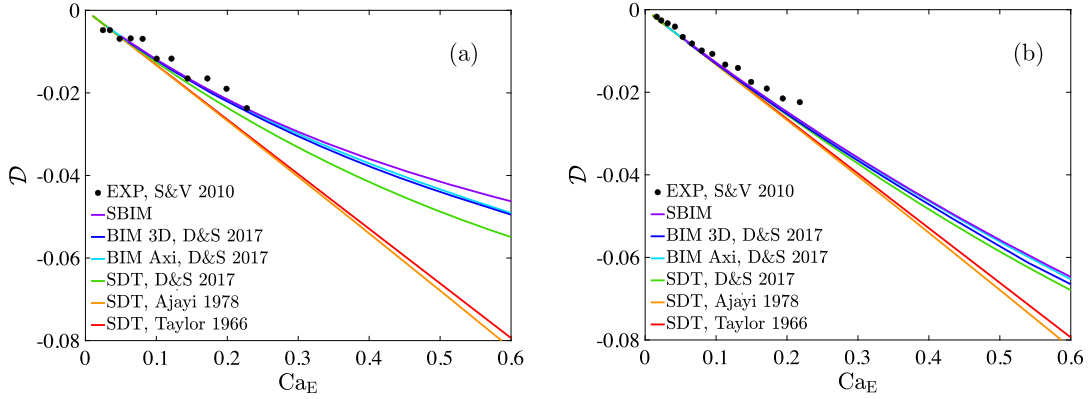


Fig. 6. Steady drop deformation as a function of electric capillary number Ca_E for systems **S1** in (a) and **S2** in (b). Results from our spectral boundary integral method (SBIM) are compared with the experimental data (EXP) of Salipante and Vlahovska [55], along with axisymmetric and three-dimensional boundary integral simulations (BIM) by Das and Saintillan [21,16] and various small-deformation theories (SDT) [8,9,16]. We use $N = 8$ and $M = 3N$ in the simulations shown here.

4.1. Axisymmetric regime

Under weak electric fields ($Ca_E \ll 1$), the drop adopts a steady axisymmetric shape and charge distribution, as shown in Fig. 5. Following the pioneering work of Taylor [8], various analytical models have been proposed to improve the predictions of the steady shape of a drop subject to weak electric fields and small deformations [34,16]. We characterize the deviation from the spherical shape using Taylor’s deformation parameter \mathcal{D} defined as

$$\mathcal{D} = \frac{l - b}{l + b}, \tag{54}$$

where l and b are the lengths of the major axes of the drop in the directions parallel and perpendicular to the applied electric field (see Fig. 1). Fig. 6 shows the steady-state deformation \mathcal{D} from our simulations compared to the predictions of different small-deformation theories (SDT), prior numerical simulations, as well as experimental data. As shown in Fig. 6, our simulation results closely match the experimental data and past predictions. In agreement with past studies [20,16, 21], interfacial charge convection by the fluid flow plays a significant role in the dynamics and tends to weaken drop deformations.

4.2. Quincke regime

Upon increasing the intensity of the electric field, the drop can undergo a symmetry-breaking bifurcation to a dynamical regime characterized by a rotational component to the EHD flow, as shown in Fig. 7 [56,57,55]. When $RQ > 1$, the induced flow is from the poles to the equator, and the induced electric dipole is anti-parallel to the applied electric field. This condition is unfavorable for stability as perturbations to the system result in a destabilizing electric torque that can drive spontaneous rotation beyond a critical electric field strength E_c . Quincke first observed a similar phenomenon in solid spheres subject to uniform electric fields [58]. In solid spheres, the threshold for Quincke electroration is given by [59]

$$E_c = \sqrt{\frac{2\mu^+}{3\epsilon^+ \tau_{MW} (\bar{\epsilon} - \bar{\sigma})}}, \quad \text{where } \bar{\sigma} = \frac{\sigma^- - \sigma^+}{\sigma^- + 2\sigma^+}, \quad \bar{\epsilon} = \frac{\epsilon^- - \epsilon^+}{\epsilon^- + 2\epsilon^+}. \tag{55}$$

Fig. 7 shows the velocity field and charge distribution in a leaky dielectric drop during Quincke electroration. Following the onset of instability, the drop tilts away from its initial orientation. The tilt angle α and deformation oscillate until they reach their steady-state values. Fig. 8 shows the steady-state tilt angle and deformation as functions of the applied electric field E_∞/E_c for a drop of the system **S3**. For a given set of material properties, different values of Ca_{MW} correspond to different drop sizes, and the effect of capillary forces is stronger for smaller drops (small Ca_{MW}). Therefore, we expect the drop tilt angle α to approach that of a solid sphere β [59,55]:

$$\beta = \frac{\pi}{2} - \arctan\left(\frac{E_\infty^2}{E_c^2} - 1\right)^{-\frac{1}{2}}, \quad \text{for } E_\infty \geq E_c, \tag{56}$$

for sufficiently large λ and small Ca_{MW} . This is verified in Fig. 8(a), where the green curve for $Ca_{MW} = 0.44$ closely follows the black line corresponding to β .

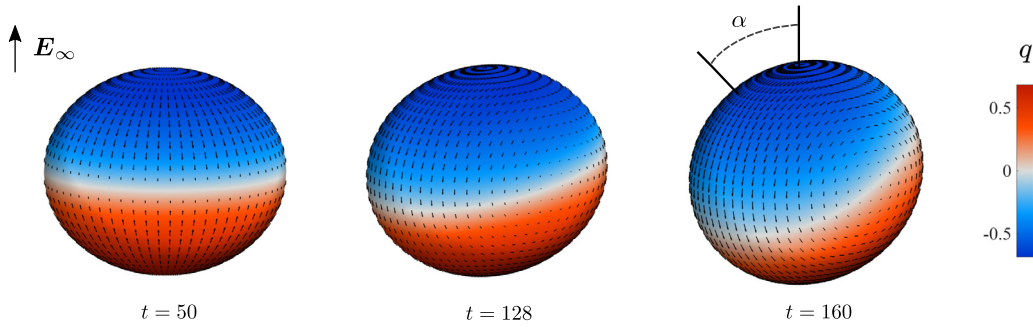


Fig. 7. Quincke regime: spontaneous rotation and tilt of a drop of system **S3** subject to an applied electric field with $E_\infty/E_c = 1$ and $Ca_{MW} = 1.32$ ($N = 10$, $M = 3N$). The tilt angle α is measured from the initial orientation of the drop as depicted. See Video 2 in the Supplemental Material, which shows the evolution of the drop in the Quincke regime.

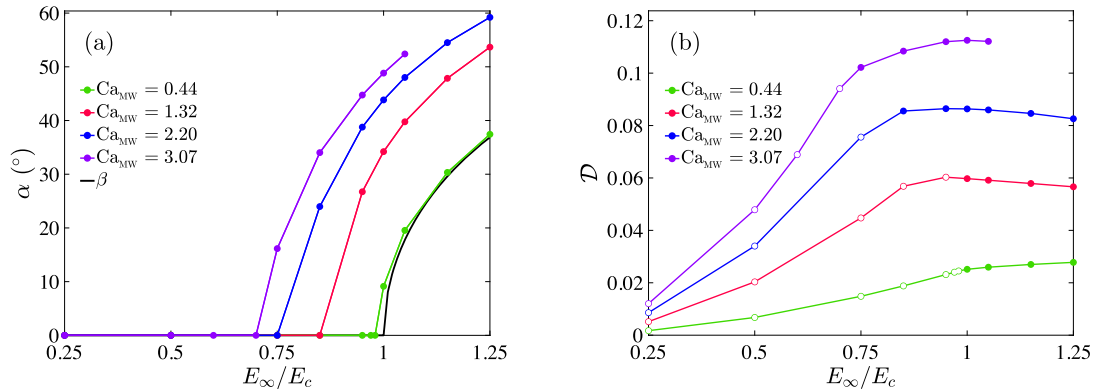


Fig. 8. Steady tilt angle α (a), and deformation \mathcal{D} (b) as functions of the applied electric field E_∞/E_c for different values of Ca_{MW} (different r_0) for system **S3**. The solid black line in (a) shows the tilt angle β of a rigid sphere. Filled markers in (b) show cases where Quincke electroration occurs. We used $N = 10$ and $M = 3N$ in all simulations.

4.3. Low-viscosity drops

When a leaky dielectric drop with $RQ < 1$ is subject to a uniform electric field, it compresses along the field direction. The induced quadrupolar flow is from the poles to the equator at the interface. Under weak electric fields, the characteristic interfacial velocity is:

$$u_T = \frac{9}{10} \frac{R(RQ - 1)}{(2R + 1)^2 (1 + \lambda)} r_0 \tau_{EHD}^{-1} \quad (57)$$

based on Taylor's classic solution [8]. According to Eq. (57), we expect the EHD flow to be stronger in low-viscosity drops ($\lambda < 1$) as $u_T \propto (1 + \lambda)^{-1}$. As we increase the intensity of the applied electric field, low-viscosity drops may undergo different types of EHD instabilities, such as dimpling, equatorial streaming, or Quincke rotation, depending on their material properties [60]. In other experiments, colloidal particles adsorbed on a drop interface were observed to accumulate at the equator and form a belt, which broke into vortices of particles [61,62]. In this section, we study the dynamics of an oblate drop with a small viscosity ratio $\lambda < 0.1$, represented by **S4**, which is similar to the systems studied in [62].

Our results show that charge convection significantly affects the system's dynamics in this regime. As a result of the nonlinear coupling between the flow and charge dynamics, strong charge gradients build up around the equator, as shown in Fig. 9. The nonlinear steepening disappears when we switch off the convective term in the charge conservation equation (5). Therefore, we can conclude that the charge convection is responsible for the nonlinear behavior mentioned. This can be verified in Fig. 9(c) by comparing the charge profiles for simulations with and without convection. The convergent EHD flow sweeps charges of opposite signs towards the equator, which is the stagnation line of the flow where the electric field is locally tangent to the interface. It has been shown that in systems where an interface is subject to a tangential electric field and a converging flow, strong charge gradients develop around the stagnation line due to the effect of charge convection [44].

The emergence of sharp local features in the charge profile poses a fundamental challenge for our numerical method, as spectral methods are most efficient for smooth field variables. It is known that large gradients or discontinuities in a function result in Gibbs phenomenon or ringing artifacts in its spherical harmonic representation [63]. Following the emergence of

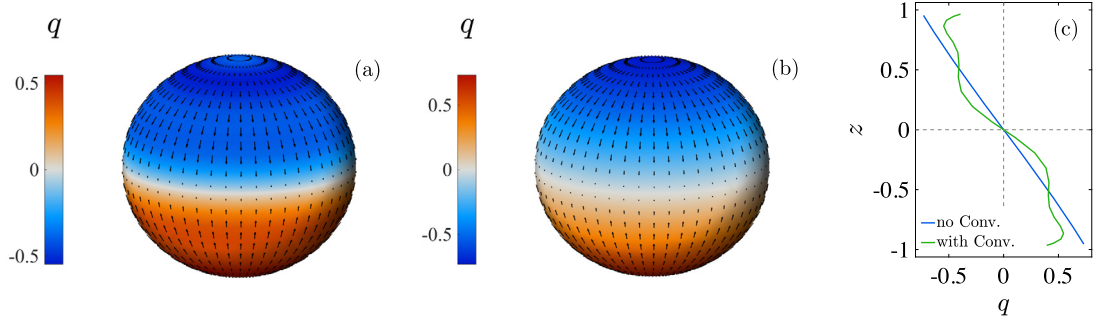


Fig. 9. Charge dynamics in a low-viscosity drop of **S4** with $(Ca_E, Ma) = (0.3, 1.43)$: steady-state velocity and charge density fields in the presence of charge convection (a), and with no convection (b), and the corresponding charge profiles along the direction of the applied electric field (c). We used $N = 8$, $M = 3N$ in both simulations.

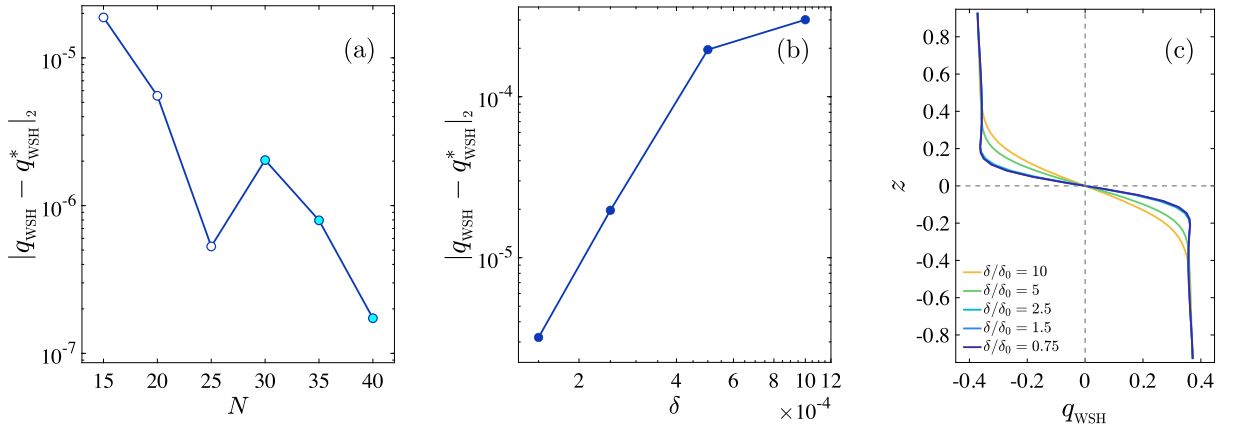


Fig. 10. WSH convergence study on a drop of **S4** with $(Ca_E, Ma) = (1.0, 0.43)$: (a) Error in the steady-state charge density distribution as a function of mesh resolution N for a fixed value of $\delta = 10^{-3}$ ($N^* = 42$). Open markers correspond to cases with ringing artifacts. (b) Error in the steady-state charge density distribution as a function of relaxation factor δ ($N^* = 80$, $\delta/\delta_0 = 0.75$ and $\delta_0 = 10^{-4}$ for q_{WSH}^*). (c) Interfacial charge profiles corresponding to (b). See Video 3 for $N^* = 80$, $\delta/\delta_0 = 0.75$ and $\delta_0 = 10^{-4}$ in the Supplemental Material.

ringing artifacts, spurious oscillations cannot be contained locally as spherical harmonics are global basis functions. This phenomenon occurs due to the inability of the finite spherical harmonic expansions to properly represent the infinite (or very large) derivatives in the sharp regions of a discontinuous (or nearly discontinuous) function.

For relatively small electric capillary numbers $Ca_E < 0.2$, we can suppress ringing artifacts by using a finer mesh corresponding to a larger number of spherical harmonics modes, a more restricted time step size, and, therefore, a higher computational cost. However, due to the stronger effect of charge convection at larger values of Ca_E , refining the resolution is insufficient to remove the ringing artifacts when Ca_E is increased.

To resolve this problem, we propose a weighted spherical harmonic expansion (WSH) for the charge density distribution, where the contribution from high-frequency components is exponentially relaxed [64]. For a given charge density distribution $q(\theta, \phi)$, the spherical harmonic expansion (30) gives coefficients $\tilde{s} = \{\tilde{a}_{nm}, \tilde{b}_{nm}\}$ where $n = 0, 1, 2, \dots, N-1$ and $m \leq n$. Accordingly, we define WSH: $q \mapsto q_{WSH}$ as:

$$q_{WSH}(\theta, \phi) = \sum_{n=0}^{N-1} \sum_{m=0}^n e^{-n(n+1)\delta} \left(\tilde{a}_{nm} \cos m\phi + \tilde{b}_{nm} \sin m\phi \right) \bar{P}_n^m(\cos \theta), \quad (58)$$

where δ is the relaxation factor. Based on (58), we note that WSH does not affect the mean, and q_{WSH} converges to q in the limit of $\delta \rightarrow 0$. In Fig. 10(a), we demonstrate that increasing the mesh resolution at a fixed value of δ eliminates the ringing artifacts. In addition, Fig. 10(b) shows that the steady-state charge profile at a fixed resolution converges as the relaxation factor δ decreases. We combine both approaches and increase N while decreasing δ to refine the numerical solution while avoiding the emergence of ringing artifacts. The numerical solution with the finest resolution, $N^* = \max(N)$, is then chosen as a surrogate to the exact solution in Fig. 10(a,b). To compare the charge profiles computed on different mesh resolutions, we upsample them to a fine grid with $N^u > N^*$. Convergence of the charge profile is illustrated in Fig. 10(c), where the coupling of the flow and charge dynamics is found to result in the formation of strong gradients in the charge density profile near the equator, consistent with previous studies [16]. We note that our simulations for system **S4** under electric

capillary numbers of up to $Ca_E \sim O(1)$ reach a steady axisymmetric profile and do not exhibit any electrohydrodynamic instability, suggesting that the equatorial vortices observed by Ouriemi and Vlahovska [62] could be a consequence of the colloidal particles present at the interface in their experiments.

5. Conclusions

We developed a spectral boundary integral method for simulating electrohydrodynamic flows in leaky dielectric viscous drops. The drop surface, charge density, and all other variables are represented using truncated series of spherical harmonic expansions. In addition to the excellent accuracy of the spectral representation, it enables us to develop a nondissipative dealiasing method required for numerical stability. The charge transport is modeled using a conservation equation for Ohmic conduction from the bulk and surface charge convection by the flow and finite charge relaxation. We employ a reparametrization technique which allows us to explore regimes where drops undergo significant deformations. Our results closely match the existing experimental data and analytical predictions in the axisymmetric Taylor and Quincke electrorotation regimes. The code used in this article is publicly available at https://github.com/mfirouzn/EHD_Drop_3D [42].

Moreover, our simulations confirmed that the dynamics of low-viscosity drops are strongly influenced by interfacial charge convection. In this regime, the interplay between the flow and charge dynamics results in steep gradients in the interfacial charge density. The development of these sharp features in the charge density profile results in the formation of ringing artifacts due to the Gibbs phenomenon. Increasing the mesh resolution would eliminate the Gibbs phenomenon under relatively small electric capillary numbers $Ca_E < 0.2$. However, the effect of nonlinear steepening by the flow is sufficiently strong at larger Ca_E that increasing the resolution does not prevent ringing. To solve this problem, we introduced a weighted spherical harmonic transformation that serves as a relaxation method to exponentially damp high-wave-number coefficients. The convergence of our numerical solution with decreasing level of relaxation shows that it provides a close approximation to the true solution in this regime. Further characterizing the system's dynamics in this regime will require more sophisticated numerical methods for solutions containing discontinuities.

CRedit authorship contribution statement

Mohammadhossein Firouznia: Formal analysis, Investigation, Software, Validation, Visualization, Writing – original draft. **Spencer H. Bryngelson:** Methodology, Software, Supervision, Writing – review & editing. **David Saintillan:** Conceptualization, Funding acquisition, Supervision, Writing – review & editing.

Declaration of competing interest

The authors declare the following financial interests/personal relationships which may be considered as potential competing interests:

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Data availability

The code used for simulations is available on GitHub (link provided in the manuscript).

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Appendix A. Supplementary material

Supplementary material related to this article can be found online at <https://doi.org/10.1016/j.jcp.2023.112248>.

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