# Deformation and stability of a viscous electrolyte drop in a uniform electric field

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# Abstract

We study the deformation and breakup of an axisymmetric electrolyte drop which is freely suspended in an infinite dielectric medium and subjected to an imposed electric field. The electric potential in the drop phase is assumed small, so that its governing equation is approximated by a linearized Poisson-Boltzmann or modified Helmholtz equation (the Debye-Hückel regime). An accurate and efficient boundary integral method is developed to solve the low-Reynolds-number flow problem for the time-dependent drop deformation, in the case of arbitrary Debye layer thickness. Extensive numerical results are presented for the case when the viscosity of the drop and surrounding medium are comparable. Qualitative similarities are found between the evolution of a drop with a thick Debye layer (characterized by the parameter  $\chi \ll 1$ , which is an inverse dimensional drop with a thick Debye layer (characterized by the parameter  $\chi \ll 1$ , which is an inverse dimensional drop with a thick Debye layer (characterized by the parameter  $\chi \ll 1$ , which is an inverse dimensional drop with a thick Debye layer (characterized by the parameter  $\chi \ll 1$ ). sionless Debye layer thickness) and a perfect dielectric drop in an insulating medium. In this limit, a highly elongated steady state is obtained for sufficiently large imposed electric field, and the field inside the drop is found to be well approximated using slender body theory. In the opposite limit  $\chi \gg 1$ , when the Debye layer is thin, the drop behaves as a highly conducting drop, even for moderate permittivity ratio  $Q = \epsilon_1/\epsilon_2$ , where  $\epsilon_1, \epsilon_2$  is the dielectric permittivity of drop interior and exterior, respectively. For parameter values at which steady solutions no longer exist, we find three distinct types of unsteady solution or breakup modes. These are termed conical end formation, end splashing, and open end stretching. The second breakup mode, end splashing, resembles the breakup solution presented in a recent paper [R. B. Karyappa et al., J. Fluid Mech. 754, 550-589 (2014). We compute a phase diagram which illustrates the regions in parameter space in which the different breakup modes occur.

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#### I. INTRODUCTION

The behavior of a viscous liquid drop immersed in a viscous surrounding fluid and acted on by an imposed electric field is a classical problem which has been extensively studied for over one hundred years. It is known that a mismatch in electrical properties between the fluids results in a jump in electric stress at the drop interface. In the case of a drop subjected to a uniform far-field electric field, non-uniform tractions at the drop surface lead to deformation of the interface and, for sufficiently large imposed field, breakup of the drop. This problem arises in a number of important applications, including electrosprays, electrohydrodynamic atomization, breakup of droplets in thunderstorms, microfluidic processes, and others. A thorough review of the topic can be found in [1–3].

For either the case of a perfect dielectric or a perfectly conducting drop in an insulating medium, in which free charge is absent, the electric field only modifies the normal stress balance. This leads to elongated (prolate) steady-state drop shapes, as shown in a number of early studies [4–9]. Criteria on the critical electric field strength beyond which no steady shape exists have been established based on a spheroidal approximation [8, 10] (exact solutions for the electric field in a spheroidal (solid) geometry can be found in [11, 12]). For either a charged or uncharged drop in a uniform electric field, these theoretical results are in fair agreement with experiments. In addition to steady prolate shapes, various unsteady shapes have been observed. For example, the work of Taylor [8], together with experiments, elucidated the well-known 'Taylor-cone' phenomenon for a conducting drop which includes predictions of the angles for a drop or bubble with conical ends (see also related work in [13– 16). It is worth noting that Taylor's solution [8] is in fact based on a local analysis that assumes a steady or equilibrium cone, while in experiments the dynamics is often observed to be unsteady, with a thin fluid jet emitted from the end of the conical or pointed tip. The results in this paper, using a simple electrokinetic model, also show conical behavior in the conducting drop limit. A recent review on topic of Taylor cones in two-phase flow can be found in [17].

Experiments [7] in weakly conducting fluids show oblate drop shapes, in addition to the prolate profiles mentioned above. To explain the oblate drop shapes, Taylor proposed the so-called leaky dielectric [18] or Taylor-Melcher (TM) model (see [2]) for weakly conducting fluids by incorporating surface charge. Tangential electric stresses generated by surface

charge are found to be important for predicting steady state shapes and unsteady drop behavior that is in qualitative agreement with experiment, although quantitative agreement is sometimes lacking. A few drop shapes other than simply prolate or oblate spheroidal, such as an elongated shape with multiple lobes or unsteady profiles with small blobs of fluid connected by a thin thread to a larger fluid mass [7, 9, 19, 20] have been predicted numerically based on the leaky dielectric model [21, 22]. However, various experiments also report discrepancies between the TM model and experimental results (see for example Vizika and Saville [23] and Ha and Yang [19]), which suggest the necessity of further modeling work.

Despite the lack of quantitative agreement with experiments on drop deformation in electrified fluids, the TM and related models have been widely and successfully applied to many problems in electrohydrodynamics. This includes electrified thin liquid films [24, 25] and liquid threads or layers [26–30]. The phenomenon in which a thin thread or small droplets are ejected from a Taylor-cone-like drop end is predicted using the TM model (interfacial surfactant can produce similar 'tip-streaming' phenomena in the absence of an electric field, see e.g. [31–33]). The recent study of Collins et al. [34] based on the TM model demonstrated that fluid ejection from Taylor cones is due to the effect of surface charge convection in the model (see also the work [35] for a charged inviscid drop with surface charge convection). Our results show the possibility of liquid ejection of a different type for an electrokinetic fluid model, when intrinsic surface charge is not explicitly represented in the model. Similar electrohydrodynamic tip-streaming is numerically observed in Pillai et al. [36, 37], who include ion dynamics to describe the origin of the charges in the TM model. A thorough review on the TM model can be found in Melcher and Taylor |2|, Saville |38|, and Vlahovska [39], which makes further attempts to explain the origin of surface charges in the TM model by employing the Poisson-Nernst-Planck equation of electrokinetics (see also the theoretical studies of [40–42]).

In recent years there has been renewed interest in more detailed electrokinetic models which incorporate equations governing the dynamics of (bulk) ionic charge. Of particular interest has been the relationship of Taylor's leaky dielectric model to the thin Debye layer limit of the electrokinetic models. Following the early work by Baygents and Saville [40], Schnitzer and Yariv [43] derived a macro-scale model based on asymptotic analysis of electrokinetic equations. The TM model without surface charge convection is shown to be recovered in the thin-Debye-layer and large electric field limit (it is argued that effective sur-

face charge convection emerges from higher order terms). As in [40], the model of Schnitzer and Yariv assumes surface adsorption-desorption of charge from the bulk. An alternative analysis is given by Mori and Young [44], who derived Taylor's full leaky-dielectric model in the thin Debye layer limit by including a weak reaction term in the ion transport equations as a model for a weak electrolyte, as suggested in Saville [38]. They assume that no charges accumulate at the genuine interface, so that interface charge in the macro-scale model results solely from the layer structure.

In this paper, we modify the traditional model for a perfect dielectric drop in an insulating medium by assuming that the fluid inside the drop is an electrolyte, while keeping the exterior fluid as nonionic. We avoid the difficulty of solving the full nonlinear problem for the ion dynamics, as in [36], by employing the Debye-Hückel approximation for the electric potential inside the drop, which results a linearized Poisson-Boltzmann or a modified Helmholtz equation. The model presented here is formally valid for small potential, or more precisely, when  $e\phi/(k_BT) \ll 1$ . However, we will sometimes also apply it to cases with large deformation (e.g. Figure 9) which may be beyond its range of formal validity. In this case, our model gives behavior that is in qualitative agreement with experiments, and with other models. A discussion of the Debye-Hückel theory and its limitations is given in [45]. A similar model was previously presented in Hua et al. [46], but there the focus is on analytical theory for small drop deformation. We go beyond this and carry out a more comprehensive numerical investigation by reformulating the problem as a system of boundary-integral equations for the coupled electrostatic field and fluid flow (Stokes flow) problem. Numerical computations based on a boundary integral formulation for the problem of freely suspended drops in an electric field have been popular due to their accuracy and simplicity [14, 21, 22, 47–52]. When the electric potential is governed by the Laplace equation, there is an analytical expression for the axisymmetric version of the Green's function (i.e., the azimuthal part of the surface integral can be done analytically). However, this is not the case for the modified Helmholtz equation that arises here from the linearized Poisson-Boltzmann equation. Although accurate numerical schemes to solve boundary integral formulations of Laplace and Helmholtz equations in axisymmetric geometries have been developed (see [53] and references therein), one of the main contributions in this paper is to develop a scheme to accurately and efficiently compute the Green's function for the modified Helmholtz equation and apply it to a moving boundary problem. In this way, we extend previous studies to assess the effect of ions on drop deformation in the case of *arbitrary* Debye-layer thickness. For numerical efficiency, our results are specialized to the case when the viscosity of the drop and surrounding medium are comparable.

Along with the numerical simulations, we also carry out a slender-body analysis (in the case of highly elongated drops) starting from the boundary-integral equations to approximate the electric field inside the drop. A correction term to the results of [13] that takes into account the presence of ions is derived, and the result is shown to agree reasonably well with simulations of the full problem.

The analysis and numerics are used to find steady state solution branches or response curves over a wide range of parameter values. For parameter values at which steady solutions no longer exist, we find three distinct types of unsteady solution or breakup modes, which are termed conical end formation, end splashing, and open end stretching (see Section IV B for examples of these breakup modes). We compare our results with similar behavior observed in other studies [15, 34–37, 50] based on different models.

The paper is organized as follows. We begin in Section II with a complete description of the equations governing the electric field, viscous flow and the boundary conditions. In Section III, the problem is reformulated as a system of boundary integral equations and the numerical method is introduced. Numerical results are presented in Section IV. We summarize the effect of ions on the drop's steady deformation. Meanwhile, the nonlinear instability is also discussed. Closing remarks are provided in Sec. V. In Appendix A, we present the formulations of the Green's function and its derivatives for the modified Helmholtz equation, as well as results demonstrating the accuracy of our numerical scheme in computing the Green's functions. In Appendix B, we present a brief derivation of the small deformation theory for our problem. This is used to compare with and partially validate our numerical results. Finally, details of the slender-body analysis are presented in Appendix C.

#### II. MATHEMATICAL FORMULATION

#### A. Electrokinetic equations

We consider the dynamics of an electrolyte fluid drop with viscosity  $\lambda\mu$  (region 1) immersed in a dielectric (nonionic) medium with viscosity  $\mu$  (region 2), as shown in Figure 1.

Cylindrical polar coordinates  $\mathbf{x} = r\mathbf{e}_r + z\mathbf{e}_z$  are used with the z-axis alligned with the drop's axis of symmetry. The surrounding medium is considered as a perfect dielectric and the electric potential  $\phi_2$  satisfies Laplace equation with a far field condition  $\phi_2 \to -\mathcal{E}z$  due to the applied electric field  $\mathbf{E} = \mathcal{E}\mathbf{e}_z$ .

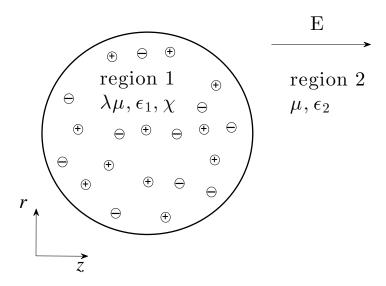


FIG. 1. An electrolyte fluid drop with viscosity  $\lambda \mu$  is surrounded by a nonionic fluid with viscosity  $\mu$ . A constant electric field, directed along the z-axis, is imposed in the far field.

In the drop phase, the electric potential is governed by Poisson's equation,

$$-\epsilon_1 \nabla^2 \phi_1 = \rho = \sum_{i=1}^N e z_i c_i, \tag{1}$$

with  $\epsilon_1$  the dielectric constant and  $\rho$  the bulk volume charge density. We assume that the ions are in thermo-equilibrium and that their concentration follows a Boltzmann-distribution [54–56],

$$c_i = c_i^0 e^{-z_i e(\phi_1 - \phi_0)/(k_B T)},$$
 (2)

where  $c_i^0$  is the constant equilibrium concentration of ion species i,  $k_B$  is the Boltzmann constant, and T is temperature. Here, the neutral bulk condition that  $\sum_{i=1}^n z_i c_i^0 = 0$  is applied. Here  $\phi_0$  is a reference potential, which is set to zero without loss of generality. We introduce

$$\beta = \frac{\mathscr{E}eR}{k_B T} \tag{3}$$

which measures the ratio of the imposed field potential over the thermal potential. Using R,  $\mathscr{E}R$  and  $c^*$  as the characteristic length, electric potential and ion concentration in equation (1), where R is the unperturbed drop radius and  $\mathscr{E}$  is the uniform electric field at infinity, we obtain the Poisson-Boltzmann equation in dimensionless form

$$-\nabla^2 \phi_1 = \frac{eRc^*}{\epsilon_1 \mathscr{E}} \sum_{i=1}^N z_i c_i^0 e^{-\beta z_i \phi_1}.$$
 (4)

In the case of small applied (drop phase) potential  $|\beta\phi_1| \ll 1$  this equation simplifies at leading order to the linearized Poisson-Boltzmann or modified Helmholtz equation,

$$\nabla^2 \phi_1 = \chi^2 \phi_1,\tag{5}$$

where

$$\chi^2 = \frac{e^2 R^2 c^*}{\epsilon_1 k_B T} \sum_{i=1}^N z_i^2 c_i^0.$$
 (6)

This linearized equation, also known as Debye-Hückel approximation, is widely used in various problems involving electrolyte solutions [41, 46, 54–58]. At the drop interface, we have the boundary conditions

$$\phi_1 = \phi_2, \quad Q\phi_{1n} = \phi_{2n}, \tag{7}$$

where  $Q = \epsilon_1/\epsilon_2$ . This specifies that no ionic charge accumulates at the interface, a condition used in the electrolyte-drop model of [46].

## B. Fluid motion and stress boundary conditions

The fluid motion is approximated by Stokes equations, which are nondimensionalized using the spherical drop radius R for lengths,  $\gamma/R$  for pressure, and  $\gamma/\mu$  for velocities, with  $\gamma$  the surface tension,

$$-\nabla p_1 + \alpha \phi_1 \nabla \phi_1 + \lambda \nabla^2 \boldsymbol{u}_1 = 0, \quad \nabla \cdot \boldsymbol{u}_1 = 0,$$
 (8)

$$-\nabla p_2 + \nabla^2 \boldsymbol{u}_2 = 0, \quad \nabla \cdot \boldsymbol{u}_2 = 0. \tag{9}$$

In the above,

$$\alpha = \chi^2 E_b Q \text{ with } E_b = \epsilon_2 \mathcal{E}^2 R / \gamma,$$
 (10)

where the latter quantity is an electric bond number which measures the ratio of Maxwell or electric stress to capillary pressure. The stress balance on the interface is written as

$$[\mathbf{T} \cdot \mathbf{n}]_2^1 = [\mathbf{\sigma} \cdot \mathbf{n}]_2^1 - \Delta \mathbf{f}^e = -\kappa \mathbf{n}, \tag{11}$$

where  $\sigma = -p\mathbf{I} + 2\lambda_i \mathbf{e}$  is the hydrodynamic stress tensor with  $\lambda_1 = \lambda$  and  $\lambda_2 = 1$ ,  $\mathbf{e}$  is the symmetric part of the velocity gradient,  $\mathbf{n}$  is the outward unit normal, and  $\kappa$  is the interface curvature, taken as positive for a convex surface. Here  $[\cdot]_2^1$  denotes the jump across the interface, with the convention that it is the limit as the interface is approached from the interior domain (region 1) minus the limit from the exterior domain (region 2). The Maxwell stress, or electric contribution to the stress balance, is directed normally to the interface and is given by

$$\Delta \mathbf{f}^e = \frac{E_b(Q-1)}{2} \left( Q E_{1n}^2 + E_t^2 \right) \mathbf{n}, \tag{12}$$

where  $E_n$  and  $E_t$  are normal and tangential derivatives of electric field respectively (see also Sherwood [21], Lac and Homsy [22], Miksis [47]). Far from the drop,  $\phi_2 \to -z$  as  $|\mathbf{x}| \to \infty$ .

#### III. BOUNDARY INTEGRAL METHOD

#### A. Integral equations

We reformulate the electrostatic problem as a system of boundary integral equations using classical potential theory [59]. Denote the Green's function for the modified Helmholtz equation by  $G^{\chi}$ ; expressions for this Green's function are presented in Appendix A. The electric potentials  $\phi_1$  and  $\phi_2$  in regions 1 and 2 satisfy

$$\frac{1}{2}\phi_{1}(\boldsymbol{x}_{0}) + \int_{S}\phi_{1}(\boldsymbol{x})\frac{\partial G^{\chi}}{\partial n_{x}}(\boldsymbol{x},\boldsymbol{x}_{0}) dS(\boldsymbol{x}) = \int_{S}\frac{\partial \phi_{1}}{\partial n}(\boldsymbol{x})G^{\chi}(\boldsymbol{x},\boldsymbol{x}_{0}) dS(\boldsymbol{x}), \qquad (13)$$

$$-\frac{1}{2}\left(\phi_{2}(\boldsymbol{x}_{0}) - \phi_{\infty}(\boldsymbol{x}_{0})\right) + \int_{S}\left(\phi_{2}(\boldsymbol{x}) - \phi_{\infty}(\boldsymbol{x})\right)\frac{\partial G^{0}}{\partial n_{x}}(\boldsymbol{x},\boldsymbol{x}_{0}) dS(\boldsymbol{x}), \qquad (14)$$

$$= \int_{S}\frac{\partial\left(\phi_{2} - \phi_{\infty}\right)}{\partial n}(\boldsymbol{x})G^{0}(\boldsymbol{x},\boldsymbol{x}_{0}) dS(\boldsymbol{x}), \qquad (14)$$

where  $\phi_{\infty} = -z$  is the imposed far-field electric potential. We note that setting  $\chi = 0$  in (13) recovers the case of a perfect dielectric drop, in which electrolyte is not present in the interior (see [47]).

The standard boundary integral formulation of the Stokes flow problem for the fluid velocity is modified to include the electrostatic forcing. Starting from the Lorentz reciprocal relation [60], we obtain

$$\mathbf{u}_{1}(\mathbf{x}_{0}) = \frac{\alpha}{8\pi\lambda} \int_{V} \phi_{1}(\mathbf{x}) \nabla \phi_{1}(\mathbf{x}) \cdot \mathbf{J}(\mathbf{x}, \mathbf{x}_{0}) \ dV(\mathbf{x}) + \frac{1}{8\pi\lambda} \int_{S} \mathbf{n}(\mathbf{x}) \cdot \mathbf{\sigma}_{1}(\mathbf{x}) \cdot \mathbf{J}(\mathbf{x}, \mathbf{x}_{0}) \ dS(\mathbf{x}) 
+ \frac{1}{8\pi} \int_{S} \mathbf{n}(\mathbf{x}) \cdot \mathbf{K}(\mathbf{x}, \mathbf{x}_{0}) \cdot \mathbf{u}(\mathbf{x}) \ dS(\mathbf{x}),$$

$$\mathbf{u}_{2}(\mathbf{x}_{0}) = -\frac{1}{8\pi} \left( \int_{S} \mathbf{n}(\mathbf{x}) \cdot \mathbf{\sigma}_{2}(\mathbf{x}) \cdot \mathbf{J}(\mathbf{x}, \mathbf{x}_{0}) \ dS(\mathbf{x}) + \int_{S} \mathbf{n}(\mathbf{x}) \cdot \mathbf{K}(\mathbf{x}, \mathbf{x}_{0}) \cdot \mathbf{u}(\mathbf{x}) \ dS(\mathbf{x}) \right),$$
(15)

where J and K are the Stokeslet and Stresslet Green's functions for Stokes flow. The first term on the right hand side of (15) is transformed into a surface integral by using the divergence free property of the Stokeslet, namely,

$$\nabla \cdot \boldsymbol{J} = 0, \tag{17}$$

which leads to

$$\int_{V} \phi_{1}(\boldsymbol{x}) \nabla \phi_{1}(\boldsymbol{x}) \boldsymbol{J}(\boldsymbol{x}, \boldsymbol{x_{0}}) \ dV(\boldsymbol{x}) = \frac{1}{2} \int_{V} \nabla_{\boldsymbol{x}} \cdot \left( \phi_{1}^{2}(\boldsymbol{x}) \boldsymbol{J}(\boldsymbol{x}, \boldsymbol{x_{0}}) \right) \ dV(\boldsymbol{x}), \qquad (18)$$

$$= \frac{1}{2} \int_{S} \phi_{1}^{2}(\boldsymbol{x}) \boldsymbol{J}(\boldsymbol{x}, \boldsymbol{x_{0}}) \cdot \boldsymbol{n}(\boldsymbol{x}) \ dS(\boldsymbol{x}). \qquad (19)$$

As  $x_0$  approaches interface, the integral equations can be combined to one equation by using (11),

$$u(\mathbf{x}_0) = -\frac{1}{4\pi(1+\lambda)} \int_S \mathbf{J}(\mathbf{x}, \mathbf{x}_0) \cdot \triangle \mathbf{F}^e(\mathbf{x}) \ dS(\mathbf{x})$$
$$-\frac{1-\lambda}{4\pi(1+\lambda)} \int_S \mathbf{n}(\mathbf{x}) \cdot \mathbf{T}(\mathbf{x}, \mathbf{x}_0) \cdot \mathbf{u}(\mathbf{x}) \ dS(\mathbf{x}), \quad (20)$$

where  $\phi_s$  is the electric potential on the interface S and

$$\Delta \mathbf{F}^e = \left(\kappa - \frac{\alpha}{2}\phi_s^2\right)\mathbf{n} - \Delta \mathbf{f}^e,\tag{21}$$

where  $\alpha$  is given in (10) and  $\Delta f^e$  is given by (12). Similar integral equation formulations for a viscous drop in an electric field have appeared in [21, 22, 47], and [60, 61] provides more details in the derivation, as well as numerical implementations. Note that the additional term  $\alpha \phi_s^2$  in (21) is a consequence of the electric body force in the Stokes equation. This term can also be understood via a modified or effective pressure  $\hat{p}_1 = p_1 - \frac{\alpha}{2}\phi_1^2$  in (8), which

similarly implies (21). In the limit of a perfect conductor  $(\chi \to \infty)$ ,  $\phi_1 = 0$ , whereas in the limit of a perfect dielectric drop,  $\chi \to 0$ , hence  $\alpha \to 0$ . Therefore, this additional term vanishes in both extreme cases.

# B. Computation of Green's functions

The axisymmetric version of the free space Green's function for Laplace's equation (defined as the azimuthal integral of the 3D Green's function) can be expressed in closed form, see for example, [61],

$$G^{0}(z, z_{0}, r, r_{0}) = \int_{0}^{2\pi} G^{3D}(z, z_{0}, r, r_{0}, \phi, \phi_{0}) d\varphi = \frac{K(k)}{\pi \sqrt{(z - z_{0})^{2} + (r + r_{0})^{2}}},$$
 (22)

where  $G^{3D}(z, z_0, r, r_0, \phi, \phi_0) = ((z - z_0)^2 + r^2 + r_0^2 - 2rr_0\cos(\phi - \phi_0))^{-1/2}/(4\pi)$  is the Green's function for the 3D Laplace's equation in cylindrical coordinates,  $k^2 = 4rr_0/\left[(z - z_0)^2 + (r + r_0)^2\right]$  and  $k \leq 1$ . Here K(k) is the complete elliptical integral function of first kind,

$$K(k) = \int_0^{\pi/2} \frac{d\theta}{\sqrt{1 - k^2 \cos^2 \theta}}.$$
 (23)

The axisymmetric Green's function for the modified Helmholtz equation, however, does not have an analytical expression. Starting from the Green's function for the modified Helmholtz equation in cylindrical coordinates, we write the axisymmetric version as follows:

$$G^{\chi}(z, z_{0}, r, r_{0}) = \frac{1}{4\pi} \int_{0}^{2\pi} \frac{\exp\left(-\chi \left[(z - z_{0})^{2} + r^{2} + r_{0}^{2} - 2rr_{0}\cos u\right]^{1/2}\right)}{((z - z_{0})^{2} + r^{2} + r_{0}^{2} - 2rr_{0}\cos u)^{1/2}} du$$

$$= \frac{1}{4\pi} \int_{0}^{2\pi} \frac{\exp\left(-\chi \left[(z - z_{0})^{2} + (r + r_{0})^{2} - 4rr_{0}\cos^{2}(u/2)\right]^{1/2}\right)}{((z - z_{0})^{2} + (r + r_{0})^{2} - 4rr_{0}\cos^{2}(u/2))^{1/2}} du$$

$$= \frac{k}{2\pi (rr_{0})^{1/2}} \int_{0}^{\pi/2} \frac{\exp\left(-\Lambda \left[1 - k^{2}\cos^{2}\theta\right]^{1/2}\right)}{(1 - k^{2}\cos^{2}\theta)^{1/2}} d\theta, \tag{24}$$

where  $\Lambda = 2\chi(rr_0)^{1/2}/k$ . When  $|\Lambda| \ll 1$  or  $|\Lambda| \gg 1$  in (24), the Green's function is expanded in an appropriate series for the numerical calculations in [62]. In the present study, we focus on the direct evaluation of (24) by proper quadrature.

Substitution of  $t = \cos \theta$  into (24) results in

$$G^{\chi}(z, z_0, r, r_0) = \frac{k}{2\pi (rr_0)^{1/2}} \int_0^1 \frac{\exp\left(-\Lambda \left[1 - k^2 t^2\right]^{1/2}\right)}{(1 - k^2 t^2)^{1/2}} \frac{dt}{(1 - t^2)^{1/2}}.$$
 (25)

Gauss-Chebyshev quadrature would seem a natural choice to integrate (25), treating  $(1 - t^2)^{-1/2}$  as the weight function. However, our experience shows that Alpert quadrature [63] gives faster convergence and better performance. By recognizing  $(1 - t)^{-1/2}$  as a singular function inside the integrand and setting t = 1 - x, (25) can be rewritten as

$$G^{\chi}(z, z_0, r, r_0) = \frac{k}{2\pi (rr_0)^{1/2}} \int_0^1 \frac{\exp\left(-\Lambda \left[1 - k^2 (1 - x)^2\right]^{1/2}\right)}{\left(1 - k^2 (1 - x)^2\right)^{1/2} (2 - x)^{1/2}} x^{-1/2} dx.$$
 (26)

Alpert quadrature uses a hybrid Gauss-trapezoidal quadrature rule for the integration  $\int_0^1 f(x)dx$  where  $f(x) = g(x)x^{-1/2}$  and g(x) is regular. The quadrature follows the formula (see [63] for more details)

$$T_n^{jkab}(f) = h\left(\sum_{i=1}^j u_i f(v_i h) + \sum_{i=0}^{m-1} f(ah + ih) + \sum_{i=1}^k w_i f(1 - x_i h)\right), \tag{27}$$

where the nodes  $v_1, ..., v_j, x_1, ..., x_k$  and weights  $u_1, ..., u_j, w_1, ..., w_k$  are given for known j, k, a, b which are related to the convergence order, here chosen as fourth order. The total number of nodes is denoted by n = j + m + k. This method is of high order accuracy when the Green's function is regular.

As  $k \to 1$ , i.e. Green's function (24) is close to singular and exhibits the same singular behavior as (22) for Laplace's equation. Simple calculation shows that the singular behavior of the normal gradient of  $G^{\chi}$  is also the same as that of  $G^0$ . We add and subtract the singular Laplace kernel to obtain

$$G^{\chi}(z, z_0, r, r_0) = \frac{k}{2\pi (rr_0)^{1/2}} \int_0^1 \frac{\exp\left(-\Lambda \left[1 - k^2 t^2\right]^{1/2}\right) - 1}{(1 - k^2 t^2)^{1/2}} \frac{dt}{(1 - t^2)^{1/2}} + G^0(z, z_0, r, r_0),$$
(28)

so that the first term is regular, and we use the hybrid quadrature method as described above. Meanwhile, the singularity in  $G^0$  is treated in a standard way, via Gauss-log quadrature [28, 64]. Expressions for the gradient of the Green's function are given in Appendix A and the method of treating the singularity in derivatives of  $G^{\chi}$  is the same as for  $G^{\chi}$ . Furthermore, in Appendix A we provide sample calculations of both the Green's function and its derivatives, which demonstrates the accuracy and of our numerical method. We note that for thin Debye layers  $\Lambda \gg 1$ , and the main contribution to the integral in (28) is localized near t=1 and k=1. Resolution studies show that our computation of the Green's function has error of about  $10^{-6}$  for  $\Lambda$  up to 1414, when N=2048 is employed in the Alpert quadrature.

### C. Numerical procedure

In this paper, we focus on the deformation of axisymmetric drops. The azimuthal part of the surface integrations in each of the integral equations is carried out analytically, except for the ones with Green's functions from the modified Helmholtz equations, for which Alpert quadrature is implemented as described above. The drop interface is discretized by N+1 points, which divide it into N segments. The discretized equations assume the unknown 'densities'  $\phi_s$  and u vary linearly between node points along the interface. On each boundary element, this gives an integral of the product of a linear (density) function and the Green's function. Integrations of this product are carried out in double precision using six-point Gaussian quadrature when the element is regular. As  $x_0 \to x$  the integrand is logarithmically singular, and Gauss-Log quadrature is used to handle the singularity (see also the implementation in [28]). The normal and curvature along the drop interface are calculated by fitting cubic splines, which is similar to Stone and Leal [65].

The linear system that results from discretizing an integral equation is solved by using the Fortran subroutine dgesv in LAPACK. After obtaining surface velocities, the drop interface is advanced by Euler's method via the kinematic condition. The full method is second order accurate in space and first order accurate in time. A solution is deemed to be in a steady state when  $\max |u_n| < 10^{-4}$  along the drop interface. For the steady state calculations reported here,  $N=40\sim70$  is typically enough to resolve the interface. The code for the Stokes droplet without an electric field has been extensively tested and used in earlier work [33, 66]. When electrostatic fields are included in both the drop and surrounding phases in the simpler case of  $\chi=0$ , the code has been validated against small deformation theory as well as the results in [21, 22], where good agreement is obtained. For example, in Figure 2 we compare the numerically computed electric potential  $\phi_s$  for a spherical drop with an analytical prediction from Hua et al. [46], for a uniform imposed field with  $\phi \to -z$  as  $x \to \infty$ . Parameter values are Q=10 and  $\chi=0.1,1,10$ . Excellent agreement is obtained.

If the drop deforms into a highly elongated spheroidal shape or a spindle shape with conical ends, an adaptive regridding scheme is employed. In particular, grid points are redistributed using cubic spline interpolation to be inversely proportional to local curvature, so that the density of points is high near conical ends. One check on the overall method is to compare our calculated results with those in Sherwood [21] (e.g. their Figure 3). The

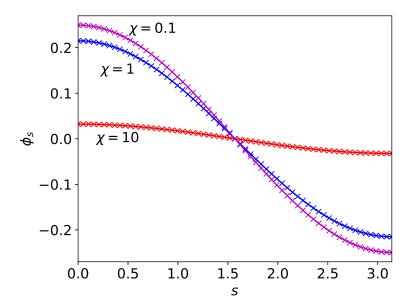


FIG. 2. Comparison of numerically computed results (solid lines) for the surface potential  $\phi_s$  and analytical solutions (cross symbols) for a spherical particle, from [46]. Here Q = 10 and  $\chi$  is indicated in the figure.

results are in excellent agreement (see also our Figure 5). If the drop becomes elongated and exhibits end pinching or other breakup modes, larger N is used (typically  $N \sim 160$  to 320) together with adaptive time stepping. No special adaptive spatial regridding is applied in this case.

## IV. RESULTS AND DISCUSSION

We focus the discussion on the case where the interior and exterior fluids have equal viscosity, i.e.,  $\lambda = 1$ . The third term in equation (20) is then absent, which greatly simplifies the numerics, but also allows for a rich bifurcation diagram and wide variety of unsteady shapes. Results for  $\lambda \neq 1$  will be reported elsewhere.

## A. Steady state drops

In this section we show the computed steady states of drops in a uniform imposed electric field. We note that unless specified otherwise, the reported simulation time is rescaled following [67] as  $t = \tilde{t}\gamma/2\pi R\mu(1+\lambda)$ , where  $\tilde{t}$  is dimensional time. Following other work we

measure drop deformation using the parameter

$$D_f = \frac{l-b}{l+b},\tag{29}$$

where b and l are a typical half width and length of the drop at steady state. To compute the steady response curve we use continuation in the parameter  $E_b$ : once a steady state solution is obtained, we increase  $E_b$  to a larger value and use the steady solution at the previous  $E_b$  as initial data.

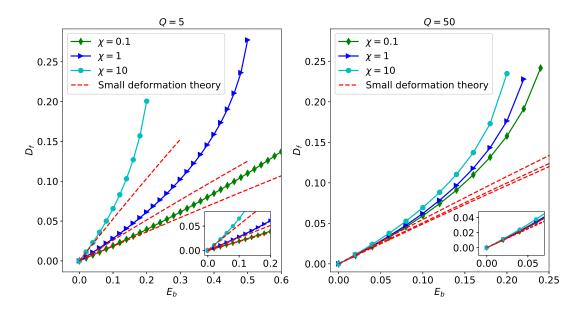


FIG. 3. Comparison between computed results (solid lines with filled symbols) and small deformation theory from [46] and Appendix B (dashed lines) with Q and  $\chi$  indicated in the figure. Insets: Magnification of results near  $E_b = 0$ .

Figure 3 compares the drop deformation for fixed permittivity ratios Q = 5 and Q = 50 and a range of  $\chi$ . The deformation curves follow small deformation theory (see (B6) in Appendix B) when  $E_b$  is relatively small. The deformation is seen to be greater when  $\chi$  is larger, with the same imposed electric field  $E_b$ . This is because capillary pressure has a reduced effect, relative to electrostatic stresses, as  $\chi$  is increased, per (21). This permits a more deformed surface before a local force balance between the capillary force and Maxwell (electrostatic) traction is reached. Increasing Q also tends to increase the deformation at a fixed imposed field strength.

# Conducting drops

In our model, there are two ways to approach the conducting drop limit: either  $\chi \to \infty$  or  $Q \to \infty$ . The surface potential  $\phi_s$  in either case tends to zero, as can be seen by taking the appropriate limit in the small deformation theory (see Appendix B). This theory also shows that the deformation for  $\chi \gg 1$  and  $Q \gg 1$  is given by  $D_f \approx (9/16)E_b$ , for  $E_b \ll 1$ . Figure 4 shows the steady deformation curves for various values of Q and  $\chi$  which all correspond to a highly conducting liquid drop. As expected, the deformation curves nearly overlap each other. The critical value of  $E_b$  at which point steady solutions no longer exist is roughly the same for each branch and is about 0.21, which is consistent with the value reported for a perfectly conducting drop in Karyappa et al. [50] and Dubash and Mestel [48]. The maximum interface potential over all the steady solutions represented in the figure is less than 0.03.

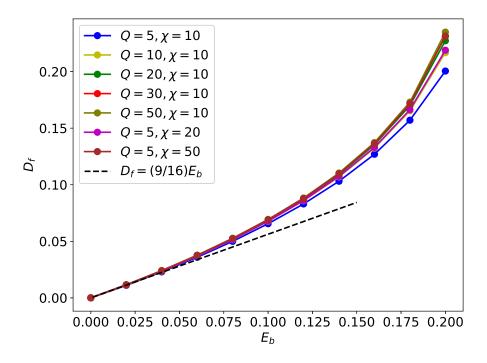


FIG. 4. Steady state deformation curves for various Q and  $\chi$ , corresponding to a highly conducting drop, compared with the small deformation result  $D_f = (9/16)E_b$ .

# Dielectric drops

When  $\chi$  is small and Q is not too large (i.e.,  $\chi \lesssim 1$  and  $Q \lesssim 10^1$ ), the drop is close to a perfect dielectric suspended in an insulating medium (e.g., see the case  $\chi = 0.1$  and Q = 5 in the left panel of Figure 3). Steady solution branches for such nearly dielectric drops are shown in Figure 5, which extends the plot in Figure 3 to smaller values of  $\chi$  and larger imposed field strength  $E_b$ . Instead of plotting the deformation as in Figure 3, we plot the aspect ratio b/l to which is better suited to the wide range of  $E_b$  used here. We also overlay both the analytical solution using the spheroidal approximation at  $\chi = 0$  and the boundary integral solution for a perfect dielectric drop in an insulating medium (the analytical expression for the spheroidal approximation is available in Pillai et al. [36]). It is seen that the computed deformation curves for  $\chi = 0$  and 0.01 almost exactly lie on top of the analytical curve. For  $\chi = 0.1$ , differences between the curves only occur when  $E_b$  is sufficiently large. Consistent with previous observations at small deformation, increasing  $\chi$  promotes larger deformation for a given  $E_b$ . For  $\chi = 0.2$  and 0.25, deviation from the

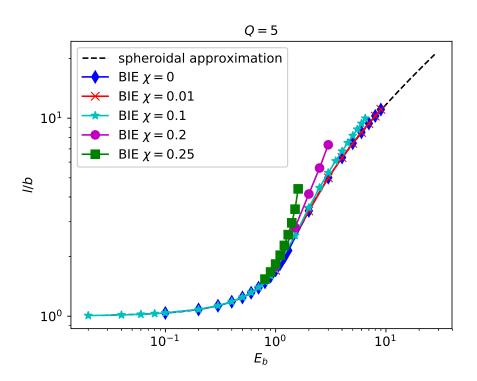


FIG. 5. Comparison of drop deformation for various  $\chi$  and Q=5 together with the analytical results based on a spheroidal approximation for  $\chi=0$ .

insulating drop limit occurs at smaller  $E_b$ , as expected. An aspect ratio of  $l/b \approx 4.4$  is quickly reached at about  $E_b = 1.6$  for  $\chi = 0.25$ , beyond which the drop is found to be unstable and steady solutions no longer exist.

It is natural to carry out a slender body analysis when a highly elongated drop is obtained. In Appendix C, we assume a highly deformed drop and obtain an asymptotic approximation of the electric field for a spheroidal drop, see (C15). The field is shown to be almost uniform. Our results serve as a correction to the result in [13] by taking into account the presence of ions. Integration of the field gives the electric potential, and the theoretical drop shape and electrostatic potential are compared to our numerically computed solutions in Figure 6. In the upper panels of the figure, the drop is highly elongated with aspect ratio about 10 (i.e. slenderness parameter  $\epsilon = b/l \approx 0.1$ ). Both the shape and electric potential are in excellent agreement with theory. In the bottom panels, the drop shape is shown to deviate from a

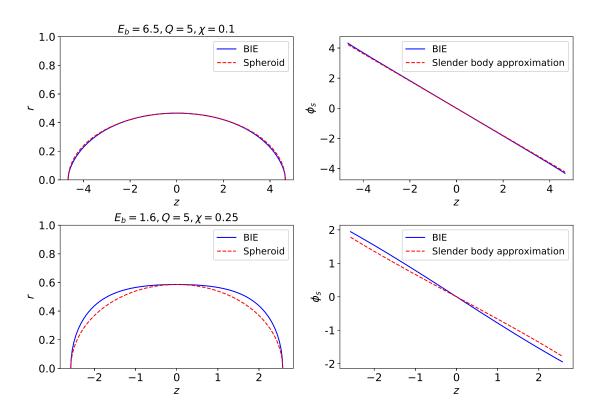


FIG. 6. Comparison of drop shape with a spheroid that has the same aspect ratio. Top left: aspect ratio  $l/b \approx 10$  for a slightly conducting drop with  $\chi = 0.1$ ,  $E_b = 6.5$ , Q = 5. Bottom left: aspect ratio  $l/b \approx 4.4$  with  $\chi = 0.25$ ,  $E_b = 1.6$ , Q = 5. Right: comparisons between calculated surface potentials and slender body approximation, for same parameter values as in left panels.

spheroid with the same aspect ratio (about 4.4). The interfacial potential is also slightly different from the prediction of slender-body analysis but the agreement is still reasonably good.

# B. Breakup behavior

In this section, we investigate drop deformation for parameter values in which steady state solutions do not exist. Several different types of unsteady solution are observed (depending on parameter values), which are classified into three groups: (i) conical end formation, (ii) end-splashing, and (iii) open end stretching: A few case studies are presented before a summary is given.

## Conical end formation

In Figure 7, conical end formation is shown for conducting drops with Q=50 and  $\chi=0.1,1,10$ . In the upper panel of the figure, a time-sequence of unsteady drop shapes are

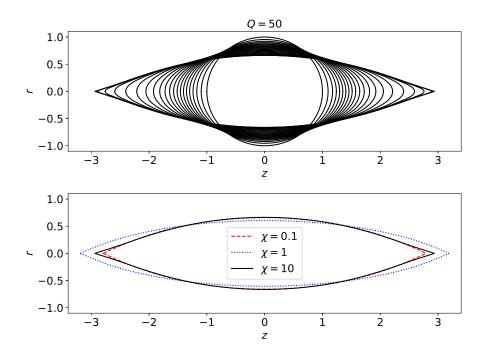


FIG. 7. Upper panel: Evolution of drop with Q = 50,  $\chi = 10$  and  $E_b = 0.26$ . Bottom panel: Drop shapes at breakdown of the numerical scheme for Q = 50 and  $E_b = 0.26$ .

shown for  $\chi = 10$ , Q = 50, starting from an initially spherical shape. In the lower panel, drop shapes at the point at which the simulation is terminated are shown for  $\chi = 0.1, 1, 10$ , which all show the formation of unsteady conical drop tips that are similar to the shapes reported in [13, 14, 21]. The simulation is terminated when the tip curvature  $\kappa_{tip}$  becomes sufficiently large that the number of grid points and time-step required to resolve the interface make the simulation too computationally costly. It is argued in Appendix C 2, based on equation (C16), that conical end formation can occur only for sufficiently large Q, or more precisely  $Q \gtrsim 15$ , regardless of  $\chi$ .

Fontelos et al. [15] present an analysis of conical singularity formation for a charged, perfectly conducting drop in an insulating medium. They find that the singularity formation is self similar with  $\kappa_{tip} = O(\tau^{-\delta})$  and  $U_{n,tip} = O(\tau^{\delta-1})$ , where  $\tau = t_s - t$  is the time to singularity formation (i.e., the singularity occurs at  $t = t_s$ ) and  $\delta$  is a similarity exponent that depends on the opening angle of the cone. Although our model incorporating electrokinetic effects in the drop interior via the linearized Poisson-Boltzmann equation is different from Fontelos et al. [15], this form of the similarity scalings seems to remain unchanged based on our numerical results. We will determine  $\delta$  from numerical data, but one difficulty in doing so is that the singularity time  $t_s$  is unknown. While it can be estimated from numerical data, we take a different approach. Assuming the above self-similar scalings, then  $\tau \sim 1/(\kappa_{tip} \cdot U_{n,tip})$ , and since accurate values for  $\kappa_{tip}$  and  $U_{n,tip}$  are provided by the numerical data, we can replace  $\tau$  by  $1/(\kappa_{tip} \cdot U_{n,tip})$  in a log-log plot to determine  $\delta$ . Such a plot of the time evolution of tip curvature  $\kappa_{tip}$  versus  $1/(\kappa_{tip} \cdot U_{n,tip})$  for a drop which forms conical tips is shown in Figure 8. The figure shows linear behavior for  $\log \kappa_{tip}$  with a slope that is very slightly dependent on  $\chi$ . We estimate the slope magnitude or similarity exponent to be very near  $\delta = 0.71$ , which is close to the value of 0.72 reported in Fontelos et al. [15]. Our simulations give slightly different results for the cone angles than Betelú et al. [14] and Fontelos et al. [15]. We find the semi-angles are between 21  $\sim$  24 degrees for the different  $\chi$  values. In the work of Betelú et al. [14], the semi-angle is shown to be dependent on the viscosity ratio and is about 25 degrees for  $\lambda = 1$ . A slightly different result of about 27.5 degrees is reported in Fontelos et al. [15] (for a different model). While we cannot completely rule out numerical error as a source of the variation of cone angle with  $\chi$ found here, resolution studies suggest that the computed angles are well resolved. We note that the current numerical method is only able to resolve about 2 decades of scaling in the space-time neighborhood of the singularity, which is similar to the other cited studies. For much more than this, and for a more detailed investigation of cone angles, it is anticipated that a specialized numerical treatment of the emerging singularity is needed.

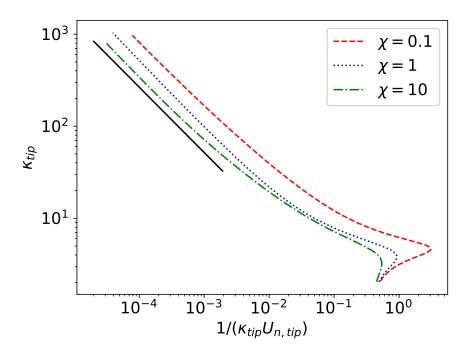


FIG. 8. Evolution of interface tip curvature  $\kappa_{tip}$  versus  $1/(\kappa_{tip}U_{n,tip})$  for the three cases in Figure 7. Here,  $U_{n,tip}$  is the normal velocity at the tip. Following the scaling law  $\kappa_{tip} = O(\tau^{-\delta})$  and  $U_{n,tip} = O(\tau^{\delta-1})$  from Fontelos et al. [15], we plot  $\kappa_{tip}$  versus the time to singularity  $\tau \sim 1/(\kappa_{tip} \cdot U_{n,tip})$ . The estimated or average slope  $\delta \approx 0.71$  is shown as a black solid line.

#### End-splashing mode

When  $\chi$  is large enough (see Figure 12 for precise values), conical end formation is replaced by a small finger that is emitted from the tip, nearly perpendicular to the axis of symmetry or z-axis. This behavior persists even in the highly conducting drop limit of large  $\chi$ . The interface shape near the ends eventually evolves into a 'snail-head' that forms in the vertical direction. We call this the 'end-splashing mode'. Representative examples of end-splashing are plotted in the top two panels of Figure 9. For these simulations, N=320 and the profile is well resolved up to the point when bulbous ends first form, which marks the onset of snail-head finger formation (see bottom panel). Resolution studies of the fully

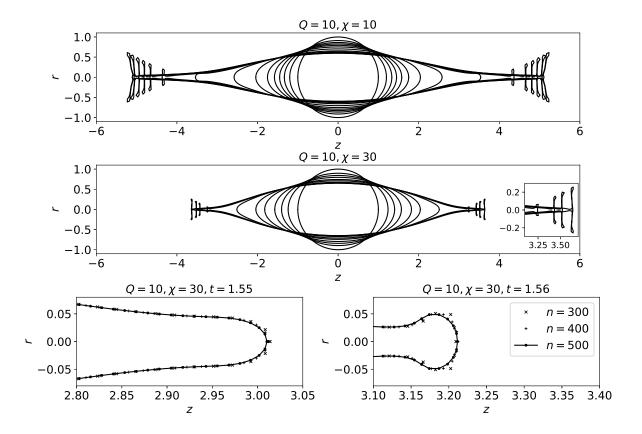


FIG. 9. Breakup of a viscous drop for Q=10 and  $E_b=0.26$  with  $\chi=10$  in upper panel (t=0,0.25,0.49,0.74,0.98,1.23,1.47,1.66,1.77,1.83,1.86,1.88,1.90,1.92,1.94) and  $\chi=30$  in the middle panel (t=0,0.30,0.60,0.90,1.20,1.35,1.50,1.56,1.58,1.59,1.61). The inset of the middle panel show local finger formation before breakup for  $\chi=30$ . The lower two panels show tip profiles at different resolution N for  $\chi=30$  at times before (left) and after the snail head is formed (right). The profiles are well-resolved, at least up to the onset of snail formation.

developed snail head profile, e.g. at the final time in the upper panel of Figure 9, show similar shapes, but slightly decreased snail head length, as resolution is increased. In a 3D view, the drop end looks like a disk or nearly flat cone with a ring rim. This is similar to the so-called dimple formation and lobe-breakup solution reported in Karyappa et al. [50] (see their figure 11 for example). This interface morphology is distinctly different from that observed in Taylor [8], Betelú et al. [14], Grimm and Beauchamp [68] and Fontelos et al. [15], where a Taylor-cone-like solution first develops, then is followed by the ejection of a thin fluid thread in the axial direction.

For highly conducting fluids, similar breakup behavior has been reported in the experiments and numerical simulations of Karyappa et al. [50] which have NaCl added to the drop

phase, suggesting that ions in the drop may contribute to the fingering instability. More recently, Mohamed et al. [69] observed a similar end-splashing mode in experiments on conducting fluids for a pendant drop problem, although there it was attributed to the effect of a more viscous fluid in the surrounding medium. In the current study, this behavior is also obtained when the viscosity ratio is one. The middle panel of Figure 9 shows a similar fingering instability for the larger value  $\chi = 30$ , except that a narrower finger is formed. Our numerical results show a trend of decreasing finger width with increasing  $\chi$ .

# Open end stretching

When both  $\chi$  and Q are both moderate in size (roughly of the order  $10^0$  to  $10^1$ ) our model exhibits relatively long drops, i.e., with  $\epsilon = b/l < 0.01$ . One such example is presented in Figure 10 for  $E_b$  slightly larger than the critical value for nonexistence of steady sates. The top two panels show the drop shape and interface potential, respectively, while the bottom panel plots the maximum normal velocity. As it evolves, the drop forms a cylindrical central thread and the electrostatic field is nearly uniform and directed along the axis of symmetry.

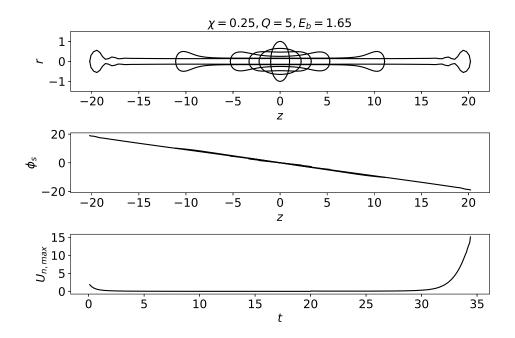


FIG. 10. Breakup of a viscous drop for Q = 5 and  $E_b = 1.65$  with  $\chi = 0.25$ . Top: Drop profiles at times t = 0, 1.73, 29.31, 32.14, 33.56, 34.41. Middle: interfacial potential  $\phi_s$  versus s, at the same t as top. Bottom: maximum normal velocity versus t.

During the evolution, the normal velocity decreases and the drop nearly settles into an elongated steady state. However, the elongational velocity is reestablished after about  $t \approx 30$  as the electric traction overwhelms surface tension. Two 'blobs' develop at the drop ends in a manner similar to the initial stages of the end-splashing mode in Figure 9. However, due to its highly elongated cylindrical thread, we classify this to be a third breakup mode, 'open end stretching'. The simulation in Figure 10 was stopped when the drop aspect ratio exceeded 100.

# Summary of breakup modes

We summarize our results in Figures 11 and 12. Figure 11 shows numerically determined curves in  $E_b - \chi$  space that separate regions where steady drop shapes exist (or S-regions) from those with purely unsteady dynamics (U-regions). Toward the conducting drop limit, i.e., for sufficiently large Q, the U-region is roughly independent of  $\chi$  and occurs above  $E_b \approx 0.22$ . For moderate or small Q, a narrow S-region occurs when  $\chi$  is sufficiently small, i.e., as the perfect dielectric limit is approached. Alternatively, Figure 12 plots the phase diagram in  $Q - \chi$  space with a fixed electric field strength  $E_b = 0.5$ . The behavior in the

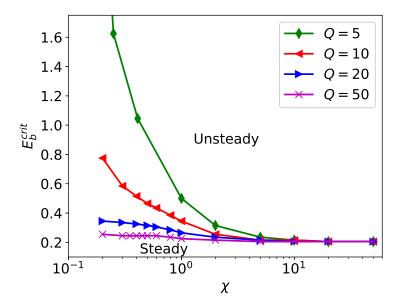


FIG. 11. Boundaries separating steady (S-region) and unsteady (U-region) solutions in the  $E_b - \chi$  plane for various Q.

U-region is further classified by breakup mode. When  $Q \lesssim 15$ , we find the S-region for small  $\chi$  and end-splashing breakup modes (Figure 12(a)) for large  $\chi$ , while for  $\chi$  in between we find open end stretching modes as shown at example point Figure 12(c). When  $Q \gtrsim 15$ , conical end solutions are found for sufficiently small  $\chi$  (Figure 12(b)), consistent with the asymptotic theory in Appendix C2, whereas for larger  $\chi$  we find end-splashing breakup modes.

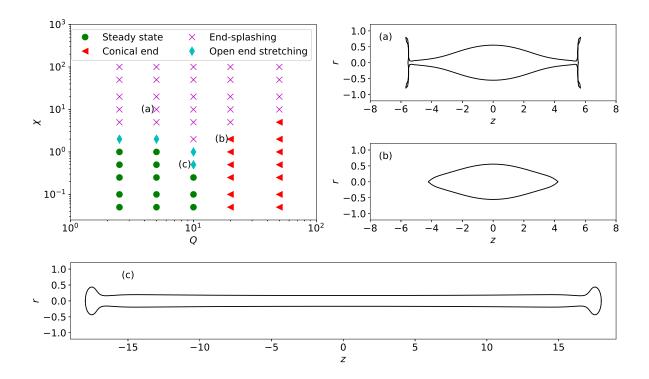


FIG. 12. Phase diagram of steady shapes and breakup modes in  $Q - \chi$  space with  $E_b = 0.5$ . Interface shapes at markers (a)-(c) on the phase diagram are shown in panels at right and below. (a) End-splashing  $(Q = 5, \chi = 10)$ , (b) conical end formation  $(Q = 20, \chi = 2)$  and (c) open end stretching  $(Q = 10, \chi = 0.5)$ 

We conclude this section by noting that only the three aforementioned breakup modes are observed in the current study when the viscosity ratio is unity, while the well-known pinch-off breakup mode is expected to also exist for  $\lambda < 1$ , which is outside of the scope of this work.

# V. CONCLUSION

We have developed a robust and accurate numerical method to evaluate Green's functions for the linearized Poisson-Boltzmann equation and applied it to solve the moving boundary problem for Stokes flow, including electrostatic forces. The method is used to investigate the steady deformation of an electrolyte drop suspended in an insulating fluid medium, for viscosity ratio  $\lambda = 1$ . We demonstrate that the conducting drop limit can be approached through either  $Q \to \infty$  or  $\chi \to \infty$ , and in these limits, the electric field inside the drop vanishes. For large but finite Q, however, the electric field inside the drop is nonzero and can contribute to the deformation. The perfect dielectric drop limit is  $\chi \to 0$ , and for small  $\chi$  we find that a highly elongated steady drop is possible. For given imposed field strength  $E_b$ , the presence of electrolyte enhances the deformation, and increasing  $\chi$  leading to more deformed drops. When the drop is long and slender, results from our full numerical simulations agree well with approximate solutions based on slender body theory. Finally, we study drop breakup behavior by choosing parameters in the regime where steady solutions no longer exist. In addition to conical end formation, we find other two breakup modes, which we call 'end-splashing' and 'open end stretching'. The type of break up depends on parameter values and an example phase diagram is presented which illustrates the dependence on two of those parameters (Q and  $\chi$ ), for fixed  $E_b$ .

## ACKNOWLEDGMENTS

MS acknowledges the financial support from National Science Foundation through the grant DMS-1412789, MM acknowledges the financial support from Natural Science Foundation of China (No. 11701428) and "Chen Guang" project supported by Shanghai Municipal Education Commission and Shanghai Education Development Foundation.

# Appendix A: Computation of axisymmetric Green's functions and their derivatives

In this section, we present the derivatives of the axisymmetric Green's functions (24) and (22). Gradients of the axisymmetric Green's function for Laplace's equation are given by

$$\frac{\partial G^0}{\partial z} = -\frac{(z - z_0)k^3 E(k)}{8\pi (r r_0)^{3/2} (1 - k^2)},\tag{A1}$$

$$\frac{\partial G^0}{\partial r} = -\frac{k^3}{8\pi (rr_0)^{3/2}} \left[ \left( \frac{r - r_0}{1 - k^2} - \frac{2r_0}{k^2} \right) E(k) + \frac{2r_0}{k^2} K(k) \right], \tag{A2}$$

where K(k) is the complete elliptic integral of the first kind (see (23)) and

$$E(k) = \int_0^{\pi/2} \sqrt{1 - k^2 \cos^2 \theta} d\theta \tag{A3}$$

is the complete elliptic integral of the second kind. For the axisymmetric Green's function  $G^{\chi}$  of the modified Helmholtz equation, we find that the Green's function gradient is given by

$$\frac{\partial G^{\chi}}{\partial z} = \frac{k_z}{2\pi (rr_0)^{1/2}} \int_0^{\pi/2} \frac{1 + \Lambda \left(1 - k^2 \cos^2 \theta\right)^{1/2}}{\left(1 - k^2 \cos^2 \theta\right)^{3/2}} \exp\left(-\Lambda \left[1 - k^2 \cos^2 \theta\right]^{1/2}\right) d\theta, \quad (A4)$$

$$\frac{\partial G^{\chi}}{\partial r} = \frac{1}{2\pi (rr_0)^{1/2}} \int_0^{\pi/2} \frac{1 + \Lambda \left(1 - k^2 \cos^2 \theta\right)^{1/2}}{\left(1 - k^2 \cos^2 \theta\right)^{1/2}} \left(\frac{k_r}{1 - k^2 \cos^2 \theta} - \frac{k}{2r}\right) \times \exp\left(-\Lambda \left[1 - k^2 \cos^2 \theta\right]^{1/2}\right) d\theta. \tag{A5}$$

For the general calculations of Green's functions for Laplace equation, details can be found in Pozrikidis [61].

In table I to III, we show data from a computation of  $G^{\chi}$  comparing the Gauss-Chebyshev method and Gauss-Trapezoidal method (or Alpert quadrature), for  $\chi = 0.1, 1, 10$ , and different n or number of quadrature points. The points of evaluation are  $(z_0, r_0) = (\cos(\pi/4), \sin(\pi/4)), (z, r) = (\cos(\pi/4 + \pi/4096), \sin(\pi/4 + \pi/4096))$ , so that (z, r) is slightly different from  $(z_0, r_0)$ . When  $(1 - t^2)^{-1/2}$  is treated as a weight function in (25), the integration, for a regular function f(x), can be done by Gauss-Chebyshev quadrature

$$\int_{-1}^{1} \frac{f(t)}{(1-t^2)^{1/2}} dt = \frac{\pi}{n} \sum_{j=1}^{n} f(t_{j,n}) + \frac{2\pi}{2^{2n}(2n)!} f^{(2n)}(\eta), \tag{A6}$$

for some  $-1 < \eta < 1$  and

$$t_{j,n} = \cos\left(\frac{2j-1}{2n}\pi\right). \tag{A7}$$

TABLE I.  $\chi=0.1,$  Gauss-Trapezoidal parameters j=7, k=6, n=j+k+m

$\overline{n}$	Gauss-Trapezoidal			Gauss-Chebyshev		
	$G^{\chi}$	$G_z^{\chi}$	$G_r^{\chi}$	$G^{\chi}$	$G_z^{\chi}$	$G_r^{\chi}$
16	1.95602152	207.50604021	-208.60234657	1.95602514	207.50604303	-208.60234683
32	1.95602152	207.50604021	-208.60234657	1.95602243	207.50604261	-208.60234833
64	1.95602152	207.50604021	-208.60234657	1.95602175	207.50604218	-208.60234839
128	1.95602152	207.50604021	-208.60234657	1.95602158	207.50604176	-208.60234809
256	1.95602152	207.50604021	-208.60234657	1.95602154	207.50604134	-208.60234770
512	1.95602152	207.50604021	-208.60234657	1.95602153	207.50604093	-208.60234730
1024	1.95602152	207.50604021	-208.60234657	1.95602153	207.50604057	-208.60234693

TABLE II.  $\chi = 1$ 

$\overline{n}$	Gauss-Trapezoidal			Gauss-Chebyshev		
	$G^{\chi}$	$G_z^{\chi}$	$G_r^\chi$	$G^{\chi}$	$G_z^{\chi}$	$G_r^{\chi}$
16	1.66552053	207.50556904	-208.52289051	1.66588210	207.50585107	-208.52291721
32	1.66552054	207.50556904	-208.52289050	1.66561085	207.50580881	-208.52306613
64	1.66552056	207.50556904	-208.52289048	1.66554306	207.50576654	-208.52307180
128	1.66552057	207.50556904	-208.52289048	1.66552612	207.50572433	-208.52304161
256	1.66552057	207.50556904	-208.52289048	1.66552191	207.50568240	-208.52300272
512	1.66552057	207.50556904	-208.52289048	1.66552087	207.50564156	-208.52296267
1024	1.66552057	207.50556904	-208.52289048	1.66552063	207.50560466	-208.52292600

For  $\chi=0.1$ , both quadrature methods work well, however, Gauss-Trapezoidal quadrature is more accurate than Gauss-Chebyshev for moderate and large  $\chi$ , as seen in Table II and III. In particular, when  $\chi=10$ , it is seen that convergence is poor for the Gauss-Chebyshev method. For example,  $G_r^{\chi}$  obtains only one digit of precision at the largest n. In this paper, we therefore employ Gauss-trapezoidal quadrature since it has performed well in our tests. We have not performed an extensive investigation of quadrature methods, as this is beyond the scope of current paper.

TABLE III.  $\chi = 10$ 

$\overline{n}$	Gauss-Trapezoidal			Gauss-Chebyshev		
	$G^{\chi}$	$G_z^{\chi}$	$G_r^\chi$	$G^{\chi}$	$G_z^{\chi}$	$G_r^{\chi}$
16	1.12249190	207.47248568	-208.10737309	1.15507371	207.50023604	-208.11656235
32	1.12249346	207.47248566	-208.10737197	1.13128003	207.49634418	-208.12532268
64	1.12249500	207.47248564	-208.10737086	1.12472887	207.49220608	-208.12550479
128	1.12249602	207.47248562	-208.10737012	1.12305031	207.48800774	-208.12247820
256	1.12249628	207.47248562	-208.10736994	1.12262964	207.48382016	-208.11859296
512	1.12249621	207.47248562	-208.10736999	1.12252599	207.47973719	-208.11458941
1024	1.12249620	207.47248562	-208.10737000	1.12250164	207.47604760	-208.11092268

#### Appendix B: Small deformation theory

Small deformation analysis of an electrolyte droplet immersed in a dielectric fluid and deformed by a nonuniform electric field is provided by [46], from which the following solution for a uniform field can be easily recovered. When the drop is spherical, the general solution for the electric potential is given in spherical radial and polar coordinates  $\bar{r}, \bar{\theta}$  by

$$\phi_1 = -A_1(\bar{r})\cos(\bar{\theta}), \quad \phi_2 = -\left(1 - \frac{A_2}{\bar{r}^3}\right)\bar{r}\cos(\bar{\theta})$$
 (B1)

with

$$i_1(x) = \frac{x \cosh(x) - \sinh(x)}{x^2},\tag{B2}$$

$$i_2(x) = \frac{(x^2+3)\sinh(x) - 3x\cosh(x)}{x^3},$$
 (B3)

$$A_1(\bar{r}) = \frac{3i_1(\chi \bar{r})}{(Q+2)i_1(\chi) + Q\chi i_2(\chi)},$$
(B4)

$$A_2 = \frac{(Q-1)i_1(\chi) + Q\chi i_2(\chi)}{(Q+2)i_1(\chi) + Q\chi i_2(\chi)}.$$
(B5)

When deformability is included, first order perturbation can be used to approximate the shape of the drop. Assuming the shape is perturbed slightly when  $E_b \ll 1$ , [46] derived an expression for the deformation

$$D_f \approx \frac{3E_b h(\chi, Q)}{4 + E_b h(\chi, Q)} \approx \frac{3}{4} E_b h(\chi, Q) + O(E_b^2)$$
(B6)

where

$$h(\chi, Q) = \frac{1}{12Q} \left[ (Q - 1)(1 + 2A_2)^2 + (\chi^2 + 1 - Q)A_1^2(r = 1)Q \right].$$
 (B7)

It is instructive to note some limits in these formulas. First consider the limit of a conducting drop: as  $Q \to \infty$  or  $\chi \to \infty$ , we have  $\phi_1(\bar{r}) = 0$  and  $\phi_2(\bar{r}) = (r^{-2} - r)\cos\theta$ . Inside the drop, both  $\phi_{1\bar{r}}$  and  $\phi_{1\bar{\theta}}$  tend to zero, that is, the electric field is zero for  $\bar{r} < 1$ , but on the drop surface there is a nonzero normal component given by  $Q\phi_{1\bar{r}} \mid_{\bar{r}=1} = \phi_{2\bar{r}} \mid_{\bar{r}=1} = -3\cos\bar{\theta}$ . The surface deformation function satisfies  $\lim_{\chi \to \infty} h(\chi, Q) = (12Q)^{-1}(9(Q-1) + 9Q^{-1})$ , and  $\lim_{Q \to \infty} h(\chi, Q) = 3/4$ . In the limit  $\chi \to 0$  of a dielectric drop,  $\phi_1(\bar{r}) = -(3r/(Q+2))\cos\theta$  and  $\phi_2(\bar{r}) = [(Q-1)/(Q+2)](1/r^2 - r)\cos\theta$ , from which the electric field is easily obtained by differentiation. The surface deformation function satisfies  $\lim_{\chi \to 0} h(\chi, Q) = (3/4)[(Q-1)/(Q+2)]^2$ .

## Appendix C: Slender body analysis

We carry out slender-body analysis on the boundary-integral equations, following Stone et al. [13],. Define a slenderness parameter  $\epsilon = b/l$ , where l and b are the half length and half width of drop, respectively. The existing dimensionless equations are adapted for the slender-body scales by making the substitution

$$S \to \frac{\epsilon}{\nu} \hat{S}, \quad z \to \frac{1}{\nu} \hat{z},$$
 (C1)

where  $\nu = R/l$  and variables with a hat are O(1).

For a slender drop, i.e.,  $\epsilon = b/l \ll 1$ , the electric field inside the drop is to leading order in the axial direction, i.e.,  $E_t \approx E(z) = \nu \hat{E}(\hat{z})$ . As a result, the normal stress balance (11) simplifies to

$$\frac{E_b(Q-1)}{2}(QE_{1n}^2 + E_t^2) + \Delta p = \frac{E_b(Q-1)\nu^2}{2\epsilon^2}(Q\hat{E}_{1n}^2 + \epsilon^2\hat{E}^2) + \Delta\hat{p}$$

$$\approx \frac{\nu}{\epsilon\hat{S}} - \epsilon\nu\hat{S}_{\hat{z}\hat{z}} - \frac{\alpha}{2}\phi_1^2, \tag{C2}$$

where  $\hat{r} = \hat{S}(z)$  is the drop surface shape and  $\triangle \hat{p}$  is the constant pressure difference between drop interior and exterior. The second term on the right hand side of (C2) is the contribution to surface tension from the axial curvature, which is commonly retained despite being higher

order. Since  $\nabla \cdot E = -\chi^2 \phi_1$ , the internal field can be estimated to leading order as

$$\hat{E}_{\hat{r}} \approx -\frac{1}{2} \epsilon^2 \hat{S} \left( \hat{E}_{\hat{z}} + \nu^{-2} \chi^2 \phi_1 \right), \tag{C3}$$

hence (see also [13]),

$$\hat{E}_{1n} \approx -\frac{\epsilon^2}{2\hat{S}} \left( \left( \hat{S}^2 \hat{E} \right)_{\hat{z}} + \nu^{-2} \chi^2 \hat{S}^2 \phi_s \right). \tag{C4}$$

The potential exterior to the slender drop is approximated following [67]. Starting from the boundary integral representation of the exterior potential, we subtract (14) from (13) to obtain

$$\phi_1(\boldsymbol{x}_0) - \phi_{\infty} + \int_S \phi_1 \left( \frac{\partial G^{\chi}}{\partial n} - \frac{\partial G^0}{\partial n} \right) dS = \int_S \frac{\partial \phi_1}{\partial n} \left( G^{\chi} - QG^0 \right) dS, \tag{C5}$$

which is equivalent to the integral-equation used in [13, 22] when  $\chi = 0$ . Following [70], we focus on the contribution of the integral from  $\epsilon \hat{S} \ll |\hat{z} - \hat{z}_0| \ll 1$ . For a point on the drop centerline  $(0, z_0)$ , the Green's function (24) is expanded before evaluating along drop surface  $\hat{r} = \hat{S}$ ,

$$G^{\chi} = \nu \frac{e^{\left(-\frac{\chi}{\nu}\left[(\hat{z}-\hat{z}_{0})^{2}+\epsilon^{2}\hat{r}^{2}\right]^{1/2}\right)}}{2\left((\hat{z}-\hat{z}_{0})^{2}+\epsilon^{2}\hat{r}^{2}\right)^{1/2}} = \frac{\nu-\chi\left[(\hat{z}-\hat{z}_{0})^{2}+\epsilon^{2}\hat{r}^{2}\right]^{1/2}+\frac{\chi^{2}}{2\nu}\left[(\hat{z}-\hat{z}_{0})^{2}+\epsilon^{2}\hat{r}^{2}\right]}{2\left((\hat{z}-\hat{z}_{0})^{2}+\epsilon^{2}\hat{r}^{2}\right)^{1/2}} + \cdots$$
(C6)

Substituting into (C5) yields

$$\phi_{1}(\hat{z}_{0}) - \phi_{\infty} + \frac{\chi^{2}}{4\nu^{2}} \int_{-1}^{1} \phi_{1} \frac{\epsilon^{2} \hat{S}^{2}}{\left((\hat{z} - \hat{z}_{0})^{2} + \epsilon^{2} \hat{S}^{2}\right)^{1/2}} d\hat{z}$$

$$= -(1 - Q) \int_{-1}^{1} \frac{\hat{S} \hat{E}_{1n}}{2\left((\hat{z} - \hat{z}_{0})^{2} + \epsilon^{2} \hat{S}^{2}\right)^{1/2}} d\hat{z} - \frac{\chi}{2\nu} \int_{-1}^{1} \hat{E}_{1n} \hat{S} d\hat{z} + \cdots$$
(C7)

which is further evaluated to be

$$\phi_{1} - \phi_{\infty} + \frac{\epsilon^{2} \chi^{2} \ln(1/\epsilon)}{2\nu^{2}} \phi_{1} \hat{S}^{2}$$

$$= (1 - Q) \frac{\epsilon^{2} \ln(1/\epsilon)}{2} \left( (\hat{S}^{2} \hat{E})_{\hat{z}} + \nu^{-2} \chi^{2} \hat{S}^{2} \phi_{1} \right) - \frac{\chi}{2\nu} \int_{-1}^{1} \hat{E}_{1n} \hat{S} d\hat{z} + \cdots$$
(C8)

This is coupled with the equation for drop volume,

$$\int_{-1}^{1} \hat{S}^2 d\hat{z} = \frac{4\nu^3}{3\epsilon^2},\tag{C9}$$

which readily yields  $\nu = \epsilon^{2/3}$ .

#### 1. Electric field inside a spheroid

For a spheroid,  $\hat{S}^2 + \hat{z}^2 = 1$  and equation (C8) with  $\chi = 0$  is satisfied by a uniform electric field  $E_{1D}$ ,

$$E_{1D} = \frac{1}{1 + \epsilon^2 \ln(1/\epsilon) (Q - 1)} \sim 1 - \epsilon^2 \ln(1/\epsilon) (Q - 1) + \cdots$$
 (C10)

which agrees with the approximation in Stone et al. [13]. For order one  $\chi > 0$  and  $\chi/\nu \gg 1$ , (C8) at leading order (after taking one derivative with respect to  $\hat{z}$ ) becomes

$$-\hat{E}_{1D} + \frac{1}{\nu} \approx -Q \frac{\epsilon^2 \chi^2 \ln(1/\epsilon)}{2\nu^2} \left(\phi_1 \hat{S}^2\right)_{\hat{z}}.$$
 (C11)

Assume  $\hat{E}_{1D} \sim \nu^{-1} + \Psi(\hat{z})$  so that  $\phi_1 \sim -\nu^{-1}\hat{z} - \int^{\hat{z}} \Psi(s)ds$ . After denoting  $F = \int^{\hat{z}} \Psi(s)ds$  we arrive at

$$-F = Q \frac{\epsilon^2 \chi^2 \ln(1/\epsilon)}{2\nu^2} \left[ (1 - \hat{z}^2) \left( \nu^{-1} \hat{z} + F \right) \right].$$
 (C12)

This is rewritten as

$$F = -K\frac{\hat{z}}{\nu} \frac{1 - \hat{z}^2}{1 + K(1 - \hat{z}^2)} \sim -K\frac{\hat{z}}{\nu} (1 - \hat{z}^2)$$
 (C13)

where  $K=Q\frac{\chi^2\epsilon^2\ln(1/\epsilon)}{2\nu^2}\ll 1$ . After some algebra, we arrive at an approximation for the electric field inside the drop, which holds for  $Q\ll \frac{\nu^2}{\chi^2\epsilon^2\ln(1/\epsilon)}$ 

$$\hat{E}_{1D} = \nu^{-1} + F_{\hat{z}} - \epsilon^2 \ln(1/\epsilon) (Q - 1) \nu^{-1} + \cdots$$

$$= \frac{1}{\nu} - Q \frac{\epsilon^2 \chi^2 \ln(1/\epsilon)}{2\nu^3} (1 - 3\hat{z}^2) - \frac{\epsilon^2 \ln(1/\epsilon) (Q - 1)}{\nu} + \cdots$$
(C14)

After using  $E = \nu \hat{E}$ , the electric field is recovered under the original scaling to yield

$$E_{1D} \approx 1 - \epsilon^2 \ln(1/\epsilon) (Q - 1) - Q \frac{\epsilon^{4/3} \chi^2 \ln(1/\epsilon)}{2} (1 - 3(\epsilon^{2/3} z)^2).$$
 (C15)

This is consistent with Stone et al. [13] for  $\chi = 0$ , and the term with  $\chi$  provides a correction due to the presence of ions. The field is used to compare to the full boundary-integral simulation when the drop is elongated.

# 2. Drop with conical ends

For a drop with conical end, locally  $\hat{S} \sim 1 - \hat{z}$  and  $\epsilon = \tan \theta_0$ . It is seen in normal stress balance (C2) that  $E \sim (1 - \hat{z})^{-1/2}$ , then (C8) shows the terms with  $\chi$  serve as higher order

corrections, and we still have the same equation from Stone et al. [13]

$$Q = 1 - \frac{8}{3\tan^2\theta_0\ln(\tan\theta_0)},\tag{C16}$$

which reflects a local balance of force contributions from the electric field and surface curvature. Therefore, to leading order, the formation of a conical drop is independent of  $\chi$ , i.e. the influence of ions. Thus, as for a drop without electrolyte, a conical end is only expected when Q is sufficiently large regardless of  $\chi$ . To be specific, Stone et al. [13] gives a minimum Q around 15.5, above which conical tip is possible.

Finally, we note that the slender drop shape can be analyzed by coupling (C8), (C2) and (C9) (see Stone et al. [13], Sherwood [71], Rhodes and Yariv [72]). Our preliminary results show a mild singularity between a conical and a rounded end exists (same in [71, 72]) and  $E_b(Q-1)\epsilon^{7/3}\ln(1/\epsilon) \sim O(1)$ . However, we do not pursue this further in current study as the predicted shape using slender-body is usually in poor comparison with the full simulation.

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