



HAL
open science

DROPLET IN MICRO-CHANNELS: A NUMERICAL APPROACH USING AN ADAPTIVE TWO PHASE FLOW SOLVER

Jose-Maria Fullana, Yue Ling, Stéphane Popinet, Christophe Josserand

► **To cite this version:**

Jose-Maria Fullana, Yue Ling, Stéphane Popinet, Christophe Josserand. DROPLET IN MICRO-CHANNELS: A NUMERICAL APPROACH USING AN ADAPTIVE TWO PHASE FLOW SOLVER. 1st Pan-American Congress on Computational Mechanics - PANACM 2015, 2015, Buenos Aires, Argentina. hal-01176070

HAL Id: hal-01176070

<https://hal.sorbonne-universite.fr/hal-01176070>

Submitted on 15 Jul 2015

HAL is a multi-disciplinary open access archive for the deposit and dissemination of scientific research documents, whether they are published or not. The documents may come from teaching and research institutions in France or abroad, or from public or private research centers.

L'archive ouverte pluridisciplinaire **HAL**, est destinée au dépôt et à la diffusion de documents scientifiques de niveau recherche, publiés ou non, émanant des établissements d'enseignement et de recherche français ou étrangers, des laboratoires publics ou privés.

DROPLET IN MICRO-CHANNELS: A NUMERICAL APPROACH USING AN ADAPTIVE TWO PHASE FLOW SOLVER

Jose-Maria Fullana, Yue Ling, Stéphane Popinet, Christophe Josserand

¹ Sorbonne Universités,
UPMC Univ Paris 06, UMR 7190, Institut Jean Le Rond d'Alembert,
F-75005, Paris, France

² CNRS, UMR 7190,
Institut Jean Le Rond d'Alembert,
France 4 Place Jussieu, Paris, France
jose.fullana@upmc.fr

Key words: Computational Mechanics, microfluids, two-phase flows

Abstract. We propose a numerical approach to study the mechanics of a flowing bubble in a constraint micro channel. Using an open source two phase flow solver (Gerris, gfs.sourceforge.net) we compute solutions of the bubble dynamics (i.e. shape and terminal velocity) induced by the interaction between the bubble movement, the Laplace pressure variation, and the lubrication film near the channel wall. Quantitative and qualitative results are presented and compared against both theory and experimental data for small Capillary numbers. We discuss the technical issues of explicit integration methods on small Capillary numbers computations, and the possibility of adding Van der Waals forces to give a more precise picture of the Droplet-based microfluidic problem.

1 INTRODUCTION

Droplet-based microfluidics is a very promising tool for performing biochemical or chemical assays. Droplets are unit systems of controlled volume and content, within which mixing can be easily achieved. Several physical phenomena (mechanics, thermocapillarity, solutocapillarity, thermomechanics) either in cumulative or compensative ways appears when we develop microfluidic setups. It is of prime importance to characterize, under controlled experimental conditions, within which range each contribution is the dominant phenomena regarding element migration. Rationalizing these various effects would have important consequences for lab-on-a-chips, and numerical studies are an interesting way to understand each contribution separately.

In microfluidic setups we often have measurements of shape deformations, bubble or drop velocities, and pressures or flow rates at the input/exit conditions, but the knowledge local values of these variables are not easily available, because principally of the small length scales involved in the system. Numerical approaches are then an interesting way for acceding to small length dynamical fields. The validity of numerical approaches requires the validation and the confrontation against theories and experimental data. Theoretical results exist in few academic microfluidic configurations like bubble in cylinders or plates, and it is then necessary to be able to compare positively these academic configurations before extend the prediction to more complex situations which are actually common in microfluidic devices.

In this communication we present an open source two phase flow tool (gfs.sourceforge.net) for computing solutions of the bubble dynamics (i.e. shape and terminal velocity) in microfluidic channel and; numerical results are presented and compared, quantitatively and qualitatively, against both theory and experimental data for small Capillary numbers. We discuss the possibility of adding Van der Waals forces to give a more precise picture of the Droplet-based microfluidic problem.

2 EQUATIONS AND NUMERICAL SCHEME

We use the incompressible, two-dimensional variable-density, NavierStokes equations with surface tension which can be written

$$\rho\left(\frac{\partial U}{\partial t} + U\nabla U\right) = -\nabla p + \mu\nabla^2 U + \sigma\kappa\delta\bar{n} \quad (1)$$

$$\nabla U = 0$$

with $U = (u, v)$ the fluid velocity, $\rho = \rho(x, t)$ the fluid density, $\mu = \mu(x, t)$ the dynamic viscosity. The Dirac distribution function δ expresses the fact that the surface tension term is concentrated on the interface; σ is the surface tension coefficient, κ and \bar{n} the curvature and normal to the interface.

For two-phase flows we introduce the volume fraction $c(x, t)$ of the first fluid and define the density and viscosity as a function of c , i.e. $\rho = \rho(c(x, t))$ and $\mu = \mu(c(x, t))$. The advection equation for the density can then be replaced with an equivalent advection equation for the volume fraction

$$\partial_t c + \nabla(Uc) = 0$$

The Navier-Stokes equations are solved using a finite volume approach based into a projection method. The numerical problem is solved using the open-source package Gerris [6]. A staggered in time discretisation of the volume-fraction/density and pressure leads to the a formally second-order accurate time discretisation. The interface between the different fluids are tracked and followed using a VOF (Volume of Fluid) method. The spatial discretisation is done using a quad-tree square cells which give a very important

flexibility allowing dynamical costless grid refinement into user-defined regions. Finally a powerful discretisation scheme was developed to capture accurately the surface tension term. More information can be found in reference [1]

3 BUBBLE IN A CHANNEL

The studied configuration is presented in Figure 1: a bubble is pushed into a microchannel of width H by a mean flow velocity U_f . The typical length of the bubble is larger than the width H , the bubble is then constrained by the channel. In the stationary regime the velocity bubble is U_b . Using the width H as characteristic length and the mean flow velocity U_f as characteristic velocity the dimensionless Navier-Stokes is then written

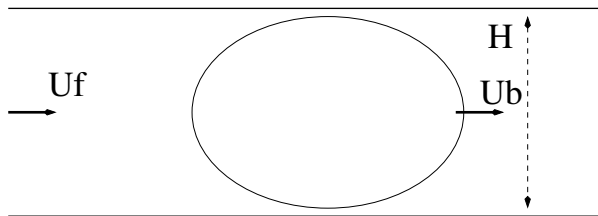


Figure 1: Microfluid configuration : channel width is H , velocities are U_f the mean flow velocity and U_b the bubble velocity.

$$\frac{\partial U}{\partial t} + U \nabla U = -\nabla p + \frac{1}{Re} \nabla^2 U + \frac{1}{Re C_a} \kappa \delta \bar{n} \quad (2)$$

where all dynamical variables are dimensionless and we define the Reynolds and Capillary numbers as

$$Re = \frac{\rho U H}{\mu}, \quad (3)$$

$$C_a = \frac{\mu U_f}{\sigma}. \quad (4)$$

Over this communication we do not discriminate densities and viscosities for liquid and bubble as long as in the subsequent computations we fix the density and the viscosity ratios between the fluid and the bubble to one, then $\rho_f/\rho_b = 1$ and $\mu_f/\mu_b = 1$. Bretherton [2] studied theoretically and experimentally the dynamic of a bubble on a cylindrical configuration. In the limit of small capillarity number based on the bubble velocity, it is shown that the ratio between the gap of the thin film of lubrication h (between the wall and the bubble) and the typical height H of the channel as well as the ratio between the bubble and mean fluid velocity scale both as $C_a^{2/3}$:

$$\frac{h}{H} \sim \frac{U_b}{U_f} \sim C_a^{2/3} \quad (5)$$

the proportionality constant depends, at least, on the geometrical configuration (planar, squared or cylindrical) and on the viscosity ratio. By inspecting the h/H relation and its dependency into the capillary number C_a , it appears that the grid refinement plays an important role if we want to be able to capture the dynamics of the thin film. The Bretherton theory stands that thin film is very important as long as is the key point determining the bubble shape. The Figure 2 (left) present the computed shape for a capillary number of 0.01, which is in fully agreement with the theory. The Figure 2 (right) shows the detail of the grid refinement at the rear of the bubble where the film is thinner. We can also note the specific grid refinement along the interface.

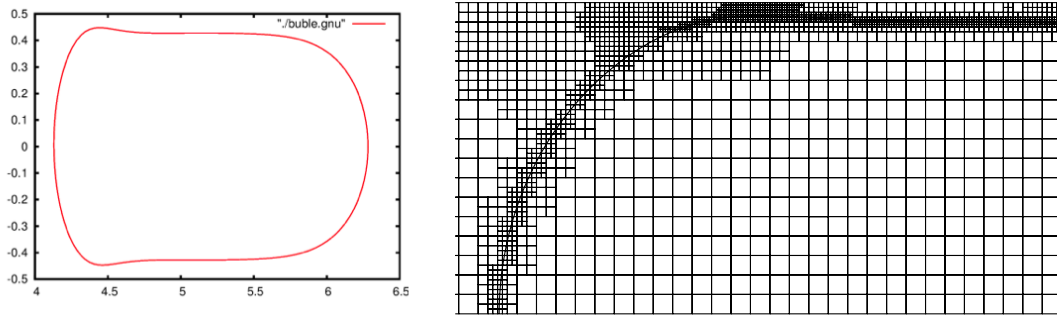


Figure 2: (left) Typical bubble in the stationary state. (right) Grid refinement at the rear of the bubble.

The bubble velocity is evaluated by computing the x position of the center mass of the bubble along the channel, $\frac{\int_S x c(x,t) dS}{\int_S dS}$ where x is the spatial position of each fraction c .

We present now some quantitative and qualitative numerical results on a bubble flowing on a micro channel. For a bubble between parallel plates the analytical solutions are

$$\frac{U_b}{U_f} \sim 1 + 0.643 (3C_a)^{2/3} \quad \text{or} \quad \sim 1 + 0.51 (3C_a)^{2/3} \quad (6)$$

the second relation valid for very viscous drops (liquid-liquid interfaces). In our simulations we impose a Reynolds number of 0.1 and the only variable parameter is the capillary number C_a . The Figure 3 shows the log-log scaling of the excess of velocity $\frac{U_b}{U_f} - 1$ as function of the capillary number C_a up to a capillary number of $0.5 \cdot 10^{-4}$. The solid lines shows both limits from the later relations. These results are consistent with those of reference [3] where capillary number are indeed greater.

Experimental observations in a microfluidic setup [5] show

- that below a capillary number of around 10^{-5} the film gap and the bubble velocity remain constant giving no dependency of the ratios h/H and U_b/U_f on the capillary number, and
- the bubble lost its Bretherton shape becoming more symmetric, like a pancake.

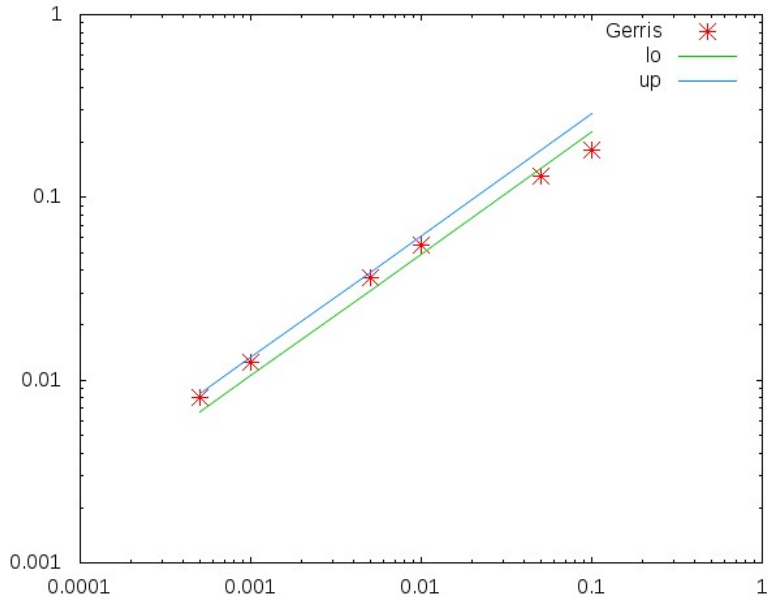


Figure 3: Log-log scaling : $\frac{U_b}{U_f} - 1$ as function of C_a .

This phenomenon appears for h of the order of tens of nanometers, an argument advanced as explanation is that in this region, near of the wall, the Van der Waals forces are not negligible pushing away the bubble by the apparition of an equilibrium film of constant width. The main difficult for including the Wan der Waals forces in this continuum approach is that's necessary resolving numerically a lot of scales, from the small, few nanometers for the equilibrium film, to the large ones, the channel width H . Including Van der Waals forces in a continuum approach was recently done using Gerris to impose the macroscopic contact angles from microscopic physics [4], the numerical computations were done locally and the wide scale range problem was not matter of fact. Starting from the Lennard-Jones potential of two particles and doing some approximations we can add to the r.s.h. of the Navier-Stokes momentum equation (equation (2)) the force $F(d)$ per unit of volume which depends only on the distance d between the bubble interface and the wall

$$F(d) = \frac{K}{d^*} \left[m \left(\frac{h^*}{d} \right)^{m+1} - n \left(\frac{h^*}{d} \right)^{n+1} \right] \quad (7)$$

where K is a constant and h^* is the equilibrium film thickness, $m = 3$ and $n = 2$. (details in reference [4]).

To compare qualitatively this approach without resolving all the spatial scales we impose a large value of h^* which is indeed not physical but the mechanism is still the same, pushing the bubble away from the wall. The Figure 4 present two numerical

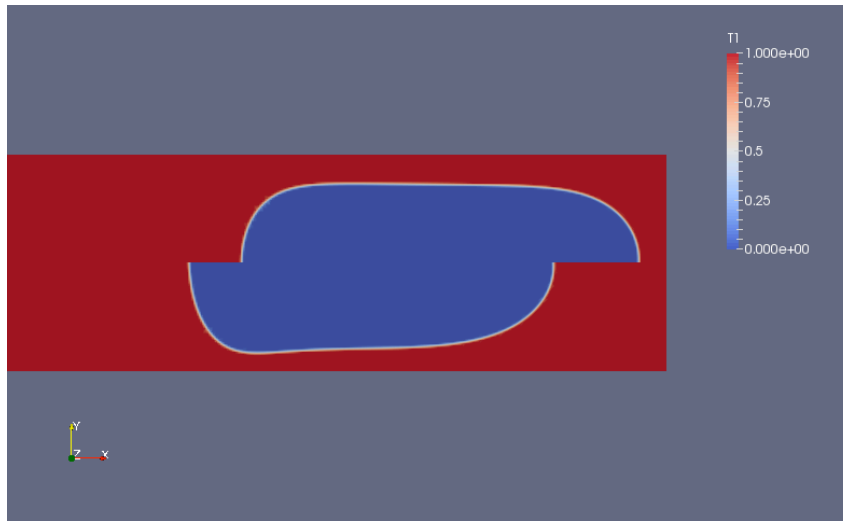


Figure 4: Steady bubble shape and final position including (upper side) or not (lower side) the Van der Waals forces.

simulations of a bubble flowing into a channel for $C_a = 0.01$ and $Re = 0.1$ for the same final time with (upper side) and without (lower side) Van der Waals forces from equation (7). We observe that (i) imposing the Van der Waals forces we found a larger gap h and consequently a faster bubble velocity U_b , (ii) the bubble shape becomes a pancake like.

4 CONCLUSIONS

We have presented numerical simulations of a bubble into a channel, the well behavior of the numerical implementation of the Navier-Stokes equations with a surface tension model was demonstrated by a comparison with Bretherton theory for very small capillary numbers where the scaling law in $C_a^{2/3}$ was validated. The quality of the numerical results are, in particular, a consequence of the grid refinement approach which allows computing the very thin films of liquid between the bubble and the wall. We have also implemented a Van der Waals like force and the imposed minimum gap gives numerical prediction in accordance with experimental observations.

REFERENCES

- [1] Popinet, S. *An accurate adaptive solver for surface-tension-driven interfacial flows*, Journal of Computational Physics 228.16 (2009): 5838-5866.
- [2] Bretherton, F. P. *The motion of long bubbles in tubes*, Journal of Fluid Mechanics 10.02 (1961): 166-188.
- [3] Afkhami, S., A. M. Leshansky, and Y. Renardy. *Numerical investigation of elongated drops in a microfluidic T-junction*, Physics of Fluids (1994-present) 23.2 (2011): 022002.

- [4] Mahady K., Afkhami S., and Kondic L. *A volume of fluid method for simulating fluid/fluid interfaces in contact with solid boundaries*. arXiv preprint arXiv:1409.1276 (2014).
- [5] Project ANR (Agence Nationale de la Recherche) - 2014-2017 - TheRmo-Actuated Migration in a microsystem.
- [6] Gerris is an open-source two phase solver *gfs.sourceforge.net* and Basilisk *basilisk.fr* is another tool, more stand alone, using the methods of Gerris.