

A numerical model for temporal variations during explosive central vent eruptions

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Abstract. An axisymmetrical numerical model has been developed in order to find the temporal evolution of pressure, the position of the exsolution level, the velocity field, the eruption rate, and the amount of erupted material of a shallow, volatile-rich, felsic magma chamber during a Plinian central vent eruption. The overpressure necessary to trigger the eruption is assumed to result from crystallization-driven volatile oversaturation. We solve the resulting set of equations using a finite element method. The results obtained show that the pressure at the conduit entrance decreases exponentially as the eruption proceeds. This produces a shifting of the exsolution level, so that deeper parts of the chamber become progressively volatile oversaturated during the eruption. We assess the influence of chamber geometry and the physical properties of the magma on the computed parameters using several numerical examples. The results are also compared with those predicted by previous models from the literature and are found to be in good agreement with documented eruptions. The model constitutes a first attempt to numerically model the dynamics and the temporal evolution of the most relevant physical parameters during withdrawal from a closed magma chamber.

1. Introduction

Silicic magma is commonly stored in shallow magma chambers located a few kilometers below the surface of the Earth [Smith, 1979]. When the chamber overpressure exceeds the tensile strength of the surrounding rocks a volcanic eruption may be triggered through fractures that propagate to the surface. The increase in magma pressure necessary to trigger the eruption can be produced either by the addition of new magma into the chamber [Blake, 1981] or by the exsolution of volatiles associated with the cooling of the magma during fractional crystallization [Blake 1984; Tait *et al.*, 1989]. We define a system as open when the eruption is due to the addition of new magma into the chamber and as closed when the triggering mechanism is crystallization-driven volatile exsolution.

A good knowledge of the formation and dynamics of magma chambers is crucial to understanding how magmas evolve and how they subsequently erupt at the Earth surface. The dynamics of magma chambers prior to and during volcanic eruptions has become a subject of increasing interest during the recent years. Several analytical [Blake, 1981; 1984; Tait *et al.*, 1989; Bower and Woods, 1997], experimental and numerical [Spera, 1984; Spera *et al.*, 1986; Trial *et al.*, 1992] models have been developed to study the withdrawal of magma from a shallow magma chamber. However, the dynamics of magma chambers during volcanic eruptions are still far from well-understood.

The application of numerical simulations has become a useful tool in such studies, as they allow the variation of some of the

main physical parameters that control the dynamics of magma chambers during volcanic eruptions to be constrained. This has provided an opportunity to understand the significance of some compositional and stratigraphic variations found in pyroclastic deposits in terms of preeruptive gradients in the chamber and dynamics of the magma withdrawal process [Spera, 1984; Spera *et al.*, 1986; Trial *et al.*, 1992]. So far, all these numerical simulations consider the magma chamber as an open system, in which the magma withdrawal process is only driven by pressure forces associated with magma replenishment. Spera [1984] simulated a central vent eruption assuming that magma behaves as a homogeneous, Newtonian, incompressible fluid of constant viscosity. Trial *et al.* [1992] improved this model considering a preeruptive compositional gradient in order to account for the variations observed in erupted deposits. Overpressure produced by oversaturation of volatiles during fractional crystallization has also been proposed as a plausible mechanism of triggering explosive eruptions in felsic systems [Blake, 1984; Tait *et al.*, 1989]. However, no attempts have been made to simulate numerically the dynamics of an erupting magma chamber in such conditions.

The aim of this paper is to characterize the evolution of the physical properties of magma inside the chamber during a Plinian eruption triggered by oversaturation of volatiles in a closed system. We develop an axisymmetrical numerical model of the interior of a chamber which is filled with a homogeneous, volatile-rich, felsic magma in which vesiculation is considered to have taken place prior to the eruption. The driving force that leads to the eruption is thus the volatile overpressure produced during magma cooling and crystallization. In order to find the pressure evolution, the evolution of the exsolution level, the velocity field in the chamber, and the mass discharge rate at the

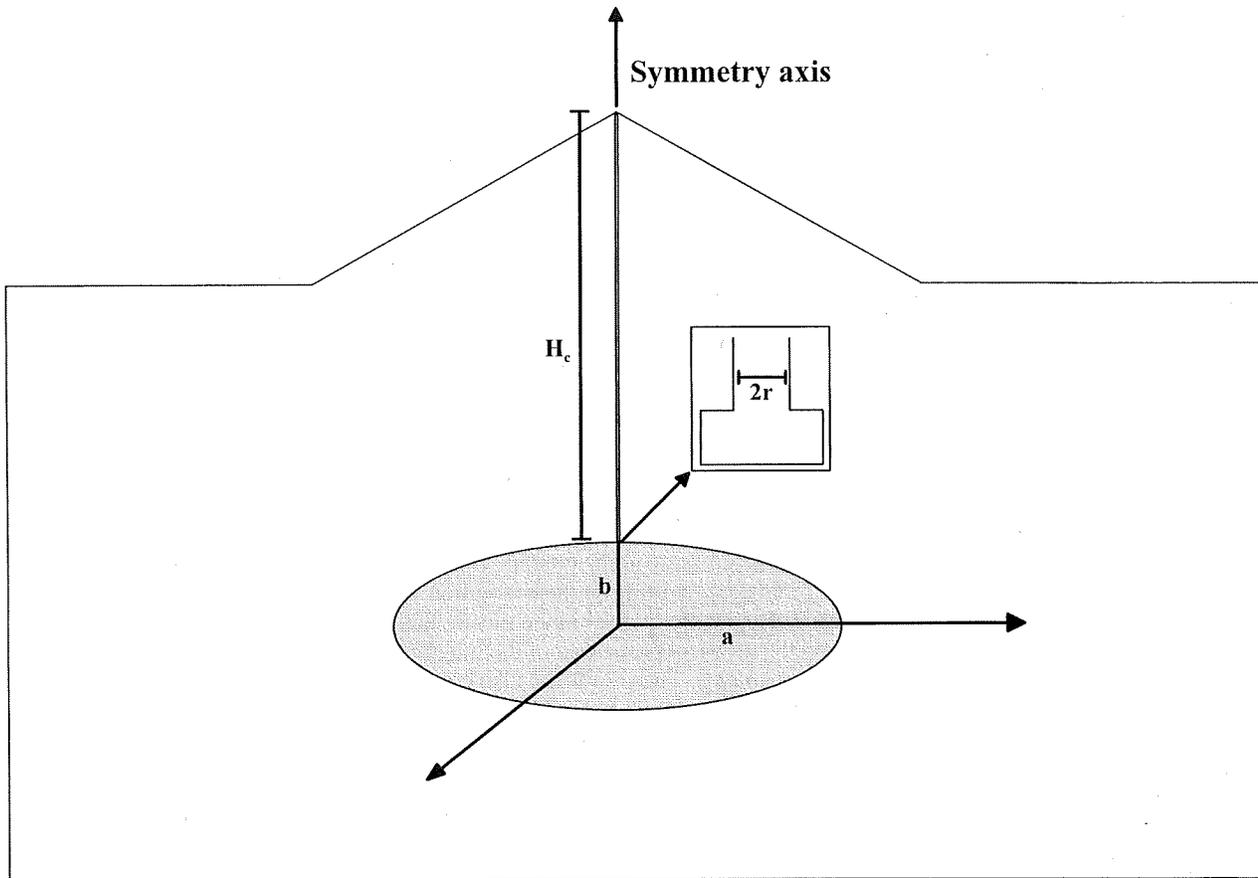


Figure 1. Central vent eruption with axial symmetry. An elliptical magma chamber with semimajor axis a and semiminor axis b is located at depth H_c below the surface of the Earth. Magma flows through a conduit of radius r that coincides with the symmetry axis.

conduit entrance, we solve the corresponding Navier-Stokes equations using a finite element method (FEM). Magma is treated as an incompressible Newtonian fluid below the exsolution level but is compressible once is vesiculated. In the domain above the exsolution level, magma is considered as a homogeneous two-phase flow, and we develop a barotropic state law, where density depends only on pressure, depending on the solubility law.

All the results obtained depend on several parameters (geometry of the chamber, chemical composition of magma, magma properties, etc.). The application of the model allows us to discuss the relative significance of each of these parameters in the eruption process. As we compute the pressure evolution inside the chamber during the withdrawal process, we can estimate the variation of the exsolution level with time. To know the temporal pressure variation at the conduit entrance it is also important to constrain models for the conduit, which generally assume either a constant pressure or a linear decrease with time. In contrast, our model indicates that pressure at the entrance of the conduit decreases exponentially with time.

2. The Model

2.1. Physical Model

We consider a central vent eruption with axial symmetry as shown in Figure 1. The domain comprises a magma chamber located at a depth H_c and a central conduit of radius r along the

symmetry axis. Assuming axial symmetry, we can solve a three-dimensional problem in two dimensions by using cylindrical coordinates.

Plinian eruptions involve conduit erosion during their initial stages. To avoid the coupling with this mechanical problem, we will assume that the magma chamber has rigid walls and that the conduit is not eroded by the flow. Neither the chamber nor the conduit geometries change during the eruption. We consider that the magma behaves as a Newtonian incompressible fluid below the exsolution level but is compressible above this level, where magma is assumed to be a barotropic fluid.

The problem may be approached without considering temperature because magmas have a large thermal capacity and the host rock has a low thermal conductivity so that temperature variations inside the chamber occur on timescales which are orders of magnitude greater than the duration of an eruption [Dobran, 1992]. Therefore we assume that the chamber behaves as a thermally isolated system during the eruption. This assumption would not be so valid in modeling the conduit, where cooling of the magmatic mixture by up to 100°C may occur [Buresti and Casarosa, 1989].

As water is by far the most abundant gas species in felsic magmas, it is considered as the only gas phase. Water content is a parameter of the model, and its value depends upon the magma chamber depth and the lithostatic pressure. We choose a water content which ensures that the chamber is partially vesiculated prior to eruption, in order to reproduce a situation where the triggering mechanism of the eruption could be the

volatile oversaturation. Finally, we assume a chemically homogeneous magma composition for the whole chamber. It should be said that this homogeneous composition is a major limitation of the model, because most magma chambers are zoned systems with compositional gradients [Huppert and Sparks, 1984].

2.2. Governing Equations

The governing equations for this problem are the Navier-Stokes equations, which are derived in continuum mechanics from general principles of mass, energy, and momentum conservation. The energy conservation equation is not needed in our model because temperature is assumed to be a constant inside the chamber during the withdrawal process. A barotropic state law can be therefore established in this region. However, even for barotropic flows, the energy conservation equation remains coupled to the mechanical ones due to the Joule effect. If this contribution is not considered in the energy equation, the use of a barotropic state law will uncouple the thermal equation from the mechanical ones. Therefore only mass and momentum equations are required in our case. Using a spatial notation and Cartesian coordinates, they can be written as follows :

Continuity equation (mass conservation)

$$\frac{\partial \rho}{\partial t} + \frac{\partial (\rho u_k)}{\partial x_k} = 0 \quad (1)$$

Momentum equation (momentum conservation)

$$\frac{\partial \sigma_{ij}}{\partial x_i} + \rho g_j = \rho \left(\frac{\partial u_j}{\partial t} + u_k \frac{\partial u_j}{\partial x_k} \right) \quad (2)$$

Navier-Poisson law (constitutive equation)

$$\sigma_{ij} = -P \delta_{ij} + \mu \left(\frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} - \frac{2}{3} \frac{\partial u_k}{\partial x_k} \delta_{ij} \right) \quad (3)$$

Barotropic state law

$$\rho = \rho (P) \quad (4)$$

where (see notation section) ρ is density, μ is viscosity, P is pressure, u_i is the i th component of the velocity vector, σ_{ij} are the components of the Cauchy stress tensor, and g_j is the j th component of the gravity. In all the equations of this paper, indices run from 1 to 2 or 3 (space dimensions) and repeated indices always imply summation. The proposed barotropic state law for magmas inside the chamber (4) is discussed in the section 2.3. Defining the viscous stress tensor τ (excluding pressure isotropic term in σ) as

$$\tau_{ij} = \sigma_{ij} - \frac{1}{3} \sigma_{kk} \delta_{ij} = \sigma_{ij} + P \delta_{ij} \quad (5)$$

the set of equations can be expressed in its pure conservative form

$$\frac{\partial \vec{V}}{\partial t} + \frac{\partial \vec{C}_i}{\partial x_i} + \frac{\partial \vec{D}_i}{\partial x_i} + \vec{S} = 0 \quad (6)$$

where the conservative variables (mass and momentum)

$$\vec{V}^T = (\rho, \rho u_1, \rho u_2, \rho u_3) \quad (7)$$

are transported by means of convection through the convective fluxes

$$\vec{C}_i^T = (\rho u_i, \rho u_i u_1 + \delta_{i1} P, \rho u_i u_2 + \delta_{i2} P, \rho u_i u_3 + \delta_{i3} P) \quad (8)$$

and by means of diffusion through the diffusive fluxes

$$\vec{D}_i^T = (0, -\tau_{i1}, -\tau_{i2}, -\tau_{i3}) \quad (9)$$

The source term is

$$\vec{S}^T = -(0, \rho g_1, \rho g_2, \rho g_3) \quad (10)$$

where g_i is the i th component of the gravity acceleration vector. Then, in this conservative form, the equations to solve are

Continuity equation

$$\frac{\partial \rho}{\partial t} + \frac{\partial}{\partial x_i} (\rho u_i) = 0 \quad (11)$$

Momentum equation

$$\frac{\partial U_j}{\partial t} + \frac{\partial}{\partial x_i} (u_i U_j) + \frac{\partial}{\partial x_i} (\delta_{ij} P - \tau_{ij}) - \rho g_j = 0 \quad (12)$$

Barotropic state law

$$\rho = \rho (P) \quad (13)$$

where $U = \rho \vec{u}$ is the momentum vector. In a two-dimensional space (or in a three-dimensional space with axial symmetry), equations (11)-(13) form a set of four equations with four unknowns: density (ρ), pressure (P), and the velocity components.

2.3. The State Law

In order to close the problem, a state law (4) is required. For high-viscosity magmas above the exsolution level, the level where exsolution due to melt oversaturation begins, and below the fragmentation level, where the flow is characterized by a mixture of gas with pyroclasts and lithics, the mixture can be considered as a bubbly flow, characterized by a liquid continuum with dispersed gas bubbles and crystals in thermal and mechanical equilibrium [Papale, 1996]. In this bubbly flow regime, the pressure difference between both phases (liquid and the gas bubbles within) can be ignored [Sparks, 1978]. The gas bubbles and the liquid are in mechanical equilibrium and can be considered to move at the same velocity along the conduit because buoyancy effects are negligible compared to viscous drag, even for centimetric bubble sizes (notwithstanding this, the formation of large bubbles in high-viscosity magmas is unlikely because growth by coalescence is impeded). We aim to find a state law for the mixture inside the magma chamber and in the lowermost part of the conduit, that is, in the bubbly flow regime domain, where the mixture is a continuous medium with density:

$$\rho = \alpha \rho_g + (1 - \alpha) \rho_s \cong \alpha \rho_g + (1 - \alpha) \rho_l \quad (14)$$

with α being the gas volume fraction, ρ_g being the gas density, ρ_s being the density of the melt (containing dissolved volatiles) and ρ_l being the liquid density. In (14) it is assumed that the dissolved volatiles do not change the bulk density of the liquid. This homogeneous approximation for the mixture is commonly assumed in the literature [Wilson *et al.*, 1980; Giberti and Wilson, 1990], at least in the lowermost part of the conduit and for high-viscosity magmas. The gas phase is assumed to behave as a perfect gas [Vergnolle and Jaupart, 1986], with

$$P = \rho_g Q T \quad (15)$$

where T is temperature, ρ_g is the gas density, and Q is a gas constant ($Q=461.66 \text{ J}^\circ\text{K kg}$ for water). For the liquid component, a linearized expression in terms of the compressibility modulus β can be assumed [Blake, 1984]:

$$\rho_l = \rho_0(1 + \beta(P - P_0)) \quad (16)$$

where ρ_0 is the density at the pressure reference P_0 . In order to find the state law, we should consider that the volatile species can be either dissolved in the melt or exsolved as a gas phase if the magma is above the exsolution level. Therefore a solubility law is also needed to determine the mass fraction of volatiles dissolved in the melt χ at given thermodynamic conditions. The simplest relationship under magmatic conditions is a Henry's law equation, where the solubility of the silicate melt depends only on a power of the pressure:

$$\chi = sP^m \quad (17)$$

where s and m are constants depending on magma and volatile compositions. Following Tait *et al.*, [1989], values of $s = 4.11 \times 10^{-6} \text{ Pa}^{-1/2}$ and $m = 0.5$ for H_2O in a rhyolitic melt

can be assumed. Using this solubility law, the critical pressure P_c

$$P_c = \left(\frac{W\%}{s} \right)^{\frac{1}{m}} \quad (18)$$

below which gas exsolution takes place depends only on the volatile content ($W\%$). For simplicity, we assume that the exsolution occurs instantaneously if pressure descends below this critical value, and no energetic contributions are considered during the vesiculation process. Let m_d be the mass of dissolved gas in the melt, m_l be the mass of the liquid, m_g be the mass of gas, m_v be the mass of volatiles (dissolved plus exsolved, where $m_v = m_g + m_d$), m_s be the mass of the melt (liquid plus dissolved gas, where $m_s = m_l + m_d$), and M be the total mass ($M = m_l + m_v = m_s + m_g$). Then

$$\chi = \frac{m_d}{m_d + m_l} = \frac{m_d}{m_s}$$

is the gas mass fraction dissolved in the melt,

$$W\% = \frac{m_v}{M} = \frac{m_v}{m_v + m_l}$$

is the volatile mass fraction (dissolved plus exsolved),

$$n_g = \frac{m_g}{M} = \frac{m_g}{m_g + m_s}$$

is the gas mass fraction, and

$$W\% = n_g + (1 - n_g)\chi \quad (19)$$

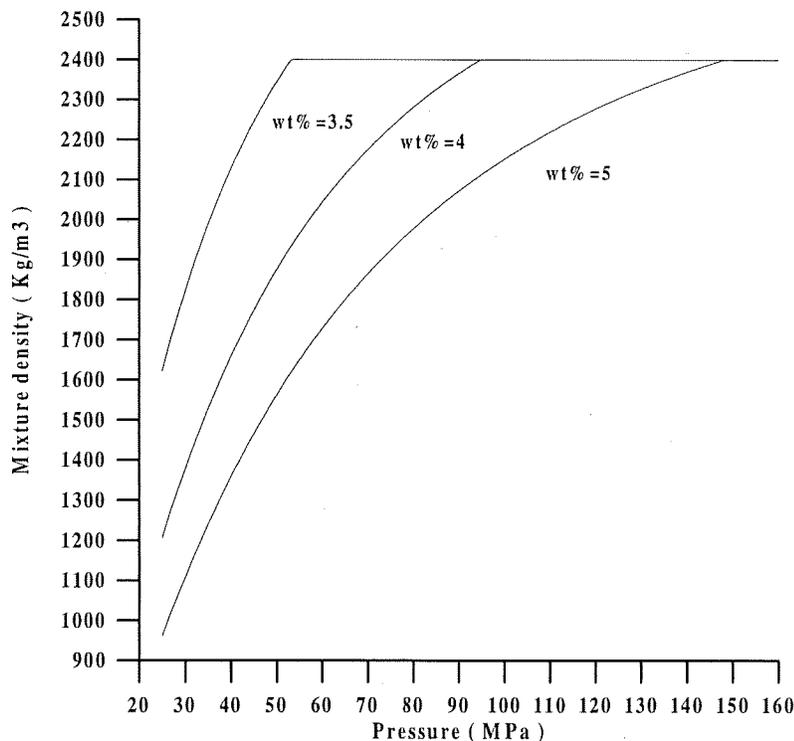


Figure 2. Magma density versus pressure given by the state law (20). The results are for a rhyolitic magma with $T=850^\circ\text{C}$, $\rho_l=2400\text{kg/m}^3$ and having water contents of 3.5, 4.0, and 5.0 in weight, respectively. Note how the exsolution level (the level where the mixture density becomes a constant equal to ρ_l) depends on the water content.

$$sP^m = \chi = \left. \begin{aligned} & \frac{m_d}{m_d + m_l} < W\% = \frac{m_v}{m_v + m_l} \\ & n_g > 0 \end{aligned} \right\}$$

above exsolution level

$$sP^m = \chi > \left. \begin{aligned} & \frac{m_d}{m_d + m_l} = W\% = \frac{m_v}{m_v + m_l} \\ & n_g = 0 \end{aligned} \right\}$$

below exsolution level. Using equations (14) to (19) and after some rearrangement, one arrives at a final expression for the state law:

$$\rho = \frac{\rho_l}{1 + \left(\frac{\rho_l QT}{P} \right) \left(\frac{W\% - sP^m}{1 - sP^m} \right) \theta (P_c - P)} \quad (20)$$

where θ is the step function (defined as $\theta=0$ for $P \geq P_c$ and $\theta=1$ for $P \leq P_c$). Note that (20) is a general state law, not only applicable to vesiculated magmas. In general, the critical pressure P_c at which gas exsolution begins will be attained inside the chamber depending on the chamber depth (lithostatic pressure) and on the volatile content. If this critical pressure is not attained, that is, if magma is not vesiculated, we have $\theta=0$ in (20), and the density is that of the liquid. In turn, liquid density ρ_l can be considered in general either as a constant or, depending also on the pressure, as indicated in (16). As the magma has values of the compressibility modulus β of 10-100 GPa [Touloukian *et al.*, 1981], we will assume that the magma behaves as an incompressible flow (of constant density) below the exsolution level. Hence liquid density ρ_l is set as a constant in all computations, and the state law (20) becomes

$$\rho = \rho_l = \text{const} \quad (20')$$

below the exsolution level

$$\rho = \frac{\rho_l}{1 + \left(\frac{\rho_l QT}{P} \right) \left(\frac{W\% - sP^m}{1 - sP^m} \right)} \quad (20'')$$

above the exsolution level.

Figure 2 shows the density profile of the magma predicted by the equation (20.b') for a rhyolitic magma with $\rho_l = 2400 \text{ kg/m}^3$ at 850°C . Note how the magma density is constant below the exsolution level and descends progressively above this level due to the volatile exsolution induced by the pressure decrease. Equation (20') has been also recently proposed by Bower and Woods [1997] as a suitable state law for the magma-volatile mixture.

3. Numerical Method

3.1. Procedure

The set of equations (11), (12), and (20) are numerically solved using a finite element method (FEM). The characteristics of the method are briefly summarized here. A full description of the algorithm is given by Codina *et al.* [1998]. Let us write the conservative equations (11) and (12) as

$$\frac{\partial \rho}{\partial t} + \frac{\partial (\rho u_k)}{\partial x_k} = 0 \quad (21)$$

and

$$\frac{\partial U_i}{\partial t} = M_i - \frac{\partial p}{\partial x_i} \equiv R_i \quad (22)$$

where R_i is the i th component of the spatial operator and the body force and

$$M_i \equiv - \frac{\partial}{\partial x_j} (\rho u_i u_j - \tau_{ij}) + \rho g_i \quad (23)$$

First, transport equations are time discretized using a traditional finite differences scheme. Second, the space discretization is done using a Galerkin FEM method. The momentum equation has a convective contribution appearing in (23) that could lead to numerical oscillations when discretizing the space using a standard Galerkin formulation. In this algorithm, the equations with convective terms are time discretized along the characteristics, thus providing a consistent artificial diffusion when the spatial discretization is done. This leads to the following time-discretized equations:

$$\frac{\Delta \tilde{U}_i^n}{\Delta t} = M_i^{n+\vartheta_3} - \frac{\Delta t}{2} u_k^n \frac{\partial R_i^n}{\partial x_k} \quad (24)$$

$$\frac{\Delta p^n}{\Delta t} = - \frac{\partial}{\partial x_i} \left(U_i^n + \vartheta_1 \Delta \tilde{U}_i^n - \vartheta_1 \Delta t \frac{\partial p^{n+\vartheta_2}}{\partial x_i} \right) \quad (25)$$

$$\frac{\Delta U_i^n}{\Delta t} = \frac{\Delta \tilde{U}_i^n}{\Delta t} - \frac{\partial p^{n+\vartheta_2}}{\partial x_i} \quad (26)$$

where Δt is the time step size, the superscripts denote time step level, and ϑ_1 , ϑ_2 , and $\vartheta_3 \in [0,1]$. Equations (24) and (26) come from the splitting of the momentum equation (discretized in time along the characteristics). We use a fractional step method thus introducing a new unknown to the problem, the fractional momentum \tilde{U} . This technique was first proposed by Chorin [1967] and Temam [1969] for incompressible flows and allows the use of equal interpolation spaces for pressure and velocity fields. The splitting of the linear momentum equation produces a stabilizing effect on the pressure, eliminating the need for compatibility conditions, the Babuška-Brezzi conditions, when the incompressible limit is reached. This is an important point: in our model magma behaves either as a compressible flow or as a incompressible flow depending on pressure, so that we need an algorithm able to solve for both kind of flows. Once time discretization is done, the weak form of the equations is obtained, that is, time-discretized equations are projected into the usual space of test functions and integrated over the domain Ω . Integrals with spatial second-order derivatives are integrated by parts. Finally, we can discretize the space using a standard Galerkin method where the test functions are equal to the shape functions. In order to stabilize the numerical scheme at high Reynolds number (convection dominated flows), the time derivatives are discretized along the characteristics [Codina *et al.*, 1998], thus introducing the required artificial diffusion. This allows one to deal with high Reynolds number flows. Once the spatial discretization has been

performed, the equations can be written in matrix form:

$$\mathbf{M} \frac{\Delta \bar{\mathbf{U}}^n}{\Delta t} = \bar{\mathbf{F}}_1 - \mathbf{K} \bar{\mathbf{U}}^{n+\vartheta_3} \quad (27)$$

$$\mathbf{M}_\alpha \frac{\Delta \bar{\mathbf{P}}^n}{\Delta t} + \vartheta_1 \Delta t \mathbf{L} \bar{\mathbf{P}}^{n+\vartheta_2} = \bar{\mathbf{F}}_2 \quad (28)$$

$$\mathbf{M} \frac{\Delta \bar{\mathbf{U}}^n}{\Delta t} = \mathbf{M} \frac{\Delta \bar{\mathbf{U}}^n}{\Delta t} - \mathbf{G} \bar{\mathbf{P}}^{n+\vartheta_2} + \bar{\mathbf{F}}_3 \quad (29)$$

where \mathbf{U} , \mathbf{P} , and \mathbf{U} are the vectors of nodal unknowns and \mathbf{M} , \mathbf{K} , \mathbf{G} , and \mathbf{L} are the standard mass matrix, the matrix coming from the viscous and convective terms in the fractional momentum equation, the matrix coming from the gradient operator, and the matrix coming from the Laplacian operator respectively. The matrix \mathbf{M}_α in the discrete version of the continuity equation (28) is defined as

$$\mathbf{M}_{\alpha,ij} \equiv \int_{\Omega} \alpha N_i N_j d\Omega \quad (30)$$

where N_i is the shape function associated to the i th node of the FE mesh and α depends on the state law considered. Using the state law (20), one gets

$$\alpha = \frac{\rho_l}{p^n + \rho_l Q T \left(\frac{W\% - s(p^n)^m}{1 - s(p^n)^m} \right) \theta (p_c - p^n)} \quad (31)$$

Finally, the nodal vectors \mathbf{F}_1 , \mathbf{F}_2 , and \mathbf{F}_3 denote a vector which is known when solving each time step in any particular equation. These vectors contain terms with boundary integrals in which some boundary conditions are prescribed. Each time step is solved as follows: (1) Calculate a time increment Δt ; (2) calculate the fractional momentum using equation (27); (3) calculate the pressure using equation (28); (4) calculate the density using the state law (20), and (5) calculate the momentum using equation (29). Some remarks concerning the implementation of the algorithm are required. Three parameters θ_1 , θ_2 , and θ_3 , with $\theta_i \in [0,1]$ have been introduced when discretizing in time the momentum equation (θ_3), the continuity equation (θ_1), and during the splitting (θ_2). Depending on its values the algorithm becomes explicit, semi-implicit or fully implicit. If $\theta_2 = \theta_3 = 0$, the algorithm is explicit. In this case, the incompressible regime cannot be dealt with, since it can be shown that (28) becomes an identity when $\alpha = \rho_l / \rho$ in (31). This reflects the fact that pressure must be treated implicitly for incompressible flows. If $0 < \theta_2 \leq 1$ and $\theta_3 = 0$, the scheme is semi-implicit, and either pressure or density can be considered as unknowns except in the incompressible limit. The semi-implicit scheme can be applied either in compressible flows or in incompressible flows and using the same interpolation functions for velocity and pressure (density) fields. In our case, the implicit scheme with the pressure as unknown is used for two reasons. First, because the state law (20) gives density as a function of pressure directly, while pressure as a function of density cannot easily be obtained, and second, because neither the boundary conditions at the outlet or the initial conditions are known in terms of density. The critical time step Δt for a convection-diffusion equation can be computed as [Hirsch, 1990]

$$\Delta t = \frac{F_s}{\frac{4\varepsilon}{h^2} + \frac{|\bar{u}|}{h}} \quad (32)$$

where F_s is a safety factor (less than 1 in the explicit scheme) and h is the characteristic element dimension. For the continuity equation, $\varepsilon = 0$, while for the momentum equation, $\varepsilon = \mu / \rho$, μ being the viscosity. If a stationary problem were considered, discrete time advance could be different for each equation and could be local or global. In our case, the time advance must be uniform because a transient problem is studied. Substituting typical magmatic values in (32), one gets $\Delta t \approx F_s / 10$, so that the maximum time step able to ensure the stability of the algorithm is extremely small. Due to this limitation, we performed all simulations using the fully implicit option ($\theta_1 = \theta_3 = 0$), where F_s can achieve values up to 20 without any stability problem. Due to the small critical time step, one has to solve the whole set of equations to advance just 2s in real time. Thousands of time steps are required to simulate the entire withdrawal process. Examples of the reliability of the code are given in Zienkiewicz *et al.* [1995].

3.2. Boundary and Initial Conditions

Magma inside the chamber is always in a subsonic regime [Spera, 1984]. Under these circumstances, either pressure or density must be prescribed as a boundary condition at the computational outlet [e.g., Hirsch, 1990]. The coupling between the modeling of magma ascent, magma chamber processes, and the mechanical problem of host rock behavior should be considered because the processes which occur in each of these domains affect the dynamics of the others. When each of these domains (chamber, conduit, and host rock) is modeled as an isolated system, some unknown boundary condition must be assumed. Thus, for instance, most of the conduit models found in the literature assume a pressure at the conduit entrance when in fact it is an unknown of the problem. In our computations, the outlet is set at 500 m above the entrance of the conduit, and therefore the computational outlet does not coincide with the physical one (the vent). We cut the computational domain in the middle of the conduit in order to ensure the validity of equation (20), that is, in order to keep the flow within the bubbly flow regime (see the appendix). Pressure at the computational outlet is prescribed to the lithostatic, despite the fact that conduit models predict a magma pressure drop below the lithostatic value inside the conduit [Papale and Dobran, 1993, 1994]. However, this assumption can be justified in gross terms because the average stress field within the conduit cannot be too far from lithostatic in order for it to remain open against compression of the country rocks. Another possibility would be to not cut the conduit and therefore to set the pressure at the vent. This would probably produce a more realistic boundary condition. However, as discussed in the appendix, the state law (20) would not be realistic physically because the assumption of homogeneity would not be valid above the fragmentation level.

Boundary and initial conditions considered in our problem are schematically illustrated in Figure 3. The no-slip condition (velocity vector equal to zero) is imposed at the chamber walls ($u = v = 0$), while only the horizontal component of the velocity is set to zero at the symmetry axis ($x = 0$). Initial values are assigned to the velocity field and pressure. Magma is

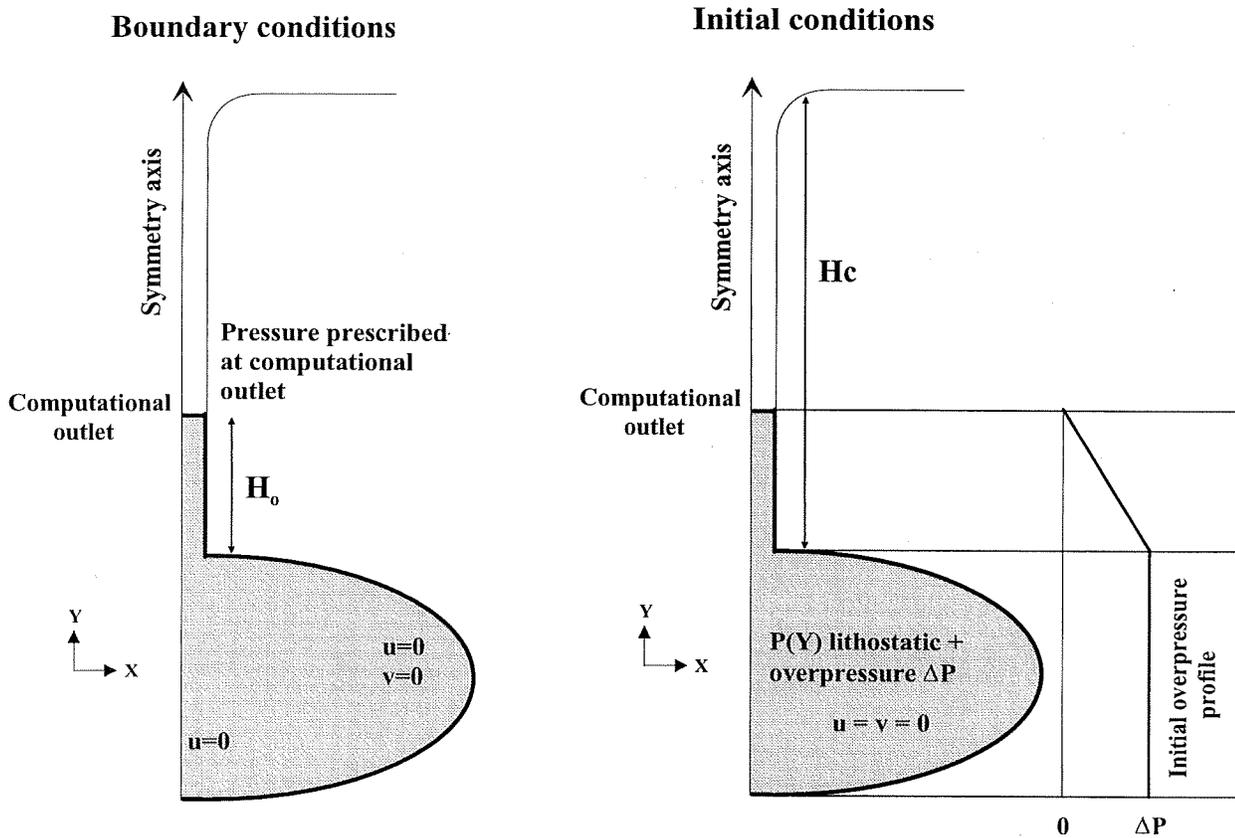


Figure 3. Schematic cartoon showing boundary and initial conditions. The computational domain is shown in grey. Pressure is the lithostatic value at the computational outlet (located at depth H_0 above the conduit entrance), and a no-slip condition ($u = v = 0$) is imposed at the chamber walls. The initial pressure profile is the lithostatic plus an overpressure ΔP . This overpressure is the driving force in triggering the eruption and is assumed to be equal to the tensile strength of the host rock inside the chamber and to decrease linearly along the conduit.

assumed to be at rest before the eruption, so that the initial velocity field is set to zero in the whole domain. The initial pressure distribution is assumed to be lithostatic with an overpressure inside the chamber equal to the tensile strength of the host rock. The conduit has an initial overpressure that decreases linearly from a value equal to the tensile strength of the rock at the inlet to zero at the computational outlet. It should be pointed out that the flow is induced by the pressure gradient in the conduit and is therefore not a departure from a state of equilibrium unstable, as would, for example, be the case for thermal convection. The magnitude of the pressure gradient at the beginning of the conduit depends on how the overpressure in the chamber is dissipated along the conduit, and this in turn depends on its degree of obturation. We have assumed that this overpressure is dissipated across 500 m. It might be thought that the initial pressure gradient is balanced by a porosity-like force with a small porosity coefficient and a negligible initial velocity. It is in this state that the sudden opening of the conduit takes place. A constant overpressure within the conduit with a sudden drop to zero at the computational outlet could also be assumed for the sudden conduit opening. However, the high-pressure gradient in such situation at $t = 0$ gives rise to numerical instabilities.

4. Numerical Results

We have focused on modeling magma chambers during Plinian central vent eruptions in order to find how pressure, density, and mass discharge rate at the conduit entrance evolve with time. Many numerical simulations have been done considering different values for the parameters involved. Any numerical simulation is done by defining a chamber-conduit configuration and the magma properties (initial overpressure, viscosity, volatile content, etc.). The relevant quantities are computed over time as the simulated eruption proceeds. The eruption rate is obtained by computing the integral

$$\int_S \rho \bar{v} \cdot d\bar{s}$$

over the conduit section S , where ρ is the mixture density given in (20') and \bar{v} is the velocity. The integral of the eruption rate with respect to time gives the total mass erupted. The position of the exsolution level is given by the critical pressure P_c , found using (18). In order for the overpressure to exist in a closed system, the exsolution level must occur inside the chamber before the eruption. Therefore the gas content is determined from an assumed lithostatic pressure or a magma chamber depth.

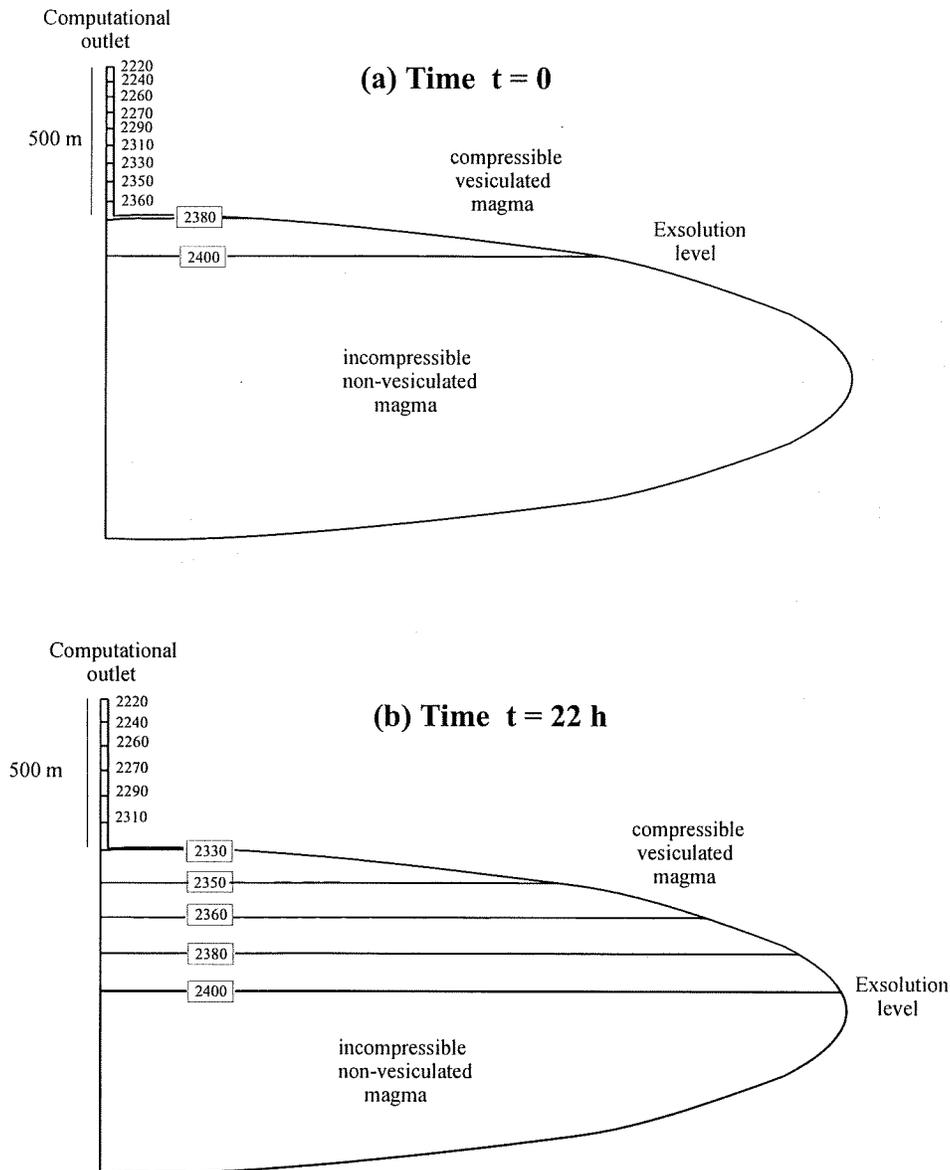


Figure 4. Density profiles for an ellipsoidal magma chamber with a major axis $a=2.5$ km, a minor axis $b=0.5$ km, a volume $V=4/3\pi a^2 b \cong 14$ km³, and a conduit radius of 25 m. The density is given in kg/m³. Volatile content is set to 4.3 in wt% and the lithostatic pressure at the computational outlet is 105 MPa. The nonvesiculated magma has a density $\rho = \rho_l = 2400$ kg/m³ (a) Density profile at the onset of the eruption ($t = 0$) when only the uppermost part of the chamber is vesiculated. (b) Density profile after 22 hours of simulated eruption. Note the drop in the exsolution level due to the pressure decrease as the eruption proceeds.

4.1. A General Overview

In this section, the results for a particular numerical simulation are shown as an example of applicability of the model. Although the results presented are obtained by applying the model to a particular case, we stress those characteristics which are expected to be found in other simulations. A rhyolitic magma having $\rho_l = 2400$ kg/m³, $T = 850^\circ\text{C}$, $s = 4.11 \times 10^{-6}$ Pa^{-1/2}, $m = 0.5$ and a water content of 4.3 in wt% is assumed. The hypothetical shallow magma chamber is elliptical and has a major axis $a = 2.5$ km, a minor axis $b = 0.5$ km, a volume $V=4/3\pi a^2 b \cong 14$ km³, and a conduit radius of 25 m. The chamber has an initial pressure equal to the lithostatic pressure plus an overpressure $\Delta P = 10$ MPa. This overpressure is a typical value of the tensile strength of the host rocks [Touloukian *et al.*,

1981]. Note that this is an approximation because there is no straightforward relation between the overpressure within an ellipsoidal reservoir and the tensile stress induced in the elastic medium surrounding. The effect of the initial overpressure is discussed below. The lithostatic pressure at the computational outlet is set to 105 MPa. Note that we set the lithostatic pressure rather than the chamber depth H_c in order to avoid assumptions about the host rock mean density. Thus, if the host rock density were 2500 kg/m³, the computational outlet should be located at 4280 m below the Earth surface ($H_c=4780$ m); if the host rock density were 2800 kg/m³, the computational outlet should be located at 3820 m below the Earth surface ($H_c=4320$ m).

All the numerical experiments yield similar results: the initial overpressure decreases exponentially as the eruption proceeds until it becomes zero and the eruption ceases. This pressure

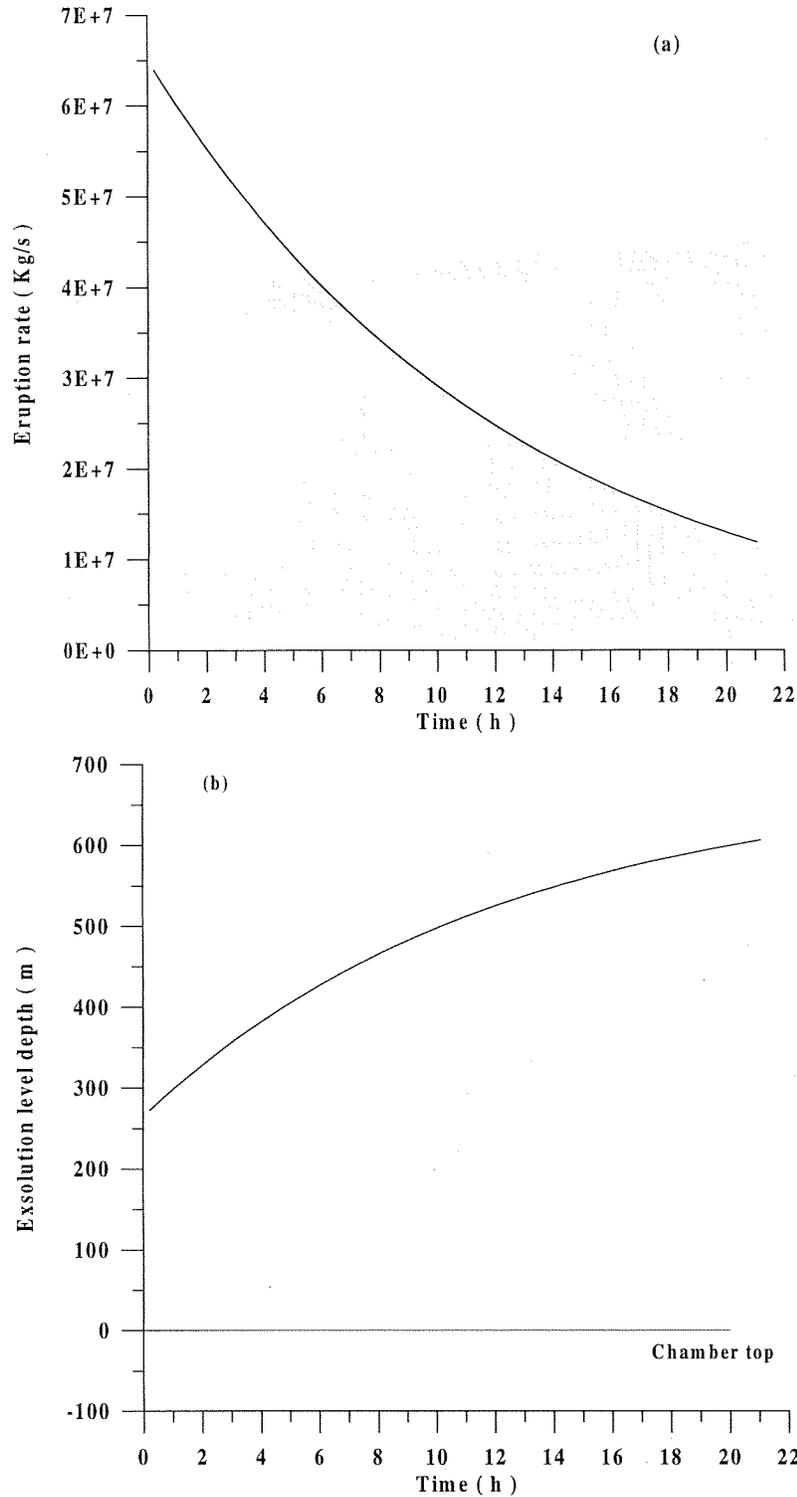


Figure 5. Results using the same chamber geometry and magma properties as in Figure 4. (a) Eruption rate as a function of time. (b) Position of the exsolution level as the eruption proceeds. The origin of the vertical axis is set at the top of the chamber (4 km below the surface) with the positive direction being downward.

decrease produces a drop in the exsolution level so that deeper parts of the chamber become progressively volatile oversaturated as the eruption proceeds. The upper oversaturated magma layer thus becomes progressively thicker and less dense. Figure 4 shows the density distribution for this elliptical magma chamber at the onset of the eruption and after 22 hours, when the initial overpressure has decreased by an order of magnitude.

Note that due to the pressure variations the position of the exsolution level descends hundreds of meters as the eruption proceeds. Figure 5 shows the eruption rate and the position of the exsolution level versus time for this particular simulation. A common characteristic of all the experiments is that the eruption rate reaches a peak immediately after the onset of the eruption and then decreases exponentially as the eruption proceeds.

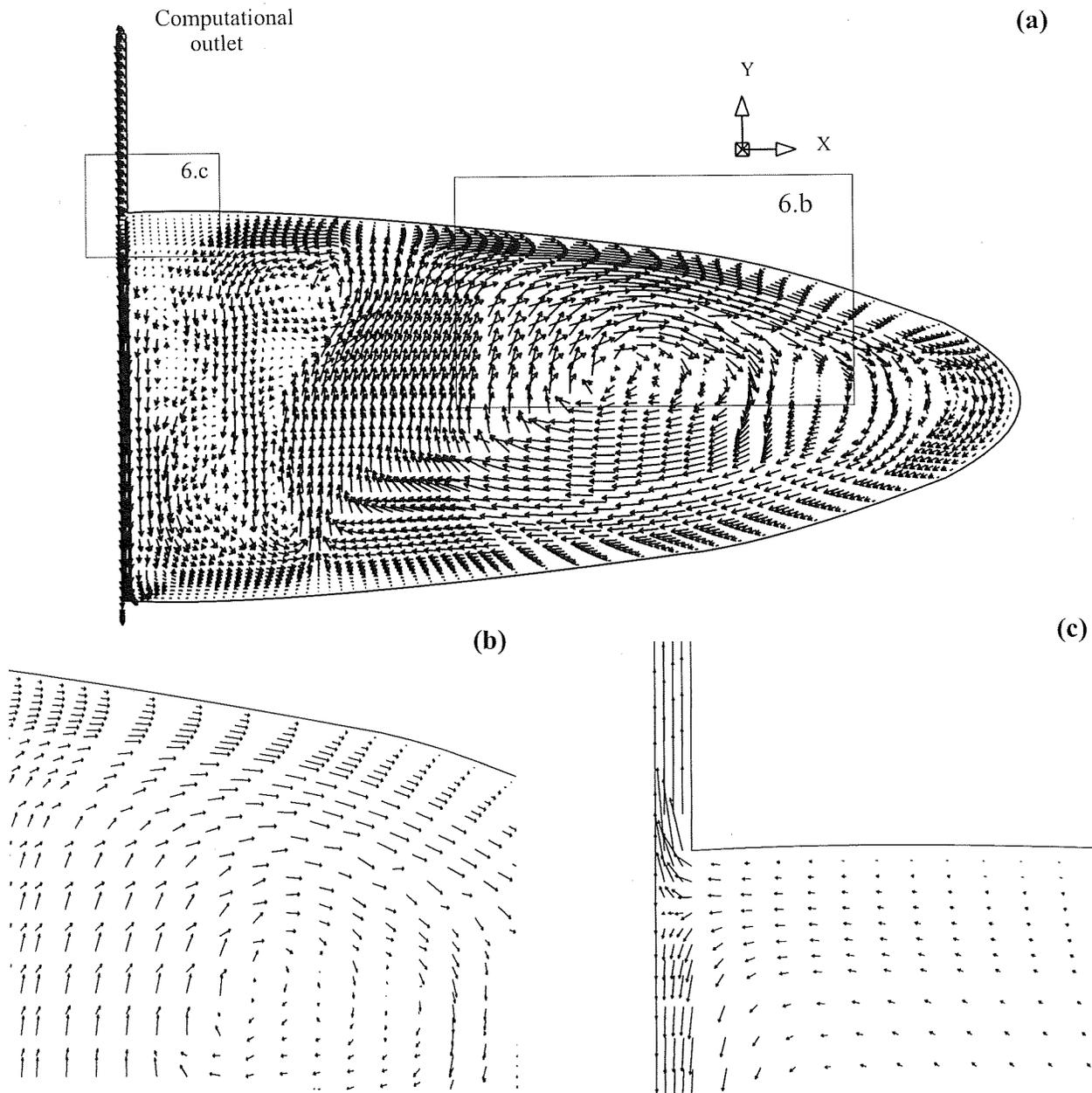


Figure 6. Velocity field inside the chamber after 22 hours of simulated eruption when using the same chamber geometry and magma properties as in Figures 4 and 5. Although the velocity field is unsteady, its temporal variations are slow so that trajectories approach streamlines. Thus Figure 6c illustrates how a parcel of magma reaches the conduit entrance laterally and is then either erupted or reinjected back into the chamber.

Nevertheless, it should be said that many Plinian eruptions present phases in which the eruption rate descends progressively only to then reach a new peak of intensity. These Plinian phases are commonly associated with tapping effects [*e.g.*, Dobran, 1992], where the magma that reaches the conduit entrance changes its properties (viscosity, gas content, etc.) as the eruption proceeds, as a consequence of preeruptive heterogeneities. As we have assumed a chemically homogeneous magma chamber composition, our model is not able to reproduce such oscillations.

Although we do not model any specific volcanic system, the numerical results obtained are in good agreement with values

observed in natural systems with comparable characteristics. For instance, in this example, we obtain a total erupted mass of approximately 3×10^{12} kg. This value is comparable with those found in explosive volcanic eruptions: 0.6×10^{12} kg in the 1980 Mount St. Helens eruption [Scandone and Malone, 1985], $5-7 \times 10^{12}$ kg in the Nisyros eruption [Barberi *et al.*, 1988] or 8.6×10^{12} kg in the 79 A.D. Vesuvius eruption [Barberi *et al.*, 1981].

Figure 5b shows the depth of the exsolution level versus time. As also illustrated in Figure 4, the position of the exsolution level descends as the chamber is depressurized. In this particular example, the exsolution level is located initially at 275 m below

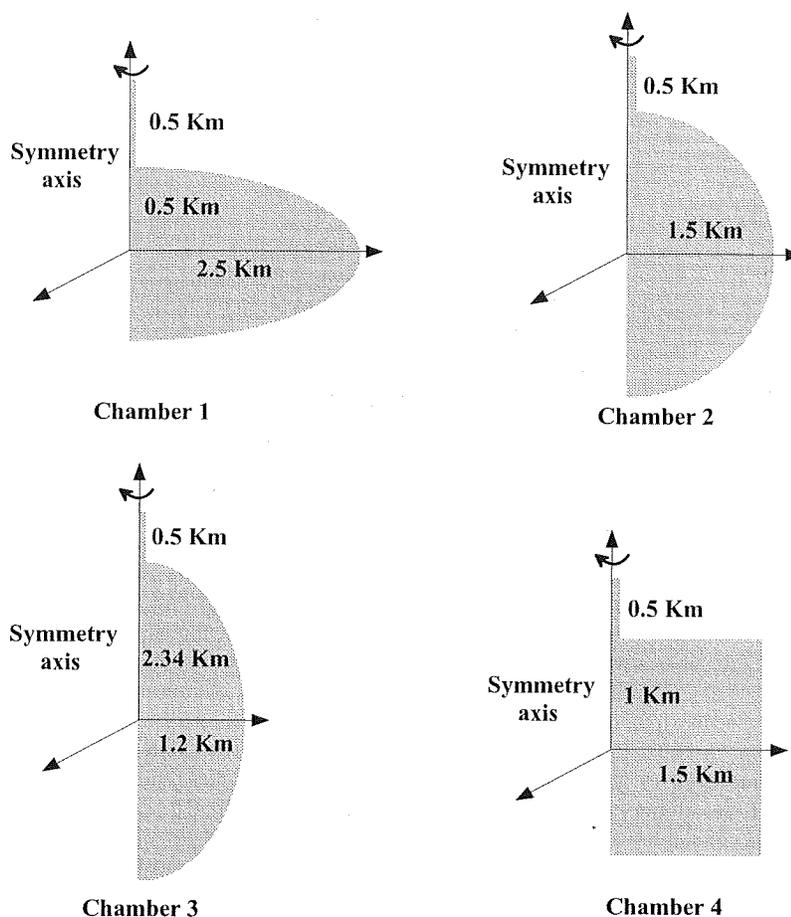


Figure 7. Schematic cartoon showing the magma chamber geometries considered: chamber 1 is an oblate ellipsoid, chamber 2 is spherical, chamber 3 is a prolate ellipsoid, and chamber 4 is cylindrical. All the chambers have equal volume $V \cong 14 \text{ km}^3$ and equal conduit radius $r = 25 \text{ m}$. The conduit has been cut 500 m above the chamber top to ensure the validity of (20').

the conduit entrance and moves 300 m after 22 hours, reaching depths of up to 600 m below the chamber top. Figure 6 shows the velocity field inside the chamber after 22 hours of eruption. The magnitude of the velocity field is determined directly by the initial assumed overpressure. Convection could also be driven by thermal effects, although in the present simulation these have not been taken into account. To account for these effects, buoyancy forces would have had to be accounted for in the region where the flow is incompressible.

A common feature found during the withdrawal of a chemically homogeneous magma chamber is that the magma reaches the conduit entrance laterally as reflected in Figure 6c. Note that although the velocity field is unsteady, its variations are given in a characteristic timescale of hours. This means that we can safely approach the streamlines given by the trajectories of the particles. A volume of magma reaching this point is partly erupted through the conduit and partly reinjected into the chamber. Thus, a parcel of magma initially emplaced at the chamber top moves first laterally into the vicinity of the conduit entrance and is then either erupted or forced to descend before again moving up. As the parcel descends, exsolved volatiles within it are progressively redissolved back into the magma due to the pressure effect on the volatile solubility in the melt. This

phenomenon has an important consequence: as we have assumed that magma behaves as an incompressible fluid below the exsolution level, any inflow into the undersaturated incompressible layer must produce an equal outflow in order to ensure the mass conservation. This movement is progressively transmitted to the neighborhood of the parcels through viscous stresses, so that finally the whole magma chamber is affected by this induced movement, as reflected in Figure 6a. This is a common feature of all the numerical experiments, though its intensity and characteristics are obviously strongly dependent on magma viscosity and on chamber geometry. This phenomenon could have important consequences on the eruption products because magmas initially located at different depths could be simultaneously erupted. *Trial et al.* [1992] suggested that the presence of different compositions in the same Plinian deposits is not necessarily the consequence of a mixing process but may result from the withdrawal process itself when preeruptive gradients existed in the chamber. Our homogeneous flow model does not allow us to draw conclusions about stratified magma chambers with preeruptive density gradients but could explain some chemical heterogeneities, such as those involving trace element abundances, which have negligible effects on magma density.

Table 1. Properties of the FEM Meshes

	Nodes	Elements	Boundary Elements	Element Type
Chamber 1	3190	3080	218	Q1
Chamber 2	2680	2580	198	Q1
Chamber 3	2680	2580	198	Q1
Chamber 4	2900	2775	248	Q1

All the elements are type Q1, that is, bilinear Lagrangian (four nodes per element).

4.2. A Parametrical Study

Many parameters such viscosity, gas content, magma chamber geometry or initial overpressure appear in the model, and its influences on the solution are not always clear. In this section, we aim to illustrate the influence of these parameters by showing how their values affect the results obtained in the numerical example previously described.

4.2.1. Chamber Geometry. In order to assess the importance of chamber geometry we consider four different magma chambers with volume $V \cong 14 \text{ km}^3$ (Figure 7). Chamber 1 is oblate, chamber 2 spherical, chamber 3 prolate, and chamber 4 cylindrical. In all cases, the FEM meshes are structured and made of linear Lagrangian elements (four nodes per element). Their properties are summarized in Table 1. The dimensions of these chambers are of the same order of magnitude as natural shallow magma chambers but are arbitrary in representing any particular volcanic system. The conduit has a radius of 25 m and again is terminated 500 m above the chamber because the state law (20') may not be applicable in the uppermost parts of the conduit. Figure 8 shows the eruption rate at the conduit entrance versus time for this particular numerical experiment. An important result is that although shapes of the curves are similar

in all four cases, the total mass erupted increases as the chamber becomes flatter and as the volumetric fraction found above the exsolution level increases. This can be explained by the fact that magma is nearly incompressible below the exsolution level (fully incompressible in our model) but that above this level, the mixture containing exsolved gas becomes much more compressible. Thus the total mass erupted corresponds only to the volume change in the upper compressible layer as a response to the pressure variation. In consequence, those chambers containing more magma in the upper oversaturated layer (like chambers 1 and 4 in our example) will erupt a greater mass of material for a given pressure decrease. The total erupted mass is obtained by computing the area below the eruption rate curves versus time. In this example, we achieve erupted fractions of the initial mass of 0.090, 0.077, 0.064, and 0.078 for chambers 1 to 4, respectively. Recently, *Bower and Woods* [1997] have predicted that in saturated magma chambers, the chamber overpressure is only relieved when a fraction of 0.01-0.1 of the initial mass in the chamber has erupted. The numerical results are thus in excellent agreement with those predicted by the analytical model of *Bower and Woods* [1997]. Nevertheless, the use of a numerical model allows us to go one step further since we are able not only to compute the total mass erupted but also the temporal variation of the eruption rate and the influence of the chamber geometry.

4.2.2. Volatile Content. The influence of volatile content on the eruption rates and on the pressure variations is now considered. All the examples are obtained using magma chamber 1 (oblate ellipsoid), an initial overpressure of $\Delta P = 10 \text{ MPa}$, a viscosity of $\mu = 10^5 \text{ Pa s}$, and a density of $\rho_l = 2400 \text{ kg/m}^3$. The eruption rates obtained and the evolution of the exsolution level are shown in Figure 9. Although a small change in the content of volatiles does not change substantially the mass discharge rate, it has an important effect on the position of the

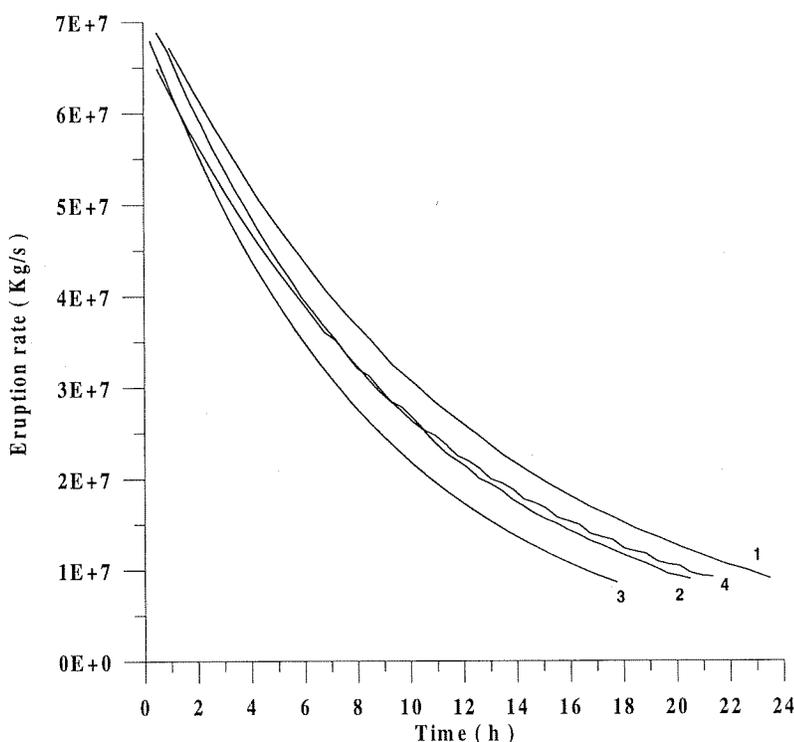


Figure 8. Eruption rate versus time for different magma chamber geometries. The chambers are those shown in Figure 7.

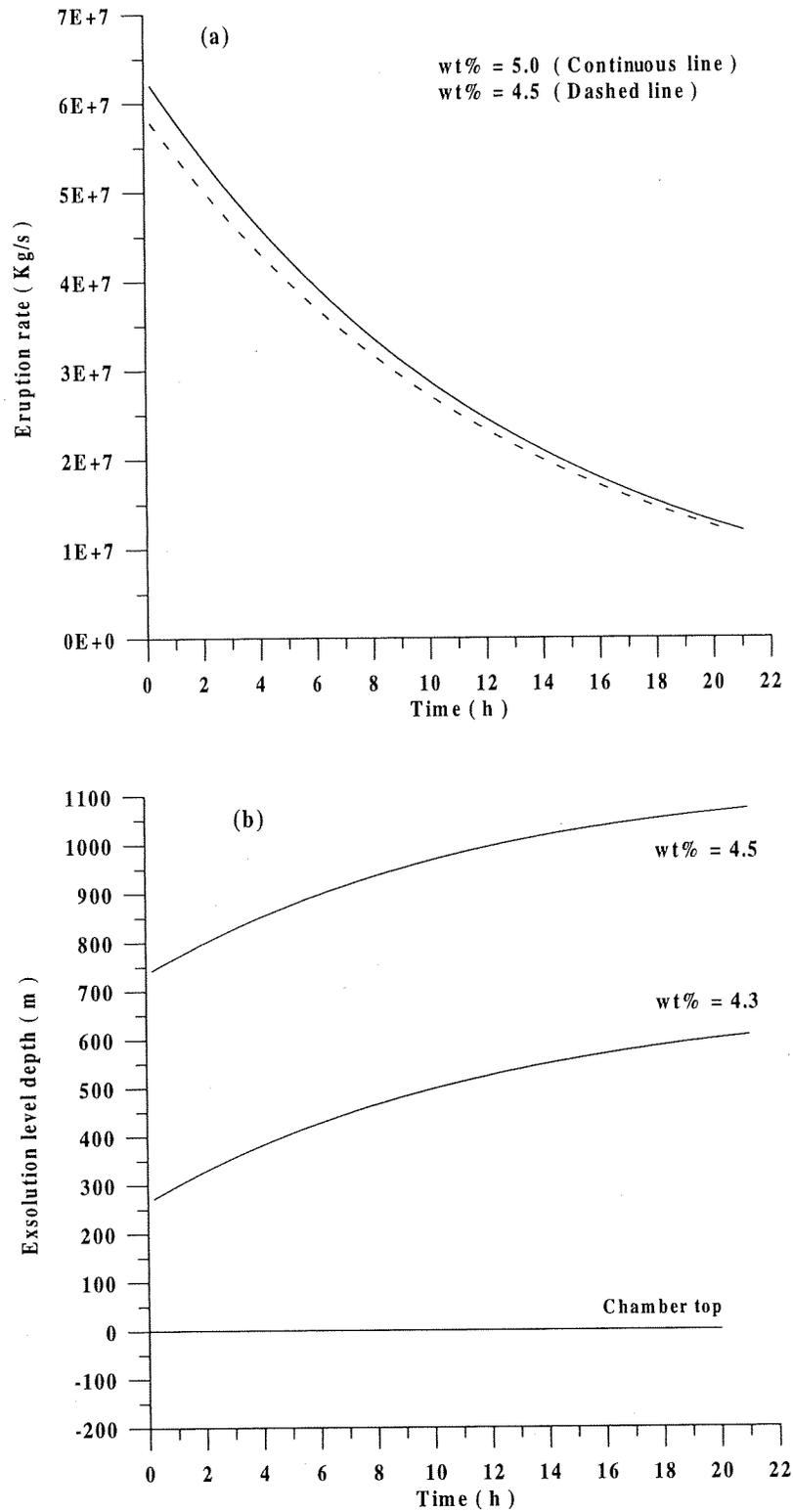


Figure 9. Results using chamber 1, an initial overpressure of $\Delta P = 10$ MPa, a viscosity $\mu = 10^5$ Pa s, and a liquid density $\rho_l = 2400$ kg/m³. (a) Eruption rate versus time for water contents of 4.5 and 5.0 in wt %. The exponential decrease is only weakly dependent on volatile content. (b) Position of the exsolution level for water contents of 4.3 and 4.5 in wt %. The origin of the vertical axis is set at the top of the chamber. Note that the shape of the function does not vary substantially when the volatile content is changed, although the position of the exsolution level is dependent on the volatile content.

exsolution level. A small increase in the volatile content produces a substantial deepening of the exsolution level. The total displacement is insensitive to the volatile content, but the initial and final positions are strongly dependent on its value.

4.2.3. Viscosity. Figure 10 shows the effect of varying viscosity on the magma overpressure, the position of the exsolution level, and the mass discharge rate when other parameters are kept constants. Figure 10 illustrates that the less viscous the magma, the faster the chamber becomes depressurized. The position of the exsolution level changes faster for less viscous magmas despite the fact that the initial and final positions are the same for a given pressure decrease ($\Delta P = 10$ MPa in this case). Although the total erupted mass is only weakly dependent on viscosity, this parameter plays a major role in controlling the duration of the eruption. The more viscous the magma, the longer the eruption and the lower the erupted mass peak, that is, chambers containing a less viscous magma are more rapidly withdrawn and have a greater intensity peak.

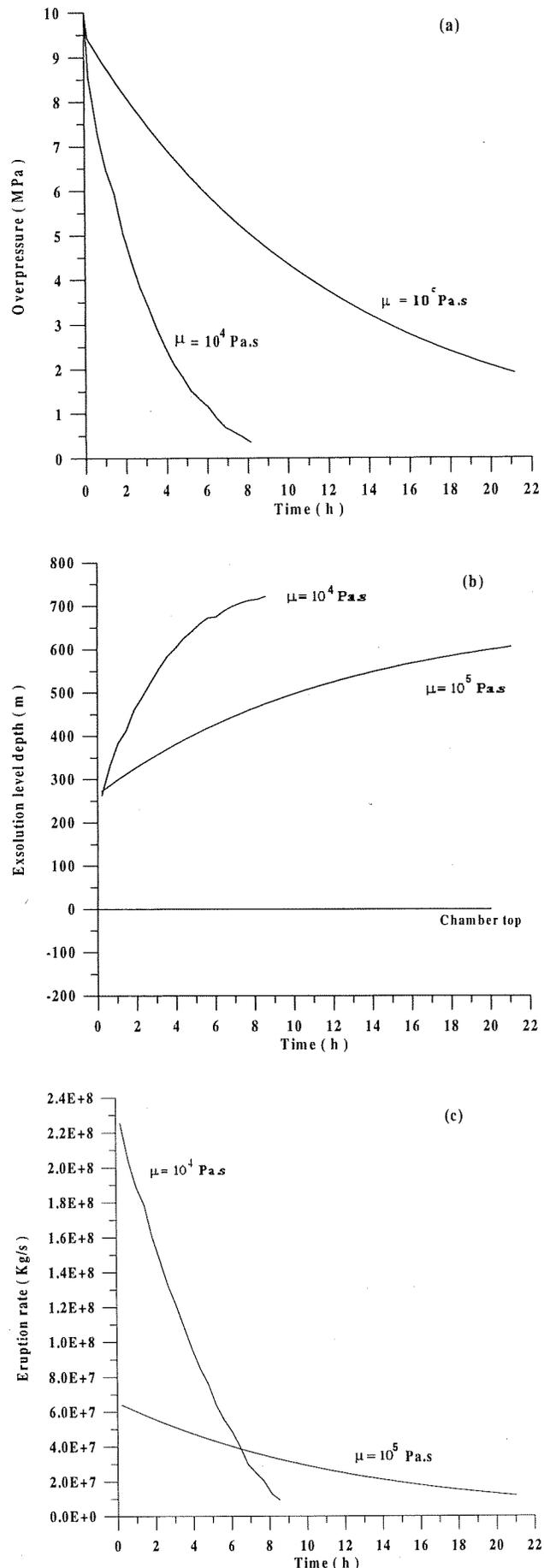
4.2.4. Initial Overpressure. The influence of the initial overpressure is now considered. All of the examples presented are for a chamber 1 geometry (Figure 7), with a viscosity of $\mu = 10^5$ Pa s, a density of $\rho_l = 2400$ kg/m³ and a water content of 4.3 in wt %, but with initial overpressures of 5, 10, and 15 MPa respectively (Figure 11). The total drop in the exsolution level is strongly dependent on the initial overpressure, as is its initial position, while its final position is invariant. Obviously, an increase in the initial overpressure yields an increase in the displacement of the exsolution level. This leads to a greater erupted mass because a significant volume of magma changes from volatile-undersaturated (incompressible) to volatile-oversaturated (compressible) during the withdrawal process.

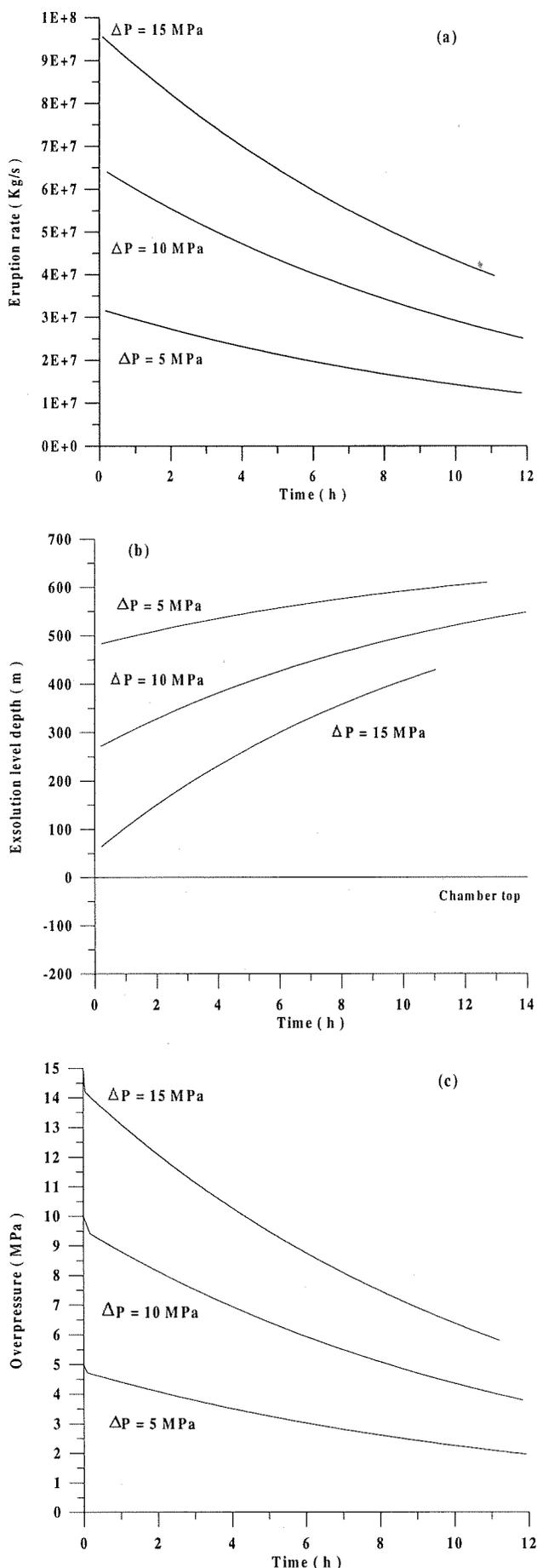
5. Conclusions

We have developed an axisymmetrical numerical model in order to find the temporal evolution of pressure, the position of the exsolution level, the velocity field, and the eruption rate during the withdrawal from a shallow volatile-rich felsic magma chamber, where the overpressure is due to volatile oversaturation. In this model, magma behaves as an incompressible fluid below the exsolution level and as a compressible gas-liquid mixture above this level. In this domain, magma is treated as a homogeneous two-phase flow, and we propose a barotropic state law (20') as suitable for the mixture under magmatic conditions.

The resulting set of equations is solved numerically using a finite element method (FEM). We propose an algorithm able to deal simultaneously with compressible and incompressible flows. This uses a fractional step method which allows the use of equal interpolation spaces for pressure and velocity fields.

Figure 10. (a) The magma overpressure, (b) the position of the exsolution level, and (c) the mass discharge rate for two different viscosity values of $\mu = 10^4$ Pa s and $\mu = 10^5$ Pa s, with the other parameters kept the same as in Figure 9. The less viscous the magma, the faster the chamber becomes depressurized and the faster is the descent of the exsolution level. Chambers containing less viscous magma are thus more rapidly withdrawn and have a greater peak intensity of erupted mass.





The artificial diffusion required in the equations with convective terms is obtained by discretizing in time along the characteristics. By assuming axial symmetry, we solve a three-dimensional problem like a two-dimensional using cylindrical coordinates.

The results obtained show that the pressure at the conduit entrance decreases exponentially as the eruption proceeds. This pressure decrease produces a shifting of the exsolution level, so that deeper parts of the chamber become progressively volatile oversaturated. As previously proposed by *Triel et al.* [1992] for open magma chambers in which the trigger for the eruption is magma inflow into the chamber, our numerical experiments suggest that in closed, homogeneous magma chambers, magma from different depths can be erupted simultaneously. This may explain some chemical heterogeneities in the erupted products, such as different abundances of trace elements, which are not necessarily related to magma density. The total erupted mass predicted by the experiments is in good agreement with the results of *Bowers and Woods* [1997], who find that in closed volatile-saturated magma chambers, a mass fraction of 0.01-0.1 of the initial mass is erupted. In addition, the numerical simulations allow us to predict not only the total mass erupted but also its temporal variation for any chamber geometry and set of magma properties.

For chambers of equal volume and magma properties, the qualitative behavior is insensitive to the chamber geometry. However, the total erupted mass increases as the chamber becomes flatter because flat chambers have a major volumetric fraction of compressible magma. It should be pointed out that geometry would play a major role when considering the coupling with the mechanical problem. In natural systems, the stress field around the chamber is directly related to its geometry. The stress field controls the opening, propagation and closure of those fractures that allow the magma to erupt. By assuming rigid walls, our model partially neglects the influence of geometry.

A small change in the magma volatile content does not substantially change the mass discharge rate but has an important influence on the position of the exsolution level. A small increase in volatile content produces a substantial deepening of the exsolution level. The exsolution level drops faster in less viscous magmas although the initial and final positions are insensitive to viscosity. The total erupted mass is weakly dependent on viscosity, although viscosity plays a major role in controlling the duration and the peak intensity of the eruption. An increase in the initial overpressure, here associated with the tensile strength of the host rock, produces an increase in the total displacement of the exsolution level. As a consequence, when the initial overpressure is high we see a greater erupted mass because a major volumetric fraction of magma changes from volatile-undersaturated (incompressible)

Figure 11. (a) The mass discharge rate, (b) the position of the exsolution level, and (c) the overpressure evolution for different initial overpressures of 5, 10, and 15 MPa. An increase in the initial overpressure can be accounted for by an increase in the tensile strength of the host rock. The total drop in the exsolution level depends strongly on the initial overpressure, although its final position is the same in all cases. Increasing the initial overpressure yields a greater erupted mass because a major volumetric fraction of magma changes from undersaturated to oversaturated during the depressurization process.

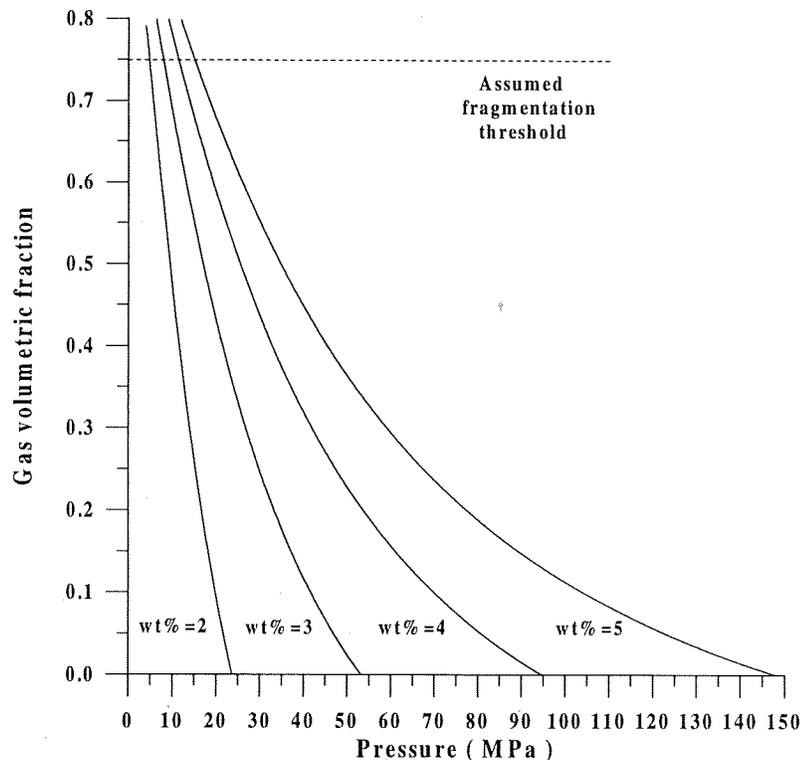


Figure 12. Gas volumetric fraction α predicted by the state law (20). Assuming a fragmentation threshold for $\alpha=0.75$ [e.g., Sparks, 1978], the domain in which the state law (20) is valid can be estimated quantitatively. Above the fragmentation level, at the uppermost part of the conduit, the bubbly flow regime cannot be maintained and the assumptions of the state law are physically unrealistic. Results for volatile contents of 2, 3, 4, and 5 in wt % respectively. The fragmentation level predicted by (20) is constrained to lie in the uppermost part of the conduit, even for high volatile contents.

to volatile-oversaturated (compressible) during the withdrawal process.

Our model constitutes a first attempt to numerically model the dynamics and temporal evolution of the most relevant physical parameters during magma withdrawal from a closed chamber. The relative importance of convective flow induced by overpressure and convective flow induced by thermal convection requires further study. Future models should also take account of chemical inhomogeneities in the magma, as well as the coupling with the mechanical problem. If a constitutive model for the host rock were coupled with this model, it should be possible to determine under which conditions of magma composition, chamber geometry, chamber depth, stress field, etc., a Plinian central vent eruption could generate a ring fault system and become a caldera-forming eruption.

Appendix. Discussion on Boundary and Initial Conditions

The barotropic state law (20) has been deduced assuming that the gas bubbles and the liquid are in mechanical equilibrium. As magma flows through the conduit, the gas bubbles grow by diffusion and by decompression. Growth by coalescence is inhibited in silicic magmas because gas bubbles are unable to come in contact by displacing surrounding high-viscosity magma. The volumetric fraction of gas progressively increases as magma ascends and eventually reaches a critical bubble packing leading to magma fragmentation. Above this

fragmentation level, the bubbly regime cannot be sustained, assumptions of mechanical equilibrium are no longer valid, and the state law (20) becomes physically unrealistic. The gas volume fraction α can be computed combining (14), (15), and (20). The result is shown in Figure 12, and reflects the need to limit the computational domain in order to ensure the validity of equation (20). We set the computational outlet at 500 m above the conduit entrance. This is an arbitrary choice but sufficient to ensure the validity of the state law even for very shallow magma chambers.

Notation

We use $f^{n+\theta} \equiv \theta f^{n+1} + (1-\theta) f^n$ and $\Delta f^n \equiv f^{n+1} - f^n$ for any function f . The superscript denotes time step level. Bold vectors stand for vectors of nodal unknowns, and \mathbf{F}_i denotes a vector which is known at the moment of solving a particular equation.

a	semimajor axis (ellipsoidal chamber).
b	semiminor axis (ellipsoidal chamber).
\vec{g}	gravity vector.
F_s	safety factor.
H_0	computational conduit height.
H_c	magma chamber depth.
m	solubility law exponent.
m_d	mass of dissolved gas.
m_g	mass of exsolved gas.
m_l	mass of liquid magma.

m_s	mass of the melt.
m_v	mass of volatiles.
M	total mass.
n	molecular mass (18 kg/kmol for H ₂ O).
n_g	exsolved gas mass fraction.
P	pressure.
P_c	critical (exsolution) pressure.
Q	gas constant.
r	conduit radius.
R	universal gas constant (8314 J ^o K kmol)
s	solubility law constant.
T	temperature.
\vec{u}	velocity vector.
u	x velocity component.
\vec{U}	momentum vector
\vec{U}	fractional momentum vector.
v	y velocity component.
$W\%$	volatile mass fraction.
α	gas volumetric fraction.
β	compressibility modulus.
ΔP	magma overpressure.
χ	solubility law ; maximum allowable mass fraction of gas into solution
μ	viscosity.
ρ	mixture density.
ρ_g	gas density.
ρ_l	liquid density.
ρ_s	melt (liquid plus exsolved gas within) density.
σ_{ij}	Cauchy stress tensor components.
τ_{ij}	viscous stress tensor components.
θ	step function.
θ_i	parameters of the algorithm.

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