

Modelling unsaturated soils with the Material Point Method. A discussion of the state-of-the-art*

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Abstract

Many natural hazards involve large deformations of unsaturated soils, e.g. rainfall-induced landslides, embankment collapses due to wetting, seepage-induced instabilities of dams and levees, etc. The study of these phenomena requires accounting for the complex hydro-mechanical interactions between solid skeleton and pore fluids and modeling large deformations to predict the post-failure behaviour, which poses significant computational challenges. In recent years, several hydro-mechanical coupled MPM formulations were developed to model saturated and unsaturated soils. These approaches are slightly different in terms of governing equations, integration schemes and have been implemented in different MPM software; thus, they benefit from various computational features. The purpose of this paper is to present an overview of the available MPM approaches to model unsaturated soils discussing differences and similarities of the formulations and their impact on the results under different conditions in a range of geotechnical appli-

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cations. In addition, the effect of partially saturated conditions on the critical time step in explicit numerical integration schemes is studied for the first time. Different analytical expressions are derived and compared with the numerical results.

Keywords: unsaturated soils, material point method, large deformations, dynamic coupled analysis, critical time step, state-of-the-art

1. Introduction

Many geotechnical applications involve multi-phase interactions and large deformations of the ground. This is the case of rainfall-induced landslides, embankment collapses due to seepage or wetting, shallow foundation stability, and penetration problems in unsaturated soils, etc. Soil is a mixture of solid grains and pore fluids (liquid and/or gas) that interact with each other. In some cases, e.g. when considering dry or saturated soils in fully drained or fully undrained conditions, the interactions among the different phases can be simplified or even neglected. However, when the soil is in partially saturated conditions (i.e. gas and liquid coexist in the pores) or in saturated conditions when simultaneous generation and dissipation of fluid pore pressure is present, the mechanical and hydraulic response of the soil is fully controlled by the coupling between pore fluids and solid skeleton controls. Hence, the hydromechanical interactions need to be considered to accurately capture the material behaviour. The numerical analysis of these problems requires implementing multi-phase formulations and the consideration of appropriate boundary conditions (BCs) that can reproduce realistic scenarios of water infiltration, seepage, ponding, and changes in the water table.

Historically, fully-coupled hydro-mechanical formulations have been developed and adapted to traditional numerical frameworks such as finite elements (FEM)

20 and finite differences (FDM) methods. However, it is well known that these tools
21 are limited to modeling relatively small deformations because of mesh tangling
22 limitations. Alternatively, the numerical tools used for the analysis of post-failure
23 runouts (i.e. propagation models) typically use rheological models that are difficult
24 to trace back to soil's pre-failure conditions and commonly rely on depth-averaged
25 integration schemes that cannot capture the failure initiation [1, 2, 3]

26 Several techniques have been proposed to overcome such difficulties e.g. A Ar-
27 bitraryLagrangian Eulerian (ALE), Coupled Eulerian – Lagrangian (CEL), Particle
28 Finite Element Method (PFEM), Finite Element Method with Lagrangian Integra-
29 tion Points (FEMLIP), Smooth Particle Hydrodynamics (SPH), Material Point
30 Method (MPM), etc [e.g. 4, 5, 6, 7, 8]. A review of these methods and their ap-
31 plicability for the study of landslide mass movements can be found in [9]. Among
32 them, the Material Point Method (MPM) has recently increased its popularity in
33 the geotechnical community. The MPM is a particle-based numerical method es-
34 pecially well suited for the simulation of large deformations in history-dependent
35 materials. It was initially developed in the framework of fluid mechanics under
36 the name of Particle-In-Cell method (PIC) by Harlow [10] in Los Alamos National
37 Laboratory. The basis of the method was to represent the flow of a continuum
38 mass by a collection of material points (MPs) that moved through a background
39 fixed mesh. It was not until 1994 that Sulsky et al. [8] extended the concept to solid
40 mechanics for single-phase materials. During the last twenty years, MPM has been
41 increasingly applied in the geotechnical engineering field for the study of soils and
42 granular earth material. In particular, slope instabilities, failure of earth retain-
43 ing structures, tunneling and underground collapses, soil penetration testing, pile
44 installation, scour, internal erosion, among others [e.g. 11, 12, 13, 14, 15, 16, 17].

45 Different MPM coupled multi-phase formulations have been developed to ad-
46 dress hydro-mechanical interactions between the solid skeleton and pore fluids.
47 These approaches are slightly different in terms of governing equations, integra-
48 tion schemes, and they benefit of different computational features. In particular,
49 this paper addresses the challenges of modeling hydro-mechanical interactions and
50 large deformations of unsaturated soils using the MPM framework. The objectives
51 are to provide an overview and discuss the available MPM approaches to model
52 solid-liquid-gas mixtures, highlighting differences and similarities of the formu-
53 lations, implementation aspects, and boundary conditions. Particular attention is
54 paid to explain the relevance and physical meaning of common simplifications and
55 their implications for the use of these approaches in practical applications. For the
56 first time, the effect of partially saturated conditions on the time step in explicit
57 MPM numerical integration schemes is studied. A set of analytical expressions for
58 the critical time step are derived and compared with numerical results. Addition-
59 ally, a discussion on existing applications is provided and the preliminary results
60 of potential future applications is presented.

61 The paper is organized into four different sections. First, the MPM formulations
62 are presented and discussed. Secondly, different implementation aspects are ex-
63 amined including boundary conditions and critical time step. Then, an overview of
64 geotechnical applications is presented. Finally, the conclusions are summarized.

65 **2. MPM formulations for unsaturated soils**

66 Unsaturated porous media consist of a combination of three phases (*ph*): solid
67 (*S*), liquid (*L*), and gas (*G*). The phases interact with each other determining the
68 mechanic and hydraulic behaviour of the material. The solid phase is made of

69 solid grains that constitute the solid skeleton while the fluid phases (i.e. liquid and
70 gas) fill the pore space.

71 In MPM, the material is considered as a continuum and is represented by a finite
72 number of material points (MPs) that move together with the reference domain
73 carrying all the information (e.g. stress, strain, mass, constitutive state variables).
74 The computational cycle is summarized in Fig. 1. At the beginning of each time
75 step, information carried by MPs is mapped to the nodes of the computational
76 mesh and nodal mass, nodal velocities, forces, and dragging terms are evaluated.
77 The main governing equations, usually the momentum balances, are solved in
78 terms of main unknowns, typically acceleration or velocity. The nodal solution
79 is interpolated using shape functions and the kinematic quantities and position
80 of the MPs are updated accordingly. Finally, strains and stresses among other
81 quantities are evaluated at the MPs using compatibility equations, mass balances,
82 and constitutive relationships. At the end of the computational cycle, the nodal
83 information is generally discarded. These features make the MPM a powerful tool
84 in geotechnics. MPM also suffers from numerical artifacts (e.g. cell-crossing
85 instability, volumetric locking, sporadic occurrence of empty elements in the
86 material domain). Many publications are available in the literature describing the
87 MPM algorithm and the numerical solutions to address its limitations (the reader
88 is directed to [18] for an overview).

89 Two MPM frameworks have been presented to study multi-phase problems [19]:
90 the single-point approach and the multi-point approach. The single-point frame-
91 work represents the multi-phase soil mixture as a unique continuum represented
92 with one set of MPs. During the calculation, all MPs move with the displacement
93 of the solid skeleton represented using Lagrangian formulation and consequently

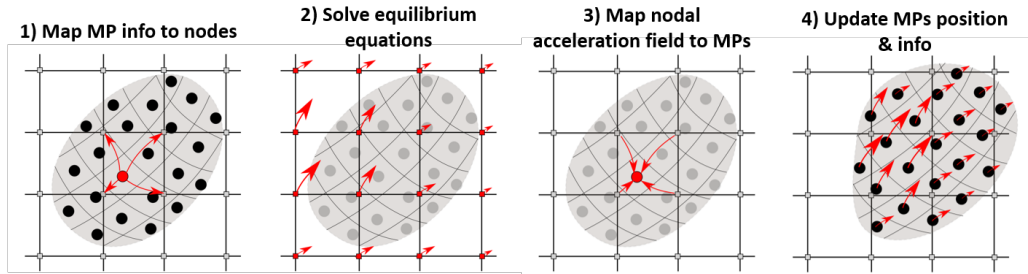


Fig. 1. Diagram of the MPM computational cycle (modified from [22]).

94 the solid mass in the MP remains constant. Fluids are allowed to flow in and
 95 out of the reference porous domain and their motion is described with respect to
 96 the moving solid skeleton. The single-point approach is the general framework
 97 selected to represent soils in unsaturated conditions as discussed herein. The
 98 multi-point framework requires more than one set of MPs to represent different
 99 phases. In particular, the double-point approach for saturated soil is available in
 100 the literature and represents solid skeleton and liquid with two completely separate
 101 sets of MPs [e.g. 20, 21]. Each set of MPs moves accordingly to the displacement
 102 of the corresponding phase and carries the information of one phase only, hence
 103 solid and fluid are represented using Lagrangian formulation. In the double-point
 104 configuration presented in [20, 21], the soil is considered either fully dry or fully
 105 saturated but this same approach can be extended to include hydro-mechanical
 106 interactions resulting from partially saturated conditions. Based on the authors
 107 knowledge, the study of unsaturated soils using a three-point approach (i.e. three
 108 sets of MP for solid, liquid and gas) has never been attempted yet.

109 The aim of following subsections is to provide an overview of the MPM formu-
 110 lations available in the literature for the simulation of unsaturated soils, proposed
 111 by Yerro et al. [23], Yerro [24], Bandara et al. [25], Wang et al. [26, 27], Lei and

112 Solowski [28], Lei et al. [16], Ceccato et al. [29, 30], Martinelli et al. [31]. A
 113 summary of the main features discussed below is provided in Tab. 1.

114 The governing equations of a comprehensive hydro-mechanically coupled formu-
 115 lation for unsaturated soils in a continuum framework should consider the balance
 116 equations of the phases and the interactions among them. Additionally, compat-
 117 ibility equations and constitutive relationships are also required to complete the
 118 equations set. General assumptions considered in all MPM formulations presented
 119 below include: (a) solid grains are incompressible, (b) fluids are weakly compress-
 120 ible, (c) fluid flows are laminar (Darcy's law is valid for gas and liquid phases), (d)
 121 fluid convective terms are negligible, and (e) isothermal conditions.

122 2.1. Momentum balance equations

123 The first MPM formulation for unsaturated soils was proposed by Yerro [24], Yerro
 124 et al. [23]. This is a three-phase single-point formulation and each MP represents
 125 the solid-liquid-gas mixture. The main governing equations posed at the nodes of
 126 the computational mesh are the linear momentum balances of the gas, liquid, and
 127 mixture (Eqs. 1, 2, and 3, respectively). The first two are posed per unit of fluid
 128 volume, while the mixture is posed per unit of total soil volume. This formulation
 129 is fully dynamic and all relative acceleration terms are fully accounted with the
 130 accelerations of each phase being the primary unknowns of the system, i.e. \mathbf{a}_S , \mathbf{a}_L
 131 and \mathbf{a}_G .

$$\rho_G \mathbf{a}_G = \nabla p_G - \mathbf{f}_G^d + \rho_G \mathbf{g} \quad (1)$$

$$\rho_L \mathbf{a}_L = \nabla p_L - \mathbf{f}_L^d + \rho_L \mathbf{g} \quad (2)$$

$$n_S \rho_S \mathbf{a}_S + n_L \rho_L \mathbf{a}_L + n_G \rho_G \mathbf{a}_G = \nabla \cdot \boldsymbol{\sigma} + \rho_m \mathbf{g} \quad (3)$$

Table 1. Summary of most relevant characteristics of currently available MPM formulations for unsaturated soils.

Formulation	Yerro et al. [23]	Yerro [24]	Bandara et al. [25]	Wang et al. [26, 27], Lee et al. [32], Ceccato et al. [30] et al. [32], Ceccato et al. et al. [16] [33, 29]	Martinelli et al. [31]
Number of phases	solid-liquid-gas	solid-liquid-gas	solid-liquid (+ suction)	solid-liquid (+ suction)	solid-liquid (+ suction)
Dynamics	Fully dynamic	Fully dynamic	Relative acceleration of the liquid is neglected	Fully dynamic	Fully dynamic
Main governing equations	momentum of the mixture, momentum of the liquid, momentum of the gas	momentum of the mixture, momentum of the liquid, momentum of the gas	momentum of the mixture, momentum of the liquid	momentum of the mixture, momentum of the liquid	momentum of the solid, momentum of the liquid
Principal unknowns	a_S, a_L, a_G	a_S, a_L, a_G	a_S, w	a_S, a_L	a_S, a_L
Mass exchange	Liquid-Gas	Liquid-Gas	No	No	No
Stress soil	Independent stress variables	Independent stress variables	Bishop effective stress	Bishop effective stress	Bishop effective stress
Incompressible solid grains	x	x	x	x	x
Weakly compressible fluids	x	x	x	x	x
Laminar fluid flow (Darcy's law)	x	x	x	x	x
Isothermal conditions	x	x	x	x	x
Solid-fluid relative acceleration is neglected		x			
Gas pressure is neglected		x			x
Gas density is neglected		x			x
Gradient of degree of saturation is negligible	x				
Gradient of porosity is negligible	x				x
Gradient of liquid density is negligible	x				x

Main hypothesis

132 Where n_S , n_L and n_G are the volumetric concentration ratio of solid, liquid, and
 133 gas; ρ_S , ρ_L , ρ_G are densities of all phases and ρ_m is density of the mixture
 134 ($\rho_m = n_S\rho_S + n_L\rho_L + n_G\rho_G$); \mathbf{f}_G^d and \mathbf{f}_L^d are gas and liquid drag forces; p_G and
 135 p_L are gas and liquid pressures; $\boldsymbol{\sigma}$ is the total stress tensor; and \mathbf{g} is the gravity
 136 vector. The porosity of the solid skeleton becomes $n = n_L + n_G$, and the volumetric
 137 concentration ratio of the fluid phases can be expressed in terms of the degree of
 138 saturation (S_L) and porosity as $n_L = nS_L$ and $n_G = n(1 - S_L)$, respectively for
 139 liquid and gas. Note that fully saturated conditions are the particular case when
 140 the degree of saturation S_L is one.

141 The fluid flow (either liquid or gas) is assumed laminar and the liquid and gas drag
 142 forces (\mathbf{f}_G^d and \mathbf{f}_L^d) are written taking into account that Darcy's law is valid (Eq. 4).
 143 The subscript f denotes fluid phase, that can be either gas G or liquid L .

$$\mathbf{f}_f^d = \frac{n_f \mu_f}{k_f} (\mathbf{v}_f - \mathbf{v}_S) \quad (4)$$

144 Where μ_f and k_f correspond to the viscosity and intrinsic permeability of the
 145 fluids (i.e. $f = G, L$). The intrinsic permeability is defined in terms of hydraulic
 146 conductivity (κ_f) as

$$k_f = \kappa_f \frac{\mu_f}{\rho_f g} \quad (5)$$

147 The three momentum balances are integrated into the domain and discretised at
 148 the nodes of the computational mesh (the detailed expressions can be found in
 149 [24, 23]). At the beginning of each computational cycle, information carried by
 150 the MPs is mapped to the mesh to calculate nodal mass, nodal velocities, internal

151 and external forces, and dragging terms. Nodal gas and liquid accelerations are
 152 obtained first from solving the discretised version of Eqs. 1 and 2 independently.
 153 Finally, the nodal acceleration of the solid skeleton is obtained by solving Eq. 3
 154 and using updated nodal fluid velocities.

155 The rest of the MPM formulations addressing unsaturated conditions assume that
 156 the density and pressure of the gas phase are negligible ($\rho_G = 0$ and $p_G = 0$). These
 157 formulations are essentially two-phase approaches that account for the suction
 158 effect. In consequence, the linear momentum balance of the gas is dropped from
 159 the system of equations and only two momentum balance equations are considered
 160 instead of three. This simplification reduces the computational cost. In the
 161 formulations proposed by Wang et al. [26, 27], Lei and Solowski [28], Lee et al.
 162 [32], Lei et al. [16], Ceccato et al. [29, 30] all dynamic terms, i.e. liquid and solid
 163 inertia, are considered and the main unknowns of the system are the accelerations
 164 of the solid and liquid phases (\mathbf{a}_S and \mathbf{a}_L). The linear momentum of the liquid is
 165 identical to Eq. 2. The linear momentum of the mixture is equivalent to Eq. 3 but
 166 removing the dynamic term corresponding to the gas phase, which yields Eq. 6.

$$n_S \rho_S \mathbf{a}_S + n_L \rho_L \mathbf{a}_L = \nabla \cdot \boldsymbol{\sigma} + \rho_m \mathbf{g} \quad (6)$$

167 Very similarly, the formulation from Martinelli et al. [31] considers all dynamic
 168 terms and the solid and liquid accelerations as primary variables, but the momen-
 169 tum balance of the solid per unit of total volume (Eq. 7) is posed instead of the
 170 momentum balance of the mixture. \mathbf{I} is the identity matrix.

$$n_S \rho_S \mathbf{a}_S = \nabla \cdot (\boldsymbol{\sigma} - n_L p_L \mathbf{I}) + (\rho_m - n_L \rho_L) \mathbf{g} + \mathbf{f}_L^d \quad (7)$$

171 Finally, the formulation by Bandara et al. [25] neglects the relative acceleration of
 172 the liquid with respect to the solid skeleton. Therefore, all phases have the same
 173 acceleration \mathbf{a} (i.e. acceleration of the system, $\mathbf{a} = \mathbf{a}_S = \mathbf{a}_L$). The momentum
 174 balances of the liquid, which is equivalent to the generalized Darcy equation
 175 (Eq. 8), and the mixture (Eq. 9) are the governing equations posed at the nodes of
 176 the computational mesh and the main unknowns are the acceleration of the system
 177 \mathbf{a} and the liquid seepage velocity ($\mathbf{w} = nS_L(\mathbf{v}_L - \mathbf{v}_S)$).

$$\mathbf{w} = -\frac{\kappa_L}{\mu_f}(\nabla p_L + \rho_L - \rho_L \mathbf{g}) \quad (8)$$

$$\rho_m \mathbf{a}_S = \nabla \cdot \boldsymbol{\sigma} + \rho_m \mathbf{g} \quad (9)$$

178 2.2. Mass balance equations

179 Additional governing equations include the mass balances. In the three-phase
 180 MPM formulation [23, 24] the liquid and gas are considered to be a mixture of
 181 water and air, and mass exchange between liquid and gas phases is allowed to
 182 account for "water vapour" in the gas and "dissolved gas" in the liquid. The
 183 process of mass exchange is not included herein to facilitate the reading of the
 184 paper. Assuming incompressible solid grains ($\frac{D^S \rho_S}{Dt} = 0$), the expressions for the
 185 mass balance of the solid and fluids (i.e. liquid and gas) phases are written as
 186 Eqs. 10 and 11, respectively. $\frac{D^S(\bullet)}{Dt}$ is the material derivative with respect to the
 187 solid motion.

$$\frac{D^S n}{Dt} - (1 - n)\nabla \cdot \mathbf{v}_S = 0 \quad (10)$$

$$n \frac{D^S(\rho_f S_f)}{Dt} + \nabla \cdot (\rho_f S_f n(\mathbf{v}_f - \mathbf{v}_S)) + \rho_f S_f \nabla \cdot \mathbf{v}_S = 0 \quad (11)$$

188 The mass balance of the solid (Eq. 10) represents the variation of porosity n
 189 caused by volumetric deformation of the solid skeleton. In those problems where
 190 the spatial variations of fluids mass in the soil are small ($\nabla(\rho_f S_f n) \approx 0$), Yerro
 191 et al. [23] simplify the term with the gradient of the advective fluxes of the fluid
 192 phases (i.e. $\nabla \cdot (\rho_f S_f n(\mathbf{v}_f - \mathbf{v}_S))$) yielding to Eq. 12. The effect of neglecting
 193 part of the advective flow gradients is further discussed in Sec. 2.5.

$$n \frac{D^S(\rho_f S_f)}{Dt} + \rho_f S_f n \nabla \cdot (\mathbf{v}_f - \mathbf{v}_S) + \rho_f S_f \nabla \cdot \mathbf{v}_S = 0 \quad (12)$$

194 The solid material derivatives from Eqs. 11 and 12 are evaluated using the chain
 195 rule and considering the fluid pressures (p_G and p_L) as state variables (Eq. 13).
 196 The fluid mass balances are solved in the MPM computational cycle in terms of
 197 the variation of liquid and gas pressure ($\frac{D^S p_L}{Dt}$ and $\frac{D^S p_G}{Dt}$).

$$\frac{D^S(\rho_f S_f)}{Dt} = \left(S_f \frac{\partial \rho_f}{\partial p_L} + \rho_f \frac{\partial S_f}{\partial p_L} \right) \frac{D^S p_L}{Dt} + \left(S_f \frac{\partial \rho_f}{\partial p_G} + \rho_f \frac{\partial S_f}{\partial p_G} \right) \frac{D^S p_G}{Dt} \quad (13)$$

198 The two-phase MPM formulations neglect the gas density and only the mass
 199 balances of the solid and liquid phases are required. All approaches consider the
 200 mass balance of the solid as presented in Eq. 10, but propose slightly different
 201 versions of the mass balance of the liquid. Ceccato et al. [30] consider the liquid
 202 mass balance with no simplification regarding the the gradient of the advective
 203 fluxes, leading to Eq. 14. Wang et al. [26, 27], Ceccato et al. [33, 29], Lee et al.
 204 [32], Martinelli et al. [31] neglect the gradients of the liquid mass ($\nabla(\rho_L S_L n) \approx 0$)
 205 and use Eq. 15 instead. Bandara et al. [25] rewrites Eq. 14 in terms of \mathbf{w} .

$$n \frac{D^S \rho_L S_L}{Dt} = -\nabla \cdot (\rho_L S_L n (\mathbf{v}_L - \mathbf{v}_S)) - \rho_L S_L \nabla \cdot \mathbf{v}_S \quad (14)$$

$$n \frac{D^S \rho_L S_L}{Dt} = -\rho_L S_L n \nabla \cdot (\mathbf{v}_L - \mathbf{v}_S) - \rho_L S_L \nabla \cdot \mathbf{v}_S \quad (15)$$

206 The solid material derivatives from Eqs. 14 and 15 are evaluated assuming a
 207 weakly-compressible liquid (see Sec. 2.4.2), and the variation of the liquid pressure
 208 ($\frac{D^S p_L}{Dt}$) is derived accordingly.

$$\frac{D^S \rho_L S_L}{Dt} = \rho_L \frac{D^S S_L}{Dt} - \frac{\rho_L}{K_L} \frac{D^S p_L}{Dt} \quad (16)$$

209 Finally, the formulation proposed by Lei and Solowski [28], Lei et al. [16] extends
 210 the previous works to account for internal erosion of the solid skeleton, based
 211 on the work from Yerro et al. [34] for saturated conditions. The mass exchange
 212 between the solid and liquid phases is allowed according to an erosion law that
 213 controls the rate of eroded mass. This implies that the mass balances of solid and
 214 liquid phases (Eq. 10 and Eq. 14) are extended with an additional term to account
 215 for the lost or gained mass, respectively. In this formulation, the gradient of the
 216 liquid density is neglected, but gradients of porosity and degree of saturation are
 217 accounted. In addition to the solid and liquid mass balances, the mass balance of
 218 the eroded grains (i.e. liquidized spices) is also required in the system.

219 2.3. Compatibility equations

220 All formulations consider the same compatibility equations to obtain the infinites-
 221 imal strain rate tensor of each phase $\boldsymbol{\varepsilon}_{ph}$ (i.e. $ph = S, L, G$) at the MPs. These
 222 can be written in terms of the corresponding phase velocities as in Eq. 17. In the

223 MPM computational cycle, the compatibility equation is evaluated at the MPs and
 224 updated nodal velocities are considered to calculate the corresponding gradients.

$$\frac{D^S \boldsymbol{\varepsilon}_{ph}}{Dt} = \frac{1}{2} \left[\nabla \cdot \mathbf{v}_{ph} + (\nabla \cdot \mathbf{v}_{ph})^T \right] \quad (17)$$

225 2.4. Constitutive laws

226 Different constitutive laws are required to complete the set of governing equations.
 227 These are evaluated at the MPs, generally at the end of the computational cycle.

228 2.4.1. Stress-strain relationship for solid skeleton

229 The relationship between the strains of the solid skeleton and the stress is defined
 230 by means of a constitutive equation. Various hydro-mechanical stress frameworks
 231 are available in the literature to describe the behaviour of unsaturated soils [35].
 232 They are inherited from the Bishop single effective stress approach [36] or from the
 233 independent stress variable approach [37]. The Bishop effective stress approach is
 234 the most commonly implemented in the MPM formulations [25, 27, 16, 32, 30, 31].
 235 The general form of the constitutive equation is presented incrementally as Eq. 18,
 236 where the effective stress essentially controls the stress state of unsaturated soil
 237 and is defined as Eq. 19. \mathbf{D} is the tangent matrix, $\boldsymbol{\varepsilon}$ is the strain vector of the
 238 solid skeleton, $\boldsymbol{\sigma}_{net}$ is the net stress ($\boldsymbol{\sigma}_{net} = \boldsymbol{\sigma} - p_G \mathbf{I}$), s is the matric suction
 239 ($s = p_G - p_L$), and χ is an effective stress parameter generally assumed equal to
 240 S_L . Note that most of the works assume $p_G = 0$.

$$d\boldsymbol{\sigma}' = \mathbf{D} \cdot d\boldsymbol{\varepsilon} \quad (18)$$

$$\boldsymbol{\sigma}' = \boldsymbol{\sigma}_{net} - \chi s \mathbf{I} \quad (19)$$

241 The independent stress variable approach is used in [23], where the net stress is
 242 defined more generally as Eq. 20. In this context, the stress state is described by
 243 two state variables (i.e. σ_{net} and s) and a double constitutive matrix is required
 244 (Eq. 21), where \mathbf{h} is a constitutive vector controlling the influence of suction.

$$\sigma_{net} = \sigma - \max\{p_G, p_L\}\mathbf{I} \quad (20)$$

$$d\sigma_{net} = \mathbf{D} \cdot d\boldsymbol{\varepsilon} + \mathbf{h}ds \quad (21)$$

245 Very limited number of constitutive models have been considered in both stress
 246 frameworks. Linear elasticity has been used for validation purposes and the stan-
 247 dard elastic-perfectly plastic Mohr Coulomb and the Mohr Coulomb with suction
 248 hardening [38] have also been used for the study of different failure problems
 249 (Sec. 4 includes a review of the applications). Finally, the Jaumann stress rate is
 250 adopted in case of large deformation to achieve objectivity (i.e. frame indifferent
 251 formulation) of the Cauchy stress rate matrix [25, 31].

252 2.4.2. Weakly-compressible fluids

253 All the currently available MPM formulations for unsaturated soils consider linear
 254 elastic weakly-compressible fluids (Eq. 22), where K_f is the bulk modulus of the
 255 reference fluid.

$$-\frac{1}{\rho_f} \frac{\partial \rho_f}{\partial t} = \frac{1}{K_f} \frac{\partial p_f}{\partial t} \quad (22)$$

256 2.4.3. Soil water retention curve

257 The soil water retention curve (SWRC) is an essential relationship required to
 258 model the behaviour of unsaturated soils that correlates suction ($s = p_G - p_L$) with

259 degree of saturation (S_L). For the purpose of simulating realistic soil behaviour,
 260 the Van Genuchten SWRC [39] (Eq. 23) is preferred among the authors, where
 261 p_0 and λ are fitting parameters. Alternatively, a linear SWRC (Eq. 24) is also
 262 employed for validation purposes, where a_v is a constant parameter, S_{min} is the
 263 residual degree of saturation, and S_{max} is the maximum degree of saturation.

$$S_L = S_{min} + (S_{max} - S_{min}) \left[1 + \left(\frac{p_G - p_L}{p_0} \right)^{\frac{1}{1-\lambda}} \right]^{-\lambda} \quad (23)$$

$$S_L = 1 - a_v(p_G - p_L) \quad (24)$$

264 2.4.4. Liquid relative hydraulic conductivity

265 It is well known that unsaturated soils are less permeable than fully saturated
 266 soils. The liquid relative hydraulic conductivity (i.e. the ratio between the actual
 267 liquid hydraulic conductivity and the saturated hydraulic conductivity, κ_L/κ_{sat})
 268 is a function of the degree of saturation. This dependence is also included in
 269 the MPM formulations with the introduction of the hydraulic conductivity curve
 270 (HCC). The most common HCC are the functions proposed by Hillel [40] (Eq. 25),
 271 and Mualem [41] (Eq. 26), where r and λ are fitting parameters.

$$\frac{\kappa_L}{\kappa_{sat}} = S_L^r \quad (25)$$

$$\frac{\kappa_L}{\kappa_{sat}} = \sqrt{S_L} \left[1 - \left(1 - S_L^{\frac{1}{\lambda}} \right)^\lambda \right]^2 \quad (26)$$

272 2.5. Discussion

273 This section discusses the effect of the gradient of advective fluxes in the fluid
 274 mass balance equations. Eq. 11 can be written as Eq. 27, where $f = L, G$, by
 275 expanding the gradient of the advective flux term.

$$n \frac{D^S(\rho_f S_f)}{Dt} + \nabla(\rho_f S_f n) \cdot (\mathbf{v}_f - \mathbf{v}_S) + \rho_f S_f n \nabla \cdot \mathbf{v}_f + \rho_f S_f (1 - n) \nabla \cdot \mathbf{v}_S = 0 \quad (27)$$

276 The first addend is related to the liquid mass change as a result of pressure and
 277 degree of saturation variation in time (Eq. 16). All other terms explain the variation
 278 of liquid mass as a result of the liquid inflow/outflow gradients resulting from
 279 different mechanisms. In particular, the second term in Eq. 27 describes the
 280 variation of the liquid mass induced by a flow triggered by the fluid mass gradient;
 281 the third term describes the variation of liquid mass due to the divergence of the
 282 fluid velocity (equivalent to a volumetric deformation of the fluid); and the fourth
 283 term describes the variation of liquid mass due to the divergence of the solid
 284 velocity (equivalent to a volumetric deformation of the solid skeleton).

285 The mass balance equation is solved at the MP level, where the computation of the
 286 last two terms of Eq. 27 is straightforward because nodal velocities are already
 287 available during the traditional computational cycle. In contrast, the evaluation of
 288 the fluid mass gradient requires the additional step of mapping the quantity $\rho_f S_f n$
 289 to the nodes and then calculate the gradient at the MP [24]. This is relatively
 290 simple and does not increase significantly the computational cost. The gradient
 291 of fluid mass can be calculated as Eq. 28. While the spatial gradient of liquid
 292 density ($\nabla \rho_f$) and porosity (∇n) can be assumed negligible in most cases, the
 293 gradient of the degree of saturation depends on the pressure gradient and the
 294 SWRC. The importance of the last term of Eq. 28 increases with pressure gradient
 295 and $\partial S_f / \partial p_f$.

$$\nabla(\rho_f S_f n) = S_f n \nabla \rho_f + S_f \rho_f \nabla n + n \rho_f \nabla S_f \quad (28)$$

296 To visualize the effect of this term, the evolution of suction in a 1m soil column
297 during an infiltration test is considered. For $t = 0$, an initial suction s_0 is applied
298 along the column. For $t > 0$, zero suction is imposed at the top boundary while the
299 bottom is impervious. Gravity is neglected. The soil permeability is constant and
300 the SWRC is linear (Eq. 24). Under these assumptions, an analytical expression
301 that describes the evolution of the normalized suction along the column with time
302 can be derived from the mass balance equation of the liquid following [42]

303 Numerical simulations are performed with the two-phase formulations for unsat-
304 urated soils with the material parameters listed in Tab. 2.

305 Figure 2 compares the results obtained with the complete mass balance equation
306 (Eq. 14, circles) and with the simplified mass balance equation (Eq. 15, cross
307 symbol) considering different values of initial suction s_0 and $\partial S_L / \partial p_L = a_v$. For
308 $s_0 = 500kPa$ and $a_v = 1 \cdot 10^{-3}$ (Fig. 2a) the simplified mass balance equation leads
309 to a delay in the evolution of suction, while the complete mass balance equation
310 gives results in very good agreement with the analytical solution. Decreasing the
311 slope of the SWRC to $a_v = 1 \cdot 10^{-5}$ (Fig. 2b) or reducing the initial suction to
312 $s_0 = 5kPa$, i.e. the pressure gradient, (Fig. 2c) the two approaches give very
313 similar results. Indeed, in these cases, the contribute of the last term of Eq. 28 is
314 very small and neglecting the gradient of fluid mass is an acceptable simplification.
315 For the range of pressure gradients and SWRC typical of many civil engineering
316 applications the error introduced using Eq. 15 is acceptably small.

317 Similar results are obtained with the three-phase formulation as discussed in [43].

Solid density [kg/m^3]	ρ_S	2700
Liquid density [kg/m^3]	ρ_L	1000
Porosity [-]	n	0.4
Liquid bulk modulus [kPa]	K_L	80000
Liquid dynamic viscosity [$kPa \cdot s$]	μ_L	$1 \cdot 10^{-6}$
Intrinsic permeability liquid [m^2/s]	κ_L	$1 \cdot 10^{-11}$

Table 2. Material parameters for 1D infiltration example with applied pressure.

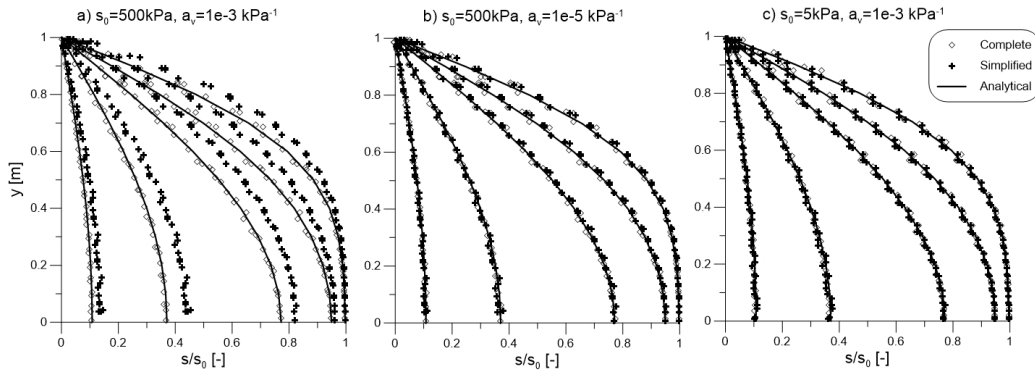


Fig. 2. Evolution of normalized suction with depth. Comparison between numerical and analytical formulations.

318 3. Implementation aspects

319 This section presents an overview of different implementation aspects concern-
 320 ing unsaturated MPM formulations such as the stability of the solution and the
 321 application of boundary conditions.

322 3.1. Time integration scheme and critical time step

323 All MPM formulations presented for the study of unsaturated soils are based on
 324 numerical schemes explicitly integrated in time. Commonly, the Euler-Cromer

325 algorithm is used. Explicit schemes are conditionally stable depending on a
 326 maximum or critical time step Δt_c , that depends on the material properties and
 327 characteristic size of the computational grid. According to the authors' knowl-
 328 edge, no studies have been performed concerning the effects of partially saturated
 329 conditions on Δt_c .

330 3.1.1. Review of stability criteria

331 In two-phase problems we can define two time scales of interest, namely that of
 332 excess pore pressure dissipation related to consolidation and infiltration and that
 333 of compression wave propagation within the solid-fluid mixture. In mathematics,
 334 consolidation (infiltration) is seen as parabolic behaviour, while wave propagation
 335 is hyperbolic behaviour. The associated critical time steps are expressed by Eq.
 336 29 and Eq. 30 respectively, where L_i is a characteristic length, i.e. the size of an
 337 element.

$$\Delta t_c = \frac{L_i^2}{2c} \quad (29)$$

$$\Delta t_c = \frac{L_i}{v_c} \quad (30)$$

338 The coefficient c can be written as Eq. 31 where E'_c is the effective oedometric
 339 modulus. Note that for $S_L = 1$ and $\frac{dS_L}{ds} = 0$ it coincides with the consolidation
 340 coefficient.

$$c = \frac{k_L}{\mu_L \left(\frac{1}{E'_c} + \frac{S_L n}{K_L} + n \frac{dS_L}{ds} \right)} \quad (31)$$

341 v_c in Eq. 30 represents the one-dimensional compressing wave speed propagation

342 and can be estimated with Eq. 32, where E_c and ρ are the oedometric modulus
 343 and the density of the considered material.

$$v_c = \sqrt{\frac{E_c}{\rho}} \quad (32)$$

344 Eq. 32 can be particularized for undrained conditions using $E_c = E'_c + n/K_L$ and
 345 $\rho = \rho_{sat}$ or for dry conditions using $E_c = E'_c$ and $\rho = \rho_{dry}$.

346 Yerro [24] and Mieremet [44] mathematically studied the stability of the two-
 347 phase MPM formulation for saturated conditions proposed in [45]. In the work of
 348 Yerro [24], the eigenvalue problem is solved to study the stability of four different
 349 equations: (a) the liquid momentum balance $\Delta t_{c,liquid}$, (b) the mixture momentum
 350 balance $\Delta t_{c,mixture}$, (c) the coupled hydro-mechanical system ($\Delta t_{c,coupled}$) (i.e.
 351 liquid and mixture momentum balances), and (d) the momentum balance of the
 352 mixture in undrained conditions ($\Delta t_{c,undrained}$). The critical time step criterion for
 353 the undrained analysis is identical to Eq. 30, while the other three criteria share
 354 the following expression

$$\Delta t_{c,Yerro} = \frac{2}{\omega} \left[-\xi + \sqrt{\xi^2 + 1} \right] \quad (33)$$

355 where ω and ξ vary depending on the set of equations considered, and provided in
 356 Tab.3 (saturated column).

357 Mieremet [44] considered the stability of the coupled hydro-mechanical system
 358 concerning liquid and mixture momentum balances. Rigorous mathematical sta-
 359 bility analysis were conducted by means of the matrix method [46], and an esti-
 360 mation of Δt_c is proposed as

$$\Delta t_{c, \text{Mieremet}} = \frac{-2a + \sqrt{4a^2 + 8(b + \sqrt{b^2 - 4d})}}{b + \sqrt{b^2 - 4d}} \quad (34)$$

361 where terms a , b and d are reported in Tab. 4 (saturated column). The performance
 362 of the proposed criterion is verified with 1D FEM-based simulations.

363 3.1.2. New stability criteria for unsaturated conditions

364 Assuming that saturation can be understood as a particular case of unsaturated
 365 conditions, one could expect that the Δt_c for unsaturated formulations is an ex-
 366 tension of those obtained for saturated conditions. In this work, the stability of
 367 the MPM formulation proposed by Ceccato et al. [30] in unsaturated conditions is
 368 studied considering the same approach followed by Yerro [24]. The stability of (a)
 369 the momentum balance of the liquid, (b) the momentum balance of the mixture,
 370 and (c) the coupled system is derived. The main passages to obtain the stability
 371 criteria are presented in detail in the Appendix (Sec. 7). The obtained expressions
 372 for the critical time step follow the same structure as Eq. 33. The new terms ω
 373 and ξ are reported in Tab. 3 (unsaturated column).

374 Given the consistent complexities in applying the same rigorous procedure by
 375 Mieremet [44] to the formulation by Ceccato et al. [30], a simplified "adaptation"
 376 of Eq. 34 to unsaturated conditions is proposed. The terms a , b and d are modified
 377 in a similar manner to ω and ξ in saturated vs. unsaturated conditions (3). In
 378 the term a , S_L^2 is added and ρ_{sat} is replaced by ρ_m . In terms b and d , the liquid
 379 bulk modulus K_L is replaced by $\frac{S_L}{\left(\frac{dS_L}{ds}\right) + \frac{S_L}{K_L}}$, based on ω_C and ω_M unsaturated
 380 expressions. Note that, consistently for all the criteria, the saturated case ($S_L = 1$)
 381 is a particular case of the unsaturated expressions.

	Saturated	Unsaturated
Coupled	$\omega_C = \frac{1}{L_i} \sqrt{\frac{K_L}{\rho_L} + \frac{E_c}{\rho_S(1-n)} + \frac{(1-n)K_L}{n\rho_S}}$ $\xi_C = \frac{n\mu_L}{2k_L\omega_C} \left[\frac{n}{(1-n)\rho_S} + \frac{1}{\rho_L} \right]$	$\omega_C = \frac{1}{L_i} \sqrt{\left[\frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)} \right] \left[\frac{1}{\rho_L} + \frac{1}{\rho_S} \left(\frac{1}{n} - S_L \right) \right]} + \frac{E_c}{\rho_S(1-n)}$ $\xi_C = \frac{S_L n \mu_L}{2k_L \omega_C} \left[\frac{S_L n}{(1-n)\rho_S} + \frac{1}{\rho_L} \right]$
Liquid	$\omega_L = \frac{1}{L_i} \sqrt{\frac{K_L}{\rho_L}}$ $\xi_L = \frac{n\mu_L}{2k_L\omega_L} \frac{1}{\rho_L}$	$\omega_L = \frac{1}{L_i} \sqrt{\frac{S_L}{\rho_L \left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)}}$ $\xi_L = \frac{S_L n \mu_L}{2k_L \omega_L} \frac{1}{\rho_L}$
Mixture	$\omega_M = \frac{1}{L_i} \sqrt{\frac{E_c}{(1-n)\rho_S} + \frac{(1-n)K_L}{n\rho_S}}$ $\xi_M = \frac{n^2\mu_L}{2(1-n)\rho_S k_L \omega_M}$	$\omega_M = \frac{1}{L_i} \sqrt{\frac{1}{\rho_S} \left[\frac{E_c}{(1-n)} + \left(-S_L + \frac{1}{n} \right) \frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)} \right]}$ $\xi_M = \frac{S_L^2 n^2 \mu_L}{2(1-n)\rho_S k_L \omega_M}$

Table 3. Parameters ω and ξ used in Δt_c expression by Yerro [24]. Left column: original version for saturated conditions. Right column: new development for unsaturated conditions.

	Saturated	Unsaturated
Mieremet	$a = \frac{n\rho_{sat}\mu}{(1-n)\rho_S\rho_L k_L}$ $b = \frac{4(n\rho_{sat}K_L + (1-2n)\rho_L K_L + n\rho_L E_c)}{n(1-n)\rho_S\rho_L L_i^2}$ $d = \frac{16E_c K_L}{(1-n)\rho_S\rho_L L_i^4}$	Mieremet _{adapted} $a = \frac{S_L^2 n \rho_m \mu}{(1-n)\rho_S \rho_L k_L}$ $b = \frac{4 \left(n \rho_m \frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)} + (1-2n)\rho_L \frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)} + n\rho_L E_c \right)}{n(1-n)\rho_S\rho_L L_i^2}$ $d = \frac{16E_c \frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_L} \right)}}{(1-n)\rho_S\rho_L L_i^4}$

Table 4. Parameters a , b and d used in Δt_c expression by Mieremet [44]. Left column: original version for saturated conditions. Right column: adapted criterion for unsaturated conditions.

382 3.1.3. Numerical analysis and discussion

383 In this section, the performances of the unsaturated criteria are evaluated. A trial
384 and error procedure is used to explore Δt_c in unsaturated conditions. A series of
385 MPM-based simulations is performed with Anura3D [47] using the unsaturated
386 formulation in [30]. The objective is to determine Δt_c for different degrees of
387 saturation. The model is a soil column, 1m high and 0.05m wide. At the top of the
388 column, an external load of 100kPa is applied, see Fig. 3 (a). An initial suction
389 is imposed to the entire column and the same suction is applied at the top of the
390 column during the rest of the calculation to ensure that the degree of saturation
391 remains as constant as possible in the model. The material is linear elastic and the
392 intrinsic permeability ($k_L = 2 \cdot 10^{-10} m^2$, equivalent to a hydraulic conductivity
393 of $\kappa \approx 2 \cdot 10^{-3} m/s$) is assumed constant. The mesh is made by linear triangular
394 elements of 0.05m side length. The bottom boundary is fully fixed and impervious.
395 The lateral boundaries only allow vertical movement and are impervious.

396 The same analysis is repeated for three soil types (sand, silt, and clay) to evaluate
397 the effect of the SWRCs on the numerical stability. The Van Genuchten model
398 is used to describe three reference SWRC from [48] (Fig. 3 (b)). The SWRC
399 parameters are provided in Tab. 5. The simulations are performed for each soil
400 type ensuring a large range of suction values, covering degrees of saturation from
401 saturated ($S_L = 1.0$) to approximately dry conditions ($S_L = 10^{-2}$).

402 Moreover, to evaluate the critical time step in diverse conditions we consider
403 different sets of analysis varying the porosity ($n = 0.4$ and $n = 0.6$) and the Young
404 modulus ($E_1 = 10^4 kPa$ and $E_2 = 5 \cdot 10^4 kPa$). The Poisson ratio is assumed
405 constant for all the simulations, $\nu = 0.2$.

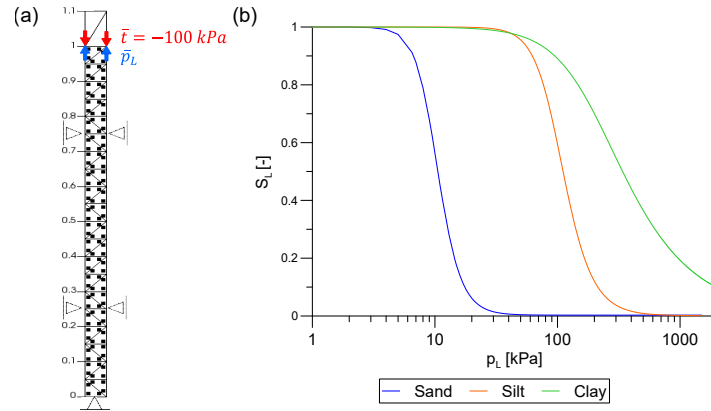


Fig. 3. Critical time step in unsaturated conditions obtained from numerical simulations: (a) geometry of the problem, (b) SWRCs tested.

406 The time step of each calculation is manually increased until the simulation doesn't
 407 converge. This iterative process is repeated for each combination of material pa-
 408 rameters. The maximum value that ensures the solution convergence is considered
 409 the critical time step.

410 The critical time steps obtained from the numerical tests performed in sand, silt,
 411 and clay are presented in Fig. 4, Fig. 5 and Fig. 6, respectively. In all figures, the
 412 numerical results plotted together with the stability criteria proposed in this work
 413 for unsaturated conditions (Tab. 3 and 4). In addition, the expressions from the
 414 literature, i.e. consolidation and infiltration (Eq. 29), and CFL in undrained and
 415 dry conditions (Eq. 30) are also presented for reference. All the expressions are
 416 evaluated considering $L_i = 0.035m$ which corresponds to the minimum altitude
 417 of the element.

418 The numerical results show a general increase of Δt_c in the unsaturated regime.

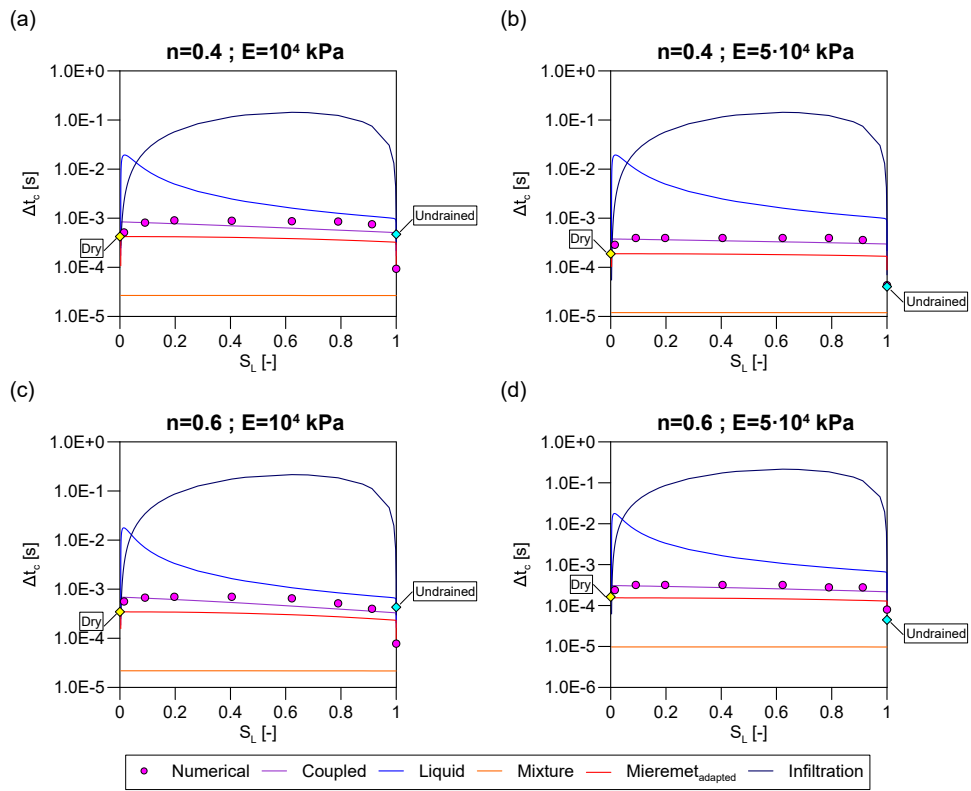


Fig. 4. Δt_c results from MPM simulations compared with mathematical expressions and literature criteria. Sand SWRC.

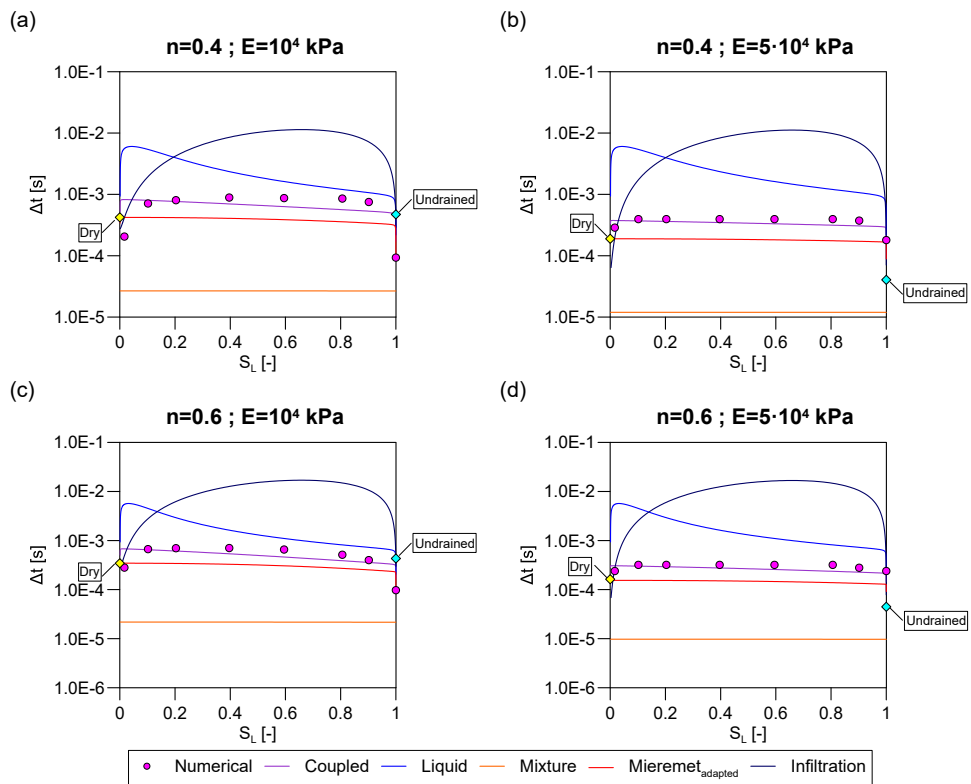


Fig. 5. Δt_c results from MPM simulations compared with mathematical expressions and literature criteria. Silt SWRC.

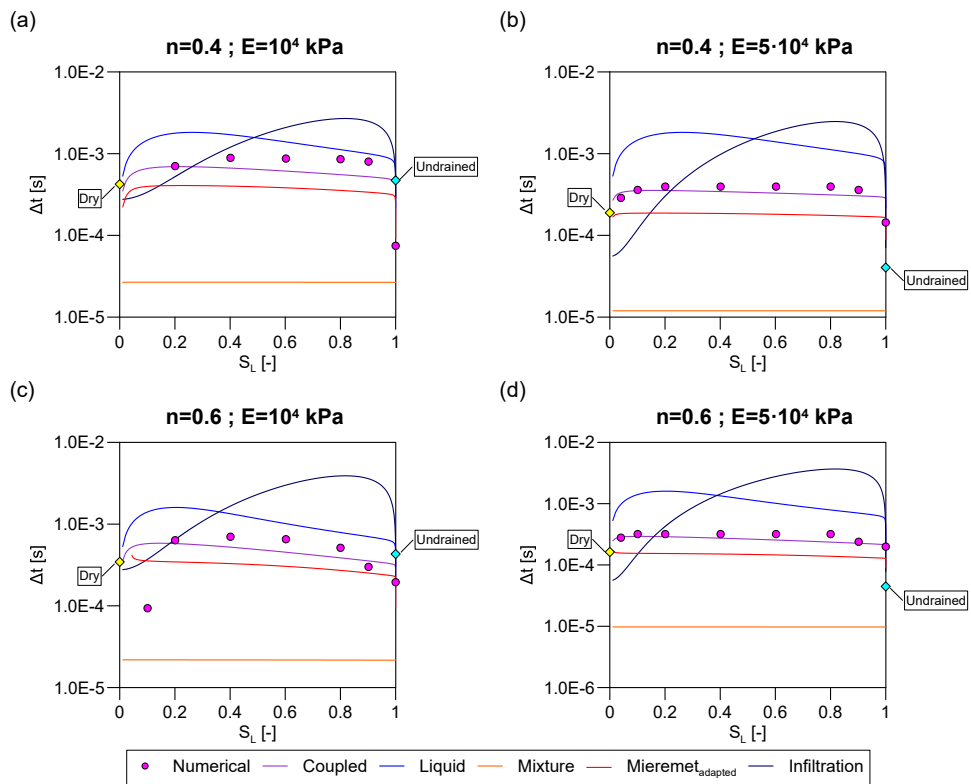


Fig. 6. Δt_c results from MPM simulations compared with mathematical expressions and literature criteria. Clay SWRC.

	Sand	Silt	Clay
S_{min} [-]	0.003	0.001	0.001
S_{max} [-]	1.0	1.0	1.0
p_{ref} [kPa]	9.9	100.0	196.0
λ [-]	0.8	0.75	0.5

Table 5. Van Genuchten parameters for the three SWRCs used in the critical time step numerical analysis.

419 The coupled criterion $\Delta t_{c,Coupled}$ appears as the one better fitting the results,
420 however, it overestimates Δt_c near the extremes (dry and saturated conditions).
421 $\Delta t_{c,Mieremetadapted}$ is the second closest criterion, and it trends to be more con-
422 servative than $\Delta t_{c,Coupled}$. It is important to note that near the dry conditions
423 ($S_L \approx 10^{-2}$) a sharp decrease of Δt_c is consistently observed. This decrease is
424 highlighted by $\Delta t_{c,Coupled}$, $\Delta t_{c,Liquid}$ and $\Delta t_{c,Mieremetadapted}$ criteria. It should be
425 noted that the use of $\Delta t_{c,Mieremetadapted}$ is limited by negative values under the
426 square root present at the numerator of this expression (Eq. 34). This situation
427 seems to occur when suction values are extremely high (i.e. S_L is very low).
428 This behaviour occurs for silt SWRC when $s > 1800kPa$ and for clay SWRC
429 when $s > 4500kPa$ (although it is difficult to appreciate in the figures because
430 of the logarithmic scale). These values of suction are hardly encountered in real
431 geotechnical scenarios.

432 To better quantify the increase of Δt_c in unsaturated conditions and to visualize the
433 effect of each material parameter on it, in Fig. 7(a) the numerical Δt_c values are
434 normalized with respect to $\Delta t_{c,Mieremet}$ (criterion for saturated conditions, Eq. 34).
435 We can observe that smaller porosity values result in bigger Δt_c , as highlighted

436 by circle and diamond symbols. On the other side, the higher Young modulus
437 results in lower Δt_c values, visible with square and cross symbols. Furthermore,
438 the higher Young modulus gives a more flat trend in the interval between $S_L =$
439 0.2 and $S_L = 0.8$. It is interesting to observe the comparable (in some cases
440 identical) increase in Δt_c among the different materials for the same combination
441 of porosity and Young modulus. The major differences occur when approaching
442 dry conditions.

443 In addition, in Fig.7(b), the outcomes using the same model scaled up 10 times, $L_i =$
444 $0.5m$ are introduced to emphasize the mesh size effect on Δt_c . This comparison is
445 related to the silt material. It is clear that Δt_c depends on the mesh size. Results
446 with the coarser mesh have a different trend of critical time step. The sharp
447 reduction of Δt_c near dry conditions is no more present, on the contrary, in this
448 normalization, Δt_c is continuously increasing and the two porosity values tested
449 results in very similar results. A more extensive investigation in this regard is
450 further necessary to understand the different behaviours.

451 These preliminary results can encourage future studies and provide an indicative
452 support to stimulate researchers using MPM unsaturated formulations in testing
453 higher values of Δt_c in situations where significant variations of S_L or extreme
454 values in the unsaturated regime, i.e. full saturation or dry conditions, are not
455 encountered. Further studies need to be performed to investigate the effect of
456 other parameters such as permeability on the stability of the solution.

457 3.2. *Boundary conditions*

458 Given the multi-phase nature of the formulation, the boundary conditions (BCs)
459 need to be applied on each phase separately (solid, liquid, and eventually gas).

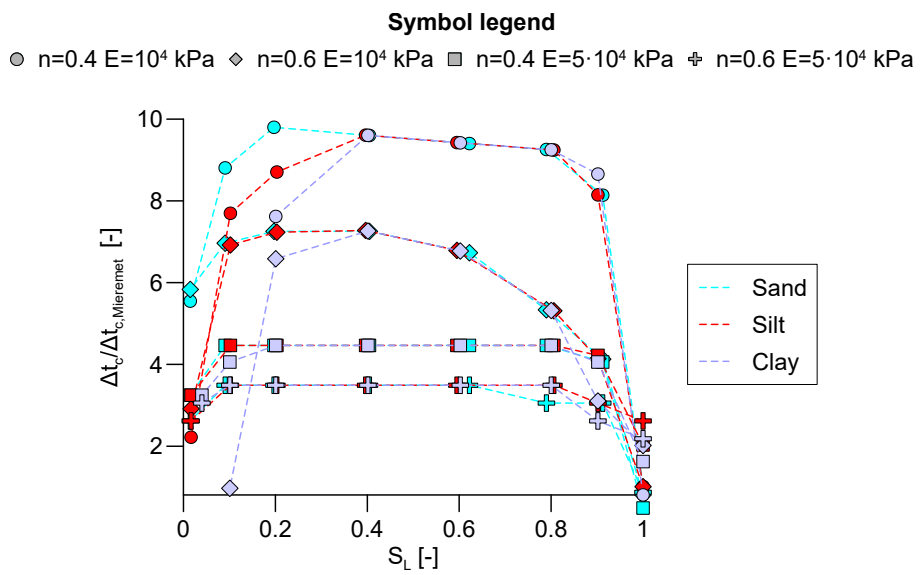


Fig. 7. Comprehensive numerical results normalized with respect to $\Delta t_{c,Mieremet}$, as function of saturation degree (a) ($L_i = 0.05m$). Comparison of normalized Δt_c for different mesh sizes using silt SWRC (b).

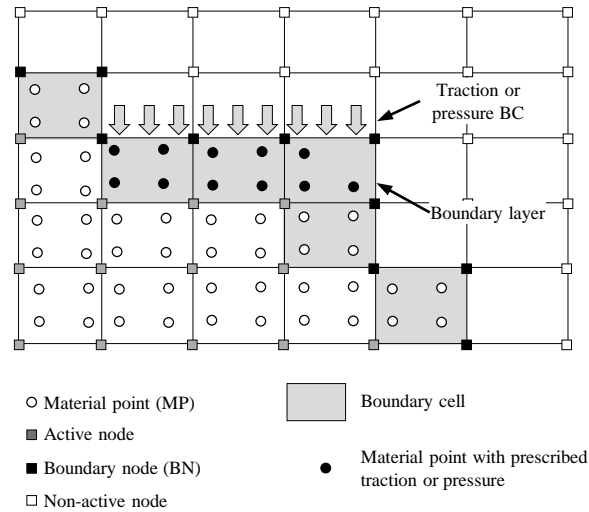


Fig. 8. BC applied with the *boundary layer* approach.

460 Two kinds of BCs can be distinguished – essential and natural. Essential BCs
 461 are imposed directly on the solution, and the degrees of freedom are eliminated
 462 from the system of equations. Fixities, prescribed accelerations, velocities, or
 463 displacements are typical examples of essential BCs. Natural BCs are imposed
 464 on a secondary variable, such as stresses or pressures, and they are included in
 465 the weak form of the governing equations. In classical FEM, the application of
 466 BCs is relatively simple as these can be specified directly on the boundary nodes,
 467 which coincide with the boundary of the continuum body, and are well defined
 468 throughout the computation. However, the computational mesh in MPM does not
 469 necessarily align with the boundary of the material making the application of the
 470 BCs more challenging.

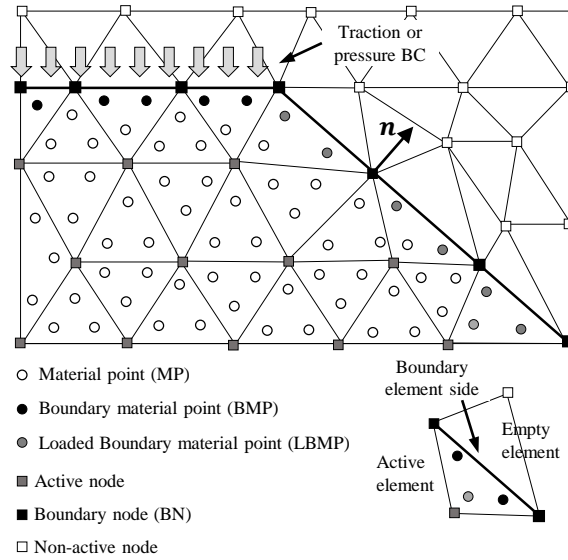


Fig. 9. BC applied at the boundary MP or at the boundary nodes.

471 Bandara et al. [25] apply the concept of *boundary layer* in which the applied
 472 traction or pressure is prescribed to all the MP inside the boundary cells. Boundary
 473 cells are those adjacent to empty cells, and they are determined at each time step
 474 (Fig. 8). The disadvantage of this approach is that traction (or pressure) is
 475 distributed over a thickness that has the size of the cell. In the Anura3D software
 476 [47] used by [23, 30, 32, 31] loads can be applied on either the boundary element
 477 side or the boundary material points (BMPs, Fig. 9) [49]. The first option
 478 is applicable only when the boundary of the body remains aligned with loaded
 479 element boundaries throughout the computation. The nodal traction force is
 480 integrated like in FEM applying Gauss quadrature and then used in the momentum
 481 balance equation. The applied load is thus integrated accurately, and the traction
 482 nodal force is non-zero only for the nodes belonging to the loaded surface. The

483 second option consists of storing the load on specific MPs (loaded BMP, LBMP)
 484 that can move through the mesh. This is mapped from LBMPs to all nodes of
 485 the element where it is located by means of the shape functions. In this way, the
 486 surface force is distributed across the layer of elements that borders the boundary,
 487 but the effect is less severe compared to the boundary layer approach. The LBMPs
 488 may move a long distance and become non-boundary material points, which may
 489 lead to unrealistic results.

490 An alternative definition of pressure BC is the *hydraulic head* BC, which is very
 491 useful to simulate, for example, water levels in reservoirs [30, 50]. A total head
 492 value \widehat{H} is prescribed on the boundary, which is related to the applied pressure \widehat{p}_L
 493 by means of Bernoulli's equation (Eq. 35) where h_g is the geometric head

$$\widehat{H} = h_g - \frac{\widehat{p}_L}{\rho_L g} \quad (35)$$

494 Essential BCs on the liquid phase include *impermeable boundaries* and *infiltration*
 495 *BC*. While the first is trivial, the second is more complex because the infiltration
 496 capacity of a soil depends on its hydraulic conductivity, and it can be solved
 497 in different ways. Bandara et al. [25] impose the infiltration velocity \widehat{w} at the
 498 surface boundary nodes (BNs), i.e. nodes of the cell side which bound active and
 499 non-active nodes (Fig. 8). The infiltration velocity is applied only if the MPs
 500 located adjacent to the BNs are unsaturated. There is no control on the maximum
 501 infiltration velocity, i.e. the user should not apply a value of \widehat{w} larger than the
 502 maximum soil infiltration capacity \widehat{w}_{max} if ponding is not allowed. The case
 503 $\widehat{w} > \widehat{w}_{max}$ will lead to pore pressures higher than zero at the boundary, meaning
 504 that a layer of free water is present above the free surface.

505 Martinelli et al. [31] and Ceccato et al. [30] apply a predictor-corrector scheme:
 506 liquid and solid velocities (\tilde{v}_L, \tilde{v}_S) are predicted assuming zero pressure at the
 507 infiltration boundary and then (eventually) corrected to ensure the prescribed
 508 infiltration rate $\widehat{w} = n_L(v_L - v_S)$. If the net infiltration discharge q_{net} (Eq. 36) is
 509 positive, ponding conditions occur, and if fluid accumulation above the boundary is
 510 not allowed (it must remain at zero pressure), no correction is necessary, meaning
 511 that the maximum soil infiltration capacity is met. If the net infiltration discharge
 512 is negative or liquid ponding is allowed above the surface, then the liquid velocity
 513 must be corrected to ensure the correct infiltration rate.

$$q_{net} = (n_L(\tilde{\mathbf{v}}_L^{t+\Delta t} - \tilde{\mathbf{v}}_S^{t+\Delta t}) - \widehat{\mathbf{w}}) \cdot \mathbf{n} \quad (36)$$

514 \mathbf{n} is the normal direction at the node that is determined by means of the gradient
 515 of mass.

516 The two approaches differ in terms of velocity correction. Martinelli et al. [31]
 517 correct both liquid and solid velocity, while Ceccato et al. [30] correct only liquid
 518 velocity. The first one accurately ensures the conservation of momentum balance
 519 of the mixture at the boundary, i.e. the total stress remains constant at the ground
 520 level. The second one implicitly assumes that the relative acceleration between
 521 solid and liquid is negligible and that the effective stress at the boundary is constant
 522 and equal to zero.

523 In order to emphasize the difference between the two implementations on the
 524 results, a 1D infiltration example is used for comparison. A 1m-high soil column
 525 is considered with the material parameters listed in Tab. 6. For simplicity,
 526 the intrinsic permeability is assumed constant and equal to $k_L = 1 \cdot 10^{-11} m^2$,

527 corresponding to an hydraulic conductivity of $\kappa = 1.0 \cdot 10^{-4}$ m/s. Van Genuchten
 528 SWRC is accounted (Eq. 23), with parameters $p_{ref} = 3 \text{ kPa}$, $\lambda = 0.7$, $S_{min} = 0.0$,
 529 $S_{max} = 1$.

Solid density [kg/m^3]	ρ_S	2700
Liquid density [kg/m^3]	ρ_L	1000
Porosity [-]	n	0.4
Liquid bulk modulus [kPa]	K_L	80000
Liquid dynamic viscosity [$\text{kPa} \cdot \text{s}$]	μ_L	$1 \cdot 10^{-6}$
Young modulus [kPa]	E	10000
Poisson ratio [-]	ν	0.2

Table 6. Material parameters for 1D infiltration example with applied infiltration velocity.

530 The column is discretised with 20 rows of 2 square triangular elements filled with
 531 3 MPs each (Fig. 10a). The bottom and lateral boundaries are impervious, while
 532 a vertical infiltration rate $\widehat{w}_y = 1.0 \cdot 10^{-4}$ m/s is applied at the top boundary. An
 533 initial suction of 2kPa is assigned along the column at $t = 0$, which corresponds
 534 to an initial degree of saturation $S_{L0} = 0.85$.

535 The liquid infiltrates from the top and flows down through the column accumu-
 536 lating at the impervious bottom. Liquid suction decreases while soil saturates.
 537 As expected, the approaches proposed by Martinelli et al. [31] and Ceccato et al.
 538 [30] give different results. Due to the different correction of solid velocity, soil
 539 displacements at the column head are different. The approach from Ceccato et al.
 540 [30] results in a non-realistic solid velocity at the beginning of the calculation.
 541 This difference reduces with time, i.e. with the reduction of relative acceleration
 542 between liquid and solid. Despite the results obtained in terms of liquid pres-

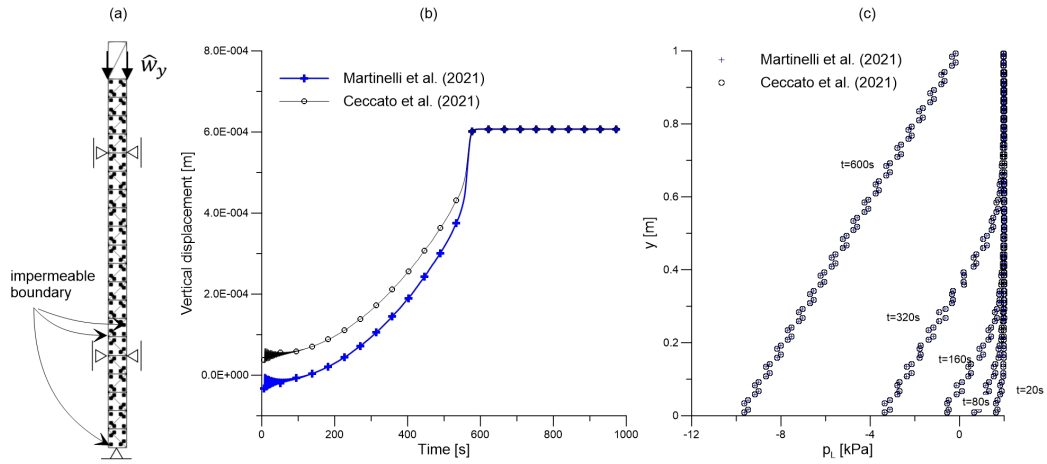


Fig. 10. 1D infiltration BC (a) geometry of the problem, (b) liquid pressure with depth, (c) displacement with time.

543 sures are very similar (Fig. 10b), the formulation from Martinelli et al. [31] is
 544 theoretically more accurate and therefore recommended.

545 In some cases, it is unknown if the boundary is an essential or a natural BC, for
 546 example, interface between soil and atmosphere where the fluid is free to exit at
 547 zero pressure when the soil is saturated. Still, it cannot enter when the soil is
 548 partially saturated. This boundary is called *potential seepage face*, and it is typical
 549 of the downstream surface of a dam or levee, or it can also arise on the reservoir
 550 side after a rapid drawdown of the water level.

551 Once more, Bandara et al. [25] solve this BC with the concept of *boundary layer*.
 552 If the boundary cell is fully saturated, i.e. $S_L = 1$ for all MP inside the cell, the
 553 pore pressure of all the MPs inside the cell is set to zero. If the boundary cell is
 554 unsaturated, i.e. there is at least 1 MP for which $S_L < 1$, then zero infiltration

555 rate is prescribed at the boundary nodes. The BC switches between natural and
556 essential depending on the saturation level of the *boundary cell*. Ceccato et al.
557 [30] solves the potential seepage face as a particular case of infiltration boundary
558 condition in which $\widehat{w} = 0$ and applies the predictor-corrector scheme described
559 previously. In this case, the BC switches between natural and essential depending
560 on the infiltration discharge (Eq. 36). Similarly, the formulation by Martinelli
561 et al. [31] can also be used for the same purpose.

562 **4. Applications**

563 All MPM formulations for partially saturated soils developed to simulate large-
564 deformation problems were first validated using applications where soil displace-
565 ment is limited. In general, the 1D infiltration problem is considered and the
566 MPM results are compared with analytical solutions [24, 29], experimental data
567 [25, 27, 16], or FEM calculations [33, 30, 31, 16].

568 For large-deformation applications, the MPM formulations for unsaturated soils
569 were mainly used to study rainfall-induced slope failures, with particular emphasis
570 on levees [30], embankments [24, 23] and landslides [25, 27, 32, 51, 16]. Ceccato
571 et al. [30] looked at the displacement evolution of a levee after failure induced
572 different boundary conditions, i.e. the rapid drawdown of the water level on the
573 river side or an infiltration due to heavy rainfall, highlighting the importance of
574 a large deformation tool for improving the accuracy of risk assessment. Yerro
575 [24], Yerro et al. [23] described the instability of a embankment slope subjected
576 to rain infiltration for a real case described in [52]. The model was able to show
577 the development of the initial failure surface, and the progressive large strain
578 deformation of the slope together with the dynamics of the motion characterised

579 by the history of displacement, velocity and acceleration of the unstable mass.
580 The first rainfall-induced landslide using MPM considering a rainfall boundary
581 condition was performed by Bandara et al. [25] for a case study inspired by the
582 full-scale field test performed in Rüdlingen, Switzerland [53]. The model was
583 able to show the negative effects on the stability of unsaturated slopes due to the
584 presence of continuous rainfall and shallow soil cover to bedrock surface. Wang
585 et al. [27] highlighted the importance of using coupled formulations for rainfall-
586 induced landslides, as large difference in retrogressive failure modes was observed
587 compared to the total stress analysis. More recently, the catastrophic event of
588 the Fei Tsui Road landslide, occurred on 13 August 1995 in Hong Kong, was
589 studied [32, 51, 54]. The first attempt was made by Lee et al. [32] where a linearly
590 distributed suction with depth was initialized and the rainfall was applied as zero
591 suction at the free surface. The evolution of the degree of saturation during the
592 failure is illustrated in Fig. 11. Later on, Liu et al. [51] revised the initial conditions
593 and simulated the failure process using a rainfall intensity equal to the saturated
594 conductivity of the soil. Lastly, Cuomo et al. [54] simulated the complete rainfall
595 event, exploring the failure process also in axisymmetric conditions.

596 Beside slope instability, other interesting applications using MPM for unsaturated
597 soils are in the field of site investigation, soil-structure interaction (e.g. pile
598 installation), excavation and trenching, and many others. For example, the role of
599 unsaturated behaviour during soil characterization by means of site investigation
600 (e.g. CPT, SPT, DMT, etc.) is relevant for soils above the water table. A new
601 example of cone penetration test (CPT) in unsaturated soil is illustrated hereafter.
602 The numerical simulation represents a penetration of a CPT, with radius $r_c = 3.6$
603 cm, in a virtual calibration chamber. The thickness of the soil domain is 1 m and

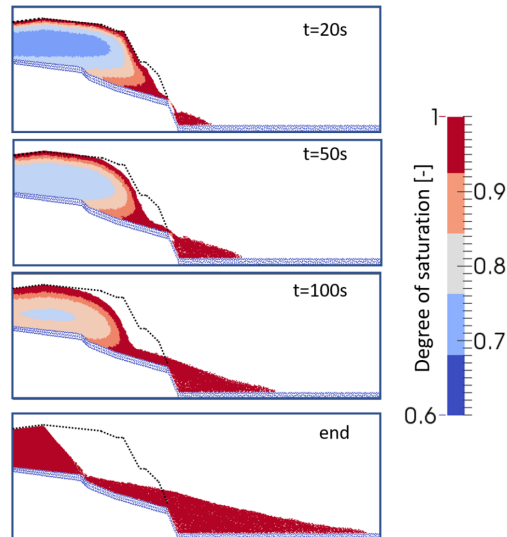


Fig. 11. Evolution of the degree of saturation during failure ([32])

604 the radius of 84 cm, loaded by 50 kPa constant in time. The computational mesh
 605 and numerical settings are equivalent to the one used by Martinelli and Galavi
 606 [55] for dry sand, but here unsaturated soil is considered. The soil behaviour
 607 is simulated using Mohr-Coulomb model with friction angle $\phi = 28^\circ$ and zero
 608 cohesion. The full list of material parameters is in Table 6 and the water retention
 609 curve is the one for "clay" in Table 5. The contact between pile and penetrometer is
 610 perfectly smooth. The simulations are performed using the unsaturated formulation
 611 described in [31], with the additional assumption that soil-water relative movement
 612 is negligible.

613 Three different values of initial suction in the liquid pressure are considered (i.e.
 614 0, 50 and 100 kPa), and the corresponding initial effective stress is then calculated
 615 to meet the equilibrium. The cone penetrates 40 cm into the soil sample and the
 616 cone resistance (q_c) is calculated as the ratio between the vertical reaction forces

617 on the cone and the area of the cross section of the penetrometer. The evolution
618 of q_c during the penetration is illustrated in Fig. 12, where the cone resistance
619 increases with the initial suction (u_0).

620 Fig.13 shows the distribution of pore water pressure in a close-up view of the
621 computational mesh, close to the cone, at the penetration of 40 cm. It is observed
622 that, if $u_0 = 0$, the soil domain is fully saturated and large pore water pressure
623 is computed nearby the cone and along the shaft. In this case, a soil region of
624 approximately 5 cone radii ($5r_c$) is affected by such a large pore water pressure
625 change. As the initial suction increases ($u_0 = 50$ or 100 kPa), the degree of
626 saturation decreases and the volumetric compressibility increases. It follows that,
627 during the penetration, large pore pressures can still be developed below the cone,
628 due to the large volumetric strains. However, the larger soil compressibility makes
629 such a pore pressure changes concentrated in the vicinity of the penetrometer,
630 whereas the rest of the domain remain almost unaffected.

631 Despite the relatively simple constitutive model used for to describe soil behaviour,
632 this example highlights the important role of the initial suction and the unsaturated
633 behaviour of soils on the measured cone resistance during site investigations.

634 As discussed in Sec. 3.1, all MPM formulations developed for unsaturated soil
635 are implemented in explicit, typically suitable for fast processes (e.g. impacts,
636 fast flows). For some of the applications which include long-term process (e.g.
637 consolidation or creep), the very small time step, required by explicit integration
638 schemes for stability criteria, makes the simulation extremely time consuming and
639 sometimes unfeasible. The implementation of implicit MPM codes will overcome
640 this issue.

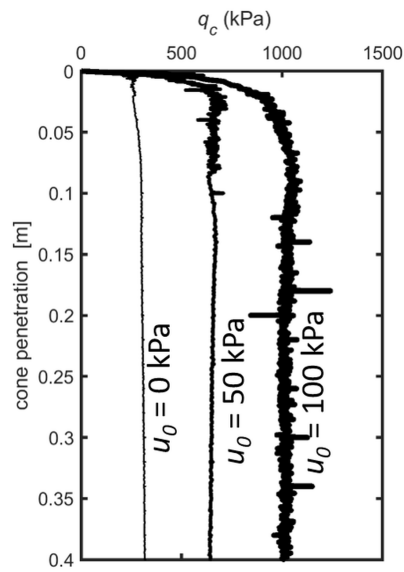


Fig. 12. Cone resistance with depth. Effect of initial suction (u_0) in water.

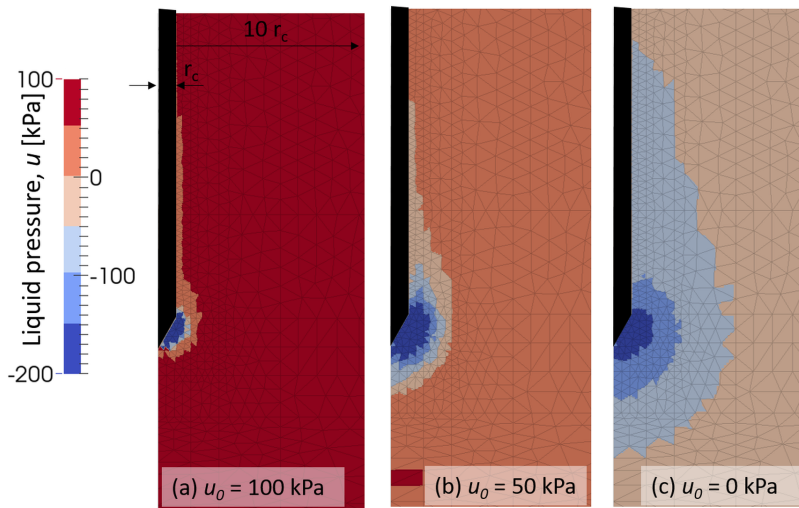


Fig. 13. Cone penetration of 40 cm. Distribution of pore water pressure in the soil nearby the cone. Effect of initial suction (u_0) in water: (a) $u_0 = 100$ kPa; (b) $u_0 = 50$ kPa; (c) $u_0 = 0$ kPa. (negative values are in compression)

641 **5. Conclusions**

642 This paper presented an overview of the MPM formulations available for the
643 simulation of unsaturated soils. Three-phase and two-phase formulations are
644 proposed in the literature. The first one accounts for solid-liquid-gas interactions,
645 while the second one neglects the pressure and density of the gas. While the three-
646 phase formulations can be essential for the simulation of problems where pressure
647 variations are important such as extremely rapid motions and explosion, the two-
648 phase formulations are adequate for the study of problems where the gas pressure
649 remains constant. The two-phase approach is computationally more efficient than
650 the three-phase one because a reduced number of governing equations is required.

651 Most of the formulations include all inertial terms, which is ideal when dealing with
652 dynamic problems. Neglecting the relative acceleration of the fluid with respect
653 to the solid is not recommended in dynamic problems where relative movement
654 between liquid and solid is expected. Finally, a few approaches neglect the gradient
655 of fluid mass when evaluating the mass balances. This assumption is acceptable
656 for most of the pressure gradients and SWRC present in many geotechnical appli-
657 cations, but it can induce errors when large gradients of pressure and degree of
658 saturation are expected.

659 All MPM formulations currently available for unsaturated soils are integrated
660 using explicit time schemes. The stability of the solution is limited by a critical
661 time step. This paper presents the first investigation on the effects of degree of
662 saturation (among other properties) on the stability of the solution. In general, the
663 critical time step increases in unsaturated conditions. Practical recommendations
664 are proposed based on the results to optimize the time step of the calculations.

665 Nevertheless, the implementation of implicit schemes is needed to study long-term
666 processes efficiently.

667 Despite the large number of geotechnical applications involving unsaturated soils,
668 a limited number of problems have been investigated, mainly focused on slope
669 stability and runout analysis. In this paper, preliminary results of a CPT model are
670 presented to illustrate the capabilities of unsaturated MPM in application involving
671 soil-structure interaction.

672 **6. Acknowledgements**

673 The second author greatly acknowledges the financial support from Fondazione
674 Ing. Aldo Gini.

675 **7. Appendix**

676 In this appendix, we present the steps followed to study the stability of the for-
677 mulation presented by Ceccato et al. [30] for the simulation of unsaturated soils.
678 Three different Δt_c criteria are obtained taking into the stability of (a) the coupled
679 system of momentum balance equations, (b) the momentum balance of the liquid
680 phase, and (c) the momentum balance of the mixture.

681 *7.1. Stability of the coupled system of momentum balance equations*

682 The governing equations in the strong form considered are the liquid momentum
683 balance and the mixture momentum balance as in

$$\rho_L \mathbf{a}_L = \nabla p_L - \mathbf{f}_L^d + \rho_L \mathbf{g} \quad (37)$$

$$n_S \rho_S \mathbf{a}_S + n_L \rho_L \mathbf{a}_L = \nabla \cdot \boldsymbol{\sigma} + \rho_m \mathbf{g} \quad (38)$$

684 Additionally, the mass balance equation is considered to express the liquid pressure
 685 as function of velocity (Eq.39). A linear elastic solid constitutive law (Eq.40) is
 686 assumed to express effective stress as function of solid displacements. The sign
 687 convention used expresses suction s as $s = -p_L$.

$$n \left(S_L \frac{\partial \rho_L}{\partial p_L} + \rho_L \frac{\partial S_L}{\partial p_L} \right) \frac{D^S p_L}{Dt} = \rho_L n S_L \operatorname{div}(\mathbf{v}_S - \mathbf{v}_L) - \rho_L S_L \operatorname{div}(\mathbf{v}_S) \quad (39)$$

$$d\boldsymbol{\sigma}' = \mathbf{E}_c \cdot d\boldsymbol{\varepsilon} \quad (40)$$

688 If a single node i is considered, the momentum balances per unit volume at time t_k
 689 can be written as a system of second-order ordinary differential equations (ODE).
 690 The homogeneous form for both equations is

$$\rho_L a_L^k + q_L v_L^k - q_L v_S^k + K_L u_L^k + \frac{(1-n)}{n} K_L u_S^k = 0 \quad (41)$$

$$(1-n) \rho_S a_S^k + n S_L \rho_L a_L^k + K_S u_S^k + \frac{(1-n)}{n} K_L u_S^k + K_L u_L^k = 0 \quad (42)$$

where

$$q_L = \frac{S_L n \mu_L}{k_L} \quad (43)$$

$$K_L = \frac{S_L}{L_i^2} \cdot \frac{1}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_w} \right)} \quad (44)$$

$$K_S = \frac{E_c}{L_i^2} \quad (45)$$

691 In this notation K_w is the bulk modulus of the liquid, while K_L is a term which
 692 includes K_w , S_L is the saturation degree, $\frac{dS_L}{ds}$ is the derivative of S_L respect to

693 suction, and L_i is the characteristic length of the element. The use of an Euler-
 694 Cromer time scheme leads to the following set of equations for the kinematic
 695 variables of liquid and solid phases.

$$v_L^{k+1} = v_L^k + \Delta t a_L^k \quad (46)$$

$$v_S^{k+1} = v_S^k + \Delta t a_S^k \quad (47)$$

$$u_L^{k+1} = u_L^k + \Delta t v_L^{k+1} \quad (48)$$

$$u_S^{k+1} = u_S^k + \Delta t v_S^{k+1} \quad (49)$$

696 From Eq.41 and Eq.42 it is possible to express accelerations as function of all the
 697 other terms as follows.

$$a_L^k = -\frac{q_L}{\rho_L} v_L^k + \frac{q_L}{\rho_L} v_S^k - \frac{K_L}{\rho_L} u_L^k - \frac{(1-n)}{n\rho_L} K_L u_S^k \quad (50)$$

$$a_S^k = -\frac{nS_L\rho_L}{(1-n)\rho_s} a_L^k - \frac{K_S}{(1-n)\rho_s} u_S^k - \frac{K_L}{n\rho_s} u_S^k - \frac{K_L}{(1-n)\rho_s} u_L^k \quad (51)$$

698 Now, the acceleration expressions (Eqs. 50 and 51) are substituted in the explicit
 699 time scheme set (Eqs. 48, 46, 49, 47).

$$u_L^{k+1} = u_L^k + \Delta t \left[v_L^k + \Delta t \left(-\frac{q_L}{\rho_L} v_L^k + \frac{q_L}{\rho_L} v_S^k - \frac{K_L}{\rho_L} u_L^k - \frac{(1-n)}{n\rho_L} K_L u_S^k \right) \right] \quad (52)$$

$$v_L^{k+1} = v_L^k + \Delta t \left(-\frac{q_L}{\rho_L} v_L^k + \frac{q_L}{\rho_L} v_S^k - \frac{K_L}{\rho_L} u_L^k - \frac{(1-n)}{n\rho_L} K_L u_S^k \right) \quad (53)$$

$$u_S^{k+1} = u_S^k + \Delta t \left[v_S^k + \Delta t \left(-\frac{nS_L\rho_L}{(1-n)\rho_s} a_L^k - \frac{K_S}{(1-n)\rho_s} u_S^k - \frac{K_L}{n\rho_s} u_S^k - \frac{K_L}{(1-n)\rho_s} u_L^k \right) \right] \quad (54)$$

$$v_S^{k+1} = v_S^k + \Delta t \left(-\frac{nS_L\rho_L}{(1-n)\rho_s} a_L^k - \frac{K_S}{(1-n)\rho_s} u_S^k - \frac{K_L}{n\rho_s} u_S^k - \frac{K_L}{(1-n)\rho_s} u_L^k \right) \quad (55)$$

700 The coefficients' matrix of the system of Eqs.52, 53, 54, 55 is evaluated to solve
701 the eigenvalues problem,

$$\det(A - \lambda I) = 0 \quad (56)$$

702 Thus, a quadratic equation for λ is obtained, where the terms $\geq \Delta t_c^3$ are neglected,

$$\lambda^2 + \lambda \left(\Delta t_c q_L Q - \frac{c\Delta t_c^2}{\rho_s} + \frac{K_L \Delta t_c^2}{\rho_L} - 2 \right) + \left(1 - \Delta t_c q_L Q \right) = 0 \quad (57)$$

703 If Eq.57 is solved for Δt_c by imposing $|\lambda| = 1$, it results in

$$\Delta t_c = \frac{-q_L Q + \sqrt{(q_L Q)^2 + 4 \left(\frac{K_L}{\rho_L} - \frac{c}{\rho_s} \right)}}{\left(\frac{K_L}{\rho_L} - \frac{c}{\rho_s} \right)} \quad (58)$$

704 Arranged in a more general form, it is equivalent to Eq.33 proposed by [24],

$$\Delta t_c = \frac{2}{\omega} \left[-\xi + \sqrt{\xi^2 + 1} \right] \quad (59)$$

705 where

$$\omega_c = \sqrt{\left(\frac{K_L}{\rho_L} - \frac{c}{\rho_s} \right)} = \frac{1}{L_i} \sqrt{\left[\frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_w} \right)} \right] \left[\frac{1}{\rho_L} + \frac{1}{\rho_s} \left(\frac{1}{n} - S_L \right) \right] + \frac{E_c}{\rho_s(1-n)}} \quad (60)$$

$$\xi_c = \frac{q_L Q}{2\omega_c} = \frac{S_L n \mu_L}{2k_L \omega_c} \left[\frac{S_L n}{(1-n)\rho_s} + \frac{1}{\rho_L} \right] \quad (61)$$

706 *7.2. Stability of the momentum balance of the liquid phase*

707 Same procedure presented for the coupled system can be applied to study the
 708 stability of the momentum balance of the liquid phase. By considering the ho-
 709 mogeneous form of Eq.41 (without the terms referring to the solid phase), we
 710 obtain

$$\rho_L a_L^k + q_L v_L^k + K_L u_L^k = 0 \quad (62)$$

711 The set of two ODEs (by considering the Euler-Cromer time scheme) is now

$$u_L^{k+1} = u_L^k + \Delta t \left[v_L^k + \Delta t \left(-\frac{q_L}{\rho_L} v_L^k - \frac{K_L}{\rho_L} u_L^k \right) \right] \quad (63)$$

$$v_L^{k+1} = v_L^k + \Delta t \left(-\frac{q_L}{\rho_L} v_L^k - \frac{K_L}{\rho_L} u_L^k \right) \quad (64)$$

712 The characteristic polynomial based on the resolution of the eigenvalues problem
 713 for Eqs. 63 and 64 is

$$\lambda^2 + \lambda \left(\frac{\Delta t_c q_L}{\rho_L} + \frac{K_L \Delta t_c^2}{\rho_L} - 2 \right) + \left(1 - \frac{\Delta t_c q_L}{\rho_L} \right) = 0 \quad (65)$$

714 Eq.65 solved for Δt_c gives

$$\Delta t_c = \frac{-\frac{q_L}{\rho_L} + \sqrt{\left(\frac{q_L}{\rho_L}\right)^2 + 4\frac{K_L}{\rho_L}}}{\frac{K_L}{\rho_L}} \quad (66)$$

715 Arranged in a more general form, it is equivalent to 33, where

$$\omega_L = \sqrt{\frac{K_L}{\rho_L}} = \frac{1}{L_i} \sqrt{\frac{S_L}{\rho_L \left(\frac{dS_L}{ds} + \frac{S_L}{K_w} \right)}} \quad (67)$$

$$\xi_L = \frac{\left(\frac{q_L}{\rho_L} \right)}{2\omega_L} = \frac{S_L n \mu_L}{2k_L \omega_L} \frac{1}{\rho_L} \quad (68)$$

716 *7.3. Stability of the momentum balance of the mixture*

717 Same procedure presented for the coupled system can be applied to study the sta-
 718 bility of the momentum balance of the mixture. By considering the homogeneous
 719 form of Eq.42, we obtain

$$(1 - n)\rho_S a_S^k + nS_L q_L v_S^k + \left[K_S + (1 - n) \left(-S_L + \frac{1}{n} \right) K_L \right] u_S^k = 0 \quad (69)$$

720 The set of 2 ODEs (by considering the Euler-Cromer time scheme) is now

$$u_S^{k+1} = u_S^k + \Delta t (v_s^k + \Delta t a_s^k) \quad (70)$$

$$v_S^{k+1} = v_S^k + \Delta t a_s^k \quad (71)$$

721 The characteristic polynomial (related to Eqs.70 and 71) is now

$$\lambda^2 + \lambda \left[P \Delta t_c^2 + \frac{S_L \Delta t_c n q_L}{(1 - n) \rho_S} - 2 \right] + \left[1 - \frac{S_L \Delta t_c n q_L}{(1 - n) \rho_S} \right] = 0 \quad (72)$$

722 where the term P is expressed as

$$P = \frac{1}{\rho_S} \left[\frac{K_S}{(1 - n)} + \left(-S_L + \frac{1}{n} \right) K_L \right] \quad (73)$$

723 If we solve Eq.72 for Δt_c we obtain

$$\Delta t_c = \frac{\left(-\frac{nS_L q_L}{(1-n)\rho_s} + \sqrt{\left(\frac{nS_L q_L}{(1-n)\rho_s}\right)^2 + 4P} \right)}{P} \quad (74)$$

724 Arranged in a more general form, it is equivalent to 33, where

$$\omega_M = \sqrt{P} = \frac{1}{L_i} \sqrt{\frac{1}{\rho_s} \left[\frac{E_c}{(1-n)} + \left(-S_L + \frac{1}{n} \right) \frac{S_L}{\left(\frac{dS_L}{ds} + \frac{S_L}{K_w} \right)} \right]} \quad (75)$$

$$\xi_M = \frac{\left(\frac{nS_L q_L}{(1-n)\rho_s} \right)}{2\omega_M} = \frac{S_L^2 n^2 \mu_L}{2(1-n)\rho_s k_L \omega_M} \quad (76)$$

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