A Manifold Learning Approach to Data-Driven Computational Mechanics

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Résumé — Standard simulation in classical mechanics is based on the use of two very different types of equations. The first one, of axiomatic character, is related to balance laws (momentum, mass, energy, ...), whereas the second one consists of models that scientists have extracted from collected, natural or synthetic data. In this work we propose a new method, able to directly link data to computers in order to perform numerical simulations. These simulations will employ universal laws while minimizing the need of explicit, often phenomenological, models. They are based on manifold learning methodologies. **Mots clés** — data-driven, model-free, manifold learning.

1 Introduction

In the present work we will assume that all the needed data is available. We will not address all the difficulties related to data generation or obtention from adequate experiments. This is a topic that, of course, remains open. On the contrary, we develop a method in which this stream of data plays the role of a constitutive equation, without the need of a phenomenological fitting to a prescribed model.

To better understand the data-driven rationale addressed in the present work, let us consider, for the sake of clarity, a very simple problem : linear elasticity. In that case the balance of (linear and angular) momentum leads to the existence of a symmetric second-order tensor σ (the so-called Cauchy's stress tensor) verifying equilibrium, expressed in the absence of body forces, as :

Find the displacement field $u \in (\mathcal{H}^1(\Omega))^3$ satisfying the essential boundary conditions $u(x \in \Gamma_D) = u_g(x)$ such that

$$\int_{\Omega} \varepsilon^* : \sigma \, d\boldsymbol{x} = \int_{\Gamma_N} \boldsymbol{u}^* \cdot \boldsymbol{t} \, d\boldsymbol{x}, \tag{1}$$

 $\forall u^* \text{ regular enough and vanishing on } \Gamma_D, \text{ i.e. } \forall u^* \in \left(\mathcal{H}_0^1(\Omega)\right)^3.$

The weak form given by Eq. (1) involves kinematic and dynamic variables from the test displacement field u^* and the stress tensor σ respectively. In order to solve it a relationship linking kinematic and dynamic variables is required, the so-called constitutive equation. The simplest one, giving rise to linear elasticity, is known as Hooke's law (even if, more than a law, it is simply a model), and writes

$$\boldsymbol{\sigma} = \lambda \mathrm{Tr}(\boldsymbol{\varepsilon}) \boldsymbol{I} + \boldsymbol{\mu} \boldsymbol{\varepsilon}, \tag{2}$$

where $Tr(\bullet)$ denotes the trace operator, and λ and μ are the Lame coefficients directly related to the Young modulus *E* and the Poisson coefficient v.

By introducing the constitutive model, Eq. (2), into the weak form of the balance of momentum, Eq. (1), a problem is obtained that can be formulated entirely in terms of the displacement field u. By discretizing it, using standard finite element approximations, for instance, and performing numerically the integrals involved in Eq. (1), we finally obtain a linear algebraic system of equations, from which the nodal displacements can be obtained.

In the case of linear elasticity there is no room for discussion : the approach is simple, efficient and has been applied successfully to many problems of interest. Today, there are numerous commercial codes making use of this mechanical behavior and nobody doubts about its pertinence in engineering practice.

However, there are other material behaviors for whom simple models fail to describe any experimental finding. These models lack of generality (universality) and for this reason a mechanical system is usually associated to different models that are progressively adapted and/or enriched from the collected data.

The biggest challenge could then be formulated as follows : can simulation proceed directly from data by circumventing the necessity of establishing a constitutive model? In the case of linear elasticity it is obvious that such an approach lacks of interest. However, in other branches of engineering science and technology it should be an appealing alternative to standard constitutive model-based simulations. In our opinion, we are at the beginning of a new era, the one of data-based or, more properly, data-driven engineering science and technology, where as much as possible data should be collected and information extracted in a systematic way by using adequate machine learning strategies. Then, simulations could proceed directly from this automatically acquired knowledge.

2 Constitutive manifold

Imagine, to begin with (more general scenarios will soon be considered) mechanical tests conducted on a perfectly linear elastic material, in a specimen exhibiting uniform stresses and strains. As previously indicated, in this paper we do not address issues related to data generation. Thus, for M randomly applied external loads, we assume ourselves able to collect M couples $(\sigma_m, \varepsilon_m), m = 1, ..., M$. These pairs could be represented as a single point X_m in a phase space of dimension D = 12 (the six distinct components of the stress and strain tensors, respectively). In the sequel Voigt notion will be considered, i.e. stress and strain tensors will be represented as vectors and the fourth-order elastic tensor reduces to a square matrix.

Each vector X_m thus defines a point in a space of dimension D and, therefore, the whole set of samples represents a set of M points in \mathbb{R}^D . We conjecture that all these points belong to a certain low-dimensional manifold embedded in the high-dimensional space \mathbb{R}^D . Imagine for a while that the M points belong to a curve, a surface or a hyper-surface of dimension $d \ll D$.

Therefore, appropriate manifold learning (or non-linear dimensionality reduction) techniques are needed to extract the underlying manifold (when it exists) in multidimensional phase spaces. A panoply of techniques exist to this end. In this work we focus on the particular choice of Locally Linear Embedding —LLE— techniques [1]. In the linear elastic behavior the application of the just described technique results, as expected, in a flat manifold of dimension two, i.e. d = 2. This is in perfect agreement to the fact that Hooke's law is completely characterized by two coefficients (either Young's modulus and Poisson coefficient, or Lame's coefficients) and is linear.

However, for the method to be useful, we need to define a strategy to solve problems stated in weak form and discretized by finite elements. Several options can be considered, which are described next.

1. **Identifying the locally linear behavior**. If we consider locally linear approximations, we can write

$$\mathbf{X}_m = \sum_{i=1}^M W_{mi} \mathbf{X}_i,$$

with $W_{mi} = 0$ if $i \notin S_m$ (set containing the *K*-nearest neighbors of \mathbf{X}_m). By minimizing the functional

$$\mathcal{H}(\mathsf{C}) = \sum_{i \in \mathcal{S}_m} (\sigma_i - \mathsf{C} \cdot \varepsilon_i)^2$$

we obtain $C(X_m) \equiv C_m$.

2. Identifying the locally linear *tangent* behavior. In order to consider Newton strategies the locally tangent linear behavior should be computed. Again, it is easy to obtain by considering $\Delta_{mi} \equiv X_m - X_i = (\sigma_m - \sigma_i, \varepsilon_m - \varepsilon_i)$ or $\Delta_{mi} = (\Delta \sigma_i^m, \Delta \varepsilon_i^m)$, $i \in S_m$. Because of the locally linear behavior around point X_m , we can write

$$\Delta \sigma_i^m = \mathsf{C}_T \cdot \Delta \varepsilon_i^m, \tag{3}$$

that allows defining the functional $\mathcal{H}_T(\mathsf{C}_T)$

$$\mathcal{H}_T(\mathsf{C}_T) = \sum_{i \in \mathcal{S}_m} (\Delta \sigma_i^m - \mathsf{C}_T \cdot \Delta \varepsilon_i^m)^2, \tag{4}$$

whose minimization results in the tangent elastic tensor $C_T(X_m) \equiv C_{T,m}$.

3. No model at all. The third level of description considers points X_m without trying to identify local behavior models at all. This third description was deeply discussed in [2].

3 Data-driven simulation

From the just identified locally linear behavior C(X) one could apply the simplest explicit linearization technique operating on the standard weak form

$$\int_{\Omega} \varepsilon^*(\boldsymbol{x}) : \boldsymbol{\sigma}^{n+1}(\boldsymbol{x}) \, d\boldsymbol{x} = \int_{\Gamma_N} \boldsymbol{u}^*(\boldsymbol{x}) \cdot \boldsymbol{t}(\boldsymbol{x}) \, d\boldsymbol{x}, \tag{5}$$

where at each point, from the stress-strain couple at position x, X(x), the locally linear behavior C(X(x)) can be obtained (in practice at the Gauss points used for the integration of the weak form) that allows us to write (using Voigt notation)

$$\int_{\Omega} \boldsymbol{\varepsilon}^*(\boldsymbol{x}) \cdot (\mathsf{C}(\boldsymbol{x}) \cdot \boldsymbol{\varepsilon}(\boldsymbol{x})) \ d\boldsymbol{x} = \int_{\Gamma_N} \boldsymbol{u}^*(\boldsymbol{x}) \cdot \boldsymbol{t}(\boldsymbol{x}) \ d\boldsymbol{x}.$$

This allows, in turn, to compute the displacement field and from it, to update the strain and stress fields, to compute again the locally linear behavior. The process continues until convergence. The discretization related to the other two descriptions just introduced were deeply considered in [2].

4 A simple nonlinear behavior : the linear-elastic perfectly-plastic behavior

we start by addressing the case of a linear-elastic-perfectly plastic 2D behavior. We assume the linear elastic contribution defined locally from $C(X^e)$ (X^e refers to the stress-elastic strain manifold) whereas the plastic contribution that involves the yield surface $f(\sigma)$ is assumed given by its own manifold.

Using again Voigt notation, the elastic behavior expressed from $\sigma = C \cdot \varepsilon^e$, where C represents the manifold-based elastic tensor and ε^e refers to the elastic component of the deformation (the reversible one). The total strain can be decomposed in its elastic and inelastic components,

$$\varepsilon = \varepsilon^e + \varepsilon^p$$
,

where we assume the plastic flow rate

$$\dot{\varepsilon}^p = \lambda \frac{\partial f(\sigma)}{\partial \sigma} = \lambda n,$$

where the yield surface $f(\sigma)$ is provided by experimental data. To generate these data *in silico*, we assume that it follows a von Mises model $f(\sigma) = \sigma_e - Y$, with Y the yield stress (no hardening is considered) and σ_e the equivalent stress related to the von Mises criterion. $f(\sigma)$ results in the surface represented in Fig. 1 where, for the sake of clarity, it is represented in the space of stresses.

The persistency condition $\dot{f}(\sigma) = 0$ when plastic flow occurs, results in the following plastic flow

$$\lambda = rac{oldsymbol{n}^T \cdot \mathsf{C} \cdot \dot{arepsilon}}{oldsymbol{n}^T \cdot \mathsf{C} \cdot oldsymbol{n}},$$

or in its incremental counterpart

$$\lambda = \frac{\boldsymbol{n}^T \cdot \mathsf{C} \cdot \Delta \varepsilon}{\boldsymbol{n}^T \cdot \mathsf{C} \cdot \boldsymbol{n}},$$

with now $\Delta \varepsilon^p = \lambda n$.

Here three fields must be considered, stress, strain and plastic strain. As soon as the last one is known, the elastic strain can be locally determined and the stresses obtained from the elastic manifold using the couple stress-elastic component of the strain.

In these expressions everything is properly defined except n, since we assume that the explicit form of the yield condition, i.e. $f(\sigma)$ is unknown and the only available data is the manifold depicted in Fig.

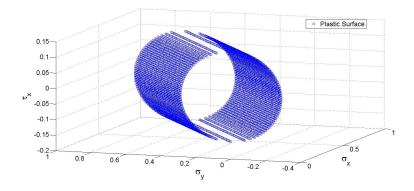


FIGURE 1 – Plastic manifold associated to the von Mises plasticity case.

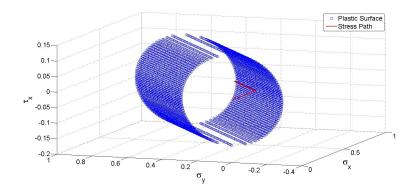


FIGURE 2 - Stress trajectory in the stress space in the elastic-perfectly plastic behavior

1. However, n is easily accessible by considering the normal vector to the plastic manifold depicted in Fig. 1.

Now one could imagine performing a standard linear elastic-perfectly plastic simulation by using a finite element explicit code where the plastic deformation is computed from the manifold that allows extracting n instead of the knowledge of function $f(\sigma)$ and its explicit derivative with respect to the stresses.

When considering the traction of a square domain along its right side, with appropriate boundary conditions on its left side (with tension-free conditions on the top and bottom boundaries) ensuring an homogeneous stress and strain fields everywhere in the domain, the stress trajectory in the stress space is depicted in Fig. 2. It can be noticed that the elastic behavior applies when the stress remains inside the plastic surface and then it remains in the surface during the plastic flow. Again, for the sake of simplicity, the results are shown in the stress domain.

Références

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