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## Article

# An information-theoretic framework for optimal design: analysis of protocols for estimating soft tissue parameters in biaxial experiments 

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#### Abstract

A new framework for optimal design based on the information-theoretic measures of mutual information, conditional mutual information, and their combination is proposed. The framework is tested on the analysis of protocols-combination of angles along which strain measurements can be acquired-in a biaxial experiment of soft tissues for the estimation of hyperelastic constitutive model parameters. The proposed framework sees information gain about the parameters from the experiment as the key criterion to be maximised which can be directly used for optimal design. Information gain is computed through $k$-nearest neighbour algorithms applied to the joint samples of the parameters and measurements produced by the forward and observation models. For biaxial experiments, the results show that low angles have relatively low information content compared to high angles. They also show that fewer number of angles with suitably chosen combinations can result in higher information gains when compared to a larger number of angles which are poorly combined. Finally, it is shown that the proposed framework is consistent with classical approaches, particularly the D-optimal design.


Keywords: optimal design; soft tissue mechanics; mutual information; biaxial experiment; inverse problems; information theory.

## 1. Introduction

Soft tissues exhibit complex biomechanical behaviour, including nonlinearity, anisotropy and heterogeneity [1]. Moreover, the tissues also demonstrate inelastic properties, such as rate-dependence, hysteresis and permanent set. The important link between the biomechanics and their physiological function have motivated a large number of ex-vivo studies aimed at characterizing their biomechanical properties. Given the complex interplay between the different aspects of their biomechanical properties, the experimental design of ex-vivo soft tissues is extremely challenging and has been a subject of investigation with a variety of experiments been proposed [2-6].

Since a variety of soft tissues are thin, e.g., blood vessels, heart valves and skin, biaxial testing is a widely used experimental technique that allows independently stretching the tissue in two orthogonal directions and measuring the corresponding forces $[7,8]$. Applying different stretches in two directions allows characterization of the in-plane anisotropic behavior of a given tissue, while a range of stretches provides us its nonlinear elastic response. However, even with this relatively simple set of options, the choices of what stretches to apply are unclear. Moreover, it is not obvious what these choices will depend on.

A variety of hyperelastic models have been developed to describe the anisotropic and nonlinear elastic properties of specific soft tissues [4,9-11]. The biaxial experimental data is commonly fit to these models in order to determine the model parameters. As the unknown parameters depend on specific model, the choice of experimental setup-the problem of optimal design-might depend on the choice of model. However, in practice, a predetermined set of experimental protocols are used.

In the present work, an optimal design problem is defined as finding the most suitable protocol in view of estimating the parameters of the material model. A comprehensive overview of the optimal design problem can be found in [12,13], and several criteria for optimal design have been proposed in the literature, often based on the minimisation of the variance of the parameters and sensitivities [14,15]. In the present work, we investigate a criterion based on information theoretic quantities, in the spirit of what has been proposed in the work [16] (from a Bayesian point of view) and [17]. The goal is to characterise the amount of information the measurements convey about the parameters we would like to estimate and then maximise that information. Estimating information theoretic quantities is in general a challenging problem, and this is especially the case in high-dimensional settings. In the present work, a model reduction method is coupled with a non-parametric sample based mutual information estimation in order to provide a pertinent estimation of the information theoretic quantities involved in the optimal design problem and then applied to the biaxial testing of soft tissues.

The structure of the work is as follows: in Section 2 the model and information theoretical aspects of the problem are introduced. In particular, in Section 2.1 we detail the mathematical model of the biaxial experiments for soft tissues: after having introduced the notation and the non-linear elasticity model, in Section 2.1 . 1 we particularise it to the biaxial testing experimental setup. In Section 2.1.2 we introduce the experimental protocol definition.

The second part of the Section is devoted to the description of the informationtheoretic framework for solving the optimal design problem. In Section 2.2.1 we introduce the problem, in Sections 2.2.2 and 2.2.3 the information-theoretic quantities and their numerical estimation are detailed. We then present the reduce order modeling method used and how to validate the results obtained by the proposed approach. The section ends with an overview of the method.

The results and the discussion are presented in Section 3, followed by the conclusion and the perspectives.

## 2. Methods

The methodological aspects are divided into two broad categories: the mathematical model of the biaxial experiments; and the information-theoretic optimal design framework.

### 2.1. Mathematical model of the biaxial experiments

We begin by defining the notation: a material point at its reference position $X \in \mathbb{R}^{3}$ moves to $x \in \mathbb{R}^{3}$ after deformation. The elastic behavior of soft tissues is described using hyperelastic strain energy density $\Psi$ which depends on the deformation gradient tensor $\mathbf{F}=\nabla_{X} \boldsymbol{x}$. The ratio of volume after deformation to that before deformation is given by $J=\operatorname{det}(\mathbf{F})$. Soft tissues are commonly regarded as incompressible due to high water content, i.e. $J$ is constrained to be unity.

We consider the hyperelastic model proposed by Gasser et al.[18] which defines the strain energy density

$$
\begin{equation*}
\Psi=\frac{k_{1}}{2 k_{2}}\left[e^{k_{2}\left(\kappa I_{1}+(1-3 \kappa) I_{4}-1\right)}-1\right]+\mu\left(I_{1}-3\right), \tag{1}
\end{equation*}
$$

where $I_{1}=\operatorname{tr}\left(\mathbf{F}^{\top} \mathbf{F}\right)$ is the first invariant of the right Cauchy-Green strain tensor $\mathbf{C}=\mathbf{F}^{\top} \mathbf{F}$ and $I_{4}=\boldsymbol{M} \cdot \mathbf{C M}$ is the fourth invariant representing stretch along fiber direction $\boldsymbol{M}$. The resulting Cauchy stress is given by

$$
\begin{equation*}
\sigma=2 \mathbf{F} \cdot \frac{\partial \Psi}{\partial \mathbf{C}} \cdot \mathbf{F}^{\top}-p \mathbf{I} \tag{2}
\end{equation*}
$$

where $p$ acts as the Lagrange multiplier to enforce incompressibility and $\mathbf{I}$ is the identity matrix.

For this model, the set of unknown parameters can be written as $\left\{k_{1}, k_{2}, \kappa, \mu\right\}$ assuming that the fiber direction $\boldsymbol{M}$ is known a priori (based on another experiment, e.g., light scattering [6]). In this paper, in order to simplify the problem, we assume that $\kappa=0.1$ and $\mu=0.1 \mathrm{kPa}$ are known and fixed. Thus, the aim of an ex-vivo biomechanical experiment is to determine parameters $k_{1} \in[5,100] \mathrm{kPa}$ and $k_{2} \in[5,80]$ robustly and with high confidence $[19,20]$. A commonly used experiment called the biaxial testing is described next.

### 2.1.1. Biaxial experiments for soft-tissues

Many of the soft tissue types are planar with small thickness. In a biaxial experiment, a square-shaped tissue sample is mounted via clamps or rakes and stretched along two orthogonal directions aligned with the sample edges (Fig. 1a). If these directions are used as the two coordinate axes and incompressibility is assumed, the stretching results in a diagonal deformation gradient tensor

$$
\begin{equation*}
\mathbf{F}=\operatorname{diag}\left[\lambda_{1}, \lambda_{2}, \frac{1}{\lambda_{1} \lambda_{2}}\right] \tag{3}
\end{equation*}
$$

where $\lambda_{1}$ is the stretch along first in-plane direction and $\lambda_{2}$ is the stretch along second in-plane direction. The fiber direction $M$ is generally aligned with the first coordinate axis, which results in only normal stress components. As no force is applied along the thickness of the tissue, $\sigma_{33}=0$ is used to determine the Lagrange multiplier $p$. Thus, we get

$$
\begin{align*}
& \sigma_{11}=2 \frac{\partial \Psi}{\partial I_{1}}\left[\lambda_{1}^{2}-\frac{1}{\lambda_{1}^{2} \lambda_{2}^{2}}\right]+2 \frac{\partial \Psi}{\partial I_{4}} \lambda_{1}^{2}  \tag{4}\\
& \sigma_{22}=2 \frac{\partial \Psi}{\partial I_{1}}\left[\lambda_{2}^{2}-\frac{1}{\lambda_{1}^{2} \lambda_{2}^{2}}\right] . \tag{5}
\end{align*}
$$

The applied stresses $\sigma_{11}, \sigma_{22}$ are controlled using load cells. The resulting strains defined as $e_{1}:=\lambda_{1}-1$ and $e_{2}:=\lambda_{2}-1$ are measured from the marker positions ${ }^{1}$. It is important to note that a homogeneous stress and strain state is assumed in the middle of the sample (Fig. 1a). Therefore, an implicit assumption is that the material properties and sample thickness are homogeneous. Moreover, these measurement techniques carry an error due to the limitations in measurement tools and/or the deviation from homogeneity, incompressiblity and material direction.

### 2.1.2. Protocol definition

In practice, there are two approaches to the biaxial experiment: 1) displacement controlled where known stretches are imposed and forces are measured and 2) force controlled where known forces are applied and stretches are measured. Generally, the force controlled approach is used as it is easier to implement. Therefore, in the force controlled approach, different values of stresses $\sigma_{11}$ and $\sigma_{22}$ can be applied.

[^0]
b)


Figure 1. a) A schematic of a biaxial experimental setup where a thin planar tissue sample (in light gray) is mounted via rakes and two orthogonal forces are applied to induce stresses $\sigma_{11}$ and $\sigma_{22}$, and the resulting strains are measured by tracking the locations of the markers (in dark gray). b) The $\sigma_{11}-\sigma_{22}$ space, where the applied stresses lie on the dotted line with a finite number of protocol angles $\phi$ used.

A single-angle biaxial protocol is defined as a straight line in the $\sigma_{11}-\sigma_{22}$ space (Fig. $1 b)$. That is, the ratio between the two stresses is kept constant while the applied forces are increased until a maximum value $\sigma^{\max }=200 \mathrm{kPa}$. Thus, for a chosen angle $\phi$, we apply:

$$
\begin{align*}
& \sigma_{11}= \begin{cases}\sigma & \text { if } \phi \leq \frac{\pi}{4} \\
\tan (\phi) \sigma & \text { else }\end{cases}  \tag{6}\\
& \sigma_{22}= \begin{cases}\cot (\phi) \sigma & \text { if } \phi \leq \frac{\pi}{4} \\
\sigma & \text { else }\end{cases} \tag{7}
\end{align*}
$$

where $\sigma \in\left[0, \sigma^{\max }\right]$. For $\sigma, 100$ linearly spaced observation points between zero and the maximum stress ( $\sigma^{\max }=200 \mathrm{kPa}$ ) are used. The resulting strains are calculated by iteratively solving equation (5) for $\lambda_{1,2}$ and thereby obtaining $e_{1,2}$. In practice, a combination of angles can be successively tested. We refer to this combination as the experimental protocol that needs to be optimally designed.

### 2.2. Information-theoretic framework for optimal design

The problem of optimal design typically refers to the choice of a design of experiments such that the design is optimal with respect to a pre-determined statistical criterion. We propose that the information-theoretic measures naturally define such statistical criteria. The central idea is that information gain [21,22] from an experiment or protocol-as quantified by the information-theoretic quantities of mutual information and conditional mutual information-can be directly used as a reasonable statistical criterion for optimal design. These quantities are described next after presenting the framework for optimal design.

### 2.2.1. Optimal design problem

Consider the following general model:

$$
\begin{equation*}
\mathbf{y}=\mathcal{M}(\boldsymbol{\theta}), \tag{8}
\end{equation*}
$$

where $\mathcal{M}$ denotes for the forward model that takes $\boldsymbol{\theta} \in \mathbb{R}^{m}$ and outputs $\mathbf{y} \in \mathbb{R}^{n}$. Note
that $\boldsymbol{\theta}$ may contain initial and boundary conditions of the model and that $\mathbf{y}$ may subsume
the output at many time-points in the case of a dynamical system. Subsequently, consider that the measurement model is as follows

$$
\begin{equation*}
\mathbf{z}=\mathcal{H}_{p}(\mathbf{y}, \boldsymbol{\theta})+\boldsymbol{\varepsilon}, \tag{9}
\end{equation*}
$$

where $\mathcal{H}_{p}$ represents the observation operator, $\mathbf{z} \in \mathbb{R}^{d}$ represents the measurement vector, and $\varepsilon$ represents the vector of measurement error/noise. Note that the the observation operator $\mathcal{H}_{p}$ depends on the design of experiments, which specifies which quantities are measured. Given a set of possible $\mathcal{H}_{p}=\left\{\mathcal{H}_{1}, \mathcal{H}_{2}, \cdots, \mathcal{H}_{h}\right\}$, and a statistical criterion $\mathcal{S}\left(\mathcal{H}_{p}\right)$ to be maximised, the optimal design is given by

$$
\begin{equation*}
\hat{\mathcal{H}}_{p}=\underset{\mathcal{H}_{p}}{\arg \max } \mathcal{S}\left(\mathcal{H}_{p}\right) . \tag{10}
\end{equation*}
$$

In the case of the biaxial experiments, the model $\mathcal{M}$ represents the model for the force controlled experiment (sections 2.1.1 and 2.1.2) and $\mathcal{H}_{p}$ essentially denotes the experimental protocol, see section 2.1.2, representing the combination of angles-each representing a straight line in the $\sigma_{11}-\sigma_{22}$ plane-along which the strain measurements of $e_{1}$ and $e_{2}$ are acquired. With the possible variation of each angle between 0 and $\pi / 2$, the set $\Phi$ of possible angles $\phi$ are constructed through a uniform discretisation of the space between 0 and $\pi / 2$ into $\alpha$ levels, thus

$$
\begin{equation*}
\Phi=\left\{\phi_{0}, \phi_{1}, \cdots, \phi_{\alpha}\right\} . \tag{11}
\end{equation*}
$$

The possible set of protocols is then given by any combination of elements in $\Phi$ with the restriction that the number of elements in a protocol be limited to $\mathcal{C}$. Thus, if $\bar{\Phi} \subset \Phi$ is a subset of angles representing a protocol, our set of protocols is given by

$$
\begin{equation*}
\mathcal{H}_{p}=\{\bar{\Phi} \subset \Phi|1 \leq|\bar{\Phi}| \leq \mathcal{C}\}, \tag{12}
\end{equation*}
$$

where $|\cdot|$ represents the number of elements in the set. In other words, we choose at least 1 and up to $\mathcal{C}$ elements from $\Phi$, with the total elements in $\mathcal{H}_{p}$ being

$$
\begin{equation*}
\left|\mathcal{H}_{p}\right|=\binom{\alpha}{1}+\binom{\alpha}{2} \cdots+\binom{\alpha}{\mathcal{C}} . \tag{13}
\end{equation*}
$$

### 2.2.2. Information-theoretic quantities for optimal design

In the framework of section 2.2.1, we propose that information-theoretic quantities of mutual information and conditional mutual information are a natural choice for the statistical criterion $\mathcal{S}$. Denoting the random variables associated with $\boldsymbol{\theta}$ and $\mathbf{z}$ as $\boldsymbol{\Theta}$ and $\mathbf{Z}$, respectively, the mutual information (MI) between the parameters $\boldsymbol{\Theta}$ and the measurements $\mathbf{Z}$ is defined as [21]

$$
\begin{equation*}
\mathcal{I}(\boldsymbol{\Theta} ; \mathbf{Z})=\int_{\mathcal{X}_{\boldsymbol{\Theta}} \times \mathcal{X}_{\mathbf{Z}}} p_{\boldsymbol{\Theta}, \mathbf{Z}}(\boldsymbol{\theta}, \mathbf{z}) \frac{p_{\boldsymbol{\Theta}, \mathbf{Z}}(\boldsymbol{\theta}, \mathbf{z})}{p_{\boldsymbol{\Theta}}(\boldsymbol{\theta}) p_{\mathbf{Z}}(\mathbf{z})} d \boldsymbol{\theta} d \mathbf{z} \tag{14}
\end{equation*}
$$

where $p_{X}(x)$ represents the probability density of a random variable $X$ with a realisation $X=x$ and support $\mathcal{X}_{X}$. The mutual information $\mathcal{I}(\boldsymbol{\Theta} ; \mathbf{Z})$ quantifies the amount of information that can be gained on average by knowing one random variable, say $\mathbf{Z}$, about the other, say $\boldsymbol{\Theta}$. Indeed, with this interpretation, MI is a good candidate for the statistical criterion $\mathcal{S}$ for optimal design. For an individual parameter, $\Theta_{i}$, or indeed for any combination of parameters $\left\{\Theta_{i}, \Theta_{j}\right\}$, the corresponding information gains can be similarly computed through $\mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)$ and $\mathcal{I}\left(\left\{\Theta_{i}, \Theta_{j}\right\} ; \mathbf{Z}\right)$, respectively. Thus, while $\mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)$ quantifies the information gain individually for the parameter $\Theta_{i}$, the quantity $\mathcal{I}\left(\left\{\Theta_{i}, \Theta_{j}\right\} ; \mathbf{Z}\right)$ quantifies information gain for the pair $\left\{\Theta_{i}, \Theta_{j}\right\}$ jointly. A measure of correlation between the parameters $\Theta_{i}$ and $\Theta_{j}$ is, however, missing, and is provided by conditional mutual information (CMI) defined as

$$
\begin{equation*}
\underbrace{\mathcal{I}\left(\Theta_{i} ; \Theta_{j} \mid \mathbf{Z}\right)}_{\mathrm{I}}=\underbrace{\mathcal{I}\left(\Theta_{i} ;\left\{\Theta_{j}, \mathbf{Z}\right\}\right)}_{\mathrm{II}}-\underbrace{\mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)}_{\mathrm{III}} . \tag{15}
\end{equation*}
$$

The CMI $\mathcal{I}\left(\Theta_{i} ; \Theta_{j} \mid \mathbf{Z}\right)$ represents the additional information gained about the parameter $\Theta_{i}$ when both $\Theta_{j}$ and $Z$ are known (term II) relative to when only the measurements $\Theta_{i}$ alone are known (term III). Note that CMI is symmetric, i.e. $\mathcal{I}\left(\Theta_{i} ; \Theta_{j} \mid \mathbf{Z}\right)=$ $\mathcal{I}\left(\Theta_{j} ; \Theta_{i} \mid \mathbf{Z}\right)$, and can be interpreted as a measure of dependence between the parameters given the measurements $\mathbf{Z}$. Also note that both MI and CMI are non-negative.

With the above background, many statistical measures can be constructed. For example:

1. The mutual information for any single parameter may be maximised, giving $\mathcal{S}=$ $\mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)$. This approach only concerns with the posterior of the parameter $\Theta_{i}$ and ignores all other parameters.
2. The joint mutual information may be maximised, giving $\mathcal{S}=\mathcal{I}(\boldsymbol{\Theta} ; \mathbf{Z})$. In the sense of classical optimal design this can be interpreted as the D -optimal design. This is because D -optimal designs minimise the determinant of the Fisher Information Matrix inverse, and $\mathcal{S}=\mathcal{I}(\boldsymbol{\Theta} ; \mathbf{Z})$ measures the information gain in the joint $\boldsymbol{\Theta}$ space.
3. The sum of individual parameter mutual information may be be maximised, giving $\mathcal{S}=\sum_{i=1}^{m} \mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)$. In the sense of classical optimal design this can be interpreted as the A-optimal design. This is because the the A-optimal design minimises the trace of the Fisher Information Matrix inverse, and $\mathcal{S}=\sum_{i=1}^{m} \mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)$ measures the sum of the information gains for all the parameters.
4. Alternatively, one may seek to maximise individual parameter information gain while minimising pairwise CMI, thus seeking both small posterior variances and minimising pairwise correlations between the parameters. In this case, the statistical criterion is

$$
\begin{equation*}
\mathcal{S}=\sum_{i=1}^{m} \mathcal{I}\left(\Theta_{i} ; \mathbf{Z}\right)-\tau \sum_{i=1}^{m} \sum_{j=i}^{m} \mathcal{I}\left(\Theta_{i} ; \Theta_{j} \mid \mathbf{Z}\right), \tag{16}
\end{equation*}
$$

where $\tau>0$ is a regularisation parameter.
Note that the above list is not exhaustive, and based on the interpretations of MI and CMI, other criteria may be constructed based on the desired sense of optimality.

### 2.2.3. Estimating mutual information

In general, the forward model in equation (8) is non-linear, and hence even if the observation operator is linear (implying linear combinations of the state are measured), the analytical computation of mutual information is intractable. Thus, the informationtheoretic quantities of MI and CMI must be estimated. A common method is to generate samples of $\boldsymbol{\Theta}$ through specification of an appropriate prior probability density $p_{\boldsymbol{\Theta}}(\boldsymbol{\theta})$. Denoting these $N_{s}$ samples as $\theta^{(i)}, i=\left\{1,2, \cdots N_{s}\right\}$, each $\boldsymbol{\theta}^{(i)}$ can be propagated through the forward and observation models of equations (8) and (9) to produce corresponding samples of $\mathbf{Z}$, denoted as $\mathbf{z}^{(i)}$. The samples of $\boldsymbol{\theta}^{(i)}$ and $\mathbf{z}^{(i)}$ can subsequently be used on non-parametric estimators of MI and CMI. Such non-parametric estimators can broadly be classified into two categories: kernel density estimators (KDE) [23] and $k$-nearest neighbour (kNN) estimators [24,25]. For an overview of such methods, we refer to [26]. While the estimator proposed by Kraskov et. al. [24] is widely used and performs very well across a range of scenarios, one of its drawbacks is that it suffers from higher errors when extreme correlations are present between the variables and/or when the the data effectively lives in a lower-dimensional manifold. Since we are working with models that specify explicit relationships between the variables through the forward and observation model, this is likely to be true for the data-set of $\left(\boldsymbol{\theta}^{(i)}, \mathbf{z}^{(i)}\right)$. Thus, in this study, we
employ the local non-uniformity correction (LNC) proposed in [27], which includes a correction term to the original estimator by Kraskov et. al. [24]. This term accounts for strong dependencies between the variables through local principle component analysis [27]. The method of [27] is used for estimation of all MIs, and CMIs are estimated from the difference of two MIs, see equation (15).

### 2.2.4. Dimensionality reduction for the biaxial experiment

One of the main difficulties in estimating information theoretic quantities is related to the data dimension. The non-parametric estimation is particularly challenging whenever the data are close to manifolds embedded in high-dimensional spaces. This is indeed the case when a physical model relates parameters and observable quantities. One of the possible ways to overcome this difficulty or, at least, to mitigate it, is (dimension or) model reduction, which aims at discovering the underlying low-dimensional structure of a set of data (a comprehensive review of the topic can be found in [28-31]). A large spectrum of methods have been proposed in the literature. In the present contribution, we adopt a local reduced-basis method (similar, in the spirit, to the methods proposed in [32,33]). Let the strains computed by the model be $e_{1,2}\left(\sigma ; \phi ; k_{1}, k_{2}\right)$, where $k_{1}$ and $k_{2}$ are the model parameters $\left(k_{1}, k_{2}\right) \in \Omega_{k} \subset \mathbb{R}^{2}$, and $\sigma \in \Omega_{\sigma} \subset \mathbb{R}$ is the variable defined in section 2.1.2. Let $n \in \mathbb{N}^{*}$, we introduce the following approximation:

$$
\begin{equation*}
e_{1,2} \approx \sum_{i=1}^{n} \eta_{i} r_{i}(\sigma, \phi) s_{i}\left(k_{1}, k_{2}, \phi\right), \tag{17}
\end{equation*}
$$

which is well defined by virtue of the Eckart-Young theorem. First, let us observe that a given protocol consists in a set of known angles $\bar{\Phi}$. An efficient way to construct the local reduced basis is hence to introduce a Proper Orthogonal Decomposition (POD) for each of the angles $\phi_{j} \in \bar{\Phi}$. This corresponds to looking for an approximation of the form:

$$
\begin{equation*}
e_{1,2}^{(j)}\left(\sigma ; \phi_{j} ; k_{1}, k_{2}\right) \approx \sum_{i=1}^{n} \eta_{i}^{(j)} r_{i}^{(j)}(\sigma) s_{i}^{(j)}\left(k_{1}, k_{2}\right) \tag{18}
\end{equation*}
$$

where $\left\langle r_{i}^{(j)}, r_{k}^{(j)}\right\rangle_{\Omega_{\sigma}}=\delta_{i k}$ and $\left\langle s_{i}^{(j)}, s_{k}^{(j)}\right\rangle_{\Omega_{k}}=\delta_{i k}\left(\langle\cdot, \cdot\rangle_{\Omega_{\sigma}, \Omega_{k}}\right.$ being the standard $L^{2}$ scalar product). The error in the approximation is related to the number $n$ of modes retained:

$$
\begin{equation*}
\left\|e_{1,2}^{(j)}-\sum_{i=1}^{n} \eta_{i}^{(j)} r_{i}^{(j)}(\sigma) s_{i}^{(j)}\left(k_{1}, k_{2}\right)\right\|_{L^{2}\left(\Omega_{\sigma} \times \Omega_{k}\right)}^{2}=\sum_{i=n+1}^{\infty} \eta_{i}^{(j)^{2}} \tag{19}
\end{equation*}
$$

In the present work, a number $n=4$ modes proved to be sufficient in order to get errors smaller than $10^{-3}$ in $L^{2}$ norm on the solution reconstruction. This means that the set of elements $e_{1,2}\left(\sigma ; \phi_{j} ; k_{1}, k_{2}\right)$ are close to the linear subspace spanned by the first $n=4$ modes $r_{i}^{(j)}$. Henceforth, instead of considering the discretised $e_{1,2}$ we will consider their coordinates in the subspace, given by:

$$
\begin{equation*}
\left.z_{1,2}^{(j)}\right|_{i}=\left\langle e_{1,2}, r_{i}^{(j)}\right\rangle_{\Omega_{\sigma}}=\eta_{i}^{(j)} s_{i}^{(j)}\left(k_{1}, k_{2}\right) \tag{20}
\end{equation*}
$$

### 2.2.5. Validation of results against existing methods

Several methods and criteria to define and reach an optimal design of experiments were proposed [12]. Among them, D-optimality criterion attempts at maximising the determinant of the information matrix. In the present case, this is equivalent to minimize the determinant of the inverse of the average Hessian of the loss function we would introduce in a classical parameter estimation method. In a noisy setting, and, in particular, when the noise is Gaussian, this cost function is equivalent to minus the logarithm of the likelihood function. Let the misfit function be $f(\theta)$ and $\mathbb{E}_{\Theta}$ denote the expectation operator. The average of the Hessian reads:

$$
\begin{equation*}
H=\mathbb{E}_{\Theta}\left[\left.\partial_{\theta}^{2} f\right|_{\theta_{*}}\right] \tag{21}
\end{equation*}
$$

where $\theta_{*}$ is the value of the parameter minimising the loss function.

### 2.2.6. Overview of approach for the biaxial experiments

In the context of the biaxial experiments, the parameters are $k_{1}$ and $k_{2}$, represented as random variables $K_{1}$ and $K_{2}$, respectively. The variability in these parameters is considered to be uniform (thus imposing a uniform prior distribution) in the following intervals: $k_{1} \in[5,100] \mathrm{kPa}$ and $k_{2} \in[5,80]$. For a single value of angle $\phi$, the measurements are the strain values $e_{1}$ and $e_{2}$ and measured at 100 points along the line defined by the angle $\phi$. Here, we consider $\alpha=16$ discrete values of possible measurement angles $\phi$ uniformly distributed between, and including, $0^{\circ}$ and $90^{\circ}$. For each angle $\phi$, separate reduced bases of 4 modes for $e_{1}$ and $e_{2}$ are constructed through POD over 400 values of ( $K_{1}, K_{2}$ ) sampled uniformly in the aforementioned parametric space. Thus, for any angle $\phi$, the dimensionality reduction approach projects $e_{1}$ and $e_{2}$ measured at 100 points along the line defined by $\phi$ to a basis of $4+4$ modes. For a given protocol consisting of multiple angles, the measurement vector $\mathbf{z}$ (with corresponding random variable $\mathbf{Z}$ ), is the collection of all the reduced basis representation of $e_{1}$ and $e_{2}$ along the angles in the protocol. Lastly, the maximum number of angles in a protocol is restricted to $\mathcal{C}=5$, giving a total of 6,884 unique combinations of the $\alpha=16$ angles.

For the estimation of MI and CMI, a total of $N=10,000$ values of $\left(K_{1}, K_{2}\right)$ are uniformly distributed in the parametric space. For each such sample $\left(k_{1}^{(i)}, k_{2}^{(i)}\right)$, the numerical model of the biaxial experiment is run to produce $e_{1}^{(i)}$ and $e_{2}^{(i)}$, which are then projected on to the reduced basis, giving $\mathbf{z}^{(i)}$. The $N$ triplets of $\left(k_{1}^{(i)}, k_{2}^{(i)}, \mathbf{z}^{(i)}\right)$ are subsequently used for estimation of MI and CMI through the LNC estimator (see section 2.2.3). In equation (16) we use $\tau=1$.

## 3. Results and discussion

For all the 6,884 combinations of angles, three statistical criteria are evaluated: (i) $\mathcal{I}(A ; \mathbf{Z}) ;\left(\right.$ ii $\mathcal{I}(B ; \mathbf{Z})$; and (iii) $\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$. While the first two criteria aim to maximise information gain about $A$ and $B$ individually, the third criterion aims to maximise information gain about $A$ and $B$ simultaneously while minimising the information dependence between them. Figures 2-4 show the variation in these three criterion when grouped by the number of angles in a protocol. In these Figures the values of information criterion when using two approaches to uniformly discretise the angular space within protocols is also presented. Observations from these plots are as follows:

1. Generally, all the three information criteria increase with increasing number of angles in the protocol. Intuitively, this is expected as higher number of angles imply more measurement data and hence a higher potential for improved estimation of the parameters. This observation is true for the maximum information gain, minimum information gain, and the mean information gain.
2. Across all the three criteria, it is observed that the uniform discretisation is not necessarily reflective of the best protocol for estimating the parameters. In fact, in most cases, the performance of uniform discretisation is close to the mean information gain observed across all the angle combinations.
3. From Figures 2 and 3 it is observed that the angular combinations that maximise information gain for $K_{1}$ are not identical—and vary significantly when more than 2 angles are simultaneously used-to those that maximise information gain for $K_{2}$. This further motivates the use of a criterion that balances information gains in both the parameters while minimising their interdependence.
4. Figure 4 shows that the best combinations that maximises a balanced criterion such as $\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$ are a trade-off between the combinations of
angles that maximises $\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)$ and $\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)$ individually. For example, when 5 angles are considered, the angles that maximise $\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)$ are $\phi_{a}=[66,72,78,84,90]$, those that maximise $\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)$ are $\phi_{b}=[30,36,42,48,54]$, while the combination that maximises $\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$ is $[30,36,48,78,90]$, which has two angles from $\phi_{a}$ and three angles from $\phi_{b}$. It should be noted that such a trade-off between maximising individual parameter gains is still significantly different than a uniform discretisation.
5. Finally, it is observed that the worst combinations are all low angles: $[0,6,12,18,24]$. This can be related to the fact that at low angles, the applied stress is largely aligned along the stiff fibers of the tissue, thus resulting in lower strain values. Thus, the lower angles provide a small range of the observations, while the larger angles provide a larger range (Fig. 9a), thereby containing more information about the parameters.


Figure 2. The variation of information criterion $\mathcal{S}=\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)$ across the 6,884 combinations grouped by the number of angles in a protocol. The vertical lines represent the variation around the mean value, which is shown in black circles. Black text shows the combinations that produce maximum and minimum values of $\mathcal{S}$. The red and blue pointers show $\mathcal{S}$ for angle combinations that follow a uniform discretisation of the angular space between 0 and 90 degrees. Red and blue texts show the associated angle combinations.


Figure 3. The variation of information criterion $\mathcal{S}=\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)$ across the 6,884 combinations grouped by the number of angles in a protocol. The vertical lines represent the variation around the mean value, which is shown in black circles. Black text shows the combinations that produce maximum and minimum values of $\mathcal{S}$. The red and blue pointers show $\mathcal{S}$ for angle combinations that follow a uniform discretisation of the angular space between 0 and 90 degrees. Red and blue texts show the associated angle combinations.


Figure 4. The variation of information criterion $\mathcal{S}=\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$ across the 6,884 combinations grouped by the number of angles in a protocol. The vertical lines represent the variation around the mean value, which is shown in black circles. Black text shows the combinations that produce maximum and minimum values of $\mathcal{S}$. The red and blue pointers show $\mathcal{S}$ for angle combinations that follow a uniform discretisation of the angular space between 0 and 90 degrees. Red and blue texts show the associated angle combinations.


Figure 5. The variation of information criterion $\mathcal{S}=\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$ across the 6,884 combinations. The vertical red lines show the groupings with respect to the number of angles in a protocol and $\mathcal{S}$ values are sorted in increasing order within each such grouping. The $x$-axis is represents the index associated with the protocol and is in logarithmic scale to capture the spread between 1 angle in a protocol ( 16 values) vs 5 angles in a protocol ( 4,368 values).

From this point onward, we present results only for the balanced information criterion $\mathcal{S}=\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$. Figure 5 shows the variation in $\mathcal{S}$ across all the combinations (x-axis, and in log-scale to capture the spread) grouped by the number of angles in a protocol and sorted in increasing order of $\mathcal{S}$ within each such group. Within each group, when observing the minimum and maximum values of $\mathcal{S}$, it shows that a better choice of angles can lead to more than $100 \%$ increase in the information gain compared to a poor choice. Furthermore, it shows that good combinations of a lower number of angles can lead to higher information gain compared to higher number of angles with bad combinations. For example, the maximum $\mathcal{S}$ when only one angle is used is higher than many combinations with $2-4$ angles. This emphasises the utility of optimal design and the proposed framework.

Figure 6 shows $\mathcal{S}$ for all the 6,884 combinations in increasing order of magnitude and Figure 7 shows a zoomed plot for the first 150 combinations along with the corresponding combinations of angles. Observing index values of 26 (red) and 28 (blue) in Figure 7 shows that even though four combinations are used in index 26 protocol, it produces a lower $\mathcal{S}$ compared to only a single angle used in index 28 protocol. Furthermore, since Figure 7 shows the first 150 out of 6,884 combinations of Figure 6 (which is sorted in creasing order of $\mathcal{S}$ ), all combinations here are relatively low $\mathcal{S}$ producing protocols. Observing the high density of angles in the region $\phi<24^{\circ}$, is indicative that lower values of angles, in particular those less than $24^{\circ}$ are relatively less informative when compared to higher values of angles. This behaviour is also apparent in Figure 8, which shows $\mathcal{S}$ values for protocols that use only one angle, and where a sharp jump can be observed when transitioning from $18^{\circ}$ to $24^{\circ}$. This peculiar behaviour may be explained by the physics of the biaxial experiment. Looking at the resulting strains $e_{1}$ and $e_{2}$ at this transition (Fig. 9b), we observe that the $e_{1}$ changes from positive to negative values. This behavior captures the important coupling between the two normal stresses and strains and is also related to the fiber dispersion in our constitutive model (equation (1), [18]). It is remarkable and encouraging that the information-theoretic framework captures the physics of the problem without explicitly considering it in the framework. While for simpler low-dimensional models the association between physics and optimal design may be relatively easy to see, inferring such behaviour is, in general, not trivial for more complex and higher-dimensional models.

Similarly to Figure 8, the results of the information-theoretic optimal design and further be analysed for higher number of angles. When two angles are considered, the $\mathcal{S}$ values in increasing order of magnitude and the corresponding angle combinations are shown in Figure 10. This figure re-iterates observations made previously: i) the choice of combinations significantly affects the information gain, the best combination gives
approximately 20 nats more information compared to the worst combination; and ii) generally speaking, higher angles are more informative compared to lower angles, in particular angles below $24^{\circ}$. While similar analysis for more than two angle combinations can be easily performed, efficient visual representation of such results is cumbersome and avoided in this manuscript.


Figure 6. The variation of information criterion $\mathcal{S}=\mathcal{I}\left(K_{1} ; \mathbf{Z}\right)+\mathcal{I}\left(K_{2} ; \mathbf{Z}\right)-\mathcal{I}\left(K_{1} ; K_{2} \mid \mathbf{Z}\right)$ across the 6,884 combinations sorted in increasing order of $\mathcal{S}$


Figure 7. Zoomed view into the first 150 protocols from Figure 6. The upper panel shows $\mathcal{S}$ and the lower panel shows the angles (by circles) in the corresponding protocol.


Figure 8. Information criterion against the angle when the protocols are restricted to a maximum of one angle.


Figure 9. Representative observations from the model with $k_{1}=40 \mathrm{kPa}$ and $k_{2}=40$. a) The observations using angles $\phi=0$ and 90 degrees, with the latter covering a significnantly larger range. b) The change in observations the angle is changed from 18, 24 , to 30 degrees shows a transition in $e_{1}$ from positive to negative values, indicating a coupling between the two directions.


Figure 10. Information criterion against the angles when the protocols are restricted to only two angles. The upper panel shows $\mathcal{S}$ and the lower panel shows the angles (by circles) in the corresponding protocol.


Figure 11. Information criterion $\mathcal{S}$ (in blue) and the $\log$ of the determinant of the Fisher Information Matrix inverse $\log _{10}\left|H^{-1}\right|$ (in red) against the angles when the protocols are restricted to a maximum of two angles. The upper panel shows $\mathcal{S}$ and $\log _{10}\left|H^{-1}\right|$, while the lower panel shows the angles (by circles) in the corresponding protocol. Note that here $e_{1}$ is shown in solid lines (left y-axis) and $e_{2}$ is shown in dashed lines (right y -axis).

To further illustrate the validity of the information-theoretic approach, comparison with a classical method, see section 2.2.5, is presented. For one and two angles in a protocol, Figure 11 shows a comparison between $\mathcal{S}$ and the $\log$ of the determinant of the Fisher Information Matrix inverse, $\log \left|H^{-1}\right|$. It is encouraging that a high correspondence between the two metrics is observed. In particular, increases in $\mathcal{S}$, implying higher information gains, are accompanied by corresponding decreases in $\log \left|H^{-1}\right|$, implying a smaller volume of the parameter posteriors. A Pearson correlation coefficient of $r=-0.76$ is observed between $\mathcal{S}$ and $\log \left|H^{-1}\right|$ implying a high similarity between the two metrics, and validating the information-theoretic approach in part. Let us remark that, when the number of the parameters become large, evaluating the Hessian would imply a non-negligible computational cost. On the contrary, the method used to evaluate the Mutual Information, being a primarily Monte Carlo based estimation, is less severely dependent on the number of parameters. Furthermore, computation of derivatives (either numerically or through adjoint based methods) may be cumbersome for certain type of models. Finally, we note that the effect of noise on information gain, and hence optimal design, can be easily assessed in the proposed framework by adding noise to the samples of $\mathbf{Z}$, see equation (9).

## 4. Conclusion

A framework for optimal design based on information-theoretic quantities of mutual information and conditional mutual information is proposed. The framework treats information gain as the central criterion for inverse problems and proposes several information-theoretic frameworks for desired sense of optimality. The capabilities of this framework are tested on the optimal design problem for biaxial experiments, where the effect of angle combinations along which the strains are measured is assessed on parameter estimation through information gain. Without including any physics based reasoning and purely through the information-theoretic measures, it is found that low angles $\leq 24^{\circ}$ are not very informative about the parameters relative to high angles.

These observations are then found to be consistent based on physics-based reasoning, thereby showing the efficacy of the proposed framework. Furthermore, it is demonstrated measurements for a low total number of angles which are carefully chosen can be more informative compared to the case when measurements along a high number of poorly chosen angles are acquired, thus highlighting both the importance of optimal design for biaxial experiments and the utility of the proposed framework in uncovering good angle combinations. Relation of the proposed framework to classical optimal design is performed and shown that the results produced by the new framework are consistent with classical frameworks.

## 5. Limitations and future work

While the proposed framework is shown to perform well on a two-parameter problem, its performance in higher parameter problems is not assessed. This assessment forms the the primary limitation and area of future assessment. In particular, the problems envisaged are largely related to the performance of the MI and CMI estimators in higher dimensions of both parameters and the measurements. While a dimensionality reduction approach was adopted in this study to minimise the adverse effects of the latter, this may not be possible in many forward and inverse problems. Thus, a large area of future work relates to the development of efficient and robust MI and CMI estimators. Lastly, a thorough comparison against classical optimal design methods (C-, E-, T-, V-optimal designs, etc.) needs to be performed, along with the construction and analysis of corresponding information-theoretic metrics.

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[^0]:    1 Although $e_{1}$ and $e_{2}$ are not the usual strain measures, we use these as our observations

