Elastic Full Procrustes Means for Sparse and Irregular Planar Curves

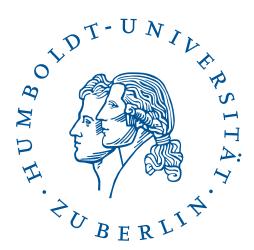
Masters Thesis

in partial fulfillment of the requirements for the degree

M.Sc. Statistics

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Figure: Example reparameterized SRV and coordinate functions				
Distance cannot be isometric				
Make duality of approach more explicit!				
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Citation, write more here!				
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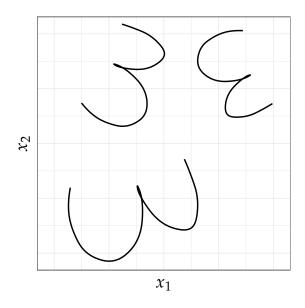
1. Introduction

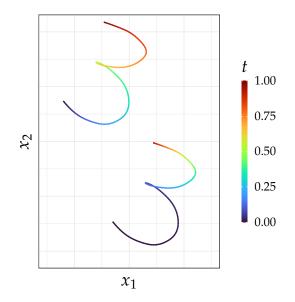
sec:intro

Statistical Shape Analysis (see e.g. DRYDEN and MARDIA 2016) is the branch of statistics concerned with modelling geometrical information. Such information might come in the form of outlines of bones or organs in a medical image, traced points along a handwritten digit, or data on the folding structure of a protein. This data is commonly captured using *landmarks*, which are characteristic points on the objects of interest that "match between and within populations" (DRYDEN and MARDIA 2016, p. 3). As an example, we might geometrically compare a set of mouse vertebrae by comparing the coordinates of prominent points along the bone outlines, which are common between all mouse vertebrae. More formally, we could say that each mouse vertebra's geometrical information is then represented by a landmark configuration $X \in \mathbb{R}^{k \times d}$, which is the stacked matrix of the k d-dimensional landmark coordinates, allowing for a multivariate treatment of shape or geometrical form.

A more flexible approach might be to treat e.g. the outline of an object as a whole, represented in the form of a continous curve $\beta:[0,1]\to\mathbb{R}^d$. Landmarks have the drawback that there is no clear way of choosing which points to include in the configuration, leaving the decision up to the subjectivity of the researcher. Furthermore, using landmarks leads to an inherently discrete treatment of the available data, which means modes of variation that lie in between landmarks may not be picked up by the analysis. By using curves, the analysis is not restricted to a fixed set of k discrete points, but instead uses all the available information. At the same time, the subjectivity in choosing the landmarks is eliminated. As each object then corresponds to one observation, this opens up a connection to the branch of statistics concerned with observations that are whole functions: Functional Data Analysis (see e.g. Ramsay and Silverman 2005).

When analyzing the geometry of objects, differences in location, rotation, and size are often not of interest. Instead, the focus lies purely on their differences in *shape*, a widely adapted definition of which was established by Kendall 1977 and which might be





- (a) The same digit with randomized rotation, scaling and translation applied three times.
- **(b)** Original (left) and re-parameterized digit (right; $t \mapsto t^5$). Color indicates the value of the parametrization $t \in [0,1]$ at point $\beta(t)$.



Figure 1.1.: Several representations of the same shape. Data: digits3.dat from the shapes package (Dryden 2019) for the R programming language (R Core Team 2021) with smoothing applied using methods discussed in Appendix A.3. Original dataset collected by Anderson 1997.

ig:1-shape

fig:1-eucl

formulated in the following way: "[A]ll the geometrical information that remains when location, scale and rotational effects are removed from an object"(DRYDEN and MARDIA 2016, p. 1). This is illustrated in Figure 1.1a, where the same shape of a handwritten digit "3" is plotted in three different orientations and sizes. When considering the shape of a curve $\beta:[0,1]\to\mathbb{R}^d$, one has to additionally take into account effects relating to re-parametrisation. As illustrated in Figure 1.1b, curves $\beta(t)$ and $\beta(\gamma(t))$, with some re-parametrisation or warping function $\gamma:[0,1]\to[0,1]$ monotonically increasing and differentiable, have the same image and therefore represent the same shape as well.

A pre-requisite for any statistical analysis of shape is the ability to calculate a distance between and to estimate a mean from observations, in a fashion that does not depend on location, rotation, scale and/or parametrization of the input. In this thesis two established approaches to shape analysis will be combined: Firstly, the *full Procrustes distance* and *mean* are widely used for translation-, rotation-, and scaling-invariant analysis of landmark data (see e.g. DRYDEN and MARDIA 2016, Chap. 4, 6). Secondly, SRIVASTAVA, KLASSEN, et al. 2011 introduced a mathematical framework for the *elastic*

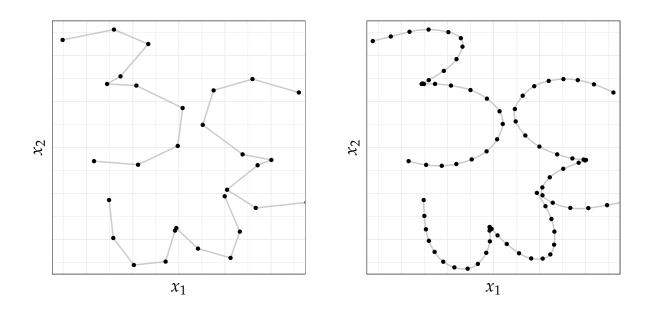
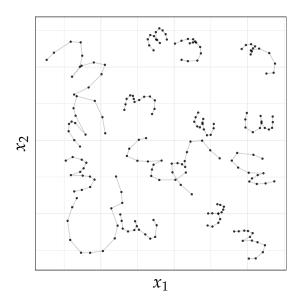


Figure 1.2.: Sparse (left) and dense (right) observations of the same three digits. Data: see Figure 1.1.

g:1-sparse

(re-parametrization invariant) shape analysis of curves, by using their square-root-velocity (SRV) representations. Taken together, both approaches allow for analysing curves in a fashion that is invariant to all four shape-preserving transformations, leading to an *elastic full Procrustes distance* and *mean*. As the full Procrustes mean has particularly nice properties in two dimensions, when identifying \mathbb{R}^2 with \mathbb{C} (see DRYDEN and MARDIA 2016, Chap. 8), this thesis will be restricted to the case of planar curves.

While we are interested in modeling a (planar) object's geometrical information as a continous curve $\beta:[0,1]\to\mathbb{R}^2$, the curve itself is usually only observed as a discrete set of points $\beta(t_1),\beta(t_2),\ldots,\beta(t_m)$. As shown on the left side of Figure 1.2 this is no problem when the number of observed points m is high and the whole length of the curve is densly observed, as we can easily interpolate any value $\beta(t)$ for $t\in[0,1]$. However, in cases where β is only observed over a small number of points (right side) and where the density and position of observed points may even vary between different curves—a setting known as *sparse* and *irregular*—more sophisticated smoothing techniques have to be applied. While the SRV framework has been combined with a Procrustes distance before, to estimate elastic shape means that also include invariance under scaling, rotation and translation (see Srivastava, Klassen, et al. 2011), these approaches have mostly focused on "Riemannian" or



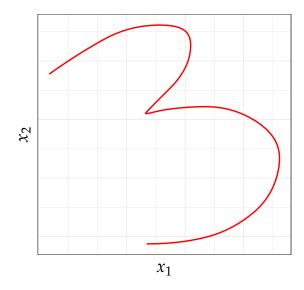


Figure 1.3.: Elastic full Procrustes mean function (right) estimated from sparse and irregular observations (left) using 13 knots and linear p-splines with a 2nd order penalty on SRV level. Data: digits3.dat with additional random rotation, scaling and translation applied.

fig:1-mean

"geodesic" mean concepts and are not specially designed with sparse or irregular observations in mind. On the other hand, as will be shown, the estimation of the *elastic full Procrustes mean* in two dimensions is related to an eigenfunction problem over the complex covariance surface of the observed curves. This offers an advantage, as we can then make use of established smoothing techniques for the estimation of covariance surfaces in the sparse and irregular setting. Here, in particular Cederbaum, Scheipl, and Greven 2018 offers a method for efficient covariance smoothing using *tensor product p-splines* (see e.g. Fahrmeier et al. 2013, Chap. 8.2).

The aim of this thesis, as illustrated in Figure 1.3, is to extend existing methods for elastic mean estimation of sparse and irregularly sampled curves, as proposed by STEYER, A. STÖCKER, and GREVEN 2021 and implemented in the package elasdics (STEYER 2021) for the R programming language (R CORE TEAM 2021), to also include invariance with respect to rotation and scaling. The later will be achieved by generalizing the concept of the full Procrustes mean from landmark to functional data and by iteratively applying full Procrustes mean estimation, rotation-alignment and parametrization-alignment, leading to the estimation of elastic full Procrustes means. Here, smoothing techniques for sparse estimation of the complex covariance surfaces as available in the R package sparseFLMM (Cederbaum, Volkmann, and A. Stöcker

2021) will be used.

The thesis is organized as follows. In Section 2 relevant background material is reviewed and the elastic full Procrustes mean is derived, in the case where curves $\beta:[0,1]\to\mathbb{R}^2$ are fully observed. Then, in Section 3, an estimation procedure for the setting of sparse and irregularly observed curves $\beta(t_1),\ldots,\beta(t_m)$ is proposed, concluding the theoretical part of this thesis. In Section 4 the proposed method will be verified and applied using simulated and empirical datasets. Finally, all results will be summarized in Section 5. Appendix A and Supplements B offer additional considerations and some reproducability guidelines.

2. Elastic Full Procrustes Means for Planar Curves

sec:theo

As a starting point, it is important to establish a notational and mathematical framework for the treatment of planar shapes. While the restriction to the 2D case might seem a major one, it still covers all shape data extracted from e.g. imagery and is therefore very applicable in practice. The outline of a 2D object may be naturally represented by a planar curve $\beta:[0,1]\to\mathbb{R}^2$ with $\beta(t)=(x_1(t),x_2(t))^T$, where $x_1(t)$ and $x_2(t)$ are the scalar-valued *coordinate functions*. Calculations in two dimensions, and in particular the derivation of the full Procrustes mean, are greatly simplified by using complex notation. As shown in 2.1, we will therefore identify \mathbb{R}^2 with \mathbb{C} and always use complex notation when representing a planar curve:

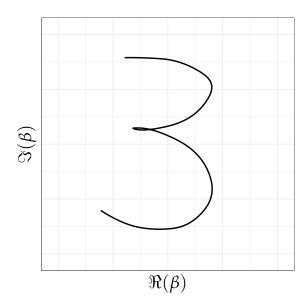
$$\beta: [0,1] \to \mathbb{C}, \quad \beta(t) = x_1(t) + i x_2(t).$$

We will assume the curves to be absolutely continuous, denoted as $\beta \in \mathcal{AC}([0,1],\mathbb{C})$, which guarantees integrable derivatives $\frac{\partial \beta(t)}{\partial t}$. This is important when working in the square-root-velocity (SRV) framework, as will be discussed in Section 2.2. All considerations will be restricted to the case of open curves, with possible extensions to closed curves $\beta \in \mathcal{AC}(\mathbb{S}^1,\mathbb{C})$ discussed in Section A.2 of the appendix.

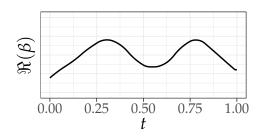
2.1. Equivalence Classes and Shape Invariance

theo:inv

As mentioned in the introduction, shape is usually defined by its invariance under the transformations of scaling, translation, and rotation. When considering the shape of curves, we additionally have to take into account invariance with respect to re-parametrisation. This can be seen, by noting that the curves $\beta(t)$ and $\beta(\gamma(t))$, with some re-parametrisation or warping function $\gamma:[0,1]\to[0,1]$ monotonically increasing and differentiable, have the same image and therefore represent the same geometrical object. We can say that the actions of translation, scaling, rotation, and



ig:2-curve



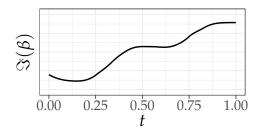


Figure 2.1.: Example of a planar curve (left) with respective coordinate functions (right) using complex notation. Data: see Figure 1.1.

re-parametrisation are *equivalence relations* with respect to shape, as each action leaves the shape of the curve untouched and only changes the way it is represented. The shape of a curve can then be defined as the respective *equivalence class*, i.e. the set of all possible shape preserving transformations of the curve. As two equivalence classes are neccessarily either disjoint or identical, we can consider two curves as having the same shape, if they are elements of the same equivalence class (see Srivastava and Klassen 2016, p. 40).

When defining an equivalence class, one has to first consider how the individual transformations act on a planar curve with complex representation $\beta:[0,1]\to\mathbb{C}$. This is usually done using the notion of *group actions* and *product groups*, with the later desciribing multiple transformations acting at once. A brief introduction to group actions may be found in Srivastava and Klassen 2016, Chap. 3.

- 1. The *translation* group $\mathbb C$ acts on β by $(\xi, \beta) \stackrel{\operatorname{Trl}}{\longmapsto} \beta + \xi$, for any $\xi \in \mathbb C$. We can consider two curves as equivalent with respect to translation $\beta_1 \stackrel{\operatorname{Trl}}{\backsim} \beta_2$, if there exists a complex scalar $\widetilde{\xi} \in \mathbb C$ so that $\beta_1 = \beta_2 + \widetilde{\xi}$. Then, for some function β , the related equivalence class with respect to translation is given by $[\beta]_{\operatorname{Trl}} = \{\beta + \xi \mid \xi \in \mathbb C\}$.
- 2. The *scaling* group \mathbb{R}^+ acts on β by $(\lambda, \beta) \stackrel{\operatorname{Scl}}{\longmapsto} \lambda \beta$, for any $\lambda \in \mathbb{R}^+$. We define $\beta_1 \stackrel{\operatorname{Scl}}{\leadsto} \beta_2$, if there exists a scalar $\widetilde{\lambda} \in \mathbb{R}^+$ so that $\beta_1 = \widetilde{\lambda} \beta_2$. An equivalence class is

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$$[\beta]_{Scl} = \{ \lambda \beta \, | \, \lambda \in \mathbb{R}^+ \}.$$

- 3. The *rotation* group $[0,2\pi]$ acts on β by $(\theta,\beta) \stackrel{\text{Rot}}{\longmapsto} e^{i\theta}\beta$, for any $\theta \in [0,2\pi]$. We define $\beta_1 \stackrel{\text{Rot}}{\leadsto} \beta_2$, if there exists a $\widetilde{\theta} \in [0,2\pi]$ with $\beta_1 = e^{i\widetilde{\theta}}\beta_2$. An equivalence class is $[\beta]_{\text{Rot}} = \{e^{i\theta}\beta \mid \theta \in [0,2\pi]\}$.
- 4. The *warping* group Γ acts on β by $(\gamma, \beta) \xrightarrow{\text{Wrp}} \beta \circ \gamma$, for any $\gamma \in \Gamma$ with Γ being the set of monotonically increasing and differentiable warping functions. We define $\beta_1 \xrightarrow{\text{Wrp}} \beta_2$, if there exists a warping function $\widetilde{\gamma} \in \Gamma$ with $\beta_1 = \beta_2 \circ \widetilde{\gamma}$. An equivalence class is $[\beta]_{\text{Wrp}} = \{\beta \circ \gamma \mid \gamma \in \Gamma\}$.

Wording and sentence structure not super clear in this paragraph.

In a next step, we can consider how these transformations act in concert and whether they *commute*, that is, whether the order of applying the transformations changes outcomes. Consider for example the actions of the rotation and scaling product group $\mathbb{R}^+ \times [0,2\pi]$ given by $((\lambda,\theta),\beta) \xrightarrow{\mathrm{Scl+Rot}} \lambda e^{i\theta} \beta$. These clearly commute, because the order of applying rotation or scaling do not make a difference, as $\lambda(e^{i\theta}\beta) = e^{i\theta}(\lambda\beta)$. However, the joint actions of scaling and translation do not commute, as $\lambda(\beta+\xi) \neq \lambda\beta+\xi$, with the same holding for the joint action of rotation and translation. As the order of translating and rotating or scaling matters, one usually takes the translation to act on the already scaled and rotated curve. The joint action defined using this ordering is usually called an *Euclidean similarity transformation*.

Definition 2.1 (Euclidean similarity transformation (DRYDEN and MARDIA 2016, p. 62)). We define an *Euclidean similarity transformation* of a curve $\beta : [0,1] \to \mathbb{C}$ as the joint action of scaling, rotation, and translation by

$$((\xi,\lambda,\theta),\beta)\mapsto \lambda e^{i\theta}\beta+\xi,$$

with
$$\xi \in \mathbb{C}$$
, $\lambda \in \mathbb{R}^+$, and $\theta \in [0, 2\pi]$.

With respect to the action of re-parametrization, we can note that it necessarily commutes with all Euclidean similarity transformations, as those only act on the image of β , while the former only acts on the parametrization. Putting everything together, we can finally give a formal definition of the shape of a planar curve.

Definition 2.2 (Shape). The *shape* of an absolutely continous planar curve $\beta \in \mathcal{AC}([0,1],\mathbb{C})$ is given by its equivalence class with respect to all Euclidean similarity transformations and re-parametrisations

$$[eta] = \left\{ \lambda e^{i heta} (eta \circ \gamma) + \xi \, | \, \xi \in \mathbb{C}, \, \lambda \in \mathbb{R}^+, \, heta \in [0, 2\pi], \, \gamma \in \Gamma
ight\}.$$

The *shape space* S is then given by the corresponding quotient space

$$\mathcal{S} = \mathcal{AC}([0,1],\mathbb{C})/\Gamma \times \mathbb{C} \rtimes \left(\mathbb{R}^+ \times [0,2\pi]\right) = \left\{ [\beta] \,|\, \beta \in \mathcal{AC}([0,1],\mathbb{C}) \right\} \right\},$$

where the symbol "×" denotes a semi-direct product, i.e. the translation group acting "after" scaling and rotation.

Is this clear?

2.2. The Elastic Full Procrustes Distance for Planar Curves

Rewrite motivation. Make approach to derivation more clear.

Let us now turn to the calculatation of distances between the shapes of curves. As shapes are represented by certain equivalence classes, and are therefore elements of a non-Euclidean quotient space, calculating their distance is not straight-forward. A common approach is to "project" the distance calculation in shape space down into the underlying functional space. For example, consider β_1 , $\beta_2 \in \mathbb{L}^2([0,1], \mathbb{C})$ with $[\beta_1]$, $[\beta_2]$ their equivalence classes with respect to all shape-preserving transformations. We might want to calculate their shape-distance as the minimal \mathbb{L}^2 -distance, when optimizing over all elements of their respective equivalence classes:

$$d([\beta_1],[\beta_2]) = \inf_{\widetilde{\beta}_1 \in [\beta_1], \, \widetilde{\beta}_2 \in [\beta_2]} d_{\mathbb{L}^2}(\widetilde{\beta}_1,\widetilde{\beta}_2) = \inf_{\widetilde{\beta}_1 \in [\beta_1], \, \widetilde{\beta}_2 \in [\beta_2]} \|\widetilde{\beta}_1 - \widetilde{\beta}_2\|_{\mathbb{L}^2}.$$

Which is equivalent to optimizing over all shape-preserving transformations:

$$\begin{split} d([\beta_{1}],[\beta_{2}]) &= \inf_{\lambda_{1,2} \in \mathbb{R}^{+},\, \theta_{1,2} \in [0,2\pi],\, \xi_{1,2} \in \mathbb{C},\, \gamma_{1,2} \in \Gamma} \, \left\| \, \lambda_{1} e^{i\theta_{1}}(\beta_{1} \circ \gamma_{1}) + \xi_{1} \right. \\ &\left. - \left(\lambda_{2} e^{i\theta_{2}}(\beta_{2} \circ \gamma_{2}) + \xi_{2} \right) \, \right\|_{\mathbb{L}^{2}}. \end{split}$$

theo:dist

However, this approach runs into problems, when considering whether all transformations act by isometries on this distance, i.e. whether equally changing the translation, rotation, scaling or re-parametrization of both curves affects their distance.

As it turns out, neither re-parametrization nor scaling are distance preserving when using the \mathbb{L}^2 -distance: For two equally re-parameterized curves $\widetilde{\beta}_{1,2} = \beta_{1,2} \circ \gamma$, their squared \mathbb{L}^2 -distance is given by $\|\beta_1 \circ \gamma - \beta_2 \circ \gamma\|^2 = \int_0^1 \|\beta_1(\gamma(t)) - \beta_2(\gamma(t))\|^2 dt = \int_0^1 \|\beta_1(s) - \beta_2(s)\|^2 \frac{1}{\dot{\gamma}(\gamma^{-1}(s))} ds$ with $s = \gamma(t)$. It follows that $\|\widetilde{\beta}_1 - \widetilde{\beta}_2\| \neq \|\beta_1 - \beta_2\|$, as in general $\dot{\gamma}(\gamma^{-1}(s)) \neq 1$. Likewise, it holds for equal re-scaling that $\|\lambda\beta_1 - \lambda\beta_2\| = \lambda \|\beta_1 - \beta_2\| \neq \|\beta_1 - \beta_2\|$. As one consequence the above optimization problem might simply be solved by $\lambda_1, \lambda_2 \to 0$, leading to $d([\beta_1], [\beta_2]) \to 0$ for any two curves β_1, β_2 . Furthermore, optimizing over re-parametrization using the \mathbb{L}^2 -distance has problems relating to the so called *pinching effect* and *inverse-inconsistency*, where the later means that aligning the parametrisation of one curve to another by $\inf_{\gamma \in \Gamma} \|\beta_1 - \beta_2 \circ \gamma\|$ may yield different results than $\inf_{\gamma \in \Gamma} \|\beta_2 - \beta_1 \circ \gamma\|$ (see Srivastava and Klassen 2016, p. 88).

Exact cita-tions!

A solution proposed in Srivastava, Klassen, et al. 2011 is to ditch the \mathbb{L}^2 -metric in favor of an *elastic metric*, which is isometric with respect to re-parametrization. Calculation of this metric, the Fisher-Rao Riemannian metric (Rao 1945), can be greatly simplified by using the *square-root-velocity* (SRV) framework, as the Fisher-Rao metric of two curves can be equivalently calculated as the \mathbb{L}^2 -distance of their respective SRV curves.

Definition 2.3 (SRV function of a planar curve (Srivastava, Klassen, et al. 2011)). The *SRV function* (SRVF) of an absolutely continuous planar curve $\beta \in \mathcal{AC}([0,1],\mathbb{C})$ is given by

$$q(t) = egin{cases} rac{\dot{eta}(t)}{\sqrt{\|\dot{eta}(t)\|}} & \dot{eta}(t)
eq 0 \\ 0 & \dot{eta}(t) = 0 \end{cases}$$

where $q \in \mathbb{L}^2([0,1],\mathbb{C})$ and $\dot{\beta}(t) = \frac{\partial \beta(t)}{\partial t}$.

More detail about bijective between AC and L^2 !

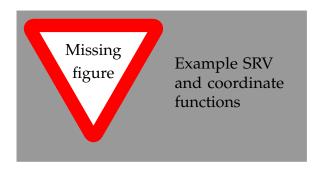


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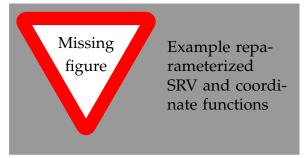


Figure 2.2.: SRV function of the planar curve from Figure 2.1 in original (left) and reparameterized (right; $t \mapsto t^5$) form, with respective coordinate functions on SRV level. Data: see Figure 1.1.

Remark. The original curve β can be re-constructed from its SRVF, up to translation, by $\beta(t) = \beta(0) + \int_0^t q(s) ||q(s)|| ds$.

As this representation makes use of derivatives, any curve β that has a SRVF must fulfill some kind of differentiability constraint. Here it is enough to consider only curves that are absolutely continuous $\beta \in \mathcal{AC}([0,1],\mathbb{C})$. In particular, this means that the original curves do not have to be smooth but might also be piecewise linear (see Srivastava and Klassen 2016, p. 91). The SRVFs are considered elements of a Hilbert space, which is given by $\mathbb{L}^2([0,1],\mathbb{C})$ equipped with the complex inner product $\langle \cdot, \cdot \rangle$ and corresponding norm $\|\cdot\|$. The complex inner product of $q, q' \in \mathbb{L}^2([0,1],\mathbb{C})$ is defined as

$$\langle q, q' \rangle = \int_0^1 \overline{q(t)} q'(t) dt$$
,

with $\bar{z} = \text{Re}(z) - i \, \text{Im}(z)$ denoting the complex conjugate.

As we can always recover the original curve up to translation, the SRV representation holds all relevant information about the shape of a curve. Furthermore, because of the use of derivatives, the SRV representation is invariant under changes in translation of the original curve. As a consequences, all shape-preserving transformations commute on SRV level.

Lemma 2.1. The actions of the translation, scaling, rotation, and re-parametrization groups commute on SRV level.

Proof. The SRVF $\widetilde{q}(t)$ of $\widetilde{\beta}(t) = \lambda e^{i\theta} \beta(\gamma(t)) + \xi$ is given by

$$\widetilde{q}(t) = \frac{\lambda e^{i\theta} \dot{\beta} \left(\gamma(t) \right) \dot{\gamma}(t)}{\sqrt{\left| \left| \lambda e^{i\theta} \dot{\beta} \left(\gamma(t) \right) \dot{\gamma}(t) \right| \right|}} = \sqrt{\lambda} e^{i\theta} \frac{\dot{\beta} \left(\gamma(t) \right)}{\sqrt{\left| \left| \dot{\beta} \left(\gamma(t) \right) \right| \right|}} \sqrt{\dot{\gamma}(t)} = \sqrt{\lambda} e^{i\theta} \left(q \circ \gamma \right) \sqrt{\dot{\gamma}(t)}.$$

The result is irrespective of the order of applying the transformations.

Remark. It follows, that the individual transformations translate to SRV level by

i.)
$$(\xi, q) \xrightarrow{\text{Trl}} q$$
, ii.) $(\lambda, q) \xrightarrow{\text{Scl}} \sqrt{\lambda} q$, iii.) $(\theta, q) \xrightarrow{\text{Rot}} e^{i\theta} q$, iv.) $(\gamma, q) \xrightarrow{\text{Wrp}} (q \circ \gamma) \sqrt{\dot{\gamma}}$.

Going forward, we will now work in the SRV framework and use the elastic metric for distance calculations between shapes. This means, instead of optimizing over the \mathbb{L}^2 distance between the original curves, we optimize over the \mathbb{L}^2 distance between their respective SRVFs. For two absolutely continuous curves $\beta_1, \beta_2 \in \mathcal{AC}([0,1], \mathbb{C})$ with respective SRV curves $q_1, q_2 \in \mathbb{L}^2([0,1], \mathbb{C})$ we might define the *elastic* distance between their shapes as

$$d([\beta_1], [\beta_2]) = \inf_{\widetilde{q}_1 \in [q_1], \widetilde{q}_2 \in [q_2]} \|\widetilde{q}_1 - \widetilde{q}_2\|_{\mathbb{L}^2}$$

or when equivalently optimizing over all possible transformations as

$$d([\beta_1], [\beta_2]) = \inf_{\lambda_{1,2} \in \mathbb{R}^+, \, \theta_{1,2} \in [0,2\pi], \, \gamma_{1,2} \in \Gamma} \|\sqrt{\lambda_1} e^{i\theta_1} (q_1 \circ \gamma_1) \sqrt{\dot{\gamma}_1} - \sqrt{\lambda_2} e^{i\theta_2} (q_2 \circ \gamma_2) \sqrt{\dot{\gamma}_2} \|_{\mathbb{L}^2}.$$

We can simplify the optimization, by considering that both rotation and warping act by isometries on the elastic metric, which means in practice it is enough to optimize over the rotation and warping of only one of the curves. For warping, we can reformulate the optimization as a problem over the relative parametrisation between the curves, as $\inf_{\gamma_1,\gamma_2\in\Gamma}\|(q_1\circ\gamma_1)\sqrt{\dot{\gamma}_1}-(q_2\circ\gamma_2)\sqrt{\dot{\gamma}_2}\|=\inf_{\gamma_1,\gamma_2\in\Gamma}\|q_1-(q_2\circ(\gamma_2\circ\gamma_1^{-1}))\sqrt{(\gamma_2\circ\gamma_1^{-1})}\|=\inf_{\gamma\in\Gamma}\|q_1-(q_2\circ\gamma)\sqrt{\dot{\gamma}}\|$. And smilarly for rotation, we can optimize over the relative rotation between both curves, as $\inf_{\theta_1,\theta_2\in[0,2\pi]}\|e^{i\theta_1}q_1-e^{i\theta_2}q_2\|=\inf_{\theta_1,\theta_2\in[0,2\pi]}\|q_1-e^{i(\theta_2-\theta_1)}q_2\|)=\inf_{\theta\in[0,2\pi]}\|q_1-e^{i\theta}q_2\|$. Taken together these lead to

$$d([\beta_1], [\beta_2]) = \inf_{\lambda_{1,2} \in \mathbb{R}^+, \, \theta \in [0,2\pi], \, \gamma \in \Gamma} \|\sqrt{\lambda_1} q_1 - \sqrt{\lambda_2} e^{i\theta} (q_2 \circ \gamma) \sqrt{\dot{\gamma}} \|,$$

which, however, still has the problem of not being isometric with respect to scaling.

A possible solution, mirroring the definition of the *full Procrustes distance* for landmark data, is to work with the normalized representations $z = \frac{q}{\|q\|}$, while only aligning

Distance cannot be isometric

the scaling of one of the curves. This leads to a distance that is isometric with respect to scaling, as for $\widetilde{q}_{1,2}=\lambda q_{1,2}$ with $\lambda\in\mathbb{R}^+$ it holds that $\|\widetilde{z}_1-\widetilde{z}_2\|=\|\frac{\lambda q_1}{\|\lambda q_1\|}-\frac{\lambda q_2}{\|\lambda q_2\|}\|=\|z_1-z_2\|$, while also being inverse consistent, as $\inf_{\lambda\in\mathbb{R}^+}\|z_1-\lambda z_2\|^2=\inf_{\lambda\in\mathbb{R}^+}\|z_1\|^2+\lambda^2\|z_2\|^2-\lambda(\langle z_1,z_2\rangle+\langle z_2,z_1\rangle)=\inf_{\gamma\in\Gamma}\|z_2\|^2+\lambda^2\|z_1\|^2-\lambda(\langle z_1,z_2\rangle+\langle z_2,z_1\rangle)=\inf_{\lambda\in\mathbb{R}^+}\|z_2-\lambda z_1\|^2$. We can finally take everything together and define an *elastic full Procrustes distance*.

def:dist

Definition 2.4 (Elastic full Procrustes distance). The *elastic full Procrustes distance* between the shapes $[\beta_1]$, $[\beta_2]$ of two continuously differentiable planar curves β_1 , $\beta_2 \in \mathcal{AC}([0,1],\mathbb{C})$ is given by

$$d_{EF}([\beta_1], [\beta_2]) = \inf_{\lambda \in \mathbb{R}^+, \theta \in [0, 2\pi], \gamma \in \Gamma} ||z_1 - \lambda e^{i\theta}(z_2 \circ \gamma) \sqrt{\dot{\gamma}}||$$

with normalized SRV representation $z_i = \frac{q_i}{\|q_i\|} \in \mathbb{L}^2([0,1],\mathbb{C})$, where q_i is the SRVF of β_i , i = 1, 2.

Remark. If the original curve β is of unit length $L[\beta] = \int_0^1 |\dot{\beta}(t)| dt = 1$, the SRV curve $q = \frac{\dot{\beta}}{\|\dot{\beta}\|}$ will be automatically normalized, as $\|q\| = \sqrt{\int_0^1 |\dot{q}(t)|^2 dt} = \sqrt{\int_0^1 |\dot{\beta}(t)| dt} = 1$.

Make duality of approach more explicit!

To calculate the distance, we need to solve the joint optimization problem over $\Gamma \times \mathbb{R}^+ \times [0,2\pi]$. For a fixed $\gamma \in \Gamma$, the optimization problem in Definition 2.4 mirrors the full Procrustes distance for landmark data, where an explicit solution is known in the planar case (see Dryden and Mardia 2016, Chapter 8). Likewise, for fixed rotation $\theta \in [0,2\pi]$ and scaling $\lambda \in \mathbb{R}^+$, there are known optimization techniques dealing with re-parametrisation.

$$(\lambda^*, \theta^*, \gamma^*) = \underset{\lambda \in \mathbb{R}^+, \theta \in [0, 2\pi], \gamma \in \Gamma}{\operatorname{argmin}} \|z_1 - \lambda e^{i\theta}(z_2 \circ \gamma) \sqrt{\dot{\gamma}}\|$$

Here, using one parameter $\omega=\lambda e^{i\theta}\in\mathbb{C}$ for rotation and scaling can simplify notation. The rotation and scaling parameters can always be recovered by $\lambda=|\omega|$ and $\theta=\arg(\omega)$.

$$(\omega^*, \gamma^*) = \underset{\omega \in \mathbb{C}, \gamma \in \Gamma}{\operatorname{argmin}} \|z_1 - \omega(z_2 \circ \gamma) \sqrt{\dot{\gamma}}\|$$

$$(2.1) \quad \boxed{\text{eq:argmin}}$$



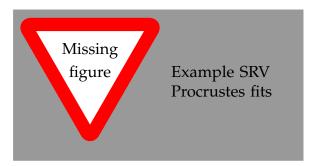


fig:2-pfit

Figure 2.3.: Procrustes fit of two example curves on original (left) and SRV (right) curve level. Data: see Figure 1.1.

The usual strategy is to optimize over the sets of parameters individually and then to iterate through both solutions. Let us first consider the optimization with respect to rotation and scaling. For fixed $\gamma \in \Gamma$ with $\tilde{z}_2 = (z_2 \circ \gamma) \sqrt{\dot{\gamma}}$ we can write Eq. 2.1 as

Citation write more ed argmi

$$\omega^* = \underset{\omega \in \mathbb{C}}{\operatorname{argmin}} \|z_1 - \omega \widetilde{z}_2\|, \tag{2.2}$$

which can be solved analytically.

lem:dist

Lemma 2.2. i.) For a fixed $\gamma \in \Gamma$, the optimal scaling and rotation solving Eq. 2.2 is

$$\omega^* = \langle \widetilde{z}_2, z_1 \rangle = \langle (z_2 \circ \gamma) \sqrt{\dot{\gamma}}, z_1 \rangle$$

ii.) The optimization problem in Definition 2.4 can then be reduced to

$$d_{EF}([\beta_1], [\beta_2]) = \inf_{\gamma \in \Gamma} \sqrt{1 - \langle z_1, (z_2 \circ \gamma) \sqrt{\dot{\gamma}} \rangle \langle (z_2 \circ \gamma) \sqrt{\dot{\gamma}}, z_1 \rangle}$$

Proof. See A.1.1 in the appendix.

Remark. For fixed $\gamma \in \Gamma$, we can use the first part of Lemma 2.2 to calculate the optimal rotation and scaling alignment of $\tilde{z}_2 = (z_2 \circ \gamma) \sqrt{\dot{\gamma}}$ onto z_1 . The alignment $\widetilde{z}_2^p = \langle \widetilde{z}_2, z_1 \rangle \widetilde{z}_2$ is called the *Procrustes fit* of \widetilde{z}_2 onto z_1 . The second part of Lemma 2.2 will be useful for mean calculation in the next section.

Given the current $\omega \in \mathbb{C}$ with $z_2^P = \omega z_2$, the optimization with respect to reparametrization can be written as

$$\gamma^* = \underset{\gamma \in \Gamma}{\operatorname{argmin}} \| z_1 - (z_2^P \circ \gamma) \sqrt{\dot{\gamma}} \|. \tag{2.3}$$

This is a well known problem and usually solved numerically by minimizing a cost function $H[\gamma] = \int_0^1 ||z_1(t) - z_2^p(\gamma(t)) \sqrt{\dot{\gamma}(t)}|| \, dt$ using a dynamic programming algorithm (DPA) or gradient based methods (see Srivastava, Klassen, et al. 2011). In this thesis we will use the methods laid out in Steyer, A. Stöcker, and Greven 2021, for solving Eq. 2.3 in the setting of sparse and irregularly sampled curves.

Anmerkung für Lisa Ich habe mich noch gefragt ob man anstatt Iteration über Eq. 2.2 und Eq. 2.3 die Distanz in Lemma 2.2 ii.) auch direkter optimieren kann.

$$\begin{split} \gamma^* &= \underset{\gamma \in \Gamma}{\operatorname{argmin}} \sqrt{1 - \langle z_1, \widetilde{z}_2 \rangle \langle \widetilde{z}_2, z_1 \rangle} \\ &= \underset{\gamma \in \Gamma}{\operatorname{argmax}} \langle z_1, \widetilde{z}_2 \rangle \langle \widetilde{z}_2, z_1 \rangle \\ &= \underset{\gamma \in \Gamma}{\operatorname{argmax}} \int_0^1 \int_0^1 \overline{z_1(t)} \underbrace{\widetilde{z}_2(t)}_{:=\widetilde{C}(s,t)} z_1(s) \, dt ds \\ &= \underset{\gamma \in \Gamma}{\operatorname{argmax}} \langle \widetilde{C} \, z_1, z_1 \rangle \\ &= \underset{\gamma \in \Gamma}{\operatorname{argmax}} \langle \widetilde{C} \, z_1, z_1 \rangle \end{split}$$

Mit $\widetilde{C}(s,t)$ der Kovarianz-Funktion von $\widetilde{z}_2=(z_2\circ\gamma)\sqrt{\dot{\gamma}}$. Habe das jetzt nicht weiter verfolgt (weil Zeit), aber vielleicht ist das ganz interessant? Man kann das glaube ich auch umschreiben mit $\breve{z}_1=(z_1\circ\gamma^{-1})\sqrt{\dot{\gamma}^{-1}}$ als $\mathrm{argmax}_{\gamma\in\Gamma}\langle C\,\breve{z}_1,\breve{z}_1\rangle$.

Write iterative procedure as algorithm! Basically as in Srivastava, Klassen, et al. 2011.

2.3. The Elastic Full Procrustes Mean for Planar Curves

Nummerierung der Equations überarbeiten

We now want to calculate shape means for a random sample of planar curves. Again, we assume all curves to be absolutely continous $\beta_i \in \mathcal{AC}([0,1], \mathbb{C})$ with corresponding SRV curves $q_i \in \mathbb{L}^2([0,1], \mathbb{C})$, $i=1,\ldots,N$. As we want to take into account shape invariance in the mean calculation, we cannot simply use sums or integrals to calculate a sample mean shape. Instead, we can use a more general concept for mean calculation, where the mean is defined as a minimizer over the sum of squared distances to each observation, for choice of a sensible distance in the space of interest. If the resulting

mean is a global minimum, it is usually called a "sample Fréchet mean" (Fréchet 1948), if it is a local minimum a "sample Karcher mean" (Какснек 1977) (see Dryden and Mardia 2016, p. 111).

def:mean

Definition 2.5 (Sample elastic full Procrustes mean). For a set of curves $\beta_i \in \mathcal{AC}([0,1], \mathbb{C})$, i = 1, ..., N, their *sample elastic full Procrustes mean* is given by the minimizing shape $[\hat{\mu}]$ with

$$[\hat{\mu}] = \underset{[\mu] \in \mathcal{S}}{\operatorname{arginf}} \sum_{i=1}^{N} d_{EF}([\mu], [\beta_i])^2,$$

where $S = \{ [\beta] : \beta \in \mathcal{AC}([0,1], \mathbb{C}) \}$ is the shape space.

Instead of working with equivalence classes as in Def. 2.5, it is often simpler to work with a specific element $\hat{\mu} \in [\hat{\mu}]$, that acts as a "representation" of the sample mean shape. One possibility is to use a representation that is of unit-length and starts at the origin, so that $L[\hat{\mu}] = \int_0^1 ||\dot{\hat{\mu}}(t)|| dt = 1$ and $\hat{\mu}(0) = 0$. This is an attractive choice when working in the SRV framework, as we do not have to worry about reconstructing translation when calculating $\hat{\mu}$ from its respective SRV curve $\hat{\mu}_q$ by $\hat{\mu}(t) = \hat{\mu}(0) + \int_0^t \hat{\mu}_q(s) \|\hat{\mu}_q(s)\| \, ds$. Another possibility would be to use a unit-length representation that is centered, so that $\left|\int_0^1 \hat{\mu}(t) dt\right| = 0$, which may be achieved by setting $\hat{\mu}(0) = \int_0^1 \int_0^t \hat{\mu}_q(s) \|\hat{\mu}_q(s)\| ds dt$ when reconstructing $\hat{\mu}$ from $\hat{\mu}_q$. From the point of shape analysis, the choice of representation does not make a difference, as both mean curves are elements of $[\hat{\mu}]$ and therefore have the same shape. However, the distinction becomes important when the estimated mean curve $\hat{\mu}$ is used in concert with other curves, for example when visualizing multiple curves or when comparing multiple class mean shapes, as those do not typically share the same center or starting point. Differences between both representations will be explored using empirical data of tounge shapes in Section ??.

Note that we can always construct $\hat{\mu}$ (and therefore $[\hat{\mu}]$) from $\hat{\mu}_q$ by integration. Turning back to the actual mean calculation, we can therefore use the Def. 2.4 to reformulate the optimization problem in Def. 2.5 on SRV level, where we now optimize

over possible normalized SRV representations μ_q of the mean shape $[\mu]$.

$$\hat{\mu}_q = \operatorname*{argmin}_{\mu_q \in \mathbb{L}^2: \, \|\mu_q\| = 1} \, \sum_{i=1}^N \, \left(\inf_{\omega_i \in \mathbb{C}, \gamma_i \in \Gamma} \|\mu_q - \omega_i(z_i \circ \gamma_i) \sqrt{\dot{\gamma}_i} \| \right)^2$$

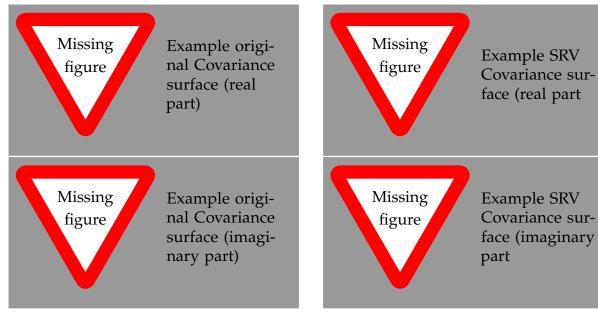
Here, $z_i = \frac{q_i}{\|q_i\|}$ are the observed normalized SRV curves, ω_i the rotation and scaling alignment to μ_q and γ_i the re-parametrisation alignment to μ_q . We can further simplify this by solving the optimization over ω_i using Lemma 2.2 ii.

$$\begin{split} \hat{\mu}_{q} &= \underset{\mu_{q} \in \mathbb{L}^{2}: \, \|\mu_{q}\| = 1}{\operatorname{argmin}} \sum_{i=1}^{N} \inf_{\gamma_{i} \in \Gamma} \left(1 - \langle \mu_{q}, \, (z_{i} \circ \gamma_{i}) \sqrt{\dot{\gamma}_{i}} \rangle \langle (z_{i} \circ \gamma_{i}) \sqrt{\dot{\gamma}_{i}}, \, \mu_{q} \rangle \right) \\ \hat{\mu}_{q} &= \underset{\mu_{q} \in \mathbb{L}^{2}: \, \|\mu_{q}\| = 1}{\operatorname{argmax}} \sum_{i=1}^{N} \sup_{\gamma_{i} \in \Gamma} \langle \mu_{q}, \, (z_{i} \circ \gamma_{i}) \sqrt{\dot{\gamma}_{i}} \rangle \langle (z_{i} \circ \gamma_{i}) \sqrt{\dot{\gamma}_{i}}, \, \mu_{q} \rangle \\ \hat{\mu}_{q} &= \underset{\mu_{q} \in \mathbb{L}^{2}: \, \|\mu_{q}\| = 1}{\operatorname{argmax}} \sum_{i=1}^{N} \sup_{\gamma_{i} \in \Gamma} \langle \mu_{q}, \, (\gamma_{i}, z_{i}) \rangle \langle (\gamma_{i}, z_{i}), \mu_{q} \rangle \end{split}$$

We end up with a two step optimization problem consisting of an outer optimization over μ_q and an inner optimization over the set $\{\gamma_i\}_{i=1,\dots,N}$. Similarly to the approaches discussed in Srivastava and Klassen 2016 and to Steyer, A. Stöcker, and Greven 2021, we solve this by *template based alignment* (see e.g. Srivastava and Klassen 2016, p. 271). In a first step the mean $\hat{\mu}_q$ is estimated while keeping the parametrisations fixed, after which the γ_i are updated by pairwise warping-alignment between z_i and $\hat{\mu}_q$, which is usually achieved by DPA or a gradient based approach. Both steps are iterated until the mean shape has converged. [TODO: Noch genauer auf warping alignment step eingehen -> muss ja erstmal auch procrustes fit berechnen um warping alignen zu können.]

Let us now consider the outer optimization problem for a fixed set of warping function $\{\gamma_i^*\}_{i=1,\dots,N}$ where we denote the warping aligned normalized SRV curves (γ_i^*, z_i) as \widetilde{z}_i . Note that if no warping alignment has happened yet, we can always set $\gamma_i^*(t) = t$ for all $i = 1, \dots, N$ as a starting value. The problem we have to solve is:

$$\begin{split} \hat{\mu}_{q} &= \underset{\mu_{q} \in \mathbb{L}^{2}: \, \|\mu_{q}\| = 1}{\operatorname{argmax}} \sum_{i=1}^{N} \, \left\langle \mu_{q}, \, \widetilde{z}_{i} \right\rangle \left\langle \widetilde{z}_{i}, \, \mu_{q} \right\rangle \\ \hat{\mu}_{q} &= \underset{\mu_{q} \in \mathbb{L}^{2}: \, \|\mu_{q}\| = 1}{\operatorname{argmax}} \sum_{i=1}^{N} \, \int_{0}^{1} \int_{0}^{1} \overline{\mu_{q}(s)} \widetilde{z}_{i}(s) \overline{\widetilde{z}_{i}(t)} \mu_{q}(t) \, ds \, dt \end{split}$$



(a) Covariance surface on original curve level.

(b) Covariance surface on SRV curve level.

Figure 2.4.: Complex covariance surface on original and SRV curve level. Data: see Figure 1.1.

$$\hat{\mu}_q = \operatorname*{argmax}_{\mu_q \in \mathbb{L}^2: \, \|\mu_q\| = 1} \ \int_0^1 \int_0^1 \overline{\mu_q(s)} \left(\sum_{i=1}^N \, \widetilde{z}_i(s) \overline{\widetilde{z}_i(t)} \right) \mu_q(t) \, ds \, dt$$

We can identify the inner term as proportional to a sample estimator $\check{C}(s,t) = \frac{1}{N} \sum_{i=1}^{N} z_i(s) \overline{z_i(t)}$ of the population covariance surface $C(s,t) = \mathbb{E}[z(s) \overline{z(t)}]$, when noting that $\mathbb{E}[z(t)] = 0$ for all $t \in [0,1]$ due to rotational symmetry. [TODO: Hier wird die Notation etwas unsauber. Man hat hier ja einen Cov. Estimator aus den "gewarpten" Kurven. Vielleicht müsste man sich hier nicht nochmal kurz Gedanken machen wie C(s,t) eigentlich aussieht, wenn man berücksichtigt, dass die z(t) auch random gewarpt sind. Vielleicht ist's auch egal.]

$$\hat{\mu}_q = \underset{\mu_q \in \mathbb{L}^2: \|\mu_q\| = 1}{\operatorname{argmax}} N \cdot \int_0^1 \int_0^1 \overline{\mu_q(s)} \check{\mathsf{C}}(s, t) \mu_q(t) \, ds \, dt$$

By replacing $\check{C}(s,t)$ by its expectation C(s,t), we can analogously formulate an estimator on the population level. **[TODO: Notation?]**

$$\hat{\mu}_q = \underset{\mu_q \in \mathbb{L}^2: \|\mu_q\| = 1}{\operatorname{argmax}} \int_0^1 \int_0^1 \overline{\mu_q(s)} C(s, t) \mu_q(t) \, ds \, dt$$

We can rewrite this again as a functional scalar product by considering the covari-

vvc can

fig:2-cov

ance operator C with $(C\mu_q)(s) = \int_0^1 C(s,t)\mu_q(t)dt$ (see Ramsay and Silverman 2005, p. 153).

$$\hat{\mu}_q = \underset{\mu_q \in \mathbb{L}^2, \|\mu_q\| = 1}{\operatorname{argmax}} \left\langle \mu_q, C\mu_q \right\rangle \tag{2.4} \quad \text{eq:quadr_c}$$

This is a well known problem in the context of functional principal component analysis (FPCA). From $\overline{C(s,t)} = \overline{\mathbb{E}[z(s)\overline{z(t)}]} = \mathbb{E}[z(t)\overline{z(s)}] = C(t,s)$ it follows that $\langle \mu_q, C\mu_q \rangle = \langle C\mu_q, \mu_q \rangle$ and therefore that C is a *self-adjoint* operator. The optimization problem then reduces to an eigenfunction problem

$$Cu = \lambda u \quad \Leftrightarrow \quad \int_0^1 C(s,t)u(t) \, dt = \lambda u(s) \,,$$
 (2.5) eq:funceign

where $\lambda = \langle \mu_q, C\mu_q \rangle$ is the target function to maximize. For normalized eigenfunctions u_1, u_2, \ldots and corresponding eigenvalues $\lambda_1 \geq \lambda_2 \geq \ldots$ of C(s,t), the solution $\hat{\mu}_q(t)$ is given by the leading normalized eigenfunction $u_1(t)$ of C(s,t) (see Ramsay and Silverman 2005, pp. 153, 397).

Algorithm 2.1 (Sample elastic full Procrustes mean). Let $\{\beta_i\}_{i=1,...,N}$ be a sample of planar curves with corresponding SRV curves $\{q_i\}_{i=1,...,N}$. Let $z_i = \frac{q_i}{\|q_i\|}$. Set $\gamma_i^0 = t$ for all i=1,...,N as the initial parametrisation alignment. Set k=0.

1. For
$$i = 1, ..., N$$
: Set $z_i^k(t) = z_i \left(\gamma_i^k(t) \right) \cdot \sqrt{\dot{\gamma}_i^k(t)}$.

- 2. Estimate $C(s,t) = \mathbb{E}[z(s)\overline{z(t)}]$ from $\{z_i^k\}_{i=1,\dots,N}$. Call this estimate $\hat{C}^k(s,t)$.
- 3. Set μ_q^k as the leading normalized eigenfunction of $\hat{C}^k(s,t)$. Stop if μ_q^k is close to μ_q^{k-1} .
- 4. For $i=1,\ldots,N$: Calculate the optimal rotation and scaling alignment $\omega_i^k=\left\langle z_i^k,\,\mu_q^k\right\rangle$.
- 5. For $i=1,\ldots,N$: Solve $\gamma_i^{k+1}=\operatorname{argmin}_{\gamma\in\Gamma}\|\mu_q^k-(\omega_i^k\cdot z_i\circ\gamma)\sqrt{\dot{\gamma}}\|$.
- 6. Set k = k + 1 and return to Step 1.

3. Estimation Strategy for Sparse and Irregular Observations

sec:est

Create figures for everything.

Alg. 2.1 shows an idealized version of the elastic full Procrustes mean estimation, where it is assumed that each curve β_i is fully observed. This is not the case in practice, as each observation β_i is usually itself only observed at a finite number of discrete points $\beta_i(t_{i1}), \ldots, \beta_i(t_{im_i})$. Additionally, the number of observed points per curve m_i might be quite small and the points do not need to follow a common sampling scheme across all curves, a setting which is respectively known as *sparse* and *irregular*.

Following the steps laid out in Alg. 2.1, this section proposes a mean estimation strategy for dealing with sparse and irregular observations. In a first step, the construction of SRV and warped SRV curves from discrete (and possibly sparse) observations will be shown in Section 3.1. Section 3.2 discusses efficient estimation of the complex covariance surface C(s,t) from sparse observations. In Section 3.3 the calculation of the leading eigenfunction u_1 of C(s,t) in a fixed basis will be shown. Section 3.4 deals with the estimation of the scalar product $\omega = \langle z, \mu_q \rangle$, which gives the optimal rotation and scaling alignment. Note that the final warping alignment step in Alg. 2.1 can be solved by using methods for warping alignment of sparse and irregular curves provided in Steyer, A. Stöcker, and Greven 2021.

c:discrete

3.1. Discrete Treatment of SRV Curves

A natural first consideration might be how to calculate SRV curves from sparse observations. We defined the SRV curve of a function $\beta \in \mathcal{AC}([0,1],\mathbb{C})$ as $q=\frac{\dot{\beta}}{\sqrt{\|\dot{\beta}\|}}$ (for $\dot{\beta}\neq 0$). This means that if we want to calculate the SRV curve, we have to calculate the derivative of β . As we never observe the whole function β but only a set discrete points $\beta(t_1),\ldots,\beta(t_m)$, this is already not straight forward, as we cannot simply calculate a pointwise derivative. However, following STEYER, A. STÖCKER, and

Greven 2021, we treat a discretely observed curve β as piecewise linear between its observed corners $\beta(t_1), \ldots, \beta(t_m)$, which allows us to calculate a piecewise constant derivative on the intervalls $[t_j, t_{j+1}], j = 1, \ldots, m-1$. Let us first consider the case of unwarped observations.

Initial parameterization Usually only the image $\beta(t_1), \ldots, \beta(t_m)$, but not the parametrisation t_1, \ldots, t_m , is observed. Therefore it is first necessary to construct an initial parameterisation. A common choice is the *arc-length-parametrisation*, where we set $t_j = \frac{l_j}{l}$ with $l_j = \sum_{k=1}^{j-1} |\beta(t_{k+1}) - \beta(t_k)|$ the polygon-length up to point j for $j \leq 2$, $l_1 = 0$ and $l_m = l$.

Piecewise-constant SRV curve Consider the discrete derivative $\Delta \beta \big|_{[t_j,t_{j+1}]} = \frac{\beta(t_{j+1})-\beta(t_j)}{t_{j+1}-t_j}$, which assumes that β is linear between its observed corners. The corresponding SRV curve q can then be treated as piecewise constant $q\big|_{[t_j,t_{j+1}]} = q_j$ with $q_j = \Delta \beta \big|_{[t_j,t_{j+1}]} / \sqrt{\|\Delta \beta \big|_{[t_j,t_{j+1}]}\|} = \frac{\beta(t_{j+1})-\beta(t_j)}{\sqrt{t_{j+1}-t_j}\cdot\sqrt{\|\beta(t_{j+1})-\beta(t_j)\|}}$ the discrete square-root-velocity of β between the corners $\beta(t_j)$ and $\beta(t_{j+1})$.

Approximate discrete SRV curve As shown in Steyer, A. Stöcker, and Greven 2021 (cf. Fig. 3), treating the SRV curves as piecewise-constant functions can lead to overfitting, where the mean shape is estimated too polygon-like. As an alternative they propose to approximate the derivative, by assuming that it attains the value of the discrete derivative $\Delta \beta \big|_{[t_j,t_{j+1}]}$ at the center $s_j = \frac{t_{j+1}-t_j}{2}$ of the interval $[t_j,t_{j+1}]$. Using this, we can construct "approximate observations" $q(s_j) \approx q_j$ of the SRV curve q.

Normalization We can approximate the normalized SRV curve z using the polygon-length l of β by $z_i = q_i/\sqrt{l}$.

[TODO: Eventuell den ganzen Part mit den piecewise-constant curves streichen. Am ende wird sowieso nur die Mittelwert-Approximation benutzt.]

[TODO: Nochmal überlegen mit γ^{-1} und γ . Gerade ist das hier glaube ich falsch.] Let us now consider a warping function γ . The warped discrete derivative is given by $\Delta(\beta\circ\gamma)\big|_{[\gamma(t_j),\gamma(t_{j+1})]}=\frac{\beta(\gamma(t_{j+1}))-\beta(\gamma(t_j))}{\gamma(t_{j+1})-\gamma(t_j)}$. The corresponding warped SRV curve is then given by $(q\circ\gamma)\sqrt{\dot{\gamma}}\big|_{[\gamma(t_j),\gamma(t_{j+1})]}=\frac{\beta(\gamma(t_{j+1}))-\beta(\gamma(t_j))}{\sqrt{\gamma(t_{j+1})-\gamma(t_j)}}.$





fig:3-disc

Figure 3.1.: A sparsely observed planar curve (left) with sparse approximate SRV curve (right). Data: see Figure 1.1.

[TODO: Das hier zuende formulieren.]

3.2. Efficient Estimation using Hermitian Covariance Smoothing

sec:cov

We want to estimate $C(s,t) = \mathbb{E}[z(s)\overline{z(t)}]$ given approximate observations of the normalized SRV curves $z_i(s_{ij})$, with $j=1,\ldots,m_i-1$ and $i=1,\ldots,N$, where m_i denotes the number of observed points per curve. Following **[TODO: Cite]**, we can treat this estimation as a smoothing problem, by constructing responses $y_{ijk} = z_i(s_{ij})\overline{z_i(s_{ik})}$ and treating the pairs s_{ij} , s_{ik} as covariates s and t. Smoothing the responses y_{ijk} gives an estimate $\hat{C}(\cdot,\cdot)$ of C(s,t), as each response has expectation $\mathbb{E}[y_{ijk}] = C(s_{ij},s_{ik})$. A popular approach **[TODO: Cite]** is to carry out the smoothing in a *tensor product spline* basis

$$C(s,t) = b(s)^T \Xi b(t)$$

where $b(s) = (b_1(s), ..., b_K(s))$ denotes the vector of a spline basis and Ξ is a $K \times K$ coefficient matrix to be estimated. As C(s,t) is complex, we restrict the spline basis to be real-valued with $b_k : \mathbb{R} \to \mathbb{R}$ for k = 1, ..., K and the coefficient matrix to be complex-valued with $\Xi \in \mathbb{C}^{K \times K}$, without loss of generality.

Considering the symmetry properties of the covariance surface allows for more efficient estimation, as shown in Cederbaum, Scheipl, and Greven 2018 for real valued, symmetric covariance surfaces, by considering every unique pair s_{ij} , s_{ik} only once. In the complex case, the covariance surface is hermitian with $C(s,t) = \overline{C(t,s)}$, which means we can decompose the estimation into two separate regression problems

over the symmetric real and skew-symmetric imaginary parts of C(s, t).

$$\mathbb{E}[\Re(y)] = b(s)^T \Xi_{\Re} b(t)$$

$$\mathbb{E}[\Im(y)] = b(s)^T \Xi_{\Im} b(t)$$

with $\Xi_{\Re}, \Xi_{\Im} \in \mathbb{R}^{K \times K}$ and $\Xi = \Xi_{\Re} + i\Xi_{\Im}$, under the constraints that $\Xi_{\Re}^T = \Xi_{\Re}$ and $\Xi_{\Im}^T = -\Xi_{\Im}$. In this thesis I estimate Ξ_{\Re} and Ξ_{\Im} using the R (**Rcore**) package mgcv (Wood 2017). For efficient estimation, two mgcv smooths from the package sparseFLMM (Cederbaum, Volkmann, and A. Stöcker 2021) are used, which implement and generalize the approach proposed by Cederbaum, Scheipl, and Greven 2018 for symmetric and skew-symmetric tensor product p-splines. [**TODO: Erklärung p-splines?**] [**TODO: Motivate b-spline/p-spline basis using Lisa's paper?**] [**TODO: REML?**]

3.3. Estimating the Elastic Full Procrustes Mean in a Fixed Basis

To estimate the elastic full Procrustes Mean, we have to solve a functional eigenvalue problem on the estimated covariance surface $\hat{C}(s,t) = b(s)^t \hat{\Xi}b(t)$. This may be achieved by evaluating $\hat{C}(s,t)$ on a dense grid and performing an eigendecomposition on the matrix of evaluations [**TODO: Cite**]. Alternatively, we can estimate the mean directly in some basis $b(s) = (b_1(s), \ldots, b_K(s))$, where a natural choice might be to evaluate mean and covariance surface in the same basis, i.e. $\mu_q(s) = b(s)^T \theta$.

Remember that the elastic full Procrustes mean (for fixed warping) is given by the solution to the optimization problem

$$\hat{\mu}_q = \underset{\mu_q \in \mathbb{L}^2: \|\mu_q\|=1}{\operatorname{argmax}} \int_0^1 \int_0^1 \overline{\mu_q(s)} C(s, t) \mu_q(t) \, ds \, dt .$$

Given an estimate of the covariance surface $\hat{C}(s,t) = b(s)^T \hat{\Xi}b(t)$, the mean estimation then reduces to estimating the vector of coefficients $\theta = (\theta_1, \dots, \theta_K) \in \mathbb{C}^K$ with

$$\hat{\theta} = \underset{\theta \in \mathbb{C}^K: \|b^T \theta\| = 1}{\operatorname{argmax}} \int_0^1 \int_0^1 \theta^H b(s) b(s)^T \hat{\Xi} b(t) b(t)^T \theta \, ds \, dt$$

nean_basis

$$\begin{split} &= \underset{\theta \in \mathbb{C}^K: \, \|b^T\theta\| = 1}{\operatorname{argmax}} \ \theta^H \left(\int_0^1 b(s)b(s)^T \, ds \right) \hat{\Xi} \left(\int_0^1 b(t)b(t)^T \, dt \right) \theta \\ &= \underset{\theta \in \mathbb{C}^K: \, \theta^H G\theta = 1}{\operatorname{argmax}} \ \theta^H G \hat{\Xi} G \theta \end{split}$$

where $(\cdot)^H = \overline{(\cdot)}^T$ denotes the conjugate transpose and G is the $K \times K$ Gram matrix with entries given by the basis products $g_{ij} = \langle b_i, b_j \rangle$. In the special case of an orthonormal basis with $\langle b_i, b_j \rangle = \delta_{ij}$ the Gram matrix is an identity matrix, however, this is not the case for many basis representations such as the b-spline basis.

We have reduced the functional eigenvalue problem to a multivariate eigenvalue problem over the covariance coefficient matrix. We might solve this using Lagrange optimization with the following Langrangian:

$$\mathcal{L}(\theta, \lambda) = \theta^H G \hat{\Xi} G \theta - \lambda (\theta^H G \theta - 1)$$

Taking into account that we identified \mathbb{R}^2 with \mathbb{C} we can split everything into real and imaginary parts and optimize with respect to $\Re(\theta)$ and $\Im(\theta)$ seperately, to avoid having to take complex derivatives. Using $\theta = \theta_{\Re} + i\theta_{\Im}$ and $\hat{\Xi} = \hat{\Xi}_{\Re} + i\hat{\Xi}_{\Im}$ we can write

$$\begin{split} \mathcal{L}(\theta_{\Re},\theta_{\Im},\lambda) &= (\theta_{\Re}^T - i\theta_{\Im}^T)G(\hat{\Xi}_{\Re} + i\hat{\Xi}_{\Im})G(\theta_{\Re} + i\theta_{\Im}) - \lambda \left((\theta_{\Re}^T - i\theta_{\Im}^T)G(\theta_{\Re} + i\theta_{\Im}) - 1\right) \\ &= \theta_{\Re}^TG\hat{\Xi}_{\Re}G\theta_{\Re} + i\theta_{\Re}^TG\hat{\Xi}_{\Im}G\theta_{\Re} + \theta_{\Im}^TG\hat{\Xi}_{\Im}G\theta_{\Re} - \theta_{\Re}^TG\hat{\Xi}_{\Im}G\theta_{\Im} \\ &+ \theta_{\Im}^TG\hat{\Xi}_{\Re}G\theta_{\Im} + i\theta_{\Im}^TG\hat{\Xi}_{\Im}G\theta_{\Im} + \lambda \left(\theta_{\Re}^TG\theta_{\Re} + \theta_{\Im}^TG\theta_{\Im} - 1\right) \end{split}$$

using $\hat{\Xi}_{\Re}^T = \hat{\Xi}_{\Re}$ and $\hat{\Xi}_{\Im}^T = -\hat{\Xi}_{\Im}$. Differentiation w.r.t. θ_{\Re} and θ_{\Im} yields

$$\frac{\partial \mathcal{L}}{\partial \theta_{\Re}} = 2G \hat{\Xi}_{\Re} G \theta_{\Re} - 2G \hat{\Xi}_{\Im} G \theta_{\Im} - 2\lambda G \theta_{\Re} \stackrel{!}{=} 0 \tag{3.1}$$
 [eq:lagrRe]

eq:lagrIm

$$\frac{\partial \mathcal{L}}{\partial \theta_{\Im}} = 2G \hat{\Xi}_{\Re} G \theta_{\Im} + 2G \hat{\Xi}_{\Im} G \theta_{\Re} - 2\lambda G \theta_{\Im} \stackrel{!}{=} 0 \tag{3.2}$$

with the additional constraint $\theta_{\Re}^T G \theta_{\Re} + \theta_{\Im}^T G \theta_{\Im} = 1$. TWe can simplify this further and multiply Eq. 3.2 by i, leading to

$$\hat{\Xi}_{\Re}G\theta_{\Re} - \hat{\Xi}_{\Im}G\theta_{\Im} = \lambda\theta_{\Re} \tag{3.3}$$

$$i\hat{\Xi}_{\Re}G\theta_{\Im} + i\hat{\Xi}_{\Im}G\theta_{\Re} = i\lambda\theta_{\Im}. \tag{3.4}$$

Adding both equations finally leads to

$$(\hat{\Xi}_{\Re} + i\hat{\Xi}_{\Im})G\theta_{\Re} + i(\hat{\Xi}_{\Re} + \hat{\Xi}_{\Im})G\theta_{\Im} = \lambda(\theta_{\Re} + i\theta_{\Im})$$

or likewise, using θ and $\hat{\Xi}$

$$\hat{\Xi}G\theta = \lambda\theta$$

which is an eigenvalue problem on the product of the complex coefficient matrix and the Gram matrix. Multiplying by $\theta^H G$ from the left yields $\lambda = \theta^H G \hat{\Xi} G \theta$, i.e. the eigenvalues correspond to the target function to maximize. It follows that the estimate for the coefficient vector of the elastic full Procrustes mean is given by the eigenvector of the leading eigenvalue of $\hat{\Xi} G$. [TODO: Cite Reiss.]

3.4. Numerical Integration of the Procrustes Fits

sec:pfit

[TODO: Mean value theorem in the integral, etc.]

 $\hat{p}(t) = b(t)^T \hat{\theta}$ estimated Procrustes mean function, q piecewise constant SRV transform with $q|_{[t_j,t_{j+1}]} = q_j \in \mathbb{C}$.

For estimation of the Procrustes fits we need to estimate two scalar products:

$$\hat{q}_p = \frac{\langle q, \hat{p} \rangle}{\langle q, q \rangle} q$$

So far I treat *q* as piecewise constant (as in Lisa's paper):

$$\langle q, \hat{p} \rangle = \int_0^1 \langle q(t), \hat{p}(t) \rangle dt \approx \sum_{j=0}^{m-1} \int_{t_j}^{t_{j+1}} \langle q_j, \hat{p}(t) \rangle dt$$

$$\langle q, q \rangle = \int_0^1 \langle q(t), q(t) \rangle dt \approx \sum_{j=0}^{m-1} (t_{j+1} - t_j) \langle q_j, q_j \rangle$$

4. Empirical Applications

sec:app

- 4.1. Mean Estimation for Simulated Spirals
- 4.2. Classification of Hand-written Digits
- 4.3. Mean Differences of Tounge Shapes in a Phonetics Study

5. Summary

sec:sum

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A. Appendix

app:app

A.1. Proofs and Derivations

-proc-dist

A.1.1. Derivation of Lemma 2.2

[TODO: Make dependence on γ explicit!]

Proof. Start with ii.). Let $\tilde{z}_2 = (\gamma, z_2) = (z_2 \circ \gamma) \sqrt{\dot{\gamma}}$, then optimise $d_{EF}([\beta_1], [\beta_2])^2$ over rotation and scaling, keeping γ fixed.

$$\begin{split} d_{EF}([\beta_{1}], [\beta_{2}])^{2} &= \inf_{\lambda \in \mathbb{R}^{+}, \theta \in [0, 2\pi], \gamma \in \Gamma} \|z_{1} - \lambda e^{i\theta} \widetilde{z}_{2}\|^{2} \\ &= \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^{+}, \theta \in [0, 2\pi]} \langle z_{1} - \lambda e^{i\theta} \widetilde{z}_{2}, z_{1} - \lambda e^{i\theta} \widetilde{z}_{2} \rangle \right) \\ &= \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^{+}, \theta \in [0, 2\pi]} \|z_{1}\|^{2} + \lambda^{2} \|\widetilde{z}_{2}\|^{2} - \lambda (e^{i\theta} \langle z_{1}, \widetilde{z}_{2} \rangle + e^{-i\theta} \langle \widetilde{z}_{2}, z_{1} \rangle) \right) \end{split}$$

As $\langle z_1, \widetilde{z}_2 \rangle \in \mathbb{C}$, define $\langle z_1, \widetilde{z}_2 \rangle = \kappa_{\gamma} e^{i\phi_{\gamma}}$, where $(\cdot)_{\gamma}$ denotes the dependence on γ . Furthermore, use $||z_{1,2}|| = 1$, which implies $||\widetilde{z}_2|| = 1$, as re-parametrisation is norm-preserving, when using the elastic metric.

$$\begin{split} \dots &= \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^+, \theta \in [0, 2\pi]} 1 + \lambda^2 - \lambda (e^{i\theta} \kappa_{\gamma} e^{i\phi_{\gamma}} + e^{-i\theta} \kappa_{\gamma} e^{-i\phi_{\gamma}}) \right) \\ &= \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^+, \theta \in [0, 2\pi]} 1 + \lambda^2 - \lambda \kappa_{\gamma} \left(e^{i(\theta + \phi_{\gamma})} + e^{-i(\theta + \phi_{\gamma})} \right) \right) \\ &= \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^+} 1 + \lambda^2 - \sup_{\theta \in [0, 2\pi]} 2\lambda \kappa_{\gamma} \cos\left(\theta + \phi_{\gamma}\right) \right) \\ &\stackrel{\theta = -\phi_{\gamma}}{=} \inf_{\gamma \in \Gamma} \left(\inf_{\lambda \in \mathbb{R}^+} 1 + \lambda^2 - 2\lambda \kappa_{\gamma} \right) \end{split}$$

From $\frac{\partial}{\partial \lambda} \left(1 + \lambda^2 - 2\lambda \kappa_{\gamma} \right) = 2\lambda - 2\kappa_{\gamma} \stackrel{!}{=} 0$ it follows that $\lambda = \kappa_{\gamma}$.

$$\ldots = \inf_{\gamma \in \Gamma} \left(1 + \kappa_{\gamma}^2 - 2\kappa_{\gamma}^2 \right) = \inf_{\gamma \in \Gamma} \left(1 - \kappa_{\gamma}^2 \right)$$

[TODO: Make this part clearer!] Lemma 2.2 ii.) follows from
$$\kappa_{\gamma}^2 = |\langle z_1, \widetilde{z}_2 \rangle|^2 = \langle z_1, \widetilde{z}_2 \rangle \langle \widetilde{z}_2, z_1 \rangle$$
 and $\widetilde{z}_2 = (z_2 \circ \gamma) \sqrt{\dot{\gamma}}$. Lemma 2.2 i.) follows from $\lambda e^{i\theta} = \kappa_{\gamma} e^{-i\phi_{\gamma}} = \overline{\langle z_1, \widetilde{z}_2 \rangle}$.

app:closed

A.2. Discussion of Possible Extensions to Closed Curves

app:smooth

A.3. Shape-Smoothing Using the Estimated Covariance-Surface

Idea: Smooth q in mean basis b(t), so that $\hat{q}(t) = b(t)^T \hat{\theta}_q$. Then the scalar products simplify to

$$\langle q, \hat{p} \rangle \approx \hat{\theta}_q^H G \hat{\theta}$$

$$\langle q,q\rangle \approx \hat{\theta}_q^H G \hat{\theta}_q$$

As q can be sparse, we may want to use the estimated covariance matrix $\hat{\Xi}$ when estimating θ_q in the form of a Normal prior $\theta_q \sim \mathcal{N}_{\mathbb{C}^k}(0,\hat{\Xi})$ leading to

$$\hat{\theta}_q = (B^T B + \Xi^{-1})^{-1} B^T q$$

or equivalently using eigendecompositon $\Xi = V\Lambda V^{-1}$ with $V^HV = \mathcal{I}$ and $\Lambda = diag(\lambda_1, \ldots, \lambda_k)$

$$\hat{\theta}_q = V(V^H B^T B V + \Lambda^{-1})^{-1} V^H B^T q$$

In general how well this works tends to depend on how close the curve is to the estimated mean.

Maybe using a penalty parameter - controling the strength of the regularization - is also thinkable? Something like this:

$$\hat{\theta}_q(\lambda) = V(V^H B^T B V + \lambda \cdot \Lambda^{-1})^{-1} V^H B^T q$$

Thought: A normal prior is probably not appropriate, because while its true that $\mathbb{E}[\theta] = 0$ (due to rotational symmetry) I would think that in general $\mathbb{E}[\|\theta\|] \neq 0$. A distribution more like a Normal distribution lying on a ring around 0 with radius $r = \|\theta\|$ is probably better? Then estimates wouldn't be pressed to zero so strongly, and instead would be pressed to curve of size equal to the mean.

B. Supplementary Materials

app:sup

- **B.1.** Dataset Replication Guide
- **B.2.** Implementation Notes

—Discarded—

Math-Basics Recap

Scalar Products

V n-dimensional vector space with basis $B = (b_1, \ldots, b_n)$, then any scalar product $\langle \cdot, \cdot \rangle$ on V can be expressed using a $(n \times n)$ matrix G, the Gram matrix of the scalar product. Its entries are the scalar products of the basis vectors:

$$G = (g_{ij})_{i,j=1,...,n}$$
 with $g_{ij} = \langle b_i, b_j \rangle$ for $i, j = 1,...,n$

When vectors $x, y \in V$ are expressed with respect to the basis B as

$$x = \sum_{i=1}^{n} x_i b_i \quad \text{and} \quad y = \sum_{i=1}^{n} y_i b_i$$

the scalar product can be expressed using the Gram matrix, and in the complex case it holds that

$$\langle x,y\rangle = \sum_{i,j=1}^{n} \bar{x}_i y_j \langle b_i, b_j \rangle = \sum_{i,j=1}^{n} \bar{x}_i g_{ij} y_j = x^{\dagger} G y$$

when $x_i, y_i \in \mathbb{C}$ for i = 1, ..., n with x^{\dagger} indicating the conjugate transpose of $x = (x_1, ..., x_n)^T$. If B is an *orthonormal* basis, that is if $\langle b_i, b_j \rangle = \delta_{ij}$, it further holds that $\langle x, y \rangle = x^{\dagger}y$ as $G = \mathbb{1}_{n \times n}$.

Functional Scalar Products

This concept can be generalized for vectors in function spaces. Define the scalar product of two functions f(t), g(t) as:

$$\langle f, g \rangle = \int_a^b \bar{f}(t) w(t) g(t) dt$$

with weighting function w(t) and [a,b] depending on the function space. The scalar product has the following properties:

1.
$$\langle f, g + h \rangle = \langle f, g \rangle + \langle f, h \rangle$$

2.
$$\langle f, g \rangle = \overline{\langle g, f \rangle}$$

3.
$$\langle f, cg \rangle = c \langle f, g \rangle$$
 or, using (2), $\langle cf, g \rangle = \bar{c} \langle f, g \rangle$ for $c \in \mathbb{C}$

If we have a functional basis $\{\phi_1, \dots, \phi_n\}$ (and possibly $n \to \infty$) of our function space we can also write the function f as an expansion

$$f = \sum_{i=1}^{n} a_i \phi_i$$
 so that $f(t) = \sum_{i=1}^{n} a_i \phi_i(t)$

Additionally, if we have a *orthogonal* basis, so that $\langle \phi_i, \phi_j \rangle = 0$ for $i \neq j$, we can take the scalar product with ϕ_k from the left

$$\langle \phi_k, f \rangle = \sum_{i=1}^n a_i \langle \phi_k, \phi_i \rangle = a_k \langle \phi_k, \phi_k \rangle$$

which yields the coefficients a_k :

$$a_k = \frac{\langle \phi_k, f \rangle}{\langle \phi_k, \phi_k \rangle}$$

For an *orthonormal* basis it holds that $\langle \phi_i, \phi_j \rangle = \delta_{ij}$. Suppose that two functions f, g are expanded in the same orthonormal basis:

$$f = \sum_{i=1}^{n} a_i \phi_i$$
 and $g = \sum_{i=1}^{n} b_i \phi_i$

We can then write the scalar product as:

$$\langle f, g \rangle = \langle \sum_{i=1}^n a_i \phi_i, \sum_{i=1}^n b_i \phi_i \rangle = \sum_{i=1}^n \sum_{j=1}^n \hat{a}_i b_j \langle \phi_i, \phi_j \rangle = \sum_{i=1}^n \bar{a}_i b_i = a^{\dagger} b$$

for coefficient vectors $a, b \in \mathbb{C}^n$. This means that the functional scalar product reduces to a complex dot product. Additionally it holds that for the norm $||\cdot||$ of a function f:

$$||f|| = \langle f, f \rangle^{\frac{1}{2}} = \sqrt{a^{\dagger}a} = \sqrt{\sum_{i=1}^{n} |a_i|^2}$$

FDA-Basics Recap

As discussed in the last section we can express a function f in its basis function expansion using a set of basis functions ϕ_k with k = 1, ..., K and a set of coefficients $c_1, ..., c_K$ (both possibly $\mathbb C$ valued e.g. in the case of 2D-curves)

$$f = \sum_{k=1}^{K} c_k \phi_k = c' \phi$$

where in the matrix notation c and ϕ are the vectors containing the coefficients and basis functions.

When considering a sample of N functions f_i we can write this in matrix notation as

$$f = C\phi$$

where *C* is a $(N \times K)$ matrix of coefficients and *f* is a vector containing the *N* functions.

Smoothing by Regression

When working with functional data we can usually never observe a function f directly and instead only observe discrete points (x_i, t_i) along the curve, with $f(t_i) = x_i$. As we don't know the exact functional form of f, calculating the scalar products $\langle \phi_k, f \rangle$ and therefore calculating the coefficients c_k of a given basis representation is not possible.

However, we can estimate the basis coefficients using e.g. regression analysis an approach motivated by the error model

$$f(t_i) = c' \phi(t_i) + \epsilon_i$$

If we observe our function n times at t_1, \ldots, t_n , we can estimate the coefficients from a least squares problem, where we try to minimize the deviation of the basis expansion from the observed values. Using matrix notation let the vector f contains the observed values $f(t_i)$, $i = 1, \ldots, n$ and $(n \times k)$ matrix $\mathbf{\Phi}$ contains the basis function values $\phi_k(t_i)$. Then we have

$$f = \Phi c + \epsilon$$

with the estimate for the coefficient vector *c* given by

$$\hat{c} = (\mathbf{\Phi'\Phi})^{-1} \mathbf{\Phi'} f.$$

Spline curves fit in this way are often called *regression splines*.

Common Basis Representations

Piecewise Polynomials (Splines) Splines are defined by their range of validity, the knots, and the order. Their are constructed by dividing the area of observation into subintervals with boundaries at points called *breaks*. Over any subinterval the spline function is a polynomial of fixed degree or order. The term *degree* refers to the highest power in the polynomial while its *order* is one higher than its degree. E.g. a line has degree one but order two because it also has a constant term. [...]

Polygonal Basis [...]

Bivariate Functional Data

The analogue of covariance matrices in MVA are covariance surfaces $\sigma(s,t)$ whose values specify the covariance between values f(s) and f(t) over a population of curves. We can write these bivariate functions in a *bivariate basis expansion*

$$r(s,t) = \sum_{k=1}^{K} \sum_{l=1}^{K} b_{k,l} \phi_k(s) \psi_l(t) = \boldsymbol{\phi}(s)' \boldsymbol{B} \boldsymbol{\psi}(t)$$

with a $K \times K$ coefficient matrix B and two sets of basis functions ϕ_k and ψ_l using *Tensor Product Splines*

$$B_{k,l}(s,t) = \phi_k(s)\psi_l(t).$$