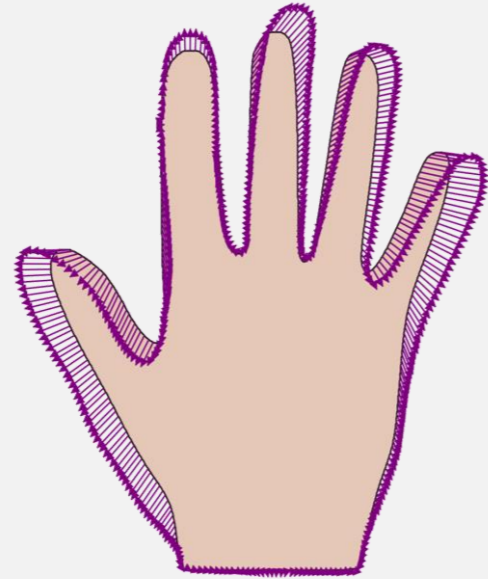


**University
of Basel**

Finite rank representation of a GP (advanced topic)

Finite observations revisited

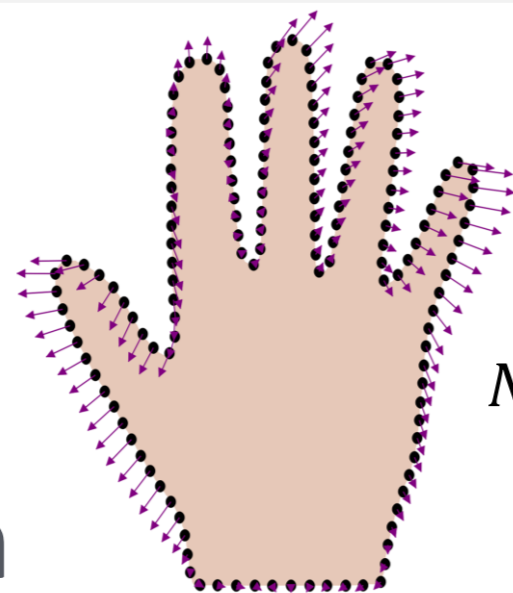
Continuous representation



$GP(\mu, k)$

Computer implementation

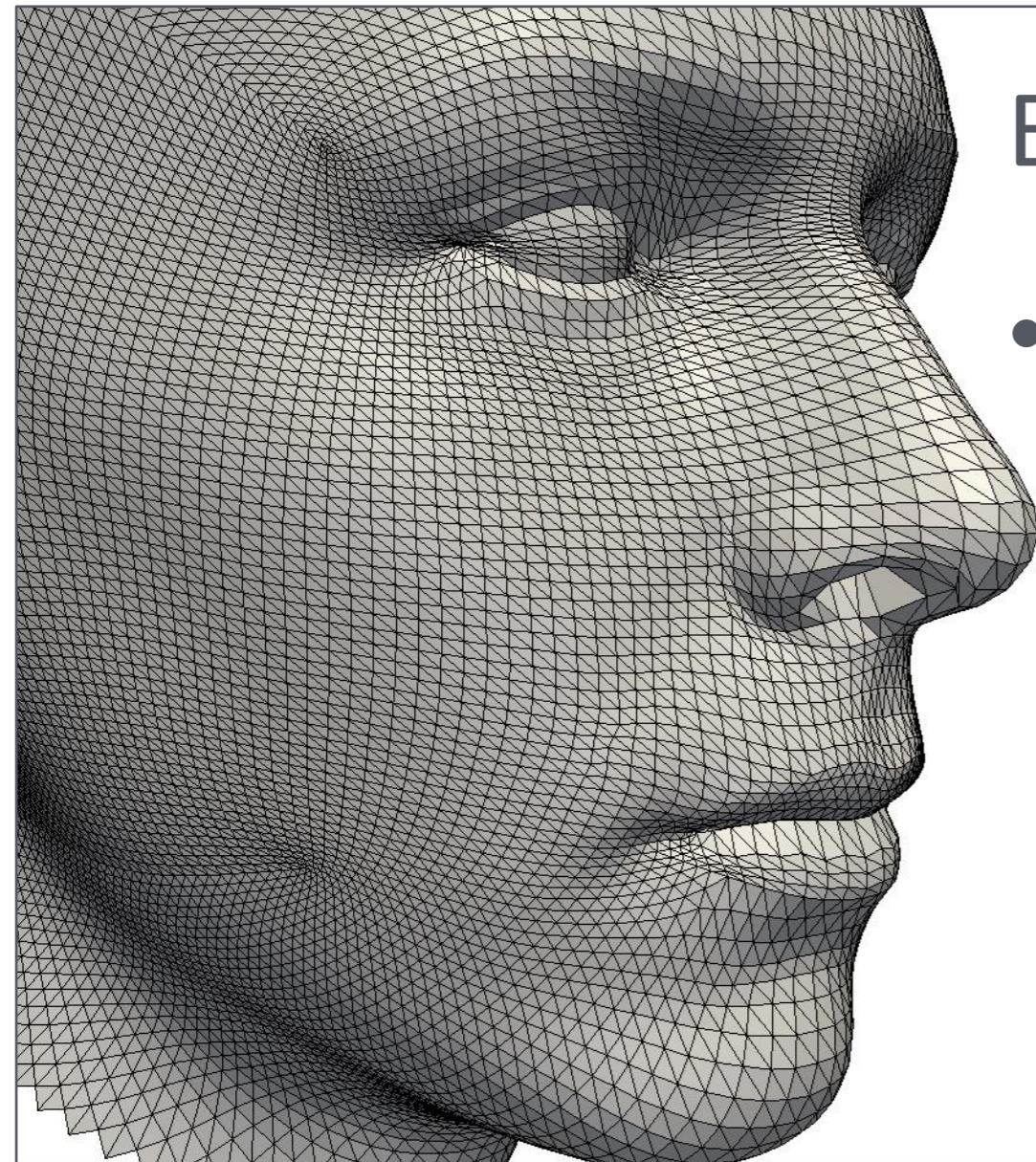
Discrete representation



$$N \left(\begin{pmatrix} \mu_1(x_1) \\ \mu_2(x_1) \\ \vdots \\ \mu_1(x_n) \\ \mu_2(x_n) \end{pmatrix}, \begin{pmatrix} k_{11}(x_1, x_1) & k_{12}(x_1, x_1) & \dots & k_{11}(x_1, x_n) & k_{12}(x_1, x_n) \\ k_{21}(x_1, x_1) & k_{22}(x_1, x_1) & \dots & k_{21}(x_1, x_n) & k_{22}(x_1, x_n) \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ k_{11}(x_n, x_1) & k_{12}(x_n, x_1) & \dots & k_{11}(x_n, x_n) & k_{12}(x_n, x_n) \\ k_{21}(x_n, x_1) & k_{22}(x_n, x_1) & \dots & k_{21}(x_n, x_n) & k_{22}(x_n, x_n) \end{pmatrix} \right)$$

A problem with GPs

- The resulting covariance matrix can be large



Example: Basel face model

- Mesh contains 53490 vertices

$$N \left(\begin{pmatrix} \mu(x_1) \\ \vdots \\ \mu(x_n) \end{pmatrix}, \underbrace{\begin{pmatrix} k(x_1, x_1) & \cdots & k(x_1, x_n) \\ \vdots & \ddots & \vdots \\ k(x_n, x_1) & \cdots & k(x_n, x_n) \end{pmatrix}}_{(53490 \cdot 3) \times (53490 \cdot 3)} \right)$$

The Karhunen-Loève expansion

We can write

$$u \sim GP(\mu, k)$$

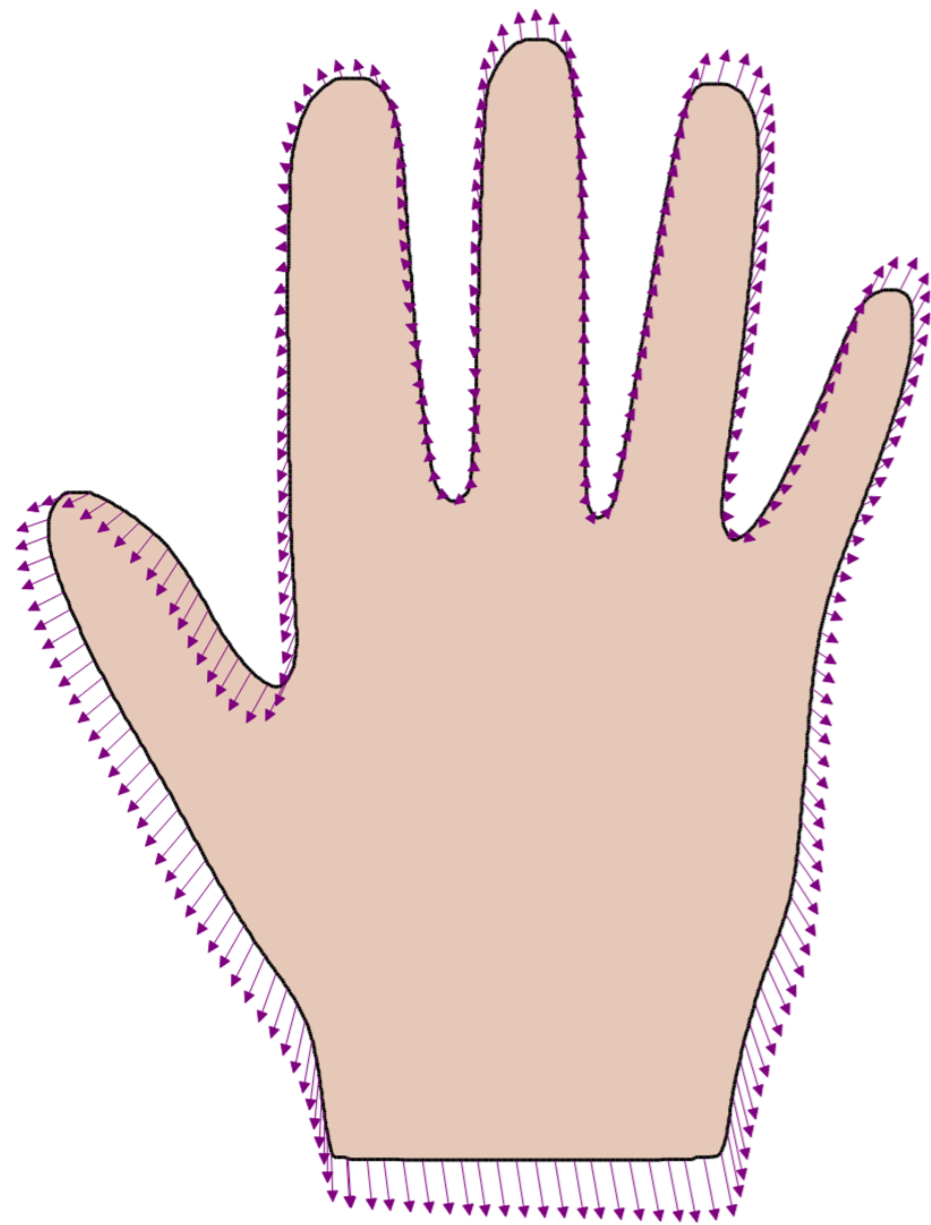
as

$$u \sim \mu + \sum_{i=1}^{\infty} \alpha_i \sqrt{\lambda_i} \phi_i, \quad \alpha_i \sim N(0, 1)$$

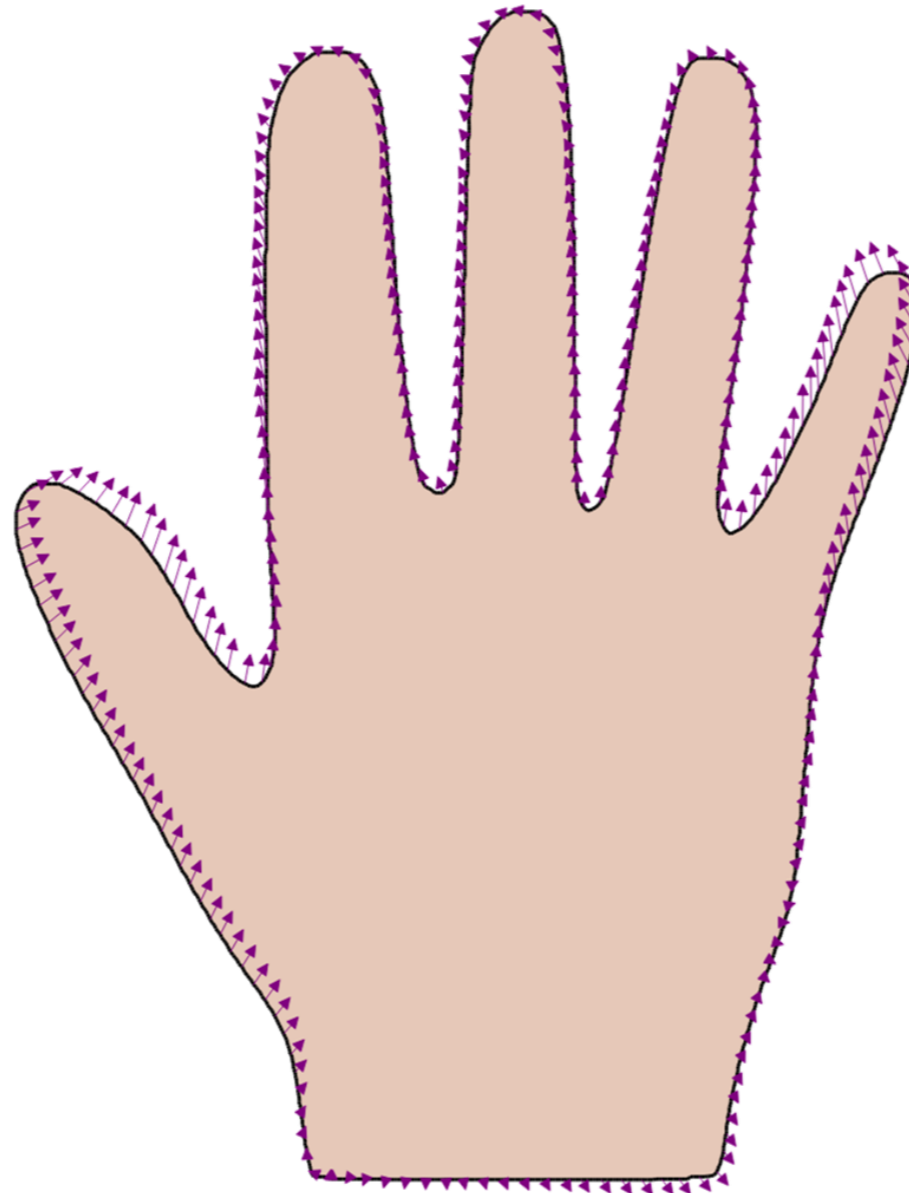
- ϕ_i is the eigenfunction with associated eigenvalue λ_i of the linear operator

$$[T_k u](x) = \int k(x, s) u(s) ds$$

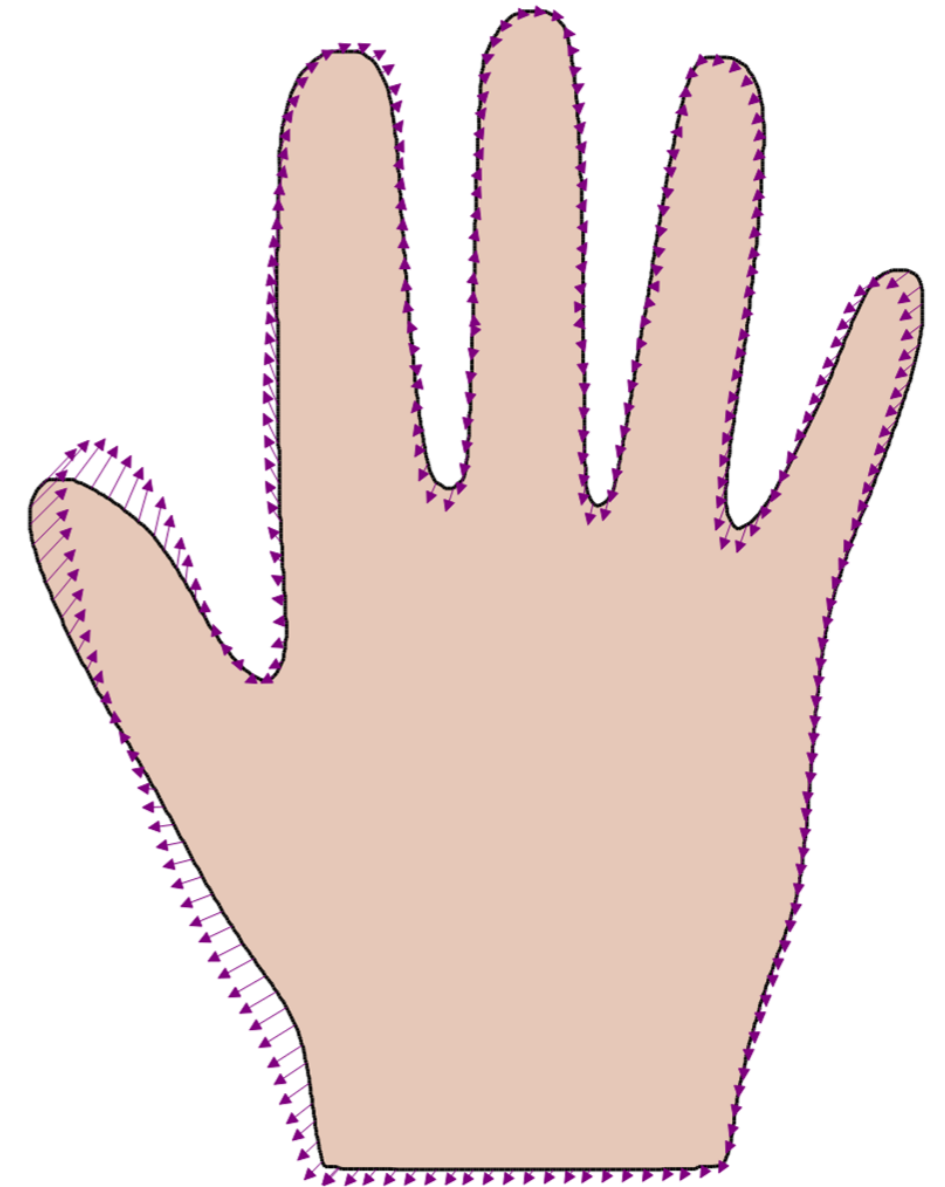
Eigenfunctions of the hand model



$$\sqrt{\lambda_1}\phi_1$$

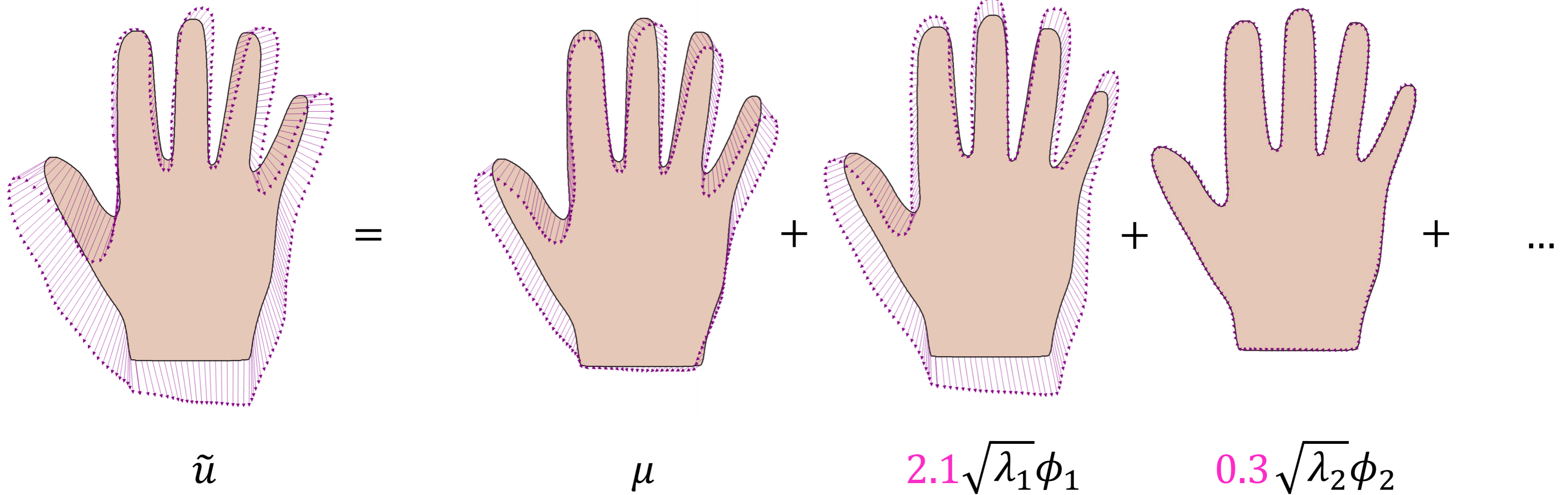


$$\sqrt{\lambda_2}\phi_2$$

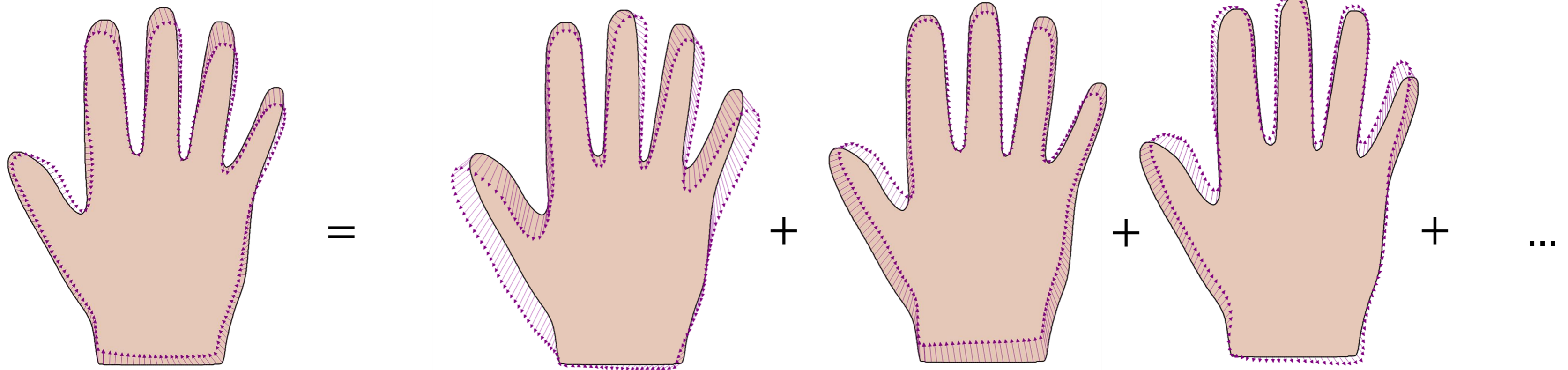


$$\sqrt{\lambda_3}\phi_3$$

Karhunen-Loève expansion



Karhunen-Loève expansion



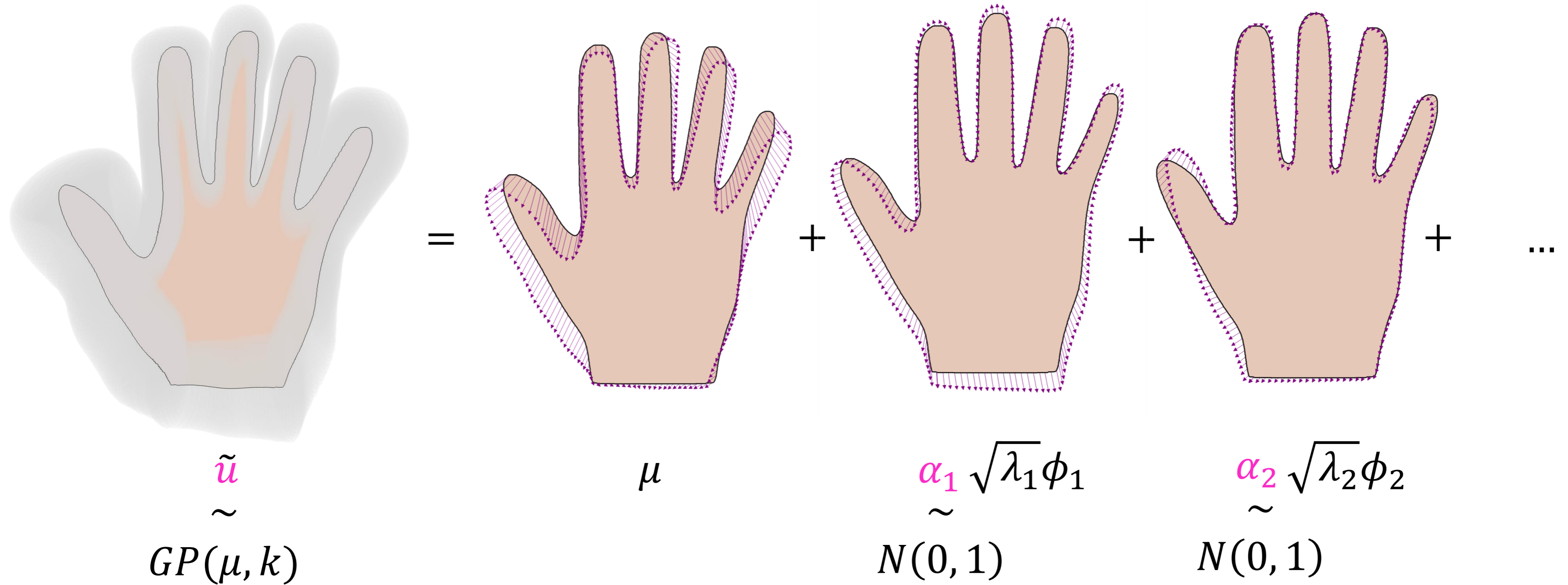
\tilde{u}

μ

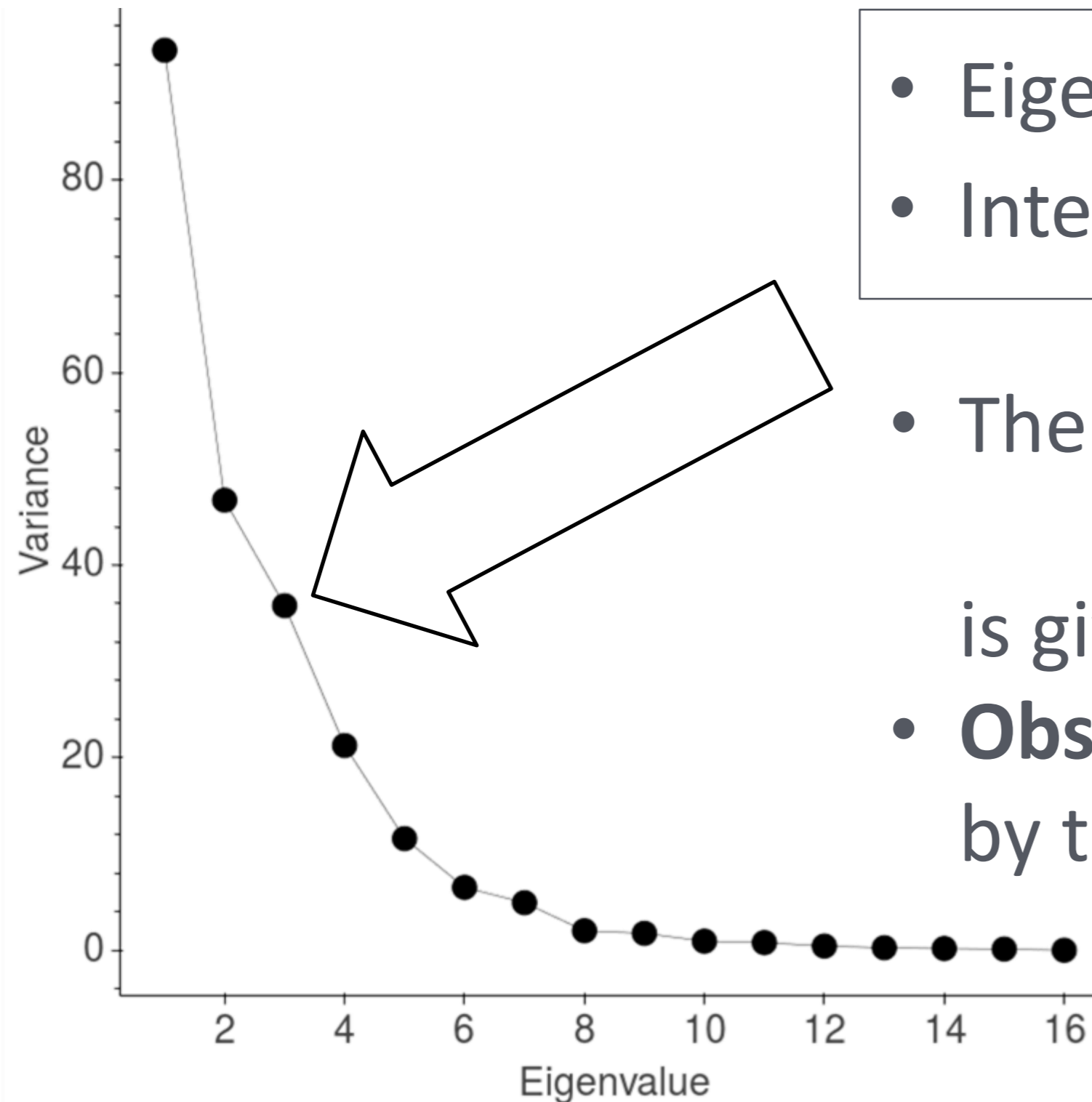
$-1.2\sqrt{\lambda_1}\phi_1$

$1.8\sqrt{\lambda_2}\phi_2$

Karhunen-Loève expansion



Eigenvalues and variance



- Eigenvalue λ_i
- Interpretation: Variance of $\alpha_i \sqrt{\lambda_i} \phi_i$
- The total variance of the process $u \sim GP(\mu, k)$ is given by $\sum_{i=1}^{\infty} \lambda_i$.
- **Observation:** Most variance is explained by the first eigenfunctions

Low-rank approximation

$$u = \mu + \sum_{i=1}^r \alpha_i \sqrt{\lambda_i} \phi_i, \quad \alpha_i \sim N(0, 1)$$

Main idea: Represent process using only the first r components

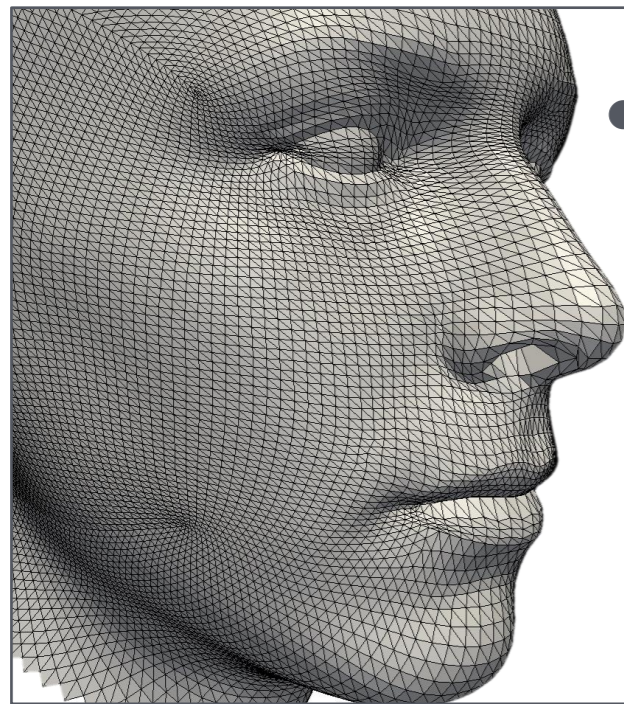
- We have a **finite**, parametric representation of the process.
- Any deformation u is determined by the coefficients

$$\alpha = (\alpha_1, \dots, \alpha_r)$$

$$p(u) = p(\alpha) = \prod_{i=1}^r \frac{1}{\sqrt{2\pi}} \exp(-\alpha_i^2 / 2)$$

Computational aspects

- Numerical methods are used for computing (λ_i, ϕ_i)
 - Very efficient for models that are learned from examples
 - Possible for general models
 - The full covariance matrix is never computed



- If rank r is low, the representation is efficient
 - can represent shapes with millions of points