

# Inferring Human Body Parts and Correlations from Images, Pointclouds and Meshes

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# Abstract

This thesis addresses the development of a novel sample thesis. We analyze the requirements of a general template, as it can be used with the  $\text{\LaTeX}$  text processing system. (And so on...) The abstract should not exceed half a page in size!



# Zusammenfassung

Diese Arbeit beschäftigt sich mit der Entwicklung einer neuartigen Beispielausarbeitung. Wir untersuchen die Anforderungen, die sich für eine allgemeine Vorlage ergeben, die innerhalb der  $\text{\LaTeX}$ -Textverarbeitungsumgebung verwendet werden kann. (Und so weiter und so fort...) Die Zusammenfassung sollte nicht länger als eine halbe Textseite sein!



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# Introduction

The price for breast enhancement surgery was estimated to be \$3718 in 2017 according to the American Society of Plastic Surgery<sup>1</sup>. Next to the financial aspects there are also risks connected to undergoing surgery as well as not knowing exactly what the result will look like. Before committing to this kind of operation, it should be possible to generate a preview of the outcome from a few images. This thesis aims to design a method that is able to predict a 3D model of the outcome by learning a mapping between parametric models. Additionally, the idea is explored if it is possible to generate a parametric model from a character modelling software.

## 1.1. Parametric Model

A previous implementation by Biland [Bil17] was used to create parametric models. A parametric model can describe all data that went into the model with its parameters. For example, the physical appearance of a person can be roughly described by their height, skin tone and hair color, where these three are the parameters of this parametric model. This is, of course, only an approximation as the accuracy of the description of the person would increase with more parameters. It is also possible that one parameter influences multiple features. In the previous example, when the height of a person is raised, the length of the arms are also proportionally increased.

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<sup>1</sup><https://www.plasticsurgery.org/cosmetic-procedures/breast-augmentation/cost>

## 1. Introduction

### 1.2. Mapping

A mapping between sets associates each element in the first set with one or more elements of the second set. An example for a simple mapping could be the numbers one to twenty-six as the first set and the letters of the alphabet as the second. In the case of two parametric models, the goal is to find a mapping that describes the relationship between the parameters of the first and second model.

This mapping can either be linear or non-linear. The difference between a linearity and a non-linearity can be described with a simple example. The time it takes to drive  $10km$  in a car at  $10\frac{km}{h}$  is  $1h$ . If the distance to drive is doubled to  $20km$ , so will the time it takes. That is because distance to drive and time it requires are in a linear relationship. On the other hand, given that the braking distance while travelling with  $10\frac{km}{h}$  is  $1m$ , the braking distance while travelling with  $20\frac{km}{h}$  is  $4m$ . This is due to the fact, that speed and braking distance are related in a quadratic, non-linear manner.

### 1.3. Applications

The applications for this method could be plenty. While the most obvious use for this implementation would be plastic surgery - creating 3D models from images of possible breast enhancements, it could also be used in oncology, to show affected patients what their body might look like after a mastectomy and even further, after a possible reconstruction. This approach could also be used in an adapted version for commercial use, such as tailor fitting clothes by using an image of the customer.

## Related Work

In the field of modelling the human body, early on, a lot of work was done concerning the face. Jeng et al. [JLH<sup>+</sup>98] proposed a geometric face model to localize the face in an image and detect facial features efficiently. The reason for the need of a functioning model for faces may have been the upcoming importance of facial recognition.

Parametric models have also been applied to reshaping the human body in images using semantic parameters. [ZFL<sup>+</sup>10] Their approach included reshaping a morphable 3D model and applying the reshaping effects using an image warping approach. Other work was done by Allen et al. [ACP03] on parametrizing full body scans and using the parameters in a variety of applications. One application was to find features connected to certain parameters, such that for example the weight could be increased and the resulting mesh would increase in size in a meaningful way.

Work about specifically modelling the breast has been done by Galle et al. [GGC10]. They proposed a parameter space based on MRI scans to describe the shape of a human breast using PCA. Ciechomski et al. [dHCCG<sup>+</sup>12] proposed a web approach to create 3D models from images where participants uploaded images and selected keypoints. These 3D models can later be accessed by surgeons to visualize different types and sizes of implants applied.





## Methods

In this chapter the methods needed to create a parametric model and a mapping between parametric models are introduced and explained. First off, the data gathering and preprocessing are outlined. Multiple variants of the parametric model are discussed. Different learning approaches for the mapping are reviewed. Lastly, a character editor is presented that is used to generate data for a parametric model.

### 3.1. Data Acquisition and Preprocessing

Optimally it would be necessary to have a large data set of images of women before and after breast enhancement surgery, where the patients pose topless. It is very unlikely though that such a database exists, due to the fact that having breast surgery is a very personal topic and people generally don't enjoy posing naked. Therefore the images used were downloaded from a website<sup>1</sup> that offered to simulate various plastic surgical procedures including breast enhancement. For each user a 3D model of their torso was displayed side by side with different enhancements varying in size. Each model was made up of a sequence of 24 images displaying the torso from different angles. This dataset fit the requirements nicely as images are available for *before* and *after*, except the *after* is generated and based on their model. Additionally, each after image sequence had a short label, usually describing how much silicon was added, that was also saved for further evaluation. In total 2'937 examples were retrieved and preprocessed. This dataset included images from 748 subjects of which each one was comprised of one *before* and at least one *after* image sequence.

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<sup>1</sup><https://my.crisalix.com/>

### 3. Methods

Term	R:	L:	LSC	73	cc	LSA	350	Custom	300
Frequency	205	205	100	85	54	73	73	70	60

**Table 3.1.:** The frequency of the top 9 terms can be seen in the upper row. The corresponding frequency is stated below.

In a next step these image sequences needed to be transformed into point clouds. This was done using a general-purpose Structure-from-Motion (SfM) [SF16] and Multi-View Stereo (MVS) [SZPF16] pipeline called COLMAP. This generated point clouds spanning from 5'000 to 15'000 points. Some of the images needed to be discarded, due to the fact that SfM created a point cloud with less than 1'000 points or the point clouds had holes, such that certain areas had no points and were not defined at all. The remaining point clouds were cleaned using a C++ implementation by Biland [Bil17] that removed white points around the point clouds.

The mapping required to have one set of point clouds of *before* examples and the corresponding<sup>2</sup> *after* examples. Therefore the data was split into sets of *before* and *after* point clouds. As all of the *after* point clouds were based on various breast enhancements, all the corresponding labels of given were processed and analyzed. This resulted in a table of terms sorted by the frequency of the term. The 9 most frequent terms can be found in table 3.1.

The first few terms are not informative, as they don't describe the breast themselves. The terms "LSC" and "73" are descriptions of a certain implant<sup>3</sup> and "cc" is a measure of volume. The first meaningful term is "350", which means that an implant with a volume of 350ml was inserted. Therefore only the *after* examples that were labelled "350" were included, to create a better mapping. After removing some of the point clouds that were not reconstructed well enough, this resulted in 57 examples in the *before* and 57 in the *after* set. All of these point clouds were further processed in a MATLAB implementation by Biland [Bil17] to generate mesh files.

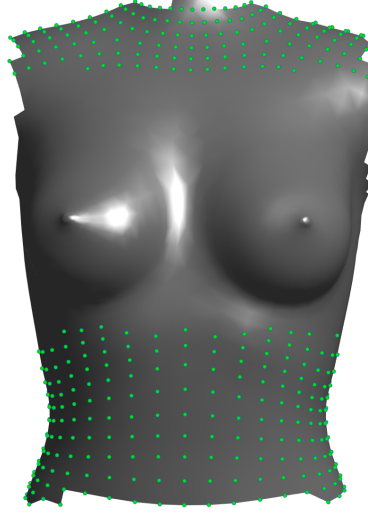
#### 3.1.1. Alignment

Eventhough the meshes were aligned in the MATLAB implementation by Biland [Bil17] in a general fashion, each corresponding *before* and *after* should be pairwise aligned to avoid that the mapping also learns rotations. This was achieved by using an implementation of Horn's method [Hor87]. Given two sets of vertices, Horn's method computes the translation, rotation and possibly scale change from one set to the other. As it is expected that for the same subject, only points defining the breasts should vary from before to after, only a subset of vertices should be used. The points used in the alignment can be seen in figure 3.1.

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<sup>2</sup>Corresponding meaning, based on the same subject.

<sup>3</sup>[http://www.simplybreastimplants.com/breast\\_implant\\_sizes/sebbin\\_firm\\_high\\_profile.html](http://www.simplybreastimplants.com/breast_implant_sizes/sebbin_firm_high_profile.html)



**Figure 3.1.:** The mesh is depicted as the grey surface. The points drawn in green are used for the alignment.

## 3.2. Parametric Model from Meshes

In this section it is described how a parametric model is obtained using principle component analysis (PCA). Given  $n$  meshes  $m_i \in \mathbb{R}^{k \times 3}$ , where  $k$  describes the amount of vertices  $m$  has, each mesh needs to be transformed to be of shape  $\mathbb{R}^{1 \times 3k}$ . Then, all transformed meshes are stacked into a matrix  $M \in \mathbb{R}^{n \times 3k}$ . As the differences over each column isn't significant, the mean  $\bar{m}$  of the matrix  $M$  is subtracted from each row of  $M$ .

$$\mathbf{A} := \begin{bmatrix} m'_1 - \bar{m} \\ m'_2 - \bar{m} \\ \vdots \\ m'_n - \bar{m} \end{bmatrix} \in \mathbb{R}^{n \times 3k} \quad (3.1)$$

Next, PCA is run with matrix  $A$  as the input. PCA is able to reduce the dimensionality of the data while retaining most of the information from the initial data set. This is achieved by finding orthogonal basis vectors, where the first basis vector is responsible for the largest variance in the data. The second basis vector needs to be orthogonal to the first and is responsible for the second largest variance of the data. This holds for each following basis vector. These basis vectors in PCA are also known as principle components.

The output of the PCA function is:

$$\mathbf{coeff} := \begin{bmatrix} c_1 & c_2 & \cdots & c_{n-1} \end{bmatrix} \in \mathbb{R}^{3k \times n-1} \quad (3.2)$$

$$\mathbf{score} := \begin{bmatrix} s_1 & s_2 & \cdots & s_{n-1} \end{bmatrix} \in \mathbb{R}^{n \times n-1} \quad (3.3)$$

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where coefficient  $c_i$  is the  $i$ -th principle component and score  $s_i$  is the  $i$ -th parameter vector corresponding to the  $i$ -th input. Therefore the input data can be reproduced by computing  $\mathbf{A} = \text{score} \times \text{coeff}^T$ . Instead of using all  $n - 1$  coefficients it is possible to only use the first  $q$  coefficients resulting in an approximation of the space. The parameters  $\mathbf{p}_{q,\text{new}}$  for a new input mesh  $m_{\text{new}}$  can be easily computed for  $q$  coefficients by first reshaping the mesh to be of the form  $\mathbb{R}^{3k \times 1}$  and subtracting the mean  $\bar{m}$ . This is done the same way as above, adding an additional step to transpose. Then the pseudoinverse of  $\text{coeff}_{(q)}$  is multiplied from the left to compute the corresponding parameters  $\mathbf{p}_{q,\text{new}}$ :

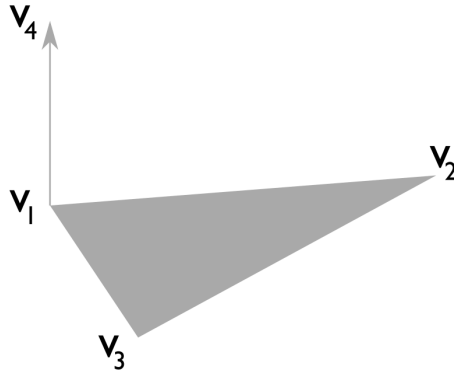
$$\mathbf{p}_{q,\text{new}} = \text{coeff}_{(q)}^+ \cdot (m'_{\text{new}} - \bar{m})^T \text{ where } \text{coeff}_{(q)} := \begin{bmatrix} c_1 & c_2 & \dots & c_q \end{bmatrix} \in \mathbb{R}^{3k \times q} \quad (3.4)$$

In this example the parametric model is based on the vertices of the mesh. In the next subsections, two other variants are explored and explained.

#### 3.2.1. Face Deformations

Instead of defining a mesh by its vertices, it is possible to describe it by the deformations of the faces. This is done by defining one source mesh and computing how each face is deformed in relation to the source mesh. One way to quantify a deformation of a face was described by Sumner [SP04], where the idea was to transfer triangle deformations between similar meshes. First, a new vertex needs to be computed such that an affine transformation can be determined. The fourth vertex is defined as follows:

$$\mathbf{v}_4 = \mathbf{v}_1 + (\mathbf{v}_2 - \mathbf{v}_1) \times (\mathbf{v}_3 - \mathbf{v}_1) / \sqrt{|(\mathbf{v}_2 - \mathbf{v}_1) \times (\mathbf{v}_3 - \mathbf{v}_1)|} \quad (3.5)$$



**Figure 3.2.:** A vertex is added to the triangle that lies in the direction of the normal.

The crossproduct is scaled by the reciprocal of the square root of its length, to keep the distance proportional to the length of the triangle edges. See figure 3.2 for a visualization of adding a vertex on the normal. Given both faces with one additional vertex, the following equations can be posed:

$$\mathbf{Q}\mathbf{v}_i + \mathbf{d} = \tilde{\mathbf{v}}_i, i \in 1 \dots 4 \quad (3.6)$$

where  $\mathbf{Q} \in \mathbb{R}^{3 \times 3}$  and translation vector  $\mathbf{d}$  describe the affine transformation. By subtracting the first equation from the following three equations and rewriting the resulting system in matrix form, the problem can be defined as  $\mathbf{Q}\mathbf{V} = \tilde{\mathbf{V}}$  where

$$\mathbf{V} = \begin{bmatrix} \mathbf{v}_2 - \mathbf{v}_1 & \mathbf{v}_2 - \mathbf{v}_1 & \mathbf{v}_2 - \mathbf{v}_1 \end{bmatrix} \quad (3.7)$$

$$\tilde{\mathbf{V}} = \begin{bmatrix} \mathbf{v}_2 - \mathbf{v}_1 & \mathbf{v}_2 - \mathbf{v}_1 & \mathbf{v}_2 - \mathbf{v}_1 \end{bmatrix} \quad (3.8)$$

The closed form expression for  $\mathbf{Q}$  is defined as

$$\mathbf{Q} = \tilde{\mathbf{V}}\mathbf{V}^{-1} \quad (3.9)$$

To fully describe a mesh, this  $\mathbf{Q}$  matrix needs to be computed for each face of the mesh. Finally, the mesh is represented by  $\mathbf{D} \in \mathbb{R}^{3 \times 3h}$  where  $h$  is the number of faces the mesh has. To be able to run PCA, each mesh needs to be reshaped to be of form  $\mathbf{D}' \in \mathbb{R}^{1 \times 9h}$  and the rest of the process is the same as above in section 3.2.

### 3.2.2. Point Normals

This method is very much similar to one described in section 3.2. In addition to the vertices of the mesh, one vertex per face is computed as described in equation 3.5 and added to the list of vertices of the mesh. Therefore the mesh matrix will be of form  $\mathbb{R}^{k+h \times 3}$  where  $k$  is the number of original vertices and  $h$  is the number of faces of the mesh.

## 3.3. Mapping

The following segment describes how a mapping can be computed in a linear or a non-linear way. The data available is the parameters returned by both the parametric models for the before and after examples. The matrices of the data are

$$\mathbf{P}_{\text{before}} = \begin{bmatrix} \mathbf{p}_{b,1} \\ \mathbf{p}_{b,2} \\ \vdots \\ \mathbf{p}_{b,n} \end{bmatrix} \in \mathbb{R}^{n \times n-1}, \mathbf{P}_{\text{after}} = \begin{bmatrix} \mathbf{p}_{a,1} \\ \mathbf{p}_{a,2} \\ \vdots \\ \mathbf{p}_{a,n} \end{bmatrix} \in \mathbb{R}^{n \times n-1} \quad (3.10)$$

where each parameter pair  $(\mathbf{p}_{b,i}, \mathbf{p}_{a,i}) \forall i$  is related, as the after was generated from the before.

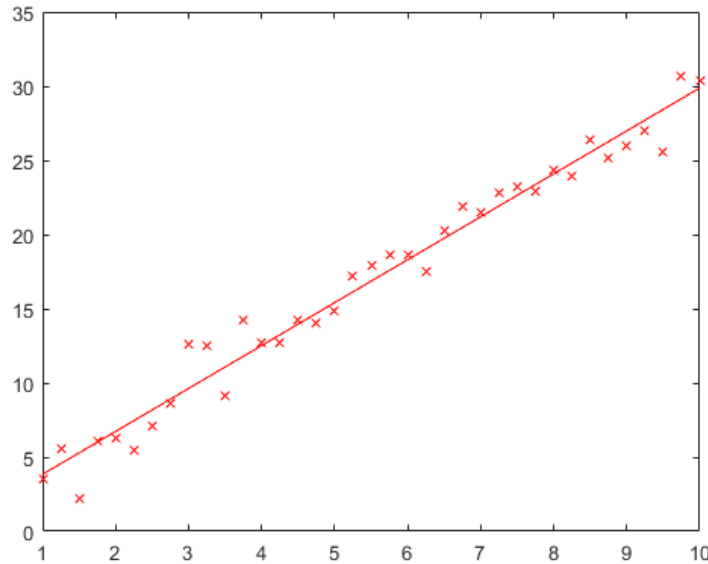
### 3.3.1. Linear Method

The first method used, was the linear system solver by MATLAB (also known as backslash solver) that solved the equation

$$\mathbf{P}_{\text{before}} \cdot \mathbf{M}_{\text{linear}} = \mathbf{P}_{\text{after}} \text{ where } \mathbf{M}_{\text{linear}} \in \mathbb{R}^{n-1 \times n-1}. \quad (3.11)$$

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It is clear that an explicit linear transformation could be found for each parameter pair, but the goal is to have one mapping that works for each parameter pair. This is also known as linear regression. One basic example in linear regression is the task of fitting a line to a point cloud. As all the points of the point cloud do not lie on one line, it isn't possible to define a straight line. Therefore a line is sought after, such that all points are as close as possible on average. This is depicted in figure 3.3.



**Figure 3.3.:** Fitting a line to a point cloud using linear least-squares.

#### 3.3.2. Non-Linear Variants

This section deals with non-linear methods to solve the regression problem described above. All the methods used are part of the scikit-learn library [PVG<sup>+</sup>11]. It is common in a machine learning setting to split data into two sets. One set is used to train the model and the other is used to evaluate how well the method does on data, that was not part of the first set. Both data sets should be drawn from the same distribution. This is necessary because the model shouldn't be penalized for performing badly on problems it didn't encounter in the training phase.

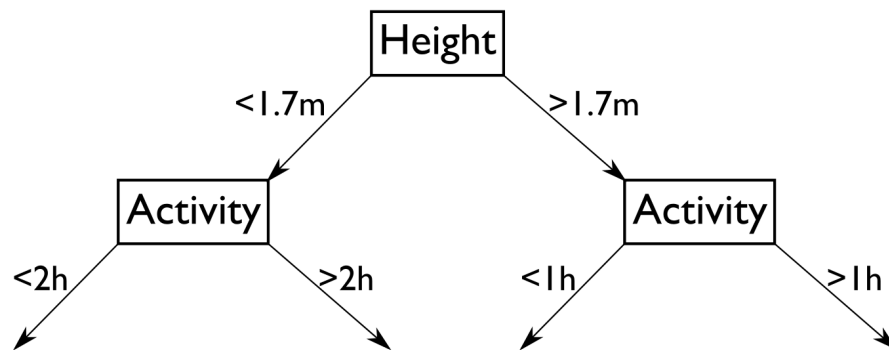
During the training phase it is standard to run grid search. Grid search varies the parameters of the model to find the best set of parameters for a specific task. It needs to be considered that grid search can lead to overfitting, this occurs when the error on the training set is reduced at the cost of increasing the test error. This can be avoided by adding regularization terms or by running grid search along with cross-validation.

#### Decision Tree

One category of machine learning approaches is based on so called decision trees (DT). A decision tree uses a tree-like approach, where nodes of the tree are defined to split the data set

into subsets corresponding to groups. As an example, the goal is to have a model that can predict the weight of a person, based on height, daily activity and daily calorie intake as features. Using a decision tree, the value of height could split the data set, as taller people tend to be heavier than small people. This is done over all features until the complete data set is divided into smaller subsets. Leafs make up the end of a decision tree. Each leaf has a certain value connected to it, in the case mentioned before it would be a certain weight range or value.

When the weight of a new person would need to be predicted, one would follow the tree from top to bottom until a leaf was met. The corresponding weight value of the leaf would be equal to the prediction the model returns. An example of a decision tree can be seen in figure 3.4.



**Figure 3.4.:** Example of a decision tree following example of section 3.3.2

## Random Forest

The random forest (RF) regressor is also a tree-like approach similar to the decision tree. The main difference is that the decision tree considers all features when splitting and thus can be prone to overfitting. Random forest chooses features at random and builds decision trees based on those features. This is done for multiple trees until all subtrees are combined.

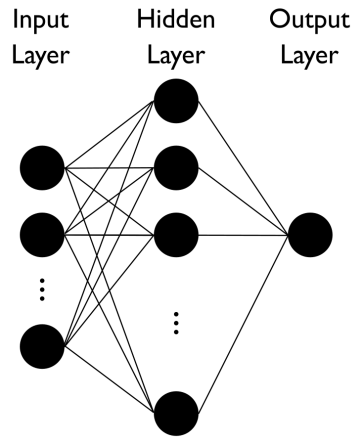
## Multilayer Perceptron

The multilayer perceptron (MLP) is a type of neural network. It uses a feedforward approach to propagate information through the network. MLP have at least three layers where the first layer is the input layer and the last layer is the output layer. Each layer is made up of neurons, that receive information from the previous layer, apply some operation followed by an activation function and pass the result on to the next layer. These activations can either be linear or non-linear. It doesn't make sense to have multiple intermediate layers with linear activation functions, as each linear combination of linear functions, can be described one other linear function.

Each layer is made up of multiple neurons, each with their own weights and biases, and an activation function. All weights and biases are initialized at random. In the training phase a technique known as backpropagation is applied. Given an example, input and outcome, the input data is passed through all the layers. The resulting output is compared to the actual outcome. Given the difference between the two values, weights and biases are modified to

### 3. Methods

reduce the difference, starting from the output layer, moving back through the hidden layers up to the input layer. A basic scheme of a MLP is shown in figure 3.5.



**Figure 3.5.:** Scheme of a multilayer perceptron for regression.

## 3.4. Parametric Model from Editor

In all of the previous sections the parametric model was based off of real world data. In this section the same procedure is done as in section 3.2, except that the meshes used for PCA were generated by software called MakeHuman<sup>4</sup>. MakeHuman is a free and open source character modelling tool. It starts out with a gender neutral base mesh that can be modelled by varying sliders. These sliders apply linear interpolations to the base mesh between the extreme positions. There are over 40 sliders in MakeHuman and the sliders are categorized by what areas of the body they modify. The graphical user interface of MakeHuman can be seen in figure 3.6. Using the scripting feature in MakeHuman multiple meshes could be generated, exported and later imported into MATLAB.

## 3.5. Fitting parametric models to point clouds

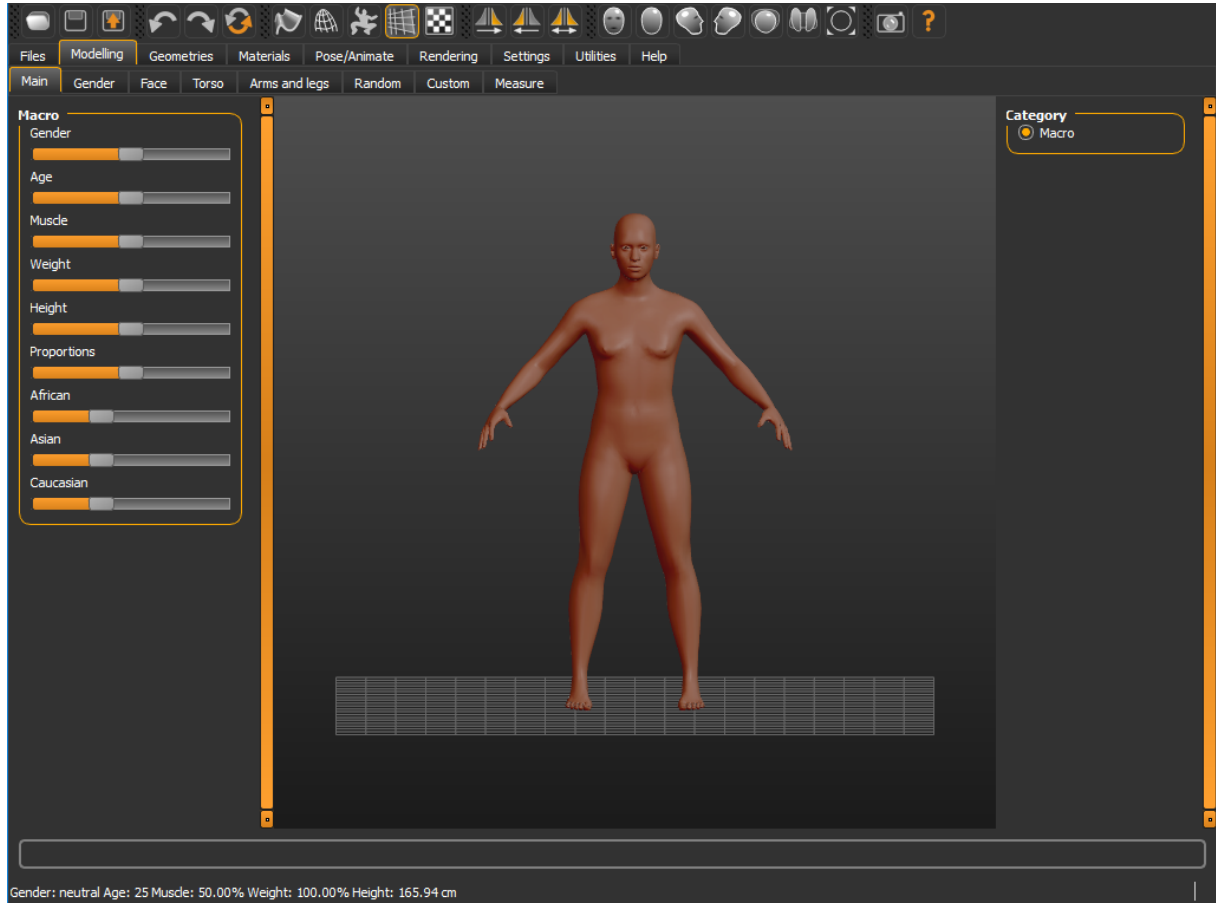
Given a parametric model, it is possible to create a large diversity of possible outputs by varying the parameters. But finding the specific parameters for a new point cloud is a non-linear optimization problem. Therefore a C++ implementation by Biland [Bil17], using an open source C++ library for optimization problems called the Ceres solver [AMO], was slightly modified and used. This implementation allows to fit a parametric model to a point cloud given a few keypoints of the point cloud.

In a first step the keypoints are used to find the parameters of translation, rotation and scaling. Afterwards a select number of points are chosen as correspondences between the model and

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<sup>4</sup><http://www.makehumancommunity.org/>





**Figure 3.6.:** Graphical user interface of MakeHuman.

the point cloud. Iteratively this implementation tries to minimize the distance between all the correspondences by varying the parameters of the model. Additionally, constraints are applied to the values the parameters are allowed to take. This prevents one parameter becoming too large and also possibly forcing the model to use multiple parameters than to rely on a single parameter. After a certain number of iterative steps has been reached, a certain threshold in accuracy is passed or the change in parameters is vanishingly small, the resulting mesh and parameters are returned.



## Results

In the next section two error metrics are briefly described, followed by evaluations of each experiment. At the beginning of each section a small description of the problem setting is given, followed by the results and discussion.

### 4.1. Error Metrics

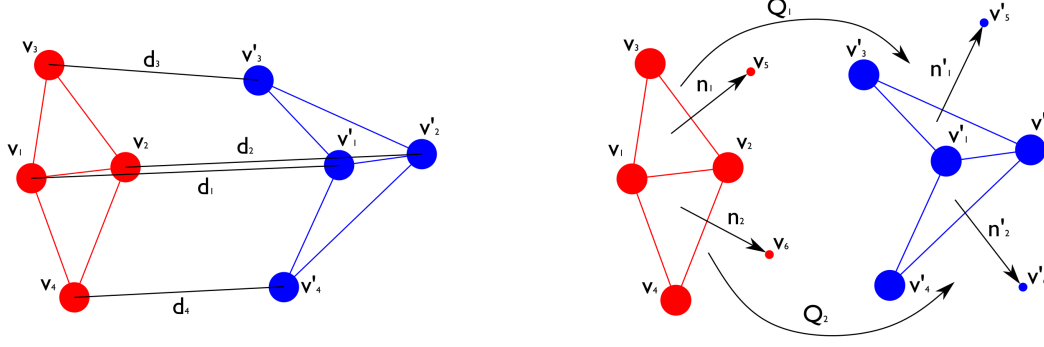
For this thesis two error metrics were considered:

- **Point Distance:** Given two meshes with equal topology, meaning both meshes have the same amount of vertices and layout, the error computed is based on the distance between corresponding points. In general, the mean distance is computed over all point correspondences.
- **Face orientations:** The idea behind this method is based on the  $Q$  matrix mentioned in section 3.2.1. Given two topologically equal meshes, the  $Q$  matrix is computed for all corresponding faces. As it is of interest to investigate the difference in breast shapes, this error tries to quantify the variation of the  $Q$  matrices.

One drawback of the point distance method occurs when the two meshes are not aligned properly. This can be solved by applying the same algorithm mentioned in section 3.1.1 before computing the error. One advantage of this method is the unit of the error, as it could be converted into a unit of distance, given the distance between two point is known in real world lengths. This error can also be computed when two meshes are not topologically equal by first finding the correspondences between vertices.

One advantage of the face orientation method is the invariance of translation and global rotation. This means that moving or rotating the mesh doesn't affect the error. On the other hand, the

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**Figure 4.1.:** The left image depicts distances between corresponding points of the mesh. The right image shows a representation of the information used to compute the  $Q$  matrix.

values of this error don't have a real world meaning and it is a lot harder to apply this method to topologically different meshes, as finding correspondences between faces is a lot more challenging.

The methods are depicted in figure 4.1. Both methods have their advantages and disadvantages, but due to the fact that alignment isn't a problem, the error metric used in this thesis is the point distance method.

## 4.2. Evaluation

In the following section the evaluation methods, that were used, are explained. Optimally, when evaluating a prediction some representation of the true outcome, often called the "ground truth", is given. This is useful to quantify how accurate the prediction was.

### 4.2.1. Mesh Point Error

For this method the comparison is done as explained in section 4.1. Before evaluating the point distances between a prediction and the ground truth, the meshes are aligned as mentioned in section 3.1.1. Additionally, the mean point distance error is computed by dividing by the number of vertices in the mesh and also by dividing by the diagonal of the largest bounding box over all examples to scale the error to a meaningful unit. Given a prediction of a mesh  $M^{(p)}$  and the ground truth  $M^{(g)}$  the problem can be stated as:

$$error = \frac{1}{nd_{max}} \sum_{i=1}^n \sqrt{\sum_{j=1}^3 (M_{i,j}^{(p)} - M_{i,j}^{(g)})^2} \quad (4.1)$$

where  $n$  is the number of vertices of the mesh,  $d_{max}$  is the diagonal length of the largest bounding box over all examples and the final result is a scalar value.

### 4.2.2. Error Heat Map

The idea of this method is to generate a visualization of the error, indicating where the prediction failed to represent the ground truth the most. This can be accomplished by computing the error described in section 4.1 but instead of adding up the errors keeping them separate for each vertex. This error per vertex is accumulated over multiple samples. Given the  $k$ -th prediction of a mesh  $M^{(p_k)}$  and the corresponding ground truth  $M^{(g_k)}$  the problem can be described as:

$$e'_i = \sum_{k=1}^m \sqrt{\sum_{j=1}^3 (M_{i,j}^{(p_k)} - M_{i,j}^{(g_k)})^2} \quad \forall i = 1 \dots n \text{ with } e = \frac{e'}{\max(e')} \quad (4.2)$$

where  $m$  is the number of samples,  $e'$  is a vector of the accumulated errors over all samples and  $e$  is a vector of the errors scaled by the largest entry in  $e'$ . This is done to visualize the error on a range, such that, for example the error can be color coded where red is equivalent to a large error and green is a small error.

## 4.3. Input Data for PCA

In this experiment three different variations of data inputs are tested for PCA as described in section 3.2 and the following two subsections. Based on the resulting parameters of each model, a linear mapping is created and compared based on the error of the predictions.

### 4.3.1. Problem Setting

Given a data set comprised of 57 pairs of meshes where one mesh is based on the before point cloud and the other is based on the after point cloud. Both sets are split into equally sized sets of 42 training pairs and 15 testing pairs. Each variation is trained on the training set and afterwards evaluated on the test set. The evaluation is done by computing a linear mapping over the parameters of each model. Then the mean mesh point error over all 15 testing pairs is computed and a error heat map is generated.

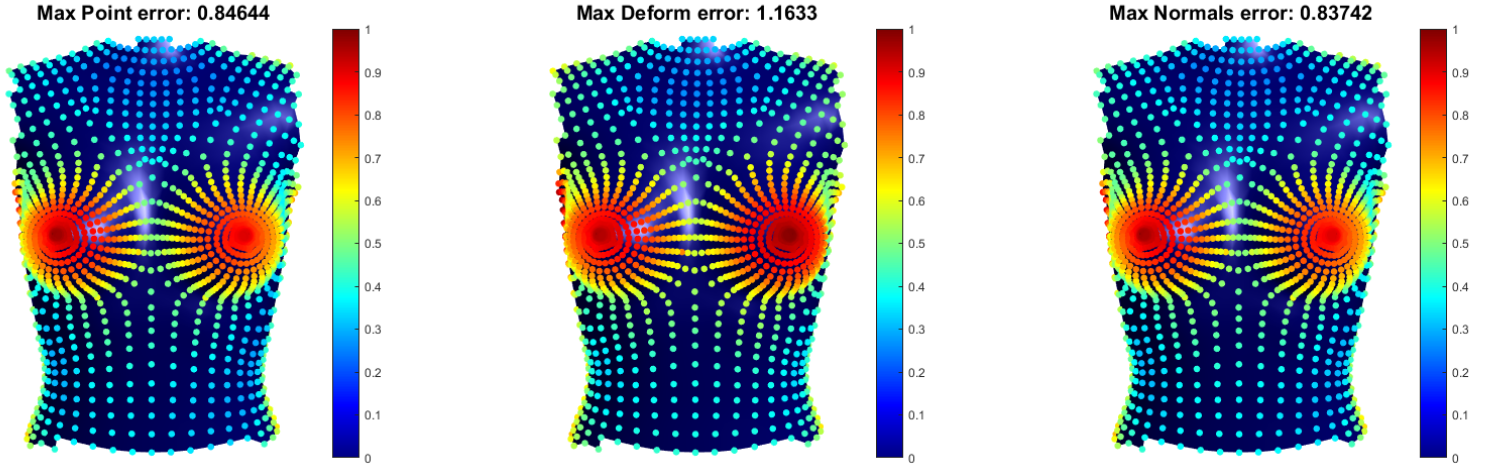
### 4.3.2. Results

In the table 4.1 the time column is the amount of time each method needed to compute the parameteric model. The training mean error is computed as mentioned in 4.2.1 for each case. The mean is computed over all cases. The same is done for the test mean error. In each subfigure in figure 4.2 the values that are listed above are the respective  $\max(e')$  as explained in 4.2.2. Additionally, one vertex on the border doesn't have a colored point as the error wasn't being computed correctly for it.

#### 4. Results

	Time	Training Mean Error	Test Mean Error
<b>Point</b>	0.06s	0.00733	0.01845
<b>Deformation</b>	6.81s	0.0090	0.02795
<b>Points and Normals</b>	2.92s	0.0071	0.01841

**Table 4.1.:** Results in terms of mean mesh point error are listed for the training and for the test set. Also the time each function took to compute the parametric model is listed in seconds.



**Figure 4.2.:** The left image shows the error heat map when using the point method. The middle one uses the deformation method. The right one uses the point normals method. The color of each point reflects the amount of the error. Red color indicates large error, blue small error.

### 4.3.3. Discussion

The amount of time it takes to compute the parametric model based on deformations takes a lot longer due to the fact that additional normals need to be computed followed by the computation of the  $Q$  matrix. One can see that the additional time to compute normals is also reflected in the point and normals method. Therefore roughly 40% of the time is accounted for by the normal computation.

All three methods worsen going from training to test case, but that is expected. This is caused by the models having seen the training set before and are already familiar with those examples. Of the three methods the points and normals method achieves the best results over test and training. The reason this occurs is due to the fact that the normals carry some information of the size of the faces.

Figure 4.2 looks similar for all three cases. The error for the deformation method is higher than for the other two methods. Also around the right breast the red area seems larger and darker than for the other two. Visually no difference can be made out between the point method and the point normals method.

## 4.4. Learning Mapping

The next step was to investigate, if non-linear mappings could perform better than a linear mapping. Additionally, it is checked how the error changes, when PCA is done on the complete set, instead of, only the training set. The reason, this is done, is because the mapping shouldn't be penalized or perform badly because of the parametric model. For each of the approaches the best parameters need to be found first by running cross-validation grid search. Each of the learnt mappings can be applied to the test set and compared against the ground truth. The mean point error is computed to measure the error.

### 4.4.1. Problem Setting

The parameters for each of the 57 pair meshes are computed according to two parametric model. These will be referred to as the "complete" and "incomplete" model, where the "complete" model was based on the training and test data, where as the "incomplete" model only received the training data. The non-linear approaches used are a decision tree regressor, a random forest regressor, a multilayer perceptron with a rectified linear unit (ReLU) as its activation function and a multilayer perceptron without a ReLU.

For each approach the best parameters are evaluated by running cross-validation grid search for either cases. The parameters varies are listed in a table that can be found in the appendix A.1. Finally, the training and test errors are computed based on the best parameters and a ratio between test and training error is calculated. Optimally, this ratio should be one, as the test error should be equally low as the training error.

#### 4. Results

PCA Model	Method	Training	Test	Ratio
		Mean Error	Mean Error	Test/Train
Incomplete	Linear	0.0073	0.0184	2.52
	Random Forest	0.0068	0.0196	2.88
	Decision Tree	0.0127	0.0237	1.86
	MLP	0.0154	0.0186	1.21
	MLP ReLU	0.0173	0.0203	1.17
Complete	Linear	0.0076	0.0357	4.70
	Random Forest	0.0077	0.0196	2.55
	Decision Tree	0.0125	0.0218	1.74
	MLP	0.0152	0.0193	1.27
	MLP ReLU	0.0161	0.0193	1.20

**Table 4.2.:** This tables shows the mean errors computed on the training and test set. The first five rows show the results for the "incomplete" model, the last five for the "complete" model. The first two columns are the training and test error, followed by the ratio of test to training error.

#### 4.4.2. Results

The results for the different cases can be seen in 4.2. The table is divided into two sections, the first half for the results of the "incomplete" model and the second for the "complete" model. The first row of results are the mean errors for the training set, the second row for the test set. The error for a prediction is computed by comparing to the ground truth with the mesh point error as explained in 4.2.1. The resulting error is equivalent to the mean over all errors. The last column is the ratio between test and training.

#### 4.4.3. Discussion

It is interesting to see that the training error of the linear method and the random forest increases slightly going from the "incomplete" to the "complete" model. This must be caused by some examples in the test set being similar to some of the examples in the training set and forcing the parameters to adjust. The other three methods improve slightly in terms of training error due to the "complete" model. Independant of case, the test error is always larger than the training error. This is expected and was already discussed above.

The behaviour of the test error going from "incomplete" to "complete" is similar to the training error. The linear method worsens gravely, perhaps due to the "complete" model describing a wider space, making it harder for the linear method to find an appropriate mapping. All of the non-linear methods perform similarly in the "complete" case compared to the "incomplete" case



in terms of mean test error.

Even though the best performing method in terms of error is the MLP, the MLP with ReLU has similar error values but has a slightly better test/train ratio. This could indicate that this method is more robust, meaning that the error behaves similarly when applied to new data.

## 4.5. Parametric Models

Given two parametric models and a test set of point clouds, both models process the point clouds as described in section 3.5.

### 4.5.1. Problem Setting

### 4.5.2. Results

### 4.5.3. Discussion



# 5

## **Conclusion**

### **5.1. Outlook**



# A

## Appendix

## A. Appendix

PCA	Method	Possible Parameters	Best
Incomplete	Random Forest	'estimator__n_estimators':[5,10,20,30],	30
		'estimator__max_features': ( 'auto', 'sqrt', 'log2' ),	auto
		'estimator__max_depth':[2,4,8,16]	8
	DT	'estimator__criterion':('mse', 'friedman_mse', 'mae'),	mse
		'estimator__splitter':('random', 'best'),	best
		'estimator__max_features':('auto', 'sqrt', 'log2'),	auto
	MLP	'estimator__max_depth':[2,4,8,16]	2
		'estimator__activation':('identity', 'tanh', 'logistic'),	identity
		'estimator__solver':('lbfgs', 'adam', 'sgd'),	sgd
	MLPrelu	'estimator__hidden_layer_sizes': [(15,),(30,),(45,),(60,),(45,30,),(60,45,)],	(60,45,)
		'estimator__activation':( 'relu' ),	relu
		'estimator__solver':('lbfgs', 'adam', 'sgd'),	adam
		'estimator__hidden_layer_sizes': [(15,45,),(45,30,),(60,45,),(60,45,30),(60,45,15)],	(45,30,)
		early_stopping: true	true
Complete	RF	'estimator__n_estimators':[5,10,20,30],	30
		'estimator__max_features':('auto', 'sqrt', 'log2'),	auto
		'estimator__max_depth':[2,4,8,16]	4
	DT	'estimator__criterion':('mse', 'friedman_mse', 'mae'),	friedman_mse
		'estimator__splitter':('random', 'best'),	best
		'estimator__max_features':('auto', 'sqrt', 'log2'),	auto
	MLP	'estimator__max_depth':[2,4,8,16]	2
		'estimator__activation':('identity', 'tanh', 'logistic'),	tanh
		'estimator__solver':('lbfgs', 'adam', 'sgd'),	sgd
	MLPrelu	'estimator__hidden_layer_sizes': [(15,),(30,),(45,),(60,),(45,30,),(60,45,)],	(60,45,)
		'estimator__activation':( 'relu' ),	relu
		'estimator__solver':('lbfgs', 'adam', 'sgd'),	sgd
		'estimator__hidden_layer_sizes': [(15,45,),(45,30,),(60,45,),(60,45,30),(60,45,15)],	(60,45,30,)
		early_stopping: true	true

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