



Bavarian Graduate School of Computational Engineering

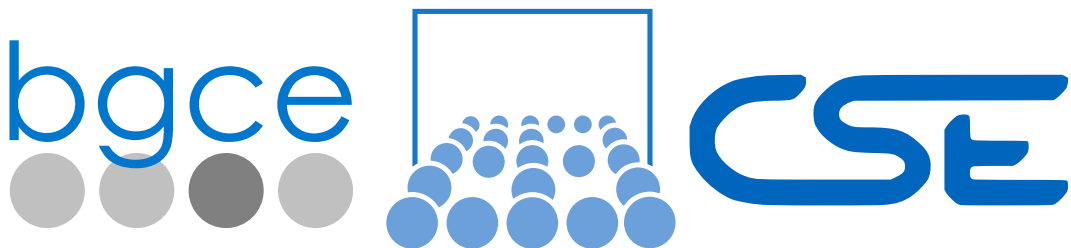
Technische Universität München

BGCE Honours project report

CAD-integrated Topology Optimization

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Preface

The Bavarian Graduate School of Computational Engineering (BGCE) honours project at the Computational Science and Engineering (CSE) Institute of Technische Universität München (TUM) is a 10-month project where students conduct research on cutting-edge topics in the field of Computational Engineering, in cooperation with a partner in industry or academia. The 2015-16 project is titled *CAD-Integrated Topology Optimization* and is initiated and supervised in a cooperation between TUM and Siemens AG in Munich.

Acknowledgments

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Outline and Overview

Purpose of the document

The purpose of this document is to describe the *CAD-integrated Topology Optimization* software tool along with the theoretical background it relies on.

Document overview

The document is arranged in chapters, covering introductions to the field and the project, background theory, and parts of implementation. The chapters are described in more detail below.

CHAPTER 1: INTRODUCTION

This chapter presents an overview of the motivation behind *CAD-integrated Topology Optimization*, including the current state of the field. It also provides general organizational information about the project execution, timeline and structure.

CHAPTER 2: BACKGROUND THEORY

This chapter introduces the theoretical background for the implementation of the *CAD-integrated Topology Optimization* tool. It consists of four parts, each describing essential background of the topology optimization pipeline. Furthermore, detailed description of selected algorithms used in each step is given.

CHAPTER 3: IMPLEMENTATION

This chapter provides details on the implementation and structure of the *CAD-integrated Topology Optimization* tool itself. The different parts of the topology optimization pipeline are presented along with underlying implementation details.

1 Introduction

1 Introduction

In this chapter, the motivation for the project and a short description of the problem task is provided, together with a brief introduction to topology optimisation and the project structure.

1.1 Motivation

A common problem in product design is to create a functional structure using as few material as possible. Three decades ago engineering design versions were drawn, prototypes created and experimental test performed. Nowadays, the field of topology optimisation simplifies this process and has become a great help in all fields of engineering.

Topology optimisation tackles the problem of material distribution in a structure in order to fulfil certain target loads. Several topology optimisation open-source tools exist that are ready to use; however, it is still a challenge to incorporate these tools handily in the design process. The idea of this project is to allow these tools to work directly from CAD files and to transfer the obtained mesh based solution of the topology optimization tool back to the CAD world. Unfortunately, at the moment, there is no open source solution for the conversion mesh based geometry to the spline-based CAD format (where in our case *Non-Uniform Rational B-Splines*, commonly referred to as NURBS, are used). The would-be straightforward approach – to convert each triangle of mesh geometry directly to CAD format – results in enormous file sizes. One of the biggest challenges of this project is thus to develop a conversion tool that feasibly provides a useful CAD-representation of the optimized surface.

1.2 Project structure

The aim of the project is to provide a tool that allows to utilize topology optimisation without leaving the CAD-framework. Therefore, the main goals of the project are as follows:

- Implementation of a CAD format-accepting topology optimisation framework by using and extending available open source libraries
- Development of a flexible tool for conversion of an optimized surface back to the CAD format.

The duration of the project has been set to 10 months. Hence, this project has been divided into 4 phases:

Phase 1: Getting familiar with the topic and agreement on the project specification.

Phase 2: Implementation of the first part of the pipeline (Topology Optimisation from CAD surface using existing tools); investigating the tools and algorithms available for the conversion

of the geometry generated after topology optimisation back to CAD format (later referred as *NURBS fitting pipeline*); prototyping (using MATLAB) and evaluating of found results.

Phase 3: Implementing the prototypes developed on the previous stage, using non-proprietary languages; extension of the NURBS fitting pipeline to more complex cases; finalising the first part of the pipeline.

Phase 4: Implementation of the extended NURBS fitting pipeline; integration with the topology optimisation part and delivering the final product to costumer.

2 Background Theory

2 Background Theory

In this chapter, the theoretical background for the implementation of the *CAD-integrated Topology Optimization* tool is presented. The chapter is divided into four sections: *CAD overview* (section 2.1), *Topology Optimisation* (section 2.2), *From Voxels to a surface representation* (section 2.3), and *From a surface representation to NURBS* (section 2.4).

2.1 CAD overview

Computer Aided Design (CAD) refers to the process of designing a product using a computer system. Before CAD applications were used, products were designed using a sketch board. It was a challenge to incorporate changes in the construction drafts as well as to keep documentations up to date; hence, it was no surprise that CAD systems spread rapidly across all design development branches. Computer aided design is now of irreplaceable use in architecture, mechanical, electrical and civil engineering.

Depending on the discipline, different requirements are set on the virtual model. One may imagine that in a civil engineering model of a building a 2D floor plan is often sufficient; however, in the design of a mechanical motor a 3D model is always necessary. Given these circumstances, various CAD software bundles evolved in the different disciplines with completely different modelling approaches. Besides the geometry representation additional parameters, such as material properties or manufacturing information, are stored. In order to move between different data structures standardized exchange interfaces are commonly used.

This section presents the relevant geometrical and computational aspects of CAD for the project; for a more thorough introduction we refer to [1].

2.1.1 Geometry representations

In general, two different ways of describing a geometry are used in CAD systems: a *constructive solid geometry* (CSG) or a *boundary representation* (BREP). Other approaches, such as a complete voxelised geometry are not common due to extensive memory consumption.

Constructive solid geometry

One way of representing a geometry in CAD is the approach of CSG. The basic idea is to start from a set of primitives, e.g. spheres, cylinders and/or cubes. Basic Boolean operations link these primitives towards a complex geometry, as illustrated in Figure 2.1.

Key advantages of this format is the precise representation using very little storage memory. However, not all desired forms can be represented by CSG and hence, a second type of geometry description is needed.

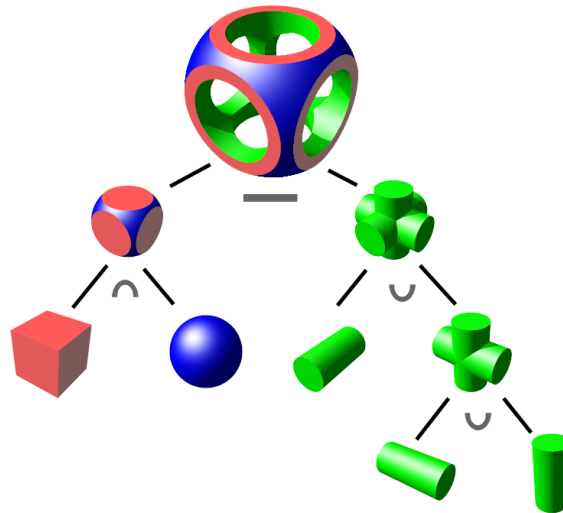


Figure 2.1: CSG object tree. The picture shows the construction of a complex object from a cube, a sphere, and a set of cylinders. Figure taken from [2].

Boundary representation

A different kind of modelling approach is the BREP. Instead of storing the geometry information as geometrical objects, BREP formats only save the boundary surface of the body. The interior is assumed to be uniformly filled. Especially in complex geometries, this approach simplifies the model to such an extent, that the amount of data becomes much easier to handle. Surfaces can then be for example stored as a set of triangles (as in STL files, see section 2.1.2 below) or in NURBS patches (see section 2.4). Furthermore, holes in the body are made possible by saving the surface normal of the respective boundary.

By the boundary representation arbitrary geometries can be created. While the data sizes are commonly larger than in CSG representation, BREP files are usually easier to work with. One also has to keep in mind, that non-physical geometries can result from BREP formats through a not closed surface.

2.1.2 Data exchange interfaces

CAD software programs usually use their own data formats; in order to exchange models standardized interface formats have been developed. Geometric models are compressed to certain geometry descriptions; transferring additional information, such as material properties or manufacturing information, is in general a difficult task and in some exchange file formats even prohibited. A few common exchange file types are described below, as also compared in [3].

STL file format

The STL (from *ST*ereo*L*ithography) file format describes the model only by its boundary and is thus a BREP format. The idea behind its files is simple. The geometric model is discretized into a cloud of points, where sets of three vertices form a triangle; hence, a connected surface of

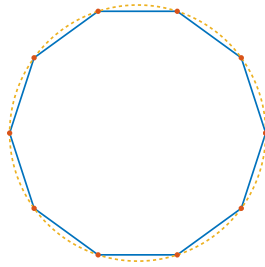


Figure 2.2: STL discretization for a circle, self-made in MATLAB [4]. Note that the circle cannot be exactly represented by the vertices and edges.

triangles emerges which describes the geometry. The procedure is shown in Figure 2.2 for a two dimensional circle. The aforementioned triangles boil down to lines in two dimensions. The advantages and disadvantages of this approach become clear: It can be applied to an arbitrary geometry, but accuracy causes difficulties. In order to transfer high precision geometries many vertices are necessary, resulting in big files. Still, as is illustrated in the figure, a perfect circle can never be represented.

ASCII STL files begin with a name and the data on the triangles is constructed as follows:

- a facet normal pointing outward
- a sequence of vertex coordinates

As this is the only information provided, no additional data such as material properties are transferred through STL files, reducing file size but also range of usage.

STEP and IGES file formats

To overcome issues of insufficient precision also more elaborate exchange formats exist; these save e.g. a circle as a parameter where no discretization step is involved. Also, the possibility of passing additional parameter information (e.g. density, manufacturing information) is required by certain users. Popular file types that offer these two functionalities are STEP and IGES files.

The STEP file format is a newly developed CAD data exchange standard and documented in the ISO 10303 norm. On the contrary to STL files it uses a combination of CSG and BREP to store the geometry. Additional information (e.g. density) are passed through attribute sets that are stored besides geometry instances (e.g. a circle). A key disadvantage, however, is that STEP files carry much redundant information [3].

The Initial Graphics Exchange Specification (IGES) is an American National Standard since 1981 to exchange graphics information. Similar to the STEP format it uses a combination of CSG and BREP for the geometry representation. Instead of storing a set of manufacturing information as done in the STEP file, the IGES is build only to exchange graphics information. For example, the step file transfers a physical density information; in the IGES format the only additional parameter store node coloring information. Consequentially, IGES file sizes are significantly smaller compared to the STEP file format [3].

The IGES file format contains five different sections: a *Start*, *Global*, *Directory Entry*, *Parameter Data* and *Terminate* section. The *Start* and *Global* section are used for naming and part information. In the *Directory Entry* section additional information like the node color is saved. The

Parameter Data section is used for storing the coordinate points; the *Terminate* section signals the end of the file [1].

2.2 Topology Optimisation

Topology optimization describes the process of finding the optimal distribution of a limited amount of material for a given area or volume based on a predefined constraint/minimization problem. Possible optimization goals are for example [5]:

- **Minimum compliance**, in which one seeks to find the optimal distribution of material that returns the stiffest possible structure. The structure is thereby subjected to loads (forces) and supports (boundary conditions). By maximizing the stiffness, the compliance is minimized. This is also analogous to minimizing the stress energy stored by the applied loads.
- **Heat conduction**, where one tries to optimize the domain of a conductive material with respect to conductivity for the purpose of heat transfer. This maximization problem is the same as minimizing the temperature gradient over the domain— a poor conductor will create a large gradient.
- **Mechanism synthesis**, where the objective is to obtain a device that can convert an input displacement in one location to an output displacement in another location. Thus, one hereby seeks the optimal design which maximizes the output force for a given input, or respectively, minimizes the input force for a given output.

As one can already imagine by this short list of optimization goals, topology optimization has a wide field of possible applications. Hence, it has become a well established technology used by engineers in the fields of aeronautics, civil, materials, mechanical and structural optimization. Furthermore, due to the rising significance of additive manufacturing techniques in industry, the realisation of complex optimized designs is now much easier. For the rest of this section, and the rest of the document, we will concentrate on the *Minimum compliance* problem. Note however, that almost all parts in the *CAD-Integrated Topology Optimization* tool could just as well be applied to any other topology optimization problem.

2.2.1 Minimum compliance: Problem formulation

In order to constrain the resulting structure as little as possible, the formulation of the topology optimization problem is generally given as follows: for a given set of external fixture points, external loads and/or body forces, the distribution of material within the reference domain should be found such that the structure has maximum stiffness. This is obtained when the structure has the minimum energy stored by external work for the applied forces. The problem is also usually formed to allow for regions in the domain to be specified as filled or empty of material (see Figure 2.3).

The formulation allows the problem to be cast as finding a displacement field u and a stiffness tensor field E that is in equilibrium with the applied loads, and that minimizes the external work done by these external loads to reach that equilibrium.

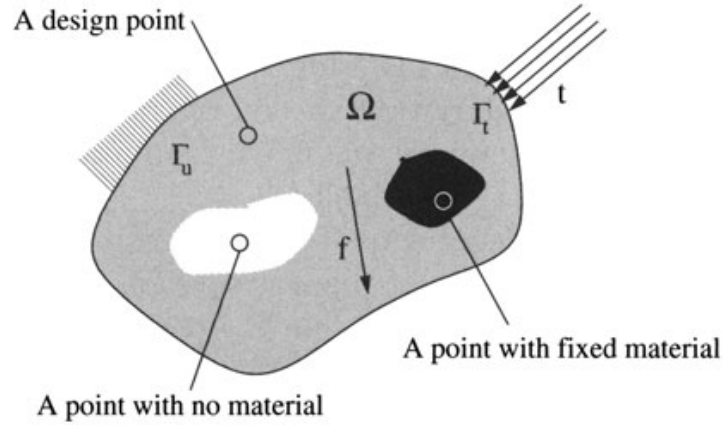


Figure 2.3: The reference domain Ω for the minimum compliance problem. The problem is formulated such that for a set of external loads t on boundaries λ_t , body forces f and a set of fixed support points λ_u , the material distribution within Ω is such that the stiffness with regards to these loads and forces is maximal and the energy stored by the application of those forces is minimal. The problem also allows defining areas which either cannot or must be filled with material. Figure taken from [6].

2.2.2 Physical and mathematical simplifications

To turn this into a more tractable mathematical problem, a few physical assumptions are also typically made: the material be isotropic and linearly elastic. From the assumptions of isotropy and linear elasticity of the material, the stiffness field becomes a constant of the material, defined where there is material in the domain.

The problem is also easy to cast into a weak form. First of all, we compute the integrated internal virtual work and external work. The former is the work of deforming the elastic material from equilibrium by an admissible displacement. The latter is done by the loads and forces to bring out this displacement. Having computed these, we set them equal to one another in order to conserve energy. As a result we obtain an equation that relates the equilibrium displacement, stiffness tensor, and the forces and loads. We then cast this into the weak form, which can be solved using Finite Element Methods (FEM). These can also incorporate the calculation of the external work done.

2.2.3 SIMP: Solid Isotropic Material with Penalization

As described in the previous section, we aim to minimise the external work done by looking at different material distributions. However, the usual problem of finding an optimum arises: the search space is vast. After discretising the domain with FEM, the possibilities of where to put material at least are not infinite – but they still grow exponentially with the number of elements; hence, trying out one-by-one is not going to prove efficient. One popular way of recasting the problem to allow for easier solving is the SIMP model. Here, instead of either being present or not at a point, the material presence can take a continuous set of values between one and zero. The total final volume is then obtained and fixed by integrating this presence variable over the domain, instead of constraining the allowed occupied space. This allows for the interpretation

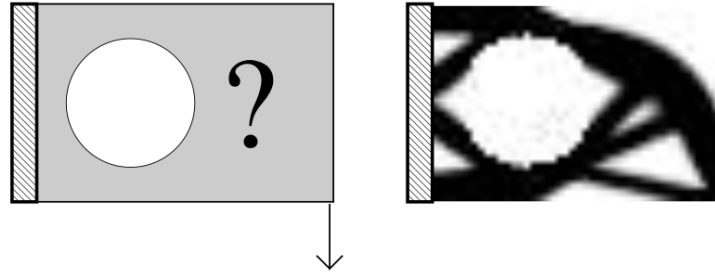


Figure 2.4: Topology optimisation of end-loaded cantilever with fixed hole. The optimisation of a loaded cantilever is one of the model problems in topology optimisation, due to its simplicity and its multitude of used solutions throughout the history of engineering. The picture is taken from a known paper by Sigmund [7] where a 99-line Matlab code for topology optimisation is introduced.

as some kind of density.

In order to still obtain topologies where material is predominant in certain areas– of densities one, with the rest being empty at densities close to zero– a “penalty” is applied to the intermediate values. This is effected by raising the density to a power > 1 in the elastic energy calculation, but not in the volume calculation. That way, an intermediate density value provides less elastic support, but still “costs” as much volume, and will thus be suboptimal.

2.2.4 Solution and implementation

In typical implementations, a heuristic iterative scheme is then used for finding a solution. The optimal solution is assumed to have all present parts stressed (as they would otherwise be unnecessary, not providing any support). Thus, at places where the elastic energy is high, material is added if possible, and where it is low, material is likewise removed, with the values “high” and “low” being determined dynamically to keep the total volume constraint.

This whole scheme is one of the simpler topology optimisation schemes to implement, and has been done so in several pieces of open-source software, including a known 99-line Matlab code by Sigmund [7] and ToPy described in section 3.2. An example optimised topology is shown in Figure 2.4. For an extended explanation and discussion, as well as further alternative methods for topology optimisation, the interested reader is referred to [6].

2.3 From Voxels to a surface representation

In order to fit NURBS and other curves to an optimized geometry, a *mesh-based geometry*, that is, a representation of the object at a set of (connected) points, is typically needed, this as opposed to the volumetric representation of density in each voxel, that results from the topology optimization process.

In order to produce the mesh-based geometry, the data can be represented by a contour at a value of a smooth function in space, that is, an isosurface. Below, we describe two methods that solve this problem, the *Marching Cubes* and the *Dual Contouring* methods.

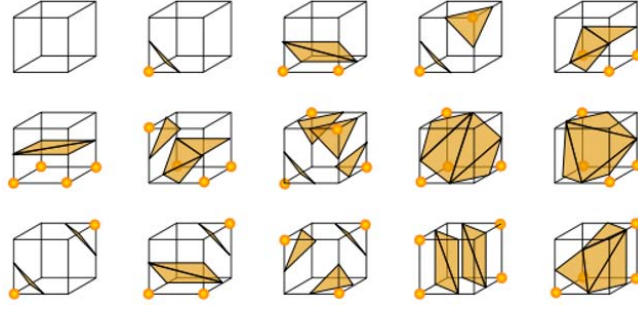


Figure 2.5: The base cases of Marching Cubes. These are drawn with each polygon vertex intercepting its edge in the middle between the cube’s corners, as in the case when the isovalue is exactly halfway between the function values at the vertices. Figure taken from [8].

2.3.1 Marching Cubes

The Marching Cubes method [8] takes as an input a set of scalar function values on a cartesian mesh and extracts an approximate isosurface in the form of a mesh of triangles. The method starts by dividing the space into cubes with the set of points as cube vertices. On these points, the value is determined to be above or below the desired isovalue. According to which corners are set to be above or below, the corner configuration is then mapped to a polygon inside the cube, with vertices on the cube’s edges. On an edge between a vertex above and a vertex below the desired isovalue, the exact location of the surface is then determined via linear interpolation and set as the polygon’s vertex on that edge.

Since there are 8 vertices on each cube, each above or below the isovalue (with equality falling to one of these categories), there are $2^8 = 256$ possible polygon configurations. However, many of these can be constructed by rotating or reflecting other polygons. As such, there are therefore 15 base cases which represent all the surface polygons of the marching cubes. To illustrate what the polygons may look like, these are shown in Figure 2.5, where one can also see that they are composed of triangles. The original algorithm presents two main problems. Firstly, it does not guarantee neither correctness nor topological consistency, which means that holes may appear on the surface due to inaccurate base case selection. Second, another problem is ambiguity, which appears when two base cases are possible and the algorithm chooses the incorrect one. There are many extended Marching Cubes algorithms that tackle the problems of the original one, getting rid of the ambiguities and providing correctness (see for example [9]).

2.3.2 Dual Contouring

The idea of *Dual algorithms*, to which Dual contouring belongs, is similar to Marching Cubes. However, instead of generating polygon vertices on the edges of the cubes, it locates them inside the cubes that have vertex values both above and below the isovalues. After locating the vertices, the ones associated with four contiguous cubes are joined to form a quadrilateral face, commonly referred to as a *quad*. The approach can be seen in Figure 2.6, with a similar Marching Cubes illustration for comparison.

The relevant question is now where in the cube the ideal place for the vertex is, and here is

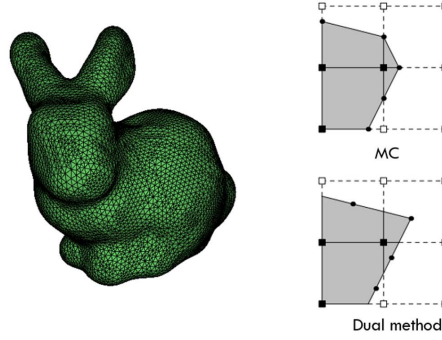


Figure 2.6: *Left*: The famous Stanford Bunny, a popular computer graphics test object, here after application of Marching Cubes. *Right*: Main difference between MC and Dual methods. Figures taken from [10].

where different dual algorithms are distinguished. Dual Contouring in particular generates a vertex positioned at the minimizer of a certain quadratic function, which depends on the (interpolated) isosurface intersection points, as well as the gradient – or just the normal of the isosurface – at these points, together the first-order *Hermite data* of the set. The quadratic function for Dual Contouring is defined in [10] as follows:

$$E(x) = x^T A^T A x - 2x^T A^T b + b^T b$$

where A is a matrix whose rows are the isosurface normals at the intersection points, and b is a vector whose entries are the product of normals and the intersection points. This system can be solved numerically, for example as proposed in [10] by computing the singular value decomposition of A and forming the pseudo-inverse, truncating its small singular values.

The main advantage of this method over Marching Cubes is the acquisition of better aspect ratios [10]. On the other hand the need of gradient data represents a disadvantage.

2.4 From a surface representation to NURBS

Parametrised geometries are often given in terms of *Non-Uniform Rational B-Spline* (NURBS) curve patches (see for example, documentation of FreeCAD software [11]). To define NURBS from a mathematical standpoint, we first define so-called *Bezier curves* and use them later for the definition of NURBS. For these two sections, we refer to [12] for a more in-depth introduction and further material.

2.4.1 Bezier Curves

A Bezier curve is a *parametric* curve, which is often used for producing a smooth approximation of a given set of data points.

An analytical expression for the Bezier curve parametrized by the variable u is given by:

$$\vec{B}(u) = \sum_{i=0}^n b_i^n(u) \vec{P}_i$$

where \vec{P}_i is the i^{th} control point, $i \in 0, 1, \dots, n$ ($n+1$ control points in total), and

$$b_i^n(u) = \binom{n}{i} (1-u)^{n-i} u^i$$

with $\binom{n}{i}$ being a binomial coefficient, is the i^{th} *Bernstein polynomial* (see [13]) of degree n .

Additionally to the expression with the Bernstein polynomials, one can use a recursion formula (so-called *de Casteljau Algorithm*) for the construction of the Bezier curve, which we will not cover here.

Analogically to Bezier curves, but with $n \cdot m$ points $\vec{P}_{i,j}$, one can define a *Bezier surface*, given by the analytical expression

$$\vec{S}(u, v) = \sum_{i=0}^n \sum_{j=0}^m b_i^n(u) b_j^m(v) \vec{P}_{i,j}$$

Note that Bezier curves and surfaces may be unstable – minor changes in control points might lead to major global changes.

2.4.2 NURBS basis functions

Extending the idea described in previous section, one could use *B-spline basis functions* (see below) instead of the Bernstein polynomial basis.

Unlike with Bezier curves, for the B-splines a parameter domain is subdivided by so-called *knots*. For the one-dimensional parameter domain $[u_0, u_m]$, the *knot vector* will be given by $u_0 \leq u_1 \leq \dots \leq u_m$. In most cases $u_0 = 0, u_m = 1$ is chosen, so that we get the unit interval for our parameter values. For the case of NURBS, the knots u_0, \dots, u_m need not be equidistant – hence the Non-Uniform in the beginning NU of NURBS.

Given a knot vector $[u_0, u_m]$ and a degree of B-spline p , the i -th B-spline basis function is then defined recursively as follows:

$$N_{i,0}(u) = \begin{cases} 1, & \text{if } u_i \leq u < u_{i+1} \\ 0, & \text{otherwise} \end{cases} \quad (2.1)$$

$$N_{i,p}(u) = \frac{u - u_i}{u_{i+p} - u_i} N_{i,p-1}(u) + \frac{u_{i+p+1} - u}{u_{i+p+1} - u_{i+1}} N_{i+1,p-1}(u) \quad (2.2)$$

For $p = 0$ we get just step functions, and $p = 1$ familiar hat functions, whereas the quadratic basis functions look more complicated (Figure 2.7).

By giving each of these basis functions a weight ω_i and normalizing them at each point by dividing by the total sum, we get the rational basis functions. Writing them out explicitly, in terms of B-spline basis functions $N_{i,p}$, the n^{th} -degree NURBS curve with k control points P_i is finally given by:

$$C(u) = \frac{\sum_{i=1}^k N_{i,n} \omega_i P_i}{\sum_{i=1}^k N_{i,n} \omega_i}. \quad (2.3)$$

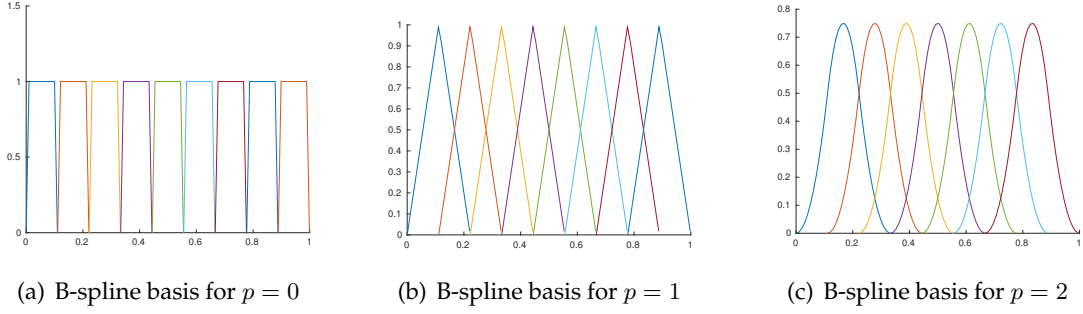


Figure 2.7: B-spline basis functions, of degree $p = 0$ (left), $p = 1$ (middle) and $p = 2$ (right).

B-splines have the following properties, which are useful for our problem:

- Degree n and number of control points $\vec{P}_{i \dots m}$ are independent.
- B-Splines only change locally (depending on the degree n) when a control point is changed.

Analogously to the Bezier curve surfaces, one can define B-spline or NURBS surfaces. For more information about NURBS see [14].

2.4.3 Fitting problem

In order to represent a given line or surface by Bezier curves or surfaces respectively, just as with NURBS, one has to solve a fitting problem. The goal of it is to fit in a parametric curve to the set of given data points. In the case of interest, the given set of points is a mesh, obtained from surface contouring.

Fitting problem: Bezier curve

First we want to find a Bezier-curve $\vec{B}_n(u)$ of degree n which is approximating a given spline $\vec{s}_m(u)$ defined by m points in an optimal way. For this purpose we want to minimize the L2-error. This leads to the minimization problem

$$\text{find } \min_{\vec{B}_n \in \mathbb{B}_n} \left\| \vec{B}_n - \vec{s}_m \right\|_{L_2}. \quad (2.4)$$

Minimizing the L2-norm is equal to minimizing the functional:

$$F(\vec{B}_n) = \int_{u=0}^{u=1} \left(\vec{B}_n - \vec{s}_m \right)^2 du \quad (2.5)$$

Using the variational principle we get the system of linear equations

$$Aa = b \quad (2.6)$$

where A is the matrix of pair-wise scalar products of basis functions (Bernstein polynomials), b is the vector of scalar products of the given spline s_m and basis functions and a is a required vector of coefficients for Bezier curve representation.

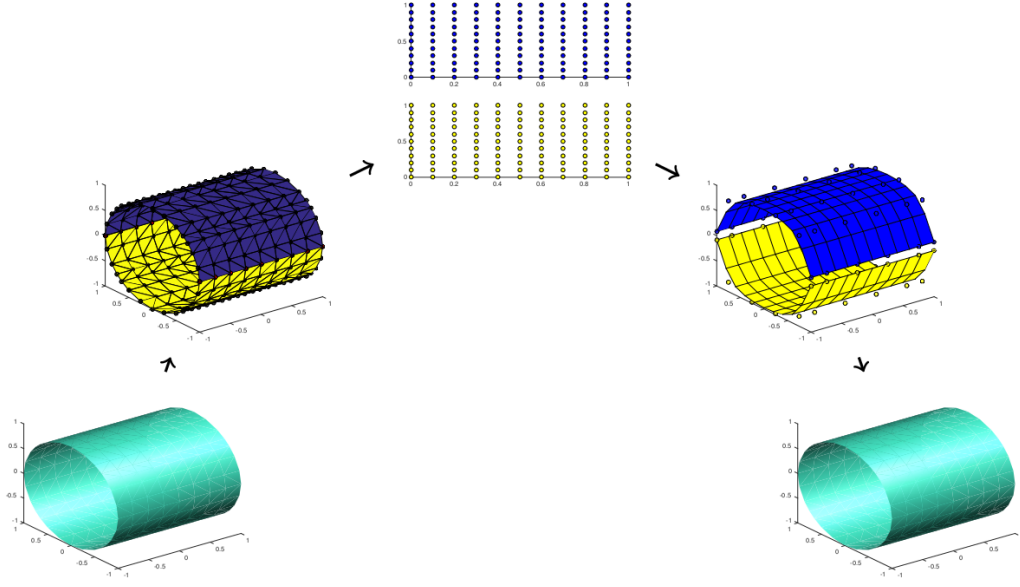


Figure 2.8: NURBS fitting pipeline

Fitting problem (least squares): NURBS

Although the approach used above showed good results, it appears to be very computationally intensive. Analogically, one could reduce the original problem to a *regression problem*, which allows us to reduce computational costs. Also, to improve the locality of our solution, from now on we are going to use NURBS basis functions with weights $\omega_j = 1$ instead of Bezier curves. For this purpose, we adopted an algorithm, provided in [15]:

Let X^0 be the $n \times 2$ matrix of the given set of points, N^p the basis functions of degree p ($n \times (n + p)$ matrix, where n is the number of points), and P^0 the control points ($(n + p) \times 2$ matrix).

The original problem can be written as:

$$X_i^0 = \sum_{j=1}^{n+p} P_j^0 N_{i,j}^p, \quad i \in \{1, \dots, n\} \quad (2.7)$$

Or, in short:

$$X^0 = N^p P^0 \quad (2.8)$$

The above system needs to be solved for the unknown P^0 . This can be done numerically, for example through singular value decomposition.

2.4.4 Fitting pipeline

Since the geometry obtained after topology optimization can be arbitrary complex, we might not be able to find a good fit using only one patch. We seek a multi step algorithm, allowing us to break the overall big problem into smaller problems, which can be handled relatively easy. Based on the algorithm described in [16], our overall fitting pipeline looks as follows (see Figure 2.8):

- Patch selection (breaking our problem in small pieces which can be solved using least squares)
- Parametrization of obtained patches
- B-spline fitting using least squares
- Smooth connection of patches
- Conversion back to CAD

The pipeline given above, once implemented, will provide us with a flexible algorithm for converting an arbitrary complex mesh based geometry into NURBS and, hence, CAD-representation.

3 Implementation

3 Implementation

3.1 From CAD to Voxels

One of the hurdles with most state-of-the-art open source topology optimization tools is their input format, where many of them (including ToPy, our topology optimizer of choice) require input to be specified as a 3-dimensional voxel grid. Presence (or absence) of material in these voxels is defined by a boolean variable, and boundary conditions are imposed on the appropriate locations. An example of voxelized data can be seen in Figure 3.1.

In the variety of toolboxes available for voxelization, we decided on the *Common Versatile Multi-purpose Library for C++* (CVMLCPP). This is a collection of mathematical algorithms whose objective is "to eliminate this redundancy by offering high-quality implementations of commonly needed functionality" [18]. The library offers an easy-to-use voxelizer, which we use for conversion of CAD input to a boolean voxel grid.

Another very popular open-source choice is the 3D modeling toolbox is OpenCascade [17], a versatile library with huge amounts of functionality. Although this also offers a voxelizer, we decided that its additional functionality did not justify its disadvantages in size and cumbersome installation requirements on for example a linux system.

In terms of implementation, the only thing required is the installation of the CVMLCPP library. The voxelizer is then included as a callable binary, that takes STL-file input (section 2.1.2) and converts it to a .dat binary file with dimension and voxel information, with specifiable voxel size. An example result of using the voxelizer is shown in Figure 3.2.

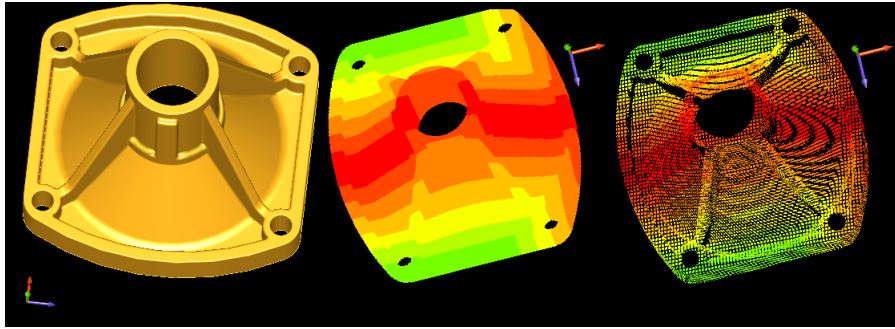


Figure 3.1: A shape and its voxel representation. *Leftmost picture:* The original parametrized shape. *Rightmost picture:* The voxel representation. Picture from OpenCascade [17].

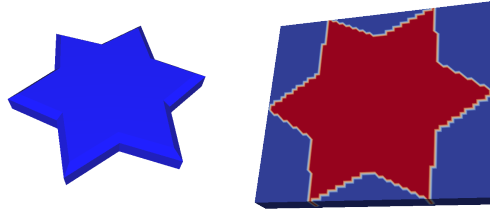


Figure 3.2: The STL geometry of a star (left) and its voxelized form (right) obtained via the CVMLCPP voxelizer, visualized by Paraview [19].

3.2 Topology Optimization

Due to the good range of topology optimization software available, we decided to adapt an open-source topology optimizer to our needs, namely *ToPy*.

ToPy [20] is a python library/program, written by William Hunter and documented in [5], implementing the SIMP model and method described in section 2.2. It is based on the 99-line Matlab code by Sigmund's for minimum compliance [7]. The program can optimize the previously named problem types: minimum compliance, heat conduction and mechanism synthesis—in 2D as well as 3D. It uses highly optimized open source python libraries such as Pysparse [21] and Numpy [22], leading to improved speed, porta- and scalability.

We use ToPy as a black-box topology optimizer. This means, we launch the program with an input file based on our scenario and let ToPy run. Next, as soon as the output is available, we proceed with the next step. The intention is to create separate modules to be able to plug in different solvers later on. In order to specify our input, we wrote an auxiliary program taking a voxelized CAD design provided by CVMLCPP as input, outputting a .tpd file with geometry, load and fixture information in the format required by ToPy.

Results of the topology optimization process can be seen in figure 3.3. Here, a star was given as input from a stl-file. We set the voxels in the star's points as fixtures, while we set a load in the middle, in the direction normal to the plane of the star. As can be seen, the optimization process "cuts" away unnecessary material in-between the corners and even in the middle of the material, returning an optimally stiff structure for the chosen remaining volume fraction.

3.3 From optimized voxels to surface representation

As was mentioned above, surface extraction is an intermediate step after Topology Optimization and NURBS representation in order to facilitate the conversion. In terms of implementing the surface extraction, we used VTK [23].

The VTK Toolbox is an open-source tool, providing algorithms for "3D computer graphics, image processing, and visualization" [23]. Among the variety of tools, VTK offers algorithms that allow us to obtain a surface representation from voxel data. Among these algorithms, we could find Marching Cubes, Dual Contouring and also a Decimation tool, the last of which can be useful for reducing the data size further for the NURBS-representation step.

The Marching Cubes implementation in VTK is however not applicable to the VTK *Unstructured Grid* type data that we obtain from ToPy, since it only works with *Image Data* and *PolyData*,

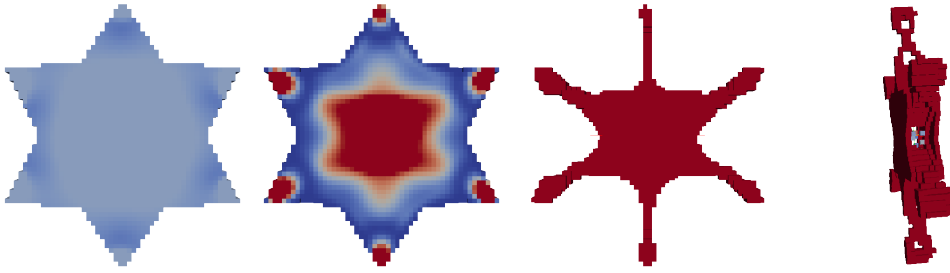


Figure 3.3: Topology Optimization by ToPy [20], with minimum compliance. *From left to right:* increasing number of SIMP iterations until convergence. The star-shaped structure was given by an STL-file which was processed into input readable by ToPy, with fixtures in the corners, and a load in the middle. Throughout the SIMP iterations, one can see how material from the less dense regions (blue) is concentrated into denser regions (red) that carry the load. The last picture gives a rotated view, to illustrate how material has been eliminated even from the inside of the star.

the main data types in VTK. For structured and unstructured grids the tool to render the iso-surface is the *Contour Filter* tool. Unfortunately, the documentation does not present which algorithm the tool uses. It can be inferred that it is an extended Marching Cubes algorithm. Although an implementation with Contour Filtering worked, the visualization of the data was still not possible making an intermediate step needed. Here, we used the *Implicit Modelling* tool, a filter that computes the distance from the input geometry to the points of an output structured point set. This distance function can then be "contoured" to generate new, offset surfaces from the original geometry. Although this approach allowed the visualisation, some crucial information was lost, as for example holes were not represented in the final model. This can be seen in Figure 3.4.

In order to reduce computational costs of the following NURBS fitting process (see section 2.4) we also need to create a coarser mesh of polygons from the fine one. The number of triangles that represent the isosurface can here be reduced with the VTK Decimation tool mentioned above. A smoothing step is however necessary in-between to get the new connections right. As can be seen in Figure 3.5, a 50 % reduction of the triangles barely provides a noticeable difference, and even with a 90 % reduction it is difficult to see a difference. Triangle meshes can be easily coarsened since there are many open source algorithms that reduce the number of triangles.

3.4 From extracted surface to NURBS representation: an algorithmic journey

As of yet, there is no open source software which provides the conversion from a *mesh-based* geometry to NURBS representation. Hence, one of the main challenges of both the algorithmic and implementation part of this project is to develop one from scratch. Due to a variety of possible approaches to tackle this problem (e.g. [15],[16]), we have conducted a profound prototyping work. In order to avoid a cumbersome and time-consuming implementation overheads during the prototyping phase, we have used MATLAB [4]. Once the algorithms to be used are finalized,

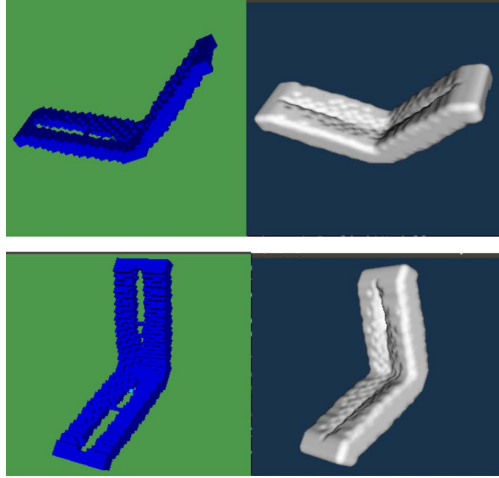


Figure 3.4: Contour Filtering tool after Implicit Modelling. The original geometry (left) is an optimized geometry output from ToPy (see section 3.2). Note that some essential features of the geometry after the process (right) are lost, such as the holes through the structure.



Figure 3.5: Decimation of triangles. *Top*: 50% reduction of triangle number from the object in Figure 3.4. *Lower*: 90% reduction. The difference is barely noticeable in both cases.

the prototypes are to be implemented using a non-proprietary language, such as Python or C++.

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