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Orthotropic $k$-nearest foams for additive manufacturing

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1 INTRODUCTION

Additive manufacturing (AM) enables the fabrication of objects embedding metamaterials. By creating fine-scale structures, the object’s physical properties can be graded (e.g., elasticity, porosity), even though a single base material is used for fabrication. Designing the fine and detailed geometry of a metamaterial while attempting to achieve specific properties is difficult. In addition, the structures are intended to fill comparatively large volumes, which quickly leads to large data structures and intractable simulation costs. Thus, most metamaterials are defined as periodic structures repeated in regular lattices. The periodicity simplifies modeling, simulation, and reduces memory costs—however it limits the possibility to smoothly grade properties along free directions.

In this work, we propose a novel metamaterial with controllable, freely orientable, orthotropic elastic behavior—orthotropy means that elasticity is controlled independently along three orthogonal axes, which leads to materials that better adapt to uneven, directional load scenarios, and offer a more versatile material design primitive. The fine-scale structures are generated procedurally by a stochastic process, and resemble a foam. The absence of global organization and periodicity allows the free gradation of density, orientation, and stretch, leading to the controllable orthotropic behavior. The procedural nature of the synthesis process allows it to scale to arbitrarily large volumes at low memory costs.

We detail the foam structure synthesis, analyze and discuss its properties through numerical and experimental verifications, and finally demonstrate the use of orthotropic materials for the design of 3D printed objects.

CCS Concepts: • Computing methodologies → Shape modeling.

Additional Key Words and Phrases: 3D printing, additive manufacturing, procedural modeling, material design, metamaterials

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of designer Lilian van Daal\(^1\). Unfortunately, such metamaterials are much more challenging to synthesize and analyze, and to the best of our knowledge no works have yet focused on automatically synthesizing orientable, orthotropic graded materials that can be fabricated.

We propose a novel technique to produce microstructures that result in a controlled, freely-orientable orthotropic elastic behavior (see Figure 1). The structures are efficiently generated by a procedural function, without the need for storing or simulating their entire geometry.

Our main contributions are:

- The definition of an orientable, orthotropic foam well suited for modeling and additive manufacturing.
- An efficient procedural evaluation of the microstructures, which is remarkably simple to implement.
- The analysis of the relationship between procedure parameters and elastic properties.

We demonstrate several applications of orthotropic graded foams on 3D printed objects.

2 RELATED WORK

We start by discussing techniques which optimize for material properties within objects in Section 2.1. The output of these methods can be used as a control field for the parameters of a (meta)material defined by a microstructure. We discuss in Section 2.2 microstructures for material design, and in Section 2.3 methods that consider modeling and computational challenges of infilling 3D shapes with microstructures. Finally, we discuss in Section 2.4 approaches that optimize internal structures within objects with similar objectives of rigidity and weight reduction.

2.1 Optimizing for material properties

The field of topology optimization is dedicated to the optimization of material properties within a domain [Eschenauer and Olhoff 2001; Rozvany 2001]. The methods optimizing for continuously spatially-varying material properties (e.g. density) are of direct interest to us [Allaire 2012; Bendse and Kikuchi 1988]. After optimization, the question of how to map the continuous, homogenized properties to an actual material arises. Microstructures are one possible approach to this problem [Bendse and Kikuchi 1988; Suzuki and Kikuchi 1991]. It is worth noting that topology optimization via homogenization optimizes for a full elasticity tensor at every point (e.g. free material optimization [Bendse et al. 1994; Kočvara and Stingl 2007], laminate optimization [Allaire and Francfort 1993]), and thus anisotropy information is available. In practice, it is difficult to physically produce materials with this type of control, and this information is often unexploited. One exception is the work of Pantz and Trabelsi [2008], which maps non-regular lattices onto the homogenized solution of a compliance minimization problem. This however requires a global optimization over the domain. Our goal is to provide a metamaterial capable of producing an orthotropic elastic behavior while being simple to conform to a control field – e.g. without having to solve for a global mapping problem.

In computer graphics there is also a strong interest in fabricating deformable, animated objects, as well as interactive editing of such designs. Bickel et al. [2010] design materials with prescribed deformation from a set of predefined base materials. Chen et al. [2013] propose a framework for the optimization of heterogenous multi-material arrangements achieving specific properties. Skouras et al. [2013] optimize for a bi-material distribution to achieve a desired deformation under imposed displacements. Xu et al. [2015] optimize for distributions of different isotropic materials to meet displacement and internal force constraints. Ion et al. [2016] provide an advanced interface to allow users to paint material properties and create mechanisms embedding microstructures.

2.2 Microstructures for (meta)material design

The main approach to make microstructure design tractable is to rely on periodic structures [Sigmund 1994, 1995; Sigmund and Torquato 1999]. This offers two significant advantages. First, the periodicity allows storing in memory a single tile, that is implicitly repeated in a regular grid covering the object. This enables compact storage, efficient display and processing by the AM device (see Section 2.3). Second, the theory of homogenization affords for the computation of the parameters of an equivalent homogeneous material [Allaire 2012]. Given these, the object can be simulated while abstracting away fine scale heterogeneities.

The base tile of a periodic microstructure can be optimized through an inverse homogenization problem to target a specific material [Andressen et al. 2014; Radman et al. 2013; Sigmund 1994; Xia and Breitkopf 2015; Zhou and Li 2008]. Schumacher et al. [2015] extended this methodology to optimize families of periodic tiles with varying properties. Panetta et al. [2015] proposed a family of isotropic tiles based on periodic truss structures. In both these works, tiles with different properties can be spatially arranged to grade properties, such as obtaining varying degrees of elasticity. This requires a special treatment of the boundaries, either ensuring the tile borders are compatible across the entire tile set [Panetta et al. 2015], or performing a global optimization step to choose tiles with best matching borders [Schumacher et al. 2015]. Whenever using a periodic grid care must be taken where the grid intersects the object surface, as tiles are cut by the surface [Robbins et al. 2016].

While extremely efficient and widely adopted, periodic microstructures have a number of disadvantages. The underlying regular grid makes it difficult to smoothly grade the structure properties along arbitrary fields. In particular, when the material is orthotropic, the grid makes it difficult to follow an orientation field seamlessly across grid cells.

Martinez et al. [2016] deviate from periodic microstructures and propose to generate aperiodic, stochastic open-cell foams with prescribed isotropic elasticity, through a procedure akin to solid procedural textures. This builds upon well known properties of open-cell foams [Gibson and Ashby 1997; Roberts and Garboczi 2002], the Young’s modulus of which is strongly correlated to their volume (e.g. denser foams produce stiffer materials following a quasi-linear relationship). We discuss in Figure 3 possible ways to obtain an

\(^1\)Chair by designer Lilian van Daal: [url]

The chair was manually modeled in Rhinoceros, which took hundreds of hours, see [here](#) and [here](#).
orthotropic behavior from such foams, however none proved efficient. Another possibility, not in the Figure, is to rely on anisotropic Voronoi diagrams (e.g. [Lévy and Bonneel 2013]). For instance, Nervous System uses such diagrams on surfaces to model 3D printed jewelry. However, extraction of curved cell boundaries (edges in 3D) of anisotropic diagrams remains computationally expensive. Disconnected islands may also appear under varying metrics. In comparison, our approach is simple to implement, procedural, and produces structures made of straight beams that exhibit a controlled orthotropic behavior.

2.3 Efficient infilling with microstructures

Given a specific microstructure geometry the question of how to fill a target shape with this detailed geometry arises. Wang et al. [2005] conform truss lattices in a thick shell below an object’s surface through an efficient procedure that directly outputs a STL model. Rosen [2007] considers the challenges of designing parts with internal lattices and proposes a slicing procedure which works directly from the truss skeleton. Chen [2007] extends texture mapping and signal specialized parameterization to infill volumes with density-varying truss lattices. Brennan-Craddock [2011] studies several approaches to apply microstructures within objects, and proposes a slicing algorithm exploiting the periodicity of the structures for efficiency. Pasko et al. [2011] and Fryazinov et al. [2013] explore procedural formulations of periodic microstructures. This enables efficient visualization through ray-tracing, and efficient fabrication by streaming slice images to the printer. Li et al. [2015] rely on such microstructures to optimize the internal cross-sections of an object. Vidimé et al. [2016; 2013] explore a voxel approach to define complex procedural structures within objects, with an emphasis on multi-materials, usability, and slicing efficiency.

Our approach answers the requirements that provide efficient procedural synthesis and direct slicing for scalability of manufacturing. Compared to lattices, we propose a significantly different viewpoint by considering stochastic structures, that do not require solving for a mapping to conform to a vector field. Despite its randomized nature, our approach defines a metamaterial in a principled way, and affords for a precise control of the orthotropic behavior.

2.4 Globally optimized internal structures

A number of approaches optimize internal structures globally. Lu et al. [2014] optimize closed-cell foams formed by the faces of a Voronoi diagram. Wu et al. [2016a] solve for a global, high-resolution topology optimization problem under local material density constraints, which produces porous structures. Wu et al. [2016b] subdivide rhombic structures in a global optimization process, using smaller and denser cells in regions of high compliance. Other approaches optimize sparse truss structures within the volume of objects, such as to obtain rigid but lightweight 3D prints [Wang et al. 2013; Zhang et al. 2015]. Recent software for AM, such as Autodesk Netfabb, Autodesk Within or nTopology Element, provide...
In this work we always use beams of minimal printable thickness. For the same volume this lets us maximize foam density.

Point distribution. Our structures rely on an isotropic point distribution similarly to procedural Voronoi foams [Martínez et al. 2016]. We generate one random point in each cell of a virtual grid covering the space, which provides a crude but efficient approximation of a Poisson disc distribution [Wolsey 1996]. The subdivision scheme of Martínez et al. [2016] is used to locally increase or decrease the point density.

Edge connections. In absence of anisotropy, we generate the edges of the graph by connecting each point to its k closest neighbors. For sufficiently large values of k (as discussed later), this produces fully connected isotropic structures.

The edges of the graph are the essential component in inducing anisotropy in the elastic behavior of the structures. As seen in Figure 3 (top), a global stretch of an isotropic graph introduces an excellent orthotropy in its elastic response. The effect of the stretch is to bias the angular distribution of edge lengths, making the equivalent material stiffer along the direction with longest edges. Our key insight is to produce a similar bias, and hence a similar elastic response, without having to rely on a global stretch. Instead we introduce local stretches that can be easily controlled. In the limit our approach converges towards an ideal laminate (like a global stretch), which reach orthotropic elasticity bounds [Lipton 1994].

We produce the angular edge length bias by stretching the distance computation when selecting the k nearest neighbors. This encourages longer edges to appear in the direction of largest stretch, as k-nearest neighbors are selected further away (see the inset below and Figure 4). Section 5 provides an in-depth analysis of the impact of the stretch of metric on the orthotropic elastic behavior.

More precisely, let us consider two points pi, pj in the embedded graph in dimension d. We denote the frame orientation at point p as Θ(p) : R^d → [0, π]^d, with α = {1, 3} for respectively the 2D and 3D cases (Euler angles). We denote the stretch at p as H(p) : R^d → R^d, which, without loss of generality is assumed to be an ordered length vector. In 2D (h_u, h_v) with h_u ≥ h_v and h_v = 1. In 3D (h_u, h_v, h_w) with h_u ≥ h_v ≥ h_w and h_w = 1.

We define the metric tensor as [Du and Wang 2005]:

\[ M(p) = E^T U E, \quad E = \text{diag}(H(p)^{-2}) \]  

where R(Θ) is the n dimensional clockwise rotation matrix.

The asymmetric anisotropic distance from pi to pj is:

\[ d_{pi}(p_1, p_2) = \sqrt{(p_1 - p_2)^T M(p_1)(p_1 - p_2)} \]  

In order to simplify the synthesis of the structure, we define a symmetric distance d(pi, pj) as:

\[ d(p_i, p_j) = \frac{d_{pi}(p_1, p_2) + d_{pj}(p_j, p_i)}{2} \]  

This distance is used to select the k nearest neighbors, producing an anisotropic distribution of edges. We here conjecture that this translates to an orthotropy in the elasticity tensor. This, however, is
not obvious. We verify it is the case in Section 5. We also quantify the relationship between the metric anisotropy and the impact on the Young’s moduli and Poisson’s ratios in each direction.

Connectivity. A major concern for fabrication is to obtain fully connected structures. Studies in graph theory indicate that $k$-nearest random graphs are connected with very high probability for even small values of $k$ [Balister et al. 2005]. Our case is, in fact, more favorable: when using a point distribution similar to Worley [1996] one can easily guarantee that at least one sample exists in each virtual grid cell. In practice we use $k = 6$, which with uniform density always achieves connectivity even with strong anisotropy.

We experimentally evaluate the impact of anisotropy on connectivity by producing random instances of the foam in a finite, large domain. We assume the structure to be attached to the boundary (i.e. the surface of the object), and track for any disconnected component inside. We use a square domain of unit size covered by a grid of $32^2$ points, and a maximum tested stretch ratio of $h_u = 80$. The maximum edge length was measured at 0.7. Running $7k$ tests on stretch ratios from 1 to 80 did not produce any disconnected component.

While we did not observe any detrimental effect of anisotropy on connectivity, disconnections may occur when varying density. In fact, following the first steps of the proof in Xue and Kumar [2004], we can easily setup a “trap” as illustrated in the inset. Such a trap consists in sandwiching a narrow band of low density in between two bands of much higher density. As a result, the nodes from the coarse band will take their $k$-nearest neighbors in the high-density bands only, creating a gap in connectivity. Figure 5 reveals this behavior on our foams. Such cases seldom occur in practice as the density control fields are typically smooth; if needed the connectivity of the structure can be checked efficiently through a sweeping scheme (line in 2D, slice in 3D) as the structure synthesis is procedural. Larger values of $k$ also alleviate this issue.

Locality. Locality is an important property for the efficient evaluation of the foam geometry: we seek for an algorithm where a sub-set of the foam can be generated without having to globally synthesize the result. To achieve this, the search radius (with the usual $L_2$ distance) to find the $k$-nearest neighbors of a point has to be bounded.

Since we guarantee one sample per jittered grid cell, such an upper bound exists. Finding a tight upper bound is however difficult. We need to find the smallest radius that will contain at least $k$ points – this relates to the Gauss circle problem generalized for ellipses [Bentkus and Götze 1997; Hardy 1915], but is not strictly the same as we rely on a jittered grid. Nevertheless, from these works we can derive a lower bound (the radius that must at least be searched) that is given in 2D by $c \cdot h_u \sqrt{k/(\pi \cdot h_u \cdot h_v)}$, and in 3D by $c \cdot h_u \sqrt{k/(\pi \cdot h_u \cdot h_v \cdot h_w)}$, where $c$ is the grid cell size length. In practice, we observe that the lower bound follows a trend similar to the experimental upper bound, see Figure 6. We therefore set the upper bound proportional to the lower bound, in practice using a factor of 2. As shown Figure 6 the 2D upper bound remains well above the measured edge lengths for the stretch ratios we use. During evaluation we use the bound obtained from the maximum stretch in the control field.

4.2 Procedural evaluation of orthotropic foams

Given the properties of connectivity and locality, we devise a procedural evaluation scheme. A procedural solid texture function, which returns either empty of solid for any coordinate is possible. However
it would be computationally wasteful due to the repeated $k$-nearest searches. Instead, we propose a per-block synthesis scheme that groups computations in a local region of space.

Let us assume our point distribution algorithm is pseudo-random and can generate point-sets deterministically in any box in space. This is the case of the jitted grid approach we favor. Let us consider an axis aligned box with min corner $c_{\text{min}}$ and max corner $c_{\text{max}}$. To produce the foam geometry within the box $[c_{\text{min}}, c_{\text{max}}]^d$, we generate the point distribution in an enlarged box $[c_{\text{min}}, c_{\text{max}}]^d @ L_{\text{max}}$ where $@$ is the morphological dilation and $L_{\text{max}}$ is the bound on maximum edge length, computed from the maximal stretch in the field. This guarantees that all the points required for the $k$-nearest search within the smaller box $[c_{\text{min}}, c_{\text{max}}]^d$ are available within the enlarged box. Thus, the set of edges matches the set that would have been generated globally. In other words, this evaluation is deterministic: generating the foam in overlapping boxes produces the same geometry within the overlapping regions. This process and its results are shown in Figure 7.

Using this approach, we can efficiently generate the foam in any spatial box, in a time that only depends on the box extent (the foam domain being infinite). In particular, this can be used for front-to-back rendering during visualization, or during slicing for AM where only a thin slab of foam has to be computed. Thus, the set of edges matches the set that would be computed using correct computations inside. Right: Overlapping the two blocks produces a seamless result: their content is deterministic and exactly matches in the overlap region. The space spawned by the procedure is infinite.

5 ANALYSIS OF ORTHOTROPIC FOAMS

For the foams to be useful as a metamaterial it is important that the equivalent material properties correlate to the synthesis parameters. If this relationship exists and can be characterized, one can easily fill a volume with a foam that produces a target elastic behavior.

We use numerical homogenization (Section 5.1) to verify the link between the anisotropy of the foam geometry and the orthotropy of the elasticity tensor (Section 5.2). We then analyze the material space spawned by the foams (Section 5.3) for varying synthesis parameters. We conclude by mechanical tests on printed samples (Section 5.4). We provide additional analysis in supplemental material for homogenization versus detailed simulation and for the isotropic case.

In the remainder, we consider an isotropic base material of Young’s modulus $E = 1$ (without loss of generality since the Young’s modulus values are linearly proportional to that of the base material). We consider a base Poisson’s ratio of $\nu = 0.3$ which is representative of plastic materials.

5.1 Homogenization

We analyze the macroscale behavior using numerical homogenization in grids of hexahedral (3D) or quad (2D) linear elements. This computes the tensor of an equivalent homogeneous material for a given sample of our foam. We provide additional details on numerical homogenization in the supplemental material.

Similarly to Martinez et al. [2016], we generate periodic versions of the foams in a grid of varying spatial extent, and apply numerical periodic homogenization to each. We fit the homogenized elasticity tensor $C_{\text{hom}}$ to an ideal orthotropic tensor $C_{\text{ortho}}$ of orientation $O$ (see Section 3), which retrieves both the orthotropic parameters and a fitting error. We increase the spatial extent until the fitting error is negligible, ensuring enough foam is observed to obtain a reliable result (grid of $300^2$ elements in 2D, $80^3$ in 3D). For now, let us assume the orthotropic orientation $O$ is known. We discuss its selection in Section 5.2.

To perform the fitting we consider the logarithmic Euclidean distance. This has the desirable property of evaluating to the same value for both the elasticity and compliance tensors, and provides more accurate results for elasticity compared to the Frobenius norm [Moakher and Norris 2006]. The logarithmic Euclidean distance for the elasticity tensor is:

$$d_L(C_{\text{ortho}}, C_{\text{hom}}) = \|\log(C_{\text{ortho}}) - \log(C_{\text{hom}})\|_F$$

The optimization is constrained such that $C_{\text{ortho}}$ is positive definite [Ting 1996]. We use a gradient-based optimization with nonlinear constraints [Johnsen 2016]. Since the objective function involves matrix logarithms, its gradient is computed with numerical differentiation.

5.2 Orientation conjecture

We generate anisotropic foam geometries by locally stretching the metric used to connect a point to its $k$-nearest neighbors (Section 4.1). In this section we verify that the geometric stretch translates to an orthotropy of same angle in the elasticity tensor. We perform the analysis in 2D.

To observe this, in particular on low stretches, we setup the following numerical experiment. We generate 6800 foam samples for stretch values $h_{\text{u}}$ in $[1.5, 2.5]$, where $h_{\text{u}}$ aligns with the horizontal axis, $h_{\text{v}}$ with the vertical axis. We fit an orthotropic tensor to each, rotating the expected angle of orthotropy $O$ (Section 5.1), and considering the fitting error. If the material is indeed orthotropic with direction $O$, then the fitting error is expected to be small. Otherwise, the fitting error is expected to increase as the directions of orthotropy do not match.
Rather than using absolute values of fitting error, which are difficult to interpret, we consider the normalized measure proposed in [Moakher and Norris 2006]:

\[ d_N = \frac{d_1(C_{\text{ortho}}, C_{\text{iso}})^2}{d_1(C_{\text{hom}}, C_{\text{iso}})^2} \in [0 \ldots 1] \quad (7) \]

where \( C_{\text{iso}} \) is a fitted isotropic tensor (obtained by the fitting process of Section 5.1, using an idealized isotropic tensor, see supplemental material), and \( C_{\text{ortho}} \) is the fitted orthotropic tensor. A value of one indicates that the fitted tensor is indeed close to ideally orthotropic, while zero indicates the opposite. A value above 0.9 is considered to be a good agreement [Moakher and Norris 2006].

We perform this analysis for three different uniform point distributions: uniform random coordinates, Poisson disc, and the jitted grid approach we use in practice. Results are reported in Figure 8. As can be seen, for angles where \( O \) does not align with the metric stretch the error increases (lower values of \( d_N \)), while the agreement is excellent when axes align. This verifies that the stretch in the metric translates to a similarly aligned orthotropy in the material.

Another interesting observation is to consider how randomness impacts the properties (i.e. the spread of points around the average behavior). The spread is wider for a uniform random coordinate distribution, and narrower for Poisson disc, while the jitted grid lies in between. The spread of the jitted grid actually narrows on larger stretch values, and becomes similar to that of the Poisson disc distribution. This indicates that there is little difference in quality between using a jitted grid or a Poisson disc distribution, especially with high orthotropy.

### 5.3 Material space

We now analyze the space of materials covered by our technique. We generate a large number of orthotropic 2D samples, and evaluate each with numerical homogenization. We vary the stretch \( h_u \) (up to 15 to cover a wide range) and the density. The orientation is fixed to zero since we verified numerically in Section 5.2 the orientation conjecture. We exploit symmetries in orthotropy when exploring the material space. The samples all have the same spatial extent but are randomized: the dataset may contain samples with similar parameters, but they are different random realizations.

The results for the 2D foams are summarized in Figure 9. The variations in normalized Young’s modulus along each axis are shown in the top part. An isotropic material would only cover a diagonal, as the orthotropic axes would have the same Young’s modulus. Our structures cover the full spectrum of Young’s moduli and ratios between them. The main factor in controlling the ratio between the Young’s modulus is, as expected, the metric stretch. The main factor in controlling the absolute Young’s modulus values is the density.

The plot at the bottom of Figure 9 reveals the behavior of the Poisson’s ratio together with the Young’s modulus. Figure 9 plots only \( \nu_{xy} \), as \( \nu_{yx} \) follows from Equation (2). The Poisson’s ratio remains positive, and sharply increases on small volumes and high stretches. Figure 9 contains a foam representative of this case (\( \nu_{xy} = 5.16 \)). Since the Poisson’s ratio of orthotropic materials is unbounded [Ting and Chen 2005] large values are to be expected.

#### Figure 8. Numerical verification of the material orthotropy for uniform random coordinates, jitted grid, and Poisson disc point distributions. The normalized measure \( d_N \) is plotted against stretch angle \( O \) (polar axis in degrees), and used as color code (bottom colorbar).

Figure 10 presents the analysis of Young’s moduli on 3D foam samples (cubes). As can be seen, the foams provide a good coverage of possible ratios between \( E_x, E_y, E_z \). We limited the exploration to a stretch of at most 14 in all directions, but the space could be explored further. The Poisson’s ratio behaves as with 2D foams. Please refer to the supplemental material for the plots of the 2D and 3D shear modulus, and 3D Poisson’s ratio.

In summary, our foams cover the Young’s modulus space densely. However, for a fixed orientation of orthotropy, our foams have only two (2D) or three (3D) independently controlled elasticity parameters, so they do not offer a complete control. For instance, in 2D, choosing both Young’s moduli fixes the Poisson’s ratio and shear modulus. Varying \( k \) changes the stiffness for a same density, but does not significantly change the covered material space.

The user selects a foam by specifying a desired pair (2D) or triplet (3D) of Young’s modulus. The foam parameters are obtained by interpolating from the three closest samples in our dense datasets (distance weighted interpolation). If any of the three is further than a threshold derived from sampling density, the point is rejected as not covered by the foam pre-computed material space. Projecting to a closest match is left for future work. In 2D, for 13000 randomly tests we obtain an average relative Young’s modulus error of 0.84%, and in 3D for 350 tests an error of 2.06%. The higher error in 3D is

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**Table 1:**

<table>
<thead>
<tr>
<th>Uniform</th>
<th>Jittered grid</th>
<th>Poisson disc</th>
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<td>( h_u = 1.5 )</td>
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<td>( h_u = 2.5 )</td>
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5.4 Measurements on printed samples

We verify our numerical analysis on actual 3D printed samples. We focus this analysis on 2D versions of the foam, which are vertically extruded to obtain a 3D solid object. We print the structures using SemiFlex filament on an Ultimaker 2+ with a 0.25 mm nozzle, with a custom slicer. Each beam is exactly one filament wide. Please refer to the supplemental material for details.

We printed a family of samples with varying degrees of metric stretches. We additionally print a full sample, which acts as a reference for the (3D printed) full base material (cross-hatched, 100% infill). The test specifications and results are summarized in Figure 11. Figure 11a shows the raw data measured from the sensor. For fitting the linear material parameters we only use the first half millimeter of displacement (from 0 to 0.5 mm), as the material enters non-linear elasticity beyond. Note that the non-linear behavior is only a very gently bent, quasi linear curve.

Figure 11b compares the normalized predicted Young’s modulus to the measured one, relative to the reference. That is, we divide each measured sample Young’s modulus by the Young’s modulus measured on the reference specimen. The overall behavior matches remarkably well, as the shapes of the curves closely correspond. The vertical offset between the measure and the prediction indicates that our samples are overall stiffer than expected. This is most likely explained by the fabrication process, and in particular by the imperfect infilling of the fabricated reference (changes to the reference stiffness moves the curve up/down).

We provide in the graph both numerical predictions (homogenized and detailed simulations). It is expected that the homogenized material is slightly different since it considers an infinite periodic medium. The close agreement between them shows that the samples are sufficiently large to capture the macroscale foam behavior.

We report the result of repeated compression tests in Figure 12. Compared to tensile tests, under compression the foams enter a non-linear regime after less deformation, which is due to the buckling of elongated edges. Nevertheless, all specimens exhibit a similar non-linear behavior that follows the same proportions as their Young’s modulus ratios. Repeated testing of the same specimens reveal that after the first few tests the behavior changes in the non-linear regime — while it remains very similar in the linear one. Interestingly, subsequent tests behave more stably, indicating a burn-in behavior that converges towards a stable response. This could be explained by failures in places of high local stresses, which no longer influence the behavior after the first few compressions.
Fig. 11. Uniaxial tensile tests on a MTS® 4/ML testing machine, with a force sensor of 100 N, displacement speed of 1 mm/ min. The specimens have a dimension of 18 × 108 mm and thickness 3.2 mm. Nine of them correspond to varying ratios of stretch $h_x / h_z$, one is the reference base material. (a) Measured force/displacement curves (raw data). (b) Predicted (homogenized and detailed simulation) vs measured normalized Young’s modulus. (c) Specimen with $h_x / h_z = 2$ being tested.

Fig. 12. Repeated compression tests with three prints having the same dimensions (50 × 50 × 7 mm) and using the same point density. The stretch varies in the vertical direction ($h_x = 1, 3, 5$). All the tests are performed on an Instron 3345 testing machine, with a force sensor of 500 N. The maximum compressive strain is 8% (4 mm), attained at a speed of 1 mm/ min. Each print is repeatedly tested five times, waiting for two minutes between each test. Please refer to the text for discussion. Left: The curves revealing the measured forces as the compressive strain increases. Short dashes correspond to first tests, while longer dashes indicate later tests. The vertical dashed lines corresponds to 1.6% and 4.8% strain. Right: From left to right, photos for increasing compressive strains.

We report in Figure 13 compression tests on 3D foam specimens.

6 APPLICATION TO ADDITIVE MANUFACTURING

In this section we discuss the use of our foams for the design of 3D objects with orthotropic elasticity. Figure 1 (left) illustrates an orthotropic chair design inspired from the work of designer Lilian van Daal. The initial foam design is two dimensional and generated within the chair profile. It is then extruded into the final object, producing orthotropy in the seat and the back.

Figure 1 (right) is an illustrative model of a gear made lighter by an orthotropic foam oriented against the surface, keeping it resistant to pressure. Figure 14 is a tree model with a foam oriented vertically, along the inner distance field, showing how longitudinal strength can be preserved. Figure 10 (right) illustrates a 3D printed cube of orthotropic foam, printed on a B9 Creator resin printer. The cube is 35 × 35 × 35 millimeters with Young’s moduli ratios of $E_z / E_y = 20.5, E_y / E_x = 8.4$, for stretch $h_x = 1.0, h_y = 2.0, h_z = 6.0$. Figure 15 show two cylinders filled with our foams and having different orthotropic behaviors. In each cylinder, the local stretch rigidifies one direction while making the other two more flexible.

Regarding results with deformations, please keep in mind our analysis is performed on linear elasticity (small deformations), and
Fig. 13. Compression tests of 3D $k$-nearest foams with a beam radius of 0.25 mm within cubes of $20 \times 20 \times 20$ mm, printed on a B9Creator (DLP resin printer). To minimize errors due to the fabrication process all specimens were printed together and measured in one session. The stretch along the compression direction is different for each specimen ($h_u = 1, 1.5, 3, 6$), all other parameters are equal ($h_v = 1$). Tests are performed on an Instron 3345 testing machine. The compression speed is 0.5 mm/min, the maximum compressive strain is 1.5% using a high precision 50 N sensor. (a) Force measured as a function of relative compressive strain. (b) Pictures of the foams (front view, initial position). (c) The first two columns report the Young’s modulus ($E_z$) for each specimen, for simulation (relative) and experiment (MPa). For comparison the two last columns show the ratios over the average value, for each column (i.e. how much more rigid/flexible each is with respect to the average). As can been seen the ratios agree well overall.

Fig. 14. 3D printed tree with internal orthotropic foam. The foam is oriented vertically along the surface.

Fig. 15. 3D printed cylinders with orthotropic material. Top: The center segment is made more rigid horizontally and flexible vertically. Bottom: The extremities are isotropic, the center is rigid along the main axis but flexible radially.

While this case is essentially illustrative, we believe future integration of our foams within free material optimization methods will allow to further exploit the material space.

---

$\text{Stretch } h_u$
- 6.0
- 3.0
- 1.5
- 1.0

Force (N)
- Compressive strain (%)
- Stretch $h_u$
  - 6.0
  - 3.0
  - 1.5
  - 1.0

Simulation (relative)  Experimental (MPa)  Simul. (ratio)  Exp. (ratio)

<table>
<thead>
<tr>
<th>$h_u$</th>
<th>Simulation (relative)</th>
<th>Experimental (MPa)</th>
<th>Simul. (ratio)</th>
<th>Exp. (ratio)</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>0.006</td>
<td>0.366</td>
<td>x0.16</td>
<td>x0.13</td>
</tr>
<tr>
<td>1.5</td>
<td>0.014</td>
<td>0.707</td>
<td>x0.36</td>
<td>x0.26</td>
</tr>
<tr>
<td>3</td>
<td>0.041</td>
<td>3.429</td>
<td>x1.08</td>
<td>x1.25</td>
</tr>
<tr>
<td>6</td>
<td>0.091</td>
<td>6.469</td>
<td>x2.39</td>
<td>x2.36</td>
</tr>
</tbody>
</table>

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Therefore we cannot predict large deformation behaviors accurately. Nevertheless, the designed structures trigger the expected global deformation. Under small deformations the analysis is accurate (see Figure 11 and 13).

Figure 16 illustrates how our foams can be used in conjunction with the results of topology optimization. We run the SIMP method [Bendsøe 1989] using a penalty of $p = 1$ (continuous material distribution) and a volume constraint of 40%. We then extract the stress field from the FEM simulation. Material density, principal stress direction and magnitude are mapped to the foam parameters. While this case is essentially illustrative, we believe future integration of our foams within free material optimization methods will allow to further exploit the material space.
Density (SIMP)

while quite large, remains positive. Whether or not stochastic structures (extruded 2D foams) and powder based systems (SLS). While we did interest materials.

porous structures in a multi-material setting, with beams in different angles, metric stretch, and density. The foam scales trivially through procedural generation.

While our foams can be used for a wide range of design tasks and to interpret the results of topology optimization, we believe a tighter integration can be achieved with free material optimization methods [Bendse et al. 1994; Kočvara and Stingl 2007], constraining the elasticity tensor to remain in the covered material space.

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