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## Hierarchical mesh-to-points as-rigid-as-possible registration

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#### ABSTRACT

Surface registration is a fundamental problem in computer graphics and computer-aided design. The problem consists in finding a deformation from one surface to another that preserves some properties. For instance, in our inverse engineering context, we aim at finding the best, as isometric as possible, map between an input triangular model, and a large point cloud acquired on the actual mechanical part being processed. Existing solutions are not able to handle very large models with a good level of precision. We propose a method which is accurate and fast. Our solution combines an efficient iterative energy minimization scheme on a hierarchical decomposition of the problem geometry. Our experiments show that we obtain a fast and efficient algorithm compared to the state-of-the art method, while keeping its numerical accuracy.

#### 1. Introduction

Surface registration is a fundamental problem in computer graphics and computer-aided design. Given two surfaces, called source and target, the objective of the registration is to find a deformation that maps from the *source* into the *target* satisfying some properties (isometric, conformal, one-to-one, partial...). 6 This is a common problem, for example in 3D scanning, where multiple datasets captured from different viewpoints must be registered. In geometry processing, a mapping between the source and the target is for example needed in post-processing 10 such as reconstruction, morphing or information transfer. 11

This is also a key problem in inverse engineering where CAD 12 models are compared to measure their real counterparts. In 13 this context, real objects are captured live using laser range or 14 probing devices. Hence, the source is a CAD model (e.g. a 15 clean triangular mesh), and the target a large, possibly noisy, 16 point cloud. In sheet metal manufacturing, which motivates 17 the present research, the springback problem [1], caused by the elastoplastic material behavior of sheet metals, falls into reg-19 istration problems where the CAD model source must be reg-20 istered into the measure of the real model *target* before being 21 processed by the device (e.g. for machining). The springback 22 being mainly bending, a classical assumption is that tangential 23 shearing or stretching are neglected and therefore the deforma-24 tion can be considered as non-rigid and almost isometric. This 25 defines the class of maps between the source and the target we 26 are looking for. 27

The computer graphics literature contains several non-rigid registration techniques, see *e.g.* [2] for a survey. Despite the 34 quality of the results of such approaches, they may not be suited 35 to handle very large models as it is the case in the production 36 engineering field. The CAD literature proposes fewer solu-37 tions (see Section 2). Even though they are more suited for 38 our problem, they are still not fast enough to deal with very 39 large datasets. One common remark about the non-rigid isomet-40 ric registration solutions is almost all of them register mesh to 41 mesh instead of mesh to point cloud. Therefore they all require 42 a reconstruction preprocessing on the point cloud to extract a 43 mesh. 44

In this paper, we extend computer graphics techniques and 45 propose an accurate and efficient algorithm to compute an as-46 rigid-as-possible map from a 3D CAD model to a large (possi-47 bly noisy) point cloud. This approach combines an efficient it-48 erative energy minimization scheme and a multiscale optimiza-49 tion using a hierarchical mesh representation of the source.

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The paper is organized as follows: the next section reviews 51 the state-of-the-art of non-rigid isometric registration. In Sec-52 tion 3 we introduce the concepts used in the paper. In Section 4 53 we present our registration algorithm. The method is validated 54 on our dataset in Section 5. 55

Since time efficiency is central for engineering production, 28 the registration method must be time efficient on very large 29 datasets. In the following, we will assume that the source is 30 a mesh with more than a million of vertices and the *target* is a 31 point cloud with more than 5 million samples. 32

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#### 2. Related works

Almost isometric, also known as as-rigid-as-possible (arap 2 for short), and non-rigid approaches have been studied in many fields of computer graphics: Sorkine and Alexa [3] Chao and al. [4] and Levi and al. [5] proposed surface manipulation tech-5 niques based on the arap energy, which we detail in Section 3.1. 6 Gotsmam and al. [6] introduced an arap parametrization method and more recently Smith and al. [7] proposed optimization so-8 lutions for distortion energy based problems (such as the ones 9 previously cited). Functional maps have been widely used to 10 map scalar or vector functions from a given mesh onto another 11 one while preserving some arap properties [8]. In this paper, 12 we focus on the explicit construction of an arap map between a 13 model (mesh) and a target (point cloud). 14

The NI-ICP algorithm introduced by Sacharow and al. [9] 15 is a Non-rigid Isometric variant of ICP that registers a mesh 16 onto another one. The method deforms the source into the tar-17 get using an alternate scheme that first projects the source ver-18 tices onto their nearest neighbor in the *target*, and then enforces 19 the isometry of those projections (gradient mesh editing step). 20 Given the correspondences, the algorithm fits, for each source 21 face, the best rigid transformation from the non-projected face 22 to the projected one, transforming the source mesh into a soup 23 of faces where each face is isometric to its original version. Be-24 cause this last step breaks the mesh connectivity, a global stitch-25 ing step is applied following the mesh editing proposed by Yu 26 27 and al. [10]. A final step NURBS-based technique is applied from the resulting deformation to compute a continuous map. 28

While the need for this final step is questionable in our con-29 text, the method provides an actual solution to the non-rigid 30 isometric registration for CAD models under springback de-31 formations. However, it requires a well-reconstructed surface 32 mesh from the target point cloud and is still not fast enough on 33 34 large datasets. Authors say in the paper they had to prematurely abort their experiments on full-resolution models. 35

Schweinoch and al. [11] have proposed a similar non-rigid 36 isometric registration algorithm based on a combination of ICP 37 and arap mesh deformation. However, they use a hierarchical 38 divide-and-conquer approach: the deformed source is consid-39 ered as a collection of segments initially set to a single segment. 40 The core step of the algorithm consists of subdividing each seg-41 ment according to its distance from the *target* and rigidly align 42 all those segments onto the target. Similarly to the NI-ICP ap-43 proach, the rigid alignment of all segments results in a loss of 44 connectivity at the segment endpoints. This connectivity is sub-45 sequently reestablished using the arap mesh deformation pro-46 posed by Sorkine and al. [3] that we detail in Section 3. Simi-47 larly to Sacharow and al. [9] algorithm, this method is an actual 48 solution for CAD models under springback deformation. How-49 ever it was not tested on full-resolution models. Although the 50 method uses a hierarchical approach, the whole set of source 51 vertices is used at each iteration. All vertices of each segment 52 are used to compute the rigid alignment and only the bound-53 ary vertices are used for the stitching step. So over the itera-54 tions, the total number of vertices used for alignment remains 55 the same and only the number of vertices used for the stitching 56 step increases. Therefore the method does not benefit from the 57

speed up of classical hierarchical approaches.

Klein and al. [12] propose a new method by exactly formulating the isometry registration with a proper objective function:

min 
$$\sum_{ij} |||\mathbf{p}_i - \mathbf{p}_j||^2 - ||\mathbf{z}_i - \mathbf{z}_j||^2 |$$
, (1)

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with  $\mathbf{p}_i$  and  $\mathbf{p}_i$  being vertices of the *source* surface mesh,  $\mathbf{z}_i$ 61 and  $\mathbf{z}_i$  being their associated vertices on the *target*, and *ij* a set 62 of vertex pairs not specified yet. The first step of this method 63 consists in computing a neighborhood structure of the source 64 mesh for the objective function such that close-by situated ver-65 tex pairs are conserved in the energy and distant pairs are re-66 moved. The second step is a global registration where the (non-67 convex) energy minimization is treated as a Quadratic Assign-68 ment Problem (QAP for short) from the neighborhood structure 69 and the downsampled *target*. The final step is a local reopti-70 mization where the energy is minimized by a gradient descent 71 method initialized by the QAP solutions. The method provides 72 an actual isometric registration energy and works on a point 73 cloud target. However the computation of the neighborhood 74 structure and the energy minimization being based on a QAP 75 problem (whose optimal solution is NP-hard) make the method 76 not suited for large datasets. 77

Finally, Huang and al. [13] proposed a non-rigid registration 78 method under isometric assumptions that alternates between a correspondence and a deformation optimization step. The cor-80 respondences are computed by matching source and target us-81 ing the nearest neighbor approach in a metric space defined 82 by the Euclidean distance and a feature metric defined from the principal curvatures of multi-level quadric patch fitted for 84 each vertex according to Cazals and Pouget formulation [14]. 85 Then the correspondences are pruned based on geodesic distance consistency and the remaining ones are propagated to as-87 sure uniform correspondences. The source is segmented into 88 clusters whose deformations can be described as single rigid 89 deformations and then registration of the source onto the target is computed by minimizing a combination of point-to-point 91 and point-to-plane energy according to the optimized correspondences. This method focuses on handling complex computer graphics geometries, leaving open its numerical accuracy on controlled shapes in the engineering context.

#### 3. Preliminaries

In this section, we introduce the concepts used in the following.

#### 3.1. Notations and As-Rigid-As-Possible deformation

We denote by S a source triangle mesh and by  $\mathcal{T}$  a target 100 point cloud. The mesh S has N vertices  $\{s_i\}$  in  $\mathbb{R}^3$ , its vertex 101 normals are denoted  $\{\mathbf{n}_i\}$  and its faces are denoted  $\{\mathbf{f}_k\}$ . The 102 neighboring vertices to a vertex  $\mathbf{s}_i$  is given by  $\mathcal{N}(i)$  and  $\Delta_{cotan}$  is 103 the discrete Laplace-Beltrami operator on S following the cotan 104 approach of Meyer and al. [15]. The *target* point cloud  $\mathcal{T}$  has 105 *M* vertices  $\{\mathbf{t}_i\}$  and its normal vectors are denoted  $\{\mathbf{m}_i\}$ . When 106 not specified otherwise, we will assume that N < M (the CAD 107 model has fewer vertices than the acquired point cloud). 108

We denote by  $\pi : \{0 \dots N - 1\} \rightarrow \{0 \dots M - 1\}$  an assignment map from vertices of S to points of  $\mathcal{T}$ . We further elaborate below on the definition of such assignment.

Our objective is to determine the deformation S onto T resulting in the registered surface mesh S' such that S and S' are isometric and S' and T are aligned. Because true isometric deformation cannot be achieved, the deformation between S and S' must be arap instead of isometric. This objective implies two different energies: a first one to account for the proximity between S' and T, and a second one for the arap evaluation.

The proximity energy is a simple quadratic fitting energy between S' and T from the assignment:

$$E^{prox}(\mathcal{S}',\mathcal{T}) := \sum_{i} \|\mathbf{s}'_{i} - \mathbf{t}_{\pi(i)}\|^{2}.$$
 (2)

The arap energy is a local rigidity energy between S and S':

$$E^{arap}(\mathcal{S}, \mathcal{S}', \mathcal{R}) := \sum_{i} w_i \sum_{j \in \mathcal{N}(i)} w_{ij} \|\mathbf{e}'_{ij} - R_i \mathbf{e}_{ij}\|^2, \qquad (3)$$

where  $\mathbf{e}_{ij} := (\mathbf{s}_j - \mathbf{s}_i)$ , resp.  $\mathbf{e}'_{ij} := (\mathbf{s}'_j - \mathbf{s}'_i)$ , denotes the edge vec-14 tor,  $\mathcal{R}$  denotes a set of rotation with  $R_i$  the rotation of the tangent 15 space at  $s_i$ . Weights  $w_i$  and  $w_{ij}$  are the cotangent weights of the 16  $\Delta_{cotan}$  operator attached to S. A more generic and continuous 17 elasticity model following this principle can be found in [4]. In 18 Sorkine and al. [3] arap deformation framework they define  $R_i$ 19 as the rotation between  $s_i$  and  $s'_i$  neighborhoods and derive  $R_i$ 20 from the singular value decomposition  $S_i = U_i \Sigma_i V_i^T$  of 21

$$S_i := \sum_{j \in \mathcal{N}(i)} w_{ij} \mathbf{e}_{ij} \mathbf{e}_{ij}^{\prime T}, \tag{4}$$

leading to  $R_i := V_i U_i^T$  (up to changing the last column of  $U_i$  if det $(R_i) < 0$ ).

From the arap energy, one can define several inverse problems. For instance, given a dense set of rotations  $\mathcal{R}$  and  $\mathbf{s}$ , find the  $\mathbf{s}'$  minimizing  $E^{arap}$ . The new positions of a given vertex  $\mathbf{s}'_i$ minimizing  $E^{arap}$  is given by solving the (sparse) linear system:

$$\sum_{j \in \mathcal{N}(i)} w_{ij}(\mathbf{s}'_i - \mathbf{s}'_j) = \sum_{j \in \mathcal{N}(i)} \frac{w_{ij}}{2} (R_i + R_j) \mathbf{e}_{ij} \,. \tag{5}$$

<sup>28</sup> If **s**' denotes the  $3 \times N$  matrix where each row *i* is **s**'<sub>*i*</sub>, and **b** the  $3 \times$ <sup>29</sup> *N* matrix defined from the right-hand side of (5), optimizing the <sup>30</sup> positions amounts to solving the following Poisson problem:

$$\Delta_{cotan} \, \mathbf{s}' = \mathbf{b} \,. \tag{6}$$

#### 31 3.2. Mesh simplification

Our approach relies on a hierarchical structure of the source 32 mesh S. Although many solutions exist in geometry processing 33 and computer graphics, we rely on the simple, yet very effi-34 cient, edge collapsing simplification proposed by Garland and 35 al. [16]. This method combines edge collapsing and quadric 36 error metric that estimates the squared distance to the mesh. 37 Initially an error quadric matrix  $Q_i$  is assigned to each  $s_i$  that 38 encodes the total squared distance a point **p** to  $s_i$ 's neighboring 39 faces which is given by computing  $\mathbf{p}^T Q_i \mathbf{p}$ . The cost of each 40

edge (i, j) is computed by minimizing the energy  $\mathbf{p}^T (Q_i + Q_j) \mathbf{p}$ where  $\mathbf{p}$  is the point that minimizes the cost (see Figure 1).

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Fig. 1. Edge collapsing: the edge formed by the two green vertices in (a) is collapsed and the green vertices are merged into the red vertex in (b).

The edge with the minimal cost is collapsed into a new vertex 43 whose position is the point **p** that minimizes the edge cost and 44 tis quadric error matrix is the sum of the quadrics of the edge extremities. 46

From this elementary edge collapse step, the overall hierarchical representation  $\mathcal{H} : S \to S^{h-1} \to S^{h-2} \to \ldots \to S^0$  of Swith *h* layers can be given by considering disjoint sequences of minimal edge collapses per layer (see 4.2 for details).

#### 4. Fast and accurate hierarchical As-Rigid-As-Possible registration

In this section we present our solution to compute an arap registration of S onto T and then its hierarchical optimization. <sup>53</sup>

#### 4.1. Single level registration

Let us first detail the arap registration for a single layer of 56 the hierarchical reconstruction. Algorithm 1 presents the over-57 all algorithm to construct S' with the same topology as S, such 58 that S' is close to  $\mathcal{T}$  and arap with respect to S. This algorithm 59 alternates between an assignment and an arap optimization step 60 close to the one proposed by Sacharow and al. [9]. We extend 61 this preminilary formulation with two main contributions: First, we propose a hierarchical optimization detailed in Section 4.2. 63 Second, Sacharow and al. method requires that the *target* is a 64 triangulated surface to define per face rotations needed in their 65 formulation, leading to a mandatory preliminary surface reconstruction step from the point cloud. Our formulation has per 67 vertex rotation information allowing us to directly work on the input point cloud stored in an efficient associated data structure 69 (see below). 70

Assignment step (line 4 to 6). This step consists in assigning 71 each vertex  $\mathbf{s}'_i$  to a point  $\mathbf{t}_{\pi(i)}$  of  $\mathcal{T}$  and therefore results in a set 72 of pairs  $(i, \pi(i))$ . Although the nearest neighbor operator from 73  $\mathbf{s}'$  to  $\mathbf{t}$  is not injective and may produce inconsistent pairs, our 74 experiments and the results of Sacharow and al. [9] and Klein 75 and al. [12] has shown that its use under small or simple de-76 formations is a good trade-off between assignment quality and 77 computation efficiency. We use a kdTree structure to perform 78 the nearest neighbor requests. Since T is static, the kdTree 79 is computed only once before the optimization loop. During 80 the minimization process, geometrical information from  $\mathcal{T}$  will 81 only be given through nearest-neighbor queries on the kdTree. 82



Fig. 2. Overall workflow: Starting from a *source* mesh S and a *target* point cloud T(a), we first start with assignation  $\pi : S \to T(b)$ . From rotations  $\mathcal{R}$  computed from the assignment, we solve a first arap matching (c). In (d) we illustrate the final surface after 3 iterations.

We provide more technical details in Section 5.1, efficient NN
 queries allow us to consider very large point clouds as targets.

Algorithm 1: arap registration algorithm							
$\epsilon$ : a stopping criterion							
<b>Output:</b> $S' = ({s'_i}, {f_k})$ : as-rigid-as possible deformed							

<sup>3</sup> *Optimization step (line 7 to 13).* We could use the  $\{\mathbf{t}_{\pi(i)}\}$  as our <sup>4</sup> mapping solution, but since they may not respect the arap con-<sup>5</sup> straint, we need to refine those pairs to ensure that  $\mathbf{s}'$  minimizes <sup>6</sup>  $E^{arap}$ . The solution of the arap energy minimization (5) requires <sup>7</sup> the original *source* vertices  $\mathbf{s}$ , it's cotangent weights expressed <sup>8</sup> in the Laplacian operator  $\Delta_{cotan}$  and rotation matrices  $\mathcal{R}$ . From <sup>9</sup> the assignment  $\pi$ , rotation matrices  $R_i$  are given by:

$$R_i := I + [\mathbf{n}_i \times \mathbf{m}_{\pi(i)}]_{\times} + \frac{1}{1 + \mathbf{n}_i \cdot \mathbf{m}_{\pi(i)}} [\mathbf{n}_i \times \mathbf{m}_{\pi(i)}]_{\times}^2.$$
(7)

with *I* is the identity  $3 \times 3$  matrix and  $[\mathbf{v}]_{\times}$  is the map from a 3D vector  $\mathbf{v}$  to a  $3 \times 3$  skew-symmetric matrix such that  $[\mathbf{v}]_{\times}\mathbf{q} = \mathbf{v} \times \mathbf{q}$  for any  $\mathbf{q} \in \mathbb{R}^3$ . Because we assume the deformation of S onto  $\mathcal{T}$  is isometric,  $R_i$  is the rotation occurring between  $\mathbf{s}_i$  and  $\mathbf{s}'_i$  13 local plane, thus  $R_i$  can be computed from  $\mathbf{s}_i$  and  $\mathbf{s}'_i$  normals according to (7). With a rotation for each  $\mathbf{s}_i$  (5) can be solved to get new vertices  $\mathbf{s}'$  that respect the arap constraint. Solving (5) handles the rotational part of the mapping, but because it also writes off its translation part, we move the barycenter of  $\mathbf{s}'$  to  $\mathcal{T}$ barycenter.

Solver (line 12). The  $\Delta_{cotan}$  being sparse, (5) can be efficiently solved either using factorized solvers (*e.g.* LU decomposition) or iterative schemes (*e.g.* conjugate gradient). The first one would lead to an expensive preprocessing overhead while the second approach also allows us to reuse results from previous iterations in Algorithm 1, *i.e.* vertices s' (see line 11), as warm start. This warm start could be problematic for the very first iteration as S' may be far from the target, but it becomes more and more efficient for the remaining iterations.

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#### 4.2. Hierarchical registration

When the source mesh S is large, the main bottleneck of Algorithm 1 is the linear system solve on line 12 (see Section 5). To improve the overall performances we use a geometrical multigrid approach which constructs a hierarchical representation of S, and a coarse-to-fine minimization of the arap energy.

As first sketched in Section 3.2, H denotes a hierarchical 35 view of S in which  $S^h$  denotes the *h*-th layer with vertices  $\mathbf{s}_i^h$ 36 (resp. normal vectors  $\mathbf{n}_{i}^{h}$ , faces  $\mathbf{f}_{k}^{h}$  and Laplace-Beltrami oper-37 ator  $\Delta_{cotan}^{h}$ ). Layers are sorted in ascending order (with respect 38 to the number of vertices), the last layer being S. Connections 39 between two consecutive layers  $S^h$  and  $S^{h+1}$  (see Figure 4) are given by linking each vertex  $\mathbf{s}_i^{h+1}$  to a face  $\mathbf{f}_k^h$  and  $\mathbf{s}_i^{h+1}$  local po-sition within  $\mathbf{f}_k^h$  is represented by the barycentric coordinates 40 41 42  $(\alpha_i^{h+1}, \beta_i^{h+1}, \gamma_i^{h+1})$  of its projection onto  $\mathbf{f}_k^h$  along the face normal and the (possibly negative) height  $d_i^{h+1}$  from its projection such 43 44 that: 45

$$d_i^{h+1} = (\mathbf{s}_i^{h+1} - \mathbf{p}_0) \cdot \mathbf{q}$$
(8)

$$\mathbf{s}_{i}^{h+1} - d_{i}^{h+1}\mathbf{q} = \alpha_{i}^{h+1}\mathbf{p}_{0} + \beta_{i}^{h+1}\mathbf{p}_{1} + \gamma_{i}^{h+1}\mathbf{p}_{2}, \qquad (9)$$

where **p** is one of the three vertices of  $\mathbf{f}^h$ , and **q** the face normal vector. With  $\mathbf{s}_i^{h+1}$  local coordinates to  $\mathbf{f}_k^h$  and given  $\mathbf{f}_k^h$  updated vertices **p**' and normal **q**',  $\mathbf{s}_i^{h+1}$  new position is simply:

$$\mathbf{s}_{i}^{h+1} = \alpha_{i}^{h+1}\mathbf{p}_{0}' + \beta_{i}^{h+1}\mathbf{p}_{1}' + \gamma_{i}^{h+1}\mathbf{p}_{2}' + d_{i}^{h+1}\mathbf{q}' \,. \tag{10}$$



Fig. 3. Overall hierarchical workflow: Starting from a coarse mesh  $S^0(a)$ , we solve a first arap matching leading to the mesh  $S'^0(b)$ . We then consider the mesh  $S^1$  with updated positions from  $S'^0(c)$ , and repeat the minimization+optimization process to obtain the final surface S' after 3 iterations (d).

Since there is no inclusion between  $\{\mathbf{s}^{h}\}$  (resp.  $\{\mathbf{f}^{h}\}$ ) and  $\{\mathbf{s}^{h+1}\}$ (resp.  $\{\mathbf{f}^{h+1}\}$ ), the sparse operator  $\Delta_{cotan}^{h+1}$  is simply reconstructed from  $S^{h+1}$  (linear time algorithm in the number of edges/vertices).



Fig. 4. Exploded view of the hierarchical structure: Vertex  $s_i^{h+1}$  of upper layer  $S^{h+1}$  is linked to face  $f_k^h$  of lower layer  $S^h$ . Vertex  $s_i^{h+1}$  position within  $f_k^h$  is encoded the barycentric coordinates  $(\alpha_i^{h+1}, \beta_i^{h+1}, \gamma_i^{h+1})$  of its projection and its signed distance  $d_i^{h+1}$  from  $f_k^h$ .

Algorithm 2: Hierarchical arap registration	
<b>Input:</b> $\mathcal{H} = \{S^0 \dots S^h\}$ : hierarchical <i>source</i> surface	
mesh	
$\mathcal{T} = (\{\mathbf{t}_i\}, \{\mathbf{m}_i\}): target \text{ point cloud}$	
$\epsilon$ : a stopping criterion	
1 foreach $S^h \in \mathcal{H}$ with h ascending do	
/* Register $\mathcal{S}^h$ to $\mathcal{T}$ */	
2 $\mathcal{S}^{h\prime} \leftarrow \text{Arap}_{\text{Registration}}(\mathcal{S}^{h}, \mathcal{T}, \epsilon)$	
3 <b>if</b> $S^h$ is last layer <b>then</b>	
4 return $S^{h'}$	
5 else	
/* Propagate $\mathcal{S}^{h\prime}$ to $\mathcal{S}^{h+1}$ */	
6 foreach $\mathbf{s}_i^{h+1} \in \mathcal{S}^{h+1}$ do	
7 compute $\mathbf{s}_i^{h+1}$ according to (10)	
8 end	
9 end	
10 end	

Given a set of desired vertices sizes, Garland and Heckbert's edge collapsing method [16] is applied to the original *source* S until all desired sizes are reached and results in  $\mathcal{H} = \{S^0 \dots S^h\}$ without connection between layers. Because this method minimizes the distance between the vertices from collapsed edges and their neighboring faces, each vertex  $s^{h+1}$  of  $S^{h+1}$  is linked to S closest face  $\mathbf{f}_{h}^{h}$ .

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Our hierarchical optimization of the arap registration of Sonto T is presented in Algorithm 2. Given a hierarchy H of the *source* surface mesh S and the *target* point cloud T, each layer  $S^h$  from the coarsest to the finest is registered to T (line 2) and the current registration result  $S'^h$  is propagated to the next layer  $S^{h+1}$  (line 7). As  $S^h$  has fewer vertices than  $S^k$  (for k > h), fewer nearest neighbor queries are required and the size of the  $\Delta^h_{cotan}$  is smaller. The last iteration of Algorithm 2 is performed on the full resolution mesh S. Note thanks to the hierarchical construction, very few optimization steps of Algorithm 1 are necessary to obtain a correct registration of this last step.

In the next section, the hierarchical optimization is experimented on data to evaluate the quality and the efficiency of the method.

#### 5. Experiments

#### 5.1. Dataset

The dataset (see Figure 5) is composed of three analytical 28 models that mimic real-application deformations (springback, 29 torsion and complex deformation), a car hood model (exhibiting 30 high curvature features), and a CAD model corresponding to a 31 real test case. To illustrate the *springback*, a hat shape whose parametrization is detailed in Appendix A is registered onto 33 another hat. The hat shape is controlled by a bending factor and given its parametrization, two hats with different bending 35 factor are isometric. For the torsion, a rectangle is mapped onto 36 a helicoid twisted 90° whose parametrization is also detailed 37 in Appendix A. The sheet case corresponds to the registration 38 of a square to a deformed one generated by applying several 39 rotation fields according to (5) and (6). The same method is 40 used to generate the deformed car hood. Because the rotation 41 fields used are manually chosen, there is no reason for them to be actual solutions of  $E^{arap}$  minimization. 43

Input scans for our analytical models and the hood have been generated by uniformly sampling each object and by adding Gaussian noise on the samples positions and normals. Given a triangular mesh, a face is randomly selected with an importance function proportional to the face area and the point lying on the



Fig. 5. Dataset: Yellow meshes are *source* models, blue point clouds correspond the *target* scans. Finally, red meshes are outputs we obtain after registration (see Table 1 for the settings). From left to right columns: two hats (a, f) under *springback* deformation, a mapping between a rectangle (b) and a twisted helicoid (g) (90° to mimic a torsion), the sheet (c, h) for a complex deformation and a car hood model (d) with simulated scan (i). The last column (e, j, o) corresponds to a real test case of a CAD model and a scan of the real metal sheet. Meshes and point clouds are downsampled for visualization purposes.

face is selected by uniformly sampling its barycentric coordinates. The normal vector of the point is set to the normal vector of the face it belongs to, and this process is repeated until the desired number of points *M* has been sampled. The positions are perturbed by adding to their coordinates a noise sampled from a Gaussian distribution ( $\mu = 0, \sigma_{coord}$ ) and the normals are perturbed by a rotation ( $\theta, \phi$ )<sup>1</sup> with  $\theta$  sampled from an uniform distribution [0,  $\pi$ [ and  $\phi$  sampled from a Gaussian distribution ( $\mu = 0, \sigma_{angle}$ ).

Finally, the scan of the CAD model is a measure of the real piece which undergoes a *springback* deformation. Because of the machining context, the actual *target* surface (which corresponds to the *source* after deformation) is yet unknown. To ensure that no part is missing, the scanned surface is bigger and so the arap registration is performed with a *source* being a sub-part of the *target* point cloud.

For all experiments, the dataset has been normalized such that the diagonal of *source* bounding box measures 1. Our algorithm has been implemented in C++ with the libraries eigen [17] and libigl [18] for the data structures and geometry processing, nanoflann [19] for the kdTree and polyscope [20] for the visualization. All tests were run on a 4.0 GHz PC with 16 GB RAM. For each experiment (*e.g.* Table 1) we measure the total num-

<sup>24</sup> ber of iterations  $n_{iter}$  which is the sum of Algorithm 1 itera-<sup>25</sup> tions for all layers, the final  $E^{prox}$  and  $E^{arap}$ , the initialization step duration  $T_{init}$  (construction of  $\mathcal{H}$  and the Laplace-Beltrami 26 operator  $\Delta_{cotan}$  computations), the iterative step duration  $T_{core}$ 27 which is the execution time of Algorithm 2. The assignment 28 step duration  $T_{NN}$  is the timing of all nearest-neighbor queries 29 on  $\mathcal{T}$  (accumulated for all layers),  $T_{opt}$  denotes the remain-30 ing optimization timing of S (also accumulated for all layers). 31 Hence,  $T_{core} = T_{NN} + T_{opt}$  and the final timing is given by 32  $T_{total} = T_{init} + T_{core}$ . We denote  $a_{total}$  (resp.  $a_{core}$ ), the speedup 33 between the hierarchical and the nonhierarchical variants of the 34 algorithms (in total and for the core part only). 35

Hat and helicoid sources have both  $N = 10^6$  vertices and their targets have  $M = 5 \times 10^6$  points. The sheet source has  $N = 4 \times 10^5$  vertices and its target has  $M = 2.5 \times 10^6$  points. The CAD model has  $N = 8 \times 10^5$  vertices and its target has  $M = 1.4 \times 10^6$  points. The number of layers for the hierarchical registration is set to 3 with the first, second and last layer having respectively N/100, N/10 and N vertices. The stopping criterion d and the maximum number of iterations of Algorithm 1 per layer are set to  $10^{-6}$  and 100.

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#### 5.2. Accuracy and Efficiency

In this experiment the hierarchical optimization is evaluated by comparing for each dataset case, the registration results with and without the hierarchy. The results for each case are displayed in Table 1.

For the hat object, both global and hierarchical approaches produce similar results in terms of energy. However, the hierarchical registration is 43 times faster in total and 115 times 52

 $<sup>^{1}\</sup>theta$  being the azimuthal angle and  $\phi$  the polar angle.

Case	sizes (N, M)	<i>n<sub>iter</sub></i>	Eprox	Earap	T <sub>total</sub>	T <sub>init</sub>	$\begin{array}{c c} T \\ T_{NN} \end{array}$	core T <sub>opt</sub>	a <sub>total</sub>	<i>a<sub>core</sub></i>
hat	$N = 10^{6}$	17	$1.3  imes 10^{-2}$	$5.1 \times 10^{-6}$	1734	2 (0.1%)	1732 ( 1373	(99.9%) 361	42	115
hat hierarchical	$M = 5 \times 10^6$	12	$1.9 \times 10^{-2}$	$4.6  imes 10^{-6}$	40	25 (62.5%)	<b>15 (3</b> 12	3 <b>7.5%</b> )	43	115
helicoid	$N = 10^{6}$	100	$2.5 \times 10^{-1}$	$5.0 \times 10^{-3}$	514	2 (0.4%)	512 ( 288	99.6%) 224	7	10
helicoid hierarchical	$M = 5 \times 10^6$	136	$1.0  imes 10^{-1}$	$2.0 imes10^{-6}$	72	25 (34.7%)	<b>47 (6</b> 33	5.3%) 14		10
sheet	$N = 4 \times 10^5$	7	1.3	$1.8 \times 10^{-2}$	107	1 (0.9%)	106 ( 51	99.1%) 55		0
sheet hierarchical	$M = 2.5 \times 10^6$	40	$1.3  imes 10^{-1}$	$2.7  imes 10^{-3}$	24	11 (45.8%)	<b>13 (5</b> 3	<b>4.2%</b> ) 10	4	0
hood	$N = 8 \times 10^5$	8	$3.7 \times 10^{-1}$	$2.5 \times 10^{-3}$	311	1 (0.3%)	310 ( <sup>1</sup> 55	99.7%) 255	15	17
hood hierarchical	$M = 5 \times 10^6$	18	$3.7 \times 10^{-2}$	$1.0  imes 10^{-3}$	201	21 (10.4%)	<b>180 (8</b>	<b>89.6%</b> ) 173	1.3	1.7
CAD	$N = 8 \times 10^5$	100	$6.7 \times 10^{-1}$	$4.5 \times 10^{-4}$	2065	1 (0.04%)	2065 (	99.96%)		
						(	38	2027	21	28
CAD hierarchical	$M = 1.4 \times 10^6$	150	$6.3 \times 10^{-1}$	$3.8\times10^{-6}$	94	21 (22.3%)	3	70		

Table 1. Metrics of nonhierarchical and hierarchical registration on the dataset after convergence (with stopping criterion  $d = 10^{-6}$  – see Algorithm 1 – and number of iterations per layer limited to 100). Metrics are n<sub>iter</sub> the sum of Algorithm 1 iterations for all layers, the final E<sup>prox</sup> and E<sup>arap</sup>, the total duration  $T_{total}$ , the initialization step and iterative step duration  $T_{init}$  and  $T_{core}$  ( $T_{total} = T_{init} + T_{core}$ ), the assignment steps and optimization step  $T_{NN}$  and  $T_{opt}$  ( $T_{core} = T_{NN} + T_{opt}$ ) and the total and iteration step speedups  $a_{total}$  and  $a_{core}$ . Times are expressed in seconds.

faster if we only consider the iterative part. More detailed measures show that iterations with a bad alignment onto  $\mathcal{T}$  (*i.e.* with large  $E^{prox}$ ) were performed on the smaller layers. Even though the hierarchical registrations of the helicoid, the sheet and the CAD models have smaller speed ups and required more iter-5 ations than their nonhierarchical versions, they both produce mappings with lower arap energies. This can be due to the fact that optimizing the energies in a coarse-to-fine scheme make the process unlikely to get stuck on local minima.

For the sheet and hood cases for which the deformations are 10 not fully arap, the hierarchical method produces better  $E^{prox}$ 11 than the nonhierarchical one. The speedup for the hierarchical 12 car hood (1.5) is less favorable than for the other shapes as sev-13 eral high resolution iterations are required to cope with the high 14 curvature features. Figure 6 shows  $E^{prox}$  and  $E^{arap}$  distributions 15 on the sheet case. 16

Results in Table 2 show the impact on the *target* point cloud 17 size to the arap registration computation time. For all analytical 18 cases, we thus increase the sample count of the target. The 19 assignment time  $T_{NN}$  (which is a function of N and M) scales up 20 with M thanks to the kdTree structure used for nearest neighbor 21 queries, while the optimization time  $T_{opt}$  which is a function of 22 N is constant as expected. 23

#### 5.3. Robustness to noise 24

In this experiment, we evaluate the robustness to perturba-25 tions on the positions and normal vectors of  $\mathcal{T}$  on the hat case 26 (hierarchical approach). To evaluate the robustness with respect 27

to the perturbations, we consider the  $l_2$ -norm between the de-28 formed surface  $S^*$  vertices for the noise-free target T (the 29 groundtruth for this test), and the surface obtained on the per-30 turbed  $\mathcal{T}$ : 31

$$E^{*}(\mathcal{S}, \mathcal{S}^{*}) := \sum_{i} \|\mathbf{s}_{i} - \mathbf{s}_{i}^{*}\|^{2}.$$
(11)

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When perturbing the normal vectors, experiments (Figure 7) 32 show that Algorithm 1 is stable for small perturbations (closer 33 to our use-case). For extreme cases with strong perturbations  $(30^{\circ} (c) \text{ and } 60^{\circ})$ , the output surface becomes highly impacted. 35 When perturbing the sample positions, (Table 3 and Figure 8), 36 the algorithm produces very stable solutions.

Finally, in Figure 9, we provide additional experiments when 38 we consider non-uniform point distributions for the target point 30 cloud  $\mathcal{T}$ . Note that the last case (c - d) exactly corresponds to 40 the kind of point clouds we are facing in inverse engineering 41 when  $\mathcal{T}$  is given by a laser range or a probing device. Our 42 method can efficiently handles these cases. 43

#### 5.4. Failure example

In this experiment we apply our registration method on data 45 that are slightly out-of-scope of the reverse engineering con-46 text. As illustrated in Figure 10, our method does not produce 47 satisfactory results on complex geometries, due to the assign-48 ment step. In the bunny case (see Figure 10(c)) the two sides of 49 the source left ear are assigned on the opposite sides of the tar-50 get two ears. With this assignment the two sides of the source 51 ear should be flipped. The two sides of the source right ear 52



Fig. 6. Sheet registration results: Top row displays nonhierarchical results, bottom row the hierarchical results. From left to right the results are registration output (with T in blue and S' in red), the  $E^{prox}$  per vertex and the  $E^{arap}$  per vertex.  $E^{prox}$  (resp.  $E^{arap}$ ) share the same colormap for both versions with purple (resp. yellow) as lower (resp. upper) bound.

Case	N	М	$1 \times 10^{6}$	$5 \times 10^{6}$	$10 \times 10^{6}$	$15 \times 10^{6}$	$20 \times 10^{6}$	$25 \times 10^{6}$
hat	$N = 10^{6}$	$T_{NN}$	3	13	20	28	34	40
		T <sub>opt</sub>	4	10	11	12	11	11
helicoid	$N = 10^{6}$	$T_{NN}$	23	38	56	70	81	96
		T <sub>opt</sub>	17	16	15	17	17	18
sheet	$N = 4 \times 10^5$	$T_{NN}$	2	6	8	11	11	12
		T <sub>opt</sub>	16	17	17	16	18	17

Table 2. Scan size impact on hierarchical registration duration: For all cases the *source* vertex number N is fixed, M is the number of points in the *target* scan and  $T_{NN}$  and  $T_{opti}$  the total amounts of time spent on assignment and optimization steps (see Section 4.1 for step details).



Fig. 7. Registration results with perfect position and noisy normals: The normals are perturbed using a Gaussian noise as specified in Section 5.1 with  $\sigma_{angle}$  equals from left to right 3° (*a*), 6° (*b*), 30° (*c*) and 60° (*d*). Top row displays the normal distribution and bottom row the registration output S'. With small enough noises (*e.g.* 3° (*a*) and 6° (*b*)) the mapping is not smooth but consistent while with too big noises (*e.g.* 30° (*c*) and 60° (*d*)) the algorithm produces degenerated results.



Fig. 8. Registration results with noisy positions and perfect normals: The positions are perturbed using a Gaussian noise as specified in Section 5.1 with  $\sigma_{coord}$  equals from left to right 0.1% (*a*), 0.4% (*b*), 0.7% (*c*) and 1% (*d*) of the diagonal length of the point cloud bounding box. Top row displays the position distribution and bottom row the registration output S' colored by  $E^*$  per vertex,  $E^*$  share the same colormap for all  $\sigma_{coord}$  with purple (resp. yellow) as lower (resp. upper) bound. See Table 1 for detailed metrics.

$\sigma_{coord}$	<i>n<sub>iter</sub></i>	Eprox	Earap	$E^*$	$T_{total}$ (sec)
0%	16	$1.3 \times 10^{-2}$	$3.6 \times 10^{-6}$	0	48
0.1%	23	$1.5 \times 10^{-2}$	$7.0 imes10^{-6}$	$4.5 \times 10^{-3}$	43
0.4%	81	$3.2 \times 10^{-2}$	$4.7  imes 10^{-6}$	$3.0 \times 10^{-2}$	100
0.7%	174	$4.7 \times 10^{-2}$	$1.2 \times 10^{-4}$	$2.3 \times 10^{-2}$	179
1%	290	$5.9 \times 10^{-2}$	$2.5 \times 10^{-4}$	$8.5 \times 10^{-2}$	297

Table 3. Registration results with noisy positions and perfect normals after convergence:  $\sigma_{coord}$  is expressed as percent of the diagonal length of the point cloud bounding box (see Table 1 for column details). See Figure 8 for visualizations of  $E^*$ . Under  $\sigma_{coord} = 0.4\%$  the algorithm produces solutions with similar  $E^{arap}$  and  $E^{prox}$  (which is affected by  $\sigma_{coord}$ ). Under  $\sigma_{coord} = 0.1\%$  the algorithm produces a solution quasi-equal to  $S^*$  in the same amount of time and iterations.



Fig. 9. Robustness to different point distributions: When the sampling density is not uniform (*a*) or when its has high aliasing structure when simulating probing/laser based acquisition devices (*c*), our approach still provides stable and accurate outputs (*b*) and (*c*).

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Fig. 10. Failed bunny registration: Registration between two Stanford bunnies (a) with one bunny having bent ears. The algorithm returns a bunny with degenerated ears (b). Bad assignments (c) produce inconsistent rotations leading to an inconsistent solution (d).

are assigned on the same side of one *target* ear, so one side of the source ear should also be flipped. Theses assignments pro-2

duce areas where two neighboring vertices could have opposite

3 normals leading to a inconsistent rotation field  $\mathcal{R}$ , and incon-

sistent **b** and a degenerated (6) solution (e.g. bunny result and 5

Figure 7(h)). Note that the overall algorithm (assignment/arap alternate steps, hierarchical construction...) perfectly handles this case but providing a better assignment for these complex 8 shapes (mimicking the one in [13] without scarfing the speed of q

our approach) is a challenging future work. 10

#### 6. Conclusion 11

In this paper, we have presented a new method of non-rigid 12 isometric registration of a source triangle mesh onto a target 13 point cloud that is fast, accurate and can handle large geome-14 tries. Our solution is based on a hierarchy of meshes built from 15 the source triangle mesh. Our algorithm alternates between 16 an assignment and an arap optimization step, starting from the 17 coarsest mesh of the hierarchy, and going up progressively each 18 time the solution with the current level has converged. Our ex-19 periments show a very highest speed up compared to the non-20 hierarchical approach, while ensuring the same accuracy. 21

Several future works exist. As mentioned in Section 5.4, the 22 assignment step could be improved to be able to handle more 23 complex shapes in terms of geometry or topology. We would 24 also consider alternative schemes for the hierarchical construc-25 tion (for instance to keep some user-specified features, or to 26 preserve some spectral properties similarly to [21]). The chal-27 lenge in these future works would be to keep a solution that 28

- scales up with respect to the size of the source and the target, 29
- while keeping the accuracy of the reconstruction as required in 30 our inverse engineering context.

#### Appendix A. Hat and helicoid parametrization

By definition a ruled surface can be described by a 33 parametrization p(u, v) = c(u) + v \* r(u) with p the surface ver-34 tices, c its directrix curve and r its generator. Hat directrix  $c_{hat}$ is controlled by a bending factor b and its piecewise definition 36 is given on each piece *i* by:

$$c_{hat}(u,b) := \begin{bmatrix} L_i(b) \cos(\omega_i(b) \, u + \phi_i(b)) \\ L_i(b) \sin(\omega_i(b) \, u + \phi_i(b)) \\ 0 \end{bmatrix} + P_i(b)$$
(A.1)

with  $L_i$ ,  $\omega_i$ ,  $\phi_i$  and  $P_i$  chosen such that  $c_{hat}$  is at least  $C^1$  and isometric for any *b*. Hat generator  $r_{hat}$  is a simple constant [0, 0, W].

The helicoid is a well known ruled surface, like with the hat we introduce a bending factor b to control its deformation such that:

$$c_{helicoid}(u) := \begin{bmatrix} 0\\0\\W u \end{bmatrix} \text{ and } r_{helicoid}(u) := \begin{bmatrix} L\cos(b u)\\L\sin(b u)\\0 \end{bmatrix}. \quad (A.2)$$

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