

# Hands-on AI based 3D Vision Summer Semester 25

Lecture 5\_0 – Surface Reconstruction

Prof. Dr.-Ing Gerard Pons-Moll

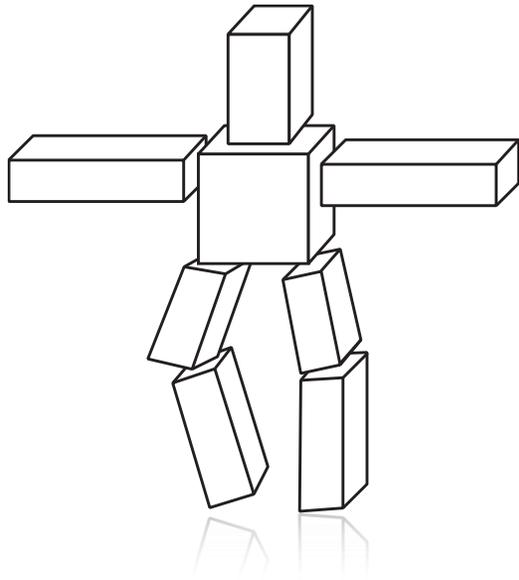
University of Tübingen / MPI-Informatics

EBERHARD KARLS  
UNIVERSITÄT  
TÜBINGEN



# Increasing complexity of our models

**Transformations**



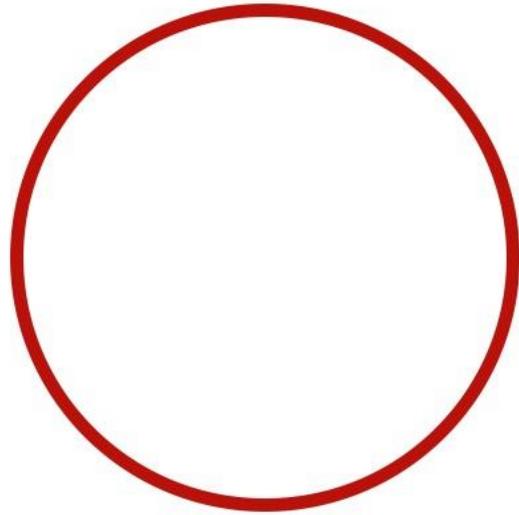
**Geometry**



**Materials, lighting, ...**



# How can we describe geometry?



# How can we describe geometry?

**IMPLICIT**

$$x^2 + y^2 = 1$$

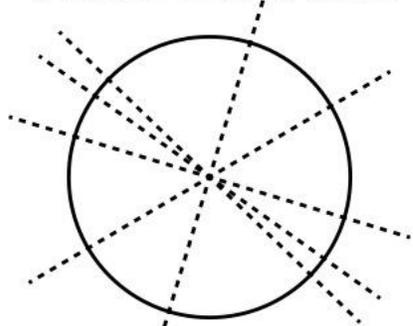
**LINGUISTIC**

“unit circle”

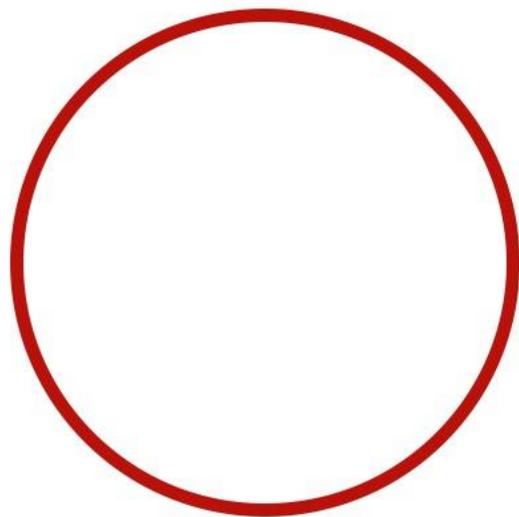
**EXPLICIT**

$$\left( \underbrace{\cos \theta}_x, \underbrace{\sin \theta}_y \right)$$

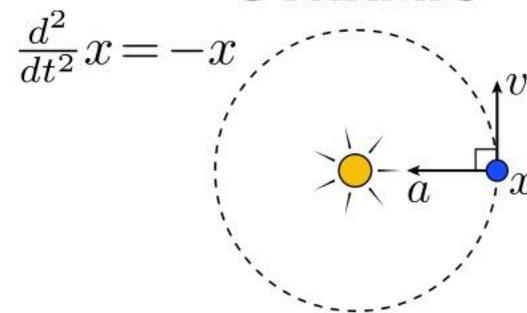
**TOMOGRAPHIC**



(constant density)



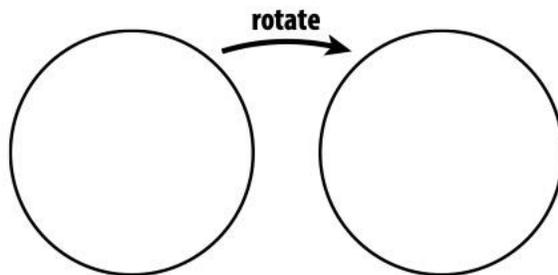
**DYNAMIC**



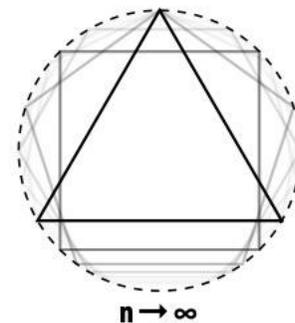
**CURVATURE**

$$\kappa = 1$$

**SYMMETRIC**



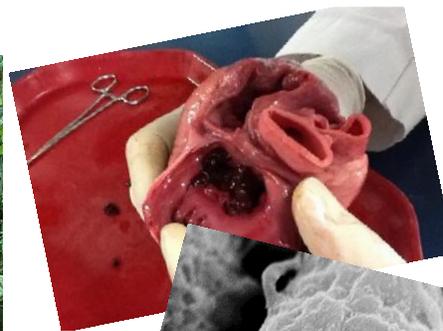
**DISCRETE**



$n \rightarrow \infty$

Given all these options, what is the best way to encode geometry on a computer?

# It's a jungle out there!



# No one “best” choice—geometry is hard!

*“I hate meshes.  
I cannot believe how hard this is.  
Geometry is hard.”*

**—David Baraff**

**Senior Research Scientist  
Pixar Animation Studios**

# Many ways to digitally encode geometry

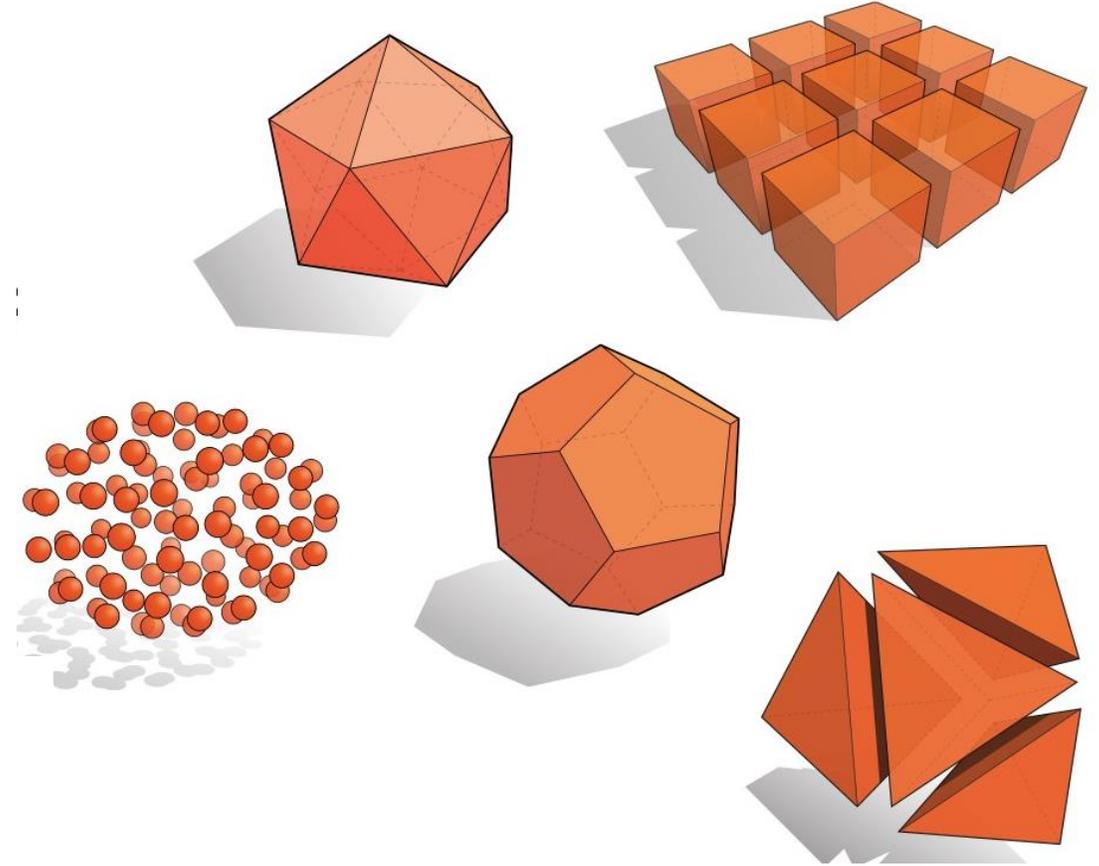
- EXPLICIT

- point cloud
- polygon mesh
- subdivision, NURBS
- ...

- IMPLICIT

- level set
- algebraic surface
- L-systems
- ...

- Each choice best suited to a different task/type of geometry



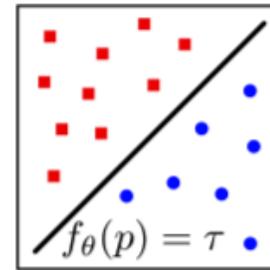
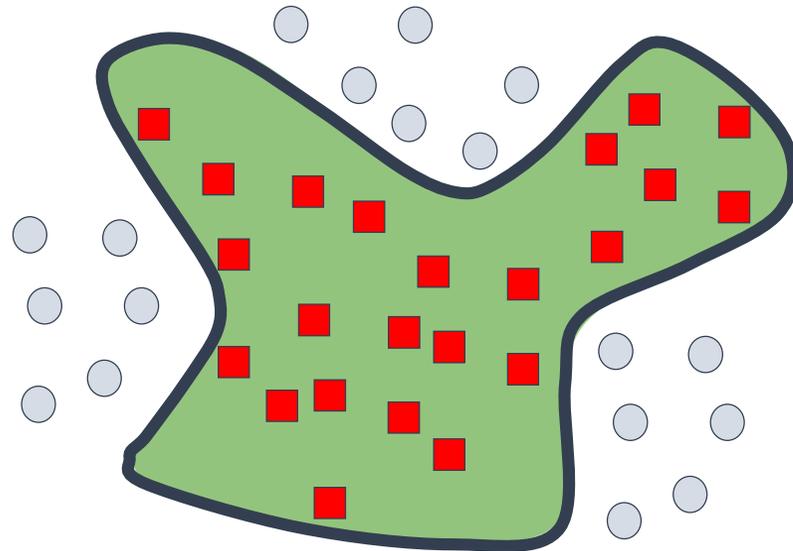
# "Implicit" Representations of Geometry

- Points aren't known directly, but satisfy some relationship
- E.g., unit sphere is all points such that  $x^2+y^2+z^2=1$
- More generally,  $f(x,y,z) = 0$

# Surfaces as an Implicit Function

$$\mathbf{p} = (x, y, z) \in \mathbb{R}^3$$

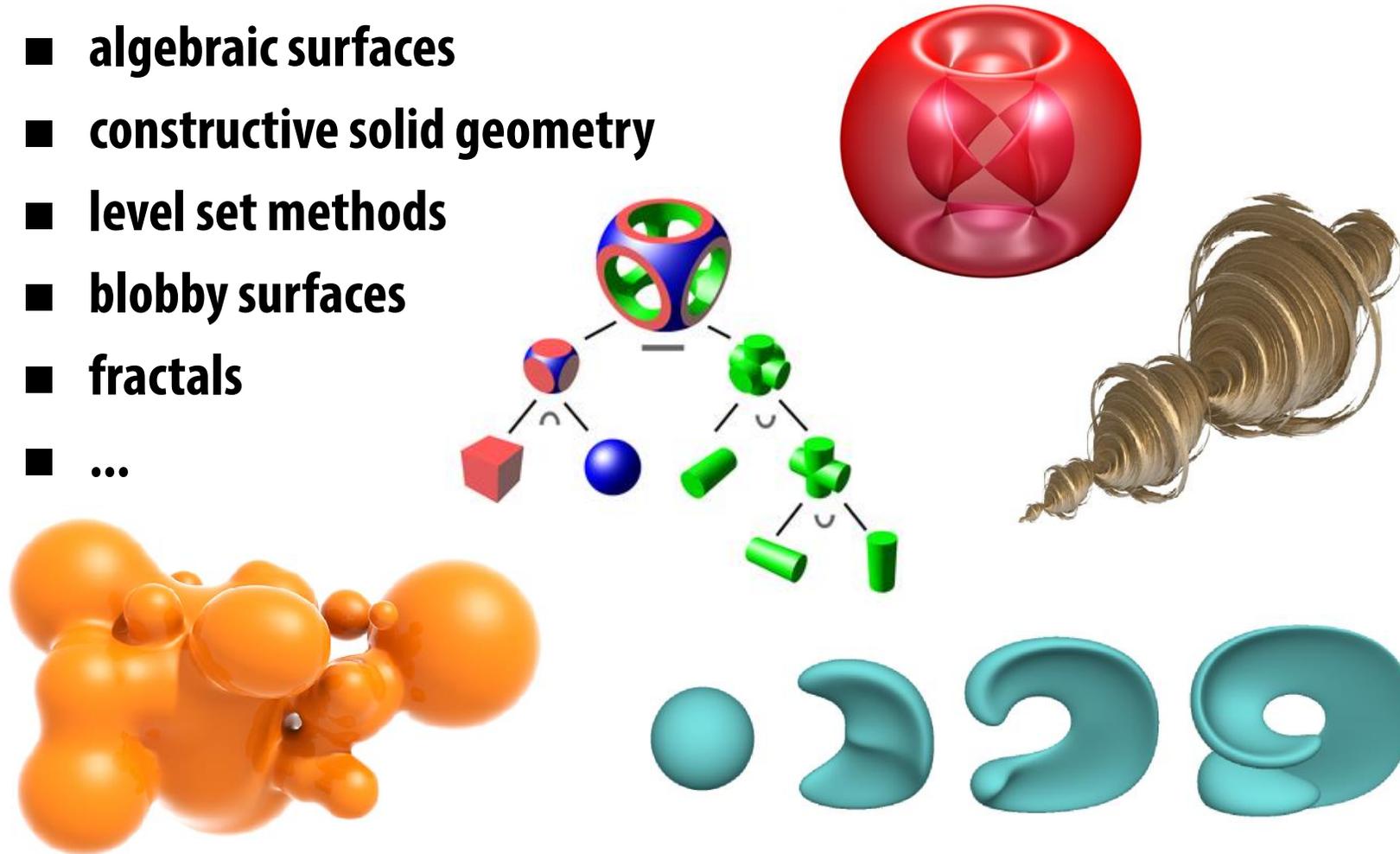
$$f(\mathbf{p}) = \begin{cases} 0, & \text{if } \mathbf{p} \in \text{outside } \circ \\ 1, & \text{if } \mathbf{p} \in \text{inside } \blacksquare \end{cases}$$



$$\mathcal{S} = \{\mathbf{p}, f(\mathbf{p}) = \tau\}$$

# Many implicit representations in graphics

- algebraic surfaces
- constructive solid geometry
- level set methods
- blobby surfaces
- fractals
- ...



(Will see some of these a bit later.)

**But first, let's play a game:**

**I'm thinking of an implicit surface  $f(x,y,z)=0$ .**

**Find *any* point on it.**

**Let's play another game.**

**I have a new surface  $f(x,y,z) = x^2 + y^2 + z^2 - 1$ .**

**I want to see if a point is *inside* it.**

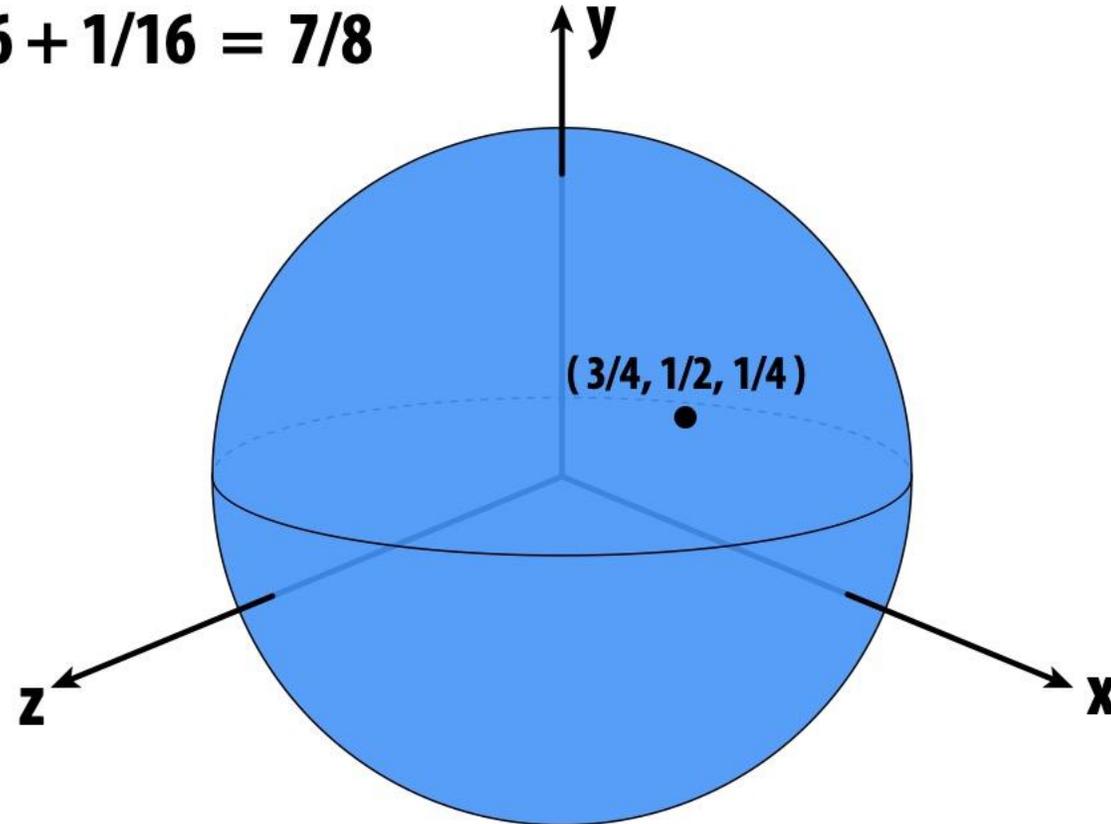
# Check if this point is inside the unit sphere

How about the point  $(3/4, 1/2, 1/4)$ ?

$$9/16 + 4/16 + 1/16 = 7/8$$

$$7/8 < 1$$

**YES.**



**Implicit surfaces make other tasks easy (like inside/outside tests).**

# Algebraic Surfaces (Implicit)

- Surface is zero set of a polynomial in  $x, y, z$
- Examples:



$$x^2 + y^2 + z^2 = 1$$



$$(R - \sqrt{x^2 + y^2})^2 + z^2 = r^2$$



$$\left(x^2 + \frac{9y^2}{4} + z^2 - 1\right)^3 = x^2 z^3 + \frac{9y^2 z^3}{80}$$

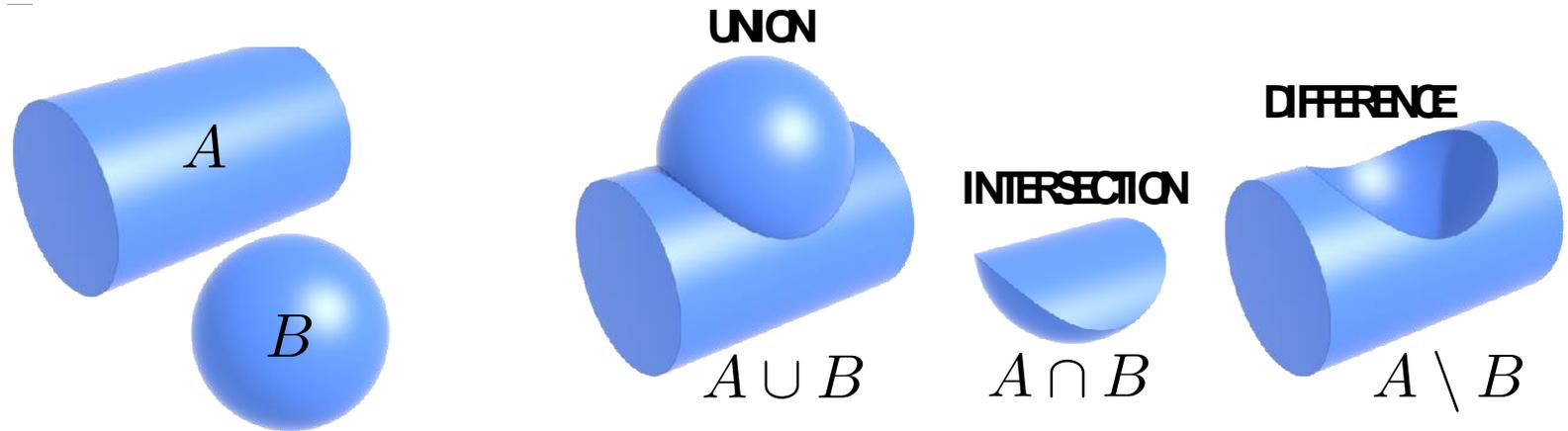
- What about more complicated shapes?
- Very hard to come up with polynomials!



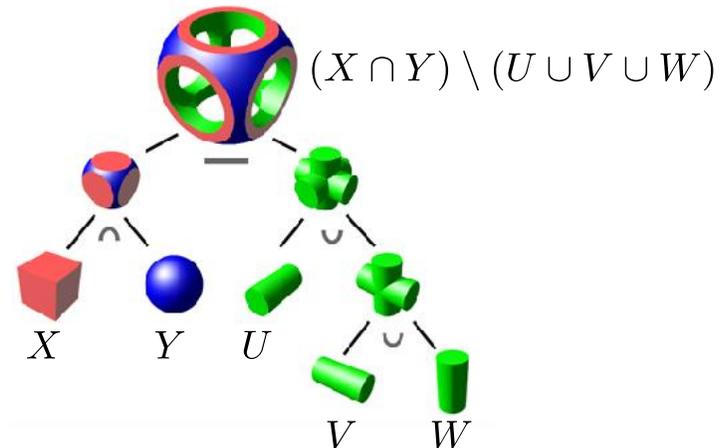
# Constructive Solid Geometry (Implicit)

- Build more complicated shapes via Boolean operations

- Basic operations:

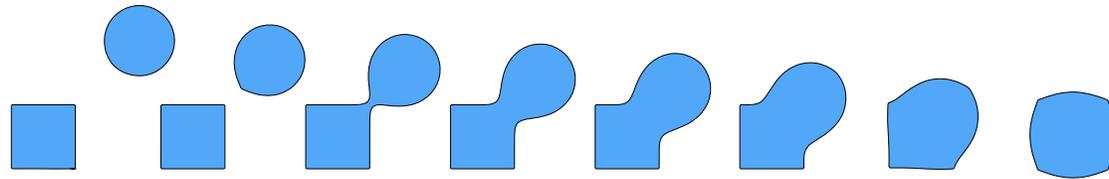


- Then chain together expressions:



# Blending Distance Functions (Implicit)

- A distance function gives distance to closest point on object
- Can blend any two distance functions  $d_1, d_2$ :



- Similar strategy to points, though many possibilities. E.g.,

$$f(x) := e^{d_1(x)^2} + e^{d_2(x)^2} - \frac{1}{2}$$

- Appearance depends on how we combine functions
- Q: How do we implement a Boolean union of  $d_1(x), d_2(x)$ ?
- A: Just take the minimum:  $f(x) = \min(d_1(x), d_2(x))$

# Scene made of pure signed distance functions

## Art with math -- really hard!

Shadertoy  Browse New Sign In



19.59 39.7 fps 1280 x 720

### Slisesix (2008)

Views: 3904, Tags: procedural, 3d, raymarching, distancefield, sdf, demoscene

Created by iq in 2021-07-22

A Shader I made in 2008. While not my first raymarched SDF, this was my first SDF based content creation exercise. Lots of the usual SDF techniques are used, although soft shadows and smooth minimum were not fully evolved yet.

Comments (15)

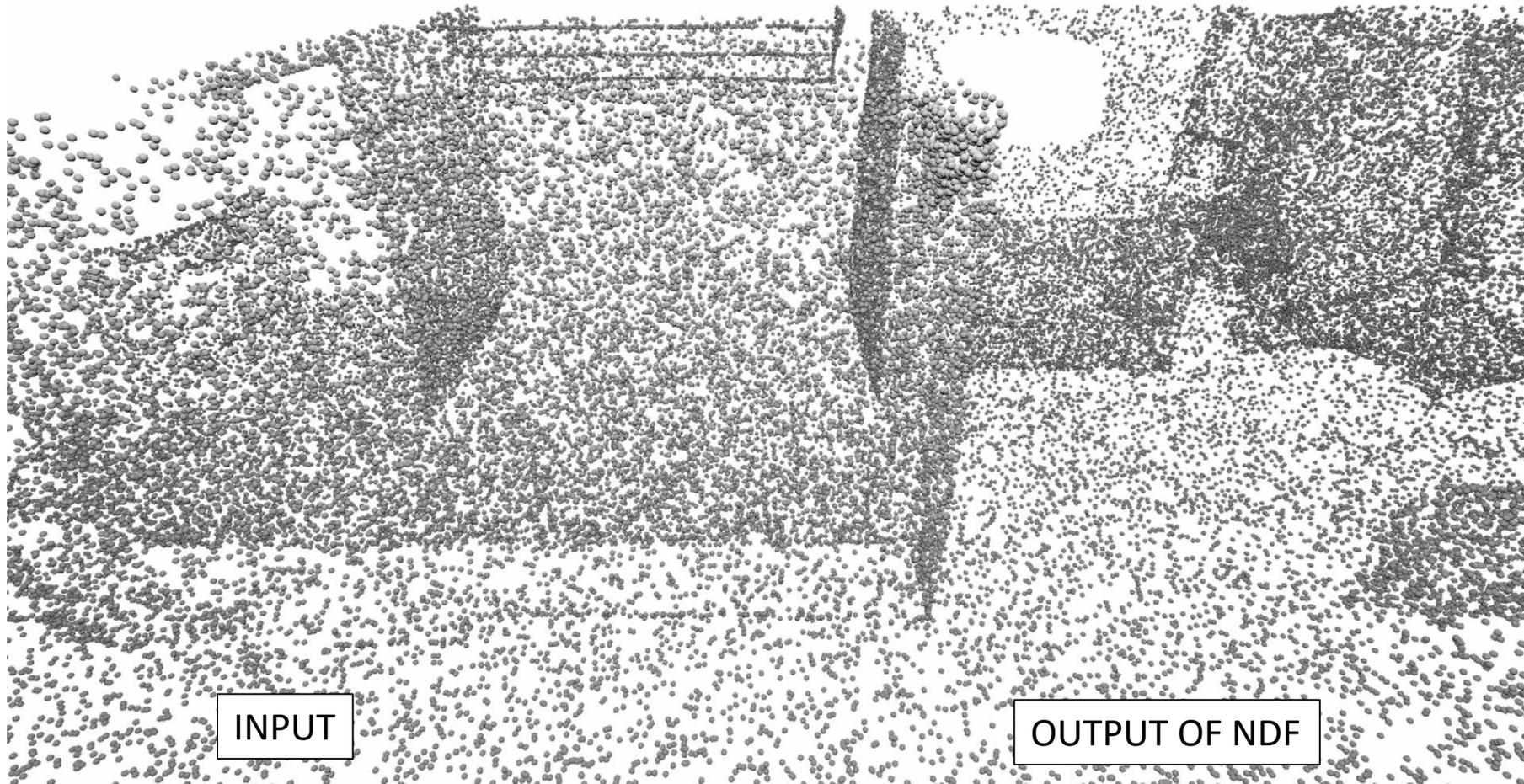
[Sign in](#) to post a comment.

```
127
128 float columna( vec3 pos, float offx )
129 {
130     float x = pos.x;
131     float y = pos.y;
132     float z = pos.z;
133
134     float y2=y-0.40;
135     float y3=y-0.35;
136     float y4=y-1.00;
137
138     float di = udSqBox( vec3(x, y, z), vec3(0.10, 1.00, 0.10) );
139     di = min( di, udSqBox( vec3(x, y, z), vec3(0.12, 0.40, 0.12) ) );
140     di = min( di, udSqBox( vec3(x, y, z), vec3(0.05, 0.35, 0.14) ) );
141     di = min( di, udSqBox( vec3(x, y, z), vec3(0.14, 0.35, 0.05) ) );
142     di = min( di, udSqBox( vec3(x, y4, z), vec3(0.14, 0.02, 0.14) ) );
143     di = min( di, udSqBox( vec3(x-y2)*0.7071, (y2+x)*0.7071, z), vec3(0.10*0.7071, 0.10*0.7071, 0.10*0.7071) );
144     di = min( di, udSqBox( vec3(x, (y2+z)*0.7071, (z-y2)*0.7071), vec3(0.12, 0.10*0.7071, 0.10*0.7071) );
145     di = min( di, udSqBox( vec3(x-y3)*0.7071, (y3+x)*0.7071, z), vec3(0.10*0.7071, 0.10*0.7071, 0.10*0.7071) );
146     di = min( di, udSqBox( vec3(x, (y3+z)*0.7071, (z-y3)*0.7071), vec3(0.14, 0.10*0.7071, 0.10*0.7071) );
147
148     #if 1
149     float fb = fb( vec3(10.1-w+offx, 10.1-w, 10.1-w) );
```

Compiled in 0.1 secs 8868 chars

iChannel0 iChannel1 iChannel2 iChannel3

# Nowadays -- Neural Distance Fields (NDF)



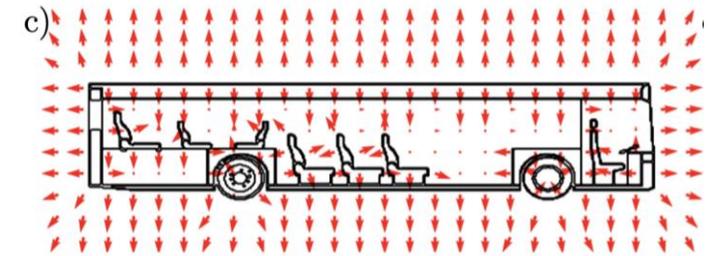
# Distance field, field normals and closest points

- **Distance to surface** is given by the distance field itself  $f(\mathbf{p})$

$$f(\mathbf{p}) = \min_{\mathbf{q} \in \mathcal{S}} \|\mathbf{p} - \mathbf{q}\| \quad \mathcal{S} = \{\mathbf{p} \in \mathbb{R}^3 \mid f(\mathbf{p}) = 0\}$$

- **Surface normal** are the gradients of the function

$$n(\mathbf{p}) = \nabla f(\mathbf{p})$$

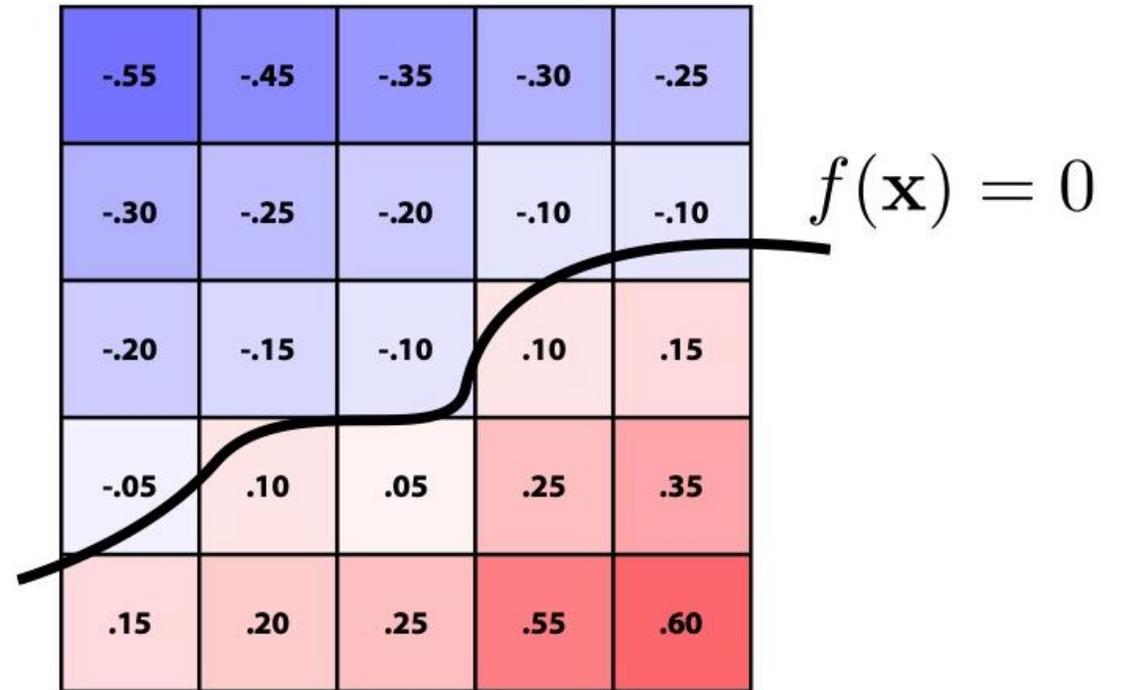


- **Closest points** are found trivially as:

$$\mathbf{q} = \mathbf{p} - f(\mathbf{p}) \nabla_{\mathbf{p}} f(\mathbf{p})$$

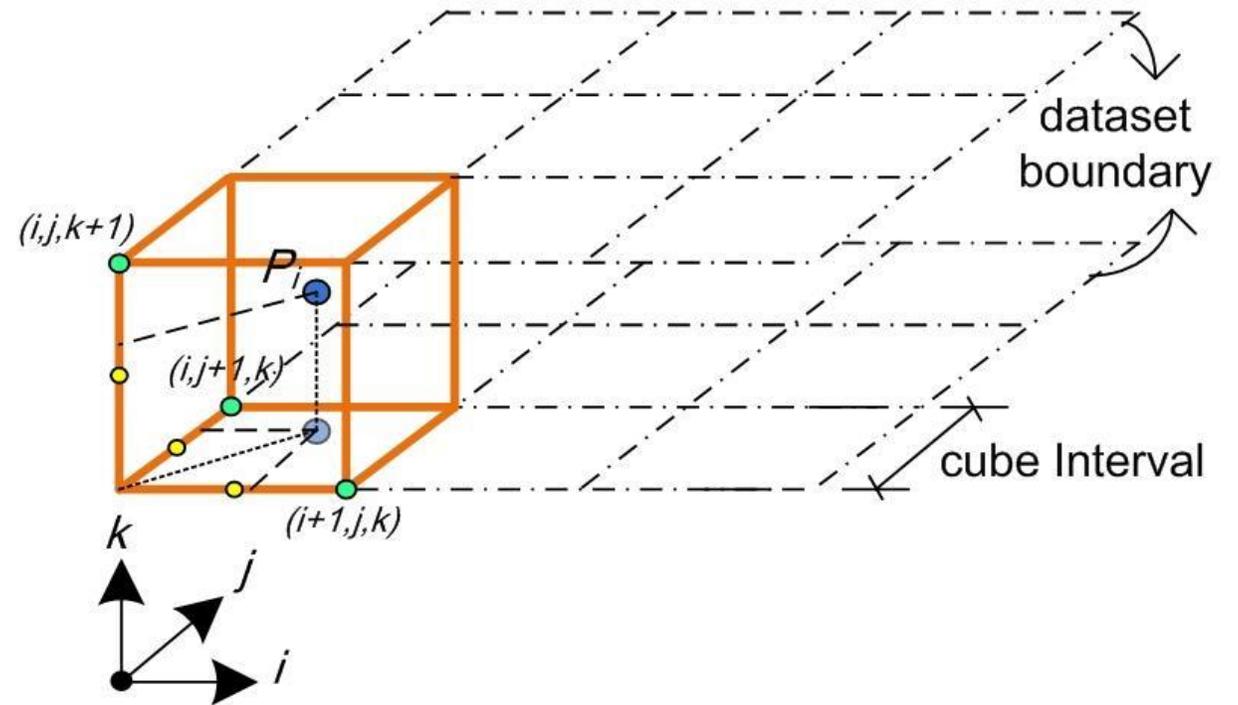
# Surface Reconstruction from Implicit Representation

- Implicit surfaces have some nice features (e.g., merging/splitting)
- But, hard to describe complex shapes in closed form
- Alternative: store a grid of values approximating function
- Surface is found where interpolated values equal zero



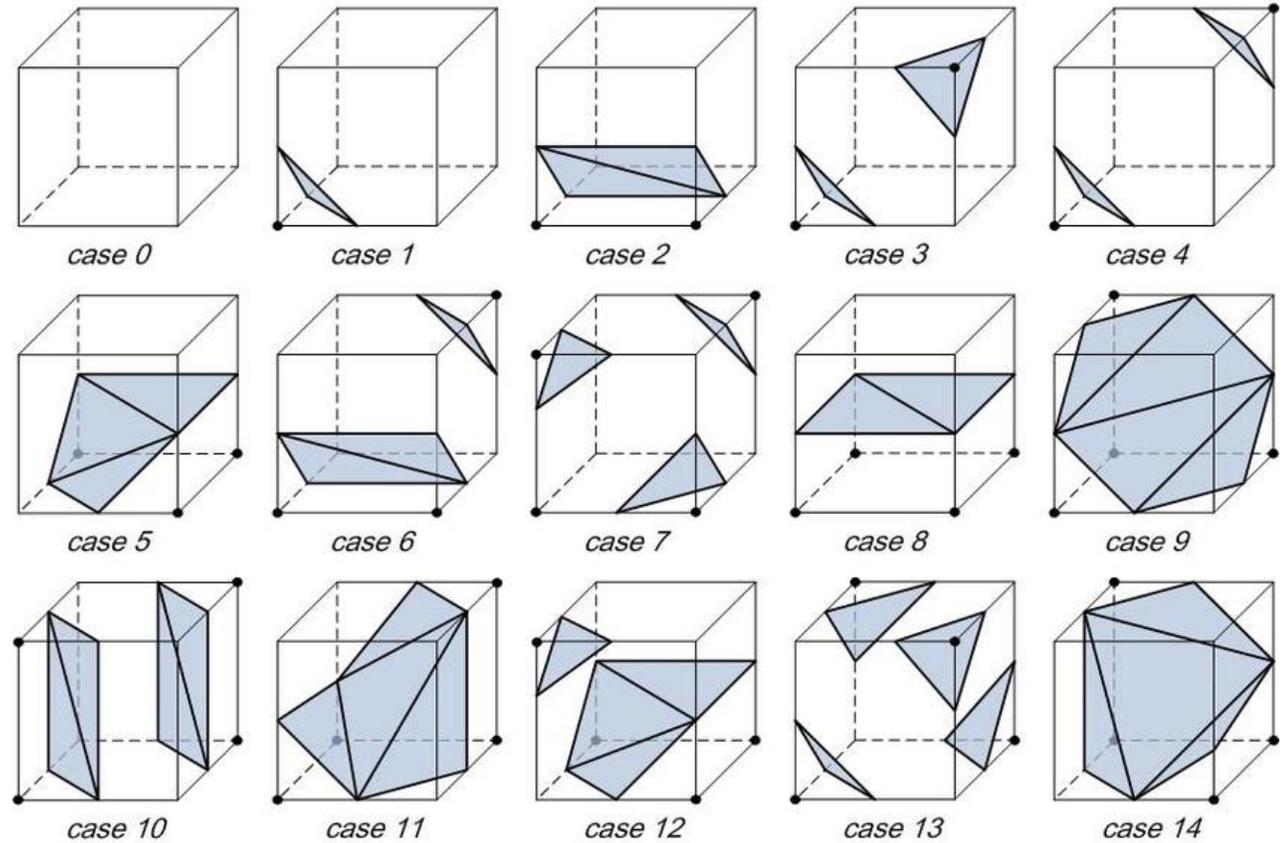
# Marching Cubes

1. Read four slices into memory.
2. Scan two slices and create a cube from four neighbors on one slice and four neighbors on the next slice.
3. ...
4. ...
5. ...
6. ...
7. ...



# Marching Cubes

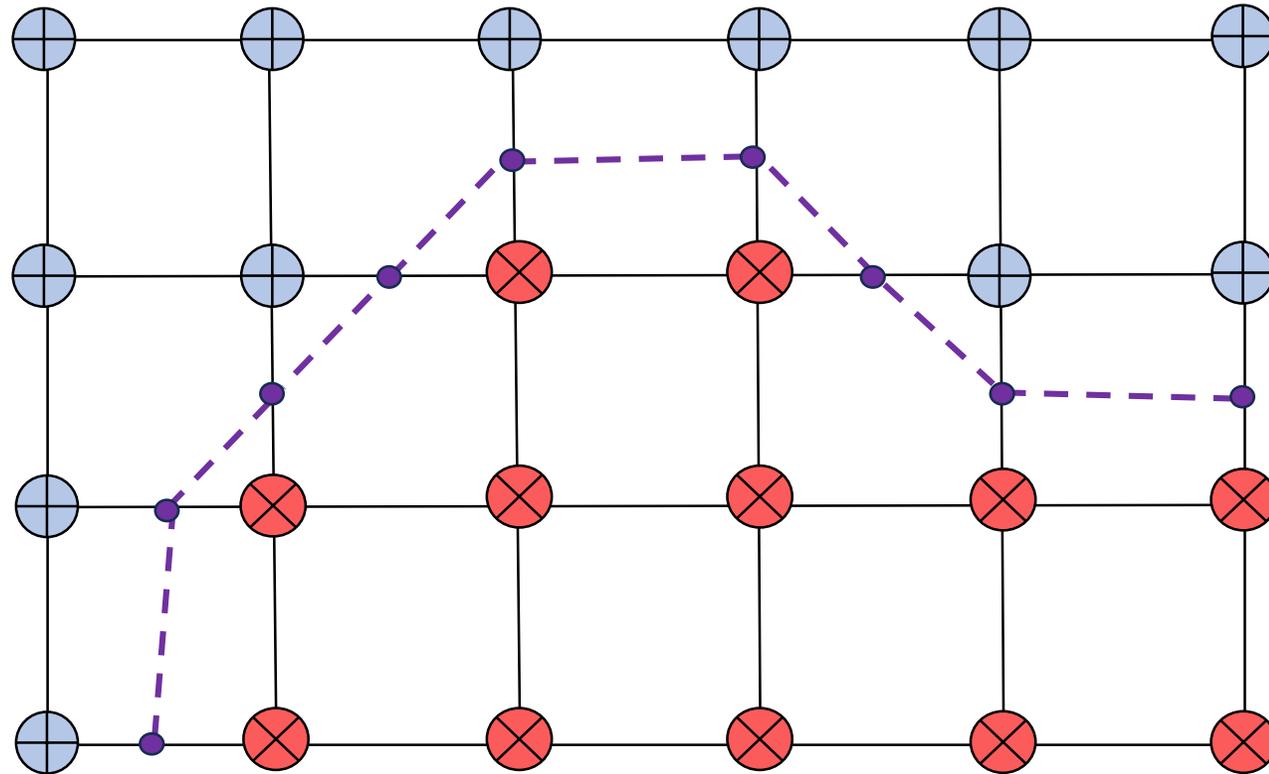
1. ...
2. ...
3. Calculate an index for the cube by comparing the eight density values at the cube vertices with the surface constant.
4. Using the index, look up the list of edges from a precalculated table.



# Marching Cubes

1. ...
2. ...
3. ...
4. ...
5. Using the densities at each edge vertex, find the surface-edge intersection via linear interpolation
6. Calculate a unit normal at each cube vertex using central differences. Interpolate the normal to each triangle vertex
7. Output the triangle vertices and vertex normals.

# Marching Cubes



# Challenges with Marching Cubes

- Generating one polygon for each portion of the contour
- Can only approximate the restriction of the contour
- Fails to capture the sharp detail of the surface
- Difficult to simplify the generated surface mesh
- Becoming challenging to extract surface for higher grid sizes (3D uniform grid induces cubic cost in resolution)

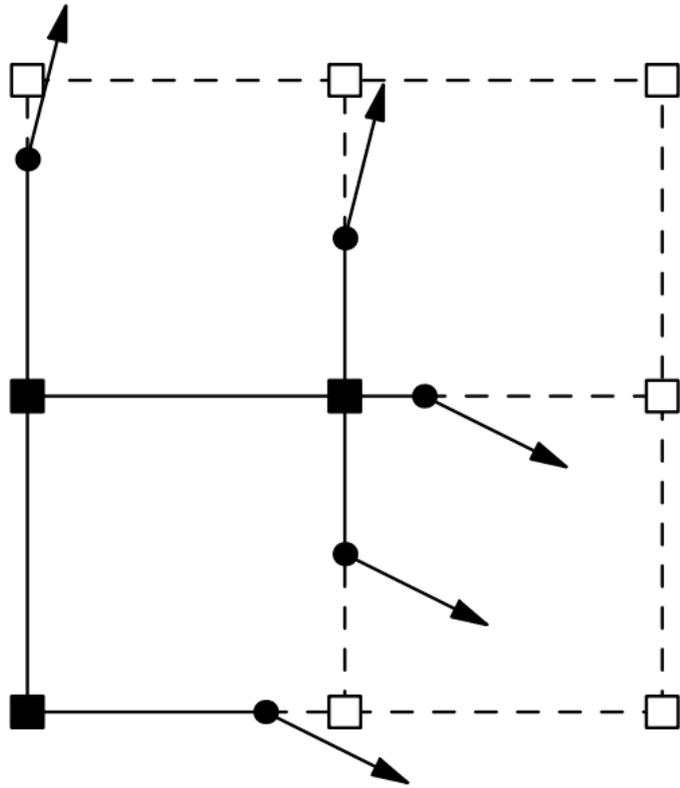
# Dual Contouring

- Can capture sharp edges along corners of surface
- Can leverage gradient information along edge intersections to reproduce a wide class of polyhedral shapes as well as curved or sharp edges
- Utilizes octree in place of a 3D uniform grid for better efficiency
- Uses normals to define Quadratic Error Function for each leaf of the octree, inspired by original Extended Marching Cubes (EMC)

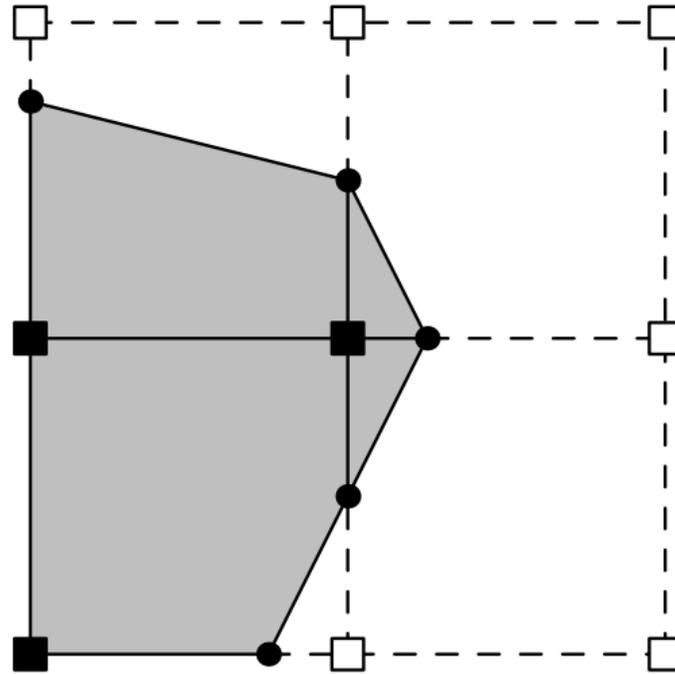
$$E[x] = \sum_i (n_i \cdot (x - p_i))^2$$

where  $p_i$  and  $n_i$  correspond to the intersections and normals of the contour with the edges of cube

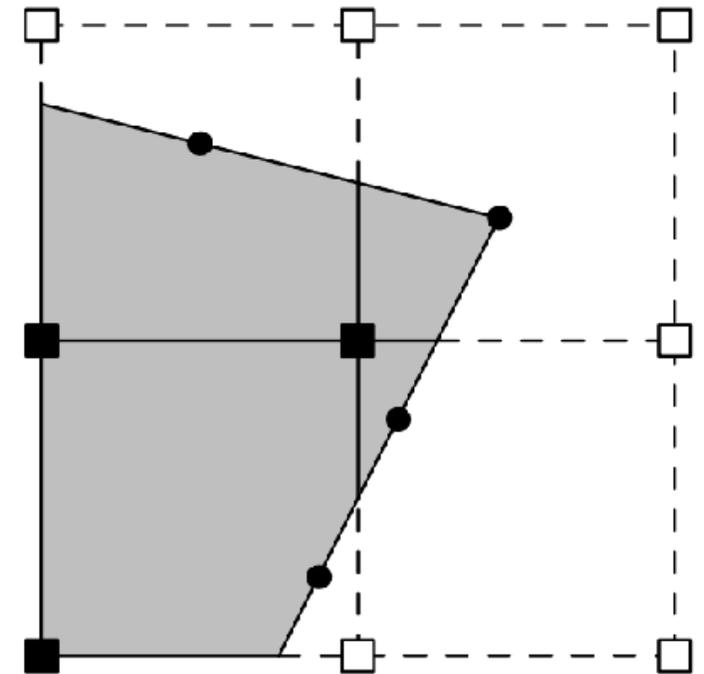
# Dual Contouring



Signed Grid with edges  
tagged by Hermite Data



Marching Cubes Contour



Dual Contour

# Challenges with Dual Contouring

- Limited expressivity due to reliance on structured grids
- Cannot faithfully extract very thin surfaces and features (otherwise requiring very fine grid resolution)
- Limited by the amount of surface simplification

# Dual Marching Cubes

- Align features of the implicit function with the features of the structured grid; enabling extraction of sharp features
- Can generate adaptive polygonalizations that are crack-free and topologically manifold
- Can reproduce thin features in surface without excessive subdivision of the octree

# Dual Marching Cubes

DMC's QEF is calculated by computing a tangent planes to the implicit surface representation  $f$  on a grid of points  $(x_i, y_i, z_i)$  sampled over  $c$ .

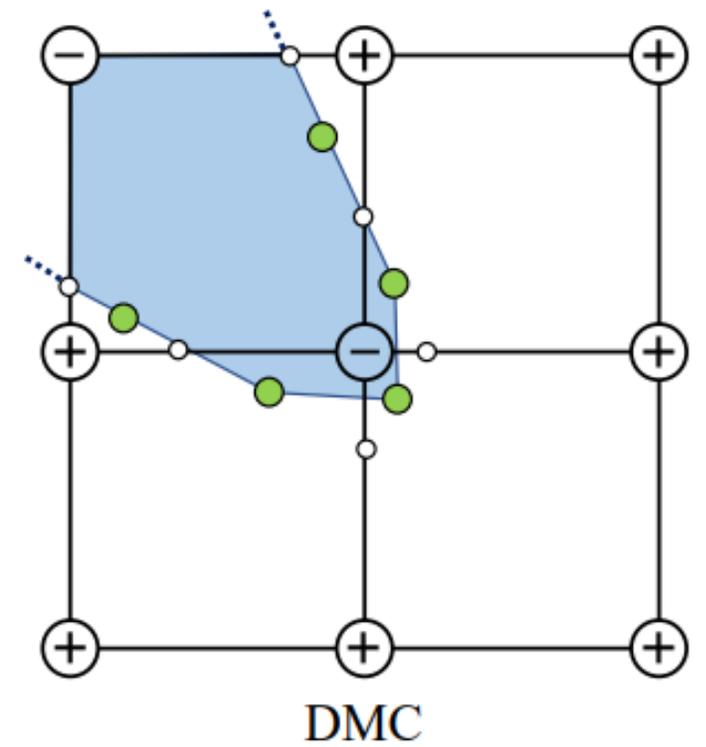
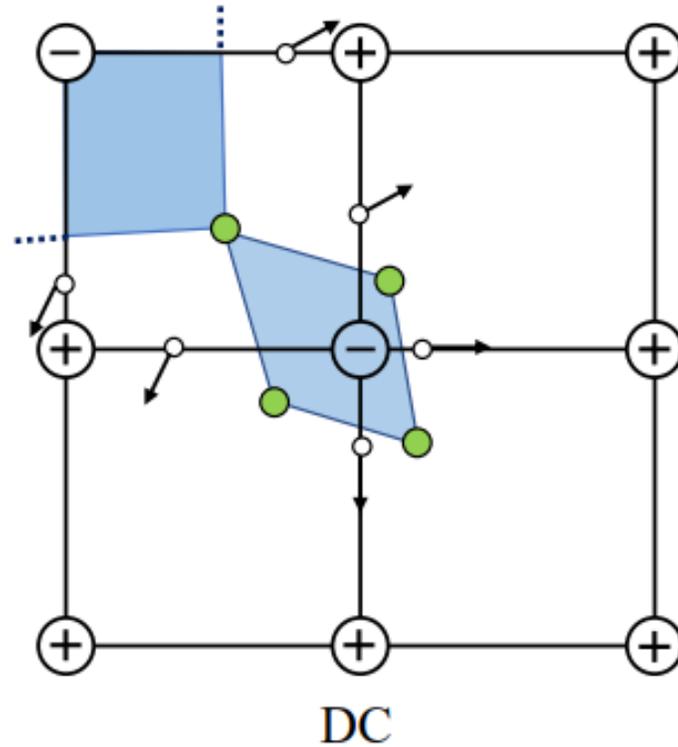
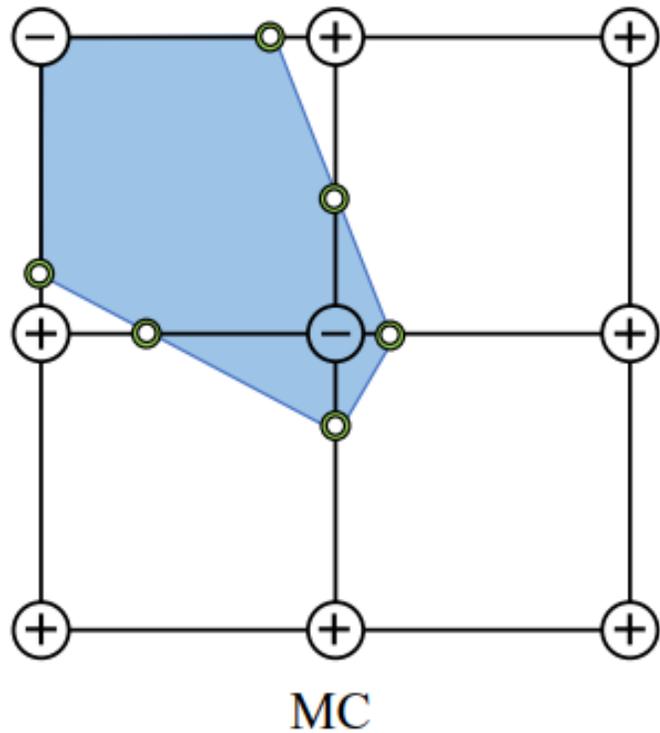
Tangent plane's equation:  $T_i(x, y, z) = \nabla f(x_i, y_i, z_i) \cdot ((x, y, z) - (x_i, y_i, z_i))$

DMC's QEF sums over all sample points yielding:

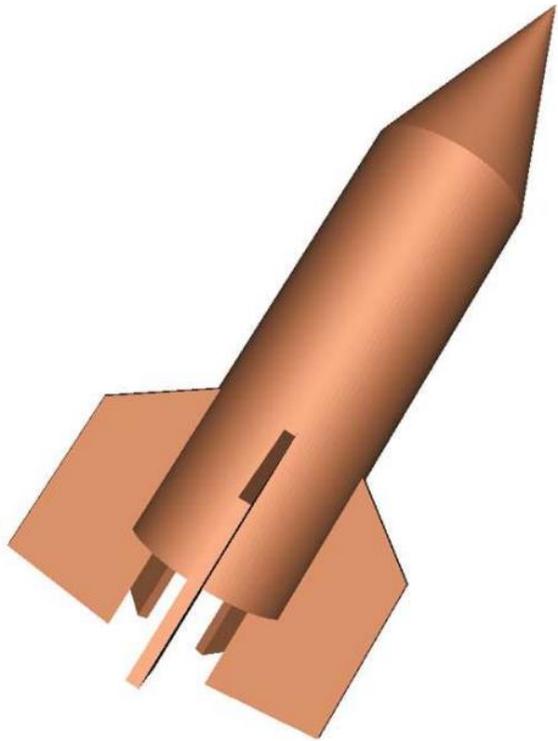
$$E(f(x, y, z), x, y, z) = \sum_i (f(x, y, z) - T_i(x, y, z))^2 / (1 + |\nabla f(x_i, y_i, z_i)|^2)$$

where the denominator normalizes the contribution of each tangent plane

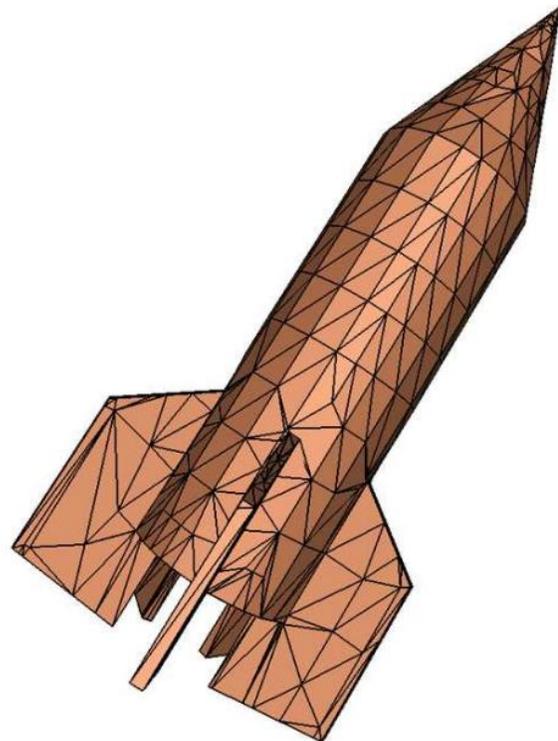
# Comparison between Marching Cubes & Dual Contouring & Dual Marching Cubes



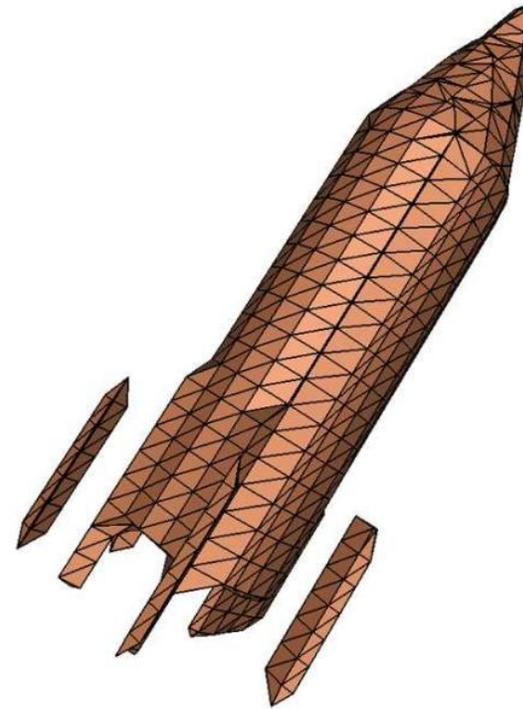
# Comparison between Marching Cubes & Dual Contouring & Dual Marching Cubes



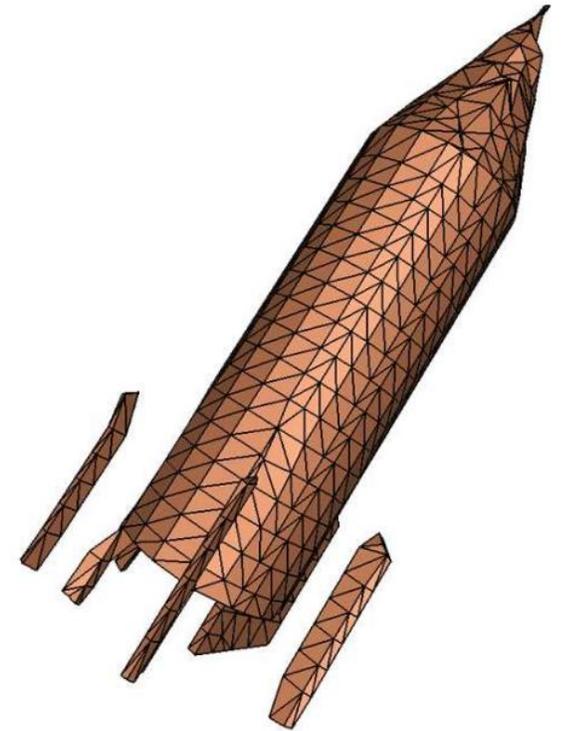
Ground Truth  
Model



Dual Marching  
Contour



Marching  
Contour

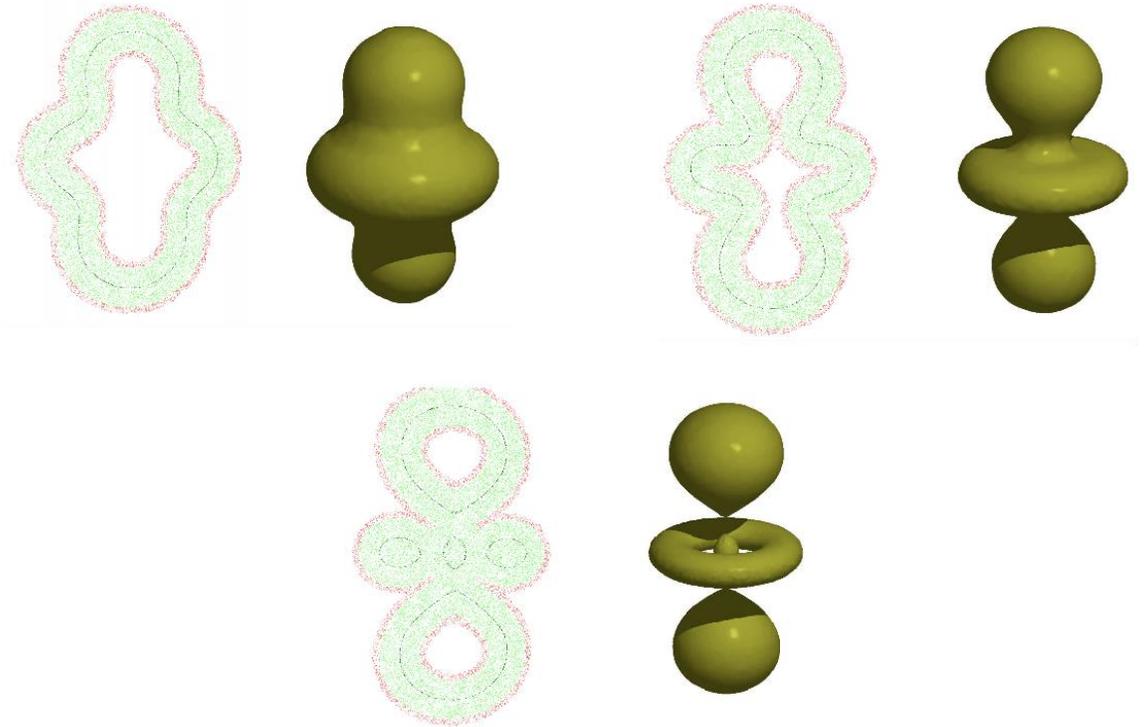


Dual  
Contour

# Challenges faced by Level Sets method

Although Level Set representation provides much more explicit control over shape (like a texture), it

- runs into problems of aliasing! (Unlike closed-form expressions)
- induces  $O(n^3)$  storage costs for 3D surface extraction
- can save space by only storing a narrow band around the surface.



# Implicit Representations - Pros & Cons

## Pros:

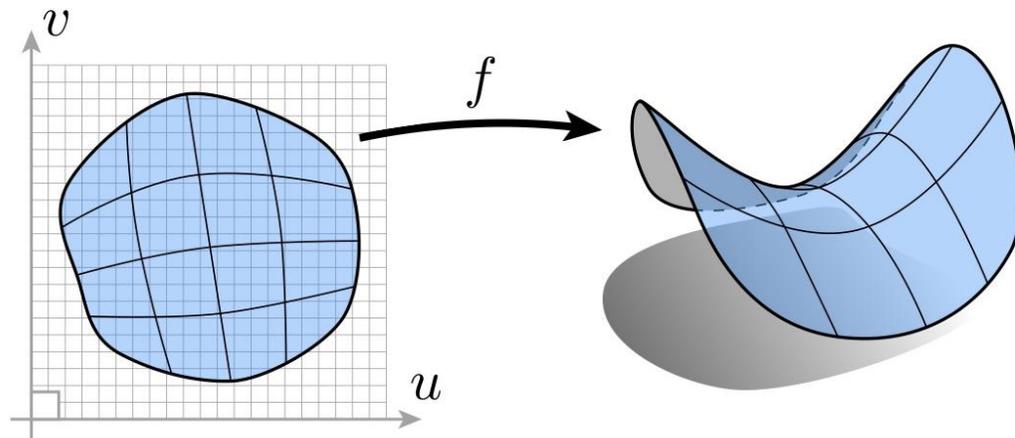
- description can be very compact (e.g., a polynomial)
- easy to determine if a point is in our shape (just plug it in!)
- other queries may also be easy (e.g., distance to surface)
- for simple shapes, exact description/no sampling error
- easy to handle changes in topology (e.g., fluid)

## Cons:

- expensive to find all points in the shape (e.g., for drawing)
- very difficult to model complex shapes

# "Explicit" Representations of Geometry

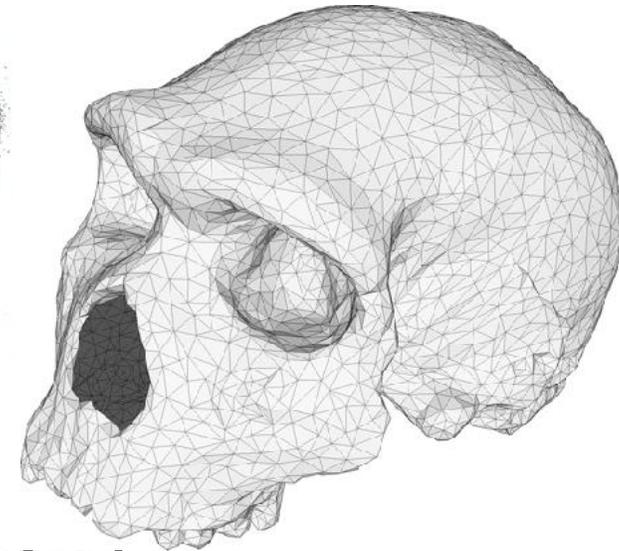
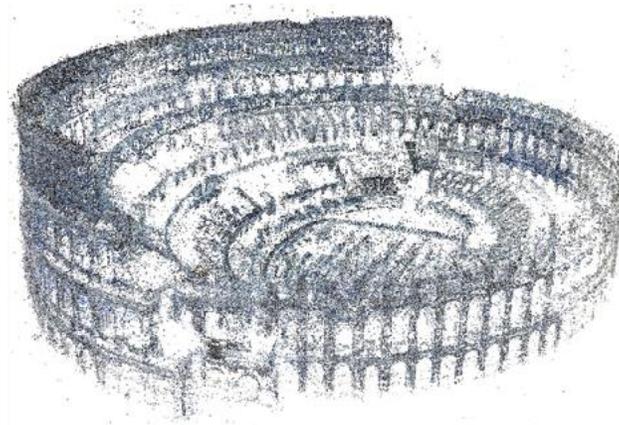
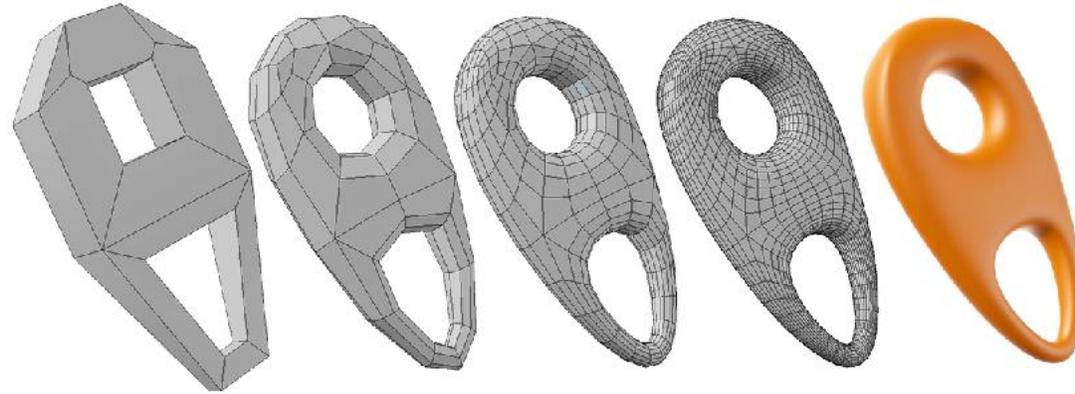
- All points are given directly
- E.g., points on sphere are  $(\cos(u) \sin(v), \sin(u) \sin(v), \cos(v))$ ,  
for  $0 \leq u < 2\pi$  and  $0 \leq v \leq \pi$
- More generally:  $f : \mathbb{R}^2 \rightarrow \mathbb{R}^3; (u, v) \mapsto (x, y, z)$



- (Might have a bunch of these maps, e.g., one per triangle!)

# Many explicit representations in graphics

- triangle meshes
- polygon meshes
- subdivision surfaces
- NURBS
- point clouds
- ...



(Will see some of these a bit later.)

**But first, let's play a game:**

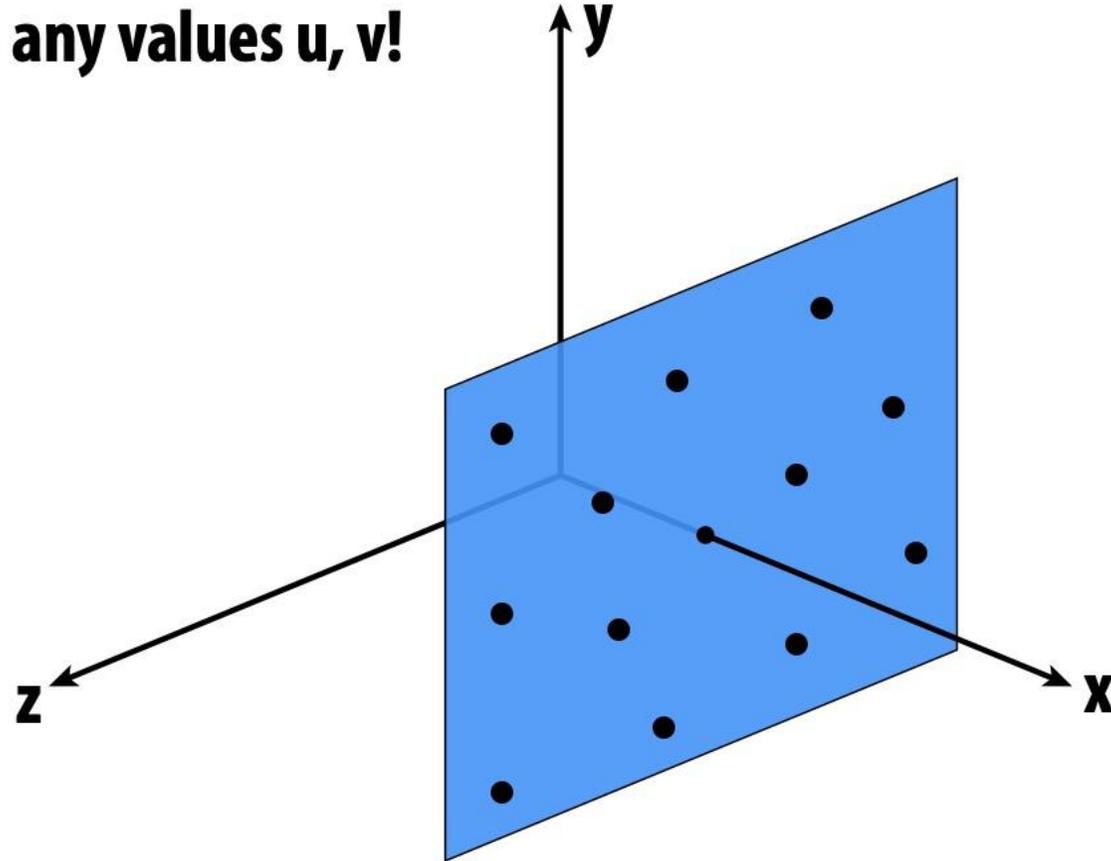
**I'll give you an explicit surface.**

**You give me some points on it.**

# Sampling an explicit surface

My surface is  $f(u, v) = (1.23, u, v)$ .

Just plug in any values  $u, v$ !



Explicit surfaces make some tasks easy (like sampling).

**Let's play another game.**

**I have a new surface  $f(u,v)$ .**

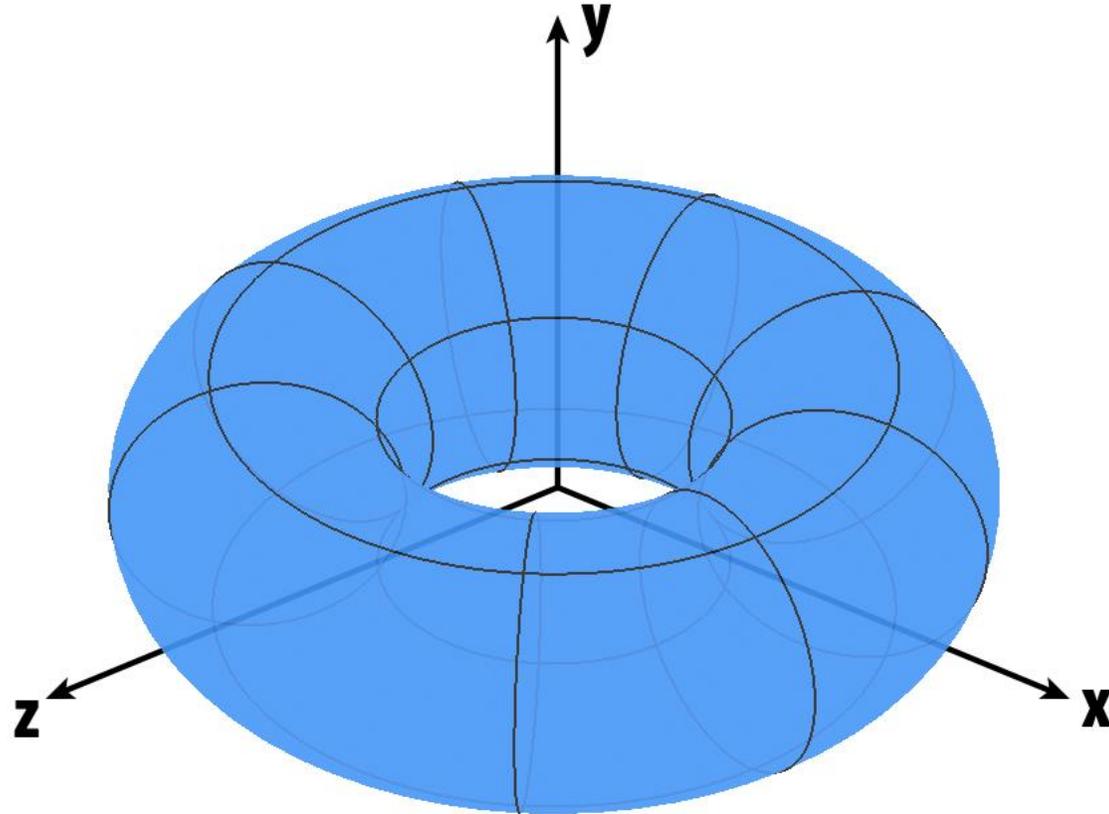
**I want to see if a point is *inside* it.**

# Check if this point is inside the torus

My surface is  $f(u,v) = ( (2+\cos u)\cos v, (2+\cos u)\sin v, \sin u )$

How about the point  $(1.96, -0.39, 0.9)$ ?

**...NO!**



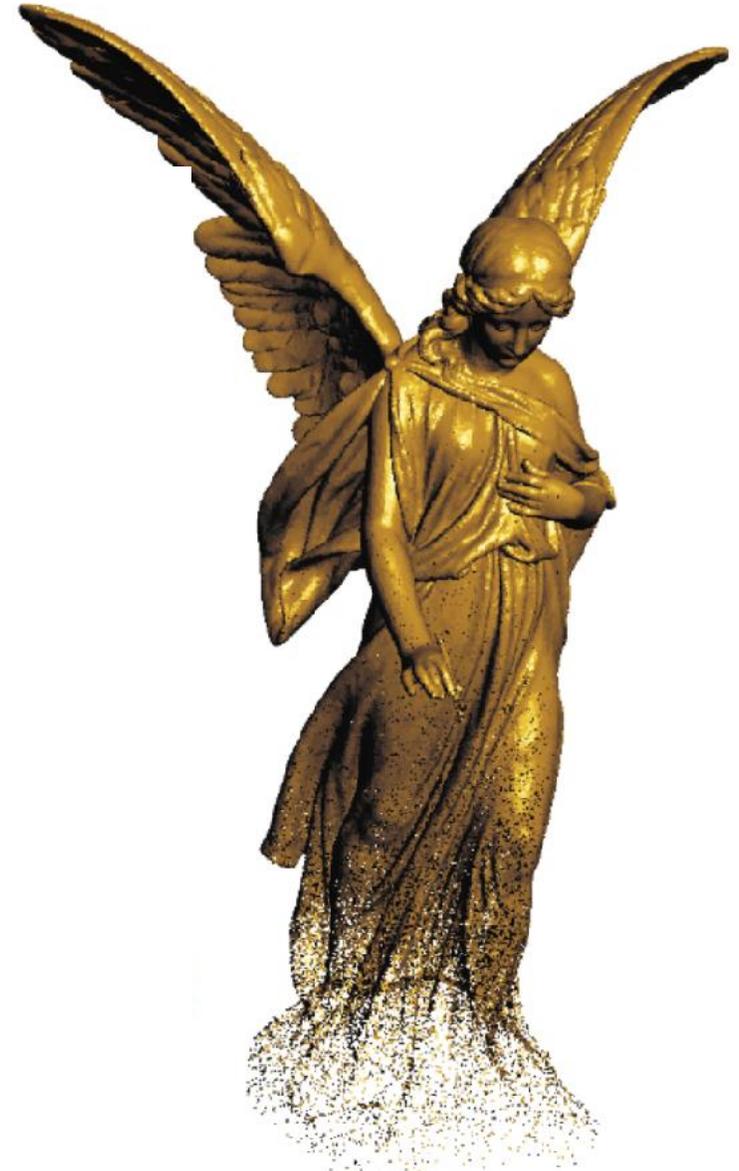
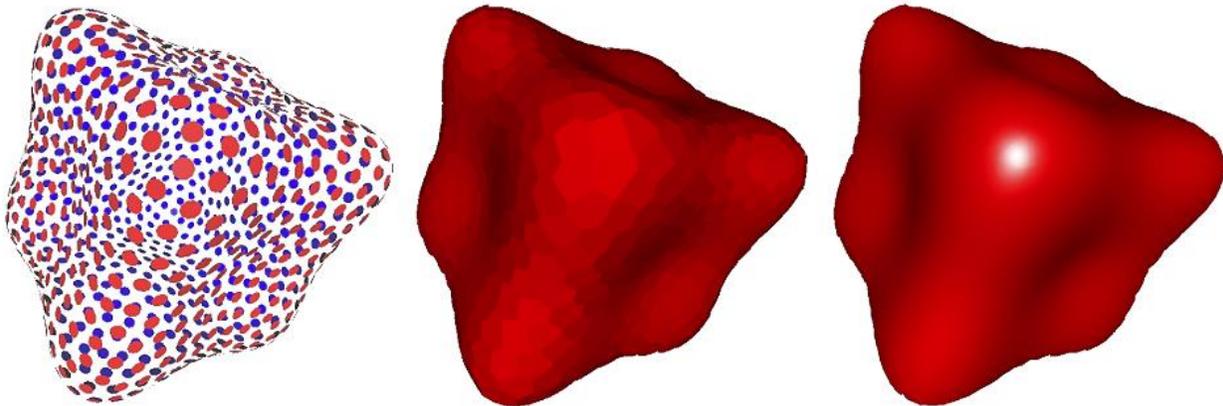
**Explicit surfaces make other tasks hard (like inside/outside tests).**

## **CONCLUSION:**

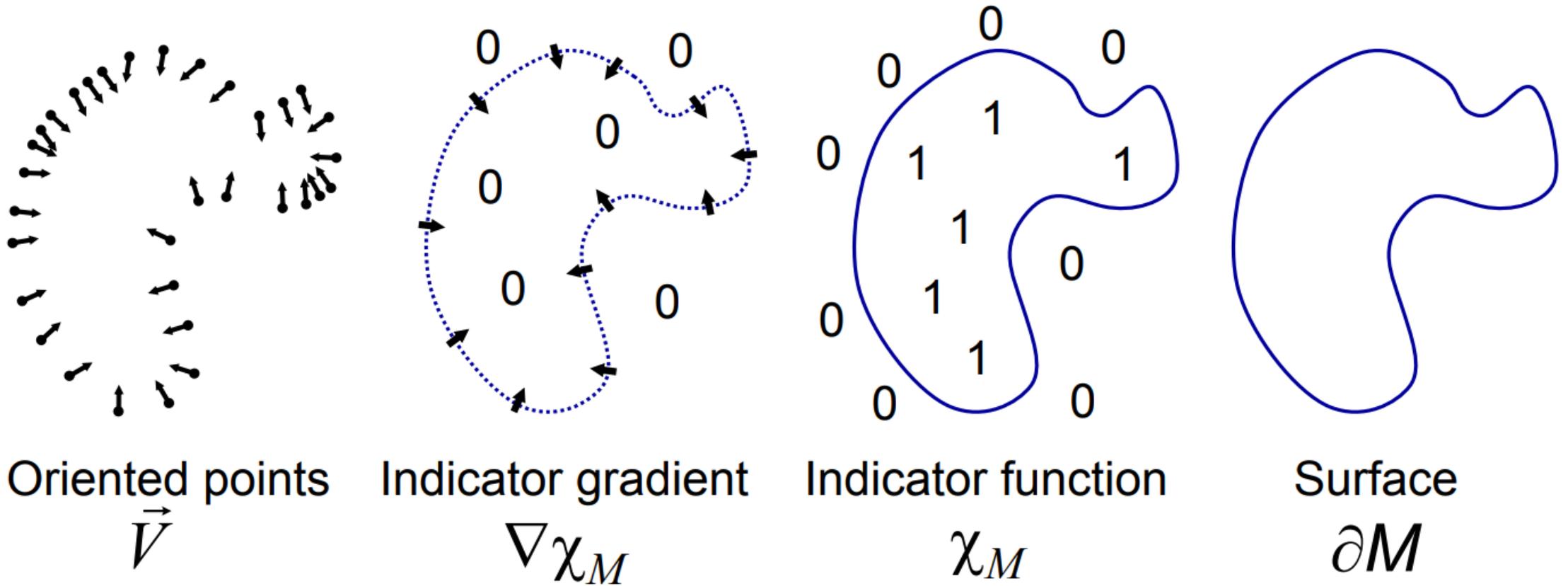
**Some representations work better than others—depends on the task!**

# Point Clouds

- Easiest representation: list of points  $(x,y,z)$
- Often augmented with normals
- Easily represent any kind of geometry
- Easy to draw dense cloud ( $\gg 1$  point/pixel)
- Hard to interpolate under-sampled regions
- Hard to do processing / simulation/...



# Poisson Surface Reconstruction



# Poisson Surface Reconstruction

Given an oriented point cloud  $P$  with points  $p_1, \dots, p_n$  and corresponding (outward-facing) normal observations  $n_1, \dots, n_n$ , we consider the implicit reconstruction task of finding a function  $f : \mathbb{R}^d \rightarrow \mathbb{R}$  such that

$$f(p_i) = 0, \nabla f(p_i) = n_i, \forall i \in \{1, \dots, n\}.$$

PSR first utilizes convolutional kernel  $F$  to interpolate  $n_i$  into a vector field  $V : \mathbb{R}^d \rightarrow \mathbb{R}^d$  defined in box  $B$  containing  $P$  as

$$V(x) = \sum_i F(x, x_i) n_i$$

# Poisson Surface Reconstruction

$f : \mathbb{R}^d \rightarrow \mathbb{R}$  is then defined as a function whose gradient best matches  $V$

$$f = \operatorname{argmin}_g \int_B \|V(x) - \nabla g(x)\|_2 dx$$

Variational problem above is equivalent to the Poisson Equation

$$\Delta f = \nabla \cdot V(x)$$

which is solved by PSR via discretization using the Finite Element Method on an octree and solve using a purpose-built multigrid algorithm

# Poisson Surface Reconstruction

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which is solved by PSR via discretization using the Finite Element Method on an octree and solve using a purpose-built multigrid algorithm. In Screened PSR (Kazhdan et.al, 2013) authors use

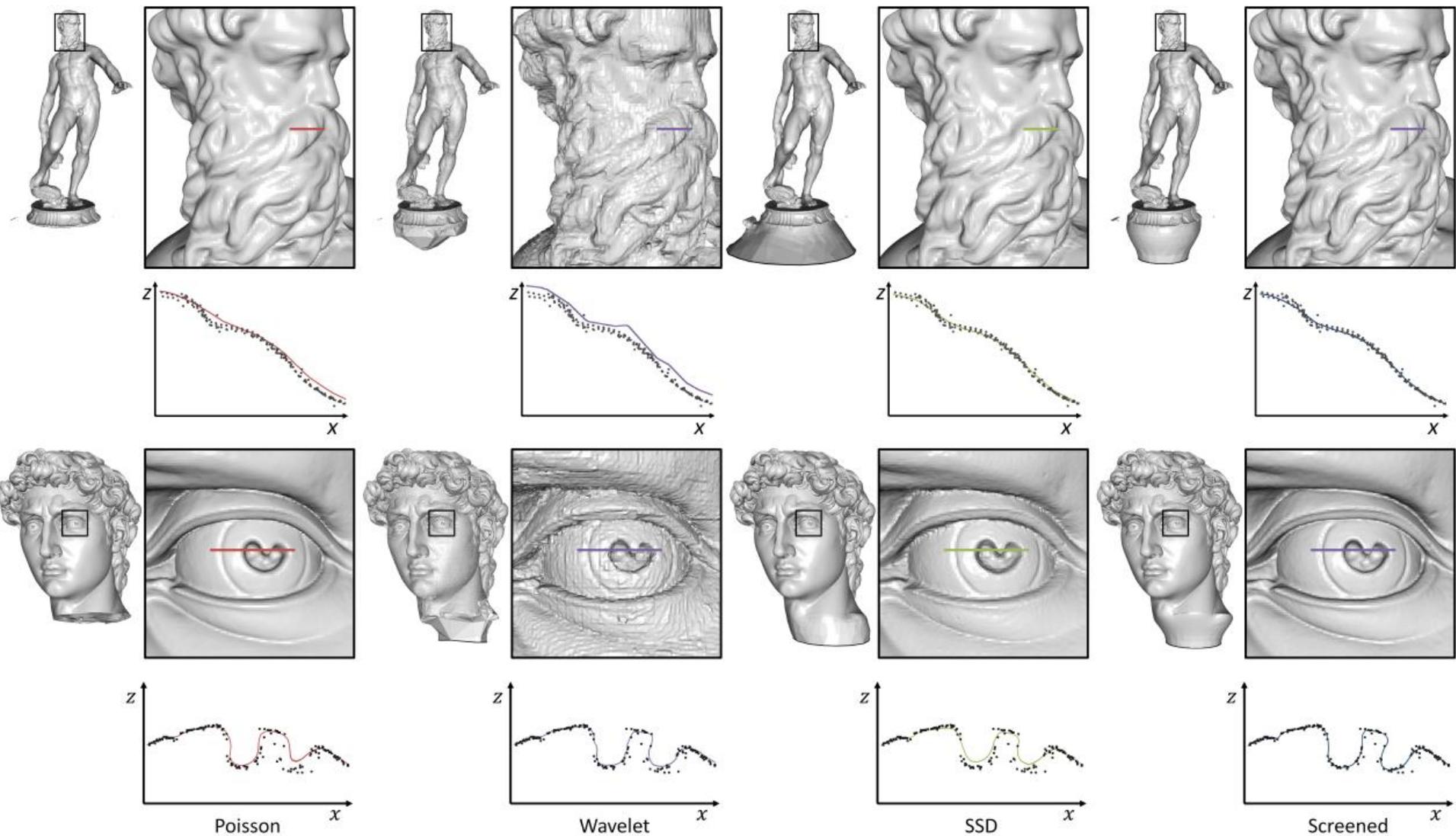
$$f = \operatorname{argmin}_g \int_B \|V(x) - \nabla g(x)\|_2 dx + \alpha * \sum_i g(p_i)^2$$

# (Screened) Poisson reconstruction

- Can extract the surface from oriented point clouds
- (Screened PSR) efficiently implements sparsity constraints
- Can better capture the details of the underlying surface
- (Screened PSR) implements efficient octree structure and multigrid algorithm to reduce the time complexity

However, these methods limit themselves to outputting the likeliest surface given the point cloud observations and their assumed prior. Whereas in reality there could be theoretically infinite number of possible surfaces representing same set of Points with normals

# (Screened) Poisson reconstruction

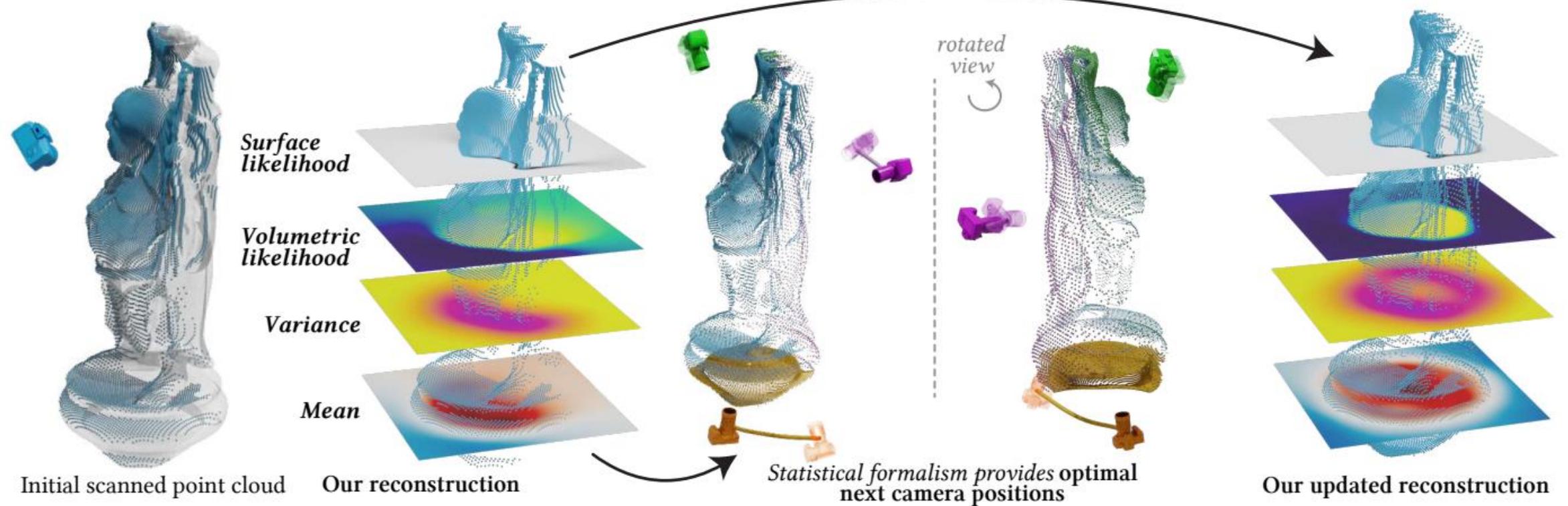


# Neural Stochastic Screened Poisson Surface Reconstruction

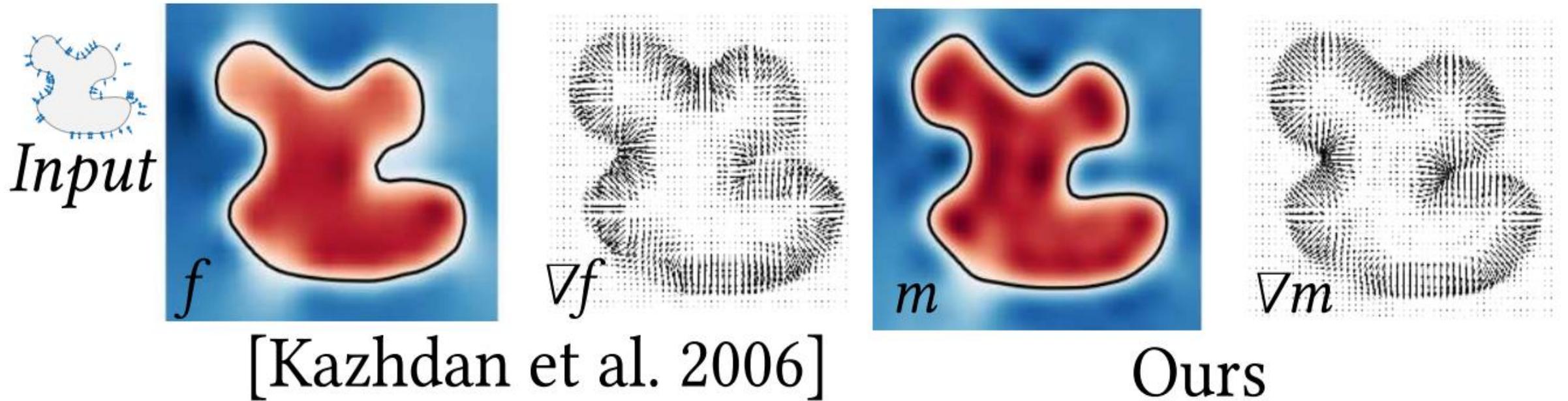
- Take a step even further and models the posterior multivariate Gaussian distributions representing the likelihood of all possible reconstructions for the given Point Set with normals
- Proposes the parametrization of mean and covariance functions using neural networks, optimizing them through gradient-based methods for a more efficient and flexible approach
- Enabling computation of the statistical quantities over the reconstruction

# Neural Stochastic Screened Poisson Surface Reconstruction

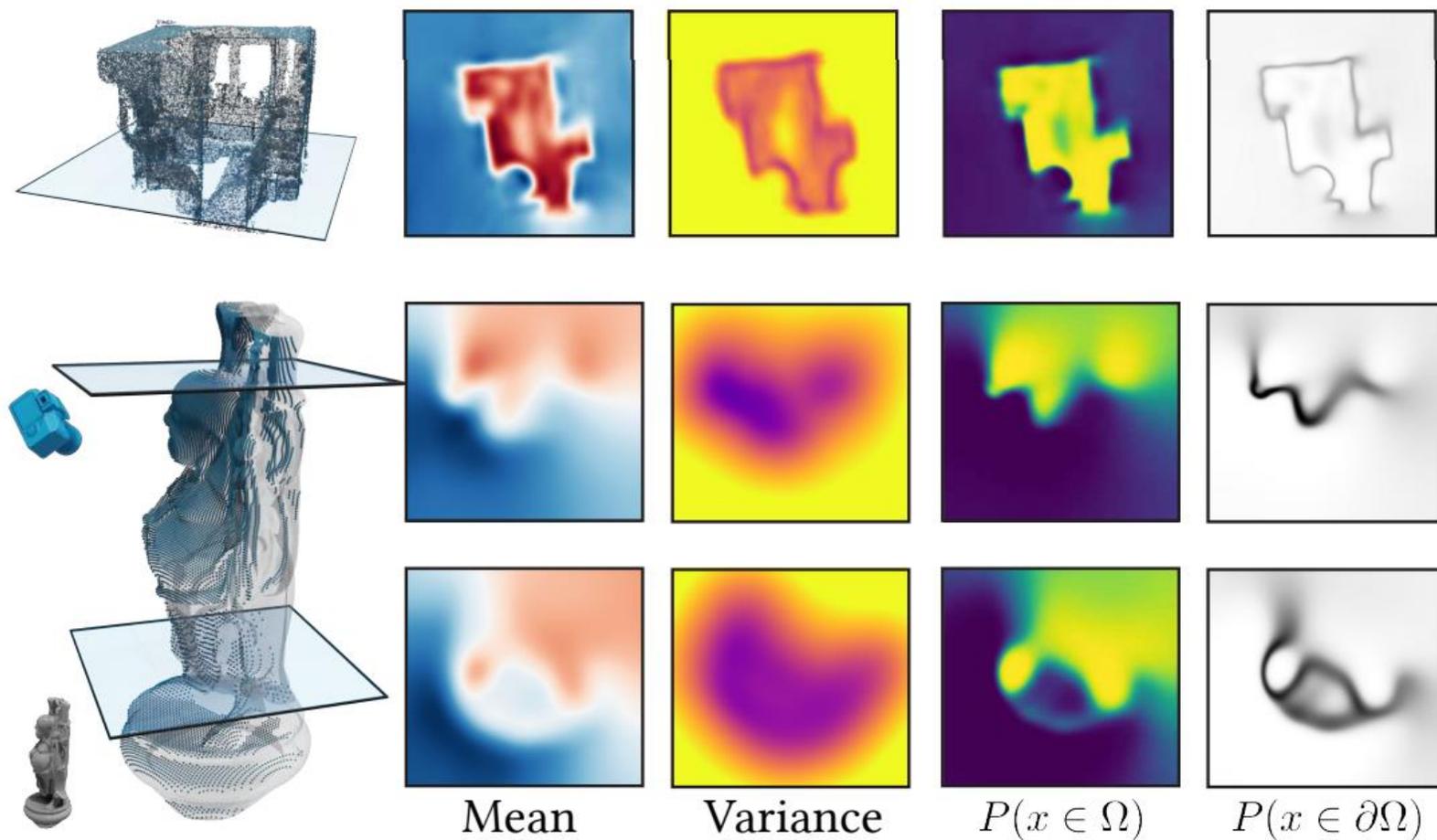
*Fine-tune our pretrained model to obtain new reconstruction*



# Neural Stochastic Screened Poisson Surface Reconstruction

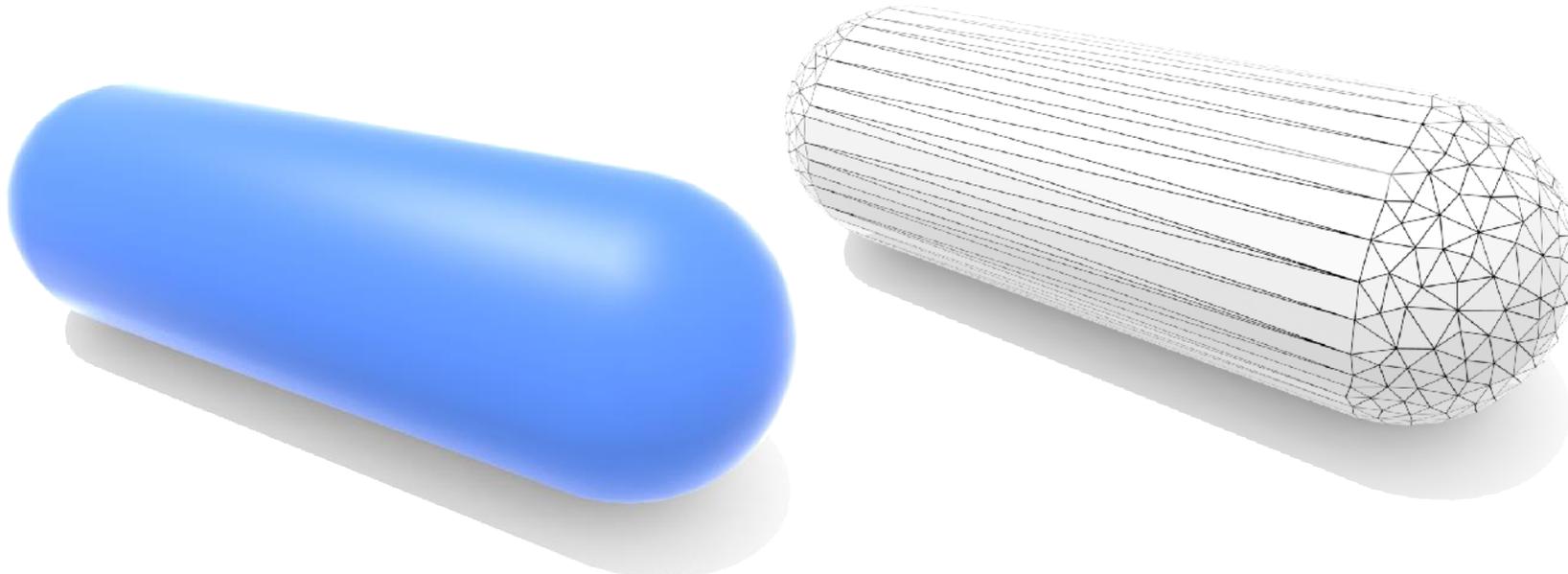


# Neural Stochastic Screened PSR



# Polygon Mesh

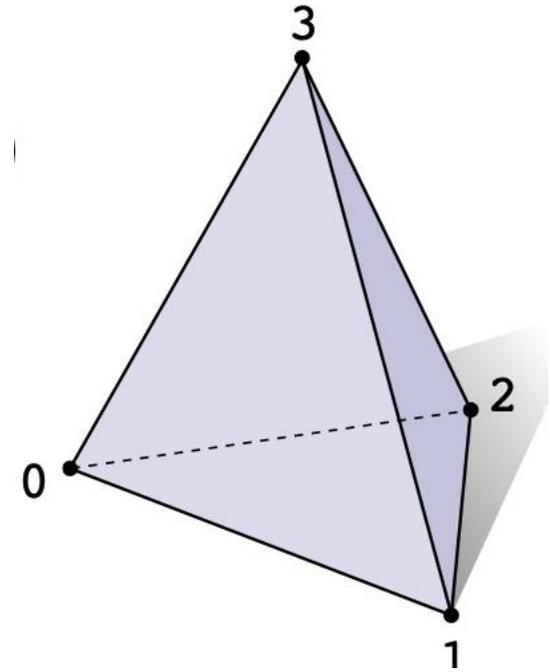
- Store vertices and polygons (most often triangles or quads)
- Easier to do processing/simulation, adaptive sampling
- More complicated data structures
- Irregular neighbourhoods



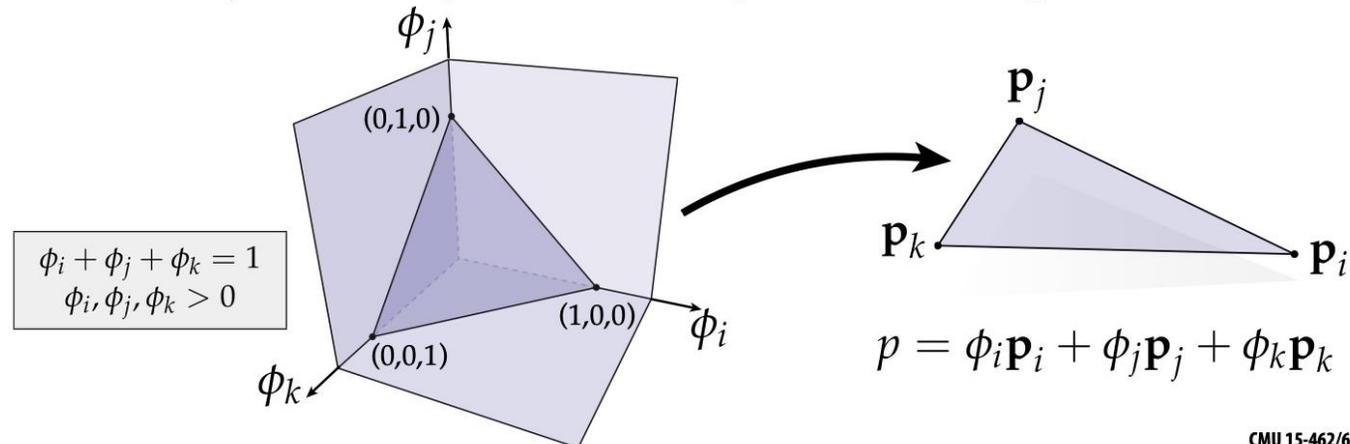
# Triangle Mesh (Explicit)

- Store vertices as triples of coordinates (x,y,z)
- Store triangles as triples of indices (i,j,k)
- E.g., tetrahedron:

	VERTICES			TRIANGLES		
	x	y	z	i	j	k
0:	-1	-1	-1	0	2	1
1:	1	-1	1	0	3	2
2:	1	1	-1	3	0	1
3:	-1	1	1	3	1	2



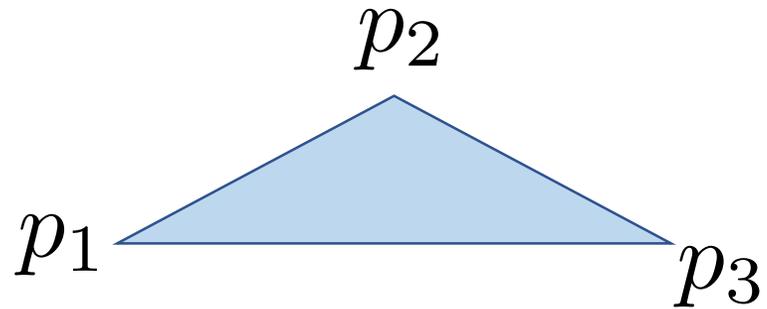
- Use barycentric interpolation to define points inside triangles:



# Triangle Mesh – Normals (Explicit)

- Option1: Normals per face (Phong)

$$n = (p_2 - p_1) \times (p_3 - p_1)$$



- Option2: Interpolate incident face normal at each vertex, and interpolate with barycentric interpolation for points inside the triangle

$$n(\phi_1, \phi_2, \phi_3) = \phi_1 n_1 + \phi_2 n_2 + \phi_3 n_3$$

# Distance queries and closest points

- Smooth parametric surface

- Evaluation  $f(u, v)$

- Normals  $\mathbf{n}(u, v) = f_u(u, v) \times f_v(u, v)$

- Distance and closest point. Find surface point  $f(u, v)$  such that:

$$[\mathbf{q} - f(u, v)] \times \mathbf{n}(u, v) = \mathbf{0}$$

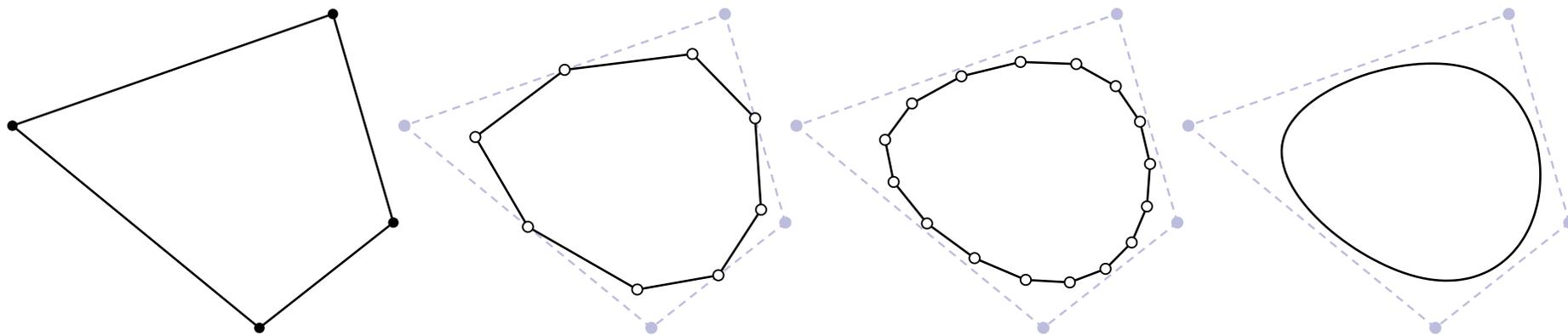
- Mesh

- Distance and closest point:

- Find closest triangle and then find closest point in triangle, and then finding point in triangle
    - Use Kd-tree, AAB Tree, to find closest triangles

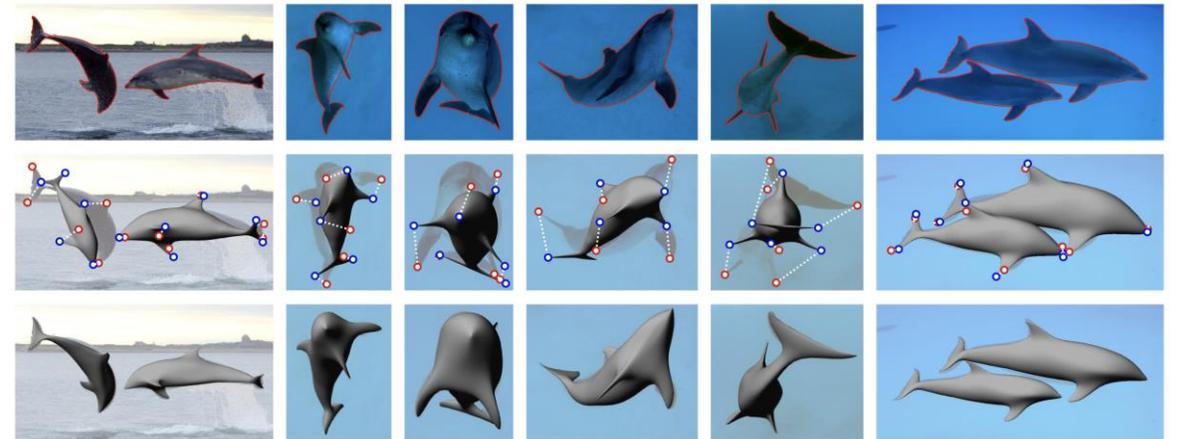
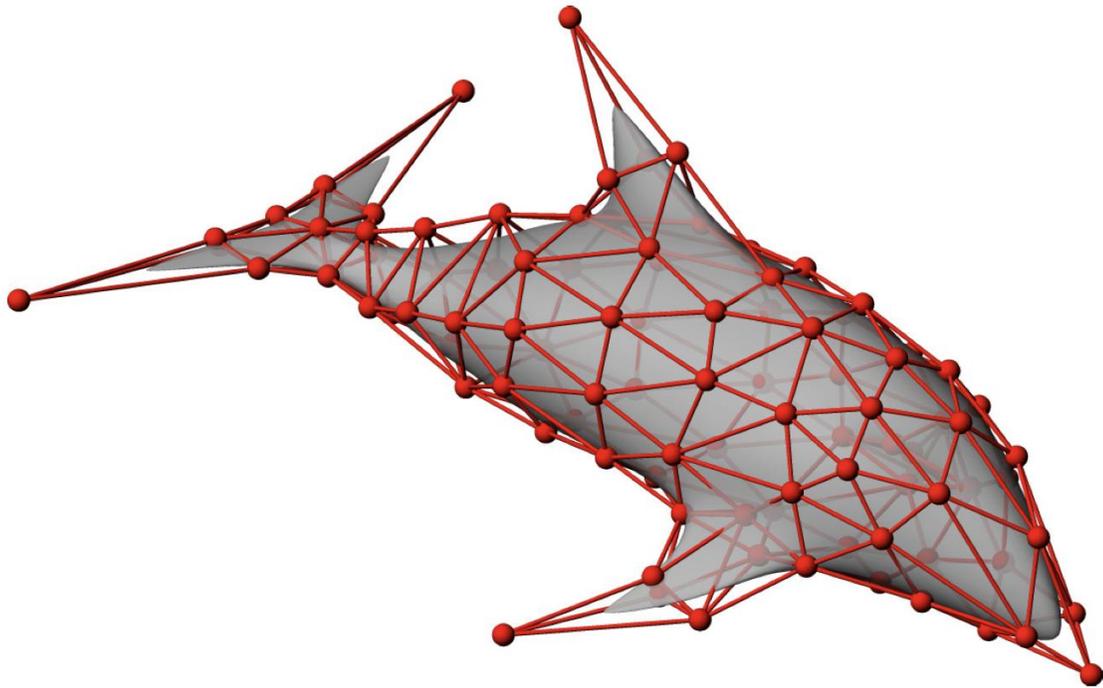
# Subdivision

- Alternative starting point for curves/surfaces: subdivision
- Start with "control curve"
- Repeatedly split, take weighted average to get new positions
- For careful choice of averaging rule, approaches nice limit curve
  - Often exact same curve as well-known spline schemes!



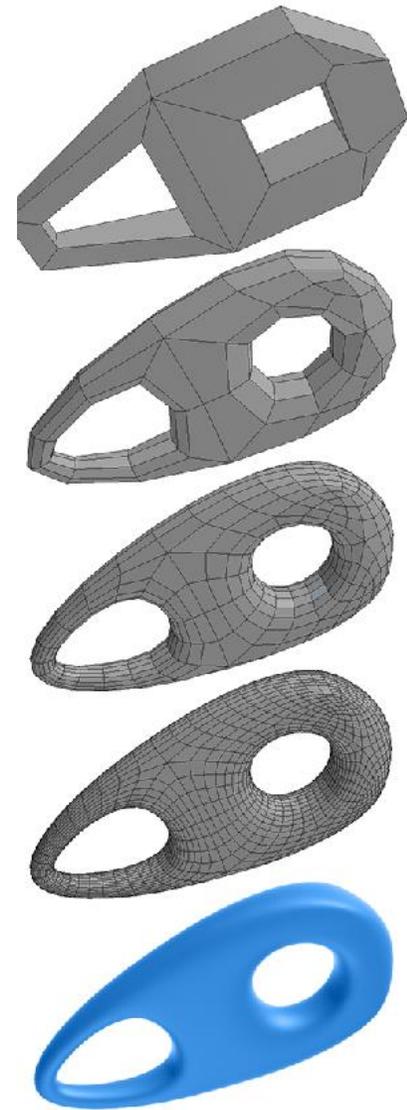
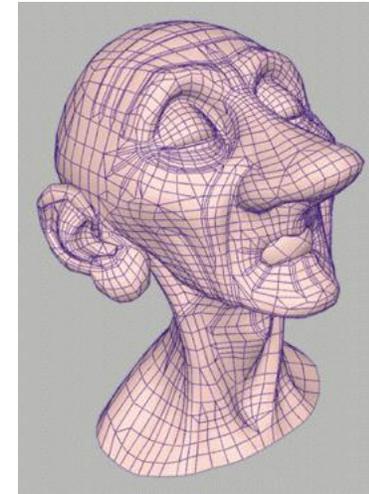
**Q: Is subdivision an explicit or implicit representation?**

# Subdivision Surfaces in Computer Vision



# Subdivision Surfaces (Explicit)

- Start with coarse polygon mesh ("control cage")
- Subdivide each element
- Update vertices via local averaging
- Many possible rules:
  - Catmull-Clark (quads)
  - Loop (triangles)
  - ...
- Common issues:
  - interpolating or approximating?
  - continuity at vertices?
- Easier than splines for modeling; harder to evaluate pointwise
- Widely used in practice (2019 Academy Awards!)



# Subdivision in Pixar (Pixar's "Geri's Game")



see: de Rose et al, "Subdivision Surfaces in Character Animation"

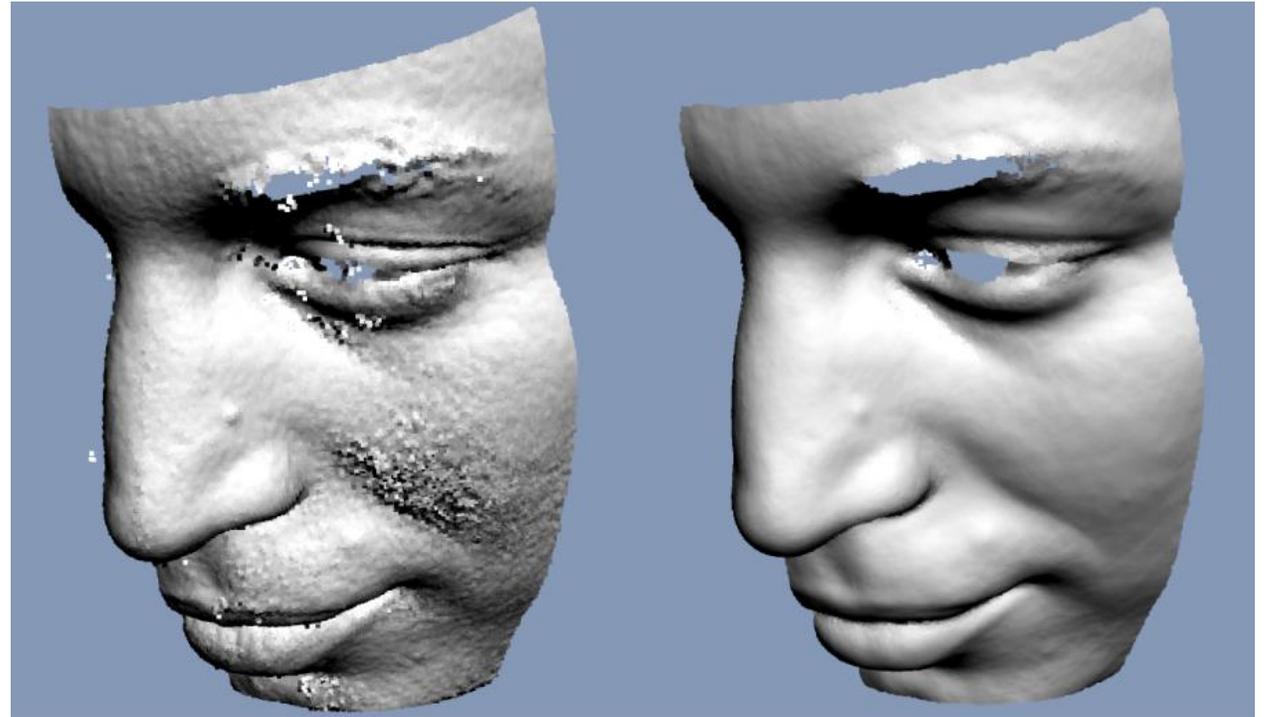
# Slide credits and further reading

- Keenan Crane – Computer Graphics Lecture CMU 15-462/662.  
Lecture 09: Introduction to Geometry.

# Point Cloud Denoising

Problems with Poisson  
Surface Reconstruction:

- cannot handle noise in input point cloud
- uneven distribution of points can make gradient estimation unreliable



# Locally Optimal Projection Operator

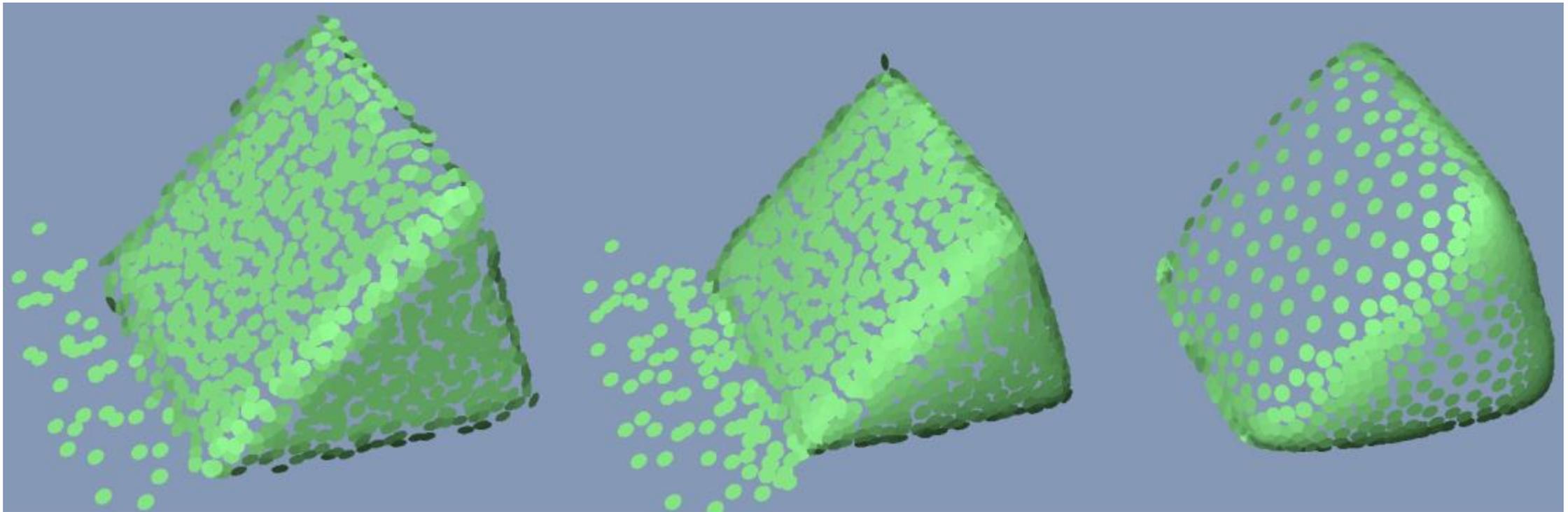
Given an unorganized set of points  $P = \{p_j\}$  for  $j \in J \subset \mathbb{R}^3$ , we define a set of projected points  $X = \{x_i\}$  for  $i \in I \subset \mathbb{R}^3$  by a fixed point iteration to minimize

$$\sum_{i \in I} \sum_{j \in J} \|x_i - p_j\|_2 * \theta(\|\xi_{ij}^k\|) + \\ + \lambda_i \sum_{i' \in I \setminus \{i\}} \eta(\|x_i - x_{i'}^k\|) \theta(\|\delta_{ii'}^k\|)$$

where  $\xi_{ij}^k = x_i^k - p_j$ ,  $\delta_{ii'}^k = x_i^k - x_{i'}^k$ ,  $\theta(r) = e^{-r^{**2}/(h/4)^{**2}}$  and  $\eta(r) = (1/3r^3)$ .

In practice,  $n = |I|$  is often significantly smaller than  $m = |J|$ .

# Locally Optimal Projection Operator



Contaminated  
Point Set

MLS Projections

LOP Projections

# Locally Optimal Projection Operator

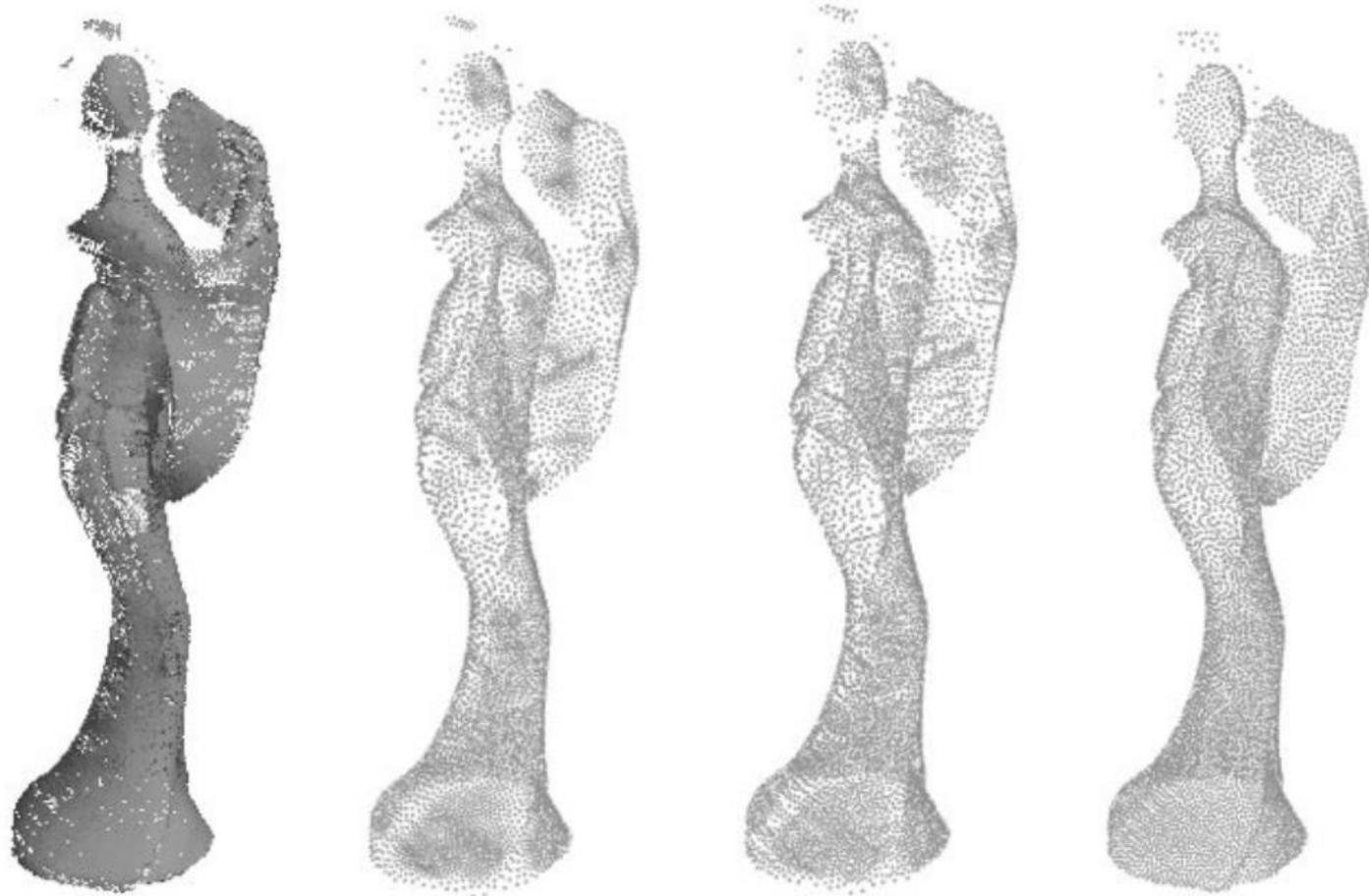
However, LOP has the following disadvantages:

- Repulsion term  $\eta(r)$  decreases too drastically
- Imposes strong attraction between point clusters

Weighted LOP improves over original LOP by

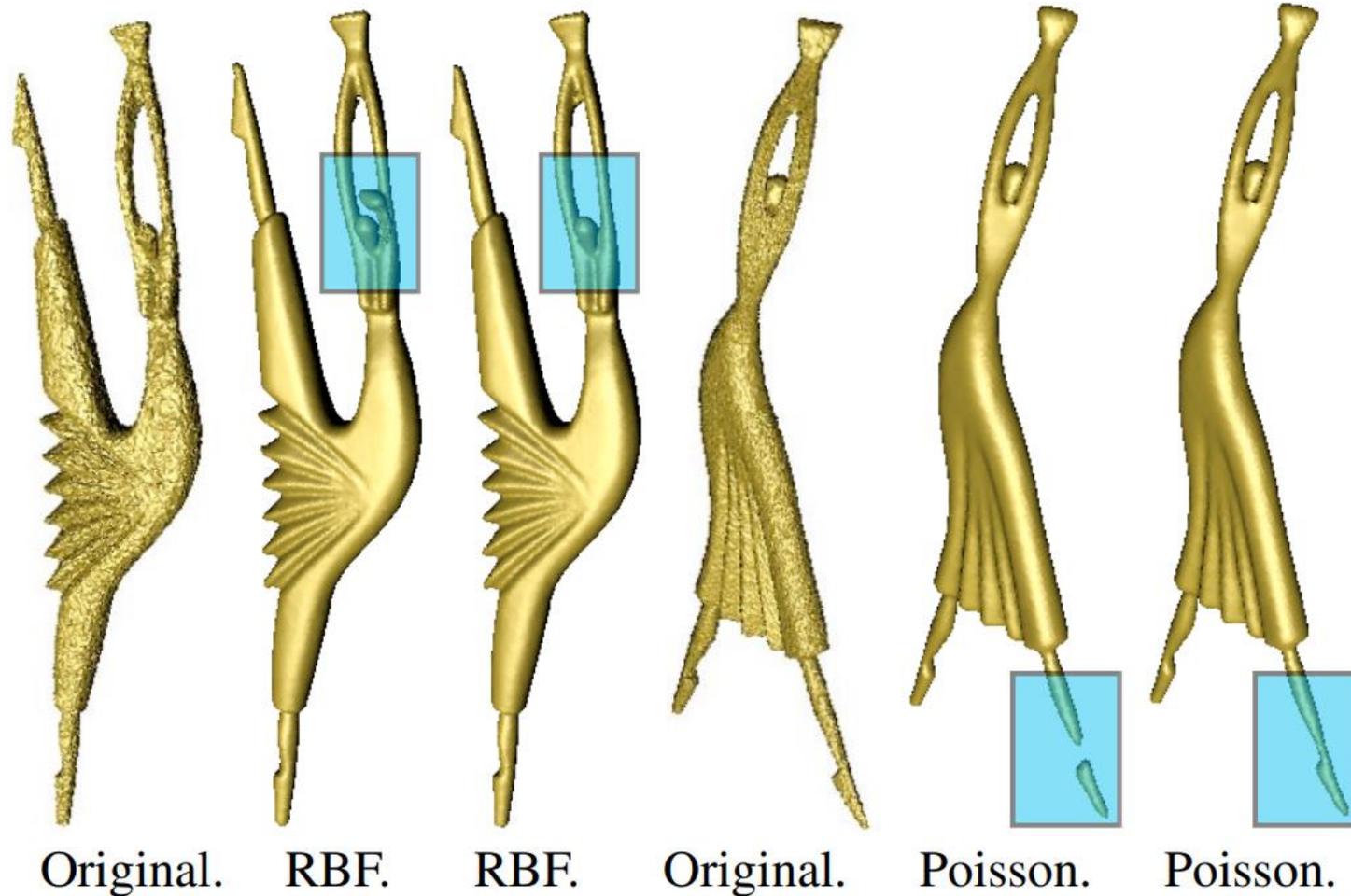
- Replacing the original repulsion term  $\eta(r) = 1/(3*r^3)$  with  $\eta(r) = -r$  which decreases more gently and penalizes more at a large  $r$
- Using weighted local density for each point in a Point Set that enables the attraction of point clusters to be relaxed and the repulsion force from points in dense areas to be strengthened

# Weighted Locally Optimal Projection Operator



(a) Raw scan. (b) LOP (old  $\eta$ ). (c) LOP (new  $\eta$ ). (d) WLOP.

# Weighted Locally Optimal Projection Operator

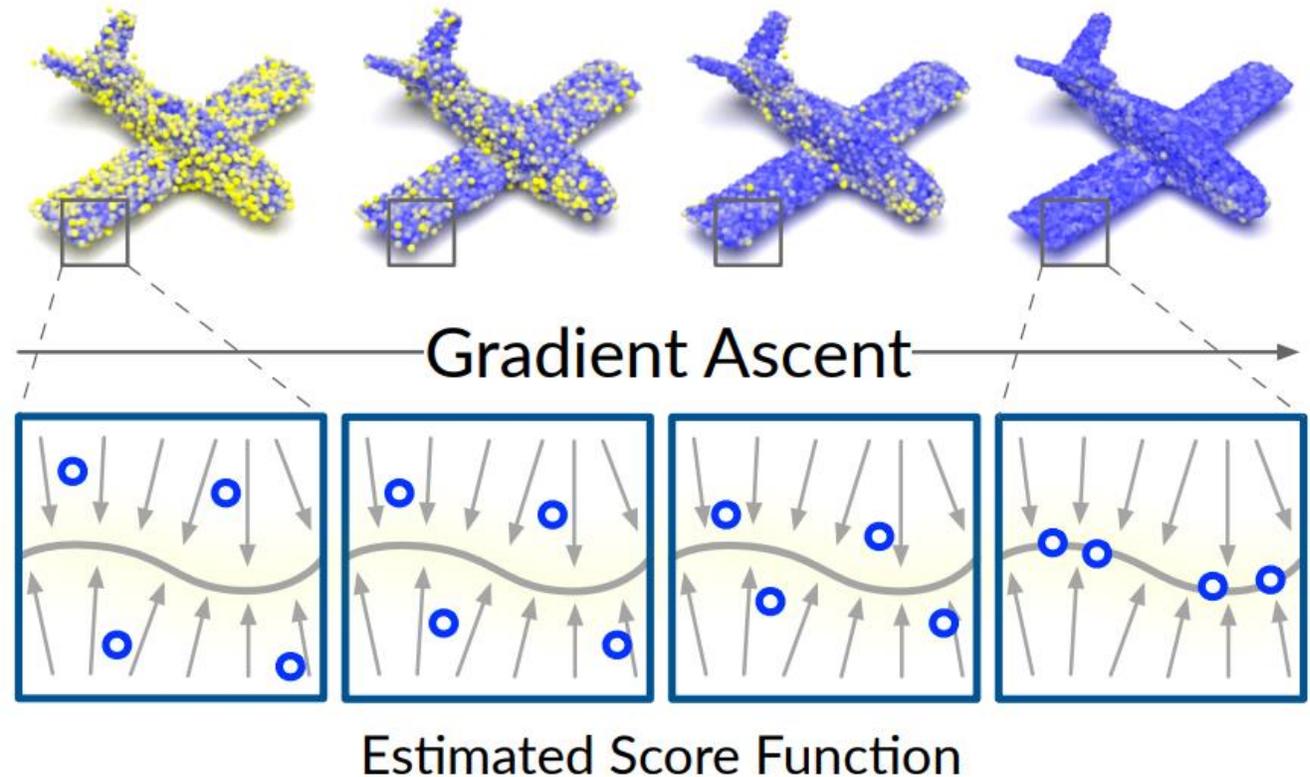


# (Weighted) Locally Optimal Projection Operator

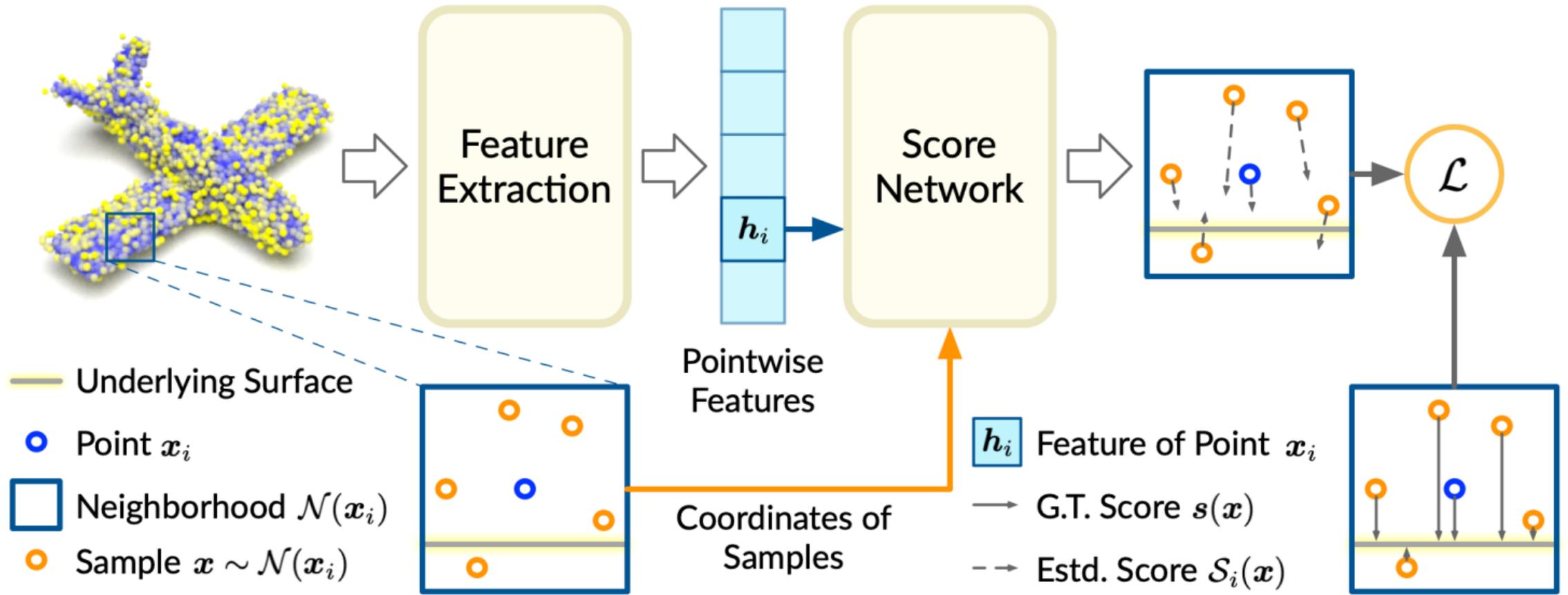
- not learning patterns from data
- mainly relied on parameters set heuristically
- applies Local PCA for estimation of point cloud normals
- cannot recover sharp features and handle missing data

# Score-Based Point Cloud Denoising

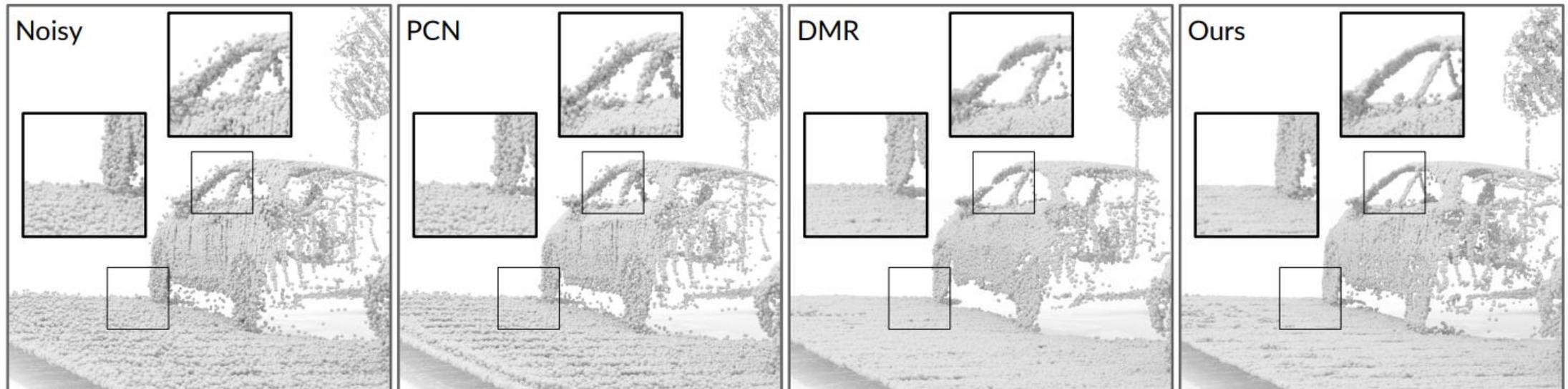
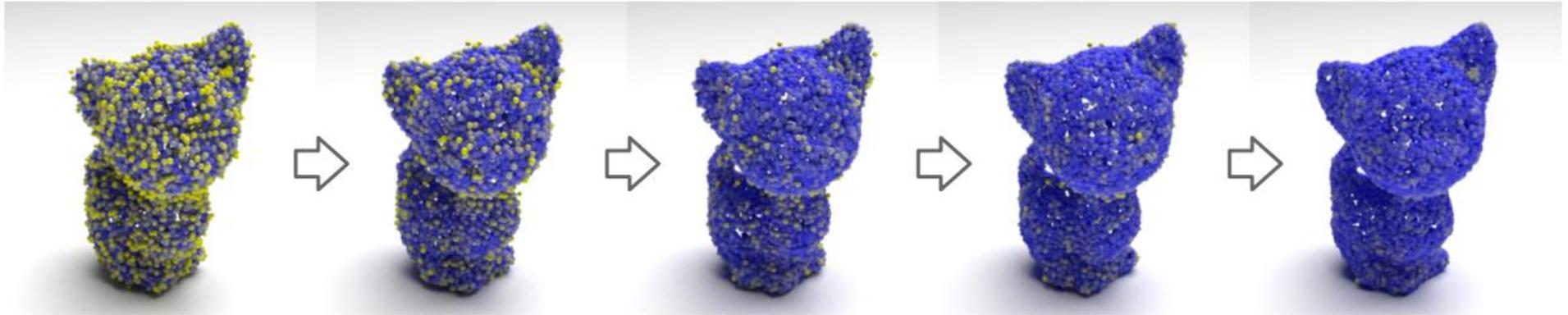
- Learns to estimate the score of the noisy point cloud
- Iteratively updating each point to increase log-likelihood via gradient descent
- Can learn variety of noisy models



# Score-Based Point Cloud Denoising



# Score-Based Point Cloud Denoising



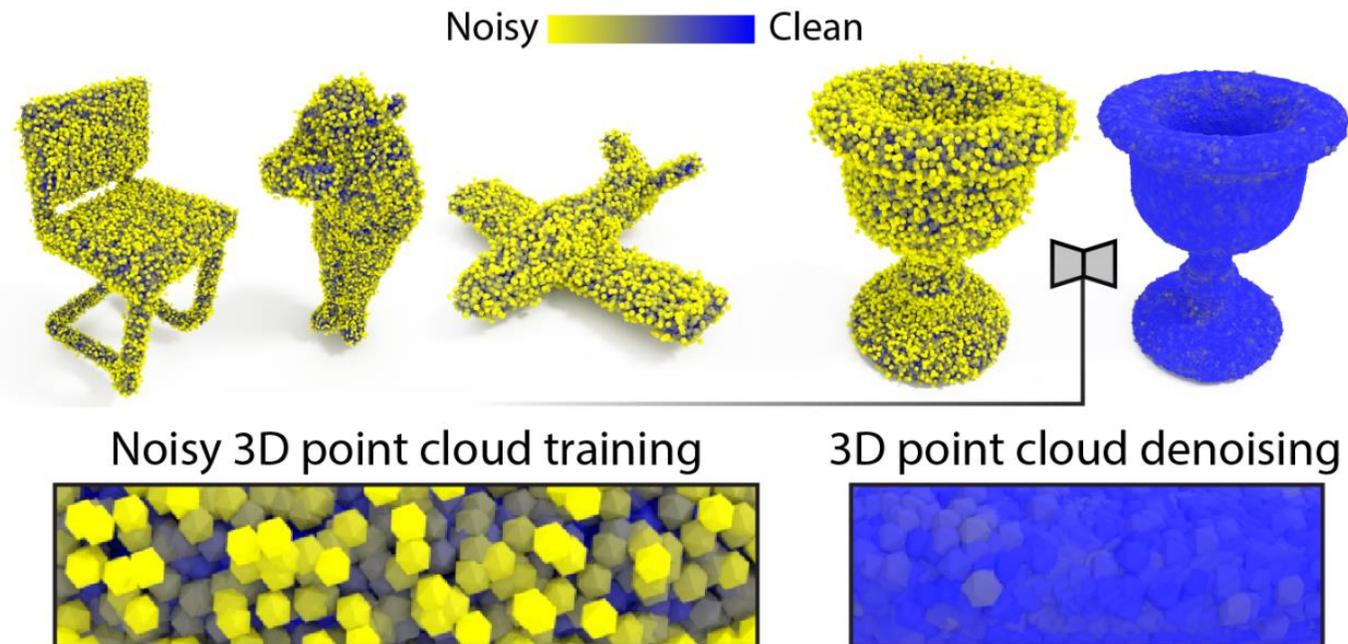
# Score-Based Point Cloud Denoising

Although performing

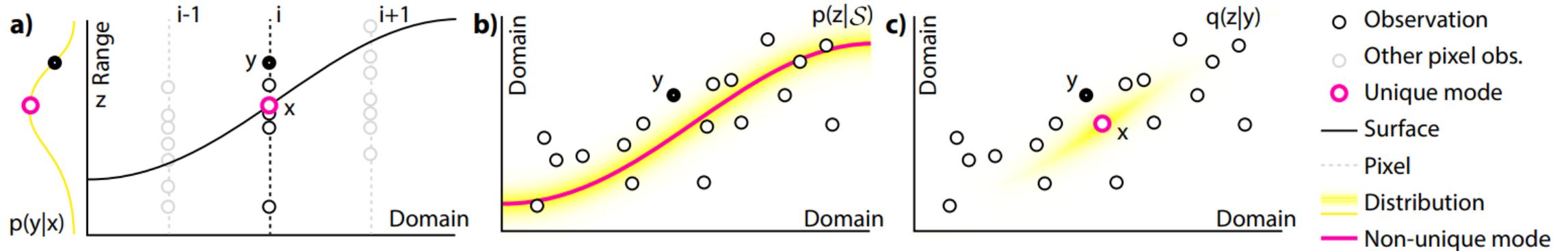
- implements synthetic datasets generation by perturbing ground truth point clouds
- thus not scalable with the real world data
- restricts evaluation to pre-determined set of noise models, including Gaussian noise, simulated LiDAR noise, non-isotropic Gaussian noise, uni-directional noise, Laplace noise, uniform noise, and discrete noise

# Total Denoising: Unsupervised Learning of 3D Point Cloud Cleaning

- Allows learning point cloud cleaning from noisy examples alone
- Enforce convergence to the unique closest out of many possible modes on a manifold of point clouds



# Total Denoising: Unsupervised Learning of 3D Point Cloud Cleaning



To enforce convergence onto a unique mode, we enforce regularization over denoising process by imposing prior distribution  $q(z|y)$  that captures the probability that a given observation  $y$  is a realization of the clean point  $z$

# Total Denoising: Unsupervised Learning of 3D Point Cloud Cleaning

Using spatial and appearance proximity:

$$q(z|y) = p(z|S) * k(z - y)$$

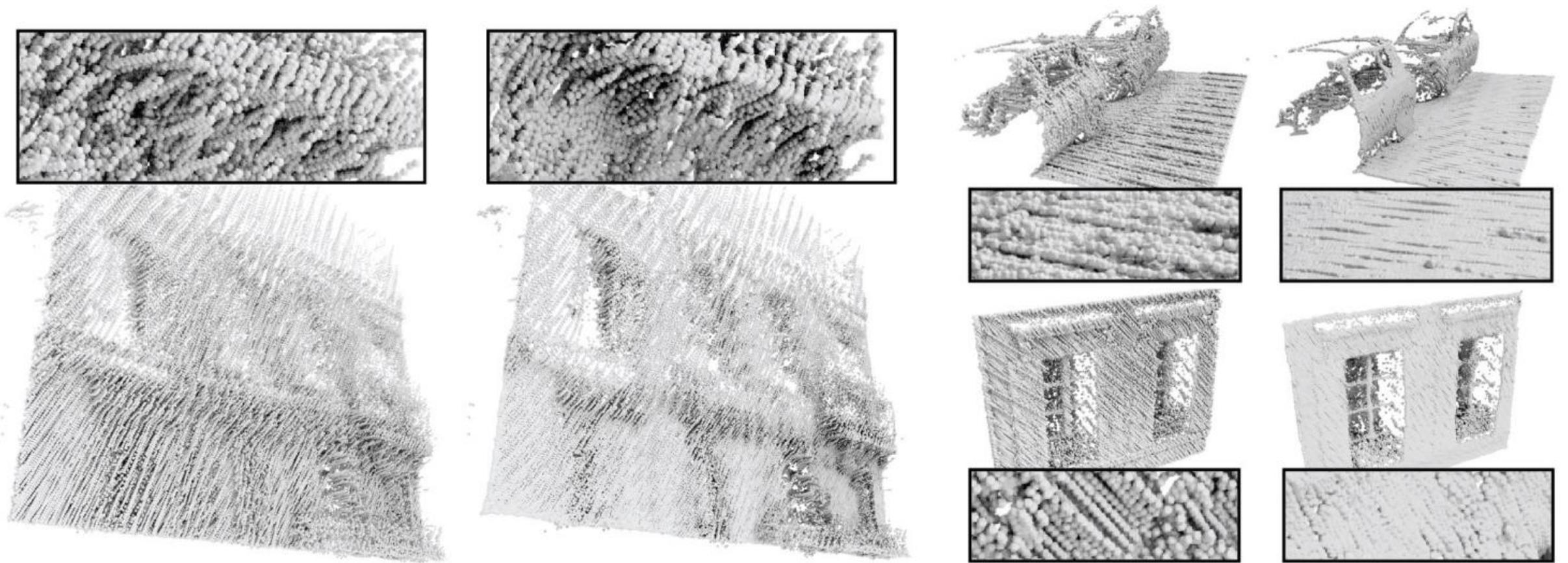
$$k(d) = (1 / \sigma \sqrt{(2\pi)}) * \exp\{- ||Wd||_2^2 / (2*\sigma^2)\}$$

where  $\sigma$  is the bandwidth of  $k$  and  $W = \text{diag}(w)$  is a diagonal weight matrix trading spatial and appearance locality. Optimizing

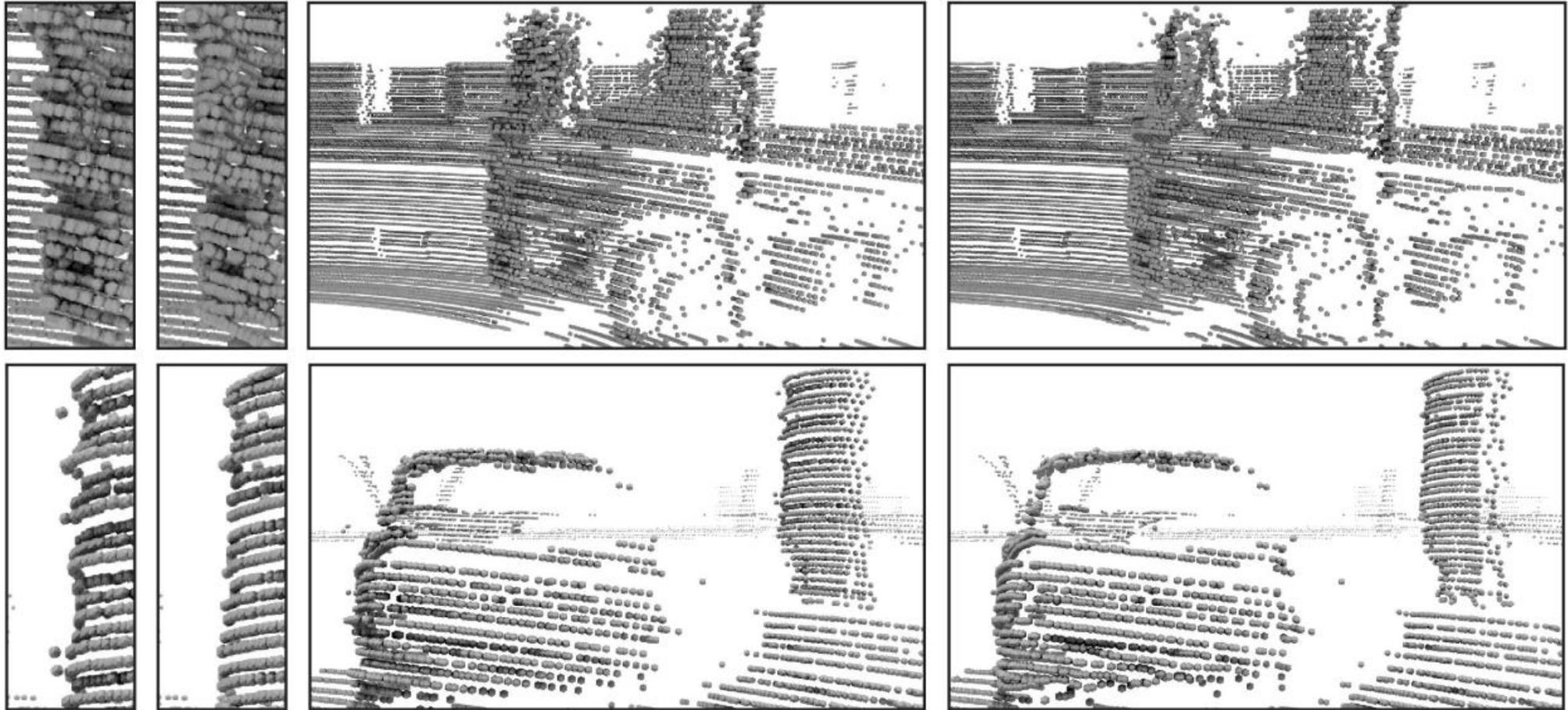
$$\arg \min_{\Theta} \mathbb{E}_{y \sim p(z|S)} \mathbb{E}_{q \sim q(z|y)} L(f_{\Theta}(y), q)$$

where  $L$  can be chosen to be a standard Chamfer distance and  $f$  is implemented using Monte Carlo Convolution

# Total Denoising: Unsupervised Learning of 3D Point Cloud Cleaning



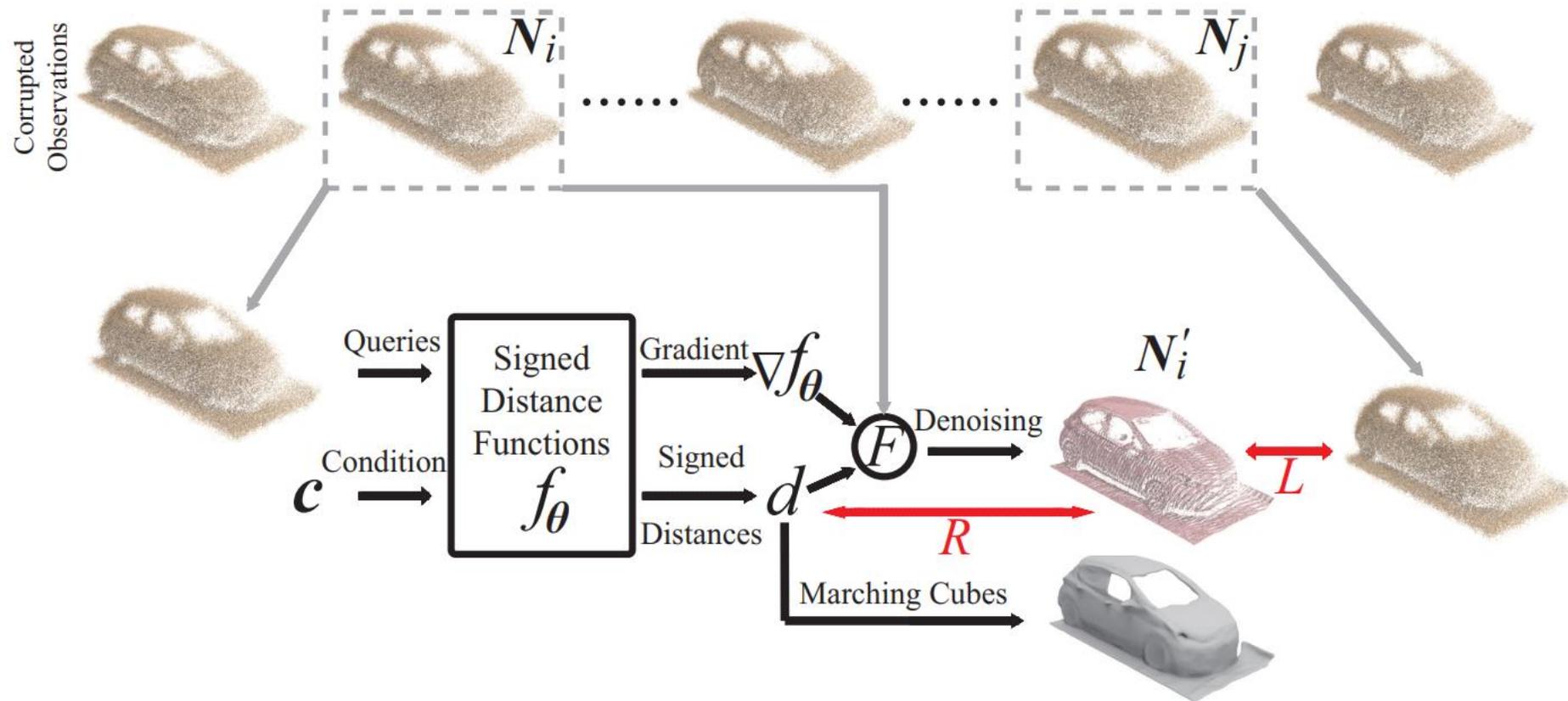
# Total Denoising: Unsupervised Learning of 3D Point Cloud Cleaning



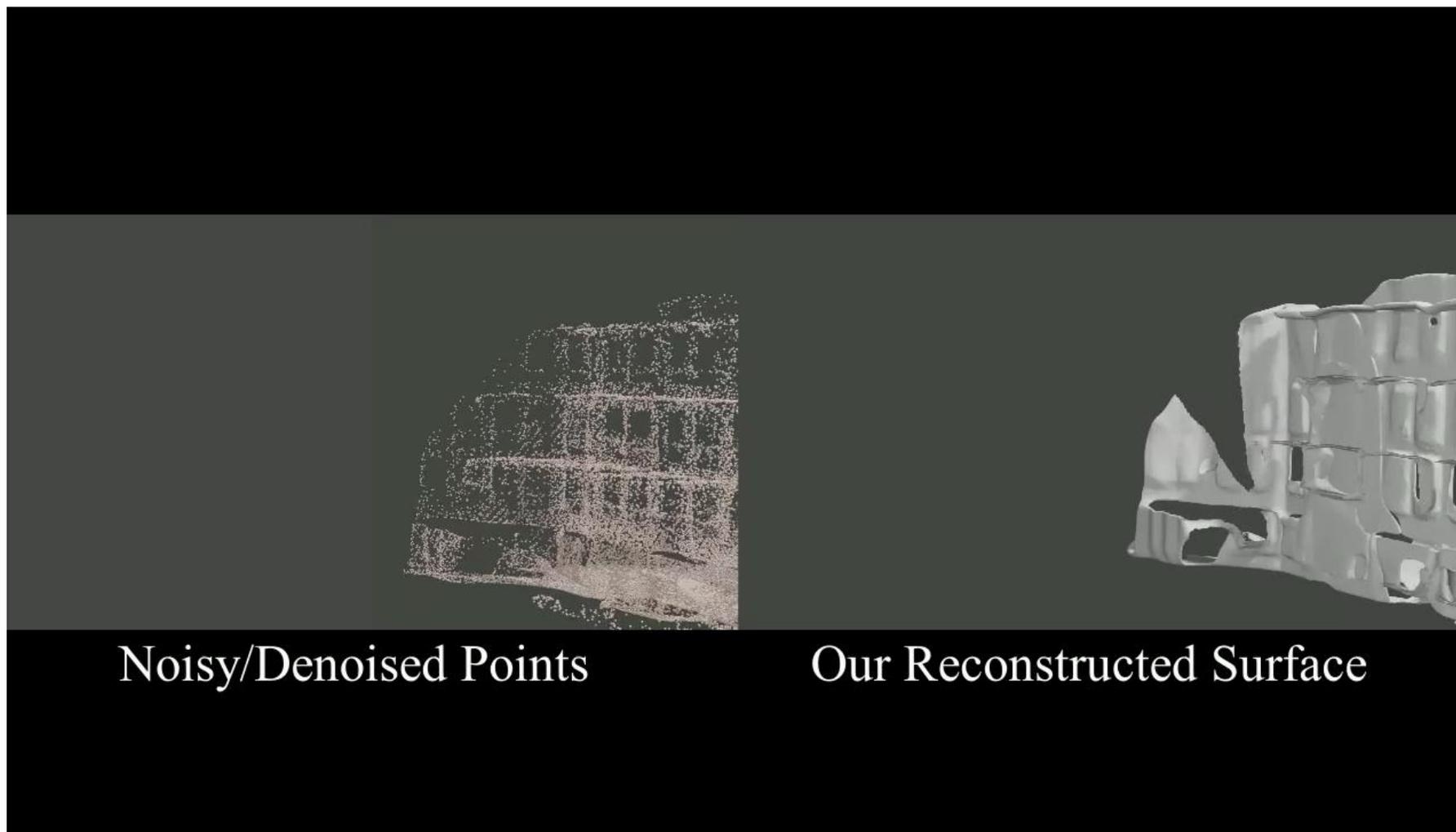
# Fast Learning of Signed Distance Functions from Noisy Point Clouds via Noise to Noise Mapping

- Doesn't require any clean point cloud or ground truth supervision
- Can infer highly accurate distance field from multiple or even a single noisy observation
- Further, introduce a novel schema to improve multi-view reconstruction by estimating SDFs as a prior

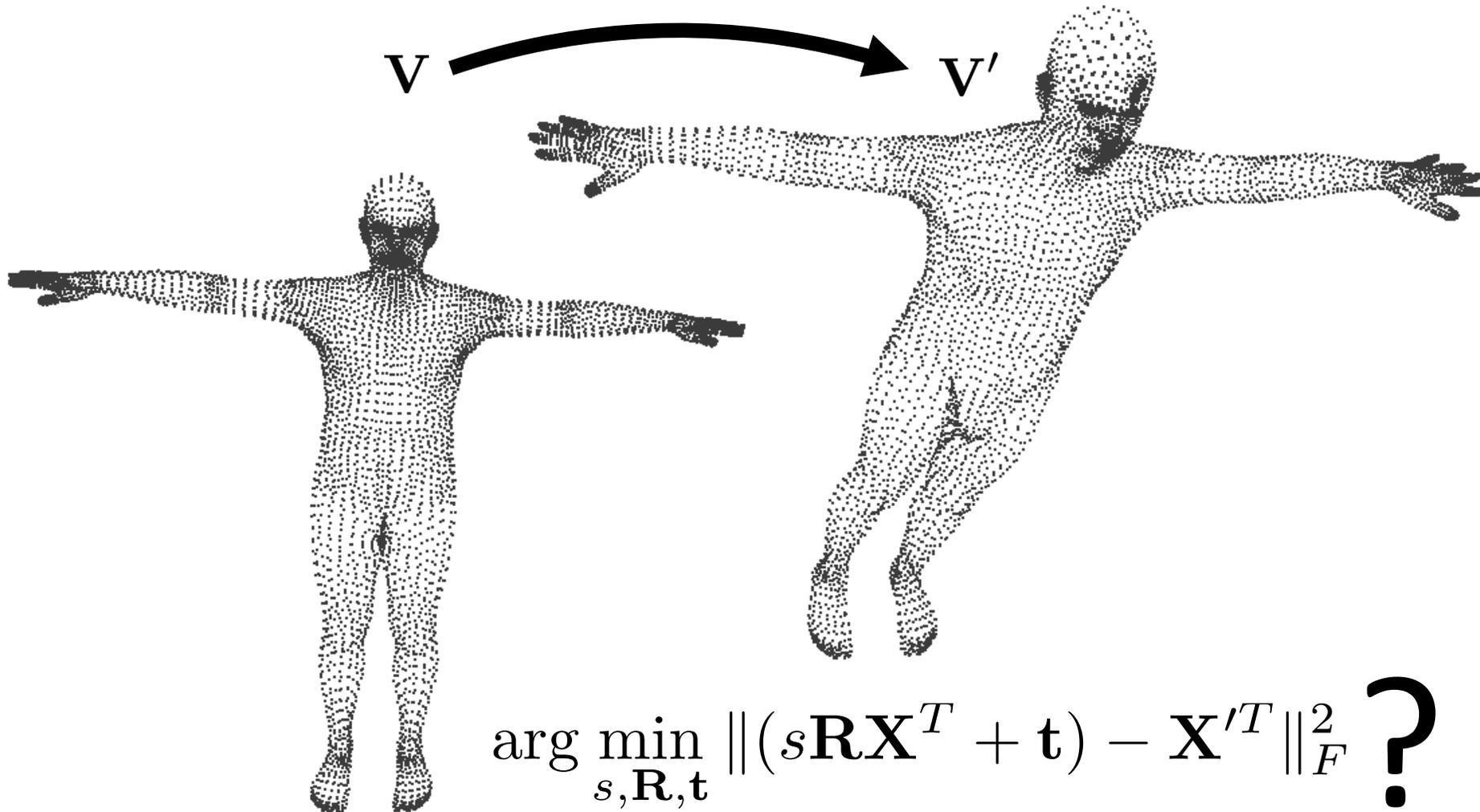
# Fast Learning of Signed Distance Functions from Noisy Point Clouds via Noise to Noise Mapping



# Fast Learning of Signed Distance Functions



# Procrustes alignment problem definition



# Rigid Alignment

How can we rigidly transform a set of points ?

- Translate it
- Rotate it
- Today we'll consider also
- Scale it

$$\mathbf{X}' = \mathbf{X} + \mathbf{t}, \mathbf{t} \in \mathbb{R}^3$$

$$\mathbf{X}'^T = \mathbf{R} \cdot \mathbf{X}^T, \mathbf{R} \in \mathbf{SO}(3)$$

$$\mathbf{X}' = s\mathbf{X}, s \in \mathbb{R}$$

# Optimisation Problem

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

# Quick recap of SVD

in general, applied to a real matrix:

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$$

$$\mathbf{A} \equiv (M \times N) \text{ real}$$

$$\mathbf{U} \equiv (M \times M) \text{ orthogonal, unit norm}$$

$$\mathbf{V} \equiv (N \times N) \text{ orthogonal, unit norm}$$

$$\mathbf{\Sigma} \equiv (M \times N) \text{ diagonal}$$

warning: this is not the vertex matrix!

\* See also clarification slide 1 at the end of slide deck

# Procrustes alignment steps

$$\mathbf{X} = [\mathbf{x}_1 \dots \mathbf{x}_n]^T$$

$$\mathbf{Y} = [\mathbf{y}_1 \dots \mathbf{y}_n]^T$$

Matrices of points

$$\mathbf{X}, \mathbf{Y} \in \mathbb{R}^{(N \times 3)}$$

$$\mathbf{x}_i, \mathbf{y}_i \in \mathbb{R}^{(3 \times 1)}$$

$$\bar{\mathbf{Y}}^T \mathbf{X} = \mathbf{U} \Sigma \mathbf{V}^T$$

$$\mathbf{R} = \mathbf{U} \mathbf{V}^T$$

Optimal **rotation** obtained by computing SVD on the point cross-covariance

$$\mathbf{t} = s \mathbf{R} \bar{\mathbf{x}} - \bar{\mathbf{y}}$$

**Translation** is the centroid difference

$$s = \frac{\text{tr}(\Sigma)}{\|\bar{\mathbf{X}}\|_F^2}$$

**Scale** is a quotient of eigenvalue sums

# Proof

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

# Procrustes derivation: translation

$$\mathbf{X}, \mathbf{Y} \in \mathbb{R}^{(N \times 3)}$$

$$\mathbf{x}_i, \mathbf{y}_i \in \mathbb{R}^{(3 \times 1)}$$

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} E$$

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

Minimize the L2 distance between transformed source points and target points

$$= \sum_i (s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i)^\top (s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i)$$

$$= \sum_i s^2 \mathbf{x}_i^\top \mathbf{x}_i + \mathbf{t}^\top \mathbf{t} + \mathbf{y}_i^\top \mathbf{y}_i + 2s\mathbf{x}_i^\top \mathbf{R}^\top \mathbf{t} - 2s\mathbf{x}_i^\top \mathbf{R}^\top \mathbf{y}_i - 2\mathbf{t}^\top \mathbf{y}_i$$

# Procrustes derivation: translation

We remove the elements that do not depend on the translation and solve for  $\mathbf{t}$ .

$$\mathbf{t} = \arg \min_{\mathbf{t}} E = \arg \min_{\mathbf{t}} \sum_i s^2 \mathbf{x}_i^T \mathbf{x}_i + \mathbf{t}^T \mathbf{t} + \mathbf{y}_i^T \mathbf{y}_i + 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{t} - 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{y}_i - 2\mathbf{t}^T \mathbf{y}_i$$

$$= \arg \min_{\mathbf{t}} \sum_i \mathbf{t}^T \mathbf{t} + 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{t} - 2\mathbf{t}^T \mathbf{y}_i$$

$$\bar{\mathbf{x}} \equiv \frac{\sum_i \mathbf{x}_i}{N}, \bar{\mathbf{y}} \equiv \frac{\sum_i \mathbf{y}_i}{N}$$

Compute the centroid of the point clouds

$$\mathbf{t} = \arg \min_{\mathbf{t}} E = \arg \min_{\mathbf{t}} (\mathbf{t}^T (2s \mathbf{R} \bar{\mathbf{x}} + \mathbf{t} - 2\bar{\mathbf{y}})) = \bar{\mathbf{y}} - s \mathbf{R} \bar{\mathbf{x}}$$

So given  $s$  and  $R$ , we can compute the translation  $\mathbf{t}$

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

subject to  $\mathbf{R} \in SO(3)$

Optimal rotation does not depend on scale

$$\mathbf{R} = \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2$$

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

$$\mathbf{R} = \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2$$

$$= \arg \min_{\mathbf{R}} \langle \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T, \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \min_{\mathbf{R}} \|\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \max_{\mathbf{R}} \langle \mathbf{R}, \bar{\mathbf{Y}}^T \bar{\mathbf{X}} \rangle_F$$

Optimal rotation does not depend on scale

Rotation does not change norm!

Inner product should be maximum!

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

$$\begin{aligned} \mathbf{R} &= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2 \\ &= \arg \min_{\mathbf{R}} \langle \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T, \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \min_{\mathbf{R}} \|\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle \mathbf{R}, \bar{\mathbf{Y}}^T \bar{\mathbf{X}} \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle \mathbf{R}, U\Sigma V^T \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle U^T \mathbf{R} V, \Sigma \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V \end{aligned}$$

Optimal rotation does not depend on scale

Rotation does not change norm!

Inner product should be maximum!

Cross-covariance matrix

SVD decomposition

# Procrustes derivation: Rotation

$$= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V$$

What kind of matrix is  $S$  ?

What kind of matrix is  $\Sigma$  ?

# Procrustes derivation: Rotation

$$= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V$$

What kind of matrix is  $S$  ?

Orthogonal

What kind of matrix is  $\Sigma$  ?

Diagonal

Hence the quantity above is maximised when  $S$  equals the identity, hence:

$$I = U^T \mathbf{R} V$$

$$\mathbf{R} = UV^T$$

SVD of cross-covariance of pointsclds

$$\bar{\mathbf{Y}}^T \mathbf{X} = U \Sigma V^T$$

# Procrustes derivation: scale

Optimize **scale** given the rotation

$$s = \arg \min_s E = \arg \min_s \left( \sum_i s^2 \bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i + \bar{\mathbf{y}}_i^T \bar{\mathbf{y}}_i - 2s \bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i \right)$$

$$= \arg \min_s \left( s^2 \sum_i (\bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i) - 2s \sum_i (\bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i) \right)$$

$$= \arg \min_s (s^2 a - 2sb) = \frac{b}{a} = \frac{\sum_i (\bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i)}{\sum_i (\bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i)}$$

$$= \frac{\text{tr}(\bar{\mathbf{X}} \mathbf{R}^T \bar{\mathbf{Y}}^T)}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\mathbf{R}^T \bar{\mathbf{Y}}^T \bar{\mathbf{X}})}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\mathbf{V} \mathbf{U}^T \mathbf{U} \Sigma \mathbf{V}^T)}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\Sigma)}{\|\bar{\mathbf{X}}\|_F^2}$$

We used:

- 1) Trace invariance with shifts
- 2)  $\bar{\mathbf{Y}}^T \bar{\mathbf{X}} = \mathbf{U} \Sigma \mathbf{V}^T$ ,  $\mathbf{R} = \mathbf{U} \mathbf{V}^T$
- 3) Trace equals sum of eigenvals for square matrices

# Slide credits and further reading

# Hands-On AI Based 3D Vision

## Winter 24/25

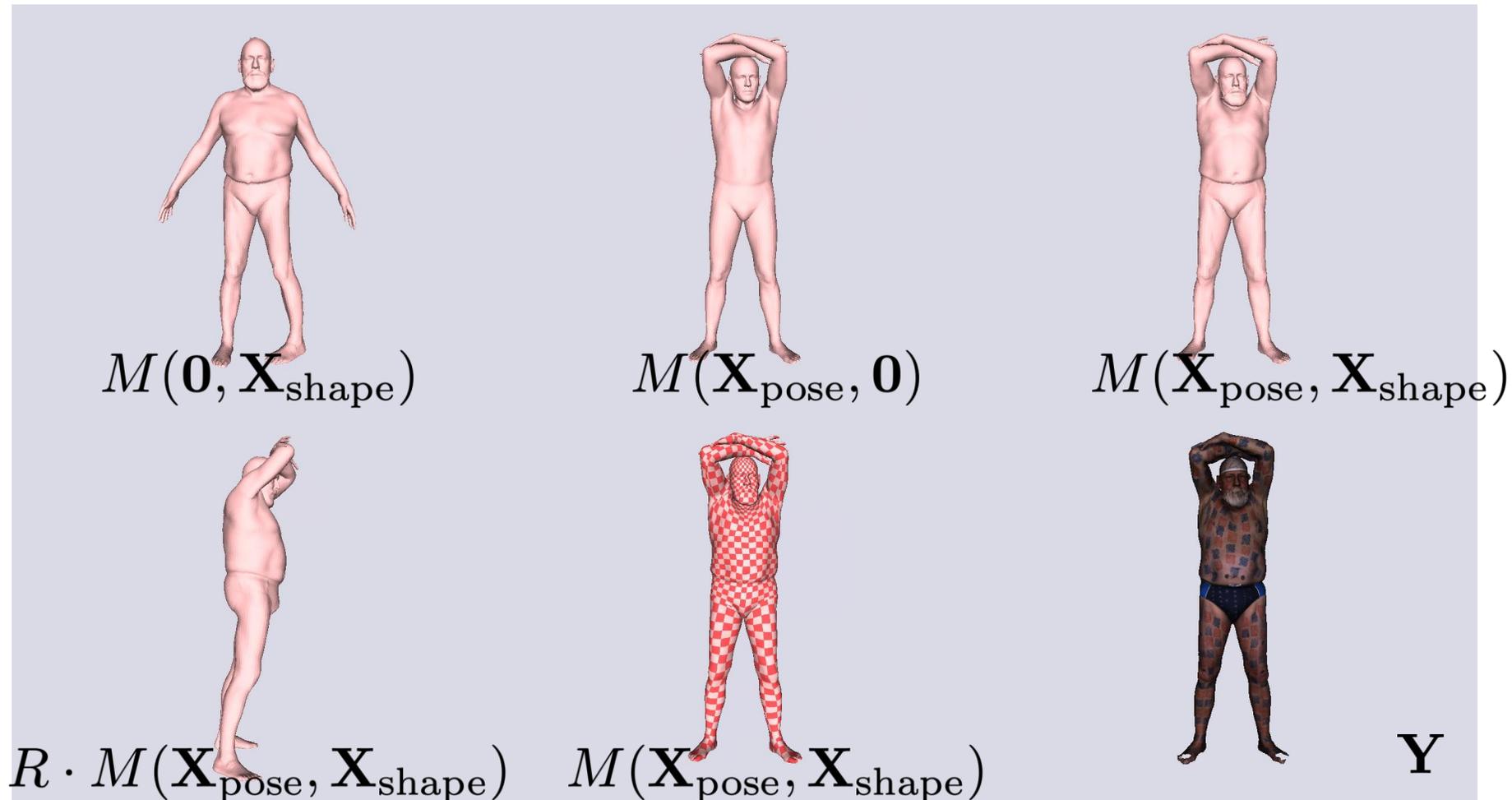
Lecture 5\_2 – Procrustes Alignment

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EBERHARD KARLS  
UNIVERSITÄT  
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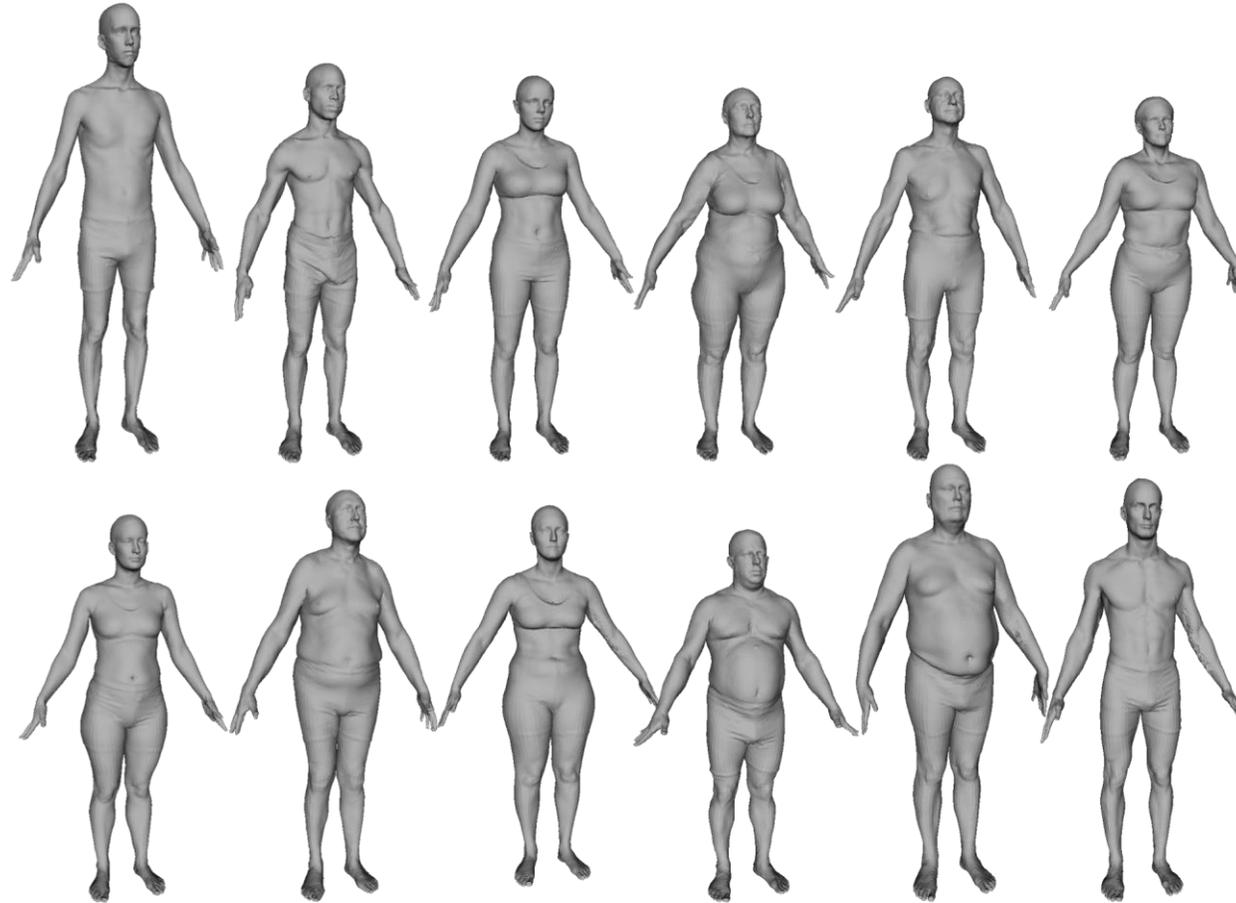


# Goal: learn a model of pose and shape



$$\mathbf{X} = \{\mathbf{X}_{\text{pose}}, \mathbf{X}_{\text{shape}}\}$$

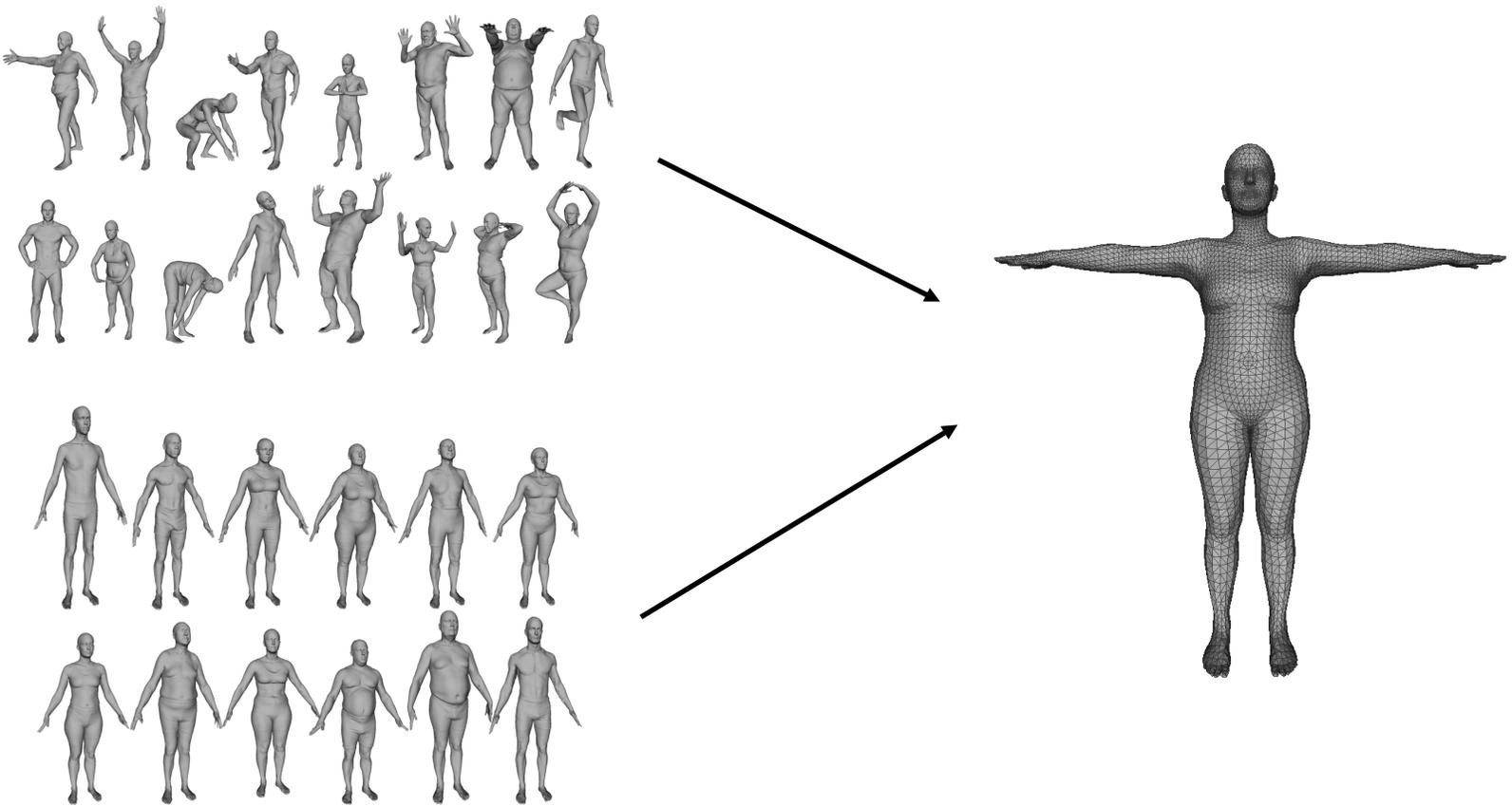
# Scan a Lot of People



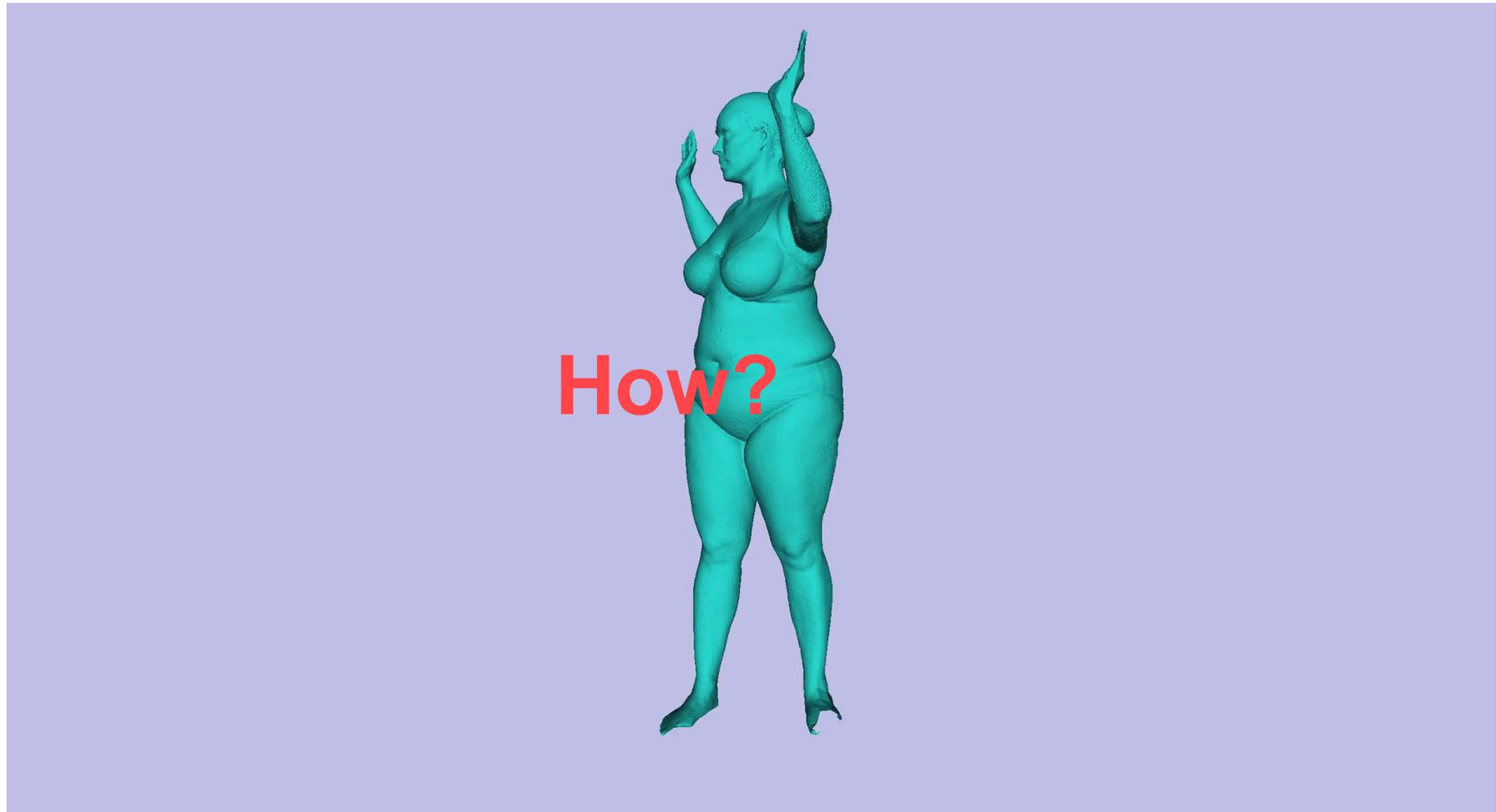
# In lots of poses



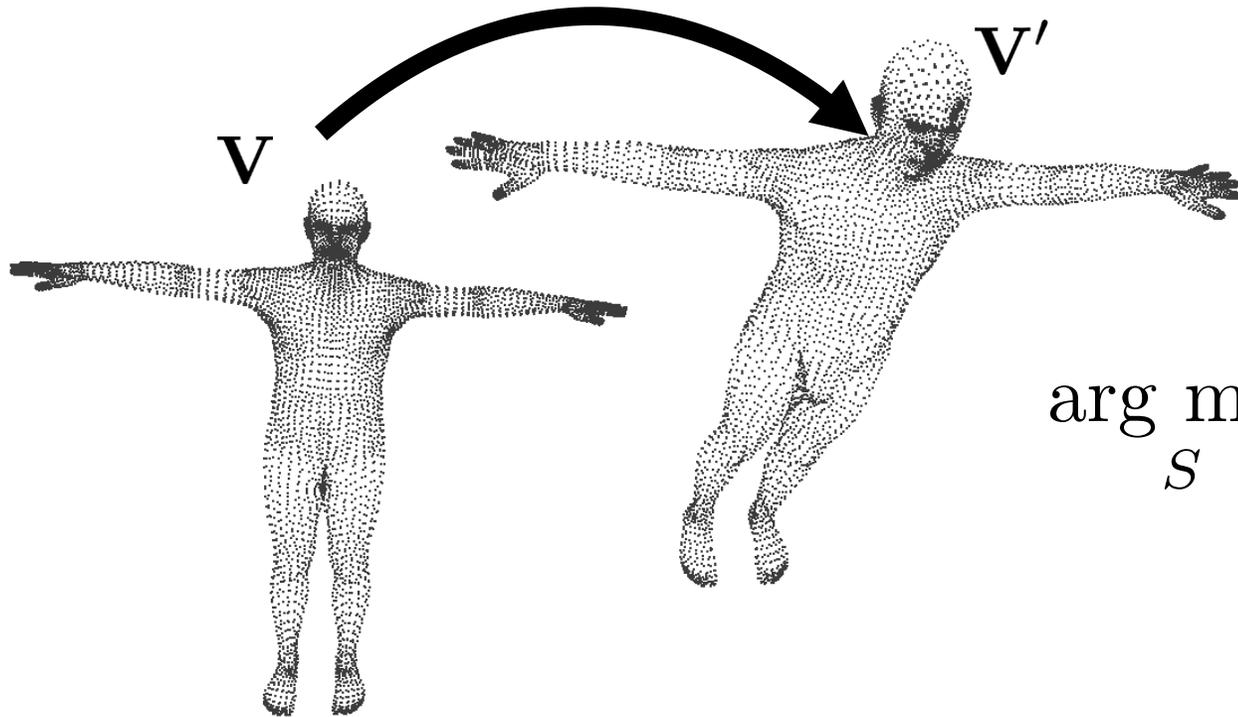
# To learn a model we need correspondence



# Non-rigid Articulated Registration



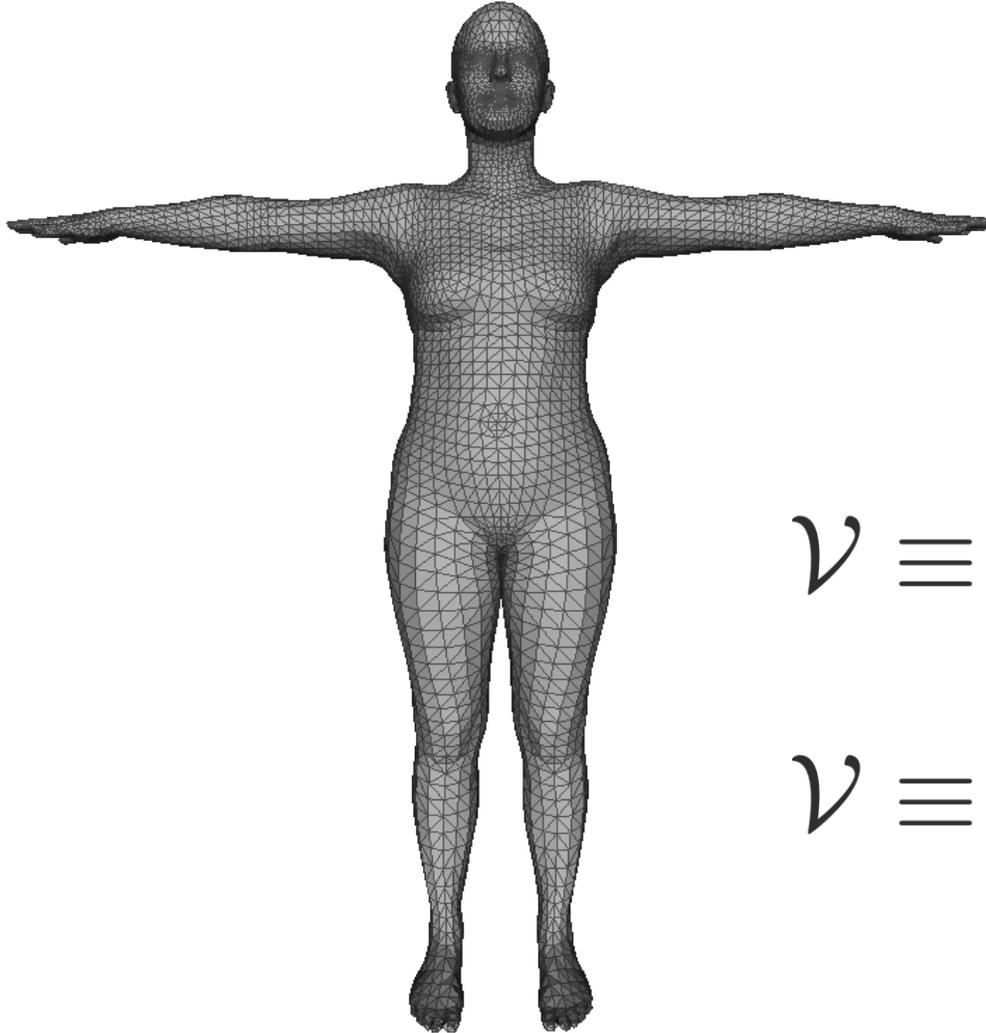
# Today: Rigid Alignment (Procrustes)



$$\arg \min_S \| (s\mathbf{R} \cdot \mathbf{X}^T + \mathbf{t}) - \mathbf{X}'^T \|$$

?

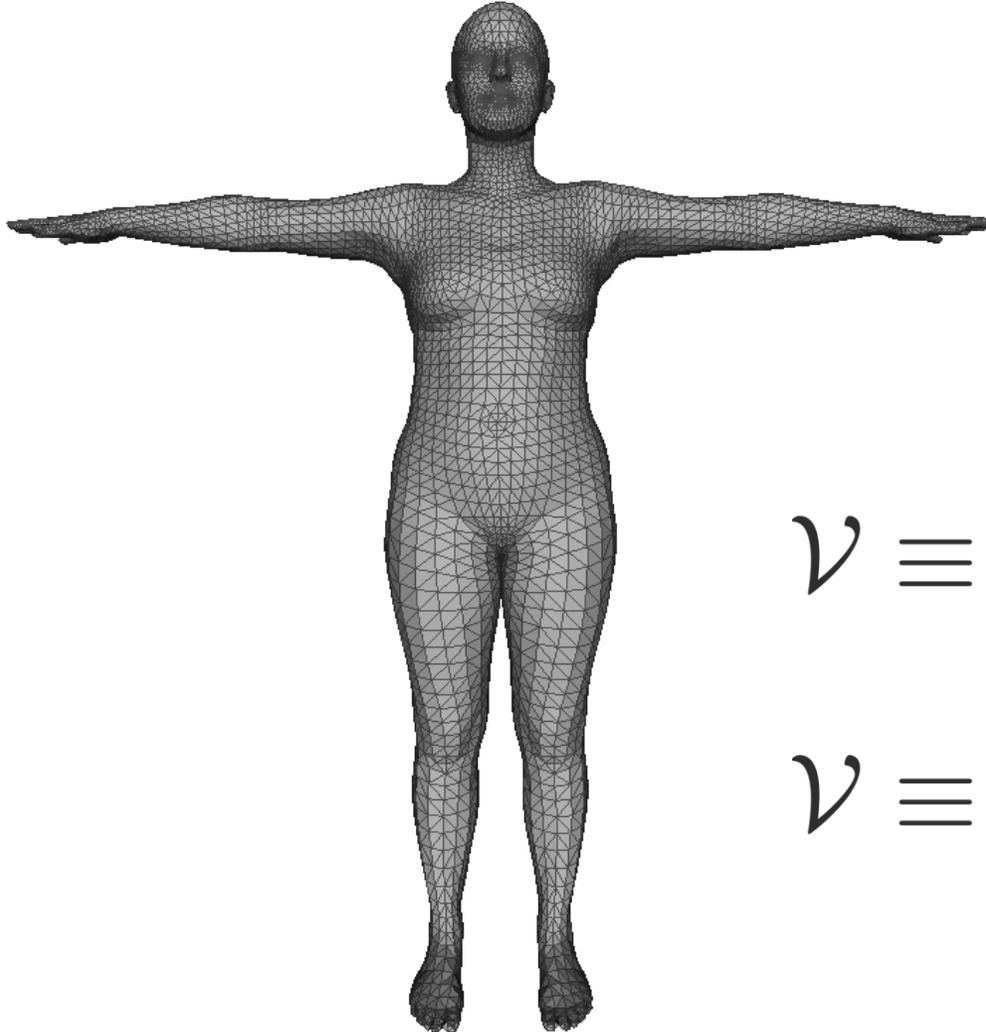
# Surface representation: Mesh



$$\mathcal{V} \equiv \begin{cases} \mathbf{F} \in \mathbb{N}^{M \times 3} \\ \mathbf{X} \in \mathbb{R}^{N \times 3} \end{cases}$$

$$\mathcal{V} \equiv \mathbb{R}^2 \rightarrow \mathbb{R}^3$$

# Surface representation: Mesh



$$\mathcal{V} \equiv \begin{cases} \mathbf{F} \in \mathbb{N}^{M \times 3} \\ \mathbf{X} \in \mathbb{R}^{N \times 3} \end{cases}$$

$$\mathcal{V} \equiv \mathbb{R}^2 \rightarrow \mathbb{R}^3$$

# Surface representation: Mesh

How can we rigidly transform a set of points ?

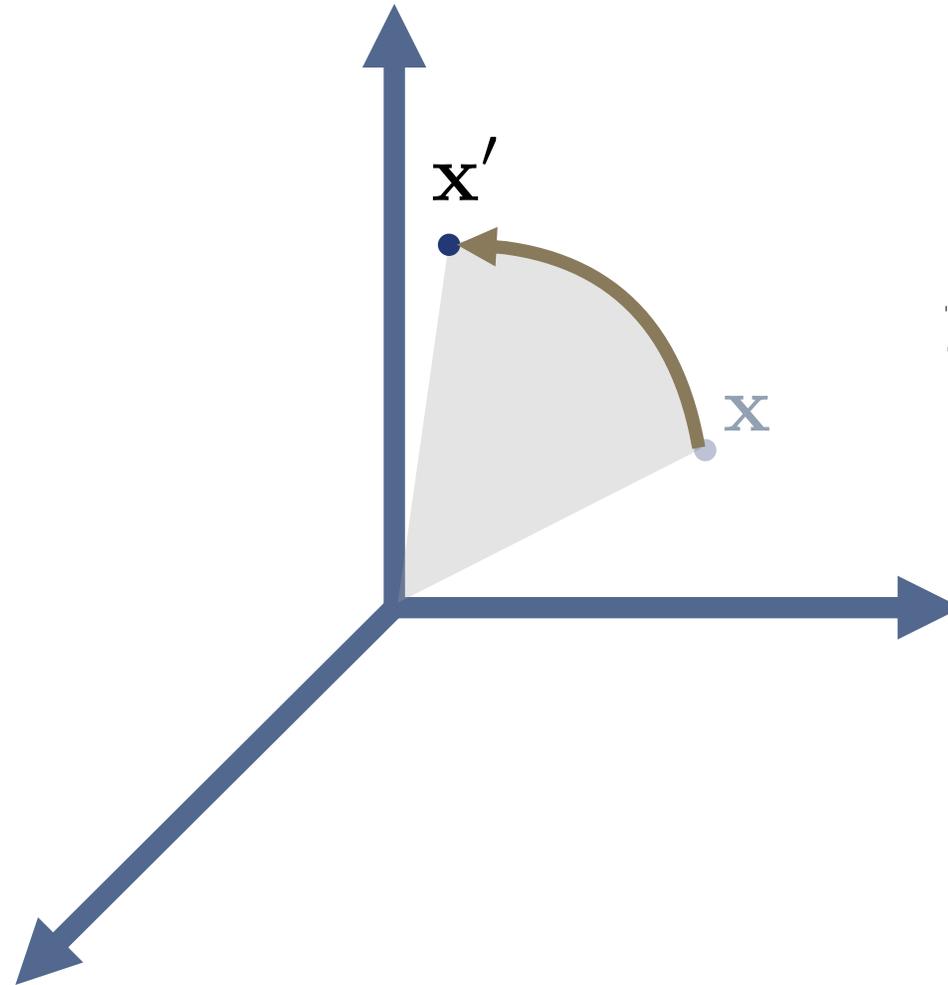
- Translate it
- Rotate it
- Today we'll consider also
- Scale it

$$\mathbf{X}' = \mathbf{X} + \mathbf{t}, \mathbf{t} \in \mathbb{R}^3$$

$$\mathbf{X}'^T = \mathbf{R} \cdot \mathbf{X}^T, \mathbf{R} \in \mathbf{SO}(3)$$

$$\mathbf{X}' = s\mathbf{X}, s \in \mathbb{R}$$

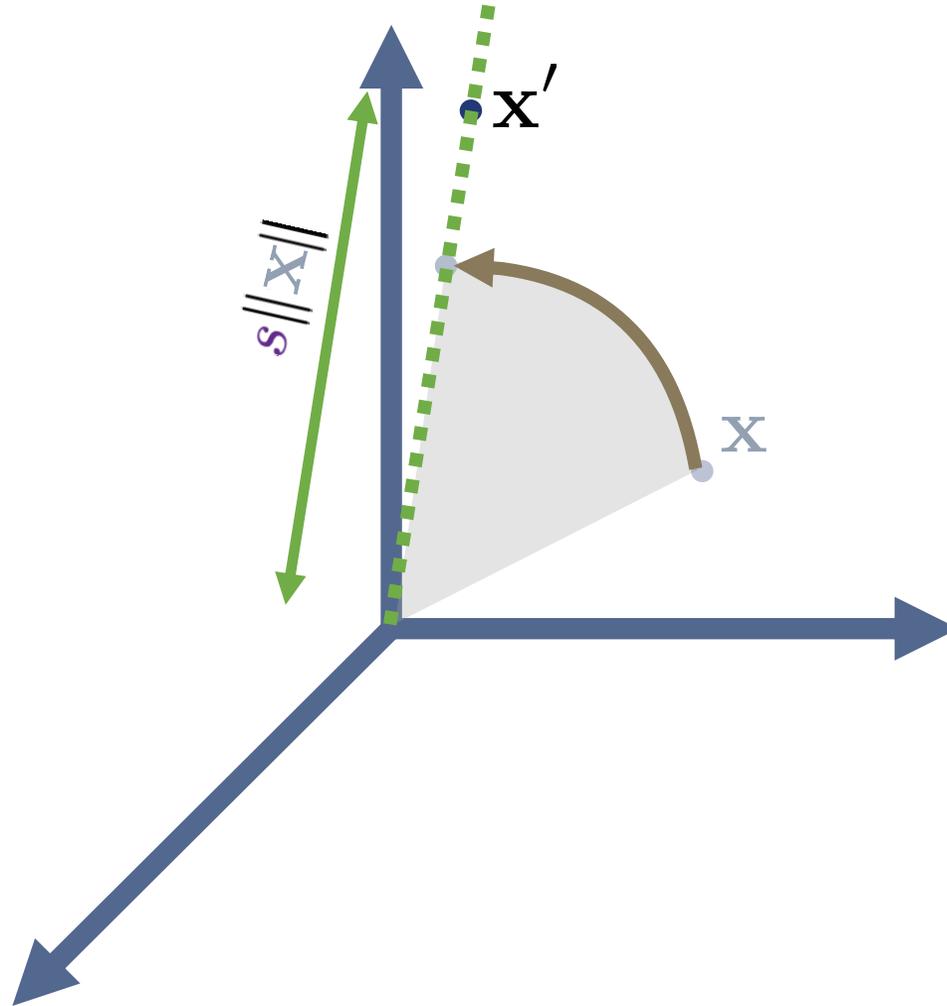
# Rigid transformation + scale



$$R : \mathbb{R}^3 \rightarrow \mathbb{R}^3$$

$$\mathbf{x}' = R\mathbf{x}$$

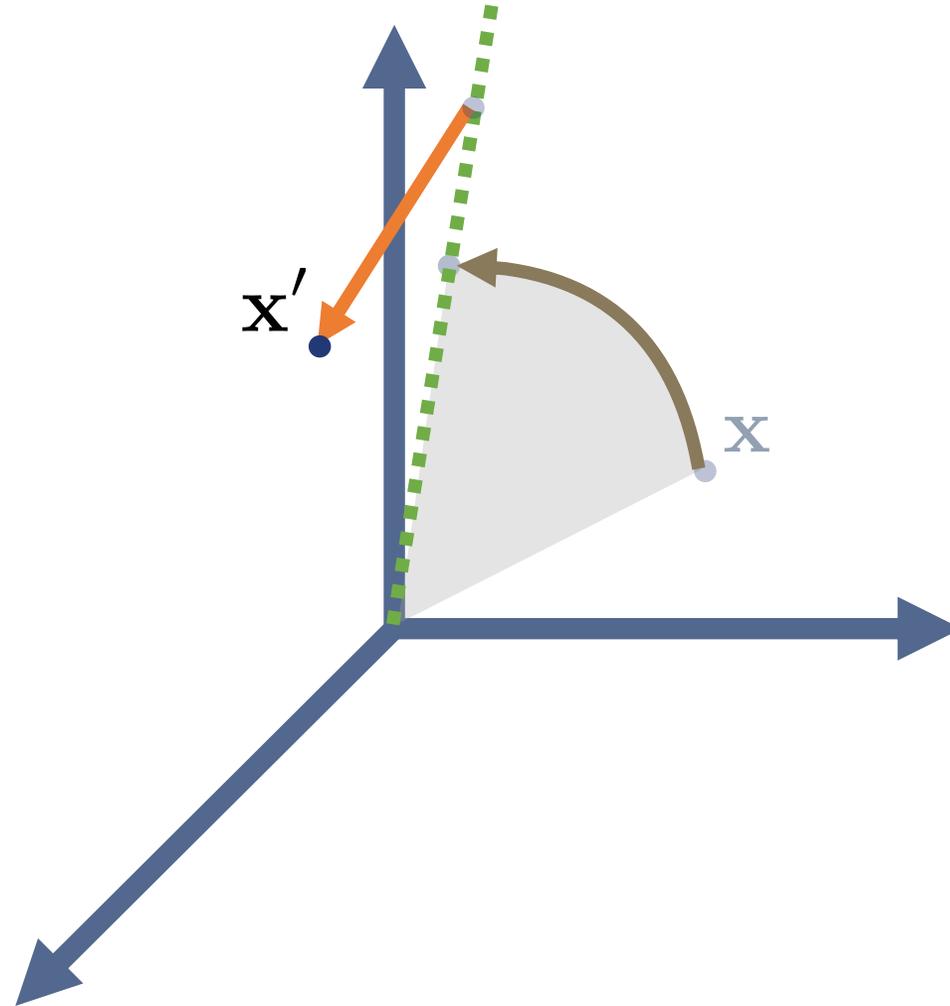
# Rigid transformation + scale



$$R : \mathbb{R}^3 \rightarrow \mathbb{R}^3$$

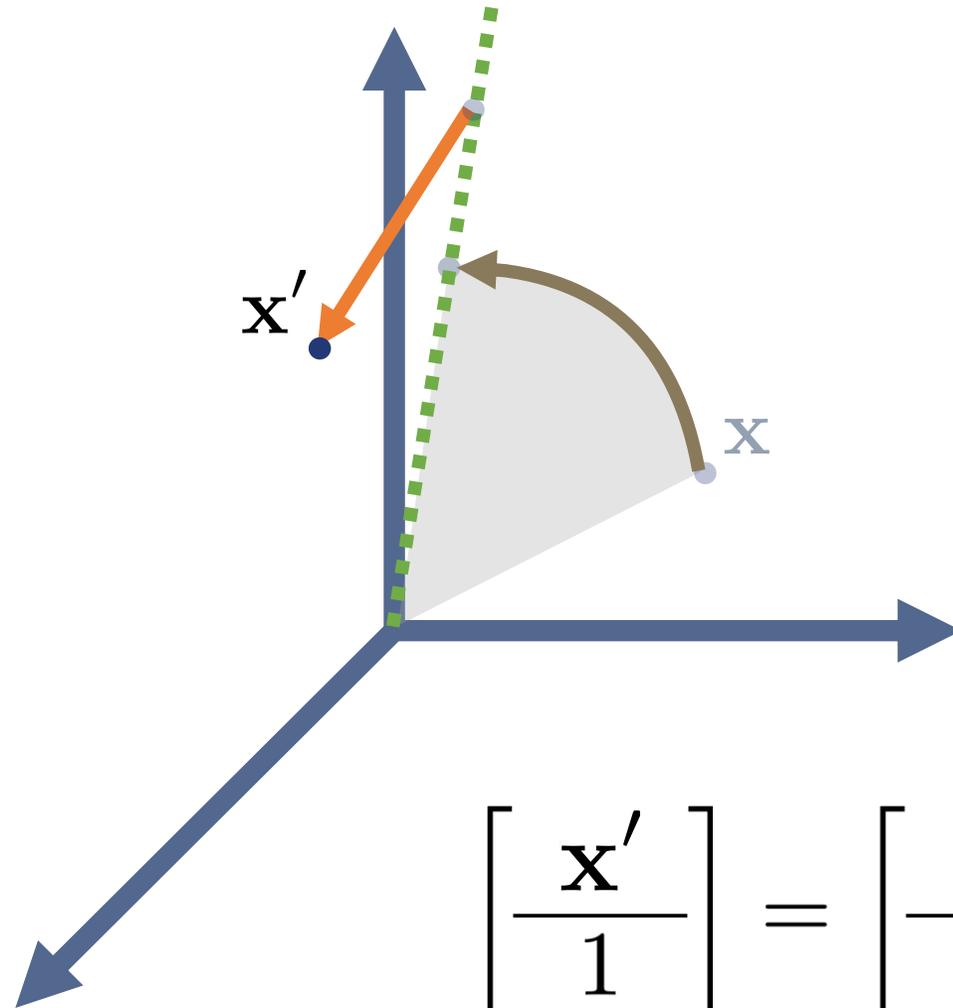
$$x' = sRx$$

# Rigid transformation + scale



$$R : \mathbb{R}^3 \rightarrow \mathbb{R}^3$$
$$\mathbf{x}' = sR\mathbf{x} + \mathbf{t}$$

# Rigid transformation + scale



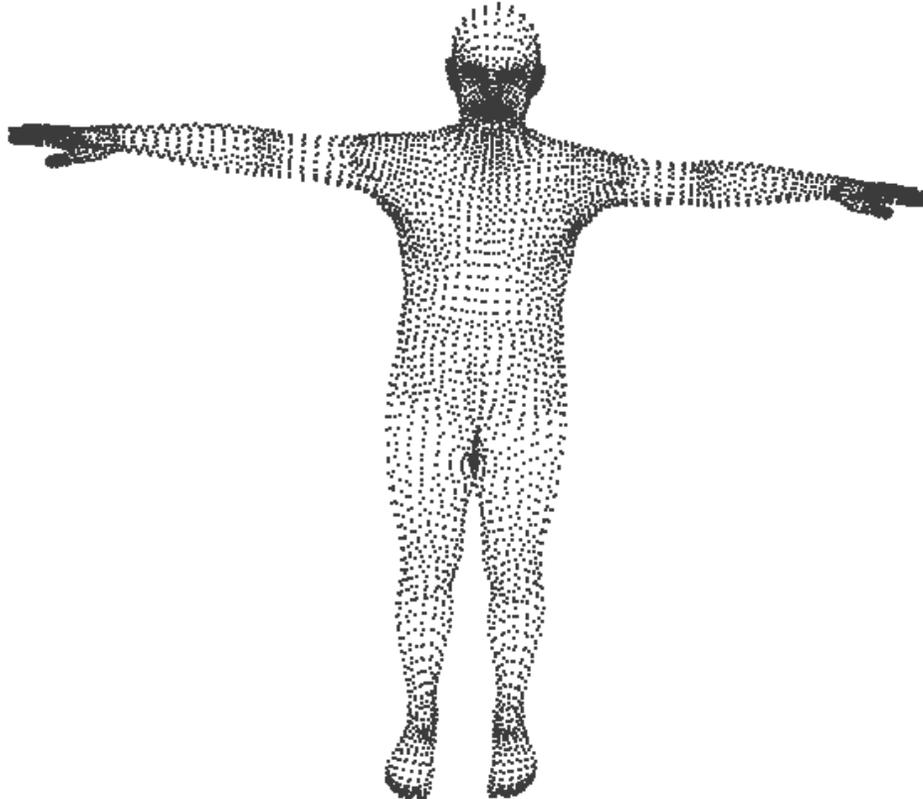
$$R : \mathbb{R}^3 \rightarrow \mathbb{R}^3$$

$$\mathbf{x}' = s\mathbf{R}\mathbf{x} + \mathbf{t}$$

$$\begin{bmatrix} \mathbf{x}' \\ 1 \end{bmatrix} = \begin{bmatrix} s\mathbf{R} & \mathbf{t} \\ 0 & 1 \end{bmatrix} \begin{bmatrix} \mathbf{x} \\ 1 \end{bmatrix}$$

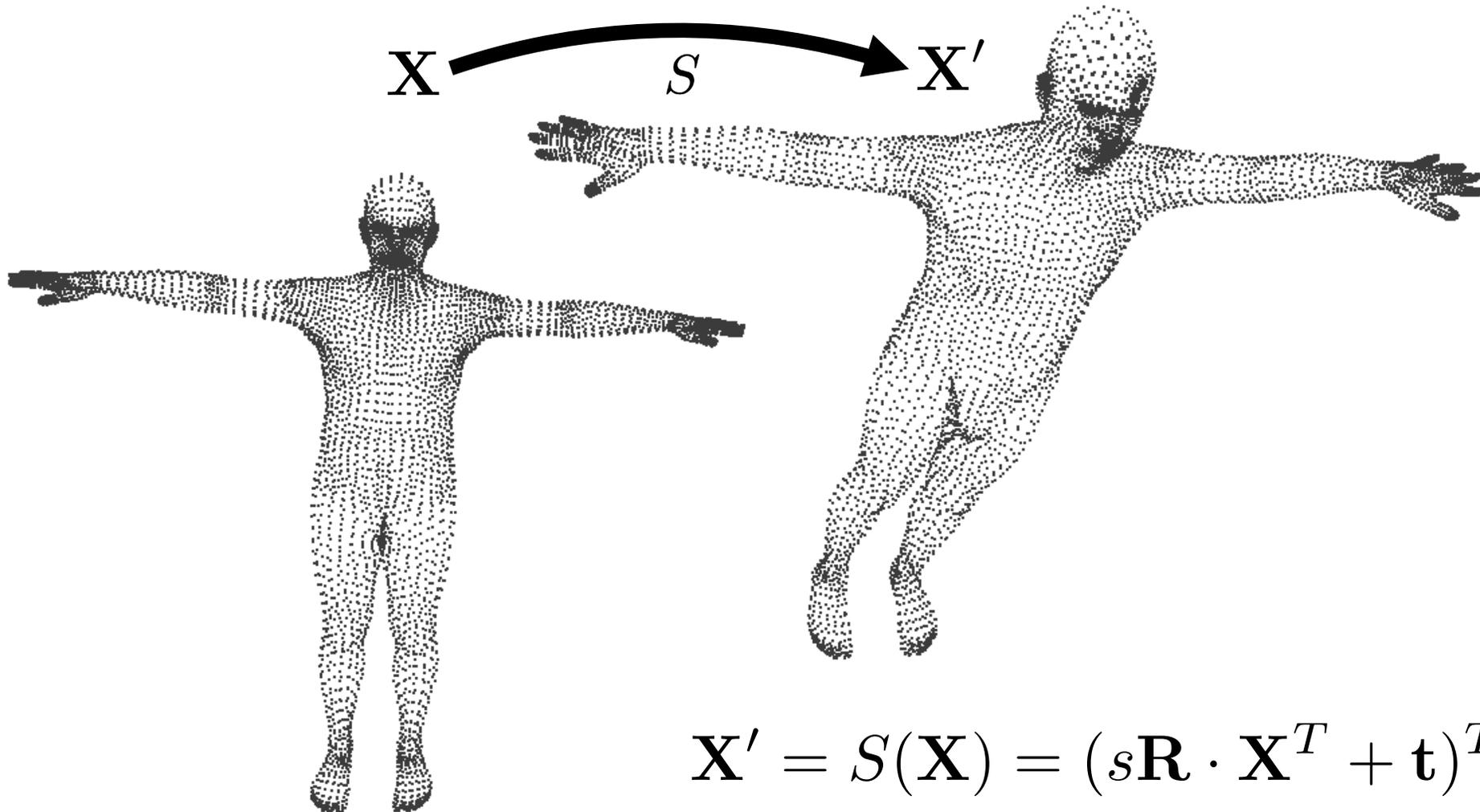
# Procrustes alignment problem definition

$$X \xrightarrow{S} X'$$



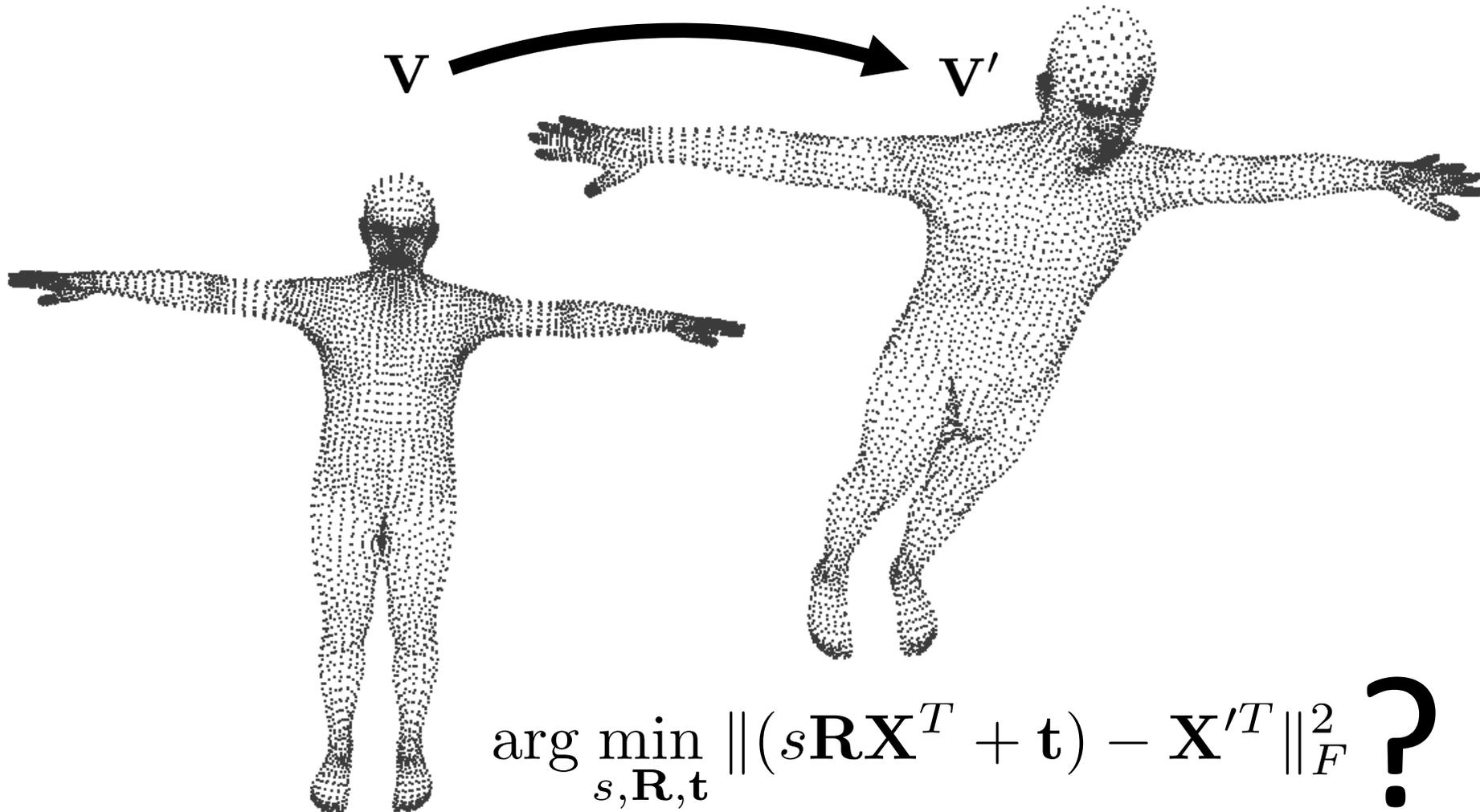
?

# Procrustes alignment problem definition



$$\mathbf{X}' = S(\mathbf{X}) = (s\mathbf{R} \cdot \mathbf{X}^T + \mathbf{t})^T$$

# Procrustes alignment problem definition



# Why Procrustes?



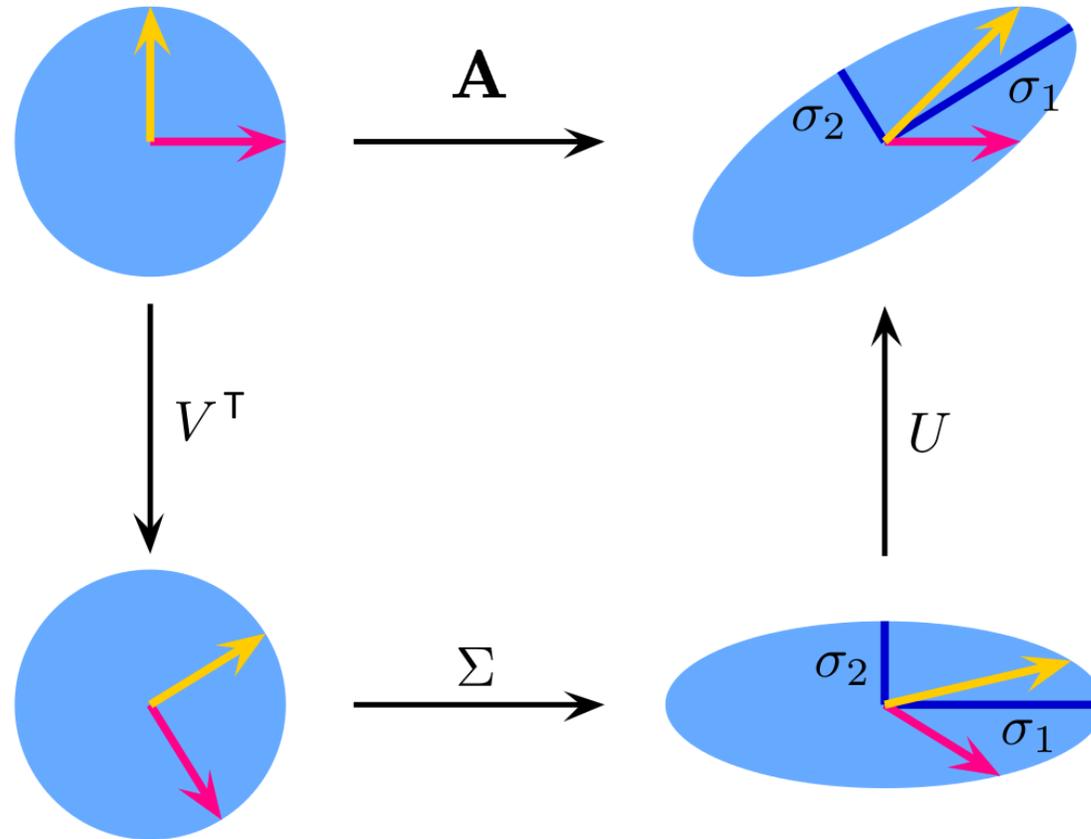
The name [Procrustes](#) ([Greek](#): Προκρούστης) refers to a bandit from Greek mythology who made his victims fit his bed either by stretching their limbs or cutting them off.

Procrustes means “he who stretches”

# Optimisation Problem

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

# Quick recap of SVD



$$A = U\Sigma V^T$$

# Quick recap of SVD

in general, applied to a real matrix:

$$\mathbf{A} = \mathbf{U}\mathbf{\Sigma}\mathbf{V}^T$$

$$\mathbf{A} \equiv (M \times N) \text{ real}$$

$$\mathbf{U} \equiv (M \times M) \text{ orthogonal, unit norm}$$

$$\mathbf{V} \equiv (N \times N) \text{ orthogonal, unit norm}$$

$$\mathbf{\Sigma} \equiv (M \times N) \text{ diagonal}$$

warning: this is not the vertex matrix!

\* See also clarification slide 1 at the end of slide deck

# Procrustes alignment steps

$$\mathbf{X} = [\mathbf{x}_1 \dots \mathbf{x}_n]^T$$

$$\mathbf{Y} = [\mathbf{y}_1 \dots \mathbf{y}_n]^T$$

Matrices of points

$$\mathbf{X}, \mathbf{Y} \in \mathbb{R}^{(N \times 3)}$$

$$\mathbf{x}_i, \mathbf{y}_i \in \mathbb{R}^{(3 \times 1)}$$

$$\bar{\mathbf{Y}}^T \mathbf{X} = \mathbf{U} \Sigma \mathbf{V}^T$$

$$\mathbf{R} = \mathbf{U} \mathbf{V}^T$$

Optimal **rotation** obtained by computing SVD on the point cross-covariance

$$\mathbf{t} = s \mathbf{R} \bar{\mathbf{x}} - \bar{\mathbf{y}}$$

**Translation** is the centroid difference

$$s = \frac{\text{tr}(\Sigma)}{\|\bar{\mathbf{X}}\|_F^2}$$

**Scale** is a quotient of eigenvalue sums

# Proof

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

# Procrustes derivation: translation

$$\mathbf{X}, \mathbf{Y} \in \mathbb{R}^{(N \times 3)}$$

$$\mathbf{x}_i, \mathbf{y}_i \in \mathbb{R}^{(3 \times 1)}$$

$$s, \mathbf{R}, \mathbf{t} = \arg \min_{s, \mathbf{R}, \mathbf{t}} E$$

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

Minimize the L2 distance between transformed source points and target points

$$= \sum_i (s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i)^\top (s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i)$$

$$= \sum_i s^2 \mathbf{x}_i^\top \mathbf{x}_i + \mathbf{t}^\top \mathbf{t} + \mathbf{y}_i^\top \mathbf{y}_i + 2s\mathbf{x}_i^\top \mathbf{R}^\top \mathbf{t} - 2s\mathbf{x}_i^\top \mathbf{R}^\top \mathbf{y}_i - 2\mathbf{t}^\top \mathbf{y}_i$$

# Procrustes derivation: translation

We remove the elements that do not depend on the translation and solve for  $\mathbf{t}$ .

$$\mathbf{t} = \arg \min_{\mathbf{t}} E = \arg \min_{\mathbf{t}} \sum_i s^2 \mathbf{x}_i^T \mathbf{x}_i + \mathbf{t}^T \mathbf{t} + \mathbf{y}_i^T \mathbf{y}_i + 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{t} - 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{y}_i - 2\mathbf{t}^T \mathbf{y}_i$$

$$= \arg \min_{\mathbf{t}} \sum_i \mathbf{t}^T \mathbf{t} + 2s \mathbf{x}_i^T \mathbf{R}^T \mathbf{t} - 2\mathbf{t}^T \mathbf{y}_i$$

$$\bar{\mathbf{x}} \equiv \frac{\sum_i \mathbf{x}_i}{N}, \bar{\mathbf{y}} \equiv \frac{\sum_i \mathbf{y}_i}{N}$$

Compute the centroid of the point clouds

$$\mathbf{t} = \arg \min_{\mathbf{t}} E = \arg \min_{\mathbf{t}} (\mathbf{t}^T (2s\mathbf{R}\bar{\mathbf{x}} + \mathbf{t} - 2\bar{\mathbf{y}})) = \bar{\mathbf{y}} - s\mathbf{R}\bar{\mathbf{x}}$$

So given  $s$  and  $\mathbf{R}$ , we can compute the translation  $\mathbf{t}$

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

subject to  $\mathbf{R} \in SO(3)$

Optimal rotation does not depend on scale

$$\mathbf{R} = \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2$$

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

$$\mathbf{R} = \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2$$

$$= \arg \min_{\mathbf{R}} \langle \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T, \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \min_{\mathbf{R}} \|\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F$$

$$= \arg \max_{\mathbf{R}} \langle \mathbf{R}, \bar{\mathbf{Y}}^T \bar{\mathbf{X}} \rangle_F$$

Optimal rotation does not depend on scale

Rotation does not change norm!

Inner product should be maximum!

# Procrustes derivation: Rotation

Subtract the centroid from the points to obtain a simpler expression for E

$$\bar{\mathbf{x}}_i \equiv \mathbf{x}_i - \bar{\mathbf{x}}, \bar{\mathbf{y}}_i \equiv \mathbf{y}_i - \bar{\mathbf{y}}$$

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 = \sum_i \|s\mathbf{R}\bar{\mathbf{x}}_i - \bar{\mathbf{y}}_i\|^2$$

$$\begin{aligned} \mathbf{R} &= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T\|_F^2 \\ &= \arg \min_{\mathbf{R}} \langle \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T, \mathbf{R}\bar{\mathbf{X}}^T - \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \min_{\mathbf{R}} \|\mathbf{R}\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \min_{\mathbf{R}} \|\bar{\mathbf{X}}^T\|_F^2 + \|\bar{\mathbf{Y}}^T\|_F^2 - 2\langle \mathbf{R}\bar{\mathbf{X}}^T, \bar{\mathbf{Y}}^T \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle \mathbf{R}, \bar{\mathbf{Y}}^T \bar{\mathbf{X}} \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle \mathbf{R}, U\Sigma V^T \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle U^T \mathbf{R} V, \Sigma \rangle_F \\ &= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V \end{aligned}$$

Optimal rotation does not depend on scale

Rotation does not change norm!

Inner product should be maximum!

Cross-covariance matrix

SVD decomposition

# Procrustes derivation: Rotation

$$= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V$$

What kind of matrix is  $S$  ?

What kind of matrix is  $\Sigma$  ?

# Procrustes derivation: Rotation

$$= \arg \max_{\mathbf{R}} \langle S, \Sigma \rangle_F \quad \text{where } S = U^T \mathbf{R} V$$

What kind of matrix is  $S$  ?

Orthogonal

What kind of matrix is  $\Sigma$  ?

Diagonal

Hence the quantity above is maximised when  $S$  equals the identity, hence:

$$I = U^T \mathbf{R} V$$

$$\mathbf{R} = UV^T$$

SVD of cross-covariance of pointsclds

$$\bar{\mathbf{Y}}^T \mathbf{X} = U \Sigma V^T$$

# Procrustes derivation: scale

Optimize **scale** given the rotation

$$s = \arg \min_s E = \arg \min_s \left( \sum_i s^2 \bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i + \bar{\mathbf{y}}_i^T \bar{\mathbf{y}}_i - 2s \bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i \right)$$

$$= \arg \min_s \left( s^2 \sum_i (\bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i) - 2s \sum_i (\bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i) \right)$$

$$= \arg \min_s (s^2 a - 2sb) = \frac{b}{a} = \frac{\sum_i (\bar{\mathbf{x}}_i^T \mathbf{R}^T \bar{\mathbf{y}}_i)}{\sum_i (\bar{\mathbf{x}}_i^T \bar{\mathbf{x}}_i)}$$

$$= \frac{\text{tr}(\bar{\mathbf{X}} \mathbf{R}^T \bar{\mathbf{Y}}^T)}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\mathbf{R}^T \bar{\mathbf{Y}}^T \bar{\mathbf{X}})}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\mathbf{V} \mathbf{U}^T \mathbf{U} \Sigma \mathbf{V}^T)}{\|\bar{\mathbf{X}}\|_F^2} = \frac{\text{tr}(\Sigma)}{\|\bar{\mathbf{X}}\|_F^2}$$

We used:

- 1) Trace invariance with shifts
- 2)  $\bar{\mathbf{Y}}^T \bar{\mathbf{X}} = \mathbf{U} \Sigma \mathbf{V}^T$ ,  $\mathbf{R} = \mathbf{U} \mathbf{V}^T$
- 3) Trace equals sum of eigenvals for square matrices

# Virtual Humans – Winter 24/25

Lecture 4\_1 – ICP: Iterative Closest Points

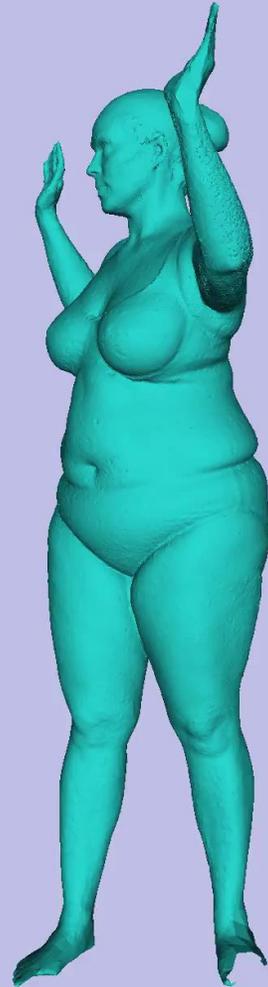
Prof. Dr.-Ing. Gerard Pons-Moll

University of Tübingen / MPI-Informatics

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# Non-rigid Articulated Registration



# What is missing?

Given correspondences, we can find the optimal rigid alignment with Procrustes.

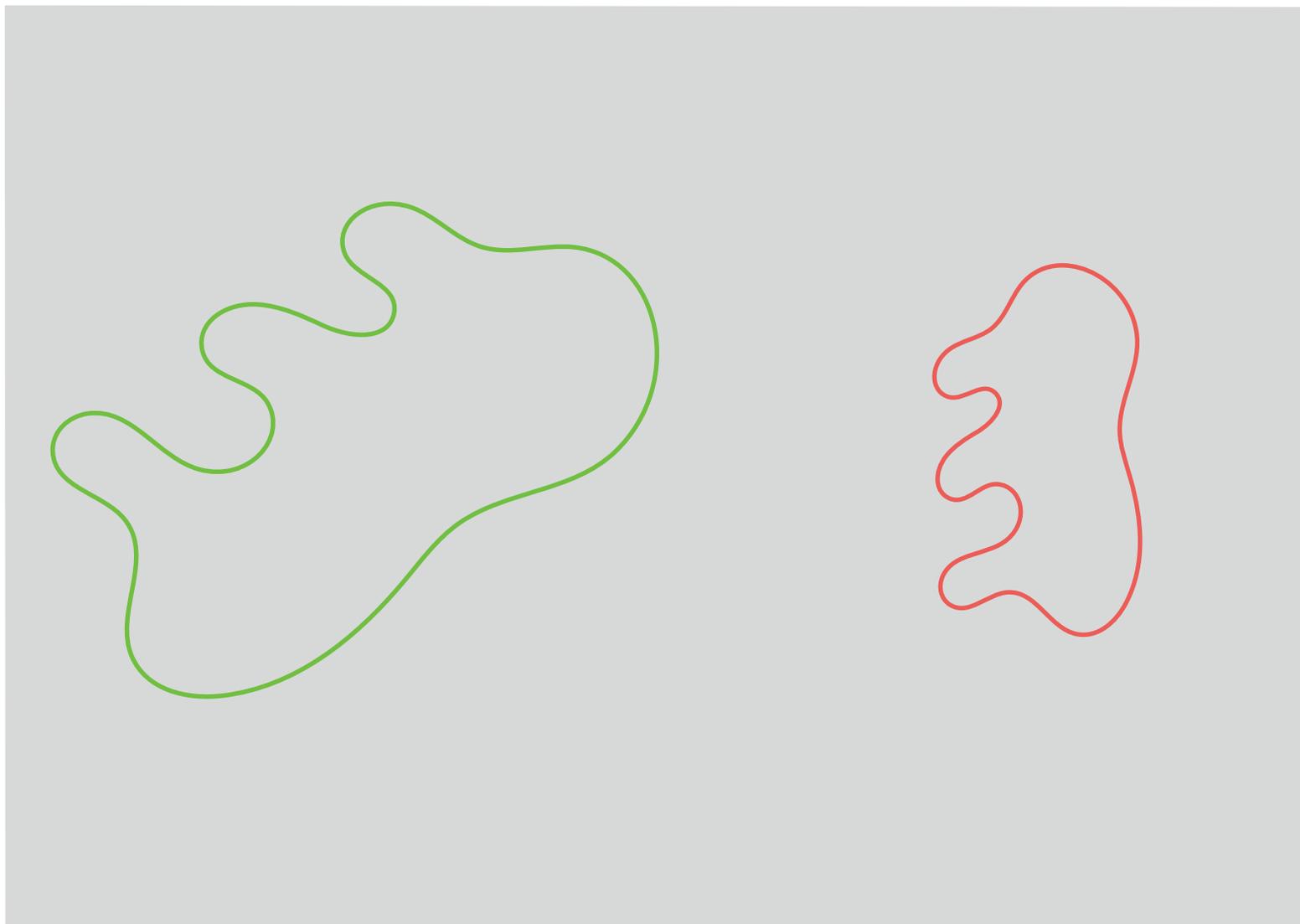
## PROBLEMS:

- How do we find the **correspondences** between shapes ?
- How do we **align** shapes **non-rigidly** ?

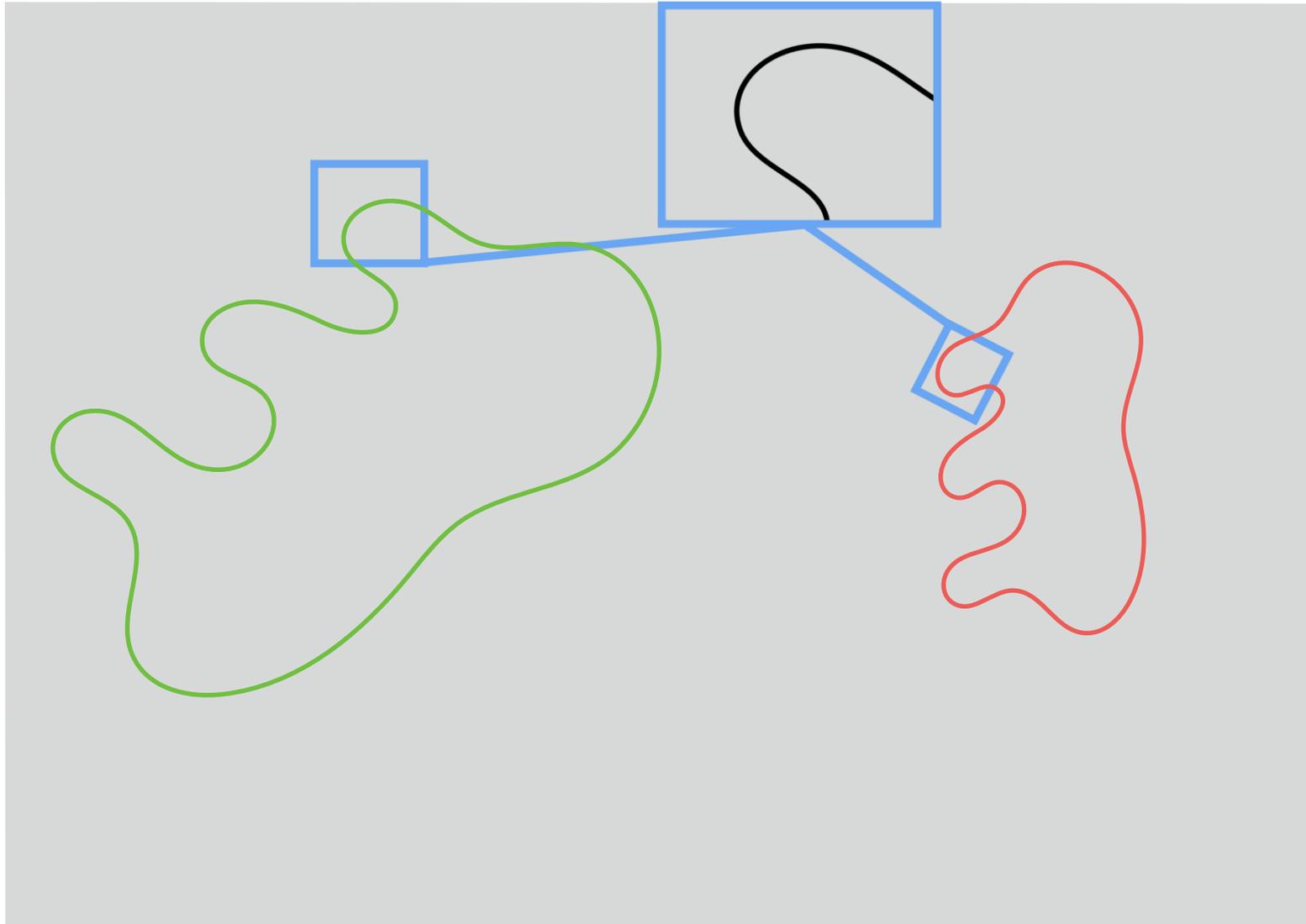
# ICP and alignment based on optimisation

- Optimising alignment and correspondences using *Iterative Closest Point (ICP)*.
- Alignment through *continuous* optimisation.

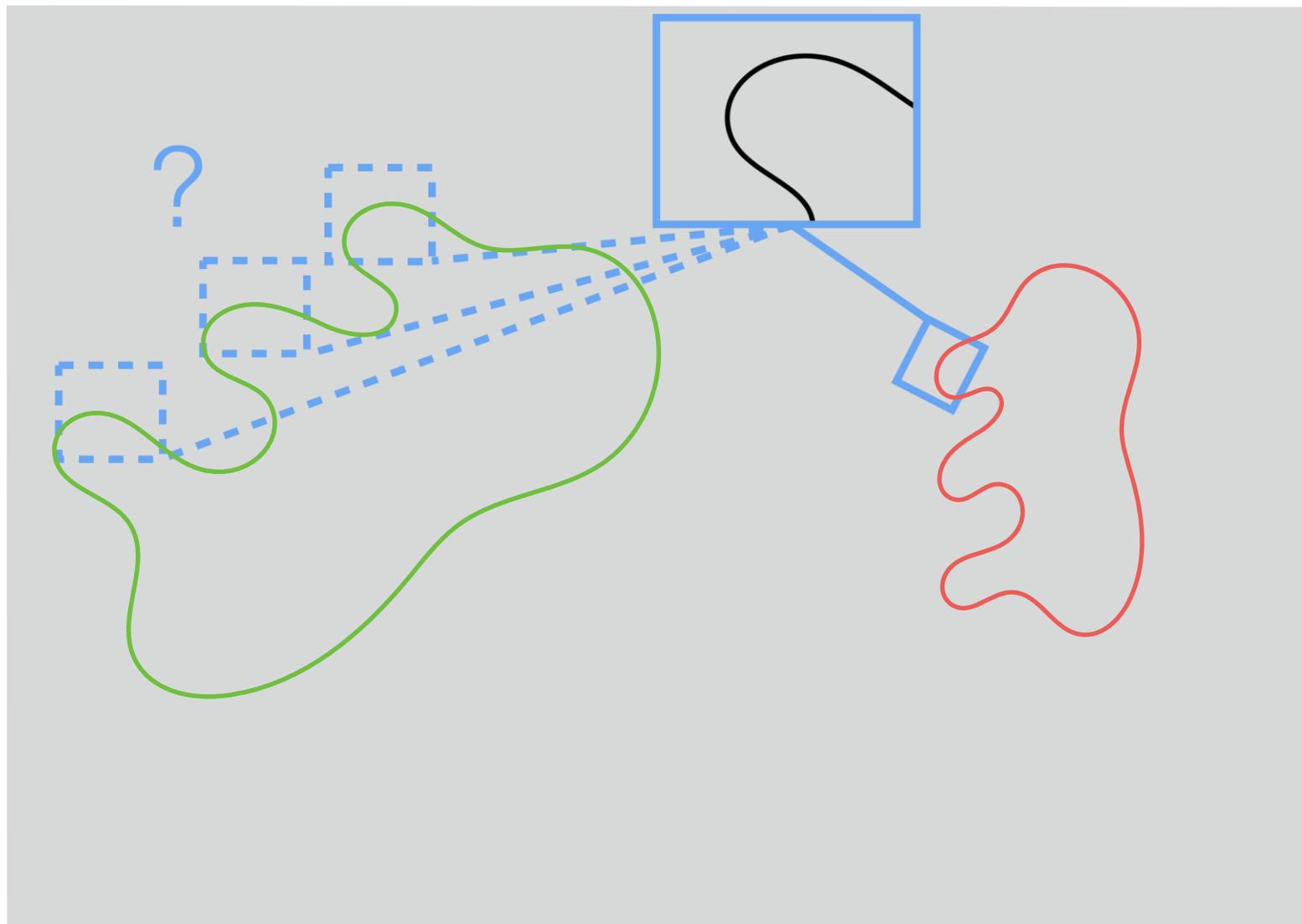
# How do we find correspondences?



# How do we find correspondences?



# How do we find correspondences?



# How do we find correspondences?

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 \equiv \sum_i \|f(\mathbf{x}_i) - \mathbf{y}_i\|^2$$

compact notation:  $f$  contains translation, rotation and isotropic scale

$\mathbf{X}_i$  Closest point to target shape point  $\mathbf{y}_i$

The optimisation is over:

- the transform  $f$
- the correspondences  $\mathcal{C} = \{(\mathbf{x}_i, \mathbf{y}_i)\}_i^N$

$$E(\mathcal{C}, f) = \sum_i \min_{\mathbf{x} \in \mathbf{X}} \|f(\mathbf{x}) - \mathbf{y}_i\|^2$$

# How do we find correspondences?

The idea was to minimise the sum of distances between the one set of points and the other set, transformed

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 \equiv \sum_i \|f(\mathbf{x}_i) - \mathbf{y}_i\|^2$$

compact notation:  $f$  contains translation, rotation and isotropic scale

# Ideas

The idea was to minimise the sum of distances between the one set of points and the other set, transformed

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 \equiv \sum_i \|f(\mathbf{x}_i) - \mathbf{y}_i\|^2$$

compact notation:  $f$  contains translation, rotation and isotropic scale

What if we estimate the correspondences?

# Solution: Iteratively find correspondences

The idea was to minimise the sum of distances between the one set of points and the other set, transformed

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 \equiv \sum_i \|f(\mathbf{x}_i) - \mathbf{y}_i\|^2$$

compact notation:  $f$  contains translation, rotation and isotropic scale

What if we estimate the correspondences?

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$
$$f^{j+1} = \arg \min_f \sum_i \|f(\mathbf{x}_i^{j+1}) - \mathbf{y}_i\|^2$$

# Alternate between finding correspondences and finding the optimal transformation

The idea was to minimise the sum of distances between the one set of points and the other set, transformed

$$E \equiv \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2 \equiv \sum_i \|f(\mathbf{x}_i) - \mathbf{y}_i\|^2$$

compact notation:  $f$  contains translation, rotation and isotropic scale

What if we estimate the correspondences?

iteration

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

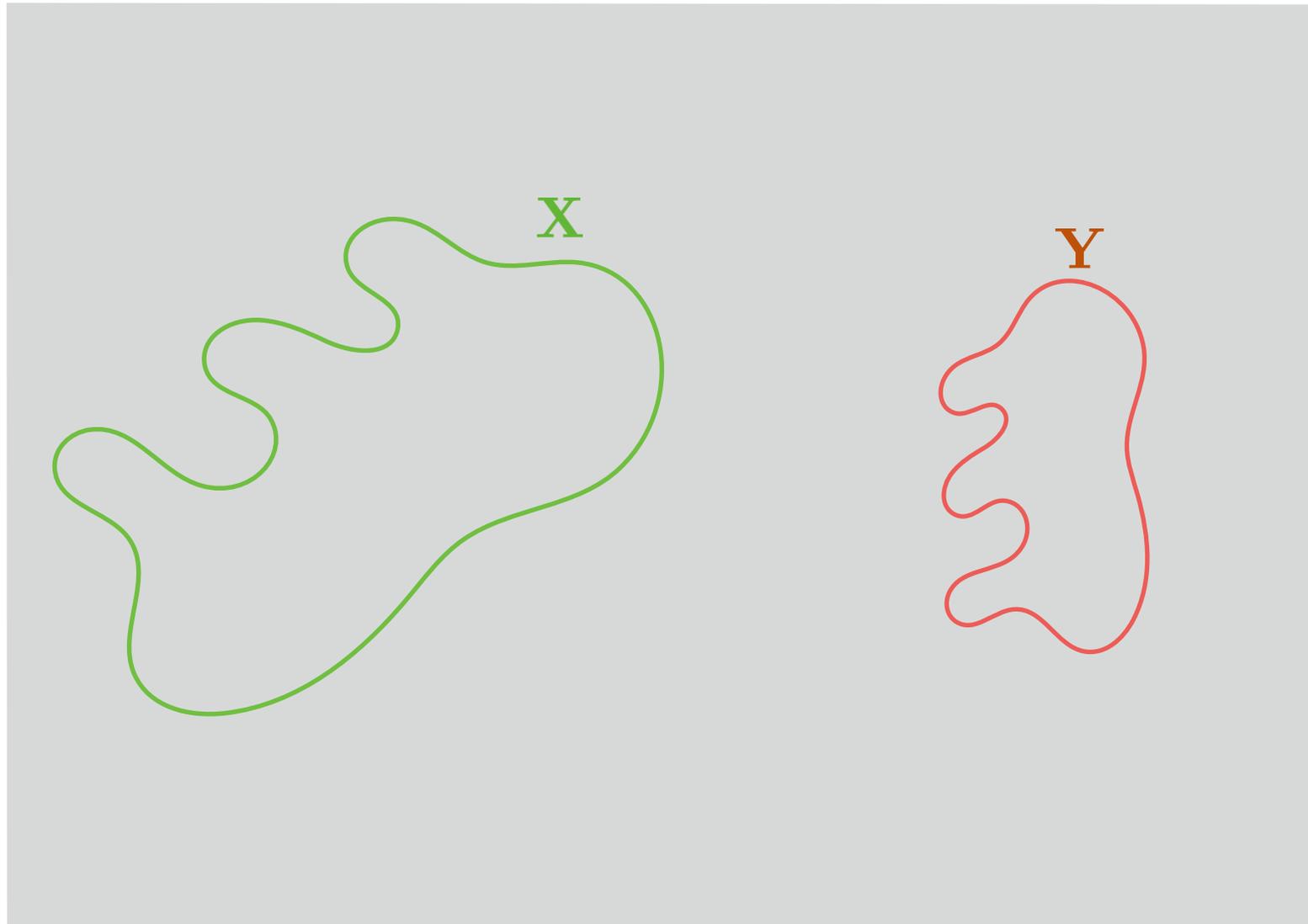
original unsorted points

$$f^{j+1} = \arg \min_f \sum_i \|f(\mathbf{x}_i^{j+1}) - \mathbf{y}_i\|^2$$

Given current best transformation, which are the closest correspondences?

Given current best correspondences, which is the best transformation?

# Make up reasonable correspondences



# Make up reasonable correspondences

$\mathbf{X} \equiv f^0(\mathbf{X})$

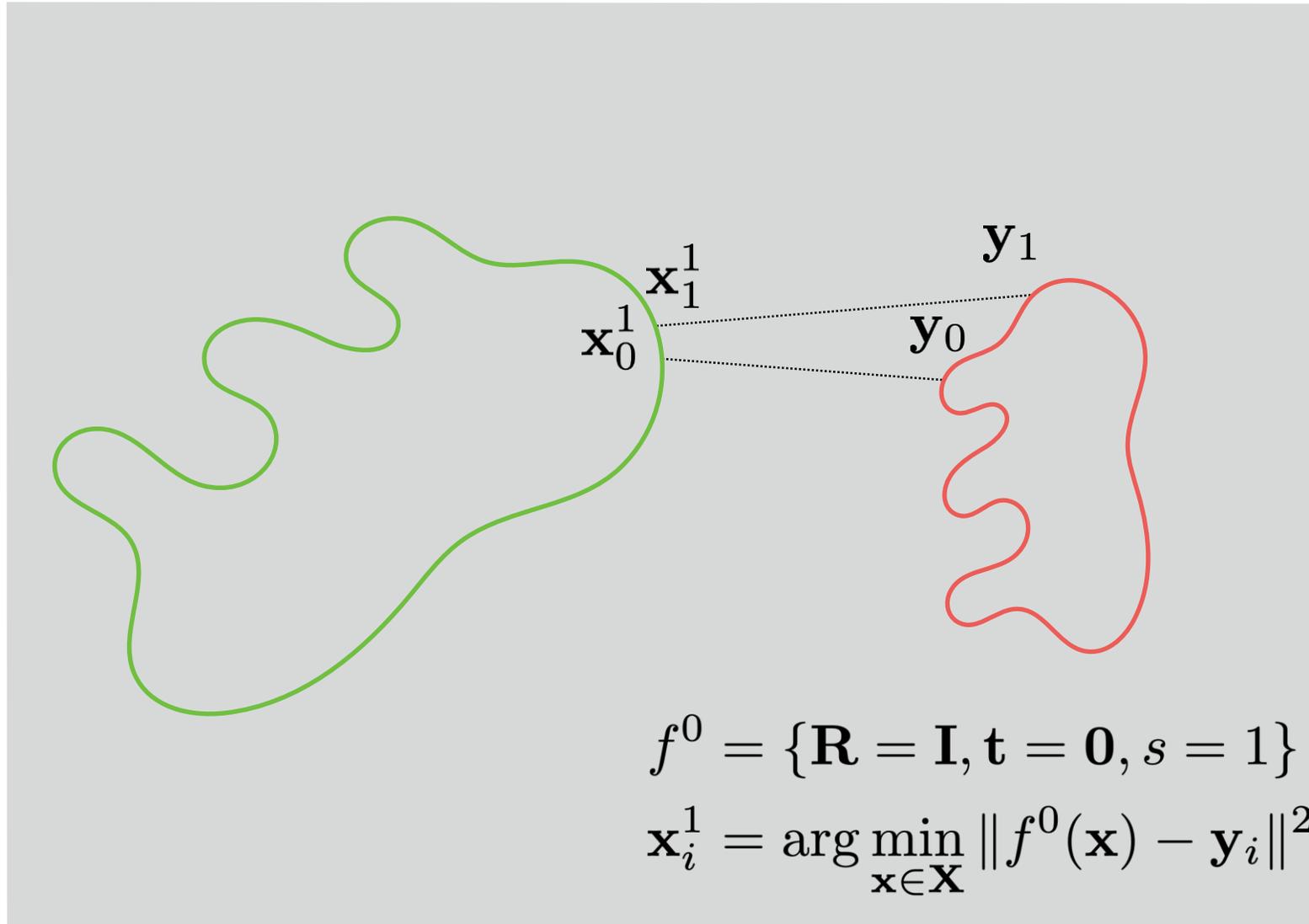
$\mathbf{x}_0^1$   $\mathbf{y}_0$

Neutral initialization.  
Initializing  $\mathbf{t}$  to align centroids  
should work better!

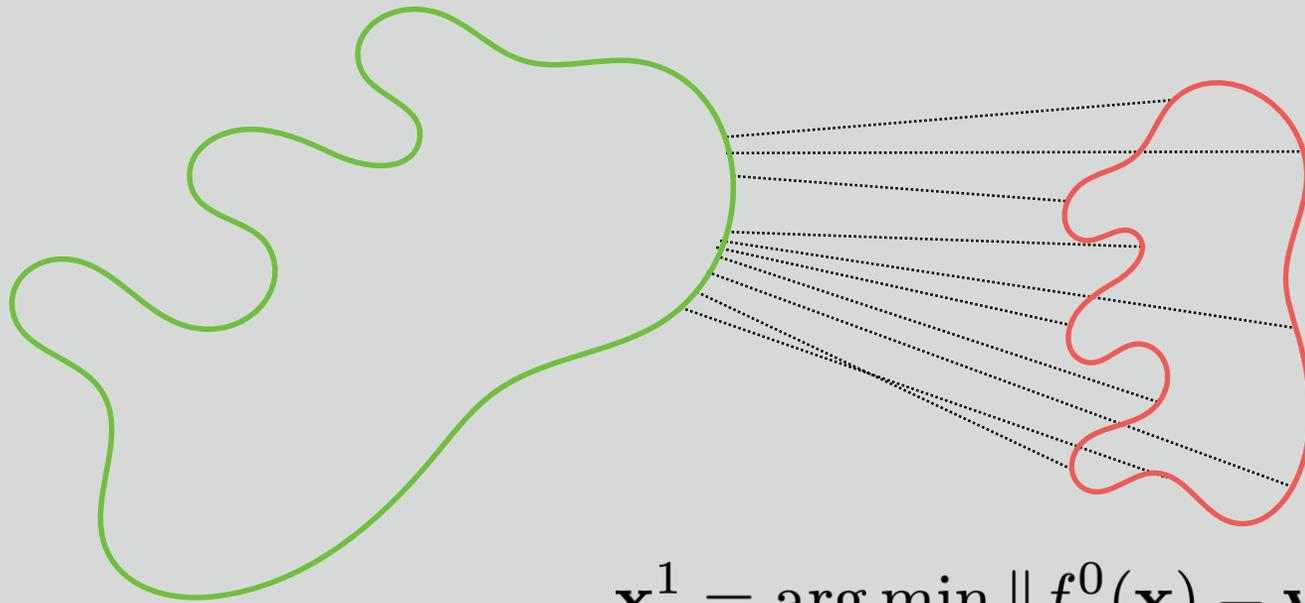
$\longrightarrow f^0 = \{\mathbf{R} = \mathbf{I}, \mathbf{t} = \mathbf{0}, s = 1\}$

$\mathbf{x}_0^1 = \arg \min_{\mathbf{x} \in \mathbf{X}} \|\mathbf{f}^0(\mathbf{x}) - \mathbf{y}_0\|^2$

# Make up reasonable correspondences



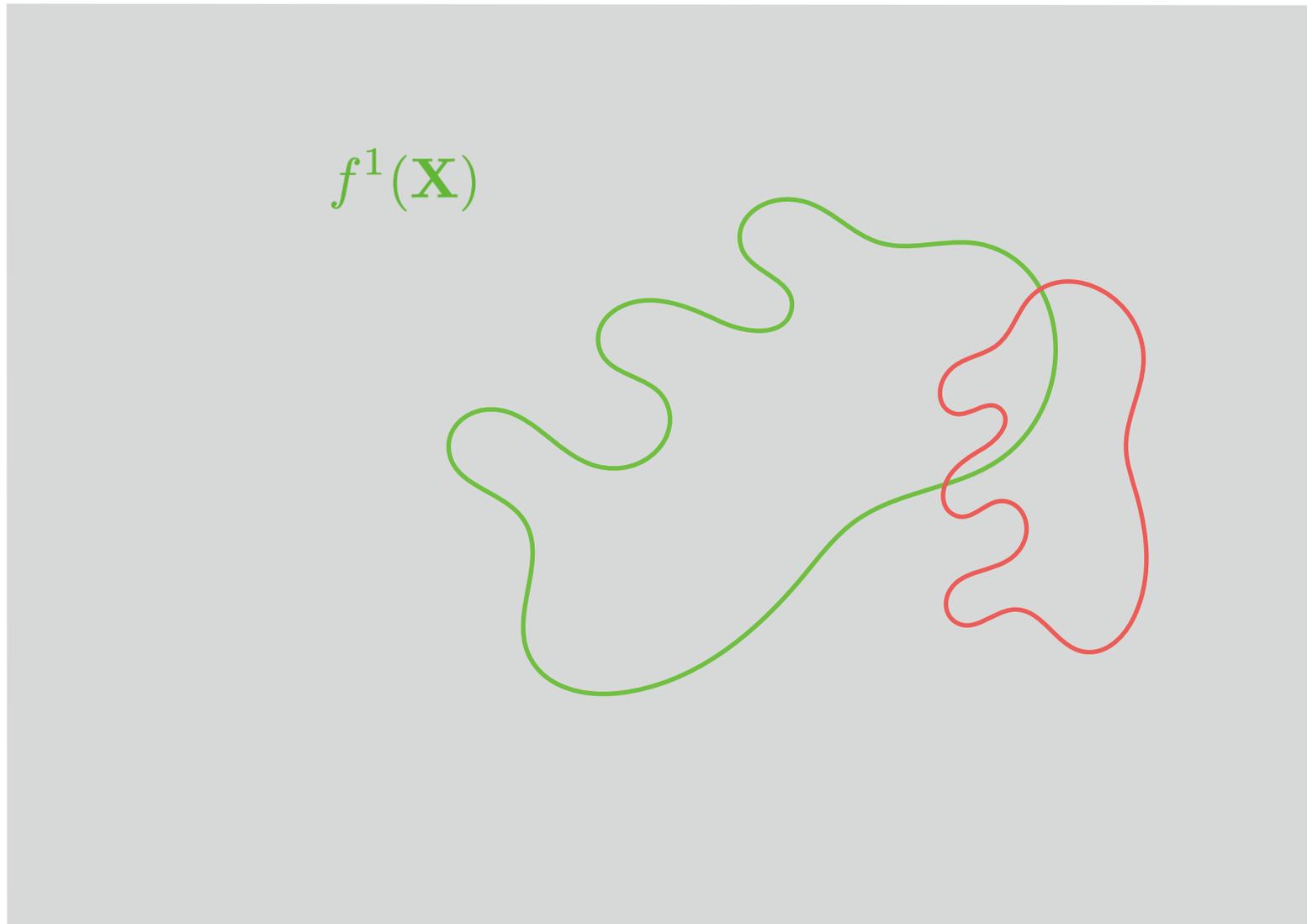
# Solve for the best transformation



$$\mathbf{x}_i^1 = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^0(\mathbf{x}) - \mathbf{y}_i\|^2$$

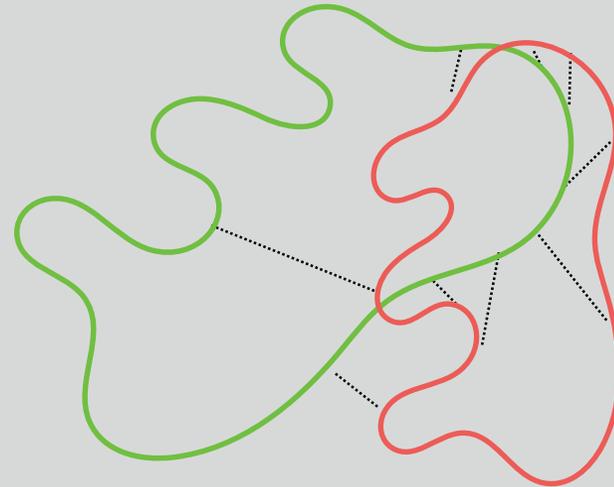
solve with procrustes  $\longrightarrow f^1 = \arg \min_f \sum_i \|f(\mathbf{x}_i^1) - \mathbf{y}_i\|^2$

Apply it ...



and iterate!

$f^1(\mathbf{X})$

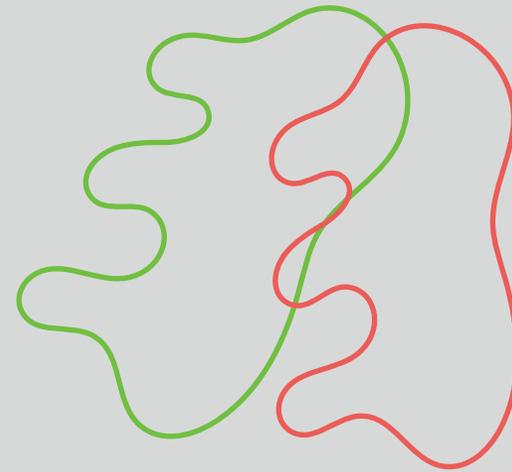


$$f^1 = \arg \min_f \sum_i \|f(\mathbf{x}_i^1) - \mathbf{y}_i\|^2$$

$$\mathbf{x}_i^2 = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^1(\mathbf{x}) - \mathbf{y}_i\|^2$$

and iterate!

$f^j(\mathbf{X})$

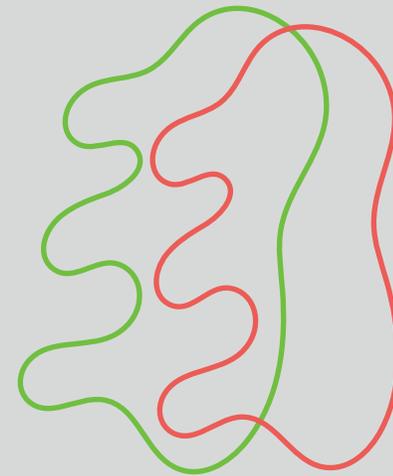


$$f^j = \arg \min_f \sum_i \|f(\mathbf{x}_i^j) - \mathbf{y}_i\|^2$$

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

and iterate!

$f^j(\mathbf{X})$



$$f^j = \arg \min_f \sum_i \|f(\mathbf{x}_i^j) - \mathbf{y}_i\|^2$$

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

and iterate!

$f^j(\mathbf{X})$



$$f^j = \arg \min_f \sum_i \|f(\mathbf{x}_i^j) - \mathbf{y}_i\|^2$$

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

and iterate!

$f^j(\mathbf{X})$



$$f^j = \arg \min_f \sum_i \|f(\mathbf{x}_i^j) - \mathbf{y}_i\|^2$$

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

# Iterative Closest Point (ICP)

1. Initialize

$$f^0 = \{\mathbf{R} = \mathbf{I}, \mathbf{t} = \frac{\sum \mathbf{y}_i}{N} - \frac{\sum \mathbf{x}_i}{N}, s = 1\}$$

typically better than 0

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2. Compute correspondences according to current best transform

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5. Converges to local minima

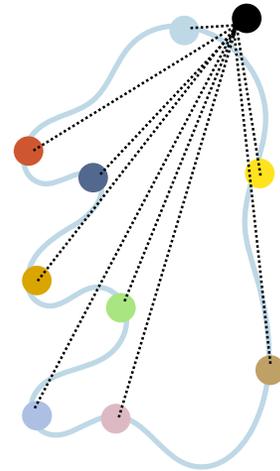
# Is ICP the best we can do?

Iteration j:

- compute closest points
- compute optimal transformation with Procrustes
- apply transformation
- terminate if converged, otherwise iterate

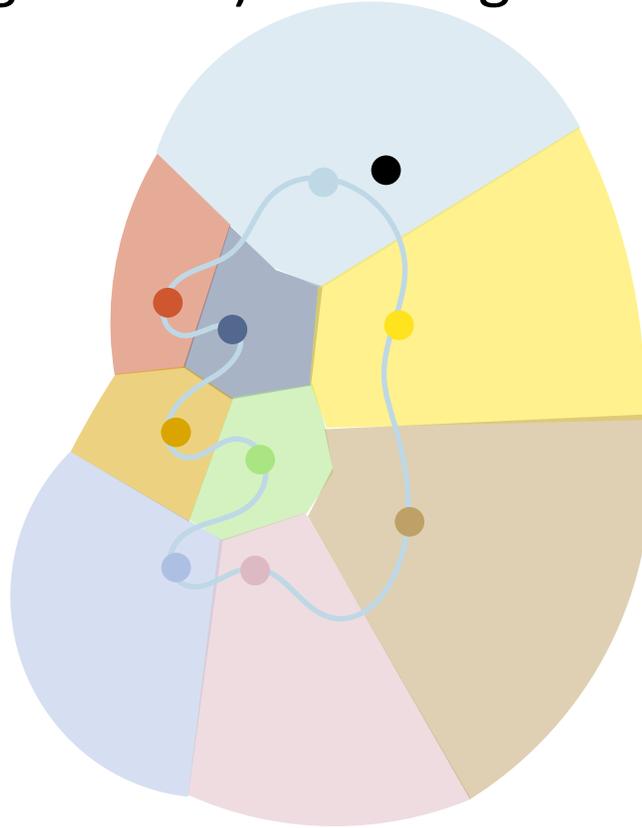
# Closest points

- Brute force is  $O(n^2)$
- For every source point find a neighbor point on the source shape



# Closest points

- Tree based methods (e.g. kdtree) have avg. complexity  $\log(n)$

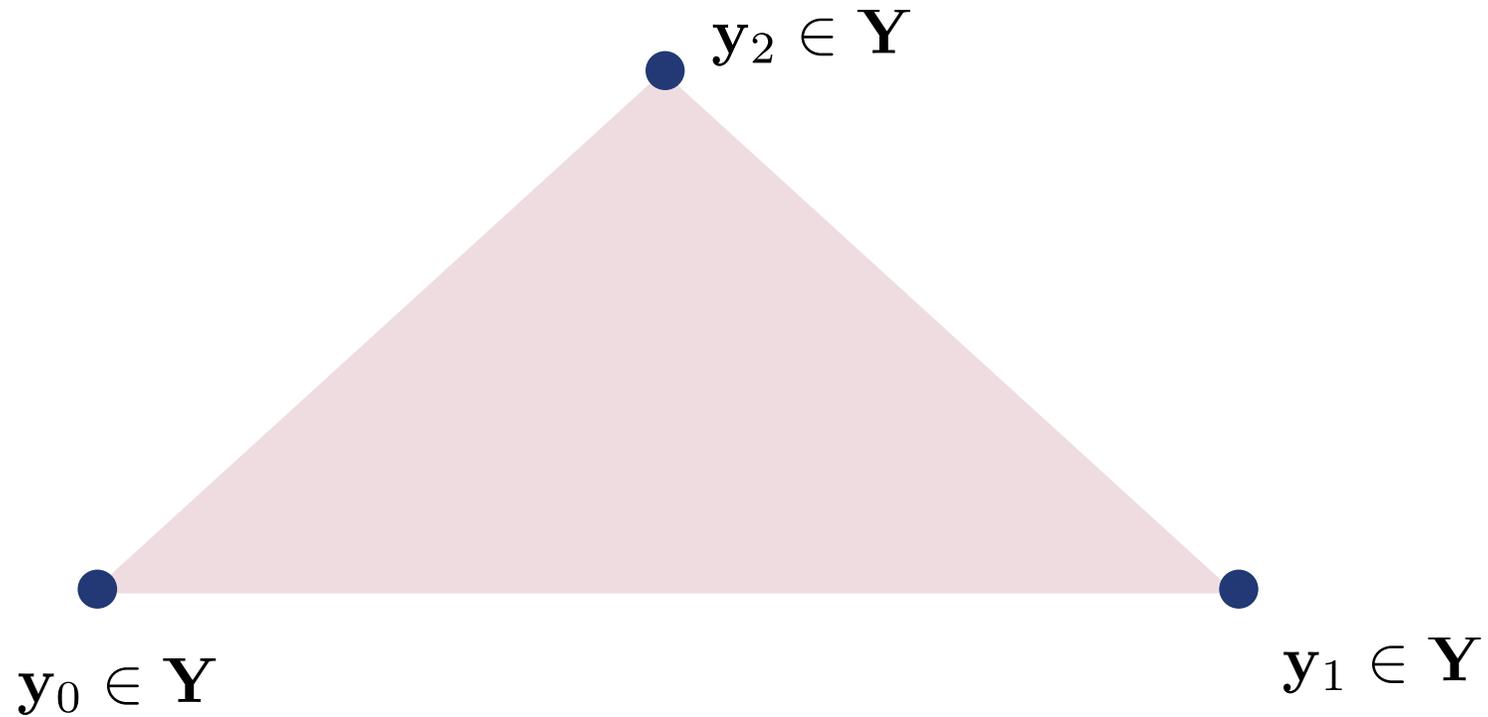


- Random point sampling also reduces the running time

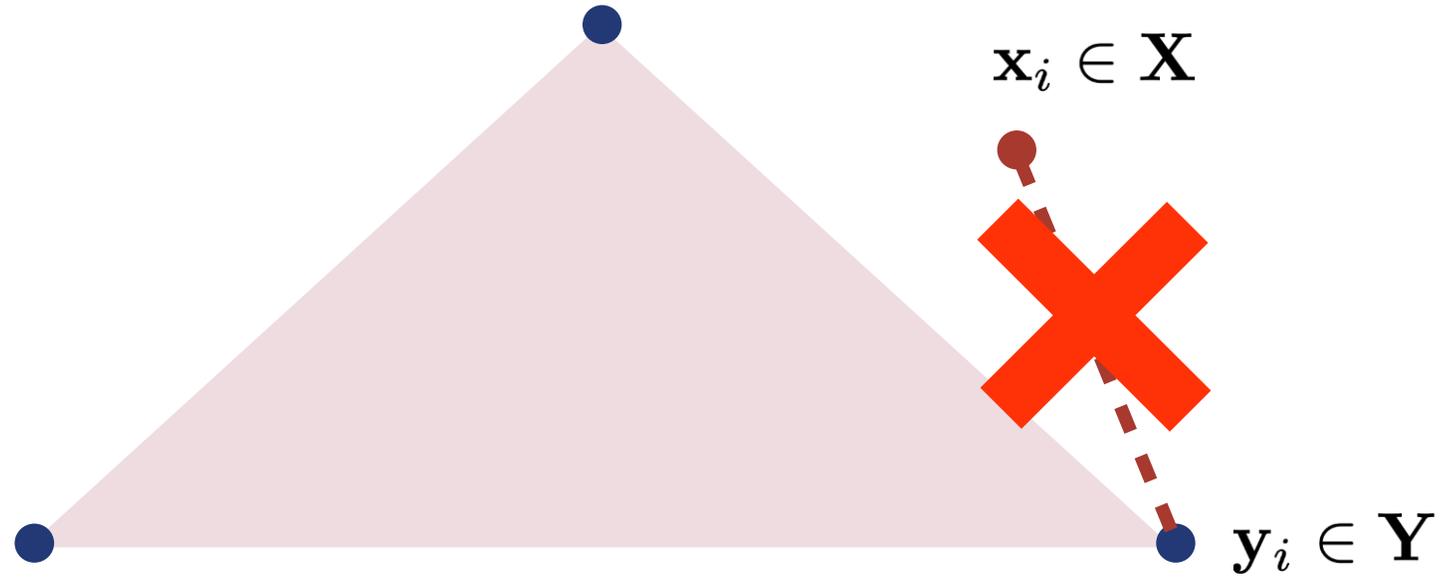
# ICP: Tips to avoid local minima

- Always find correspondences from target to source!  
Proper **data term**
- Outliers —> Robust cost functions
- Use additional information (e.g. normals)
- Compute transformation based on greedy subsets of points: RANSAC

A much better objective: Point-to-surface distance

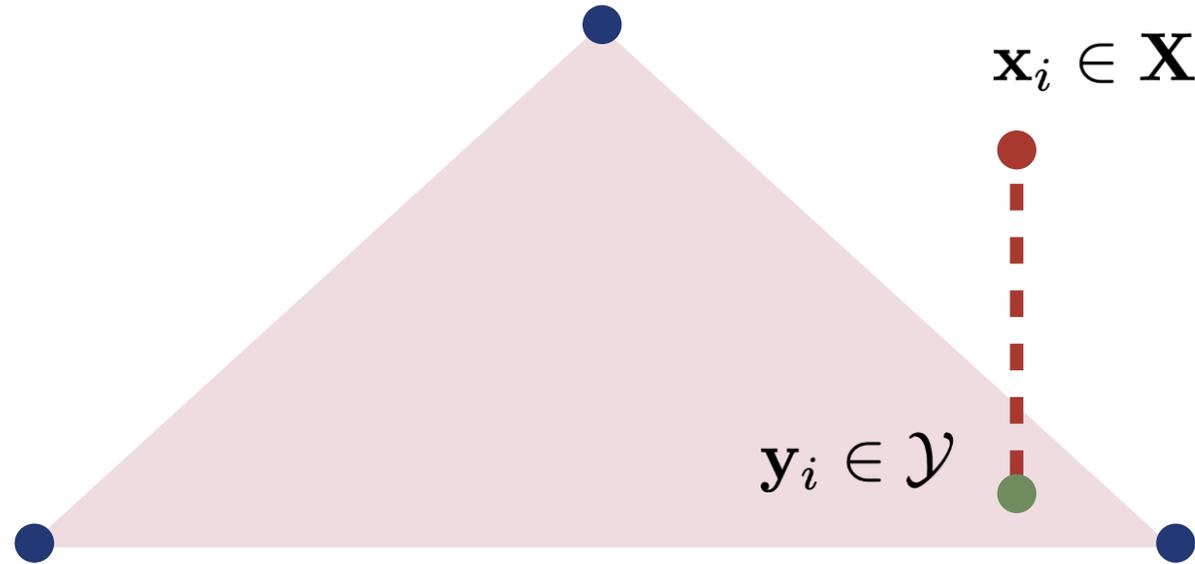


# Closest points: avoid local minima



**Point-to-point** distance

# Closest points: avoid local minima



**Point-to-surface distance**

# Is ICP the best we can do?

Iteration j:

- compute closest points
- compute optimal transformation with Procrustes
- apply transformation
- terminate if converged, otherwise iterate

# Best transformation?

- Procrustes gives us the optimal **rigid** transformation and scale given correspondences
- What if the deformation model is **not rigid** ?
- Can we generalise ICP to non-rigid deformation ?

# Iterative Closest Point (ICP)

Iteration j:

- compute closest points → Which direction to move?
- compute optimal transformation with Procrustes
- apply transformation
- terminate if converged, otherwise iterate

# Iterative Closest Point (ICP)

Iteration j:

- compute closest points → Which direction to move?
- compute optimal transformation with Procrustes →  
Compute a transform that reduces the error
- apply transformation
- terminate if converged, otherwise iterate

# Gradient-based ICP

Iteration j:

- compute closest points → Which direction to move?
- ~~compute optimal transformation with Procrustes~~ →  
Compute descent step by linearising the energy  
Jacobian of distance-based energy
- apply transformation
- terminate if converged, otherwise iterate

# Gradient-based ICP

$$\arg \min_f E(f) = \arg \min_f \sum_i \|f(\mathbf{x}_i^{j+1}) - \mathbf{y}_i\|^2$$

- If  $f$  is a rigid transformation we can solve this minimisation using Procrustes
- If  $f$  is a general non-linear function ?
- Gradient descent:  $f^{k+1} = f^k - \lambda \nabla_f E(f)$
- For least squares, is there a better optimisation method ?  
yes: Gauss-Newton based methods.

# Gradient-based ICP

1. Energy:

$$E \equiv \sum_i \left\| \min_{\mathbf{x}} f(\mathbf{x}) - \mathbf{y}_i \right\|^2$$

2. Consider the correspondences fixed in each iteration  $j+1$

$$\mathbf{x}_i^{j+1} = \arg \min_{\mathbf{x} \in \mathbf{X}} \|f^j(\mathbf{x}) - \mathbf{y}_i\|^2$$

3. Compute gradient of the energy around current estimation

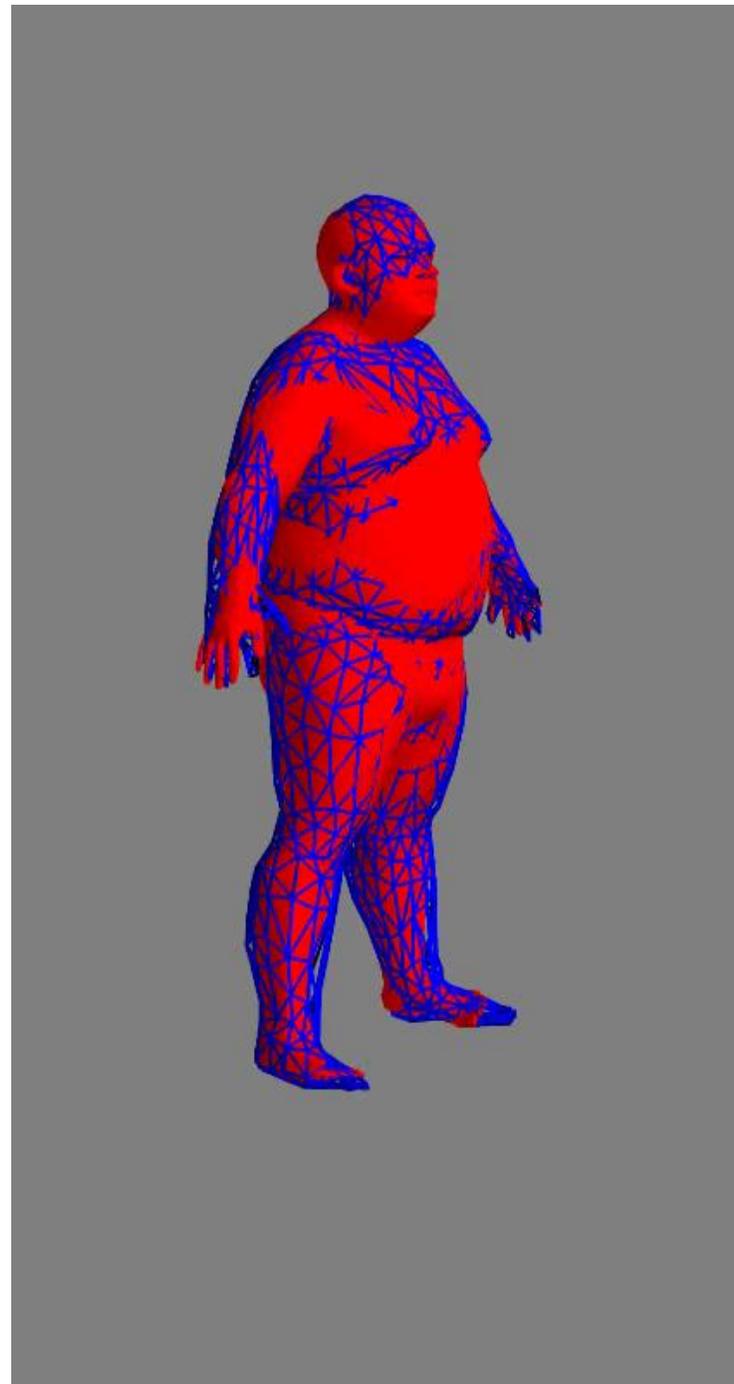
$$g^{j+1} = \nabla E(f^j)$$

4. Apply step (gradient descent, dogleg, LM, BFGS...)

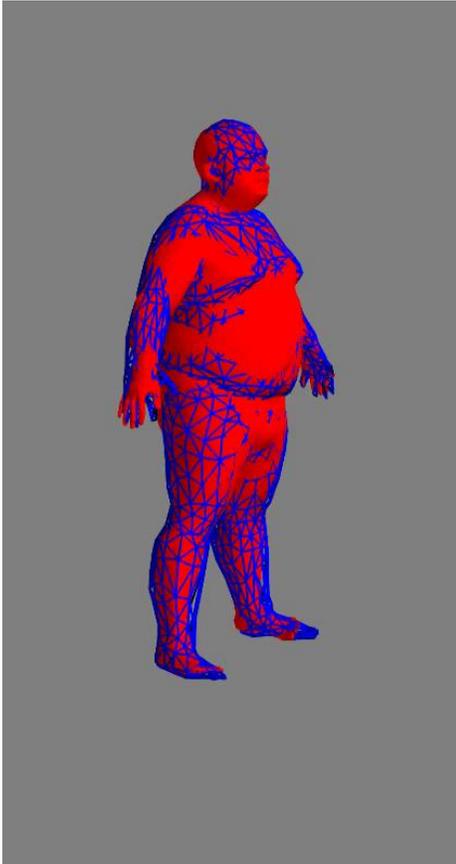
$$f^{j+1} = k_{step}(g^{0\dots j+1}, f^{0\dots j}) \longleftarrow \text{(for example } f^{j+1} = f^j - \alpha g^{j+1}\text{)}$$

5. terminate if converged, otherwise iterate (go to step 2)

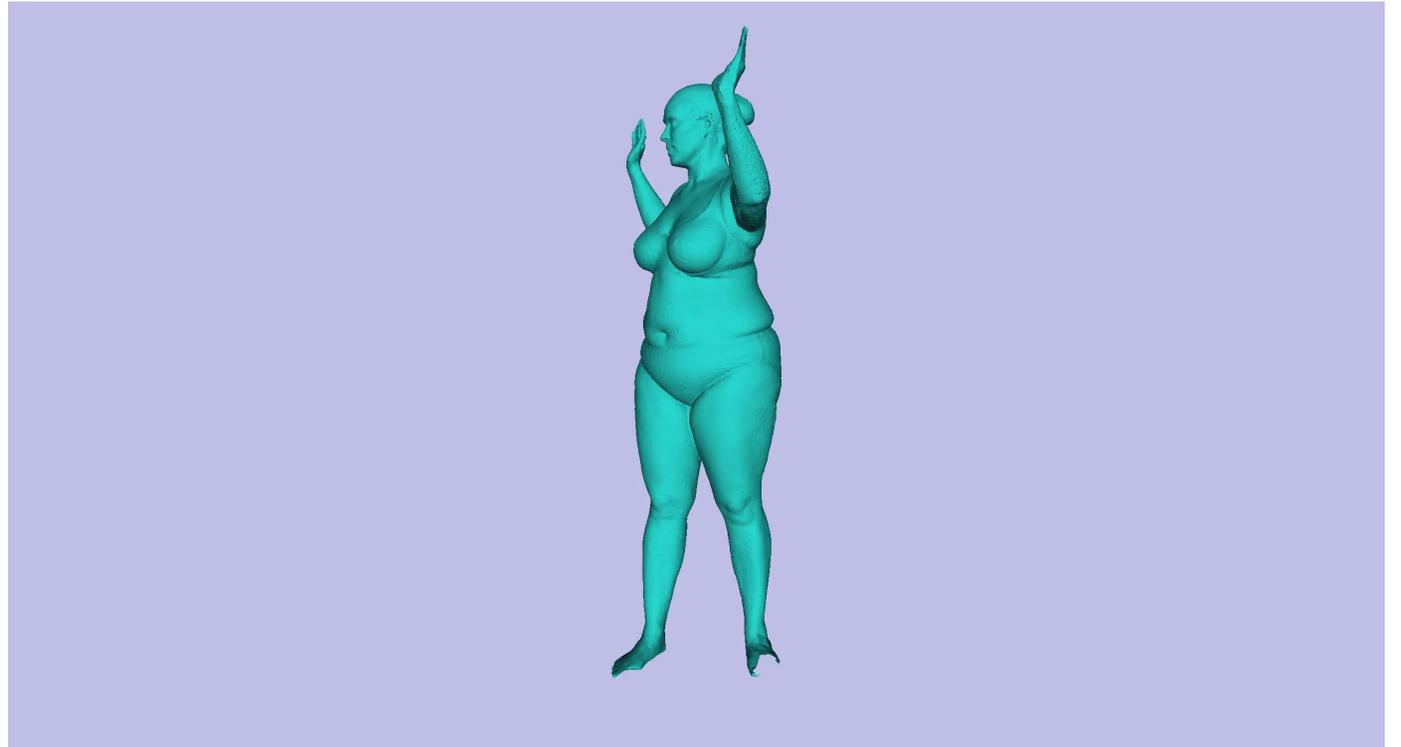
# Gradient-based ICP



# Why is convergence on the left less smooth?



Point to point objective



Point to surface objective

# Gradient-based ICP

- Energy:
- Consider the correspondences fixed in each iteration  $j+1$
- Compute gradient of the energy around current estimation
- Apply step (gradient descent, dogleg, LM, BFGS...)
- terminate if converged, otherwise iterate  $f^{j+1} = k_{step}(g^{0\dots j+1}, f^{0\dots j})$

# Gradient-based ICP

- Gradient: derivative of the sum of squared distances with respect to transformation  $f$  parameters

$$E \equiv \sum_i \left\| \min_{\mathbf{x}} f(\mathbf{x}) - \mathbf{y}_i \right\|^2$$
$$g^{j+1} = \nabla E(f^j)$$

# Gradient-based ICP

- Gradient: derivative of the sum of squared distances with respect to transformation  $f$  parameters
- Each derivative is easy
  - Who wants to writes it down?

$$E \equiv \sum_i \left\| \min_{\mathbf{x}} f(\mathbf{x}) - \mathbf{y}_i \right\|^2$$
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# Gradient-based ICP

- Gradient: derivative of the sum of squared distances with respect to transformation  $f$  parameters
- Each derivative is easy
  - Who wants to writes it down?
- **Chain rule and automatic differentiation!**

$$E \equiv \sum_i \left\| \min_{\mathbf{x}} f(\mathbf{x}) - \mathbf{y}_i \right\|^2$$
$$g^{j+1} = \nabla E(f^j)$$

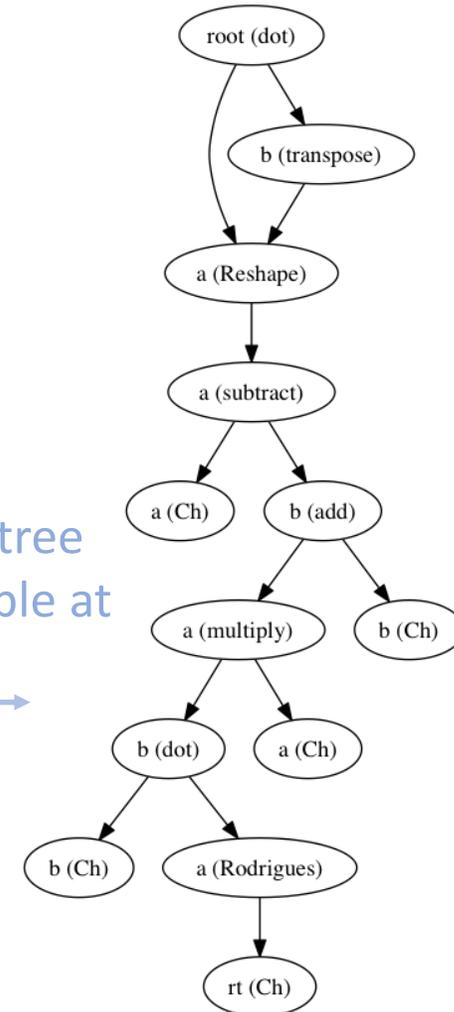
# Automatic differentiation

$$E = \sum_i \|s\mathbf{R}\mathbf{x}_i + \mathbf{t} - \mathbf{y}_i\|^2$$

write as if it was numpy code

```
1 import numpy as np
2 from numpy.linalg import norm
3 rot = np.zeros((3,3))
4 scale = np.zeros(3)
5 trans = np.zeros((3,1))
6 x = np.random.randn(3, 100)
7 y = np.random.randn(3, 100)
8 s = (y - (scale*rodriguez(rot).dot(x) + trans)).norm()
9 rot = s.dot(x.T)
10 rot = rot.dot(x)
11 import sys; sys.exit(0)
```

results in expression tree with jacobians available at each step



# Gradient-based ICP

- Energy:
- Consider the correspondences fixed in each iteration  $j+1$
- Compute gradient of the energy around current estimation
- Apply step (gradient descent, dogleg, LM, BFGS...)
- terminate if converged, otherwise iterate  $f^{j+1} = k_{step}(g^{0\dots j+1}, f^{0\dots j})$

# Why Gradient-based ICP?

- Formulation is much more generic: the energy can incorporate other terms, more parameters, etc
- A lot of available software for solving this least squares problem (cvx, ceres, ...)
- **However**, the resulting energy is non-convex for general deformation models. Optimisation can get trapped in local minima.

# Take-home message

- **Procrustes** is **optimal** for **rigid alignment problems** with *known* correspondences. For other problems:
- We can compute **correspondences** and solve for the best **transformation** iteratively with Iterative Closest Point (**ICP**)

# Slide credits

- Javier Romero