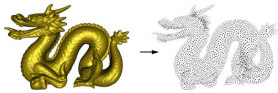


Efficient Simplification of Point-sampled Surfaces



Overview

- Introduction
- Local surface analysis
- Simplification methods
- Error measurement
- Comparison

Introduction

- Point-based models are often sampled very densely
- Many applications require coarser approximations, e.g. for efficient
 - Storage
 - Transmission
 - Processing
 - Rendering

⇒ we need simplification methods for reducing the complexity of point-based surfaces

Introduction

- We transfer different simplification methods from triangle meshes to point clouds:
 - Incremental clustering
 - Hierarchical clustering
 - Iterative simplification
 - Particle simulation
- Depending on the intended use, each method has its pros and cons (see comparison)

Local Surface Analysis

- Cloud of point samples describes underlying (manifold) surface
- We need:
 - mechanisms for locally approximating the surface ⇒ MLS approach
 - fast estimation of tangent plane and curvature ⇒ principal component analysis of local neighborhood

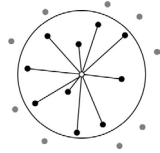
Neighborhood

- No explicit connectivity between samples (as with triangle meshes)
- Replace geodesic proximity with spatial proximity (requires sufficiently high sampling density!)
- Compute neighborhood according to Euclidean distance

Neighborhood



- k-nearest neighbors

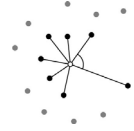


- can be quickly computed using spatial data-structures (e.g. kd-tree, octree, bsp-tree)
- requires isotropic point distribution

Neighborhood



- Improvement: angle criterion (Linsen)



- project points onto tangent plane
- sort neighbors according to angle
- include more points if angle between subsequent points is above some threshold

Neighborhood



- Local Delaunay triangulation (Floater)



- project points into tangent plane
- compute local Voronoi diagram

Covariance Analysis



- Covariance matrix of local neighborhood N:

$$\mathbf{C} = \begin{bmatrix} \mathbf{p}_i - \bar{\mathbf{p}} \\ \dots \\ \mathbf{p}_i - \bar{\mathbf{p}} \end{bmatrix}^T \cdot \begin{bmatrix} \mathbf{p}_i - \bar{\mathbf{p}} \\ \dots \\ \mathbf{p}_i - \bar{\mathbf{p}} \end{bmatrix}, \quad i_j \in N$$

- with centroid $\bar{\mathbf{p}} = \frac{1}{|N|} \sum_{i \in N} \mathbf{p}_i$

Covariance Analysis



- Consider the eigenproblem:

$$\mathbf{C} \cdot \mathbf{v}_l = \lambda_l \cdot \mathbf{v}_l, \quad l \in \{0,1,2\}$$

- C is a 3x3, positive semi-definite matrix
 - ⇒ All eigenvalues are real-valued
 - ⇒ The eigenvector with smallest eigenvalue defines the least-squares plane through the points in the neighborhood, i.e. approximates the surface normal

Covariance Analysis



- The total variation is given as:

$$\sum_{i \in N} |\mathbf{p}_i - \bar{\mathbf{p}}|^2 = \lambda_0 + \lambda_1 + \lambda_2$$

- We define surface variation as:

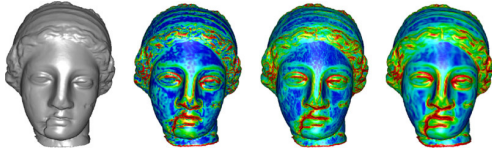
$$\sigma_n(\mathbf{p}) = \frac{\lambda_0}{\lambda_0 + \lambda_1 + \lambda_2}, \quad \lambda_0 \leq \lambda_1 \leq \lambda_2$$

- measures the fraction of variation along the surface normal, i.e. quantifies how strong the surface deviates from the tangent plane ⇒ estimate for curvature

Covariance Analysis



- Comparison with curvature:



original mean curvature variation n=20 variation n=50

Surface Simplification



- Incremental clustering
- Hierarchical clustering
- Iterative simplification
- Particle simulation

Incremental Clustering

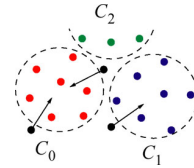


- Clustering by region-growing:
 - Start with random seed point
 - Successively add nearest points to cluster until cluster reaches maximum size
 - Choose new seed from remaining points
- Growth of clusters can also be bounded by surface variation
 - ⇒ Curvature adaptive clustering

Incremental Clustering



- Incremental growth leads to internal fragmentation
 - ⇒ assign stray samples to closest cluster



- Note: this can increase maximum size and variation bounds!

Incremental Clustering



- Replace each cluster by its centroid



original model with color-coded clusters (34,384 points)



simplified model (1,000 points)

Hierarchical Clustering

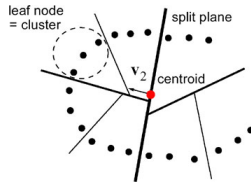


- Top-down approach using binary space partition:
 - Split the point cloud if:
 - Size is larger than user-specified maximum or
 - Surface variation is above maximum threshold
 - Split plane defined by centroid and axis of greatest variation (= eigenvector of covariance matrix with largest associated eigenvector)
- Leaf nodes of the tree correspond to clusters

Hierarchical Clustering



- 2D example



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Hierarchical Clustering



- Adaptive clustering



original model with color-coded clusters (34,384 points)



simplified model (1,000 points)

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Iterative Simplification



- Iteratively contracts point pairs
 - ⇒ Each contraction reduces the number of points by one
- Contractions are arranged in priority queue according to quadric error metric (Garland and Heckbert)
- Quadric measures cost of contraction and determines optimal position for contracted sample
- Equivalent to QSlm except for definition of approximating planes

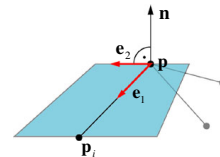
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Iterative Simplification



- Quadric measures the squared distance to a set of planes defined over *edges* of neighborhood
 - plane spanned by vectors $e_1 = p_i - p$ and $e_2 = e_1 \times n$



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Iterative Simplification



original model (187,664 points)



simplified model (1,000 points)



remaining point pair contraction candidates

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Particle Simulation



- Resample surface by distributing particles on the surface
- Particles move on surface according to inter-particle repelling forces
- Particle relaxation terminates when equilibrium is reached (requires damping)
- Can also be used for up-sampling!

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Particle Simulation

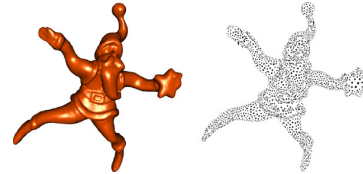


- Initialization
 - randomly spread particles
- Repulsion
 - linear repulsion force $F_i(\mathbf{p}) = k(r - \|\mathbf{p} - \mathbf{p}_i\|) \cdot (\mathbf{p} - \mathbf{p}_i)$
 - ⇒ only need to consider neighborhood of radius r
- Projection
 - keep particles on surface by projecting onto tangent plane of closest point
 - apply full MLS projection at end of simulation

Particle Simulation



- Adaptive simulation
 - Adjust repulsion radius according to surface variation
 - ⇒ more samples in regions of high variation



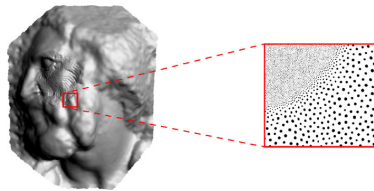
original model
(75,781 points)

simplified model
(6,000 points)

Particle Simulation



- User-controlled simulation
 - Adjust repulsion radius according to user input



Measuring Error

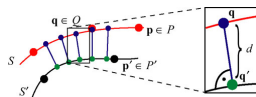


- Measure the distance between two point-sampled surfaces using a sampling approach
- Maximum error: $\Delta_{\max}(S, S') = \max_{\mathbf{q} \in Q} d(\mathbf{q}, S')$
- ⇒ Two-sided Hausdorff distance
- Mean error: $\Delta_{\text{avg}}(S, S') = \frac{1}{|Q|} \sum_{\mathbf{q} \in Q} d(\mathbf{q}, S')$
- ⇒ Area-weighted integral of point-to-surface distances
- Q is an up-sampled version of the point cloud that describes the surface S

Measuring Error



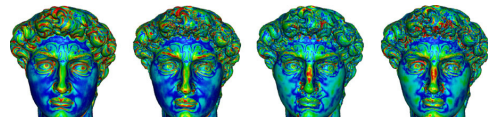
- $d(\mathbf{q}, S')$ measures the distance of point \mathbf{q} to surface S' using the MLS projection operator with linear basis functions



Comparison



- Error estimate for Michelangelo's David simplified from 2,000,000 points to 5,000 points

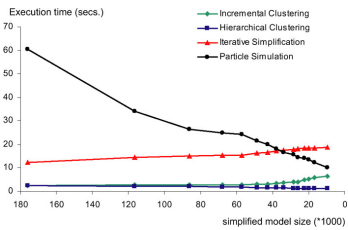


$\Delta_{\max} = 0.0049$ $\Delta_{\text{avg}} = 0.12 \cdot 10^{-4}$ $\Delta_{\max} = 0.0049$ $\Delta_{\text{avg}} = 0.14 \cdot 10^{-4}$ $\Delta_{\max} = 0.0052$ $\Delta_{\text{avg}} = 2.43 \cdot 10^{-4}$ $\Delta_{\max} = 0.0061$ $\Delta_{\text{avg}} = 2.68 \cdot 10^{-4}$
 (a) uniform incremental clustering (b) adaptive hierarchical clustering (c) iterative simplification (d) particle simulation

Comparison



- Execution time as a function of target model size (input: dragon, 535,545 points)



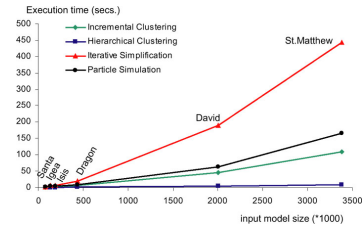
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Comparison



- Execution time as a function of input model size (reduction to 1%)



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Comparison



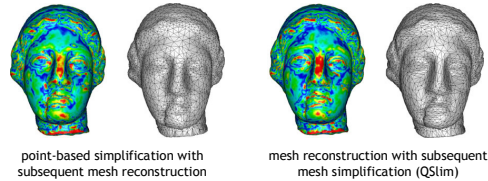
- Summary

	Efficiency	Surface Error	Control	Implementation
Incremental Clustering	+	-	-	+
Hierarchical Clustering	+	-	-	+
Iterative Simplification	-	+	o	o
Particle Simulation	o	+	+	-

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Point-based vs. Mesh Simplification



⇒ point-based simplification saves an expensive surface reconstruction on the dense point cloud!

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References



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- Garland, Heckbert: *Surface Simplification using Quadric Error Metrics*, SIGGRAPH 1997
- Turk: *Re-Tiling Polygonal Surfaces*, SIGGRAPH 1992
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