

3D Spherical based Segmentation and Registration

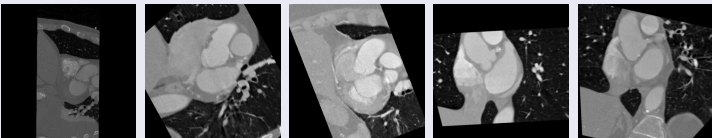
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December 2, 2008



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Comparing CT datasets

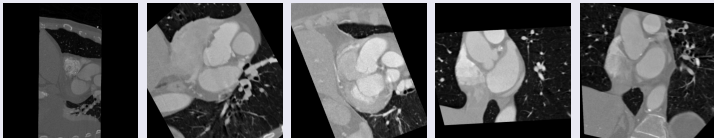


Selected CT slices from different patients showing similar areas
(Source: Nuklearmed. Klinik der TU Muenchen)

- Comparing different CT datasets taken at different time
- Matching of 3D datasets performed by hand takes a lot of time
- Existing algorithms work on projective matrices...
 - - matching a large amount of points created by edge detection
 - - computing the difference of every data domain voxel
 - - ...
- ⇒ Use (blood) vessels as (more) characteristic data with **less data for representation**



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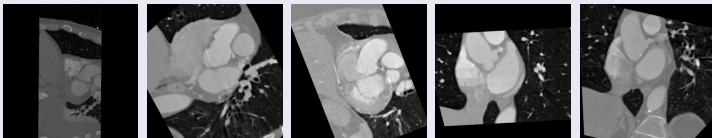


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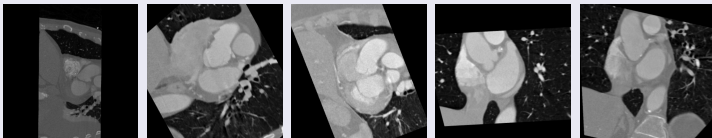
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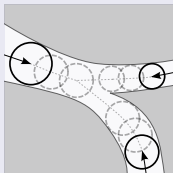


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Simulation of spheres moving through vessels

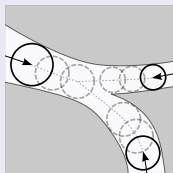


Simulated movement of sphere within the blood vessel

- **Simulate the movement** of spheres through vessels touching the borders like a chimney-sweeper
- **Sphere radii are variable** and grow/shrink to touch the vessel borders
- Spheres stay in the **center of the vessels**
- The radius of **the sphere representing the blood vessels** is stored at each **center point**
- **Direct implementation would be too inefficient** due to collision tests, realignment of sphere, etc.



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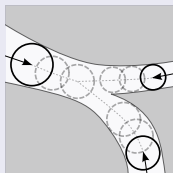


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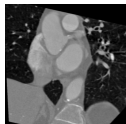


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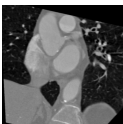


- **Raw** CT data
- Growing **Spheres**
- Computation of **Gradient**
- **Particle** emission
- ...

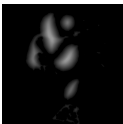


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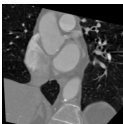
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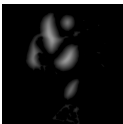


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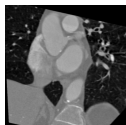
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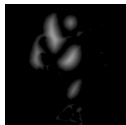


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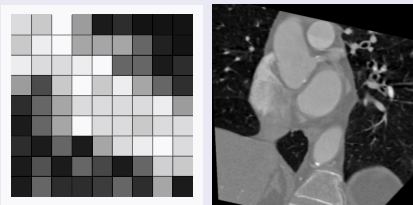
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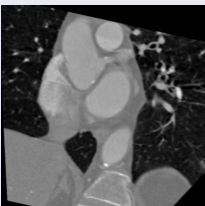
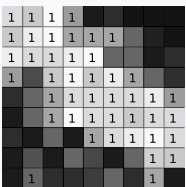
Spheres - Window



- Values for coronary contrast media are usually **within a specific window**



Spheres - Threshold flag field

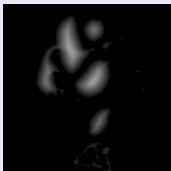
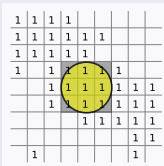


- Thresholding CT data by the window range $[win_{min}; win_{max}]$
 - Flag field **speeds up computations**
 - Important for **convolution** in frequency room (later)
 - Typical values for coronary contrast media: $[150; 1000]$

$$FlagData_{pos} = \begin{cases} 1 & win_{min} < value_{pos} < win_{max} \\ 0 & else \end{cases}$$



Spheres - First spherical test



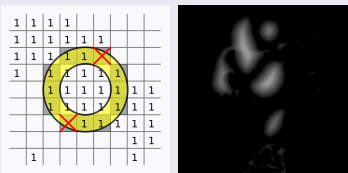
- **Avoiding early stop** of spherical growing on **noisy data**:
 - Start with a radius *StartRadius*
 - Abort if there are **too many mismatching flags** within the sphere
 - Output value $SphereData_{pos}$ of current voxel is set to **0** if **first spherical test was not successful**

Start sphere abort criteria

$$\frac{\sum_{pos \in StartSphere} FlagData_{pos}}{|Voxels \text{ in Sphere}|} < MaxMismatch$$



Spheres - Growing spheres, radius 3



- If there was no output data set continue growing the sphere
 - Growing is stopped if too many **mismatching voxels on the sphere surface exceed a specific error value**
 - Output value $SphereData_{pos}$ is set to the **current sphere radius if the abort criteria is met**

Sphere growing abort criteria

$$\frac{\sum_{pos \in SphereSurface} FlagData_{pos}}{|Sphere\ Surface|} < MaxMismatch$$



Spheres - Growing spheres, radius 4



Left: Mismatch value 9/16 exceeds the allowed rate *MaxMismatch*

Middle: Stored radius values after applying spherical filter

Right: Spherical dataset created by the spherical filter

- Sphere growing for every voxel returns data set with the following **properties**:
 - The sphere radii **represent the blood vessels** with a diameter of at least $2 \cdot \textit{StartRadius}$
 - Blood vessels could be **reconstructed** with the spherical dataset by **joining the sphere volumes**
 - Spheres totally covered by larger spheres can be dropped if we are only interested in a representative data for blood vessels

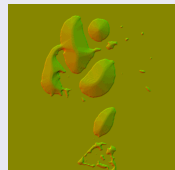


Gradient - Computation

5	3	3	2				
3	3	3	2				
2	2	3	2	2			
		2	3	3	2		
			2	3	2	2	
			2	2	3	2	2
				2	2	2	2
					2	2	3
						2	2



0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0
0.0	0.1	-1.0	-3.0	-2.2	0.0	0.0	0.0
0.0	1.3	0.1	-1.1	-2.3	-2.2	0.0	0.0
0.0	2.2	3.3	1.0	-1.1	-3.2	-2.2	0.0
0.0	0.0	2.2	3.1	0.1	-1.1	-2.2	-2.2
0.0	0.0	2.0	2.2	1.1	0.0	-1.0	0.2
0.0	0.0	0.0	2.2	2.2	0.3	0.0	0.0
0.0	0.0	0.0	0.0	0.2	2.2	2.2	1.2
0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0



Gradient computation with central differences

$$\text{GradientData}_{pos} = \begin{pmatrix} \frac{\delta \text{SphereData}_{pos}}{\delta x} \\ \frac{\delta \text{SphereData}_{pos}}{\delta y} \\ \frac{\delta \text{SphereData}_{pos}}{\delta z} \end{pmatrix}$$

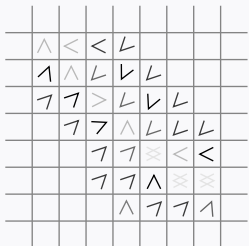
$$= \begin{pmatrix} \text{SphereData}_{pos+(1,0,0)} - \text{SphereData}_{pos-(1,0,0)} \\ \text{SphereData}_{pos+(0,1,0)} - \text{SphereData}_{pos-(0,1,0)} \\ \text{SphereData}_{pos+(0,0,1)} - \text{SphereData}_{pos-(0,0,1)} \end{pmatrix} \cdot 0.5$$



Gradient - Meaning

Gradient vectors scaled by 2

0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0
0,0	0,1	-1,0	-3,0	-2,-2	0,0	0,0	0,0	0,0	0,0
0,0	1,3	0,1	-1,-1	-2,-3	-2,-2	0,0	0,0	0,0	0,0
0,0	2,2	3,3	1,0	-1,-1	-3,-2	-2,-2	0,0	0,0	0,0
0,0	0,0	2,2	3,1	0,1	-1,-1	-2,-2	-2,-2	0,0	0,0
0,0	0,0	2,0	2,2	1,1	0,0	-1,0	0,-2	0,0	0,0
0,0	0,0	0,0	2,2	2,2	0,3	0,0	0,0	0,0	0,0
0,0	0,0	0,0	0,0	0,2	2,2	2,2	1,2	0,0	0,0
0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0	0,0

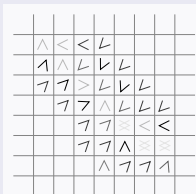


- Gradients aim to the local center of largest sphere in neighborhood
 - ⇒ **Can be used for efficient simulation of origin problem**
 - Growing** the sphere **forces movement to the center** of vessel
 - Movement **direction** is **given by gradient**
- Gradient can be **smoothed** if sphere data has a high frequency



Particles - Emission and movement

Particle emission and movement along the gradient

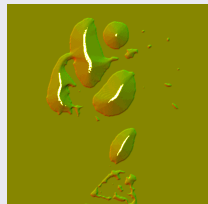
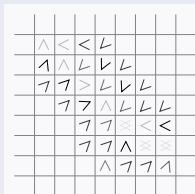


- **Emit particles** starting on voxels with $sphere > MinEmissionRadius$
- $MinEmissionRadius$ **avoids emitting particles in small vessels** (for registration unnecessary) and positive-false segmented areas like bones
- Particles **follow the local gradient vector**
- Length of gradient vector is **small at the center of vessels**
- Particle **stops** if the length of gradient vector is **below a specific value**



Particles - Emission and movement

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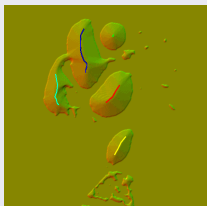
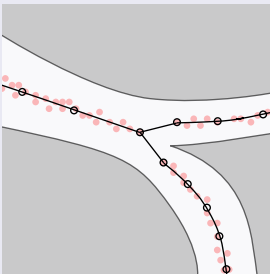
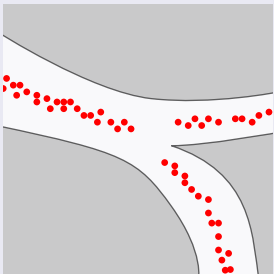


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Graph - Construction

Graph reconstruction

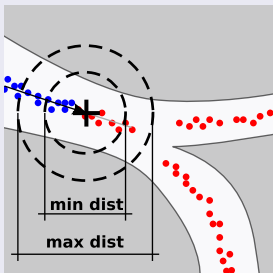


- Representing blood vessels by graphs
- Radius is also stored for each node for advanced registration
- **Reduces** matching of the **large particle amount** to matching of **sparse graph nodes**



Graph - Construction - Step 1 of 3

Connecting neighbored particles to strips

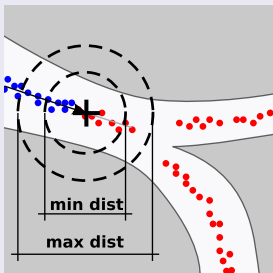


- Each particle has a **flag used** which is set if the particle is already **represented by an edge**
- Search for **neighbored particle within a specific range** $[min_dist, max_dist]$ where the used flag is not yet set
- Take particle which is **furthest away** as node *NextNode*
- **Set used flag** for all particles within the range *max_dist*
- Continue at the node *NextNode*



Graph - Construction - Step 1 of 3

Connecting neighbored particles to strips

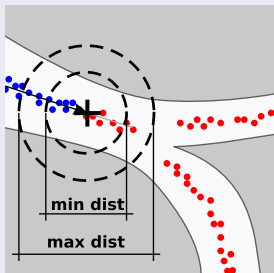


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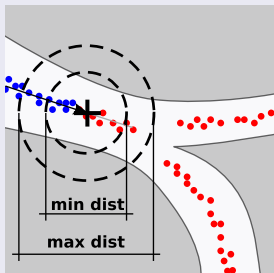


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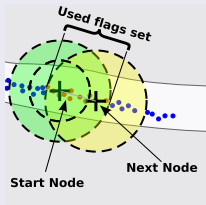


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Graph - Construction - Step 2 of 3

Handling of first strip node within non-forking area

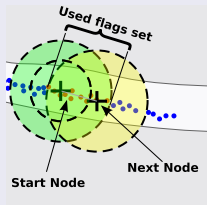


- **Start** each edge construction at unused particle with **maximum radius** \Rightarrow First node can have **2 neighbors**
- Setting all used flags of first node within the range *max_dist* **avoids creating edges in the opposite direction**
- \Rightarrow Set *used* flags only for particles which are **also in the range *max_dist* of NextNode**
- After creation of a stripline "in one direction", **restart again at first node** to extend stripline in opposite direction



Graph - Construction - Step 2 of 3

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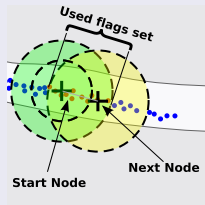


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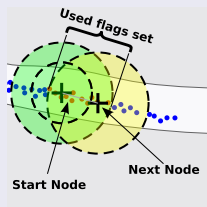


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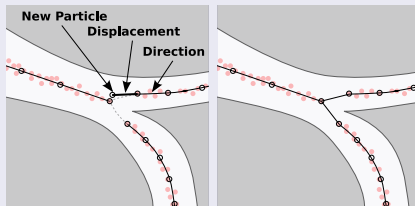


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Graph - Construction - Step 3 of 3

Connecting strips to graphs

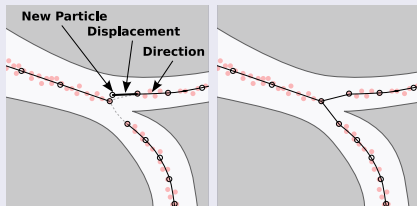


- The Gradient is very high on vessel **forkings** \Rightarrow strips are **disconnected**
- Use particle emission to create a connection at forkings
 - Particle is emitted with a **displacement** in the direction described by the **two corner nodes**
 - This particle **follows the gradient** until the **gradient value is below a specific value**
 - If there's a node N_n within the range max_dist : **connect corner node with node N_n**



Graph - Construction - Step 3 of 3

Connecting strips to graphs

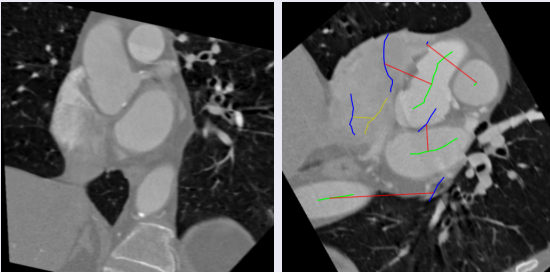


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Matching - Terminology

Matching graphs

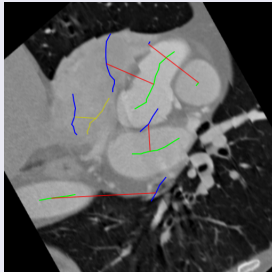
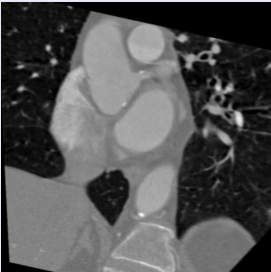


- Finding **transformation matrix M**
- Projecting **nodes P** from matching graph (applying matrix M for each point) results in **points P'** in "destination space"
- **Minimize distance** between nodes P' and nearest edge of destination graph



Matching - Terminology

Matching graphs



- Finding **transformation matrix M**
- Projecting **nodes P from matching graph** (applying matrix M for each point) results in **points P' in "destination space"**
- **Minimize distance** between **nodes P'** and **nearest edge** of destination graph

Matching

Transformation Matrix M projecting point P to P' :

$$M \cdot P = \begin{pmatrix} m_{1,1} & m_{1,2} & m_{1,3} & m_{1,4} \\ m_{2,1} & m_{2,2} & m_{2,3} & m_{2,4} \\ m_{3,1} & m_{3,2} & m_{3,3} & m_{3,4} \\ m_{4,1} & m_{4,2} & m_{4,3} & m_{4,4} \end{pmatrix} \cdot P = P'$$

Transformation matrix M for 3D CT data can be computed with **4 nodes** P^i, P^j, P^k, P^l of source graph and **4 points** O^i, O^j, O^k, O^l of destination graph (given in homogeneous form)

$$\begin{pmatrix} P_x^i & P_y^i & P_z^i & 1 \\ P_x^j & P_y^j & P_z^j & 1 \\ P_x^k & P_y^k & P_z^k & 1 \\ P_x^l & P_y^l & P_z^l & 1 \end{pmatrix} \begin{pmatrix} m_{1,1} \\ m_{1,2} \\ m_{1,3} \\ m_{1,4} \end{pmatrix} = \begin{pmatrix} O_x^i \\ O_x^j \\ O_x^k \\ O_x^l \end{pmatrix}$$

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Matching - Matrix decomposition

- Creating a matrix based on all possible node combinations can produce **unlikely mappings**
 - large shearings
 - large scalings / negative scalings
 - large rotations
 - large translations
- Even an unlikely mapping can produce **wrong matching with best computed matching**
- **Decompose matrix** to basic transformations and **restrict transformations**



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Matching - Matrix decomposition (cont')

- Matrix M is decomposed in
 - 3 Translation components t_x, t_y, t_z by matrix T
 - 3 Rotation matrices R_n around axis n : R_x, R_y, R_z
 - 3 Scaling components s_x, s_y, s_z with matrix S
 - 3 Shearing components $sh_{1,2,3}$ in shearing matrix H
- $M = H \cdot S \cdot R_z \cdot R_x \cdot R_y \cdot T$
- Decompositions "simulate" different **basic transformations**
 - First the translation is done to align both datasets
 - Secondly the translated dataset is rotated around y axis for better matching
 - ...
- Decomposition to basic transformations give the information for an **"early drop"** (omit the current matrix)



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Matching - Translation

- Translation Matrix $T = \begin{pmatrix} 1 & \cdot & \cdot & t_x \\ \cdot & 1 & \cdot & t_y \\ \cdot & \cdot & 1 & t_z \\ \cdot & \cdot & \cdot & 1 \end{pmatrix}$

- Translation decomposition:

$$M = M' \cdot T = \begin{pmatrix} m_{1,1} & m_{1,2} & m_{1,3} & 0 \\ m_{2,1} & m_{2,2} & m_{2,3} & 0 \\ m_{3,1} & m_{3,2} & m_{3,3} & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix} \cdot \begin{pmatrix} 1 & 0 & 0 & t_x \\ 0 & 1 & 0 & t_y \\ 0 & 0 & 1 & t_z \\ 0 & 0 & 0 & 1 \end{pmatrix}$$

(Search values for T to eliminate the rightmost column)

- Computation of the rightmost column values of M gives the implicit solution for T

$$\begin{pmatrix} m_{1,1} & m_{1,2} & m_{1,3} \\ m_{2,1} & m_{2,2} & m_{2,3} \\ m_{3,1} & m_{3,2} & m_{3,3} \end{pmatrix} \cdot \begin{pmatrix} t_1 \\ t_2 \\ t_3 \end{pmatrix} = \begin{pmatrix} m_{1,4} \\ m_{2,4} \\ m_{3,4} \end{pmatrix}$$



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Matching - Rotation

Rotation matrices (example for given y axis):

- Use rotation matrices (from QR decomposition) to set **values below diagonal to 0**

$$R_y = \begin{pmatrix} \cos(\alpha) & 0 & \sin(\alpha) & 0 \\ 0 & 1 & 0 & 0 \\ -\sin(\alpha) & 0 & \cos(\alpha) & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}$$

- Eliminating entry $m''_{3,1}$:

$$M' = M'' \cdot R_y \iff M' \cdot R_y^{-1} = M''$$

Assuming $m''_{3,1}$ should be set to 0:

$$0 = m'_{3,1} \cos(\alpha) - m'_{3,3} \sin(\alpha) \iff \alpha = \text{atan} \left(\frac{m'_{3,1}}{m'_{3,3}} \right)$$

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Matching - Scaling and Shearing

- Result **after translation and rotation** elimination is an **upper diagonal matrix** $M^{(3)}$
- Final decomposition returns the **shearing and scaling matrix**

$$M^{(3)} = H \cdot S \iff$$

$$\begin{pmatrix} m_{1,1}^{(3)} & m_{1,2}^{(3)} & m_{1,3}^{(3)} & \cdot \\ \cdot & m_{2,2}^{(3)} & m_{2,3}^{(3)} & \cdot \\ \cdot & \cdot & m_{3,3}^{(3)} & \cdot \\ \cdot & \cdot & \cdot & 1 \end{pmatrix} = \begin{pmatrix} 1 & h_1 & h_2 & \cdot \\ \cdot & 1 & h_3 & \cdot \\ \cdot & \cdot & 1 & \cdot \\ \cdot & \cdot & \cdot & 1 \end{pmatrix} \begin{pmatrix} s_x & \cdot & \cdot & \cdot \\ \cdot & s_y & \cdot & \cdot \\ \cdot & \cdot & s_z & \cdot \\ \cdot & \cdot & \cdot & 1 \end{pmatrix}$$

$$s_x = m_{1,1}^{(3)} \quad s_y = m_{2,2}^{(3)} \quad s_z = m_{3,3}^{(3)}$$

$$h_1 \cdot s_y = m_{1,2}^{(3)} \iff h_1 = \frac{m_{1,2}^{(3)}}{s_y}$$

$$h_2 \cdot s_z = m_{1,3}^{(3)} \iff h_2 = \frac{m_{1,3}^{(3)}}{s_z}$$

$$h_3 \cdot s_z = m_{2,3}^{(3)} \iff h_3 = \frac{m_{2,3}^{(3)}}{s_z}$$



Matching

- To work with minimal transformations of permitted matrices, **pretranslate the center of image** to $(0, 0, 0)$
- Empirical values to match $256 \times 256 \times 200$ CT scans:
 - maximum relative rotation angle: 90°
 - scale factor $\in [0.7; 1.3]$
 - shear factor $\in [-0.3; 0.3]$
 - maximum relative translation: 100
 - pretranslation matrix: $P = \begin{pmatrix} 1 & \cdot & \cdot & -128 \\ \cdot & 1 & \cdot & -128 \\ \cdot & \cdot & 1 & -100 \\ \cdot & \cdot & \cdot & 1 \end{pmatrix}$

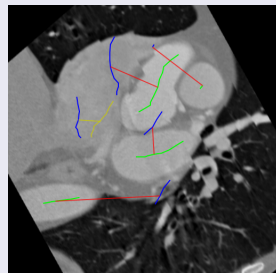
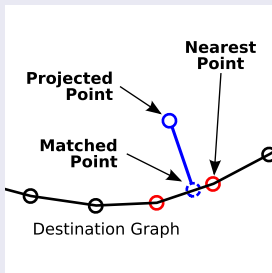


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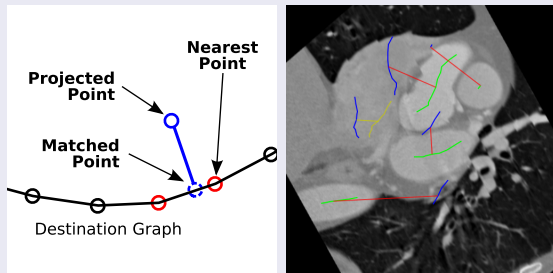
Matching - Computation of matching difference



- If the matrix is valid, **nodes P are projected to the destination graph space giving P'**
- For every projected node P' , **the nearest node P'' of the destination graph is searched**
- For both adjacent nodes of P'' the **smallest distance to the edge is taken**
- Best matching matrix M : Projection with the **smallest sum of distances for all nodes P' to the corresponding edges**



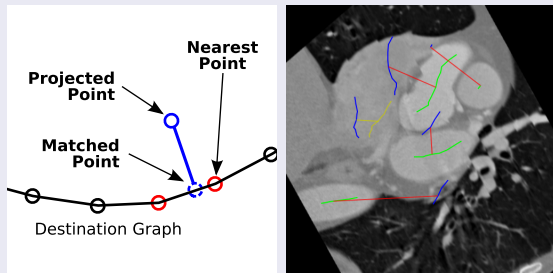
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Efficient implementation

- Direct implementation would take **more than half an hour** to match a 256x256x200 CT dataset to reference data
- **Sphere filter:**
 - Sphere filter has to be applied **for every point!**
- **Graph filter:**
 - Storing a few particles with a full grid would be a **waste of memory**
 - Computation time for nearest points with $O(r^d)$ increases with resolution n^d
- **Matching:**
 - Using naive approach for a graph with N nodes there would be $O\left(\left(\frac{N!}{(N-4)!}\right)^2\right)$ **matching possibilities** taking hours to compute



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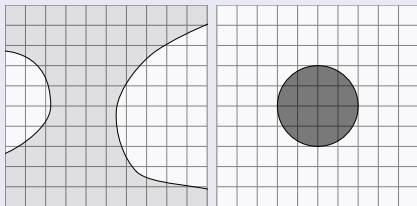


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Sphere filter using FFT

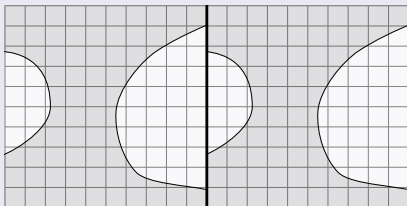


- Applying sphere filter is a **convolution for each sphere size**
- Convolution can be done very efficiently in **frequency space**, specially for large kernels



Sphere filter using FFT

FFT based on periodical data

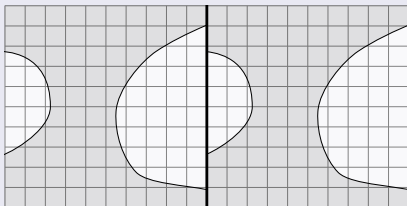


- **Boundary conditions:** Data fields have to be **padded with extra data** to work with existing FFT libraries (e.g. FFTW)
- Applying the **standard FFT** is based on a **periodical function**
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- **Large spheres at borders don't represent the local data**



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FFT based on periodical data

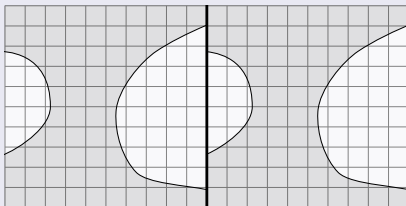


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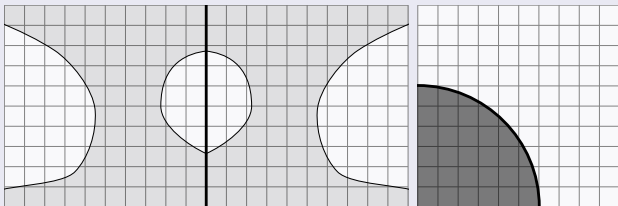


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FFT based on symmetric data



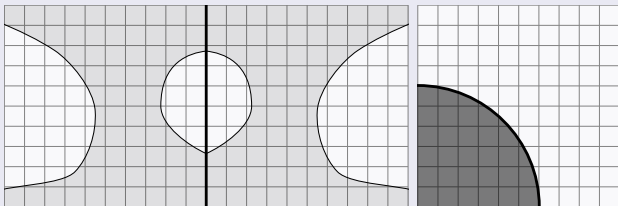
Left: Symmetric data - Right: Kernel

- Using **real data FFT** assuming domain is **symmetric on borders**
- Introduced errors just depend on the local data
- **Kernel** has to be **initialized for only $\frac{1}{8}$ of the data domain**
- Reduces the computation from > 20 minutes to a few seconds



Sphere filter using FFT

FFT based on symmetric data



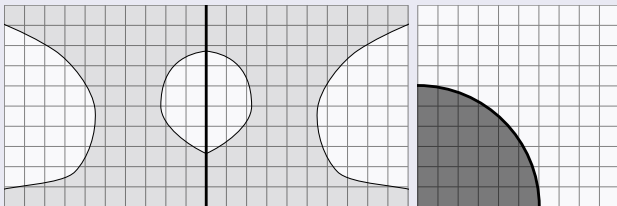
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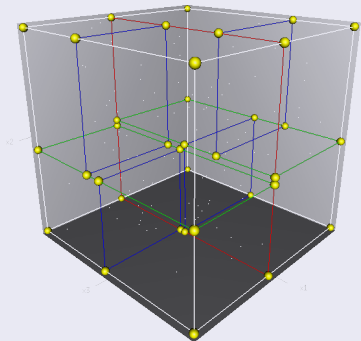
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Handling Particles with KD-Trees

KD-Tree



<http://en.wikipedia.org/wiki/Image:3dtree.png>

- KD Trees store arbitrary points using a tree like structure
- Efficient operations to **find points within a given radius**
- \Rightarrow Graph can be constructed within a second



Restricting Transformation Matrices

- $O\left(\left(\frac{N!}{(N-4)!}\right)^2\right)$ matching possibilities
- Discretization of points introduce **error** ϵ_1
- **Small anatomical differences** of points introduce **errors** ϵ_2
 - Error in matrix after construction becomes **less for far distant points**
 - Use only nodes with a **distance of at least δ** to create a **better conditioned problem**
- \Rightarrow **avoids the computation of the transformation matrix for many points**
- Omitting **impossible node combinations** and **nodes producing a bad conditioned problem** decreases computation time to a few seconds

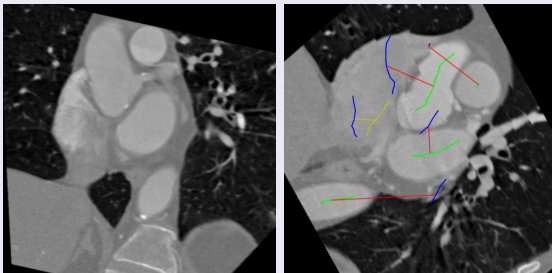


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Results



Matching 2D CT slices using only midpoints of edges as graph matching nodes

3D CT heart blood segmentation and registration:

- Can be handled in an **efficient and fast way**
- Takes just a **few seconds** on recent quad-core-systems
- Using translation matrices offers **registration of invisible areas** (yellow line in right image)



Possible improvements

- Use **sparser representation** of graph using **extrapolation or spline curves**
- Using **interpolation** with spherical filter (aliased kernel) for more accurate sphere radii
- Graph construction: include possible omitted nodes at strip endings
- **Randomized/hierarchical matching points** selection (maybe using hints of graph)
- Matching graphs
 - Using **heuristics from graphs** for matching
 - Comparing **edge slopes**
- Use **matching positive abort** if computed **overall distance is below a certain value** (assuming this is the correct matching)



Thank you for your attention



Any questions?



References and Links

- CT Datasets: Nuklearmed. Klinik der TU Muenchen, Germany
- FFTW: <http://www.fftw.org/>
- KD-Tree: <http://libkdtree.alioth.debian.org/>
- DICOM-Toolkit: <http://dicom.offis.de/dcmtoolkit.php.de>

