

Delineating objects in images via minimization of ℓ_p energies: Fuzzy Connectedness, Graph Cut, and Random Walk algorithms

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Outline

- 1 The problem of object delineation in a digital image: translating intuition to energy minimization setup
- 2 Object cost as a function of object boundary; ℓ_p cost
- 3 Delineation algorithms associated with ℓ_p energies
- 4 Comparison of GC and FC image segmentations
- 5 Spanning forests, Dijkstra algorithm, IRFC and PW objects

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Example 1, 2D, of object segmentation/delineation



An image of peppers



Delineation version 1

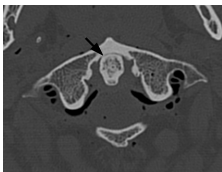


Delineation version 2

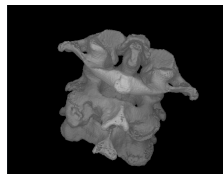


Delineation version 3

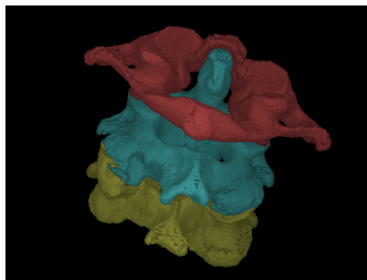
Example 2, 3D: a CT image of patient's cervical spine



A slice of an original 3D image

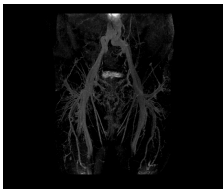


Surface rendition of segmented
three vertebrae, together

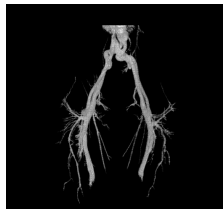


Color surface rendition of the segmented three vertebra

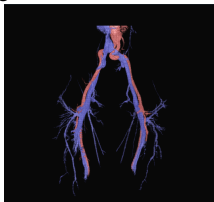
Example 3: An MR angiography image of the body region from belly to knee.



Rendition of an original 3D, contrast enhanced, image



A surface rendition of the entire vascular tree



Color surface rendition of segmented arterial (red) and venous (blue) trees

Image segmentation — formal setting

- An *(n -dimensional) image* is a map f from $C \subset \mathbb{R}^n$ into \mathbb{R}^k
The value $f(c)$ represents **image intensity at c** , a k -dimensional vector each component of which indicates a measure of some aspect of the signal, like color.
- *Segmentation problem*: Given an image $f: C \rightarrow \mathbb{R}^k$,
find a “**desired**” family $\mathcal{S}(f) = \{P_1, \dots, P_m\}$ of subsets of C .
- *Delineation problem* (on which we concentrate)
is when $m = 1$, i.e., when $\mathcal{S}(f) = P \subset C$.

Delineation of an image $f: C \rightarrow \mathbb{R}^k$ — formal setting

How to express that $\mathcal{S}(f) = P \subset C$ is **desired**?

There is no magic formulation that expresses all **desires**!

Several practical “solutions” exist. We use here the following:

- Fix seed sets:
 S indicating the foreground object P , (i.e., $S \subset P$), and
 T indicating the background (i.e., $T \cap P = \emptyset$).

This restricts the search space for $\mathcal{S}(f)$ to the family

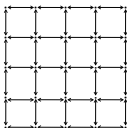
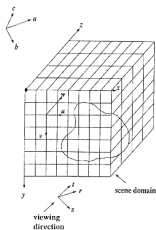
$$\mathcal{P}(S, T) = \{P \subset C \setminus T : S \subset P\}$$

- Define an **energy/cost function** $\varepsilon: \mathcal{P}(\emptyset, \emptyset) \rightarrow [0, \infty)$
- Declare $\mathcal{S}(f)$ to be **desired** when it minimizes ε on $\mathcal{P}(S, T)$.

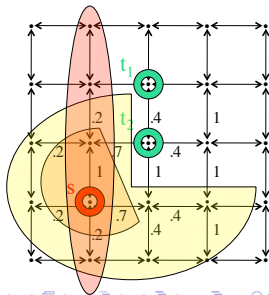
Digital vs “continuous” image $f: C \rightarrow \mathbb{R}^k$

The above set-up makes sense and was studied for the images with scene $C \subset \mathbb{R}^n$ being open bounded region.

We discuss only **digital images**, with finite rectangular scenes:



Example of sets in $\mathcal{P}(S, T)$
with $S = \{s\}$, $T = \{t_1, t_2\}$



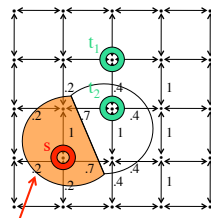
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Heuristic and the definition of boundary

Heuristic: The objects boundary areas should be identifiable in the image, as the areas of sharp image intensity change.

What constitutes **boundary** $\text{bd}(P)$ of P ?



Desired object

Need graph (or topological) structure $G = \langle V, E \rangle$ on C :

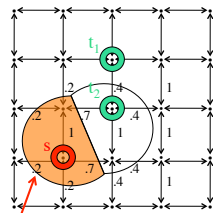
- Pixels $c \in C$ are its vertices, $V = C$;
- Edges $\{c, d\} \in E$ are “nearby” vertices (e.g. as in figure).

$\text{bd}(P)$ is the set of all edges $\{c, d\} \in E$ with $c \in P$ and $d \notin P$

Weighted graphs and ℓ_p cost functions, $1 \leq p \leq \infty$

Assume that with every edge $e = \{c, d\} \in E$ of an image f we have associated its **weight/cost** $w(e) \geq 0$, which is low, for big $\|f(c) - f(d)\|$.

Typically, $w(e) = e^{-\|f(c) - f(d)\|/\sigma^2}$, see fig.



Desired object

If $F_P: E \rightarrow [0, \infty)$, $F_P(e) = w(e)$ for $e \in \text{bd}(P)$ and $F_P(e) = 0$ for $e \notin \text{bd}(P)$, then **ℓ_p cost is defined** as

$$\varepsilon_p(P) \stackrel{\text{def}}{=} \|F_P\|_p = \begin{cases} \left(\sum_{e \in \text{bd}(P)} w(e)^p \right)^{1/p} & \text{if } p < \infty \\ \max_{e \in \text{bd}(P)} w(e) & \text{if } p = \infty. \end{cases}$$

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FC and GC algorithms as minimizers of ε_p

$$\varepsilon_p(P) \stackrel{\text{def}}{=} \|F_P\|_p = \begin{cases} \left(\sum_{e \in \text{bd}(P)} w(e)^p \right)^{1/p} & \text{if } p < \infty \\ \max_{e \in \text{bd}(P)} w(e) & \text{if } p = \infty. \end{cases}$$

$$p = 1: \varepsilon_1(P) = \sum_{e \in \text{bd}(P)} w(e);$$

Optimization solved by classic **min-cut/max-flow algorithm**.

Graph Cut, GC, delineation algorithm optimizes ε_1 .

$$p = \infty: \varepsilon_\infty(P) = \max_{e \in \text{bd}(P)} w(e);$$

Optimization solved by (versions of) **Dijkstra algorithm**.

ε_∞ optimized objects are returned by the algorithms:

Relative Fuzzy Connectedness, RFC, **Iterative RFC, IRFC**,
and **Power Watershed, PW** [C. Couprie *et al*, 2011].

$p = 2$: related to **Random Walker, RW**, algorithm [Grady, 2006],
see next slides.

Fuzzy sets

A map $x: C \rightarrow [0, 1]$ (i.e., $x \in [0, 1]^C$) can be considered as a **fuzzy set**, with $x(c)$ giving the degree of membership of c in it.

A hard set $P \subset C$ is identified with a fuzzy set (binary image) $\chi_P \in \{0, 1\}^C \subset [0, 1]^C$, $\chi_P(c) = 1$ iff $c \in P$.

For $x \in [0, 1]^C$ let $\hat{\varepsilon}_p(x) = \|F_x\|_p$, where $F_x: E \rightarrow [0, \infty)$,

$F_x(\{c, d\}) = |x(c) - x(d)|w(\{c, d\})$ for $\{c, d\} \in E$.

Then $\varepsilon_p(P) = \hat{\varepsilon}_p(\chi_P)$. We can minimize $\hat{\varepsilon}_p$ on

$\hat{\mathcal{P}}(S, T) = \{x: x(c) = 1 \text{ for } c \in S \text{ \& } x(c) = 0 \text{ for } c \in T\}$

instead of ε_p on $\mathcal{P}(S, T) = \hat{\mathcal{P}}(S, T) \cap \{0, 1\}^C$.

Random Walker, RW, algorithm

- RW finds (the unique) $\hat{\varepsilon}_2$ minimizer on $\hat{\mathcal{P}}(S, T)$.
- Defines its output as $P = \{c: x(c) \geq .5\}$.

Problems with RW:

- 1 Output need not be connected (even when S and T are).
- 2 P need not minimize ε_2 on $\mathcal{P}(S, T)$.

Neither of this happens for ε_1 (i.e. GC) or ε_∞ (i.e. RFC or PW):

Thm: For $p \in \{1, \infty\}$, any minimizer of $\hat{\varepsilon}_p$ on $\hat{\mathcal{P}}(S, T)$ actually belongs to $\mathcal{P}(S, T)$.

(Non)-uniqueness of the minimizers for ε_1 and ε_∞

Let $\mathcal{P}_p(S, T) = \{P \in \mathcal{P}(S, T) : P \text{ minimizes } \varepsilon_p \text{ on } \mathcal{P}(S, T)\}$.

Both $\mathcal{P}_1(S, T)$ and $\mathcal{P}_\infty(S, T)$ may have more than one element.

However, the **outputs** of the standard versions **of the algorithms**:

- **GC**, from $\mathcal{P}_1(S, T)$,
- **RFC**, from $\mathcal{P}_\infty(S, T)$, and
- **IRFC**, from $\mathcal{P}_\infty(S, T)$

are unique in the sense of the next theorem.

GC & FC segmentations — comparison theorem 1

Theorem (Argument minimality)

For $p \in \{1, \infty\}$, $\mathcal{P}_\varepsilon(S, T)$ contains the \subset -smallest object.

- **GC** algorithm returns the smallest set in $\mathcal{P}_1(S, T)$.
- **RFC** algorithm returns the smallest set in $\mathcal{P}_\infty(S, T)$.
- **IRFC** algorithm returns the smallest set in a refinement $\mathcal{P}_\infty^*(S, T)$ of $\mathcal{P}_\infty(S, T)$.

Moreover, if n is the size of the image (scene), then

- GC runs in time of order $O(n^3)$ (the best known algorithm) or $O(n^{2.5})$ (the fastest currently known algorithm)
- Both RFC and IRFC run in time of order $O(n)$ (for standard medical images — the intensity range size not too big) or $O(n \ln n)$ (the worst case scenario)

GC & FC — asymptotic equivalence

Theorem (Asymptotic equivalence of GC and FC)

Let $\mathcal{P}_p^m(S, T)$ be the family $\mathcal{P}_p(S, T)$ for the edge weight function w replaced by its m -th power w^m . Then

- $\mathcal{P}_\infty^m(S, T) = \mathcal{P}_\infty(S, T)$ and similarly for IRFC algorithm.
So, the outputs of RFC and IRFC are unchanged by m .
- $\mathcal{P}_1^m(S, T) \subseteq \mathcal{P}_\infty(S, T)$ for m large enough.

In particular, if $\mathcal{P}_\infty(S, T)$ has only one element, then

the output of GC coincides with the outputs of RFC and IRFC for m large enough.

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Advantages of FC over GC — theoretical angle

Speed: FC algorithms run a lot faster than GC algorithms:
 $O(n)$ (or $O(n \ln n)$) versus $O(n^3)$ (or $O(n^{2.5})$).

Robustness: RFC & IRFC are unaffected by small seed changes.
GC is sensitive for even small seed changes.

Shrinking: GC chooses objects with small size boundary
(often with edges with high weights);
No such problem for RFC & IRFC

Multiple objects: FC framework handles easily the segmentation of
multiple objects, same running time and robustness.
GC in such setting leads to NP-hard problem,
so (for precise delineation) it runs in exponential time

Iterative approach: RFC has an iterative approach refinement;
No such refinement methods exist for GC at present.

Advantages of GC over FC

Boundary smoothness: GC chooses small boundary, so it naturally smooths it; in many (but not all) medically important delineations, this is a desirable feature.

Basic FC framework has no boundary smoothing; if desirable, smoothing requires post processing

Combining image homogeneity info with known object intensity:

GC naturally combines information on image homogeneity (binary relation on voxels) with information on expected object intensity (unary relation on voxels);

Combining such informations is difficult to achieve in the FC framework.

Setup of experiments:

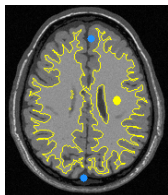
- In each experiment we used 20 MR BrainWeb phantom images (simulated T1 acquisition); graphs show averages.
- Sets of seeds were generated, from known true binary segmentations, by applying erosion operation: the bigger erosion radius, the smaller the seed sets.
- The weight map $w(c, d)$, same for FC and GC, was defined from the image intensity function f as $w(c, d) = -|G(f(c)) - G(f(d))|$, where G is an appropriate Gaussian.

Setup of experiments:

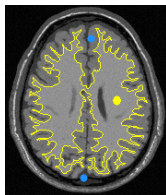
Data parameters: the simulated T1 acquisition were as follows: spoiled FLASH sequence with TR=22ms and TE=9.2ms, flip angle = 30° , voxel size = $1 \times 1 \times 1 \text{ mm}^3$, noise = 3%, and background non-uniformity = 20%.

Computer: Experiments were run on PC with an AMD Athlon 64 X2 Dual-Core Processor TK-57, 1.9 GHz, 2×256 KB L2 cache, and 2 GB DDR2 of RAM.

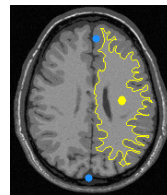
Robustness & shrinking for FC & GC: White Matter



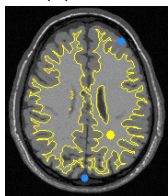
(a) RFC



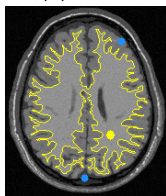
(b) IRFC



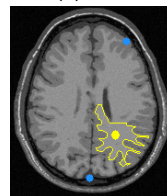
(c) GC



(d) RFC



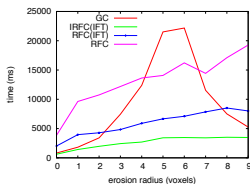
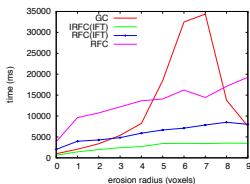
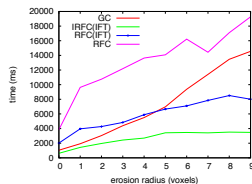
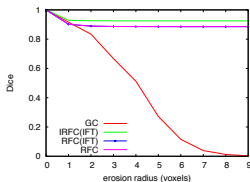
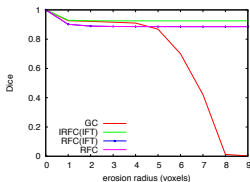
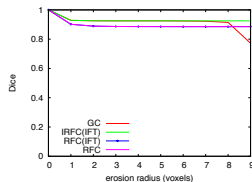
(e) IRFC



(f) GC

Figure: (a)&(d) and (b)&(e): same outputs for different seeds; (c)&(f) GC: dramatic change of output; seeds choice same as in the FC case

Time & accuracy of FC & GC: segmentation of WM

(a) Time for w^1 (b) Time for w^5 (c) Time for w^{30} (d) Accuracy for w^1 (e) Accuracy for w^5 (f) Accuracy for w^{30}

FC vs GC: Conclusions

- FC and GC quite similar,
yet FC has many advantages over GC:
 - FC runs considerably faster than GC
 - FC is robust (seed), while GC has shrinkage problem
 - FC, unlike GC, easily handles multiple-object segmentation
- unless the application requires, in an essential way, the **simultaneous** use of
 - homogeneity (binary) info on image intensity;
 - expected object intensity (unary) info on image intensity;

it makes sense to use FC (more precisely IRFC)
segmentation algorithm, rather than GC algorithm

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Forests: the powerhouse behind Dijkstra algorithm

Fix weighted graph $G = \langle C, E, w \rangle$ and $\emptyset \neq W \subset C$.

Definition (Spanning Forest w.r.t. W)

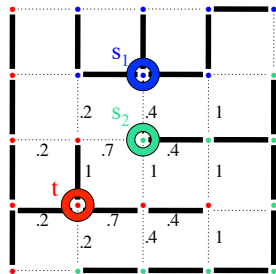
A *forest* for G is any subgraph $\mathbb{F} = \langle C, E' \rangle$ of G free of cycles.
 $\mathbb{F} = \langle C, E' \rangle$ is *spanning with respect to W* when any connected component of \mathbb{F} contains precisely one element of W .

Example of a spanning

forest w.r.t. $W = \{s_1, s_2, t\}$

Each component

marked by different color



Forest-generated (IRFC and PW) objects

$G = \langle C, E, w \rangle$ – weighted graph, $\emptyset \neq W \subset C$, $S \subset W$

Definition (Forest-generated object)

For a spanning forest \mathbb{F} w.r.t. W and $S \subset W$,

$P(S, \mathbb{F})$ is a union of all components of \mathbb{F} intersecting S .

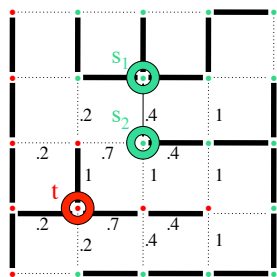
Note that $P(S, \mathbb{F}) \in \mathcal{P}(S, T)$ for $T = W \setminus S$.

Example (green vertices) of

$P(S, \mathbb{F})$ with $S = \{s_1, s_2\}$.

Outputs of the algorithms we will discuss, GC_{sum} and PW,

are in the $P(S, \mathbb{F})$ format.



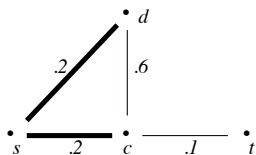
Optimal Path Forest, OPF

Definition (Optimal Path Forest, OPF)

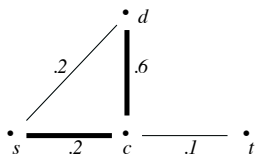
For a path $p = \langle c_1, \dots, c_k \rangle$ in G let $\mu(p) = \min_{i < k} W(\{c_i, c_{i+1}\})$,
 the *weakest link* of p .

A forest \mathbb{F} w.r.t. W is *path-optimal* provided for every $c \in C$,
 the unique path p_c in \mathbb{F} from W to c is μ -optimal in G , i.e.,
 $\mu(p_c) \geq \mu(p)$ for any path p in G from W to c .

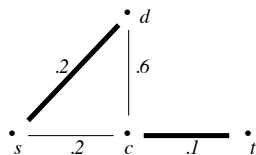
For OPF \mathbb{F} w.r.t. W , $\mu(p_c) = \mu^C(c, W)$ for every $c \in C$
 (with μ^C in the **Fuzzy Connectedness** sense)



(g) OPF, $W = \{s, t\}$



(h) another OPF



(i) not OPF

GC_{max} algorithm and IRFC

Theorem ([KC *et al.*] OPF object minimizing ε_∞)

There exists the smallest $P_{\min} \in \mathcal{P}(S, T)$ in form $P(S, \mathbb{F})$, where \mathbb{F} is an OPF w.r.t. $S \cup T$.

\mathbb{F} is found by GC_{max}, a version of Dijkstra's shortest path algorithm, in a linear time w.r.t. $|C| + M$,
where M is the size of the range of w .

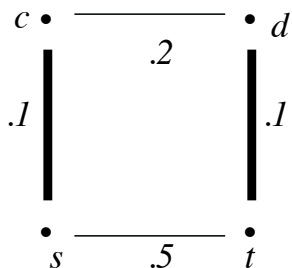
In practice, $O(|C| + M) = O(|C|)$.

The object P_{\min} , returned by GC_{max}, coincides with the Iterative Relative Fuzzy Connectedness, IRFC, object.

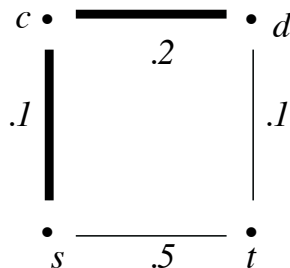
Maximal Spanning Forest, MSF

Definition (Maximal Spanning Forest, MSF)

A forest $\mathbb{F} = \langle C, E' \rangle$ w.r.t. W is *maximal spanning* provided $\sum_{e \in E'} w(e) \geq \sum_{e \in \hat{E}'} w(e)$ for every forest $\hat{\mathbb{F}} = \langle C, \hat{E}' \rangle$ w.r.t. W



(j) OPF w.r.t. $\{s, t\}$, not MSF



(k) MSF and OPF

Theorem ([Audigier & Lotufo], [Cousty et al.])

Every MSF is OPF, but not the other way around.

MSF and Power Watershed, PW, algorithm

Theorem ([C. Couprie *et al.*] PW output as MSF)

PW algorithm returns $P(S, \mathbb{F})$ for a MSF \mathbb{F} w.r.t. $S \cup T$.

\mathbb{F} is found by PW via a complicated version of Kruskal's algorithm and, locally, Random Walker algorithm.

Since

- IRFC object is indicated by OPF,
- PW object is indicated by MSF, and
- every MSF is OPF

What is the relation between IRFC and PW objects?

New results on GC_{\max} , MSF, and OPF

Theorem ([KC *et al.*] MSF vs OPF)

If P_{\min} is the output of GC_{\max} (the smallest $P(S, \mathbb{F})$, with \mathbb{F} being OPF w.r.t. $S \cup T$), then $P_{\min} = P(S, \hat{\mathbb{F}})$ for some MSF $\hat{\mathbb{F}}$.

If \mathbb{F} is a MSF w.r.t. $S \cup T$, then $P(S, \mathbb{F})$ minimizes energy ε_{∞} (in $\mathcal{P}(S, T)$).

$P(S, \mathbb{F})$, with \mathbb{F} being OPF w.r.t. $S \cup T$, need not minimize ε_{∞} .

In other words

$$P_{\min} \in \mathcal{P}_{MSF}(S, T) \subset \mathcal{P}_{OPF}(S, T) \cap \mathcal{P}_{\varepsilon_{\infty}}(S, T),$$

where $\mathcal{P}_{MSF}(S, T) = \{P(S, \mathbb{F}) : \mathbb{F} \text{ is MSF}\}$, similarly for OPF, and $\mathcal{P}_{\varepsilon_{\infty}}(S, T)$ is the set of all ε_{∞} -optimizers.

Thank you for your attention!