Workshop on image processing, computer vision and graphics Graph cut optimization class:35 36, 02.07.2010, time: 10am to 1.15 pm

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Outline

- Motivation of selecting MAP-MRF framework
- MAP-MRF approach
- Mapping of MAP-MRF to graph-cut problem
- What Energy functions can be minimized with Graph-cuts?
- Some examples of MAP-MRF framework using graph-cut
- ▶ MAP-MRF framework for Super Resolution
- Summary

Motivation of selecting MAP-MRF framework

- Bayesian framework suitable for problems in Computer Vision
- MAP-MRF with Gibbs gives easy implementation and formulation.
- Problems: High computational cost or Standard methods used are very slow.
- Boykov et.al proposed methods to solve MAP-MRF using graph-cut algorithms -MAP-MRF estimation is equivalent to min-cut problem on a graph
- Applied to many vision problems

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MAP-MRF framework

- MRF framework: Given set of pixels $S = \{s_1 \dots s_m\}$ and set of labels $\Lambda = \{l_1 \dots l_L\}$ and neighborhood system N, Find mapping of S to Λ .
- Let F be the configuration for labels $F = \{f_i \dots f_N\}, f_i \in \Lambda$ is the label for s_i
- ▶ *F* is MRF with respect to *N* iff
 - Positivity: $P(F = f) \ge 0 \ \forall \ f \in F$
 - Markovianity:

$$P(F_s = f_s/F_r = f_r, \forall r \neq s) = P(F_s = f_s/F_r = f_r, \forall r \in N_s)$$

Easy Implementability

MAP-MRF framework contd...

Let *G* be the observed image

- $G = \phi(H(F)) + N$ where H =Camera Transfer Function and $\phi =$ Recorder distortion H is assumed to be LSI and ϕ is invertible nonlinear function and N is additive noise assumed to be iid
- In the framework of Restoration : Given G What is F?
 - P(F = f/G = g), Maximum likelihood of F = f given G = g
 - From Bays rule $P(F=f/G=g) \propto P(G=g/F=f)P(F=f)$ where P(G=g/F=f)=Data model, P(F=f)=Prior and P(F=f/G=g)=Aposteriori distribution
 - Need to maximize aposteriori(MAP) distribution

Gibbs Distribution

Geman and Geman proved equivalence between MRF and Gibbs distribution

$$P(f) = \frac{1}{Z}exp(-U(f)/T)$$

where $U(f)=\sum\limits_{c\in N}V_c(f)$ = Energy function, V_c =Clique potential, $Z=\sum\limits_f exp(-U(f)/T)$ =Partition function and T=Temperature

Hammersely Clifford Theorem:

F is MRF on S with respect to N if and only if F is Gibbs random field on S with respect to N

Relates the conditional distribution(local characteristics) and joint distribution(Gibbs measure)
Workshop on CVG

MAP-MRF

- $\oint \hat{f} = \underset{f}{\operatorname{argmax}} P(f/g)$
- From Bays Rule

$$\hat{f} = \underset{f}{\operatorname{argmax}} P(g/f)\dot{P}(f)$$

I-term: Likelihood function and II-term: Prior Model

- $\hat{f} = \underset{f}{\operatorname{argmax}} \ exp \big\{ \underset{p}{\Sigma} \in (p(g/f_p)) \underset{p,q \in N}{\sum} V_{p,q}(f_p,f_q) \big\}$
- ▶ MAP estimate of *f* given *g* is equivalent to minimizing energy function with prior and data model

$$\hat{f} = \underset{f}{\operatorname{argmin}} \left\{ - \sum_{p} ln(p(g/f_p)) + \sum_{p,q \in N} V_{p,q}(f_p, f_q) \right\}$$

MAP-MRF

The Energy Function has data term and regularization term

$$U(f) = \left\{ \sum_{i} (g_i - \phi(H(f_i)) + \sum_{i,j \in N} V_{i,j}(f_i, f_j) \right\}$$

- Different ways of defining Clique potential which defines the regularization term or smoothness term in the energy function and describes the prior probability of a particular realization of the elements of the clique.
- Data model should capture the cost of assigning the label
- MAP-MRF is usually solved using SA which is very slow but guarantees the global minima for any arbitrary energy function
- Boykov et.al suggested max-flow/min-cut graph algorithms to solve some class of energy functions with MAP-MRF framework within a known factor of global minimum

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Mapping of MAP-MRF to graph-cut

- Vision problems as image labeling:Depth(stereo), Object Index(Segmentation), Original Intensity(Restoration)
- Labeling problem can be cast in terms of energy minimization
 - Labeling of pixels
 - Penalty for pixel labeling
 - Interaction between neighboring pixels:Smoothness term
- All pixels and labels are considered as vertices, edge and edge weights are calculated dynamically
- Min-cut on G has unique binary segmentation
- Segmentation associated with min-cut that satisfies user defined constraints minimizes the energy function

Energy Minimization

- Global minimum can be found in polynomial time if the energy function
 - is convex,
 - or has only two labels, eg. Icing model.
- Discontinuity- preserving energy function is not convex, eg. Potts model. Thus global minimization is NP- hard, takes exponential time,
- Thus global minimization Approximation algorithm to find local minimum
 - EM, Belief Propagation, Graph-Cuts
- What is Graph-Cuts?
 - Minimize an energy function with non binary variables by repeatedly minimizing an energy function with binary variables using Max- flow/ min- cut method

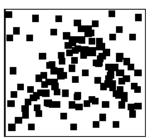
s-t graph-cuts for Binary Energy Minimization

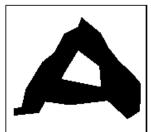
- Posterior energy (MRF)
- Complete characterization of binary energies that can be minimized with s-t graph cuts.

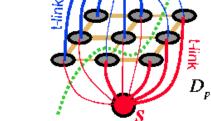
$$U(f) = \sum_{p} (D_p(f_p)) + \sum_{pq \in N} V(f_p, f_q)$$

 $lackbox{}{lackbox{}{lackbox{}{}}} U(f)$ can be minimized by graph-cuts

$$\iff V(s,s) + V(t,t) \leqslant V(s,t) + V(t,s)$$



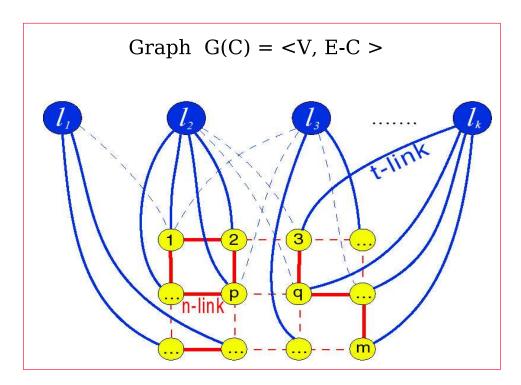




a cut

s-t graph-cuts for multi label problems

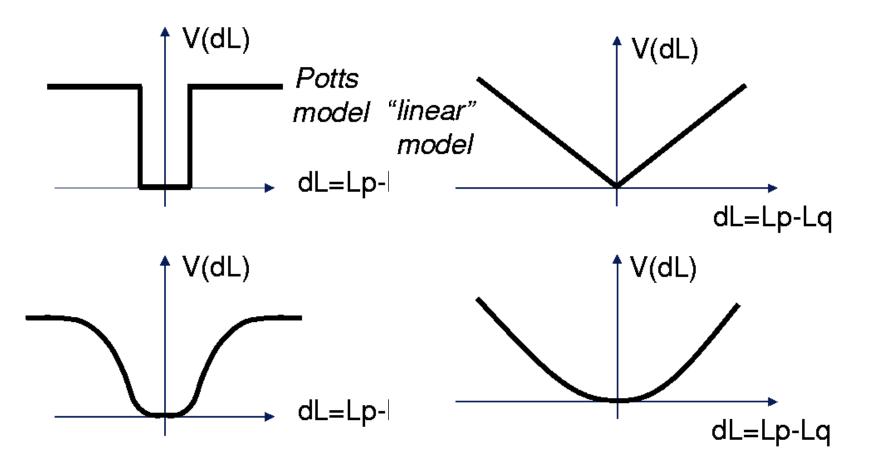
- Class of Energy that can be minimized exactly: Energies with convex interactions
 - excludes robust discontinuity-preserving interactions
- Guaranteed quality approximation algorithms for multi-label energies with discontinuity-preserving interactions like Potts model of interactions and Metric interactions



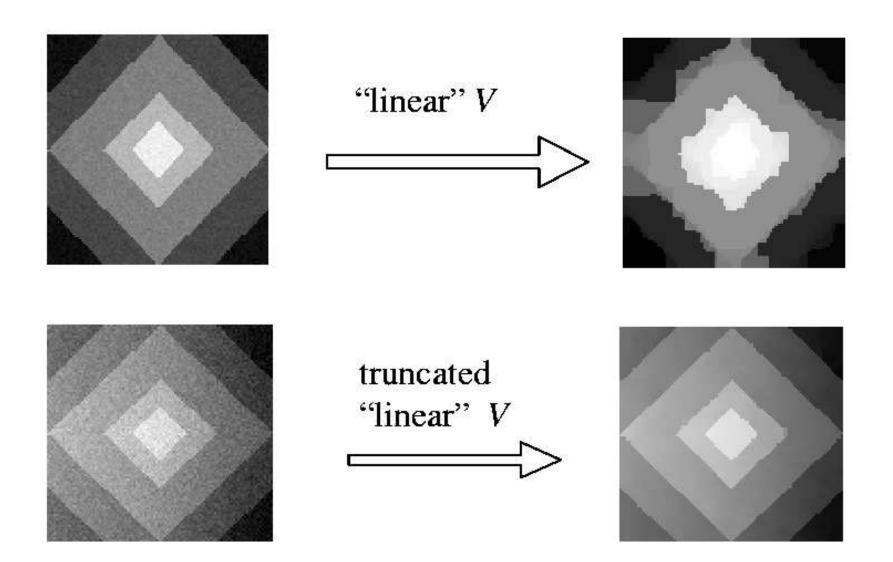
Different types of Pixel Interactions

Discontinuity preserving interactions:

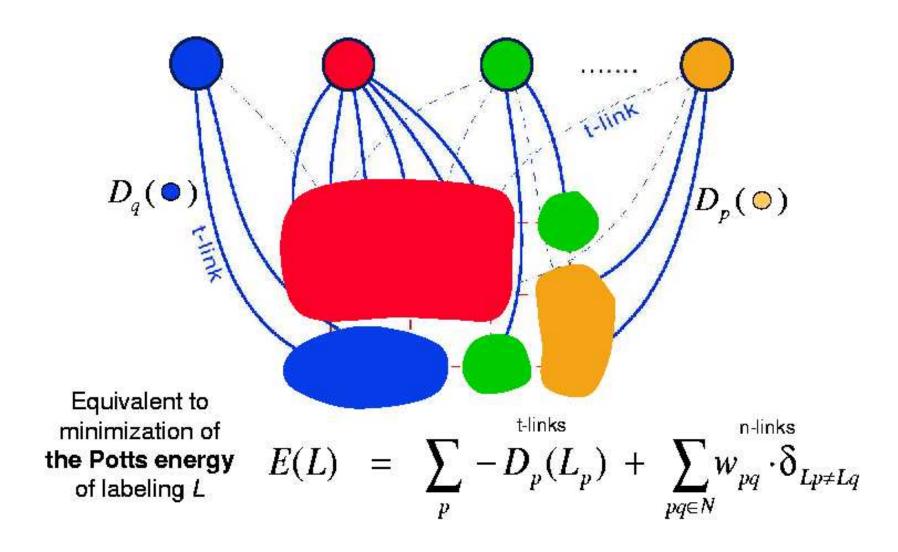
Convex interactions: Linear Models



Convex vs. Discontinuity-preserving



Multi way Graph-cut



Multi way Graph-cut algorithms by Boykov et.al

- Equivalent to Potts energy minimization
- ▶ NP-hard problem (3 or more labels)
 - two labels can be solved via s-t cuts (Greig et. al., 1989)
- Two approximation algorithms (Boykov et.al 1998,2001) Basic Idea:break multi-way cut computation into a sequence of binary s-t cuts.
 - α- Expansion
 Each label competes with the other labels for space in the image
 - $\alpha-\beta$ Swap : Define a move which allows to change pixels from α to β and β to α

α-Expansion approximation algorithm

Guaranteed quality approximation

- within a factor of 2 from Global minimum (Potts Model)
- applies to a wide class of energies with robust interactions
- ▶ Potts model (BVZ 1989), Metric interactions (BVZ 2001), Sub modular interactions (KZ 2002,2004)

Algorithm

- 1. Start with any arbitrary labeling f
- 2. Set success = 0
- 3. For each label $\alpha \in L$ (random order)
 - (a) find $\hat{f} = argmin \ U(f^1)$ among f^1 within one α -expansion f
 - (b) If $U(\hat{f}) < U(f)$, set $f = \hat{f}$ and success = 1
- 4. If success = 1 go to step 2
- 5. return f

$\alpha - \beta$ Swap approximation algorithm

Handles more general energy functions

Experimentally proved results

Algorithm

- 1. Start with any arbitrary labeling f
- 2. Set success = 0
- 3. For each pair of labels $\{\alpha, \beta\} \in L$ (random order)
 - (a) find $\hat{f} = argmin \ U(f^1)$ among f^1 within one $\alpha \beta$ swap of f
 - (b) $IfU(\hat{f}) < U(f)$, set $f = \hat{f}$ and success = 1
- 4. If success = 1 go to step 2
- 5. return f

Finding optimal expansion move

3.a step in algo. The structure of the graph is dynamically determined by the current position P and label α .

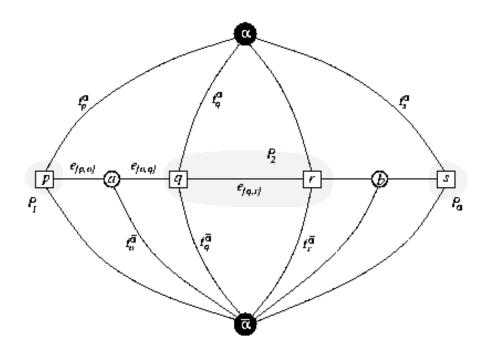


Figure 6: An example of \mathcal{G}_{α} for a 1D image. The set of pixels in the image is $\mathcal{P} = \{p, q, r, s\}$ and the current partition is $\mathbf{P} = \{\mathcal{P}_1, \mathcal{P}_2, \mathcal{P}_{\alpha}\}$ where $\mathcal{P}_1 = \{p\}$, $\mathcal{P}_2 = \{q, r\}$, and $\mathcal{P}_{\alpha} = \{s\}$. Two auxiliary nodes $a = a_{\{p,q\}}$, $b = a_{\{r,s\}}$ are introduced between neighboring pixels separated in the current partition. Auxiliary nodes are added at the boundary of sets $\mathcal{P}_{l \text{ ind IP}-p. 21/40}$

Expansion move-assignment of weights

edge	weight	for	
$t_p^{ar{lpha}}$	∞	$p \in \mathcal{P}_{\alpha}$	
$t_p^{ar{lpha}}$	$D_p(f_p)$	$p ot\in \mathcal{P}_{lpha}$	
t_p^{lpha}	$D_p(\alpha)$	$p \in \mathcal{P}$	
$e_{\{p,a\}}$	$V(f_p, \alpha)$		
$e_{\{a,q\}}$	$V(\alpha, f_q)$	$\{p,q\}\in\mathcal{N},\;f_p eq f_q$	
$t_a^{ar{lpha}}$	$V(f_p, f_q)$		
$e_{\{p,q\}}$	$V(f_p, lpha)$	$\{p,q\}\in\mathcal{N},\;f_p=f_q$	

Finding optimal swap move

3.1 step in algo. The structure of the graph is dynamically determined by the current position P and labels α, β .

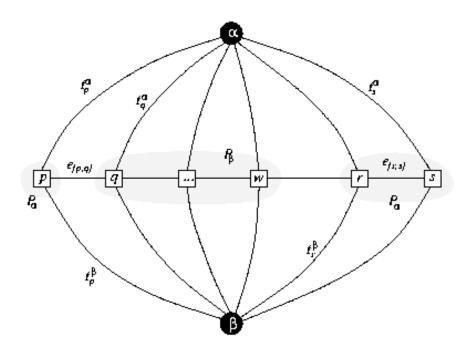


Figure 4: An example of the graph $\mathcal{G}_{\alpha\beta}$ for a 1D image. The set of pixels in the image is $\mathcal{P}_{\alpha\beta} = \mathcal{P}_{\alpha} \cup \mathcal{P}_{\beta}$ where $\mathcal{P}_{\alpha} = \{p, r, s\}$ and $\mathcal{P}_{\beta} = \{q, \ldots, w\}$.

Optimal swap move-assignment of weights

edge	weight	for
t_p^{lpha}	$D_p(lpha) + \sum_{\substack{q \in \mathcal{N}_p \ q ot\in \mathcal{P}_{lphaeta}}} V(lpha, f_q)$	$p \in \mathcal{P}_{\alpha\beta}$
t_p^eta	$D_p(eta) + \sum_{\substack{q \in \mathcal{N}_p \ q ot\in \mathcal{P}_{m{lpha}eta}}} V(eta, f_q)$	$p \in \mathcal{P}_{\alpha\beta}$
$e_{\{p,q\}}$	V(lpha,eta)	$\{p,q\}{\in}\mathcal{N} \ p,q{\in}\mathcal{P}_{m{lpha}m{eta}}$

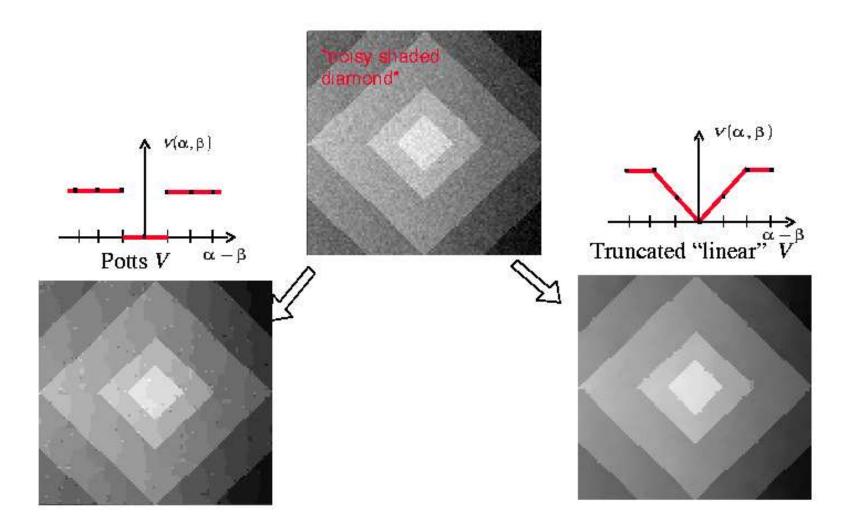
α-Expansion move



initial solution

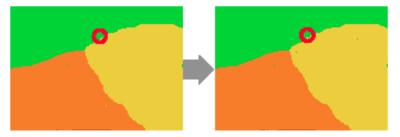
- -expansion

Example for Metric Interactions



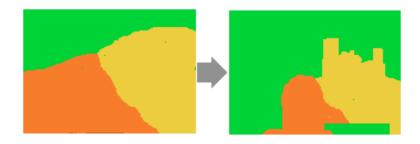
Comparison

single "one-pixel" move (simulated annealing, ICM,...)



- Only one pixel can change its label at a time
- Finding an optimal move is computationally trivial

single a-expansion move



- Large number of pixels can change their labels simultaneously
- Finding an optimal move is computationally intensive O(2ⁿ) (s-t cuts)

Comparisons contd..

simulated annealing

- Finds local minimum of energy with respect to small one-pixel moves
- Initialization is important
- solution could be arbitrarily far from the global minima
- May not know when to stop. Practical complexity may be worse than exhaustive search
- Can be applied to anything

α-Expansion

- Finds local minimum of energy with respect to very strong moves
- In practice, results do not depend on initialization
- solution is within the factor of 2 from the global minima
- In practice, one cycle through all labels gives sufficiently good results
- Applies to a restricted class of energies

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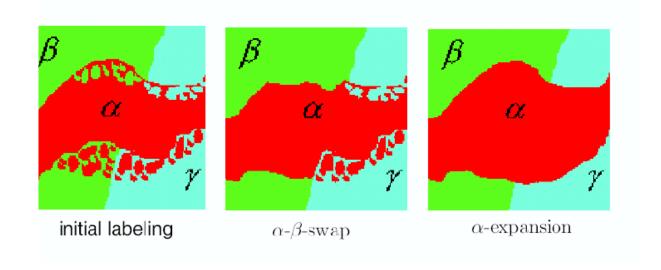
What Energy functions can be minimized with Graph-cuts?

- α-expansion algorithm can be applied to pairwise interactions that are metric on the space of labels
 - V(a,a) = 0
 - $V(a,b) \ge 0$
 - $V(a,b) \le V(a,c) + V(c,b)$
- Any truncated metric is also a metric(includes robust interactions)
- α-expansion algorithm further generalizes to submodular pair-wise interactions
- $V(c,c) + V(a.b) \le V(a,c) + V(c,b)$
- $\alpha \beta$ swap can be applied to pairwise interactions which are semi-metric on the space of labels
- Let *E* be a function of binary variables. If *E* is not regular, then *E* is not graph-representable.

Regular functions

- All functions of one variable are regular
- A function V of two variables is called regular if $V(0,0)+V(1,1) \leq V(0,1)+V(1,0)$
- A function V of more than two variables is called regular if all projections of V of two variables are regular.
- Let V be a function of n binary variables from F^3 , ie. $V(x_1, \ldots x_n) = \sum_i V^i(x_i) + \sum_{i < j} V^{i,j}(x_i, x_j) + \sum_{i < j < k} V^{i,j,k}(x_i, x_j, x_k).$ Then, V is graph-representable if and only if V is regular
- Any projection of a graph-representable function is graph-representable.

Moves



If V is Metric, then each expansion move is regular

$$E(0,0)+E(1,1)=V(\beta,\gamma)+V(\alpha,\alpha) \leq V(\beta,\alpha)+V(\alpha,\gamma)=E(0,1)+E(1,0)$$

If V is Semi-metric, then each swap move is regular

$$E(0,0) + E(1,1) = V(\beta,\beta) + V(\alpha,\alpha) \le V(\beta,\alpha) + V(\alpha,\beta) = E(0,1) + E(1,0)$$

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Some examples of MAP-MRF using graph-cut

- Image segmentation 1: Jiangjian Xiao and Mubarak Shah CVPR 2005
 - Motion cue to segment using graph-cut
 - refine the segmentation by alpha matte
 - Example 1
- Image segmentation/Object Extraction: Yuri Boykov and Vladimir Kolmogorov 2004
 - Combine both active contours and graph-cuts
 - Reduces the metrification error
 - Example 2

Examples contd...

- Texture Synthesis :Image quilting by Efros and Freeman, 2001Example
- Video Texture :3D generalization of Image quilting by Kwatra, Schodl, Essa, Bobick 2003 Process Source Synthesized
- Stereo: Boykov et.al 98, 2002 Example
- Multiview reconstruction by Boykov et.al 2004 Example
- Interactive Digital photomontage by microsoft research lab Example

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MAP-MRF framework for Super Resolution

- ▶ The Energy function is given by
- $E(f) = \sum_{p} D_p(f_p) + \sum_{p,q \in N_p} V_{p,q}(f_p, f_q) + \sum_{(p,q) \downarrow d} V_{(p,q) \downarrow d}(g_{p \downarrow d}, g_{q \downarrow d})$

where D = g - DHf is a data model term and next two terms are regularization terms. g=observed image, D=decimation function and H= Camera transfer function.

- f is the Super Resolution Image needs to estimate
- Regularization terms are truncated linear models
- $V_{p,q}(f_p, f_q) = min(K, |f_p f_q|)$
- $V_{(p,q)_{\downarrow d}}(g_{p\downarrow d},g_{q\downarrow d}) = \min(K,|g_{p\downarrow d}-g_{q\downarrow d}|)$
- Once we have the MAP-MRF framework for SR, we can apply Graph-cuts to estimate f.

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- What energy functions can be minimized with Graph-cuts
- lacktriangle α —Expansion and α β Swap algorithms in Graph-cuts
- Examples using Graph-cuts
- Framework of SR

THANK YOU