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Variational Methods

Continuous convex models and optimization



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1 Introduction to variational methods

- **2** Convex optimization primer: proximal gradient methods
- **3** Variational inverse problems
- 4 Segmentation and labeling problems
- **5** 3D reconstruction







Introduction to variational methods

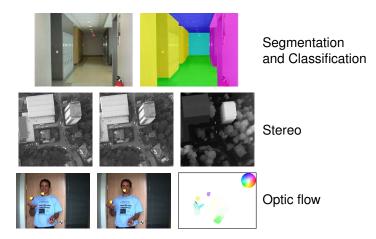
- Convex optimization primer: proximal gradient methods
- **3** Variational inverse problems
- 4 Segmentation and labeling problems





Fundamental problems in computer vision

Image labeling problems

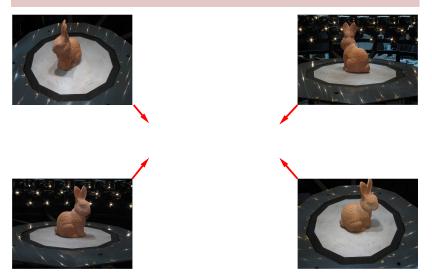






Fundamental problems in computer vision

3D Reconstruction



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Variational Methods, Eurographics 2014 Tutorial "Optimization Techniques in Computer Graphics"





Unknown object: Vector-valued function

$$u:\Omega \to \mathbb{R}^m$$

on a **continuous** domain Ω .

• Problem solution: minimizer of an energy functional

$$\underset{u \in \mathcal{V}}{\operatorname{argmin}} \{ \underbrace{J(u)}_{\text{regularizer}} + \underbrace{F(u)}_{\text{data term}} \}.$$

on an infinite dimensional (function) space $\ensuremath{\mathcal{V}}.$





Continuous vs. discrete methods

Advantages:

- Realistic physical models of the world naturally continous
- Exact modeling of geometric quantities (length, curvature) ...
- No grid bias (rotation invariance)
- Solvers easy to parallelize, fast on GPU

Disadvantages:

- · Less flexible than probabilistic graphical models
- Iterative solvers (usually no runtime guarantee)
- Slow on single core





Rudin-Osher-Fatemi (ROF), or TV- \mathcal{L}^2 model (1992)

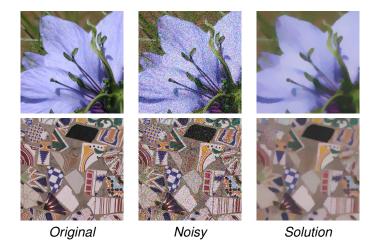
Given an image *f* and a smoothness parameter $\lambda > 0$, compute a **denoised** image as the minimizer

$$\underset{u \in L^{2}(\Omega)}{\operatorname{argmin}} \left\{ \left\| \nabla u \right\|_{1} + \frac{1}{2\lambda} \left\| u - f \right\|_{2}^{2} \right\}.$$

- Can be interpreted as maximum a posteriori (MAP) estimate under assumption of Gaussian noise on f.
- The energy is convex, but not differentiable, in particular gradient based methods do not work







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The TV-L² model

$$\underset{u}{\operatorname{argmin}} \left\{ \left\| \nabla u \right\|_{1} + \frac{1}{2\lambda} \left\| u - f \right\|_{2}^{2} \right\}.$$

is a building block for several more general algorithms. It is therefore important to fully understand it.

- Note that the total variation is not differentiable, so it cannot be exactly minimized by simple gradient descent, although it is convex.
- In this section, we discuss discretization and the most straight-forward exact solver I know of.





In the continuous setting, an image is a function or scalar field $u: \Omega \to \mathbb{R}$ with $\Omega \subset \mathbb{R}^2$. For simplicity, we assume in this course that the domain is always two-dimensional.

Discretization of an image

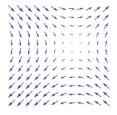
- the domain Ω is a regular grid of size $M \times N$
- an image u is a matrix $u \in \mathbb{R}^{M \times N}$
- the gray value or intensity at a pixel (*i*, *j*) is written as *u*_{*i*,*j*}



Color images have several components, usually red, green and blue for the primary colors. Thus, they are functions $u : \Omega \to \mathbb{R}^3$. In the discretization, each component u^1 is represented by its own matrix of intensity values.



A vector field is a map $\mathbf{p}: \Omega \to \mathbb{R}^2$. One can imagine it as an assignment of a little arrow to each point:



Discretization of a vector field

- a vector field **p** consists of two components p¹, p², each of which is a matrix of size M × N
- the vector at a pixel (i, j) is written as $\begin{bmatrix} k \\ j \end{bmatrix}$

s
$$\begin{bmatrix} p_{i,j}^1\\ p_{i,j}^2 \end{bmatrix}$$
.



In the continuous setting, one of the basic operators acting on a function is the derivative. The vector field of the two partial derivatives of a function $u: \Omega \to \mathbb{R}$ is called the **gradient** of u,

$$\nabla u = \begin{bmatrix} \partial_x u \\ \partial_y u \end{bmatrix}.$$

Discretization of the gradient with forward differences For i < N and j < M:

$$(\nabla u)_{i,j} = \begin{bmatrix} u_{i+1,j} - u_{i,j} \\ u_{i,j+1} - u_{i,j} \end{bmatrix}.$$

Neumann boundary conditions $\frac{\partial u}{\partial \mathbf{n}} = \mathbf{0}$ are used for functions, i.e.

$$(\nabla u)_{M,j}^1 = 0$$
 and $(\nabla u)_{i,N}^2 = 0$.



In the continuous setting, the divergence of a vector field **p** is the scalar field or function div(**p**) : $\Omega \to \mathbb{R}$ defined as

$$\operatorname{div}(\mathbf{p}) = \partial_x p^1 + \partial_y p^2.$$

Discretization of the divergence with backward differences

Derivatives on vector fields use Dirichlet boundary conditions ${\bm p}|_{\partial\Omega}{=}$ 0. Thus,

$$(\operatorname{div}(\mathbf{p}))_{i,j} = \begin{cases} p_{i,j}^1 - p_{i-1,j}^1 & \text{if } 1 < i \le M \\ p_{i,j}^1 & \text{if } i = 1 \\ + \begin{cases} p_{i,j}^2 - p_{i,j-1}^2 & \text{if } 1 < j \le N \\ p_{i,j}^2 & \text{if } j = 1 \end{cases}$$



Let $u \in \mathbb{R}^{M \times N}$ be a matrix, then its **total variation** is defined as

$$\mathsf{TV}(u) := \sum_{i=1}^{M} \sum_{j=1}^{N} \left| (\nabla u)_{i,j} \right|_2,$$

where the gradient is computed with forward differences and Neumann boundary conditions.

Note: Actually computing the discrete TV requires to compute the sum over all elements in a matrix, which is a not fully parallelizable type of operation called a **reduction**.





The Euclidean norm in \mathbb{R}^n can be written for any vector v as

$$|\boldsymbol{v}|_2 = \sup_{\boldsymbol{p} \in \mathbb{R}^n, |\boldsymbol{p}|_2 \leq 1} (\boldsymbol{v}, \boldsymbol{p}),$$

where (v, p) denotes the usual inner product.

We can use this fact to rewrite the total variation of a matrix as follows:

$$\mathsf{TV}(u) = \max_{\mathbf{p} \in K} \sum_{i,j} \left((\nabla u)_{i,j}, \, \mathbf{p}_{i,j} \right),$$

where the (convex) set K consists of all vector fields such that their vectors in each point have length at most one,

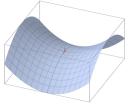
$$\mathcal{K} := \left\{ \mathbf{p} \in \mathbb{R}^{M \times N} \times \mathbb{R}^{M \times N} \; : \; \left| \mathbf{p}_{i,j} \right|_2 \leq 1 \text{ for all } i,j \right\}.$$





Instead of finding a minimizer for ROF, we can look for a saddle point (u, \mathbf{p}) of

$$\min_{\boldsymbol{u}\in\mathbb{R}^{M\times N}}\max_{\mathbf{p}\in\mathcal{K}}\sum_{i,j}\left((\nabla\boldsymbol{u})_{i,j}\,,\,\mathbf{p}_{i,j}\right)+\frac{1}{2\lambda}(\boldsymbol{u}_{i,j}-\boldsymbol{f}_{i,j})^{2}.$$



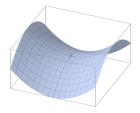
- Note that the energy is now differentiable, so as a simple method to find a saddle point we can iterate gradient descent in *u* and ascent in **p** starting from an initial point (*u*₀, **p**₀).
- We will formulate a slightly faster method, which is based on the observation that for given **p**, one can solve the minimization in *u* explicitly.





The energy of the saddle point problem is

$$E(u,\mathbf{p}) = \sum_{i,j} \left((\nabla u)_{i,j}, \mathbf{p}_{i,j} \right) + \frac{1}{2\lambda} (u_{i,j} - f_{i,j})^2.$$



Its derivative with respect to **p** is just

$$\frac{d}{d\mathbf{p}}E(u,\mathbf{p})=\nabla u,$$

so a (discrete) gradient ascent step for ${\boldsymbol{p}}$ is given by setting

$$\mathbf{p}^{n+1} = \mathbf{p}^n + \tau \nabla u^n,$$

where (\mathbf{p}^n, u^n) is the *n*th iterate for the saddle point and $\tau > 0$ is a small enough step size (to be determined later).





The allowed range for **p** is the set *K*. After the update step, \mathbf{p}^n might lie outside *K*, so we have to back-project it onto *K*. This can be done by projecting each single vector of the field **p** back onto the unit disk,

$$\Pi_{\mathcal{K}}(\mathbf{p}^{n})_{i,j} = \frac{\mathbf{p}_{i,j}^{n}}{\max\left(1, \left|\mathbf{p}_{i,j}^{n}\right|_{2}\right)}.$$





The reason the finite difference operators introduced in the first lecture were defined in this way is that they need to satisfy

$$\sum_{i,j} \left((
abla u)_{i,j} \,,\, \mathbf{p}_{i,j}
ight) = - \sum_{i,j} u_{i,j} \,(\operatorname{div}(\mathbf{p}))_{i,j}$$

like their continuous counterparts.

This means we can rewrite the saddle point energy as

$$E(u, \mathbf{p}) = \sum_{i,j} -u_{i,j} (\operatorname{div}(\mathbf{p}))_{i,j} + \frac{1}{2\lambda} (u_{i,j} - f_{i,j})^2.$$



The gradient with respect to *u* is

$$rac{d}{du} E(u,\mathbf{p}) = - {
m div}(\mathbf{p}) + rac{1}{\lambda}(u-f),$$

At a minimum for u, this must be zero. This leads to the update equation

$$u^{n+1} = f + \lambda \operatorname{div}(\mathbf{p}^{n+1})$$

which computes the exact solution for u^{n+1} explicitly, given the previously computed \mathbf{p}^{n+1} .



Algorithm to solve ROF (Bermudez-Moreno 1981, Chambolle 2005)

- Given an image *f* to be denoised and a smoothing parameter λ > 0.
- Set the step size to $\tau := \frac{1}{2\lambda \|\nabla\|} = \frac{1}{8\lambda}$.
- Start with initial *u*⁰, **p**⁰, where *u*⁰ is a scalar field and **p**⁰ is a vector field. Initial values can be arbitrary, choose e.g. zero for everything.
- Iterate until convergence:

$$\mathbf{p}^{n+1} = \Pi_{\mathcal{K}}(\mathbf{p}^n + \tau \nabla u^n)$$

 $u^{n+1} = f + \lambda \operatorname{div}(\mathbf{p}^{n+1})$



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- A Hilbert space \mathcal{V} with
 - inner product (\cdot, \cdot) and
 - norm $\|\cdot\| = \sqrt{(\cdot, \cdot)}$,

in this tutorial, this will be $\mathcal{V} = R^N$ or $\mathcal{V} = L^2(\Omega)$.

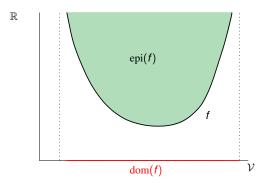
- The dual space \mathcal{V}^* in our setting, you can substitute $\mathcal{V}^*=\mathcal{V}$ if you are not familiar with the concept
- We consider "proper" functionals *f* : *V* → ℝ ∪ {∞}, i.e. they can take the value infinity, but we exclude the (boring) functional *f* = ∞ for technical reasons.





The **epigraph** epi(*f*) of a functional $f : \mathcal{V} \to \mathbb{R} \cup \{\infty\}$ is the set "above the graph", i.e.

$$\mathsf{epi}(f) := \{(x, \mu) : x \in \mathcal{V} \text{ and } \mu \ge f(x)\}.$$





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Definition

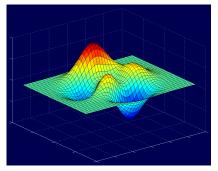
- A functional f : V → ℝ ∪ {∞} is called convex if epi(f) is a convex set.
- The set of all proper and convex functionals on V is denoted conv(V).

We choose the geometric definition of a convex function here because it is more intuitive, the usual algebraic property is a simple consequence.

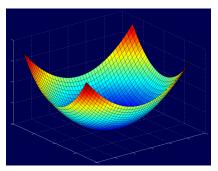




Convex versus non-convex energies



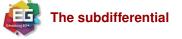
non-convex energy



convex energy

Convex energies can be globally minimized for non-convex energies this is usually impossible.





Definition

• A vector $\varphi \in \mathcal{V}^*$ is a **subgradient** of *f* at $x \in \mathcal{V}$ if

$$f(y) \ge f(x) + \langle y - x, \varphi \rangle$$
 for all $y \in \mathcal{V}$.

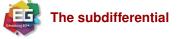
• The set of all subgradients of *f* at *x* is the **subdifferential** $\partial f(x)$.

Geometrically: the function on the right hand side is an affine function

$$h(y) = f(x) + \langle y - x, \varphi \rangle$$

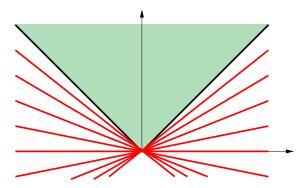
with slope φ , which touches the epigraph of *f* in *x*. The condition says that *h* needs to lie below *f*.





Example: the subdifferential of $f : x \mapsto |x|$ in 0 is

$$\partial f(\mathbf{0}) = [-1, 1].$$







The subdifferential is a generalization of the Fréchet derivative (or the gradient in finite dimension), in the following sense.

Theorem

Let $f \in conv(\mathcal{V})$ be Fréchet differentiable at $x \in \mathcal{V}$. Then

 $\partial f(x) = \{ df(x) \}.$

The proof of the theorem is surprisingly involved - it requires to relate the subdifferential to one-sided directional derivatives. We will not explore this here, see e.g. [Rockafeller 1970].



Theorem

Let $f \in conv(\mathcal{V})$. Then \hat{x} is a global minimum of f if and only if

 $0 \in \partial f(\hat{x}).$

In particular, a convex functional does not have minima which are local but not global.

Geometrically intuitive, but the proof is also surprisingly simple. Rewriting the subgradient definition, one sees that $0 \in \partial f(\hat{x})$ just means that

 $f(y) - \langle y, 0 \rangle \ge f(\hat{x}) - \langle \hat{x}, 0 \rangle$ for all $y \in \mathcal{V}$.

This already implies the equivalence.

Proximity operator

 The proximity operator for a closed convex functional H and step size parameter τ > 0 is defined as

$$\operatorname{prox}_{\tau H}(x) = \operatorname{argmin}_{y} \left\{ \frac{1}{2\tau} \left\| y - x \right\|^{2} + H(y) \right\}.$$

• At the minimizer
$$\hat{x} = \text{prox}_{\tau H}(x)$$
,

$$0 \in rac{1}{ au}(\hat{x}-x) + \partial H(\hat{x})$$

 $\Leftrightarrow \quad \hat{x} \in x - au \, \partial H(\hat{x}).$

The proximity operator computes an implicit subgradient descent for the convex functional H with step size τ .





• Let H = TV, then

$$\operatorname{prox}_{\lambda TV}(x) = \operatorname{argmin}_{y} \left\{ TV(y) + \frac{1}{2\tau} \left\| x - y \right\|^{2} \right\}$$

is an instance of ROF-denoising.

• Solve e.g. with Bermudez-Moreno.





- Let $\mathcal{C} \subset \mathcal{V}$ be a convex set, and

$$\delta_{\mathcal{C}}(x) := egin{cases} \mathsf{0} & ext{if } x \in \mathcal{C}, \ \infty & ext{otherwise} \end{cases}$$

be its indicator function.

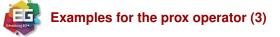
• Then

$$\operatorname{prox}_{\delta_{\mathcal{C}}}(x) = \operatorname{argmin}_{y \in \mathcal{V}} \left\{ \frac{1}{2\tau} \|y - x\|^2 + \delta_{\mathcal{C}}(y) \right\}$$
$$= \operatorname{argmin}_{y \in \mathcal{C}} \left\{ \|y - x\|^2 \right\}$$

is the projection onto C.

Adding an indicator function to a functional is equivalent to an optimization constraint.





 Let *H* = ∥·∥₁ (sparsity inducing norm), e.g. interesting in compressive sensing or dictionary learning, next talk.

• Then (component-wise)

$$\operatorname{prox}_{\tau \parallel \cdot \parallel_1}(x)_i = (|x_i| - \tau)\operatorname{sgn}(x_i),$$

which is called the shrinkage operator.



Applicability: Minimization problems of the form

 $\operatorname*{argmin}_{u\in\mathcal{V}}\left\{G(u)+F(u)\right\}$

such that

- G is convex
- F is convex and differentiable
- *dF* is Lipschitz continuous, i.e. there exist *L* > 0 such that

 $\|dF(u) - dF(v)\| \le L \|u - v\|$ for all $u, v \in \mathcal{V}$.

• The proximation for *G* is (comparatively) easy to evaluate.





Iterative shrinkage and thresholding (ISTA)

Algorithm: Initialize with $u_0 = 0$,

then compute alternating "forward-backward steps":

• gradient descent in F:

$$v_{n+1}=u_n-\frac{1}{L}dF(u_n).$$

• implicit subgradient descent in G:

$$u_{n+1} = \operatorname{prox}_{\frac{1}{L}G}(v_{n+1}).$$

only O(1/n) convergence (slow).

Bruck 1977, Passty 1979





Algorithm: Initialize with $u_0 = 0$, $\overline{u}_0 = 0$, $\mathbb{R} \ni t_0 = 1$, then iterate

• gradient descent in F:

$$v_{n+1}=\bar{u}_n-\frac{1}{L}dF(\bar{u}_n).$$

• implicit subgradient descent in G:

$$u_{n+1} = \operatorname{prox}_{\frac{1}{L}G}(v_{n+1}).$$

• extragradient step:

$$t_{n+1} = \frac{1}{2}(1 + \sqrt{1 + 4t_n^2}),$$

$$\bar{u}_{n+1} = u_n + \frac{t_n - 1}{t_{n+1}}(u_{n+1} - u_n).$$

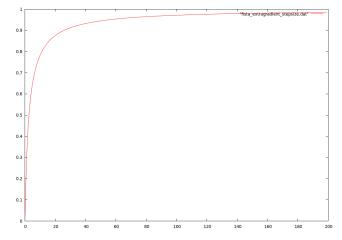
Beck and Teboulle SIIMS 2009







FISTA extragradient step size development



- Leads to $O(1/n^2)$ algorithm (optimal for a first-order method)
- Good general-purpose solver, easy to implement (in particular if proximation is already available)

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Let $C \subset V$ be a convex constraint set, i.e. the solution shall be restricted to lie in C. Thus, we want to solve

 $\operatorname*{argmin}_{x\in\mathcal{V}}\left\{ F(x)\right\} \text{ subject to }x\in\mathcal{C},$

which could be e.g. a constrained least squares problem as in the first talk. This is equivalent so solving

$$\operatorname*{argmin}_{x\in\mathcal{V}}\left\{F(x)+\delta_{\mathcal{C}}(x)\right\}.$$

When applying FISTA, $\delta_{\mathcal{C}}$ is the non-differentiable part taken care of via the prox operator, which equals the projection onto \mathcal{C} .





Many other proximal gradient methods ...

- Alternating projection (projection onto convex sets)
- Douglas-Rachford splitting, alternating direction method of multipliers ADMM (popular e.g. in learning)
- Split-Bregman (basis pursuit problems, i.e. constrained *L*¹-minimization)
- Chambolle-Pock (swiss army knife if both *F* and *G* are non-differentiable, but convex with simple prox-operators).

All of these use the same basic building blocks as the ones shown in the previous section.





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The convolution b * u with a kernel b of total mass 1 can be interpreted as a blurring operation.

Example: Gaussian blur (isotropic)





Example: Motion blur for diagonal motion (anisotropic)





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Inverse problem: We observe the image f, which was created from an unknown original via

f = b * u + Gaussian noise.

More general, f = Au + noise (linear inverse problem).

TV deblurring model

Recover a deblurred image u as the minimizer of

$$\underset{u}{\operatorname{argmin}}\left\{\left\|\nabla u\right\|_{1}+\frac{1}{2\lambda}\left\|b*u-f\right\|_{2}^{2}\right\},$$

which is again the MAP estimate for Gaussian noise with TV prior.

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Original



Blurred + Noisy



Solution





• The differentiable part is

$$F(u) = \frac{1}{2\lambda} \|Au - f\|_2^2,$$

which has functional derivative

$$dF(u) = \frac{1}{\lambda}A^{T}(Au - f),$$

which is Lipschitz-continuous with $L = \frac{1}{\lambda} \| A^T A \|_2$.

• The non-differentiable convex part is the regularizer, e.g. TV, with the proximation given by solving an instance of ROF.







Inpainting problem

- Restoration problem: recover missing regions of an image
- Input:
 - damaged region $\Gamma\subset \Omega$
 - partial image $f: \Omega \setminus \Gamma \to \mathbb{R}$
- TV inpainting: Solve

 $\min_{u} \|\nabla u\|_1 \text{ such that } u = f \text{ on } \Omega \setminus \Gamma.$



Damaged image f



Recovered image u





TV inpainting results ("textured" images)



Original

Damaged

Inpainted

- TV inpainting is unconvincing for highly textured images if the missing regions are larger. The reason is that no structure is inferred from surrounding regions, and only boundary values of Γ are taken into account.
- A better variational model for inpainting can be found e.g. in papers on **non-local TV** by Osher et al., or check methods based on **dictionary learning**.





Once we have an inpainting algorithm, we can employ it to remove unwanted regions in an image by marking them as damaged.



Uncaging a bird







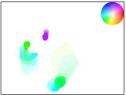
 I_t at time t

Input

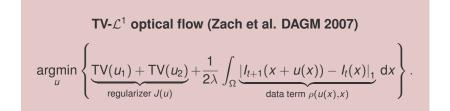


 I_{t+1} at time t+1

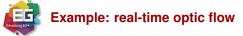
Unknown



flow field $u: \Omega \to \mathbb{R}^2$



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Idea: introduce auxiliary variable v

- Decouples regularizer from the data term.
- Let θ > 0 be a coupling parameter, define family of energies in the variables u, v,

$$E_{\theta}(u,v) = \mathcal{J}(u) + \frac{1}{2\theta} \left\| u - v \right\|_{2}^{2} + \int_{\Omega} \rho(x,v(x)) \,\mathrm{d}x.$$

For θ → 0, the quadratic term forces u to be close to v, so the solution of E_θ approximates the solution of E.





Algorithm

Set $\theta > 0$, start with initial u_0 , v_0 . Then iterate the following steps until convergence.

• Solve an instance of TV-L²:

$$u_{k+1} = \operatorname*{argmin}_{u} E_{\theta}(u, v_k).$$

Solve the point-wise problem

$$v_{k+1} = \operatorname*{argmin}_{v} E_{\theta}(u_{k+1}, v).$$

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same technique (just reduced target dimension): real-time disparity (depth) maps



Stuehmer et al. DAGM 2010

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Wedel et al. ECCV 2008

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The MAP estimate for binary segmentation

Given:

- An observed image f on a domain Ω, from which one computes (e.g. via a trained classifier)
- a foreground probability distribution

 $P_{fg}(x) =$ Probability that the point x is in the foreground

• A background probability distribution

 $P_{bg}(x) =$ Probability that the point *x* is in the background

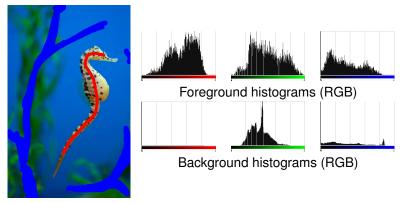
Wanted:

 The MAP estimate for a binary segmentation u : Ω → {0,1}, where u(x) = 1 means a point is in the foreground, and u(x) = 0 means it is in the background:

$$\underset{u:\Omega \to \{0,1\}}{\operatorname{argmax}} P(u \mid f) = \underset{u:\Omega \to \{0,1\}}{\operatorname{argmax}} P(f \mid u)P(u).$$



Example: histograms of user input (naive, illustration only)



 $P_{fg}(x) = rac{\text{number of marked foreground pixels with color } f(x)}{\text{total number of marked foreground pixels}}$

 $P_{bg}(x) = rac{\text{number of marked background pixels with color } f(x)}{\text{total number of marked background pixels}}$





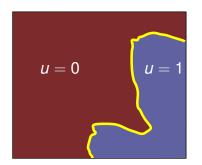
Under the above assumptions, the MAP estimate for the segmentation is

$$\operatorname*{argmin}_{u:\Omega \to \{0,1\}} \left\{ J(u) + \int_{\Omega} u \cdot \log \left[\frac{p_{bg}}{p_{fg}} \right] \mathrm{d}x \right\}.$$

Here, J(u) is the **prior**. A typical prior is the length of the interface.







Length of interface equals total variation of u,

$$\mathsf{TV}(u) = \int_{\Omega} |Du|$$

Binary segmentation problem

$$\operatorname*{argmin}_{u:\Omega\to\{0,1\}}\mathsf{TV}(u)+\int_{\Omega}cu\,\mathsf{d}x,$$

with local assigment costs c.

Space of binary functions $u : \Omega \to \{0, 1\}$ not convex but: globally optimal solution possible by relaxation to $u : \Omega \to [0, 1]$ and subsequent thresholding.

Chan, Esedoglu and Nikolova 2006

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Straightforward:

• Since F(u) = (c, u), the derivative is constant,

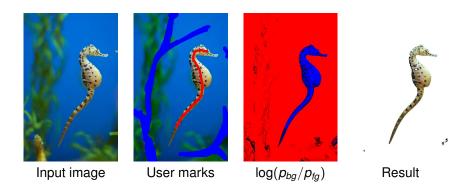
$$dF(u) = a.$$

- It is Lipschitz-continuous with arbitrary Lipschitz constant L > 0, one can e.g. just choose L = 1.
- Thus, the gradient descent step in *F* just subtracts *c* from the solution.
- Proximation in the TV-regularizer is an instance of ROF.





Segmentation example



Simple model works pretty well with easy, clean input ...



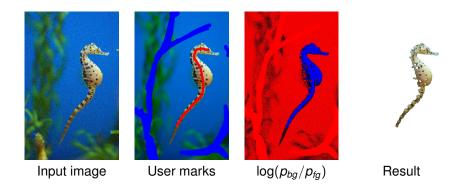
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Variational Methods, Eurographics 2014 Tutorial "Optimization Techniques in Computer Graphics"



Segmentation example: noisy input



... but is not very robust against noise (better classifiers requires, i.e. random forests)

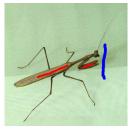


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Segmentation example: elongated structures







Input with User marks $\log(p_{bg}/p_{fg})$ (normalized) TV (best result)

Length does not work well with elongated structures (curvature regularity is better in that case, but complex)

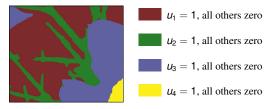
Goldluecke and Cremers ICCV 2011



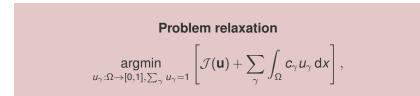


Convex relaxation of the multilabel problem

Indicator function $u_{\gamma} : \Omega \to \{0, 1\}$ assigned to each label γ :



 $\sum_{\gamma} u_{\gamma}$ must be one !



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The regularization penalty is proportional to the **label distance** times the **length of the interface**.

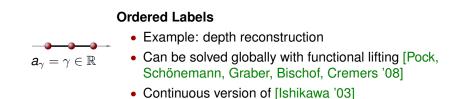
In this example $d(\gamma_1, \gamma_2) \cdot L(\Sigma)$



Euclidean representation of the label distance:

- Each label γ is *represented* by a point $a_{\gamma} \in \mathbb{R}^{k}$.
- Label distance $d(\gamma, \mu) = |a_{\gamma} a_{\mu}|_2$.





 $a_{\gamma} = e_{\gamma} \in \mathbb{R}^{N}$

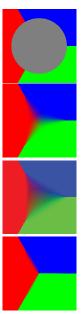
Potts model

- Example: segmentation
- No globally optimal solution possible if N > 2
- Continuous version of [Potts '52]





Different Multilabel Regularizers



A comparison using the "triple-junction" problem

Zach, Gallup, Frahm, Niethammer '08

$$\mathcal{J}_1(\mathbf{u}) = rac{1}{2}\sum_\gamma \int_\Omega |Du_\gamma|$$

Lellmann, Kappes, Yuan, Becker, Schnörr '08

$$\mathcal{J}_{2}(\mathbf{u}) = \int_{\Omega} \sqrt{\sum_{\gamma} \|Du_{\gamma}\|_{A}^{2}}, \ \|v\|_{A} = \sqrt{v^{T}A^{T}Av}$$

Chambolle, Cremers, Pock '08

$$\mathcal{J}_{3}(\mathbf{u}) = \int_{\Omega} \Psi(D\mathbf{u}), \Psi(oldsymbol{q}) = \sup_{oldsymbol{p}} \left\{ \langle oldsymbol{q}, oldsymbol{p}
ight\} \; : \; |oldsymbol{p}_{\gamma} - oldsymbol{p}_{\mu}| \leq oldsymbol{d}_{\gamma\mu}
ight\}$$

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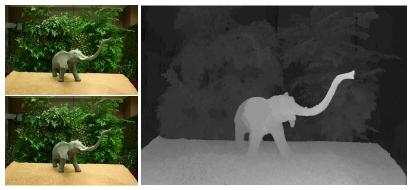
Variational Methods, Eurographics 2014 Tutorial "Optimization Techniques in Computer Graphics"





Results: indoor stereo pair

- Stereo assignment cost $c_{\gamma}(x) = |I_{\text{left}}(x) I_{\text{right}}(x + \gamma)|$
- Linear discontinuity penalty \Rightarrow globally optimal solution.



Images from UCSD lightfield repository

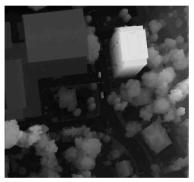




Results: aerial images (1400 x 1500 pixels)



One of two input images (Courtesy of Microsoft Graz)



Depth reconstruction



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Potts segmentation with k = 10 labels





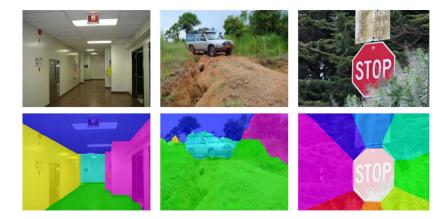
Potts segmentation with k = 16 labels











Labeling regions should have certain spatial relationships, *i.e. heaven is always above ground.*





Strekalovskiy and Cremers, ICCV 2011

The labeling penalty may depend also on the **normal n of the interface** between two regions,

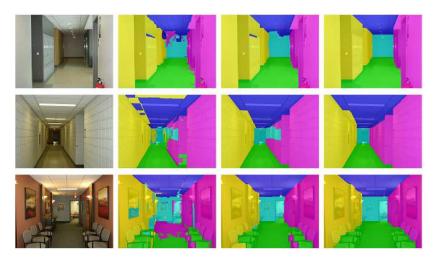
$$\boldsymbol{d}: \boldsymbol{\Gamma} \times \boldsymbol{\Gamma} \times \mathbb{S} \to \mathbb{R}^+.$$

New direction-aware regularizer:

$$\begin{split} \mathcal{J}(\mathbf{u}) &= \sup_{\mathbf{p} \in \mathcal{C}} \sum_{\gamma} \int_{\Omega} \langle \mathbf{p}_{\gamma}, \nabla u_{\gamma} \rangle \, \mathrm{d}x, \\ \text{with } \mathcal{C} &= \{ (\mathbf{p}_{\gamma} : \Omega \to \mathbb{R}^{n}) : \langle \mathbf{p}_{\mu} - \mathbf{p}_{\gamma}, \mathbf{n} \rangle \leq \mathcal{d}(\gamma, \mu, \mathbf{n}) \quad \forall \gamma, \mu, \mathbf{n} \} \, . \end{split}$$

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Input

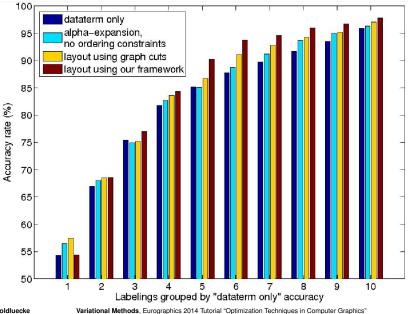
Data term

Potts

Ordering









1 Introduction to variational methods

- 2 Convex optimization primer: proximal gradient methods
- **3** Variational inverse problems
- Segmentation and labeling problems

5 3D reconstruction

6 Summary



Given *n* images $\mathcal{I}_i : \Omega_i \to \mathbb{R}^3$ with projections $\pi_i : \mathbb{R}^3 \to \mathbb{R}^2$











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Given *n* images $\mathcal{I}_i : \Omega_i \to \mathbb{R}^3$

with projections $\pi_i : \mathbb{R}^3 \to \mathbb{R}^2$



Find surface $\Sigma \subset \mathbb{R}^3$ with texture $T : \Sigma \to \mathbb{R}^3$ which optimally matches the input images.



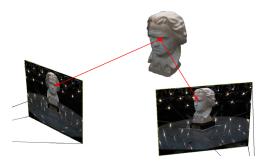


Variational 3D reconstruction

Classical variational formulation (Faugeras and Keriven, 1998)

Find a surface $\Sigma \subset \mathbb{R}^3$ which minimizes the photo-consistency error,

$$\underset{\Sigma}{\operatorname{argmin}} \int_{\Sigma} \rho(s) \, \mathrm{d}s.$$



- Point on surface
- projections look similar in different views
- small value of ρ



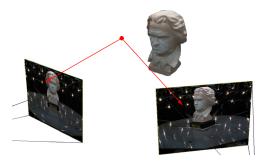


Variational 3D reconstruction

Classical variational formulation (Faugeras and Keriven, 1998)

Find a surface $\Sigma \subset \mathbb{R}^3$ which minimizes the photo-consistency error,

$$\underset{\Sigma}{\operatorname{argmin}} \int_{\Sigma} \rho(s) \, \mathrm{d}s.$$



- Point not on surface
- projections look different in different views
- large value of ρ

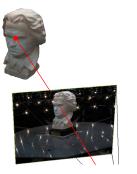




Convex functional minimizes photo-consistency error:

 $\operatorname*{argmin}_{u:\Omega \to \{0,1\}} \int_{\Omega} \rho \left| Du \right|$

Silhouette constraints to avoid constant solutions



A ray through the **silhouette** must **intersect** the surface

Kolev, Klodt, Brox, Cremers IJCV'09



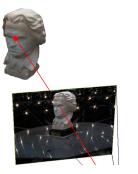


3D reconstruction as a convex variational problem

Convex functional minimizes photo-consistency error:

 $\operatorname*{argmin}_{u:\Omega \to \{0,1\}} \int_{\Omega} \rho \left| \mathcal{D} u \right|$

Silhouette constraints to avoid constant solutions



A ray through the **background** must **miss** the surface

Kolev, Klodt, Brox, Cremers IJCV'09







Kolev and Matiouk, Akademisches Kunstmuseum Bonn 2010

Bastian Goldluecke





Superresolution texture maps

Given



n images $\mathcal{I}_i : \Omega_i \to \mathbb{R}$ projections $\pi_i : \mathbb{R}^3 \to \mathbb{R}^2$



approximate surface Σ (assumed Lambertian)

Find



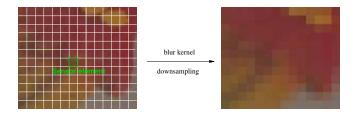
high-res color texture map $\mathcal{T}:\Sigma\to\mathbb{R}^3$ and accurate geometry

Idea:

Use information from multiple cameras for superresolution.



Image formation in a sensor



- Each sensor element samples incoming light over its area
- Sampling modeled by blur kernel b
- · Leads to image formation model

$$\mathcal{I}_{ ext{observed (low-res)}} = b * \mathcal{I}_{ ext{incoming (high-res)}}$$



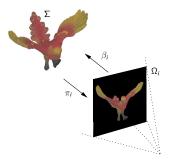
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Data term: squared difference between input images and downsampled high-resolution rendering

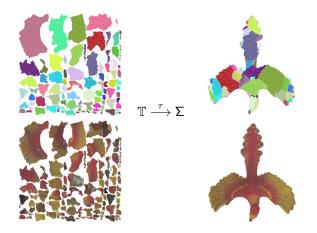
$$E(T) := \sum_{i=1}^{n} \int (\underbrace{b * (T \circ \beta_i)}_{\text{Subsampled hires rendering}} - \underbrace{\mathcal{I}_i}_{\text{Input image}})^2 \, \mathrm{d}x$$

Regularizer: total variation on the surface.



Goldluecke and Cremers, ICCV 2009





Conformal atlas: transforms energy to 2D texture space

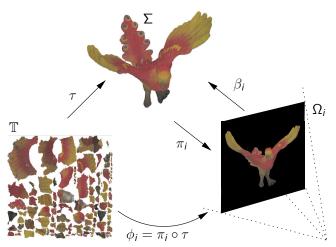
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Euler-Lagrange equation on texture space

$$\frac{1}{\lambda} \mathsf{div}\left(\sqrt{\lambda} \frac{\nabla \mathcal{T}}{\|\nabla \mathcal{T}\|}\right) + \sum_{i=1}^{n} \frac{v_i}{\sigma} \left(\left(\mathcal{J}_i \mathcal{E}_i\right) \circ \phi_i \right) = 0 \text{ on } \mathbb{T}$$



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- 30 cameras, input image resolution 768×576
- Initial geometry: Kolev and Cremers, ECCV 2008
- Texture resolution 2048×2048

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Rendered model



Input image

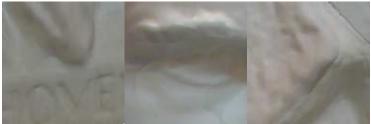


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Rendered model



Input image



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Additional dependance on **displacement map** $D : \mathbb{T} \to \mathbb{R}$,

$$E(T, \mathbf{D}) := \sum_{i=1}^{n} \int_{\hat{S}_{i}} \left(b * (T \circ \beta_{i}^{\mathbf{D}}) - \mathcal{I}_{i} \right)^{2} \mathrm{d}x + E_{\mathrm{tv}}(T, \mathbf{D}).$$

- In *T*: energy is convex
- In *D*: multilabel problem with convex regularizer, global optimization in *D* possible

Geometry

Estimated Texture

Input

Estimated displacement

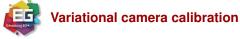
Without displacement

With displacement

Goldlücke and Cremers DAGM 2009



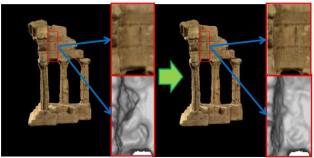
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Idea: assume the **projection parameters** π **are unknowns** in the superresolution energy,

$$E(T,\pi) := \sum_{i=1}^N \int_{\mathcal{S}_i} \|b*(T\circ\beta_i) - \mathcal{I}_i\| dx.$$

Optimize alternatingly for texture and projection. In a way, this can be thought of as a continuous version of bundle adjustment.



Goldluecke, Aubry, Kolev, Cremers, IJCV 2013.







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- Variational methods: a powerful tool for realistic modeling
- Non-convex energies: gradient descent, local minimum, usually requires good initialization
- Convex energies: no local minima, powerful optimization tools (e.g. proximal gradient methods)
- For argmin_u{*F*(*u*) + *G*(*u*)} with convex *G* and differentiable and convex *F*, FISTA is a good general-purpose method with optimal convergence rate.
- If *F* is only convex, check out e.g. [Chambolle-Pock 2012] or [Nesterov 2005].
- Numerous appliations: variational inverse problems, segmentation and labeling, 3D reconstruction ...

