

Practical Course: GPU Programming in Computer Vision Mathematics 4: Variational Methods

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Outline

Variational methods

2 Image denoising

Energy Minimization



Image denoising

Energy Minimization

Energy minimization

An established approach to model numerous computer vision problems.

Energy

Every possible candidate solution u is assigned an energy E(u). Idea: E(u) measures the costs of u: The smaller the costs the better the solution.

Minimizers

Candidates *u* with *least* energy are considered solutions to the problem.

- Clear mathematical correspondence between input data and result
- Extensive mathematical theory, optimality conditions
- Can describe sophisticated problems with only a few parameters
- Lots of algorithms to compute the minimizers

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Typical form

$$E(u) = D(u) + R(u)$$

- **Data term** D(u) measures how well the solution u fits input data.
- **Regularizer** R(u) enforces regularity and smoothness of u.

Minimizing *E* will give a solution *u* which fits to the inputs and is smooth

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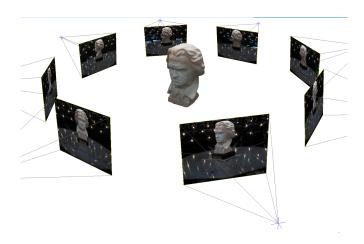
Minimizing *E* will give a solution *u* which fits to the inputs and is smooth!





Example: 3D reconstruction

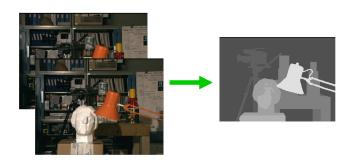
Input: views of an object from different cameras. Find: the 3D-object.





Example: Depth reconstruction

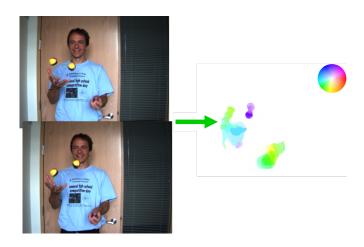
Input: a pair of stereo images. Find: the depth in every pixel





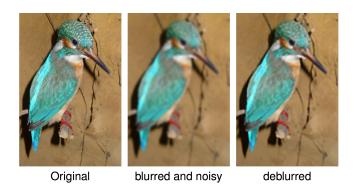
Example: Optical flow

Input: a pair of images. Find: displacement vector for each pixel





Input: a blurry image. Find: a deblurred image.



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Example: Segmentation

Input: a color image. **Find:** object with certain given characteristics (colors distribution etc.).







Example: Multilabel Segmentation

Input: a color image. **Find:** a meaningful decomposition into several regions.



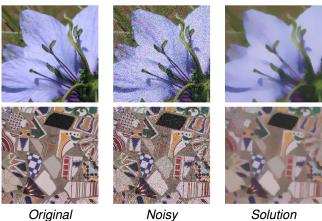






Image Denoising: The Problem

Input: a noisy image $f: \Omega \to \mathbb{R}^n$. **Find:** denoised $u: \Omega \to \mathbb{R}^n$.



Solution





Outline

2 Image denoising

3 Energy Minimization



Data term

■ The clean image *u* must be *similar* to the noisy image *f*:

$$D(u) := \int_{\Omega} (u(x,y) - f(x,y))^2 dx dy$$

■ Minimize D(u) to guarantee that $u \approx f$.

- Solution *u* must be noise-free, so we look for *smooth* images *u*.
- Colors in neighboring pixels must be similar, i.e. $|\nabla u|$ must be small:

$$R(u) := \lambda \int_{\Omega} \phi(|(\nabla u)(x, y)|) dx dy$$

- $\phi: \mathbb{R} \to \mathbb{R}$ is an increasing function, $\lambda > 0$ is a weighting parameter.
- Minimize R(u) to guarantee that $|\nabla u|$ is small, and u noise-free.

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Denoising energy

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If $\mu = f$:

Perfect fit for data: D(u) = 0. But u noisy: $R(u) \gg 1$.

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If u = const:

Bad fit for data: $D(u) \gg 1$. But u smooth: R(u) = 0.

True solution

Will be a *trade-off* between data fitting and smoothness λ controls the desired degree of smoothness of u.

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Energy Minimization



Denoising Energy

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How to find the minimizer *u* in practice?

There are many methods. The most common ones are

- Gradient descent: Go along the negative "gradient" of the energy.
- Euler-Lagrange equation: Necessary condition for the minimizers.
- 3 Primal-dual methods: Very flexible iterative algorithms



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Gradient Descent: Gradient of the Energy

Intuitively: $(\nabla E)(u)$ is the gradient w.r.t. values u(x, y) at each (x, y).

Analogy with finite $e: \mathbb{R}^k \to \mathbb{R}$:

- For $z \in \mathbb{R}^k$: $(\nabla e)(z)$ has $(\dim \mathbb{R}^k)$ -many components.
- If the position z is changed slightly to z + h,

$$e(z+h) \approx e(z) + \sum_{i=1}^{k} ((\nabla e)(z))_{i} \cdot h_{i}$$

- For $u: \Omega \to \mathbb{R}$: $(\nabla E)(u)$ has $(\dim \{\hat{u}: \Omega \to \mathbb{R}\})$ -many components.
- If the image u is changed slightly in each pixel to u(x, y) + h(x, y), then

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$$E(u+h) \approx E(u) + \int_{\Omega} ((\nabla E)(u))(x,y) \cdot h(x,y) dx dy$$

Idea

- The gradient is the direction of steepest increase of *E*.
- The *negative* gradient is the direction is *steepest descent*.

$$\partial_t u = -(\nabla E)(u)$$

$$(u_{\text{new}})(x,y) = u(x,y) + \tau \left(-\left(\nabla E(u)\right)(x,y)\right)$$

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Gradient descent equation

$$\partial_t u = -(\nabla E)(u)$$

So, having computed some candidate u with energy E(u), we can construct a better candidate u_{new} with a *potentially lower* energy $E(u_{\text{new}})$:

$$(u_{\text{new}})(x,y) = u(x,y) + \tau \left(-\left(\nabla E(u)\right)(x,y)\right)^{T}$$



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Functional derivative

$$(\nabla E)(u) = 2(u - f) - \lambda \operatorname{div}\left(\frac{\phi'(|\nabla u|)}{|\nabla u|}\nabla u\right)$$

Gradient descent equation

$$\partial_t u = -(\nabla E)(u) = 2(f - u) + \lambda \operatorname{div}\left(\frac{\phi'(|\nabla u|)}{|\nabla u|}\nabla u\right)$$

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Gradient Descent: Quadratic Regularizer Example

Quadratic regularizer: Set $\phi(s) := \frac{1}{2}s^2$.

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Using this regularizer leads to oversmoothing, solutions are too blurry.

Gradient descent equation

We have $\frac{\phi'(s)}{s} = 1$, therefore

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$$\phi(s) := h_{\varepsilon}(s) := \begin{cases} \frac{s^2}{2\varepsilon} & \text{if } s < \varepsilon \\ s - \frac{\varepsilon}{2} & \text{else} \end{cases}$$
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Idea

Setting the gradient to zero, i.e. considering $(\nabla E)(u) = 0$, yields a *necessary* optimality condition for the minimizers u.

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Euler-Lagrange Equation: Discretization

Forward differences for the diffusivity $g := \widehat{g}(|\nabla^+ u|), \ \widehat{g}(s) := \frac{\phi'(s)}{s}$. Forward differences for ∇ , backward differences for div :

$$2(\mathbf{u} - \mathbf{f}) - \lambda \operatorname{div}^{-} (\mathbf{g} \nabla^{+} \mathbf{u}) = 0.$$

$$2(u-f) - \lambda \left(g_r u(x+1,y) + g_l u(x-1,y) + g_u u(x,y+1) + g_d u(x,y-1) - (g_r + g_l + g_d) u(x,y) \right) = 0$$

$$g_r := \mathbf{1}_{x+1 < W} \cdot g(x, y), \qquad g_l := \mathbf{1}_{x>0} \cdot g(x-1, y), \\ g_u := \mathbf{1}_{y+1 < H} \cdot g(x, y), \qquad g_d := \mathbf{1}_{y>0} \cdot g(x, y-1).$$



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This is a nonlinear equations system. Use a fixed point iteration scheme.

- 1 Start with an image u^0 .

$$(2 + \lambda(g_r + g_l + g_u + g_d)) u^{k+1}(x, y)$$

$$- \lambda g_r u^{k+1}(x+1, y) - \lambda g_l u^{k+1}(x-1, y)$$

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Iterate until convergence.



Linear Equation Systems: Jacobi Method

Jacobi Method

To solve Az = b: split A = D + R with diagonal D and off-diagonal R:

$$D = \begin{pmatrix} a_{11} & 0 & \cdots & 0 \\ 0 & a_{22} & & \vdots \\ \vdots & & \ddots & 0 \\ 0 & \cdots & 0 & a_{nn} \end{pmatrix}, R = \begin{pmatrix} 0 & a_{12} & \cdots & a_{1n} \\ a_{21} & 0 & & \vdots \\ \vdots & & \ddots & a_{n-1,n} \\ a_{n1} & \cdots & a_{n,n-1} & 0 \end{pmatrix}$$

$$\mathbf{z}_{i}^{k+1} = \frac{1}{\mathbf{a}_{ii}} \Big(\mathbf{b}_{i} - \sum_{i \neq i} \mathbf{a}_{ij} \mathbf{z}_{j}^{k} \Big)$$

$$U^{k+1}(X,Y) = \frac{2f(x,y) + \lambda g_r u^k(x+1,y) + \lambda g_l u^k(x-1,y) + \lambda g_u u^k(x,y+1) + \lambda g_u u^k(x,y-1)}{2 + \lambda (g_r + g_l + g_u + g_d)}$$



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Update for the Euler-Lagrange equation

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Split $A = L_* + U$, with L_* lower triangular and U upper triangular:

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To parallelize the Gauss-Seidel update: *First*: update only at pixels (x, y) with (x + y)%2 = 0. *Then*: only with (x + y)%2 = 1.



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Accelerates the Gauss-Seidel method by linear extrapolation.

SOR update step

Let \bar{z}^{k+1} be the result of one Gauss-Seidel iteration applied to z^k .

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where $\theta \in [0, 1)$ is a fixed parameter.

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SOR converges for any $\theta \in [0, 1)$. The optimal θ depends on A. In practice, one uses values near 1, typically 0.5-0.9, or 0.9-0.98

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