Local, Unconstrained Function Optimization

COMPSCI 527 — Computer Vision

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Outline

- Motivation and Scope
- 2 First Order Methods
- Gradient Descent
- 4 Step Size Selection Methods
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- Stochastic Gradient Descent

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Motivation and Scope

- Most estimation problems are solved by optimization
- Machine learning:
 - Parametric predictor: $h(\mathbf{x}; \mathbf{v}) : \mathbb{R}^d \times \mathbb{R}^m \to Y$
 - Training set $T = \{(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_N, y_N)\}$ and $loss = \ell(y_n, y)$ Risk: $L_T(\mathbf{v}) = \frac{1}{N} \sum_{n=1}^N \ell(y_n, h(\mathbf{x}_n; \mathbf{v})) : \mathbb{R}^m \to \mathbb{R}$

 - Training: $\hat{\mathbf{v}} = \arg\min_{\mathbf{v} \in \mathbb{R}^m} L_T(\mathbf{v})$
- 3D Reconstruction:
 - Computer Graphics: $I = \pi(C, S)$ where I are (multiple) images, C are the camera positions and orientations, S is scene shape
 - Computer Vision: Given I, find $\hat{C}, \hat{S} = \arg\min_{C,S} \|I - \pi(C,S)\|$
- In general, "solving" the system of equations E(z) = 0 can be viewed as

 $\hat{\mathbf{z}} = \arg\min_{\mathbf{z}} \|E(\mathbf{z})\|$

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Only Local Minimization

 $\hat{\mathbf{z}} = \arg\min_{\mathbf{z}\in\mathbf{?}} f(\mathbf{z})$

- All we know about *f* is a "black box" (think Python function)
- For many problems, f has many local minima
- Start somewhere (z₀), and take steps "down"
 f(z_{k+1}) < f(z_k)
- When we get stuck at a local minimum, we declare success
- · We would like global minima, but all we get is local ones
- For some problems, f has a unique minimum...
- ... or at least a single connected set of minima

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Gradient

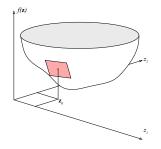
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- We worked with gradients for the case $\mathbf{z} \in \mathbb{R}^2$ (images)
- Now $\mathbf{z} \in \mathbb{R}^m$ with *m* possibly very large
- If ∇f(z) exists everywhere, the condition ∇f(z) = 0 is necessary and sufficient for a stationary point (max, min, or saddle)
- Warning: only *necessary* for a minimum!
- Reduces to first derivative when $f : \mathbb{R} \to \mathbb{R}$

First Order Taylor Expansion

 $f(\mathbf{z}) \approx g_1(\mathbf{z}) = f(\mathbf{z}_0) + [\nabla f(\mathbf{z}_0)]^T (\mathbf{z} - \mathbf{z}_0)$

approximates $f(\mathbf{z})$ near \mathbf{z}_0 with a (hyper)plane through \mathbf{z}_0



 $\nabla f(\mathbf{z}_0)$ points to direction of steepest *increase* of *f* at \mathbf{z}_0

- If we want to find z₁ where f(z₁) < f(z₀), going along
 −∇f(z₀) seems promising
- This is the general idea of gradient descent

A Template

Gradient descent methods fit the following template:

Design Decisions

k = 0

while \mathbf{z}_k is not a minimum compute the gradient \mathbf{g}_k compute a *learning rate* $\alpha_k > \mathbf{0}$

$$\mathbf{z}_{k+1} = \mathbf{z}_k - \alpha_k \mathbf{g}_k$$
$$k = k + 1$$

end

- In what direction to proceed (-g_k)
- How long a step to take in that direction (α_k ||g_k||)
- When to stop ("while z_k is not a minimum")
- Different decisions lead to different methods

Gradient Descent

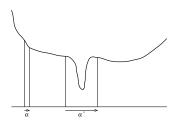
- In what direction to proceed: $-\mathbf{g}_k = -\nabla f(\mathbf{z}_k)$
- "Gradient descent"
- Problem reduces to one dimension:
 h(α) = f(z_k αg_k)
- $\alpha = \mathbf{0} \Leftrightarrow \mathbf{z} = \mathbf{z}_k$
- Find $\alpha = \alpha_k > 0$ such that $f(\mathbf{z}_k \alpha_k \mathbf{g}_k) < f(\mathbf{z}_k)$
- How to find α_k ?

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Step Size

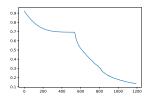
- Simplest idea: $\alpha_k = \alpha$ (fixed learning rate)
 - Small α leads to slow progress
 - Large α can miss minima



- Scheduling α :
 - Start with α relatively large (say $\alpha = 10^{-3}$)
 - Decrease α over time
 - Determine decrease rate of α by trial and error

Momentum

• Sometimes **z**_k meanders around in shallow valleys



$$f(\mathbf{z}_k)$$
 versus k

- α is too small, direction is still promising
- Add momentum

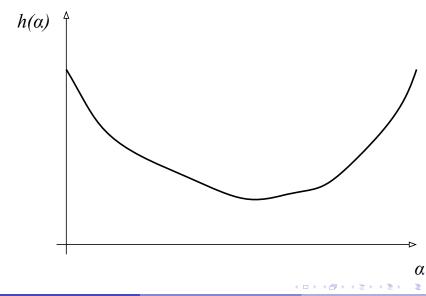
$$\begin{aligned} \mathbf{v}_0 &= \mathbf{0} \\ \mathbf{v}_{k+1} &= \mu_k \mathbf{v}_k - \alpha \nabla f(\mathbf{z}_k) \\ \mathbf{z}_{k+1} &= \mathbf{z}_k + \mathbf{v}_{k+1} \end{aligned} \qquad (\mathbf{0} \leq \mu_k < \mathbf{1} \end{aligned}$$

Line Search

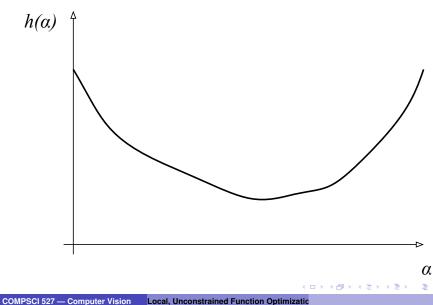
- Find a local minimum in the search direction **p**_k = -**g**_k
 h(α) = f(**z**_k + α**p**_k), a one-dimensional problem
- Bracketing triple:
- a < b < c, $h(a) \ge h(b)$, $h(b) \le h(c)$
- Contains a (local) minimum!
- Split the bigger of [a, b] and [b, c] in half with a point u
- Find a new, narrower bracketing triple involving *u* and two out of *a*, *b*, *c*
- Stop when the bracket is narrow enough (say, 10⁻⁶)
- Pinned down a minimum to within 10⁻⁶

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Phase 1: Find a Bracketing Triple



Phase 2: Shrink the Bracketing Triple



if
$$b - a > c - b$$

 $u = (a + b)/2$
if $h(u) > h(b)$
 $(a, b, c) = (u, b, c)$
otherwise
 $(a, b, c) = (a, u, b)$
end
otherwise
 $u = (b + c)/2$
if $h(u) > h(b)$
 $(a, b, c) = (a, b, u)$
otherwise
 $(a, b, c) = (b, u, c)$
end
end

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Termination

- Are we still making "significant progress"?
- Check $f(\mathbf{z}_{k-1}) f(\mathbf{z}_k)$? (We want this to be strictly positive)
- Check $\|\mathbf{z}_{k-1} \mathbf{z}_k\|$? (We want this to be large enough)
- Second is more stringent close the the minimum because ∇f(z) ≈ 0

• Stop when
$$\|\mathbf{z}_{k-1} - \mathbf{z}_k\| < \delta$$

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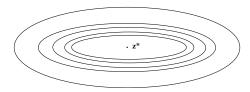
Is Gradient Descent a Good Strategy?

- "We are going in the direction of fastest descent"
- "We choose an optimal step size by line search"
- "Must be good, no?" Not so fast!
- An example for which we know the answer:

$$f(\mathbf{z}) = \mathbf{c} + \mathbf{a}^T \mathbf{z} + \frac{1}{2} \mathbf{z}^T Q \mathbf{z}$$

 $Q \succcurlyeq 0$ (convex paraboloid)

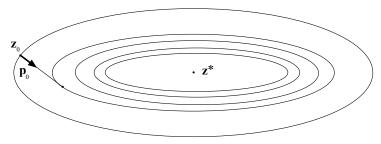
All smooth functions look like this close enough to z*



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Skating to a Minimum



- Many 90-degree turns slow down convergence
- There are methods that take fewer iterations, but each iteration takes more time and space
- We will stick to gradient descent
- See appendices in the notes for more efficient methods for problems in low-dimensional spaces

Stochastic Gradient Descent

• A special case of gradient descent, SGD works for *averages* of many terms (*N* very large):

$$f(\mathbf{z}) = \frac{1}{N} \sum_{n=1}^{N} \phi_n(\mathbf{z})$$

- Computing $\nabla f(\mathbf{z}_k)$ is too expensive
- Partition B = {1,..., N} into J random mini-batches B_j each of about equal size

$$f(\mathbf{z}) \approx f_j(\mathbf{z}) = rac{1}{|B_j|} \sum_{n \in B_j} \phi_n(\mathbf{z}) \quad \Rightarrow \quad \nabla f(\mathbf{z}) \approx \nabla f_j(\mathbf{z}) \;.$$

Mini-batch gradients are correct on average

SGD and Mini-Batch Size

- SGD iteration: $\mathbf{z}_{k+1} = \mathbf{z}_k \alpha_k \nabla f_j(\mathbf{z}_k)$
- Mini-batch gradients are correct on average
- One cycle through all the mini-batches is an epoch
- Repeatedly cycle through all the data (Scramble data before each epoch)
- *Asymptotic* convergence can be proven with suitable step-size schedule
- Small batches \Rightarrow low storage but high gradient variance
- Make batches as big as will fit in memory for minimal variance
- In deep learning, memory is GPU memory

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