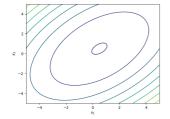
Optimization in Machine Learning

First order methods GD on quadratic forms





Learning goals

- Eigendecomposition of quadratic forms
- GD steps in eigenspace

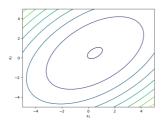
QUADRATIC FORMS & GD

- We consider the quadratic function $q(\mathbf{x}) = \mathbf{x}^{\top} \mathbf{A} \mathbf{x} \mathbf{b}^{\top} \mathbf{x}$.
- We assume that Hessian $\mathbf{H} = 2\mathbf{A}$ has full rank
- We assume q is convex, so A psd.
- Optimal solution is $\mathbf{x}^* = \frac{1}{2}\mathbf{A}^{-1}\mathbf{b}$
- As $\nabla q(\mathbf{x}) = 2\mathbf{A}\mathbf{x} \mathbf{b}$, iterations of gradient descent are

$$\mathbf{x}^{[t+1]} = \mathbf{x}^{[t]} - \alpha (2\mathbf{A}\mathbf{x}^{[t]} - \mathbf{b})$$



QUADRATIC FORMS & GD / 2

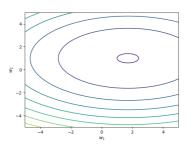




The following slides follow the blog post "Why Momentum Really Works", Distill, 2017. http://doi.org/10.23915/distill.00006

EIGENDECOMPOSITION OF QUADRATIC FORMS

- We want to work in the coordinate system given by q
- ullet Recall: Coordinate system is given by the eigenvectors of ${f H}=2{f A}$
- ullet Eigendecomposition of ${f A} = {f V} {f \Lambda} {f V}^{ op}$
- ullet V contains eigenvectors $oldsymbol{v}_i$ and $oldsymbol{\Lambda}= ext{diag}(\lambda_1,...,\lambda_n)$ eigenvalues
- ullet Change of basis: $\mathbf{w}^{[t]} = \mathbf{V}^{\top} (\mathbf{x}^{[t]} \mathbf{x}^*)$





GD STEPS IN EIGENSPACE

With
$$\mathbf{w}^{[t]} = \mathbf{V}^{\top}(\mathbf{x}^{[t]} - \mathbf{x}^*)$$
, a single GD step

$$\mathbf{x}^{[t+1]} = \mathbf{x}^{[t]} - \alpha (2\mathbf{A}\mathbf{x}^{[t]} - \mathbf{b})$$

becomes

$$\mathbf{w}^{[t+1]} = \mathbf{w}^{[t]} - 2\alpha \mathbf{\Lambda} \mathbf{w}^{[t]}.$$



GD STEPS IN EIGENSPACE / 2

Proof (for $\mathbf{w}^{[t+1]} = \mathbf{w}^{[t]} - 2\alpha \Lambda \mathbf{w}^{[t]}$):

A single GD step means

$$\mathbf{x}^{[t+1]} = \mathbf{x}^{[t]} - \alpha (2\mathbf{A}\mathbf{x}^{[t]} - \mathbf{b})$$

Then:

$$\mathbf{V}^{\top}(\mathbf{x}^{[t+1]} - \mathbf{x}^*) = \mathbf{V}^{\top}(\mathbf{x}^{[t]} - \mathbf{x}^*) - \alpha \mathbf{V}^{\top}(2\mathbf{A}\mathbf{x}^{[t]} - \mathbf{b})$$

$$\mathbf{w}^{[t+1]} = \mathbf{w}^{[t]} - \alpha \mathbf{V}^{\top}(2\mathbf{A}\mathbf{x}^{[t]} - \mathbf{b})$$

$$\mathbf{w}^{[t+1]} = \mathbf{w}^{[t]} - \alpha \mathbf{V}^{\top}(2\mathbf{A}(\mathbf{x}^{[t]} - \mathbf{x}^*) + 2\mathbf{A}\mathbf{x}^* - \mathbf{b})$$

$$= \mathbf{w}^{[t]} - 2\alpha \Lambda \mathbf{V}^{\top}(\mathbf{x}^{[t]} - \mathbf{x}^*)$$

$$= \mathbf{w}^{[t]} - 2\alpha \Lambda \mathbf{w}^{[t]}$$



GD STEPS IN EIGENSPACE

Therefore:

$$w_{i}^{[t+1]} = w_{i}^{[t]} - 2\alpha\lambda_{i}w_{i}^{[t]}$$

$$= (1 - 2\alpha\lambda_{i})w_{i}^{[t]}$$

$$= \cdots$$

$$= (1 - 2\alpha\lambda_{i})^{t+1}w_{i}^{[0]}$$



GD ERROR IN ORIGINAL SPACE

• Move back to original space:

$$\mathbf{x}^{[t]} - \mathbf{x}^* = \mathbf{V}\mathbf{w}^{[t]} = \sum_{i=1}^d (1 - 2\alpha\lambda_i)^t w_i^{[0]} \mathbf{v}_i$$

- This gives us the difference in x-space in closed form
- The difference vector is written in the eigen-basis, the initial error components $w_i^{[0]}$ decay with rate $1 2\alpha\lambda_i$,
- The smaller this term (in abs value), the faster we decay, the closer to 1, the slower we are
- For most step-sizes, eigenvectors with largest eigenvalues converge fastest. Works especially well in the first iterations.

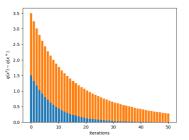


GD ERROR IN ORIGINAL SPACE / 2

We can also analyze the regret in y-space:

$$q(\mathbf{x}^{[t]}) - q(\mathbf{x}^*) = \frac{1}{2} \sum_{i}^{d} (1 - 2\alpha \lambda_i)^{2t} \lambda_i (w_i^{[0]})^2$$

- This is not hard to show, formulas become only somewhat lengthy
- Rates become somewhat equalized at the end





OPTIMAL STEPSIZE

- In order to converge, must have all : $|1 2\alpha\lambda_i| < 1$
- So for all stepsizes $0 < \alpha \lambda_i < 1$
- Overall rate determined by slowest error component:

$$\max_{i=1,...n}|1-2\alpha\lambda_i|=\max(|1-2\alpha\lambda_1|,|1-2\alpha\lambda_n|)$$

- Minimized for $\alpha = \frac{1}{\lambda_1 + \lambda_n}$
- Optimal rate $\frac{\lambda_n/\lambda_1-1}{\lambda_n/\lambda_1+1}$
- We again see that convergence is determined by condition

$$\kappa = \frac{\lambda_n}{\lambda_1}$$

• $\kappa = 1$ Is ideal, convergence in 1 step

