Introduction to Deep Learning



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Regularization

Introduction

- In machine learning, target is to make an algorithm performs well not only on training data but also on new data
- Many strategies exist to reduce test error at the cost of training error
- Any modification we make to a learning algorithm that is intended to reduce its generalization error but not its training error
- Objectives
 - To encode prior knowledge
 - Constraints and penalties are designed to express generic preference for simpler model

Regularization in DL

- In DL regularization works by trading increased bias for reduced variance
- Consider the following scenario
 - Excluded the true data generating process
 - Underfitting, inducing bias
 - Matched the true data generating process
 - Desired one
 - Included the generating process but also many other generating process
 - Overfitting, variance dominates
 - Goal of regularizer is to take an model overfit zone to desired zone

Norm penalties

- Most of the regularization approaches are based on limiting the capacity of the model
- Objective function becomes $\tilde{J}(\boldsymbol{\theta}; \boldsymbol{X}, \boldsymbol{y}) = J(\boldsymbol{\theta}; \boldsymbol{X}, \boldsymbol{y}) + \alpha \Omega(\boldsymbol{\theta})$
 - α Hyperparameter denotes relative contribution
 - Minimization of \tilde{J} implies minimization of J
 - $\boldsymbol{\Omega}$ penalizes only the weight of affine transform
 - Bias remain unregularized
 - Regularizing bias may lead to underfitting

- Weights are closer to origin as $\Omega(oldsymbol{ heta}) = rac{1}{2} \|oldsymbol{w}\|_2^2$
 - Also known as ridge regression or Tikhonov regression
- Objective function $\widetilde{J}(\boldsymbol{w}; \boldsymbol{X}, \boldsymbol{y}) = \frac{\alpha}{2} \boldsymbol{w}^{T} \boldsymbol{w} + J(\boldsymbol{w}; \boldsymbol{X}, \boldsymbol{y})$

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- New weights

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$$\boldsymbol{w} = \boldsymbol{w} - \epsilon(\alpha \boldsymbol{w} + \nabla_{\boldsymbol{w}} J(\boldsymbol{w}; \boldsymbol{X}, \boldsymbol{y})) \\ = \boldsymbol{w}(1 - \epsilon\alpha) - \epsilon \nabla_{\boldsymbol{w}} J(\boldsymbol{w}; \boldsymbol{X}, \boldsymbol{y})$$

- Assume quadratic nature of curve in the neighborhood of $w^* = \arg \min J(w)$
 - **J(w)** unregularized cost
 - Perfect scenario for linear regression with MSE

Jacobian & Hessian

- Derivative of a function having single input and single output $-\frac{dy}{dx}$
- Derivative of function having vector input and vector output that is, $f : \mathbb{R}^m \to \mathbb{R}^n$
 - Jacobian $\mathbf{J} \in \mathbb{R}^{n \times m}$ of f defined as $J_{i,j} = \frac{\partial}{\partial x_i} f(\mathbf{x})_i$
- Second derivative is also required sometime
 - For example, $f : \mathbb{R}^n \to \mathbb{R}, \ \frac{\partial^2}{\partial x_i \partial x_i} f$
 - If second derivative is 0, then there is no curvature
- Hessian matrix $H(f)(x)_{ij} = \frac{\partial^2}{\partial x_i \partial x_j} f(x)$

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 - Jacobian of gradient
 - Symmetric

Directional derivative

The directional derivative of a scalar function f(x) = f(x₁, x₂,..., x_n) along a vector
 v = (v₁,..., v_n) is given by

$$\nabla_{\mathbf{v}}f(\mathbf{x}) = \lim_{h \to 0} \frac{f(\mathbf{x} + h\mathbf{v}) - f(\mathbf{x})}{h}$$

• If f is differentiable at point x then

 $abla_{\mathbf{v}}f(\mathbf{x}) =
abla f(\mathbf{x}) \cdot \mathbf{v}$

Taylor series expansion

• A real valued function differentiable at point x_0 can be expressed as

$$f(x) = f(x_0) + \frac{f'(x_0)}{1!}(x - x_0) + \frac{f''(x_0)}{2!}(x - x_0)^2 + \frac{f^{(3)}(x_0)}{3!}(x - x_0)^3 + \cdots$$

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When input is a vector

$$f(\pmb{x}) pprox f(\pmb{x}^{(0)}) + (\pmb{x} - \pmb{x}^{(0)})^{ op} \pmb{g} + rac{1}{2} (\pmb{x} - \pmb{x}^{(0)})^{ op} \pmb{H}(\pmb{x} - \pmb{x}^{(0)})$$

• g — gradient at $x^{(0)}$, H — Hessian at $x^{(0)}$

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• If ϵ is the learning rate, then $f(x^{(0)} - \epsilon g) = f(x^{(0)}) - \epsilon g^T g + \frac{1}{2} \epsilon^2 g^T H g$

Quadratic approximation

- Let $w^* = \arg \min_{w} J(w)$ optimum weights for minimal unregularized cost
- If the objective function is quadratic then $\hat{J}(\theta) = J(w^*) + \frac{1}{2}(w w^*)^T H(w w^*)$
 - **H** is the Hessian matrix of **J** with respect to **w** at **w***
 - No first order term as w^{*} is minimum
 - **H** is positive semidefinite
- Minimum of \hat{J} occurs when $\nabla_{w}\hat{J}(w) = H(w w^{*}) = 0$
- With weight decay we have

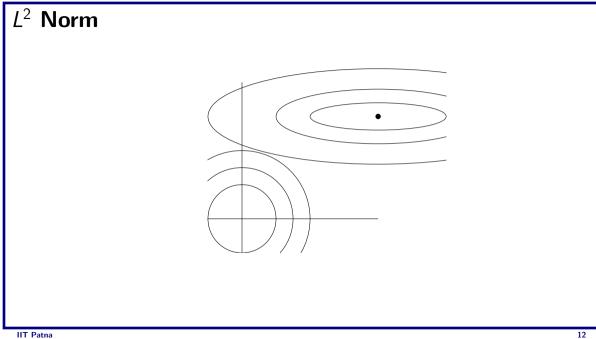
 $\alpha \tilde{\boldsymbol{w}} + \boldsymbol{H}(\tilde{\boldsymbol{w}} - \boldsymbol{w}^*) = 0$ $(\boldsymbol{H} + \alpha \boldsymbol{I}) \tilde{\boldsymbol{w}} = \boldsymbol{H} \boldsymbol{w}^*$ $\tilde{\boldsymbol{w}} = (\boldsymbol{H} + \alpha \boldsymbol{I})^{-1} \boldsymbol{H} \boldsymbol{w}^*$

Quadratic approximation (contd)

- As $\alpha \rightarrow 0$, regularized solution \hat{w} approaches to w^*
- As $\alpha \to \infty$
 - **H** is symmetric, therefore $H = Q \Lambda Q^T$. Now we have

$$\widetilde{\boldsymbol{w}} = (\boldsymbol{Q}\boldsymbol{\Lambda}\boldsymbol{Q} + \alpha\boldsymbol{I})^{-1}\boldsymbol{Q}\boldsymbol{\Lambda}\boldsymbol{Q}^{\mathsf{T}}\boldsymbol{w}^{*} \\ = [\boldsymbol{Q}(\boldsymbol{\Lambda} + \alpha\boldsymbol{I})\boldsymbol{Q}^{\mathsf{T}}]^{-1}\boldsymbol{Q}\boldsymbol{\Lambda}\boldsymbol{Q}^{\mathsf{T}}\boldsymbol{w}^{*} \\ = \boldsymbol{Q}(\boldsymbol{\Lambda} + \alpha\boldsymbol{I})^{-1}\boldsymbol{\Lambda}\boldsymbol{Q}^{\mathsf{T}}\boldsymbol{w}^{*}$$

- Weight decay rescale w^* along the eigen vector of H
 - Component of w^* that is aligned to i-th eigen vector, will be rescaled by a factor of $\frac{\lambda_i}{\lambda_i + \alpha}$
 - $\lambda_i \gg \alpha$ regularization effect is small



Linear regression

- For linear regression cost function is $(Xw y)^T (Xw y)$
- Using L^2 regularization we have $(\mathbf{X}\mathbf{w} \mathbf{y})^T (\mathbf{X}\mathbf{w} \mathbf{y}) + \frac{1}{2}\alpha \mathbf{w}^T \mathbf{w}$

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L¹ regularization

- Formally it is defined as $\Omega(\boldsymbol{\theta}) = \|\boldsymbol{w}\|_1 = \sum |w_i|$
- Regularized objective function will be $\tilde{J}(w; X, y) = \alpha ||w||_1 + J(w; X, y)$

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- Regularized objective function will be $\tilde{J}(w; X, y) = \alpha ||w||_1 + J(w; X, y)$
- The gradient will be $\nabla_{w} \widetilde{J}(w; X, y) = \alpha \operatorname{sign}(w) + \nabla_{w} J(w; X, y)$
 - Gradient does not scale linearly compared to L^2 regularization
- Taylor series exapansion with approximation provides $\nabla_{w} \hat{J}(w) = H(w w^{*})$
- Simplification can be made by assuming *H* to be diagonal
 - Apply PCA on the input dataset

L¹ regularization

- Quadratic approximation of L^1 regularization objective function becomes $\hat{J}(w; \mathbf{X}, \mathbf{y}) = J((w^*; \mathbf{X}, \mathbf{y}) + \sum_i \left[\frac{1}{2}H_{i,i}(w_i w_i^*)^2 + \alpha |w_i|\right]$
- So, analytical in each dimension will be $w_i = \operatorname{sign}(w_i^*) \max \left\{ |w_i^*| \frac{\alpha}{H_{i,i}}, 0 \right\}$
- Consider the situation when $w_i^* > 0$
 - If $w_i^* \leq \frac{\alpha}{H_{i,i}}$, optimal value for w_i will be 0 under regularization
 - If $w_i^* > \frac{\alpha}{H_{i,i}}$, w_i moves towards 0 with a distance equal to $\frac{\alpha}{H_{i,i}}$

Constrained optimization

• Cost function regularized by norm penalty is given by

 $\widetilde{J}(oldsymbol{ heta};oldsymbol{X},oldsymbol{y}) = J(oldsymbol{ heta};oldsymbol{X},oldsymbol{y}) + lpha \Omega(oldsymbol{ heta})$

• Let us assume f(x) needs to be optimized under a set of equality constraints $g^{(i)}(x) = 0$ and inequality constraints $h^{(j)}(x) \le 0$, then generalized Lagrangian is then defined as

$$\mathcal{L}(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{lpha}) = f(\mathbf{x}) + \sum_{i} \lambda_{i} g^{(i)}(\mathbf{x}) + \sum_{j} \alpha_{i} g^{(i)}(\mathbf{x})$$

• If there exists a solution then

$$\min_{x} \max_{\boldsymbol{\lambda}} \max_{\boldsymbol{\alpha} \geq 0} = L(x, \boldsymbol{\lambda}, \boldsymbol{\alpha}) = \min_{x} f(x)$$

• This can be solved by $\nabla_{\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\alpha}} L(\mathbf{x}, \boldsymbol{\lambda}, \boldsymbol{\alpha}) = 0$

Constraint optimization (contd.)

• Suppose $\Omega(\theta) < k$ needs to be satisfied. Then regularization equation becomes

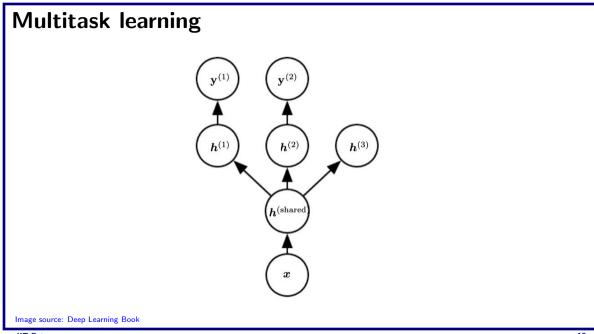
 $L(\boldsymbol{\theta}, \alpha; \boldsymbol{X}, \boldsymbol{y}) = J(\boldsymbol{\theta}; \boldsymbol{X}, \boldsymbol{y}) + \alpha(\Omega(\boldsymbol{\theta}) - \boldsymbol{k})$

• Solution to the constrained problem

$$oldsymbol{ heta}^* = rg\min_{oldsymbol{ heta}} \max_{lpha > 0} L(oldsymbol{ heta}, lpha)$$

Dataset augmentation

- If data are limited, fake data can be added to training set
 - Computer vision problem
 - Speech recognition
- Easiest for classification problem
- Very effective in object recognition problem
 - Translating
 - Rotating
 - Scaling
 - Need to be careful for 'b' and 'd' or '6' and '9'
- Injecting noise to input data can be viewed as data augmentation



Early stopping

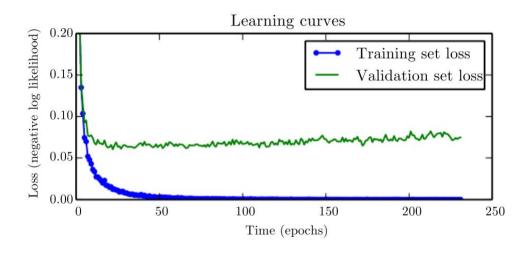


Image source: Deep Learning Book

Early stopping approach

- Intialize the parameters
- Run training algorithm for *n* steps and update i = i + n
- Compute error on the validation set (v')
- If v' is less than previous best, then update the same. Start step 2 again
- If v' is more than the previous best, then increament the count that stores the number of such occurrences. If the count is less than a threshold go to step 2, otherwise exit.

Early stopping (contd)

- Number of tarining step is a hyperparameter
 - Most hyperparameters that control model capacity have U-shaped curve
- Additional cost for this approach is to store the parameters
- Requires a validation set
 - It will have two passes
 - First pass uses only training data for update of the parameters
 - Second pass uses both training and validation data for update of the parameters
 - Possible strategies
 - Intialize the model again, retrain on all data, train for the same number of steps as obtained by early stopping in pass 1
 - Keep the parameters obtained from the first round, continue training using all data until the validation loss is than the training loss at the early stopping point
- It reduces computational cost as it limits the number of iteration
- Provides regularization without any penalty

- Let us assume au training iteration, ϵ learning rate
 - $\epsilon \tau$ measures effective capacity
- We have, $\hat{J}(\theta) = J(w^*) + \frac{1}{2}(w w^*)H(w w^*)$ and $\nabla_w \hat{J}(w) = H(w w^*)$
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$$Q^{T}w^{(\tau)} = [I - (I - \epsilon \Lambda)^{\tau}]Q^{T}w^{*}$$

Early stopping as regularizer (contd)

- Assuming $w^{(0)} = 0$ and ϵ is small value such that $|1 \epsilon \lambda_i| < 1$
- From L^2 regularization, we have

$$Q^{\mathsf{T}} \tilde{w} = (\Lambda + \alpha I)^{-1} \Lambda Q^{\mathsf{T}} w^*$$
$$Q^{\mathsf{T}} \tilde{w} = [I - (\Lambda + \alpha I)^{-1} \alpha] Q^{\mathsf{T}} w^*$$

- Therefore we have, $(I \epsilon \Lambda)^{\tau} = (\Lambda + \alpha I)^{-1} \alpha$
- Hence, $\tau \approx \frac{1}{\epsilon \alpha} \ \alpha \approx \frac{1}{\tau \epsilon}$

Bagging

- Also known as Bootstrap aggregating
- Reduces generalization error by combining several models
- Train multiple models then vote on output for the test example
 - Also known as model averaging, ensemble method
- Suppose we have k regression model and each model makes an error ϵ_i such that $\mathbb{E}(\epsilon_i) = 0$, $\mathbb{E}(\epsilon_i^2) = v$, $\mathbb{E}(\epsilon_i \epsilon_j) = c$
- Error made by average prediction $\frac{1}{k}\sum \epsilon_i$
- Expected mean square error

$$\mathbb{E}\left[\left(\frac{1}{k}\sum_{i}\epsilon_{i}\right)^{2}\right] = \frac{1}{k^{2}}\mathbb{E}\left[\sum_{i}\left(\epsilon_{i}^{2}+\sum_{i\neq j}\epsilon_{i}\epsilon_{j}\right)\right] = \frac{v}{k} + \frac{k-1}{k}\epsilon_{i}\epsilon_{j}$$

Dropout

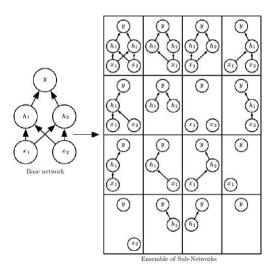
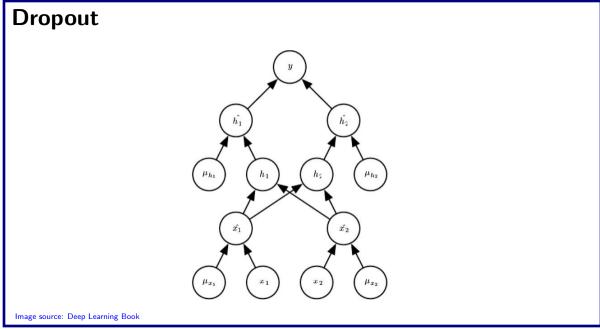


Image source: Deep Learning Book



Adversarial training

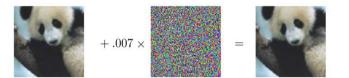


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