

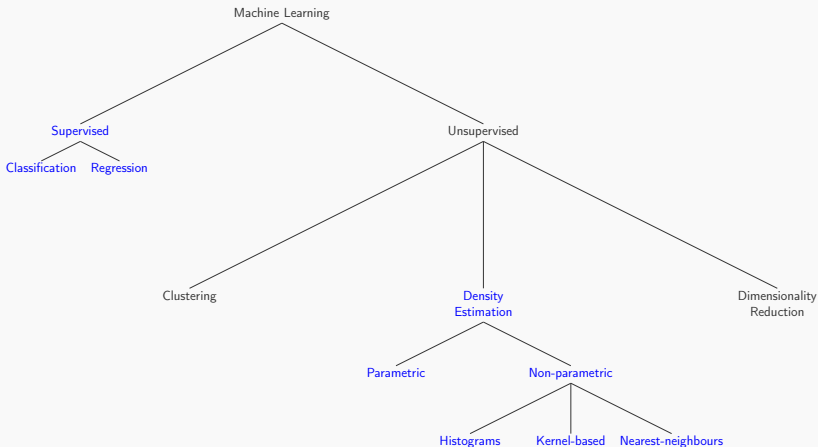
CS-667 Advanced Machine Learning

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Neural Networks

Machine Learning So Far ...



Assignment 1

Iterative Reweighted Least Squares for Logistic Regression

- ▶ Implement the IRLS algorithm for logistic regression.
 - ▶ Code up a generic implementation.
 - ▶ Train it on the first 2 classes of MNIST digits training data.
 - ▶ Relevant material has been placed on `\\printsrv`.
 - ▶ Each sample is a 784×1 vector that represents a 28×28 image. To visualise the k -th training sample as an image, you may use the following commands:

```
imagesc(reshape(train_x(k,:),28,28)');  
axis image;  
colormap gray;
```
 - ▶ Report classification accuracy and confusion matrix on the testing data for the relevant classes.
- ▶ Submit your `_roll_number_LR.zip` containing code and `report.txt/pdf` explaining your results.
- ▶ Due next Tuesday (March 07, 2017 before 5:30 pm) on `\\printsrv`.

Assignment 2

SGD for Multiclass Logistic Regression

- ▶ Implement the SGD algorithm for multiclass logistic regression.
 - ▶ Code up a generic implementation.
 - ▶ Train it on the MNIST digits training data.
 - ▶ Report classification accuracy and confusion matrix on the testing data.
- ▶ Submit your `_roll_number_MLR.zip` containing code and `report.txt/pdf` explaining your results.
- ▶ Due on Tuesday (March 14, 2017 before 5:30 pm) on `\\printsrv`.

Logistic Regression Tips

- ▶ In case you are having memory issues or training takes a lot of time, you might want to use the following tips:
 - ▶ Type in the command 'doc spdiags'
 - ▶ Do not use the `inv()` function to for matrix inverse. Use the `\` operator. For more help, consult Google or Matlab documentation.
- ▶ Also, don't forget to homogenise the inputs by appending a 1 at the end of each input. This will absorb the bias term.
- ▶ Lastly, if you start getting a warning message like "**Warning: Matrix is close to singular or badly scaled**"
 - ▶ First look at the difference between Exercises 1.1 and 1.2 from Chapter 1 and their solutions.
 - ▶ Then look at the programming solutions to both problems. We have done both the exercises as well as their programming solutions in CS 567.

Project

- ▶ Explore an interesting ML problem using a real-world data set.

- ▶ 4 deliverables

Proposal	Tuesday, March 28	1 page
Milestone	Tuesday, May 2	3-4 pages
Poster	Monday, May 29	
Final report	Monday, June 5	6-8 pages (NIPS format)

- ▶ Project ideas

- ▶ <http://cs229.stanford.edu/projects2016.html>
- ▶ <http://cs229.stanford.edu/projects2015.html>
- ▶ <http://cs229.stanford.edu/projects2014.html>
- ▶ www.cs.cmu.edu/~10701/projects.html
- ▶ www.kaggle.com
- ▶ Discuss with me

Proposal

- ▶ Maximum 1 page description containing
 - ▶ Project title
 - ▶ Data set
 - ▶ Project idea (approximately two paragraphs).
 - ▶ Software you will need to write or tool/libraries you will need to learn.
 - ▶ Papers to read. Include 1-3 relevant papers. You will probably want to read at least one of them before submitting your proposal.
 - ▶ Milestone description. What experimental results will you complete before May 2nd?
- ▶ Proposal will be rejected if you do not have the dataset available already.

Milestone

- ▶ Short report of 3-4 pages.
- ▶ Same sections as the final report (introduction, related work, method, experiment, conclusion), with a few sections "under construction".
- ▶ Specifically,
 - ▶ the introduction and related work sections should be in their final form
 - ▶ the section on the proposed method should be almost finished
 - ▶ the sections on the experiments and conclusions will have whatever results you have obtained, as well as "place-holders" for the results you plan/hope to obtain.

Final report

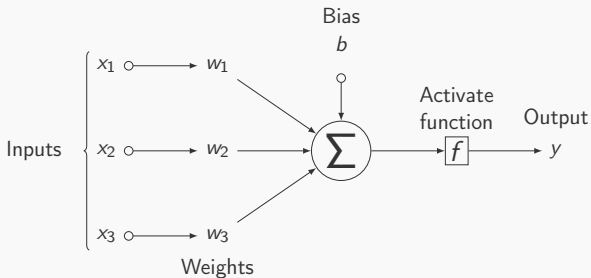
- ▶ Must be in NIPS format and page limit. Use [this](#) link.
- ▶ Think of it as a research paper being submitted to NIPS.

Neural Networks

- ▶ So far, we have learned \mathbf{w}^* for mapping inputs $\mathbf{x} \in \mathbb{R}^D$ to targets \mathbf{t} .
- ▶ Often, working in a transformed space $\phi \in \mathbb{R}^M$ makes it easier to learn the mapping.
- ▶ However, not all mappings are useful for the problem at hand. Is there an optimal mapping ϕ^* ?
- ▶ Neural networks learn the optimal mapping ϕ^* and also the optimal parameters \mathbf{w}^* .

Neural Networks

The Neuron



- ▶ The function of a biological neuron can be modelled as
$$y = f \left(\sum_j w_j x_j + b \right).$$

Neural Networks

The Neuron

- ▶ Model the output of the k -th neuron as
$$y_k = f(a_k) = f(\mathbf{w}_k^T \mathbf{x}) = f\left(\sum_j w_{kj} x_j\right)$$
 where
 - ▶ The x_j constitute values of input signals feeding into the neuron.
 - ▶ The w_{kj} are weights determining the importance given to input x_j by this neuron.
 - ▶ Dot-product $a_k = \sum_j w_{kj} x_j$ is called the *activation*.
 - ▶ $f(\cdot)$ is called the *activation function*. Determines behaviour of the neuron in response to its activation.
- ▶ The perceptron that we studied earlier is a very simple neuron model with f being the step function.

$$f(a) = \begin{cases} 1 & \text{if } a \geq 0 \\ 0 & \text{if } a < 0 \end{cases} \quad (1)$$

Neural Networks

- ▶ The linear models that we covered can be represented as

$$y_k(\mathbf{x}, \mathbf{W}) = f(\mathbf{w}_k^T \phi(\mathbf{x})) = f\left(\sum_{j=0}^M w_{kj} \phi_j(\mathbf{x})\right)$$

where y_k is the k -th output and index j starts from 0 to reflect bias inclusion.

$$f = \begin{cases} \text{identity} & \text{for regression} \\ \text{logistic sigmoid} & \text{for binary classification} \\ \text{softmax} & \text{for multiclass classification} \end{cases}$$

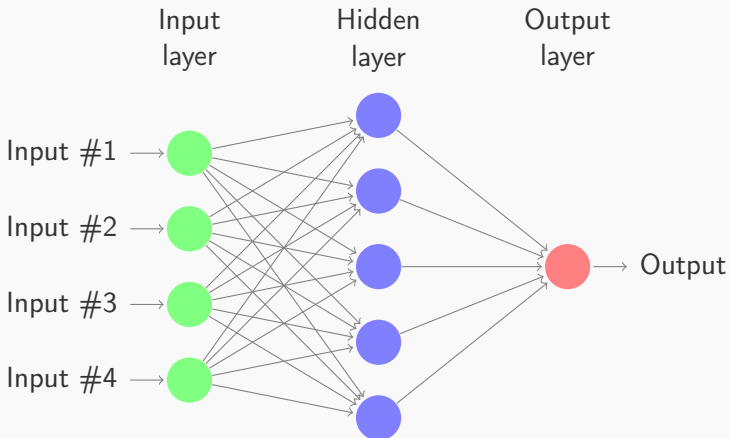
- ▶ Each $\phi_j(\mathbf{x})$ can be seen as a *basis function*.
- ▶ So far, the basis functions were fixed. Now we *adapt* them to the problem.

Neural Networks

- ▶ Model $\phi_j(\mathbf{x})$ as a non-linear function $h(a_j)$ where *activation* $a_j = \mathbf{w}_j^T \mathbf{x}$ with adjustable parameters \mathbf{w}_j .
- ▶ So the k -th output can be written as

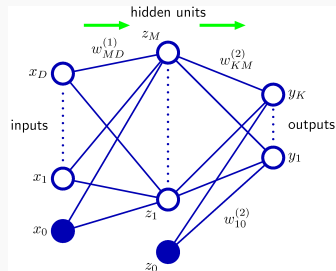
$$\begin{aligned}y_k(\mathbf{x}, \mathbf{W}) &= f(\mathbf{a}_k) = f(\mathbf{w}_k^T \boldsymbol{\phi}(\mathbf{x})) = f\left(\sum_{j=0}^M w_{kj} \phi_j(\mathbf{x})\right) \\ &= f\left(\sum_{j=0}^M w_{kj} h(a_j)\right) \\ &= f\left(\sum_{j=1}^M w_{kj} h\left(\sum_{i=0}^D w_{ji} x_i\right)\right)\end{aligned}$$

Neural Networks



These computations can be visualised graphically as *forward propagation of information* through the so-called *neural network*.

Neural Networks



A two-layer neural network.

- ▶ 3 neuron types:
 - ▶ input x_i
 - ▶ hidden z_j
 - ▶ output y_k
- ▶ 2 weight layers:
 - ▶ hidden-input $w_{ji}^{(1)}$
 - ▶ output-hidden $w_{kj}^{(2)}$
- ▶ To differentiate between different layer parameters, we can write

$$y_k(\mathbf{x}, \mathbf{W}) = f \left(\sum_{j=1}^M w_{kj}^{(2)} h \left(\sum_{i=0}^D w_{ji}^{(1)} x_i \right) \right)$$

Neural Networks

As Multilayer Perceptrons

- ▶ The computation $y_k(\mathbf{x}, \mathbf{W}) = f\left(\sum_{j=1}^M w_{kj}^{(2)} h\left(\sum_{i=0}^D w_{ji}^{(1)} x_i\right)\right)$ can be viewed in two stages:
 1. Compute $z_j = h(\mathbf{w}_j^T \mathbf{x})$, followed by
 2. $y_k = f(\mathbf{w}_k^T \mathbf{z})$.
- ▶ Both stages resemble the perceptron model.
- ▶ Therefore, another name for such neural networks is *multilayer perceptrons* or simply *MLP*.
- ▶ However, there is a key difference:
 - ▶ Perceptron uses a non-differentiable step-function non-linearity.
 - ▶ MLP uses a differentiable sigmoidal non-linearity. So we can train via gradient based approaches.
- ▶ Therefore, despite the name, MLPs never use perceptrons!

Neural Networks

As Universal Approximators

- ▶ Neural networks are considered to be *universal approximators*.
- ▶ A two-layer network with linear outputs can uniformly approximate any continuous function on a compact input domain to arbitrary accuracy.
 - ▶ Provided that the network has a sufficiently large number of hidden units.

Neural Networks for Regression

Univariate

- ▶ In the polynomial fitting example from Chapter 1, given inputs and targets $\{\mathbf{x}_n, t_n\}$, we wanted to find the *optimal parameters \mathbf{w}^* of the polynomial* that best fits the data.
- ▶ Assuming i.i.d data and $t_n \sim \mathcal{N}(y(\mathbf{x}_n, \mathbf{w}), \beta^{-1})$, we wrote the likelihood function whose maximisation corresponded to minimisation of the SSE function

$$E(\mathbf{w}) = \frac{1}{2} \sum_{n=1}^N (y(\mathbf{x}_n, \mathbf{w}) - t_n)^2$$

- ▶ By replacing the polynomial in function $y(\mathbf{x}_n, \mathbf{w})$ by the neural network function, we can minimise $E(\mathbf{w})$ to find *optimal parameters \mathbf{w}^* of the neural network*.

Neural Networks for Regression

Multivariate

- ▶ Similarly, for multivariate targets, assuming multivariate Gaussian density leads to the SSE function

$$E(\mathbf{w}) = \frac{1}{2} \sum_{n=1}^N \|\mathbf{y}(\mathbf{x}_n, \mathbf{w}) - \mathbf{t}_n\|^2$$

where $y_k = a_k = \mathbf{w}_k^T \mathbf{x}$.

- ▶ Notice and prove that

$$\frac{\partial E_n}{\partial a_k} = \underbrace{(y_{kn} - t_{kn})}_{\text{error}_n}$$

Neural Networks for Classification

Binary

- ▶ Similarly, for binary classification, we can assume a Bernoulli distribution on targets which leads to minimisation of the *cross-entropy* function

$$E(\mathbf{w}) = - \sum_{n=1}^N (t_n \ln y(\mathbf{x}_n, \mathbf{w}) + (1 - t_n) \ln(1 - y(\mathbf{x}_n, \mathbf{w})))$$

where $y(\mathbf{x}_n, \mathbf{w}) = P(C_1|\mathbf{x}_n) = \sigma(a) = \sigma(\mathbf{w}^T \mathbf{x}_n)$.

- ▶ Notice (and prove) that

$$\frac{\partial E_n}{\partial a} = \underbrace{(y_n - t_n)}_{\text{error}_n}$$

Neural Networks for Classification

Multiclass

- ▶ For multiclass classification, we can minimise the *multiclass cross-entropy* function

$$E(\mathbf{w}) = - \sum_{n=1}^N \sum_{k=1}^K (t_{kn} \ln y_k(\mathbf{x}_n, \mathbf{w}))$$

where $y_k(\mathbf{x}_n, \mathbf{w}) = P(C_k|\mathbf{x}_n) = \frac{e^{a_k}}{\sum_{j=1}^K e^{a_j}}$ and $a_k = \mathbf{w}_k^T \mathbf{x}$.

- ▶ Notice (and prove) that

$$\frac{\partial E_n}{\partial a_k} = \underbrace{(y_{kn} - t_{kn})}_{\text{error}_n}$$

- ▶ In the following we will denote the error $\frac{\partial E}{\partial a_k}$ as δ_k .

Neural Networks for Classification

- ▶ Note that we can learn classifiers via SSE minimisation also, but the cross-entropy formulations
 1. can be derived probabilistically,
 2. train faster, and
 3. generalise better.

Optimisation

- ▶ For optimisation, we notice that \mathbf{w}^* must be a *stationary point* of $E(\mathbf{w})$.
 - ▶ Minimum, maximum, or saddle point.
 - ▶ A saddle point is where gradient vanishes but point is not an extremum (Example).
- ▶ The goal in neural network minimisation is to find a local minimum.
- ▶ A global minimum, even if found, cannot be verified as globally minimum.
- ▶ Due to symmetry, there are multiple equivalent local minima. Reaching any suitable local minimum is the goal of neural network optimisation.
- ▶ Since there are no analytical solutions for \mathbf{w}^* , we use iterative, numerical procedures.

Optimisation

- ▶ Options for iterative optimisation
 - ▶ Online methods
 - ▶ Stochastic gradient descent
 - ▶ Stochastic gradient descent using mini-batches
 - ▶ Batch methods
 - ▶ Batch gradient descent
 - ▶ Conjugate gradient descent
 - ▶ Quasi-Newton methods
- ▶ Online methods
 - ▶ converge faster since parameter updates are more frequent, and
 - ▶ have greater chance of escaping local minima because stationary point w.r.t to whole data set will generally not be a stationary point w.r.t an individual data point.
- ▶ Batch methods: Conjugate gradient descent and quasi-Newton methods
 - ▶ are more robust and faster than batch gradient descent, and
 - ▶ decrease the error function at each iteration until arriving at a minimum.

Backpropagation

- ▶ For all gradient based methods, however, we must first compute the gradient $\nabla_{\mathbf{w}}E(\mathbf{w})$.
- ▶ We have seen that many error functions of practical interest can be written as a sum of terms

$$E(\mathbf{w}) = \sum_{n=1}^N E_n(\mathbf{w})$$

- ▶ So the essential gradient is $\nabla_{\mathbf{w}}E_n(\mathbf{w})$ which we write in its complete form $\nabla_{\mathbf{w}}E(\mathbf{y}(\mathbf{x}_n, \mathbf{w}))$.

Multivariate Chain Rule

- ▶ The chain rule of differentiation states

$$\frac{df(u(x))}{dx} = \frac{df}{du} \frac{du}{dx}$$

- ▶ The *multivariate* chain rule of differentiation states

$$\frac{df(u(x), v(x))}{dx} = \frac{\partial f}{\partial u} \frac{du}{dx} + \frac{\partial f}{\partial v} \frac{dv}{dx}$$

- ▶ Backpropagation is just an application of the multivariate chain rule.

Backpropagation

- ▶ For the output layer weights

$$\frac{\partial E(y_k(a_k(w_{kj})))}{\partial w_{kj}} = \frac{\partial E}{\partial a_k} \frac{\partial a_k}{\partial w_{kj}} = \delta_k z_j$$

- ▶ For the hidden layer weights, using the multivariate chain rule

$$\begin{aligned} & \frac{\partial}{\partial w_{ji}} E(y_1(a_1(z_j(a_j(w_{ji}))))), y_2(a_2(z_j(a_j(w_{ji}))))), \dots, y_k(a_k(z_j(a_j(w_{ji})))))) \\ &= \frac{\partial E}{\partial a_j} \frac{\partial a_j}{\partial w_{ji}} = \underbrace{\sum_{k=1}^K \underbrace{\frac{\partial E}{\partial a_k}}_{\delta_k} \underbrace{\frac{\partial a_k}{\partial z_j}}_{w_{kj}} \underbrace{\frac{\partial z_j}{\partial a_j}}_{h'(a_j)} \underbrace{\frac{\partial a_j}{\partial w_{ji}}}_{x_i}}_{\delta_j} = \delta_j x_i \end{aligned}$$

- ▶ For each layer, notice the familiar form

$$\text{gradient} = \text{error} \times \text{input}$$

Backpropagation

- ▶ It is important to note that

$$\delta_j = h'(a_j) \sum_{k=1}^K \delta_k w_{kj}$$

yields the error δ_j at hidden neuron j by *backpropagating* the errors δ_k from all output neurons that use the output of neuron j .

- ▶ More generally, compute error δ_j at a layer by *backpropagating* the errors δ_k from next layer.
- ▶ Hence the names *error backpropagation*, *backpropagation*, or simply *backprop*.
- ▶ Very useful machine learning technique that is not limited to neural networks.

Backpropagation

$$\delta_j^{(1)} = h'(a_j) \sum_{k=1}^K \delta_k^{(2)} w_{kj}$$

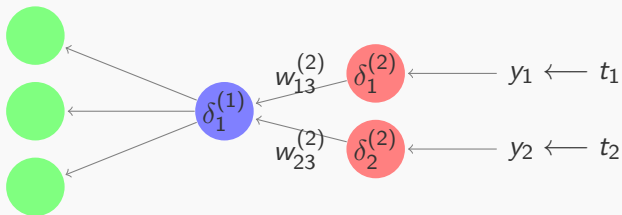


Figure: Visual representation of backpropagation of delta values of layer $l + 1$ to compute delta values of layer l .

Backpropagation

Learning Algorithm

1. Forward propagate the input vector \mathbf{x}_n to compute activations and outputs of every neuron in every layer.
2. Evaluate δ_k for every neuron in output layer.
3. Evaluate δ_j for every neuron in every hidden layer via backpropagation.
4. Compute derivative of each weight $\frac{\partial E}{\partial w}$ via $\delta \times \text{input}$.
5. Update each weight via gradient descent $w^{\tau+1} = w^\tau - \eta \frac{\partial E}{\partial w}$.

Background Math

A $(-1, 1)$ sigmoidal function

- ▶ Since range of logistic sigmoid $\sigma(a)$ is $(0, 1)$, we can obtain a function with $(-1, 1)$ range as $2\sigma(a) - 1$.
- ▶ Another related function with $(-1, 1)$ range is the [tanh](#) function.

$$\tanh(a) = 2\sigma(2a) - 1 = \frac{e^a - e^{-a}}{e^a + e^{-a}}$$

where σ is applied on $2a$.

- ▶ Preferred over logistic sigmoid as activation function $h(a)$ of hidden neurons. (Read Yann LeCun's "Efficient Backprop" paper to understand why.)
- ▶ Just like the logistic sigmoid, derivative of $\tanh(a)$ is simple: $1 - \tanh^2(a)$. ([Prove it.](#))

Backpropagation

A Simple Example

- ▶ Two-layer MLP for multivariate regression from $\mathbb{R}^D \rightarrow \mathbb{R}^K$.
- ▶ Linear outputs $y_k = a_k$ with SSE $E_n = \frac{1}{2} \sum_{k=1}^K (y_k - t_k)^2$.
- ▶ M hidden neurons with $\tanh(\cdot)$ activation functions.

Forward propagation

$$a_j = \sum_{i=0}^D w_{ji}^{(1)} x_i$$

$$z_j = \tanh(a_j)$$

$$y_k = \sum_{j=0}^M w_{kj}^{(2)} z_j$$

$$\delta_k = y_k - t_k$$

- ▶ Compute derivatives $\frac{\partial E_n}{\partial w_{ji}^{(1)}} = \delta_j x_i$ and $\frac{\partial E_n}{\partial w_{kj}^{(2)}} = \delta_k z_j$.

Backpropagate

$$\delta_j = (1 - z_j^2) \sum_{k=1}^K w_{kj}^{(2)} \delta_k$$

Backpropagation

Verifying Correctness

- ▶ *Numerical derivatives* can be computed via finite *central differences*

$$\frac{\partial E_n}{\partial w_{ji}} = \frac{E_n(w_{ji} + \epsilon) - E_n(w_{ji} - \epsilon)}{2\epsilon} + O(\epsilon^2)$$

- ▶ *Analytical derivatives* computed via backpropagation **must be compared** with numerical derivatives for a few examples to verify correctness.
- ▶ Any implementation of analytical derivatives (not just backpropagation) must be compared with numerical derivatives.
- ▶ Notice that we could have avoided backpropagation and computed all required derivatives numerically.
 - ▶ But cost of numerical differentiation is $O(W^2)$ while that of backpropagation is $O(W)$ where W is the total number of weights (and biases) in the network. (Why?)

Assignment 3

Backpropagation for MLPs

- ▶ Implement the backpropagation algorithm for training an MLP.
 - ▶ Code up a generic implementation.
 - ▶ Verify correctness of analytical derivatives.
 - ▶ Understand the experiment and network used for Figure 5.3 in Bishop's book.
 - ▶ Regenerate Figure 5.3 using your implementation.
- ▶ Submit your `_roll_number_MLP.zip` containing
 - ▶ code,
 - ▶ generated image, and
 - ▶ `report.txt/pdf` explaining your results.
- ▶ Due Tuesday (March 21, 2017 before 5:30 pm) on `\\printsrv`.

Neural Network Tips

- ▶ Modularity (`fprop()`, `bprop()`, `check_gradients()`, `update_weights()`, ...).
- ▶ Use a global structure `net.x`, `net.w{1}`, `net.w{2}`, `net.a{1}`, `net.y{1}`, `net.y{2}`, ..., `net=net.fprop(net)`, `net=net.bprop(net)`, ...
- ▶ $\mathbf{a} = W^T \mathbf{x}$.

Regularization in Neural Networks

- ▶ Recall that over-fitting can be lessened via regularization.
 1. Penalise magnitudes of weights: $\tilde{E}(\mathbf{w}) = E(\mathbf{w}) + \frac{\lambda}{2} \mathbf{w}^T \mathbf{w}$.
 2. Separately penalise magnitudes of weights of each layer:

$$\tilde{E}(\mathbf{w}) = E(\mathbf{w}) + \sum_{l=1}^L \frac{\lambda_l}{2} \mathbf{w}^{(l)T} \mathbf{w}^{(l)}.$$
 3. *Early stopping* by checking $E(\mathbf{w})$ on a validation set. Stop when error on validation set starts increasing.
 4. *Tangent propagation*: $\tilde{E}(\mathbf{w}) = E(\mathbf{w}) + \underbrace{\frac{\partial y}{\partial x}}_{\text{rough idea}}$.
 5. Training with transformed data.
 6. Building invariance into the network structure. We cover this in the next lecture on Convolutional Neural Networks.