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## Talk Announcement

Yann LeCun (NYU & FaceBook AI) 28.11. 15:00-16:30h, SuperC 6th floor (Ford Saal)

The rapid progress of AI in the last few years are largely the result of advances in deep learning and neural nets, combined with the availability of large datasets and fast GPUs. We now have systems that can recognize images with an accuracy that rivals that of humans. This will lead to revolutions in several domains such as autonomou transportation and medical image analysis. But all of these systems currently use supervised learning in which the machine is trained with inputs labeled by humans. The challenge of the next several years is to let machines learn from raw, unlabeled data such as video or text. This is known as predictive (or unsupervised) learning. Intelligen systems today do not possess "common sense", which humans and animals acquire by observing the world, by acting in it, and by understanding the physical constraints of it. I will argue that the ability of machines to learn predictive models of the world is a key component of that will enable significant progress in AI. The main technical difficulty is that the world is only partially predictable. A general formulation of unsupervised learning that deals with partial predictability will be presented. The formulation connects many well-known approaches to unsupervised learning, as well as new and exciting ones such as adversarial training.

· No lecture next Monday - go see the talk!

 $f: \mathcal{X} \to \mathbb{R}$ 

## This Lecture: Advanced Machine Learning

- · Regression Approaches
  - > Linear Regression
  - Regularization (Ridge, Lasso)
  - Kernels (Kernel Ridge Regression)
  - Gaussian Processes
- Approximate Inference
  - Sampling Approaches
  - > MCMC
- Deep Learning
  - > Linear Discriminants
  - Neural Networks
  - > Backpropagation
  - CNNs, RNNs, ResNets, etc.



- - > Transform vector  ${\bf x}$  with M nonlinear basis functions  $\phi_i({\bf x})$ :

$$y_k(\mathbf{x}) = g \left( \sum_{j=1}^{M} w_{kj} \phi_j(\mathbf{x}) + w_{k0} \right)$$

- » Basis functions  $\phi_i(\mathbf{x})$  allow non-linear decision boundaries.
- > Activation function  $g(\,\cdot\,)$  bounds the influence of outliers.
- > Disadvantage: minimization no longer in closed form.

$$y_k(\mathbf{x}) = g\left(\sum_{j=0}^M w_{kj}\phi_j(\mathbf{x})\right) \qquad \quad \text{with } \phi_0(\mathbf{x}) = 1$$

## **Recap: Gradient Descent**

- Iterative minimization
  - > Start with an initial guess for the parameter values  $w_{\iota,i}^{(0)}$  .
  - Move towards a (local) minimum by following the gradient.
- Basic strategies
- "Batch learning"

$$w_{kj}^{(\tau+1)} = w_{kj}^{(\tau)} - \eta \left. \frac{\partial E(\mathbf{w})}{\partial w_{kj}} \right|_{\mathbf{w}^{(\tau)}}$$

"Sequential updating" 
$$w_{kj}^{(\tau+1)}=w_{kj}^{(\tau)}-\eta\left.\frac{\partial E_n(\mathbf{w})}{\partial w_{kj}}\right|_{\mathbf{w}^{(\tau)}}$$
 where  $E(\mathbf{w})=\sum_{n=1}^N E_n(\mathbf{w})$ 

## **Recap: Gradient Descent**

• Example: Quadratic error function 
$$E(\mathbf{w}) = \sum_{n=1}^{N} \left(y(\mathbf{x}_n; \mathbf{w}) - \mathbf{t}_n\right)^2$$

Sequential updating leads to delta rule (=LMS rule)

$$w_{kj}^{(\tau+1)} = w_{kj}^{(\tau)} - \eta (y_k(\mathbf{x}_n; \mathbf{w}) - t_{kn}) \phi_j(\mathbf{x}_n)$$
$$= w_{kj}^{(\tau)} - \eta \delta_{kn} \phi_j(\mathbf{x}_n)$$

$$\delta_{kn} = y_k(\mathbf{x}_n; \mathbf{w}) - t_{kn}$$

⇒ Simply feed back the input data point, weighted by the classification error.

## Recap: Probabilistic Discriminative Models

· Consider models of the form

$$p(C_1|\boldsymbol{\phi}) = y(\boldsymbol{\phi}) = \sigma(\mathbf{w}^T \boldsymbol{\phi})$$

with

$$p(\mathcal{C}_2|\boldsymbol{\phi}) = 1 - p(\mathcal{C}_1|\boldsymbol{\phi})$$

- · This model is called logistic regression.
- Properties
  - > Probabilistic interpretation
  - > But discriminative method; only focus on decision hyperplane
  - > Advantageous for high-dimensional spaces, requires less parameters than explicitly modeling  $p(\phi | C_k)$  and  $p(C_k)$ .

## Recap: Logistic Regression

• Let's consider a data set  $\{\phi_n,t_n\}$  with  $n=1,\ldots,N$ , where  $\phi_n = \phi(\mathbf{x}_n)$  and  $t_n \in \{0,1\}$ ,  $\mathbf{t} = (t_1,\ldots,t_N)^T$  .

• With 
$$y_n=p(\mathcal{C}_1|\pmb{\phi}_n)$$
, we can write the likelihood as 
$$p(\mathbf{t}|\mathbf{w})=\prod_{n=1}^Ny_n^{t_n}\left\{1-y_n\right\}^{1-t_n}$$

· Define the error function as the negative log-likelihood  $E(\mathbf{w}) = -\ln p(\mathbf{t}|\mathbf{w})$ 

$$= -\sum_{n=1}^{N} \{t_n \ln y_n + (1 - t_n) \ln(1 - y_n)\}\$$

> This is the so-called cross-entropy error function.

## **Softmax Regression**

- · Multi-class generalization of logistic regression
  - > In logistic regression, we assumed binary labels  $t_n \in \{0,1\}$
  - Softmax generalizes this to K values in 1-of-K notation.

$$\mathbf{y}(\mathbf{x}; \mathbf{w}) = \begin{bmatrix} P(y = 1 | \mathbf{x}; \mathbf{w}) \\ P(y = 2 | \mathbf{x}; \mathbf{w}) \\ \vdots \\ P(y = K | \mathbf{x}; \mathbf{w}) \end{bmatrix} = \frac{1}{\sum_{j=1}^{K} \exp(\mathbf{w}_{j}^{\top} \mathbf{x})} \begin{bmatrix} \exp(\mathbf{w}_{1}^{\top} \mathbf{x}) \\ \exp(\mathbf{w}_{2}^{\top} \mathbf{x}) \\ \vdots \\ \exp(\mathbf{w}_{K}^{\top} \mathbf{x}) \end{bmatrix}$$

> This uses the softmax function

$$\frac{\exp(a_k)}{\sum_j \exp(a_j)}$$

> Note: the resulting distribution is normalized.

## Softmax Regression Cost Function

Logistic regression

> Alternative way of writing the cost function

$$\begin{split} E(\mathbf{w}) &= & -\sum_{n=1}^{N} \left\{ t_n \ln y_n + (1 - t_n) \ln(1 - y_n) \right\} \\ &= & -\sum_{n=1}^{N} \sum_{k=0}^{1} \left\{ \mathbb{I} \left( t_n = k \right) \ln P \left( y_n = k | \mathbf{x}_n; \mathbf{w} \right) \right\} \end{split}$$

Softmax regression

ightarrow Generalization to K classes using indicator functions

$$E(\mathbf{w}) = -\sum_{n=1}^{N} \sum_{k=1}^{K} \left\{ \mathbb{I}(t_n = k) \ln \frac{\exp(\mathbf{w}_k^{\top} \mathbf{x})}{\sum_{j=1}^{K} \exp(\mathbf{w}_j^{\top} \mathbf{x})} \right\}$$

$$E(\mathbf{w}) = \sum_{n=1}^{N} \left[ \mathbb{I}(t_n = k) \ln P(\mathbf{w} - k|\mathbf{x} \cdot \mathbf{w}) \right]$$

 $\nabla_{\mathbf{w}_k} E(\mathbf{w}) \ = \ -\sum_{n=1}^N \left[ \mathbb{I}\left(t_n = k\right) \ln P\left(y_n = k | \mathbf{x}_n; \mathbf{w}\right) \right]$ 

## Optimization

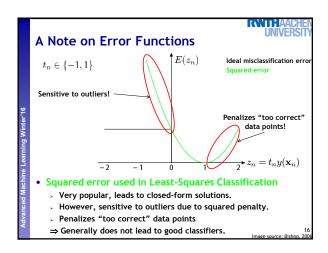
- · Again, no closed-form solution is available
  - Resort again to Gradient Descent
  - Gradient

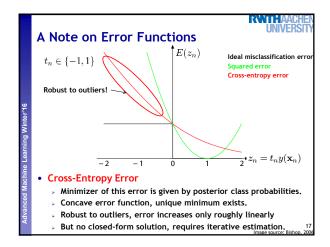
$$\nabla_{\mathbf{w}_k} E(\mathbf{w}) = -\sum_{n=1}^{N} \left[ \mathbb{I}\left(t_n = k\right) \ln P\left(y_n = k | \mathbf{x}_n; \mathbf{w}\right) \right]$$

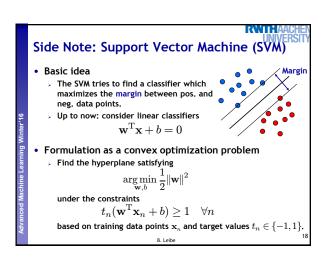
- Note
  - $abla_{\mathbf{w}^k} \, E(\mathbf{w})$  is itself a vector of partial derivatives for the different components of  $\mathbf{w}_k$ .
  - > We can now plug this into a standard optimization package.

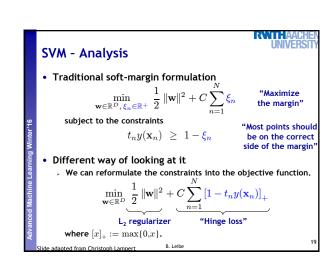
## A Note on Error Functions $E(z_n)$ Ideal misclassification error $t_n \in \{-1, 1\}$ Not differentiable! $z_n = t_n y(\mathbf{x}_n)$ Ideal misclassification error function (black)

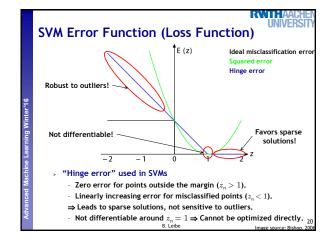
- > This is what we want to approximate,
- > Unfortunately, it is not differentiable.
- The gradient is zero for misclassified points.
- ⇒ We cannot minimize it by gradient descent.

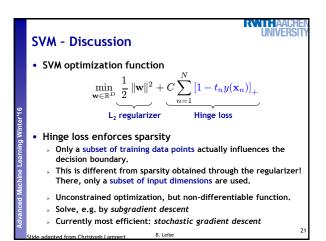


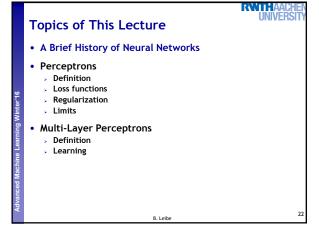


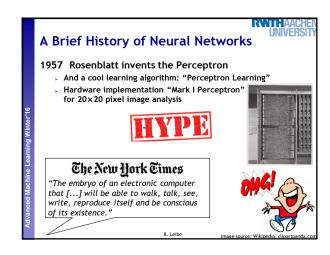


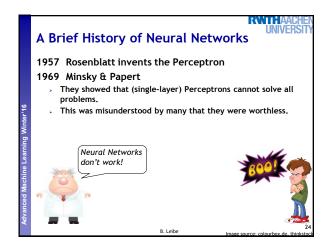


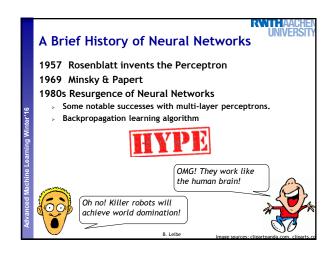


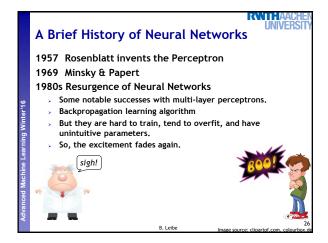


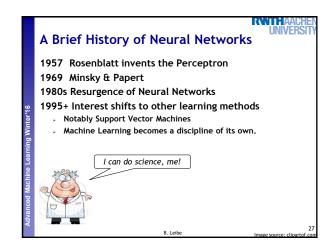


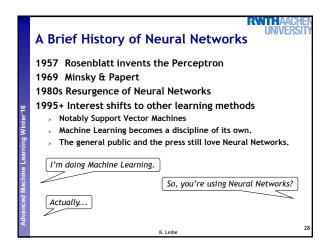


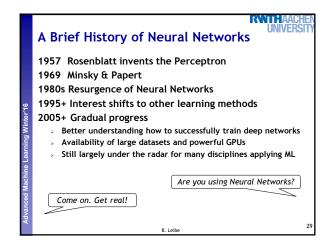


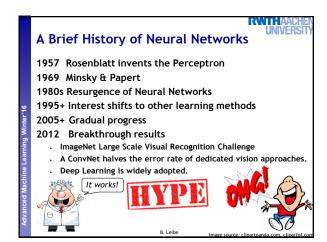


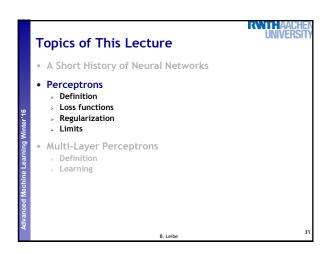


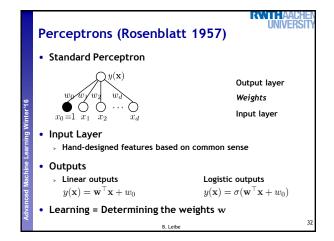


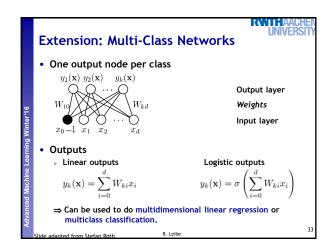


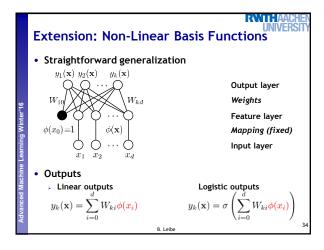


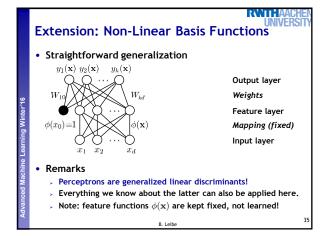












## Perceptron Learning • Very simple algorithm

- Process the training cases in some permutation
  - > If the output unit is correct, leave the weights alone.
  - > If the output unit incorrectly outputs a zero, add the input vector to the weight vector.
  - If the output unit incorrectly outputs a one, subtract the input vector from the weight vector.
- This is guaranteed to converge to a correct solution if such a solution exists.

such a solution exists.

Slide adapted from Geoff Hinton B. Leibe

# Perceptron Learning • Let's analyze this algorithm... • Process the training cases in some permutation • If the output unit is correct, leave the weights alone. • If the output unit incorrectly outputs a zero, add the input vector to the weight vector. • If the output unit incorrectly outputs a one, subtract the input vector from the weight vector. • Translation $w_{kj}^{(\tau+1)} = w_{kj}^{(\tau)}$

## Perceptron Learning Let's analyze this algorithm... Process the training cases in some permutation If the output unit is correct, leave the weights alone. If the output unit incorrectly outputs a zero, add the input vector to the weight vector. If the output unit incorrectly outputs a one, subtract the input vector from the weight vector. Translation w<sub>kj</sub><sup>(τ+1)</sup> = w<sub>kj</sub><sup>(τ)</sup> − η (y<sub>k</sub>(x<sub>n</sub>; w) − t<sub>kn</sub>) φ<sub>j</sub>(x<sub>n</sub>) This is the Delta rule a.k.a. LMS rule!

⇒ Perceptron Learning corresponds to 1st-order (stochastic) Gradient Descent of a quadratic error function!

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## Functions • We can now also apply other loss functions • L2 loss $\Rightarrow$ Least-squares regression $L(t,y(\mathbf{x})) = \sum_n (y(\mathbf{x}_n) - t_n)^2$ • L1 loss: $\Rightarrow$ Median regression $L(t,y(\mathbf{x})) = \sum_n |y(\mathbf{x}_n) - t_n|$ • Cross-entropy loss $\Rightarrow$ Logistic regression $L(t,y(\mathbf{x})) = -\sum_n \{t_n \ln y_n + (1-t_n) \ln(1-y_n)\}$ • Hinge loss $\Rightarrow$ SVM classification $L(t,y(\mathbf{x})) = \sum_n [1-t_ny(\mathbf{x}_n)]_+$ • Softmax loss $\Rightarrow$ Multi-class probabilistic classification $L(t,y(\mathbf{x})) = -\sum_n \sum_k \left\{ \mathbb{I}(t_n = k) \ln \frac{\exp(y_k(\mathbf{x}))}{\sum_j \exp(y_j(\mathbf{x}))} \right\}_{39}$

## Regularization

- In addition, we can apply regularizers
  - E.g., an L2 regularizer

$$E(\mathbf{w}) = \sum_{n} L(t_n, y(\mathbf{x}_n; \mathbf{w})) + \lambda ||\mathbf{w}||^2$$

- $\succ$  This is known as  $\stackrel{n}{\textit{weight decay}}$  in Neural Networks.
- We can also apply other regularizers, e.g. L1 ⇒ sparsity
- > Since Neural Networks often have many parameters, regularization becomes very important in practice.
- > We will see more complex regularization techniques later on...

## **Limitations of Perceptrons**

- · What makes the task difficult?
  - Perceptrons with fixed, hand-coded input features can model any separable function perfectly...
  - ...given the right input features.
  - > For some tasks this requires an exponential number of input features.
    - E.g., by enumerating all possible binary input vectors as separate feature units (similar to a look-up table).
    - But this approach won't generalize to unseen test cases!
  - $\Rightarrow$  It is the feature design that solves the task!
  - Once the hand-coded features have been determined, there are very strong limitations on what a perceptron can learn.
    - Classic example: XOR function.

Wait...

- · Didn't we just say that...
  - > Perceptrons correspond to generalized linear discriminants
  - And Perceptrons are very limited...
  - Doesn't this mean that what we have been doing so far in this lecture has the same problems???
- Yes, this is the case.
  - > A linear classifier cannot solve certain problems (e.g., XOR).
  - $\circ^{C_{_{\! 1}}}$ However, with a non-linear classifier based on the right kind of features, the problem becomes solvable.
  - ⇒ So far, we have solved such problems by hand-designing good features  $\phi$  and kernels  $\phi^{T}\phi$ .
  - ⇒ Can we also learn such feature representations?

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**Topics of This Lecture** 

## A Short History of Neural Networks

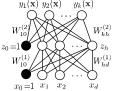
- Perceptrons
  - Definition
  - Loss functions
- Regularization
- Limits

## • Multi-Layer Perceptrons

- Definition
- Learning

## Multi-Layer Perceptrons

· Adding more layers



Output layer

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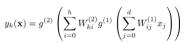
Hidden layer

Input layer

Output

$$y_k(\mathbf{x}) = g^{(2)} \left( \sum_{i=0}^h W_{ki}^{(2)} g^{(1)} \left( \sum_{j=0}^d W_{ij}^{(1)} x_j \right) \right)$$

## **Multi-Layer Perceptrons**



• Activation functions  $g^{(k)}$ :

- For example:  $g^{(2)}(a)=\sigma(a)$ ,  $g^{(1)}(a)=a$
- · The hidden layer can have an arbitrary number of nodes
  - There can also be multiple hidden layers.
- · Universal approximators
  - A 2-layer network (1 hidden layer) can approximate any continuous function of a compact domain arbitrarily well! (assuming sufficient hidden nodes)

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## Learning with Hidden Units

- Networks without hidden units are very limited in what they can learn
  - More layers of linear units do not help ⇒ still linear
  - > Fixed output non-linearities are not enough.
- We need multiple layers of adaptive non-linear hidden units. But how can we train such nets?
  - Need an efficient way of adapting all weights, not just the last layer.
  - Learning the weights to the hidden units = learning features
  - $\,\succ\,$  This is difficult, because nobody tells us what the hidden units should do.
  - ⇒ Next lecture

Slide adapted from Geoff Hintor

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## References and Further Reading

 More information on Neural Networks can be found in Chapters 6 and 7 of the Goodfellow & Bengio book

> lan Goodfellow, Aaron Courville, Yoshua Bengio Deep Learning MIT Press, in preparation



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https://goodfeli.github.io/dlbook/

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