

COS 402 – Machine Learning and Artificial Intelligence Fall 2016

Lecture 8: Introduction to Deep Learning: Part 2

(More on backpropagation, and ConvNets)

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Recap: Structure of a deep net

- "Circuit" of gates connected by wires.
- Each wire has a weight on it.
- Each gate computes a simple nonlinear function, which is applied to weighted sum of incoming signals.



Each gate first computes weighted sum of incoming signals, then applies nonlinear function on it.

Basic structure of a deep net (contd)

- "Circuit" of gates connected by wires.
- Each wire has a weight on it.
- Each gate computes a simple nonlinear function, which is applied to weighted sum of incoming signals.

More popular nonlinearities:

- Rectifier Linear Unit ("RELU").
- Sigmoid (soft threshold)



The optimization problem

- N inputs x₁, x₂,.., x_N in R^d, labeled with values y₁, y₂,..., y_N in {0,1}
- Experimenter decides on # of layers, # of nodes in each, and the nonlinearity type.
- (W, A) = Vector of unknowns.
 (Weight of each wire, and "bias" of each node.) input layer

 $f_{W,A}(x)$ = output of this net on input x.

Minimize over (W, A): $\sum_{i} (f_{W, A}(x_i) - y_i)^2 + \text{Regularizer}(W, A)$

Typical choice of regularizer = sum of squares of entries of W.



Minimize over (W, A): OBJECTIVE(W, A) = $\sum_{i} (f_{W, A}(x_i) - y_i)^2$



Gradient calculation as message passing

Imagine: On each node, one little green man doing some computation.

Desired: At the end, each edge knows $\frac{\partial f}{\partial w}$ where w is its weight, and f is the function at the last layer.

Goal:

Work per node = O(# of adjacent edges).
→ Total work by all green men = O(Network Size).



(NB: Green men = Inner loop of some program)

Main idea: Message passing (each message is a real #)



Simple inductive algorithm to compute $\partial y / \partial w_1$ for all nodes y in the network.

Work per node = O(# incoming wires); → Total work is O(network size).

Repeat for all $w_1 \rightarrow Overall$ work becomes O((network size)²)



Improve work to O(Network size) ?

Pattern of operations is identical for different w_i's: consolidate! (i.e., dynamic programming)

Backprop Algorithm:



 $\begin{array}{l} \text{Compute} \\ \text{this locally.} \end{array}$ $\begin{array}{l} \text{Message from } v_1 \text{ to } u_1 = \\ \text{Sum of all its incoming messages} \times \frac{\partial v_1}{\partial u_1} \end{array}$

Green man

at v₁ can

(NB: Amount of work at $v_1 = O(\# \text{ of nodes it is adjacent to}))$

Backprop. Lemma: This rule satisfies for any node v_i : Sum of messages received by $v_i = \frac{\partial f}{\partial v_i}$

Next few slides: Proof of this claim: By induction on (#layers –k) (NOT k)







How to connect to training objective (notes from tablet)

Loss =
$$\sum Loss(x_i, W, A)$$
.
 $\nabla Loss = \sum \nabla Loss(x_i, W, A)$.
 $SGD =) Pick random i, and compute$
 $gradient of Loss(x_i, W, A)$.
 $Via Back - Prop!$
 $x_i \left\{ = \underbrace{\bigcup_{w \in V} }_{w \in V} \circ f - \underbrace{\bigcup_{w \in$

(more tablet notes: how to get partial derivative wrt network parameters)



Can similarly compute partial derivative wrt bias parameter

Some Implementation details

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Minimize over (W, A):

\sum_{i} (f_{W, A}(x_i) - y_i)^2 + \text{Regularizer}(W, A)
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- Recall SGD: use random index i and use that x_i to compute estimate of gradient
- Better: In each iteration, estimate gradient using random sample of d inputs. "Batch" (typically power of 2 in [16, 256]).

Motivations:

(Theoretical) Avg. of a few samples gives more accurate sample of

gradient (lower variance)

(Practical) On many architectures ---eg GPUs– doing $d=2^k$ identical operations is not much more expensive than a single one.

• Let the "learning rate" drop a bit each iteration $w \leftarrow w - \nu \cdot \nabla F$ (*F* is objective) (Rule of thumb: Inversely proportional to # of iterations.)

The output layer

- If desired output is 0/1 then use sigmoid gate at the output
- If output is in {1, 2, 3,.., m} (i.e., multiclass classification) then use m sigmoids at the output layer.

Now output of deep net = $f_{W,A}(x)$ = a vector in [0,1]^m

0000000000 111111 22222222 333333 444444 555555555 66666666 7777777 888888888 999999999

Possible training objective: $\sum_{j} (f_{W,A}(x_{j}) - y_{j})^{2}$ Where y_{i} is a unary representation of output. (Number i is represented as 000010...000) i
Better : $\sum_{j} cross-entropy(f_{W,A}(x_{j}), y_{j})$ cross-entropy(z, y) = $\sum_{j} (y_{j} \ln z_{j} + (1 - y_{j}) \ln(1 - z_{j}))$ y is 0/1 vector; z is real-valued



Convolutional Neural Nets (aka "Convnet")

----- useful in image recognition, language models, etc.

Generic way to reduce # of parameters in the neural net (leverages special structure of images, text etc.) (Motivated by neuroscience studies of Visual Cortex V1)

Key component in nearly all successful deep learning in recent years.

[LeCun et al'98]

Sanity Check: Why is it good to reduce # of network parameters?

Better generalization! (Training uses fewer samples)





Correct model with Sun at center; planets In elliptical orbits. [Copernicus, Kepler]

MORAL: WITH SUFFICIENT # PARAMETERS, INCORRECT MODEL FITS DATA TOO.

Main Idea in Convolution Net: A Local Filter/Feature





"Multiply each pixel value by -4 and add to it values of neighboring pixels."



Nonzeroes whereever neighboring pixels have v. different values ("Edge Detector")!

Many other useful filters were designed ("AI by introspection")

ConvNets try to learn filters from data directly. (Above filter is 3x3 matrix; only 9 parameters!)

The philosophy

- A layer consists of M types of filters, each of which is k x k (e.g., k =5) Filter is applied in every kxk window. (Sometimes, every 2nd or every 3rd window; determined by "stride" parameter.)
- Inputs to each layer are outputs of filters of prev. layer.
- And so on...



At the end of the day, it is just a deep net with special connection structure; trained using backprop.



Convolution v. fast on GPUs.

Convolution: two dimensional case



(Stride =2 \rightarrow Apply filter every 2nd pixel.)



Same architecture at all layers. (Sometimes throw in some fully connected layers at the top.)

Clarification on how to use backprop to train convnets



This formula shows how to "pool" the gradient from the multiple occurences of this parameter.

Going further in deep nets.

- Mechanisms to allow many layers; even 100+. (Gradient gets noisier as it is backpropagated through more layers!)
- Modifications of gradient descent that allow deep nets with feedback connections (output of higher layer feeding into lower layers) "Recurrent Neural Nets"
- Deep nets with memory: e.g. Long-Short Term Memory nets (LSTMs)
- Ways to compose different deep nets automatically; use backprop to propagate gradient across the interface.
- Practical tools such as autograd, tensorflow, Caffe,...

(Look for ugrad deep nets course in spring.)