Lecture 2

Introduction to Deep Neural Networks

OUTLINE

- Introduction to multi-layered neural network
- Optimization (back-propagation)
- Regularization and Dropout
- The vanishing gradient issue
- Toolkit

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FEATURE ENGINEERING

Most current machine learning works well because of human-designed representations and input features



- Time consuming and task/domain dependant
- Features are often both over-specified and incomplete
- \bullet Machine learning \Leftrightarrow optimizing parameters to make the best prediction

REPRESENTATION LEARNING AND DEEP NETWORKS

Representation learning attempts to automatically learn useful features



- Learning a hierarchical and abstract representation
- That can be shared among tasks
- Almost all data is unlabeled \Rightarrow unsupervised learning

THE CURSE OF DIMENSIONALITY

In high-dimensional space, training data becomes sparse



http://www.edupristine.com/blog/curse-dimensionality

To generalize :

- Use the distance to define some sort of "near-ness"
- Spread the probability mass around training examples (smooth the empirical distribution)

THE CURSE OF DIMENSIONALITY (DIMENSION=3)

In 2-dimensions, two points are near if one falls within a certain radius of another.



http://www.edupristine.com/blog/ curse-dimensionality In 2-d, which proportion of uniformly spaced points within black square fall inside the red circle?

$$\frac{\pi r^2}{4r^2} = \frac{\pi}{4} \approx 78\%$$

This proportion drops to 52% in 3-d, and to 0.24% in 10-d.

Consequence

In high-dimensional space, the distance does not define a useful similarity.

THE CURSE OF DIMENSIONALITY (DIMENSION=3)

Smoothing distribution

- The mass is spread around the examples.
- While plausible in this 2-dimensional case, in higher dimensions, the balls will leave holes or be too large in high probability regions.

Manifold

- If we can discover a representation of the probability concentration,
- a lower dimensional (non-linear) manifold,
- we can "flatten" it by changing the representation
- for which the distance is useful for density estimation, interpolation,



NEURAL NETWORKS



ILLUSTRATION AT DIFFERENT LAYERS?



(Lee et al. 2009)

DEEP LEARNING AND NEURAL NETWORKS: A SUCCESS STORY

Since 2009, deep learning approaches won several challenges

- ImageNet since 2012 (Krizhevsky et al.2012)
- Traffic signs recognition : superhuman performance in 2011 (Ciresan et al.2012) based on (LeCun et al.1989)
- Handwritting recognition since 2009 (Graves and Schmidhuber2009) based on (Hochreiter and Schmidhuber1997)
- Automatic Speech recognition (Hinton et al.2012)





Facebook: millions of faces

2015: human-level performance



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DEEP LEARNING AND NEURAL NETWORKS: A LONG STORY

The breakthrough of 2006

The expression *Deep Learning* was coined around 2006 with papers on unsupervised pre-training of neural nets (Hinton et al.2006; Hinton and Salakhutdinov2006; Bengio et al.2007)

And before? (just a few dates)

- 1958 Rosenblatt proposed the perceptron (Rosenblatt1958), following the work of McCulloch and Pitts in 1943 and Hebb in 1949.
- 1980 Neocognitron (Fukushima1980) or the multilayered NNets
- 1982 Hopfield network with memory and recurrence (Hopfield1982), the unsupervised SOM (Kohonen1982), Neural PCA (Oja1982)
- 1986 Multilayer perceptrons and backpropagation (Rumelhart et al.1986b)
- 1989 Autoencoders (Baldi and Hornik1989), Convolutional network (LeCun et al.1989)
- 1993 Sparse coding (Field1993)

BUT, WHAT IS NEW?

Why today?

- The huge amount of data and the growth of computational power.
- Regularization
- and ...



BUT, WHAT IS NEW?

Why today?

- We have connected the dots, e.g. (Probabilistic) PCA / Neural PCA / Autoencoder
- We understand learning better (regularization, architecture)
- No need to be scared of non-convex optimization (initialization)
- The huge amount of data and the growth of computational power.

What is the difference between a NNet and a Deep Network?

An intensive empirical exploration of different issues



LOGISTIC REGRESSION

Logistic regression (binary classification)



$$f(a = \boldsymbol{w_1}^t \boldsymbol{x}) = \frac{1}{1 + e^{-a}}$$



A single artificial neuron



 y_1 0.25-10 -5 0 5 1 $w_1^t x$

LOGISTIC REGRESSION

From binary classification to K classes (Maxent)



$$f(a_{k} = \boldsymbol{w_{k}}^{t} \boldsymbol{x}) = \frac{e^{a_{k}}}{\sum_{k'=1}^{K} e^{a_{k'}}} = \frac{e^{a_{k}}}{Z(\boldsymbol{x})}$$

A simple neural network

$$\boldsymbol{x} \begin{cases} \mathbf{w}_{1}^{\mathbf{w}_{1}} & y_{1} = f(\boldsymbol{w}_{1}^{t}\boldsymbol{x}) \\ \vdots & \vdots \\ \mathbf{w}_{K}^{\mathbf{v}_{K}} & y_{K} = f(\boldsymbol{w}_{K}^{t}\boldsymbol{x}) \end{cases}$$

- x : input layer
- y : output layer
- each y_k has its parameters \boldsymbol{w}_k
- f is the softmax function

LOGISTIC REGRESSION



- f is usually a non-linear function
- f is a component wise function
- *e.g* the softmax function :

$$y_k = P(c = k | \boldsymbol{x}) = \frac{e^{\boldsymbol{w_k}^t \boldsymbol{x}}}{\sum_{k'} e^{\boldsymbol{w_{k'}}^t \boldsymbol{x}}} = \frac{e^{\boldsymbol{W_{k,:}} \boldsymbol{x}}}{\sum_{k'} e^{\boldsymbol{W_{k',:}} \boldsymbol{x}}}$$

 $\bullet\,$ tanh, sigmoid, relu, ...

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MAXENT CLASSIFIERS



Word representation

For each word in the context

- surface form (one-hot vector)
- prefix
- suffix

• ...



of the input for a better generalization.

NEURAL NETWORKS WITH A HIDDEN LAYER

 \boldsymbol{x} : raw input representation



the internal and tailored representation

Intuitions

- Learn an internal representation of the raw input
- Apply a non-linear transformation
- The input representation \boldsymbol{x} is transformed/compressed in a new representation \boldsymbol{h}
- Adding more layers to obtain a more and more abstract representation

HOW DO WE LEARN THE PARAMETERS ?

For a supervised single layer neural net

Just like a maxent model :

- Calculate the gradient of the objective function and use it to iteratively update the parameters.
- Conjugate gradient, L-BFGS, ...
- In practice : Stochastic gradient descent (SGD)

With one hidden layer

- The internal ("hidden") units make the function non-convex ... just like other models with hidden variables :
 - hidden CRFs (Quattoni et al.2007), ...
- But we can use the same ideas and techniques
- Just without guarantees \Rightarrow **backpropagation** (Rumelhart et al.1986b)

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OPTIMIZATION OF A SINGLE LAYER NETWORK FOR CLASSIFICATION



The set of parameters is denoted $\boldsymbol{\theta}$, in this case :

$$\boldsymbol{\theta} = (\boldsymbol{W})$$

The log-loss (conditional log-likelihood)

Assume the dataset $\mathcal{D} = (\boldsymbol{x}_{(i)}, c_{(i)})_{i=1}^N, c_{(i)} \in \{1, 2, \dots, C\}$

$$\mathcal{L}(\boldsymbol{\theta}) = \sum_{i=1}^{N} l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)}) = \sum_{i=1}^{N} \left(-\sum_{c=1}^{C} \mathbb{I}\{c = c_{(i)}\} \log(P(c|\boldsymbol{x}_{(i)})) \right)$$
(1)
$$l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)}) = -\sum_{k=1}^{C} \mathbb{I}\{k = c_{(i)}\} \log(y_k)$$
(2)

OPTIMIZATION USING SGD

Stochastic Gradient Descent (Bottou2010)

For (t = 1; until convergence; t + +):

- Pick randomly a sample $(\boldsymbol{x}_{(i)}, c_{(i)})$
- Compute the gradient of the loss function w.r.t the parameters (∇_{θ})
- Update the parameters : $\boldsymbol{\theta} = \boldsymbol{\theta} \eta_t \nabla_{\boldsymbol{\theta}}$

Questions

- convergence : what does it mean?
- what do you mean by η_t ?
 - convergence if $\sum_t \eta_t = \infty$ and $\sum_t \eta_t^2 < \infty$
 - $\eta_t \propto t^{-1}$
 - and lot of variants like Adagrad (Duchi et al.2011), Down scheduling, ... see (LeCun et al.2012)





Inference chain :

$$\boldsymbol{x}_{(i)} \longrightarrow (\boldsymbol{a} = \boldsymbol{W} \boldsymbol{x}_{(i)}) \longrightarrow (\boldsymbol{y} = f(\boldsymbol{a})) \longrightarrow l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})$$

The gradient for w_{kj}

$$\nabla_{w_{kj}} = \frac{\partial l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})}{\partial w_{kj}} = \frac{\partial l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})}{\partial \boldsymbol{y}} \times \frac{\partial \boldsymbol{y}}{\partial \boldsymbol{a}} \times \frac{\partial \boldsymbol{a}}{\partial w_{kj}}$$
$$= -(\mathbb{I}\{k = c_{(i)}\} - y_k)x_j = \delta_k x_j$$

GRADIENT COMPUTATION AT SECOND LAYER



Generalization

$$\nabla_{\boldsymbol{W}} = \boldsymbol{\delta} \boldsymbol{x}^{t}$$
$$\delta_{k} = -(\mathbb{I}\{k = c_{(i)}\} - y_{k})$$

with δ the gradient at the pre-activation level.

ITERATIONS



Inference : a forward step

- matrice multiplication with the input $oldsymbol{x}$
- Application of the activation function

One training step : forward and backward steps

- Pick randomly a sample $(\boldsymbol{x}_{(i)}, c_{(i)})$
- Compute $\boldsymbol{\delta}$
- Update the parameters : $\boldsymbol{\theta} = \boldsymbol{\theta} \eta_t \boldsymbol{\delta} \boldsymbol{x}^t$

FEED-FORWARD NETWORKS WITH A MULTI-LAYERS

One layer, indexed by l



EXAMPLE OF TWO LAYERS



$$\boldsymbol{\theta} = (\boldsymbol{W}^{(1)}, \boldsymbol{W}^{(2)})$$

Gradient for the output layer As in the Ex. 1 :

$$\begin{split} \boldsymbol{y} &\to \boldsymbol{y}^{(2)} \\ \boldsymbol{W} &\to \boldsymbol{W}^{(2)} \\ \boldsymbol{x} &\to \boldsymbol{x}^{(2)} = \boldsymbol{y}^{(1)} \\ \nabla_{\boldsymbol{W}^{(2)}} &= \boldsymbol{\delta}^{(2)} \boldsymbol{x}^{(2)^{t}}, \text{ with} \\ \delta_{k}^{(2)} &= -\mathbb{I}\{k = c_{(i)}\} - y_{k} \end{split}$$

BACK-PROPAGATION OF THE LOSS GRADIENT



Inference chain part 1 :

$$\boldsymbol{y}^{(1)} = f^{(1)}(\boldsymbol{a}^{(1)}) \rightarrow \left(\boldsymbol{a}^{(2)} = \boldsymbol{W}^{(2)} \boldsymbol{y}^{(1)}\right) \rightarrow \left(\boldsymbol{y}^{(2)} = f^{(2)}(\boldsymbol{a}^{(2)})\right) \rightarrow l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})$$

$$\nabla_{a_{j}^{(1)}} = \frac{\partial l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})}{\partial a_{j}^{(1)}} = \frac{\partial l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)})}{\partial \boldsymbol{y}^{(2)}} \times \frac{\partial \boldsymbol{y}^{(2)}}{\partial \boldsymbol{a}^{(2)}} \times \frac{\partial \boldsymbol{a}^{(2)}}{\partial y_{j}^{(1)}} \times \frac{\partial \boldsymbol{y}_{j}^{(1)}}{\partial a_{j}^{(1)}}$$
$$= \sum_{k} \left(\mathbb{I} \left\{ k = c_{(i)} \right\} - y_{k}^{(2)} \right) w_{kj}^{(2)} f'^{(1)}(a_{j}) = f'^{(1)}(a_{j}) \left(\boldsymbol{W}_{:,j}^{(2)} \boldsymbol{\delta}^{(2)} \right)^{t} \right)$$

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BACK-PROPAGATION OF THE LOSS GRADIENT (SECOND HIDDEN LAYER)



$$\nabla_{\boldsymbol{y}^{(1)}} = \boldsymbol{W}^{(2)}{}^{t} \boldsymbol{\delta}^{(2)}, \text{ then}$$
$$\boldsymbol{\delta}^{(1)} = \nabla_{\boldsymbol{a}^{(1)}} = f^{(1)'}(\boldsymbol{a}^{(1)}) \circ (\boldsymbol{W}^{(2)}{}^{t} \boldsymbol{\delta}^{(2)})$$

BACK-PROPAGATION OF THE LOSS GRADIENT (THIRD LAYER)



As for the output layer, the gradient is :

$$\nabla_{\boldsymbol{W}^{(1)}} = \boldsymbol{\delta}^{(1)} \boldsymbol{x}^{(1)}{}^{t}, \text{ with}$$
$$\delta_{j}^{(1)} = \nabla_{a_{j}^{(1)}}$$
$$\boldsymbol{\delta}^{(1)} = f'^{(1)}(\boldsymbol{a}^{(1)}) \circ (\boldsymbol{W}^{(2)}{}^{t}\boldsymbol{\delta}^{(2)})$$

The term $(\boldsymbol{W}^{(2)}{}^{t}\boldsymbol{\delta}^{(2)})$ comes from the upper layer.

BACK-PROPAGATION : GENERAL CASE

For a hidden layer l :

• The gradient at the pre-activation level :

$$\boldsymbol{\delta}^{(l)} = f'^{(l)}(\boldsymbol{a}^{(l)}) \circ \left(\boldsymbol{W}^{(l+1)^{t}} \boldsymbol{\delta}^{(l+1)} \right)$$

• The update is as follows :

$$oldsymbol{W}^{(l)} = oldsymbol{W}^{(l)} - \eta_t oldsymbol{\delta}^{(l)} oldsymbol{x}^{(l)}^t$$

The layer should keep :

- $\boldsymbol{W}^{(l)}$: the parameters
- $f^{(l)}$: its activation function
- $x^{(l)}$: its input
- $\boldsymbol{a}^{(l)}$: its pre-activation associated to the input
- $\boldsymbol{\delta}^{(l)}$: for the update and the back-propagation to the layer l-1

BACK-PROPAGATION : ONE TRAINING STEP

Pick a training example : $\boldsymbol{x}^{(1)} = \boldsymbol{x}_{(i)}$

Forward pass

For l = 1 to (L - 1)

• Compute $\boldsymbol{y}^{(l)} = f^{(l)}(\boldsymbol{W}^{(l)}\boldsymbol{x}^{(l)})$

•
$$\boldsymbol{x}^{(l+1)} = \boldsymbol{y}^{(l)}$$

$$y^{(L)} = f^{(L)}(W^{(L)}x^{(L)})$$

Backward pass

Init :
$$\boldsymbol{\delta}^{(L)} = \nabla_{\boldsymbol{a}^{(L)}}$$

For $l = L$ to 2 // all hidden units
• $\boldsymbol{\delta}^{(l-1)} = f'^{(l-1)}(\boldsymbol{a}^{(l-1)}) \circ (\boldsymbol{W}^{(l)t}\boldsymbol{\delta}^{(l)})$
• $\boldsymbol{W}^{(l)} = \boldsymbol{W}^{(l)} - \eta_t \boldsymbol{\delta}^{(l)} \boldsymbol{x}^{(l)t}$
 $\boldsymbol{W}^{(1)} = \boldsymbol{W}^{(1)} - \eta_t \boldsymbol{\delta}^{(1)} \boldsymbol{x}^{(1)t}$

INITIALIZATION RECIPES

A difficult question with several empirical answers.

One standard trick

$$\boldsymbol{W} \sim \mathcal{N}(0, \frac{1}{\sqrt{n_{in}}})$$

with n_{in} is the number of inputs

A more recent one

$$\boldsymbol{W} \sim \mathcal{U} \Big[-\frac{\sqrt{6}}{\sqrt{n_{in}+n_{out}}}, \frac{\sqrt{6}}{\sqrt{n_{in}+n_{out}}} \Big]$$

with n_{in} is the number of inputs

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REGULARIZATION L2 OR GAUSSIAN PRIOR OR WEIGHT DECAY

The basic way :

$$\mathcal{L}(\boldsymbol{\theta}) = \sum_{i=1}^{N} l(\boldsymbol{\theta}, \boldsymbol{x}_{(i)}, c_{(i)}) + \frac{\lambda}{2} ||\boldsymbol{\theta}||^2$$

- The second term is the regularization term.
- Each parameter has a gaussian prior : $\mathcal{N}(0, 1/\lambda)$.
- λ is a hyperparameter.
- The update has the form :

$$\boldsymbol{\theta} = (1 + \eta_t \lambda) \boldsymbol{\theta} - \eta_t \nabla_{\boldsymbol{\theta}}$$

DROPOUT – A NEW REGULARIZATION SCHEME





(a) Standard Neural Net

(b) After applying dropout.

- For each training example : randomly turn-off the neurons of hidden units (with p = 0.5)
- At test time, use each neuron scaled down by p
- Dropout serves to separate effects from strongly correlated features and
- prevents co-adaptation between units
- It can be seen as averaging different models that share parameters.
- It acts as a powerful regularization scheme.

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EXPERIMENTAL OBSERVATIONS (MNIST TASK) – ONE LAYER

The MNIST database

Comparison of different depth for feed-forward architecture



- Hidden layers have a sigmoid activation function.
- The output layer is a softmax.

EXPERIMENTAL OBSERVATIONS (MNIST TASK) – TWO LAYERS

Varying the depth

- Without hidden layer : $\approx 88\%$ accuracy
- 1 hidden layer (30) : $\approx 96.5\%$ accuracy
- 2 hidden layer (30) : $\approx 96.9\%$ accuracy
- 3 hidden layer (30) : $\approx 96.5\%$ accuracy
- 4 hidden layer (30) : $\approx 96.5\%$ accuracy





INTUITIVE EXPLANATION

Let consider the simplest deep neural network, with just a single neuron in each layer.



 w_i, b_i are resp. the weight and bias of neuron *i* and *C* some cost function. Compute the gradient of *C w.r.t* the bias b_1

$$\frac{\partial C}{\partial b_1} = \frac{\partial C}{\partial y_4} \times \frac{\partial y_4}{\partial a_4} \times \frac{\partial a_4}{\partial y_3} \times \frac{\partial y_3}{\partial a_3} \times \frac{\partial a_3}{\partial y_2} \times \frac{\partial y_2}{\partial a_2} \times \frac{\partial a_2}{\partial y_1} \times \frac{\partial y_1}{\partial a_1} \times \frac{\partial a_1}{\partial b_1} \quad (3)$$

$$= \frac{\partial C}{\partial y_4} \times \sigma'(a_4) \times w_4 \times \sigma'(a_3) \times w_3 \times \sigma'(a_2) \times w_2 \times \sigma'(a_1) \quad (4)$$
(5)

INTUITIVE EXPLANATION

The derivative of the activation function : σ'



The different layers in our deep network are learning at vastly different speeds :

- when later layers in the network are learning well,
- early layers often get stuck during training, learning almost nothing at all.

SOME HEURISTICS

Change the activation function (Rectified Linear Unit or ReLU)



- Avoid the vanishing gradient
- Some units can "die"

See (Glorot et al.2011) for more details

Do pre-training when it is possible

See (Hinton et al.2006; Bengio et al.2007) :

when you cannot really escape from the initial (random) point, find a good starting point.

More details

See (Hochreiter et al.2001; Glorot and Bengio2010; LeCun et al.2012)

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USEFUL TOOLS

Theano:http://deeplearning.net/software/theano/

- in python, works on CPU and GPU and several wrappers
- http://lasagne.readthedocs.org/
- http://keras.io/
- https://www.cs.cmu.edu/~ymiao/pdnntk.html
- http://deeplearning.net/software/pylearn2/
- http://blocks.readthedocs.org/

Torch7: http://torch.ch/

Lua interface to C/CUDA

TensorFlow https://www.tensorflow.org API in C++ and Python

Basics of deep learning

- Intro to deep learning: <u>http://www.deeplearningbook.org/contents/intro.htm</u>
- Feedforward multi-layer nets: <u>http://www.deeplearningbook.org/contents/r</u>
- •
- Learning deep architectures for AI
- Practical recommendations for gradient-based training of deep architecture
- <u>Quick'n'dirty introduction to deep learning: Advances in Deep Learning</u>
- <u>A fast learning algorithm for deep belief nets</u>
- Greedy Layer-Wise Training of Deep Networks
- <u>Stacked denoising autoencoders: Learning useful representations in a dee</u> network with a local denoising criterion
- <u>Contractive auto-encoders: Explicit invariance during feature extraction</u>
- <u>Why does unsupervised pre-training help deep learning?</u>
- An Analysis of Single Layer Networks in Unsupervised Feature Learning
- <u>The importance of Encoding Versus Training With Sparse Coding and Vector</u> <u>Quantization</u>
- Representation Learning: A Review and New Perspectives
- Deep Learning of Representations: Looking Forward
- Measuring Invariances in Deep Networks
- Neural networks course at USherbrooke [youtube]

Feedforward nets

- <u>http://www.deeplearningbook.org/contents/mlp.html</u>
- "Improving Neural Nets with Dropout" by Nitish Srivastava
- Batch Normalization
- "Fast Drop Out"
- <u>"Deep Sparse Rectifier Neural Networks"</u>
- <u>"What is the best multi-stage architecture for object recognition?"</u>
- <u>"Maxout Networks</u>"

MCMC

- Iain Murray's MLSS slides
- <u>Radford Neal's Review Paper</u> (old but still very comprehensive)
- Better Mixing via Deep Representations
- Bayesian Learning via Stochastic Gradient Langevin Dynamics

Restricted Boltzmann Machines

- Unsupervised learning of distributions of binary vectors using 2-layer networks
- A practical guide to training restricted Boltzmann machines
- <u>Training restricted Boltzmann machines using approximations to the likelihood</u> gradient

• <u>Tempered Markov Chain Monte Carlo for training of Restricted Boltzmann</u> <u>Machine</u>

- How to Center Binary Restricted Boltzmann Machines
- Enhanced Gradient for Training Restricted Boltzmann Machines
- Using fast weights to improve persistent contrastive divergence
- Training Products of Experts by Minimizing Contrastive Divergence

Boltzmann Machines

- <u>Deep Boltzmann Machines</u> (Salakhutdinov & Hinton)
- Multimodal Learning with Deep Boltzmann Machines
- <u>Multi-Prediction Deep Boltzmann Machines</u>
- <u>A Two-stage Pretraining Algorithm for Deep Boltzmann Machines</u>

Regularized Auto-Encoders

- <u>The Manifold Tangent Classifier</u>
- DL book chapter on autoencoders:

http://www.deeplearningbook.org/contents/autoencoders.html

• DL book chapter on representation learning:

http://www.deeplearningbook.org/contents/representation.html

• <u>Representation Learning: A Review and New Perspectives</u>, in particular section 7.

Regularization

Stochastic Nets & GSNs

- Estimating or Propagating Gradients Through Stochastic Neurons for Conditional Computation
- Learning Stochastic Feedforward Neural Networks
- Generalized Denoising Auto-Encoders as Generative Models
- Deep Generative Stochastic Networks Trainable by Backprop

Others

- Slow, Decorrelated Features for Pretraining Complex Cell-like Networks
- What Regularized Auto-Encoders Learn from the Data Generating Distribution
- Generalized Denoising Auto-Encoders as Generative Models
- <u>Why the logistic function?</u>

Recurrent Nets

- DL book chapter on recurrent nets
- Learning long-term dependencies with gradient descent is difficult
- Advances in Optimizing Recurrent Networks
- Learning recurrent neural networks with Hessian-free optimization
- On the importance of momentum and initialization in deep learning.
- Long short-term memory (Hochreiter & Schmidhuber)
- <u>Generating Sequences With Recurrent Neural Networks</u>
- Long Short-Term Memory in Echo State Networks: Details of a Simulation Study
- <u>The "echo state" approach to analysing and training recurrent neural networks</u>
- Backpropagation-Decorrelation: online recurrent learning with O(N) complexity
- <u>New results on recurrent network training:Unifying the algorithms and accelerating</u>

convergence

- Audio Chord Recognition with Recurrent Neural Networks
- Modeling Temporal Dependencies in High-Dimensional Sequences: Application to
 Polyphonic Music Generation and Transcription

Convolutional Nets

• DL book chapter on convolutional nts:

http://www.deeplearningbook.org/contents/convnets.html

- <u>Generalization and Network Design Strategies</u> (LeCun)
- ImageNet Classification with Deep Convolutional Neural Networks, Alex

Krizhevsky, Ilya Sutskever, Geoffrey E Hinton, NIPS 2012.

On Random Weights and Unsupervised Feature Learning

Optimization issues with DL

- <u>Curriculum Learning</u>
- Evolving Culture vs Local Minima
- Knowledge Matters: Importance of Prior Information for Optimization
- Efficient Backprop
- Practical recommendations for gradient-based training of deep architectures
- Batch Normalization
- Natural Gradient Works Efficiently (Amari 1998)
- Hessian Free
- Natural Gradient (TONGA)
- Revisiting Natural Gradient

READINGS (BOOKS)

- Deep Learning Book by Yoshua Bengio, Ian Goodfellow and Aaron Courville (http://www.iro.umontreal.ca/~bengioy/dlbook/)
- Deep Learning for Natural Language Processing by Stanford University http://cs224d.stanford.edu/
- Deep Learning Methods and Applications by Li Deng and Dong Yu http://research.microsoft.com/pubs/219984/BOOK2014.pdf
- Reading lists for new LISA students by University of Montreal.





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