

10707

Deep Learning

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Deep Belief Networks

Multilayer Neural Net

- Consider a network with L hidden layers.

- layer pre-activation for $k > 0$

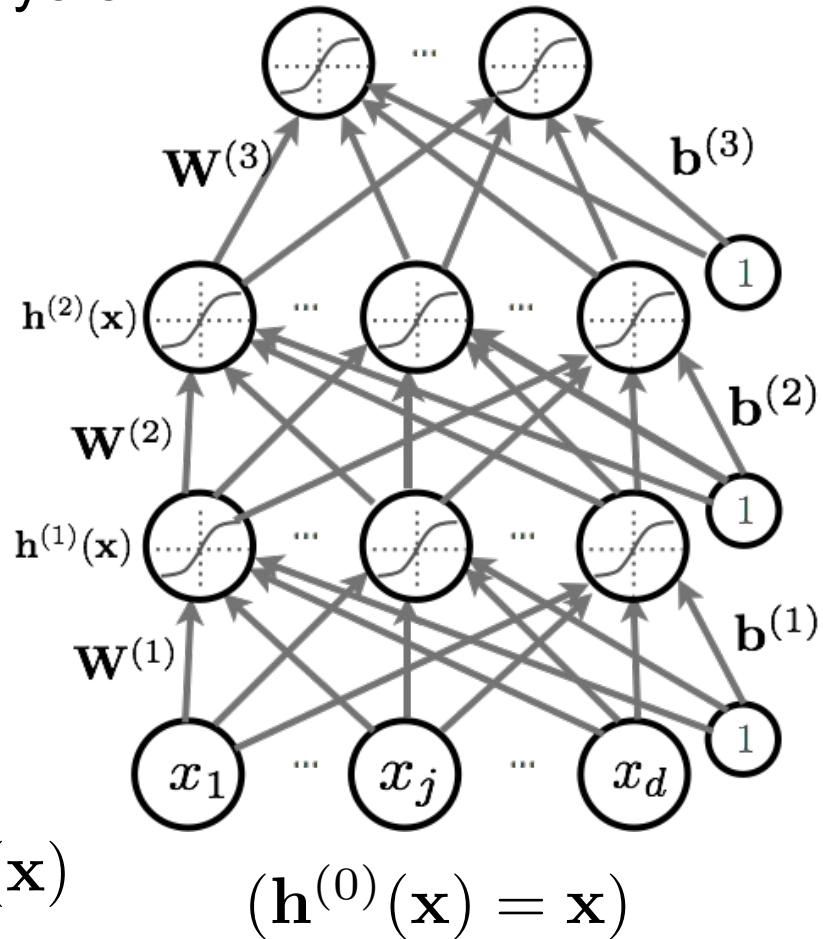
$$\mathbf{a}^{(k)}(\mathbf{x}) = \mathbf{b}^{(k)} + \mathbf{W}^{(k)} \mathbf{h}^{(k-1)}(\mathbf{x})$$

- hidden layer activation from 1 to L:

$$\mathbf{h}^{(k)}(\mathbf{x}) = \mathbf{g}(\mathbf{a}^{(k)}(\mathbf{x}))$$

- output layer activation ($k=L+1$):

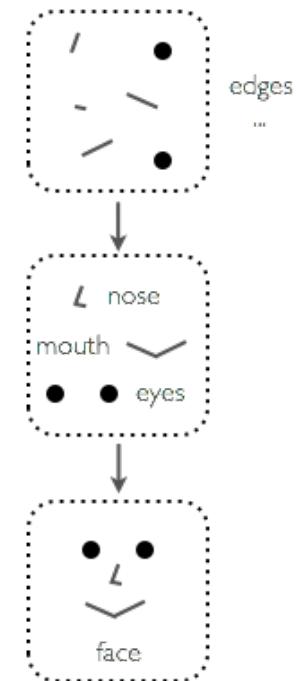
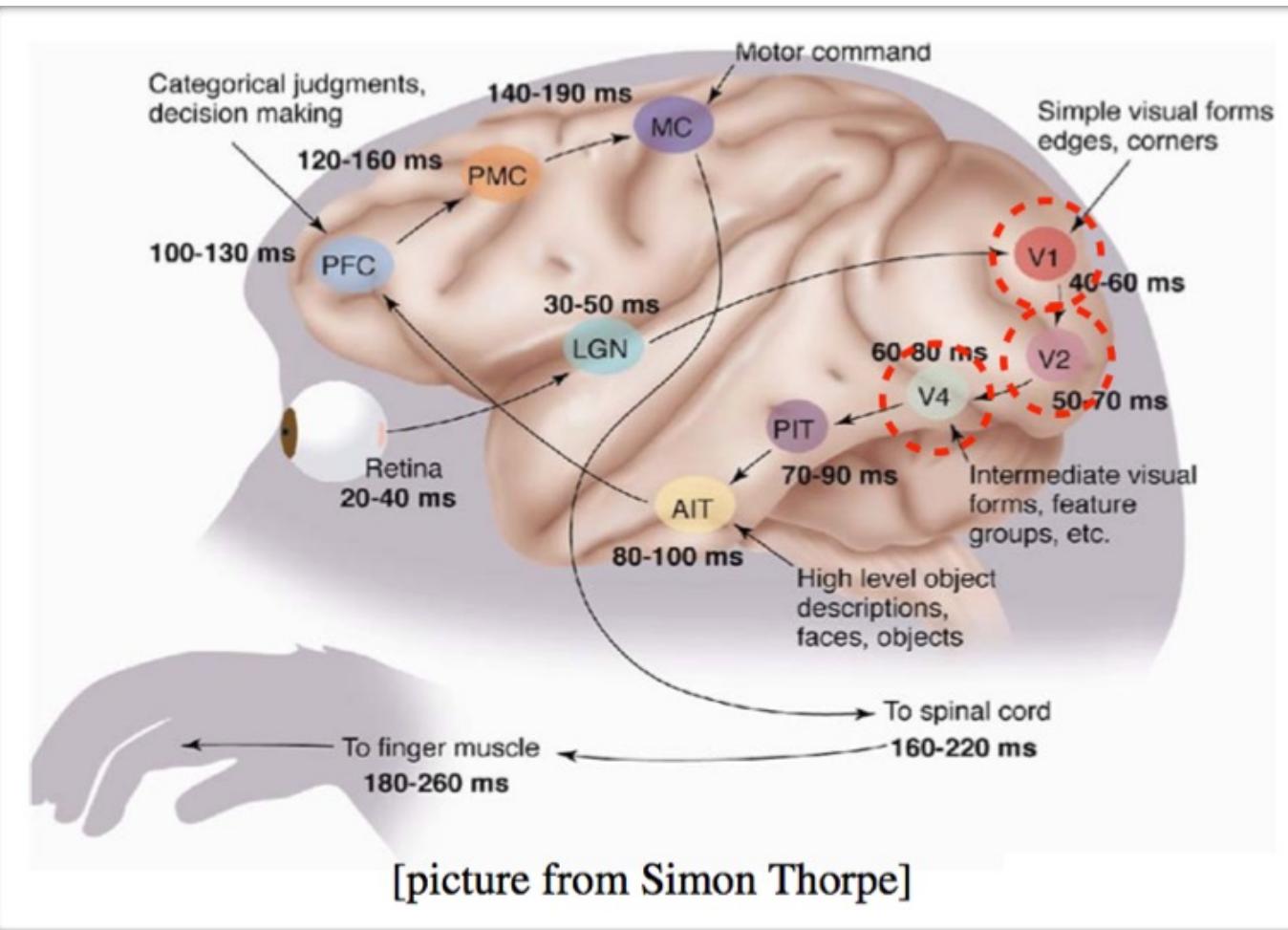
$$\mathbf{h}^{(L+1)}(\mathbf{x}) = \mathbf{o}(\mathbf{a}^{(L+1)}(\mathbf{x})) = \mathbf{f}(\mathbf{x})$$



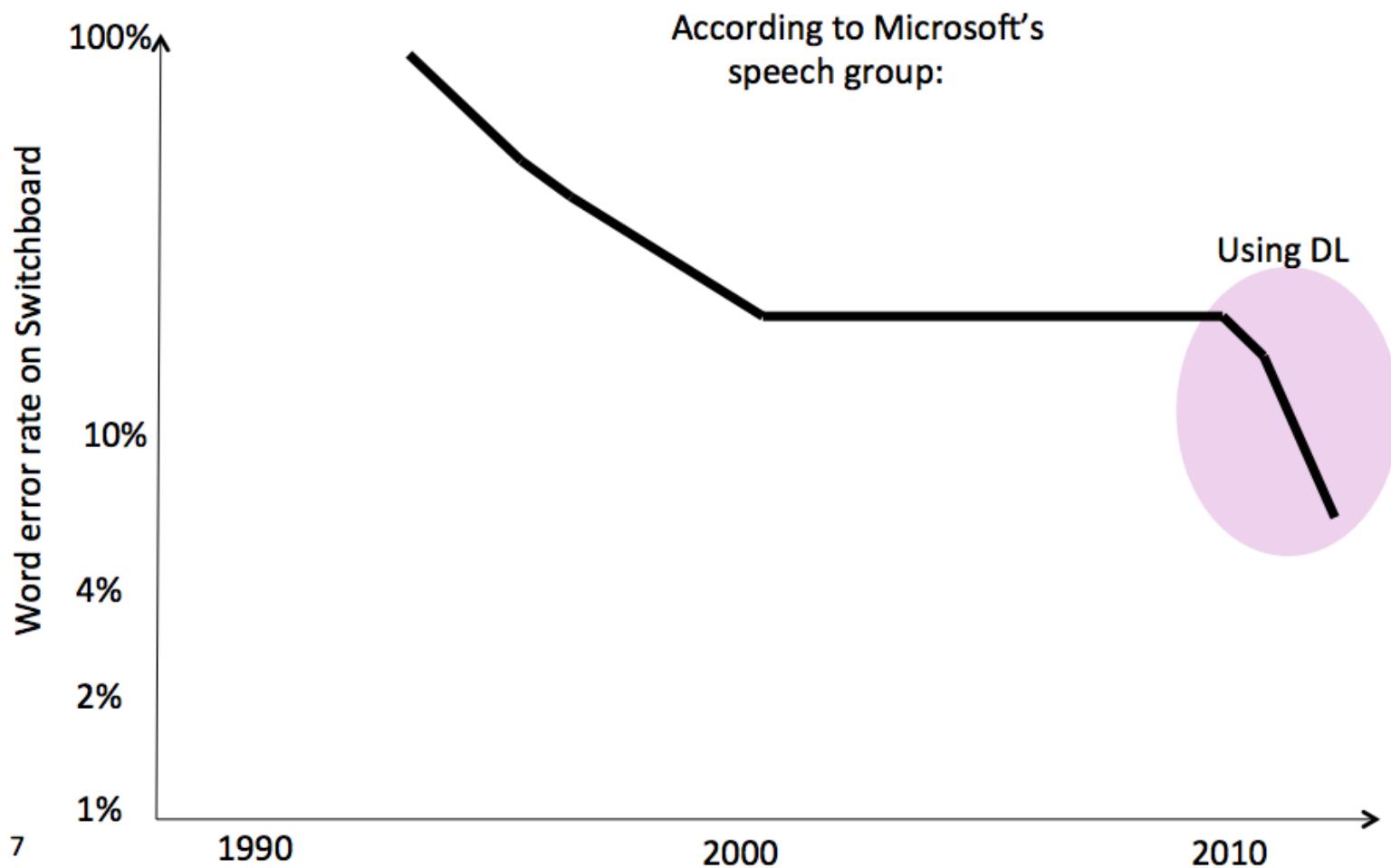
Learning Distributed Representations

- Deep learning is research on learning models with **multilayer representations**
 - multilayer (feed-forward) neural networks
 - multilayer graphical model (deep belief network, deep Boltzmann machine)
- Each layer learns “**distributed representation**”
 - Units in a layer are not mutually exclusive
 - each unit is a separate feature of the input
 - two units can be “active” at the same time
 - Units do not correspond to a partitioning (clustering) of the inputs
 - in clustering, an input can only belong to a single cluster

Inspiration from Visual Cortex

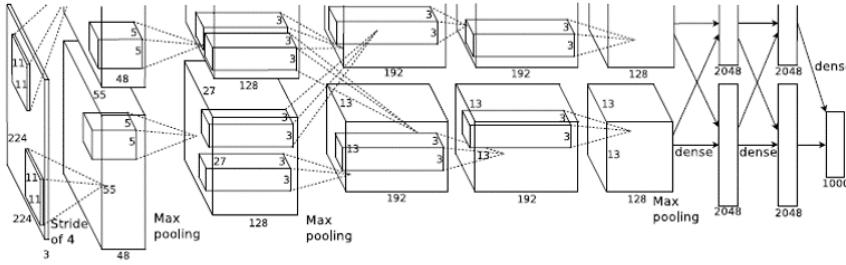


Success Story: Speech Recognition

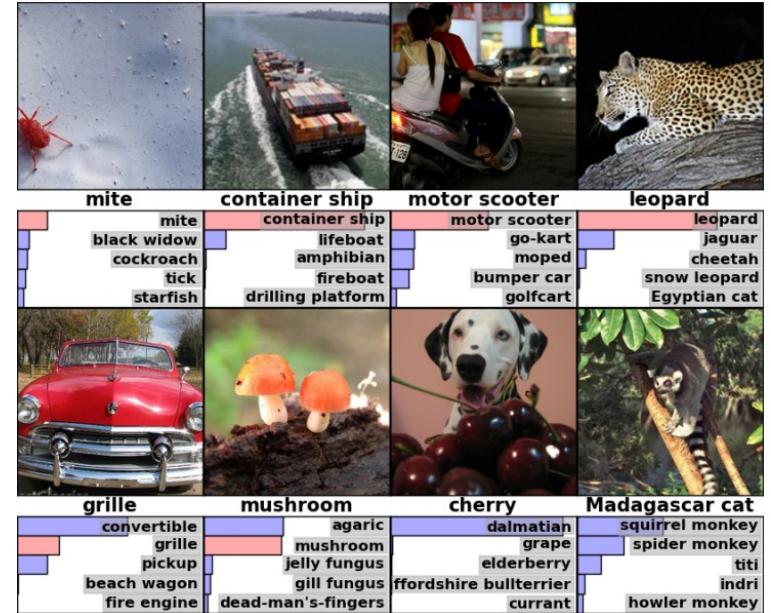


Success Story: Image Recognition

- Deep Convolutional Nets for Vision (Supervised)

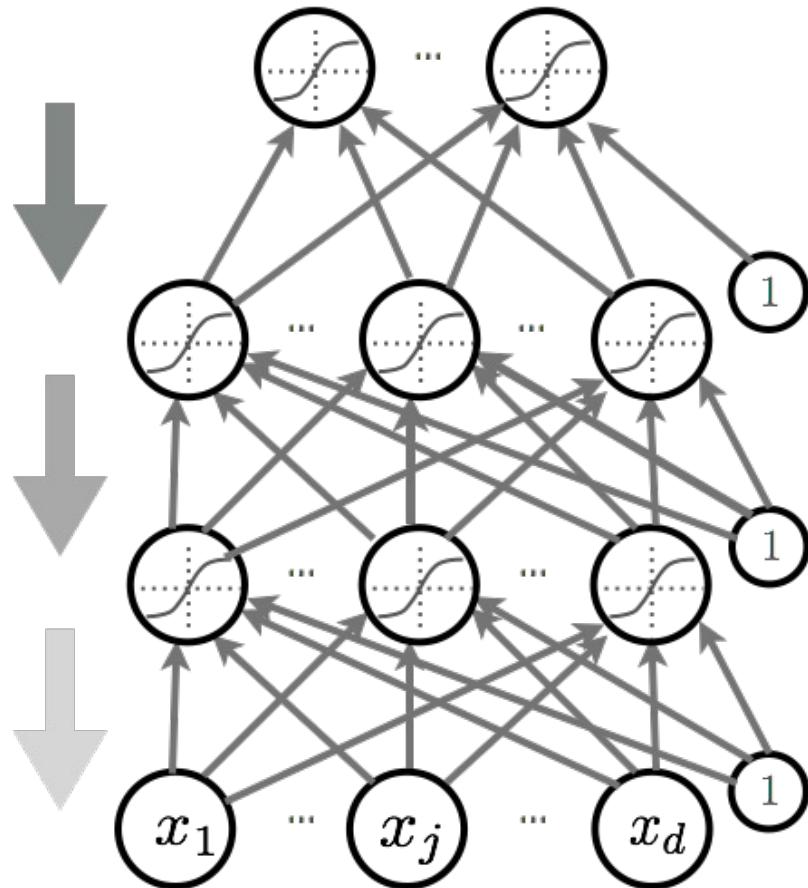


1.2 million training images
1000 classes



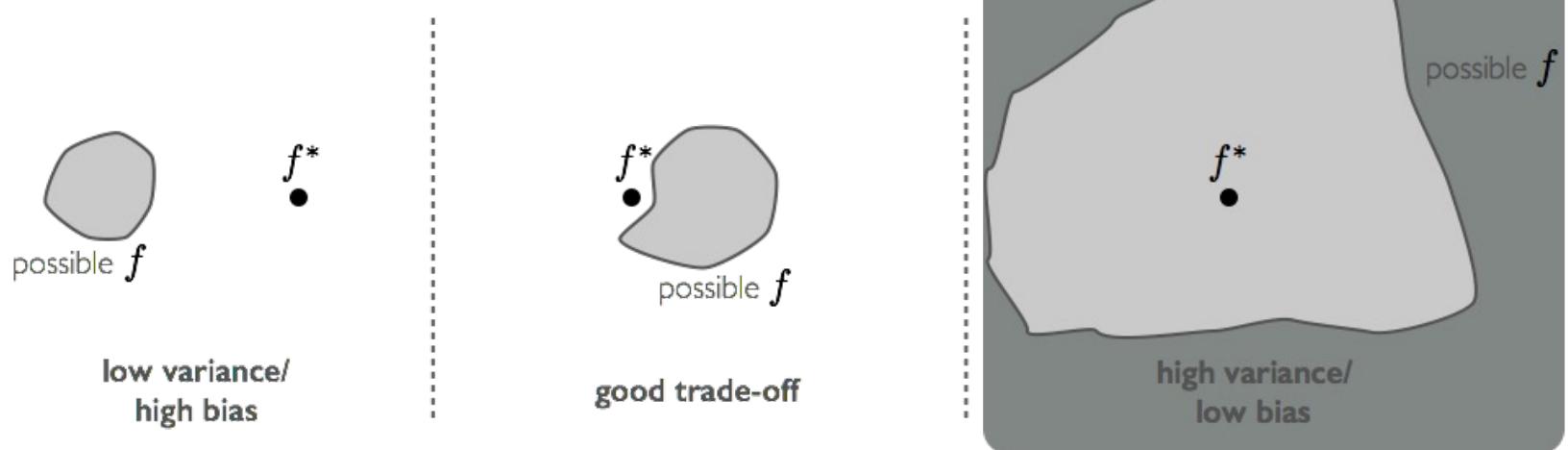
Why Training is Hard

- First hypothesis: Hard optimization problem (underfitting)
 - vanishing gradient problem
 - saturated units block gradient propagation
- This is a well known problem in recurrent neural networks



Why Training is Hard

- Second hypothesis: Overfitting
 - we are exploring a space of complex functions
 - deep nets usually have lots of parameters
- Might be in a high variance / low bias situation

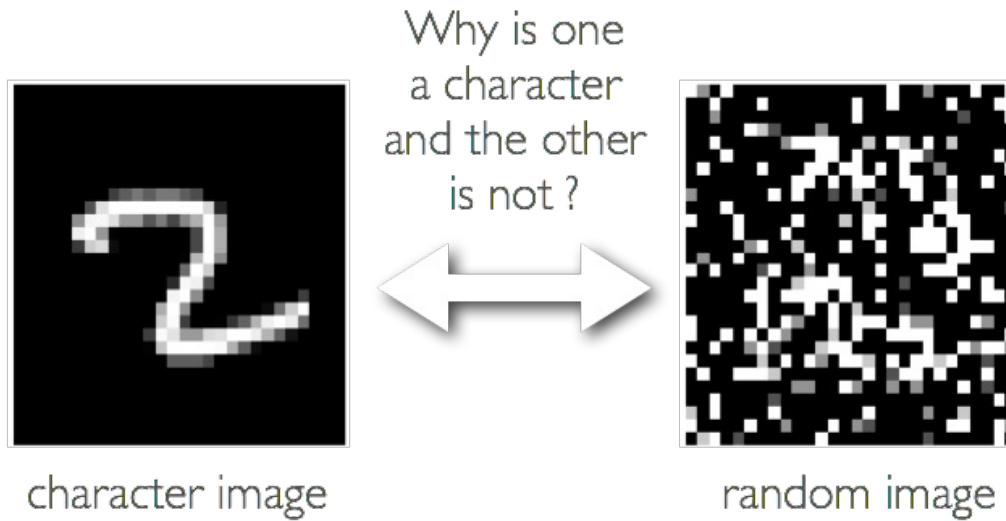


Why Training is Hard

- First hypothesis (underfitting): better optimize
 - Use better optimization tools (e.g. batch-normalization, second order methods, such as KFAC)
 - Use GPUs, distributed computing.
- Second hypothesis (overfitting): use better regularization
 - Unsupervised pre-training
 - Stochastic drop-out training
- For many large-scale practical problems, you will need to use both: better optimization and better regularization!

Unsupervised Pre-training

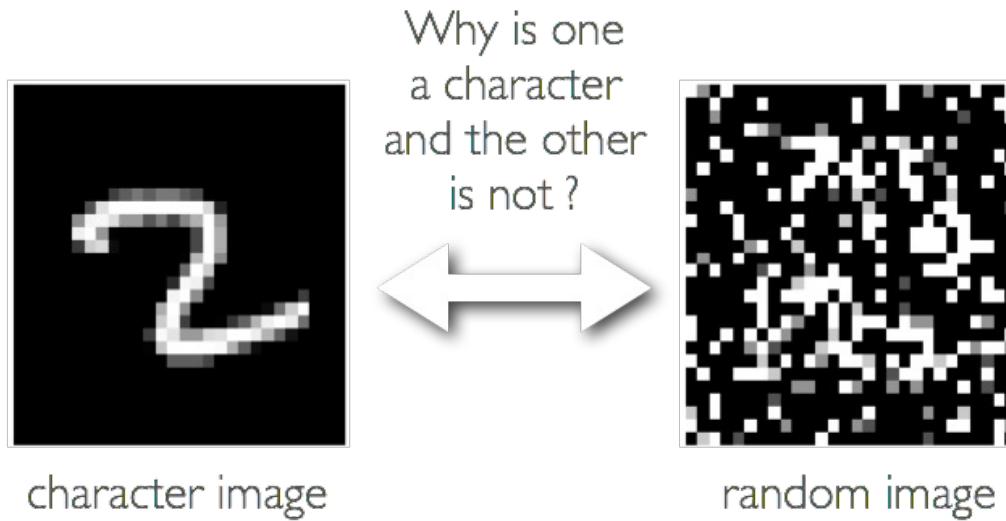
- Initialize hidden layers using unsupervised learning
 - Force network to represent latent structure of input distribution



- Encourage hidden layers to encode that structure

Unsupervised Pre-training

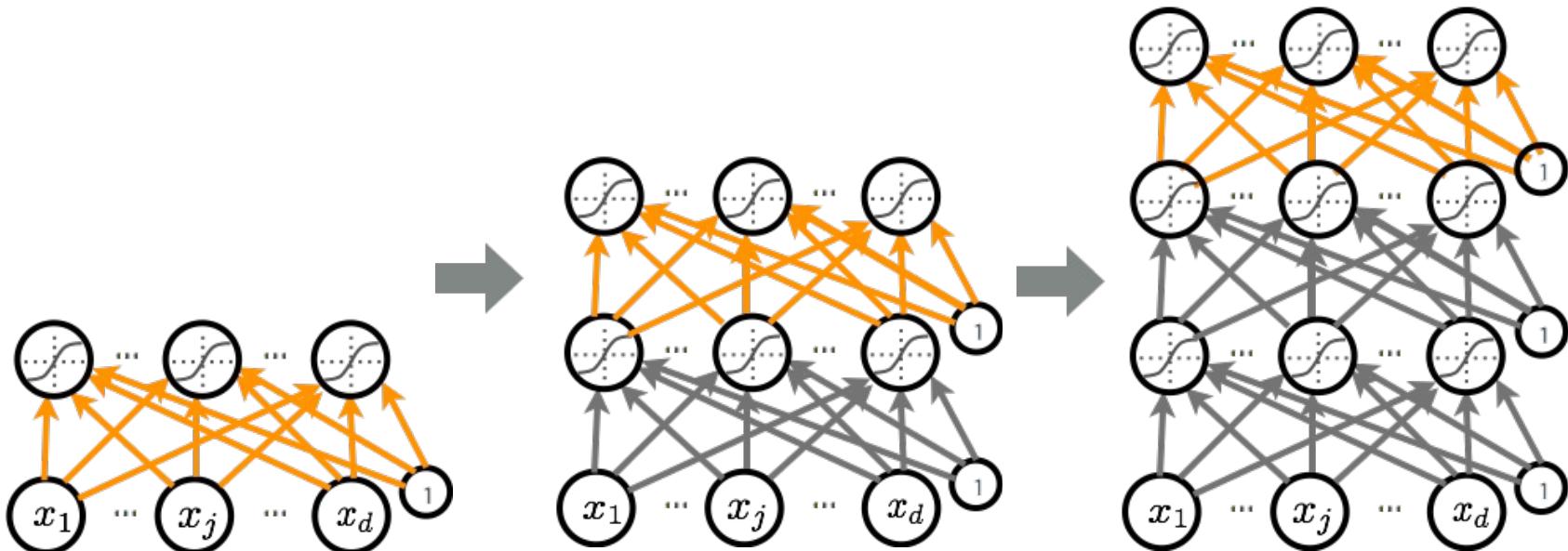
- Initialize hidden layers using unsupervised learning
 - This is a harder task than supervised learning (classification)



- Hence we expect less overfitting

Pre-training

- We will use a greedy, layer-wise procedure
 - Train one layer at a time with unsupervised criterion
 - Fix the parameters of previous hidden layers
 - Previous layers viewed as feature extraction

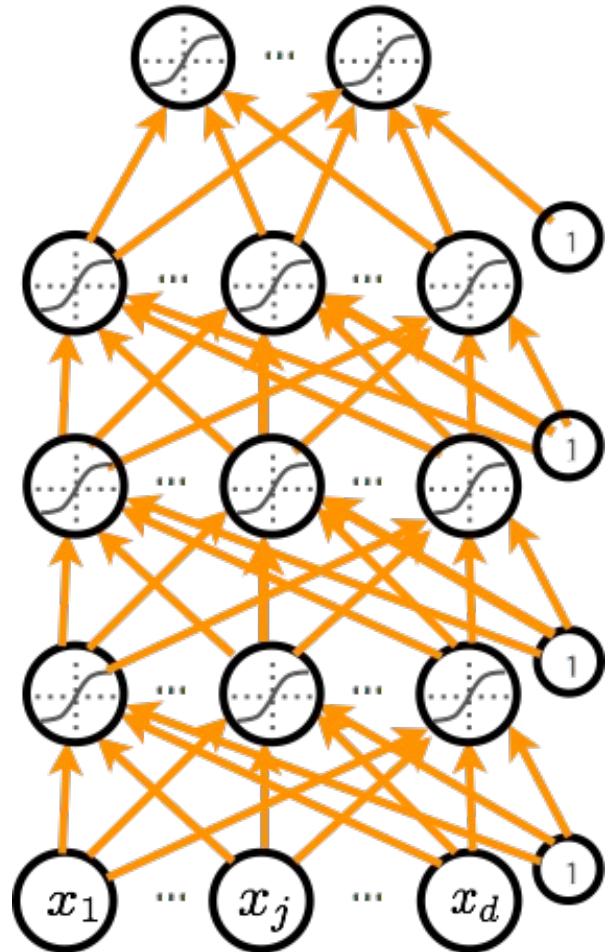


Pre-training

- Unsupervised Pre-training
 - first layer: find hidden unit features that are more common in training inputs than in random inputs
 - second layer: find combinations of hidden unit features that are more common than random hidden unit features
 - third layer: find combinations of combinations of ...
- Pre-training initializes the parameters in a region such that the near local optima overfit less the data

Fine-tuning

- Once all layers are pre-trained
 - add output layer
 - train the whole network using supervised learning
- Supervised learning is performed as in a regular network
 - forward propagation, backpropagation and update
- We call this last phase **fine-tuning**
 - all parameters are “tuned” for the supervised task at hand
 - representation is adjusted to be more discriminative



Stacking RBMs, Autoencoders

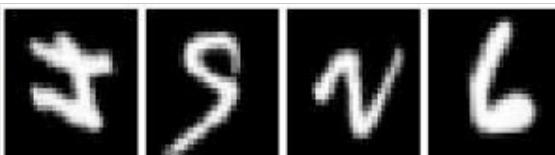
- Stacked Restricted Boltzmann Machines:
 - Hinton, Teh and Osindero suggested this procedure with RBMs,:
A fast learning algorithm for deep belief nets.
 - To recognize shapes, first learn to generate images. Hinton, 2006.
- Stacked autoencoders, sparse-coding models, etc.
 - Bengio, Lamblin, Popovici and Larochelle (stacked autoencoders)
 - Ranzato, Poultney, Chopra and LeCun (stacked sparse coding models)
- Lots of others started stacking models together.

Example

- Datasets generated with varying number of factors of variations

Variations on MNIST

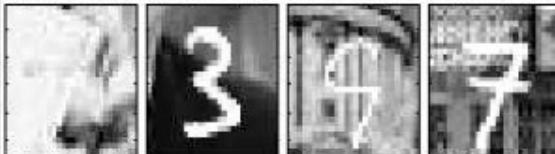
MNIST-rotation



MNIST-random-
background



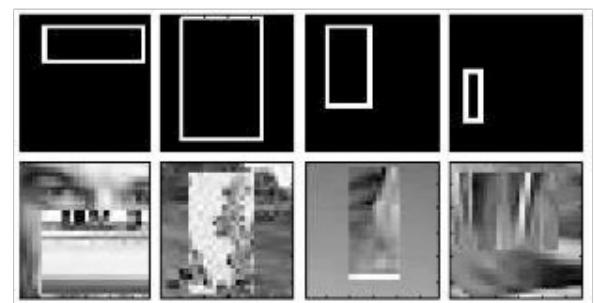
MNIST-image-
background



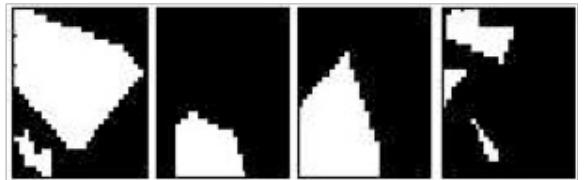
MNIST-
background-
rotation



Tall or wide?



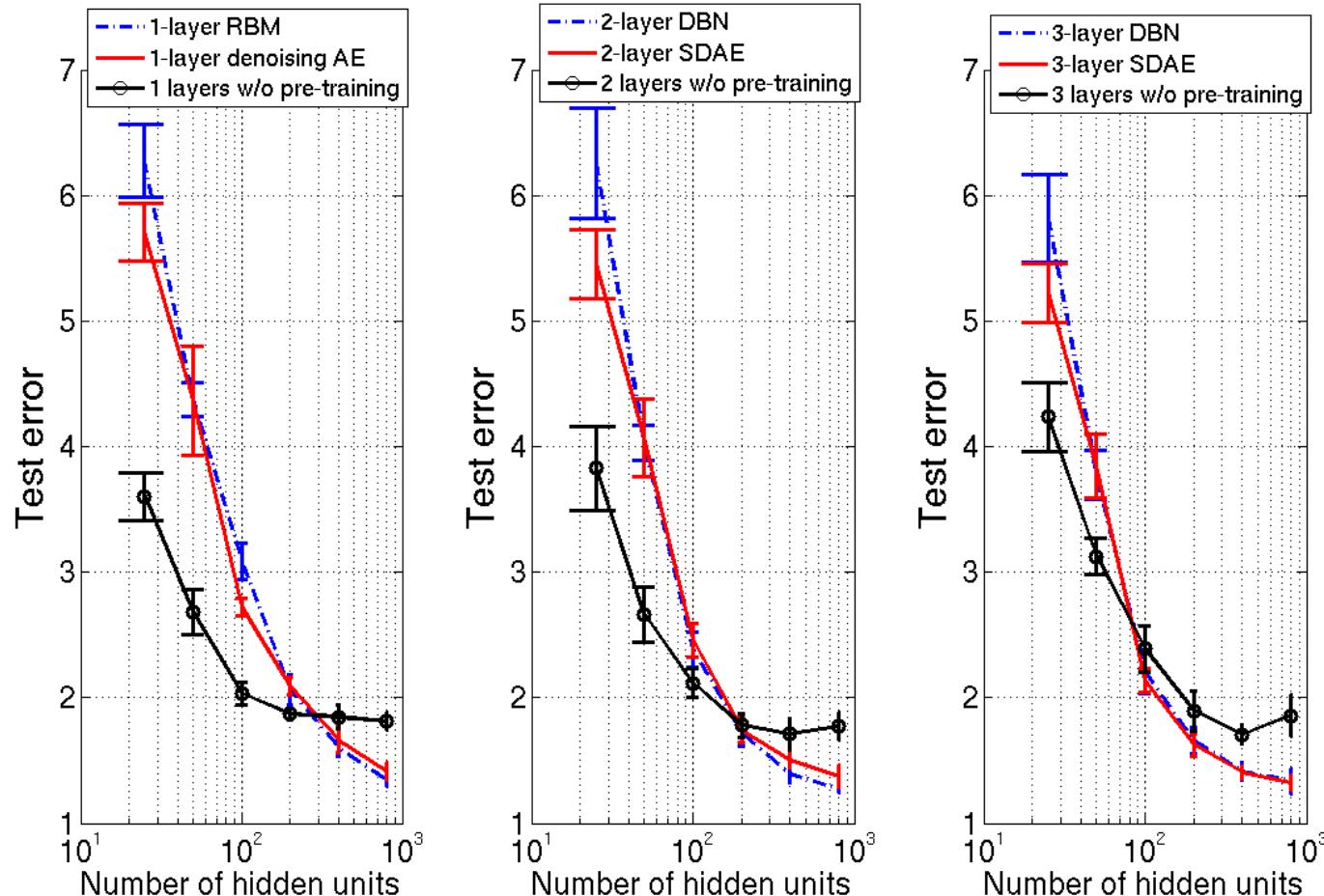
Convex shape or not?



Impact of Initialization

| Network | | MNIST-small classif. test error | MNIST-rotation classif. test error |
|---------------------------------------|-------|------------------------------------|---------------------------------------|
| Type | Depth | | |
| Neural network Deep net | 1 | 4.14 % ± 0.17 | 15.22 % ± 0.31 |
| | 2 | 4.03 % ± 0.17 | 10.63 % ± 0.27 |
| | 3 | 4.24 % ± 0.18 | 11.98 % ± 0.28 |
| | 4 | 4.47 % ± 0.18 | 11.73 % ± 0.29 |
| Deep net + autoencoder | 1 | 3.87 % ± 0.17 | 11.43% ± 0.28 |
| | 2 | 3.38 % ± 0.16 | 9.88 % ± 0.26 |
| | 3 | 3.37 % ± 0.16 | 9.22 % ± 0.25 |
| | 4 | 3.39 % ± 0.16 | 9.20 % ± 0.25 |
| Deep net + RBM | 1 | 3.17 % ± 0.15 | 10.47 % ± 0.27 |
| | 2 | 2.74 % ± 0.14 | 9.54 % ± 0.26 |
| | 3 | 2.71 % ± 0.14 | 8.80 % ± 0.25 |
| | 4 | 2.72 % ± 0.14 | 8.83 % ± 0.24 |

Impact of Pretraining



Acts as a regularizer: overfits less with large capacity, underfits with small capacity

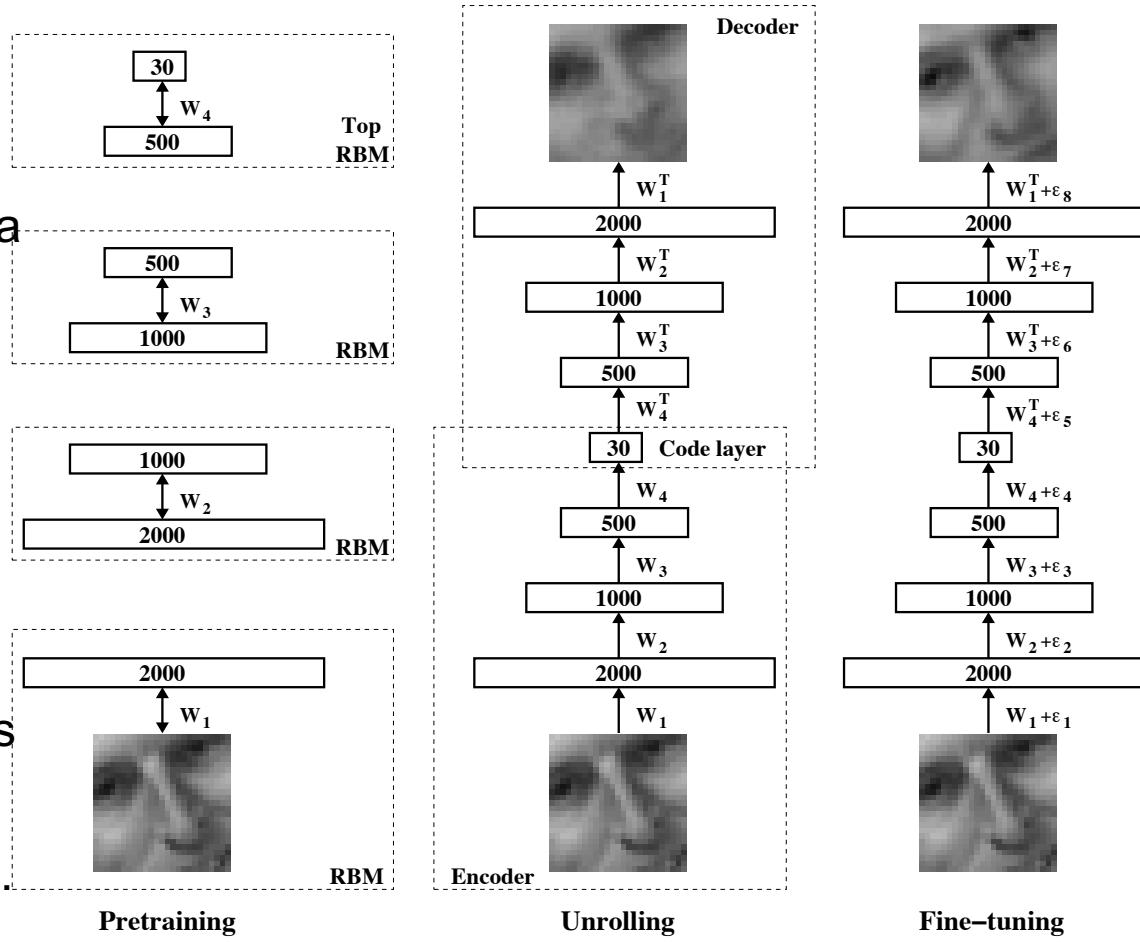
Performance on Different Datasets

| Stacked Autoencoders | Stacked RBMS | Stacked Denoising Autoencoders |
|-------------------------|-------------------|-----------------------------------|
| SAA-3 | DBN-3 | SdA-3 (ν) |
| 3.46±0.16 | 3.11±0.15 | 2.80±0.14 (10%) |
| 10.30±0.27 | 10.30±0.27 | 10.29±0.27 (10%) |
| 11.28±0.28 | 6.73±0.22 | 10.38±0.27 (40%) |
| 23.00±0.37 | 16.31±0.32 | 16.68±0.33 (25%) |
| 51.93±0.44 | 47.39±0.44 | 44.49±0.44 (25%) |
| 2.41±0.13 | 2.60±0.14 | 1.99±0.12 (10%) |
| 24.05±0.37 | 22.50±0.37 | 21.59±0.36 (25%) |
| 18.41±0.34 | 18.63±0.34 | 19.06±0.34 (10%) |

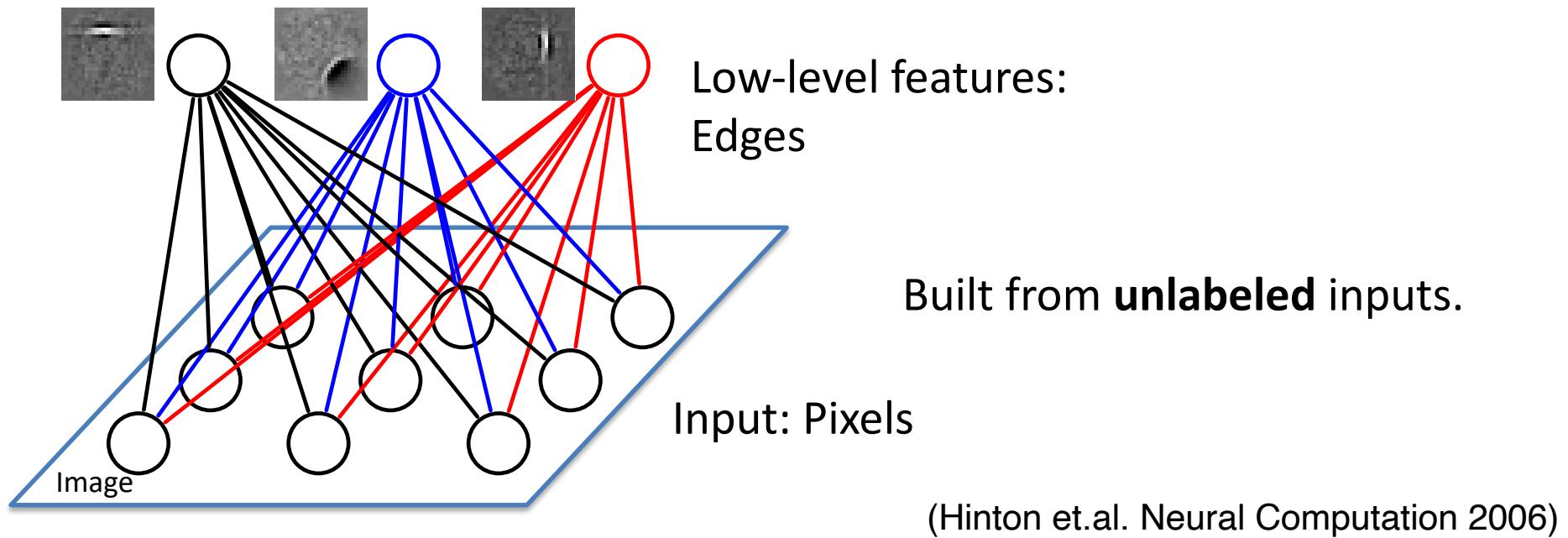
Deep Autoencoder

- Pre-training can be used to initialize a deep autoencoder

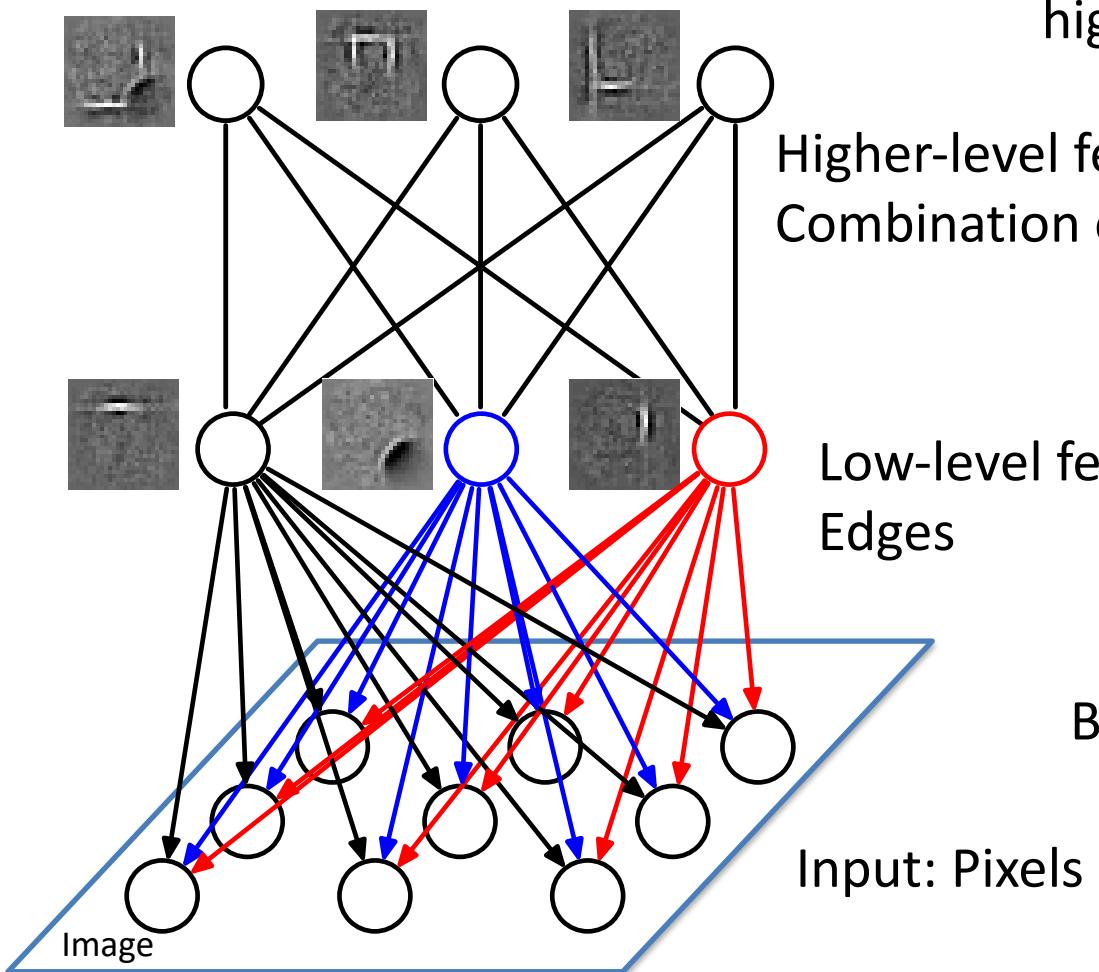
➤ Pre-training initializes the optimization problem in a region with better local optima of the training objective



Deep Belief Network



Deep Belief Network



Internal representations capture
higher-order statistical structure

Higher-level features:
Combination of edges

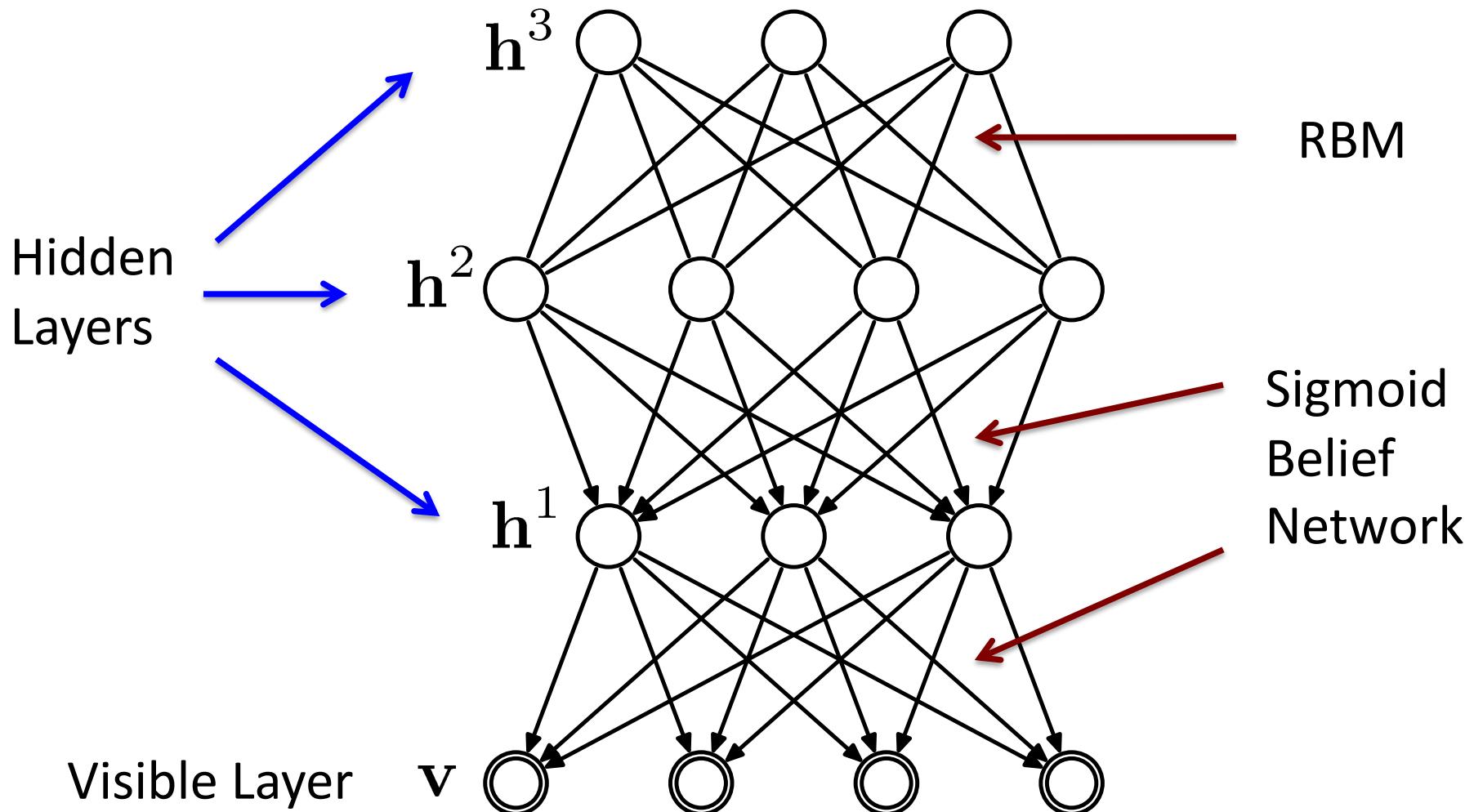
Low-level features:
Edges

Built from **unlabeled** inputs.

Input: Pixels

(Hinton et.al. Neural Computation 2006)

Deep Belief Network



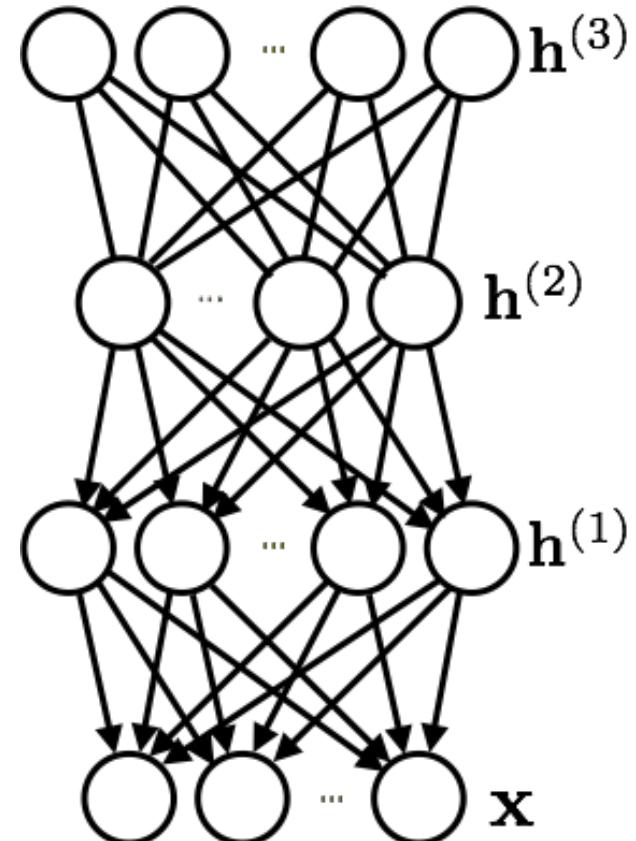
Deep Belief Network

- Deep Belief Networks:

- it is a **generative model** that mixes undirected and directed connections between variables
- top 2 layers' distribution $p(\mathbf{h}^{(2)}, \mathbf{h}^{(3)})$ is an RBM!
- other layers form a **Bayesian network** with conditional distributions:

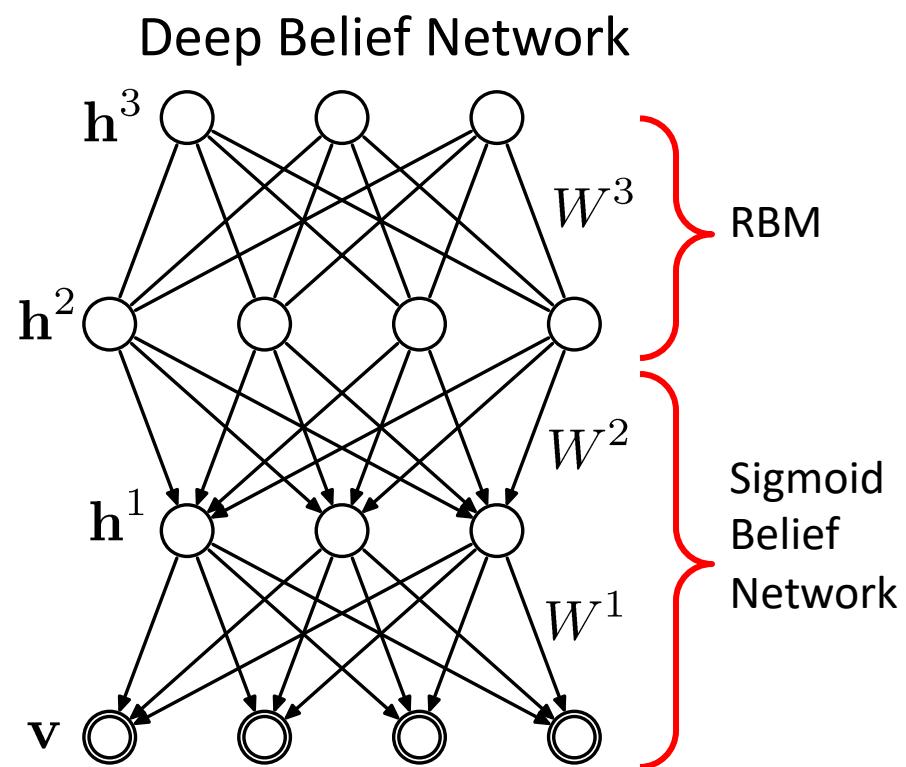
$$p(h_j^{(1)} = 1 | \mathbf{h}^{(2)}) = \text{sigm}(\mathbf{b}^{(1)} + \mathbf{W}^{(2)^\top} \mathbf{h}^{(2)})$$

$$p(x_i = 1 | \mathbf{h}^{(1)}) = \text{sigm}(\mathbf{b}^{(0)} + \mathbf{W}^{(1)^\top} \mathbf{h}^{(1)})$$



- This is **not a feed-forward neural network**

Deep Belief Network



- top 2 layers' distribution $p(\mathbf{h}^{(2)}, \mathbf{h}^{(3)})$ is an RBM
- other layers form a **Bayesian network** with conditional distributions:

$$p(h_j^{(1)} = 1 | \mathbf{h}^{(2)}) = \text{sigm}(\mathbf{b}^{(1)} + \mathbf{W}^{(2)^\top} \mathbf{h}^{(2)})$$

$$p(x_i = 1 | \mathbf{h}^{(1)}) = \text{sigm}(\mathbf{b}^{(0)} + \mathbf{W}^{(1)^\top} \mathbf{h}^{(1)})$$

Deep Belief Network

- The joint distribution of a DBN is as follows

$$p(\mathbf{x}, \mathbf{h}^{(1)}, \mathbf{h}^{(2)}, \mathbf{h}^{(3)}) = p(\mathbf{h}^{(2)}, \mathbf{h}^{(3)}) p(\mathbf{h}^{(1)} | \mathbf{h}^{(2)}) p(\mathbf{x} | \mathbf{h}^{(1)})$$

where

$$p(\mathbf{h}^{(2)}, \mathbf{h}^{(3)}) = \exp \left(\mathbf{h}^{(2)^\top} \mathbf{W}^{(3)} \mathbf{h}^{(3)} + \mathbf{b}^{(2)^\top} \mathbf{h}^{(2)} + \mathbf{b}^{(3)^\top} \mathbf{h}^{(3)} \right) / Z$$

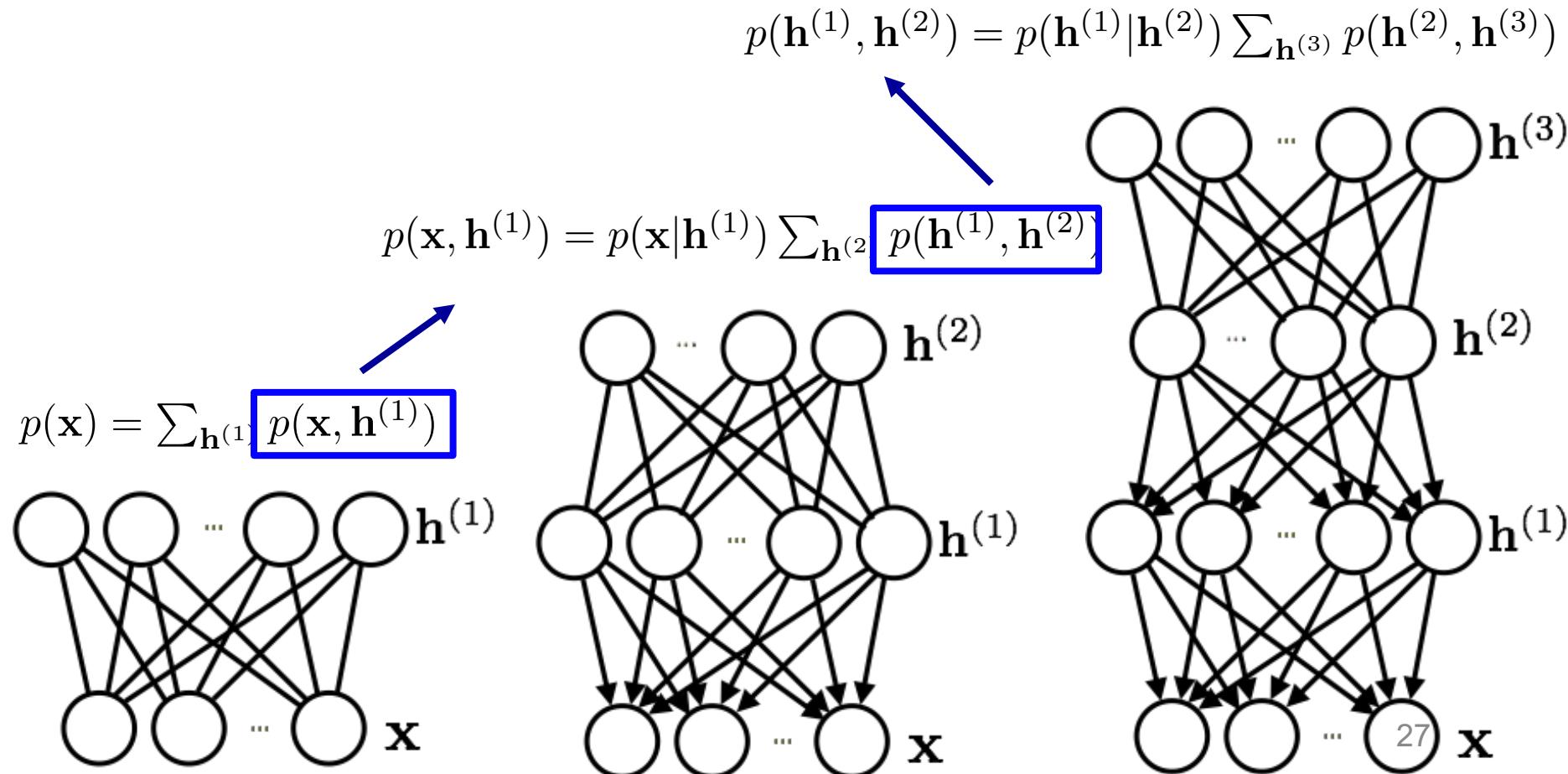
$$p(\mathbf{h}^{(1)} | \mathbf{h}^{(2)}) = \prod_j p(h_j^{(1)} | \mathbf{h}^{(2)})$$

$$p(\mathbf{x} | \mathbf{h}^{(1)}) = \prod_i p(x_i | \mathbf{h}^{(1)})$$

- As in a deep feed-forward network, training a DBN is hard

Layer-wise Pretraining

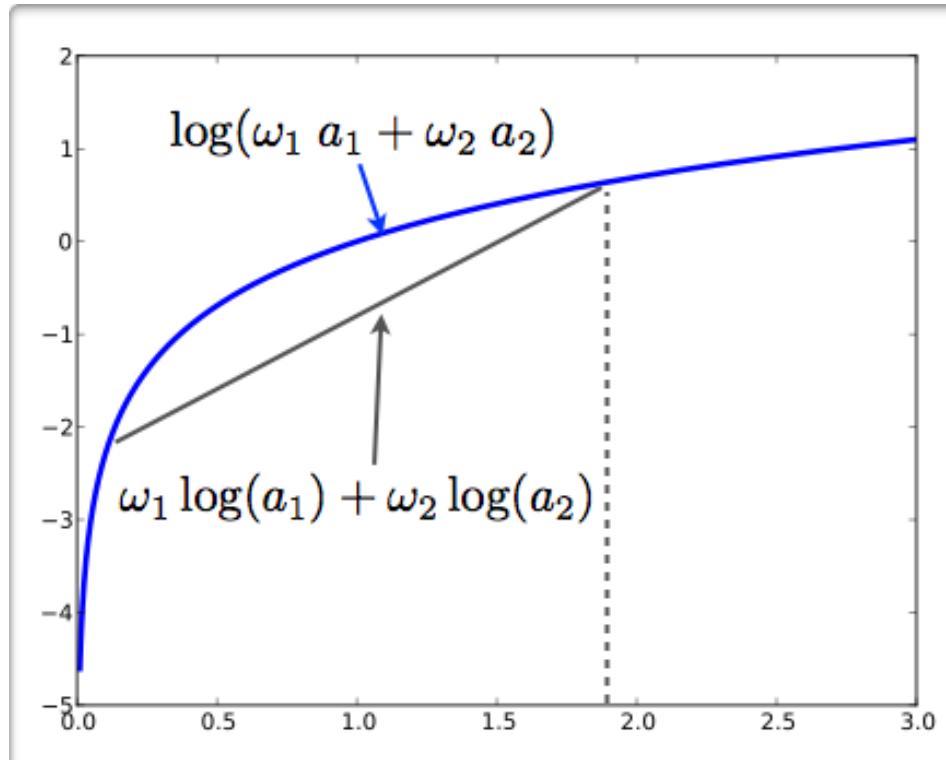
- This is where the RBM stacking procedure comes from:
 - **idea:** improve prior on last layer by adding another hidden layer



Concavity

$$\log(\sum_i \omega_i a_i) \geq \sum_i \omega_i \log(a_i)$$

(where $\sum_i \omega_i = 1$ and $\omega_i \geq 0$)



Variational Bound

- For any model $p(\mathbf{x}, \mathbf{h}^{(1)})$ with latent variables $\mathbf{h}^{(1)}$ we can write:

$$\begin{aligned}\log p(\mathbf{x}) &= \log \left(\sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \frac{p(\mathbf{x}, \mathbf{h}^{(1)})}{q(\mathbf{h}^{(1)} | \mathbf{x})} \right) \\ &\geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log \left(\frac{p(\mathbf{x}, \mathbf{h}^{(1)})}{q(\mathbf{h}^{(1)} | \mathbf{x})} \right) \\ &= \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log p(\mathbf{x}, \mathbf{h}^{(1)}) \\ &\quad - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log q(\mathbf{h}^{(1)} | \mathbf{x})\end{aligned}$$

where $q(\mathbf{h}^{(1)} | \mathbf{x})$ is any **approximation** to $p(\mathbf{h}^{(1)} | \mathbf{x})$

Variational Bound

- This is called a **variational bound**

$$\begin{aligned}\log p(\mathbf{x}) &\geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log p(\mathbf{x}, \mathbf{h}^{(1)}) \\ &\quad - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log q(\mathbf{h}^{(1)} | \mathbf{x})\end{aligned}$$

- if $q(\mathbf{h}^{(1)} | \mathbf{x})$ is equal to the true conditional $p(\mathbf{h}^{(1)} | \mathbf{x})$, then we have an equality – **the bound is tight!**
- the more $q(\mathbf{h}^{(1)} | \mathbf{x})$ is different from $p(\mathbf{h}^{(1)} | \mathbf{x})$ the less tight the bound is.

Variational Bound

- This is called a variational bound

$$\begin{aligned}\log p(\mathbf{x}) &\geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log p(\mathbf{x}, \mathbf{h}^{(1)}) \\ &\quad - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log q(\mathbf{h}^{(1)}|\mathbf{x})\end{aligned}$$

- In fact, difference between the left and right terms is the **KL divergence** between $q(\mathbf{h}^{(1)}|\mathbf{x})$ and $p(\mathbf{h}^{(1)}|\mathbf{x})$:

$$\text{KL}(q||p) = \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log \left(\frac{q(\mathbf{h}^{(1)}|\mathbf{x})}{p(\mathbf{h}^{(1)}|\mathbf{x})} \right)$$

Variational Bound

- This is called a variational bound

$$\begin{aligned}\log p(\mathbf{x}) \geq & \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \left(\log p(\mathbf{x}|\mathbf{h}^{(1)}) + \log p(\mathbf{h}^{(1)}) \right) \\ & - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log q(\mathbf{h}^{(1)}|\mathbf{x})\end{aligned}$$

- for a single hidden layer DBN (i.e. an RBM), both **the likelihood** $p(\mathbf{x}|\mathbf{h}^{(1)})$ and **the prior** $p(\mathbf{h}^{(1)})$ depend on the parameters of the first layer.
- we can now improve the model by building a better prior $p(\mathbf{h}^{(1)})$

Variational Bound

- This is called a variational bound

adding 2nd layer means
untying the parameters

$$\log p(\mathbf{x}) \geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \left(\log p(\mathbf{x} | \mathbf{h}^{(1)}) + \log p(\mathbf{h}^{(1)}) \right) - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)} | \mathbf{x}) \log q(\mathbf{h}^{(1)} | \mathbf{x})$$

- When adding a second layer, we model $p(\mathbf{h}^{(1)})$ using a separate set of parameters

- they are the parameters of the RBM involving $\mathbf{h}^{(1)}$ and $\mathbf{h}^{(2)}$
- $p(\mathbf{h}^{(1)})$ is now the marginalization of the second hidden layer

$$p(\mathbf{h}^{(1)}) = \sum_{\mathbf{h}^{(2)}} p(\mathbf{h}^{(1)}, \mathbf{h}^{(2)})$$

Variational Bound

- This is called a variational bound

adding 2nd layer means
untying the parameters

$$\log p(\mathbf{x}) \geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \left(\log p(\mathbf{x}|\mathbf{h}^{(1)}) + \log p(\mathbf{h}^{(1)}) \right) - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log q(\mathbf{h}^{(1)}|\mathbf{x})$$

- we can train the parameters of the bound. This is equivalent if other terms are constant:

Layerwise pretraining
improves variational
lower bound

$$- \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log p(\mathbf{h}^{(1)})$$

- this is like training an RBM on data generated from $q(\mathbf{h}^{(1)}|\mathbf{x})$!

Variational Bound

- This is called a variational bound

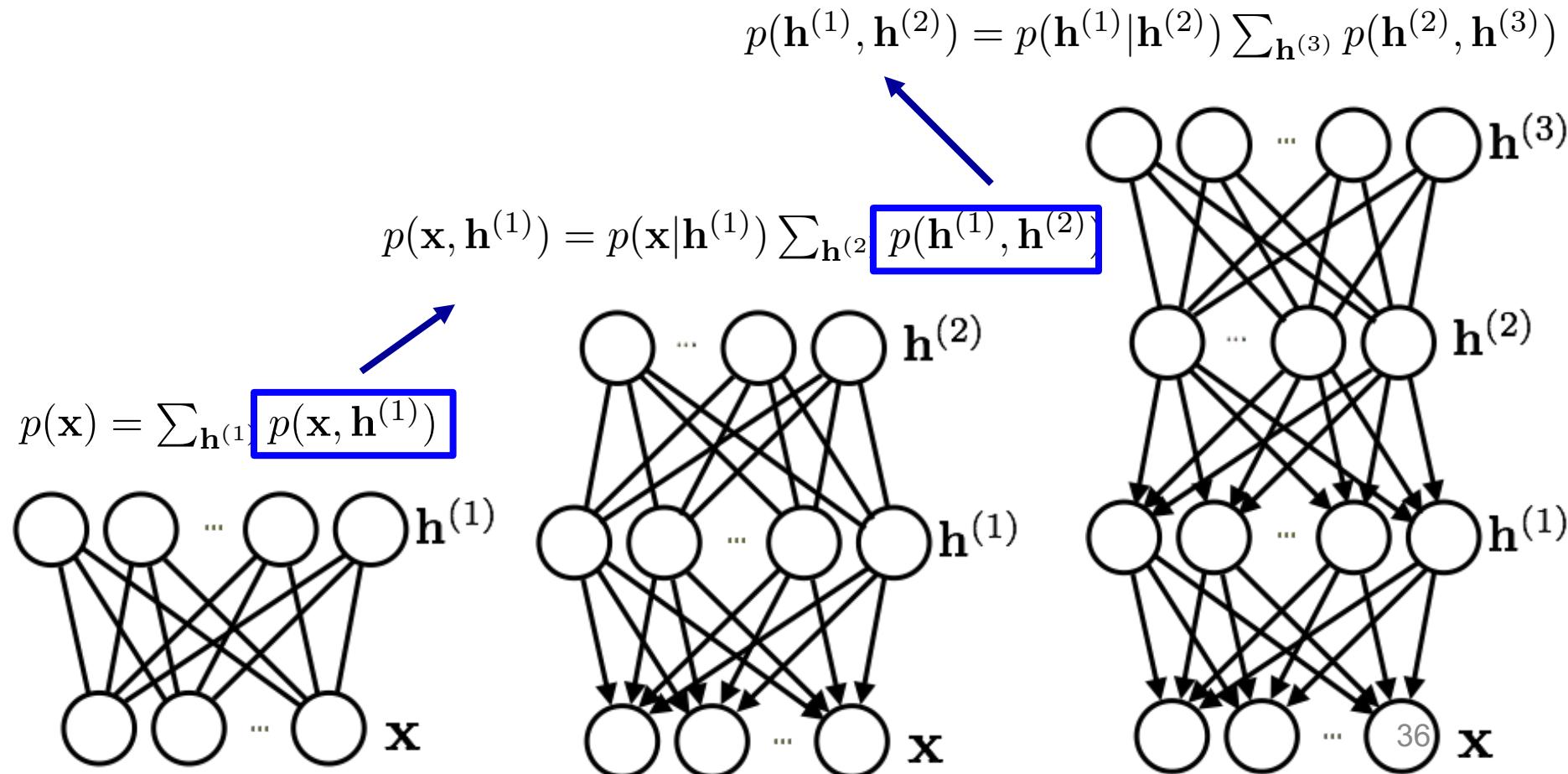
adding 2nd layer means
untying the parameters

$$\log p(\mathbf{x}) \geq \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \left(\log p(\mathbf{x}|\mathbf{h}^{(1)}) + \log p(\mathbf{h}^{(1)}) \right) - \sum_{\mathbf{h}^{(1)}} q(\mathbf{h}^{(1)}|\mathbf{x}) \log q(\mathbf{h}^{(1)}|\mathbf{x})$$

- for $q(\mathbf{h}^{(1)}|\mathbf{x})$ we use **the posterior of the first layer RBM**. This is equivalent to a feed-forward (sigmoidal) layer, followed by sampling
- by initializing the weights of the second layer RBM as the transpose of the first layer weights, **the bound is initially tight!**
- a 2-layer DBN with tied weights is equivalent to a 1-layer RBM

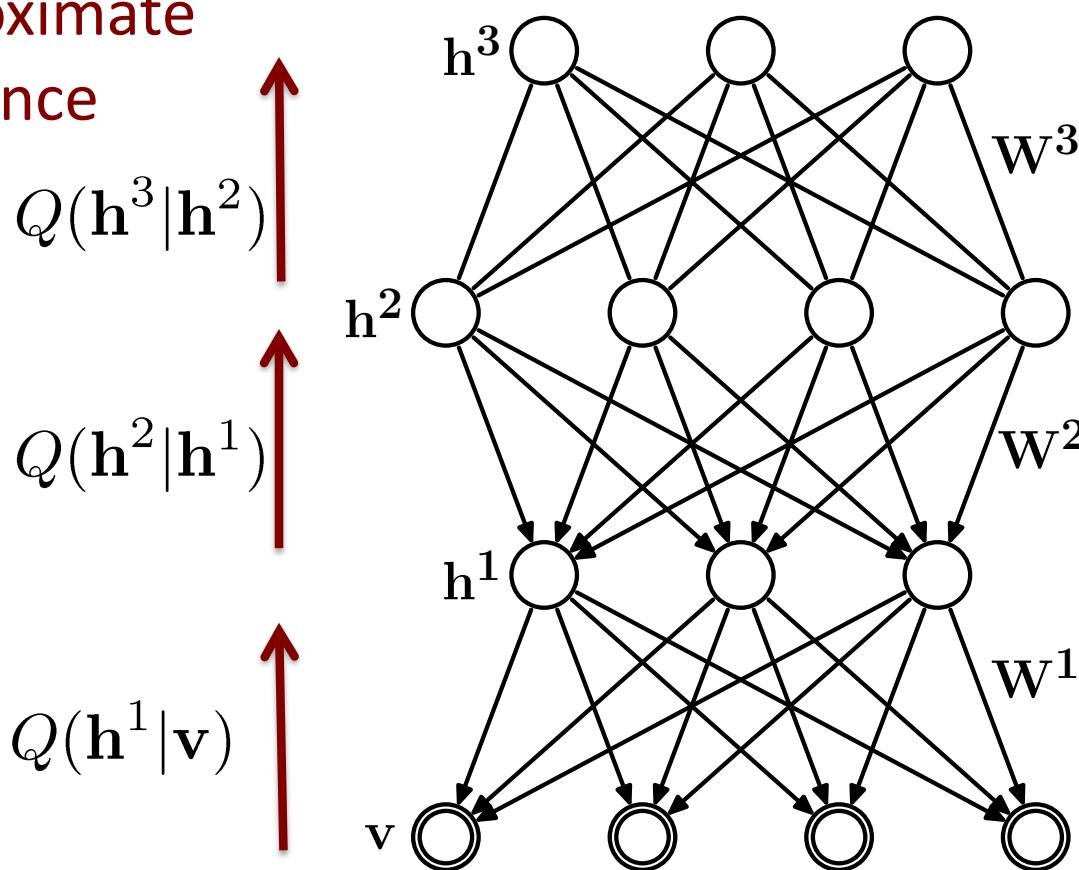
Layer-wise Pretraining

- This is where the RBM stacking procedure comes from:
 - **idea:** improve prior on last layer by adding another hidden layer



Deep Belief Network

Approximate
Inference



Generative
Process

$$P(h^2, h^3)$$

$$P(h^1|h^2)$$

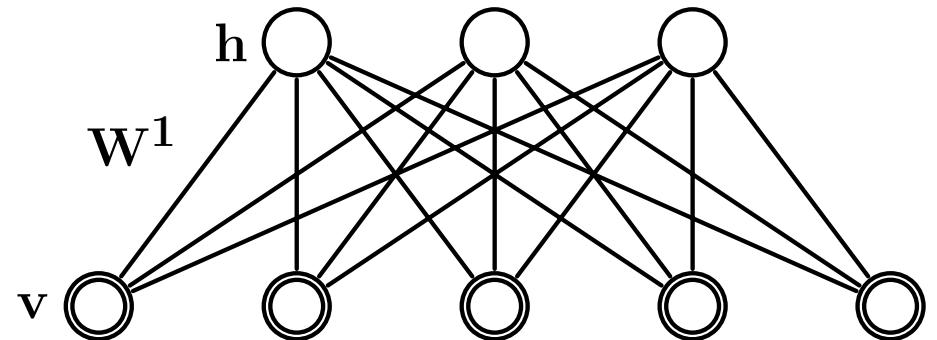
$$P(v|h^1)$$

$$Q(h^t|h^{t-1}) = \prod_j \sigma \left(\sum_i W^t h_i^{t-1} \right)$$

$$P(h^{t-1}|h^t) = \prod_j \sigma \left(\sum_i W^t h_i^t \right)$$

DBN Layer-wise Training

- Learn an RBM with an input layer $v=x$ and a hidden layer h .



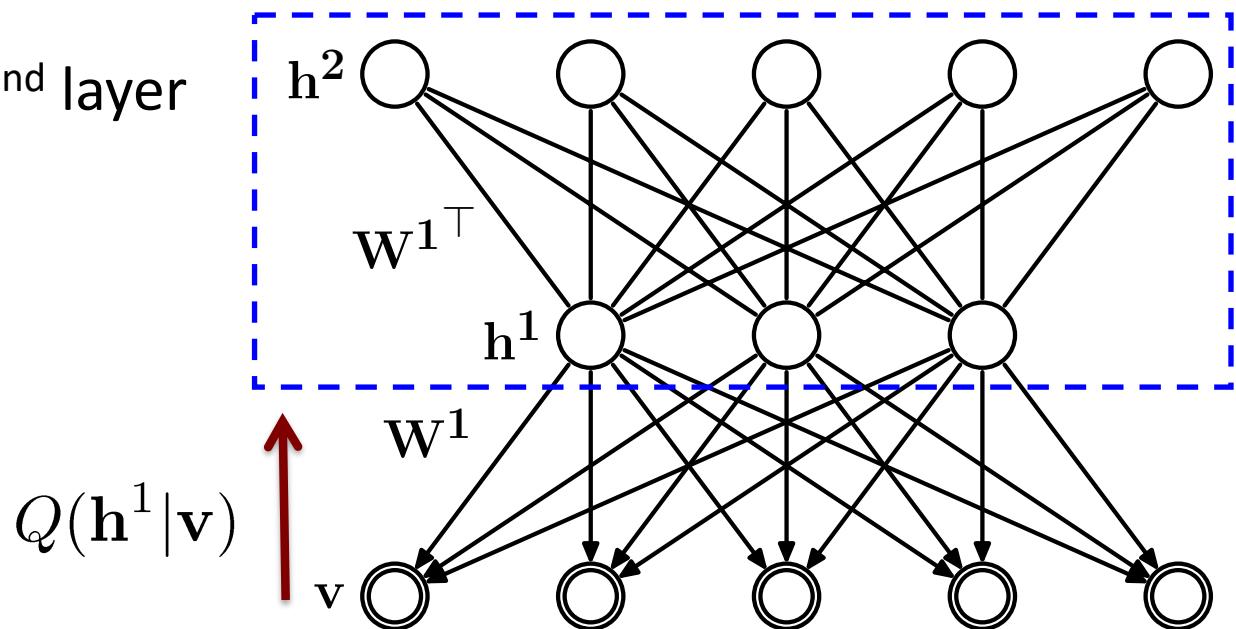
DBN Layer-wise Training

- Learn an RBM with an input layer $v=x$ and a hidden layer h .

- Treat inferred values

$Q(h^1|v) = P(h^1|v)$ as the data for training 2nd-layer RBM.

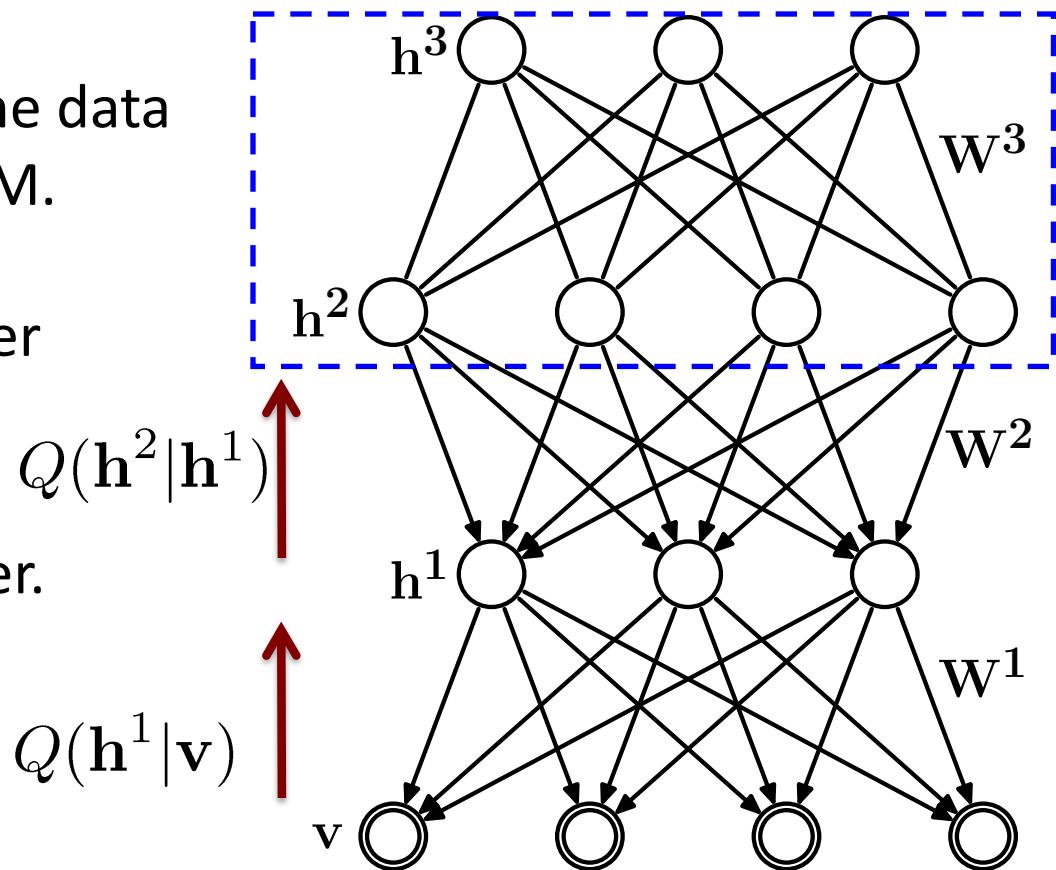
- Learn and freeze 2nd layer RBM.



DBN Layer-wise Training

- Learn an RBM with an input layer $v=x$ and a hidden layer h .
- Treat inferred values $Q(h^1|v) = P(h^1|v)$ as the data for training 2nd-layer RBM.
- Learn and freeze 2nd layer RBM.
- Proceed to the next layer.

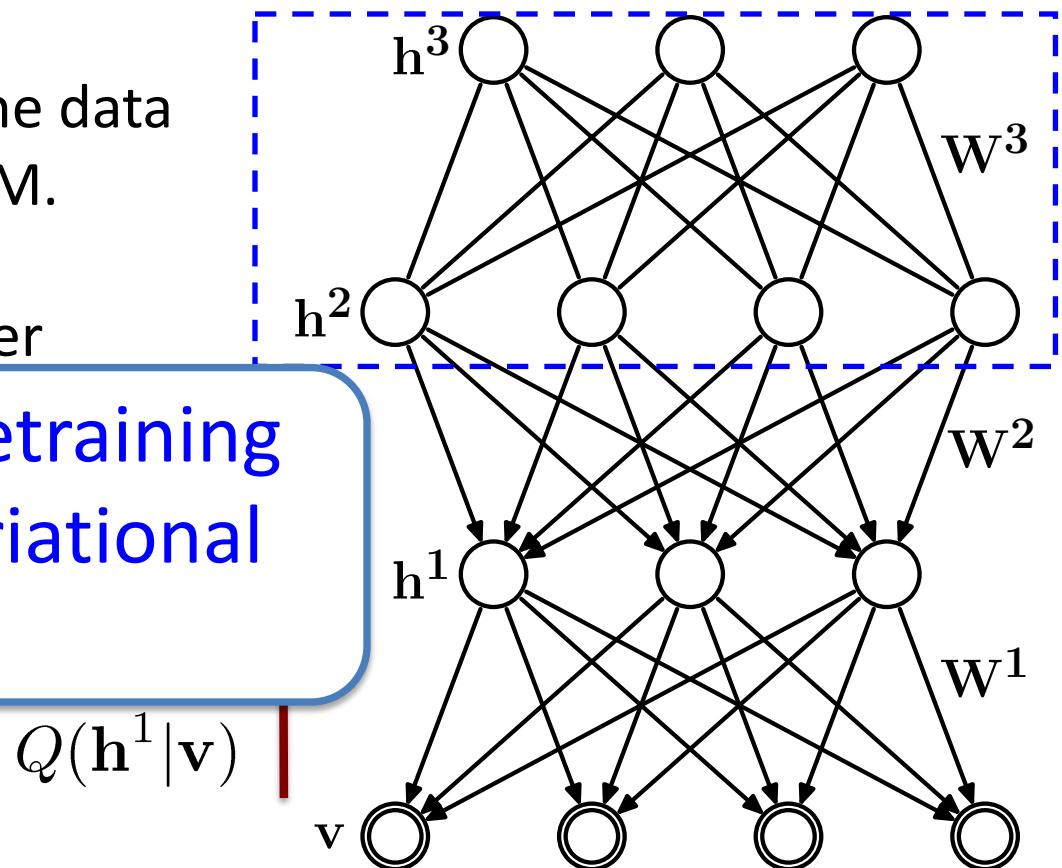
Unsupervised Feature Learning.



DBN Layer-wise Training

- Learn an RBM with an input layer $v=x$ and a hidden layer h .
- Treat inferred values $Q(h^1|v) = P(h^1|v)$ as the data for training 2nd-layer RBM.
- Learn and freeze 2nd layer RBM
- Proc

Unsupervised Feature Learning.



Deep Belief Networks

- This process of adding layers can be repeated recursively
 - we obtain **the greedy layer-wise pre-training** procedure for neural networks
- We now see that this procedure corresponds to **maximizing a bound on the likelihood of the data** in a DBN
 - in theory, if our approximation $q(\mathbf{h}^{(1)} | \mathbf{x})$ is very far from the true posterior, the bound might be very loose
 - this only means we might not be improving the true likelihood
 - we might still be extracting better features!
- Fine-tuning is done by the Up-Down algorithm
 - A fast learning algorithm for deep belief nets. Hinton, Teh, Osindero, 2006.

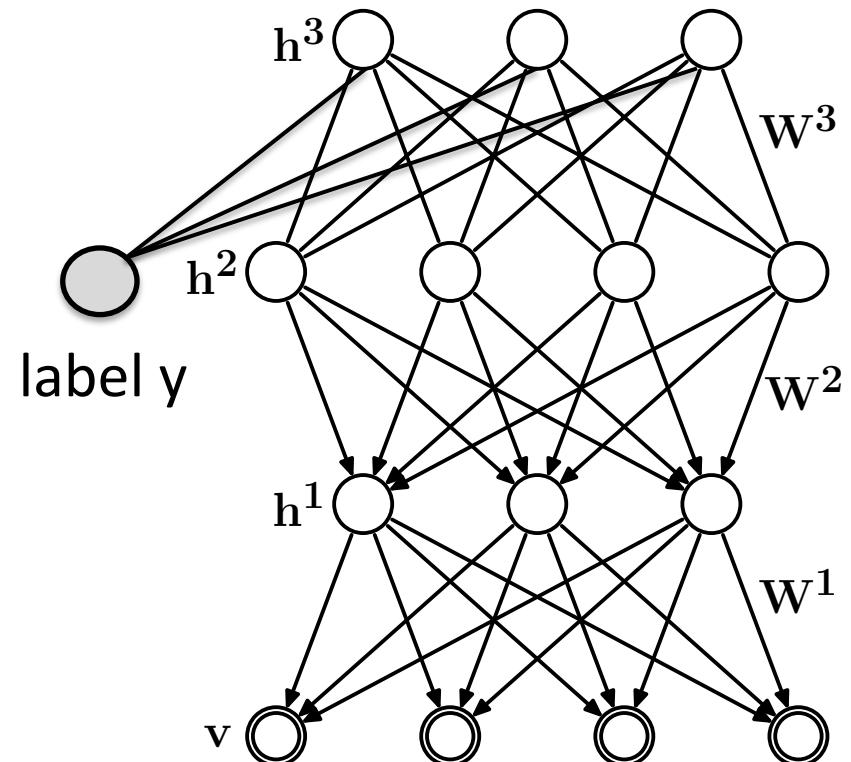
Supervised Learning with DBNs

- If we have access to label information, we can train **the joint generative model** by maximizing the joint log-likelihood of data and labels

$$\log P(\mathbf{y}, \mathbf{v})$$

- Discriminative fine-tuning:
 - Use DBN to initialize a multilayer neural network.
 - Maximize **the conditional distribution**:

$$\log P(\mathbf{y}|\mathbf{v})$$

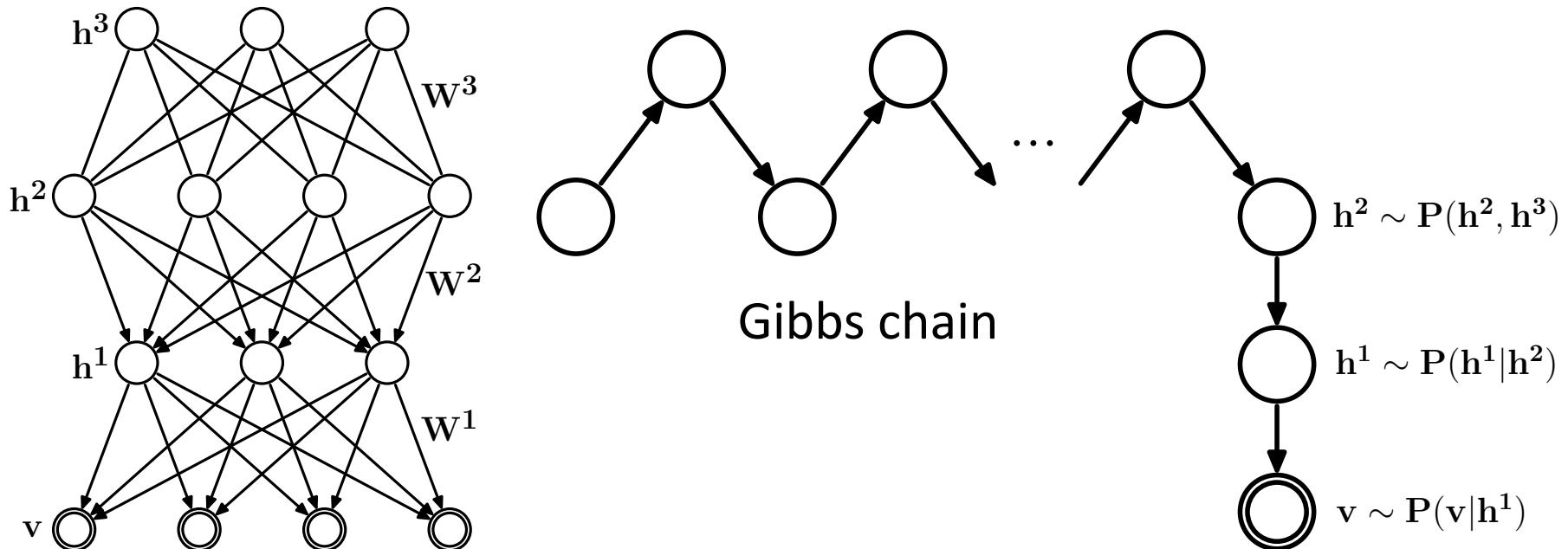


Sampling from DBNs

- To sample from the DBN model:

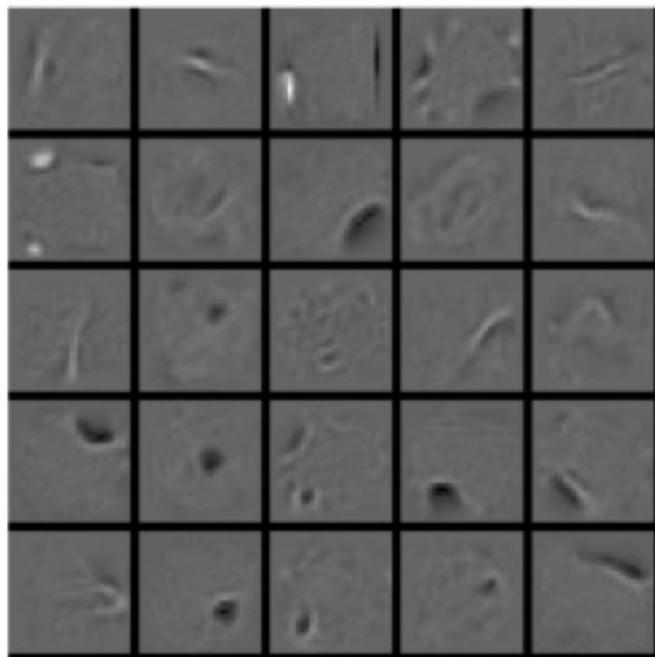
$$P(\mathbf{v}, \mathbf{h}^1, \mathbf{h}^2, \mathbf{h}^3) = P(\mathbf{v}|\mathbf{h}^1)P(\mathbf{h}^1|\mathbf{h}^2)P(\mathbf{h}^2, \mathbf{h}^3)$$

- Sample \mathbf{h}^2 using alternating Gibbs sampling from RBM.
- Sample lower layers using sigmoid belief network.

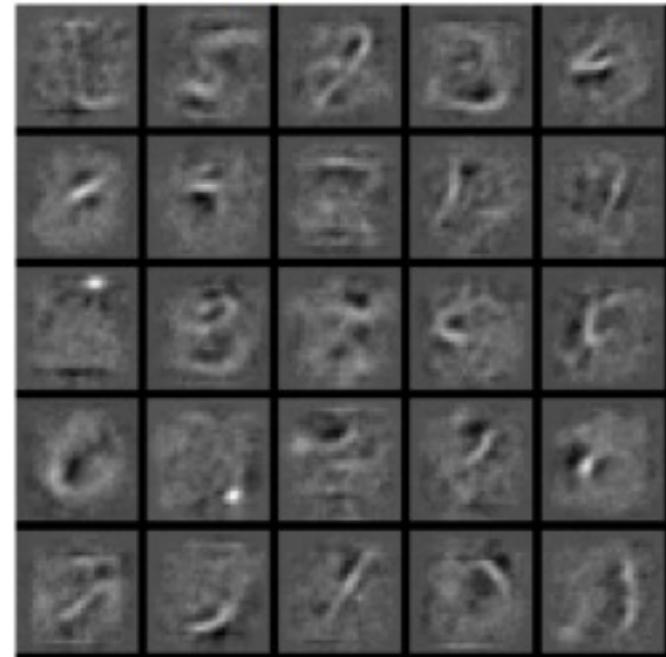


Learned Features

1st-layer features

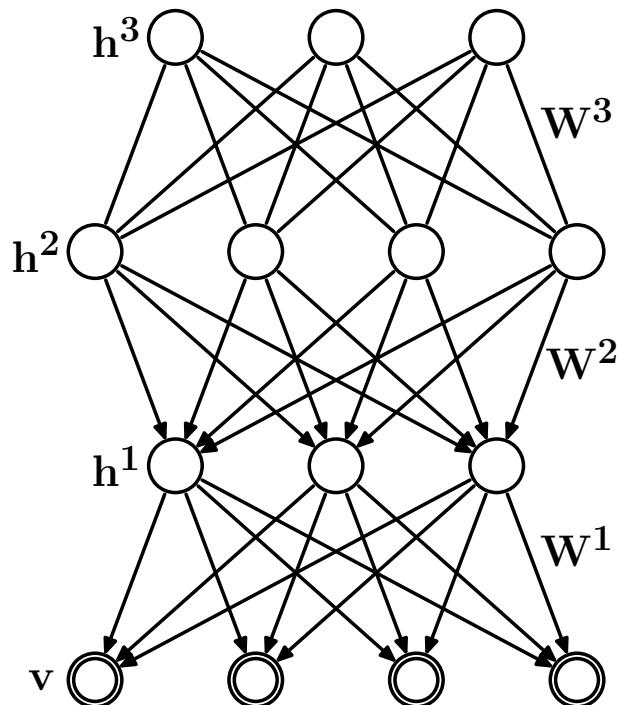


2nd-layer features

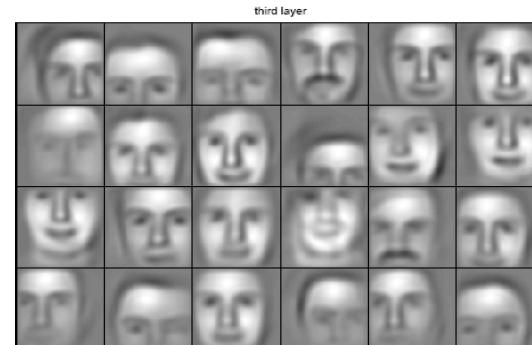


Learning Part-based Representation

Convolutional DBN



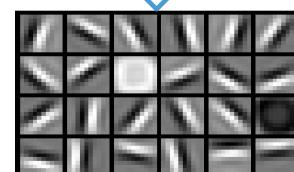
Faces



Groups of parts.



Object Parts



Trained on face images.

Learning Part-based Representation

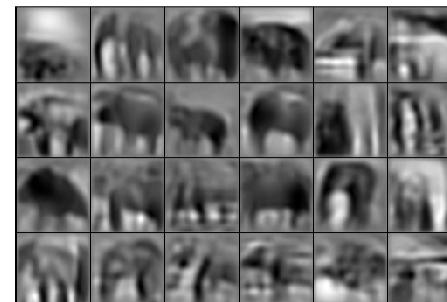
Faces



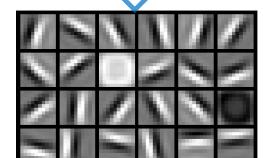
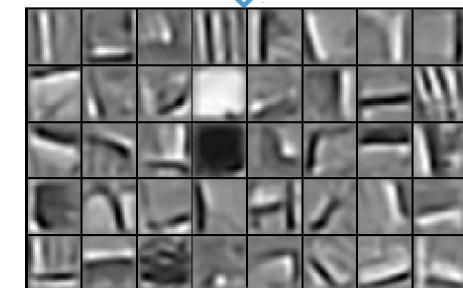
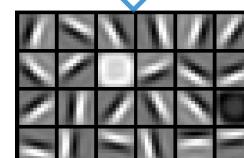
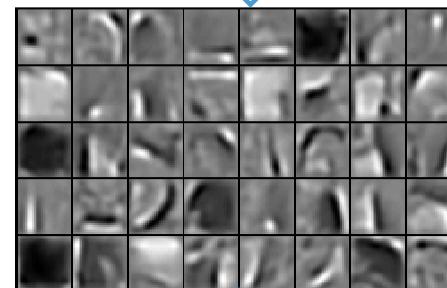
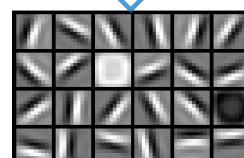
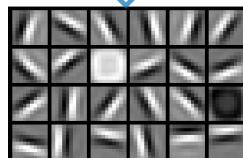
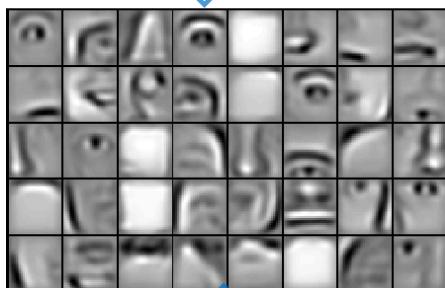
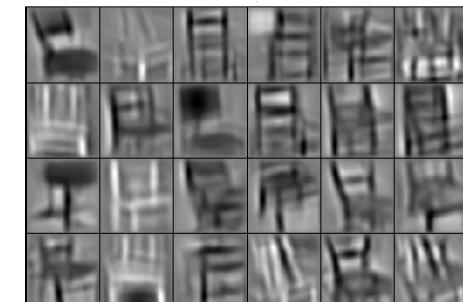
Cars



Elephants



Chairs

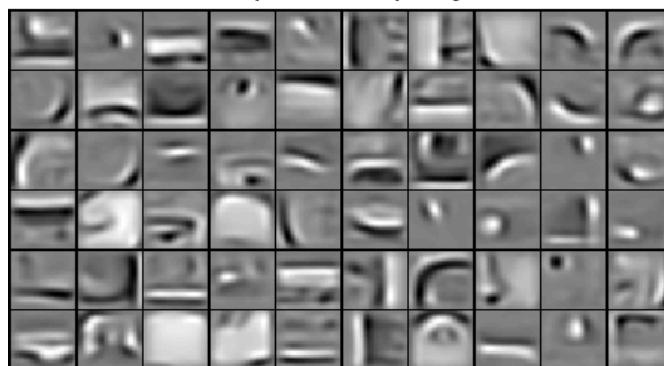


Learning Part-based Representation

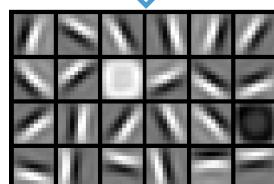


Groups of parts.

second layer learned from 4 object categories

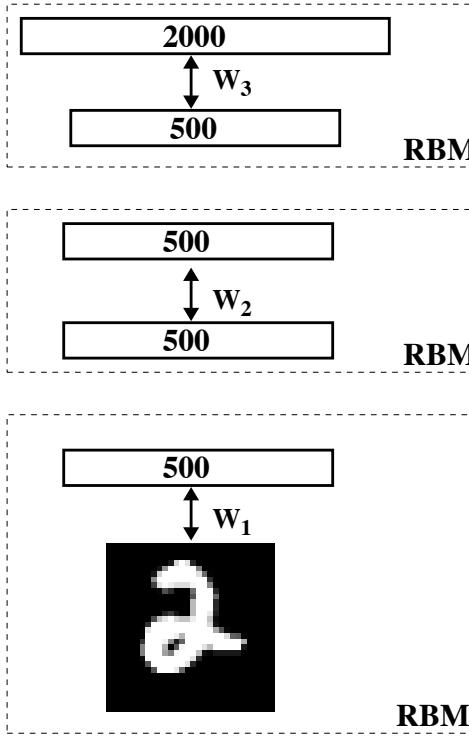


Class-specific object
parts



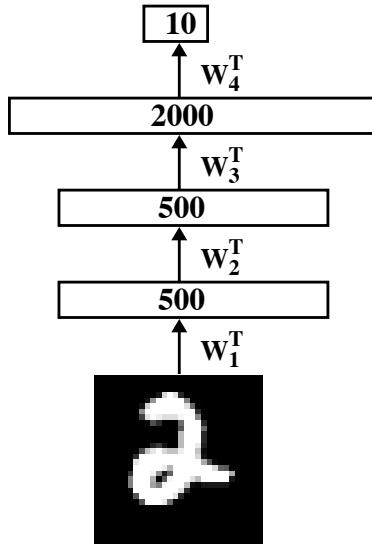
Trained from multiple
classes (cars, faces,
motorbikes, airplanes).

DBNs for Classification

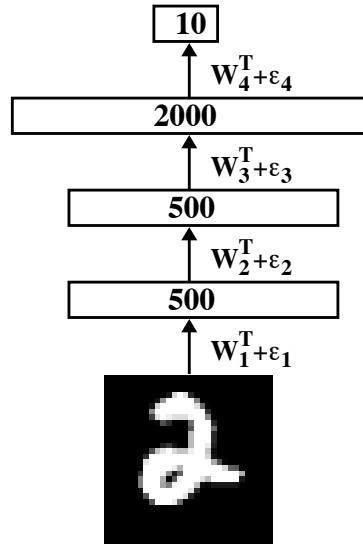


Pretraining

Softmax Output



Unrolling



Fine-tuning

- After layer-by-layer **unsupervised pretraining**, discriminative fine-tuning by backpropagation achieves an error rate of 1.2% on MNIST. SVM's get 1.4% and randomly initialized backprop gets 1.6%.
- Clearly unsupervised learning helps generalization. It ensures that most of the information in the weights comes from modeling the input data.

DBNs for Regression

Predicting the orientation of a face patch



Training Data: 1000 face patches of 30 training people.

Test Data: 1000 face patches of **10 new people.**

Regression Task: predict orientation of a new face.

Gaussian Processes with spherical Gaussian kernel achieves a RMSE (root mean squared error) of 16.33 degree.

DBNs for Regression

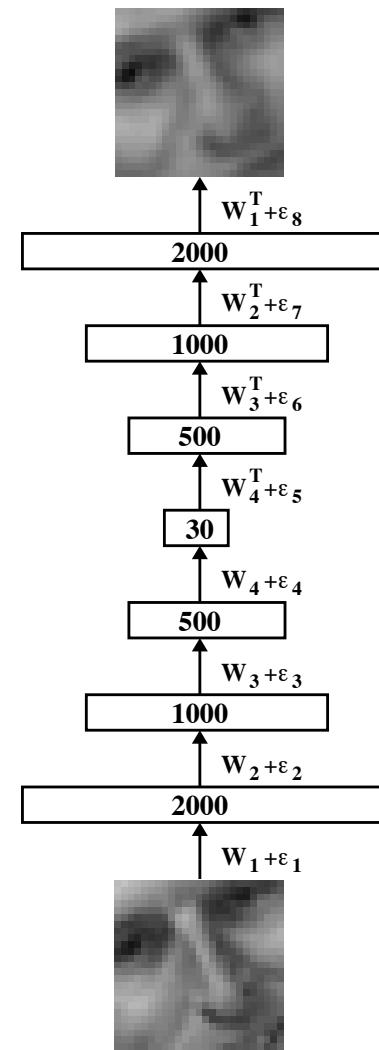
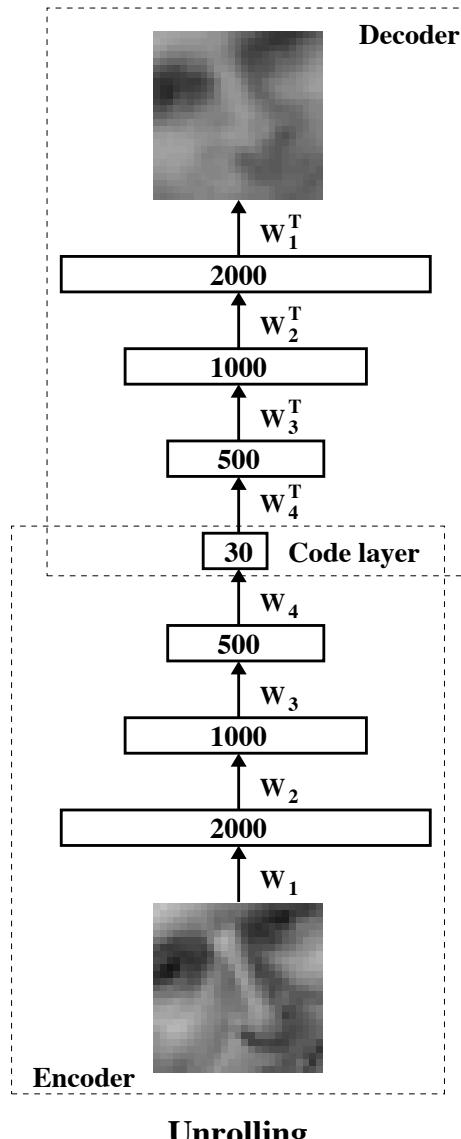
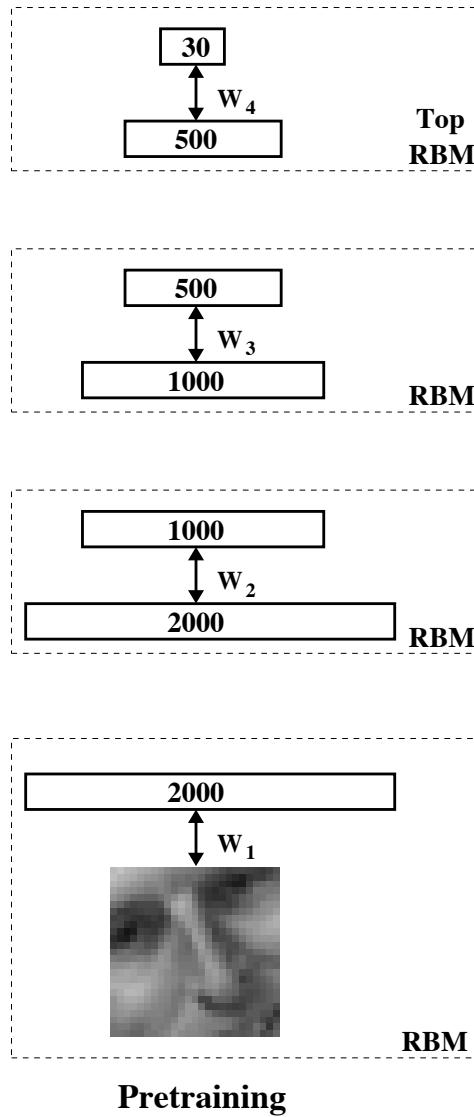


Additional Unlabeled Training Data: 12000 face patches from 30 training people.

- Pretrain a stack of RBMs: 784-1000-1000-1000.
- **Features were extracted with no idea of the final task.**

| | |
|--|-------------|
| The same GP on the top-level features: | RMSE: 11.22 |
| GP with fine-tuned covariance Gaussian kernel: | RMSE: 6.42 |
| Standard GP without using DBNs: | RMSE: 16.33 |

Deep Autoencoders



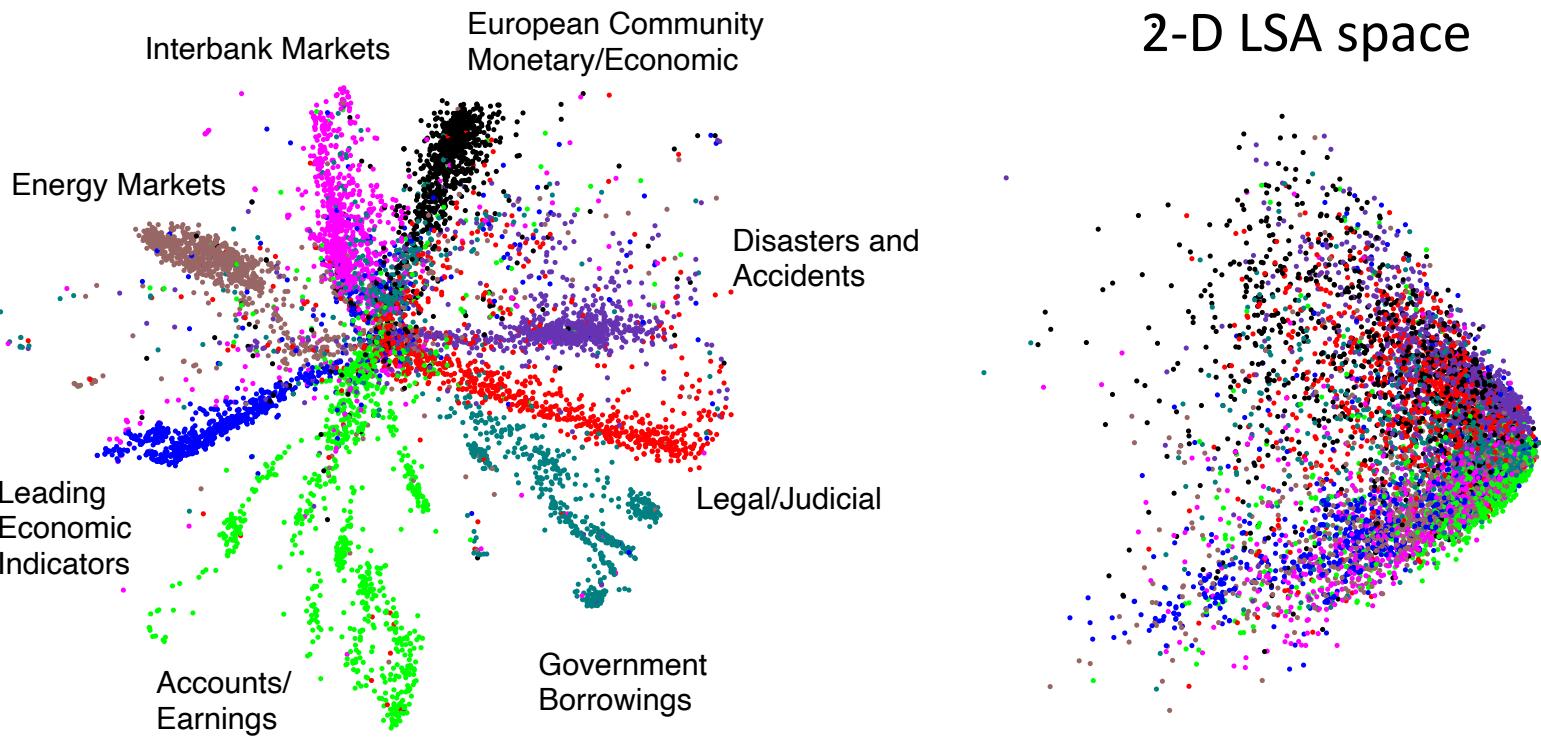
Deep Autoencoders

- We used $25 \times 25 - 2000 - 1000 - 500 - 30$ autoencoder to extract 30-D real-valued codes for Olivetti face patches.



- **Top:** Random samples from the test dataset.
- **Middle:** Reconstructions by the 30-dimensional deep autoencoder.
- **Bottom:** Reconstructions by the 30-dimensional PCA.

Information Retrieval



- The Reuters Corpus Volume II contains 804,414 newswire stories (randomly split into **402,207 training** and **402,207 test**).
- “Bag-of-words” representation: each article is represented as a vector containing the counts of the most frequently used 2000 words in the training set.