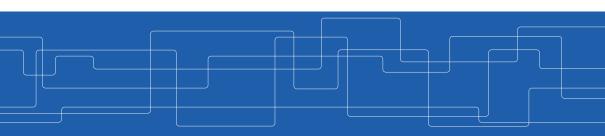


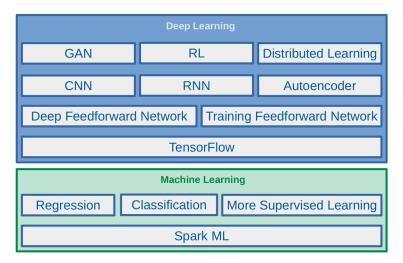
Autoencoders and Restricted Boltzmann Machines

Amir H. Payberah payberah@kth.se 27/11/2019

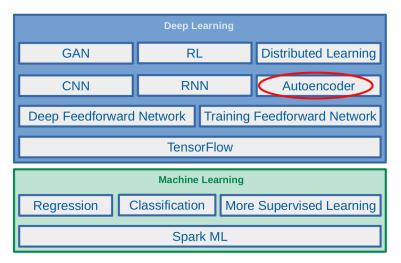


https://id2223kth.github.io











Let's Start With An Example



► Which of them is easier to memorize?



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► Seq1: 40,27,25,36,81,57,10,73,19,68



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► Seq1: 40,27,25,36,81,57,10,73,19,68

► Seq2: 50, 25, 76, 38, 19, 58, 29, 88, 44, 22, 11, 34, 17, 52, 26, 13, 40, 20



 $\mathtt{Seq2} : 50, 25, 76, 38, 19, 58, 29, 88, 44, 22, 11, 34, 17, 52, 26, 13, 40, 20$





Seq2: 50, 25, 76, 38, 19, 58, 29, 88, 44, 22, 11, 34, 17, 52, 26, 13, 40, 20

► Seq1 is shorter, so it should be easier.





Seq2: 50, 25, 76, 38, 19, 58, 29, 88, 44, 22, 11, 34, 17, 52, 26, 13, 40, 20

▶ Seq1 is shorter, so it should be easier.

► But, Seq2 follows two simple rules:





Seq2: 50, 25, 76, 38, 19, 58, 29, 88, 44, 22, 11, 34, 17, 52, 26, 13, 40, 20

- ▶ Seq1 is shorter, so it should be easier.
- ▶ But, Seq2 follows two simple rules:
 - Even numbers are followed by their half.
 - Odd numbers are followed by their triple plus one.





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- ► You don't need pattern if you could quickly and easily memorize very long sequences





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- ▶ But, Seq2 follows two simple rules:
 - Even numbers are followed by their half.
 - Odd numbers are followed by their triple plus one.
- ► You don't need pattern if you could quickly and easily memorize very long sequences
- ▶ But, it is hard to memorize long sequences that makes it useful to recognize patterns.





- ▶ 1970, W. Chase and H. Simon
- ► They observed that expert chess players were able to memorize the positions of all the pieces in a game by looking at the board for just 5 seconds.





► This was only the case when the pieces were placed in realistic positions, not when the pieces were placed randomly.





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- ► Chess experts don't have a much better memory than you and I.
- ► They just see chess patterns more easily due to their experience with the game.
- ▶ Patterns helps them store information efficiently.

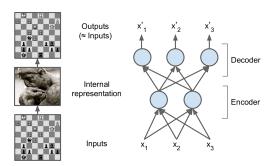




Autoencoders

Autoencoders (1/5)

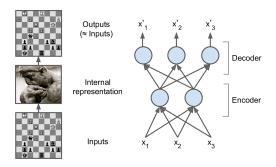
▶ Just like the chess players in this memory experiment.





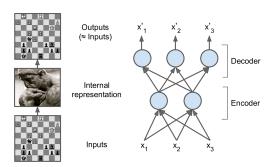
Autoencoders (1/5)

- ▶ Just like the chess players in this memory experiment.
- ▶ An autoencoder looks at the inputs, converts them to an efficient internal representation, and then spits out something that looks very close to the inputs.



Autoencoders (2/5)

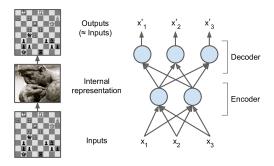
▶ The same architecture as a Multi-Layer Perceptron (MLP).





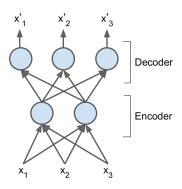
Autoencoders (2/5)

- ► The same architecture as a Multi-Layer Perceptron (MLP).
- Except that the number of neurons in the output layer must be equal to the number of inputs.



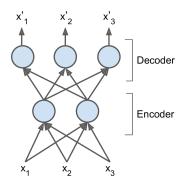


► An autoencoder is always composed of two parts.





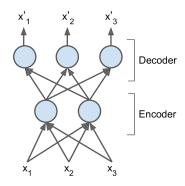
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- ► An encoder (recognition network), h = f(x)
 Converts the inputs to an internal representation.





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- ► An encoder (recognition network), **h** = **f**(**x**)

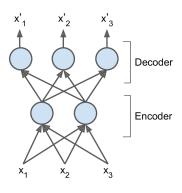
 Converts the inputs to an internal representation.
- A decoder (generative network), r = g(h)
 Converts the internal representation to the outputs.





Autoencoders (3/5)

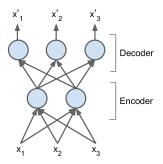
- ► An autoencoder is always composed of two parts.
- An encoder (recognition network), h = f(x)
 Converts the inputs to an internal representation.
- A decoder (generative network), r = g(h)
 Converts the internal representation to the outputs.
- ▶ If an autoencoder learns to set g(f(x)) = x everywhere, it is not especially useful, why?





Autoencoders (4/5)

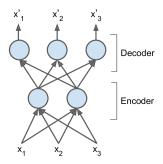
▶ Autoencoders are designed to be unable to learn to copy perfectly.





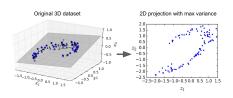
Autoencoders (4/5)

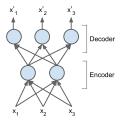
- Autoencoders are designed to be unable to learn to copy perfectly.
- ► The models are forced to prioritize which aspects of the input should be copied, they often learn useful properties of the data.





► Autoencoders are neural networks capable of learning efficient representations of the input data (called codings) without any supervision.

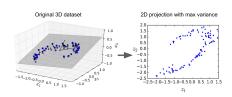


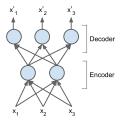




Autoencoders (5/5)

- ► Autoencoders are neural networks capable of learning efficient representations of the input data (called codings) without any supervision.
- ▶ Dimension reduction: these codings typically have a much lower dimensionality than the input data.







Different Types of Autoencoders

- ► Stacked autoencoders
- ► Denoising autoencoders
- ► Sparse autoencoders
- ► Variational autoencoders



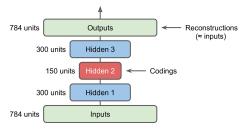
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Stacked Autoencoders (1/3)

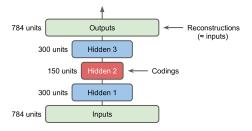
► Stacked autoencoder: autoencoders with multiple hidden layers.





Stacked Autoencoders (1/3)

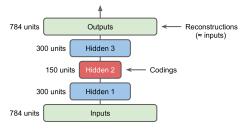
- ► Stacked autoencoder: autoencoders with multiple hidden layers.
- ▶ Adding more layers helps the autoencoder learn more complex codings.





Stacked Autoencoders (1/3)

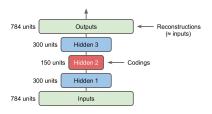
- ▶ Stacked autoencoder: autoencoders with multiple hidden layers.
- ▶ Adding more layers helps the autoencoder learn more complex codings.
- ► The architecture is typically symmetrical with regards to the central hidden layer.





Stacked Autoencoders (2/3)

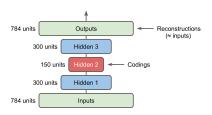
▶ In a symmetric architecture, we can tie the weights of the decoder layers to the weights of the encoder layers.





Stacked Autoencoders (2/3)

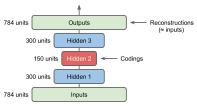
- ▶ In a symmetric architecture, we can tie the weights of the decoder layers to the weights of the encoder layers.
- ▶ In a network with N layers, the decoder layer weights can be defined as $w_{N-1+1} = w_1^T$, with $1 = 1, 2, \dots, \frac{N}{2}$.





Stacked Autoencoders (2/3)

- ▶ In a symmetric architecture, we can tie the weights of the decoder layers to the weights of the encoder layers.
- ▶ In a network with N layers, the decoder layer weights can be defined as $w_{N-1+1} = w_1^T$, with $1 = 1, 2, \dots, \frac{N}{2}$.
- ► This halves the number of weights in the model, speeding up training and limiting the risk of overfitting.



Stacked Autoencoders (3/3)

```
stacked_encoder = keras.models.Sequential([
    keras.layers.Flatten(input_shape=[28, 28]),
    keras.layers.Dense(100, activation="relu"),
    keras.layers.Dense(30, activation="relu"),
])
stacked_decoder = keras.models.Sequential([
    keras.layers.Dense(100, activation="relu", input_shape=[30]),
    keras.layers.Dense(28 * 28, activation="sigmoid"),
    keras.layers.Reshape([28, 28])
])
model = keras.models.Sequential([stacked_encoder, stacked_decoder])
```



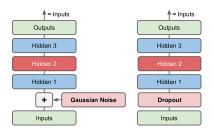
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Denoising Autoencoders (1/4)

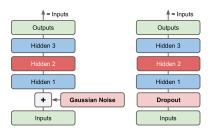
▶ One way to force the autoencoder to learn useful features is to add noise to its inputs, training it to recover the original noise-free inputs.





Denoising Autoencoders (1/4)

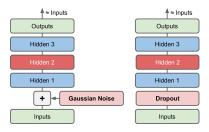
- ► One way to force the autoencoder to learn useful features is to add noise to its inputs, training it to recover the original noise-free inputs.
- ► This prevents the autoencoder from trivially copying its inputs to its outputs, so it ends up having to find patterns in the data.





Denoising Autoencoders (2/4)

► The noise can be pure Gaussian noise added to the inputs, or it can be randomly switched off inputs, just like in dropout.



Denoising Autoencoders (3/4)

```
denoising_encoder = keras.models.Sequential([
    keras.layers.Flatten(input_shape=[28, 28]),
    keras.layers.Dropout(0.5),
    keras.layers.Dense(100, activation="relu"),
    keras.layers.Dense(30, activation="relu")
])
denoising_decoder = keras.models.Sequential([
    keras.layers.Dense(100, activation="relu", input_shape=[30]),
    keras.layers.Dense(28 * 28, activation="sigmoid"),
    keras.layers.Reshape([28, 28])
])
model = keras.models.Sequential([denoising_encoder, denoising_decoder])
```

Denoising Autoencoders (4/4)

```
denoising_encoder = keras.models.Sequential([
    keras.layers.Flatten(input_shape=[28, 28]),
    keras.layers.GaussianNoise(0.2),
    keras.layers.Dense(100, activation="relu"),
    keras.layers.Dense(30, activation="relu")
])
denoising_decoder = keras.models.Sequential([
    keras.layers.Dense(100, activation="relu", input_shape=[30]),
    keras.layers.Dense(28 * 28, activation="sigmoid"),
    keras.layers.Reshape([28, 28])
])
model = keras.models.Sequential([denoising_encoder, denoising_decoder])
```



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Sparse Autoencoders (1/2)

Adding an appropriate term to the cost function to push the autoencoder to reducing the number of active neurons in the coding layer.



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- ► This forces the autoencoder to represent each input as a combination of a small number of activations.



Sparse Autoencoders (1/2)

- Adding an appropriate term to the cost function to push the autoencoder to reducing the number of active neurons in the coding layer.
- ► This forces the autoencoder to represent each input as a combination of a small number of activations.
- ► As a result, each neuron in the coding layer typically ends up representing a useful feature.

Sparse Autoencoders (2/2)

```
sparse_l1_encoder = keras.models.Sequential([
    keras.layers.Flatten(input_shape=[28, 28]),
    keras.layers.Dense(100, activation="selu"),
    keras.layers.Dense(300, activation="sigmoid", activity_regularizer=keras.regularizers.l1(1e-3))
])

sparse_l1_decoder = keras.models.Sequential([
    keras.layers.Dense(100, activation="selu", input_shape=[300]),
    keras.layers.Dense(28 * 28, activation="sigmoid"),
    keras.layers.Reshape([28, 28])
])

model = keras.models.Sequential([sparse_l1_encoder, sparse_l1_decoder])
```



Different Types of Autoencoders

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▶ Variational autoencoders are probabilistic autoencoders.



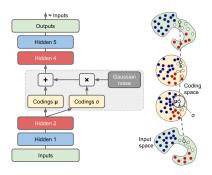
- ▶ Variational autoencoders are probabilistic autoencoders.
- ► Their outputs are partly determined by chance, even after training.
 - As opposed to denoising autoencoders, which use randomness only during training.



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- ► Their outputs are partly determined by chance, even after training.
 - · As opposed to denoising autoencoders, which use randomness only during training.
- ► They are generative autoencoders, meaning that they can generate new instances that look like they were sampled from the training set.



▶ Instead of directly producing a coding for a given input, the encoder produces a mean coding μ and a standard deviation σ .

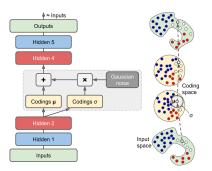




▶ Instead of directly producing a coding for a given input, the encoder produces a mean coding μ and a standard deviation σ .

► The actual coding is then sampled randomly from a Gaussian distribution with mean

 μ and standard deviation σ .



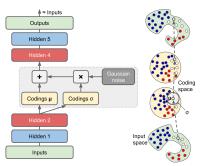


▶ Instead of directly producing a coding for a given input, the encoder produces a mean coding μ and a standard deviation σ .

► The actual coding is then sampled randomly from a Gaussian distribution with mean

 μ and standard deviation σ .

After that the decoder just decodes the sampled coding normally.



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 - Using the KL divergence between the target distribution (the Gaussian distribution) and the actual distribution of the codings.

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 - Pushes the autoencoder to reproduce its inputs.
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 - Pushes the autoencoder to have codings that look as though they were sampled from a simple Gaussian distribution.
 - Using the KL divergence between the target distribution (the Gaussian distribution) and the actual distribution of the codings.
 - latent_loss = $-\frac{1}{2}\sum_{i=1}^{K}(1+\log(\sigma_{i}^{2})-\sigma_{i}^{2}-\mu_{i}^{2})$



► Encoder part

```
inputs = keras.layers.Input(shape=[28, 28])
z = keras.layers.Flatten()(inputs)
z = keras.layers.Dense(150, activation="relu")(z)
z = keras.layers.Dense(100, activation="relu")(z)
codings_mean = keras.layers.Dense(10)(z)
codings_log_var = keras.layers.Dense(10)(z)
codings = Sampling()([codings_mean, codings_log_var]) # normal distribution
variational_encoder = keras.models.Model(inputs=[inputs], outputs=[codings])
```



Decoder part

```
decoder_inputs = keras.layers.Input(shape=[codings_size])
x = keras.layers.Dense(100, activation="relu")(decoder_inputs)
x = keras.layers.Dense(150, activation="relu")(x)
x = keras.layers.Dense(28 * 28, activation="sigmoid")(x)
outputs = keras.layers.Reshape([28, 28])(x)
variational_decoder = keras.models.Model(inputs=[decoder_inputs], outputs=[outputs])
```

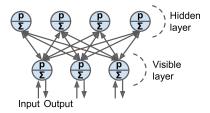






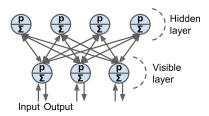


► A Restricted Boltzmann Machine (RBM) is a stochastic neural network.



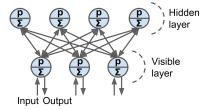


- ► A Restricted Boltzmann Machine (RBM) is a stochastic neural network.
- ► Stochastic meaning these activations have a probabilistic element, instead of deterministic functions, e.g., logistic or ReLU.





- ► A Restricted Boltzmann Machine (RBM) is a stochastic neural network.
- ► Stochastic meaning these activations have a probabilistic element, instead of deterministic functions, e.g., logistic or ReLU.
- ► The neurons form a bipartite graph:
 - One visible layer and one hidden layer.
 - A symmetric connection between the two layers.
 - There are no connections between neurons within a layer.





Let's Start With An Example



▶ We have a set of six movies, and we ask users to tell us which ones they want to watch.















- ▶ We have a set of six movies, and we ask users to tell us which ones they want to watch.
- ► We want to learn two latent neurons (hidden neurons) underlying movie preferences, e.g., SF/fantasy and Oscar winners







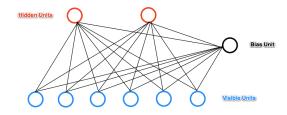








► Our RBM would look like the following.

















► Alice: (HP=1, Avatar=1, LOTR=1, Glad=0, Titan=0, Sep=0), Big SF fan.















- ► Alice: (HP=1, Avatar=1, LOTR=1, Glad=0, Titan=0, Sep=0), Big SF fan.
- ▶ Bob: (HP=1, Avatar=0, LOTR=1, Glad=0, Titan=0, Sep=0), SF fan, but not Avatar.















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- ► Eric: (HP=0, Avat=0, LOTR=1, Glad=1, Titan=0, Sep=1), Oscar winners fan, but not Titanic.















- ► Alice: (HP=1, Avatar=1, LOTR=1, Glad=0, Titan=0, Sep=0), Big SF fan.
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- ► Eric: (HP=0, Avat=0, LOTR=1, Glad=1, Titan=0, Sep=1), Oscar winners fan, but not Titanic.
- ► Fred: (HP=0, Avat=0, LOTR=1, Glad=1, Titan=1, Sep=1), Big Oscar winners fan.







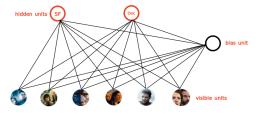








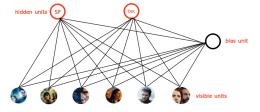
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- Assume the given input x_i is the 0 or 1 for each visible neuron v_i .
 - 1: like a movie, and 0: dislike a movie
- ► Compute the activation energy at hidden neuron h_i:







▶ For each hidden neuron h_j , we compute the probability $p(h_j)$.

$$\begin{aligned} \mathbf{a}(\mathbf{h}_j) &= \sum_i \mathbf{w}_{i\,j} \mathbf{v}_i \\ p(\mathbf{h}_j) &= \text{sigmoid}(\mathbf{a}(\mathbf{h}_j)) = \frac{1}{1 + e^{-\mathbf{a}(\mathbf{h}_j)}} \end{aligned}$$





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We turn on the hidden neuron h_j with the probability $p(h_j)$, and turn it off with probability $1 - p(h_j)$.



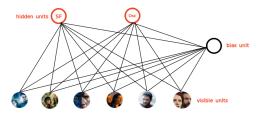


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- ▶ But it will turn on with a high probability.
 - In reality, if you want to watch all three of those movies makes us highly suspect you like SF/fantasy in general.
 - But there's a small chance you like them for other reasons.



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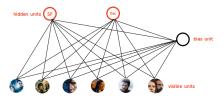
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- ▶ Being on the SF/fantasy neuron doesn't guarantee that we'll always recommend all three of Harry Potter, Avatar, and LOTR.
 - For example not everyone who likes science fiction liked Avatar.



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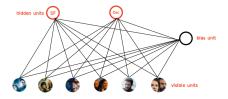


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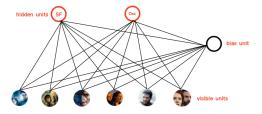


- ▶ How do we learn the connection weights w_{ij} in our network?
- ► Assume, as an input we have a bunch of binary vectors **x** with six elements corresponding to a user's movie preferences.
- ▶ We do the following steps in each epoch:
- ▶ 1. Take a training instance **x** and set the states of the visible neurons to these preferences.



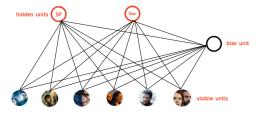


▶ 2. Update the states of the hidden neurons.





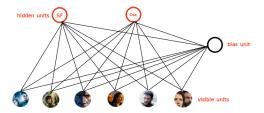
- ▶ 2. Update the states of the hidden neurons.
 - Compute $a(h_{\rm j}) = \sum_{\rm i} w_{\rm i\, j} v_{\rm i}$ for each hidden neuron $h_{\rm j}.$



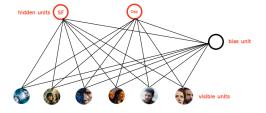


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- ▶ 3. For each edge e_{ij} , compute positive $(e_{ij}) = v_i \times h_j$
 - I.e., for each pair of neurons, measure whether they are both on.





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 - We denote the updated visible neurons with v'_i .
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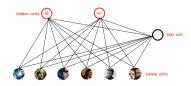


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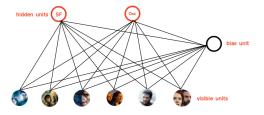
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- ▶ 6. For each edge e_{ij} , compute $negative(e_{ij}) = v_i' \times h_j'$





▶ 7. Update the weight of each edge e_{ij}.

$$\mathtt{w_{ij}} = \mathtt{w_{ij}} + \eta(\mathtt{positive}(\mathtt{e_{ij}}) - \mathtt{negative}(\mathtt{e_{ij}}))$$



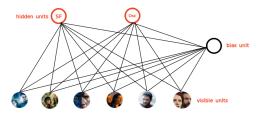


RBM Example (11/11)

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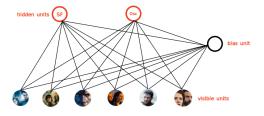


RBM Example (11/11)

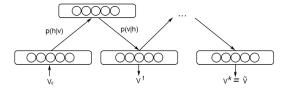
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- ▶ 8. Repeat over all training examples.
- ▶ 9. Continue until the error between the training examples and their reconstructions falls below some threshold or we reach some maximum number of epochs.

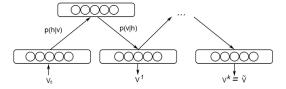


▶ Step 1, Gibbs sampling: what we have done in steps 1-6.



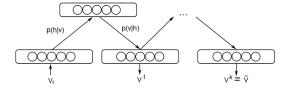


- ▶ Step 1, Gibbs sampling: what we have done in steps 1-6.
- ▶ Given an input vector \mathbf{v} , compute $p(\mathbf{h}|\mathbf{v})$.



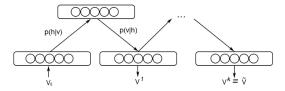


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- ► This process is repeated k times.



- ▶ Step 2, contrastive divergence: what we have done in step 7.
 - Just a fancy name for approximate gradient descent.

$$\mathbf{w} = \mathbf{w} + \eta(\text{positive}(\mathbf{e}) - \text{negative}(\mathbf{e}))$$



More Details about RBM

► Energy a quantitative property of physics.

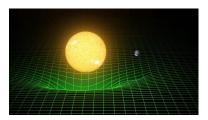


- ► Energy a quantitative property of physics.
 - E.g., gravitational energy describes the potential energy a body with mass has in relation to another massive object due to gravity.





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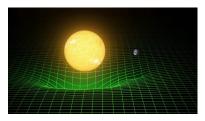


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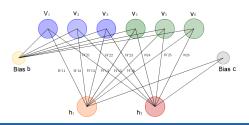
- ▶ One purpose of deep learning models is to encode dependencies between variables.
- ► The capturing of dependencies happen through associating of a scalar energy to each state of the variables.
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- ► A high energy means a bad compatibility.
- ► An energy based model tries always to minimize a predefined energy function.





► The energy function for the RBMs is defined as:

$$\mathtt{E}(\textbf{v},\textbf{h}) = -(\sum_{\mathtt{i}\mathtt{j}} \mathtt{w}_{\mathtt{i}\mathtt{j}} \mathtt{v}_{\mathtt{i}} \mathtt{h}_{\mathtt{j}} + \sum_{\mathtt{i}} \mathtt{b}_{\mathtt{i}} \mathtt{v}_{\mathtt{i}} + \sum_{\mathtt{j}} \mathtt{c}_{\mathtt{j}} \mathtt{h}_{\mathtt{j}})$$

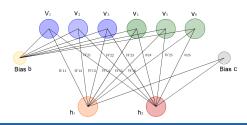




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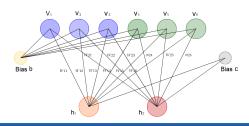




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- w represents the weights connecting visible and hidden units.

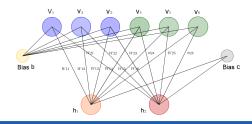




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- ▶ v and h represent the visible and hidden units, respectively.
- w represents the weights connecting visible and hidden units.
- **b** and **c** are the biases of the visible and hidden layers, respectively.





RBM is a Probabilistic Model (1/2)

► The probability of a certain state of **v** and **h**:

$$p(\boldsymbol{v},\boldsymbol{h}) = \frac{e^{-E(\boldsymbol{v},\boldsymbol{h})}}{\sum_{\boldsymbol{v},\boldsymbol{h}} e^{-E(\boldsymbol{v},\boldsymbol{h})}}$$



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- ▶ In physics, the joint distribution $p(\mathbf{v}, \mathbf{h})$ is known as the Boltzmann Distribution or Gibbs Distribution.
- ▶ At each point in time the RBM is in a certain state.
 - The state refers to the values of neurons in the visible and hidden layers **v** and **h**.



RBM is a Probabilistic Model (2/2)

▶ It is difficult to calculate the joint probability due to the huge number of possible combination of v and h.

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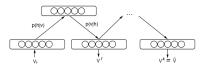
▶ It is difficult to calculate the joint probability due to the huge number of possible combination of **v** and **h**.

$$p(\mathbf{v}, \mathbf{h}) = \frac{e^{-E(\mathbf{v}, \mathbf{h})}}{\sum_{\mathbf{v}, \mathbf{h}} e^{-E(\mathbf{v}, \mathbf{h})}}$$

► Much easier is the calculation of the conditional probabilities of state **h** given the state **v** and vice versa (Gibbs sampling)

$$p(\mathbf{h}|\mathbf{v}) = \Pi_{i}p(h_{i}|\mathbf{v})$$

 $p(\mathbf{v}|\mathbf{h}) = \Pi_{i}p(v_{i}|\mathbf{h})$





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- Use the maximum-likelihood estimation.
- ► For a model of the form p(v) with parameters w, the log-likelihood given a single training example v is:

$$\log p(\boldsymbol{v}|\boldsymbol{h}) = \log \frac{\sum_{\boldsymbol{h}} e^{-E(\boldsymbol{v},\boldsymbol{h})}}{\sum_{\boldsymbol{v},\boldsymbol{h}} e^{-E(\boldsymbol{v},\boldsymbol{h})}} = \log \sum_{\boldsymbol{h}} e^{-E(\boldsymbol{v},\boldsymbol{h})} - \log \sum_{\boldsymbol{v},\boldsymbol{h}} e^{-E(\boldsymbol{v},\boldsymbol{h})}$$



► The log-likelihood gradients for an RBM with binary units:

$$\frac{\partial \log p(\textbf{v}|\textbf{h})}{\partial \textbf{w}_{\texttt{ij}}} = positive(\textbf{e}_{\texttt{ij}}) - negative(\textbf{e}_{\texttt{ij}})$$



► The log-likelihood gradients for an RBM with binary units:

$$\frac{\partial \log p(\textbf{v}|\textbf{h})}{\partial w_{ij}} = positive(e_{ij}) - negative(e_{ij})$$

► Then, we can update the weight w as follows:

$$\mathbf{w}_{\mathtt{ij}}^{(\mathtt{next})} = \mathbf{w}_{\mathtt{ij}} + \eta(\mathtt{positive}(\mathbf{e}_{\mathtt{ij}}) - \mathtt{negative}(\mathbf{e}_{\mathtt{ij}}))$$







Summary



- Autoencoders
 - Stacked autoencoders
 - Denoising autoencoders
 - Variational autoencoders
- ► Restricted Boltzmann Machine
 - Gibbs sampling
 - Contrastive divergence

Reference

- ▶ Ian Goodfellow et al., Deep Learning (Ch. 14, 20)
- ► Aurélien Géron, Hands-On Machine Learning (Ch. 17)



Questions?