Feedback — Lecture 14 Quiz

Help Center

You submitted this quiz on **Fri 1 Jul 2016 2:03 AM CEST**. You got a score of **6.00** out of **8.00**. However, you will not get credit for it, since it was submitted past the deadline.

Question 1

Why is a Deep Belief Network not a Boltzmann Machine?

| Your Answer | Score | Explanation |
|--|----------------|---|
| A DBN is not a probabilistic model of the data. | | |
| Some edges in a DBN are directed. | ✓ 1.00 | In a Boltzmann Machine, all edges must be undirected. A DBN has directed edges from the top-level RBM to each subsequent layer below. |
| A DBN does not have hidden units. | | |
| All edges in aDBN are directed. | | |
| Total | 1.00 / 1.00 | |

Question 2

Brian looked at the direction of arrows in a DBN and was surprised to find that the data is at the "output". "Where is the input ?!", he exclaimed, "How will I give input to this model and get all those cool features?" In this context, which of the following statements are true? Check all that apply.

| Your Answer | | Score | Explanation |
|---|----------|-------|--|
| ■ A DBN is a generative model of the data and cannot be used to generate features for any | ~ | 0.50 | This is not true. A DBN can be used to generate features for any given data vector by doing inference. |

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| given input. It can only be used to get features for data that was generated by the model. | | |
| A DBN is a generative model of the data, which means that, its arrows define a way of generating data from a probability distribution, so there is no "input". | x 0.00 | Unlike a feed-forward neural net, a DBN is not a mapping from inputs to outputs. It is a probabilistic generative model of the data. |
| In order to get features h given some data v , he must perform inference to find out $P(h v)$. There is an easy approximate way of doing this, just traverse the arrows in the opposite direction. | x 0.00 | Traversing arrows in the opposite direction is an approximate inference procedure. |
| In order to get features h given some data v , he must perform inference to find out $P(h v)$. There is an easy exact way of doing this, just traverse the arrows in the opposite direction. | ✔ 0.50 | Traversing arrows in the opposite direction is an approximate inference procedure. |
| Total | 1.00 / 2.00 | |

Question 3

In which of the following cases is pretraining likely to help the most (compared to training a neural net from random initialization)?

| Your Answer | Score | Explanation |
|--|-------|-------------|
| A dataset of images is to be classified into 100 semantic classes. Fortunately, there are 100 million labelled training examples. | | |
| A dataset of binary pixel images which are to be classified based on | | |

| parity, i.e., if the sum of pixels is even the image has label 0, otherwise it has label 1. | | |
|--|----------------|--|
| A speech dataset with 10 billion labelled training examples. | | |
| A dataset of images is to be classified into 100 semantic classes. There are only 1,000 labelled images but 100 million unlabelled ones are available from the internet. | ✔ 1.00 | The small labelled set of images is unlikely to have enough information to learn good features. The unlabelled images can be used to do pretraining for learning features. |
| Total | 1.00 / 1.00 | |

Question 4

Why does pretraining help more when the network is deep?

| Your Answer | | Score | Explanation |
|--|---|-------|--|
| As nets get deeper, contrastive divergence objective used during pretraining gets closer to the classification objective. | • | 0.50 | Contrastive divergence has nothing to do with the classification objective. |
| During backpropagation in very deep nets, the lower level layers get very small gradients, making it hard to learn good low-level features. Since pretraining starts those low-level | ~ | 0.50 | Lower level layers can get very small gradients, especially saturating hidden units are used (such as logistic or tanh units). Pretraining can initialize the weights in a proper region of weight space so that the features don't have to start learning from scratch. |

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| features off at a good point, there is a big win. | | | |
| Deeper nets have more parameters than shallow ones and they overfit easily. Therefore, initializing them sensibly is important. | × | 0.00 | More parameters means that the model can find ingenious ways of overfitting by learning features that don't generalize well. Pretraining can initialize the weights in a proper region of weight space so that the features learned are not too bad. |
| Backpropagation algorithm cannot give accurate gradients for very deep networks. So it is important to have good initializtions, especially, for the lower layers. | • | 0.50 | Backpropgation gives exact gradients. |
| Total | | 1.50 / 2.00 | |

Question 5

The energy function for binary RBMs goes by

$$E(\mathbf{v},\mathbf{h}) = -\sum_i v_i b_i - \sum_j h_j a_j - \sum_{i,j} v_i W_{ij} h_j$$

When modeling real-valued data (i.e., when ${f v}$ is a real-valued vector not a binary one) we change it to

$$E(\mathbf{v},\mathbf{h}) = \sum_i rac{\left(v_i - b_i
ight)^2}{2\sigma_i^2} - \sum_j h_j a_j - \sum_{i,j} rac{v_i}{\sigma_i} W_{ij} h_j$$

Why can't we still use the same old one?

| our Answer | Score | e Explanation |
|---|--------|---|
| If we use the old ne, the real-valued ectors would end up eing constrained to binary. | ✔ 0.50 | The energy function does not impose any such constraints. |

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|--|---------------|--|
| Probability distributions over real-valued data can only be modeled by having a conditional Gaussian distribution over them. So we have to use a quadratic term. | ✔ 0.50 | This is not true. Distributions over real-valued data can be modelled by a multitude of distributions or combinations thereof. The choice of Gaussian distribution is arbitrary. |
| If the model assigns an energy e_1 to state $\mathbf{v_1}$, \mathbf{h} , and e_2 to state $\mathbf{v_2}$, \mathbf{h} , then it would assign energy $(e_1+e_2)/2$ to state $(\mathbf{v_1}+\mathbf{v_2})/2$, \mathbf{h} . This does not make sense for the kind of distributions we usually want to model. | ✓ 0.50 | Suppose v_1 and v_2 represent two images. We would like e_1 and e_2 to be small. This makes the energy of the average image low, but the average of two images would not look like a natural image and should not have low energy. |
| If we continue to use the same one, then in general, there will be infinitely many \mathbf{v} 's and \mathbf{h} 's such that, $E(\mathbf{v},\mathbf{h})$ will be infinitely small (close to $-\infty$). The probability distribution resulting from such an energy function is not useful for modeling real data. | × 0.00 | If some $b_i<0$, then if $v_i\to-\infty$, then $E\to-\infty$. Similarly for $b_i>0$ if $v_i\to\infty$, then $E\to-\infty$. So the Boltzmann distribution based on this energy function would behave in weird ways. |
| Total | 1.50 / 2.00 | |