MACHINE LEARNING IN BIOINFORMATICS ANN Architectures

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OUTLINE

- Part of this lecture:
 - Embeddings
 - Auto-encoders
 - Convolutions on images and graphs
 - Attention mechanism
- Other important architectures not covered here:
 - Generative adversarial networks (GANs)
 - Deep tensor factorization
 - Recurrent neural networks (LSTM/GRU)

Embeddings

ONE-HOT ENCODING

- Assume we want to work with categorical data, e.g.
 - DNA or protein sequences
 - Text (vectors of words)
- Traditionally, we would use one-hot encoding, which use a dimension for each category
- For example, a DNA sequence ACGTTA could be represented as

[1	0	0	0
0	1	0	0
0	0	1	0
0	0	0	1
0	0	0	1
1	0	0	0

ONE-HOT ENCODING

- One-hot encodings have several problems
- For data with many categories, we obtain very high-dimensional feature vectors, e.g.
 - Protein sequences would already require 20 dimensions
 - Text would require one dimension per word type
- One-hot encodings should be used for purely categorical data, where we have no similarity between categories
- However, for most data we have certain similarities, e.g.
 - Amino acid replacements have different effects, which suggests that some amino acids are more similar in function than others

EMBEDDINGS



- We assign each category k a feature vector $x_k \in \mathbb{R}^p$
- The representations x_k are randomly initialized and optimized during training
- After training we often observe that similar categories cluster together

CONVOLUTIONAL NEURAL NETWORKS FOR IMAGES

IMAGE PATTERN DETECTION

Conway's Game of Life - glider gun:





Glider pattern:



























IMAGE PATTERN DETECTION



IMAGE PATTERN DETECTION - CONVOLUTION



• Let $\tilde{x}_j \in \mathbb{R}^r$ denote the *j*-th image patch of image X, e.g.

 $\tilde{X}_2 = (0, 0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 1, 0, 1, 0, \dots)^\top$

Let $w_k \in \mathbb{R}^r$ denote the *k*-th glider pattern or kernel, e.g.

 $W_1 = (0, 0, 0, 0, 0, 0, 0, 0, 1, 0, 0, 0, 0, 0, 1, 0, \dots)^\top$

The output y_i at position j is given by

$$y_j = \tilde{x}_j^\top W_k$$

Image pattern detection - Convolution

■ Let $\tilde{X} \in \mathbb{R}^{q \times r}$ denote the matrix of *q* image patches from image *X* and *W* ∈ $\mathbb{R}^{r \times p}$ the matrix of kernels, i.e.

$$\tilde{X} = \begin{bmatrix} \tilde{X}_1^\top \\ \tilde{X}_2^\top \\ \vdots \\ \tilde{X}_q^\top \end{bmatrix}, \quad W = [w_1, w_2, \dots, w_p]$$

The result $Y \in \mathbb{R}^{q \times p}$ of applying the kernel matrix W to image X is given by

$$Y = \tilde{X}W = X * W$$

where " * " is called convolution¹

¹Technically, we are computing a cross-correlation and not a convolution

EQUIVARIANCE

- Let X be an image and W a filter
- $\varphi(X) = X * W$ denotes a convolution with W
- $\tau(X)$ is a translation of an image
- \blacksquare The following diagram shows that φ is equivariant with respect to τ



Exception are the borders of images

WHY EQUIVARIANCE AND NOT INVARIANCE?



IMAGE PATTERN DETECTION



- Stack multiple convolutions
- Case 1: All images have the same dimension
- \Rightarrow Feed into neural network
 - Case 2: Images have variable dimension
- ⇒ Compute summary statistics (global pooling)
 - ▶ mean
 - ► max

POOLING LAYERS

- Applying kernels leads to *translation-equivariant* features
- Pooling layers add (limited amount of) translation invariance
- Average pooling
- Max pooling

1	3	3	1	1	0	0	1	1
2	1	1	0	0	2	1	1	2
0	1	1	0	3	3	0	1	6
1	5	1	1		2	1	0	1
T	5	Т	T	U	5	1	U	-
$\frac{1}{1}$	2	$\frac{1}{1}$	$\frac{1}{1}$	2	4	0	1	3

3	3	6
5	4	3

GRAPH CONVOLUTIONAL NEURAL NETWORKS (GCNNS)



- Convolutions are not only restricted to image and time-series data
- Graph convolutions are used to model the interaction between nodes
- Let G = (N, E) denote a graph with nodes N and edges E
- How could we implement a convolution of G with a weight matrix W?
- The result of a convolution is again a graph², i.e.

$$G' = G * W$$

²Remember that convolution on images also returns an image

Graph G with 5 nodes and 5 edges:



- We assign a feature vector $x_i \in \mathbb{R}^p$ to the *i*-th node
- The feature vector can depend on the type of the node
- Nodes of the same type might share the same feature vector

■ Graph *G* with 5 nodes and 5 edges:



- We assign a feature vector $x_i \in \mathbb{R}^p$ to the *i*-th node
- The feature vector can depend on the type of the node
- Nodes of the same type might share the same feature vector

- Let $A = (a_{ij})_{ij} \in \mathbb{R}^{k \times k}$ denote the adjacency matrix of a graph with *k* nodes
- The strength of the connection between node *i* and *j* is given by a_{ij}
- Self-connections $a_{ii} \neq 0$ allow to incorporate the features of the nodes itself
- The convolution operation updates the feature vector of node *i* by summing over the contributions of all neighbor nodes, i.e.

$$\mathbf{x}_{i}' = \sigma \left(\sum_{j \neq i} \mathbf{a}_{ij} \mathbf{W} \mathbf{x}_{j} \right)$$

where $W \in \mathbb{R}^{p \times p}$ and σ is the activation function³

³Graph convolutions are *permutation equivariant*

■ For the full graph we obtain



where $X \in \mathbb{R}^{k \times p}$ is the matrix of *k* feature vectors

- Note that the weight matrix W does not depend on the size and connectivity of the graph
- W can be applied to multiple graphs and optimized during training of the graph convolutional neural network (GCNN)
- GCNNs typically apply multiple convolutions and afterwards compute summary statistics of the feature vectors, the result can then be used in a conventional neural network

³Many extensions and generalizations exist [Battaglia et al., 2018, Dwivedi et al., 2020]

AUTO-ENCODERS

AUTO-ENCODERS

- Embeddings implicitly group categories by their similarity
- Auto-encoders [Kramer, 1991] learn hidden representations for non-categorical data:



- During training, the error between X and X' is minimized
- The embedding or latent space should have lower dimension than the input space

AUTO-ENCODERS - FORMAL DEFINITION

- The encoder $f_W : \mathbb{R}^p \to \mathbb{R}^q$ is a neural network with weights W that maps a sample $x \in \mathbb{R}^p$ into a q-dimensional feature space
- The decoder $g_V : \mathbb{R}^q \to \mathbb{R}^p$ takes a point in feature space and maps it back to input space
- Given a set of training points {x_i}_i we train the auto-encoder by minimizing the error between the input and output of the network, i.e.

$$W, V = \argmin_{W,V} \|x_i - (g_v \circ f_W)(x_i)\|_2^2$$

AUTO-ENCODERS - PURPOSE

- Dimensionality reduction and visualization (similar to PCA and t-SNE)
- Compression to most important features (encoder output)
- Denoising and image restauration (decoder output), by adding noise to images before sending it to the encoder



Clustering and outlier detection on the latent space

VARIATIONAL AUTO-ENCODERS (VAES)

- Can we use auto-encoders for generating data? I.e. we could sample a point from the latent space and decode the corresponding data point
- Practice has shown that this appproach does not work
- The latent space has many holes where the decoder generates garbage
- Variational auto-encoders (VAEs) [Kingma and Welling, 2013] are a probabilistic formuation of auto-encoders, that regularize the latent space

VARIATIONAL AUTO-ENCODERS (VAES)



- Instead of learning latent representations directly, VAEs learn the parameters of given distributions
- The encoder learns the parameters λ of the distribution $q_{\lambda}(z \mid x)$
- The decoder learns the parameters θ of the distribution $p_{\theta}(x \mid z)$
- Training is more complicated, i.e. minimize the KL-divergence

ATTENTION

SEQUENTIAL DATA



Translations require special architectures that can deal with:

- Variable sentence lengths, i.e. variable n
- Long-range dependencies

RECURRENT NEURAL NETWORKS



- Recurrent neural networks (RNNs) are sequentially applied to each input x_i
- The architecture and weights are the same for all steps i.e. for RNN(0), RNN(1), ..., RNN(n)
- At each step *i*, RNNs take the input x_i and the state of the previous step *i* − 1 as input

ATTENTION IS ALL YOU NEED

- Recurrent neural networks (RNNs) were traditionally used for sequence data and to model long-range interactions
- Traditional RNNs have extreme vanishing / exploding gradient problem
- Long-short term memory (LSTM) [Hochreiter and Schmidhuber, 1997] solved this problem, but is still difficult to train
 - On a large input sequence it corresponds to a very deep neural network
 - Transfer learning never worked for LSTM
- Transformers with attention layer
 [Bahdanau et al., 2014, Vaswani et al., 2017] are an alternative to RNNs and show better performance

Self-attention layer



Self-attention layer

- Let $X = [x_1^{\top}, x_2^{\top}, ..., x_n^{\top}] \in \mathbb{R}^{n \times p}$ denote the data matrix, i.e. the embeddings of the input sequence
- The self-attention layer computes the *i*-th output $y_i \in \mathbb{R}^p$ as follows:

$$S_{i} = X_{i}^{\top} X^{\top}$$
$$W_{i} = \text{softmax}(S_{i}) = \left(\frac{e^{S_{ij}}}{\sum_{k=1}^{n} e^{S_{ik}}}\right)_{j=1,2,\dots,n}$$
$$y_{i} = W_{i} X$$

■ The self-attention layer computes the entire output $Y \in \mathbb{R}^{n \times p}$ as follows:

$$Y = \text{softmax}(XX^{\top})X$$

kernel

where the softmax is applied independently to each row

Self-attention maps

The self-attention map is defined as

 $\mathsf{A} = \mathsf{softmax}(XX^{\top})$

The matrix A can be visualized to inspect attention





ATTENTION LAYER

- Except for the embeddings (x_i)_i, the self-attention layer has no parameters that can be optimized
- For self-attention, the input sequence focuses attention on the input sequence itself and a linear combination of the input sequence x₁, x₂,..., x_n is returned
- The attention layer is a generalization of the self-attention layer, where
 - ▶ attention is focused on a set of *m* keys k_1, \ldots, k_m , with $k_j \in \mathbb{R}^p$
 - ▶ a linear combination of *m* values $v_1, ..., v_m$ is returned, where $v_j \in \mathbb{R}^p$
- The attention layer implements a differentiable data retrieval method for a database of *m* keys and values

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ATTENTION LAYER



ATTENTION LAYER

- Let $K \in \mathbb{R}^{m \times p}$ and $V \in \mathbb{R}^{m \times p}$ denote a set of *m* keys and values
- The attention layer computes the entire output $Y \in \mathbb{R}^{n \times p}$ as follows:

 $Y = \operatorname{softmax}(XK^{\top})V$

- Remarks:
 - There exist several variants of the attention layer
 - Transformers use a both attention and self-attention layers
 - The sequential order is lost for self-attention and attention layers
 - Transformers use another encoding for restoring relative word positions
 - Multiple attention heads are commonly used

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TRANSFER LEARNING

- Some of the most successful deep learning models:
 - Protein folding: AlphaFold [Jumper et al., 2021]
 - Vision: GoogLeNet [Szegedy et al., 2015], Squeeze-and-Excitation Networks (SENet) [Hu et al., 2018]
 - Translation: BERT [Devlin et al., 2018], Text-to-Text Transfer Transformer (T5) [Raffel et al., 2019]
- Training T5 (11B-parameter variant) costs well above \$1.3 million [Sharir et al., 2020]
- True deep neural networks are not affordable for most academics
- Transfer learning allows to adapt pre-trained models

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