
EECS 182 Deep Neural Networks
 Fall 2025 Anant Sahai and Gireeja Ranade Homework 7

This homework is due on Mar 30, at 10:59PM.

1. Implementing RNNs (Coding)

This problem involves filling out [this notebook](#).

- (a) **Implement Section 1A in the notebook**, which constructs a vanilla RNN layer. This layer implements the function

$$h_t = \sigma(W^h h_{t-1} + W^x x_t + b)$$

where W^h , W^x , and b are learned parameter matrices, x is the input sequence, and σ is a nonlinearity such as tanh. The RNN layer “unrolls” across a sequence, passing a hidden state between timesteps and returning an array of hidden states at all timesteps.

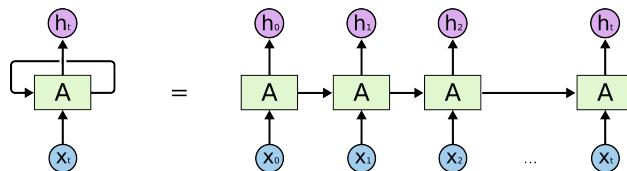


Figure 1: Source: <https://colah.github.io/posts/2015-08-Understanding-LSTMs/>

Copy the outputs of the “Test Cases” code cell and paste it into your submission of the written assignment.

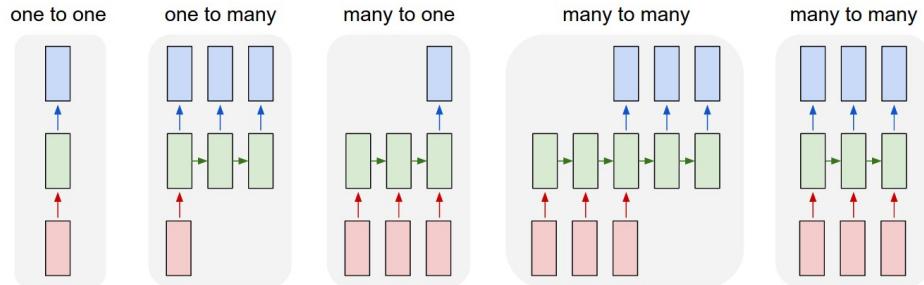
- (b) **Implement Section 1.B of the notebook**, in which you’ll use this RNN layer in a regression model by adding a final linear layer on top of the RNN outputs.

$$\hat{y}_t = W^f h_t + b^f$$

We’ll compute one prediction for each timestep.

Copy the outputs of the “Tests” code cell and paste it into your submission of the written assignment.

- (c) RNNs can be used for many kinds of prediction problems, as shown below. In this notebook we will look at many-to-one prediction and aligned many-to-many prediction.



We will use a simple averaging task. The input X consists of a sequence of numbers, and the label y is a running average of all numbers seen so far.

We will consider two tasks with this dataset:

- Task 1: predict the running average at all timesteps
- Task 2: predict the average at the last timestep only

Implement Section 1.C in the notebook, in which you'll look at the synthetic dataset shown and implement a loss function for the two problem variants.

Copy the outputs of the “Tests” code cell and paste it into your submission of the written assignment.

RNN: Computational Graph: Many to One

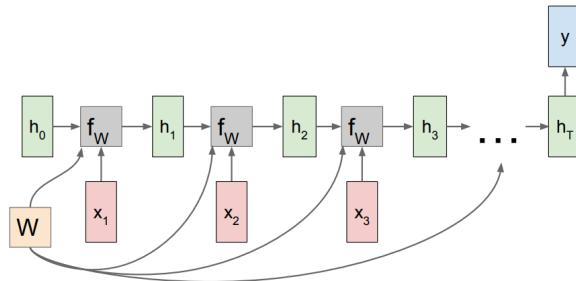


Figure 2: Image source: https://calvinfeng.gitbook.io/machine-learning-notebook/supervised-learning/recurrent-neural-network/recurrent_neural_networks

- (d) Consider an RNN which outputs a single prediction at timestep T . As shown in Figure 2, each weight matrix W influences the loss by multiple paths. As a result, the gradient is also summed over multiple paths:

$$\frac{\partial \mathcal{L}}{\partial W} = \frac{\partial \mathcal{L}}{\partial h_T} \frac{\partial h_T}{\partial W} + \frac{\partial \mathcal{L}}{\partial h_{T-1}} \frac{\partial h_{T-1}}{\partial W} + \dots + \frac{\partial \mathcal{L}}{\partial h_1} \frac{\partial h_1}{\partial W} \quad (1)$$

When you backpropagate a loss through many timesteps, the later terms in this sum often end up with either very small or very large magnitude - called vanishing or exploding gradients respectively. Either problem can make learning with long sequences difficult.

Implement Notebook Section 1.D, which plots the magnitude at each timestep of $\frac{\partial \mathcal{L}}{\partial h_t}$. Play around with this visualization tool and try to generate exploding and vanishing gradients.

Include a screenshot of your visualization in the written assignment submission.

- (e) **If the network has no nonlinearities, under what conditions would you expect the exploding or vanishing gradients with for long sequences? Why?** (Hint: it might be helpful to write out the

formula for $\frac{\partial \mathcal{L}}{\partial h_t}$ and analyze how this changes with different t). **Do you see this pattern empirically using the visualization tool in Section 1.D in the notebook** with `last_step_only=True`?

- (f) Compare the magnitude of hidden states and gradients when using ReLU and tanh nonlinearities in Section 1.D in the notebook. **Which activation results in more vanishing and exploding gradients? Why?** (This does not have to be a rigorous mathematical explanation.)
- (g) **What happens if you set `last_target_only = False` in Section 1.D in the notebook? Explain why this change affects vanishing gradients. Does it help the network's ability to learn dependencies across long sequences?** (The explanation can be intuitive, not mathematically rigorous.)
- (h) (Optional) We can create multi-layer recurrent networks by stacking layers as shown in Figure 3. The hidden state outputs from one layer become the inputs to the layer above.

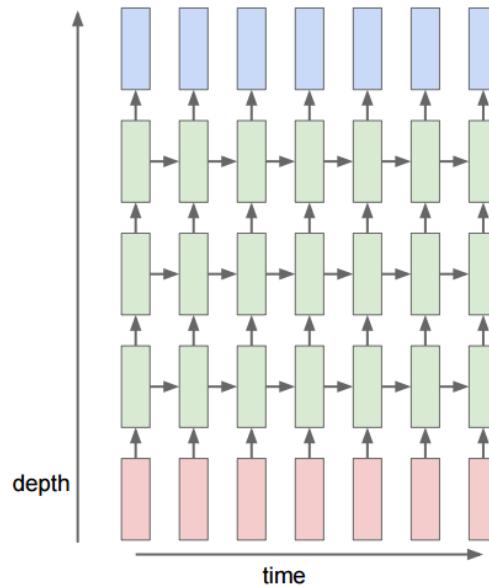


Figure 3: Image source: https://calvinfeng.gitbook.io/machine-learning-notebook/supervised-learning/recurrent-neural-network/recurrent_neural_networks

Implement notebook Section 1.K and run the last cell to train your network. You should be able to reach training loss < 0.001 for the 2-layer networks, and $<.01$ for the 1-layer networks.

2. RNNs for Last Name Classification (Coding)

Please follow the instructions in [this notebook](#). You will train a neural network to predict the probable language of origin for a given last name / family name in Latin alphabets.

- (a) Although the neural network you have trained is intended to predict the language of origin for a given last name, it could potentially be misused. **In what ways do you think this could be problematic in real-world applications?**

3. Auto-encoder : Learning without Labels

So far, with supervised learning algorithms we have used labelled datasets $\mathcal{D} = \{\mathbf{X}, \mathbf{y}\}$ to learn a mapping f_θ from input x to labels y . In this problem, we now consider algorithms for *unsupervised learning*, where we are given only inputs x , but no labels y . At a high-level, unsupervised learning algorithms leverage

some structure in the dataset to define proxy tasks, and learn *representations* that are useful for solving downstream tasks.

Autoencoders present a family of such algorithms, where we consider the problem of learning a function $f_\theta : \mathbb{R}^m \rightarrow \mathbb{R}^k$ from input x to a *intermediate representation* z of x (usually $k \ll m$). For autoencoders, we use reconstructing the original signal as a proxy task, i.e. representations are more likely to be useful for downstream tasks if they are low-dimensional but encode sufficient information to approximately reconstruct the input. Broadly, autoencoder architectures have two modules:

- An encoder $f_\theta : \mathbb{R}^m \rightarrow \mathbb{R}^k$ that maps input x to a representation z .
- A decoder $g_\phi : \mathbb{R}^k \rightarrow \mathbb{R}^m$ that maps representation z to a reconstruction \hat{x} of x .

In such architectures, the parameters (θ, ϕ) are learnable, and trained with the learning objective of minimizing the reconstruction error $\ell(\hat{x}_i, x_i) = \|x - \hat{x}\|_2^2$ using gradient descent.

$$\begin{aligned}\theta_{\text{enc}}, \phi_{\text{dec}} &= \underset{\Theta, \Phi}{\operatorname{argmin}} \frac{1}{N} \sum_{i=1}^N \ell(\hat{x}_i, x_i) \\ \hat{x} &= g_\phi(f_\theta(x))\end{aligned}$$

Note that above optimization problem does not require labels y . In practice however, we would want to use the *pretrained* models to solve the *downstream* task at hand, e.g. classifying MNIST digits. *Linear Probing* is one such approach, where we fix the encoder weights, and learn a simple linear classifier over the features $z = \text{encoder}(x)$.

(a) Designing AutoEncoders (Coding)

Please follow the instructions in [this notebook](#). You will train autoencoders, denoising autoencoders, and masked autoencoders on a synthetic dataset and the MNIST dataset. Once you finished with the notebook,

- Answer the following questions in your submission of the written assignment:
 - Show your visualization** of the vanilla autoencoder with different latent representation sizes.
 - Based on your previous visualizations, answer this question: **How does changing the latent representation size of the autoencoder affect the model's performance in terms of reconstruction accuracy and linear probe accuracy? Why?**

(b) PCA & AutoEncoders

In the case where the encoder f_θ, g_ϕ are linear functions, the model is termed as a *linear autoencoder*. In particular, assume that we have data $x_i \in \mathbb{R}^m$ and consider two weight matrices: an encoder $W_1 \in \mathbb{R}^{k \times m}$ and decoder $W_2 \in \mathbb{R}^{m \times k}$ (with $k < m$). Then, a linear autoencoder learns a low-dimensional embedding of the data $\mathbf{X} \in \mathbb{R}^{m \times n}$ (which we assume is zero-centered without loss of generality) by minimizing the objective,

$$\mathcal{L}(W_1, W_2; \mathbf{X}) = \|\mathbf{X} - W_2 W_1 \mathbf{X}\|_F^2 \tag{2}$$

We will assume $\sigma_1^2 > \dots > \sigma_k^2 > 0$ are the k largest eigenvalues of $\mathbf{X} \mathbf{X}^\top$. The assumption that the $\sigma_1, \dots, \sigma_k$ are positive and distinct ensures identifiability of the principal components, and is common in this setting. Therefore, the top- k eigenvalues of \mathbf{X} are $S = \text{diag}(\sigma_1, \dots, \sigma_k)$, with corresponding eigenvectors are the columns of $\mathbf{U}_k \in \mathbb{R}^{m \times k}$. A well-established result from ¹ shows that principal

¹Baldi, Pierre, and Kurt Hornik. "Neural networks and principal component analysis: Learning from examples without local minima" (1989)

components are the unique optimal solution to linear autoencoders (up to sign changes to the projection directions). In this subpart, we take some steps towards proving this result.

- (i) Write out the first order optimality conditions that the minima of Eq. 2 would satisfy.
- (ii) Show that the principal components U_k satisfy the optimality conditions outlined in (i).

4. Read a Blog Post: How to train your Resnet

In previous homeworks, we saw how memory and compute constraints on GPUs put limits on the architecture and the hyperparameters (e.g., batch size) we can use to train our models. To train better models, we could scale up by using multiple GPUs, but most distributed training techniques scale sub-linearly and often we simply don't have as many GPU resources at our disposal. This raises a natural question - how can we make model training more efficient on a single GPU?

The blog series [How to train your Resnet](#) explores how to train ResNet models efficiently on a single GPU. It covers a range of topics, including architecture, weight decay, batch normalization, and hyperparameter tuning. In doing so, it provides valuable insights into the training dynamics of neural networks and offers lessons that can be applied in other settings.

Read the blog series and answer the questions below.

- (a) What is the baseline training time and accuracy the authors started with? What was the final training time and accuracy achieved by the authors?
- (b) Comment on what you have learnt. (≈ 100 words)
- (c) Which approach taken by the authors interested you the most? Why? (≈ 100 words)

5. The power of the graph perspective in clustering (Coding)

Implement all the TODOs in the [notebook](#). Answer the written questions below.

- (a) We used the KMeans algorithm implementation of sklearn, and showed our attempt to cluster this dataset into 3 classes. **Comment on the output the KMeans algorithm. Did it work? If so, explain why, if not, explain why not.**

(b) adjacency_matrix = ?

- (c) As given, the data points in our dataset are represented simply with their 2D Cartesian coordinates. Let's now interpret every single point as a node in a graph. Our goal is to find a way to relate every node in the graph in such way that the points that are closer together and points that are far apart maintain that relationship explicitly.

That is, we will choose to look at every point in the dataset as a vertex in a graph where the edge connection between two vertexes is determined by the weighted distances between them.

In the notebook, implement a function that takes in the input dataset and some coefficient gamma and returns the adjacency matrix A :

$$A_{i,j} = e^{-\gamma ||x_i - x_j||^2} \quad (3)$$

where x_i and x_j represent each point in the provided dataset, γ is positive. You may find the `distance` module from `scipy.spatial` useful.

Is this a directed or an undirected graph?

- (d) The degree matrix of an undirected graph is a diagonal matrix which contains information about the degree of each vertex. In other words, it contains the number of edges attached to each vertex and it is given by:

$$D_{i,j} = \begin{cases} \deg(v_i) & \text{if } i = j, \\ 0 & \text{otherwise.} \end{cases}$$

where the degree $\deg(v_i)$ of a vertex counts the number of times an edge terminates at that vertex. Note that in the traditional definition of the adjacency matrix, this boils down to the diagonal matrix in which element along the diagonals are column-wise sum of the adjacency matrix.

Using the same idea, **write a function that takes in the adjacency matrix as an argument and returns the degree matrix.**

- (e) Using $\gamma = 7.5$, compute the adjacency matrix A , degree matrix D and the symmetrically normalized adjacency matrix matrix M ,

$$M = A^{SymNorm} = D^{-1/2}AD^{-1/2} \quad (4)$$

Note that another interpretation of the matrix M is that it shows the probability of moving/jumping from one node to another.

- (f) Applying SVD decomposition on M , **write a function that selects the top 3 vectors (corresponding to the highest singular values) in the matrix U and performs the same KMeans clustering used above on them. Show the plots. What do you observe? Did it work? If so, explain why, if not, explain why not.**

Intuition: By selecting the top 3 vectors of the U matrix, we are selecting a new representation of the data points which could be seen as a construction of a low dimension embedding of the data points as mentioned in problem 3.

- (g) Now let's think of the symmetrically normalized adjacency matrix obtained above as the transition matrix in of a Markov Chain. That is, it represents the probability of jumping from one node to another. In order to fully interpret M in such way, it needs to be a proper stochastic matrix which means that the sum of the elements in each column must add up to 1. **Write a function that takes in the matrix M and returns M_{stoch} , the stochastic version of M; compute the stochastic matrix.**

Using SVD decomposition on the newly obtained stochastic matrix M_{stoch} , **use your function in part (e) to select the top 3 vectors of the matrix U_{stoch} and perform the same KMeans clustering used above on them and show the plots. What do you observe? Did it work?**

6. Homework Process and Study Group

Citing sources and collaborators are an important part of life, including being a student!

We also want to understand what resources you find helpful and how much time homework is taking, so we can change things in the future if possible.

- (a) **What sources (if any) did you use as you worked through the homework?**
- (b) **If you worked with someone on this homework, who did you work with?**
List names and student ID's. (In case of homework party, you can also just describe the group.)
- (c) **Roughly how many total hours did you work on this homework?**

The following are old exam questions simply for your reference. You do not have to do these questions. They're coverage is redundant.

7. Machine Translation

Consider the following Machine Translation problem:

- You are learning an Encoder-Decoder model to translate sentences from Spanish into English.
- This Encoder-Decoder model consists of a single-layer RNN encoder and a single-layer RNN decoder.
- The first hidden state of the decoder is initialized with the last hidden state of the encoder.
- Words are converted into learned embeddings of length H (hidden state size), before they are passed to the model.

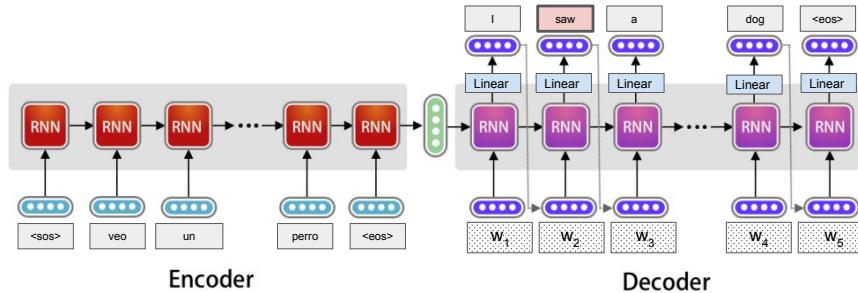


Figure 4: Translation model. The ovals with horizontal dots are learned encodings of words. The tokens $< sos >$ and $< eos >$ are “Start of Sequence” and “End of Sequence” tokens respectively. The boxes $w_1 \dots w_5$ represent the word tokens passed into the RNN decoder at each timestep (you’ll fill in which tokens go here).

- (a) (4pts) Your teammate proposes stacking the encoder and decoder vertically rather than horizontally. Instead of passing the final hidden state of the encoder h_T into the decoder’s first hidden state, at each timestep t , the encoder’s hidden state h_t gets passed as an input to timestep t of the decoder. **State one problem with this proposed design change.**

- (b) (3pts) In the example shown the correct translation is “I see a dog,” but the translation that happened to be sampled from the model incorrectly states “I saw a dog”.

What five tokens will be passed into the decoder during training for w_1, w_2, \dots, w_5 ?

(*HINT: Remember, during training we have access to correct supervision for translations. Don’t forget that you also have special tokens $< sos >$ and $< eos >$ for the beginning and end of a sentence.*)

- (c) (3pts) Continuing the previous part, **what five tokens would be passed into the decoder at evaluation time for w_1, w_2, \dots, w_5 when a translation is being generated?**

(Here, you can assume that the decoder only emits a single possibility for each word.)

8. Self-supervised Linear Autoencoders

We consider linear models consisting of two weight matrices: an encoder $W_1 \in \mathbb{R}^{k \times m}$ and decoder $W_2 \in \mathbb{R}^{m \times k}$ (assume $1 < k < m$). The traditional autoencoder model learns a low-dimensional embedding of the n points of training data $\mathbf{X} \in \mathbb{R}^{m \times n}$ by minimizing the objective,

$$\mathcal{L}(W_1, W_2; \mathbf{X}) = \frac{1}{n} \|\mathbf{X} - W_2 W_1 \mathbf{X}\|_F^2 \quad (5)$$

We will assume $\sigma_1^2 > \dots > \sigma_k^2 > \sigma_{k+1}^2 \geq 0$ are the $k + 1$ largest eigenvalues of $\frac{1}{n} \mathbf{X} \mathbf{X}^\top$. The assumption that the $\sigma_1, \dots, \sigma_k$ are positive and distinct ensures identifiability of the principal components.

Consider an ℓ_2 -regularized linear autoencoder where the objective is:

$$\mathcal{L}_\lambda(W_1, W_2; \mathbf{X}) = \frac{1}{n} \|\mathbf{X} - W_2 W_1 \mathbf{X}\|_F^2 + \lambda \|W_1\|_F^2 + \lambda \|W_2\|_F^2. \quad (6)$$

where $\|\cdot\|_F^2$ represents the Frobenius norm squared of the matrix (i.e. sum of squares of the entries).

- (a) You want to use SGD-style training in PyTorch (involving the training points one at a time) and self-supervision to find W_1 and W_2 which optimize (6) by treating the problem as a neural net being trained in a supervised fashion. **Answer the following questions and briefly explain your choice:**

- (i) **How many linear layers do you need?**

- 0
- 1
- 2
- 3

- (ii) **What is the loss function that you will be using?**

- nn.L1Loss
- nn.MSELoss
- nn.CrossEntropyLoss

- (iii) **Which of the following would you need to optimize (6) exactly as it is written? (Select all that are needed)**

- Weight Decay
- Dropout
- Layer Norm
- Batch Norm
- SGD optimizer

- (b) **Do you think that the solution to (6) when we use a small nonzero λ has an inductive bias towards finding a W_2 matrix with approximately orthonormal columns? Argue why or why not?**

(Hint: Think about the SVDs of $W_1 = U_1 \Sigma_1 V_1^\top$ and $W_2 = U_2 \Sigma_2 V_2^\top$. You can assume that if a $k \times m$ or $m \times k$ matrix has all k of its nonzero singular values being 1, then it must have orthonormal rows or columns. Remember that the Frobenius norm squared of a matrix is just the sum of the squares of its singular values. Further think about the minimizer of $\frac{1}{\sigma^2} + \sigma^2$. Is it unique?)

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