# Tricks for Training Neural Models

(Some slides by Yoav Goldberg, Graham Neubig)

# **Optimization Choices**

#### • Adaptive learning rate.

 adaptive optimizers such as Adam (<u>Kingma14</u>) because they can better handle the complex training dynamics of RNNs

#### • Gradient clipping.

- Print or plot the gradient norm to see its usual range
- then scale down gradients that exceeds this range.
- This prevents spikes in the gradients to mess up the parameters during training.
- Normalizing the loss. (To get losses of similar magnitude across datasets)
  - sum the loss terms along the sequence and divide them by the maximum seq length.
  - This makes it easier to reuse hyper parameters between experiments.
  - The loss should be averaged across the batch.

#### • Early Stopping

https://danijar.com/tips-for-training-recurrent-neural-networks/

# Network Structure (RNN)

- Use Gated Recurrent Unit.
- Layer normalization. Adding layer normalization (Ba et al 16) to all linear mappings of the recurrent network speeds up learning

#### Stacked recurrent networks.

- Recurrent networks need a quadratic number of weights in their layer size.
- More efficient to stack two or three smaller layers instead of one big one.
- Sum the outputs of all layers instead of using only the last one, similar to a ResNet or DenseNet.

# Model Parameters (RNN)

#### • Learned initial state.

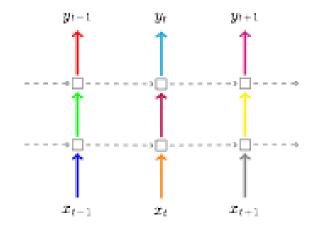
- Initializing the hidden state as zeros  $\rightarrow$  large loss initially
- Training the initial state as a variable can improve performance as described in <a href="https://r2rt.com/non-zero-initial-states-for-recurrent-neural-networks.html">https://r2rt.com/non-zero-initial-states-for-recurrent-neural-networks.html</a>

#### • Forget gate bias.

- It can take a while for a RNN to learn to remember information
- Initialize biases for LSTM's forget gate to 1 to remember more by default.
- Similarly, initialize biases for GRU's reset gate to -1.
- **Regularization.** If your model is overfitting, use dropout

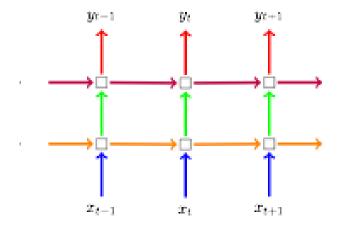
### Dropout in RNNs

- Still an open question how to perform well.
- One suggestion:



### Dropout in RNNs

- Still an open question how to perform well.
- Yarin Gal's Dropout:



# Ensembles

#### • Same model, different initialization.

- Use cross-validation to determine the best hyperparameters,
- then train multiple models with the best set of hyperparameters but with different random initialization.
- Suffers from limited variety

#### • Top models discovered during cross-validation.

- Use cross-validation to determine the best hyperparameters
- then pick the top few (e.g., 10) models to form the ensemble.
- Improves the variety of ensemble but has the danger of including suboptimal models

#### • Different checkpoints of a single model.

- If training is very expensive
- limited success in taking different checkpoints of a single network over time (for example after every epoch) and using those to form an ensemble.
- Clearly, this suffers from some lack of variety, but can still work reasonably well in practice.
- The advantage of this approach is that is very cheap.

http://lamda.nju.edu.cn/weixs/project/CNNTricks/CNNTricks.html

# Why are Neural Networks Slow and What Can we Do?

- Big operations, especially for softmaxes over large vocabularies
  - → Approximate operations or use GPUs
- GPUs love big operations, but hate doing lots of them
  - → Reduce the number of operations through optimized implementations or batching
- Our networks are big, our data sets are big
  - → Use parallelism to process many data at once

# Approximating Softmax

- Importance Sampling
- Noise Contrastive Estimation
- Simple Negative Sampling
- Hierarchical Softmax

# GPUs vs. CPUs

#### CPU, like a motorcycle

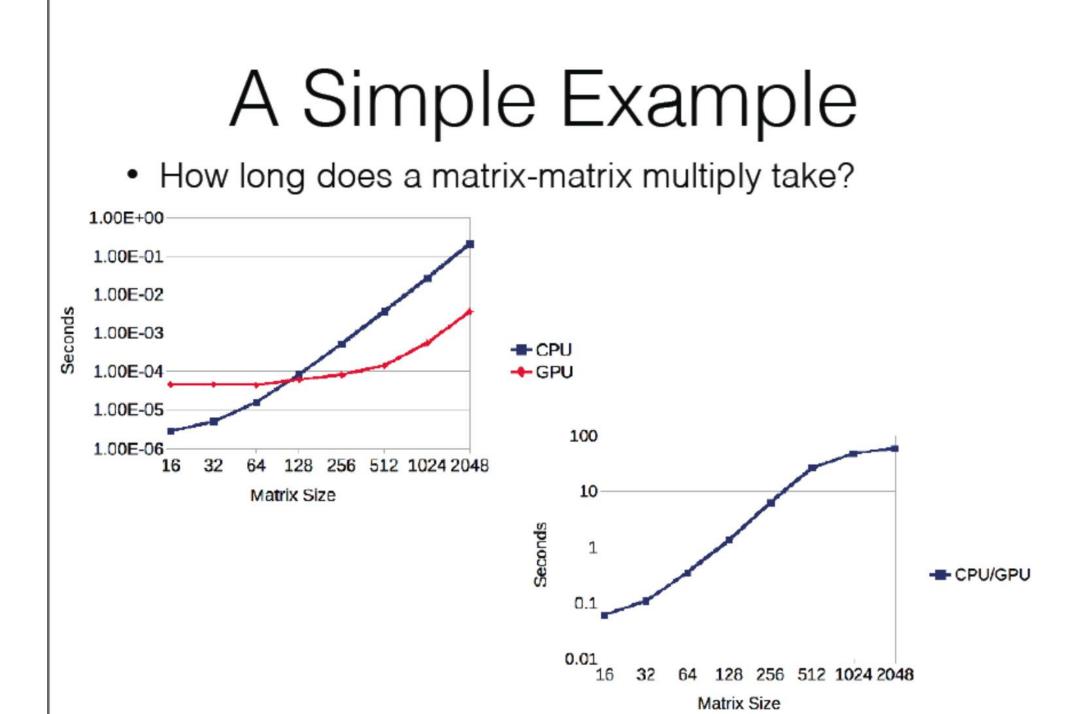


Quick to start, top speed not shabby

#### GPU, like an airplane



Takes forever to get off the ground, but super-fast once flying



# Practically

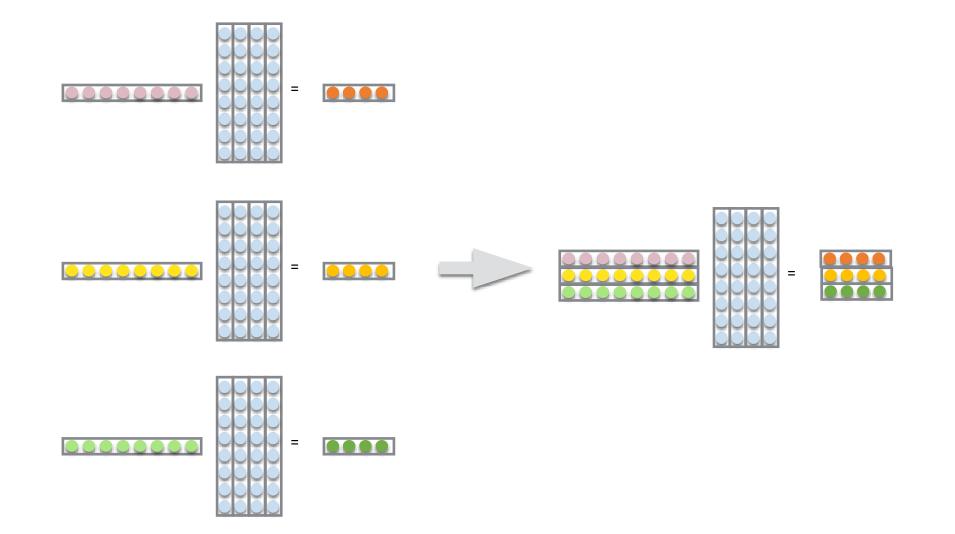
- Use CPU for profiling, it's plenty fast (esp. DyNet) and you can run many more experiments
- For many applications, CPU is just as fast or faster than GPU: NLP analysis tasks with small or complicated data/networks
- You see big gains on GPU when you have:
  - Very big networks (or softmaxes with no approximation)
  - Do mini-batching
  - Optimize things properly

# Batching (in RNNs)

- Most toolkits require a **fixed** computation graph for all examples.
- But RNNs have **different** input lengths. What do we do?
  - Option 1:
     Use a tool that does not pose this limitation.
  - Option 2:

Enforce max length + 0 padding for shorter sequences.

# Batching Reminder

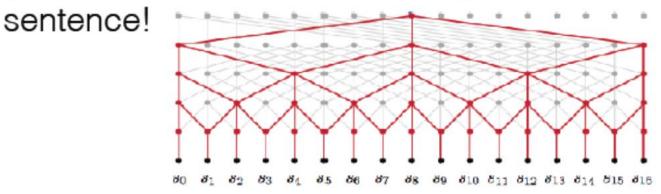


# Batching in RNNs

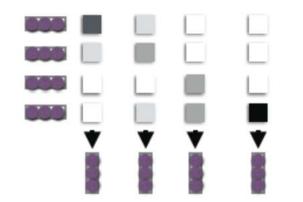
- Sequential in nature, very little parallelism.
- (Compare, e.g., to a Convolutional Network)

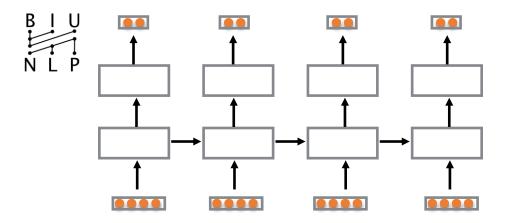
# Non-recursive Architectures

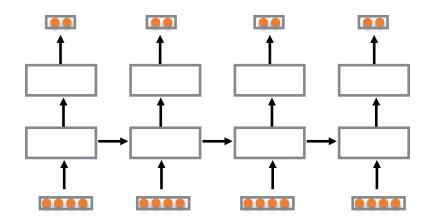
 Dilated convolutions for capturing context (Kalchbrenner et al. 2016): single GPU call for entire

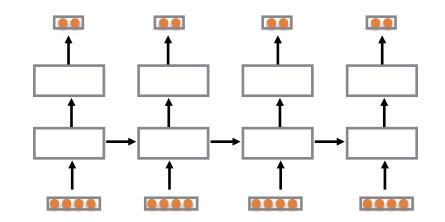


• Self-attention that decides which of previous words to focus on (Cheng et al. 2016, Vaswani et al. 2017, covered in detail a few classes): also single GPU call

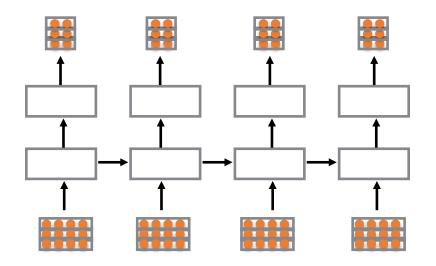






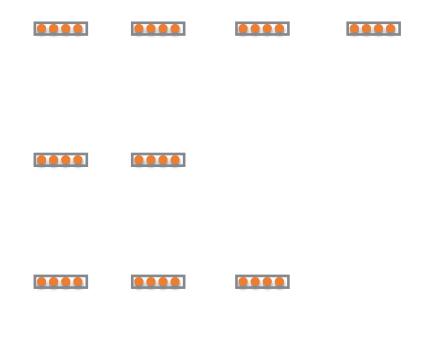




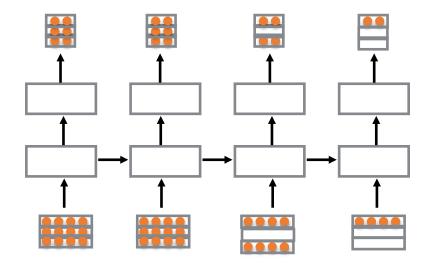




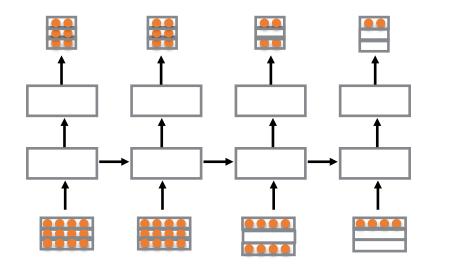
what if the sequences are different lengths?





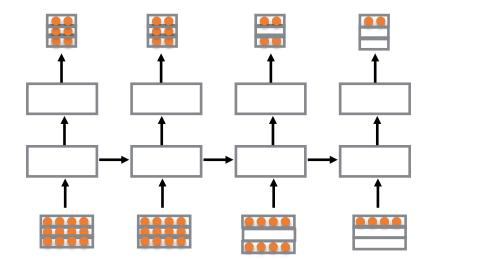






padding

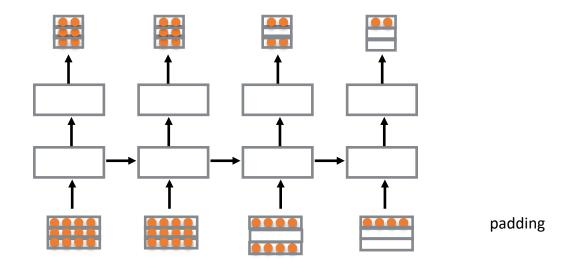




padding

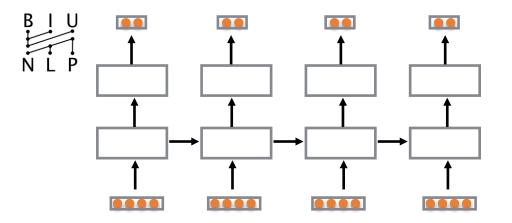
this is how its done in TF, PyTorch. supported also in DyNet, but...

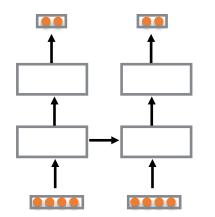




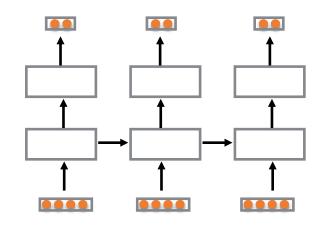
We want better

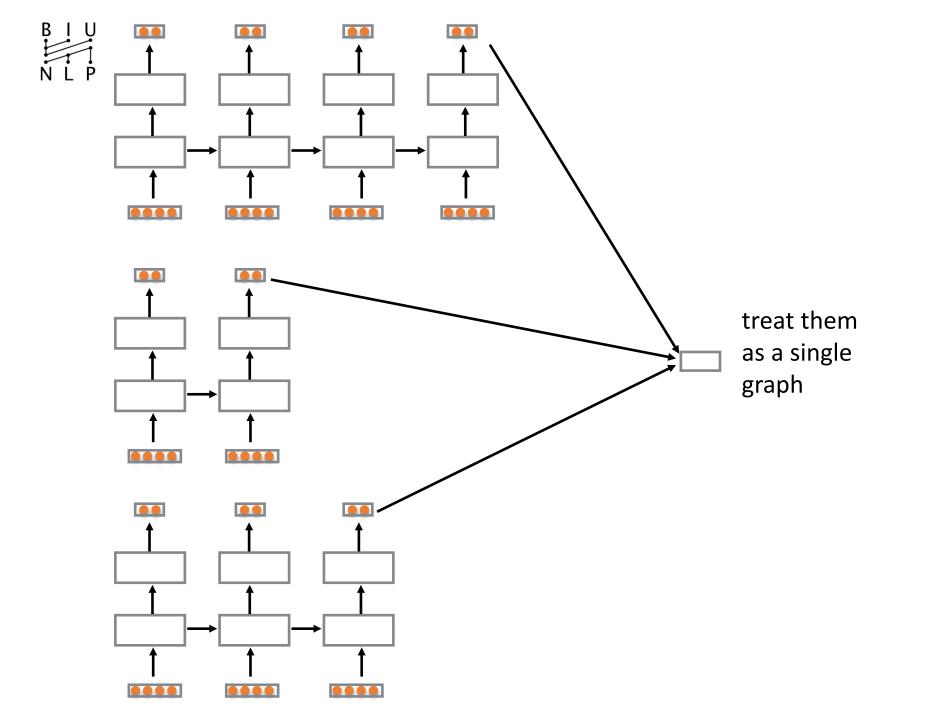
# Auto Batching

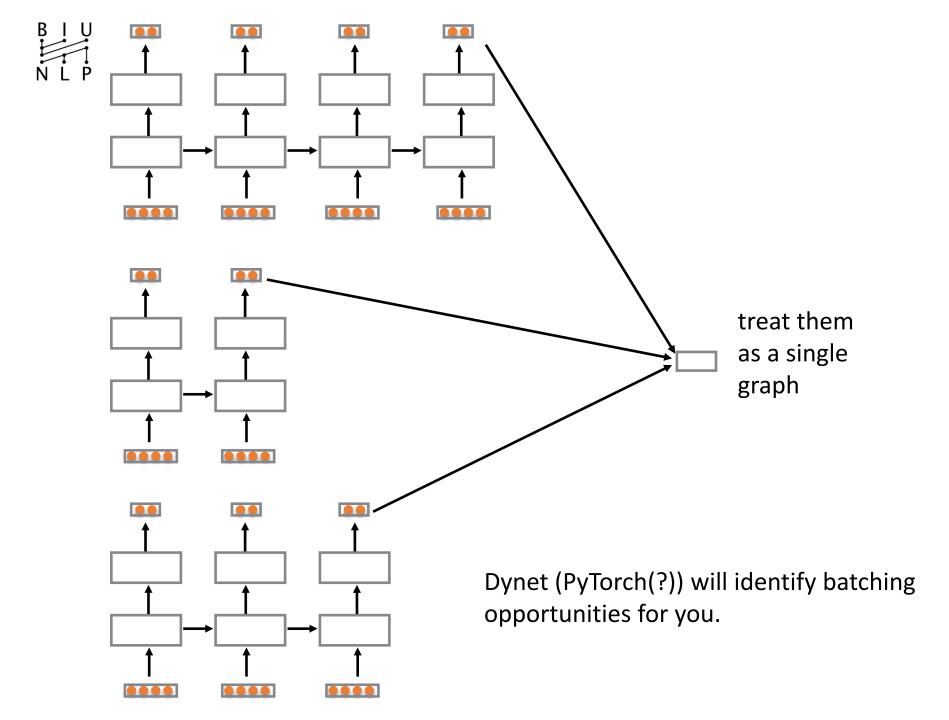


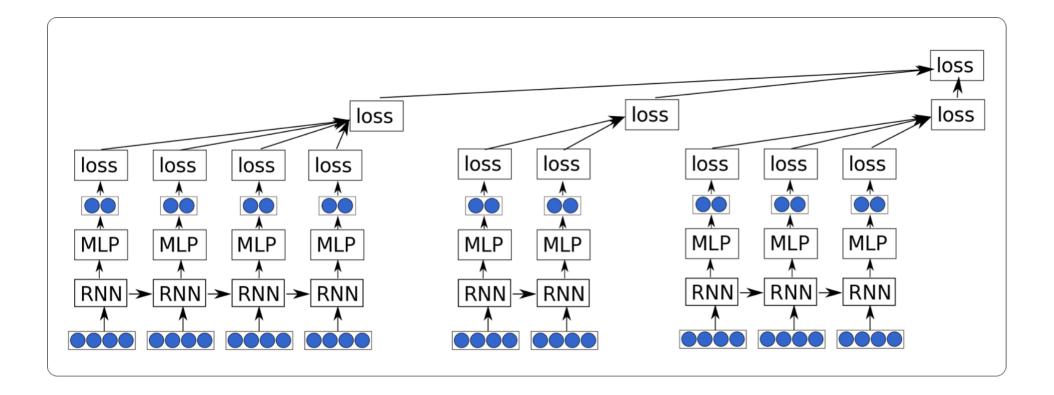


create a separate network for each (easy)

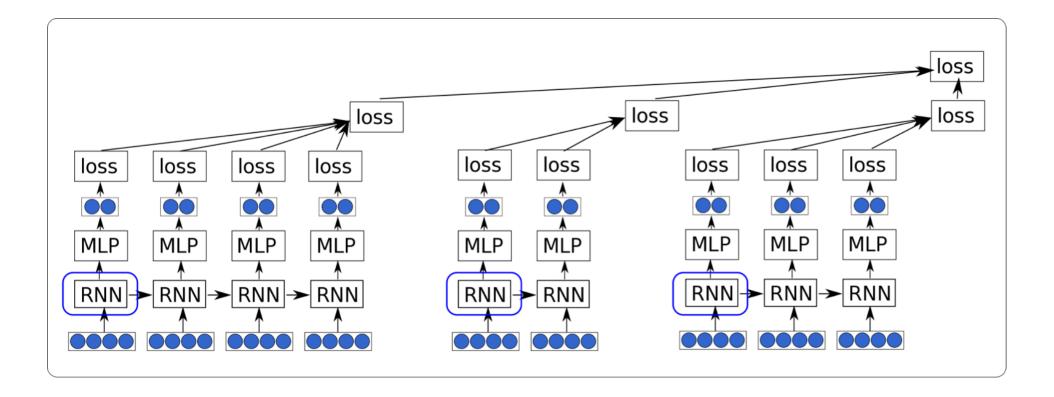




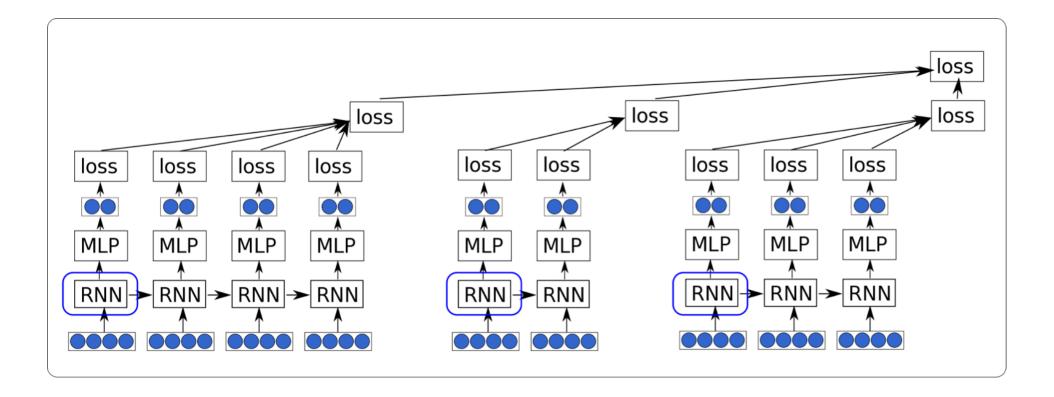


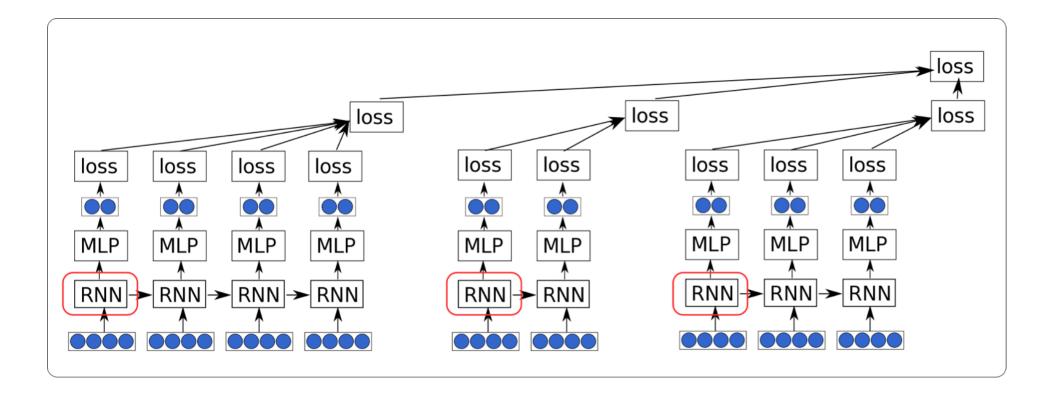


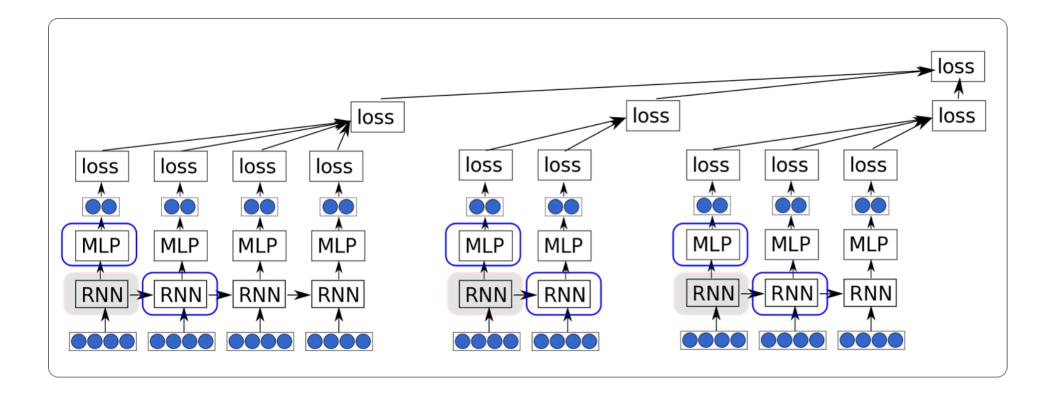
nodes in blue are ready to be executed

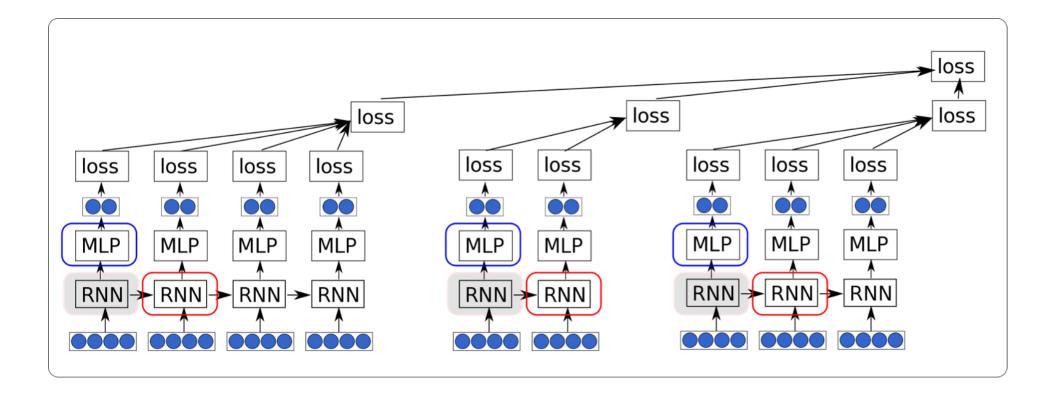


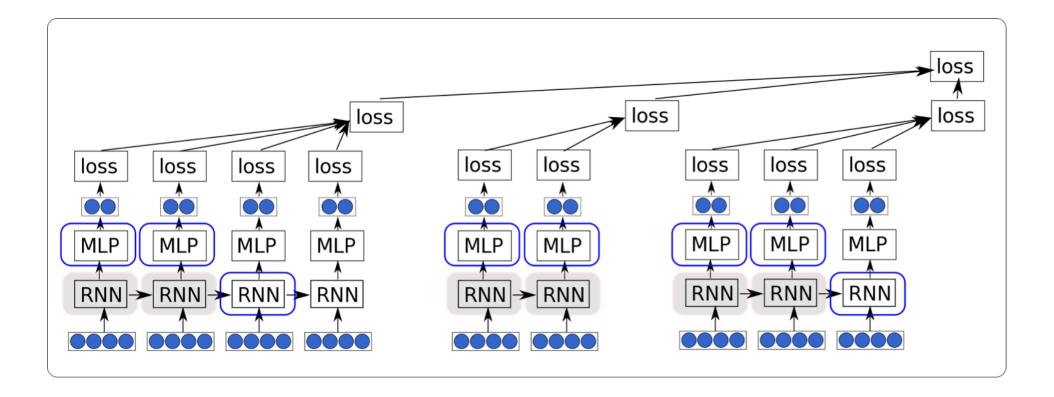
nodes in red will be executed using batch operations

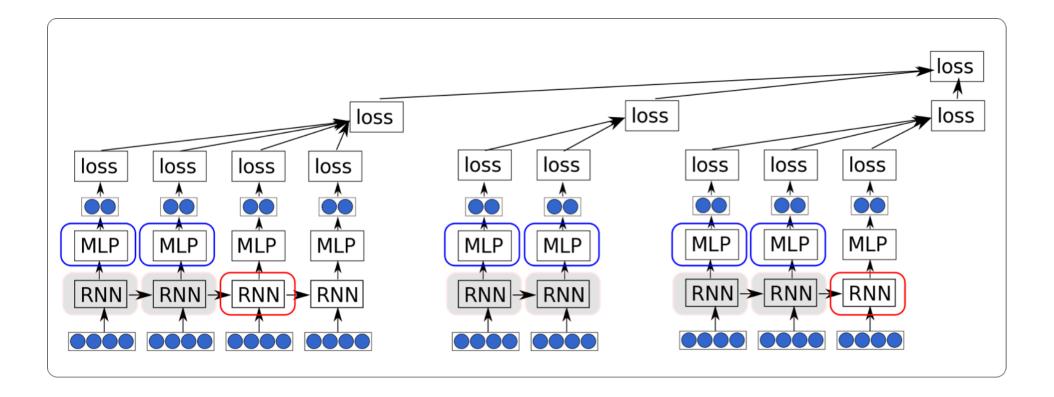


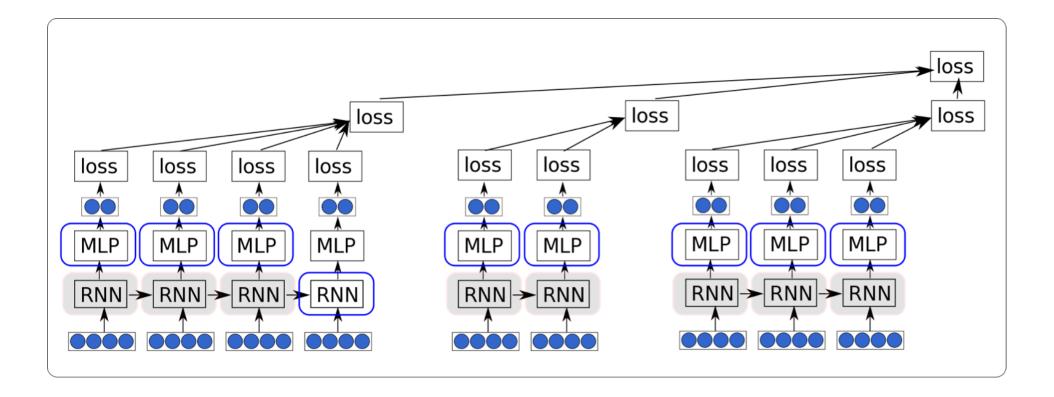


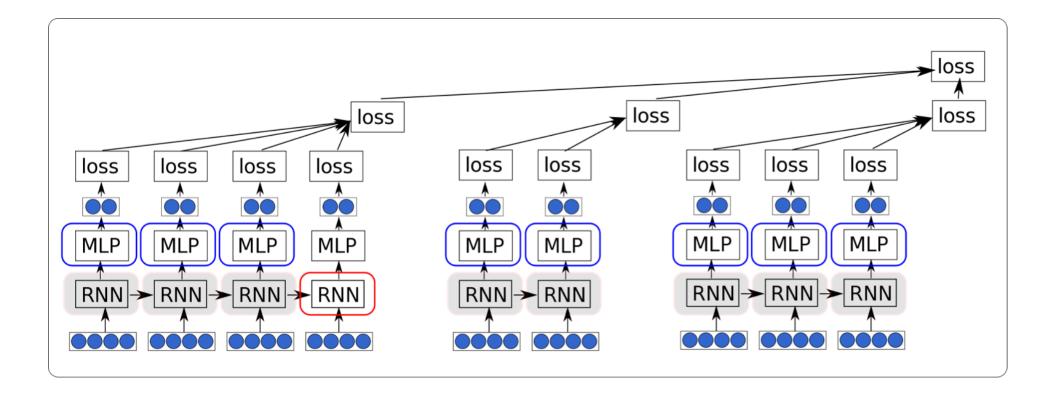


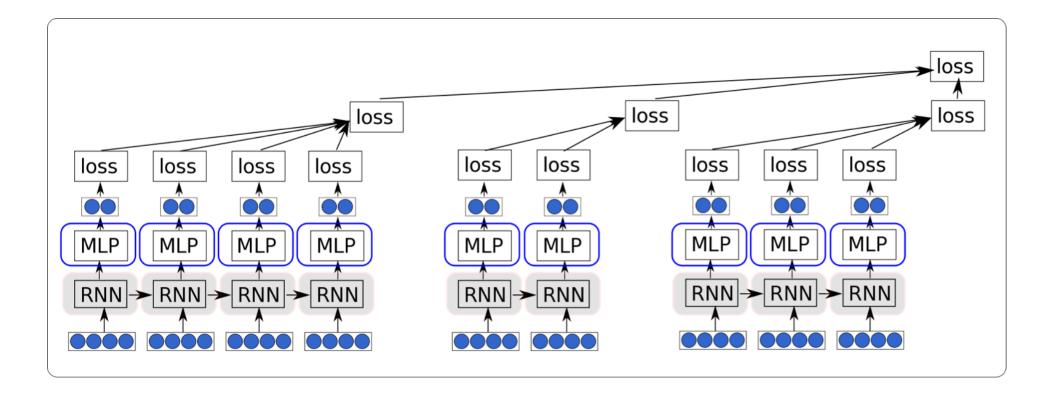


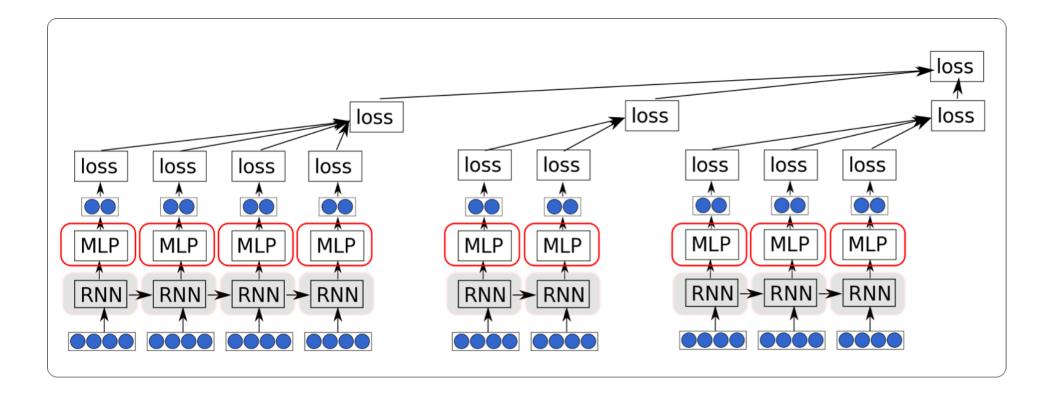


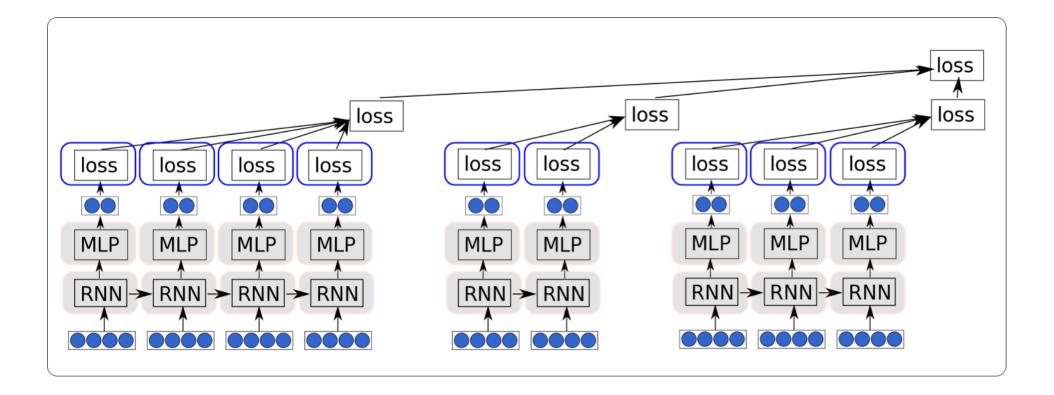


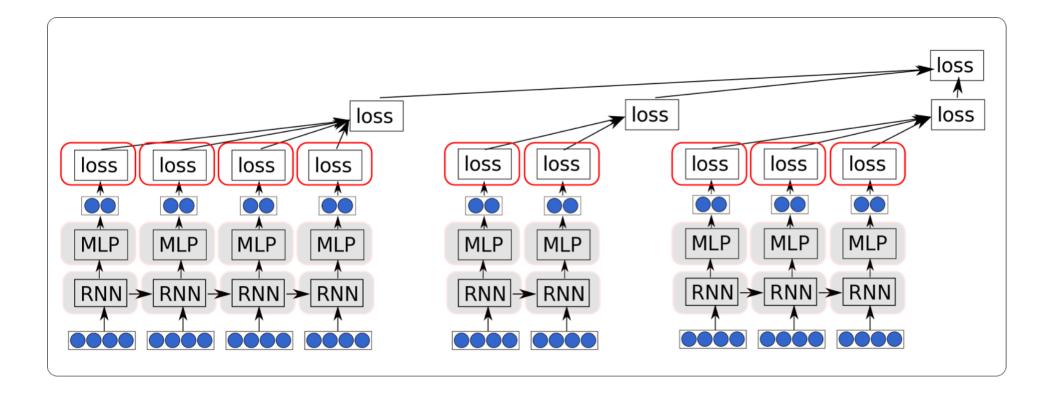


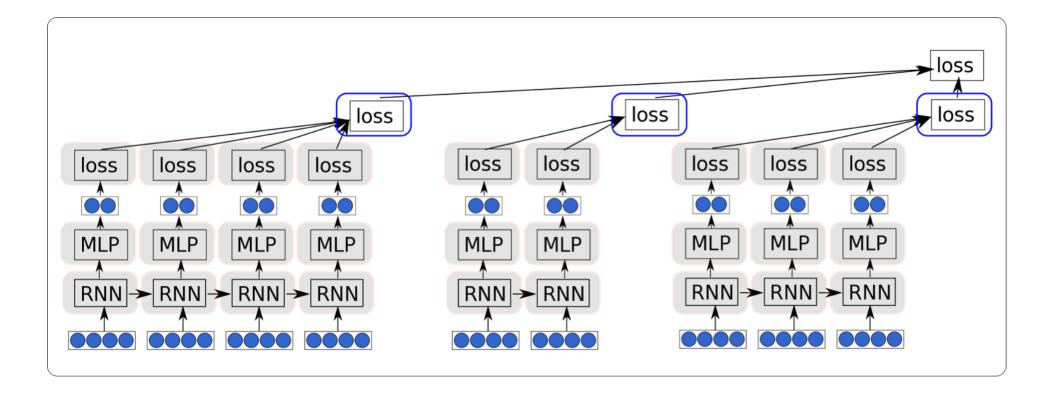


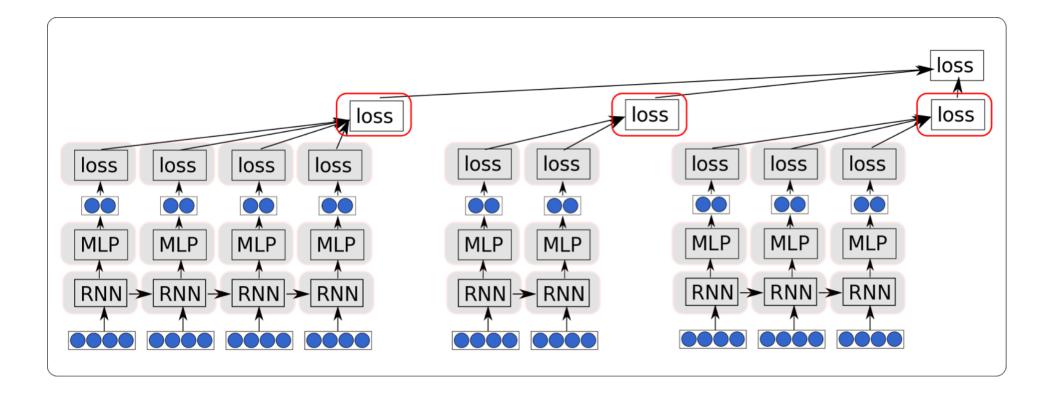


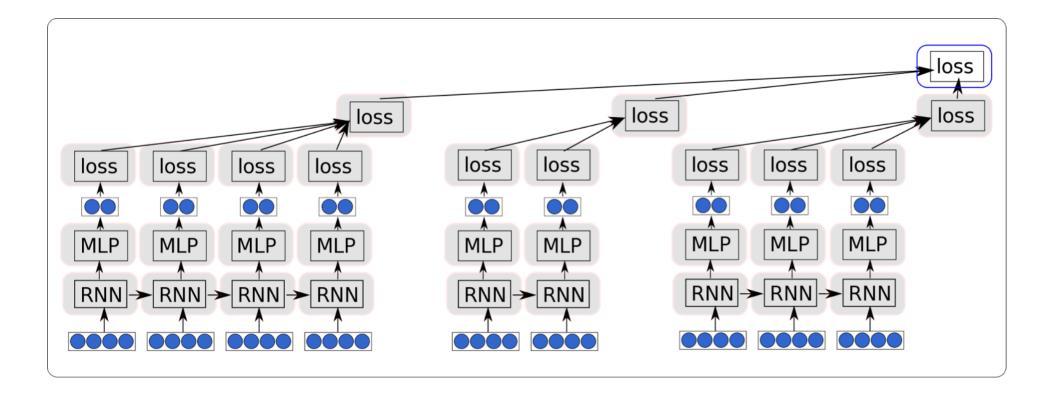


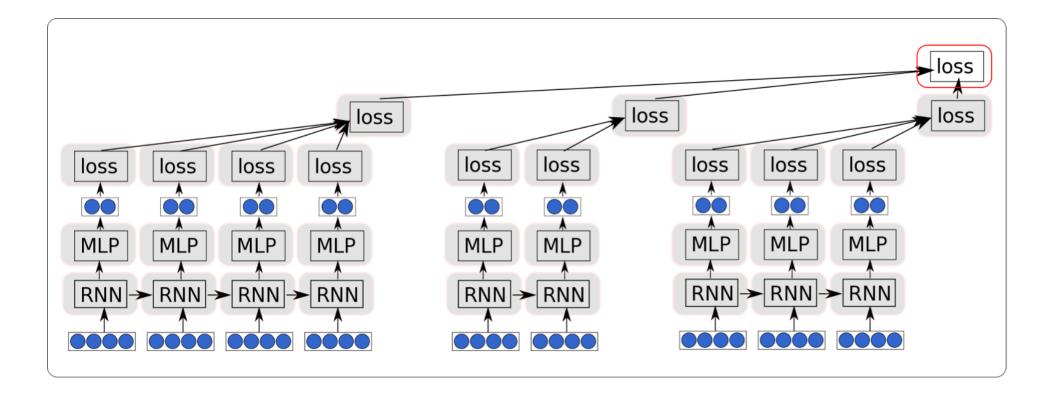


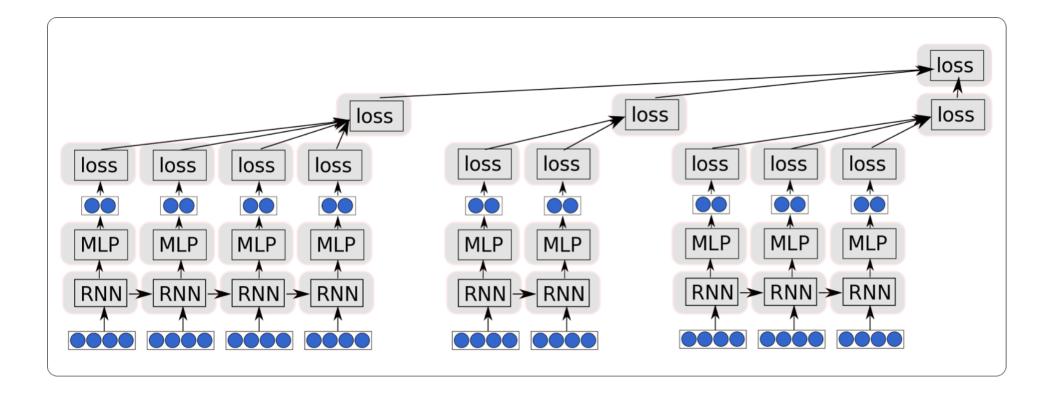


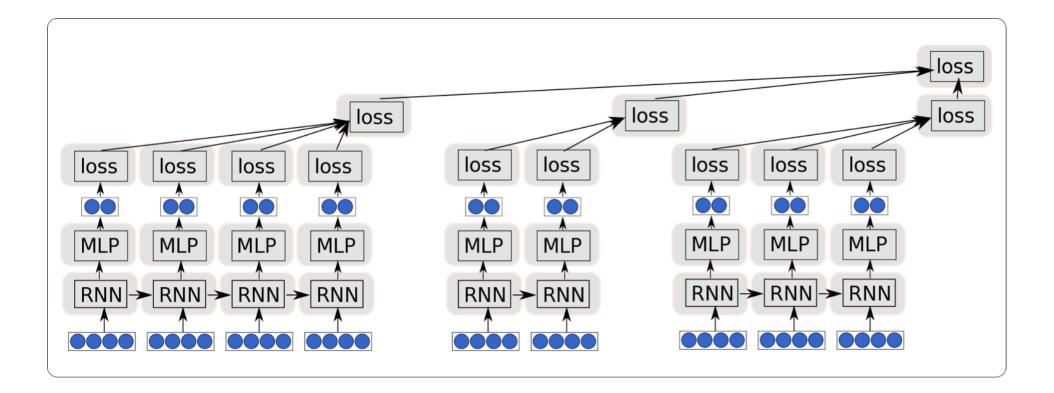












<u>note</u>: batching operations, not inputs.

# Efficiency Considerations when Implementing an LSTM

$$R_{LSTM}(\mathbf{s_{j-1}}, \mathbf{x_j}) = [\mathbf{c_j}; \mathbf{h_j}]$$

$$\mathbf{c_j} = \mathbf{c_{j-1}} \odot \mathbf{f} + \mathbf{g} \odot \mathbf{i}$$

$$\mathbf{h_j} = \tanh(\mathbf{c_j}) \odot \mathbf{o}$$

$$\mathbf{i} = \sigma(\mathbf{W^{xi}} \cdot \mathbf{x_j} + \mathbf{W^{hi}} \cdot \mathbf{h_{j-1}})$$

$$\mathbf{f} = \sigma(\mathbf{W^{xf}} \cdot \mathbf{x_j} + \mathbf{W^{hf}} \cdot \mathbf{h_{j-1}})$$

$$\mathbf{o} = \sigma(\mathbf{W^{xo}} \cdot \mathbf{x_j} + \mathbf{W^{ho}} \cdot \mathbf{h_{j-1}})$$

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all gates computations can be done in single mat-mat op.

## Speed Trick 1: Don't Repeat Operations

 Something that you can do once at the beginning of the sentence, don't do it for every word!

### Bad

for x in words\_in\_sentence:
 vals.append( W \* c + x )

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### Good

W\_c = W \* c
for x in words\_in\_sentence:
 vals.append(W\_c + x)

# Speed Trick 2: Reduce # of Operations

 e.g. can you combine multiple matrix-vector multiplies into a single matrix-matrix multiply? Do so!

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for x in words\_in\_sentence:
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```
for x in words_in_sentence:
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val = dy.concatenate(vals)
```

### Good

X = dy.concatenate\_cols(words\_in\_sentence)
val = W \* X

### Speed Trick 3: Reduce CPU-GPU Data Movement

- Try to avoid memory moves between CPU and GPU.
- When you do move memory, try to do it as early as possible (GPU operations are asynchronous)

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```
for x in words_in_sentence:
    # input data for x
    # do processing
```

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### **Bad**

for x in words\_in\_sentence:
 # input data for x
 # do processing

### Good

# input data for whole sentence
for x in words\_in\_sentence:
 # do processing