

# Deep Learning Clinic (DLC)

Lecture 5
Neural Network Architectures and Model Search

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10/8/2019

#### Today - Neural Network Architectures and Model Search

- Neural Network Architecture Design
- Hyperparameters and Network Sensitivity
- Optimize Hyperparameters Systematically
- Automatic Model Search AutoML

## Design a NN model for your problem

Things to consider:

- ☐ The characteristic of your data
- Data input/output space
- Budget on hardwares, training time...

#### Recap: Data and Neural Network Models

**Static Data** 

Convolutional Neural Networks **Dynamic Data** 

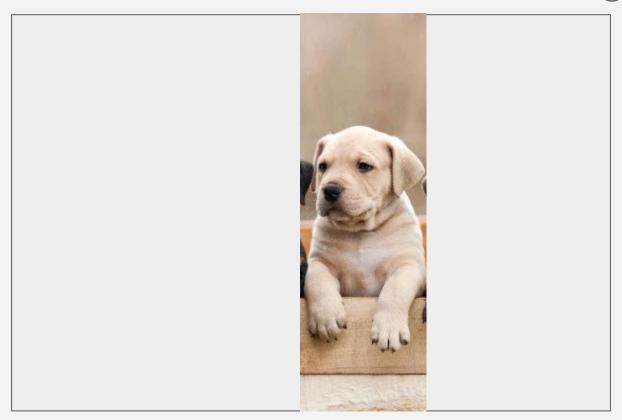
Recurrent Neural Networks **Unsupervised Data** 

Generative Neural Networks

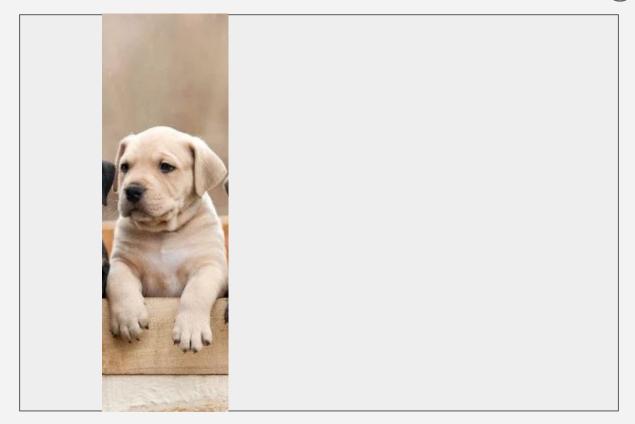
# Static Data - Image



## Static Data - Translation invariance in images

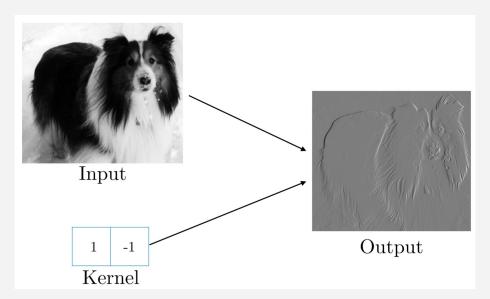


### Static Data - Translation invariance in images

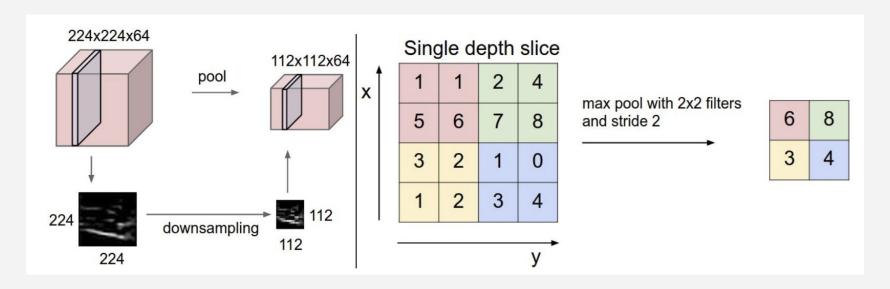


#### Convolution

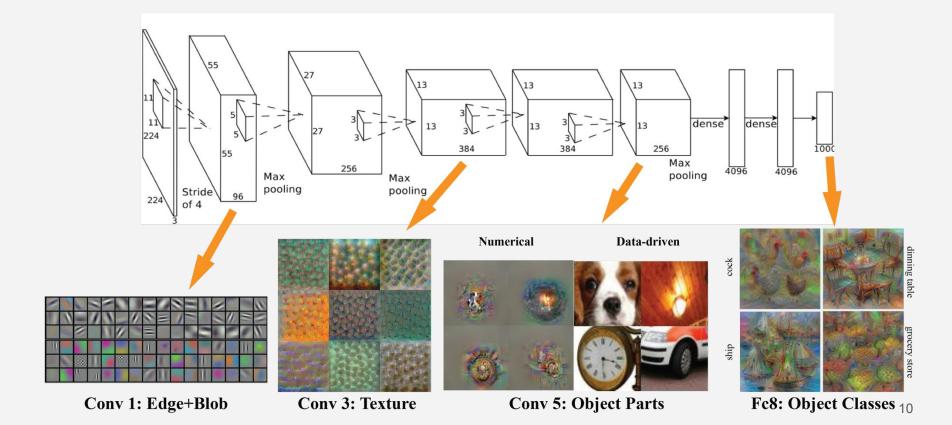
A local operation that extracts information from data.



#### **Pooling**



#### Convolutional Networks

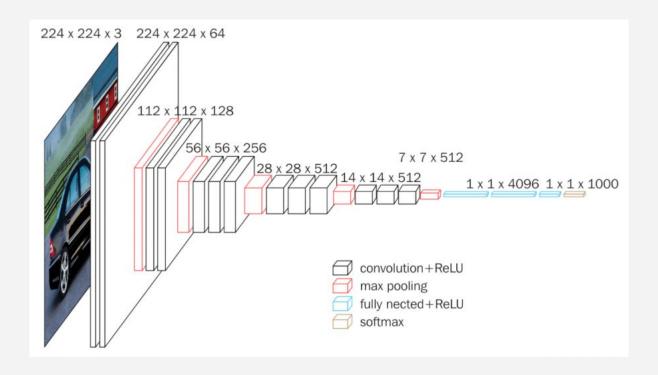


## Design a NN model for your problem

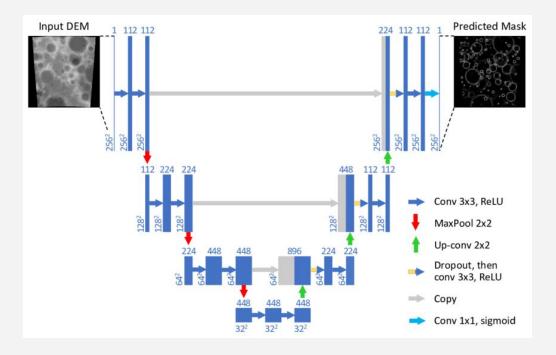
Things to consider:

- ☐ The characteristic of your data
- ☐ Data input/output space
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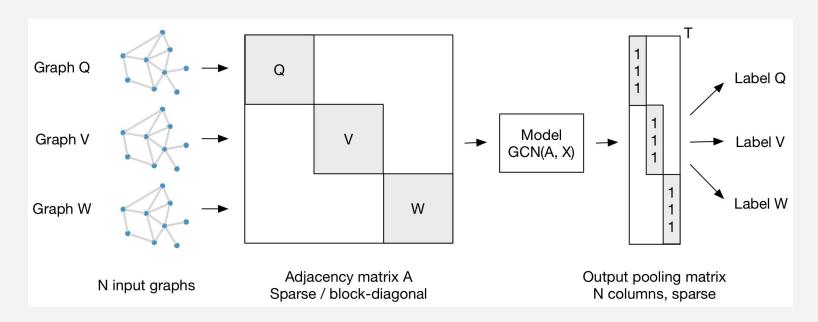
## Predicting a category



#### Dense predictions



## Input is a graph



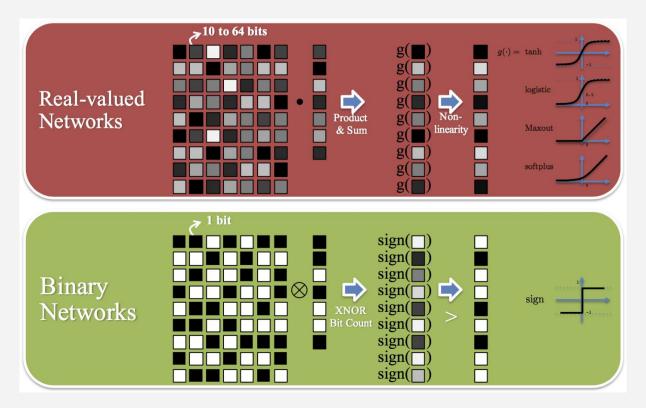
Graph Conv Net

## Design a NN model for your problem

Things to consider:

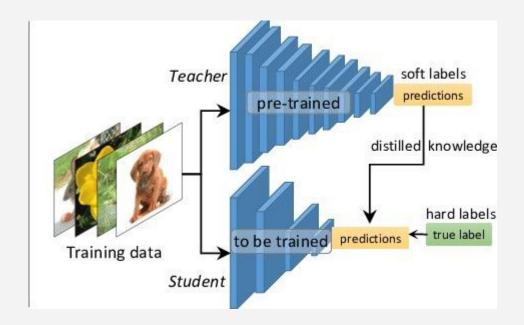
- ☐ The characteristic of your data
- ☐ Data input/output space
- **■** Budget on hardwares, training time...

## Binary Networks



## Knowledge Distillation

Train a smaller network (Student) by learning from a more powerful network (Teacher).



#### Today - Neural Network Architectures and Model Search

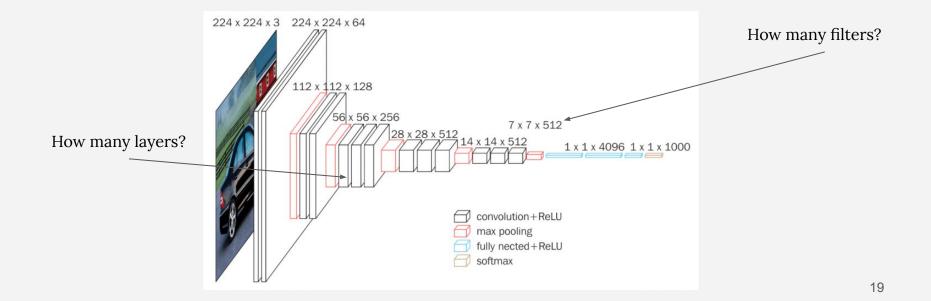
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#### A recommended read:

Bengio, Yoshua. "Practical recommendations for gradient-based training of deep architectures." *Neural networks: Tricks of the trade*. Springer, Berlin, Heidelberg, 2012. 437-478.

#### Neural Network Model Search

Even with a targeted network architecture, there are still a lot of flexibility in the exact configuration of your neural network framework.



#### Hyperparameters

Learning rate: controls how much you update the model's weights.

**Batch size**: how many data you put into the model each time to calculate the gradient.

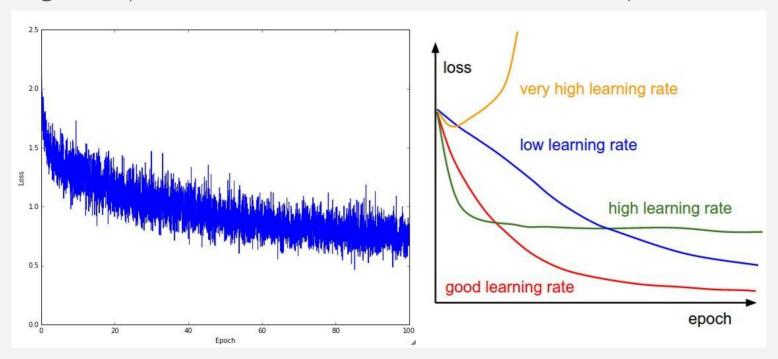
**Regularization**: e.g., weight decay, L1.

Network structure: number of layers, activation functions, etc.

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Even **Data**: it is almost always guaranteed that a model would get better with more data.

**Learning Rate** (we will cover a lot on this in Lecture - Tricks)



#### Batchsize

$$\theta = \theta - \alpha \bigtriangledown_{\theta} E[J(\theta)]$$

Decrease learning rate is to reduce the noise of each batch's gradient update.

\_

Increase batchsize to reduce the noise in each batch.

#### Batchsize

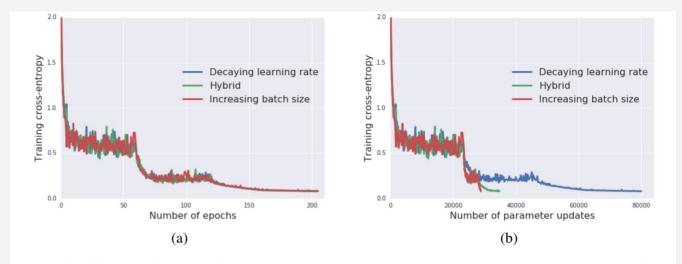
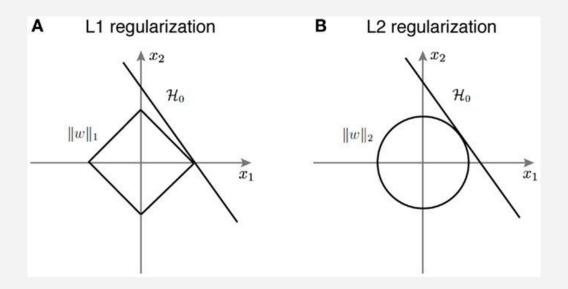


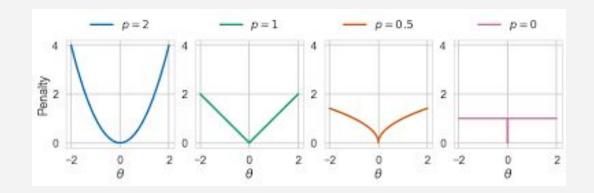
Figure 2: Wide ResNet on CIFAR10. Training set cross-entropy, evaluated as a function of the number of training epochs (a), or the number of parameter updates (b). The three learning curves are identical, but increasing the batch size reduces the number of parameter updates required.

#### Regularization



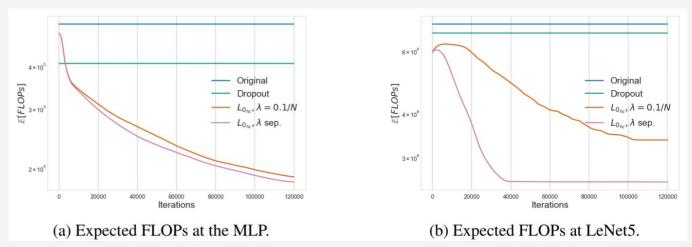
https://www.quora.com/Why-is-L1-regularization-supposed-to-lead-to-sparsity-than-L2

#### Regularization



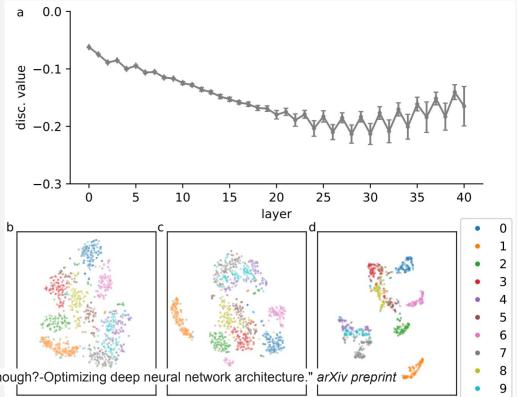
#### Regularization

Training with L0 regularization = model pruning



Louizos, Christos, Max Welling, and Diederik P. Kingma. "Learning Sparse Neural Networks through \$ L\_0 \$ Regularization." arXiv preprint arXiv:1712.01312 (2017).

Number of Layers/Depth



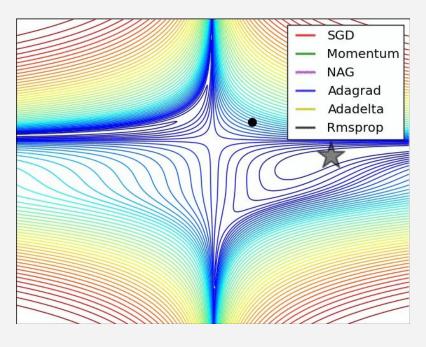
Schilling, Achim, et al. "How deep is deep enough?-Optimizing deep neural network architecture." arXiv preprint arXiv:1811.01753 (2018).

Number of Layers/Depth

Deeper networks are harder to train.

Consider add normalization (BatchNorm and its variants) and ResNet modules to the network for **stability** and **strong gradients**.

#### **Optimizer**



http://ruder.io/optimizing-gradient-descent/

#### **Optimizer**

```
torch.optim.Adadelta(params, 1r=1.0, rho=0.9, eps=1e-06, weight decay=0)
torch.optim.Adagrad (params, 1r=0.01, 1r decay=0, weight decay=0, initial accumulator value=0)
torch.optim.Adam(params, 1r=0.001, betas=(0.9, 0.999), eps=1e-08, weight decay=0, amsgrad=False)
torch.optim.ASGD(params, 1r=0.01, lambd=0.0001, alpha=0.75, t0=1000000.0, weight decay=0)
torch.optim.SGD (params, 1r=<required parameter>, momentum=0, dampening=0, weight decay=0, nesterov=False)
torch.optim.LBFGS (params, 1r=1, max iter=20, max eval=None, tolerance grad=1e-05, tolerance change=1e-09,
history size=100, line search fn=None)
```

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## Before Your Start Hyperparameter Search

Sanity Check!

Use a **small portion** of the data, remove all the fancy stuff.

Train your model, you should be able to **overfit** the model to your small data: very high accuracy on your training data (not validation data).

If not, there is some bug in your pipeline.

## This is Probably Your Current Approach

Use a set of default hyperparameter settings.

Train a model.

Check for metric.

Randomly change things.

Repeat above until satisfaction.

# Optimize Hyperparameters Systematically

#### **Master Program**

**Worker Program** 

Choose hyperparameters from a \_\_\_\_\_ Train on data preset range

Validate model on a hold out set

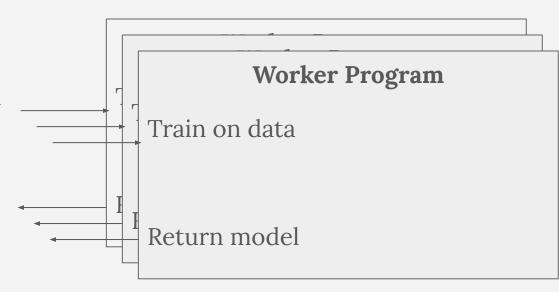
——— Return model

## Optimize Hyperparameters Systematically

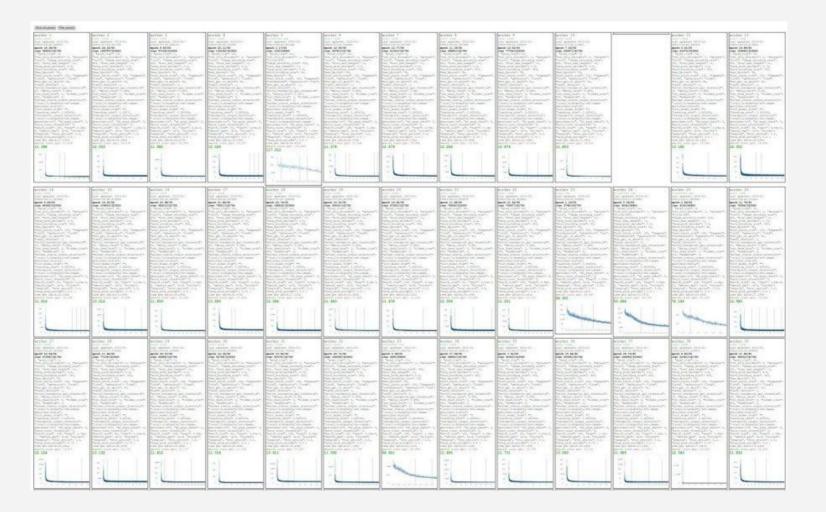
#### **Master Program**

Choose hyperparameters from a preset range

Validate model on a hold out set



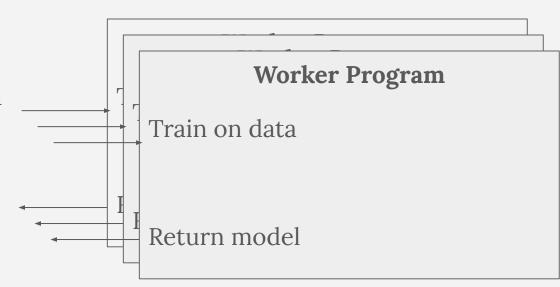
Scalable



### **Master Program**

**Choose** hyperparameters from a preset range

Validate model on a hold out set



Choose your train and validation set.

Many dataset comes with default ones, but sometimes it is good to pick a **subset for fast prototyping**.

7:3 train to validation ratio is traditionally suggested.

8:2 is also more popular now.

#### N-Fold **Cross Validation?**

In early stage of hyperparameter tuning it is enough to use a single set.

Set a **range** of your parameters:

Learning rate: [1e-3, 1e-6] with Adam, might be different for diff optimizer.

Batchsize: [16, 256], depending on the memory size.

Regularizer: {(L0), L1, L2}

• • •

Almost always good to work in **log space** for continuous values.

Almost always good to work in log space for continuous values

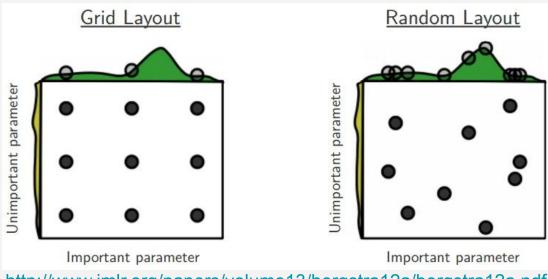
Learning rate values to be picked:

1e-3, 1e-4, 1e-5, 1e-6, ...

Batchsize:

8, 16, 32, 64, ...

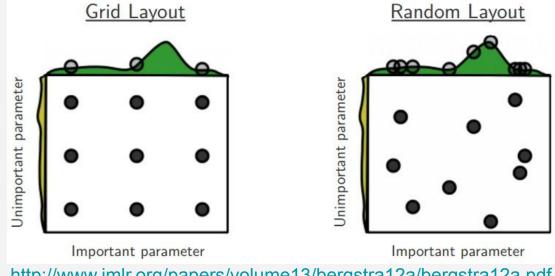
Grid search vs random search in the parameter range



http://www.jmlr.org/papers/volume13/bergstra12a/bergstra12a.pdf

Grid search vs random search in the parameter range

Because the parameters are not equally good in the parameter space



http://www.jmlr.org/papers/volume13/bergstra12a/bergstra12a.pdf

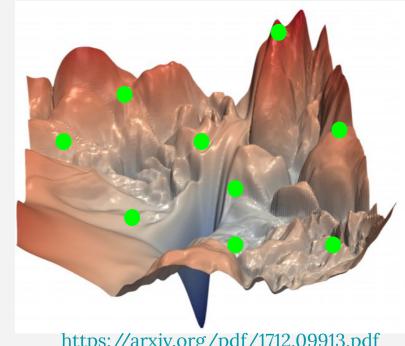
#### **Coarse to Fine Process**

Start with few epochs, on a wider range of parameters

Then use finer scale and run more epochs

#### Coarse to Fine Process

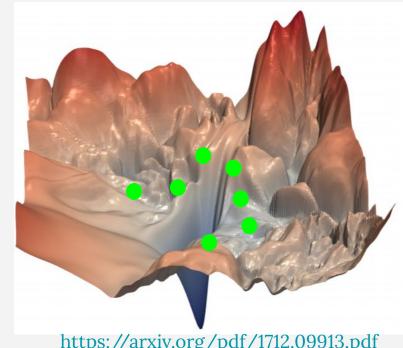
Always keep in mind this is a non-convex optimization



https://arxiv.org/pdf/1712.09913.pdf

#### Coarse to Fine Process

Always keep in mind this is a non-convex optimization

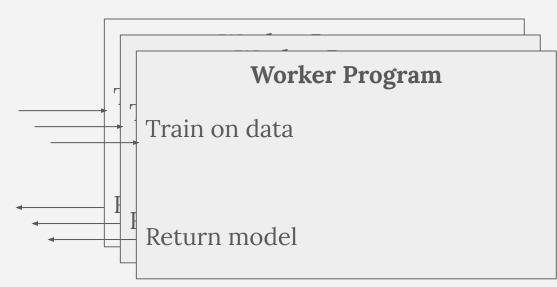


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### **Master Program**

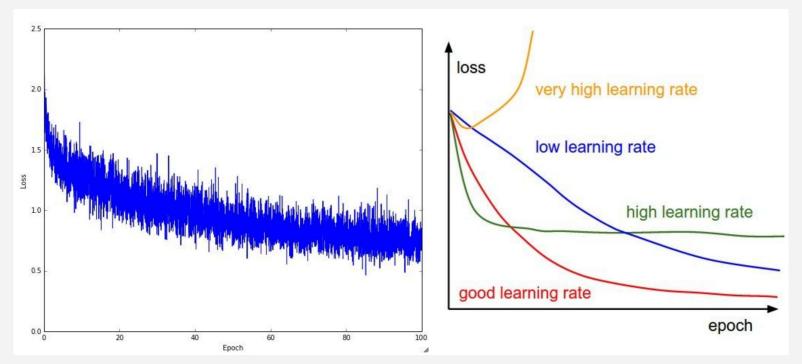
Choose hyperparameters from a preset range

Validate model on a hold out set



## What to check during the training process?

[Major] Loss values comparison between train and validation



## Look deeper

### Track the ratio of weight updates / weight magnitudes:

```
# assume parameter vector W and its gradient vector dW
param_scale = np.linalg.norm(W.ravel())
update = -learning_rate*dW # simple SGD update
update_scale = np.linalg.norm(update.ravel())
W += update # the actual update
print update_scale / param_scale # want ~1e-3
```

ratio between the updates and values:  $\sim 0.0002 / 0.02 = 0.01$  (about okay) want this to be somewhere around 0.001 or so

# What to check during the training process?

Visual inspection of predicted results might be also useful.

For example, take a look at generated images, depth maps, semantic segmentation maps.

There might be 2.3% of corrupted data in your dataset! You cannot tell this by just looking at metric numbers.

The hard way:

### **Bayesian Hyperparameter Optimization**

The Bayesian view:

Each hyperparameter is a random variable from some distribution. The goal is to find the optimal set of hyperparameters from sampling from the distribution and observe their performance.

Example methods: Spearmint, SMAC, and Hyperopt

### Remember to consult the literature

Try to take a look at state-of-the-art related work.

Our implementation for ImageNet follows the practice in [21, 41]. The image is resized with its shorter side randomly sampled in [256, 480] for scale augmentation [41]. A  $224\times224$  crop is randomly sampled from an image or its horizontal flip, with the per-pixel mean subtracted [21]. The standard color augmentation in [21] is used. We adopt batch normalization (BN) [16] right after each convolution and before activation, following [16]. We initialize the weights as in [13] and train all plain/residual nets from scratch. We use SGD with a mini-batch size of 256. The learning rate starts from 0.1 and is divided by 10 when the error plateaus, and the models are trained for up to  $60\times10^4$  iterations. We use a weight decay of 0.0001 and a momentum of 0.9. We do not use dropout [14], following the practice in [16].

In testing, for comparison studies we adopt the standard 10-crop testing [21]. For best results, we adopt the fully-convolutional form as in [41, 13], and average the scores at multiple scales (images are resized such that the shorter side is in {224, 256, 384, 480, 640}).

### Get the Most From Your Models

#### **Model Ensembles**

Instead of having a single trained model, you could generate a number of similar models and get a slightly better overall performance.

• Same model with different initializations.

Train model with the same set of hyperparameters, but with different random initialized weights.

### Get the Most From Your Models

#### **Model Ensembles**

Instead of having a single trained model, you could generate a number of similar models and get a slightly better overall performance.

• Top k models you found in cross validation.

Different snapshots of the trained model.

### Get the Most From Your Models

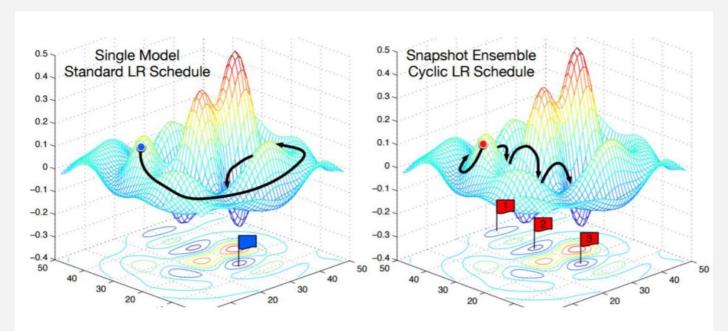
#### **Model Ensembles**

Instead of having a single trained model, you could generate a number of similar models and get a slightly better overall performance.

Final result is an averaged consensus over all models in your ensemble.

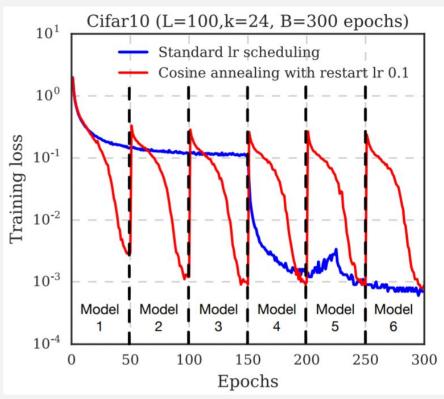
Could also take a weighted vote based on their validation accuracy.

## An Optimization View of Model Ensembles



Loshchilov and Hutter, "SGDR: Stochastic gradient descent with restarts", arXiv 2016 Huang et al, "Snapshot ensembles: train 1, get M for free", ICLR 2017 Figures copyright Yixuan Li and Geoff Pleiss, 2017. Reproduced with permission.

# Cyclic Learning Rate



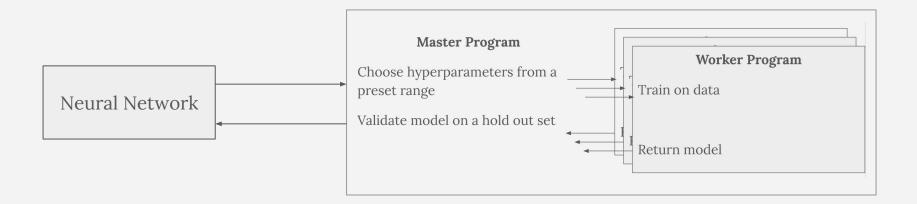
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### Automatic Model Search

The model search process is really another optimization problem.

Maybe we can train a neural network for that!



### Automatic Model Search

### Cloud AutoML<sup>BETA</sup>

Train high-quality custom machine learning models with minimum effort and machine learning expertise.

Try AutoML Translation, Natural Language, and Vision, now in beta.



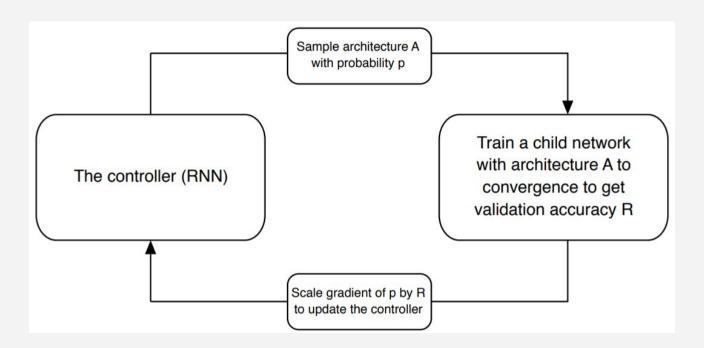
**CONTACT SALES** 

### Train custom machine learning models

Cloud AutoML is a suite of machine learning products that enables developers with limited machine learning expertise to train high-quality models specific to their business needs, by leveraging Google's state-of-the-art transfer learning, and Neural Architecture Search technology.



### Neural Architecture Search

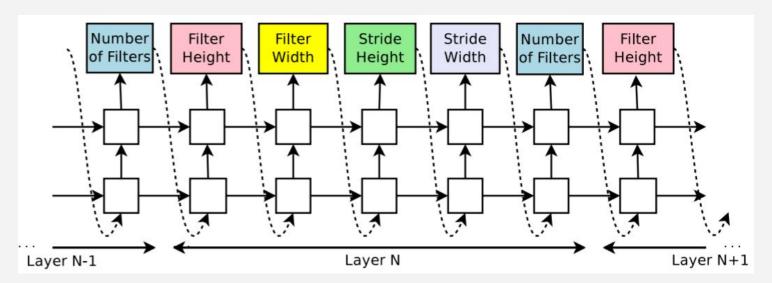


https://arxiv.org/abs/1611.01578 https://arxiv.org/pdf/1707.07012.pdf

### Generate Model Descriptions

A Controller Recurrent Neural Network

At each layer, it predicts hyperparameters to construct a network unit.



# Reinforcement Learning

### Reward signal:

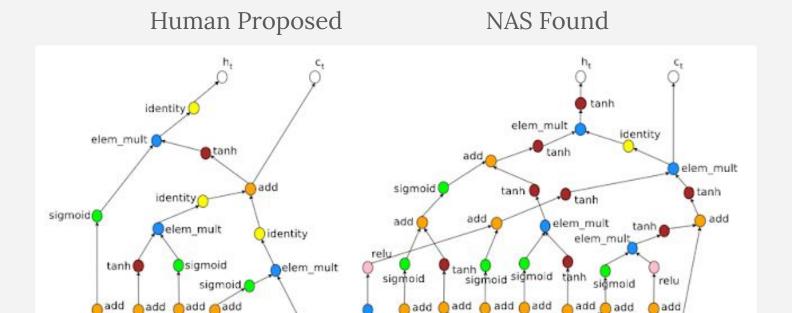
Validation accuracy

Why reinforcement learning in this task?

Supervised learning is not suitable because it doesn't make sense to set an exact target model structure: we are searching for one!

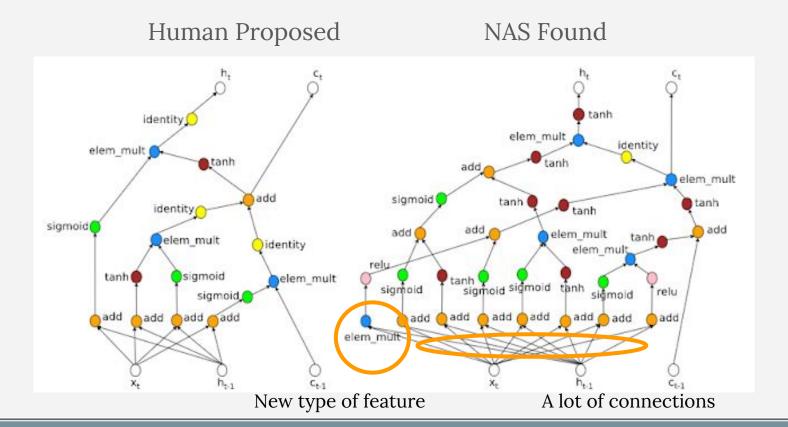
You want to optimize the model structure given some high level 'goodness' measure.

## Example of Found Structures



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## Example of Found Structures



### On Multiplicative Integration with Recurrent Neural Networks

Yuhuai Wu<sup>1,\*</sup>, Saizheng Zhang<sup>2,\*</sup>, Ying Zhang<sup>2</sup>, Yoshua Bengio<sup>2,4</sup> and Ruslan Salakhutdinov<sup>3,4</sup>
<sup>1</sup>University of Toronto, <sup>2</sup>MILA, Université de Montréal, <sup>3</sup>Carnegie Mellon University, <sup>4</sup>CIFAR
ywu@cs.toronto.edu, <sup>2</sup>{firstname.lastname}@umontreal.ca,rsalakhu@cs.cmu.edu

#### Abstract

We introduce a general and simple structural design called "Multiplicative Integration" (MI) to improve recurrent neural networks (RNNs). MI changes the way in which information from difference sources flows and is integrated in the computational building block of an RNN, while introducing almost no extra parameters. The new structure can be easily embedded into many popular RNN models, including LSTMs and GRUs. We empirically analyze its learning behaviour and conduct evaluations on several tasks using different RNN models. Our experimental results demonstrate that Multiplicative Integration can provide a substantial performance boost over many of the existing RNN models.

### On Multiplicative Integration with Recurrent Neural Networks

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The key idea behind Multiplicative Integration is to integrate different information flows  $\mathbf{W}x$  and  $\mathbf{U}z$ , by the Hadamard product " $\odot$ ". A more general formulation of Multiplicative Integration includes two more bias vectors  $\beta_1$  and  $\beta_2$  added to  $\mathbf{W}x$  and  $\mathbf{U}z$ :

$$\phi((\mathbf{W}x + \beta_1) \odot (\mathbf{U}z + \beta_2) + b) \tag{3}$$

We introduce a general and simple structural design called "Multiplicative Integration" (MI) to improve recurrent neural networks (RNNs). MI changes the way in which information from difference sources flows and is integrated in the computational building block of an RNN, while introducing almost no extra parameters. The new structure can be easily embedded into many popular RNN models, including LSTMs and GRUs. We empirically analyze its learning behaviour and conduct evaluations on several tasks using different RNN models. Our experimental results demonstrate that Multiplicative Integration can provide a substantial performance boost over many of the existing RNN models.

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### We will cover more on how to use Google AutoML later in the semester

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### Summary

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