



DARTMOUTH



# Randomization & Approximation in Machine Learning and Neural Network Pruning

Math 76 Final Project

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## Basic Randomization & Approximation in ML

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# Randomization & Approximation as we know it

The randomization techniques from class mostly involve **splitting** or **modifying data** randomly:

- splitting data into training and testing subsets

```
1 X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)
```

- bagging/bootstrapping procedures
- random parameter/weight initialization in neural networks

Approximation is of course the goal of machine learning, but we often use *exact* algorithms in ML models instead of approximation algorithms

# What else can Randomization & Approximation be used for

- Optimization:
  - trade "accuracy" for speed
  - oftentimes algorithms have  $\geq 90\%$  accuracy
- In neural networks
  - "Prune" connections (somewhat) at random to get a smaller model with higher accuracy
- Sometimes, allowing some error is more "realistic" leading to a higher accuracy
  - higher accuracy's could also be because of a reduction in over fitting

## Example: k Approximate Nearest Neighbors

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# Naive Solution

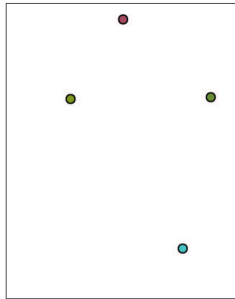
**Problem:** Given a set of vectors  $S$  in a data space and a similarity measure  $\text{DIST}$  (for example euclidean distance) find the  $k$  nearest ("most similar") neighbors to some input  $x$ .

**Brute force solution:** compare to every other point

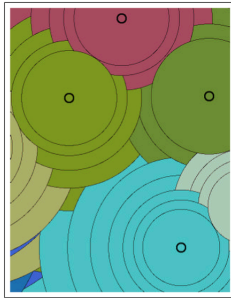
```
1: procedure BRUTE-KNN( $x, S, \text{DIST}, k = 1$ )  
2:   Compute  $\text{DIST}(x, s) \forall s \in S$   
3:   return  $k$  points ( $s \in S$ ) with smallest  $\text{DIST}$  values
```

**Issue:** expensive and **redundant** (especially in higher dimensions and when there is lots of training data);  $O(n)$  lookup

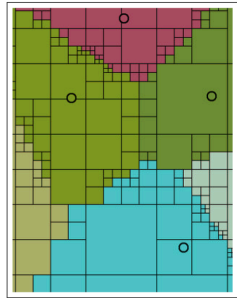
# k Approximate Nearest Neighbors (kANN)



(A)



(B)



(C)

A) data points; B) & C) approximate nearest neighbor [3]



# Locality Sensitive Hashing (LSH)

## Hashing:

- hash function:  $f: U \rightarrow [n]$  where  $|U| \gg n$  and  $n$  is (usually) prime
- typically, two inputs have a very low probability of "colliding" (their output is the same)
  - inputs will have different "hashes" even if they are very similar

## Locality Sensitive Hashing (LSH)

- two inputs which are similar have a high probability of colliding [3]
- can be used to solve kANN with high accuracy and in less time

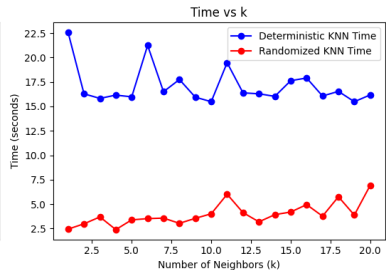
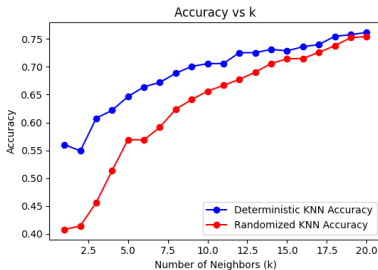
# Approximate-KNN Experiment

- using PyNNdescent package for kANN implementation [5]
  - uses modern kANN algorithm [1], written in python
  - accuracy converges to 90%
- Naive brute force implementation hand written in python
  - did not use an implementation from scikitlearn because compiled in C
  - best to have implementation we are comparing all in the same language

# Results

```
# Generate a synthetic dataset with 10,000 samples and 1,000 features
```

```
X, y = make_classification(n_samples=10000, n_features=200, n_informative=50, n_classes=3,  
    random_state=42)
```

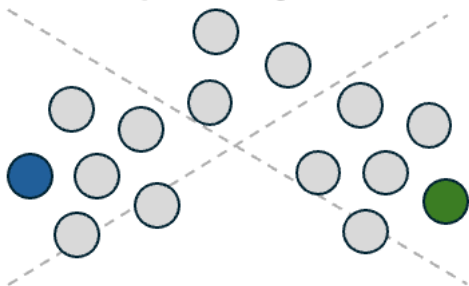


kNN vs kANN **raw** accuracy and time comparison

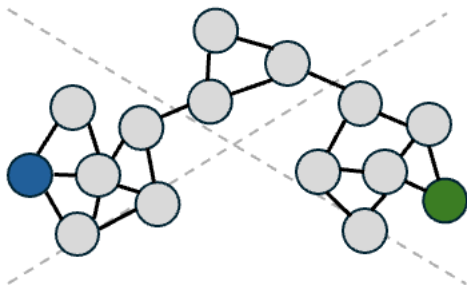
# Graph-Based Semi-Supervised Learning

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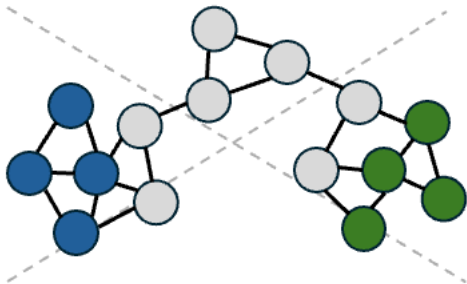
**1) Start with data in feature space;  
small percentage is labeled**



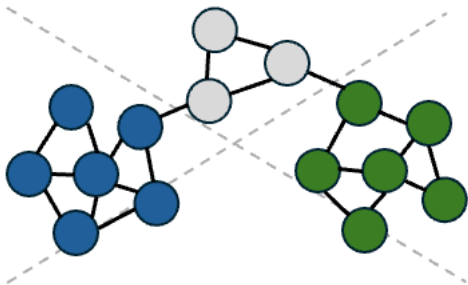
## 2) Construct $k$ nearest neighbor Graph (here $k=2$ )



## 3) "Propagate Labels"

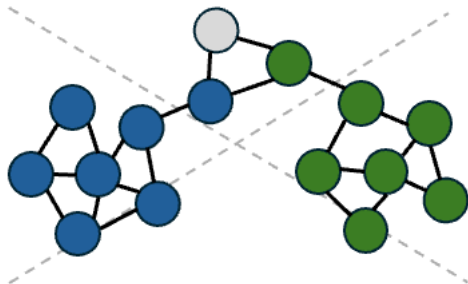


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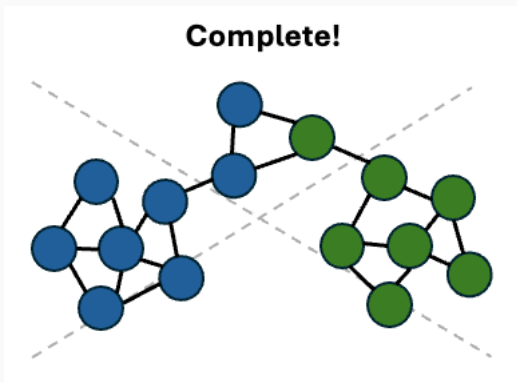




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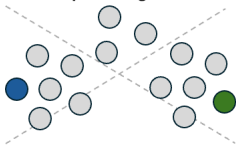


# Graph-Based Semi-Supervised Learning

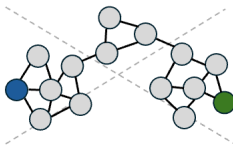


# Graph-Based Semi-Supervised Learning

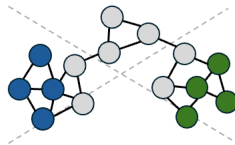
1) Start with data in feature space;  
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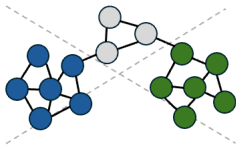
2) Construct  $k$  nearest neighbor  
Graph (here  $k=2$ )



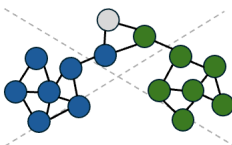
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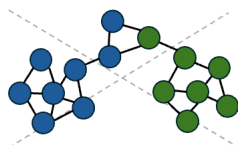
3) "Propagate Labels"



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Complete!



# Graph-Based Semi-Supervised Learning

- Constructing the  $k$  Nearest Neighbor Graph dominates runtime
- Most Label Propagation Algorithms (LPA) are **simple** and **near linear** in time

## Parameters:

- $y :=$  initial labels;  $y_i :=$  label of node  $i$ ;  $-1$  if unlabeled
- $W :=$  adjacency matrix of nearest neighbors graph
- $\alpha \in [0, 1] :=$  "clamping factor" indicating influence of initial labels **per iteration**

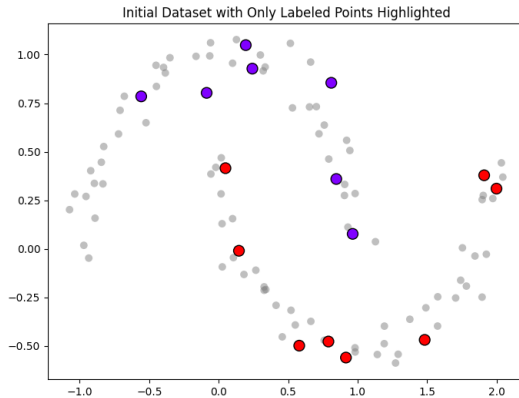
## Update Rules:

$$\bullet F_{ic}^{(0)} \leftarrow \begin{cases} 1 & \text{if } y_i = c \\ 0 & \text{otherwise} \end{cases} \quad F_{ic} \text{ is the } \mathbf{confidence} \text{ of node } i \text{ having label } c$$

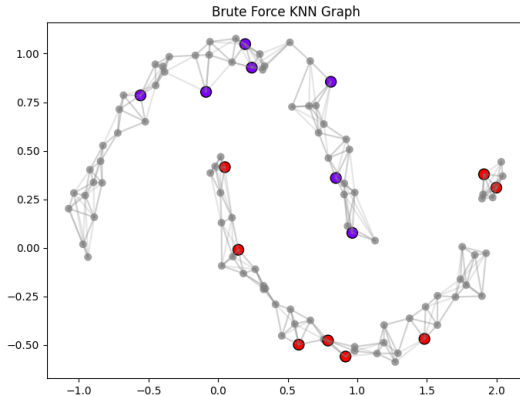
$$\bullet F_{ic}^{(t+1)} \leftarrow \underbrace{\alpha W F^{(t)}}_{\text{propagate from neighbors}} + \underbrace{(1 - \alpha) F^{(0)}}_{\text{account for initial labels}}$$

$$\bullet y_i^* \leftarrow \arg \max_c F_{ic}^{(t_{\max})} \quad y_i^* \text{ is the vector of } \mathbf{final} \text{ labels}$$

# Results I

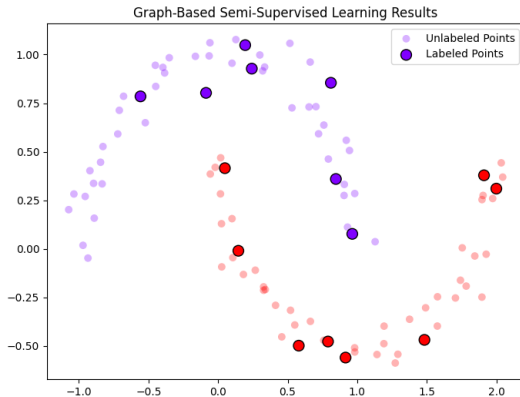


# Results I



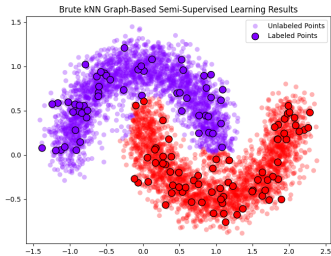
Nearest Neighbor graph with  $k = 5$

# Results I

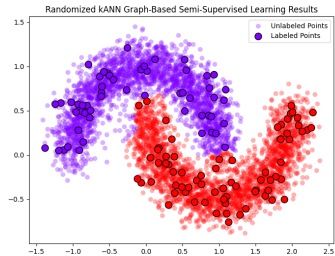


Final labels using  $\alpha = .99$

# Results II



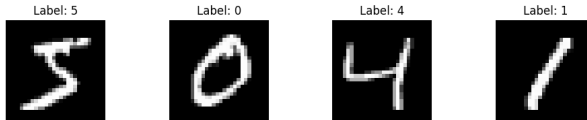
- Time: 4.20 seconds
- Accuracy: 0.976



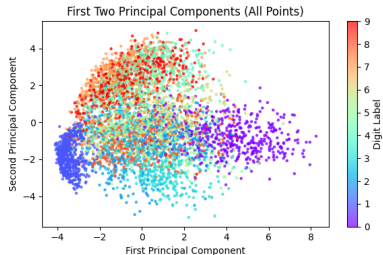
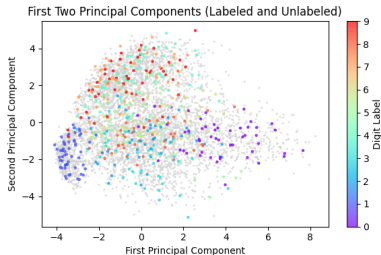
- Time: 2.89 seconds
- Accuracy: 0.977



# MNIST Dataset Preprocessing



- Normalized signal to take value  $\in [0, 1]$
- PCA to extract top 50 principle components
- masked 90% of labels to simulate semi-supervised learning; kept full label set for accuracy testing



# MNIST Dataset Results


\* Ran with A100 GPU runtime on Google Colab

\* Parameters:

## Using Brute Force kNN:

- Accuracy: 0.933
- Total Time: 18.81 minutes
- Time to construct kNN-graph: 17.14 minutes

## Using k Approximate NN:

- Accuracy: 0.955
- Time: 1.79 minutes
- Time to construct kANN-graph: 

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**Approximate NN-graph construction is over 165 times faster than the naive**

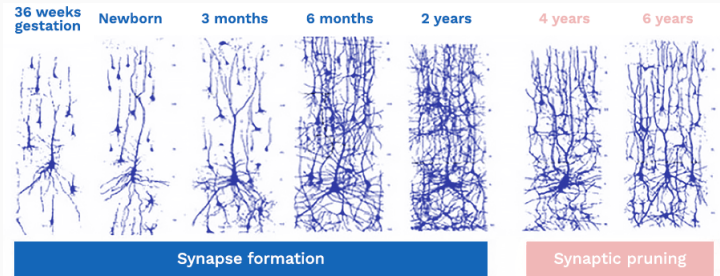
# Randomized Pruning in Neural Networks

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# Why Prune?

- "It is widely acknowledged that large and sparse models have higher accuracy than small and dense models" [4]
- pruning a model prevents over-fitting
- pruned models also require a fraction of the space a unpruned model would need in memory

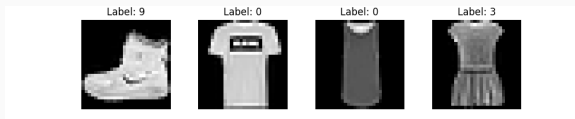
# Neurology Inspiration



Synaptic pruning in the brain [2]

# Dataset and Neural Network Architecture

- We use a simple architecture recommended by tensorflow
- We prune connections on one layer (more detail on next slide)
- pruning is done by setting pruned weights (outgoing to next layer) to zero
- Inspiration for randomized pruning: "Breaking through Deterministic Barriers: Randomized Pruning Mask Generation and Selection" [4]



MNIST Fashion Dataset



# RESULTS

Network	Pruning Strategy	Accuracy
No Pruning	No pruning is applied. All connections are retained.	87.68%
Threshold Fraction Pruning	Prune a specified fraction of connections with the smallest weights.	86.63%
Random Pruning	Prune a random subset of connections, regardless of their weights or importance.	64.56%
Importance-Weighted Random Pruning	Prune connections probabilistically based on their importance, where more important connections are less likely to be pruned.	78.06%
Fractional Random Threshold Pruning	First, select a fraction of connections with the smallest weights. Then, randomly prune a smaller fraction of these selected connections.	87.47%

Summary of Pruning Strategies for Five Neural Networks

Questions???

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W. Dong, M. Charikar, and K. Li.

**Efficient k-nearest neighbor graph construction for generic similarity measures.**

In [Proceedings of the 20th international conference on World wide web](#), pages 577–586. ACM, 2011.



J. Embrace.

**Synaptic growth, synesthesia, and savant abilities.**

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S. Har-Peled, P. Indyk, and R. Motwani.

**Approximate nearest neighbor: Towards removing the curse of dimensionality.**

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Special issue in honor of Rajeev Motwani.



J. Li, W. Gao, Q. Lei, and D. Xu.

**Breaking through deterministic barriers: Randomized pruning mask generation and selection.**

[arXiv preprint arXiv:2310.13183](#), 2023.

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