10-605/805 – ML for Large Datasets Lecture 5: Distributed Linear Regression

Front Matter

- HW1 released 8/30, due 9/13 (today!) at 11:59 PM
 - Recitation 3 on 9/16 will go over HW1 solutions
- HW2 released 9/8, due 9/22 at 11:59 PM
- Mini-project details released 9/9

Background: Big *O* Notation

- Used to describe an algorithm's time or space (storage)
 complexity in terms of the input size
- Formally:

$$f(x) = O(g(x)) \Leftrightarrow \exists C, x_0 \text{ s.t. } f(x) \leq Cg(x) \ \forall \ x \geq x_0$$

- O(1) = constant time/space, i.e., a fixed number of operations or storage regardless of input
- $O(\log(n)) = \log(n)$
- O(n) = linear time/space
- An algorithm's time and space complexity can be different
 - Example: multiplying an $a \times b$ matrix with an $b \times c$ matrix takes O(abc) time (ac dot products between b-length vectors) but the result uses O(ac) storage

Background: Empirical Risk Minimization

- A common framework for supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$
 - a loss function ℓ : $\mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$
 - $oldsymbol{\cdot}$ a hypothesis class or set of functions ${\mathcal F}$

the goal is to find

$$\hat{f} = \underset{f \in \mathcal{F}}{\operatorname{argmin}} \sum_{i=1}^{n} \ell(f(\mathbf{x}^{(i)}), y^{(i)})$$

with the hope that

$$\mathbb{E}_{p(\mathbf{x},y)}[\ell(f(\mathbf{x}),y)] \approx \sum_{i=1}^{n} \ell(f(\mathbf{x}^{(i)}),y^{(i)})$$

Background: Empirical Risk Minimization

- A common framework for supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(\mathbf{x}^{(i)}, \mathbf{y}^{(i)})\}_{i=1}^n$
 - a loss function ℓ : $\mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$
 - $oldsymbol{\cdot}$ a hypothesis class or set of functions ${\mathcal F}$

the goal is to find

$$\hat{f} = \underset{f \in \mathcal{F}}{\operatorname{argmin}} \sum_{i=1}^{n} \ell(f(\mathbf{x}^{(i)}), y^{(i)})$$

- Depending on the choice of \mathcal{F} and ℓ , this objective function may be convex (easy to optimize) or non-convex (hard)
- Our focus will be solving this problem for large n and/or k

Background: Regression

- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$
 - a loss function $\ell: \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$ where $\mathcal{Y} = \mathbb{R}$
 - $oldsymbol{\cdot}$ a hypothesis class or set of functions $oldsymbol{\mathcal{F}}$

the goal is to find

$$\hat{f} = \underset{f \in \mathcal{F}}{\operatorname{argmin}} \sum_{i=1}^{n} \ell(f(\mathbf{x}^{(i)}), y^{(i)})$$

 Fun example: predicting the year a song was released based on (a representation of) its audio (see HW2)

- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$

$$\ell(y,y') = (y-y')^2$$

• \mathcal{F} = all functions of the form $f(x) = w_0 + \sum_{d=1}^{\infty} w_d x_d$

the goal is to find

$$\hat{f} = \underset{f \in \mathcal{F}}{\operatorname{argmin}} \sum_{i=1}^{n} (f(\mathbf{x}^{(i)}) - y^{(i)})^{2}$$

- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$
 - $\ell(y,y') = (y-y')^2$
 - \mathcal{F} = all functions of the form $f(\mathbf{x}) = \mathbf{w}^T [1 \ \mathbf{x}^T]^T$

the goal is to find

$$\hat{f} = \underset{f \in \mathcal{F}}{\operatorname{argmin}} \sum_{i=1}^{n} (f(\mathbf{x}^{(i)}) - y^{(i)})^{2}$$

- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$

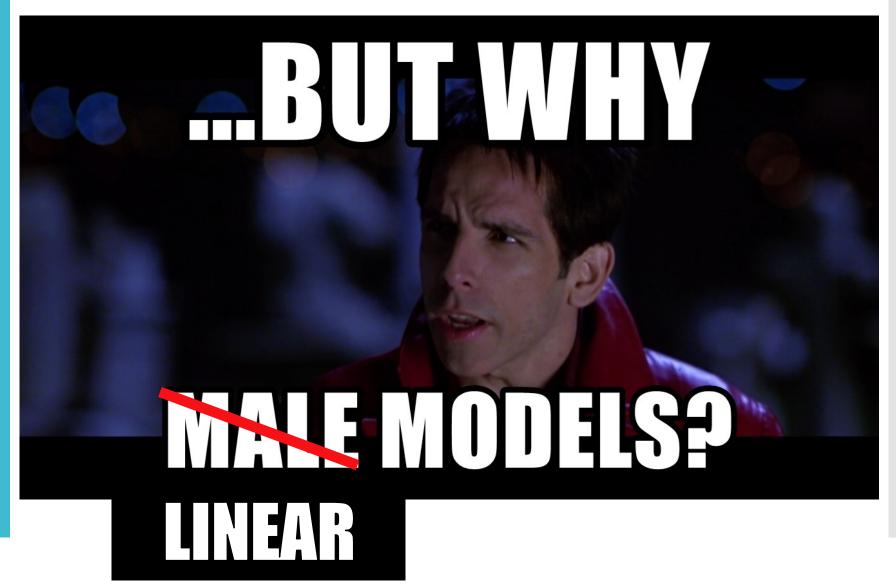
$$\cdot \ell(y, y') = (y - y')^2$$

• \mathcal{F} = all functions of the form $f(x) = \mathbf{w}^T x$ 1 implicitly

the goal is to find

prepended

$$\widehat{\boldsymbol{w}} = \underset{\boldsymbol{w}}{\operatorname{argmin}} \sum_{i=1}^{n} (\boldsymbol{w}^{T} \boldsymbol{x}^{(i)} - \boldsymbol{y}^{(i)})^{2}$$



- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$

$$\ell(y, y') = (y - y')^2$$

• \mathcal{F} = all functions of the form $f(x) = \mathbf{w}^T x$ 1 implicitly

the goal is to find

prepended

$$\widehat{\boldsymbol{w}} = \underset{\boldsymbol{w}}{\operatorname{argmin}} \sum_{i=1}^{n} (\boldsymbol{w}^{T} \boldsymbol{x}^{(i)} - \boldsymbol{y}^{(i)})^{2}$$

- A type of supervised learning
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$

$$\ell(y,y') = (y-y')^2$$

• \mathcal{F} = all functions of the form $f(x) = w^T x$ the goal is to find $f(x) = w^T x$

$$\widehat{\mathbf{w}} = \underset{\mathbf{w}}{\operatorname{argmin}} (X\mathbf{w} - \mathbf{y})^{T} (X\mathbf{w} - \mathbf{y})$$

$$\widehat{\mathbf{w}} = \underset{\mathbf{w}}{\operatorname{argmin}} \|X\mathbf{w} - \mathbf{y}\|_{2}^{2}$$

• where
$$X = \begin{bmatrix} x^{(1)}^T \\ \vdots \\ x^{(n)}^T \end{bmatrix}$$
 and $y = \begin{bmatrix} y^{(1)} \\ \vdots \\ y^{(n)} \end{bmatrix}$

$$L_{\mathcal{D}}(\mathbf{w}) = (X\mathbf{w} - \mathbf{y})^T (X\mathbf{w} - \mathbf{y})$$

Background: Regularization

- A modification to empirical risk minimization that penalizes model complexity in order to combat overfitting
- Given:
 - some labelled training dataset $\mathcal{D} = \{(\mathbf{x}^{(i)}, \mathbf{y}^{(i)})\}_{i=1}^n$
 - a loss function $\ell: \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$
 - $oldsymbol{\cdot}$ a hypothesis class or set of functions $oldsymbol{\mathcal{F}}$
 - a regularizer $R: \mathcal{W} \to \mathbb{R}$
 - a coefficient of regularization λ

the goal is to find

$$\widehat{\boldsymbol{w}} = \underset{\boldsymbol{w}}{\operatorname{argmin}} \sum_{i=1}^{n} \ell(f_{\boldsymbol{w}}(\boldsymbol{x}^{(i)}), y^{(i)}) + \lambda R(\boldsymbol{w})$$

Background: Ridge Regression

- A modification to empirical risk minimization that penalizes model complexity in order to combat overfitting
- Given:
 - some labelled training dataset $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n$
 - $\cdot \ell(y, y') = (y y')^2$
 - \mathcal{F} = all functions of the form $f(x) = \mathbf{w}^T x$
 - $R(w) = ||w||_2^2 = w^T w$
 - a coefficient of regularization λ

the goal is to find

$$\widehat{\mathbf{w}} = \underset{\mathbf{w}}{\operatorname{argmin}} (X\mathbf{w} - \mathbf{y})^T (X\mathbf{w} - \mathbf{y}) + \lambda \mathbf{w}^T \mathbf{w}$$

$$L_{\mathcal{D}}(\mathbf{w}) = (X\mathbf{w} - \mathbf{y})^T (X\mathbf{w} - \mathbf{y}) + \lambda \mathbf{w}^T \mathbf{w}$$

Background: Ridge Regression

$$\rightarrow \widehat{\boldsymbol{w}} = (X^T X + \lambda I_k)^{-1} X^T \boldsymbol{y}$$

where I_k is the $k \times k$ identity matrix

$$L_{\mathcal{D}}(\mathbf{w}) = (X\mathbf{w} - \mathbf{y})^T (X\mathbf{w} - \mathbf{y}) + \lambda \mathbf{w}^T \mathbf{w}$$

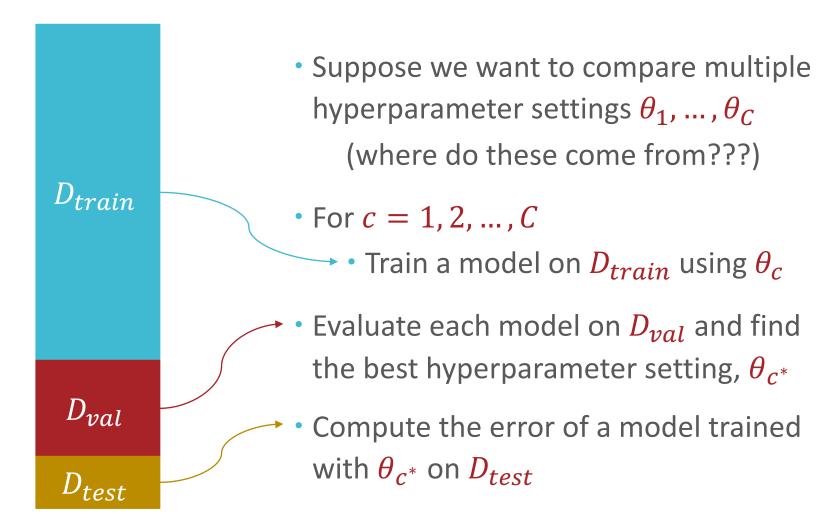
Aside: How can we set λ ?

$$\rightarrow \widehat{\boldsymbol{w}} = (X^T X + \lambda I_k)^{-1} X^T \boldsymbol{y}$$

where I_k is the $k \times k$ identity matrix

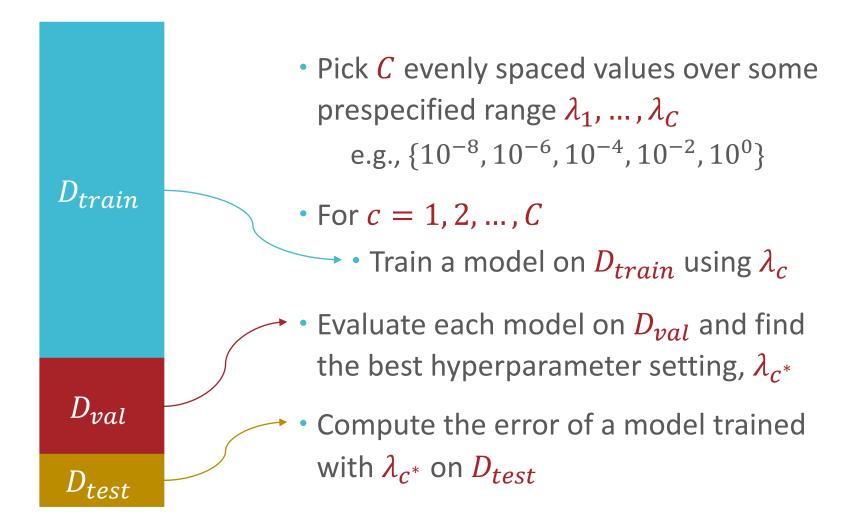
Recall: Machine Learning Pipeline

Hyperparameter optimization



HW2 Preview: Grid Search

Hyperparameter optimization



Linear Regression: Computational Cost

$$\widehat{\boldsymbol{w}} = (X^T X)^{-1} X^T \boldsymbol{y}$$

- 1. Does this quantity exist, i.e., is X^TX invertible? When $n \gg k + 1$, X^TX is (almost always) full rank and; otherwise, we can compute a *pseudoinverse* via singular value decomposition
- 2. If so, how expensive is it to compute?
 - Computing X^TX : $O(nk^2)$ time & $O(k^2)$ space
 - Inverting X^TX : $O(k^3)$ time & $O(k^2)$ space
 - Storing X: O(nk) space

Key takeaway: computational bottlenecks will change based on the relationship between n and k

Linear Regression: Large n, Small k

• Assume $O(k^3)$ computation and $O(k^2)$ storage is possible on a single machine

 \checkmark We can store and invert X^TX

We cannot compute X^TX

We cannot store X

- Idea: distribute storage of X and computation of X^TX
 - 1. Store the rows of *X* across different machines
 - 2. Compute X^TX as the sum of outer products

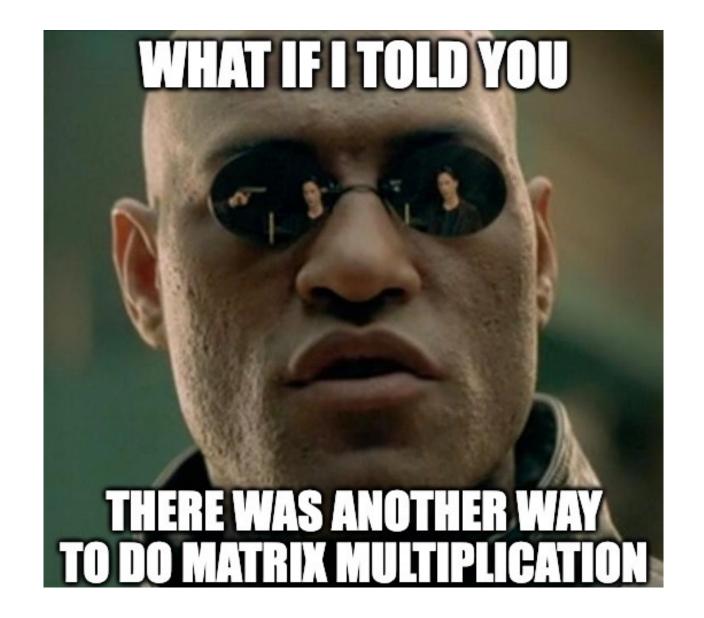
$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} \\ \end{bmatrix}$$

$$1*2+4*2+5*3=25$$

$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} 25 & 2 \\ 1 & 4 \end{bmatrix}$$

$$1*2+4*1+5*4=26$$

$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} 25 & 26 \\ 17 & 19 \end{bmatrix}$$



$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} \\ \end{bmatrix}$$

$$\begin{bmatrix} 2 & 2 \\ 6 & 6 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} 1 & 1 & 1 \\ 3 & 4 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 2 & 2 \\ 6 & 6 \end{bmatrix} + \begin{bmatrix} 8 & 4 \\ 2 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} \\ \end{bmatrix}$$

$$\begin{bmatrix} 2 & 2 \\ 6 & 6 \end{bmatrix} + \begin{bmatrix} 8 & 4 \\ 2 & 1 \end{bmatrix} + \begin{bmatrix} 15 & 20 \\ 9 & 12 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 4 & 5 \\ 3 & 1 & 3 \end{bmatrix} \begin{bmatrix} 2 & 2 \\ 2 & 1 \\ 3 & 4 \end{bmatrix} = \begin{bmatrix} 25 & 26 \\ 17 & 19 \end{bmatrix}$$

$$\begin{bmatrix} 2 & 2 \\ 6 & 6 \end{bmatrix} + \begin{bmatrix} 8 & 4 \\ 2 & 1 \end{bmatrix} + \begin{bmatrix} 15 & 20 \\ 9 & 12 \end{bmatrix}$$

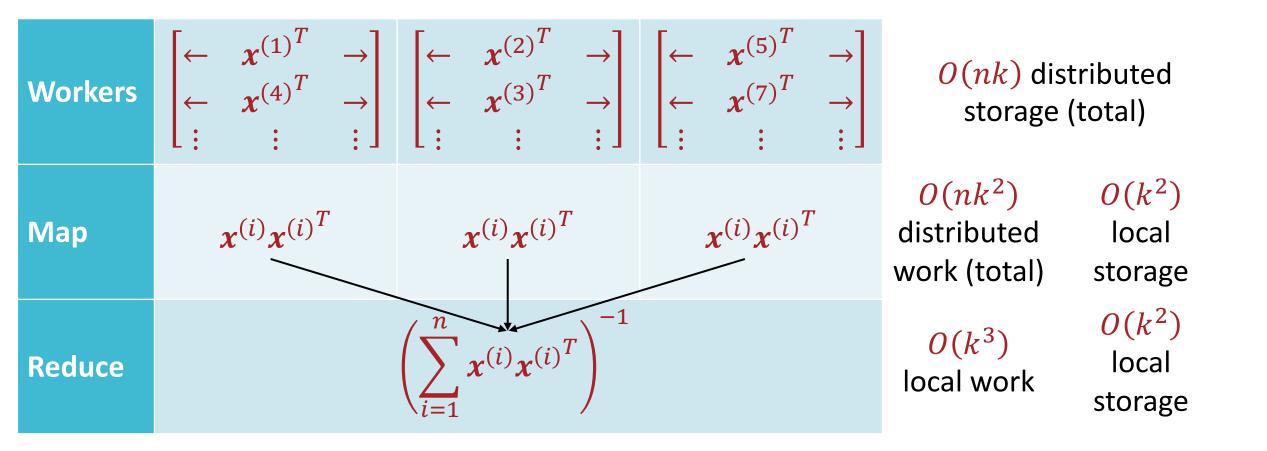
$$\begin{bmatrix} A_{11} & A_{12} & \cdots & A_{1m} \\ A_{21} & A_{22} & \cdots & A_{2m} \\ \vdots & \vdots & \ddots & \vdots \\ A_{n1} & A_{n2} & \cdots & A_{nm} \end{bmatrix} \begin{bmatrix} B_{11} & \cdots & B_{1k} \\ B_{21} & \cdots & B_{2k} \\ B_{31} & \cdots & B_{3k} \\ \vdots & \ddots & \vdots \\ B_{m1} & \cdots & B_{mk} \end{bmatrix}$$

$$= \begin{bmatrix} \sum_{i=1}^{m} A_{1i} B_{i1} & \cdots & \sum_{i=1}^{m} A_{1i} B_{ik} \\ \vdots & \ddots & \vdots \\ \sum_{i=1}^{m} A_{ni} B_{i1} & \cdots & \sum_{i=1}^{m} A_{ni} B_{ik} \end{bmatrix} = \sum_{i=1}^{m} \begin{bmatrix} A_{1i} B_{i1} & \cdots & A_{1i} B_{ik} \\ \vdots & \ddots & \vdots \\ A_{ni} B_{i1} & \cdots & A_{ni} B_{ik} \end{bmatrix}$$

Distributed Computation of $(X^TX)^{-1}$

$$X^{T}X = \begin{bmatrix} \uparrow & \uparrow & \cdots & \uparrow \\ \mathbf{x}^{(1)} & \mathbf{x}^{(2)} & \cdots & \mathbf{x}^{(n)} \end{bmatrix} \begin{bmatrix} \leftarrow & \mathbf{x}^{(1)^{T}} & \rightarrow \\ \leftarrow & \mathbf{x}^{(2)^{T}} & \rightarrow \\ \vdots & \vdots & \vdots \\ \leftarrow & \mathbf{x}^{(n)^{T}} & \rightarrow \end{bmatrix} = \sum_{i=1}^{n} \mathbf{x}^{(i)} \mathbf{x}^{(i)^{T}}$$

• Idea: distribute $x^{(i)}$ and compute summands in parallel



Distributed Computation of $(X^TX)^{-1}$

Workers
$$\begin{bmatrix} \leftarrow & \boldsymbol{x^{(1)}}^T & \rightarrow \\ \leftarrow & \boldsymbol{x^{(4)}}^T & \rightarrow \\ \vdots & \vdots & \vdots \end{bmatrix} \begin{bmatrix} \leftarrow & \boldsymbol{x^{(2)}}^T & \rightarrow \\ \leftarrow & \boldsymbol{x^{(3)}}^T & \rightarrow \\ \vdots & \vdots & \vdots \end{bmatrix} \begin{bmatrix} \leftarrow & \boldsymbol{x^{(5)}}^T & \rightarrow \\ \leftarrow & \boldsymbol{x^{(7)}}^T & \rightarrow \\ \vdots & \vdots & \vdots \end{bmatrix}$$

O(nk) distributed storage (total)

trainData.map(compute outer prods)

 $O(nk^2)$ $O(k^2)$ distributed local work (total) storage

trainData.reduce(sum and invert)

 $O(k^{3})$ local work

 $O(k^2)$ local storage

Distributed Computation of $(X^TX)^{-1}$

Linear Regression: Large n, Large k

• Now, $O(k^3)$ computation and $O(k^2)$ storage is *not* possible on a single machine

We cannot store and invert X^TX

We cannot compute X^TX

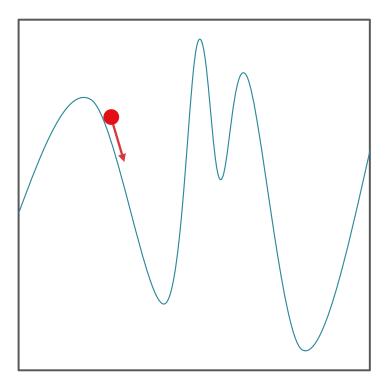
We cannot store X

10-605/805 Principle #1: computation and storage should be at most linear in n and k

Idea: use a different algorithm!

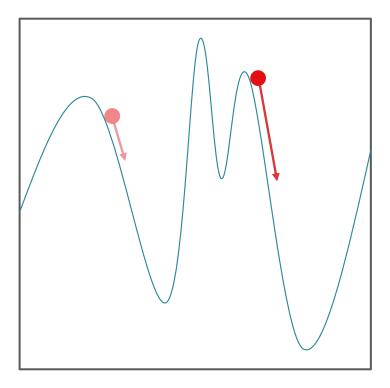
Background: Gradient Descent

- An iterative method for minimizing functions
- Requires the gradient to exist everywhere

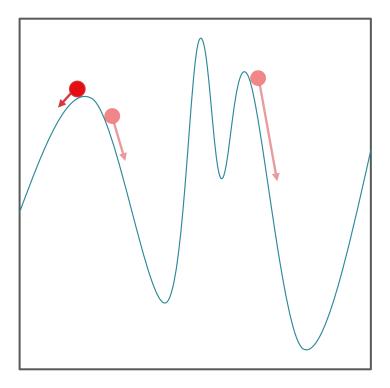


Background: Gradient Descent

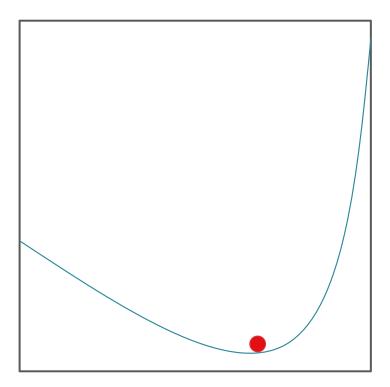
- An iterative method for minimizing functions
- Requires the gradient to exist everywhere



- An iterative method for minimizing functions
- Requires the gradient to exist everywhere



- An iterative method for minimizing functions
- Requires the gradient to exist everywhere



 Good news: the linear regression objective is convex so gradient descent will always converge to the global minimum

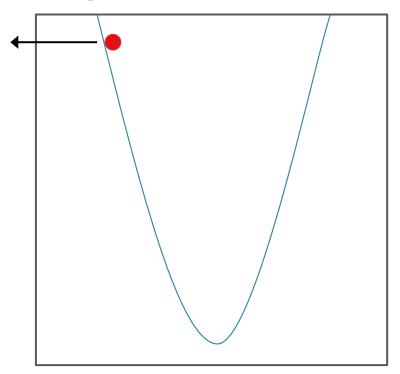
- Suppose we're trying to minimize some function L and we're currently at some location $oldsymbol{w}^{(t)}$
- Move some distance, α , in the "most downhill" direction, \boldsymbol{v} :

$$\boldsymbol{w}^{(t+1)} = \boldsymbol{w}^{(t)} + \alpha \boldsymbol{v}$$

- The gradient points in the direction of steepest increase ...
- ... so let's move in the opposite direction!

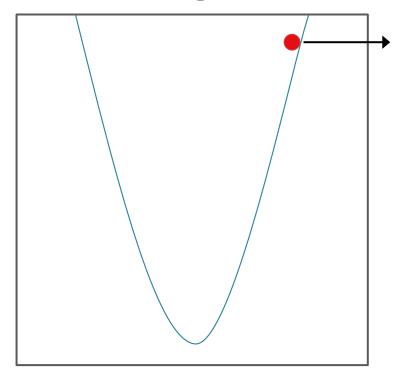
$$\mathbf{w}^{(t+1)} = \mathbf{w}^{(t)} - \eta \nabla_{\mathbf{w}} L(\mathbf{w}^{(t)})$$

Direction of gradient



$$2x < 0$$
 for $x < 0$

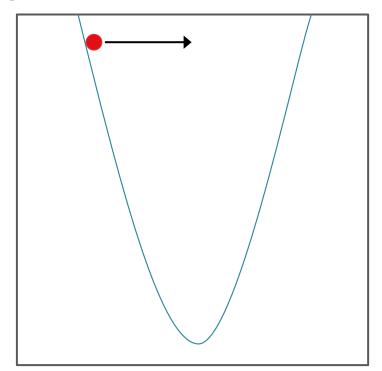
Direction of gradient



$$2x > 0$$
 for $x > 0$

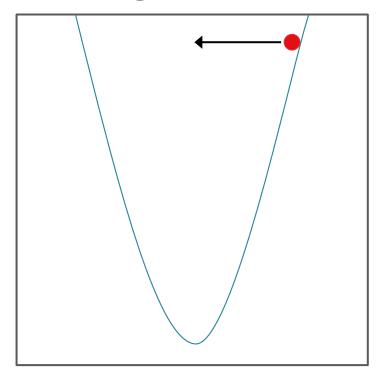
$$\frac{\partial}{\partial x}x^2 = 2x$$

Direction of global minimum



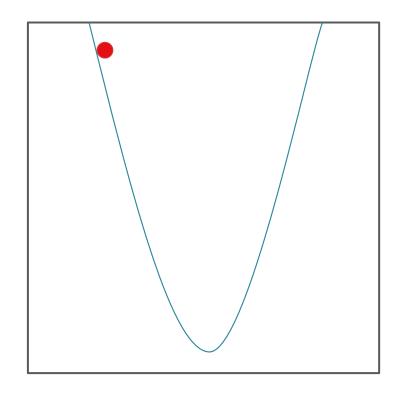
$$2x < 0$$
 for $x < 0$

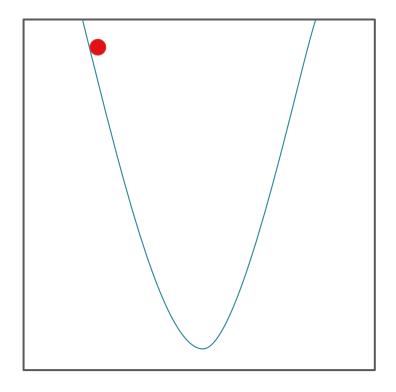
Direction of global minimum



$$2x > 0$$
 for $x > 0$

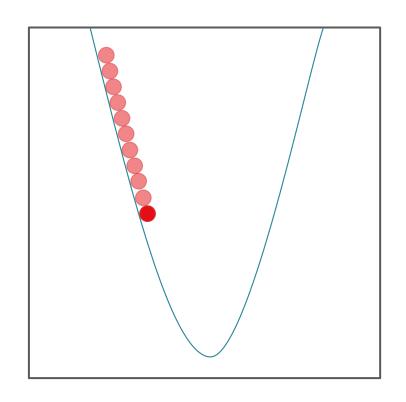
$$\frac{\partial}{\partial x}x^2 = 2x$$

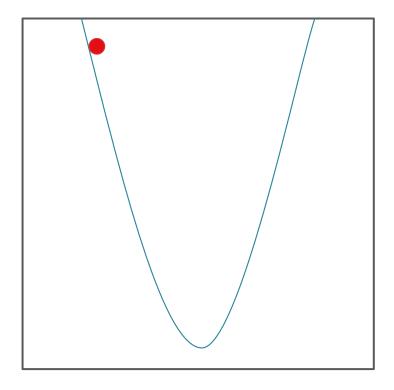




Small α

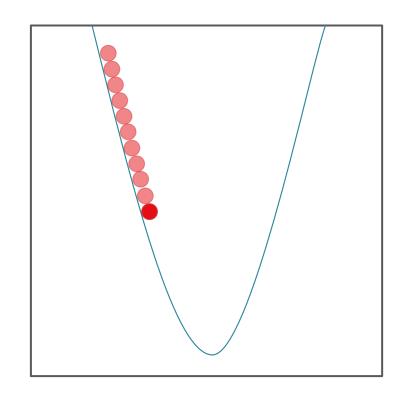
Large α

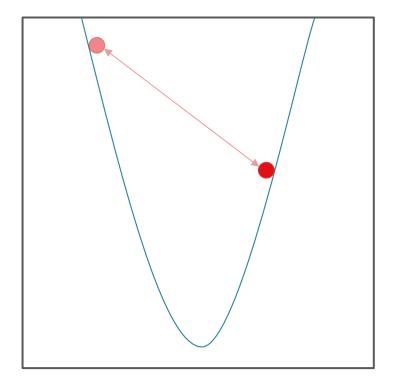




Small α

Large α

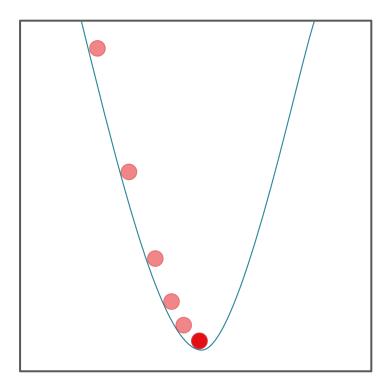




Small α

Large α

• Use a variable $\alpha^{(t)}$ instead of a fixed $\alpha!$



• Example: $\alpha^{(t)} = \frac{\alpha}{n\sqrt{t}}$

Gradient Descent for Linear Regression

• Input:
$$\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n, \alpha$$

- 1. Initialize $\mathbf{w}^{(0)}$ to all zeros and set t=0
- While TERMINATION CRITERION is not satisfied
 - a. Compute the gradient:

$$\nabla_{\boldsymbol{w}} L_{\mathcal{D}}(\boldsymbol{w}^{(t)}) = (2X^T X \boldsymbol{w}^{(t)} - 2X^T \boldsymbol{y})$$

b. Update the weights:

$$\boldsymbol{w}^{(t+1)} \leftarrow \boldsymbol{w}^{(t)} - \frac{\alpha}{n\sqrt{t}} \left(2X^T X \boldsymbol{w}^{(t)} - 2X^T \boldsymbol{y} \right)$$

c. Increment $t: t \leftarrow t + 1$

• Output: $\mathbf{w}^{(t)}$

Gradient Descent for Linear Regression

- Input: $\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n, \alpha$
- 1. Initialize $\mathbf{w}^{(0)}$ to all zeros and set t=0
- 2. While TERMINATION CRITERION is not satisfied
 - a. Compute the gradient:

$$\nabla_{\boldsymbol{w}} L_{\mathcal{D}}(\boldsymbol{w}^{(t)}) = 2 \sum_{i=1}^{n} \left(\boldsymbol{w}^{(t)^{T}} \boldsymbol{x}^{(i)} - \boldsymbol{y}^{(i)} \right) \boldsymbol{x}^{(i)}$$

b. Update the weights:

$$\mathbf{w}^{(t+1)} \leftarrow \mathbf{w}^{(t)} - \frac{\alpha}{n\sqrt{t}} \sum_{i=1}^{n} \left(\mathbf{w}^{(t)^{T}} \mathbf{x}^{(i)} - \mathbf{y}^{(i)} \right) \mathbf{x}^{(i)}$$

c. Increment $t: t \leftarrow t + 1$

• Output: $\mathbf{w}^{(t)}$

Idea: distribute $x^{(i)}$ and compute summands in parallel

• Input:
$$\mathcal{D} = \{(x^{(i)}, y^{(i)})\}_{i=1}^n, \alpha$$

- 1. Initialize $\mathbf{w}^{(0)}$ to all zeros and set t=0
- 2. While TERMINATION CRITERION is not satisfied
 - a. Compute the gradient:

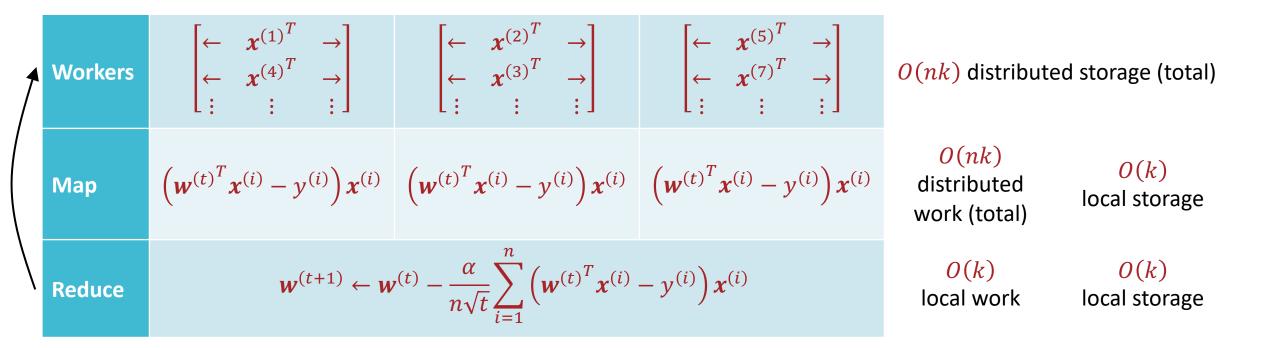
$$\nabla_{\boldsymbol{w}} L_{\mathcal{D}}(\boldsymbol{w}^{(t)}) = 2 \sum_{i=1}^{n} \left(\boldsymbol{w}^{(t)^{T}} \boldsymbol{x}^{(i)} - \boldsymbol{y}^{(i)} \right) \boldsymbol{x}^{(i)}$$

b. Update the weights:

$$\mathbf{w}^{(t+1)} \leftarrow \mathbf{w}^{(t)} - \frac{\alpha}{n\sqrt{t}} \sum_{i=1}^{n} \left(\mathbf{w}^{(t)^{T}} \mathbf{x}^{(i)} - \mathbf{y}^{(i)} \right) \mathbf{x}^{(i)}$$

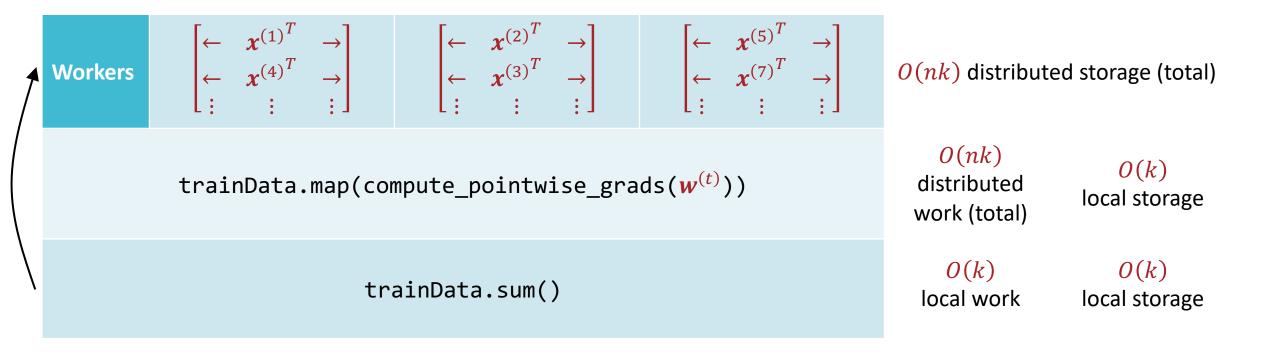
c. Increment $t: t \leftarrow t + 1$

• Output: $\mathbf{w}^{(t)}$



Issue: all workers must have the latest weight vector

Distributed Gradient Descent



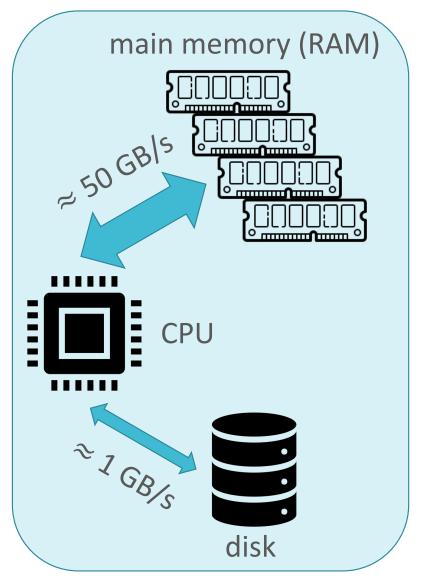
Issue: all workers must have the latest weight vector

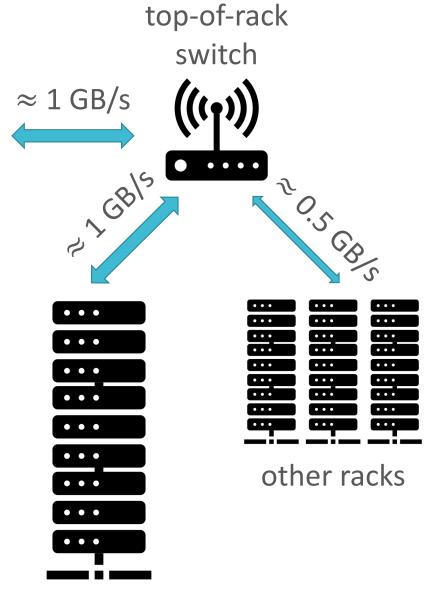
Distributed Gradient Descent

Gradient Descent

- Pros:
 - Trivially parallelizable
 - Each individual iteration is cheap
 - Can be further improved using stochastic variants
 - Guaranteed to converge on convex objective functions
- Cons:
 - Potentially slow convergence
 - Introduction of a hyperparameter
 - Network communication in each iteration

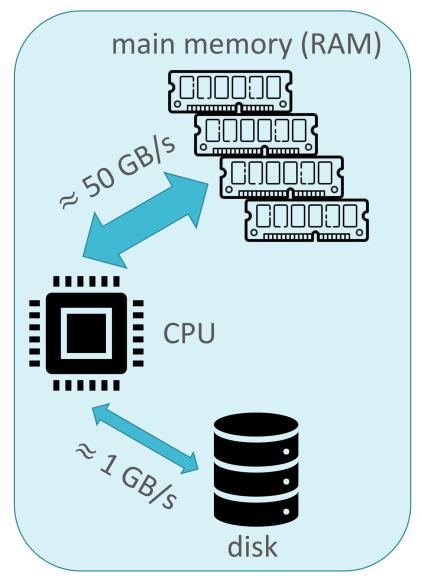
Recall: Communication Hierarchy

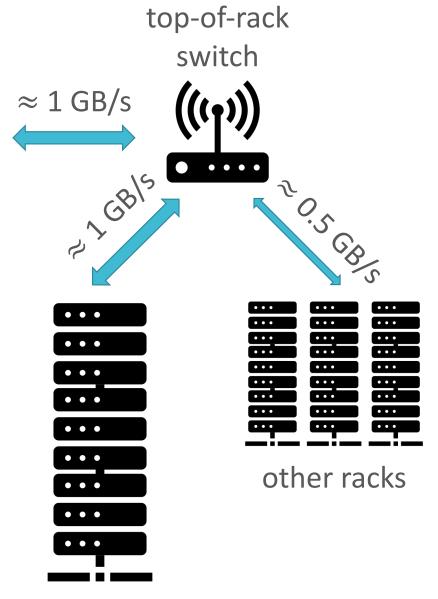




in-rack nodes

10-605/805
Principle #2:
Perform parallel and in-memory computation whenever possible





in-rack nodes

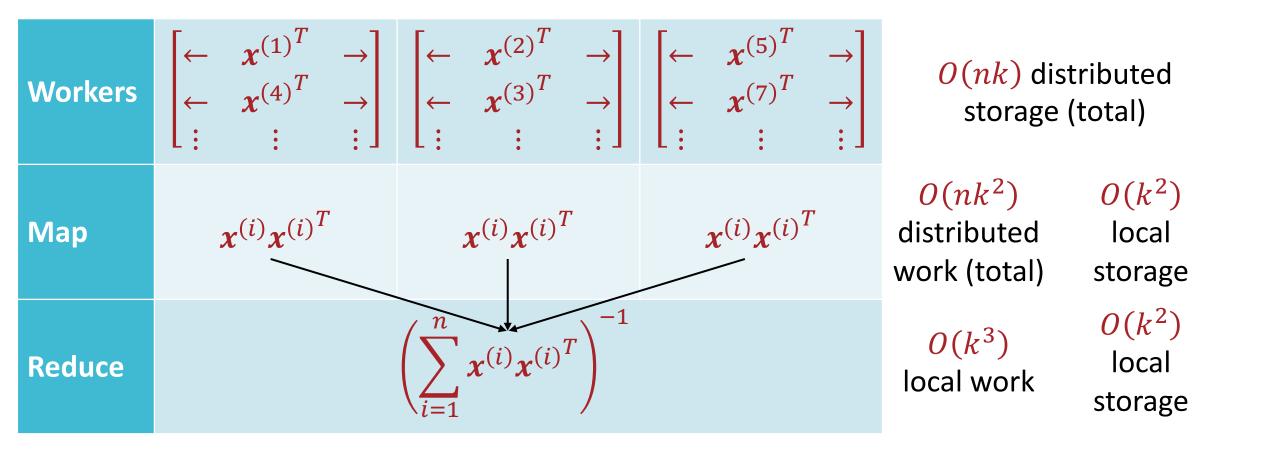
10-605/805 Principle #2: Perform parallel and in-memory computation whenever possible

 Persisting data in-memory reduces communication, especially for iterative procedures

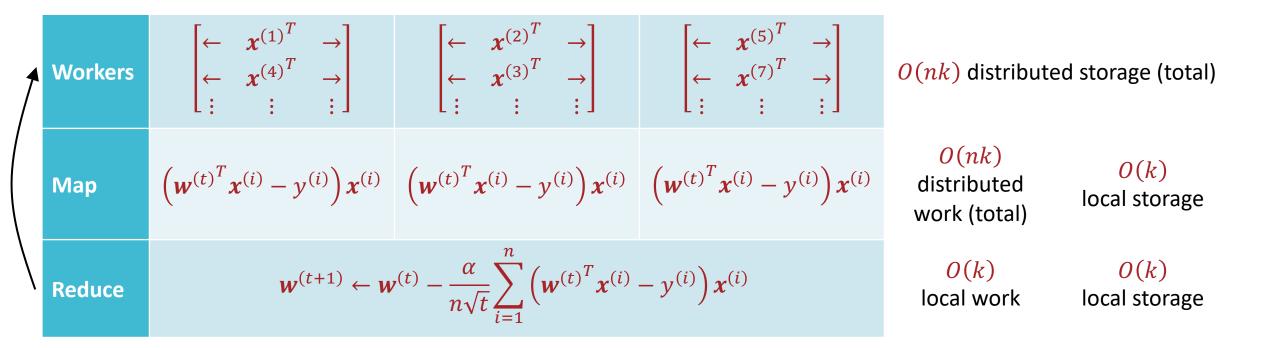
```
trainData.cache()
for t in range(num_iters):
    alpha_t = alpha / n * sqrt(t)
    grad = trainData.map(compute_pointwise_grads(w)).sum()
    w -= alpha_t * grad
```

10-605/805 Principle #3: Minimize network communication

- Inherently at odds with Principle #2 → need to tradeoff between parallelism and network communication
- Three types of objects that may need to be communicated:
 - Data
 - Models
 - Intermediate objects
- Strategies:
 - Keep large objects local
 - Reduce the number of iterations



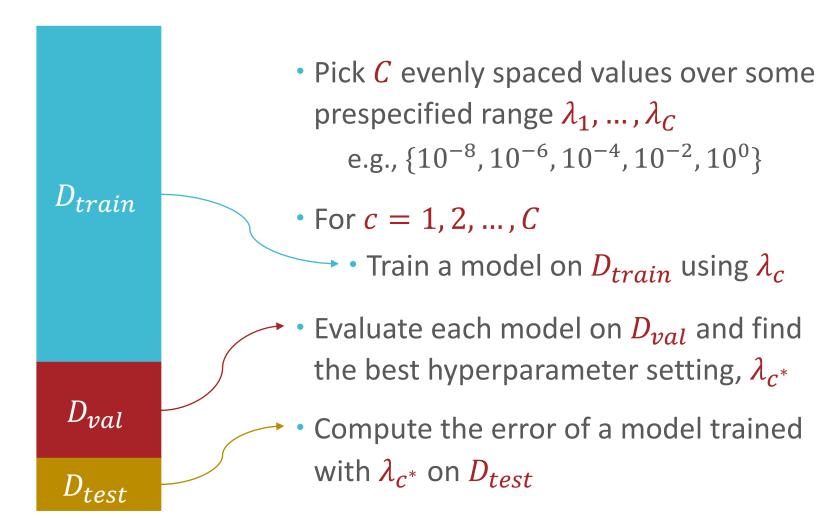
Data Parallel: Compute outer products locally



Issue: all workers must have the latest weight vector

Data Parallel: Compute pointwise gradients locally

Model Parallel: Train each hyperparameter setting on different machine(s) Hyperparameter optimization



Key Takeaways

- 10-605/805 Principles:
 - 1. Computation and storage should be linear in n and k
 - For linear regression:
 - When k is small, distribute covariance matrix computation using outer products
 - When k is large, minimize squared error via distributed gradient descent
 - 2. Perform parallel and in-memory computation whenever possible
 - Minimize network communication
 - Data vs model parallelism