#### Overview

• Previously we have seen how to learn a linear regression model using distributed closed-form solution.

• This distribution solution runs into minor issues when we try to scale.

# Limitations of the Closed-Form OLS: Solving for the $\theta$ s

- Problem is equivalent to inverting x 'x matrix.
  - Inverse does not exist if matrix is not of full rank.
    - E.g., if one column is a linear combination of another (collinearity).
    - Note that x 'x is closely related to the covariance of the x data.
      - So we are in trouble if two or more variables are perfectly correlated.
    - Numerical problems can also occur if variables are almost collinear.
- Equivalent to solving a system of *p* linear equations.
  - Many good numerical methods for doing this, e.g.,
    Gaussian elimination, LU decomposition, etc.
  - These are numerically more stable than direct inversion.
- Matrix inversion is not easily parallelized.

## **Alternative Ways of Solving OLS**

- · Closed-form solution
  - $\circ$  In this method, we will minimize RSS by explicitly taking its derivatives with respect to the  $\beta$  j 's (sometimes written as W, the weight vector) and setting them to zero.
  - Do this via calculus with matrices.

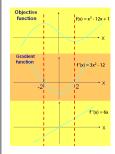
 $\theta * = (X T X) - 1 X T y \rightarrow$ 

- Gradient descent gives another way of minimizing RSS.
- Bayesian approach.
- Quadratic programming.
- · Others.

## **Linear Regression via Gradient Descent**

**Given**: Minimize f(x). Jq (W, X1m)= Minimize  $\Sigma$  i=1 m (WTXi-yi) 2

**Step 1**: Find the zeros of the gradient function f'(x) using bisection method, Newton-Raphson, or gradient descent.



RSS = Variance of  $\varepsilon$ 

$$0 = \frac{\partial \sum_{i} \hat{\varepsilon}_{i}^{2}}{\partial W} = \frac{\partial \left(\sum_{j=1}^{n} (X_{j}W - y_{i})^{2}\right)}{\partial W}$$

 $\nabla J(W) = \left(\sum_{j=1}^{n} (X_{j}W - y_{i})X_{j}\right)$ 

Gradient vector of partial derivatives

Pull model closer to examples with biggest residual.

$$W^{t+1} = W^t - \alpha^t \left( \sum_{i=1}^n \left( X_j W^t - y_i \right) X_j \right)$$

**Gradient descent** 

For another derivation see: <a href="http://cs229.stanford.edu/notes/cs229-notes1.pdf">http://cs229.stanford.edu/notes/cs229-notes1.pdf</a>.

## **OLS Using Gradient Descent**

 $J q (W, X 1 m) = Minimize \Sigma i=1 m (W T X i - y i) 2$ 

OLS objective function with decision variables W

- Initialize *W* = vector or zeros.
- Repeat until convergence.

 $W t+1 = W t - \alpha i (\sum_{j=1}^{n} n (X j W t - y i) X j)$ 

OLS batch update rule

• End repeat.

True gradient is approximated by the gradient of the cost function only evaluated at all examples; adjust parameters proportional to this approximate gradient.

Intuitively, drag weight vector closer to the incorrectly predicted examples.

#### **Batch vs. Stochastic**

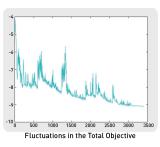
- That was batch-based descent.
- Stochastic gradient descent vs. batch.

## **OLS Using Gradient Descent**

#### Stochastic gradient descent

 $\nabla$  J wj (Wt)

Partial derivative WRT to variable w j of error function J(W) at point W t



#### Stochastic update (after each example)

Let W = (0, 0,...)

Repeat

For j in 0...n #each variable

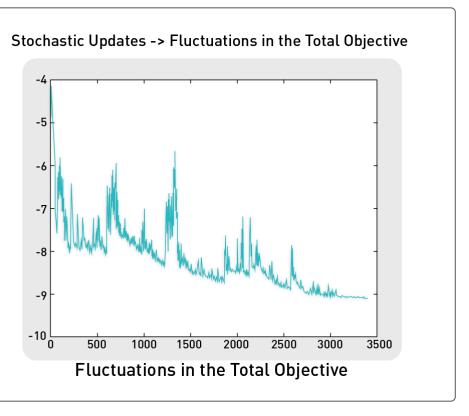
For i in 1...m #each example

$$W^{t+1} = W^t - \alpha^i \left( \sum_{j=1}^n \left( X_j W^t - y_i \right) X_j \right)$$

Until convergence (i.e., no big changes in W or error)

#### Stochastic Gradient Descent vs. Batch

- Stochastic gradient descent can start making progress right away and continues to make progress with each example it looks at.
- Often, stochastic gradient descent gets W "close" to the minimum much faster than batch gradient descent.
  - Note, however, that it may never "converge" to the minimum, and the parameters W will keep oscillating around the minimum of J(W); but in practice most of the values near the minimum will be reasonably good approximations to the true minimum.
- For these reasons, particularly when the training set is large, stochastic gradient descent is often preferred over batch gradient descent.



## **Two Approaches**

- Two gradient-descent approaches to linear regression
  - Batch gradient descent
  - Stochastic gradient descent based
- Are both amenable to parallelization? And if so, which do you expect to perform better?

# Stochastic Gradient Descent: Not Easy in MapReduce

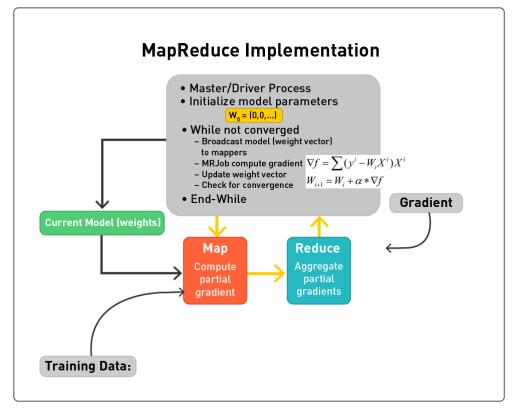
• Stochastic gradient descent requires the weight vector to be updated after each example (or minibatch of examples).

 "Some implementations of machine-learning algorithms, such as ICA, are commonly done with stochastic gradient ascent, which poses a challenge to parallelization. The problem is that in every step of gradient ascent, the algorithm updates a common set of parameters."

Some implementations of machine-learning algorithms, such as ICA, are commonly done with stochastic gradient ascent, which poses a challenge to parallelization. The problem is that in every step of gradient ascent, the algorithm updates a common set of parameters (e.g., the unmixing W matrix in ICA). When one gradient ascent step (involving one training sample) is updating W, it has to lock down this matrix, read it, compute the gradient, update W, and finally release the lock. This "lock-release" block creates a bottleneck for parallelization; thus, instead of stochastic gradient ascent, our algorithms above were implemented using batch gradient ascent.

#### **OLS via Distributed Gradient Descent**

- Master/Driver process.
- Initialize model parameters W 0 = (0,0,...).
- While not converged:
  - Broadcast model (weight vector) to mappers.
  - MRJob compute gradient ∇f = Σ ( y i W i X i ) X i.
  - Update weight vector W i+1 = W i + $\alpha * \nabla f$ .
  - · Check for convergence.
- End-While.



## **OLS via Distributed Gradient Descent: Mapper and Reducer**

- Master/Driver process.
- Initialize model parameters, W= vector of zeros: W 0 = (0,0, ...).
- While not converged:
  - Broadcast model (i.e., weight vector) to the worker nodes.
  - Mapper (MANY mappers).
    - Compute partial gradient for each training example.
    - Combine in memory  $\nabla f = \Sigma (yi WiXi) Xi$ .
    - Finally yield the partial gradient.
  - Reducer (single Reducer).
    - Aggregate partial gradients.
    - Yield full gradient  $\nabla f = \Sigma (yi WiXi)Xi$ .
  - Update weight vector W i+1 = W i +α∗∇f.
  - Check for convergence.
- End-While.

## **Divide and Conquer**

- Can we use a combiner? Yes!
  - In-memory combiner.
  - So use mapper final.
- How many reducers?
  - We can have one (should be sufficient).
  - Or many (in the case of many, the master has to aggregate the partial reducer aggregates).

