

Lecture 23

Fast Algorithms for Least Square Regression

Course: Algorithms for Big Data

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Linear Regression

- ▶ d variables (model parameters)
- ▶ n linear equations (observations)
- ▶ $n \gg d$ (over-constrained system)

$$\left\{ \begin{array}{l} 2x_1 + 3x_4 - 6x_5 + 12x_6 = 12 \\ x_1 - x_2 - 5x_3 + 12x_4 + x_5 = 1 \\ x_2 + x_3 + 6x_5 = 10 \\ 2x_1 + x_2 + x_4 + x_5 = -5 \\ -2x_1 + 3x_2 + 2x_4 - 9x_5 = 5 \\ 8x_1 - 4x_2 + x_3 + 4x_4 = 0 \\ x_2 + 10x_3 + 5x_5 = 12 \\ x_1 - 10x_4 - 2x_5 = -8 \end{array} \right.$$

$$\mathbf{Ax} = \mathbf{b}, \quad \mathbf{A} \in \mathbb{R}^{n \times d}, \quad \mathbf{b} \in \mathbb{R}^n$$

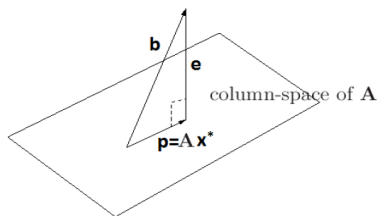
Choose $\mathbf{x} \in \mathbb{R}^d$ so that
 \mathbf{Ax} is close to \mathbf{b}

\mathbf{Ax} ranges over all linear
combinations of d columns of \mathbf{A}

Finding a closest point in the
column space of \mathbf{A} to the vector
 \mathbf{b}

Least square regression

$$\arg \min_x \|\mathbf{Ax} - \mathbf{b}\|_2 = \sum_{i=1}^n (b_i - A_{i*} \cdot \mathbf{x})^2$$



$$\mathbf{A}^T \mathbf{e} = \mathbf{0} \Rightarrow \mathbf{A}^T (\mathbf{b} - \mathbf{Ax}^*) = \mathbf{0}$$

$$\mathbf{A}^T \mathbf{Ax}^* = \mathbf{A}^T \mathbf{b} \Rightarrow \text{normal equation}$$

If \mathbf{A} is full-rank (has d independent columns), the unique solution is

$$\mathbf{x}^* = (\mathbf{A}^T \mathbf{A})^{-1} (\mathbf{A}^T \mathbf{b})$$

If \mathbf{A} is not full-rank there are multiple solutions. One solution is

$$\mathbf{x}^* = \mathbf{A}^\dagger \mathbf{b}$$

Here \mathbf{A}^\dagger is called the Moore-Penrose pseudoinverse of \mathbf{A} .

$$\mathbf{A}^\dagger = \mathbf{V} \mathbf{\Sigma}^\dagger \mathbf{U}^T, \quad \mathbf{A} = \mathbf{U}_{n \times d} \mathbf{\Sigma}_{d \times d} \mathbf{V}_{d \times d}^T$$

Here $\mathbf{U} \mathbf{\Sigma} \mathbf{V}^T$ is the SVD decomposition of \mathbf{A}

Assuming n is large, finding a solution \mathbf{x}^* takes a lot time (at least nd^2 time).

If we settle for an approximate solution, there is a faster randomized algorithm using sketching techniques.

[T. Sarlós](#). Improved approximation algorithms for large matrices via random projections. 2006.

Find \mathbf{x} where

$$\|\mathbf{Ax} - \mathbf{b}\|_2 \leq (1 + \epsilon) \|\mathbf{Ax}^* - \mathbf{b}\|_2$$

Similar to what we had in [JL lemma](#):

There is a matrix $\mathbf{S} \in \mathbb{R}^{r \times n}$ with random entries (where $r = \Theta(\frac{d}{\epsilon^2})$) such that with probability $1 - \exp(-d)$ for a $\mathbf{x} \in \mathbb{R}^d$

$$\|\mathbf{S}(\mathbf{Ax} - \mathbf{b})\|_2 \leq (1 + \epsilon)\|\mathbf{Ax} - \mathbf{b}\|_2$$

↓ (See the reference)

We can show with probability at least $1 - 1/4$:

$$\min_{\mathbf{x}} \|\mathbf{S}(\mathbf{Ax} - \mathbf{b})\|_2 \leq (1 + \epsilon)\|\mathbf{Ax}^* - \mathbf{b}\|_2$$

Strategy for the exact solution:

1. Output the exact solution x to the regression problem $\min_{\mathbf{x}} \|\mathbf{A}\mathbf{x} - \mathbf{b}\|_2$.

Time complexity: $\Omega(nd^2)$

Strategy for the approximate solution:

1. Sample a random matrix \mathbf{S} .
2. Compute $\mathbf{S} \cdot \mathbf{A}$ and $\mathbf{S} \cdot \mathbf{b}$.
3. Output the exact solution x to the regression problem $\min_{\mathbf{x}} \|(\mathbf{S}\mathbf{A})\mathbf{x} - (\mathbf{S}\mathbf{b})\|_2$.

Sarlos in his paper shows that using special random matrices \mathbf{S} one can obtain the matrix product \mathbf{SA} in time $O(nd \log d)$.

This gives the following time complexity for the approximate strategy:

Time complexity: $O(nd \log d) + \text{poly}(d/\epsilon)$

Subsequently, Clarkson and Woodruff show the following improved result.

Time complexity: $nnz(\mathbf{A}) + \text{poly}(d/\epsilon)$

Here $nnz(\mathbf{A})$ is the number of non-zero entries in matrix \mathbf{A} .