UBC MATH 307

LU Decompositions

Construction

Use Gaussian elimination to produce an upper triangular matrix, U. The lower triangular matrix L records the inverse of the row reduction operations. No pivoting of rows allowed. If A can be reduced by Gaussian elimination to row echelon form only with operations without scaling rows and without interchanging rows, then A has an LU decomposition of the form:

$$A = LU$$

Where L records the coefficients $c_{i,j}$ for each row reduction operation—add $c_{i,j}$ times row j to row i:

$$L = \begin{bmatrix} 1 & & & & \\ -c_{2,1} & 1 & & & \\ -c_{3,1} & -c_{3,2} & 1 & & \\ \vdots & \vdots & \ddots & \ddots & \\ -c_{m,1} & -c_{m,2} & \cdots & -c_{m,m-1} & 1 \end{bmatrix}$$

Back Substitution

To solve Ax = b, use the decomposition LUx = b

Solve:
$$L\boldsymbol{y} = \boldsymbol{b}$$
, then $U\boldsymbol{x} = \boldsymbol{y}$

Useful Properties of LU

$$rank(A) = rank(U)$$

$$det(A) = det(U) = \Pi(Diagonal entries U)$$

$$N(A) = N(U)$$

$$R(A) = \operatorname{span}\{\ell_1, \dots, \ell_r\} \text{ for } r = \operatorname{rank}(A),$$

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https://ubcmath.github.io/MATH307/

Matrix Norms and Error Analysis

Definitions

$$||A|| = \max_{\|\boldsymbol{x}\|=1} ||A\boldsymbol{x}||$$
 and $||A^{-1}|| = \frac{1}{\min_{\|\boldsymbol{x}\|=1} ||A\boldsymbol{x}||}$

Matrix norm generally has properties of vector norm.

$$\operatorname{cond}(A) = ||A|| ||A^{-1}||$$

Finding Matrix Norm

Use the SVD decomposition of A, the singular values of the Σ matrix determine the norm:

$$\sigma_{max} = ||A||$$
 and $\frac{1}{\sigma_{min}} = ||A^{-1}||$

Relative Error Formula

$$Ax = b \implies \frac{\|\Delta x\|}{\|x\|} \le \operatorname{cond}(A) \frac{\|\Delta b\|}{\|b\|}$$

Subspaces

Definitions

 $U \subseteq \mathbb{R}^n$ is a subspace if:

- 1. U contains the zero vector $\mathbf{0}$
- 2. $\mathbf{u}_1 + \mathbf{u}_2 \in U$ for all $\mathbf{u}_1, \mathbf{u}_2 \in U$
- 3. $c\mathbf{u} \in U$ for all $c \in \mathbb{R}$, $\mathbf{u} \in U$

 $\operatorname{span}\{u_1,\ldots,u_m\}=$ all possible linear combinations of vectors included in the brackets

Basis $\{u_1, \ldots, u_m\}$ forms basis of U if:

- 1. $\{u_1, \ldots, u_m\}$ is a linearly independent set
- 2. span $\{\boldsymbol{u}_1,\ldots,\boldsymbol{u}_m\}=U$

Range of A is $R(A) = \text{span}\{a_1, \dots, a_n\}$ column vecs

Rank-Nullity Theorem:

$$A(m \times n \text{ matrix}) \implies \operatorname{rank}(A) + \dim(N(A)) = n$$

Orthogonality of Fundamental Subspaces

$$\begin{array}{|c|c|c|c|c|}\hline R(A) = N(A^T)^{\perp} & N(A) = R(A^T)^{\perp} \\ R(A^T) = N(A)^{\perp} & N(A^T) = R(A)^{\perp} \\ \end{array}$$

Interpolation

Polynomial Interpolation

$$y = p(t) = c_0 + c_1 t + \dots + c_d t^d$$

$$\begin{bmatrix} 1 & t_0 & \cdots & t_0^d \\ 1 & t_1 & \cdots & t_1^d \\ \vdots & \vdots & \ddots & \vdots \\ 1 & t_d & \cdots & t_d^d \end{bmatrix} \quad \begin{bmatrix} c_0 \\ c_1 \\ \vdots \\ c_d \end{bmatrix} \quad = \begin{bmatrix} y_0 \\ y_1 \\ \vdots \\ y_d \end{bmatrix}$$

Vandermonde Matrix Coefficients

$$\det(\text{Vandermonde}) = \prod_{0 \le i < j \le d} (t_j - t_i)$$

Determinant is the product of the difference of each unique pair of t

Cubic Spline Interpolation

For N+1 points p(t) is a piecewise cubic function:

$$p_k(t) = a_k(t - t_{k-1})^3 + b_k(t - t_{k-1})^2 + c_k(t - t_{k-1}) + d_k$$

- 1. Interpolation at left endpoints:
- $p_k(t_{k-1}) = y_{k-1}$ for k = 1, ..., N yields N equations.
- 2. Interpolation at right endpoints:
- $p_k(t_k) = y_k$ for k = 1, ..., N yields N equations.
- 3. Continuity of p'(t):
- $p'_{k}(t_{k}) = p'_{k+1}(t_{k})$ for k = 1, ..., N-1 yields N-1 eqns.
- 4. Continuity of p''(t):
- $p_k''(t_k) = p_{k+1}''(t_k)$ for k = 1, ..., N-1 yields N-1 eqns.

Fundamental Subspaces

Space	Method of Solving
R(A)	LU decomp to find $r = rank(A)$, then pick
	the first r columns of L
N(A)	LU decomp then solve $Ux = 0$ using parame-
	ters c_1, c_2, \ldots, c_n for each free variable to get
	$x = c_1 v_1 + c_2 v_2 + \ldots + c_n v_n$ and the vectors
	v are the basis of $N(A)$
$R(A^T)$	transpose then LU, or take the orthogonal
	complement of $N(A)$
$N(A^T)$	transpose then LU, or take the orthogonal
, ,	complement of $R(A)$



Orthogonality and Projections

Orthogonal Complement

Vectors $x_1, \ldots, x_m \in \mathbb{R}^n$ are orthogonal if $\langle x_i, x_j \rangle = 0$ for all $i \neq j$, **orthonormal** if they are orthogonal and each is a unit vector, $||x_k|| = 1, k = 1, \ldots, m$.

$$U^{\perp} = \{ \boldsymbol{x} \in \mathbb{R}^n : \langle \boldsymbol{x}, \boldsymbol{y} \rangle = 0 \text{ for all } \boldsymbol{y} \in U \}$$

To find U^{\perp} place the basis vectors of U as the columns of matrix A, then:

$$U^{\perp} = \{ \boldsymbol{x} \in \mathbb{R}^n : A^T \boldsymbol{x} = 0 \} = N(A^T)$$

Projections

$$\operatorname{proj}_{\boldsymbol{u}}(\boldsymbol{x}) = \frac{\langle \boldsymbol{x}, \boldsymbol{u} \rangle}{\langle \boldsymbol{u}, \boldsymbol{u} \rangle} \boldsymbol{u} \quad \operatorname{proj}_{\boldsymbol{u}}(\boldsymbol{x}) = P \boldsymbol{x} \quad P = \frac{1}{\|\boldsymbol{u}\|^2} \boldsymbol{u} \boldsymbol{u}^T$$

Where P is a projection matrix with properties:

$$P^2 = P \quad P^T = P \quad \text{rank}(P) = 1$$

Any matrix with the first two properties is an

orthogonal projector

Gram-Schmidt Orthogonalization Algorithm

Take a basis of vectors for $U: \{u_1, \ldots, u_m\}$, then:

$$oldsymbol{v_k} = oldsymbol{u_k} - \sum_{n=1}^{k-1} \operatorname{proj}_{oldsymbol{v_n}}(oldsymbol{u_k})$$

Applied sequentially from k = 1 to k = m yields m orthogonal vectors forming a basis of U. Normalize all of the v vectors to form orthonormal basis.

Subspace Projections

 $\operatorname{proj}_{U}(\boldsymbol{x}) = \frac{\langle \boldsymbol{x}, \boldsymbol{u}_{1} \rangle}{\langle \boldsymbol{u}_{1}, \boldsymbol{u}_{1} \rangle} \boldsymbol{u}_{1} + \dots + \frac{\langle \boldsymbol{x}, \boldsymbol{u}_{m} \rangle}{\langle \boldsymbol{u}_{m}, \boldsymbol{u}_{m} \rangle} \boldsymbol{u}_{m}$ For an orthonormal basis, $\operatorname{proj}_{U}(\boldsymbol{x}) = P_{U} = AA^{T}$, where A has columns of \boldsymbol{u}

$$P_U + P_{U^{\perp}} = I$$

The projection of \boldsymbol{x} onto a subspace U gives the closest vector to \boldsymbol{x} in U

QR and Orthogonal Matrices

Orthogonal Matrices

A matrix A is **orthogonal** if $A^TA = AA^T = I$, then:

- 1. $||A\boldsymbol{x}|| = ||\boldsymbol{x}||$ for all $\boldsymbol{x} \in \mathbb{R}^n$
- 2. the columns and rows of are orthonormal
- 3. $A^T = A^{-1}$

QR Definition and Construction

A = QR where Q is an orthogonal matrix and R is an upper triangular matrix. Q by definition must be a square matrix, its columns include the orthonormal basis of R(A) via GS Algorithm for the first column entries, Q_1 . The orthonormal basis of $R(A)^{\perp}$ fills the remaining columns Q_2 .

$$\underbrace{A}_{n \times m} = QR = \underbrace{\begin{bmatrix} Q_1 & Q_2 \end{bmatrix}}_{n \times n} \underbrace{\begin{bmatrix} R_1 \\ 0 \end{bmatrix}}_{n \times m}$$

$$R_1 = egin{bmatrix} \langle oldsymbol{w}_1, oldsymbol{a}_1
angle & \langle oldsymbol{w}_1, oldsymbol{a}_2
angle & \cdots & \langle oldsymbol{w}_1, oldsymbol{a}_m
angle \ & \langle oldsymbol{w}_2, oldsymbol{a}_2
angle & \cdots & \langle oldsymbol{w}_2, oldsymbol{a}_m
angle \ & \ddots & dots \ & \langle oldsymbol{w}_m, oldsymbol{a}_m
angle \end{bmatrix}$$

The entire structure is simply a projection of the columns of A onto the orthonormal basis of A. The R vector contains the information to reverse the GS algorithm process used for Q:

$$\boldsymbol{a}_1 = \langle \boldsymbol{w}_1, \boldsymbol{a}_1 \rangle \boldsymbol{w}_1$$

$$\boldsymbol{a}_2 = \langle \boldsymbol{w}_1, \boldsymbol{a}_2 \rangle \boldsymbol{w}_1 + \langle \boldsymbol{w}_2, \boldsymbol{a}_2 \rangle \boldsymbol{w}_2$$

:

$$a_m = \langle \boldsymbol{w}_1, \boldsymbol{a}_m \rangle \boldsymbol{w}_1 + \langle \boldsymbol{w}_2, \boldsymbol{a}_m \rangle \boldsymbol{w}_2 + \dots + \langle \boldsymbol{w}_m, \boldsymbol{a}_m \rangle \boldsymbol{w}_m$$

As with any orthonormal basis:

$$\operatorname{proj}_{R(A)}(\boldsymbol{x}) = Q_1 Q_1^T \boldsymbol{x}$$

Least Squares Methods

Definition

The least squares method approximates a linear system with no solution by finding the closest vector to \mathbf{b} in R(A): $A\mathbf{x} = \operatorname{proj}_{R(A)}\mathbf{b}$

$$A x \cong b$$

We assume that A is a $m \times n$ matrix where m > n and also $\operatorname{rank}(A) = n$

Methods of Solving

$$A^T A \boldsymbol{x} = A^T \boldsymbol{b}$$
 $R_1 \boldsymbol{x} = Q_1^T \boldsymbol{b}$ $\boldsymbol{x} = A^+ \boldsymbol{b}$

Note that the QR method is generally preferred as having a lower condition number involved and not being overly computationally expensive.

Curve Fitting

For m data points (t_i, y_i) fitted to the linear combination of n functions f_k , the coefficients c_k can be found by applying least squares to $A\mathbf{c} \cong \mathbf{y}$

$$\mathbf{y} = \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_m \end{bmatrix} A = \begin{bmatrix} f_1(t_1) & f_2(t_1) & \cdots & f_n(t_1) \\ f_1(t_2) & f_2(t_2) & \cdots & f_n(t_2) \\ \vdots & & & \vdots \\ f_1(t_m) & f_2(t_m) & \cdots & f_n(t_m) \end{bmatrix} \mathbf{c} = \begin{bmatrix} c_1 \\ c_2 \\ \vdots \\ c_n \end{bmatrix}$$

PDP Decomposition and Power Method

Definitions

A $(n \times n)$ is diagonalizable if and only if A has n linearly independent eigenvectors, $A = PDP^{-1}$:

$$P = \underbrace{\begin{bmatrix} \boldsymbol{v}_1 & \cdots & \boldsymbol{v}_n \end{bmatrix}}_{\text{Eigenvectors}} D = \underbrace{\begin{bmatrix} \lambda_1 & & \\ & \ddots & \\ & & \lambda_n \end{bmatrix}}_{\text{Eigenvalues}}$$

Spectral Theorem

If A is a real symmetric matrix then the eigenvalues of A are real, the eigenvectors for distinct eigenvalues are orthogonal, and A has a

$$PDP^{-1}$$

where P is orthogonal: $A = PDP^{T}$. To make P orthogonal, make sure to normalize the eigenvectors to length 1.

Power Method

A dominant eigenvalue is a unique eigenvalue that is the largest for a given matrix A.

$$oldsymbol{x}_{k+1} = rac{Aoldsymbol{x}_k}{\|Aoldsymbol{x}_k\|_{\infty}} \qquad \|oldsymbol{x}_k\|_{\infty} = \max\{oldsymbol{x}_k\}$$

Check for convergence to $Ax_k = \lambda x_k$ with the Rayleigh quotient:

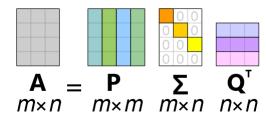
$$\lambda_k = rac{oldsymbol{x}_k^T A oldsymbol{x}_k}{oldsymbol{x}_k^T oldsymbol{x}_k}$$

The smallest eigenvalue can be found by iterating the inverse matrix (use LU with back substitution for repeated inverse operations).

SVD Decomposition

Definitions

Any $m \times n$ matrix A has a singular value decomposition $A = P\Sigma Q^T$ where P and Q are orthogonal matrices and Σ is a diagonal matrix ordered high to low $(\sigma_1 \geq \sigma_2 \geq ... \geq \sigma_r)$.



Adapted "Visualization of SVD" by Cmglee licensed under CC BY-SA 4.0

$$A = P\Sigma Q^T \quad ext{where} \quad \Sigma = \left[egin{array}{c|c} \sigma_1 & & & & \\ & \ddots & & \mathbf{0} \\ & & \sigma_r & \\ \hline & \mathbf{0} & & \mathbf{0} \end{array}
ight]_{m imes n}$$

Construction

 σ are known as the singular values of A. We find them as the $\sqrt{\lambda}$ of either AA^T or A^TA . The columns and rows of P and Q are corresponding normalized eigenvectors:

$$A^T A \mathbf{q}_i = \sigma_i^2 \mathbf{q}_i \qquad A A^T \mathbf{p}_i = \sigma_i^2 \mathbf{p}_i$$
 $\mathbf{p}_i = \frac{1}{\sigma_i} A \mathbf{q}_i \qquad \mathbf{q}_i = \frac{1}{\sigma_i} A^T \mathbf{p}_i$

Either P or Q may not have a full orthonormal basis defined based on the dimension and rank of A. In these cases simply extend the basis to a full one by finding the remaining orthogonal subspace and applying GS algorithm.

Important Properties of SVD

$$\operatorname{rank}(A) = r$$

$$\|A\| = \sigma_1$$

$$\|A^{-1}\| = 1/\sigma_r$$

$$\operatorname{cond}(A) = \sigma_1/\sigma_r$$

$$\operatorname{null}(A) = \operatorname{span}\{\boldsymbol{q}_{r+1}, \dots, \boldsymbol{q}_n\}$$

$$\operatorname{range}(A) = \operatorname{span}\{\boldsymbol{p}_1, \dots, \boldsymbol{p}_r\}$$

PCA and Applications of SVD

Principle Component Analysis

Create a data matrix X composed of row vectors, the same shape as used for least squares fitting. Normalize data so the columns have a mean value of 0. Apply SVD decomposition to the data matrix, the principle weight vectors are the q vectors. We can project X onto its principle components to reduce the dimension of the data while keeping the most significant weight vectors of dimensioning.

Pseudo Inverse

$$A^{+} = Q\Sigma^{+}P^{T}$$

$$\Sigma^{+} = \begin{bmatrix} \sigma_{1}^{-1} & & & \\ & \ddots & & \\ & & \sigma_{r}^{-1} \\ \hline & \mathbf{0} & & \mathbf{0} \end{bmatrix}_{n \times m}$$

SVD Expansion

$$A = \sum_{i=1}^r \sigma_i oldsymbol{p}_i oldsymbol{q}_i^T = \sigma_1 oldsymbol{p}_1 oldsymbol{q}_1^T + \dots + \sigma_r oldsymbol{p}_r oldsymbol{q}_r^T$$

Complex Linear Algebra

Inner Product

$$\langle \boldsymbol{v}, \boldsymbol{w} \rangle = v_0 \overline{w_0} + v_1 \overline{w_1} + \dots + v_n \overline{w_n}$$

 $\langle \boldsymbol{c}\boldsymbol{v}, \boldsymbol{w} \rangle = c \langle \boldsymbol{v}, \boldsymbol{w} \rangle \qquad \langle \boldsymbol{v}, \boldsymbol{w} \rangle = \overline{\langle \boldsymbol{w}, \boldsymbol{v} \rangle}$
 $\langle \boldsymbol{v}, c \boldsymbol{w} \rangle = \overline{c} \langle \boldsymbol{v}, \boldsymbol{w} \rangle \qquad \langle \boldsymbol{v}, \boldsymbol{v} \rangle = ||\boldsymbol{v}||$

Complex Analogs

Hermitian: $A = \overline{A}^T$ Unitary: $A^{-1} = \overline{A}^T$

If A is hermitian then diagonal entries of A are real

Roots of Unity

The Nth complex roots of 1: $\omega = e^{\frac{2\pi ik}{N}}$ for $0 < k < N, k \in \mathbb{C}^N$

Discrete Fourier Transform

Fourier Basis

The fourier basis of \mathbb{C}^N is an orthogonal basis defined as $f_0, ..., f_{N-1}$:

$$m{f}_k = egin{bmatrix} 1 & & & & \ \omega_N^k & & & \ \omega_N^{2k} & & & \ dots & & \ \omega_N^{(N-1)k} \end{bmatrix}$$

$$\overline{m{f}}_k = m{f}_{N-k}$$

Fourier Transform

$$DFT(\boldsymbol{x}) = F_N \boldsymbol{x}$$
 and $F_N =$

$$\begin{bmatrix} \overline{\boldsymbol{f}}_{0}^{T} \\ \overline{\boldsymbol{f}}_{1}^{T} \\ \vdots \\ \overline{\boldsymbol{f}}_{N-1}^{T} \end{bmatrix} = \begin{bmatrix} 1 & 1 & 1 & \cdots & 1 \\ 1 & \overline{\omega}_{N} & \overline{\omega}_{N}^{2} & \cdots & \overline{\omega}_{N}^{N-1} \\ 1 & \overline{\omega}_{N}^{2} & \overline{\omega}_{N}^{4} & \cdots & \overline{\omega}_{N}^{2(N-1)} \\ 1 & \vdots & \vdots & \ddots & \vdots \\ 1 & \overline{\omega}_{N}^{N-1} & \overline{\omega}_{N}^{2(N-1)} & \cdots & \overline{\omega}_{N}^{(N-1)^{2}} \end{bmatrix} \begin{bmatrix} \frac{1}{2} \left(\boldsymbol{f}_{k} + \overline{\boldsymbol{f}}_{k} \right) = \cos(2\pi k \boldsymbol{t}) \\ \frac{1}{2i} \left(\boldsymbol{f}_{k} - \overline{\boldsymbol{f}}_{k} \right) = \sin(2\pi k \boldsymbol{t}) \\ \vdots \\ \operatorname{For} \boldsymbol{x} = A \cos(2\pi k \boldsymbol{t} + \phi) : \\ \operatorname{DFT}(\boldsymbol{x}) = \frac{AN}{2} e^{i\phi} \boldsymbol{e}_{k} + \frac{AN}{2} e^{-i\phi} \boldsymbol{e}_{N-k} \end{bmatrix}$$

Note that $\frac{1}{\sqrt{N}}F_N$ is unitary

$$oldsymbol{x} = rac{1}{N} egin{bmatrix} oldsymbol{f}_0 & \cdots & oldsymbol{f}_{N-1} \ & oldsymbol{f}_1 \ & dots \ & oldsymbol{ar{f}}_{N-1}^T \ & dots \ & oldsymbol{ar{f}}_{N-1}^T \ \end{pmatrix} oldsymbol{x}$$

Let y = DFT(x) then:

$$\overline{\boldsymbol{y}[k]} = \boldsymbol{y}[N-k]$$

Inverse Transform

$$IDFT(\boldsymbol{y}) = \frac{1}{N} \overline{F}_N^T \boldsymbol{y}$$

DFT Analysis

Sinusoids

For N number of samples over a time period of 1, we have:

$$m{n} = egin{bmatrix} 0 \ 1 \ 2 \ dots \ N-1 \end{bmatrix} \qquad m{t} = (1/N) m{n} = egin{bmatrix} 0 \ 1/N \ 2/N \ dots \ (N-1)/N \end{bmatrix}$$

A discrete sinusoid signal is $x = A\cos(2\pi kt + \phi)$

The kth Fourier vector has the following sinusoidal properties:

$$\begin{aligned} \boldsymbol{f}_k &= \cos(2\pi k \boldsymbol{t}) + i \sin(2\pi k \boldsymbol{t}) \\ \frac{1}{2} \left(\boldsymbol{f}_k + \overline{\boldsymbol{f}}_k \right) &= \cos(2\pi k \boldsymbol{t}) \\ \frac{1}{2i} \left(\boldsymbol{f}_k - \overline{\boldsymbol{f}}_k \right) &= \sin(2\pi k \boldsymbol{t}) \end{aligned}$$

$$DFT(\boldsymbol{x}) = \frac{AN}{2}e^{i\phi}\,\boldsymbol{e}_k + \frac{AN}{2}e^{-i\phi}\,\boldsymbol{e}_{N-k}$$

Stemplots