Reviewing Linear Algebra

• Rank of a Matrix

The Rank of a matrix is defined as number of linearly independent columns in the matrix.

Recall: $x_1, x_2, x_3 \dots x_n$ n-Vectors are independent when

$$a_1x_1 + a_2x_2 + a_3x_3 \dots a_nx_n = 0$$

and $a_i = 0 \ \forall \ i \in \{1, 2, 3 \dots, n\}$

Note: Any \mathbf{r} -rank $n \times n$ matrix can be written as multiplication of 2 Matrices one containing \mathbf{r} independent column vectors(full column rank) and other containg \mathbf{r} independent row vectors(full row rank)!

$$A_{n\times n} = C_{n\times r} \times R_{r\times n}$$

Note: it is possible that Entities of original matrix are changed but the Transformation represented by this multiplication is same as original.

$$R_{1} = \begin{bmatrix} a_{11} & \dots & a_{1n} \\ \vdots & \dots & \vdots \\ a_{n1} & \dots & a_{nn} \end{bmatrix} = \begin{bmatrix} a_{11} & \dots & a_{1r} \\ \vdots & \dots & \vdots \\ a_{n1} & \dots & a_{nr} \end{bmatrix} \begin{bmatrix} a_{11} & \dots & a_{1n} \\ \vdots & \dots & \vdots \\ a_{r1} & \dots & a_{rn} \end{bmatrix}$$

• Rank-1 Matrices

An $n \times n$ matrix will be called Rank 1 Matrix when there is only 1 linearly independent columns or 1 linearly independent rows.can be represented as product of full column rank and full row rank matrices. . .

$$R_1 = \begin{bmatrix} a_{11} & \dots & a_{1n} \\ \vdots & \dots & \vdots \\ a_{n1} & \dots & a_{nn} \end{bmatrix} = \begin{bmatrix} r_1 \\ r_2 \\ \vdots \\ r_n \end{bmatrix} \begin{bmatrix} v_1 & v_2 & \dots & v_n \end{bmatrix}$$

• Rank-2 Matrices An $n \times n$ matrix will be called Rank 1 Matrix when there is only 1 linearly independent columns or 1 linearly independent rows.can be represented as product of full column rank and full row rank matrices...

$$R_{2} = \begin{bmatrix} a_{11} & \dots & a_{1n} \\ \vdots & \dots & \vdots \\ a_{n1} & \dots & a_{nn} \end{bmatrix} = \begin{bmatrix} r_{11} & r_{12} \\ r_{21} & r_{12} \\ \vdots & \vdots \\ r_{n1} & r_{n2} \end{bmatrix} \begin{bmatrix} v_{11} & v_{12} & \dots & v_{1n} \\ v_{21} & v_{22} & \dots & v_{2n} \end{bmatrix}$$

and so on we can represent any matrix as product of full column and row rank matrices

• Key Point

If we have a chain of matrix multiplication $A_1, A_2 \dots A_n$ defined as

$$A = A_1 \times A_2 \times A_3 \cdots \times A_n$$

and if we apply Transformation A on any vector \vec{x} then we can say that column space of A is subset of column space of A_1 because

$$Ax = (A_1 \times A_2 \times A_3 \cdots \times A_n)x$$
$$Ax = A_1 \times \left[(A_2 \times A_3 \cdots \times A_n) \times x \right]$$
therefore, $C(A) \subseteq C(A_1)$

• Matrix Multiplication

- Matirx with Vector(Left,Right)
- Row-Column(R-C)
- Column-Row(C-R)

Note: Any matrix can be written as sum of rank-1 matrices Intution \rightarrow as we know that we can represent any matrix(rank-r) as product of full column rank(r) and full row rank(r).now if we expand this representatin using C-R multiplication we will get sum of rank-1 matrices.

• Factorisation

-A = LU(Ellimination)

Suppose
$$L = \begin{bmatrix} L_1 & L_2 & \dots & L_k \end{bmatrix}$$
 and $U = \begin{bmatrix} U_1 \\ U_2 \\ \vdots \\ U_k \end{bmatrix}$ then

$$A = L_1 U_1 + L_2 U_2 \cdots + L_k U_k = \sum_{i=1}^k L_i U_i$$

$$L_i U_i = \begin{bmatrix} 0_{11} & 0_{12} & \dots & 0_{1(i-1)} & \dots & 0_{1n} \\ 0_{21} & 0_{22} & \dots & 0_{2(i-1)} & \dots & 0_{2n} \\ \vdots & \vdots & \ddots & \vdots & \dots & \vdots \\ 0_{(i-1)1} & 0_{(i-1)(2)} & \dots & 0_{(i-1)(i)} & & 0_{(i-1)n} \\ \vdots & \vdots & \vdots & \vdots & \vdots \\ 0_{m1} & 0_{m2} & \dots & 0_{m(i-1)} \end{bmatrix}$$

First i-1 rows and columns will be completly zero.

- -A = QR(Gram-Schimit)
- $-A = S\Lambda S^{-1}$ (Eigenvectors and Eigenvalues)
- $-A = Q\Lambda Q^T = (Symmetric Matrices)$
- $-A = u\Sigma v^T(SVD)$

- Fundamental theorem of Linear Algebra—It has two parts
 - 1. Relation between dimension of fundamental spaces of $A_{m \times n}$ (rank r).
 - dim(C(A)) = r
 - dim(R(A)) = dim(R(A)) = r
 - -dim(N(A)) = n r
 - $\dim(N(A^T)) = m r$
 - 2. Orthogonality of fundamental spaces.
 - C(A) is orthogonal to $N(A^T)$.
 - -R(A) or $C(A^T)$ is orthogonal to N(A).
- Useful Matrices
 - Rotation by $\theta \longrightarrow \begin{bmatrix} cos(\theta) & -sin(\theta) \\ sin(\theta) & cos(\theta) \end{bmatrix}$
 - Reflection about the line $y = tan(\frac{\theta}{2})x \rightarrow \begin{bmatrix} cos(\theta) & sin(\theta) \\ sin(\theta) & -cos(\theta) \end{bmatrix}$
 - Householder Matrices \rightarrow Suppose we have a unit vector $\vec{u} \in \mathbb{R}^n$ then householder matrix is defined as

$$H = I_n - uu^T, \quad u^T u = ||\vec{u}||^2 = 1$$

Note: H is Symmetric and Orthogonal.

- Hadamar Matrices \rightarrow these matrices are defined in powers of 2.

$$\mathcal{H}_{2n} = \begin{bmatrix} \mathcal{H}_n & \mathcal{H}_n \\ \mathcal{H}_n & -\mathcal{H}_n \end{bmatrix}$$
, $\mathcal{H}_2 = \begin{bmatrix} 1 & 1 \\ 1 & -1 \end{bmatrix}$

- Wavelet Matrices \rightarrow

$$H_4 = \begin{bmatrix} 1 & 1 & 1 & 0 \\ 1 & 1 & -1 & 0 \\ 1 & -1 & 0 & 1 \\ 1 & -1 & 0 & -1 \end{bmatrix}$$

$$H_8 = \begin{bmatrix} 1 & 1 & 1 & 0 & 1 & 0 & 0 & 0 \\ 1 & 1 & 1 & 0 & -1 & 0 & 0 & 0 \\ 1 & 1 & -1 & 0 & 0 & 1 & 0 & 0 \\ 1 & 1 & -1 & 0 & 0 & -1 & 0 & 0 \\ 1 & -1 & 0 & 1 & 0 & 0 & 1 & 0 \\ 1 & -1 & 0 & 1 & 0 & 0 & -1 & 0 \\ 1 & -1 & 0 & -1 & 0 & 0 & 0 & 1 \\ 1 & -1 & 0 & -1 & 0 & 0 & 0 & -1 \end{bmatrix}$$

and we can continue...

• Eigenvectors of Permutaion Matrix

suppose we have a permutation matrix with 1's shifted by 1 column right and the 1 in last column comes to 1st column example

$$P_4 = \begin{bmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \end{bmatrix}$$

Eigenvalues of such matrices are roots of $\lambda^n = 1$ let eigenvalues are $\lambda_0, \lambda_1, \ldots, \lambda_{n-1}$ then eigenvectors corrosponding to them are v_1, v_2, \ldots, v_n

$$v_i = \begin{bmatrix} 1 \\ \lambda_i \\ {\lambda_i}^2 \\ \vdots \\ {\lambda_i}^{n-1} \end{bmatrix} \quad \text{where} \quad \lambda_i = e^{(2\pi i)/n}$$

so matrix with eigenvectors as its columns would look like

$$F = \begin{bmatrix} 1 & 1 & 1 & \dots & 1 \\ 1 & \lambda_1 & \lambda_2 & & (\lambda_{n-1}) \\ 1 & {\lambda_1}^2 & {\lambda_2}^2 & & (\lambda_{n-1})^2 \\ \vdots & & & & \\ 1 & {\lambda_1}^{n-1} & {\lambda_2}^{n-1} & & (\lambda_{n-1})^{n-1} \end{bmatrix}$$

Note: It is Fourier Matrix! where $\lambda_i = \omega^i$

- **Eigen World** $A\vec{x} = \lambda \vec{x}$ where λ is eigenvalue and \vec{x} is eigenvector.
 - \circledast Eigenvalues of A^k are λ_i^k where λ_i is the eigenvalue of A.
 - * Applications
 - Difference Equations \rightarrow Solution to $v_{k+1} = Av_k$ is

$$v_k = \sum_{i=1}^n c_i (\lambda_i)^k x_i$$

– Differential Equations Solution to $\frac{dy}{dt} = Ay$ is

$$y = \sum_{i=1}^{n} c_i(e^{\lambda_i t}) x_i = e^{At} y_0$$

- Similar Matrices A, B are similar if $\exists M$ s.t $A = M^{-1}BM$.
 - \circledast Eigenvalues of A, B are same but eigenvectors may or may not!
 - \circledast Eigenvalues of PQ are same as QP as they are similar for (M=Q).
- Symmetric Matrices when $A = A^T$ we say A is symmetric!
 - * Eigenvalues of symmetric matrix are always real.
 - * Orthogonal eigenvectors can be choosen.
 - \circledast symmetric matrix can be written as $A = S \varLambda S^{-1} = S \varLambda S^T$

• Positive Definate Matrices

any symmetric matrix A is known as positive definate if its all eigenvalues are positive following are tests for positive definate matrices

- All eigenvalues $\lambda > 0$
- All pivots p > 0
- Sub-Determinants > 0
- $-A^TA$ is full rank matrix
- $\forall \vec{x} \neq \vec{0}, Energy = x^T Ax > 0 \text{ or } N(A) = \{\vec{0}\}\$

• Positive Semi-Definate Matrices

any symmetric matrix A is known as positive semi-definate if its all eigenvalues are non-negative.following are tests for positive semi-definate matrices

- All eigenvalues $\lambda > 0$
- All pivots $p \ge 0$
- Sub-Determinants ≥ 0
- $-A^TA$ may or may not full rank matrix
- $\ \forall \ \vec{x}, \ Energy = x^T A x \ge 0$
- Singular Value Decomposition(SVD) \longrightarrow for any $A_{m \times n} = u \Sigma v^T$

where u, v are left and right singular vectors respectively & Σ is diagonal matrix with non negative singular values (σ_i) .

Geometry \rightarrow Rotation $(v^T) \rightarrow$ Streching $(\Sigma) \rightarrow$ Rotation(u)

Parameters \rightarrow for $n \times n$ matrix we have toal n^2 parameters in SVD!

- $-2 \times 2 \rightarrow 2$ Parameters for streching, 1-1 for each rotations
- $-3 \times 3 \rightarrow 3$ Parameters for streching, 3-3 for each rotations
- $-4 \times 4 \rightarrow 4$ Parameters for streching,6-6 for each rotations
- $n\times n\to n$ Parameters for streching, $\frac{n^2-n}{2},\frac{n^2-n}{2}$ for each rotations

Representatin SVD form can be written in two ways but result is same

$$A = (m \times r)(r \times r)(r \times n) \tag{i}$$

$$A = (m \times m)(m \times n)(n \times n) \tag{ii}$$

in (i) we are just in row space, column space and $\Sigma_i > 0$ but in (ii) we have included both null spaces and extra zeros in diagonal of Σ

Lecture-7

• **Norm** it is a way to measure size of something in mathematics.represented by $\|.\|$ and following are some properties that a norm follows

- $-\|x\| \ge 0$, equality holds iff x = 0
- $-\|x+y\| \le \|x\| + \|y\|$ and $\|x-y\| \ge \|x\| \|y\| \ \forall \ x, y$
- $\|cx\| = abs(c)\|x\| = |c|\|x\|$ where c is a constant

• Norm of Vectors

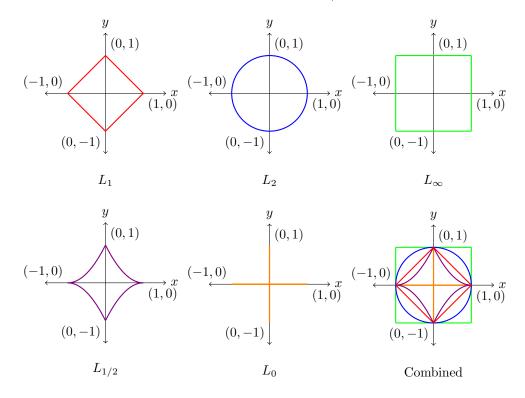
there are few kind of norms namely $L_1, L_2, \ldots, L_p, \ldots, L_{\infty}$ for any vector v with components v_1, v_2, \ldots, v_n is defined as

$$L_p = \left(\sum_{i=1}^n |v_i|^p\right)^{1/p} \tag{L_p Norm}$$

some of these are quite important such as $L_1, L_2 \& L_{\infty}$

- $-L_1 \rightarrow \text{Sum of the absolute value of components of vector, also know as Manhatten Distance or Taxicab Norm.}$
- $-L_2 \rightarrow \text{Sqrt}$ of Sum of the squares value of components of vector, also know as **Eucleadian Distance**.
- L_{∞} \rightarrow Maximum magnitude component of vector, also known as **Maximum Norm**
- $-L_0 \rightarrow$ defined as the no of non zero components of vector. Note that is does not follows the norm properties but it is useful in hamming codes this norm is also known as **Hamming Norm**

Here are the graphs in \mathbb{R}^2 for norms of $L_1, L_2, L_\infty, L_{1/2}, L_0 = 1$



• S-Norm if S is a symmetric matrix then S-Norm of a vector v is defined as following

$$||v||_S = \sqrt{(v^T S v)}$$
 (S-Norm)

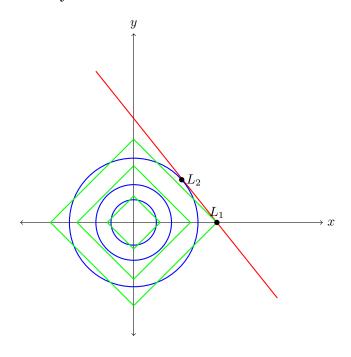
• Minising the Norm on Curve

suppose we have a curve C and we want to find the point on the tcurve that has minimum L_p norm there are two ways

- Algebrically
- Geometrically

example suppose curve ax+by=1 and we want to find the points on this line which has minimum L_1,L_2 norm

1.Geometrically



we know the picture of L_1, L_2 norm so for a particular norm starting from normvalue=0 we expand the picture the first intersection with the figure gives the point where that norm is minimum.

for L_2 norm we expanded the circle and for L_1 we expand diamond.

2. Algebrically

 L_1 norm will minimum at the closest sparsest point or where the maximum number of components are 0.

 L_2 we will write the general equation of circle with center at original and radius r, $x^2 + y^2 = r^2$ then solve with the curve for tangent.

Norm of Matrices Here are few types of matrix norms for matrix A_{m×n}
 ⊕ ||A||₂ → it is equal to maximum singular value in SVD form of the matrix.

$$||A||_2 = \max\left(\frac{||A\vec{v}||}{||\vec{v}||}\right)$$

as we know that $A\vec{v}_i = \sigma_i \vec{v}_i$ so it will be maximum for σ_1, \vec{v}_1

$$\therefore \|A\|_2 = \max\left(\frac{\|A\vec{v}\|}{\|\vec{v}\|}\right) = \sigma_1$$

 \circledast $\|\mathbf{A}\|_{\mathbf{F}} \to it$ is defined as Sqrt of sum of squares of all elements in the matrix, also known as **Frobenius Norm**.

$$||A||_F = \left(\sum_{i=1}^m \sum_{j=1}^n A_{ij}\right)^{1/2}$$

note as $A = u\Sigma v^T$ and u,v are orthogonal matrices therefore

$$||A||_F = ||\Sigma||_F = \left(\sum (\sigma_i)^2\right)^{1/2}$$

 \circledast $\|\mathbf{A}\|_{\mathbf{N}} \to \text{it}$ is defined as sum of singular values (σ_i) this norm is also known as **Nuclear Norm**.

$$||A||_N = \sum \sigma_i$$

Note: $||A||_2$, $||A||_F$, $||A||_N$ only depends on singular values $\sigma_1, \sigma_2 \dots, \sigma_n$ **Note:** Zero Norm of a matrix is defined as rank of matrix.

• Principle Component Analysis(PCA)

if we have a matrix A rank r and we have arbitrary another matrix B of rank $k \le r$, then according to **Eckart Young Theorem**

$$\boxed{\|A - B\| \ge \|A - A_k\|}$$

where A_k is diagonal matrix with largest k singular values of its σ_i 's example let we have a matrix A of order 4×4

$$\begin{bmatrix} 4 & 0 & 0 & 0 \\ 0 & 3 & 0 & 0 \\ 0 & 0 & 2 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$$

then for all matrices B of rank $k\leq 4\to \|A-B\|_k\geq \|A-A_k\|$ taking $k=2,\,\|A-B\|\geq \|A-A_2\|$ here A_2 will be

• Ways to Solve Least Square Problems $(A_{m \times n})$

$$1.\mathbf{A}^{\mathbf{T}}\mathbf{A}\mathbf{x} = \mathbf{A}^{\mathbf{T}}\mathbf{b}(m > n)$$

Basically we are trying to minimize the error, $\vec{e}=A\vec{x}-\vec{b}$ we will minimize its L_2 norm which is $\|\vec{e}\|^2$

$$\|\vec{e}\|^2 = \|A\vec{x} - \vec{b}\|^2 = (A\vec{x} - \vec{b})^T (A\vec{x} - \vec{b})$$

$$\|\vec{e}\|^2 = \vec{x}^T A^T A \vec{x} - \vec{x}^T A^T \vec{b} - \vec{b}^T A \vec{x} + \vec{b}^T \vec{b}$$

after solving it for minimum we will get $\vec{x} = (A^T A)^{-1} A^T \vec{b}$

Assumption: A has full column rank because then only A^TA is invertible

2.Pseudo Inverse $(\forall m, n)$

if a matrix is invertible then everything is fine but for non-invertible matrices we define the concept of Pseudo Inverse, which is closest to inverse as much as possible if $A_{m \times n} = u \Sigma v^T$ rank r then

$$A^{+} = v\Sigma^{+}u^{T}$$
 (Pseudo Inverse)

where Σ^+ is pseudo inverse of Σ as much as possible!

- \circledast How to solve $A\vec{x} = \vec{b}$ using pseudo inverse?
- ® When there is no inverse, our claim is that $\vec{x} = A^+ \vec{b}$ fits the best in the equation $A\vec{x} = \vec{b}$ or $||A\vec{x} \vec{b}||_2$ is minimum for $\vec{x} = A^+ \vec{b}$ as $A^+ = (A^T A)^{-1} A^T$ can be proved.

3.Undetermined(m < n)

we can use $\vec{x_1} = A^+ \vec{b}$ which is a minimizing L_2 norm but we can find $\vec{x_2}$ by minimizing L_1 norm. this case is quite useful in Deep Learning

4. $\|\mathbf{A}\mathbf{x} - \mathbf{b}\|^2 + \delta^2 \|\mathbf{x}\|^2$ when A is near singular this method can be used. Note: Here ellimination will not give satisfying results on solving for \mathbf{x}

suppose we minimize $||Ax - b||^2 + \delta^2 ||x||^2$ then we can write

$$\begin{bmatrix} A \\ \delta I \end{bmatrix} \begin{bmatrix} \vec{x} \end{bmatrix} = \begin{bmatrix} \vec{b} \\ \vec{0} \end{bmatrix}$$

say $\begin{bmatrix} A \\ \delta I \end{bmatrix} = A^*$ and $\begin{bmatrix} \vec{b} \\ \vec{0} \end{bmatrix} = \vec{b}^*$, so $A^*\vec{x} = \vec{b}^*$ and hence the solution will be

$$(A^*)^T A^* \vec{x} = (A^*)^T \vec{b}^* \rightarrow A^T A + \delta^2 I \vec{x} = A^T \vec{b}$$

Que: But what is δ and how to find it?

Ans: Later! for now let us see what will happen if $\delta \to 0$

suppose A is 1×1 matrix $A = [\sigma]$, so our solution become $\sigma^2 + \delta^2 \vec{x} = \sigma \vec{b}$

$$\vec{x} = \frac{\sigma \vec{b}}{\sigma^2 + \delta^2}$$

here we are just minimizing $(\sigma \vec{x} - \vec{b})^2 + \delta^2 x^2$ which we can do using calculus as well gives the same result *Analysising* as $\delta \to 0$

$$\vec{x} \to \begin{cases} \vec{b}/\sigma & \sigma > 0 \\ 0 & \sigma = 0 \end{cases}$$

 $Conclusion \rightarrow \text{as } \delta \rightarrow 0 \ \vec{x} \rightarrow A^+ \vec{b} \text{ where } A^+ \text{ is pseudo inverse.}$

it was not a proof but can be done using SVD form & the big result is that w.r.t L_2 norm the solution is

$$\vec{x} = A^{+}\vec{b} = (A^{T}A + \delta^{2}I)^{-1}A^{T}\vec{b}$$

Que: What if we minimize $A\vec{x} - \vec{b}$ w.r.t L_2 and $\delta^2 \vec{x}$ w.r.t L_1 norm?

$$minimize((\|A\vec{x} - \vec{b}\|_2)^2 + \delta^2(\|\vec{x}\|_1)^2)$$

its know as Lasso!

- **5.Gram Schimit** when columns are in *bad condition then first of all we will orthogonolize A (A = QR). it can we done in two ways *bad condition means it is possible that columns are nearly dependent!
 - Usual Projection way \rightarrow as new vector comes in keep elliminating projections on already calculated orthogonal vectors (ref-18.06).
 - Column Pivoting \rightarrow it involves the swapping of columns (if required).if we have $a_1, a_2 \dots, a_n$ as our columns of the matrix then
 - * choose biggest L_2 normed out of $a_1, a_2 \ldots, a_n$.
 - * next vector(a_i) out of remaining columns we will choose such that $||a_i||_2 P(a_i)$ is maximum. where $P(a_i)$ is the projection of a_i on already found orthogonal vector.
- **6.Krylov** when A is large sparse matrix

first we calculate krylov's space K_j then we use those basis to solve $A\vec{x} = \vec{b}$

$$K_i = \{\vec{b}, A\vec{b}, A^2\vec{b}, \dots, A^{j-1}\vec{b}\}$$

we can improve this method by orthogonolizing basis of K_j it is done by method which was figured out by **Arnoldi & Lanczos**

Books recomendations for numerical Linear Algebra

- Trefethen-because
- Golub and Vanloan

• Computing EigenValues

One of the method to compute eigenvalues is by setting $A - \lambda I$ to 0 then compute for eigenvalues but w.k.t as order of matrix increases (≥ 5) its very hard to find the eigenvalues.

to find the Approx eigenvalues we use following method even matlab and other maths software uses this method. here are the steps

- 1. Write, $A=A_0=Q_0R_0$, Q_0 is orthogonal & R_0 is upper triangular
- **2.** Use, $A_n = R_n Q_n \ \forall \ n \in \{1, 2 \dots, n\}$
 - $-A_0, A_1, \ldots, A_n$ will have same eigenvalues as they are similar!
 - $Accuracy \propto n$

More Improvement can be done in this method using Shifts before the step 1 and after the step 2

- **0.** Subtract sI form A or A_0
- **4.** Add sI to A_n

even more can be improved by reducing the matrix to $hessenberg\ form$ first then start with the step 0.

Note: if the matrix is symmetric then it will stay symmetric in the whole process and its hessenberg form will be a tridiagonal matrix.

• Hessenberg Matrix

when matrix is almost upper triangular except the diagonal near the main diagonal is non-zero

$$\begin{bmatrix} * & * & \dots & * & * \\ * & * & \dots & * & * \\ \vdots & \vdots & * & * & \vdots \\ 0 & 0 & * & \ddots & * \\ 0 & 0 & \dots & * & * \end{bmatrix}$$

• Computing Singular Values

w.k.t singular values are the square root of eigenvalues of A^TA so one way is to compute A^TA then calculate its eigenvalues but calculating A^TA for huge matrices will be a time consuming step (by computers) therefore another method is . . .

- 1. Write $A = u\Sigma v^T$
- **2.** Multiply by orthogonal matrices Q_1 and Q_2 s.t A becomes bidiagonal matrix $(Q_1u\Sigma v^TQ_2)$, as multiplying by orthogonal matrices does not affect the singular values.
- 3. hence A^TA will be tridiagonal 4. Use Eigenvalues method.

• Randomized Matrix Multiplication

suppose we have two matrix A with columns a_1, a_2, \ldots, a_k and B with rows b_1, b_2, \ldots, b_k then w.k.t usual multiplication of A, B is

$$A = \begin{bmatrix} \uparrow & \uparrow & \dots & \uparrow \\ a_1 & a_2 & \dots & a_k \\ \downarrow & \downarrow & \dots & \downarrow \end{bmatrix} \text{ and } B = \begin{bmatrix} \leftarrow & b_1 & \to \\ \leftarrow & b_2 & \to \\ \vdots & \vdots & \vdots \\ \leftarrow & b_k & \to \end{bmatrix}$$

$$M = A \times B = \sum_{i=0}^{k} a_i b_i$$

but in randomized matrix multiplication we consider product of columns of A & rows of B as sample space as follows

Sample Space
$$S = \{s_i = a_i b_i \mid i \in \{1, 2 \dots, k\}\}$$

with the probabilities p_i .

$$\sum_{i=1}^{k} p_i = 1$$

Prerequisites \longrightarrow Mean(μ), Varience(σ^2), Standard Deviation(σ)...

$$\mu = \sum p_i * s_i$$

$$\sigma^2 = \sum p_i * (s_i - \mu)^2 = \left(\sum p_i s_i^2\right) - (\mu^2)$$

$$\sigma = \sqrt{varience}$$

Note: In randomized matrix multiplication $\text{Mean}(\mu)$ of all the samples will always be correct we'll focus on to minimize the $\text{Varience}(\sigma^2)$ by choosing appropriate probability distribution.

Que But what should be the *Probability Distribution*? **Ans** *Norm Squared Distribution*! turns out to be the best

$$p_i = \frac{\|a_i\| \times \|b_i\|}{C}$$
, $C = \sum_{j=1}^k \|a_j\| \times \|b_j\|$ (0.1)

hence approx multiplication AB for 1 sample out of n will be

$$approx(AB) = \sum_{i=1}^{k} p_i \times \left(\frac{a_i b_i}{n p_i}\right) = \sum_{i=1}^{k} \frac{1}{n} \left(a_i b_i\right)$$

therefore $Mean(\mu)$ of whole sample space would be

$$\mu = \sum_{i=1}^{k} a_i b_i = AB$$

which is correct as it matches with original multiplication. now calculating varience of the for 1 sample

$$= \left(\sum_{i=1}^{k} \frac{p_i}{n} \left(\frac{\|a_i\| \|b_i\|}{p_i} \right)^2 \right) - \frac{1}{n} \left(\|AB\|_F \right)^2$$

$$= \left(\sum_{i=1}^{k} \frac{1}{np_i} (\|a_i\| \|b_i\|)^2\right) - \frac{1}{n} (\|AB\|_F)^2$$

substituting value of p_i from eq(0.1) we will get

$$= \frac{C}{n} \left(\sum_{i=1}^{k} (\|a_i\| \|b_i\|) \right) - \frac{1}{n} (\|AB\|_F)^2$$

$$= \frac{C^2}{n} - \frac{1}{n} (\|AB\|_F)^2 = \frac{1}{n} \left(C^2 - (\|AB\|_F)^2 \right)$$

then for the whole sample varience(σ) would be

$$\sigma^2 = C^2 - (\|AB\|_F)^2$$

Clearly $(\|AB\|_F)^2$ is a constant term so we have to minimize C.

Claim: C is minimum for Norm Squared Distribution!

Proof: It uses Lagrange Multiplier method

$$Target \rightarrow minimize \left(f(p_i) = \frac{\|a_i\|^2 \|b_i\|^2}{p_i} \right)$$

$$Constraint \rightarrow \sum p_i = 1$$

Lagrange Multiplier says that write net minimizing function as

$$f(p_i, \lambda) = Target + \lambda(Constraint)$$

$$f(p_i, \lambda) = f(p_i) + \lambda \left(\sum p_i - 1\right)$$

now use calculus and get the solution

$$\frac{\partial \left(f(p_i, \lambda)\right)}{\partial p_i} = 0 \implies -\frac{\|a_i\|^2 \|b_i\|^2}{p_i^2} + \lambda = 0 \tag{0.2}$$

$$\frac{\partial \left(f(p_i, \lambda) \right)}{\partial \lambda} = 0 \implies \sum p_i = 0 \tag{0.3}$$

from eq(0.2)

$$p_i = \sqrt{\frac{\|a_i\|^2 \|b_i\|^2}{\lambda}} = \frac{\|a_i\| \|b_i\|}{\sqrt{\lambda}} , \ \sqrt{\lambda} = C$$

• Low Rank Matrices

what will be the inverse of an Identity matrix after rank-1 perturbation or what is the inverse of $I - uv^T$ where I is order n matrix and $u, v \in \mathbb{R}^n$

Claim:

$$I = I + \frac{uv^T}{1 - v^T u}$$

Proof:

$$\begin{split} \left(I - uv^{T}\right) \left(I + \frac{uv^{T}}{1 - v^{T}u}\right) \\ &= I^{2} + \frac{uv^{T}}{1 - v^{T}u} - uv^{T} - \frac{uv^{T}uv^{T}}{1 - v^{T}u} \\ &= I + \frac{(I - uv^{T})uv^{T}}{1 - v^{T}u} - uv^{T} \\ &= I + \frac{(I - uv^{T})uv^{T} - (I - uv^{T})uv^{T}}{1 - v^{T}u} = I + 0 = I \end{split}$$

what if we do rank - k perturbation?

suppose we have an Identity matrix of order $n \times n$ and matrix U of order $n \times k$ with k independent columns also a matrix V of order ktimesn with k independent rows

Claim:

$$(I_n - UV^T)^{-1} = I_n + U((I_k - V^T U)^{-1})V^T$$

Proof:

$$(I_n - UV^T) (I_n + U ((I_k - V^T U)^{-1}) V^T)$$

$$= (I_n - UV^T) I_n + (I_n - UV^T) (U ((I_k - V^T U)^{-1}) V^T)$$

$$= I_n - UV^T + (I_n - UV^T) U (((I_k - V^T U)^{-1}) V^T)$$

$$= I_n - UV^T + U (I_k - V^T U) (I_k + V^T U)^{-1} V^T$$

$$I_n - UV^T + U I_k V^T = I_n$$

Applications:

- to solve $(I uv^T)\vec{x} = \vec{b}$
- new measurement in least squares(old was $(A\vec{x} = \vec{b})$) and normal equation was $A^T A\vec{x} = A^T \vec{b}$ now new normal equation is **dynamic**

$$\begin{bmatrix} A^T & v \end{bmatrix} \begin{bmatrix} A \\ v^T \end{bmatrix} \hat{x}_{new} = \begin{bmatrix} A^T & v \end{bmatrix} \begin{bmatrix} \vec{b} \\ b_{new} \end{bmatrix}$$
$$(A^T A + vv^T) \hat{x}_{new} = A^T b + v \vec{b}_{new}$$

Note¹: This formula was derived by Sherman, Morison, Woodbury(SMW) but it is also know as Matrix Inversion.

Note²: Kalman Filter uses this Idea of dynamic normal equations

now suppose we have a system of equation to solve Aw = b for w and then we want to solve $(A - uv^T)x = b$ for x then do find the x quickly using previous determined data we do the following.

- Solve equation Az = b for z
- & we have solution to the equation Aw = b for w
- Combine both to get x

$$x = w + \frac{wv^T z}{1 - v^T z}$$

Generalization of SMW

for any square matrix $A_{n\times n}$ and u,v be the $rank-k\leq n$ matrix with order $n\times k$ and $k\times n$ respectively

$$(A - uvT)^{-1} = A^{-1} - A^{-1}u (I - v^{T}A^{-1}u)^{-1} v^{T}A^{-1}$$

Note: rank-k perturbation in matrix will perturb its inverse by rank-k

Lecture-12

• Time Dependent Matrices

let us have a matrix A(t) is function of time then how would $A^{-1}(t)$ changes as times varies?

Que: Mathematically we are given $\frac{dA(t)}{dt}$ and we have to find $\frac{dA^{-1}(t)}{dt}$

Solution: let initially we had matrix A then it is changed to $B = A + \Delta A$ the using the following identity

$$B^{-1} - A^{-1} = B^{-1}(A - B)A^{-1}$$

$$(\Delta A)^{-1} = (A + \Delta A)^{-1} (-\Delta A)(A^{-1})$$

dividing by Δt ,

$$\left(\frac{\Delta A^{-1}}{\Delta t}\right) = \left(A + \Delta A\right)^{-1} \left(\frac{-\Delta A}{\Delta t}\right) \left(A^{-1}\right)$$

as $\Delta t \to 0$

$$\boxed{\frac{dA^{-1}}{dt} = -A^{-1} \left(\frac{dA}{dt}\right) (A^{-1})}$$

• Time Dependent Eigenvalues

Here are few facts (results) before starting

- Standard Eigenvalue Defination $\rightarrow A(t)x(t) = \lambda(t)x(t)$ --- (1)
- Transpose have same eigenvalues $\rightarrow y^T(t)A(t) = \lambda(t)y^T(t)$ ---(2)
- Normalization $\rightarrow y^T x = 1$
- $-AX = XA \& Y^TA = AY^T$ where X & Y are eigenvector matrix
- from $1.2 \rightarrow y^T(t)A(t)x(t) = \lambda(t)$

Note: When A will be symmetric $y^Ty = 1$ represents unit vector.

$$\frac{d(\lambda(t))}{dt} = \frac{d\left(y^T(t)A(t)x(t)\right)}{dt}$$

$$\frac{d(\lambda(t))}{dt} = \frac{d(y^T(t))}{dt}A(t)x(t) + y^T(t)\frac{d(A(t))}{dt}x(t) + y^TA(t)\frac{d(x(t))}{dt}$$

using (1),(2)

$$\frac{d(\lambda(t))}{dt} = \frac{d(y^T(t))}{dt}\lambda(t)x(t) + y^T(t)\frac{d(A(t))}{dt}x(t) + \lambda(t)y^T(t)\frac{d(x(t))}{dt}$$

$$\frac{d(\lambda(t))}{dt} = \lambda(t) \frac{d\left(y^{T}(t)x(t)\right)}{dt} + y^{T}(t) \frac{d(A(t))}{dt} x(t)$$

as $y^T(t)x(t) = 1$

$$\frac{d(\lambda(t))}{dt} = y^{T}(t)\frac{d(A(t))}{dt}x(t)$$

x(t) and $y^{T}(t)$ are known as right and left eigenvectors respectively at time t

*Interlacing theorem: if $S_{n\times n}$ is a symmetric matrix with eigenvalues $\gamma_1 \geq \gamma_2 \cdots \geq \gamma_n$ and $\vec{u} \in \mathbb{R}^n$ then eigenvalues of $S + uu^T$, $\lambda_1 \geq \lambda_2 \cdots \geq \lambda_n$ will be such that

$$\lambda_1 \ge \gamma_1 \ge \lambda_2 \ge \gamma_2 \cdots \ge \lambda_n \ge \gamma_n$$

if we add more rank-1 terms to S it will behave same way example let $w \in \mathbb{R}^n$ then eigenvalues of $S + uu^T + ww^T$, $\alpha_1 \ge \alpha_2 \cdots \ge \alpha_n$ then

$$\alpha_1 \ge \lambda_1 \ge \alpha_2 \ge \lambda_2 \cdots \ge \alpha_n \ge \lambda_n$$

we can relate $\alpha's$ and $\gamma's$ using both inequalities and also this theorem can be proved usign Weyl's Inequality(ref-NextLecture)

Note: uu^T will be rank - 1 positive semi-definate and one of its eigenvalue is $u^T u > 0$ and other 0's also eigenvector corresponding to $u^T u$ is u.

Que: What will happen if u is an eigenvector of S itself or $(Su = \lambda_i u)$ Ans: it means eigenvalue of $S + \theta u u^T$ corresponding to u will be $\lambda_i + \theta$

$$(S + \theta uu^T)u = Su + \theta uu^T u = (\lambda_i + \theta)u$$

clearly we can see that as θ increases $\lambda_i + \theta > \{\lambda_{i-1}, \lambda_{i-2}, \dots, \lambda_1\}$ after certain point? **No!** if θ will increase then we'll swap this $(\lambda_i + \theta)$ eigenvalue with on of appropriate $\lambda_x \in \{\lambda_{i-1}, \lambda_{i-2}, \dots, \lambda_1\}$ so that Interlace Inequlity Theorem holds true

Lecture-13

• Derivative of A²

$$\frac{dA^2}{dt} = \lim_{\Delta t \to 0} \frac{\Delta A^2}{\Delta t} = \lim_{\Delta t \to 0} \frac{(A + \Delta A)^2 - A^2}{\Delta t}$$
$$\frac{dA^2}{dt} = \lim_{\Delta t \to 0} \frac{A \cdot \Delta A + \Delta A \cdot A}{\Delta t}$$

$$\frac{dA^2}{dt} = A\frac{dA}{dt} + \frac{dA}{dt}A$$

• Time Dependent Singular Values

*What we know

 $-\sigma(t) = u^T A v$ where u, v are unit vectors and A is matrix

$$-v^Tv=1$$
 and $u^Tu=1$ therefore

$$\frac{dv^T}{dt}v + v^T\frac{dv}{dt} = 0 \quad \& \quad \frac{du^T}{dt}u + u^T\frac{du}{dt} = 0$$

as for vectors $x^T y = y^T x$ in real space(\mathbb{R}^n)

$$\frac{dv^T}{dt}v = v^T \frac{dv}{dt} = 0 \quad \& \quad \frac{du^T}{dt}u = u^T \frac{du}{dt} = 0 \tag{0.4}$$

*Derivation

$$\frac{d\sigma}{dt} = \frac{du^T}{dt}Av + u^T \frac{dA}{dt}v + u^T A \frac{dv}{dt}$$
$$\frac{d\sigma}{dt} = \frac{du^T}{dt}\sigma u + u^T \frac{dA}{dt}v + \sigma v^T \frac{dv}{dt}$$

using equation (0.4)

$$\frac{d\sigma}{dt} = 0 + u^T \frac{dA}{dt} v + 0$$

$$\boxed{\frac{d\sigma}{dt} = u^T \frac{dA}{dt} v}$$

• Weyl's Inequality

for any 2 real symmetric matrices of same order S, T

Lecture-14*

Prof: Alex Townsend

• Why Low Rank Matrices

let we have rank - k matrix X with singular values, $\sigma_1 \geq \sigma_2 \dots, \geq \sigma_n$ then we know the following facts

$$- rank(X) = k
- \sigma_i = 0, \forall i > k
- X = u_1 v_1^T + u_2 v_2^T \dots + u_k v_k^T - - -(i)
- dim(R(X)) = dim(C(X)) = k$$

• Defination of Low rank Matrix

if we can share the matrix $X_{m \times n}$ information more efficiently (space wise) then mn, using (i) we can send 2 matrix of order $m \times k$ and $k \times n$ instead of X we say $X_{m \times n}$ is low rank when

$$mk + kn < mn$$

or

$$k < \frac{mn}{m+n}$$

for square matrices

$$k < \left(\frac{n^2}{2n} = \frac{n}{2}\right)$$

in practical world we demand more i.e. $(k \ll n/2)$ to be more fast

Que: How do low rank matrices look like?

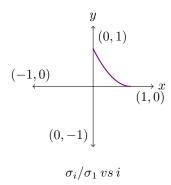
Ans: As rank decreases matrix get more aligned!

Example 1. Lower triangular matrix with all ones on lower half and diagonal elsewhere zeros then this matrix & its inverse are below

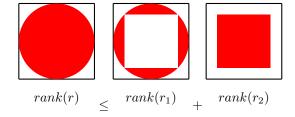
$$X = \begin{bmatrix} 1 & & & & \\ 1 & 1 & & & \\ 1 & 1 & 1 & & \\ \vdots & \vdots & & \ddots & \\ 1 & 1 & \dots & 1 & 1 \end{bmatrix} \quad \& \quad X^{-1} = \begin{bmatrix} 1 & & & & \\ -1 & 1 & & & \\ & -1 & 1 & & \\ & & \ddots & \ddots & \\ & & & -1 & 1 \end{bmatrix}$$

$$X^{T}X = \begin{bmatrix} 1 & -1 & & & & \\ -1 & 2 & -1 & & & & \\ & -1 & 2 & -1 & & & \\ & & \ddots & \ddots & \ddots & \\ & & & -1 & 2 & -1 \\ & & & & -1 & 2 \end{bmatrix} & & & \sigma_{i}(X) = \left[2sin\left(\frac{\pi(2i-1)}{2(2n+1)}\right) \right]^{-1}$$

Graph of Ratio of $(\sigma_i(X) \text{ to } \sigma_1(X)) \text{ vs } (i)$



Example2 Suppose we have following kind of matrix where non red area has 0 value



but $rank(r_2) = 1$, so $r \leq rank(r_1) + 1$ and then we can use symmetry to break r_1 as well.

• Numerical Rank

Numerical rank $(rank_{\epsilon})$ of a matrix X will be k, if for tolerence $(0 < \epsilon < 1)$

$$\sigma_{k+1}(X) \le \epsilon \sigma_1(X)$$

$$\sigma_k(X) > \epsilon \sigma_1(X)$$

assuming $\sigma_1 \geq \sigma_2 \cdots \geq \sigma_n$

Note: $rank(X) = rank_0(X)$

Eckart-Young: $\sigma_{k+1}(X) = ||X - X_k||_2$ How well X can be approximated!

• Numerical Low rank Matrices

- All low rank matrices

– Hilbert Matrices→it is full rank matrix but very low numerical rank

$$H_{ij} = \frac{1}{i+j-1}$$

- Vandermonde \rightarrow often we want V^{-1} as V is low numerical rank

$$V = \begin{bmatrix} 1 & x_1 & (x_1)^2 & \dots & (x_1)^{n-1} \\ 1 & x_2 & (x_2)^2 & \dots & (x_2)^{n-1} \\ \vdots & \vdots & \vdots & & \vdots \\ 1 & x_n & (x_n)^2 & \dots & (x_n)^{n-1} \end{bmatrix}$$

• The World is Smooth-Reade(1893)

if we take a polynomial and we form a matrix from integers sampling of this polynomial then that matrix will be

Example- P(x,y) = 1 + x + xy then matrix X with entries $X_{ij} = P(i,j)$ happens to be numerically low rank!

$$X = \begin{bmatrix} 1 & 1 & \dots & 1 \\ 1 & 1 & \dots & 1 \\ \vdots & \vdots & \ddots & \vdots \\ 1 & 1 & \dots & 1 \end{bmatrix} + \begin{bmatrix} 1 & 1 & \dots & 1 \\ 2 & 2 & \dots & 2 \\ \vdots & \vdots & \ddots & \vdots \\ n & n & \dots & n \end{bmatrix} + \begin{bmatrix} 1 & 2 & \dots & n \\ 2 & 4 & \dots & 1 \\ \vdots & \vdots & \ddots & \vdots \\ n & 2n & \dots & n^2 \end{bmatrix}$$

$$rank(X) < (1+1+1=3) \ll n$$

in general polynomial can be

$$P(x,y) = \sum_{i=0}^{m} \sum_{j=0}^{n} C_{ij} (x^{i}y^{j}), m < n$$

then rank of matrix X with $X_{ab} = P(a, b)$ will be less than equal to m^2 .

• The World is Sylvester

Matrices(X) satisfying the form AX - XB = C for some A, B, C then X is a numerical low rank

$$\sigma_{k+1} \le Z_k(E, F)\sigma_1(X)$$
, $rank(C) = r$

E, F are eigenvector matrices of A, B respectively

Example-Hilbert Matrix, Vandermonde

$$\begin{bmatrix} \frac{1}{2} & & & & \\ & \frac{3}{2} & & & \\ & & \ddots & \\ & & & \frac{2n-1}{2} \end{bmatrix} H - H \begin{bmatrix} \frac{-1}{2} & & & & \\ & \frac{-3}{2} & & & \\ & & \ddots & \\ & & & \frac{1-2n}{2} \end{bmatrix} = \begin{bmatrix} 1 & 1 & \dots & 1 \\ 1 & 1 & \dots & 1 \\ \vdots & \vdots & \ddots & \vdots \\ 1 & 1 & \dots & 1 \end{bmatrix}$$
$$\begin{bmatrix} x_1 & & & & \\ & x_2 & & \\ & & \ddots & & \\ & & & \ddots & \\ & & & & 1 & 0 \end{bmatrix} V - V \begin{bmatrix} 0 & & & & -1 \\ 1 & 0 & & & \\ & & \ddots & \ddots & \\ & & & & 1 & 0 \end{bmatrix} = \begin{bmatrix} 0 & 0 & \dots & 0 \\ 0 & 0 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & 0 \end{bmatrix}$$

• Free Parameters

Number of parameters required to know matrix are called free parameters

- Square Matrix: $A_{n\times n} \to n^2$
- Lower/Upper Triangular: $B_{n\times n} \to n(n+1)/2$
- Lower/Upper Triangular with diagonal (1's): $C_{n\times n} \to n(n-1)/2$
- Orthogonal Matrix: $Q_{n\times n} \to n(n-1)/2$
- Diagonal Matrix: $\Lambda_{n\times n} \to n$
- Symmetric Matrix: $S_{n\times n} \to n(n+1)/2$
- Eigenvector Matrix: $E_{n \times n} \to n^2 n$

different factorisation and there free parameters analysis

$$-A_{n\times n}=LU$$

$$\left(\frac{n(n-1)}{2}\right) + \left(\frac{n(n+1)}{2}\right) = n^2 \checkmark$$

$$-A_{n\times n}=QR$$

$$\left(\frac{n(n-1)}{2}\right) + \left(\frac{n(n+1)}{2}\right) = n^2 \checkmark$$

$$-A_{n\times n} = X\Lambda X^{-1}$$

$$(n^2 - n) + (n) + (0) = n^2 \checkmark$$

$$- A_{n \times n} = Q \Lambda Q^T$$

$$(n^2 - n) + (n) + (0) = n^2 \checkmark$$

$$-A_{n\times n} = QS$$

$$\left(\frac{n(n-1)}{2}\right) + \left(\frac{n(n+1)}{2}\right) = n^2 \checkmark$$

- SVD $A_{m \times n} = u \Sigma v^T$ Full rank matrix $(m \le n)$

$$\left(\frac{m(m-1)}{2}\right) + (m) + \left(mn - \frac{m(m+1)}{2}\right) = mn \checkmark$$

- SVD
$$A_{m \times n} = u \Sigma v^T$$
, $rank - r < min(m, n)$

$$\left(mr - \frac{r(r+1)}{2}\right) + (r) + \left(nr - \frac{r(r+1)}{2}\right) = mr + nr - r^2 \checkmark$$

• Saddle Points

For a function or curve $f(x_1, x_2, ..., x_n)$ point $a = (a_1, a_2, ..., a_n) \in \mathbb{R}^n$ will be a saddle point if directional derivative of in all directions is 0 & it is neither a maxima non a minima.

Que: From where saddle points arise?

Ans: There are 2 source of saddle points!

1. Suppose we have a symmetric matrix $S_{n\times n}$ and we want to minimize function f(x) on the given m constraints $A_{m\times n}(x) = b$ where f(x) is

$$f(x) = \frac{1}{2}x^T S x, \ x \in \mathbb{R}^n$$

now if we have a Lagrangian $L(x,\lambda) = f(x) - \lambda^T (Ax - b)$

Claim: $L(x,\lambda)$ do not have max or min it has saddle point when L'=0

Proof:

$$\frac{\partial L}{\partial x} = \frac{\partial}{\partial x} \left(\frac{1}{2} x^T S x - \lambda^T (A x - b) \right) = S x - A^T \lambda = 0$$

$$S x = A^T \lambda \tag{0.5}$$

$$\frac{\partial L}{\partial \lambda} = \frac{\partial}{\partial \lambda} \left(\frac{1}{2} x^T S x - \lambda^T (A x - b) \right) = A x - b = 0$$

$$A x = b \tag{0.6}$$

from equation (0.5),(0.6)

$$A'x' = b'$$

$$\begin{bmatrix} S & A^T \\ A & 0 \end{bmatrix} \begin{bmatrix} x \\ \lambda \end{bmatrix} = \begin{bmatrix} 0 \\ b \end{bmatrix}$$
(KKT)

doing ellimination on KKT matrix

$$\begin{bmatrix} S & A^T \\ A & 0 \end{bmatrix} \xrightarrow[R_2 - AS^{-1}R_1]{} \begin{bmatrix} S & A^T \\ 0 & -(AS^{-1}A^T) \end{bmatrix}$$

clearly KKT matrix (A') has n positive (as n pivots of S are positive) and m negative eigenvalues(as m pivots of $-AS^{-1}A^T$ are negative) therefore x' is saddle point of L

Note: $KKT \rightarrow Karush, Kuhn, Tucker$

2.Let S is symmetric matrix, **Rayleigh Quotient** function is defined as

$$R(x) = \frac{x^T S x}{x^T x}$$

Note¹: Max and Min values of R(x) are λ_1, λ_n for $x = q_1, q_n$ respectively. where λ_1, λ_n are the max and min eigenvalues of S and q_1, q_n are corrosponding eigenvectors

Note²: Here Saddle points are the Eigenvectors between $q_1 \& q_n$!

Example \rightarrow suppose we have a diagonal matrix D as follows

$$D = \begin{bmatrix} 5 & 0 & 0 \\ 0 & 3 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

then rayleigh quotient function R(x) will be

$$R(x) = \frac{x^T D x}{x^T x}$$

let x = (a, b, c) then

$$R(x) = \frac{5a^2 + 3b^2 + c^2}{a^2 + b^2 + c^2}$$

then R(x) will have minima and maxima when x is along the eigenvector corrosponding to 5, 1 respectively. and saddle point along the eigenvector corrosponding to 3.

Note: Saddle point means maximum of a minimum of a function

Saddle Point $\equiv max(min(f(x)))$

Lecture-16

• Probability and Statistics Basic Terminologies

- Sample Mean(\bar{x}) and Expected Mean(μ)

$$\boxed{\bar{x} = \frac{1}{n} \sum_{i=1}^{n} x_i} \boxed{\mu = \sum_{i=1}^{n} p_i x_i}$$

- Sample Varience(s^2) and Expected Varience(σ^2)

$$s^{2} = \frac{1}{n-1} \sum_{i=1}^{n} (x_{i} - \bar{x})^{2} \sigma^{2} = \sum_{i=1}^{n} p_{i} (x_{i} - \mu)^{2} = \left(\sum_{i=1}^{n} p_{i} x_{i}^{2}\right) - \mu^{2}$$

- Expected value of f(x)

$$E(f(x)) = \sum_{i=1}^{n} p_i f(x_i)$$

Note: for f(x) = x and $f(x) = (x - \mu)^2$ expected value $\mathbb{E}(f(x))$ is expected mean and varience respectively.

$$\boxed{\mu = \mathbb{E}(x)} \boxed{\sigma^2 = \mathbb{E}((x - \mu)^2) = \mathbb{E}(x^2) - \mathbb{E}(x)^2}$$

• Markov's Inequality: Probability of sample value greater than or equal to some number (a) is less than or equal to 1/a times the mean (μ) it uses an *assumption that all data points are non negative

$$P[x \ge a] \le \left(\frac{\mathbb{E}(x)}{a} = \frac{\mu}{a}\right), \ x_i \ge 0$$

• Chebyshevs's Inequality: Probability of a sample value having distance from the mean(μ) greater than some value (a) is less than or equal to $1/a^2$ times the varience(σ^2) no *assumption is used

$$(P[|x - \mu|] \ge a) \le \frac{\sigma^2}{a^2}$$

this inequality can be prooved using markov's inequality on $y_i = (x_i - \mu)^2$

- Covarience: it is defined as the measure of relationship between two random variables say x, y its formula is given by
 - Sample Covarience

$$C(x,y) = \frac{1}{n-1} \sum_{i=1}^{n} (x_i - \mu_x)(y_i - \mu_y)$$

- Population Covarience

$$C(x,y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x)(y_i - \mu_y)$$

- Expected Covarience

$$C(x,y) = \sum_{i=1}^{n} p_{ij}(x_i - \mu_x)(y_j - \mu_y)$$

• Covarience Matrix

$$V = \sum_{i,j}^{n} p_{ij} \begin{bmatrix} (x_i - \mu_x) \\ (y_i - \mu_y) \end{bmatrix} \begin{bmatrix} (x_i - \mu_x) & (y_i - \mu_y) \end{bmatrix}$$

its a symmetric matrix and its always positive semi-definate or positive definate

Example: for two random variables x,y V will be

$$V = \begin{bmatrix} \sigma_x^2 & \sigma_x \sigma_y \\ \sigma_y \sigma_x & \sigma_y^2 \end{bmatrix}$$

Lecture-17

• Taylor Series

- For 1-Variable Function $\rightarrow f(x)$

$$f(x + \Delta x) = f(x) + \Delta x \frac{df}{dx} + \frac{1}{2} (\Delta x)^2 \frac{d^2 f}{dx} \dots$$

- For n-Variable Function $\rightarrow f(x_1, x_2, \dots, x_n)$

$$f(x + \Delta x) = f(x) + (\Delta x)^T \nabla f(x) + (\Delta x)^T H(f(x))(\Delta x) \dots$$

where H(f(x)) is hessian matrix of function f(x)

Now suppose a function F(x) $x \in \mathbb{R}^n$ taking its gradient we a get vector function f(x) as follows

$$f(f_1(x), f_2(x), \dots, f_n(x)) = \nabla F(x), x \in \mathbb{R}^n$$

and

$$f(x + \Delta x) = f(x) + J\Delta x$$
, $J_{ab} = \frac{\partial f_a}{\partial x_b}$ (0.7)

J is jacobian matrix and $J = \nabla f(f_1(x), f_2(x), \dots, f_n(x))$ now to minimize F(x) we need to substitute f(x) = 0 or equivalently

$$f_1(x) = f_2(x) \cdots = f_n(x) = 0$$

Note: This is an equivalent problem to find the roots of $f_i(x)$

• Newton's Method

suppose x_k is a root of f then $f(x_k + \Delta x)$ must tends to 0 using eqⁿ(0.7)

$$0 = f(x_k) + J(x_k)(x_{k+1} - x_k)$$

or

$$x_{k+1} = x_k - (J^{-1}(x_k))f(x_k)$$

Example: let $f(x) = x^2 - 9$ then $J(x) = \lceil 2x \rceil$ and

$$x_{k+1} = x_k - \frac{1}{2x}(x_k^2 - 9) = \frac{x_k}{2} + \frac{9}{2x_k}$$

we can start with any x_0 here $x_0 \in \mathbb{R}$ after descent amount of steps either we will end up on one of roots of f(x) or it will diverge?

• Steepest Descent

at any point $x_k \in \mathbb{R}^n$ on F(x) we'll move in steepest direction at x_k

$$x_{k+1} = x_k - s_k \nabla F$$

also know as Exact Line Search!

Note¹: Hessian matrix (H) of scaler valued function F(x) is same as Jacobian matrix J(x) of its gradient $f(x) = \nabla F(x)$ so newton's method can also be written as

$$x_{k+1} = x_k - (J^{-1}(x_k))f(x_k) = x_{k+1} = x_k - (H^{-1}(x_k))f(x_k)$$

Note²: Convergence rate of newton's method is error squared e^2

Note³: Convergence rate of steepest descent method is linear wrt error!

• Convexity

- Convex Set: Set S will be called convex set if for any 2 elements $a, b \in S$ $x \in S$ where x is arbitrary element on the line joining a, b.

Note¹: Union of 2 sets is usually not convex

Note¹: Intersection of 2 sets is always convex

- Convex Function: Function $f(x_1, x_2, ..., x_n)$ is convex if $x^T H x \ge 0$ or Hessian matrix H of f is positive semi-definate over \mathbb{R}^n

• Convex Optimization

suppose we have a convex function F(x) and convex constraints set κ and we want to minimize F(x) over κ it is represented as

$$\min_{x \in \kappa} F(x)$$

Lecture-18

• Examples of convex functions Example 1.

$$f(x) = \frac{1}{2}x^T S x - a^T x - b$$
$$\nabla f(x) = S x - a$$
$$H(f(x)) = S$$

for f(x), x will be critical point when $\nabla f(x)=0$ or $x=S^{-1}a$ depending on H(f(x)), f(x) will be minimum or maximum or saddle at critical point

- if f(x) is minimum then x is known as **argmin** here $S^{-1}a$.
- if f(x) is maximum then x is known as **argmax** here $S^{-1}a$.

value of f(x) at $x = S^{-1}a$ will be

$$f(S^{-1}a) = \frac{1}{2}(S^{-1}a)^T S(S^{-1}a) - a^T (S^{-1}a) - b$$

$$= \frac{1}{2}a^T (S^{-1})^T S(S^{-1}a) - a^T (S^{-1}a) - b$$

$$= \frac{1}{2}a^T S^{-1} - a^T (S^{-1}a) - b$$

$$f(S^{-1}a) = -\frac{1}{2}a^T S^{-1} - b$$

Example 2. let $X_{n\times n}$ be a matrix then f(X) has n^2 variable $(x_{11}, x_{12}, \dots, x_{nn})$

$$f(X) = -log(|X|)$$

where |X| is determinant of X, gradient of f(X) can be computed using

$$\frac{\partial f(X)}{\partial x_{ij}} = \frac{C_{ij}}{|X|} = (X^{-1})_{ji}$$

• Gradient Descent

$$x_{k+1} = x_k - s_k \nabla f(x_k)$$

- Exact Line Search: Choose x_k move s_k distance along $-\nabla f(x)$
- Backtrakeing: Fix s then as we assect use $s = as, a^2s \dots a \in (0, 1)$

• Accelerating Gradient Descent

Gradient descent method can improvised by using an extra momentum term or daming term as follows

$$x_{k+1} = x_k - s_k Z_k$$
, $Z_k = \nabla f(x_k) + \beta Z_{k-1}$

let us apply this method to model $f(x) = \frac{1}{2}x^T Sx$ for which $\nabla f(x) = Sx$

$$x_{k+1} = x_k - s_k Z_k$$

$$Z_{k+1} - Sx_{k+1} = \beta Z_k$$

so we can write

$$\begin{bmatrix} 1 & 0 \\ -S & 1 \end{bmatrix} \begin{bmatrix} x_{k+1} \\ Z_{k+1} \end{bmatrix} = \begin{bmatrix} 1 & -s_k \\ 0 & \beta \end{bmatrix} \begin{bmatrix} x_k \\ Z_k \end{bmatrix}$$

now suppose q is the eigenvector of S corrosponding to λ i.e. $Sq = \lambda q$ also let $x_k = c_k q$ and $Z_k = d_k q$ then $Sx_k = \lambda c_k q$ hence

$$\begin{bmatrix} 1 & 0 \\ -\lambda & 1 \end{bmatrix} \begin{bmatrix} c_{k+1} \\ d_{k+1} \end{bmatrix} = \begin{bmatrix} 1 & -s_k \\ 0 & \beta \end{bmatrix} \begin{bmatrix} c_k \\ d_k \end{bmatrix}$$

$$\begin{bmatrix} c_{k+1} \\ d_{k+1} \end{bmatrix} = \begin{bmatrix} 1 & 0 \\ -\lambda & 1 \end{bmatrix}^{-1} \begin{bmatrix} 1 & -s_k \\ 0 & \beta \end{bmatrix} \begin{bmatrix} c_k \\ d_k \end{bmatrix}$$

$$\begin{bmatrix} c_{k+1} \\ d_{k+1} \end{bmatrix} = \begin{bmatrix} 1 & -s_k \\ \lambda & \beta - s_k \lambda \end{bmatrix} \begin{bmatrix} c_k \\ d_k \end{bmatrix} = R \begin{bmatrix} c_k \\ d_k \end{bmatrix}$$

now we have to choose β , s_k s.t matrix eigenvalues of R are small as possible $\forall m < \lambda < M$ where m, M are min and max eigenvalues of S.

Note¹: Condition Number κ is defined as ratio of M to m.

$$\kappa = \frac{M}{m}$$
 (Condition Number)

Note²: When $\kappa = 1$ it means S is multiple of identity matrix let eigenvalues of R are e_1, e_2 we want minimize $(max(|e_1|, |e_2|))$ Optimal Results:

$$s = \left(\frac{2}{\sqrt{M} + \sqrt{m}}\right)^2 \text{ and } \beta = \left(\frac{\sqrt{M} - \sqrt{m}}{\sqrt{M} + \sqrt{m}}\right)^2$$

• Nestrov's Idea

$$x_{k+1} = x_k + \beta(x_k - x_{k-1}) - s_k \nabla(x_k + \gamma(x_k - x_{k-1}))$$

• Linear Programming

in linear programming we try to optimize linear cost function over linear constraints as follows

cost function:
$$c^T x$$

constraints: Ax = b

where $x, c \in \mathbb{R}^n$, $x = (x_1, x_2, \dots, x_n), x_i \ge 0$ and A is $m \times n$, $b \in \mathbb{R}^m$

it can also be written as

$$\min_{Ax=b,x>0} c^T x$$

Note: Set of x s.t. Ax = b is known as **feasible set!**

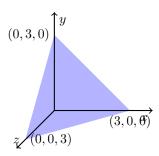
Example: Let $c^T x = x_1 + 2x_2 + 5x_3$ and we want to minimize it over the constraints $x_1 + x_2 + x_3 = 3$ then we can interpret this as

$$c = \begin{bmatrix} 1 \\ 2 \\ 5 \end{bmatrix}, x = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}, A = \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}, b = \begin{bmatrix} 3 \end{bmatrix}$$

here are 2 standard methods to solve linear programming problems

• Simplex Method Dantzig

we travel over the feasible set from corner to corner and according to that we choose optimum value as required. its complexity is Exponential!



for the above example minimum is at (3,0,0)

$\bullet \ \, \textbf{Interior Point Method} {\rightarrow} \textbf{Karmarkar} \\$

it uses steepest descent idea and calculus to travel within the feasible set and choose optimum!

• Duality LP

minization of $c^T x$ over the constraints Ax = b can also be solved as

$$\boxed{\min_{Ax=b,x>0} c^T x} \longleftrightarrow \boxed{\max_{A^T y \le c} b^T y}$$

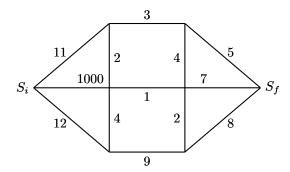
both are equivalent hence known as dual problems

– Weak Duality: $b^T y \leq c^T x, \forall x, y \implies \min(\max b^T y) = \max(\min c^T x)$

its known as minimax theorem or a saddle point at the optimum solution.

• Max Flow Problem/ Min Cut Problem

Suppose we have following graph network and some fluid flows from source S_i to sink S_f edges represent the maximum fluid can pass through it find the max fluid that can reach from S_i to S_f



Integer Flow Diagram

• Two Person Game

Suppose we have a following payoff matrix matrices for game I,II and III for persons x,y

	y_1	y_2
x_1	1	2
x_2	4	8

x is going to play row wise while y coloumn wise such that x wants to minimize the sum of rows and y wants to maximize the sum of columns

Note: it is zero sum game whatever one pays goes to other person only

Prof. Suvrit Sra

• Overview of Optimization Usecases

Large scale machine learning deals with large no. of training data points and usually in higher dimensions mathematically we writes

Training Data:
$$\{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\} \in [\mathbb{R}^d \times n]$$

in large scale machine learning d, n are quite large.

Example-1: Least squares

$$\frac{1}{n} (\|Ax - b\|_2)^2 = \frac{1}{n} \sum_{i=1}^n (a_i^T x - b_i)^2 = \frac{1}{n} \sum_{i=1}^n f_i(x)$$

Example-2: Lasso or l_1 least squares

$$\frac{1}{n} (\|Ax - b\|_2)^2 + \lambda \|x\|_1 = \frac{1}{n} \sum_{i=1}^n (a_i^T x - b_i)^2 + \lambda \sum_{j=1}^d \|x_j\|_1$$

Example-3: Support Vector Machine (SVM)

$$\frac{1}{2}(\|x\|)^{2} + \frac{C}{n} \sum_{i=1}^{n} max \left(0, 1 - y_{i}(x^{T} a_{i} + b_{i})\right)$$

all such examples can be written as finite sum problems

Example-4: Deep Neural Network(DNN)

$$\frac{1}{n} \sum_{i=1}^{n} loss(y_i, DNN(x:a)) = \frac{1}{n} \sum_{i=1}^{n} f_i(x)$$

Example-5: Maximum likelyhood estimation(MLE)

$$\frac{1}{n}\sum_{i=1}^{n} log - likelihood(x:a) = \frac{1}{n}\sum_{i=1}^{n} f_i(x)$$

• Stochastic Gradient Descent(SGD)

we had gradient descent in which one iteration used to look likelyhood

$$x_{k+1} = x_k - s_k \nabla f(x_k)$$

and as a finite sum

$$x_{k+1} = x_k - s_k \frac{1}{n} \sum_{i=1}^n \nabla f_i(x)$$

but practically calculating gradient at each point takes hours even days to complete the algorithm to overcome that we use SGD where we randomly pick an integer $i(k) \in \{1, 2, ..., n\}$ or in other words we just compute

$$x_{k+1} = x_k - s_k \nabla f_{i(k)}(x_k)$$

now each iteration become n times faster!

But: why SGD works?

Ans: in begin SGD makes rapid progress but near the solution it varies alot or fluctuations are observed in the solution domain!

Note: in ML we don't only care about solution to optimisation problem but also care about how well it works on **unseen data**.

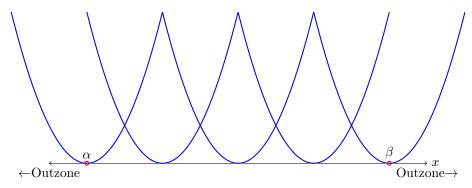
Example: Least Squares with $x \in \mathbb{R}$

$$\min f(x) = \frac{1}{2} \sum_{i=1}^{n} (a_i x - b_i)^2$$

by solving $\nabla f(x) = 0$ exact solution can be found as follows

$$x^* = \frac{\sum_{i=1}^n a_i b_i}{\sum_{i=1}^n (a_i)^2}$$

and minimum of $f_i(x)$ will be at $x_i = b_i/a_i$



Note¹: that the solution minarg, $x^* \in [\alpha, \beta]$, $\alpha = min(b_i/a_i), \beta = max(b_i/a_i)$

in the region $R = [\alpha, \beta]$ if we start the minization by picking a value of x which is in far out region from R then SGD will work kind of OK but as we enter R, SGD starts confusing

Note²: in 1-D if we are in far out region $\nabla f_i(x)$ has same sign as of $\nabla f(x)$ so using $\nabla f_i(x)$ instead of $\nabla f(x)$ we ensures our progress!

in higher dimensions the concept of sign changes to component $\nabla f_i(x)$ will have some component in the direction of true gradient $\nabla f(x)$.

Note³: SGD uses stochastic gradients g(x) such that the expection

$$\mathbb{E}(g(x)) = \nabla f(x)$$

in expection $\mathbb{E}(g(x))$ is the true gradient instead of $\nabla f(x)$ also know as **unbiased estimate** of true gradient

• Varients of SGD

1. Point wise

- Pick an index $i \in \{1, 2, ..., n\}$ without replacement OR
- Randomly pick an index $i \in \{1, 2, ..., n\}$ with replacement

then use $g_k = f_i(x_k)$ as stochastic gradient for x_{k+1}

most of us uses version 2 which is no repetation of i until we have gone through the entire training data set.

2. Using mini Batches

$$x_{k+1} = x_k - \frac{s_k}{|I_p|} \sum_{j \in I_p} \nabla f_j(x_k)$$

where $|I_p|$ is size of mini batch we have choosen

Note: Very large mini batches not favorable for DNN.because then we'll end up over fitting the neural network and it won't be able to make pridictions for unseen data.

Here are few practical challenges which we often face while training the models using in this version of stochastic gradient descent

- How to pich step size?
- Which mini batch to use?
- How to compute SGD? Backpropogation
- Gradient clipping!
- Adding momentum!

Lecture-21

• Neural Nets

suppose we have an arbitrary vector $x \in \mathbb{R}^n$ and we want to create a function that seprates the space into two parts such that

$$f(x) = \begin{cases} < 0 , 1st \text{ Part} \\ > 0 , 2nd \text{ Part} \end{cases}$$

Checkout Visual Examples → Click Here

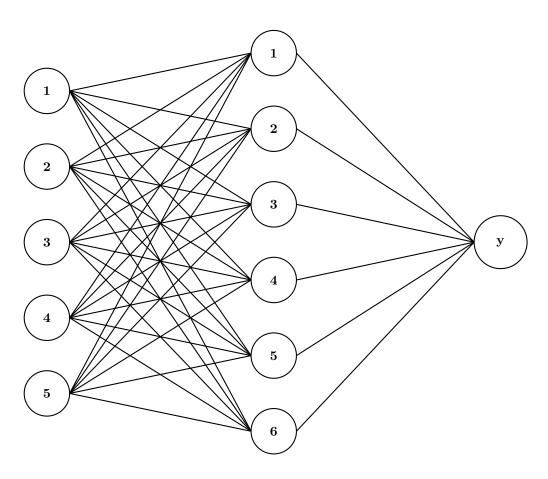
Here are few basic Terminologies

- Epoch: 1 iteration over the whole training data set
- Activation Function: Just a non linear function examples
 - * ReLu(r(x)), sigmoid $(\sigma(x))$, hyperbolic tangent (t(x))

$$r(x) = \max(0, x)$$
 , $\sigma(x) = \frac{1}{1 + e^{-x}}$, $t(x) = \tanh x = \frac{e^x - e^{-x}}{e^x + e^{-x}}$

Example of basic neural net with 1 layer

consider vector $x_0 \in \mathbb{R}^5$ is input element, and layer has 6 neurons then the output layer which here is just a single number this can be represented as



$$y = A_2(A_1x_0 + b_1) + b_2$$

where y is output of order 1×1 , A_2 is the matrix of order 1×6 between layer 1 and output layer of, A_1 is a matrix of order 6×5 between input layer and layer 1, x_0 is input vector and b_1, b_2 are constant(called affines)

But in each neuron there are non linear activation function say ReLu for now which acts upon each entity they get

$$y_1 = A_1x_0 + b_1$$

$$\downarrow \qquad \qquad \downarrow$$

$$x_1 = ReLu(y_1)$$

$$\downarrow \qquad \qquad \downarrow$$

$$y_2 = A_2x_1 + b_2$$

$$\downarrow \qquad \qquad \downarrow$$

$$y = x_2 = ReLu(y_2)$$

Note: For matrix A_1 have 30 free parameters and A_2 has 6 free parameters which we are trying to find by training the neural net In general while training we have many layers between input and output.

More abstractly if we have n layers then we have n-function composition where input vector $x\in\mathbb{R}^m$

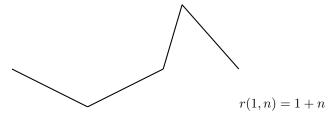
$$f = f_n (f_{n-1} \dots (f_1(x)))$$
, $f_n(x) = ReLu (A_n x_{n-1} + b_n)$

where each f_i and f is a continuous piecewise linear function!

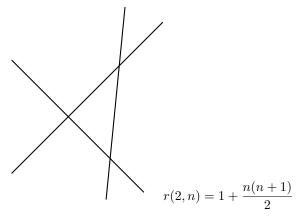
how many flat pieces we get if we are given or find r(m, n) = ?

- -r no of flat pieces
- -m- dimensions of x
- -n- no of folds

Example-1 if m = 1 then f(x) would look like where each critical point represents folds



Example-2 if m=2 then f(x) would look like where each region represent as plane and intersection lines represents folds



in General

$$r(m,n) = \binom{n}{0} + \binom{n}{1} \cdots + \binom{n}{m} = \sum_{i=1}^{m} \binom{n}{i}, m \le n$$

Recurrence Relation

$$r(m,n) = r(m,n-1) + r(m-1,n-1)$$

• Backpropogation

it is just automatic differentiation reverse mode checkout Wikipedia! and for more on reverse engineering checkout Christopher-Olah Blogs!

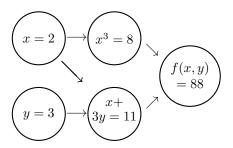
suppose we want to calculate partial differentiation of f(x, y) w.r.t x, y at (2,3) where f(x,y) is composition of f_1, f_2, f_3

$$f(x,y) = f_3(f_2(f_1(x,y))) = x^3(x+3y)$$

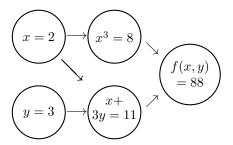
then

$$\frac{\partial f}{\partial x} = \frac{\partial f_3}{\partial f_2} \frac{\partial f_2}{\partial f_1} \frac{\partial f_1}{\partial x} = 4x^3 + 9yx^2 \; , \; \frac{\partial f}{\partial y} = \frac{\partial f_3}{\partial f_2} \frac{\partial f_2}{\partial f_1} \frac{\partial f_1}{\partial y} = 3x^3$$

Computational graph of f(x)



similarly we can draw computational graph of partial derivative and we found that reverse mode AD is much more efficient then forward mode



Note¹: Another such fact can be observed in matrix multiplication means A(BC) can be efficient then (AB)C depending on the order of A, B, C

Note²: No of operations required in $(A_{m \times n}) \times (B_{n \times p})$ are \boxed{mnp} .

• Completing Rank 1 Matrix

suppose we have to find all the entries of rank 1 matrix A of order $m \times n$ if we are given m+n-1 non zero entries of the martrix

$$A = uv^T$$

where $u \in \mathbb{R}^m$, $v \in \mathbb{R}^n$ and by choosing one of the entries of either u or v to be 1 we get m + n - 1 are free variables in the matrix

Note: Whether we can find whole matrix or not depends on the position of each given element in m + n - 1. take this example of 3×3

$$A = \begin{bmatrix} x & x & x \\ x & ? & ? \\ x & ? & ? \end{bmatrix} , B = \begin{bmatrix} x & x & ? \\ x & x & ? \\ ? & ? & x \end{bmatrix}$$

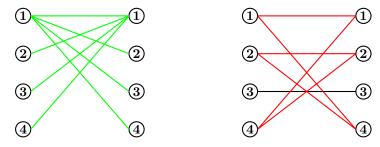
Note for matrix A we can find? setting determinant of sub matrix of 2×2 to 0.but in case of B we can't do it because determinant of matrix formed by $B_{11}, B_{12}, B_{21}, B_{22}$ may or may not be 0.if it is zero which means we are given only 4 free variables instead of 5 if its not 0 that's not possible in since it is rank-1 matrix. But how to know by positions whether we'll be able to complete the matrix or not?

Bipartite Graphs

if we make an undirected graph with source as row number and destination as coloumn number that will be a bipartite graph then if there is a cycle in graph then we can't complete the matrix! take example of 4×4

$$A = \begin{bmatrix} x & x & x & x \\ x & ? & ? & ? \\ x & ? & ? & ? \\ x & ? & ? & ? \end{bmatrix} , B = \begin{bmatrix} x & ? & ? & x \\ ? & x & ? & x \\ ? & ? & x & ? \\ x & x & ? & ? \end{bmatrix}$$

corrosponding bipartite graph of A and B would look like



in graph A there is no cycle but in graph B we does have a cycle

• Circulant Matrix:Cyclic Convulution Matrix

a general circulant matrix C of order $n \times n$ looks like this

$$\begin{bmatrix} x_1 & x_n & x_{n-1} & \dots & x_2 \\ x_2 & x_1 & x_n & \dots & x_3 \\ x_3 & x_2 & x_1 & \dots & x_4 \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ x_n & x_{n-1} & x_{n-2} & \dots & x_1 \end{bmatrix}$$

Note¹: Every circulant matrix C of order $n \times n$ can be written as sum of linear combinations powers of matrix P of order $n \times n$, where P shift one matrix of identity matrix is a permutation matrix as follows

$$P = \begin{bmatrix} 0 & 1 & 0 & 0 & \dots & 0 \\ 0 & 0 & 1 & 0 & \dots & 0 \\ 0 & 0 & 0 & 1 & \dots & 0 \\ \vdots & \vdots & \vdots & & & \vdots \\ 0 & 0 & 0 & 0 & \dots & 1 \\ 1 & 0 & 0 & 0 & \dots & 0 \end{bmatrix}$$

and P_2 is 2 shift matrix of identity and so on

$$C = c_0 I + c_1 P + c_2 P^2 \dots + c_{n-1} P^{n-1} = \left(\sum_{i=0}^{n-1} c_i P^i\right), P^0 = I$$

Note²: if C, D are circulant matrices so the matrix CD because C, D are just polynomial of P multiplication will also the polynomial of P.

• Convulution Operator

- **Non-Cyclic**: Take two vectors (a, b, c) and x, y, z then there non cyclic convulution is represented by operator (\circledast) and defined As

$$(a,b,c) \circledast (x,y,z) = (ax,ay+bx,az+by+cx,bz+cy,cz)$$

it can also be interpreted as multiplying two polynomial on LHS and writting the coefficients on RHS in vector form i.e.

$$(a + bp + cp^{2})(x + yp + zp^{2}) =$$

$$(ax) + (ay + bx)p + (az + by + cx)p^{2} + (bz + cy)p^{3} + (cz)p^{4}$$

- **Cyclic**: it is define similar to non-cyclic except the fact that $p^n = 1$ it is denoted by \otimes taking the previous example

$$(a,b,c)\otimes(x,y,z)=(ax+byz+cy,ay+bx+cz,az+by+cx)$$

it can also be interpreted as multiplying two polynomial on LHS and writting the coefficients on RHS in vector form i.e.

$$(a+bp+cp^{2})(x+yp+zp^{2}) =$$

$$(ax) + (ay+bx)p + (az+by+cx)p^{2} + (bz+cy)p^{3} + (cz)p^{4} =$$

$$(ax+by+cy) + (ay+bx+cz)p + (az+by+cx)p^{2} : p^{3} = 1, p^{4} = p$$

Note: in cyclic and non cyclic both the convulution product of sum of elements of each argument is same as the sum of elements of output taking the previous example

$$(a + b + c) * (x + y + z) = ax + ay + bx + az + by + cx + bz + cy + cz$$

in general for any 2 vector a, b convulution output c has this property

- for Cyclic

$$\left(\sum_{i=0}^{n-1} a_i\right) \times \left(\sum_{i=0}^{n-1} b_i\right) = \sum_{i=0}^{n-1} c_i$$

- for Non-Cyclic

$$\left(\sum_{i=0}^{n-1} a_i\right) \times \left(\sum_{i=0}^{n-1} b_i\right) = \sum_{i=0}^{2n-2} c_i$$

• Convulution of Continuos Functions

$$(f * g)(x) = f(x) * g(x) = \int_{-\infty}^{\infty} f(t)g(x-t)dt = \int_{-\infty}^{\infty} f(x-t)g(t)dt$$

• Eigenvalues and Eigenvectors of Circulant Matirx

Eigenvalues and eigenvectors of circulant matrix C of order $n \times n$ are same a permutation matrix Psuch that

$$C = c_0 I + c_1 P + c_2 P^2 \cdots + c_{n-1} P^{n-1}$$

but we know that eigenvalue of P are n^{th} roots of unity and eigenvectors matrix is same as fourier matrix (refer Page No.4 Reviewing Linear Algebra Section)

Note: Cyclicity of convulution in functions is seen by Periodicity!

• Toeplitz Matrix

Matirx of type below are known as Toeplitz Matirx

$$T = \begin{bmatrix} t_0 & t_1 & \dots \\ t_{-1} & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & t_1 \\ & \dots & t_{-1} & t_0 \end{bmatrix}$$

also know as Linear Shift Invarient(LSI) or Linear Time Invarient(LTI)! in machine learning images have too many feature so that we use **max pooling**(in a small part we take max of that part for all bits)

• Matrix With Orthogonal Eigenvectors

- Symmetric Matrix: $A^T = A \rightarrow \lambda = k + 0i$
- Anti-Symmetric or Skew Symmetric Matrix: $A^T = -A \rightarrow \lambda = 0 + ki$
- Orthogonal Matrix: $q_i^T q_i = 1$ only when i = j else 0
- Diagonal Matrix: Matrix itself is eigenvector matrix.
- Normal Matrix: $A^T A = AA^T$

• Recap: Convulution Formula

$$(c*d)_k = \sum_{i=0}^{n-1} c_i d_{k-i} , c \in \mathbb{R}^p, d \in \mathbb{R}^q$$

in non cyclic convulution there will be p+q-1 no. of terms in output and in case of cyclic convulution p=q=n and no of term will be n

• 2D Convulution of continuous functions

$$(f * g)(x,y) = \int_t \int_u f(x,y)g(x-t,y-u)dudt$$

• Convulution Rule

it shows the relation between point wise multiplication (.*) and convulution (*) for fast fourier transform as follows

$$F(c \circledast d) = F(c). * F(d)$$

convolve then transform OR transform then multiply component wise

Note¹: $\#\text{Cost LHS}=n^2 + nlogn \text{ and RHS}=2nlogn + n$

Note²: Eigenvalues of cd are same as product of eigenvalues of c, d

• Kronecker Product

suppose we have two matrices A, B order $n \times n$ each then

$$K = Kron(A, B) = \begin{bmatrix} a_{11}B & a_{12}B & \dots & a_{1n}B \\ a_{21}B & a_{22}B & \dots & a_{2n}B \\ \vdots & \vdots & \ddots & \vdots \\ a_{n1}B & a_{n2}B & \dots & a_{nn}B \end{bmatrix}$$

Note: K is a $n^2 \times n^2$ matrix & Kron(A, B) is Matlab Command

• Yet to complete the Lecture

- Kronecker Sum
- Laplacian

• Construction of Neural Nets:Learning Function

- Layer structure and features of input vector

$$v_l = \psi \left(A_l v_{l-1} + b_l \right)$$

where ψ is a non linear activation and $A_1, b_1, A_2, b_2, \dots, A_l, b_l$ are unknowns, together represented as x and in general number of unknowns(x) are greater than number of input features(v)

$$\#x \gg \#v$$

- Loss Function

$$L(x) = \frac{1}{n} \sum_{i=1}^{n} (F(x, v_i) - True_i)$$

- Common Loss Functions

- * Square Loss: L_2 norm, used regression.
- * Linear Loss: L_1 norms, used in lasso.
- * Hinge Loss: used in -1, 1 classification.
- * Cross Entropy Loss: used in neural nets

• Distance Matrices

Question: Given the distance matrix(D) of the points find the position matrix(X) of points in \mathbb{R}^d . (note we have to find dimension, d as well)

$$D_{ij} = ||x_i - x_j||^2$$

Solution:

$$d_{ij} = ||x_i - x_j||^2 = (x_i, x_i) - (x_i, y_i) - (y_i, x_i) + (y_i, y_i)$$

$$\uparrow \qquad \uparrow$$

Rank-1 Matrices depends only on x or y

so

$$D = d\mathbf{1}^T - 2X^TX + \mathbf{1}d^T$$

where **1** is column matrix of all 1's and now let matrix G such that G_{ij} represents dot product of X_i, X_j or $G = X^T X$

$$X^TX = G = -\frac{1}{2}D + \frac{1}{2}d\mathbf{1}^T + \frac{1}{2}\mathbf{1}d^T$$

we can find X from G or X^TX by following methods

$$X^TX=Q\Lambda Q^T\implies X=Q\sqrt{\Lambda}Q^T$$
 Singular Value Decomposition
$$X^TX=LDU=LDL^T\implies X=\sqrt{D}L^T$$
 Cholesky Factorisation

Question: Suppose $||x_1 - x_2||^2 = 1$, $||x_2 - x_3||^2 = 1$, $||x_3 - x_1||^2 = 6$ clearly triangular inequalities are failing here go and figure out the issues!

Solution:

$$D = \begin{bmatrix} 0 & 1 & 1 \\ 1 & 0 & 6 \\ 1 & 6 & 0 \end{bmatrix}$$

let position matrix of D is X then it means that $X^TX = G$ is not positive semi-definate when triangular inequalities fails

Note: G is positive semi definate $\leftarrow iff \rightarrow$ Triangular inequalities holds

• Procrustes Problem

Problem: It is a problem of finding optimal rotation from one basis to other suppose we have two set of basis vector in matrix form X, Y find an orthogonal matrix Q such that angular error is minimum

$$\min_{QQ^T=I} \left(\|YQ - X\|_F \right)^2$$

where $||A||_F$ is Frobenius norm of matrix A. Note we had following

$$||A||_F = \sum_{i=1}^m \sum_{j=1}^n (A_{ij})^2 = \sum_{i=1}^r (\sigma_i)^2 = trace(AA^T) = trace(A^TA)$$

- -trace(AB) = trace(BA)
- $trace(A^T B) = trace(B^T A) = trace(BA^T)$

Solution: write SVD of Y^TX as $u\Sigma v^T$ then $Q=uv^T$ proof yourself!

• Graph Clustring

Finding a cut in graph that divides the graph into equal set of clusters! for example to divide graph in 2 clusters C_1, C_2 we have to find x, y s.t.

$$\left(\sum \|a_i - x\|^2 + \sum \|b_i - y\|^2\right) \text{ is minimum where, } a_i \in C_1, b_i \in C_2$$

1. K-means Method: Following steps are for 2-means

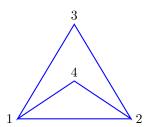
- For given a's,b's find there centroids x,y.
- Given x's,y's form best clusters
- Repeat above steps

best cluster can be formed by using the fact that arbitrary point will go in the closer cluster. Check out the internet

2. Spectral Clustering: it starts by graph Laplacian matrix(L)

$$L = A^T A = D - B$$

where A is incidence matrix, D is degree matrix, B is adjacency matrix. here are the definations of these matrices suppose we have following graph



- Incidence Matrix: -1 represents source and 1 destination

$$A = \begin{bmatrix} -1 & 1 & 0 & 0 \\ -1 & 0 & 1 & 0 \\ -1 & 0 & 0 & 1 \\ 0 & -1 & 1 & 0 \\ 0 & -1 & 0 & 1 \end{bmatrix}$$

- **Degree Matrix:** It is a diagonal matrix with degree as entries

$$D = \begin{bmatrix} 3 & 0 & 0 & 0 \\ 0 & 3 & 0 & 0 \\ 0 & 0 & 2 & 0 \\ 0 & 0 & 0 & 2 \end{bmatrix}$$

- Adjacency Matrix: It denoteds the links between nodes

$$B = \begin{bmatrix} 0 & 1 & 1 & 1 \\ 1 & 0 & 1 & 1 \\ 1 & 1 & 0 & 0 \\ 1 & 1 & 0 & 0 \end{bmatrix}$$

– Laplacian Matrix:

$$L = D - B = \begin{bmatrix} 3 & -1 & -1 & -1 \\ -1 & 3 & -1 & -1 \\ -1 & -1 & 2 & 0 \\ -1 & -1 & 0 & 2 \end{bmatrix}$$

Note matrix L is not full rank and its nullspace is 1 dimensional and

$$Lx = 0, x = \begin{bmatrix} c & c & \dots & c \end{bmatrix}^T$$

so it has 1 eigenvalue as 0 and other are positive and eigenvector y corrosponding to next eigenvalue greater than 0 is known as **Fiedler vector**

Clustering: Vector with positive components lies in one cluster and with negative components lies in other cluster

Note: x, y are orthogonal because L is symmetric P.S.D matrix