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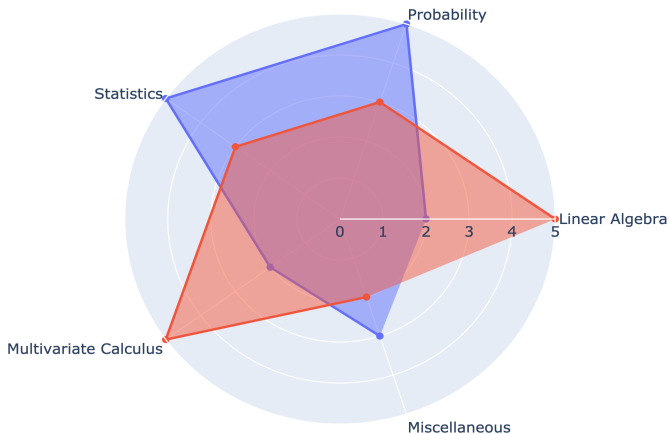
A journey from Linear Algebra to Machine Learning

Dr. Tanujit Chakraborty

Ph.D from Indian Statistical Institute, Kolkata, India.

Assistant Professor of Statistics at Sorbonne University

tanujitisi@gmail.com | <https://www.ctanujit.org>



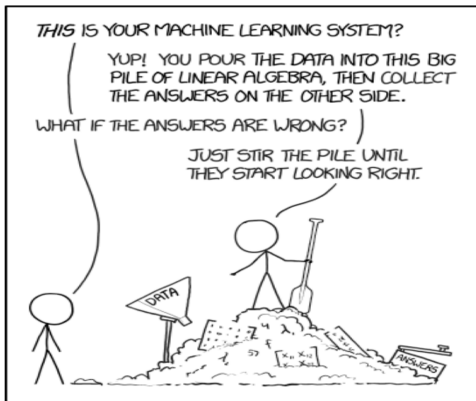
LINEAR ALGEBRA TO STATISTICAL
LEARNING

LINEAR ALGEBRA TO IMAGE PROCESSING

LINEAR ALGEBRA TO SPATIO-TEMPORAL DATA
ANALYSIS

LINEAR ALGEBRA TO STATISTICAL LEARNING

Importance in ML: We convert input vectors (x_1, \dots, x_n) into outputs by a series of linear transformations.



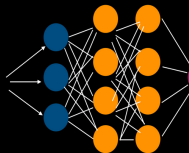
But what is RIGHT? And is that enough? (Image: [Machine Learning](#), XKCD)

ML models Leveraging Linear Algebra



[[12, 213, 44, 15],
[20, 113, 39, 10],
[13, 117, 45, 13],
[25, 120, 49, 14],
...
]]

Tensor



[[0.98, 0.45, 0.44, 0.95],
[0.23, 0.13, 0.39, 0.89],
[0.13, 0.77, 0.55, 0.23],
[0.25, 0.20, 0.49, 0.14],
...
]]

Tensor

Attend the session!

Atelactasis

Ok Google, set alarm for 5 AM.

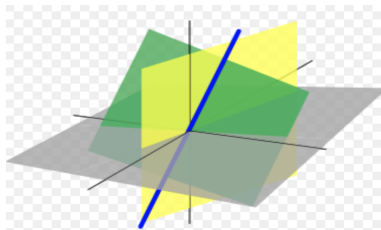
What is linear algebra?

- Linear algebra is the branch of mathematics concerning linear equations such as

$$a_1x_1 + \dots + a_nx_n = b$$

- In vector notation we say $a^T x = b$
 - Called a linear transformation of x
- Linear algebra is fundamental to geometry, for defining objects such as lines, planes, rotations.

- Linear equation
 $a_1x_1 + \dots + a_nx_n = b$ defines a
 plane in (x_1, \dots, x_n) space
 Straight lines define common
 solutions to equations.



- **Scalar:** Single number (represented by x)
 - In contrast to other objects in linear algebra, which are usually arrays of numbers.
 - They can be real-valued or be integers.
- **Vector:** An array of numbers arranged in order (represented by \mathbf{x})

$$\mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix}$$

- If each element is in \mathbb{R} then \mathbf{x} is in \mathbb{R}^n .
- We can think of vectors as points in space where each element gives coordinate along an axis.

- **Matrices:** 2-D array of numbers

- So each element identified by two indices.
- E.g.,

$$A = \begin{bmatrix} A_{1,1} & A_{1,2} \\ A_{2,1} & A_{2,2} \end{bmatrix}$$

- where $A_{i:}$ is i^{th} row of A , $A_{:j}$ is j^{th} column of A .
- If A has shape of height m and width n with real-values then $A \in \mathbb{R}^{m \times n}$.

- **Tensor:** A tensor is an array of numbers arranged on a regular grid with variable number of axes.

- Sometimes we need an array with more than two axes.
- E.g., an RGB color image has three axes.
- Element (i, j, k) of tensor denoted by $A_{i,j,k}$.

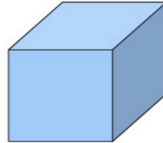
Shapes of Tensors



1d-tensor



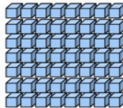
2d-tensor



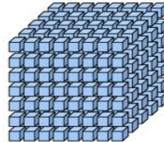
3d-tensor



4d-tensor



5d-tensor



6d-tensor

- We can add matrices to each other if they have the same shape, by adding corresponding elements
 - If A and B have same shape (height m , width n)

$$C = A + B \implies C_{i,j} = A_{i,j} + B_{i,j}$$

- A scalar can be added to a matrix or multiplied by a scalar

$$D = aB + c \implies D_{i,j} = aB_{i,j} + c$$

- Less conventional notation used in ML:

- Vector added to matrix

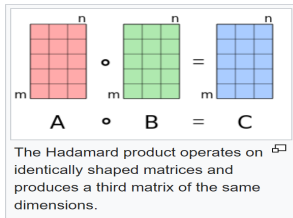
$$C = A + b \implies C_{i,j} = A_{i,j} + b_j$$

- Called broadcasting since vector b added to each row of A

- For product $C = AB$ to be defined, A has to have the same no. of columns as the no. of rows of B
 - If A is of shape $m \times n$ and B is of shape $n \times p$ then matrix product C is of shape $m \times p$

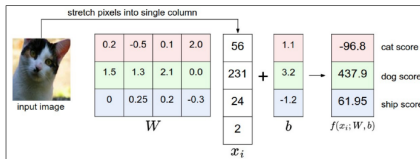
$$C = AB \implies C_{i,j} = \sum_k A_{i,k} B_{k,j}$$

- Note that the standard product of two matrices is not just the product of two individual elements.
- Such a product does exist and is called the element-wise product or the **Hadamard product** $A \odot B$.

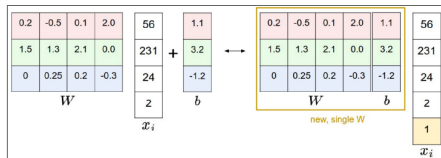


Vector x is converted
into vector y by
multiplying x by a matrix W

A linear classifier $y = Wx^T + b$



A linear classifier with bias eliminated $y = Wx^T$



Slide Credit: Sargur N. Srihari (Hari): <https://cedar.buffalo.edu/~srihari/>

- $Ax=b$
 - where $A \in \mathbb{R}^{n \times n}$ and $b \in \mathbb{R}^n$
 - More explicitly (n equations in n unknowns)

$$A_{1,1}x_1 + A_{1,2}x_2 + \dots + A_{1,n}x_n = b_1$$

$$A_{2,1}x_1 + A_{2,2}x_2 + \dots + A_{2,n}x_n = b_2$$

$$A_{n,1}x_1 + A_{n,2}x_2 + \dots + A_{n,n}x_n = b_n$$

- Consider the following:

$$A = \begin{bmatrix} A_{1,1} & \dots & A_{1,n} \\ \vdots & \vdots & \vdots \\ A_{n,1} & \dots & A_{n,n} \end{bmatrix} \quad x = \begin{bmatrix} x_1 \\ \vdots \\ x_n \end{bmatrix} \quad b = \begin{bmatrix} b_1 \\ \vdots \\ b_n \end{bmatrix}$$

Can view A as a **linear transformation of vector x to vector b**

- Sometimes we wish to solve for the unknowns $x = (x_1, \dots, x_n)$ when A and b provide constraints.

- Matrix inversion is a powerful tool to analytically solve $Ax = b$
- We can now solve $Ax = b$ as follows:

$$Ax = b$$

$$A^{-1}Ax = A^{-1}b$$

$$I_n x = A^{-1}b$$

$$x = A^{-1}b$$


- If A^{-1} exists there are several methods for finding it.
- Two closed-form solutions
 1. **Matrix inversion** $x = A^{-1}b$
 2. **Gaussian elimination**

- If A^{-1} exists, the same A^{-1} can be used for any given b
 - But A^{-1} cannot be represented with sufficient precision
 - It is not used in practice
- Gaussian elimination also has disadvantages
 - numerical instability (division by small no.)
 - $O(n^3)$ for $n \times n$ matrix
- Software solutions use value of b in finding x
 - E.g., difference (derivative) between b and output is used iteratively
- **Least squares solutions** of a $m \times n$ system

$$A x = b \implies (A^T A)x = A^T b \implies x = (A^T A)^{-1} A^T b = A^+ b.$$

$m \times n$

- A design matrix
 - N samples, D features

	# hours studied	# hours playing games	# classes missed		Grade
Student #1	10	3	0		87
Student #2	8	20	2		75
Student #3	5	1	5		63

- Feature vector has three dimensions
- This is a standard [regression problem](#).

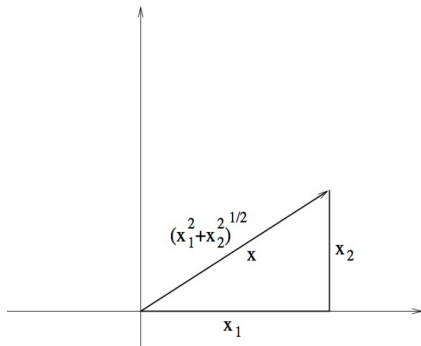
- Used for measuring the size of a vector
- Norms map vectors to non-negative values
- Norm of vector $x = [x_1, \dots, x_n]^T$ is distance from origin to x

It is any function f that satisfies:

$$f(x) = 0 \implies x = 0$$

$$f(x + y) \leq f(x) + f(y) \text{ (Triangle Inequality)}$$

$$\forall \alpha \in \mathbb{R}, f(\alpha x) = |\alpha|f(x)$$



- Definition:

$$\|x\|_p = \left(\sum_i |x_i|^p \right)^{1/p}$$

- L^2 Norm

- Called **Euclidean norm**
- Simply the Euclidean distance between the origin and the point x
- written simply as $\|x\|$
- Squared Euclidean norm is same as $x^T x$

- L^1 Norm

- Useful when 0 and non-zero have to be distinguished
- Note that L^2 increases slowly near origin, e.g., $0.1^2 = 0.01$

- L^∞ Norm

$$\|x\|_\infty = \max_i |x_i| \quad (\text{called } \mathbf{max \ norm})$$

- **Linear Regression** with x : a vector, w : weight vector

$$y(x, w) = w_0 + w_1 x_1 + \dots + w_d x_d = w^T x$$

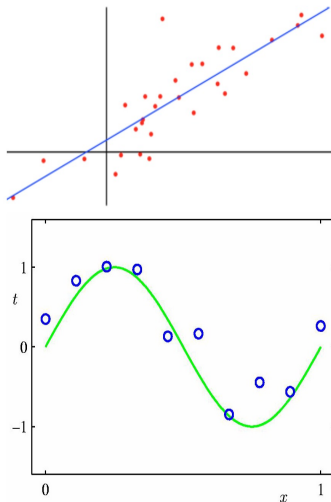
With non-linear basis functions ϕ_j

$$y(x, w) = w_0 + \sum_{j=1}^{M-1} w_j \phi_j(x)$$

- **Loss Function**

$$L(w) = \frac{1}{2} \sum_{n=1}^N \{y(x_n, w) - t_n\}^2 + \frac{\lambda}{2} \|w\|^2$$

Second term is a weighted norm called a **regularizer** (to prevent overfitting)

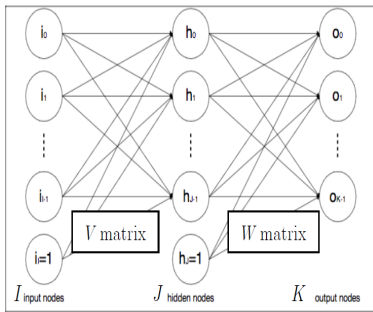


- Norm is the length of a vector
- Distance between two vectors (v, w)
 - $dist(v, w) = \|v - w\| = \sqrt{(v_1 - w_1)^2 + \dots + (v_n - w_n)^2}$
Distance to origin would be sqrt of sum of squares
- Similar to L^2 norm

$$\|A\|_F = \left(\sum_{i,j} A_{i,j}^2 \right)^{\frac{1}{2}}$$

- Frobenius in ML
 - Layers of neural network involve matrix multiplication
 - **Regularization:**
minimize Frobenius of weight matrices $\|W_{(i)}\|$ over L layers

$$J_R = J + \lambda \sum_{i=1}^L \|W^{(i)}\|_F$$



$$I_{1 \times (I+1)} \times V_{(I+1) \times J} = \text{net}_J$$

$$h_j = f(\text{net}_j) \quad f(x) = 1 / (1 + e^{-x})$$

Fig: Matrix Multiplication in Neural Net.

LINEAR ALGEBRA TO IMAGE PROCESSING

What are eigenvalues?

- Given a matrix, A , x is the eigenvector and λ is the corresponding eigenvalue if $Ax = \lambda x$.
 - A must be a square matrix, the determinant of $A - \lambda I$ must be equal to zero.

$$Ax - \lambda x = 0 \implies (A - \lambda I)x = 0$$

- Trivial solution is if $x = 0$.
 - The non trivial solution occurs when $\det(A - \lambda I) = 0$.
- Are eigenvectors unique?
 - If x is an eigenvector, then βx is also an eigenvector and $\beta\lambda$ is an eigenvalue

$$A(\beta x) = \beta(Ax) = \beta(\lambda x) = \lambda(\beta x)$$

- Expand the $\det(A - \lambda I) = 0$ for a 2×2 matrix

$$\det(A - \lambda I) = \det \left(\begin{bmatrix} a_{11} & a_{12} \\ a_{21} & a_{22} \end{bmatrix} - \lambda \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \right) = 0$$

$$\det \begin{bmatrix} a_{11} - \lambda & a_{12} \\ a_{21} & a_{22} - \lambda \end{bmatrix} = 0 \implies (a_{11} - \lambda)(a_{22} - \lambda) - a_{12}a_{21} = 0$$

$$\lambda^2 - \lambda(a_{11} + a_{22}) + (a_{11}a_{22} - a_{12}a_{21}) = 0$$

- For a 2×2 matrix, this is a simple quadratic equation with two solutions (maybe complex).
- The “characteristic equation” can be used to solve for λ .

- Consider,

$$A = \begin{bmatrix} 1 & 2 \\ 2 & 4 \end{bmatrix} \implies \begin{cases} \lambda^2 - (a_{11} + a_{22})\lambda + (a_{11}a_{22} - a_{12}a_{21}) = 0 \\ \lambda^2 - (1 + 4)\lambda + (1 * 4 - 2 * 2) = 0 \\ \lambda^2 = (1 + 4)\lambda \implies \lambda = 0, \lambda = 5 \end{cases}$$

- The corresponding eigenvectors can be computed as

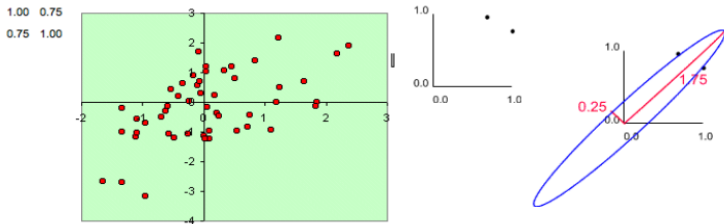
$$\lambda = 0 \implies \left[\begin{bmatrix} 1 & 2 \\ 2 & 4 \end{bmatrix} - \begin{bmatrix} 0 & 0 \\ 0 & 0 \end{bmatrix} \right] \cdot \begin{bmatrix} x \\ y \end{bmatrix} = 0 \implies \begin{bmatrix} 1 & 2 \\ 2 & 4 \end{bmatrix} \cdot \begin{bmatrix} x \\ y \end{bmatrix} = \begin{bmatrix} 1x + 2y \\ 2x + 4y \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$$

$$\lambda = 5 \implies \left[\begin{bmatrix} 1 & 2 \\ 2 & 4 \end{bmatrix} - \begin{bmatrix} 5 & 0 \\ 0 & 5 \end{bmatrix} \right] \cdot \begin{bmatrix} x \\ y \end{bmatrix} = 0 \implies \begin{bmatrix} -4 & 2 \\ 2 & -1 \end{bmatrix} \cdot \begin{bmatrix} x \\ y \end{bmatrix} = \begin{bmatrix} -4x + 2y \\ 2x - 1y \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$$

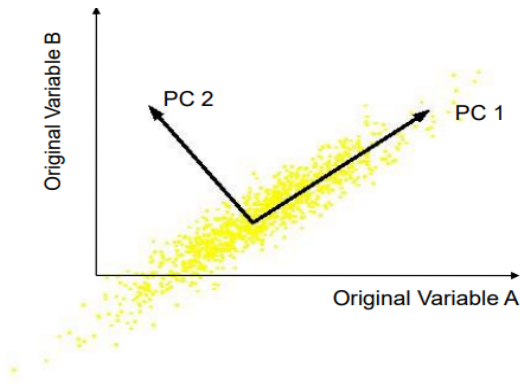
- For $\lambda = 0$, one possible solution is $x = (2, -1)$
- For $\lambda = 5$, one possible solution is $x = (1, 2)$

- Consider a correlation matrix, A

$$A = \begin{bmatrix} 1 & 0.75 \\ 0.75 & 1 \end{bmatrix} \implies \lambda_1 = 1.75, \lambda_2 = 0.25$$



- Error ellipse with the major axis as the larger eigenvalue and the minor axis as the smaller eigenvalue.



- Orthogonal directions of greatest variance in data.
- Projections along PC1 (Principal Component) discriminate the data most along any one axis.

- First principal component is the direction of greatest variability (covariance) in the data.
- Second is the next orthogonal (uncorrelated) direction of greatest variability.
 - So first remove all the variability along the first component, and then find the next direction of greatest variability.
- And so on ...
- Thus each eigenvectors provides the directions of data variances in decreasing order of eigenvalues.

- If A is symmetric and positive definite $k \times k$ matrix ($x^T A x > 0$) with λ_i ($\lambda_i > 0$) and $e_i, i = 1, 2, \dots, k$ being the k eigenvalues and eigenvectors pairs, then

$$A_{(k \times k)} = \lambda_1 \begin{matrix} e_1 & e_1^T \\ (k \times 1) & (1 \times k) \end{matrix} + \lambda_2 \begin{matrix} e_2 & e_2^T \\ (k \times 1) & (1 \times k) \end{matrix} + \dots + \lambda_k \begin{matrix} e_k & e_k^T \\ (k \times 1) & (1 \times k) \end{matrix} \implies A = \sum_{i=1}^k \lambda_i \begin{matrix} e_i & e_i^T \\ (k \times 1) & (1 \times k) \end{matrix} = P \Lambda P^T$$

$$P_{k \times k} = [e_1, e_2, \dots, e_k]; \quad \Lambda_{k \times k} = \begin{bmatrix} \lambda_1 & 0 & \dots & 0 \\ 0 & \lambda_2 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & \lambda_k \end{bmatrix}$$

- This is also called the eigen decomposition theorem.
- Any symmetric matrix can be reconstructed using its eigenvalues and eigenvectors.

- Let A be a symmetric, positive definite matrix

$$A = \begin{bmatrix} 2.2 & 0.4 \\ 0.4 & 2.8 \end{bmatrix} \implies \det(A - \lambda I) = 0$$

$$\implies \lambda^2 - 5\lambda + (6.16 - 0.16) = (\lambda - 3)(\lambda - 2) = 0$$

- The eigenvectors for the corresponding eigenvalues are $e_1^T = [\frac{1}{\sqrt{5}}, \frac{2}{\sqrt{5}}]$, $e_2^T = [\frac{2}{\sqrt{5}}, \frac{-1}{\sqrt{5}}]$.
- Consequently,

$$\begin{aligned} A &= \begin{bmatrix} 2.2 & 0.4 \\ 0.4 & 2.8 \end{bmatrix} = 3 \begin{bmatrix} \frac{1}{\sqrt{5}} \\ \frac{2}{\sqrt{5}} \end{bmatrix} \begin{bmatrix} \frac{1}{\sqrt{5}} & \frac{2}{\sqrt{5}} \end{bmatrix} + 2 \begin{bmatrix} \frac{2}{\sqrt{5}} \\ \frac{-1}{\sqrt{5}} \end{bmatrix} \begin{bmatrix} \frac{2}{\sqrt{5}} & \frac{-1}{\sqrt{5}} \end{bmatrix} \\ &= \begin{bmatrix} 0.6 & 1.2 \\ 1.2 & 2.4 \end{bmatrix} + \begin{bmatrix} 1.6 & -0.8 \\ -0.8 & 0.4 \end{bmatrix} \end{aligned}$$

- Eigendecomposition is $A = P\Lambda P^T$
 - where P is an orthogonal matrix composed of eigenvectors of A
- Decomposition is not unique when two eigenvalues are the same.
- Eigenvalues really work for square matrices. Consider a regression problem, we have input and output matrix and m and k have different dimensions. Then, this method fails. [Solution: SVD.](#)
- By convention order entries of Λ in descending order:
 - Under this convention, eigendecomposition is unique if all eigenvalues are unique

- If A is rectangular $m \times k$ matrix of real numbers, then there exists an $m \times m$ orthogonal matrix U and a $k \times k$ orthogonal matrix V such that

$$\underset{(m \times k)}{A} = \underset{(m \times m)}{U} \underset{(m \times k)}{\Lambda} \underset{(k \times k)}{V^T} \quad UU^T = VV^T = I$$

- Λ is an $m \times k$ matrix where the $(i, j)^{th}$ entry $\lambda_i, i = j = 1, 2, \dots, \min(m, k)$ and the other entries are zero. Physically, A equals rotation \times stretching \times rotation.
 - The positive constants λ_i are the singular values of A .
- If A has rank r , then there exists r positive constants $\lambda_1, \lambda_2, \dots, \lambda_r$; r orthogonal $m \times 1$ unit vectors u_1, u_2, \dots, u_r and r orthogonal $k \times 1$ unit vectors v_1, v_2, \dots, v_r such that

$$A = \sum_{i=1}^r \lambda_i u_i v_i^T$$

- If A is symmetric and positive definite then
 - SVD = Eigen decomposition
 - $\text{EIG}(\lambda_i) = \text{SVD}(\lambda_i^2)$
- Here AA^T has an eigenvalue-eigenvector pair (λ_i^2, u_i)

$$\begin{aligned} AA^T &= (U\Lambda V^T)(U\Lambda V^T)^T \\ &= U\Lambda V^T V \Lambda U^T \\ &= U\Lambda^2 U^T \end{aligned}$$

- Alternatively, the v_i are the eigenvectors of $A^T A$ with same non zero eigenvalue λ_i^2

$$A^T A = V\Lambda^2 V^T$$

- Let A be a symmetric, positive definite matrix
 - U can be computed as

$$A = \begin{bmatrix} 3 & 1 & 1 \\ -1 & 3 & 1 \end{bmatrix} \implies AA^T = \begin{bmatrix} 3 & 1 & 1 \\ -1 & 3 & 1 \end{bmatrix} \begin{bmatrix} 3 & -1 \\ 1 & 3 \\ 1 & 1 \end{bmatrix} = \begin{bmatrix} 11 & 1 \\ 1 & 11 \end{bmatrix}$$

$$\det(AA^T - \gamma I) = 0 \implies \gamma_1 = 12, \gamma_2 = 10 \implies u_1^T = \left[\frac{1}{\sqrt{2}}, \frac{1}{\sqrt{2}} \right], u_2^T = \left[\frac{1}{\sqrt{2}}, -\frac{1}{\sqrt{2}} \right]$$

- V can be computed as

$$A = \begin{bmatrix} 3 & 1 & 1 \\ -1 & 3 & 1 \end{bmatrix} \implies A^T A = \begin{bmatrix} 3 & -1 \\ 1 & 3 \\ 1 & 1 \end{bmatrix} \begin{bmatrix} 3 & 1 & 1 \\ -1 & 3 & 1 \end{bmatrix} = \begin{bmatrix} 10 & 0 & 2 \\ 0 & 10 & 4 \\ 2 & 4 & 2 \end{bmatrix}$$

$$\det(A^T A - \gamma I) = 0 \implies \gamma_1 = 12, \gamma_2 = 10, \gamma_3 = 0$$

$$\implies v_1^T = \left[\frac{1}{\sqrt{6}}, \frac{2}{\sqrt{6}}, \frac{1}{\sqrt{6}} \right], v_2^T = \left[\frac{2}{\sqrt{5}}, \frac{-1}{\sqrt{5}}, 0 \right], v_3^T = \left[\frac{1}{\sqrt{30}}, \frac{2}{\sqrt{30}}, \frac{-5}{\sqrt{30}} \right]$$

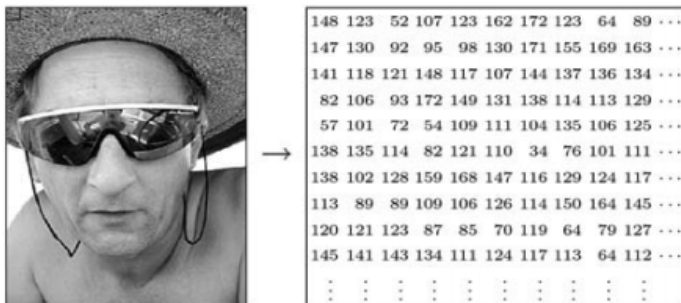
- Taking $\lambda_1^2 = 12$ and $\lambda_2^2 = 10$, the singular value decomposition of A is

$$\begin{aligned}
 A &= \begin{bmatrix} 3 & 1 & 1 \\ -1 & 3 & 1 \end{bmatrix} \\
 &= \sqrt{12} \begin{bmatrix} \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{bmatrix} \left[\frac{1}{\sqrt{6}}, \frac{2}{\sqrt{6}}, \frac{1}{\sqrt{6}} \right] + \sqrt{10} \begin{bmatrix} \frac{1}{\sqrt{2}} \\ -\frac{1}{\sqrt{2}} \end{bmatrix} \left[\frac{2}{\sqrt{5}}, \frac{-1}{\sqrt{5}}, 0 \right]
 \end{aligned}$$

- Thus, the U , V and Λ are computed by performing eigendecomposition of AA^T and A^TA .
- Any matrix has a singular value decomposition but only symmetric, positive definite matrices have an eigen decomposition.

- SVD can be used to compute optimal low-rank approximations of arbitrary matrices.
- Face recognition
 - Represent the face images as eigenfaces and compute distance between the query face image in the principal component space.
- Data mining
 - Latent Semantic Indexing for document extraction.
- Image Compression

A digital image is a representation of a real image as a set of numbers that can be stored and handled by digital computers.



The $(i, j)^{th}$ entry of the matrix comprises of $(i, j)^{th}$ pixel value, which determines the intensity of light of the image.

- **Image Size:** Number of rows * Number of columns.



- **Image Resolution:** Area covered by per pixel.



- **Number of Color Planes:** Grayscale image(1), RGB image(3).



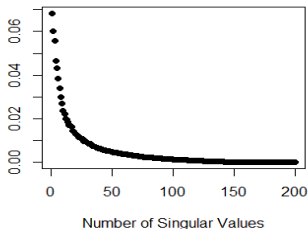
- An image is stored as a 200×200 matrix M with entries between 0 and 1. The matrix M has rank 200.
- Select $r > 200$ as an approximation to the original M .
 - As r is increased from 1 all the way to 200 the reconstruction of M would improve i.e. approximation error would reduce
- Advantage
 - To send the matrix M , need to send $200 \times 200 = 40000$ numbers.
 - To send an $r = 35$ approximation of M , need to send $35 + 35 * 200 + 35 * 200 = 14035$ numbers
 - 35 singular values.
 - 35 left vectors, each having 200 entries.
 - 35 right vectors, each having 200 entries.



Data and Code is available at <https://github.com/mad-stat/SVD-Applications>

- Obtain the singular values (Λ) and the singular vectors (U, V) of the image using SVD.
- The amount of variance explained by the singular values can be obtained by plotting $\Lambda^2 / \text{sum}(\Lambda^2)$.

Variance explained by singular values



- Most of the variance is explained by first 35 singular vectors in this image.
- After 35 vectors the variance is very low and almost steady. After about 75 vectors it's miniscule. Hence, we can compress the image without losing much of the quality by retaining the first 75 vectors and discarding others.

Compressed image

Rank= 5



Rank= 10



Rank= 20



Rank= 35



Rank= 50



Rank= 75



Rank= 100



Rank= 150



Original image



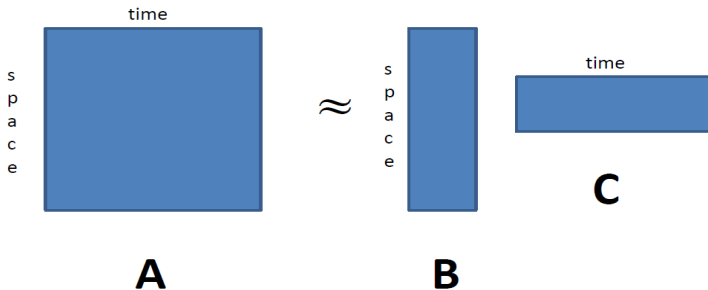
LINEAR ALGEBRA TO SPATIO-TEMPORAL DATA ANALYSIS

- Assume three matrices **A**, **B**, and **C**
- Consider equation

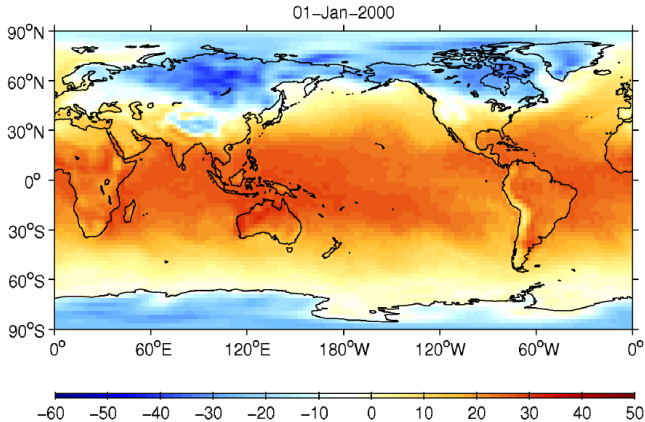
$$A = BC$$

- If any two matrices are known, the third one can be solved
- But let us consider the case when only one (say, **A**) is known.
- Then $A \cong BC$ is called a matrix decomposition for **A**.
- Some very promising machine learning techniques are based on this.

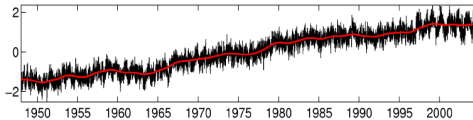
- Graphically, the situation may be like this:



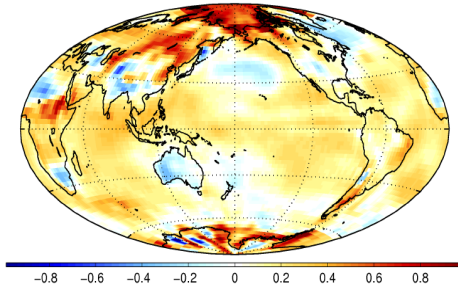
Global daily temperature



Ref: Ilin, Valpola, Oja, Neural Networks (2006).



One row of
matrix C



Correspond
ing column
of matrix B

Ref: Ilin, Valpola, Oja, Neural Networks (2006).

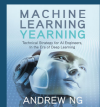
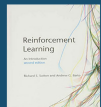
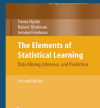
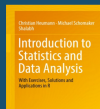
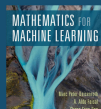
- Consider a data matrix $\mathbf{X} \in \mathbb{R}^{n \times m}$ whose m columns (or n rows) contain data vectors (signals, images, word histograms, movie ratings, distances...).
- Linear latent variable model:

$$\mathbf{X} \approx \mathbf{WH}$$

with latent (hidden) variables or sources $\mathbf{H} \in \mathbb{R}^{r \times m}$ and a weight matrix $\mathbf{W} \in \mathbb{R}^{n \times r}$.

- Typically, $r < m, n$ for compression and feature extraction, so an exact solution is not possible if \mathbf{X} is full rank.
- **Task:** To discover the "optimal" matrices \mathbf{H} and \mathbf{W} given only the observations \mathbf{X} .

Data Science, Statistics & ML Booklist



Prepared by Dr. Tanujit Chakraborty