ECE 417 Lecture 5: Eigenvectors

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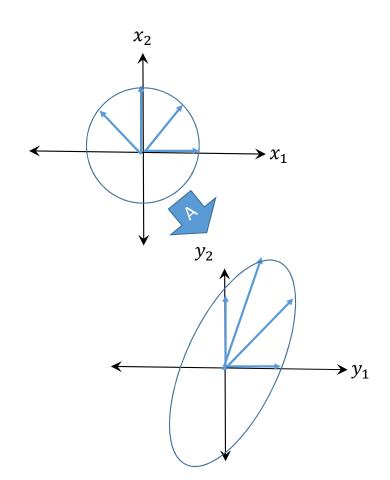
Linear Transforms

A linear transform $\vec{y} = A\vec{x}$ maps vector space \vec{x} onto vector space \vec{y} . For example: the matrix $A = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix}$ maps the vectors

$$\overrightarrow{x_1}, \overrightarrow{x_2}, \overrightarrow{x_3}, \overrightarrow{x_4} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, \begin{bmatrix} \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{bmatrix}, \begin{bmatrix} 0 \\ 1 \end{bmatrix}, \begin{bmatrix} -\frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{bmatrix}$$

to the vectors

$$\overrightarrow{y_1}, \overrightarrow{y_2}, \overrightarrow{y_3}, \overrightarrow{y_4} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, \begin{bmatrix} \sqrt{2} \\ \sqrt{2} \end{bmatrix}, \begin{bmatrix} 1 \\ 2 \end{bmatrix}, \begin{bmatrix} 0 \\ \sqrt{2} \end{bmatrix}$$



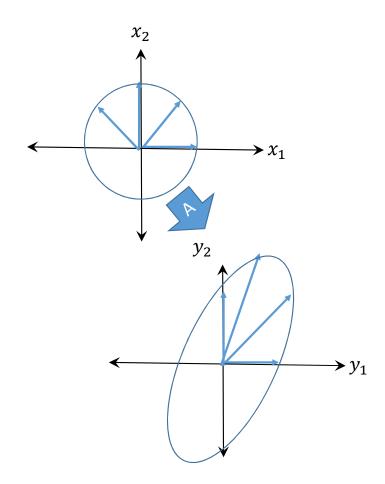
Linear Transforms

A linear transform $\vec{y} = A\vec{x}$ maps vector space \vec{x} onto vector space \vec{y} . For example: the matrix $A = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix}$ maps the vectors

$$X = \begin{bmatrix} 1 & \frac{1}{\sqrt{2}} & 0 & -\frac{1}{\sqrt{2}} \\ 0 & \frac{1}{\sqrt{2}} & 1 & \frac{1}{\sqrt{2}} \end{bmatrix}$$

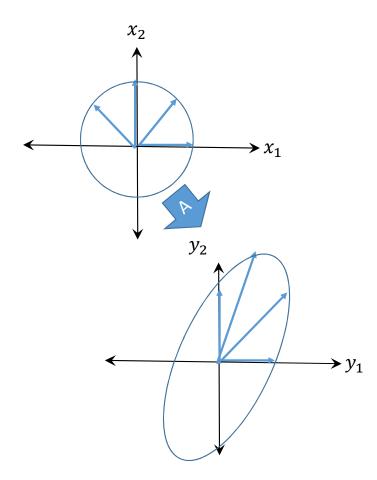
to the vectors

$$Y = \begin{bmatrix} 1 & \sqrt{2} & 1 & 0 \\ 0 & \sqrt{2} & 2 & \sqrt{2} \end{bmatrix}$$



Linear Transforms

A linear transform $\vec{y} = A\vec{x}$ maps vector space \vec{x} onto vector space \vec{y} . The determinant of A tells you how much the area of a unit circle is changed under the transformation. For example: if $A = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix}$, then the unit circle in \vec{x} (which has an area of π) is mapped to an ellipse with an area of $\pi|A| = 2\pi$.



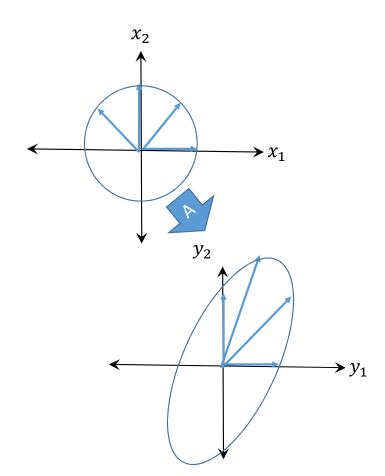
Eigenvectors

• For a D-dimensional square matrix, there may be up to D different directions $\vec{x} = \overrightarrow{v_d}$ such that, for some scalar λ_d ,

$$A\overrightarrow{v_d} = \lambda_d \overrightarrow{v_d}$$

• For example: if $A = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix}$, then the eigenvectors and eigenvalues are

$$\overrightarrow{v_1} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, \overrightarrow{v_2} = \begin{bmatrix} \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{bmatrix}, \lambda_1 = 1, \lambda_2 = 2$$

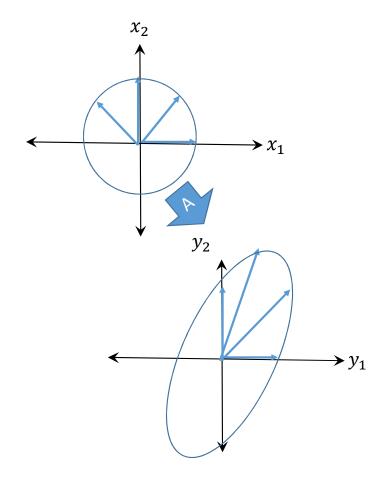


Eigenvectors

- An eigenvector is a direction, not just a vector. That means that if you multiply an eigenvector by any scalar, you get the same eigenvector: if $A\overrightarrow{v_d} = \lambda_d \overrightarrow{v_d}$, then it's also true that $cA\overrightarrow{v_d} = c\lambda_d \overrightarrow{v_d}$
- For example: the following are all the same eigenvector

$$\overrightarrow{v_2} = \begin{bmatrix} \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} \end{bmatrix}, \sqrt{2}\overrightarrow{v_2} = \begin{bmatrix} 1 \\ 1 \end{bmatrix}, -\overrightarrow{v_2} = \begin{bmatrix} -\frac{1}{\sqrt{2}} \\ -\frac{1}{\sqrt{2}} \end{bmatrix}$$

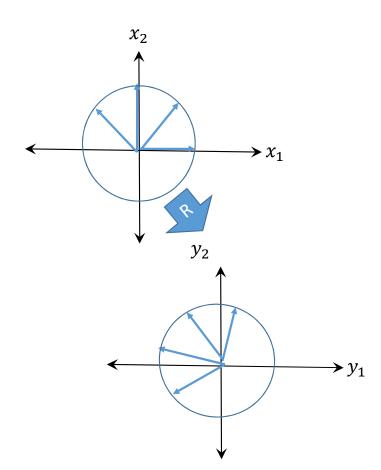
• Since scale doesn't matter, by convention, we normalize so that $||\overrightarrow{v_d}||_2=1$ and the first nonzero element is positive.



Eigenvectors

- Notice that only square matrices can have eigenvectors. For a non-square matrix, the equation $A\overrightarrow{v_d}=\lambda_d\overrightarrow{v_d}$ is impossible --- the dimension of the output is different from the dimension of the input.
- Not all matrices have eigenvectors!
 For example, a rotation matrix doesn't have any real-valued eigenvectors:

$$R = \begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix}$$



Eigenvalues

$$A\overrightarrow{v_d} = \lambda_d \overrightarrow{v_d}$$

$$A\overrightarrow{v_d} = \lambda_d I \overrightarrow{v_d}$$

$$A\overrightarrow{v_d} - \lambda_d I \overrightarrow{v_d} = \overrightarrow{0}$$

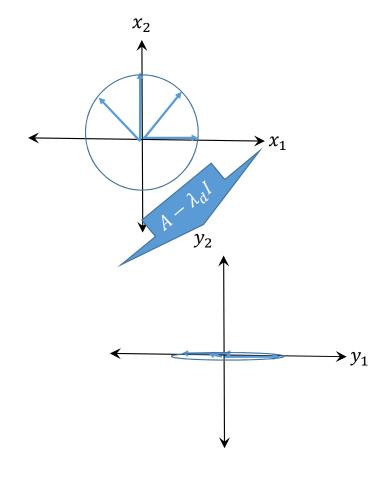
$$(A - \lambda_d I) \overrightarrow{v_d} = \overrightarrow{0}$$

That means that when you use the linear transform $(A-\lambda_d I)$ to transform the unit circle, the result has zero area. Remember that the area of the output is $\pi |A-\lambda_d I|$. So that means that, for any eigenvalue λ_d , the determinant of the matrix difference is zero:

$$|A - \lambda_d I| = 0$$

Example:

$$A - \lambda_2 I = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix} - 2 \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} = \begin{bmatrix} -1 & 1 \\ 0 & 0 \end{bmatrix}$$



Eigenvalues

Let's talk about that equation, $|A-\lambda_d I|=0$. Remember how the determinant is calculated, for example if

$$A = \begin{bmatrix} a & b & c \\ d & e & f \\ g & h & i \end{bmatrix}, \text{ then } |A - \lambda I| = 0 \text{ means that}$$

$$0 = |A - \lambda I| = \begin{vmatrix} a - \lambda & b & c \\ d & e - \lambda & f \\ g & h & i - \lambda \end{vmatrix} =$$

$$(a - \lambda)(e - \lambda)(i - \lambda) - b(d(i - \lambda) - gf) + c(dh - g(e - \lambda))$$

- We assume that a,b,c,d,e,f,g,h,i are all given in the problem statement. Only λ is unknown. So the equation $|A-\lambda I|=0$ is a D'th order polynomial in one variable.
- The fundamental theorem of algebra says that a D'th order polynomial has D roots (counting repeated roots and complex roots).

Eigenvalues

So a DxD matrix always has D eigenvalues (counting complex and repeated eigenvalues). This is true even if the matrix has no eigenvectors!! The eigenvalues are the D solutions of the polynomial equation

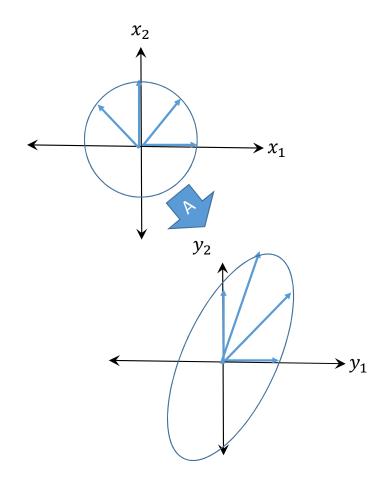
$$|A - \lambda_d I| = 0$$

Positive Definite Matrix

• A linear transform $\vec{y} = A\vec{x}$ is called "positive definite" (written A > 0) if, for any vector \vec{x} ,

 $\vec{x}^T A \vec{x} > 0$

- So, you can see that this means $\vec{x}^T \vec{y} > 0$.
- So this means that a matrix is positive definite if and only if the output of the transform, \vec{y} , is never rotated away from the input, \vec{x} , by 90 degrees or more! \leftarrow (useful geometric intuition)
- For example, the matrix $A = \begin{bmatrix} 1 & 1 \\ 0 & 2 \end{bmatrix}$ is positive-definite.



Symmetric matrices

We've been working with "right eigenvectors:"

$$A\overrightarrow{v_d} = \lambda_d \overrightarrow{v_d}$$

There may also be left eigenvectors, which are row vectors \vec{u}_d^T , and corresponding left eigenvalues μ_d :

$$\vec{u}_d^T \vec{A} = \mu_d \vec{u}_d^T$$

If A is symmetric ($A = A^T$), then the left and right eigenvectors and eigenvalues are the same, because

$$\lambda_d \vec{v}_d^T = (\lambda_d \overrightarrow{v_d})^T = (A \overrightarrow{v_d})^T = \vec{v}_d^T A^T = \vec{v}_d^T A = \mu_d \vec{u}_d^T$$

If A is symmetric ($A = A^T$), then you can do an interesting thing if you multiply the matrix by its eigenvectors both before and after:

$$\vec{v}_d^T A \vec{v}_d = \vec{v}_d^T (\lambda_d \vec{v}_d) = \lambda_d ||\vec{v}_d||_2^2 = \lambda_d$$

So if a symmetric matrix is positive definite, then all of its eigenvalues are positive real numbers. It turns out that the opposite is also true:

A symmetric matrix is positive definite if and only if all of its eigenvalues are positive.

Symmetric positive definite matrices turn out to also have one more unbelievably useful property: their eigenvectors are orthogonal.

$$\vec{v}_i^T \vec{v}_i = 0 \text{ if } i \neq j$$

If i = j then, by convention, we have

$$\vec{v}_i^T \vec{v}_i = \|\vec{v}\|_2^2 = 1$$

So suppose we create the matrix

$$V = [\vec{v}_1, \vec{v}_2, \dots, \vec{v}_D]$$

This is an orthonormal matrix:

$$V^TV = I$$

It turns out that, also, $VV^T = I$.

If A is symmetric ($A = A^T$), then

$$\vec{v}_d^T A \vec{v}_d = \vec{v}_d^T (\lambda_d \vec{v}_d) = \lambda_d ||\vec{v}_d||_2^2 = \lambda_d$$

...but also...

$$\vec{v}_i^T \vec{v}_j = \begin{cases} 1, i = j \\ 0, i \neq j \end{cases}$$

That means we can write A as

$$A = \sum_{i=1}^{D} \lambda_i \vec{v}_i \vec{v}_i^T = V \Lambda V^T$$

Because

$$\vec{v}_j^T A \vec{v}_j = \sum_{i=1}^D \lambda_i \vec{v}_j^T \vec{v}_i \vec{v}_i^T \vec{v}_j = \lambda_j$$

If A is symmetric and positive definite we can write

$$A = \sum_{i=1}^{D} \lambda_i \vec{v}_i \vec{v}_i^T = V \Lambda V^T$$

Equivalently

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

Suppose we have a dataset containing N independent sample vectors, \vec{x}_n . The true mean is approximately given by the sample mean,

$$\vec{\mu} = E[\vec{x}] \approx \frac{1}{N} \sum_{n=1}^{N} \vec{x}_n$$

Similarly, the true covariance matrix is approximately given by the sample covariance matrix,

$$\Sigma = E[(\vec{x} - \vec{\mu})(\vec{x} - \vec{\mu})^T] \approx \frac{1}{N} \sum_{n=1}^{N} (\vec{x}_n - \vec{\mu})(\vec{x}_n - \vec{\mu})^T$$

Define the "sum-of-squares matrix" to be

$$S = \sum_{n=1}^{N} (\vec{x}_n - \vec{\mu})(\vec{x}_n - \vec{\mu})^T$$

So that the sample covariance is $\Sigma \approx S/N$. Suppose that we define the centered data matrix to be the following DxN matrix:

$$\tilde{X} = [\vec{x}_1 - \vec{\mu}, \vec{x}_2 - \vec{\mu}, ..., \vec{x}_N - \vec{\mu}]$$

Then the sum-of-squares matrix is

$$S = \tilde{X}\tilde{X}^T = \left[\vec{x}_1 - \vec{\mu}, \dots, \vec{x}_N - \vec{\mu}\right] \begin{bmatrix} (\vec{x}_1 - \vec{\mu})^T \\ \dots \\ (\vec{x}_N - \vec{\mu})^T \end{bmatrix}$$

Well, a sum-of-squares matrix is obviously symmetric. It's also almost always positive definite:

$$\vec{x}^T S \vec{x} = [\vec{x}^T (\vec{x}_1 - \vec{\mu}), \dots, \vec{x}^T (\vec{x}_N - \vec{\mu})] \begin{bmatrix} (\vec{x}_1 - \vec{\mu})^T \vec{x} \\ \dots \\ (\vec{x}_N - \vec{\mu})^T \vec{x} \end{bmatrix}$$

That quantity is positive unless the new vector, \vec{x} , is orthogonal to $(\vec{x}_n - \vec{\mu})$ for every vector in the training database. As long as $N \geq D$, that's really, really unlikely.

So a sum-of-squares matrix can be written as

$$S = \sum_{i=1}^{D} \lambda_i \vec{v}_i \vec{v}_i^T = V \Lambda V^T$$

And the covariance can be written as

$$\Sigma = \frac{S}{N} = \frac{1}{N} \sum_{i=1}^{D} \lambda_i \vec{v}_i \vec{v}_i^T = V\left(\frac{\Lambda}{N}\right) V^T$$

Principal components

Suppose that

$$\Lambda = \begin{bmatrix} \lambda_1 & 0 & 0 \\ 0 & \dots & 0 \\ 0 & 0 & \lambda_D \end{bmatrix}, V = [\vec{v}_1, \dots, \vec{v}_D]$$

are the eigenvalue and eigenvector matrices of S, respectively. Define the principal components of \vec{x}_n to be $y_{dn} = \vec{v}_d^T (\vec{x}_n - \vec{\mu})$, or

$$\vec{y}_n = V^T(\vec{x}_n - \vec{\mu}) = \begin{bmatrix} \vec{v}_1^T(\vec{x}_n - \vec{\mu}) \\ ... \\ \vec{v}_D^T(\vec{x}_n - \vec{\mu}) \end{bmatrix}$$

Principal components

Suppose that Λ and V are the eigenvalue and eigenvector matrices of S, respectively. Define the principal components to be $\vec{y}_n = V^T(\vec{x}_n - \vec{\mu})$.

Then the principal components y_{dn} are not correlated with each other, and the

variance of each one is given by the corresponding eigenvalue of S.
$$E[\vec{y}\vec{y}^T] \approx \frac{1}{N} \sum_{n=1}^{N} \vec{y}_n \vec{y}_n^T = \frac{1}{N} \sum_{n=1}^{N} \begin{bmatrix} y_{1n} \\ ... \\ y_{Dn} \end{bmatrix} [y_{1n}, ..., y_{Dn}]$$

$$= \frac{1}{N} \sum_{n=1}^{N} V^{T} (\vec{x}_{n} - \vec{\mu}) (\vec{x}_{n} - \vec{\mu})^{T} V$$

$$= V^{T} S V = \Lambda = \begin{bmatrix} \lambda_{1} & 0 & 0 \\ 0 & \dots & 0 \\ 0 & 0 & \lambda_{D} \end{bmatrix}$$

Summary

- Principal component directions are the eigenvectors of the covariance matrix (or of the sum-of-squares matrix – same directions, because they are just scaled by N)
- Principal components are the projections of each training example onto the principal component directions
- Principal components are uncorrelated with each other: the covariance is zero
- The variance of each principal component is the corresponding eigenvalue of the covariance matrix

Implications

- The total energy in the signal, $E[\|\vec{x} \vec{\mu}\|_2^2]$, is equal to the sum of the eigenvalues.
- If you want to keep only a small number of dimensions, but keep most of the energy, you can do it by keeping the principal components with the highest corresponding eigenvalues.