CS9840 Learning and Computer Vision Prof. Olga Veksler

Lecture 2

Some Concepts from Computer Vision Curse of Dimensionality

PCA

Some Slides are from Cornelia, Fermüller, Mubarak Shah,

Gary Bradski, Sebastian Thrun

Optical flow





- How to estimate pixel motion from image I₁ to image I₂?
 - Solve pixel correspondence problem
 - given a pixel in I₁, look for nearby pixels of the same color in I₂
- Key assumptions
 - color constancy: a point in I₁ looks the same in I₂
 - For grayscale images, this is brightness constancy
 - small motion: points do not move very far
- This is called the optical flow problem

Outline

- Some Concepts in Image Processing/Vision
 - Optical Flow Field (related to motion field)
 - Correlation
- Curse of Dimensionality and Dimensionality reduction with PCA
- Next time:
 - "Recognizing Action at a Distance" by A. Efros, A.Berg, G. Mori, Jitendra Malik
 - Also: "80 million tiny images: a large dataset for non-parametric object and scene recognition", A. Torralba, R. Fergus, W. Freeman
 - there should be a link to PDF file on our web site

Optical Flow Field

Optical Flow and Motion Field

- Optical flow field is the apparent motion of brightness patterns between 2 (or several) frames in an image sequence
- Why does brightness change between frames?
- Assuming that illumination does not change:
 - changes are due to the RELATIVE MOTION between the scene and the camera
 - There are 3 possibilities:
 - Camera still, moving scene
 - Moving camera, still scene
 - Moving camera, moving scene

Examples of Motion Fields (a) (b) (c) (d)

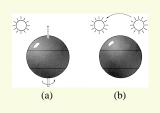
(a) Translation perpendicular to a surface. (b) Rotation about axis perpendicular to image plane. (c) Translation parallel to a surface at a constant distance. (d) Translation parallel to an obstacle in front of a more distant background.

Motion Field (MF)

- The MF assigns a velocity vector to each pixel in the image
- These velocities are INDUCED by the RELATIVE MOTION between the camera and the 3D scene
- The MF is the <u>projection</u> of the 3D velocities on the image plane

Optical Flow vs. Motion Field

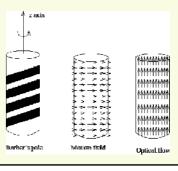
- Recall that Optical Flow is the apparent motion of brightness patterns
- We equate Optical Flow Field with Motion Field
- Frequently works, but now always:



- (a) A smooth sphere is rotating under constant illumination. Thus the optical flow field is zero, but the motion field is not
- (b) A fixed sphere is illuminated by a moving source—the shading of the image changes. Thus the motion field is zero, but the optical flow field is not

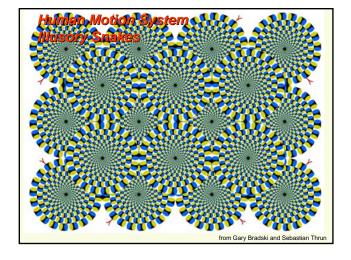
Optical Flow vs. Motion Field

 Often (but not always) optical flow corresponds to the true motion of the scene



Computing Optical Flow: Brightness Constancy Equation

- Let **P** be a moving point in 3D:
 - At time t, P has coordinates (X(t), Y(t), Z(t))
 - Let p=(x(t),y(t)) be the coordinates of its image at time t
 - Let E(x(t), y(t), t) be the brightness at p at time t.
- Brightness Constancy Assumption:
 - As P moves over time, E(x(t), y(t), t) remains constant



Computing Optical Flow: Brightness Constancy Equation

$$E(x(t), y(t), t) = Constant$$

Taking derivative wrt time:

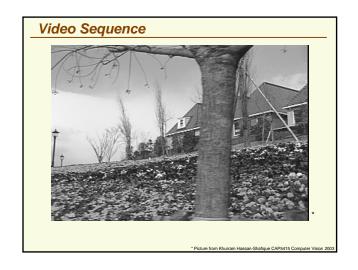
$$\frac{dE(x(t), y(t), t)}{dt} = 0$$

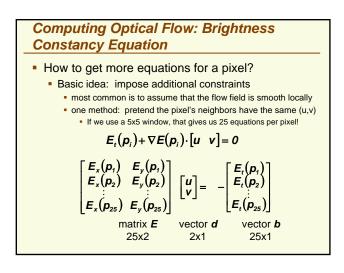
$$\frac{\partial E}{\partial x}\frac{dx}{dt} + \frac{\partial E}{\partial y}\frac{dy}{dt} + \frac{\partial E}{\partial t} = 0$$

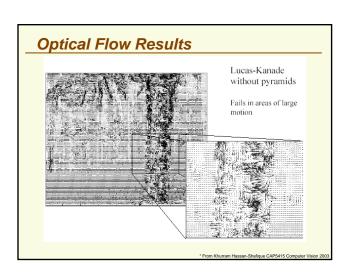
Computing Optical Flow: Brightness Constancy Equation

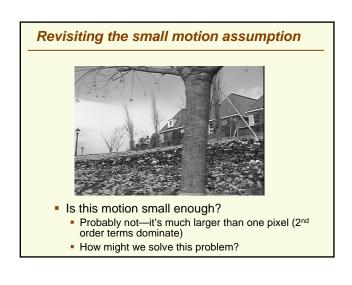
1 equation with 2 unknowns
$$\frac{\partial E}{\partial x} \frac{dx}{dt} + \frac{\partial E}{\partial y} \frac{dy}{dt} + \frac{\partial E}{\partial t} = 0$$
Let
$$\nabla E = \begin{bmatrix} \frac{\partial E}{\partial x} \\ \frac{\partial E}{\partial y} \end{bmatrix} \qquad \text{(Frame spatial gradient)}$$

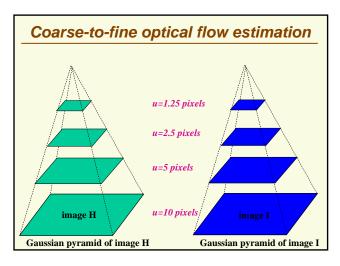
$$v = \begin{bmatrix} \frac{dx}{dt} \\ \frac{dy}{dt} \end{bmatrix} \qquad \text{(optical flow)}$$
and
$$E_t = \frac{\partial E}{\partial t} \qquad \text{(derivative across frames)}$$

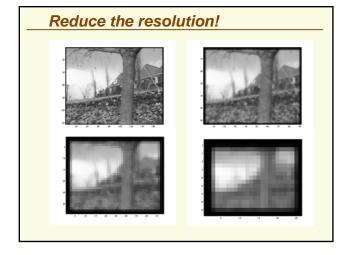






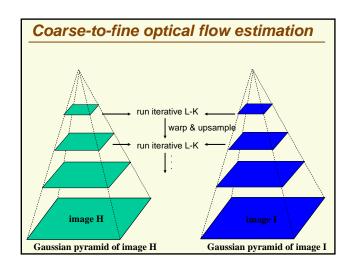


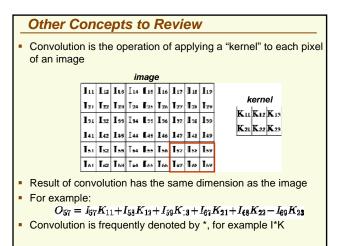


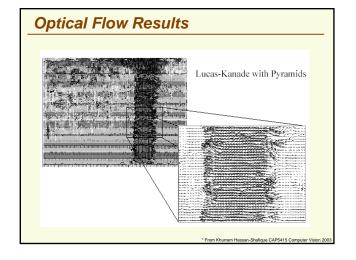


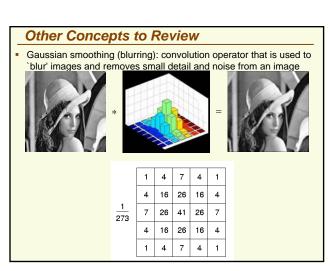
Iterative Refinement

- Iterative Lukas-Kanade Algorithm
 - Estimate velocity at each pixel by solving Lucas-Kanade equations
 - 2. Warp H towards I using the estimated flow field use image warping techniques
 - 3. Repeat until convergence

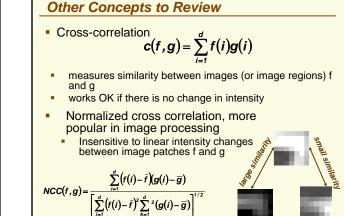


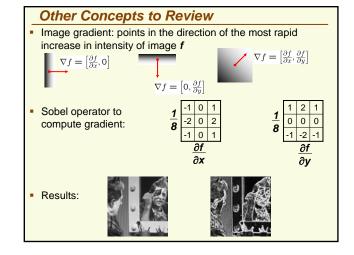










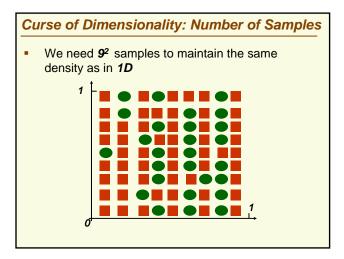


Curse of Dimensionality

- Problems of high dimensional data, "the curse of dimensionality"
 - running time
 - overfitting
 - number of samples required
- Dimensionality Reduction Methods
 - Principle Component Analysis

Curse of Dimensionality: Complexity

- Complexity (running time) increases with dimension d
- A lot of methods have at least O(nd²) complexity, where n is the number of samples
 - For example if we need to estimate covariance matrix
- So as d becomes large, O(nd²) complexity may be too costly



Curse of Dimensionality: Number of Samples

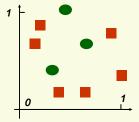
- Suppose we want to use the nearest neighbor approach with k = 1 (1NN)
- Suppose we start with only one feature



- This feature is not discriminative, i.e. it does not separate the classes well
- We decide to use 2 features. For the 1NN method to work well, need a lot of samples, i.e. samples have to be dense
- To maintain the same density as in 1D (9 samples per unit length), how many samples do we need?

Curse of Dimensionality: Number of Samples

 Of course, when we go from 1 feature to 2, no one gives us more samples, we still have 9



This is way too sparse for **1NN** to work well

Curse of Dimensionality: Number of Samples Things go from bad to worse if we decide to use 3 features: 1 0 1

If 9 was dense enough in 1D, in 3D we need 9³=729 samples!

The Curse of Dimensionality

- We should try to avoid creating lot of features
- Often no choice, problem starts with many features
- Example: Face Detection
 - One sample point is **k** by **m** array of pixels



- Feature extraction is not trivial, usually every pixel is taken as a feature
- Typical dimension is 20 by 20 = 400
- Suppose 10 samples are dense enough for 1 dimension. Need only 10⁴⁰⁰ samples

Curse of Dimensionality: Number of Samples

- In general, if n samples is dense enough in 1D
- Then in **d** dimensions we need **n**^d samples!
- And n^d grows really really fast as a function of d
- Common pitfall:
 - If we can't solve a problem with a few features, adding more features seems like a good idea
 - However the number of samples usually stays the same
 - The method with more features is likely to perform worse instead of expected better

The Curse of Dimensionality

• Face Detection, dimension of one sample point is km



- The fact that we set up the problem with km dimensions (features) does not mean it is really a km-dimensional problem
- Space of all k by m images has km dimensions
- Space of all k by m faces must be much smaller, since faces form a tiny fraction of all possible images
- Most likely we are not setting the problem up with the right features
- If we used better features, we are likely need much less than km-dimensions

Dimensionality Reduction

- High dimensionality is challenging and redundant
- It is natural to try to reduce dimensionality
- Reduce dimensionality by feature combination: combine old features x to create new features y

$$x = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_d \end{bmatrix} \rightarrow f \begin{pmatrix} \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_d \end{bmatrix} = \begin{bmatrix} y_1 \\ \vdots \\ y_k \end{bmatrix} = y \quad with \ k < d$$

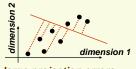
For example,

$$X = \begin{bmatrix} X_1 \\ X_2 \\ X_3 \\ X_4 \end{bmatrix} \rightarrow \begin{bmatrix} X_1 + X_2 \\ X_3 + X_4 \end{bmatrix} = y$$

 Ideally, the new vector y should retain from x all information important for classification

Principle Component Analysis (PCA)

- Main idea: seek most accurate data representation in a lower dimensional space
- Example in 2-D
 - Project data to 1-D subspace (a line) which minimize the projection error





large projection errors, bad line to project to

small projection errors, good line to project to

Notice that the the good line to use for projection lies in the direction of largest variance

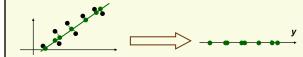
Dimensionality Reduction

- The best f(x) is most likely a non-linear function
- Linear functions are easier to find though
- For now, assume that **f**(**x**) is a linear mapping
- Thus it can be represented by a matrix **W**:

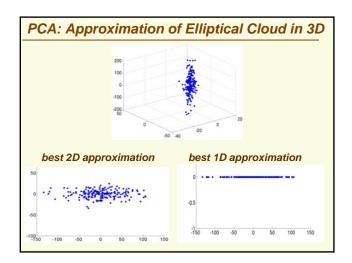
$$\begin{bmatrix} X_1 \\ X_2 \\ \vdots \\ X_d \end{bmatrix} \Rightarrow W \begin{bmatrix} X_1 \\ X_2 \\ \vdots \\ X_d \end{bmatrix} = \begin{bmatrix} W_{11} & \cdots & W_{1d} \\ \vdots & & \vdots \\ W_{k1} & \cdots & W_{kd} \end{bmatrix} \begin{bmatrix} X_1 \\ X_2 \\ \vdots \\ X_d \end{bmatrix} = \begin{bmatrix} y_1 \\ \vdots \\ y_k \end{bmatrix} \quad with \ k < 0$$

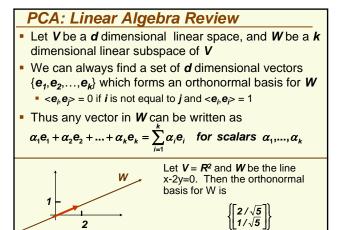
PCA

 After the data is projected on the best line, need to transform the coordinate system to get 1D representation for vector y

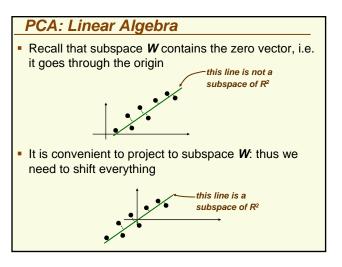


- Note that new data y has the same variance as old data x in the direction of the green line
- PCA preserves largest variances in the data



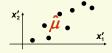


What is the direction of largest variance in data? Recall that if x has multivariate distribution N(μ,Σ), direction of largest variance is given by eigenvector corresponding to the largest eigenvalue of Σ This is a hint that we should be looking at the covariance matrix of the data (note that PCA can be applied to distributions other than Gaussian)



PCA Derivation: Shift by the Mean Vector

- Before PCA, subtract sample mean from the data $x \frac{1}{n} \sum_{i=1}^{n} x_i = x \hat{\mu}$
- The new data has zero mean: E(X-E(X)) = E(X)-E(X) = 0
- All we did is change the coordinate system





- Another way to look at it:
 - first step of getting y is to subtract the mean of x

$$x \rightarrow y = f(x) = g(x - \hat{\mu})$$

PCA: Derivation

- To find the total error, we need to sum over all x_i's
- Any \mathbf{x}_j can be written as $\sum_{i=1}^{n} \alpha_{ji} \mathbf{e}_i$
- Thus the total error for representation of all data D is: sum over all data points

$$J(\mathbf{e}_{1},...,\mathbf{e}_{k},\alpha_{11},...\alpha_{nk}) = \sum_{j=1}^{n} \left\| \mathbf{x}_{j} - \sum_{i=1}^{k} \alpha_{ji} \mathbf{e}_{i} \right\|^{2}$$
unknowns
error at one point

PCA: Derivation

- We want to find the most accurate representation of data $D=\{x_1,x_2,\ldots,x_n\}$ in some subspace W which has dimension k < d
- Let $\{e_1, e_2, ..., e_k\}$ be the orthonormal basis for **W**. Any vector in **W** can be written as $\sum_{i=1}^{\infty} \alpha_i \mathbf{e}_i$
- Thus x_1 will be represented by some vector in W $\sum_{i=1}^k \alpha_{1i} e_i$

$$\sum_{i=1}^{k} \alpha_{1i} \mathbf{e}_{i}$$

• Error this representation:

$$error = \left\| \mathbf{x}_1 - \sum_{i=1}^k \alpha_{1i} \mathbf{e}_i \right\|^2$$



PCA: Derivation

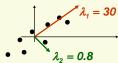
- A lot of math.....to finally get:
- Let S be the scatter matrix, it is just n-1 times the sample covariance matrix

$$\hat{\Sigma} = \frac{1}{n-1} \sum_{j=1}^{n} (x_j - \hat{\mu})(x_j - \hat{\mu})^t$$

To minimize J take for the basis of W the keigenvectors of $\bf S$ corresponding to the $\bf k$ largest eigenvalues

PCA

The larger the eigenvalue of **S**, the larger is the variance in the direction of corresponding eigenvector



- This result is exactly what we expected: project **x** into subspace of dimension **k** which has the largest variance
- This is very intuitive: restrict attention to directions where the scatter is the greatest

PCA as Data Approximation

- Let $\{e_1, e_2, ..., e_d\}$ be all d eigenvectors of the scatter matrix S, sorted in order of decreasing corresponding eigenvalue
- Without any approximation, for any sample x;

$$x_{i} = \sum_{j=1}^{d} \alpha_{j} \mathbf{e}_{j} = \alpha_{1} \mathbf{e}_{1} + \dots + \alpha_{k} \mathbf{e}_{k} + \alpha_{k+1} \mathbf{e}_{k+1} \dots + \alpha_{d} \mathbf{e}_{d}$$

$$approximation of x_{i}$$

$$coefficients \alpha_{m} = x^{t}_{i} \mathbf{e}_{m} \text{ are called } principle \ components$$

- - The larger **k**, the better is the approximation
 - Components are arranged in order of importance, more important components come first
- Thus PCA takes the first **k** most important components of x_i as an approximation to x_i

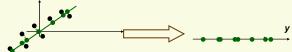
PCA

 Thus PCA can be thought of as finding new orthogonal basis by rotating the old axis until the directions of maximum variance are found



PCA: Last Step

- Now we know how to project the data
- Last step is to change the coordinates to get final **k**-dimensional vector **y**



- Let matrix $\mathbf{E} = [\mathbf{e}_1 \cdots \mathbf{e}_k]$
- Then the coordinate transformation is $y = E^t x$
- Under E^t, the eigenvectors become the standard basis:

$$E^{t}\mathbf{e}_{i} = \begin{bmatrix} \mathbf{e}_{1} \\ \vdots \\ \mathbf{e}_{i} \\ \vdots \\ \mathbf{e}_{i} \end{bmatrix} \mathbf{e}_{i} = \begin{bmatrix} \mathbf{0} \\ \vdots \\ \mathbf{1} \\ \vdots \\ \mathbf{0} \end{bmatrix}$$

Recipe for Dimension Reduction with PCA

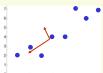
Data $D=\{x_1,x_2,...,x_n\}$. Each x_i is a **d**-dimensional vector. Wish to use PCA to reduce dimension to k

- 1. Find the sample mean $\hat{\mu} = \frac{1}{n} \sum_{i=1}^{n} x_i$
- 2. Subtract sample mean from the data $z_i = x_i \hat{\mu}$
- 3. Compute the scatter matrix $S = \sum_{i=1}^{n} z_i z_i^t$
- 4. Compute eigenvectors $\mathbf{e}_1, \mathbf{e}_2, \dots, \mathbf{e}_k$ corresponding to the k largest eigenvalues of S
- 5. Let $\mathbf{e}_1, \mathbf{e}_2, \dots, \mathbf{e}_k$ be the columns of matrix $\mathbf{E} = [\mathbf{e}_1 \cdots \mathbf{e}_k]$
- 6. The desired y which is the closest approximation to x is $y = E^t z$

PCA Example Using Matlab

Use [V,D] =eig(S) to get eigenvalues and eigenvectors of S

$$\lambda_1 = 87$$
 and $\mathbf{e}_1 = \begin{bmatrix} -0.8 \\ -0.6 \end{bmatrix}$
 $\lambda_2 = 3.8$ and $\mathbf{e}_2 = \begin{bmatrix} 0.6 \\ -0.8 \end{bmatrix}$



Projection to 1D space in the direction of e₁

Y =
$$e_1^t Z^t = \left(\begin{bmatrix} -0.8 - 0.6 \end{bmatrix} \begin{bmatrix} -3.6 & \cdots & 4.4 \\ -4.4 & \cdots & 2.6 \end{bmatrix} \right) = \begin{bmatrix} 4.3 & \cdots & -5.1 \end{bmatrix}$$

= $\begin{bmatrix} y_1 & \cdots & y_8 \end{bmatrix}$

PCA Example Using Matlab

- Let $\mathbf{D} = \{(1,2),(2,3),(3,2),(4,4),(5,4),(6,7),(7,6),(9,7)\}$
- Convenient to arrange data in array

$$X = \begin{bmatrix} 1 & 2 \\ \vdots & \vdots \\ 9 & 7 \end{bmatrix} = \begin{bmatrix} X_1 \\ \vdots \\ X_8 \end{bmatrix}$$

- Mean $\mu = mean(X) = [4.6 \ 4.4]$
- Subtract mean from data to get new data array Z

$$Z = X - \begin{bmatrix} \mu \\ \vdots \\ \mu \end{bmatrix} = X - repmat(\mu, 8, 1) = \begin{bmatrix} -3.6 - 4.4 \\ \vdots & \vdots \\ 4.4 & 2.6 \end{bmatrix}$$

Compute the scatter matrix S

$$S = 7*cov(Z) = \begin{bmatrix} -3.6 - 4.4 \end{bmatrix} \begin{bmatrix} -3.6 \\ -4.4 \end{bmatrix} + ... + \begin{bmatrix} 4.4 & 2.6 \end{bmatrix} \begin{bmatrix} 4.4 \\ 2.6 \end{bmatrix} = \begin{bmatrix} 57 & 40 \\ 40 & 34 \end{bmatrix}$$
 matlab uses unbiased estimate for covariance, so $S=(n-1)*cov(Z)$

Drawbacks of PCA

- PCA was designed for accurate data representation, not for data classification
- Preserves as much variance in data as possible
- If directions of maximum variance is important for classification, will work
- However the directions of maximum variance may be useless for classification







Next Time

- Paper: "Recognizing Action at a Distance" by A. Efros, A.Berg, G. Mori, Jitendra Malik
 - will watch the conference presentation
- Also: "80 million tiny images: a large dataset for nonparametric object and scene recognition", A. Torralba, R. Fergus, W. Freeman
- When reading papers, think about following:
 - What is the problem paper tries to solve
 - What makes this problem difficult?
 - What is the method used in the paper to solve the problem
 - What is the contribution of the paper (what new does it do)?
 - Do the experimental results look "good" to you?