#### Robust Methods in Computer Vision

CS 763 Ajit Rajwade

# Outline

- Least squares estimates
- Limitations of least squares and need for robustness
- Least median of squares method
- RANSAC method
- Application scenarios

- Consider quantity y related to quantity x in the form y = f(x;a).
- Here **a** is a vector of parameters for the function
   f. For example, y = mx + c, where **a** = (m,c).
- Now consider, we have N data points (x<sub>i</sub>,y<sub>i</sub>) where the y<sub>i</sub> could be possibly corrupted by noise, and want to estimate a.
- This is done by minimizing the following w. r. t. **a**:  $\sum_{i=1}^{N} (y_i - f(x_i; \mathbf{a}))^2$

 Why did we minimize the squared error loss? What would happen if we changed the power to 4? Or 1 or 3 (with absolute value)?

Let us assume that the noise affecting y<sub>i</sub> is (1) additive, and (2) Gaussian with mean zero and some known standard deviation σ. Then:

 $y_i = f(x_i; \mathbf{a}) + \eta_i, \eta_i \sim N(0, \sigma)$ 

- Let us also assume that the noise values affecting the different samples are independent of each other.
- Now given some value of **a** and some  $x_i$ , the probability density of  $y_i$  is: Likelihood of  $y_i$  is:

$$p(y_i | x_i, \mathbf{a}) = \frac{1}{\sigma\sqrt{2\pi}} e^{-\frac{(y_i - f(x_i; \mathbf{a}))^2}{2\sigma^2}}$$

$$p(\{y_i\}_{i=1}^N | \{x_i\}_{i=1}^N, \mathbf{a}) = \prod_{i=1}^N p(y_i | x_i, \mathbf{a}) \text{ by independence assumption on noise}$$

• The probability density of a random variable Y at a value y is defined as follows:

$$p_Y(y) = \lim_{\delta \to 0} \frac{P(y \le Y \le y + \delta)}{\delta}$$

- We want to find a value of a which maximizes this probability density. This is called the maximum likelihood estimate of a.
- ..equivalent to maximizing the log of this probability density (why?)
- ..equivalent to minimizing the negative log of this probability density, i.e.

$$J(\mathbf{a}) = \sum_{i=1}^{N} \frac{(y_i - f(x_i; \mathbf{a}))^2}{2\sigma^2} + \log\left(\sigma\sqrt{2\pi}\right)$$

- This shows us that the least squares estimate is the same as the maximum likelihood estimate under the assumption that the noise affecting different samples was *independent* and *Gaussian* distributed with *fixed* variance and mean 0.
- Why maximum likelihood estimate of a?
   Intuitively, it is the value of a that best agrees with or supports the observations {y<sub>i</sub>}, 1 ≤ i ≤ N.

#### Least squares fit of a line

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$$(m^*, c^*) = \arg \min J(m, c) = \sum_{i=1}^{N} (y_i - mx_i - c)^2$$
  

$$\therefore \left(\sum_{i=1}^{N} x_i^2 \sum_{i=1}^{N} x_i\right) \binom{m}{c} = \left(\sum_{i=1}^{N} x_i y_i\right)$$
  

$$\sum_{i=1}^{N} y_i$$
  
N = 1000;x = 2  
y = 1.5\*x + rates  
scatter(x,y,'r+1)

10

0

20

-40

-50 L -40

-30

-20

-10

N = 1000;x = 10\*randn(N,1); y = 1.5\*x + randn(N,1)\*3; scatter(x,y,'r+')



Result of a least-squares estimate under Gaussian noise:

Estimated slope: 1.5015 (versus 1.5) Estimated intercept: 0.088 (versus 0)

# Other Least-squares solutions in computer vision

- Camera calibration SVD
- Parametric motion estimation SVD or pseudo-inverse (affine, rotation, homography, etc)
- Fundamental/essential matrix estimation SVD (we will study this later in stereo-vision)
- Optical Flow (Horn-Shunck as well as Lucas-Kanade) (we will study this later)

#### **Outliers and Least-squares estimates**



**Observation:** Let *x* be a random variable with a Gaussian distribution. Then the probability that *x* takes on any value in a small range far away from the mean (typically at a distance of more than  $+/-3\sigma$ ) is **very low**. See diagram above.

#### **Outliers and Least-squares**

- The upshot of the previous observation is this: the least squares estimate assumes that most points will lie *close* to the true (unknown) model – else their probability would be very low.
- Now, suppose the given dataset contains wild outliers, i.e. stray points that simply do not obey the model.
- These outliers will skew the least squares estimate as it tries to force a solution which maximizes the likelihood of the *outliers* as well.
- Since outliers were extremely unlikely under the Gaussian probability density, the model (during maximum likelihood estimation) has to change itself to make the outliers more likely.



20% of the points are outliers. They have skewed the estimate of the slope from 1.5 to 3.6 and the intercept from 2 to 13.5.

# Examples of outliers: (1)

• Salt and pepper noise in images



Salt and pepper noise (a special case of impulse noise)

Gaussian noise

Don't worry about this example right now – we will encounter it when we study shape from shading and photometric stereo

# Examples of outliers: (2)

Shadows and specularities act as outliers in photometric stereo!



Fig. 7. Input images (4 out of 40). With all kinds of corruptions: Specularity, cast shadow, attached shadow.

Images taken from paper: perception.csl.illinois.edu/matrixrank/Files/robust v19.pdf



#### Examples of outliers: (3)

 Estimation of the spatial transformation between images (could be translation, rotation, affine or homography) requires N+ pairs of corresponding points. Some of these correspondences can be faulty.





#### Examples of outliers: (3)

- Some of these correspondences can be faulty for various reasons
- occlusions/ difference in field of view / shadows
- algorithm errors
- □ identical objects in the scene
- change in the position of some objects in the scene (even though the global motion is homography, these objects will not conform to that same motion model).



http://petapixel.com/2012/10/22/an-eerietime-lapse-of-seattle-minus-all-the-humans/

# Examples of outliers: (4)

 The motion between consecutive frames of a video in the following link may be modeled as affine. But some corresponding pairs of points (example: on independently moving objects) don't conform to that model – those are outliers.

https://www.youtube.com/watch?v=17VAuBL1Lxc



Laplacian probability density function: a distribution which has heavier tails than the Gaussian. This means that if a random variable is Laplacian distributed, the probability that it can take values far away from the mean, is higher than what it would be if the variable were Gaussian distributed.

$$p(x) = \frac{1}{2b} e^{-\frac{|x-\mu|}{b}}$$

The Laplacian (or Laplace) pdf is **not** to be confused with the Laplacian of a function f(x,y), given as  $f_{xx}(x,y) + f_{yy}(x,y)$  that we had studied in image processing.



$$p(x) = \frac{\beta}{2\alpha\Gamma(1/\beta)} e^{-\left(\frac{|x-\mu|}{\alpha}\right)^{\beta}}$$

The Laplacian probability density function is a special case of the family of **Generalized Gaussian probability density functions** with **shape parameter**  $\beta$  and **scale parameter**  $\alpha$ . As  $\beta$  reduces below 1, the density function becomes heavier tailed.

Question: Why do we care for heavier tails? Because they ensure that the wild outliers are more likely to occur (than the Gaussian pdf).

Consequently, a maximum likelihood estimation assuming heavy-tailed noise models will be less affected by outliers.



Note that x^p grows faster than x^q beyond |x| = 1, for p > q Why do these Generalized Gaussians with shape parameter  $\beta < 2$  have heavier tails than the usual Gaussian (i.e.  $\beta = 2$ )?

Consider a zero-mean Gaussian and a zeromean Laplacian (without loss of generality), i.e.  $\beta = 1$ .

The term inside the exponential in a Gaussian is  $x^2$ , where it is |x| for a Laplacian.



- Assume the noise has a Laplacian distribution.
- The maximum likelihood estimate of **a** is then given by minimizing the following:

$$J(\mathbf{a}) = \sum_{i=1}^{N} \frac{|y_i - f(x_i; \mathbf{a})|}{b} - \log(2b) \approx \sum_{i=1}^{N} \frac{\sqrt{(y_i - f(x_i; \mathbf{a})^2 + \varepsilon)}}{b} - \log(2b)$$

- Unfortunately, there is no closed-form solution (based on inverse/pseudo-inverse) in this case – unlike the case for squared error!
- One will need iterative methods like adaptive gradient descent.

• One will need iterative methods like adaptive gradient descent.

Repeat till there is no change in **a** 

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 $\alpha = \alpha_{\max}$ while  $(\alpha > \alpha_{\min})$ ł  $\mathbf{a'} = \mathbf{a} - \alpha \frac{dJ(\mathbf{a})}{d\mathbf{a}}, 0 < \alpha <<1$ if  $J(\mathbf{a'}) > J(\mathbf{a}), \alpha = \alpha_{\max} / 2$ else { $\mathbf{a} = \mathbf{a}'$ ; break }

 $\alpha$  = **step size** of gradient descent

Gradient descent converges to a local minimum of the energy function (or objective function), i.e. J(a) in this slide, if the step size  $\alpha$  is "small enough" to never.

Unfortunately, too small a step size is too expensive. A large step size may lead to increase in the energy function across iterations.

So we pick the largest possible step-size (within a given range) that reduces the energy – this is called **gradient descent** with adaptive step-size or adaptive gradient descent.

#### Robust statistics – simple example

Find "best" representative for the set of numbers

$$x_i$$

L2: 
$$J(\bar{x}) = \sum_{i} |\bar{x} - x_i|^2 \rightarrow \min$$

L1: 
$$J(\bar{x}) = \sum_{i} |\bar{x} - x_i| \rightarrow \min$$

Influence of  $x_i$  on E:  $x_i \rightarrow x_i + \Delta$ 

$$J_{new} \cong J_{old} + 2(x_i - \overline{x}) \cdot \Delta + \Delta^2$$
  
proportional to  $|\overline{x} - x_i|$ 

Outliers influence the most  $\overline{x} = \text{mean}(x_i)$ 

$$J_{new} \cong J_{old} \pm \Delta$$
 equal for all  $x_i$ 

Majority decides  $\overline{x} = \text{median}(x_i)$ 

Slide: Courtesy - Darya Frolova, Recent progress in optical flow

#### **Elections and Robust statistics**



#### New ways of defining the mean

• We know the mean as the one that minimizes the following quantity:

$$E(\mu) = \frac{\sum_{i=1}^{n} (x_i - \mu)^2}{n} \rightarrow \mu = \frac{\sum_{i=1}^{n} x_i}{n}$$

 Changing the error to sum of absolute values, we get:

$$E(\mu) = \frac{\sum_{i=1}^{n} |x_i - \mu|}{n} \rightarrow \mu = median(\{x_i\}_{i=1}^{n})$$

We will prove this in class!

# New ways of defining the mean

 We can also use errors of the following type with 0 < q ≤ 1:</li>

$$E(\mu) = \frac{\left(\sum_{i=1}^{n} |x_i - \mu|^q\right)^{1/q}}{n}$$

- Optimizing the above requires iterative methods (no closed form solutions).
- The mean computed using 0 < q ≤ 1 is quite robust to outliers – with q greater than or equal to 2, the mean is susceptible to outliers.

#### New ways of defining the mean

 The earlier definitions of the mean were for scalars. They can be extended for vectors in some d > 1 dimensions as well.

$$E(\mathbf{\mu}) = \frac{\sum_{i=1}^{n} (\mathbf{x}_{i} - \mathbf{\mu})^{2}}{n} \rightarrow \mathbf{\mu} = \frac{\sum_{i=1}^{n} \mathbf{x}_{i}}{n}$$

• For other q-norms (0 <= q < 1), we have:

$$E(\mathbf{\mu}) = \frac{\left(\sum_{i=1}^{n} |\mathbf{x}_{i} - \mathbf{\mu}|^{q}\right)^{1/q}}{n}$$

- LMedS = Least Median of Squares
- It works as follows:

 $J(\mathbf{a}) = \text{median}_{i=1:N} (y_i - f(x_i; \mathbf{a}))^2;$ select **a** for which  $J(\mathbf{a})$  is minimum.

- This has no closed form solution either and you can't do gradient descent type of techniques as the median is not differentiable.
- But it has an "algorithmic" solution.

- Step 1: Arbitrarily choose k out of N points where k is the smallest number of points required to determine a. Call this set of k points as C.
- $\circ$  Eg: If you had to do line fitting, k = ?
- $\circ$  Eg: If you were doing circle fitting, k = ?
- Eg: If you have to find the affine transformation between two point sets in 2D, you need k = ? correspondences.

- Step 2: Determine a using an inverse (say) from C.
- **Step 3:** Determine the squared residual errors for all the other *N-k* points, i.e. compute

 $\{e_i = (y_i - f(x_i; \mathbf{a}))^2\}_{i \notin C}$ 

- **Step 4:** Compute  $med_C$  = median of  $\{e_i\}$ .
- Repeat all these four steps for *S* different subsets of *k* points each.
- Pick the estimate of a corresponding to the C that has the least value of med<sub>c</sub>.
- What's the time complexity of this algorithm?

- S = number of subsets. What should be the minimum value of S?
- Let's say that some fraction p (0
- Then the probability that at least one of the S different subsets contains all inliers (i.e. yields good estimate of a) is: P = 1-(1-p<sup>k</sup>)<sup>S</sup>.
- Fix *P* to 0.99 (say) and compute *S* assuming you know *p*.
- Clearly S will increase hugely if either k is large (more parameters to determine) and/or if p is small (fewer inliers).

# Dealing with outliers: (3) RanSaC

- RanSaC = Random Sample Consensus.
- Similar in spirit to LMedS.
- Step 1: Arbitrarily choose k out of N points where k is the smallest number of points required to determine a. Call this set of k points as C.
- Step 2: Determine a using an inverse (say) from C.
- Step 3: Determine the squared residual errors for all the other *N*-*k* points, i.e. compute

 $\{e_i = (y_i - f(x_i; \mathbf{a}))^2\}_{i \notin C}$ 

# Dealing with outliers: (3) RanSaC

- Step 4: Count the number of points for which *e*<sub>i</sub> < threshold λ. These points form the "consensus set" for the chosen model.
- Repeat all 4 steps for multiple subsets and pick the subset which has maximum number of inliers and its corresponding estimate of a.
- Choice of <u>S same as LMedS.</u>

Sample result with RanSaC for line fitting.



# Dealing with outliers: (3) RanSaC

- Step 4: Count the number of points for which *e*<sub>i</sub> < threshold λ. These points form the "consensus set" for the chosen model.
- Alternatively:
- Repeat all 4 steps for multiple subsets and pick the subset C which has maximum number of inliers.
- 2. Estimate **a** using *all the points which were inliers for C*.



#### RANSAC versus LMedS

- LMedS needs no threshold to determine what is an inlier unlike RanSaC.
- But RanSaC has one advantage. What?
- LMedS will need at least 50% inliers (by definition of median).
- RanSaC can tolerate a smaller percentage of inliers (i.e. larger percentage of outliers).

#### Expected number of RanSaC iterations

 The probability that at least one point in a chosen set of k points is an outlier = 1-p<sup>k</sup>.

 The probability that the *i*-th set is the *first* set that contains no outliers = (1-p<sup>k</sup>)<sup>i-1</sup>p<sup>k</sup> to be denoted as Q(i).

#### Expected number of RanSaC iterations

 The expected number of sets to be drawn required to find the first no-outlier set =

$$= \sum_{i=1}^{k} iQ(i) = \sum_{i=1}^{k} i(1-p^{k})^{i-1} p^{k} = p^{k} \sum_{i=1}^{k} i(1-p^{k})^{i-1}$$
$$= p^{k} \sum_{i=1}^{k} \frac{d}{d(1-p^{k})} (1-p^{k})^{i} = p^{k} \frac{d}{d(1-p^{k})} \left( \sum_{i=1}^{k} (1-p^{k})^{i} \right)$$
$$= p^{k} \frac{d}{d(1-p^{k})} \left( \frac{1}{1-(1-p^{k})} \right) = \frac{p^{k}}{(p^{k})^{2}} = p^{-k}$$

#### RanSaC Variant 1

 RanSaC picks the subset C with largest number of inliers (i.e. least number of outliers), which is equivalent to picking the subset that minimizes the following cost

$$J(C) = \sum_{i \notin C} \rho(e_i);$$
$$\rho(e_i) = 1, e_i^2 \ge T$$
$$= 0, e_i^2 < T$$

#### RanSaC Variant 1: MSAC

• One could instead minimize a cost function that gives weights to inliers to see how well they fit the model:



M-estimator: an estimator that weighs inliers by their "quality", and outliers by a fixed constant

• This variant is called MSaC (M-estimator sample consensus).

#### Reminder: Planar Homography

- Given two images of a coplanar scene taken from two different cameras, how will you determine the planar homography matrix H?
- How many point correspondences will you require?

$$\mathbf{p}_{2jm} = \begin{pmatrix} u_2 \\ v_2 \\ w_2 \end{pmatrix} = \hat{\mathbf{H}} \mathbf{p}_{1jm} = \begin{pmatrix} \hat{H}_{11} & \hat{H}_{12} & \hat{H}_{13} \\ \hat{H}_{21} & \hat{H}_{22} & \hat{H}_{23} \\ \hat{H}_{31} & \hat{H}_{32} & \hat{H}_{33} \end{pmatrix} \begin{pmatrix} u_1 \\ v_1 \\ w_1 \end{pmatrix}$$

$$x_{2,im} = \frac{\hat{H}_{11}u_1 + \hat{H}_{12}v_1 + \hat{H}_{13}w_1}{\hat{H}_{31}u_1 + \hat{H}_{32}v_1 + \hat{H}_{33}w_1} = \frac{\hat{H}_{11}x_1 + \hat{H}_{12}y_1 + \hat{H}_{13}}{\hat{H}_{31}x_1 + \hat{H}_{32}y_1 + \hat{H}_{33}}, x_{1,im} = \frac{u_1}{w_1}, y_{1,im} = \frac{v_1}{w_1}$$

$$y_{2,im} = \frac{\hat{H}_{21}u_1 + \hat{H}_{22}v_1 + \hat{H}_{23}w_1}{\hat{H}_{31}u_1 + \hat{H}_{32}v_1 + \hat{H}_{33}w_1} = \frac{\hat{H}_{21}x_1 + \hat{H}_{22}y_1 + \hat{H}_{23}}{\hat{H}_{31}x_1 + \hat{H}_{32}y_1 + \hat{H}_{33}}$$

$$\begin{aligned} x_{2i}x_{1i}\hat{H}_{31} + x_{2i}y_{1i}\hat{H}_{32} + x_{2i}\hat{H}_{33} - x_{1i}\hat{H}_{11} - y_{1i}\hat{H}_{12} - \hat{H}_{13} &= 0 \\ y_{2i}x_{1i}\hat{H}_{31} + y_{2i}y_{1i}\hat{H}_{32} + y_{2i}\hat{H}_{33} - x_{1i}\hat{H}_{21} - y_{1i}\hat{H}_{22} - \hat{H}_{23} &= 0 \end{aligned} \\ \begin{pmatrix} -x_{1i} & -y_{1i} & -1 & 0 & 0 & 0 & x_{2i}x_{1i} & x_{2i}y_{1i} & x_{2i} \\ 0 & 0 & 0 & -x_{1i} & -y_{1i} & -1 & y_{2i}x_{1i} & y_{2i}y_{1i} & y_{2i} \\ \hat{H}_{13} \\ \hat{H}_{21} \\ \hat{H}_{22} \\ \hat{H}_{23} \\ \hat{H}_{31} \\ \hat{H}_{32} \\ \hat{H}_{31} \end{aligned}$$

There will be *N* such pairs of equations (i.e. totally 2*N* equations), given *N* pairs of corresponding points in the two images

=0

Ah = 0, A has size  $2N \times 9, h$  has size  $9 \times 1$ 

The equation Ah = 0 will be solved by computing the SVD of A, i.e.  $A = USV^{T}$ . The vector h will be given by the singular vector in corresponding to the null singular value (in the ideal case) or the null singular value.

#### Application: RANSAC to determine Homography between two Images

- Determine sets  $Q_1$  and  $Q_2$  of salient feature points in both images, using the SIFT algorithm.
- Q<sub>1</sub> and Q<sub>2</sub> may have different sizes!
   Determine the matching points between Q<sub>1</sub> and Q<sub>2</sub> using methods such as matching of SIFT descriptors.
- Many of these matches will be near-accurate, but there will be outliers too!



Figure 5.4.: Image pairs and detected points used in homography experiments (D and E). Inliers are marked as dots in left images and outliers as crosses in right images.

#### Application: RANSAC to determine Homography between two Images

- Pick a set of any k = 4 pairs of points and determine homography matrix H using SVD based method.
- Determine the number of inliers i.e. those point pairs for which:

$$\left\|\mathbf{q}_{1i} - \mathbf{H}\mathbf{q}_{2i}\right\|_{2}^{2} \leq \varepsilon$$

• Select the estimate of **H** corresponding to the set that yielded maximum number of inliers!





1<sup>st</sup> image: warped using estimated **H** 



#### **H** =

0.578823011557930.06780863137907 -28.33314842189324-0.060840456695420.5628359439643530.613199419103270.00002958152711-0.000031444836920.58195535780312

RANSAC result with 41% inliers (threshold on squared distance was 0.1) - point matching done using minor-Eigenvalue method with SSD based matching of 9 x 9 windows in a 50 x 50 neighborhood



1<sup>st</sup> image



2<sup>nd</sup> image

1<sup>st</sup> image: warped using estimated **H** and overlapped/merged with 2<sup>nd</sup> image – to show accuracy of alignment





Left: Result of warping 1<sup>st</sup> image using **H** estimated with simple leastsquares on the matching points (No RANSAC). Right: Result merged with 2<sup>nd</sup> image. Notice that the estimation is quite poor.

#### Some cautions with RANSAC

 Consider a dataset with a cloud of points all close to each other (degenerate set). A model created from an outlier point and any point from a degenerate set will have a large consensus set!



Image taken from Ph.D. thesis of Ondrej Chum, Czech Technical University, Prague

#### Some cautions with RANSAC

 A model created from a set of inliers need not always be optimal, i.e. it may have a very small consensus set.



Image taken from Ph.D. thesis of Ondrej Chum, Czech Technical University, Prague

#### References

- Appendix A.7 of Trucco and Verri
- Article on Robust statistics by Chuck Stewart
- Original article on RanSaC by Fischler and Bolles
- Article on RanSaC variants by Torr and Zisserman