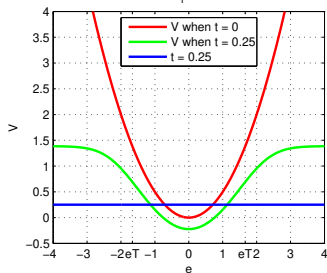


► The Action of the Robust Matching Model on Data

Example for $\hat{V}(e_{ij})$ from (26):

$$\sigma_1 = 1$$



red – the (non-robust) quadratic error

$\hat{V}(e_{ij})$ when $t = 0$

blue – the rejected match penalty t

green – robust $\hat{V}(e_{ij})$ from (26)

- if the error of a correspondence exceeds a limit, it is ignored
- then $\hat{V}(e_{ij}) = \text{const}$ and we just count outliers in (26)
- t controls the ‘turn-off’ point
- the inlier/outlier threshold is e_T – the error for which $(1 - P_0) p_1(e_T) = P_0 p_0(e_T)$: note that $t \approx 0$

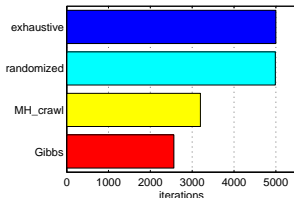
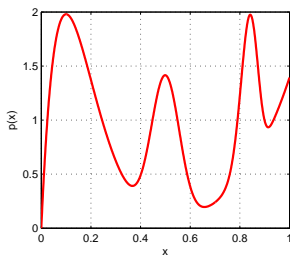
$$e_T = \sigma_1 \sqrt{-\log t^2}, \quad t = e^{-\frac{1}{2} \left(\frac{e_T}{\sigma_1} \right)^2} \quad (27)$$

The full optimization problem (23) uses (26):

$$\mathbf{F}^* = \arg \max_{\mathbf{F}} \frac{\overbrace{p(E, D | \mathbf{F})}^{\text{data model}} \cdot \overbrace{p(\mathbf{F})}^{\text{prior}}}{\underbrace{p(E, D)}_{\text{evidence}}} \approx \arg \min_{\mathbf{F}} \left[V(\mathbf{F}) + \sum_{i=1}^m \sum_{j=1}^n \log \left(e^{-\frac{e_{ij}^2(\mathbf{F})}{2\sigma_1^2}} + t \right) \right]$$

- $\pi(\mathbf{F})$ – a shorthand for the argument of the maximization
- typically we take $V(\mathbf{F}) = -\log p(\mathbf{F}) = 0$ unless we need to stabilize a computation, e.g. when video camera moves smoothly (on a high-mass vehicle) and we have a prediction for \mathbf{F}
- evidence is not needed unless we want to compare different models (e.g. homography vs. epipolar geometry)

How To Find the Global Maxima (Modes) of a PDF?



- averaged over 10^4 trials
- number of proposals before $|x - x_{\text{true}}| \leq \text{step}$

- given the function $p(x)$ at left

p.d.f. on $[0, 1]$, mode at 0.1

consider several methods:

1. exhaustive search

```
step = 1/(iterations-1);  
for x = 0:step:1  
    if p(x) > bestp  
        bestx = x; bestp = p(x);  
    end  
end
```

- slow algorithm (definite quantization)
- fast to implement

2. randomized search with uniform sampling

```
while t < iterations  
    x = rand(1);  
    if p(x) > bestp  
        bestx = x; bestp = p(x);  
    end  
    t = t+1; % time  
end
```

- equally slow algorithm
- fast to implement

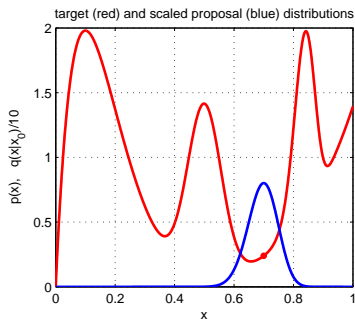
3. random sampling from $p(x)$ (Gibbs sampler)

- faster algorithm
- fast to implement but often infeasible (e.g. when $p(x)$ is data dependent (our case in correspondence prob.))

4. Metropolis-Hastings sampling

- almost as fast (with care)
- not so fast to implement
- rarely infeasible
- RANSAC belongs here

How To Generate Random Samples from a Complex Distribution?



- red: probability density function $\pi(x)$ of the toy distribution on the unit interval **target distribution**

$$\pi(x) = \sum_{i=1}^4 \gamma_i \text{Be}(x; \alpha_i, \beta_i), \quad \sum_{i=1}^4 \gamma_i = 1, \quad \gamma_i \geq 0$$

$$p(x|g) p(g) \text{Be}(x; \alpha, \beta) = \frac{1}{B(\alpha, \beta)} \cdot x^{\alpha-1} (1-x)^{\beta-1}$$

- alg. for generating samples from $\text{Be}(x; \alpha, \beta)$ is known
- \Rightarrow we can generate samples from $\pi(x)$ **how?**
- suppose we cannot sample from $\pi(x)$ but we can sample from some 'simple' proposal distribution $q(x | x_0)$, given the previous sample x_0 (blue)

$$q(x | x_0) = \begin{cases} U_{0,1}(x) & \text{(independent) uniform sampling} \\ \text{Be}(x; \frac{x_0}{T} + 1, \frac{1-x_0}{T} + 1) & \text{'beta' diffusion (crawler) } T - \text{temperature} \\ \pi(x) & \text{(independent) Gibbs sampler} \end{cases}$$

- note we have unified all the random sampling methods from the previous slide
- how to redistribute proposal samples $q(x | x_0)$ to target distribution $\pi(x)$ samples?

► Metropolis-Hastings (MH) Sampling

C – configuration (of all variable values)

e.g. $C = x$ and $\pi(C) = \pi(x)$ from →116

Goal: Generate a sequence of random samples $\{C_t\}$ from target distribution $\pi(C)$

- setup a Markov chain with a suitable transition probability to generate the sequence

Sampling procedure

1. given C_t , draw a random sample S from $q(S | C_t)$

q may use some information from C_t (Hastings)

2. compute acceptance probability

the evidence term drops out

$$a = \min \left\{ 1, \frac{\pi(S)}{\pi(C_t)} \cdot \frac{q(C_t | S)}{q(S | C_t)} \right\}$$

3. draw a random number u from unit-interval uniform distribution $U_{0,1}$

4. if $u \leq a$ then $C_{t+1} := S$ else $C_{t+1} := C_t$

'Programming' an MH sampler


1. design a proposal distribution (mixture) q and a sampler from q

2. write functions $q(C_t | S)$ and $q(S | C_t)$ that are proper distributions

not always simple

Finding the mode

- remember the best sample
- use simulated annealing
- start local optimization from the best sample


$$U(x) = \frac{1}{T}$$

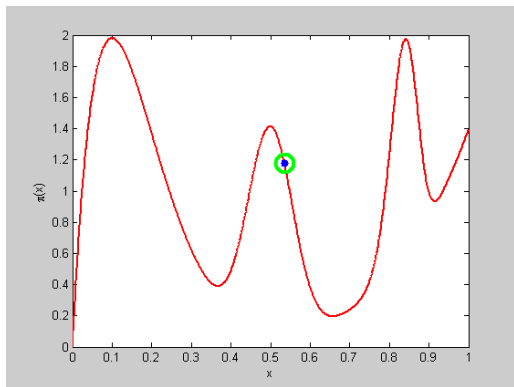
fast implementation but must wait long to hit the mode

very slow

good trade-off between speed and accuracy

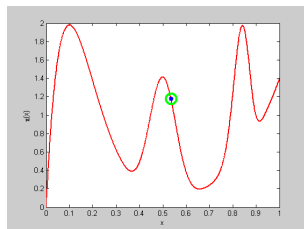
an optimal algorithm does not use just the best sample: a Stochastic EM Algorithm (e.g. SAEM)

MH Sampling Demo

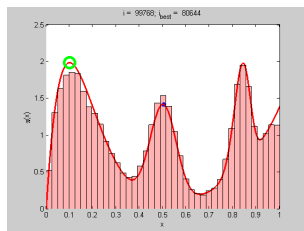


sampling process (video, 7:33, 100k samples)

- blue point: current sample
- green circle: best sample so far $\text{quality} = \pi(x)$
- histogram: current distribution of visited states
- the vicinity of modes are the most often visited states



initial sample



final distribution of visited states

Demo Source Code (Matlab)

```
function x = proposal_gen(x0)
% proposal generator q(x | x0)

    T = 0.01; % temperature
    x = betarnd(x0/T+1,(1-x0)/T+1);
end

function p = proposal_q(x, x0)
% proposal distribution q(x | x0)

    T = 0.01;
    p = betapdf(x, x0/T+1, (1-x0)/T+1);
end

function p = target_p(x)
% target distribution p(x)

% shape parameters:
a = [2 40 100 6];
b = [10 40 20 1];

% mixing coefficients:
w = [1 0.4 0.253 0.50]; w = w/sum(w);
p = 0;
for i = 1:length(a)
    p = p + w(i)*betapdf(x,a(i),b(i));
end
end
```

```
%% DEMO script

k = 10000; % number of samples
X = NaN(1,k); % list of samples

x0 = proposal_gen(0.5);
for i = 1:k
    x1 = proposal_gen(x0);
    a = target_p(x1)/target_p(x0) * ...
        proposal_q(x0,x1)/proposal_q(x1,x0);
    if rand(1) < a
        X(i) = x1; x0 = x1;
    else
        X(i) = x0;
    end
end

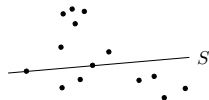
figure(1)
x = 0:0.001:1;
plot(x, target_p(x), 'r', 'linewidth',2);
hold on
binw = 0.025; % histogram bin width
n = histc(X, 0:binw:1);
h = bar(0:binw:1, n/sum(n)/binw, 'histc');
set(h, 'facecolor', 'r', 'facealpha', 0.3)
xlim([0 1]); ylim([0 2.5])
xlabel 'x'
ylabel 'p(x)'
title 'MH demo'
hold off
```

► The Elements of a Data-Driven MH Sampler

1. primitives = elementary measurements

- points in line fitting
- matches in epipolar geometry estimation

2. configuration = s-tuple of primitives minimal subsets necessary for parameter estimate



the minimization will be over a discrete set:

- of point pairs in line fitting (left)
- of match 7-tuples in epipolar geometry estimation

3. a map from configuration C to parameters θ by solving the minimal geometric problem

- line parameters \mathbf{n} from two points
- fundamental matrix \mathbf{F} from seven matches

4. target likelihood $p(E, D | \theta)$ replaces $\pi(C)$

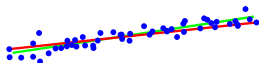
- can use log-likelihood: then it is the sum of robust errors $\hat{V}(e_{ij})$ given \mathbf{F} (26)
- robustified point distance from the line $\theta = \mathbf{n}$
- robustified Sampson error for $\theta = \mathbf{F}$
- posterior likelihood $p(E, D | \theta)p(\theta)$ can be used

MAPSAC ($\pi(S)$ includes the prior)

5. (optional) hard inlier/outlier discrimination by the threshold (27)

$$\hat{V}(e_{ij}) < e_T, \quad e_T = \sigma_1 \sqrt{-\log t^2}$$

6. parameter distribution follows the empirical distribution of s -tuples. Since the proposal is done via the minimal problem solver, it is 'data-driven',



- pairs of points define line distribution $p(\mathbf{n} | X)$ (left)
- random correspondence 7-tuples define epipolar geometry distribution $q(\mathbf{F} | M)$

7. proposal distribution $q(\cdot)$ is just a distribution of the s -tuples:

- q uniform, independent $q(S | C_t) = q(S) = \binom{mn}{s}^{-1}$, then $a = \min \left\{ 1, \frac{p(S)}{p(C_t)} \right\}$
- q dependent on descriptor similarity PROSAC (similar pairs are proposed more often)
- q dependent on the current configuration e.g. 'not far from it'

8. local optimization from promising proposals

- can use hard inliers
- cannot be used to replace C_t

9. stopping based on the probability of proposing an all-inlier sample

→122

► Data-Driven Sampler Stopping

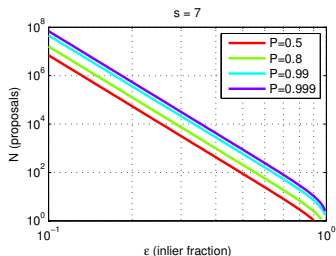
Principle: what is the number of proposals N that are needed to hit an all-inlier sample?
this will tell us nothing about the accuracy of the result

P ... probability that at least one proposal is an all-inlier $1 - P$... all previous N proposals were bad
 ε ... the fraction of inliers among primitives, $\varepsilon \leq 1$
 s ... minimal sample size (2 in line fitting, 7 in 7-point algorithm)

$$N \geq \frac{\log(1 - P)}{\log(1 - \varepsilon^s)}$$

- ε^s ... proposal does not contain an outlier
- $1 - \varepsilon^s$... proposal contains at least one outlier
- $(1 - \varepsilon^s)^N$... N previous proposals contained an outlier = $1 - P$

		N for $s = 7$	
		P	
ε	0.8	0.99	
0.5	205	590	
0.2	$1.3 \cdot 10^5$	$3.5 \cdot 10^5$	
0.1	$1.6 \cdot 10^7$	$4.6 \cdot 10^7$	



- N can be re-estimated using the current estimate for ε (if there is LO, then after LO)
the quasi-posterior estimate for ε is the average over all samples generated so far
- this shows we have a good reason to limit all possible matches to tentative matches only
- for $\varepsilon \rightarrow 0$ we gain nothing over the standard MH-sampler stopping criterion

► Stripping MH Down To Get RANSAC [Fischler & Bolles 1981]

- when we are interested in the best sample only... and we need fast data exploration...

Simplified sampling procedure

1. ~~given C_t , draw a random sample S from $q(S|C_t)$~~ $q(S)$ independent sampling
no use of information from C_t

2. ~~compute acceptance probability~~

$$a = \min \left\{ 1, \frac{\pi(S)}{\pi(C_t)} \cdot \frac{q(C_t | S)}{q(S | C_t)} \right\}$$

3. ~~draw a random number u from unit-interval uniform distribution $\mathbb{U}_{0,1}$~~
4. ~~if $u \leq a$ then $C_{t+1} := S$ else $C_{t+1} := C_t$~~
5. ~~if $\pi(S) > \pi(C_{\text{best}})$ then remember $C_{\text{best}} := S$~~

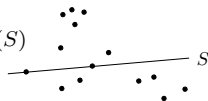
Steps 2–4 make no difference when waiting for the best sample

- ... but getting a good accuracy sample might take very long this way
- good overall exploration but slow convergence in the vicinity of a mode where C_t could serve as an attractor
- cannot use the past generated samples to estimate any parameters
- we will fix these problems by (possibly robust) 'local optimization'

► RANSAC with Local Optimization and Early Stopping

1. initialize the best sample as empty $C_{\text{best}} := \emptyset$ and time $t := 0$
2. estimate the number of needed proposals as $N := \binom{n}{s} n - \text{No. of primitives}$, $s - \text{minimal sample size}$
3. while $t \leq N$:

a) propose a minimal random sample S of size s from $q(S)$

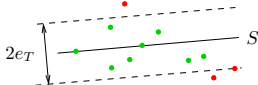


b) if $\pi(S) > \pi(C_{\text{best}})$ then

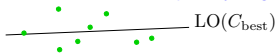
i) update the best sample $C_{\text{best}} := S$

$\pi(S)$ marginalized as in (26); $\pi(S)$ includes a prior \Rightarrow MAP

ii) threshold-out inliers using e_T from (27)



iii) start local optimization from the inliers of C_{best} LM optimization with robustified (\rightarrow 113) Sampson error possibly weighted by posterior $\pi(m_{ij})$ [Chum et al. 2003]



iv) update C_{best} , update inliers using (27), re-estimate N from inlier counts

\rightarrow 122 for derivation

$$N = \frac{\log(1 - P)}{\log(1 - \varepsilon^s)}, \quad \varepsilon = \frac{|\text{inliers}(C_{\text{best}})|}{m n},$$

c) $t := t + 1$

4. output C_{best}

• see [MPV course](#) for RANSAC details

see also [Fischler & Bolles 1981], [25 years of RANSAC]

Thank You

$s = 7$

