

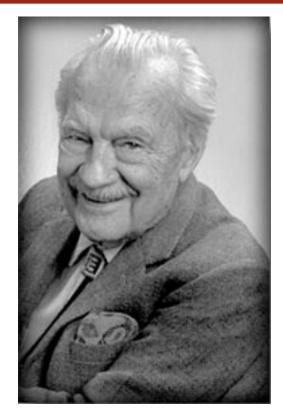
PM520: Lecture 7

Metropolis-Hastings Markov chain Monte Carlo [MCMC]



Metropolis algorithm (Metropolis, 1953)

- Define an 'energy' for a configuration of a system.
 Assume we are minimizing energy (denoted by f(x)).
- Propose moves from an old state x, to a new state x' according to a symmetric proposal kernel q()
 [So q(x->x')=q(x'->x) for all x, x'].
- Let ΔE denote the change in energy between the new and old state. i.e., we define $\Delta E(x,x')=f(x')-f(x)$.
- Move from x to x' with prob. min{1,exp(-ΔE(x,x')/T)}, where T is a 'temperature' [FIXED]. (This is for a minimizing algorithm, if you are maximizing use h=min{1,exp(ΔE(x,x')/T)}.)



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If you are minimizing:

when energy decreases, you move with prob. 1 when energy increases, you move with prob. <1.



Simulated annealing

- Kirkpatrick, S.; Gelatt, C. D.; and Vecchi, M. P. "Optimization by Simulated Annealing." Science 220, 671-680, 1983.
- Similar in spirit to the Metropolis algorithm.
- Again, uses a rule q for proposing moves to a new state x' based on the current state x. Assume we are trying to find the minimum of f.
- Define $\Delta E(x,x')=f(x')-f(x)$.
- Move to new state with prob. h=min{1,exp(-ΔE(x,x')/T)}, where T is a 'temperature'. (This is for a minimizing algorithm, if you are maximizing use h=min{1,exp(ΔE(x,x')/T)})
- e.g. T=1:
 - f(x')-f(x)=1 -> h=1/e < 1
 - f(x')-f(x)=-1 -> h=e > 1
- T is always positive, but decreases over time (What starting value? At what rate should it decrease?) so moves in the wrong direction become less likely.



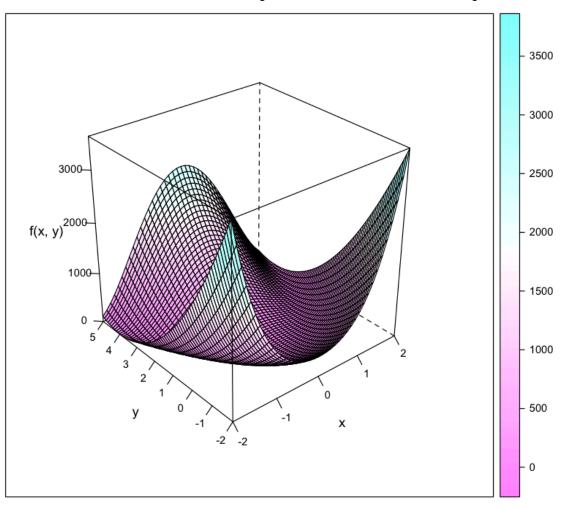
SA or Metropolis lab task #1

- Code up either the SA or Metropolis algorithm. Test it on a simple function, such as f(x,y)= x²+y²
- Then test it on the Rosenbrock function (see next slide)



Rosenbrock function

$$f(x,y)=(1-x)^2+100(y-x^2)^2$$



Commonly-used test function.

Has a global minimum at (1,1).

```
source('Rosenbrock.r')
MyRosenbrock<-function(x){
  return (Rosenbrock(x)[[1]])
}</pre>
```



SA or Metropolis task #2: regression

Regression fit aims to minimize sum of squared residuals:

$$\mathbf{x} = \{x_1, x_2, x_3, ..., x_n\}$$

$$y = \{y_1, y_2, y_3, ..., y_n\}$$

a=intercept; b=slope; so y=ax+b

SS(a,b)= Σ_i (Observed y_i - Fitted y)²

So, search through space of values for (a,b) (i.e., Intercept, slope), to find minimal (a,b) that give minimal SS(a,b).

Work in small groups if you wish to.



Sudoku

	7		1		5			
5						3		
		6	7	9		1	2	5
1	5	2	8		6		9	
				3			5	
3				5			8	
8	6			7		5	1	
				6				
		9			2		4	



Metropolis-Hastings Markov chain Monte Carlo [MCMC]



Markov Chains

- Z(t) is the state of the system at time t [t is discrete]
- The state-space, S, is the set of values that Z(t) can take.
- The transition function, $P(z_1,z_2)$, [or $P(z_2 \mid z_1)$] gives the probability that the next value of Z is z_2 , given that Z is currently in state z_1 .
- The **Markov condition**: $P(Z(t)=z(t) \mid Z(t-1)=z(t-1), Z(t-2)=z(t-2), ..., Z(0)=z(0)) = P(Z(t)=z(t) \mid Z(t-1)=z(t-1)).$
 - "Where you go depends upon where you are, but not how you got there."
 - "The future is independent of the past, except through the present state".

• Example 1:

- -X(t) = number rolled on a six-sided die
- Define Z(t) = [X(t)+X(t-1)+...+X(1)]
- So P(z_1,z_2)=1/6 if 1 ≤ z_2 - z_1 ≤ 6 (for integers z_1,z_2)



Markov Chains

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- The transition function, $P(z_1,z_2)$, [or $P(z_2 \mid z_1)$] gives the probability that the next value of Z is z_2 , given that Z is currently in state z_1 .
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 - "Where you go depends upon where you are, but not how you got there."
 - "The future is independent of the past, except through the present state".

• Example 2:

- -X(t) = -1 with prob 1/2; +1 otherwise.
- Define Z(0) = 0; Z(t)=Z(t-1)+X(t) [=X(t)+X(t-1)+...+X(0)]
- So P(z_1,z_2)=1/2 if $|z_2-z_1|$ =1 (for integers z_1,z_2)
- This is a "symmetric random walk".



Stationary distributions

The values it can take

The rule that describes how it moves

• A Markov chain, Z(), with *state-space*, S, and *transition function*, P(), has stationary distribution π () if:

$$\sum_{x \in S} \pi(x) P(x \to y) = \pi(y), \text{ for all } y \in S$$

- Intuition: if the state of Z at iteration t is distributed according to $\pi()$, then so is the state of Z at iteration t+1 (and t+2 and t+3 and.....).
- It is the long-term distribution of Z().
- Chain must also be irreducible (i.e., it can get to every state from every other state [eventually]).
- The chain must be aperiodic. (i.e., no period. Cannot have a stationary distribution otherwise)
- Fact: All irreducible, aperiodic Markov chains on finite state-spaces have a stationary distribution.



Detailed balance

• A sufficient, but not necessary condition for $\pi()$ to be the stationary distribution is the *detailed balance* requirement:

$$\pi(x)P(x \to y) = \pi(y)P(y \to x)$$
, for all $x, y \in S$

Recall previous requirement:

$$\sum_{x \in S} \pi(x) P(x \to y) = \pi(y), \text{ for all } y \in S$$



How to sample from a random variable

- Want to sample from a density f(x).
- Suppose we can calculate the cumulative density function F(x).
- 1. Generate u~Unif[0,1].
- 2. Set u=F(x).
- 3. Solve for x.
- 4. x has density f.

e.g. exponential distribution. $F(x)=1-e^{\lambda x}$, so $u=1-e^{\lambda x}$, so $x=\log(1-u)/\lambda$.

What if we cannot calculate the CDF F? (e.g. urns, population genetics models,...)



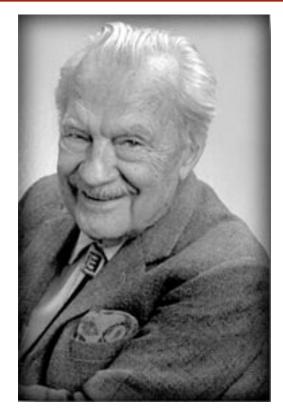
Methods for approximating a probability density f

- Monte Carlo methods:
 - find a way to sample from f,
 - do this many times and use the outcomes as an empirical approximation to f.
 - -e.g.,
 - Number of different colored balls in urn model.
- Simple, but not always very efficient.



Recall: Metropolis algorithm for 'optimization'

- Define an 'energy' for a configuration of a system.
 Assume we are minimizing energy
- Propose moves from an old state x, to a new state x' according to a symmetric proposal kernel q()
 [So q(x->x')=q(x'->x) for all x, x']
- Let ΔE=f(x')-f(x) denote the change in energy between the new and old state
- Move to new state with prob. min{1,exp(-ΔE/T)}, where T is a 'temperature' [FIXED]
- In our context, the "energy" is f(x).



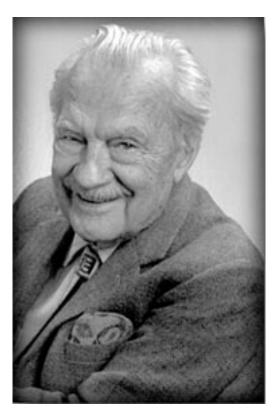
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Metropolis algorithm for **sampling** from a probability density function f...

- Assume we can calculate density f(x) (but cannot easily draw samples from it).
- Let x be the current state of the system (e.g. $x \in \mathbb{R}$).
- Iterate the following:
 - Propose move from old state x, to new state x' according to a symmetric proposal kernel q() [e.g. symmetric random walk or Normal(0, σ^2)]
 - Calculate h= $min\{1,f(x')/f(x)\}$
 - With prob. h, set x=x'; otherwise leave x unchanged.
- Let X(i) denote the state of this process at iteration i.
- X() is a Markov Chain (sequence of random variables in which the next state depends only upon the last state, not any earlier states).





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X has stationary distribution f

- Showing the following three things will show that the stationary distribution of this version of Metropolis algorithm is f:
- 1. Detailed balance equation:
 - f(x)P(x->x')=f(x')P(x'->x)
 - So, for us, f(x)q(x->x')P(accept move to x')=f(x')q(x'->x)P(accept move to x).
- 2. Ergodicity: the chain has period 1. (e.g., it does not alternate between two sets of states).
- 3. Irreducible: all states can be reached.
- Ergodicity and irreducibility will be a function of how you choose to propose changes to x, - i.e., how you define q(). So don't pick something that has a period ≠ 1, or that can't visit all possible values of x.
- But what about detailed balance?



Detailed balance

- Need to show that f(x)P(x->x')=f(x')P(x'->x)
 i.e., f(x)q(x->x')P(accept move x->x')=f(x')q(x'->x)P(accept move x'->x)
- If q is symmetric (e.g. symmetric random walk), then, by definition we will have q(x->x')=q(x'->x), so we just need to show that...
- f(x)P(accept move x->x')=f(x')P(accept move to x'->x), i.e.
- f(x')/f(x) = P(accept move x->x')/P(accept move to x'->x).
- But
 - P(accept move x->x')= min{1, f(x')/f(x)}
 - P(accept move x'->x)= min{1,f(x)/f(x')}
- At least one of f(x')/f(x) and f(x)/f(x') will be ≥ 1 by construction. Without loss of generality, suppose $f(x)/f(x')\geq 1$. So...
- f(x')/f(x) = P(accept move x->x')/P(accept move to x'->x).
 - = $\min\{1, f(x')/f(x)\}/\min\{1, f(x)/f(x')\}= f(x')/f(x)\}.$



Example: Simulating values from Normal distⁿ.

Pseudocode:

set.seed(4970) # any number will do here

Call this function as follows: MyMH<-MHNormal(0,1,5,50000)

```
MHNormal<-function(Mean,SD,EndPt,NumberOflts){ # Mean and SD are the Mean and Std. dev. of the Normal,
      #We start somewhere in the range [-EndPt, +EndPt] and run NumberOflts iterations.
      NormalXs<-mat.or.vec(1,NumberOflts) # declare a vector to store the x-values we generate during the course of the
       algorithm
      # you need to start somewhere: sample a random number, x, between +EndPt and -EndPt. Then do the following loop:
      for (i in 1:NumberOflts){ # each iteration performs one step of the Metropolis process
         # propose a new x-value (xprime) by adding a Unif(-0.5,0.5) (say) random variable to the current value
         # decide whether to move to the new value -
         # Calculate Metropolis term (the ratio of the likelihoods of the two points, i.e., H < -f(x')/f(x))
         H<-min(1,dnorm(xprime,Mean,SD)/dnorm(x,Mean,SD))
         p<-runif(1,0,1) # generate an random number
         if (p<H){
                      # move to x'
              X<-X'
         } # we don't need an else, since we just keep the existing x in that case
         # save the current value of x to the vector NormalXs
         # every now and again, plot a histogram of the accepted values
         if (i%%100==0){ # this will pot one every 100 iterations (i%%100 means the remainder when i is divided by 100)
              hist(NormalXs[0:i],breaks=HistBars)
      return (NormalXs) # returns the vector of x-values
```

On Github in repo: MetropolisForNormals



(Non-examinable) Lab exercise 1 - part 1

- Code up the Metropolis algorithm for sampling from a Normal distribution with mean 0 and std. dev. 1.
- Use a symmetric proposal kernel that adds a number y to x each time, so that x' = x+y, where, for example, y~Unif[-c,+c] or y~Normal(0, σ^2).
- Run the algorithm:
 - Start it from some value (how are you going to choose that value?)
 - Note that consecutive samples are not independent.
 - You can construct (almost) independent samples by sub-sampling your iterations, sampling every kth iteration, for some k. (How to choose k?)
 - Explore how the choice of k you need to use to get independent subsamples varies as a function of c above.



Non-examinable lab exercise 1 - part 2

- Use the code you wrote for the Metropolis algorithm for sampling from a Normal distribution with mean 0 and std. dev. 1.
- Use an asymmetric proposal kernel that at each iteration:
 - with prob. 3/4 adds a number y to x e, so that x' = x+y, where y~Unif[0,c]
 - with prob. 1/4 subtracts a number y from x, so x' = x-y, where $y\sim Unif[0,c]$.
- Run the algorithm for c=0.5 (say),
 - what happens?



For general (non-symmetric) proposal kernels - Detailed balance

- Recall, we need f(x)P(x->x')=f(x')P(x'->x)
- Rewriting, we need:
- f(x')/f(x) = P(move x->x')/P(move to x'->x).
 - = P(propose move x -> x')P(accept x -> x') / P(propose move x' -> x)P(accept x' -> x))
 - = q(x->x') P(accept x->x')/q(x'->x) P(accept x'->x)

Rearranging, we need:

[P(accept x->x') / P(accept x'->x)] = [q(x'->x)/q(x->x')][f(x')/f(x)]



Metropolis-Hastings

- We need: [P(accept x->x') / P(accept x'->x)] = [q(x'->x)/q(x->x')][f(x')/f(x)]
- Old acceptance prob: h= min{1,f(x')/f(x)}
- New acceptance probability: use h=min{1, [f(x')q(x'->x)] / [f(x)q(x->x')]} the Hastings Ratio.
- Detailed balance is satisfied, so the stationary distribution is f.



Hastings at the U. Toronto chess club



MCMC: Metropolis-Hastings

- 1. If at x, propose move to x' according to transition kernel q(x -> x').
- 2. Calculate

$$h = \min \left\{ 1, \frac{f(x')q(x' \to x)}{f(x)q(x \to x')} \right\}$$

No need for q() to be symmetric - although it often is.

- 3. Move to x' with prob. h, else remain at x.
- 4. Return to 1.

(Metropolis et al. 1953, Hastings 1970)

The resulting Markov chain has stationary distribution f.



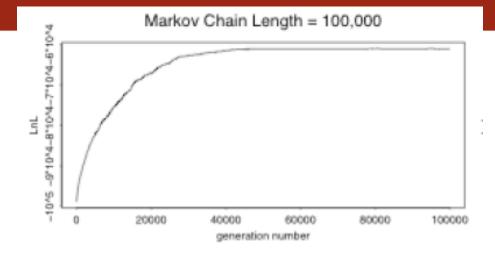
Non-examinable lab exercise 1 - part 3

- Use the code you wrote for the Metropolis algorithm for sampling from a Normal distribution with mean 2 and std. dev. 1.
- Use an **asymmetric** proposal kernel that at each iteration:
 - with prob. 3/4 adds a number y to x e, so that x' = x+y, where y~Unif[0,c]
 - with prob. 1/4 subtracts a number y from x, so x' = x-y, where $y\sim Unif[0,c]$.
- Correct it by using the Hastings Ratio [i.e. adding the term q(x'->x)/ q(x->x')]
- Run the algorithm and show that it works now.
- You are now doing Metropolis-Hastings Markov chain Monte Carlo [MH-MCMC]. This
 is perhaps the most common form of MCMC.



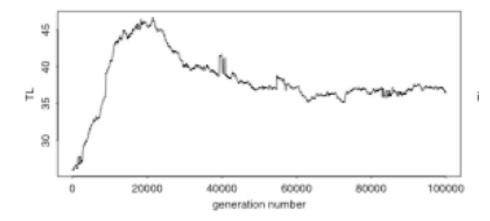
- Once stationarity is reached each iteration represents a sample from f.
- Store every iteration after 'burn-in' period.
- Consecutive outputs are correlated, so we sometimes subsample to produce 'independent' samples. (This has become less important as available computational memory has grown.)
- The art is in choosing a good transition kernel q.
- Aim for an acceptance rate (the proportion of proposals that are accepted) in the region of 20%. [If it is too low, your step size is probably too big and the algorithm won't mix very well; if it is too high, your step-size is probably too small and again the algorithm won't mix well] But this is just a rule of thumb!



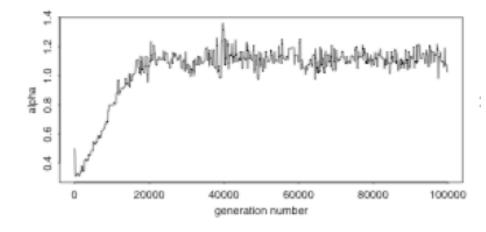


Convergence: "once stationarity has been reached"

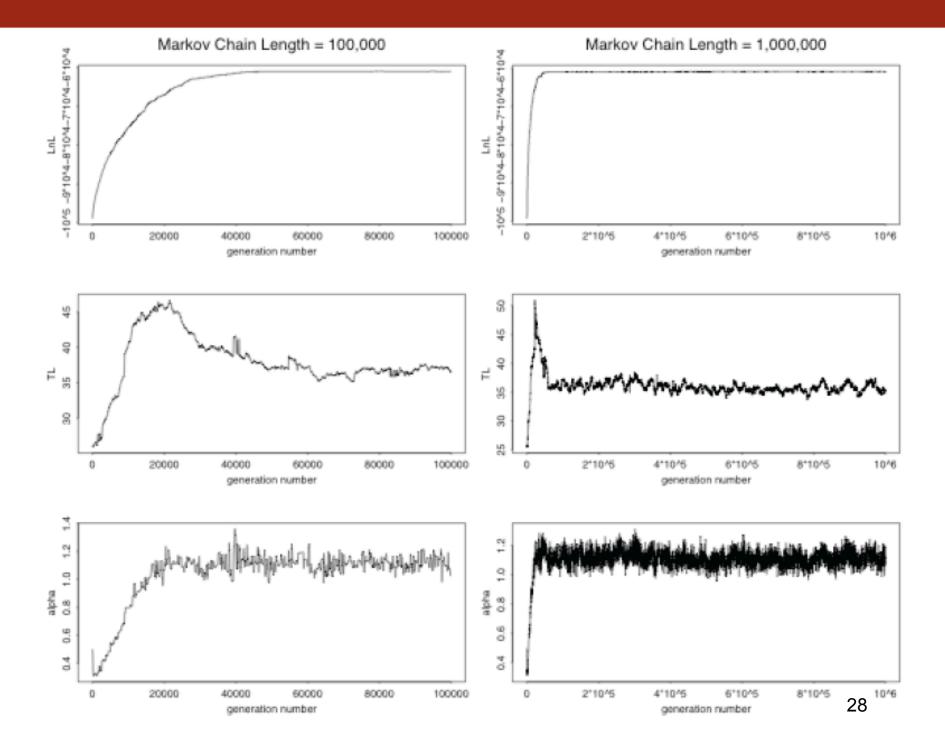
- to assess by eye, plot output against iteration number



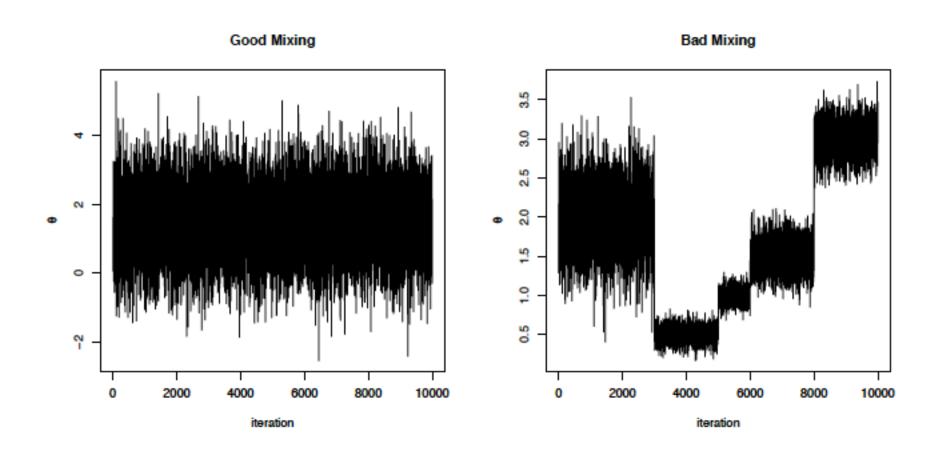
[MCMC package in R. CODA convergence diagnostics (SPLUS/R).]





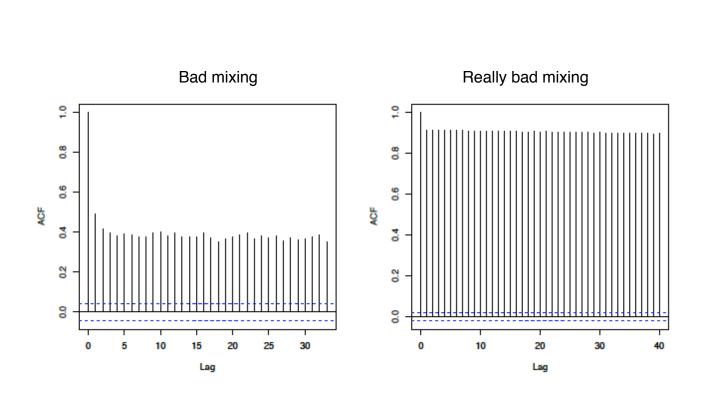


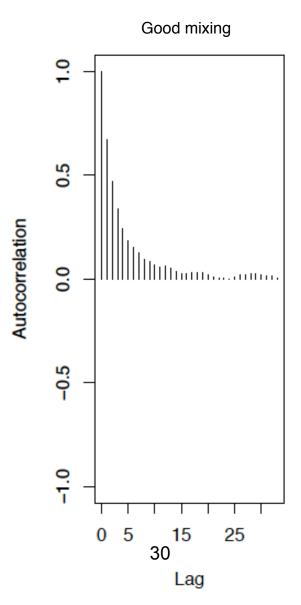






 Autocorrelation plots (ith bar shows correlation between X(t) and X(t+i))







Metropolis-Hastings: Features

• Pros:

- The algorithm 'learns'.
- Comparison is 'local'.
- MCMC package in R [next week!], Gelman/CODA convergence diagnostics, WinBUGS, OpenBUGS, JAGS, many others...

• Cons:

- Need to be able to calculate f(x) (may be a problem in some cases).
- Need to sample from stationary distribution.
- Consecutive outputs are correlated.
- May get stuck in local minima



Gelman stationarity test

- Run two (or more) replicate analyses and compare the answers.
- Compare the variance of the parameter (e.g., x) distribution in the two (or more) chains.
- If both chains have reached stationarity, the variance within each replicate analysis will be the same as that across (i.e. after combining) the two replicates.
- So the Gelman test compares "across run" variance to "within run" variance (dividing the former by the latter). Stationarity has been reach if the statistic is close to 1 ("less than 1.1" is a common, but not very strict, "rule of thumb").



Non examinable lab exercise 2

- Code up the Metropolis-Hastings MCMC algorithm for sized-biased sampling from an arbitrary distribution over the range [0,10].
 - Test your algorithm by using it to construct sized-biased samples from an $\exp(\lambda)$ distribution. Plot your results and compare them to the curve $x\lambda e^{-\lambda x}$.
 - Recall: Sized-biased samples from a density f(x) have density proportional to xf(x), rather than f(x)
 - Note that your algorithm will work for any density function f.

Pseudocode on Github in SizeBiasedMCMC



Gelman stationarity test

library("coda")

plot(SB,type="l")

Example will be posted on GitHub in the solutions repo as SizeBiasedMCMCWorking.R

convert to mcmc objects, with a burn-in

MCMC1<-mcmc(SB,start=1000)

MCMC2<-mcmc(SB2,start=1000)

combine different mcmc chain objects to an mcmc list.

Combined<-mcmc.list(list(MCMC1,MCMC2))

gelman functions are
gelman.plot(Combined) # for plots
gelman.diag(Combined) # for diagnostic values



Examinable assignment 3 - part 1

- Use the IndeptGamma.R code to use MCMC to produce samples from a Gamma(2.3,2.7) random variable
- Show how the performance deteriorates when you run the algorithm to produce samples from a Gamma(0.1,0.01).
 - Use the Gelman plot to illustrate the deterioration in performance
- Find a proposal kernel (i.e., q(x->x')), the way of producing new candidate values) that performs more efficiently.
 - Again use the Gelman plots to show how performance has changed.
 - Discuss how the performance has improved
- Due On March 25th, at 11.59pm.



Examinable assignment 3 - part 1

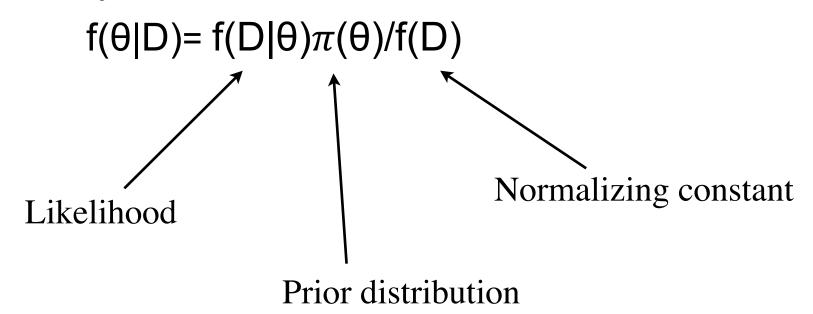
```
# metropolis-hastings independence sampler for a
# gamma based on normal candidates with the same mean and variance
set.seed(371)
gamm<-function (n, a, b)
 mu <- a/b # the mean of the gamma distribution
 sig <- sqrt(a/(b * b)) # the stadard deviation of the gamma distn
 vec <- vector("numeric", n) # this is where we are going to put the random variables we generate
 x <- a/b
 vec[1] <- x # We arbitrarily start the MCMC process at the mean
 for (i in 2:n) {
  can <- rnorm(1, mu, sig)
  hprob <- min(1, (dgamma(can, a, b)/dgamma(x,a,b))/(dnorm(can, mu, sig)/dnorm(x, mu, sig))) # where is the q term here?
  u <- runif(1)
  if (u < hprob)
   x <- can
  vec[i] <- x
 vec
```

Code on GitHub in Assignment3 repo as 'IndeptGamma.Rmd'



Sampling from Posterior Distributions

- Data, D; model parameter(s) θ. Density f(D|θ).
- Bayes theorem:





- 1. If at θ , propose move to θ according to transition kernel $q(\theta \rightarrow \theta)$.
- 2. Calculate

$$h = \min \left\{ 1, \frac{[P(\theta'|D)q(\theta' \to \theta)}{[P(\theta|D)q(\theta \to \theta')]} \right\}$$

- 3. Move to θ ' with prob. h, else remain at θ .
- 4. Return to 1.

The stationary distribution of this chain will be $P(\theta \mid D)$



- 1. If at θ , propose move to θ' according to transition kernel $q(\theta \rightarrow \theta')$.
- 2. Calculate

$$h = \min \left\{ 1, \frac{[P(D|\theta')\pi(\theta')/P(D)]q(\theta' \to \theta)}{[P(D|\theta)\pi(\theta)/P(D)]q(\theta \to \theta')} \right\}$$

- 3. Move to θ' with prob. h, else remain at θ .
- 4. Return to 1.

The stationary distribution of this chain will be $P(\theta \mid D)$



- 1. If at θ , propose move to θ' according to transition kernel $q(\theta \rightarrow \theta')$.
- 2. Calculate

$$h = \min \left\{ 1, \frac{P(D|\theta')\pi(\theta')q(\theta' \to \theta)}{P(D|\theta)\pi(\theta)q(\theta \to \theta')} \right\}$$

- 3. Move to θ' with prob. h, else remain at θ .
- 4. Return to 1.

The stationary distribution of this chain will be $P(\theta \mid D)$.



- 1. If at θ , propose move to θ' according to transition kernel $q(\theta \rightarrow \theta')$.
- 2. Calculate

$$h = \min \left\{ 1, \frac{P(D|\theta')\pi(\theta')q(\theta' \to \theta)}{P(D|\theta)\pi(\theta)q(\theta \to \theta')} \right\}$$

- 3. Move to θ' with prob. h, else remain at θ .
- 4. Return to 1.

The stationary distribution of this chain will be $P(\theta \mid D)$.

The $q(\theta \to \theta')$ in step 2. is the prob. (or prob. density) of proposing a move from θ to θ' . It must be calculated using the transition kernel you use, and it does not involve the acceptance probability at all. Similarly for $q(\theta' \to \theta)$



Transition kernel, q

- I've added some examples to Github of how to calculate the q(x -> x') term in the Hasting's ratio.
- You can find those in Week7—MCMC-calculating-q
- For those of you that use LaTeX, this is also an example of how to do that on Github.



Examinable Assignment 3 part 2: Codebreaking (due March 25th, 11.59pm)

Gzo uclfg gcpo C qhcs okof te Gollk Qoeetb zo rhf slvem ce h LtqqfLtkio Fcqxol Rlhcgz tvgfcso gzo gollhio tu Gzo Sheiolf. Gzo ahlmced qtg hggoesheg zhs yltvdzg gzo ihl tvg hes zo rhf fgcqq ztqsced gzo sttl taoe yoihvfo Gollk Qoeetbf qoug uttg rhf fgcqq shedqced tvgfcso, hf cu zo zhs utldtggoe zo zhs teo. Zo zhs h ktvedqttmced uhio yvg zcf zhcl rhf yteo rzcgo. Ktv itvqs goqq yk zcf okof gzhg zo rhf aqhfgolos gt gzo zholqceo, yvg tgzolrcfo zo qttmos qcmo hek tgzol ecio ktveddvk ce h sceeol jhimog rzt zhs yooe faoesced gtt pviz pteok ce h jtceg gzhg obcfgf utl gzhg avlatfo hes utl et tgzol. Gzolo rhf h dclq yofcso zcp. Zol zhcl rhf h qtxoqk fzhso tu shlm los hes fzo zhs h scfgheg fpcqo te zol qcaf hes txol zol fztvqsolf fzo zhs h yqvo pcem gzhg hqptfg phso gzo LtqqfLtkio qttm qcmo jvfg hetgzol hvgtptycqo. Cg scseg wvcgo. Etgzced ihe. Gzo hggoesheg rhf gzo vfvhq zhqugtvdz izhlhigol ce h rzcgo ithg rcgz gzo ehpo tu gzo lofghvlheg fgcgizos hiltff gzo ulteg tu cg ce los. Zo rhf doggced uos va. "Qttm, pcfgol," zo fhcs rcgz he osdo gt zcf xtcio, "rtvqs ktv pces h rztqo qtg avqqced ktvl qod cegt gzo ihl ft C ihe mces tu fzvg gzo sttl Tl fztvqs C taoe cg hqq gzo rhk ft ktv ihe uhqq tvg" Gzo dclq dhxo zcp h qttm rzciz tvdzg gt zhxo fgvim hg qohfg utvl ceizof tvg tu zcf yhim. Cg scseg ytgzol zcp oetvdz gt dcxo zcp gzo fzhmof. Hg Gzo Sheiolf gzok dog gzo ftlg tu aotaqo gzhg scfcqqvfcte ktv hytvg rzhg h qtg tu dtquced pteok ihe st utl gzo aolftehqcgk.



Your clues:

- It is a simple substitution code
 - -e.g. "a/A" might stand for "m/M"; "b/B" might stand for "f/F"
- The substitution is the same for the entire message.
- I have prepared a table of pairwise letter frequencies...
- We are doing this just after we learned about optimization and MCMC.



- 1 20 33 52 0.1 12 18 5 39 1 12 57 26 181 1 20 1 75 95 104 9 20 13 1 26 1
- 11 1 0.1 0.1 47 0.1 0.1 0.1 6 1 0.1 17 0.1 0.1 19 0.1 0.1 11 2 1 21 0.1 0.1 0.1 11 0.1
- 31 0.1 4 0.1 38 0.1 0.1 38 10.1 0.1 18 9 0.1 0.1 45 0.1 1 11 1 15 7 0.1 0.1 0.1 1 0.1
- 48 20 9 13 57 11 7 25 50.1 3 1 11 14 16 41 6 0.1 14 35 56 10.1 2 19 0.1 10.1 0.1
- 110 23 45 126 48 30 15 33 41 3 5 55 47 111 33 28 2 169 115 83 6 24 50 9 26 0.1
- 25 2 3 2 20 11 1 8 23 1 0.1 8 5 1 40 2 0.1 16 5 37 8 0.1 3 0.1 2 0.1
- 24 3 2 2 28 3 4 35 18 1 0.1 7 3 4 23 1 0.1 12 9 16 7 0.1 5 0.1 1 0.1
- 114 2 2 1 302 2 1 6 97 0.1 0.1 2 3 1 49 1 0.1 8 5 32 8 0.1 4 0.1 4 0.1
- 10 5 32 33 23 17 25 6 1 1 8 37 37 179 24 6 0.1 27 86 93 1 14 7 2 0.1 2
- 2 0.1 0.1 0.1 2 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 8 0.1 0.1 0.1 0.1 0.1
- 6 1 1 1 29 1 0.1 2 14 0.1 0.1 2 1 9 4 0.1 0.1 0.1 5 4 1 0.1 2 0.1 2 0.1
- 40 3 2 36 64 10 1 4 47 0.1 3 56 4 2 41 3 0.1 2 11 15 8 3 5 0.1 31 0.1
- 44 7 1 1 68 2 1 3 25 0.1 0.1 1 5 2 29 11 0.1 3 10.1 9 8 0.1 4 0.1 18 0.1
- 40 7 25 146 66 8 92 16 33 2 8 9 7 8 60 4 1 3 33 106 6 2 12 0.1 11 0.1
- 16 12 13 18 5 80 7 11 12 1 13 26 48 106 36 15 0.1 84 28 57 115 12 46 0.1 5 1
- 23 1 0.1 0.1 30 1 0.1 3 12 0.1 0.1 15 1 0.1 21 10.1 0.1 18 5 11 6 0.1 1 0.1 1 0.1
- 50 7 10.1 20 133 8 10 12 50 1 8 10 14 16 55 6 0.1 14 37 42 12 4 11 0.1 21 0.1
- 67 11 17 7 74 11 4 50 49 2 6 13 12 10 57 20 2 4 43 109 20 2 24 0.1 4 0.1
- 59 10 11 7 75 9 3 330 76 1 2 17 11 7 115 4 0.1 28 34 56 17 1 31 0.1 16 0.1
- 7 5 12 7 7 2 14 2 8 0.1 1 34 8 36 1 16 0.1 44 35 48 0.1 0.1 2 0.1 1 0.1
- 5 0.1 0.1 0.1 65 0.1 0.1 0.1 11 0.1 0.1 0.1 0.1 0.1 4 0.1 0.1 0.1 0.1 0.1 0.1 0.1 0.1 0.1 1 0.1
- 66 1 1 2 39 1 0.1 44 39 0.1 0.1 2 1 12 29 0.1 0.1 3 4 4 1 0.1 2 0.1 1 0.1
- 1 0.1 2 0.1 1 0.1 0.1 0.1 2 0.1 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1
- 18 7 6 6 14 7 3 10.1 11 1 1 4 6 3 36 4 0.1 3 19 20.1 1 1 12 0.1 2 0.1



- 1 <u>20 3</u>3 52 0.1 12 18 5 39 1 12 57 26 181 1 20 1 75 95 104 9 20 13 1 26 1
- 11 1 0.1 0.1 47 0.1 0 1 0.1 6 1 0.1 17 0.1 0.1 19 0.1 0.1 11 2 1 21 0.1 0.1 0.1 11 0.1
- 31 0.1 4 0.1 38 0.1 0.1 38 10.1 0.1 18 9 0.1 0.1 45 0.1 1 11 1 15 7 0.1 0.1 0.1 1 0.1
- 48 20 9 13 57 11 7 25 50.1 3 1 11 14 16 41 6 0.1 14 35 56 10.1 2 19 0.1 10.1 0.1
- 110 23 45 126 48 30 15 33 41 3 5 55 47 111 33 28 2 169 115 83 6 24 50 9 26 0.1
- 25 2 3 2 20 11 1 8 23 1 0.1 8 5 1 40 2 0.1 16 5 37 8 0.1 3 0.1 2 0.1
- 24 3 2 2 28 3 4 35 18 1 0.1 7 3 4 23 1 0.1 12 9 16 7 0.1 5 0.1 1 0.1
- 114 2 2 1 302 2 1 6 97 0.1 0.1 2 3 1 49 1 0.1 8 5 32 8 0.1 4 0.1 4 0.1
- 114 2 2 1 302 2 1 0 97 0.1 0.1 2 3 1 49 1 0.1 6 3 32 6 0.1 4 0.1 4 0.1
- 10 5 32 33 23 17 25 6 1 1 8 37 37 179 24 6 0.1 27 86 93 1 14 7 2 0.1 2
- 2 0.1 0.1 0.1 2 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1 8 0.1 0.1 0.1 0.1
- 6 1 1 1 29 1 0.1 2 14 0.1 0.1 2 1 9 4 0.1 0.1 0.1 5 4 1 0.1 2 0.1 2 0.1
- 40 3 2 36 64 10 1 4 47 0.1 3 56 4 2 41 3 0.1 2 11 15 8 3 5 0.1 31 0.1
- 44 7 1 1 68 2 1 3 25 0.1 0.1 1 5 2 29 11 0.1 3 10.1 9 8 0.1 4 0.1 18 0.1
- 40 7 25 146 66 8 92 16 33 2 8 9 7 8 60 4 1 3 33 106 6 2 12 0.1 11 0.1
- 16 12 13 18 5 80 7 11 12 1 13 26 48 106 36 15 0.1 84 28 57 115 12 46 0.1 5 1
- 23 1 0.1 0.1 30 1 0.1 3 12 0.1 0.1 15 1 0.1 21 10.1 0.1 18 5 11 6 0.1 1 0.1 1 0.1
- 50 7 10.1 20 133 8 10 12 50 1 8 10 14 16 55 6 0.1 14 37 42 12 4 11 0.1 21 0.1
- 67 11 17 7 74 11 4 50 49 2 6 13 12 10 57 20 2 4 43 109 20 2 24 0.1 4 0.1
- 59 10 11 7 75 9 3 330 76 1 2 17 11 7 115 4 0.1 28 34 56 17 1 31 0.1 16 0.1
- 7 5 12 7 7 2 14 2 8 0.1 1 34 8 36 1 16 0.1 44 35 48 0.1 0.1 2 0.1 1 0.1
- 66 1 1 2 39 1 0.1 44 39 0.1 0.1 2 1 12 29 0.1 0.1 3 4 4 1 0.1 2 0.1 1 0.1
- 1 0.1 2 0.1 1 0.1 0.1 0.1 2 0.1 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1
- 18 7 6 6 14 7 3 10.1 11 1 1 4 6 3 36 4 0.1 3 19 20.1 1 1 12 0.1 2 0.1

How often "a" was followed by "a"



- 1 20 <u>33</u> 52 0.1 12 18 5 39 1 12 57 26 181 1 20 1 75 95 104 9 20 13 1 26 1
- 11 1 0.1 0.1 47 0.1 0.1 0.1 6 1 0.1 17 0.1 0.1 19 0.1 0.1 11 2 1 21 0.1 0.1 0.1 11 0.1
- 31 0.1 4 0.1 38 0.1 0.1 38 10.1 0.1 18 9 0.1 0.1 45 0.1 1 11 1 15 7 0.1 0.1 0.1 1 0.1
- 48 20 9 13 57 11 7 25 50.1 3 1 11 14 16 41 6 0.1 14 35 56 10.1 2 19 0.1 10.1 0.1
- 110 23 45 126 48 30 15 33 41 3 5 55 47 111 33 28 2 169 115 83 6 24 50 9 26 0.1
- 25 2 3 2 20 11 1 8 23 1 0.1 8 5 1 40 2 0.1 16 5 37 8 0.1 3 0.1 2 0.1
- 24 3 2 2 28 3 4 35 18 1 0.1 7 3 4 23 1 0.1 12 9 16 7 0.1 5 0.1 1 0.1
- 114 2 2 1 302 2 1 6 97 0.1 0.1 2 3 1 49 1 0.1 8 5 32 8 0.1 4 0.1 4 0.1

How often "a" was followed by "b"

- 10 5 32 33 23 17 25 6 1 1 8 37 37 179 24 6 0.1 27 86 93 1 14 7 2 0.1 2
- 2 0.1 0.1 0.1 2 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1 8 0.1 0.1 0.1 0.1
- 6 1 1 1 29 1 0.1 2 14 0.1 0.1 2 1 9 4 0.1 0.1 0.1 5 4 1 0.1 2 0.1 2 0.1
- 40 3 2 36 64 10 1 4 47 0.1 3 56 4 2 41 3 0.1 2 11 15 8 3 5 0.1 31 0.1
- 44 7 1 1 68 2 1 3 25 0.1 0.1 1 5 2 29 11 0.1 3 10.1 9 8 0.1 4 0.1 18 0.1
- 40 7 25 146 66 8 92 16 33 2 8 9 7 8 60 4 1 3 33 106 6 2 12 0.1 11 0.1
- 16 12 13 18 5 80 7 11 12 1 13 26 48 106 36 15 0.1 84 28 57 115 12 46 0.1 5 1
- 23 1 0.1 0.1 30 1 0.1 3 12 0.1 0.1 15 1 0.1 21 10.1 0.1 18 5 11 6 0.1 1 0.1 1 0.1
- 50 7 10.1 20 133 8 10 12 50 1 8 10 14 16 55 6 0.1 14 37 42 12 4 11 0.1 21 0.1
- 67 11 17 7 74 11 4 50 49 2 6 13 12 10 57 20 2 4 43 109 20 2 24 0.1 4 0.1
- 59 10 11 7 75 9 3 330 76 1 2 17 11 7 115 4 0.1 28 34 56 17 1 31 0.1 16 0.1
- 7 5 12 7 7 2 14 2 8 0.1 1 34 8 36 1 16 0.1 44 35 48 0.1 0.1 2 0.1 1 0.1
- 66 1 1 2 39 1 0.1 44 39 0.1 0.1 2 1 12 29 0.1 0.1 3 4 4 1 0.1 2 0.1 1 0.1
- 1 0.1 2 0.1 1 0.1 0.1 0.1 2 0.1 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1
- 18 7 6 6 14 7 3 10.1 11 1 1 4 6 3 36 4 0.1 3 19 20.1 1 1 12 0.1 2 0.1



- 1 20 33 **5**2 0.1 12 18 5 39 1 12 57 26 181 1 20 1 75 95 104 9 20 13 1 26 1
- 11 1 0.1 0.1 47 0.1 0 1 0.1 6 1 0.1 17 0.1 0.1 19 0.1 0.1 11 2 1 21 0.1 0.1 0.1 11 0.1
- 31 0.1 4 0.1 38 0.1 0.1 38 10.1 0.1 18 9 0.1 0.1 45 0.1 1 11 1 15 7 0.1 0.1 0.1 1 0.1
- 48 20 9 13 57 11 7 25 50.1 3 1 11 14 16 41 6 0.1 14 35 56 10.1 2 19 0.1 10.1 0.1
- 110 23 45 126 48 30 15 33 41 3 5 55 47 111 33 28 2 169 115 83 6 24 50 9 26 0.1
- 25 2 3 2 20 11 1 8 23 1 0.1 8 5 1 40 2 0.1 16 5 37 8 0.1 3 0.1 2 0.1
- 24 3 2 2 28 3 4 35 18 1 0.1 7 3 4 23 1 0.1 12 9 16 7 0.1 5 0.1 1 0.1
- 114 2 2 1 302 2 1 6 97 0.1 0.1 2 3 1 49 1 0.1 8 5 32 8 0.1 4 0.1 4 0.1

How often "a" was followed by "c"

- 10 5 32 33 23 17 25 6 1 1 8 37 37 179 24 6 0.1 27 86 93 1 14 7 2 0.1 2
- 2 0.1 0.1 0.1 2 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1 8 0.1 0.1 0.1 0.1
- 6 1 1 1 29 1 0.1 2 14 0.1 0.1 2 1 9 4 0.1 0.1 0.1 5 4 1 0.1 2 0.1 2 0.1
- 40 3 2 36 64 10 1 4 47 0.1 3 56 4 2 41 3 0.1 2 11 15 8 3 5 0.1 31 0.1
- 44 7 1 1 68 2 1 3 25 0.1 0.1 1 5 2 29 11 0.1 3 10.1 9 8 0.1 4 0.1 18 0.1
- 40 7 25 146 66 8 92 16 33 2 8 9 7 8 60 4 1 3 33 106 6 2 12 0.1 11 0.1
- 16 12 13 18 5 80 7 11 12 1 13 26 48 106 36 15 0.1 84 28 57 115 12 46 0.1 5 1
- 23 1 0.1 0.1 30 1 0.1 3 12 0.1 0.1 15 1 0.1 21 10.1 0.1 18 5 11 6 0.1 1 0.1 1 0.1
- 50 7 10.1 20 133 8 10 12 50 1 8 10 14 16 55 6 0.1 14 37 42 12 4 11 0.1 21 0.1
- 67 11 17 7 74 11 4 50 49 2 6 13 12 10 57 20 2 4 43 109 20 2 24 0.1 4 0.1
- 59 10 11 7 75 9 3 330 76 1 2 17 11 7 115 4 0.1 28 34 56 17 1 31 0.1 16 0.1
- 7 5 12 7 7 2 14 2 8 0.1 1 34 8 36 1 16 0.1 44 35 48 0.1 0.1 2 0.1 1 0.1
- 66 1 1 2 39 1 0.1 44 39 0.1 0.1 2 1 12 29 0.1 0.1 3 4 4 1 0.1 2 0.1 1 0.1
- 1 0.1 2 0.1 1 0.1 0.1 0.1 2 0.1 0.1 0.1 0.1 0.1 0.1 3 0.1 0.1 0.1 3 0.1 0.1 0.1 0.1 0.1 0.1
- 18 7 6 6 14 7 3 10.1 11 1 1 4 6 3 36 4 0.1 3 19 20.1 1 1 12 0.1 2 0.1



R commands to read coded messages

read coded message

input2<- readLines("CodedMessage_Short.txt", n=1000)

read frequency table

PairsFreqs<-read.table("LetterPairFreqFrom7Novels.txt")



Assignment 3 - Part 2

- Use one of the methods you have met in the course to try to break the code.
- It may not be easy to break the code completely without resorting to manual 'tweaks' at the end, but you should be able to get to the point at which you can work out what the text is saying.
- Write it up as an Rmarkdown file and include a description of the methods you are using (for both parts of the assignment) and a discussion of your results.
- Due March 25th, 11.59pm.



Suggestion

- Need something to maximize, so let's maximize the likelihood.
- Suppose we have a sequence of letters $l_1 l_2 l_3 l_4 l_5$. Then [Recall P(A | B)=P(A \cap B)/P(B)]:
 - $P(l_1 l_2 l_3 l_4 l_5) = P(l_1) P(l_2 | l_1) P(l_3 | l_2 l_1) P(l_4 | l_3 l_2 l_1) P(l_5 | l_4 l_3 l_2 l_1)$
- Suppose the sequence of letters is Markovian. Then:
 - $-P(l_1 l_2 l_3 l_4 l_5)=P(l_1) P(l_2 | l_1) P(l_3 | l_2) P(l_4 | l_3) P(l_5 | l_4)$
 - $-P(l_1 l_2 l_3 l_4 l_5)=P(l_1) [P(l_1 l_2)/P(l_1)] [P(l_2 l_3)/P(l_2)] [P(l_3 l_4)/P(l_3)] [P(l_4 l_5)/P(l_4)]$
- Let $f_{\alpha\beta}$ denote the frequency with which the letter pair $\alpha\beta$ is observed in text (from the file on Blackboard). Let f_{α} denote the frequency with which the letter α is observed in text. Then $f_{\alpha} = \sum_{\beta} f_{\alpha\beta}$.
- Use $f_{\alpha\beta}$ as an estimate of $P(\alpha\beta)$ and f_{α} as an estimate of $P(\alpha)$.
- Now treat it as a maximization (of likelihood) or MCMC problem (in which case you get a posterior distribution over all possible decodings).

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END