CPSC 535 Metropolis-Hastings: Applications

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- Initialization: Select deterministically or randomly $\theta = \left(\theta_1^{(0)}, ..., \theta_p^{(0)}\right).$
- Iteration i; $i \ge 1$:
 - For k = 1 : p
 - Sample $\theta_k^{(i)}$ using an MH step of proposal distribution $q_k\left(\left(\theta_{-k}^{(i)}, \theta_k^{(i-1)}\right), \theta_k'\right)$ and target $\pi\left(\theta_k \mid \theta_{-k}^{(i)}\right)$ where $\theta_{-k}^{(i)} = \left(\theta_1^{(i)}, ..., \theta_{k-1}^{(i)}, \theta_{k+1}^{(i-1)}, ..., \theta_p^{(i-1)}\right)$.

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- In 1986, Challenger exploded; the explosion being the result of an O-ring failure. It was believed to be a result of a cold weather at the departure time: 31°F.
- We have access to the data of 23 previous flights which give for flight *i*: Temperature at flight time x_i and $y_i = 1$ failure and zero otherwise (Robert & Casella, p. 15).
- We want to have a model relating Y to x. Obviously this cannot be a linear model Y = α + xβ as we want Y ∈ {0, 1}.

• We select a simple logistic regression model

$$\mathsf{Pr}\left(\left.Y=1
ight|x
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• Equivalently we have

$$\log \mathsf{it} = \log \left(\frac{\mathsf{Pr} \left(\left. Y = 1 \right| x \right)}{\mathsf{Pr} \left(\left. Y = 0 \right| x \right)} \right) = \alpha + x\beta$$

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$$\Pr\left(\left|Y=1\right|x\right) = 1 - \Pr\left(\left|Y=0\right|x\right) = \frac{\exp\left(\alpha + x\beta\right)}{1 + \exp\left(\alpha + x\beta\right)}.$$

• Equivalently we have

$$\log \operatorname{it} = \log \left(\frac{\Pr(Y = 1 | x)}{\Pr(Y = 0 | x)} \right) = \alpha + x\beta.$$

• This ensures that the response is binary.

• We follow a Bayesian approach and select $\pi(\alpha, \beta) = \pi(\alpha | b) \pi(\beta) = b^{-1} \exp(\alpha) \exp(-b^{-1} \exp(\alpha))$; i.e. exponential prior on $\exp(\alpha)$ and flat prior on β .

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- *b* is selected as the data-dependent prior such that $\mathbb{E}(\alpha) = \hat{\alpha}$ where $\hat{\alpha}$ is the MLE of α (Robert & Casella).

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- *b* is selected as the data-dependent prior such that $\mathbb{E}(\alpha) = \hat{\alpha}$ where $\hat{\alpha}$ is the MLE of α (Robert & Casella).
- As a simple proposal distribution, we use

$$q\left(\left(lpha,eta
ight),\left(lpha',eta'
ight)
ight)=\pi\left(\left.lpha'
ight|b
ight)\mathcal{N}\left(eta';eta,\widehat{\sigma}_{eta}^{2}
ight)$$

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where $\hat{\sigma}_{\beta}^2$ is the variance associated to the MLE $\hat{\beta}$.

The algorithm proceeds as follows at iteration i

• Sample
$$(\alpha^*, \beta^*) \sim \pi(\alpha | b) \mathcal{N}(\beta; \beta^{(i-1)}, \hat{\sigma}_{\beta}^2)$$
 and compute

$$\zeta((\alpha^{(i-1)}, \beta^{(i-1)}), (\alpha^*, \beta^*))$$

$$= \min\left(1, \frac{\pi(\alpha^*, \beta^* | \text{data}) \pi(\alpha^{(i-1)} | b)}{\pi(\alpha^{(i-1)}, \beta^{(i-1)} | \text{data}) \pi(\alpha^* | b)}\right)$$

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• Set
$$(\alpha^{(i)}, \beta^{(i)}) = (\alpha^*, \beta^*)$$
 with probability
 $\zeta \left(\left(\alpha^{(i-1)}, \beta^{(i-1)} \right), (\alpha^*, \beta^*) \right)$, otherwise set
 $(\alpha^{(i)}, \beta^{(i)}) = (\alpha^{(i-1)}, \beta^{(i-1)})$.

Image: A match a ma



Figure: Plots of $\frac{1}{k} \sum_{i=1}^{k} \alpha^{(i)}$ (left) and $\frac{1}{k} \sum_{i=1}^{k} \beta^{(i)}$ (right).



Figure: Histogram estimates of $p(\alpha | data)$ (left) and $p(\beta | data)$ (right).

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- The response variable y is 0 for genuine and 1 for counterfeit and the explanatory variables are
 - x^1 the length,
 - x^2 : the width of the left edge
 - x³: the width of the right edge
 - x^4 : the bottom margin witdth



Figure: Left: Plot of the status indicator versus the bottom margin width. Right: Boxplots of the bottom margin width for both counterfeit status.

• Instead of selecting a logistic link, we select a probit one here

$$\Pr(Y = 1 | x) = \Phi(x^{1}\beta_{1} + ... + x^{4}\beta_{4})$$

where

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• For *n* data, the likelihood is then given by

$$f(y_{1:n}|\beta, x_{1:n}) = \prod_{i=1}^{n} \Phi\left(x_{i}^{T}\beta\right)^{y_{i}} \left(1 - \Phi\left(x_{i}^{T}\beta\right)\right)^{1-y_{i}}.$$

• We assume a vague prior where $\beta \sim \mathcal{N}(0, 100I_4)$ and we use a simple random walk sampler with $\hat{\Sigma}$ the covariance matrix associated to the MLE (estimated using simple deterministic method).

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 - Sample $eta^* \sim \mathcal{N}\left(eta^{(i-1)}, \tau^2 \widehat{\Sigma}
 ight)$ and compute

$$\alpha\left(\beta^{(i-1)},\beta^*\right) = \min\left(1,\frac{\pi\left(\beta^*\mid y_{1:n},x_{1:n}\right)}{\pi\left(\beta^{(i-1)}\mid y_{1:n},x_{1:n}\right)}\right)$$

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• Set $\beta^{(i)} = \beta^*$ with probability $\alpha \left(\beta^{(i-1)}, \beta^* \right)$ and $\beta^{(i)} = \beta^{(i-1)}$ otherwise.

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- Set $\beta^{(i)} = \beta^*$ with probability $\alpha \left(\beta^{(i-1)}, \beta^* \right)$ and $\beta^{(i)} = \beta^{(i-1)}$ otherwise.
- Best results obtained with $\tau^2 = 1$.



Figure: Traces (left), Histograms (middle) and Autocorrelations (right) for $\left(\beta_1^{(i)},..,\beta_4^{(i)}\right)$.

• One way to monitor the performance of the algorithm of the chain $\{X^{(i)}\}\$ consists of displaying $\rho_k = cov\left[X^{(i)}, X^{(i+k)}\right] / var\left(X^{(i)}\right)$ which can be estimated from the chain, at least for small values of k.

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- Sometimes one uses an effective sample size measure

$$\mathcal{N}^{\mathsf{ess}} = \mathcal{N} \left(1 + 2 \sum_{k=1}^{N_0} \widehat{
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This represents approximately the sample size of an equivalent i.i.d. samples.

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 One should be very careful with such measures which can be very misleading.
• We found for $\mathbb{E}(\beta | y_{1:n}, x_{1:n}) = (-1.22, 0.95, 0.96, 1.15)$ so a simple plug-in estimate of the predictive probability of a counterfeit bill is

$$\widehat{p} = \Phi \left(-1.22x^1 + 0.95x^2 + 0.96x^3 + 1.15x^4
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For x = (214.9, 130.1, 129.9, 9.5), we obtain $\hat{p} = 0.59$.

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A better estimate is obtained by

$$\int \Phi \left(\beta_{1} x^{1} + \beta_{2} x^{2} + \beta_{3} x^{3} + \beta_{4} x^{4}\right) \pi \left(\beta | y_{1:n}, x_{1:n}\right) d\beta$$

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- Introduce the following unobserved latent variables

$$\begin{aligned} & Z_i \quad \sim \quad \mathcal{N}\left(x_i^\mathsf{T}\beta,1\right), \\ & Y_i \quad = \quad \left\{ \begin{array}{cc} 1 & \text{if } Z_i > 0 \\ 0 & \text{otherwise.} \end{array} \right. \end{aligned}$$

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• We have now define a joint distribution

$$f(y_i, z_i | \beta, x_i) = f(y_i | z_i) f(z_i | \beta, x_i).$$

• Now we can check that

$$f(y_i = 1 | x_i, \beta) = \int f(y_i, z_i | \beta, x_i) dz_i$$

=
$$\int_0^\infty f(z_i | \beta, x_i) dz_i = \Phi(x_i^{\mathsf{T}} \beta).$$

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 \Rightarrow We haven't changed the model!

 We are now going to sample from π (β, z_{1:n} | x_{1:n}, y_{1:n}) instead of π (β | x_{1:n}, y_{1:n}) because the full conditional distributions are simple

$$\pi \left(\beta | y_{1:n}, x_{1:n}, z_{1:n} \right) = \pi \left(\beta | x_{1:n}, z_{1:n} \right) \text{ (standard Gaussian!),}$$

$$\pi \left(z_{1:n} | y_{1:n}, x_{1:n}, \beta \right) = \prod_{i=1}^{n} \pi \left(z_{k} | y_{k}, x_{k}, \beta \right)$$

where

$$z_k | y_k, x_k, \beta \sim \begin{cases} \mathcal{N}_+ \left(x_k^\mathsf{T} \beta, 1
ight) & ext{if } y_k = 1 \\ \mathcal{N}_- \left(x_k^\mathsf{T} \beta, 1
ight) & ext{if } y_k = 0. \end{cases}$$



Figure: Traces (left), Histograms (middle) and Autocorrelations (right) for $\left(\beta_1^{(i)},..,\beta_4^{(i)}\right)$.

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- Very similar were also obtained using a logistic fonction using the MH (Gibbs is feasible but more difficult).

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- Consider the following simple generalization of the previous model

$$Z_i \sim \mathcal{N}\left(x_i\beta, \sigma^2\right)$$
, $Y_i = \begin{cases} 1 & \text{if } Z_i > 0 \\ 0 & \text{otherwise.} \end{cases}$

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ight.$

• We complete the model by $\sigma^2 \sim \mathcal{IG}(1.5, 1.5)$ and $\beta | \sigma^2 \sim \mathcal{N}(0, 100)$.

• Not only the data Z_i and (eta,σ^2) are very correlated but we have

$$\Pr\left(\left|Y_{i}=1\right| x_{i},\beta,\sigma^{2}\right) = \Phi\left(\frac{x_{i}\beta}{\sigma}\right)$$

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- The likelihood only depends on β/σ and the parameters β and σ are not identifiable.
- One way to improve the mixing consists of using an additional MH step that proposes to randomly rescale the current value.
- We use a proposal distribution such that

$$\left(eta^{\prime},\sigma^{\prime}
ight)=\lambda\left(eta,\sigma
ight)$$
 with $\lambda\sim\mathcal{E} ext{xp}\left(1
ight)$

that proposes to randomly rescale the current value.

• Consider the following hidden Markov model

$$\begin{array}{rcl} X_k \left| \left(X_{k-1} = x_{k-1} \right) & \sim & f_\theta \left(\cdot \mid x_{k-1} \right), \ X_1 \sim \mu \\ Y_n \left| \left(X_k = x_k \right) & \sim & g_\theta \left(\cdot \mid x_k \right), \end{array} \right. \end{array}$$

and we set a prior $\pi\left(\theta\right)$ on the unknown hyperparameters $\theta.$

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and we set a prior $\pi(\theta)$ on the unknown hyperparameters θ . • Given *n* data, we are interested in the joint posterior

 $\pi(\theta, x_{1:n}|y_{1:n})$

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and we set a prior $\pi(\theta)$ on the unknown hyperparameters θ . • Given *n* data, we are interested in the joint posterior

$$\pi\left(\left.\theta, x_{1:n}\right| y_{1:n}\right)$$

• There is no closed-form expression for this joint distribution even if the model is linear Gaussian or for finite state-space model.

 In previous lectures, we propose sampling from π (θ, x_{1:n} | y_{1:n}) using the Gibbs sampler where the variables are updated according to

$$X_k \sim \pi(x_k | y_{1:n}, x_{-k}, \theta)$$

 In previous lectures, we propose sampling from π (θ, x_{1:n} | y_{1:n}) using the Gibbs sampler where the variables are updated according to

$$X_k \sim \pi(x_k | y_{1:n}, x_{-k}, \theta)$$

• For 2 < k < n, we have

$$\pi \left(x_{k} \middle| y_{1:n}, x_{-k}, \theta \right) \propto \pi \left(x_{1:n}, y_{1:n}, \theta \right)$$

$$\propto \underbrace{\pi \left(\theta \right) \mu \left(x_{1} \right) \prod_{i=2}^{n} f_{\theta} \left(x_{i} \middle| x_{i-1} \right) \prod_{i=1}^{n} g_{\theta} \left(y_{i} \middle| x_{i} \right)}_{\text{prior}}$$

$$\underset{\text{likelihood}}{\underset{\text{like}}{\underset{\text{like}}}}}}} x_{like}}} x_{like}} x_{like} x_{like}} x_{like}} x_{l$$

and $\theta \sim \pi(\theta | y_{1:n}, x_{1:n})$ (or by subblocks).

• It is often possible to implement the Gibbs sampler even if this can be expensive; e.g. if you use Accept/Reject to sample from $\pi(x_k | y_{1:n}, x_{-k}, \theta)$ using the proposal $\pi(x_k | x_{-k}, \theta) \propto f_{\theta}(x_k | x_{k-1}) f_{\theta}(x_{k+1} | x_k)$.

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 π (x_k | y_{1:n}, x_{-k}, θ) using the proposal
 π (x_k | x_{-k}, θ) ∝ f_θ (x_k | x_{k-1}) f_θ (x_{k+1} | x_k).
- Even if it is possible to implement the Gibbs sampler, one can expect a very slow convergence of the algorithm is successive variables are highly correlated.
- Indeed, as you update x_k with x_{k-1} and x_{k+1} being fixed, then you cannot move much into the space.

• Consider the very simple case where $heta=\left(\sigma_{v}^{2},\sigma_{w}^{2}
ight)$

$$\begin{array}{rcl} X_k & = & X_{k-1} + V_k \text{ where } V_k \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0, \sigma_v^2\right), \\ Y_k & = & X_k + W_k \text{ where } W_k \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0, \sigma_w^2\right) \end{array}$$

then we have

$$\begin{aligned} \pi\left(x_{k} \mid x_{-k}, \theta\right) &\propto \quad f_{\theta}\left(x_{k} \mid x_{k-1}\right) f_{\theta}\left(x_{k+1} \mid x_{k}\right) \\ &= \quad \mathcal{N}\left(x_{k}; \frac{x_{k-1} + x_{k+1}}{2}, \frac{\sigma_{\nu}^{2}}{2}\right) \end{aligned}$$

and

$$\begin{aligned} &\pi\left(x_{k} \mid y_{1:n}, x_{-k}, \theta\right) \\ &\propto &\pi\left(x_{k} \mid x_{-k}, \theta\right) g_{\theta}\left(y_{k} \mid x_{k}\right) \\ &= &\mathcal{N}\left(x_{k}; \frac{\sigma_{v}^{2} \sigma_{w}^{2}}{\sigma_{v}^{2} + 2\sigma_{w}^{2}} \left(\frac{x_{k-1} + x_{k+1}}{\sigma_{v}^{2}} + \frac{y_{k}}{\sigma_{w}^{2}}\right), \frac{\sigma_{v}^{2} \sigma_{w}^{2}}{\sigma_{v}^{2} + 2\sigma_{w}^{2}}\right) \end{aligned}$$

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• Assume for the time being that instead of sampling from $\pi(x_k | y_{1:n}, x_{-k}, \theta)$ directly, we use rejection sampling with $\pi(x_k | x_{-k}, \theta)$ as a proposal distribution.

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- In this case we have to bound

$$g_{\theta}\left(\left.y_{k}\right|x_{k}
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- In this case we have to bound

$$g_{\theta}\left(\left|x_{k}\right| x_{k}\right) = \frac{1}{\sqrt{2\pi}\sigma_{w}} \exp\left(-\frac{\left(y_{k}-x_{k}\right)^{2}}{2\sigma_{w}^{2}}\right) \leq \frac{1}{\sqrt{2\pi}\sigma_{w}}.$$

• We accept each proposal $X^* \sim \pi(x_k | x_{-k}, \theta)$ with probability $\exp\left(-\frac{(y_k - X^*)^2}{2\sigma_w^2}\right)$, so the (unconditional) acceptance probability is given by

$$\int \pi (x_k | x_{-k}, \theta) \exp\left(-\frac{(y_k - x_k)^2}{2\sigma_w^2}\right) dx_k \\ = \frac{\sigma_w \exp\left(-\frac{1}{2}\left(\frac{y_k^2}{\sigma_w^2} - (x_{k-1} + x_{k+1})^2 / \sigma_v^2\right)\right)}{\sqrt{\sigma_v^2 + 2\sigma_w^2}}.$$

 To improve the algorithm, we would like to be able to sample a whole block of variables simultaneously; i.e. being able to sample for 1 < k < k + L < n from

$$\pi \left(x_{k:k+L} | y_{1:n}, x_{-(k:k+L)}, \theta \right) = \pi \left(x_{k:k+L} | y_{k:k+L}, x_{k-1}, x_{k+L+1}, \theta \right)$$
$$\propto \prod_{i=k}^{k+L+1} f_{\theta} \left(x_i | x_{i-1} \right) \prod_{i=k}^{k+L} g_{\theta} \left(y_i | x_i \right).$$

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• In this case, it is typically impossible to sample from $\pi\left(x_{k:k+L} \mid y_{1:n}, x_{-(k:k+L)}, \theta\right)$ exactly as *L* is large, say 5 or 10.

 To improve the algorithm, we would like to be able to sample a whole block of variables simultaneously; i.e. being able to sample for 1 < k < k + L < n from

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In this case, it is typically impossible to sample from π (x_{k:k+L} | y_{1:n}, x_{-(k:k+L}), θ) exactly as L is large, say 5 or 10.
We are propose to use a MH step of invariant distribution π (x_{k:k+L} | y_{1:n}, x_{-(k:k+L}), θ) instead, hence we need to build a

proposal distribution $q\left((x_{1:n}, \theta), x'_{k:k+L}\right)$.

• We first propose to use the conditional prior

$$q((x_{1:n},\theta), x'_{k:k+L}) = \pi \left(x_{k:k+L} | x_{-(k:k+L)}, \theta \right)$$

= $\pi \left(x_{k:k+L} | x_{k-1}, x_{k+L+1}, \theta \right)$
 $\propto \prod_{i=k}^{k+L+1} f_{\theta}(x_i | x_{i-1}).$

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• In this case, the candidate $X'_{k:k+L} \sim \pi(x_{k:k+L}|x_{k-1}, x_{k+L+1}, \theta)$ is accepted with probability

$$\min \left(1, \frac{\pi \left(x_{k:k+L}' \middle| y_{k:k+L}, x_{k-1}, x_{k+L+1}, \theta \right) \pi \left(x_{k:k+L} \middle| x_{k-1}, x_{k+L+1}, \theta \right)}{\pi \left(x_{k:k+L} \middle| y_{k:k+L}, x_{k-1}, x_{k+L+1}, \theta \right) \pi \left(x_{k:k+L}' \middle| x_{k-1}, x_{k+L+1}, \theta \right)} \right)$$

= min $\left(1, \frac{\prod_{i=k}^{k+L} g_{\theta}(y_i | x_i')}{\prod_{i=k}^{k+L} g_{\theta}(y_i | x_i)} \right)$
• We first propose to use the conditional prior

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= min $\left(1, \frac{\prod_{i=k}^{k+L} g_{\theta}(y_i | x'_i)}{\prod_{i=k}^{k+L} g_{\theta}(y_i | x_i)} \right)$

• Simple but one cannot expect it to be too efficient when the observations are very informative compared to the prior.

• Consider the case where

$$X_{k} = AX_{k-1} + BV_{k}, V_{k} \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}(0, I)$$
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• Particular cases include

$$\begin{array}{ll} X_{k} & = & X_{k-1} + \sigma V_{k}, \text{ where } V_{k} \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0,1\right), \\ X_{k} & = & \left(\begin{array}{c} \alpha_{k} \\ \alpha_{k-1} \end{array}\right) = \left(\begin{array}{c} 2 & -1 \\ 1 & 0 \end{array}\right) X_{k-1} + \left(\begin{array}{c} \sigma \\ 0 \end{array}\right) V_{k}, \ V_{k} \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0,1\right) \end{array}$$

• In this case, it is simple to see that $\pi(x_{k:k+L}|x_{k-1}, x_{k+1}, \theta)$ is a Gaussian distribution.

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- In this case, it is simple to see that $\pi(x_{k:k+L}|x_{k-1}, x_{k+1}, \theta)$ is a Gaussian distribution.
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- In this case, it is simple to see that $\pi(x_{k:k+L}|x_{k-1}, x_{k+1}, \theta)$ is a Gaussian distribution.
- In (Knorr-Held, 1999), one samples from this distribution by computing directly the parameters of this joint distribution: complexity $O(L^2)$.
- We can derive a simpler method of complexity O(L) based on the following decomposition (omitting θ in the notation)

$$\pi (x_{k:k+L} | x_{k-1}, x_{k+L+1}) = \prod_{i=k}^{k+L} \pi (x_i | x_{k-1}, x_{k+L+1}, x_{i+1}).$$

=
$$\prod_{i=k}^{k+L} \pi (x_i | x_{k-1}, x_{i+1})$$

• Moreover it is easy to establish the expression for $\pi(x_i | x_{k-1}, x_{i+1})$

$$\pi(x_i|x_{k-1}, x_{i+1}) \propto \pi(x_i|x_{k-1}) f(x_{i+1}|x_i)$$

as

$$\pi(x_{i}|x_{k-1}) = \int \pi(x_{k:i}|x_{k-1}) dx_{k:i-1} = \mathcal{N}(x_{i}; \mu_{i}(x_{k-1}), \Sigma_{i})$$

with, for $X_n = AX_{n-1} + BV_n$, $\mu_{k-1}\left(x_{k-1}
ight) = x_{k-1}$, $\Sigma_{k-1} = 0$ and for $i \geq k$

$$\begin{split} \mu_i \left(x_{k-1} \right) &= & A \mu_{i-1} \left(x_{k-1} \right), \\ \Sigma_i &= & A \Sigma_{i-1} A^\mathsf{T} + \Sigma \text{ with } \Sigma = B B^\mathsf{T}. \end{split}$$

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• To obtain $\pi(x_i | x_{k-1}, x_{i+1})$, we combine the prior $\pi(x_i | x_{k-1})$ with the "likelihood" $f(x_{i+1} | x_i)$.

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• We have
$$\pi(x_i | x_{k-1}) = \mathcal{N}(x_i; \mu_i(x_{k-1}), \Sigma_i)$$
 and
 $f(x_{i+1} | x_i) = \mathcal{N}(x_{i+1}; Ax_i, \Sigma)$ then
 $\pi(x_i | x_{k-1}, x_{i+1}) = \mathcal{N}(x_i; \mu_i(x_{k-1}, x_{i+1}), \widetilde{\Sigma}_i)$

where

$$\widetilde{\Sigma}_{i} = \left(\Sigma_{i}^{-1} + A^{\mathsf{T}}\Sigma^{-1}A\right)^{-1},$$

$$\mu_{i}\left(x_{k-1}, x_{i+1}\right) = \widetilde{\Sigma}_{i}\left(A^{\mathsf{T}}\Sigma^{-1}x_{i+1} + \Sigma_{i}^{-1}\mu_{i}\left(x_{k-1}\right)\right).$$

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• To sample a realization of $\pi(x_{k:k+L}|x_{k-1}, x_{k+L+1})$, first compute $\mu_i(x_{k-1})$, Σ_i for i = k, ..., k + L using a forward recursion. Then sample backward $X_{k+L} \sim \pi(\cdot|x_{k-1}, x_{k+L+1})$, $X_{k+L-1} \sim \pi(\cdot|x_{k-1}, X_{k+L})$, ..., $X_k \sim \pi(\cdot|x_{k-1}, X_{k+1})$.



Figure: Number of occurences of rainfall in Tokyo for each day during 1983-1984 reproduced as relative frequencies between 0, 0.5 and 1 (n = 366)

• Consider the following model

$$X_{k} = \begin{pmatrix} \alpha_{k} \\ \alpha_{k-1} \end{pmatrix} = \begin{pmatrix} 2 & -1 \\ 1 & 0 \end{pmatrix} X_{k-1} + \begin{pmatrix} \sigma \\ 0 \end{pmatrix} V_{k}, V_{k} \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}(0,1)$$

and

$$Y_k | \, X_k \sim \left\{egin{array}{cc} B\left(2, \pi_k
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• We also use for $\sigma^2 \sim \mathcal{IG}\left(rac{\nu_0}{2}, rac{\gamma_0}{2}
ight)$.

.

 We use the block sampling strategies discussed before where candidates are sampled according to π (x_{k:k+L} | x_{k-1}, x_{k+L+1}) and accepted with proba

$$\min\left(1,\frac{\prod_{i=k}^{k+L}g\left(y_{i}\mid x_{i}'\right)}{\prod_{i=k}^{k+L}g\left(y_{i}\mid x_{i}\right)}\right).$$

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 $\bullet\,$ The parameter σ^2 is updated through a simple Gibbs step

$$\begin{aligned} \sigma^2 &\sim & \pi \left(\sigma^2 \big| \, \mathbf{x}_{1:n}, \, \mathbf{y}_{1:n} \right) = \pi \left(\sigma^2 \big| \, \mathbf{x}_{1:n} \right) \\ &= & \mathcal{IG} \left(\frac{\nu_0 + n - 1}{2}, \frac{\gamma_0 + \sum_{k=2}^n \left(\alpha_k - 2\alpha_{k-1} + \alpha_{k-2} \right)^2}{2} \right) \end{aligned}$$

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For block size L = 1, 5, 20 and 40, we compute the average trajectories of 100 parallel chains after 10, 50, 100 and 500 iterations with initialization x_k = 0 for all k, σ² = 0.1.





Figure: Average trajectories over 100 chains for L = 1, 5, 20 and 40 from top to bottom.

After 50 Iterations



Figure: Average trajectories over 100 chains for L = 1, 5, 20 and 40 from top to bottom.

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Figure: Average trajectories over 100 chains for L = 1, 5, 20 and 40 from top to bottom.



Figure: Average trajectories over 100 chains for L = 1, 5, 20 and 40 from top to bottom.



Figure: Traces of α_1 , α_{100} , α_{333} and σ^2 for L = 1 (left) and L = 20 (right).

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- For a linear Gaussian observation equation, Knorr-Held compares this strategy to a direct Gibbs sampling implementation. As expected, the conditional proposal strategy is competitive when the observations are not very informative compared to the prior.

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- For a linear Gaussian observation equation, Knorr-Held compares this strategy to a direct Gibbs sampling implementation. As expected, the conditional proposal strategy is competitive when the observations are not very informative compared to the prior.
- For more complex problems, such strategies are inefficient and we will need to use the observations to build the proposal.

• (Pitt & Shephard, 1999) propose a more efficient strategy... also more computationally intensive.

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- (Pitt & Shephard, 1999) propose a more efficient strategy... also more computationally intensive.
- Consider the log full conditional distribution

$$\begin{split} &\log \pi \left(x_{k:k+L} | y_{k:k+L}, x_{k-1}, x_{k+L+1} \right) \\ &= \sum_{i=k}^{k+L} \log g \left(y_i | x_i \right) + \sum_{i=k}^{k+L+1} \log f \left(x_{i+1} | x_i \right) \\ &\equiv \sum_{i=k}^{k+L} \log g \left(y_i | x_i \right) - \frac{1}{2} \sum_{i=k}^{k+L+1} \left(x_{i+1} - Ax_i \right)^{\mathsf{T}} \Sigma^{-1} \left(x_{i+1} - Ax_i \right) \end{split}$$

which is not quadratic in x_i hence $\pi(x_{k:k+L}|y_{k:k+L}, x_{k-1}, x_{k+1})$ is not Gaussian.

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which is not quadratic in x_i hence $\pi(x_{k:k+L}|y_{k:k+L}, x_{k-1}, x_{k+1})$ is not Gaussian.

The idea is to expand the log-likelihood part around some point estimates

$$\log g(y_i|x_i) \simeq \log g(y_i|\hat{x}_i) + \nabla \log g(y_i|\hat{x}_i) \cdot (x_i - \hat{x}_i) + \frac{1}{2} (x_i - \hat{x}_i)^{\mathsf{T}} \nabla^2 \log g(y_i|\hat{x}_i) (x_i - \hat{x}_i)$$

• By doing this, we have a Gaussian approximation of the log-likelihood and then we obtain a Gaussian proposal $q(x_{1:n}, x'_{k:k+L}) = q(x_{-(k:k+L)}, x'_{k:k+L})$

$$\log q\left(x_{-(k:k+L)}, x'_{k:k+L}\right) \equiv \sum_{i=k}^{k+L} \nabla \log g\left(y_i | \hat{x}_i\right) \cdot \left(x_i - \hat{x}_i\right) \\ + \frac{1}{2} \left(x_i - \hat{x}_i\right)^{\mathsf{T}} \nabla^2 \log g\left(y_i | \hat{x}_i\right) \left(x_i - \hat{x}_i\right) \\ - \frac{1}{2} \sum_{i=k}^{k+L+1} \left(x_{i+1} - Ax_i\right)^{\mathsf{T}} \Sigma^{-1} \left(x_{i+1} - Ax_i\right)$$

• By doing this, we have a Gaussian approximation of the log-likelihood and then we obtain a Gaussian proposal $q(x_{1:n}, x'_{k:k+L}) = q(x_{-(k:k+L)}, x'_{k:k+L})$

$$\log q\left(x_{-(k:k+L)}, x'_{k:k+L}\right) \equiv \sum_{i=k}^{k+L} \nabla \log g\left(y_{i} | \hat{x}_{i}\right) \cdot \left(x_{i} - \hat{x}_{i}\right)$$
$$+ \frac{1}{2} \left(x_{i} - \hat{x}_{i}\right)^{\mathsf{T}} \nabla^{2} \log g\left(y_{i} | \hat{x}_{i}\right) \left(x_{i} - \hat{x}_{i}\right)$$
$$- \frac{1}{2} \sum_{i=k}^{k+L+1} \left(x_{i+1} - Ax_{i}\right)^{\mathsf{T}} \Sigma^{-1} \left(x_{i+1} - Ax_{i}\right)$$

(Pitt & Shepard, 1999) propose to select

$$\hat{x}_{k:k+1} = \arg \max \pi (x_{k:k+L} | y_{k:k+L}, x_{k-1}, x_{k+L+1})$$

and a scheme to sample from $q\left(x_{-(k:k+L)}, x'_{k:k+L}\right)$ which is of complexity O(L).

• This algorithm is applied to the SV model where

$$\begin{aligned} X_k &= \phi X_{k-1} + \sigma V_k, \ V_k \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0,1\right) \\ Y_k &= \beta \exp\left(X_k/2\right) W_k, \ W_k \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}\left(0,1\right). \end{aligned}$$

Image: A math a math

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- Compared to standard single move strategies, the authors report significant improvement.



Figure: Autocorrelation plots for (ϕ, σ^2, β) with L = 1



Figure: Autocorrelation plots for (ϕ, σ^2, β) with L = 50 on average