

Further simulations issues

- Bayesian inference is based on properties of posterior.
- We have seen that MCMC (the MH algorithm/Gibbs sampler) is a general tool which will (at least in theory) explore the posterior.
- In the remainder of this lecture we will briefly touch upon cases where the methods considered so far will work very inefficiently — or not at all.

Intractable normalising constants

Setup: Consider situation where the data model

$$\pi(x|\theta) = \frac{1}{c(\theta)} f(x|\theta),$$

where $c(\theta) = \int f(x|\theta)dx$ is the normalising constant which depends on the parameter θ .

Assume that is either impossible or infeasible to calculate $c(\theta)$.

Example: The **Ising model** is one such model.

Here $c(\theta)$ is obtained by a sum over all possible pixel images. For even moderately large pixel images, this becomes infeasible.

Posterior inference and intractable normalising constants

- Assume $\pi(x|\theta) = \frac{1}{c(\theta)} f(x|\theta)$ where $c(\theta)$ is intractable.
- We want to perform Bayesian inference, i.e. we want to sample the posterior $\pi(\theta|x) \propto \pi(x|\theta)\pi(\theta)$ using a MH algorithm.
- Sampling the posterior using the MH algorithm we have:

$$H(\theta, \theta') = \frac{\pi(x|\theta')\pi(\theta')q(\theta', \theta)}{\pi(x|\theta)\pi(\theta)q(\theta, \theta')} = \frac{\frac{1}{c(\theta')} f(x|\theta')\pi(\theta')q(\theta', \theta)}{\frac{1}{c(\theta)} f(x|\theta)\pi(\theta)q(\theta, \theta')}$$

Notice that a ratio $c(\theta)/c(\theta')$ appears.

Intractable normalising constants: One solution

- Assume that $\pi(x|\theta) > 0$ implies $\pi(x|\theta') > 0$ for all pairs θ, θ' .
- Then we have the following importance sampling identity:

$$\frac{c(\theta)}{c(\theta')} = \mathbb{E}_{\theta'} \left[\frac{f(X|\theta)}{f(X|\theta')} \right],$$

where $\mathbb{E}_{\theta'}$ is expectation wrt. $\pi(x|\theta')$.

- Hence the ratio $\frac{c(\theta)}{c(\theta')}$ can be estimated using a MH algorithm (say) with $\pi(x|\theta')$ as the invariant distribution.
- That is, for each update of the main MH algorithm we need to run an addition MH algorithm to estimating the Hastings ratio in the main algorithm.
- There are number of alternative solutions to one sketched above.

Likelihood free Bayesian inference

Setup: We do not have an expression for the likelihood $\pi(x|\theta)$ — not even $f(x|\theta)$.

BUT, given θ , we can generate $x|\theta \sim \pi(x|\theta)$.

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BUT, given θ , we can generate $x|\theta \sim \pi(x|\theta)$.

We can simulate from the posterior, $\theta|x \sim \pi(\theta|x)$, as follows:

- * Repeat steps 1 and 2...
 - 1 Generate $\theta \sim \pi(\theta)$
 - 2 Generate $\tilde{x}|\theta \sim \pi(x|\theta)$
- * Until $\tilde{x} = x$.
- * Return θ

The repeat loop generates prior predictions until \tilde{x} matches the data x (exactly).

In most situations of interest the probability of $\tilde{x} = x$ is very small — maybe zero. In other words: the algorithm does not work in practise.

Approximate Bayesian Computations (ABC)

The idea is to make an approximation of the previous algorithm:

Assume we have *distance* measue $d(x, x')$ which measures the difference between two data sets.

We can simulate from the posterior, $\theta|x \sim \pi(\theta|x)$, as follows:

- * Repeat steps 1 and 2...
- 1 Generate $\theta \sim \pi(\theta)$
- 2 Generate $\tilde{x}|\theta \sim \pi(x|\theta)$
- * Until $d(\tilde{x}, x) < \epsilon$.
- * Return θ

The repeat loop generates prior predictions until \tilde{x} *approximatly* matches the data x .

If $\epsilon = \infty$: We sample the prior.

If $\epsilon = 0$: We sample the posterior.

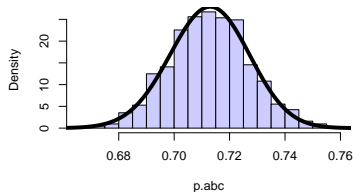
ABC: Binomial model

- **Data model:** $X \sim B(n, p)$, n known
 - **Data:** x the number of successes
 - **Prior:** $\pi(p) = Be(\alpha, \beta)$ — beta distribution
 - **Posterior:** $\pi(p|x) = Be(\alpha + x, \beta + n - x)$
-
- **Distance:** $d(x, x') = \left| \frac{x}{n} - \frac{x'}{n} \right|$

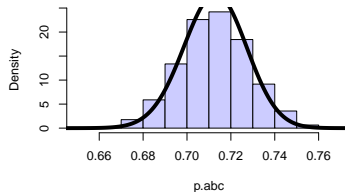
ABC: Binomial model —*cont.*

In example: $n = 1000$ and $x = 713$ and 2500 samples

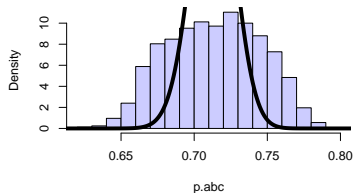
epsilon = 0 accept.prob. = 0.001



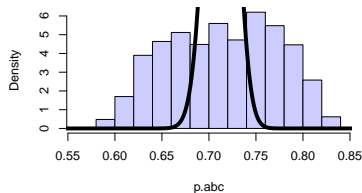
epsilon = 0.01 accept.prob. = 0.02



epsilon = 0.05 accept.prob. = 0.1018



epsilon = 0.1 accept.prob. = 0.2



A Metropolis-Hastings style ABC algorithm

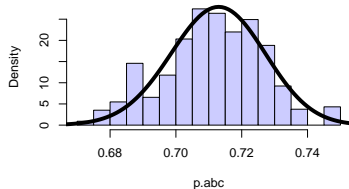
- Choose initial value $\theta^{(0)}$
- For $t = 1, \dots, T$
 - Generate $\theta' \sim q(\theta^{(t-1)}, \theta')$
 - Generate $x' | \theta' \sim \pi(x' | \theta')$
 - If $d(x, x') > \epsilon$ reject and set $\theta^{(t)} = \theta^{(t-1)}$.
 - If $d(x, x') \leq \epsilon$
 - Calculate $H(\theta^{(t-1)}, \theta') = \frac{\pi(\theta')}{\pi(\theta^{(t-1)})} \frac{q(\theta', \theta^{(t-1)})}{q(\theta^{(t-1)}, \theta')}$
 - Generate $u \sim \text{Unif}([0, 1])$.
 - If $u < H(\theta^{(t-1)}, \theta')$ set $\theta^{(t)} = \theta'$ else set $\theta^{(t)} = \theta^{(t-1)}$

Notice the Hastings ratio only involves the priors.

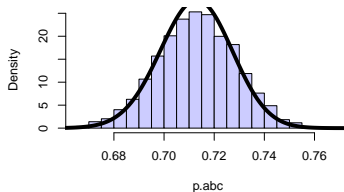
ABC: Binomial model —*cont.*

In example: $n = 1000$ and $x = 713$ and 25,000 iterations.

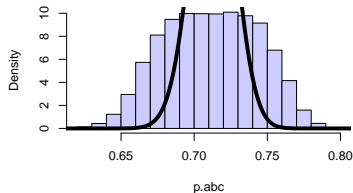
epsilon = 0 accept.prob. = 0.0141



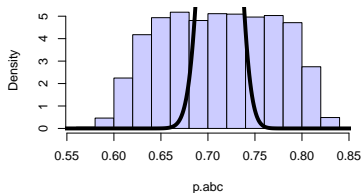
epsilon = 0.01 accept.prob. = 0.2881



epsilon = 0.05 accept.prob. = 0.7826



epsilon = 0.1 accept.prob. = 0.8906



Mixture models

The density of the data model may take the form

$$\pi(y|\theta) = \lambda_1\pi_1(y|\theta_1) + \lambda_2\pi_2(y|\theta_1) + \cdots + \lambda_k\pi_k(y|\theta_1)$$

where each $\pi_j(y|\theta_j)$ is a normalised density and $\lambda_1, \dots, \lambda_k \geq 0$ are weight with $\sum_{j=1}^k \lambda_j = 1$.

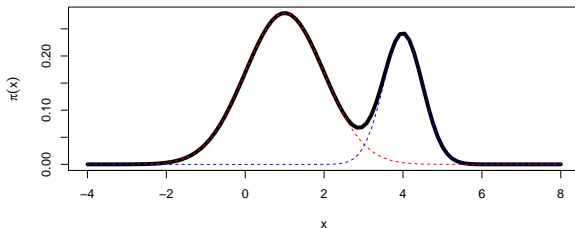
This is referred to as a (k component) mixture distribution.

Mixture model: example

Mixture distribution consisting of two normal distributions, ie.

$$\pi_j(y; \theta_j) = \mathcal{N}(y; \mu_j, \sigma_j^2):$$

- Component 1: $\lambda_1 = 0.7$ $\mu_1 = 1$ $\sigma_1^2 = 1$
- Component 2: $\lambda_2 = 0.3$ $\mu_2 = 4$ $\sigma_2^2 = 0.25$



Notice that the likelihood is *exactly* the same if we swap $(\lambda_1, \mu_1, \sigma_1^2)$ and $(\lambda_2, \mu_2, \sigma_2^2)$. In other words: the model is symmetric in the mixture components.

Dirichlet distribution

To perform a Bayesian analysis we need a prior for the unknown weights. A k dimensional random vector $\theta = (\theta_1, \dots, \theta_k)$ is said to follow a Dirichlet distribution with parameters $\alpha = (\alpha_1, \dots, \alpha_k)$, $\alpha_i > 0$, if it has density

$$\pi(\theta|\alpha) \propto \prod_{j=1}^k \theta_j^{\alpha_j-1},$$

where $\theta_j \geq 0$ and $\sum_{j=1}^k \theta_j = 1$.

As a prior for λ we typically use a *Dirichlet*(1, 1, ..., 1) prior, ie. a uniform distribution on the allowed set of weights.

The posterior

We assume a priori that the parameters for each component are independent and independent of the weight:

$$\pi(\theta, \lambda) = \prod_{j=1}^k \pi(\theta_j) \pi(\lambda)$$

The posterior is then

$$\pi(\theta, \lambda | y) \propto \pi(\theta, \lambda) \prod_{i=1}^n \sum_{j=1}^k \pi_j(y_i | \theta_j) \lambda_j$$

The full conditional for θ_j is

$$\pi(\theta_j | \theta_{-j}, \lambda, y) \propto \pi(\theta_j) \prod_{i=1}^n \sum_{j=1}^k \pi_j(y_i | \theta_j) \lambda_j$$

Problem: Even in a simple setup with normal mixtures this is a non standard distribution — which quickly becomes intractable.

Solution: Make the model more complicated!

Dummy variables

For each observation y_i , introduce dummy variable $z_i \in \{1, \dots, k\}$ which indicates which component x_i belongs to.

$$\pi(y_i | z_i = j, \theta) = \pi_l(y_i | \theta_j)$$

Let λ_j be the a priori probability that y_i belongs to the j th component. Hence $\pi(z_i = j | \lambda) = \lambda_j$.

This can be combined to

$$\begin{aligned}\pi(y_i, z_i | \lambda, \theta) &= \pi(y_i | z_i, \theta, \lambda) \pi(z_i | \theta, \lambda) \\ &= \pi(y_i | z_i, \theta) \pi(z_i | \lambda)\end{aligned}$$

Next: Verify that the introduction of the z_i s does not change the model.

Marginal distributiun of y_i

The marginal disitribution of y_i is

$$\begin{aligned}\pi(y_i|\theta) &= \sum_{j=1}^k \pi(y_i, z_i = j|\lambda, \theta) \\ &= \sum_{j=1}^k \pi(y_i|z_i = j, \theta) \pi(z_i = j|\lambda) \\ &= \sum_{j=1}^k \pi_j(y_i|\theta_j) \lambda_j\end{aligned}$$

Hence, the marginal distribution af y_i is unaffected by the introduction of the indicator variables.

Rewriting the likelihood

Notice that

$$\pi(y_i | z_i = l, \theta) = \pi_l(y_i | \theta_l) = \prod_{j=1}^k \pi_j(y_i | \theta_j)^{1_{[z_i=l]}}$$

Similarly

$$\pi(z_i = l | \lambda) = \lambda_l = \prod_{j=1}^k \lambda_j^{1_{[z_i=l]}}$$

Likelihood

$$\pi(y, z | \lambda, \theta) = \prod_{i=1}^n \prod_{j=1}^k \left(\pi_j(y_i | \theta_j) \lambda_j \right)^{1_{[z_i=j]}}$$

Full conditionals: z_i

The full conditional for the dummy variables:

$$\begin{aligned}\pi(z_s = l | y, \theta, \lambda, z_{-s}) &\propto \pi(\theta, \lambda) \prod_{i=1}^n \prod_{j=1}^k (\pi_j(y_i | \theta_j) \lambda_j)^{1[z_i=j]} \\ &\propto \pi(\theta, \lambda) \prod_{j=1}^k (\pi_j(y_s | \theta_j) \lambda_j)^{1[z_s=j]} \\ &\propto \pi_l(y_s | \theta_l) \lambda_l\end{aligned}$$

Normalising the probability we obtain

$$\pi(z_s = l | y, \theta, \lambda, z_{-s}) = \frac{\pi_l(y_s | \theta_l) \lambda_l}{\sum_{j=1}^k \pi_j(y_s | \theta_j) \lambda_j}$$

This is a simple distribution to sample from.

Full conditionals: θ_l

The full conditional for θ_j is

$$\begin{aligned}\pi(\theta_l | \theta, y, z) &\propto \pi(\theta, \lambda) \prod_{i=1}^n \prod_{j=1}^k (\pi_j(y_i | \theta_j) \lambda_j)^{1_{[z_i=j]}} \\ &\propto \pi(\theta_l) \prod_{i=1}^n \pi_l(y_i | \theta_l)^{1_{[z_i=l]}} \\ &\propto \pi(\theta_l) \prod_{i: z_i=l} \pi_l(y_i | \theta_l)\end{aligned}$$

This is equivalent to the posterior in the case of independent observations from π_l (restricted to observation for the l th component). If the mixture components are normal and we choose priors as in earlier lectures, we know how to sample this full conditional.

Full conditionals: λ

The (joint) full conditional distribution of λ is

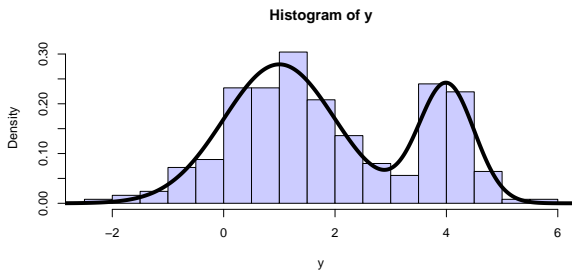
$$\begin{aligned}\pi(\lambda|\theta, y, z) &\propto \prod_{i=1}^n \prod_{j=1}^k (\pi_j(y_i|\theta_j)\lambda_j)^{1_{[z_i=j]}} \pi(\theta, \lambda) \\ &\propto \prod_{i=1}^n \prod_{j=1}^k \lambda_j^{1_{[z_i=j]}} \pi(\lambda) \\ &\propto \prod_{j=1}^k \lambda_j^{n_j(z)} \prod_{j=1}^k \lambda_j^{\alpha_j-1} \\ &\propto \text{Dirichlet}(n_1(z) + \alpha_1, \dots, n_k(z) + \alpha_k),\end{aligned}$$

where $n_j(z)$ is the number of dummy variables equal to j .

Sample model

Generate a sample of size 250 from the mixture distribution:

```
N = 250  
z = sample(size = N, x = 1:2, prob = lambda, replace = TRUE)  
y = rnorm(N, mean = mu[z], sd = sd[z])
```

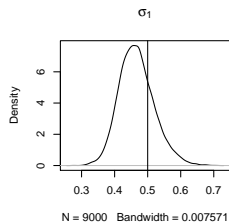
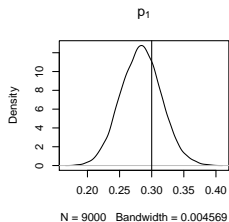
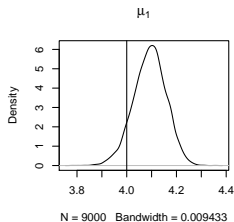


```
model{
  # Likelihood:
  for(i in 1 : N){
    y[i] ~ dnorm( mu[z[i]] , tau[z[i]] )
    z[i] ~ dcat( p[1:2])
  }
  # Prior:
  for ( j in 1:2 ) {
    tau[j] ~ dgamma( 0.001 , 0.001 )
    mu[j] ~ dnorm(0,0.001)
  }
  p ~ ddirch(alpha)
  alpha[1] <- 1
  alpha[2] <- 1
}
```

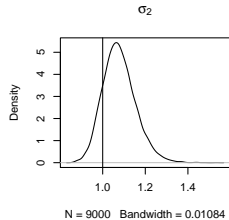
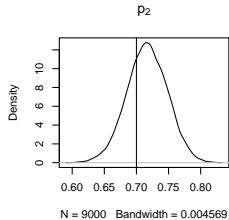
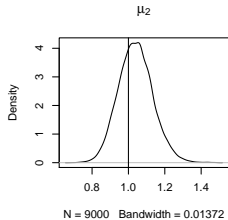
```
library(rjags)
m1 <- jags.model("mixturemodel.jag", data = list(N = length(y), y = y))
res <- coda.samples(m1, var = c("mu", "tau", "p", "z"), n.iter = 10000)
```

Posterior distributions

Summary for component 1:

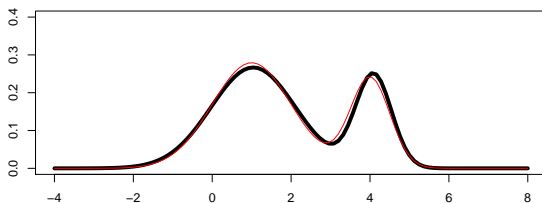


Summary for component 2:



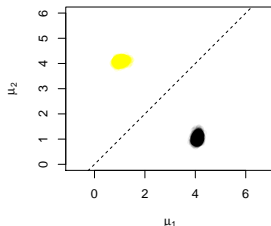
An estimate of the mixture distribution

Comparison of true mixture model (red line) and fitted mixture model (using posterior mean values).



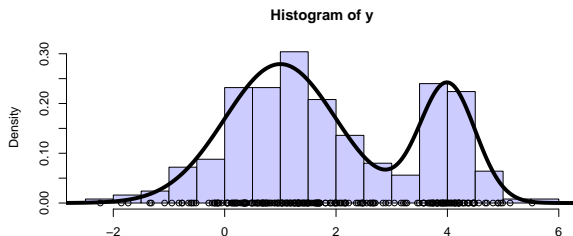
Joint posterior of mean parameters

joint posterior distribution of μ_1 and μ_2 (black) and the expected (due to symmetry) but missing part of the posterior (yellow).

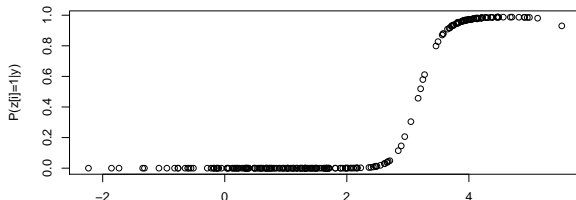


As expected the posterior simulations do not cover both modes.

Posterior probability of $z_i = 1$



Plot of posterior probability that $z_i = 1$ for each y_i :



Notice that this is not simply (a function of) the likelihood ratio.