# Lecture 2. Computational methods Markov Chain Monte Carlo

### Why computational methods?

Recall that in our target formula for posterior  $p(\theta \mid x) = \frac{p(\theta)p(x \mid \theta)}{\int_{\mathbb{R}} p(\theta)p(x \mid \theta) d\theta}$ 

where  $\theta$  are our parameters. The integral below can get really nasty!

**BUT:** this integral is just a constant! Rewrite  $p(\theta|x) = \frac{1}{Z}p(x,\theta)$ , where Z is just a normalising constant, although possibly varying over a large range.

#### What to do?

#### Monte Carlo integration.

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This is called Markov chain Monte Carlo.

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Important: We can construct an MCMC algorithm which will have  $p(\theta \mid X)$  as the stationary distribution!

At each time t the next state  $\theta_{t+1}$  is chosen by first sampling a candidate Y from a **proposal** distribution  $q(. | \theta_t)$  which **depends only on the current state**  $\theta_t$  (or not even that)

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$$\int \pi(\theta_t) P(\theta_{t+1} \mid \theta_t) d\theta_t = \pi(\theta_{t+1}) \quad \text{Meaning: if } \theta_t \text{ is from the distribution } \pi(.), \text{ then } \theta_{t+1} \text{ will be also.}$$

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$$\pi(\theta_t)P(\theta_{t+1}\,|\,\theta_t)d\theta_t=\pi(\theta_{t+1})$$
 **Meaning:** if  $\theta_t$  is from the distribution  $\pi(\,.\,)$ , then  $\theta_{t+1}$  will be also.

Hence, once sample from stationary has been obtained, all subsequent samples are going to be from it. This means MCMC has <u>converged</u>. The period before convergence is called <u>burn-in</u>

### Metropolis-Hastings: how it works in practice

- 1. Start at current position *X*.
- 2. Propose moving to a **new position** Y using proposal q(Y|X)
- 3. Accept/Reject the new position based on the position's adherence to the data and prior distributions using  $\alpha(X,Y)$ 
  - If you accept: Move to the new position Y. Return to Step 1.
  - Else: Do not move to new position, stay at X. Return to Step 1.
- 4. After a large number of iterations, return all accepted positions.

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- 2. Even if the chain converged it may **mix** slowly (move around the states). And hence one needs to **run it for longer** to obtain **reliable estimates**.
- 3. Proposal has to **explore the space efficiently**, sometimes it requires to perform experimentation and craftsmanship to construct a good one.

Jupyter notebook 2

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- 2. The larger the variance of the proposal is the lower the acceptance rate is.
- 3. This can be used during burn-in to reach the desired acceptance rate.

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Acceptance probability will then be  $\alpha(\theta_{-i}, \theta_i, Y_i) = \min\left(1, \frac{\pi(Y_i \mid \theta_{-i})q(\theta_i \mid Y_i, \theta_{-i})}{\pi(\theta_i \mid \theta_{-i})q(Y_i \mid \theta_i, \theta_{-i})}\right)$ 

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Acceptance probability will then be  $\alpha(\theta_{-i}, \theta_i, Y_i) = \min \left( 1, \frac{\pi(Y_i \mid \theta_{-i}) q(\theta_i \mid Y_i, \theta_{-i})}{\pi(\theta_i \mid \theta_{-i}) q(Y_i \mid \theta_i, \theta_{-i})} \right)$ 

Gibbs sampler:  $q(Y_i | \theta_i, \theta_{-i}) = \pi(Y_i | \theta_{-i})$ . Acceptance probability in this case is always equals to 1!

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Gibbs sampling uses the property of tractability of all *conditional* posterior distributions to get samples from the unknown *full* posterior distribution of all model variables.

### Gibbs sampling scheme

Assume we have data  $X \sim p(X | \theta_1, \theta_2)$ 

- 1. Randomly initialize  $\theta_1^{(0)}$  and sample  $\theta_2^{(0)} \sim p(\theta_2 \mid X, \theta_1^{(0)})$
- 2. For step t = 1, ..., T
  - (a) Sample  $\theta_1^{(t)} \sim p(\theta_1 | X, \theta_2^{t-1})$
  - (b) Sample  $\theta_2^{(t)} \sim p(\theta_2 \mid X, \theta_1^t)$