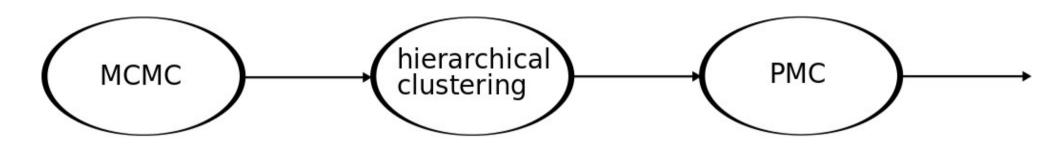
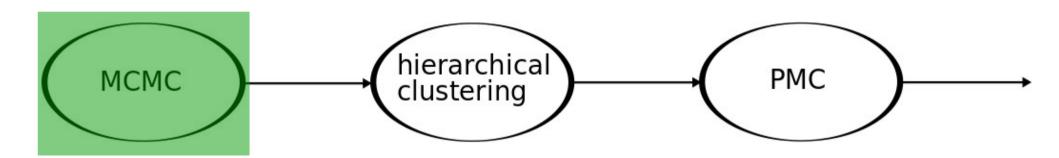
# Monty Python's Stephan's search for the holy grail

- 1. Starting point: Fred's algorithm
- 2. Improvements, where and why
- 3. Hierarchical clustering → Variational Bayes
- 4. PMC → Variational Bayes

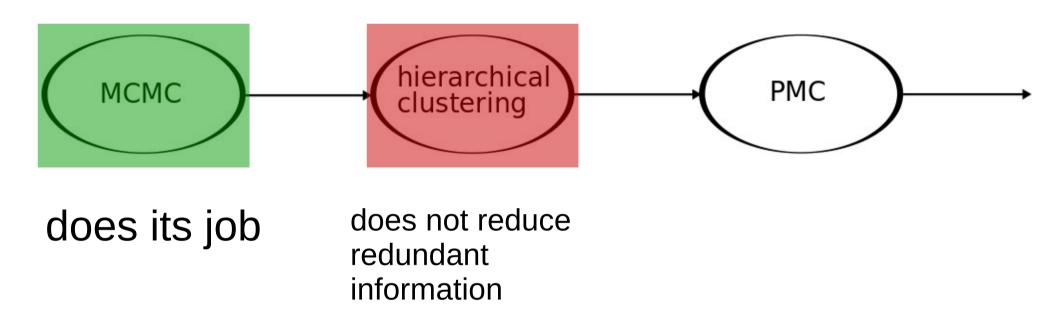


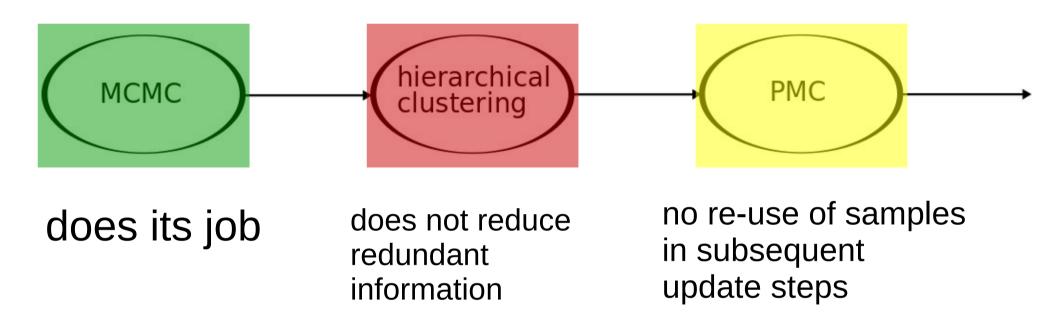
Overall goal: Find "good" proposal for Importance Sampling

$$\int P(x) = \int \frac{P(x)}{q(x)} q(x) \approx \sum_{x_i \in samples} \frac{P(x_i)}{q(x_i)} \text{ where } x_i \sim q(x)$$

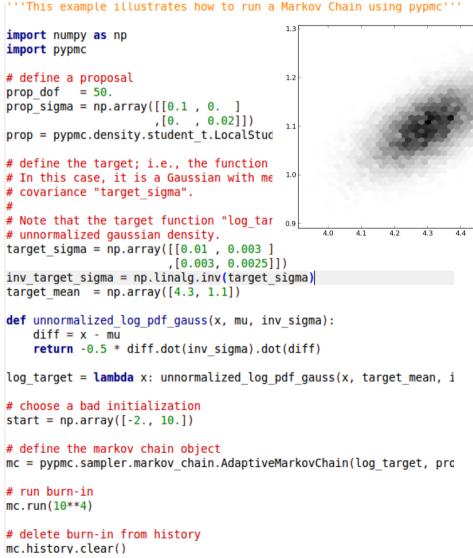


does its job





### https://github.com/fredRos/pypmc



(effective samples).

#### • rel tol-

Relative tolerance  $\epsilon$ . If two consecutive values of the log likelihood bound,  $L_t, L_{t-1}$ , are close, declare convergence. More precisely, check that

$$\left\|rac{L_t-L_{t-1}}{L_t}
ight\|<\epsilon.$$

#### • abs tol-

Absolute tolerance  $\epsilon_a$ . If the current bound  $L_t$  is close to zero, ( $L_t<\epsilon_a$ ), declare convergence if

$$||L_t - L_{t-1}|| < \epsilon_a.$$

verbose - Output status information after each update.

#### set\_variational\_parameters()

Reset the parameters to the submitted values or default.

Use this function to set the prior value (indicated by the subscript  $\theta$  as in  $\alpha_0$ ) or the initial value (e.g.,  $\alpha$ ) used in the iterative procedure to find the posterior value of the variational distribution.

Every parameter can be set in two ways:

- 1. It is specified for only one component, then it is copied to all other components.
- 2. It is specified separately for each component as a K vector.

The prior and posterior variational distributions of  $\mu$  and  $\Lambda$  for each component are given by

$$q(oldsymbol{\mu}, oldsymbol{\Lambda}) = q(oldsymbol{\mu} | oldsymbol{\Lambda}) q(oldsymbol{\Lambda}) = \prod_{k=1}^K \mathcal{N}(oldsymbol{\mu}_k | oldsymbol{m}_k, (eta_k oldsymbol{\Lambda}_k)^{-1}) \mathcal{W}(oldsymbol{\Lambda}_k | oldsymbol{W}_k, 
u_k),$$

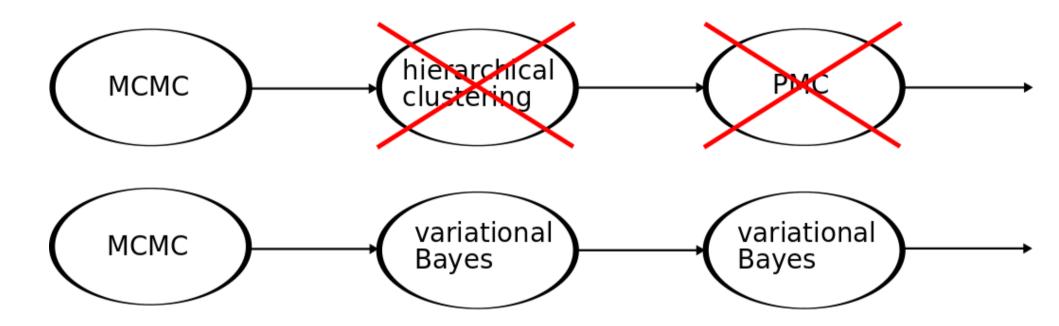
where  ${\cal N}$  denotes a Gaussian and  ${\cal W}$  a Wishart distribution. The weights  ${m \pi}$  follow a Dirichlet distribution

$$q(oldsymbol{\pi}) = Dir(oldsymbol{\pi} | oldsymbol{lpha}).$$

#### Warning

This function may delete results obtained by update().

# Suggested new algorithm



- adapt Gaussian mixture to samples
- Expectation Maximization (EM) with non trivial prior

 Very detailed derivation in "Pattern Recognition and Machine learning" (Christopher M. Bishop)

### Hierarchical clustering (hc) → VB

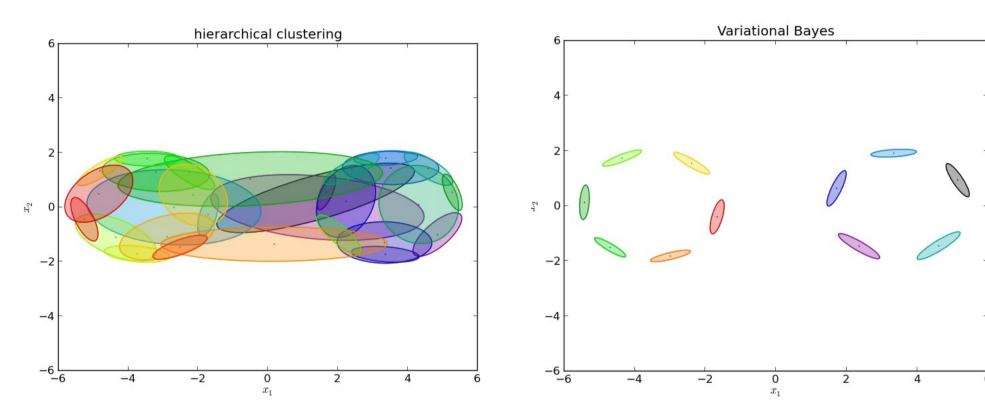
#### Advantages:

- Can at least reproduce the hc results, most runs bring better perplexity and effective sample size
- Automatically removes components with redundant information

### Disadvantage:

Computationally much more expensive

### Example: Gaussian shells, 2 dim

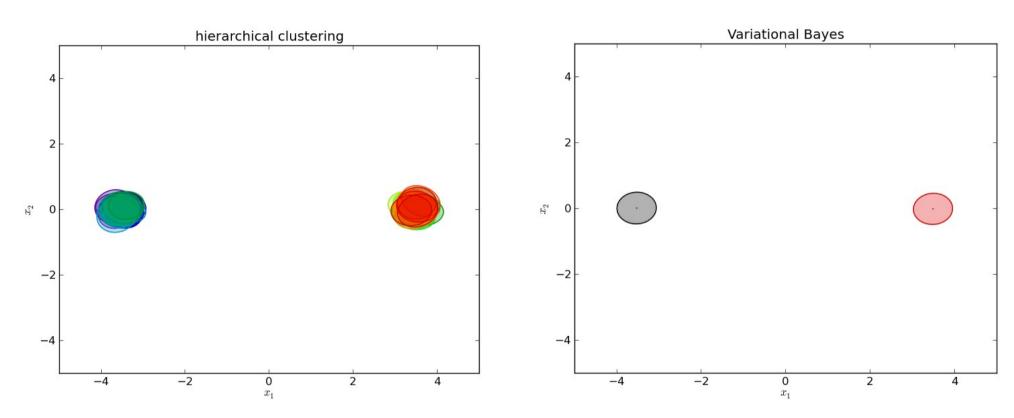


### Start clustering with 15 components per shell

- Converged with 29 components after 21 steps
- Perplexity: ~32%
- ESS: ~27%

- Converged with 11 components after 304 steps
- Perplexity: ~65%
- ESS: ~46%

### Example: Gaussian shells, 20 dim



### Start clustering with 25 components per shell

- Converged with 50 components after 5 steps
- Perplexity: ~33%
- ESS: ~23%

- Converged with 2 components after 304 steps
- Perplexity: ~39%
- ESS: ~28%

### PMC → VB (to come ...)

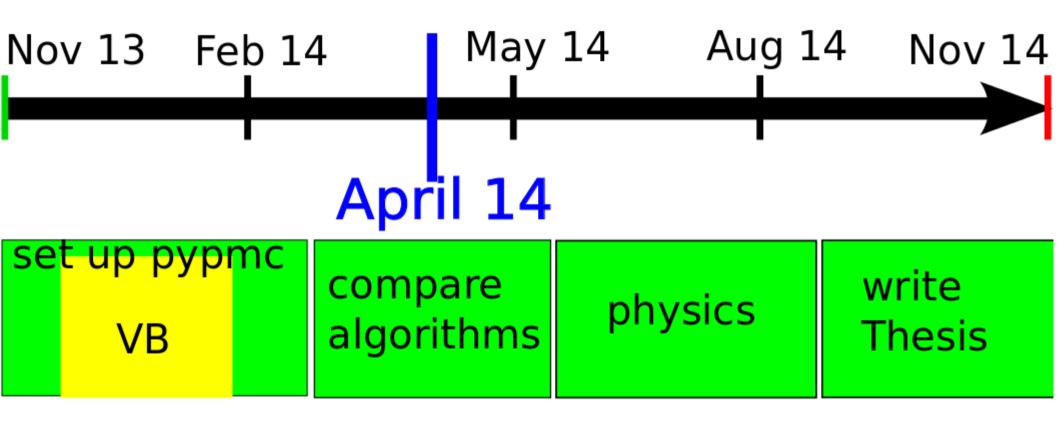
#### Hopes:

- VB allows (forces) you to define priors for the model parameters
  - → We can set the posterior from the previous step as new prior. → Use all information properly.

#### **Fears:**

Unjustifiable computational effort

### **Timeline**



Given data X, we can write for the evidence of any model:

$$\ln p(\mathbf{X}) = \mathcal{L}(q) + KL(q||p)$$

where we define:

$$\mathcal{L}(q) \equiv \int q(\mathbf{Z}) \ln \left\{ \frac{p(\mathbf{X}, \mathbf{Z}, \mathbf{\theta})}{q(\mathbf{Z}, \mathbf{\theta})} \right\} d\mathbf{Z} d\mathbf{\theta}$$

$$KL(q||p) \equiv -\int q(\mathbf{Z}) \ln \left\{ \frac{p(\mathbf{Z}, \boldsymbol{\theta}|\mathbf{X})}{q(\mathbf{Z}, \boldsymbol{\theta})} \right\} d\mathbf{Z} d\boldsymbol{\theta}$$

**\theta** are model parameters

Z are "latent variables" (next slide)

Definition: Latent (hidden) variables

Be  $\mathbf{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$  your data

 $Z = \{z_1, \dots, z_N\}$  is called latent if Z contains

information which can not uniquely be determined from **X** but inferred in a probabilistic model.

Example (Gaussian mixtures):

$$g(x) = \sum_{i} \omega_{i} G_{i}(x|\mu_{i}, \Sigma_{i})$$
$$X \sim g(x)$$

Then the index i is a latent variable

Given data X, we can write for the evidence of any model:

$$\ln p(\mathbf{X}) = \mathcal{L}(q) + KL(q||p)$$

Holds for any probability distribution q

Main idea: Restrict q to factorize as below and minimize the Kullback-Leibler divergence

$$q(\mathbf{Z},\mathbf{\theta}) = q(\mathbf{Z})q(\mathbf{\theta})$$

I.e. we look for the closest factorizing q to p in the sense of KL(q||p)

Given data X, we can write for the evidence of any model:

$$\ln p(\mathbf{X}) = \mathcal{L}(q) + KL(q||p)$$

Minimizing KL is equivalent to maximizing  $\mathcal{L}$ .

Maximizing  $\mathcal{L}$ , in the case of a Gaussian mixture model, leads to an EM-like algorithm.