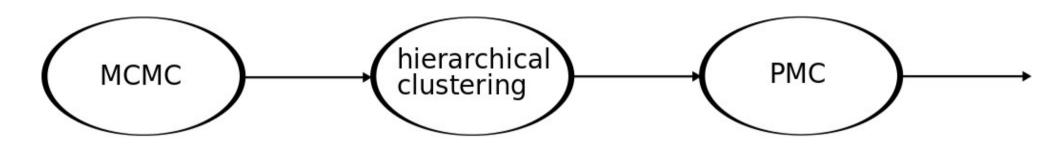
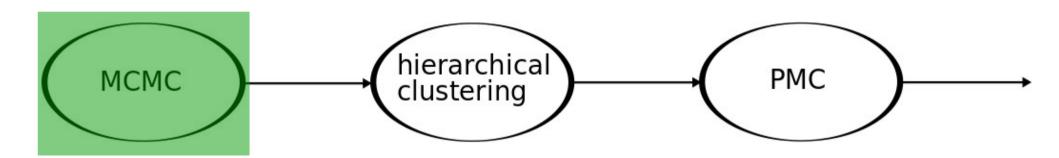
Monty Python's Stephan's search for the holy grail

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

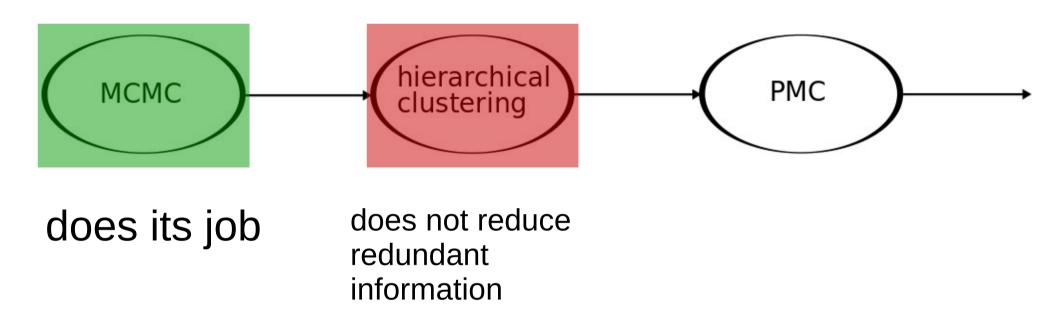


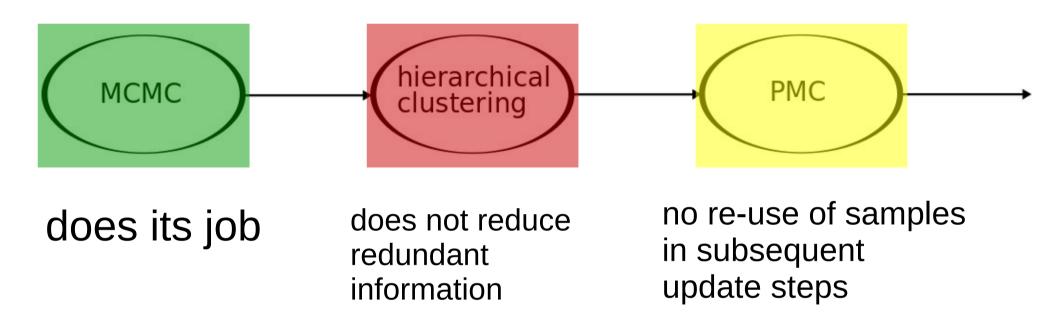
Overall goal: Find "good" proposal for Importance Sampling

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

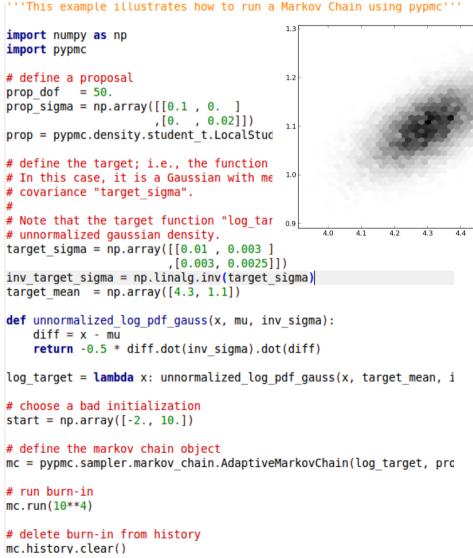


does its job





https://github.com/fredRos/pypmc



(effective samples).

• rel tol-

Relative tolerance ϵ . 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 ϵ_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 θ as in α_0) or the initial value (e.g., α) 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 μ and Λ for each component are given by

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

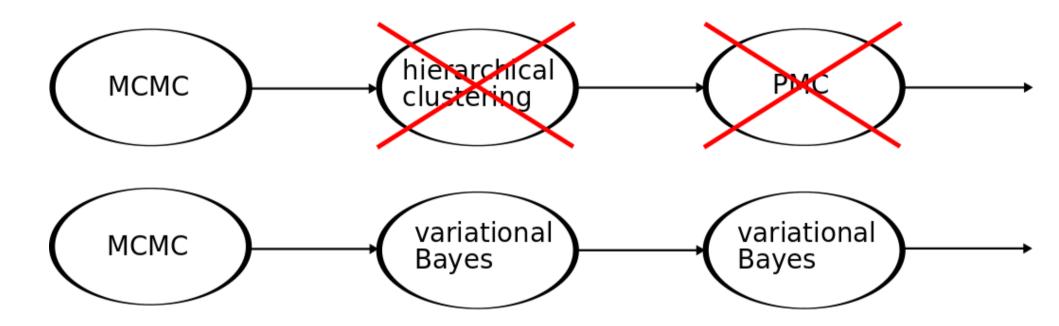
where ${\cal N}$ denotes a Gaussian and ${\cal W}$ a Wishart distribution. The weights ${m \pi}$ follow a Dirichlet distribution

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

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

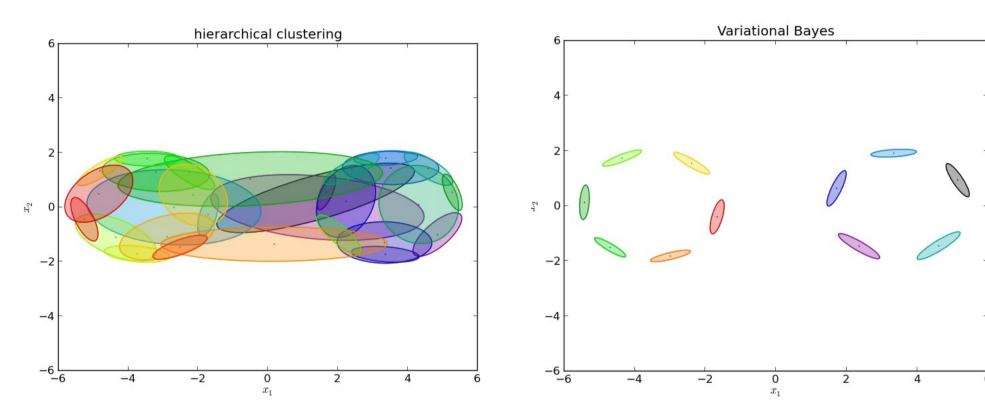
Advantages:

- 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

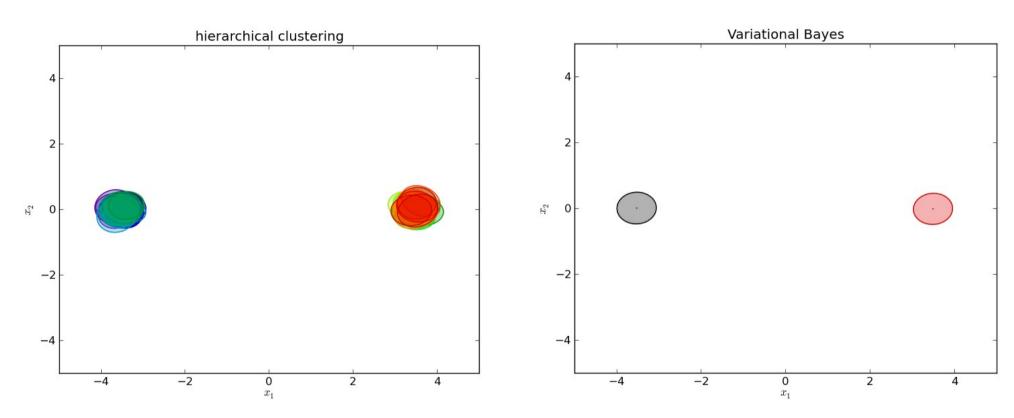


Start clustering with 15 components per shell

- 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 ...)

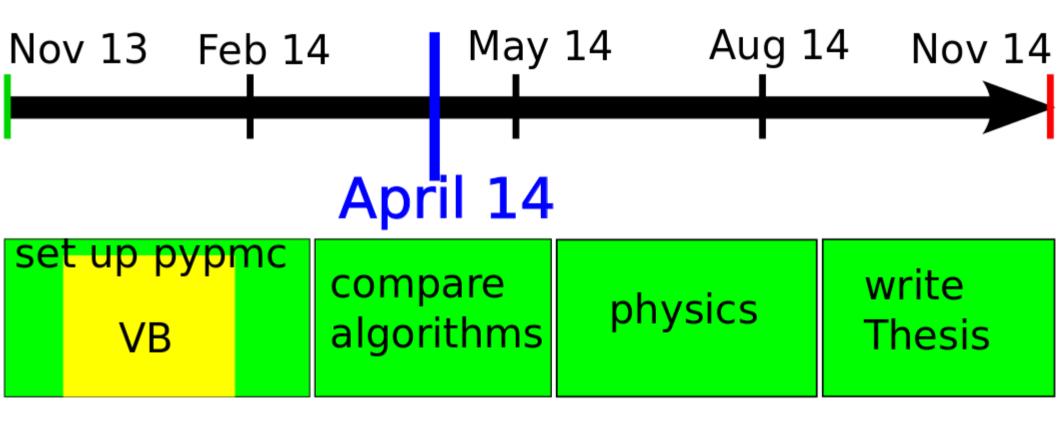
Hopes:

- 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:

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

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

\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):

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

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.