

# Bayesian Theory and Computation

## Lecture 18: Dirichlet Processes

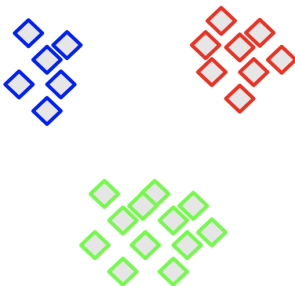


**Cheng Zhang**

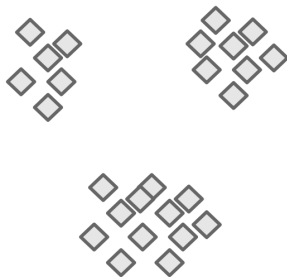
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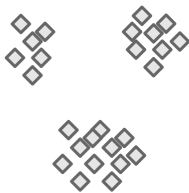
How to choose the number of clusters?



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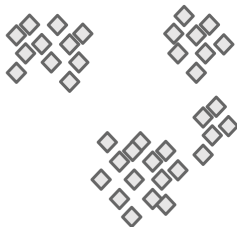


How to choose the number of clusters?



**T=1**

**Streaming Data**



**T=2**



- ▶ A generative approach to clustering
  - ▶ pick one of  $K$  clusters from a distribution  $\pi = (\pi_1, \dots, \pi_K)$
  - ▶ generate a data point from a cluster-specific probability distribution
- ▶ This yields a finite mixture model:

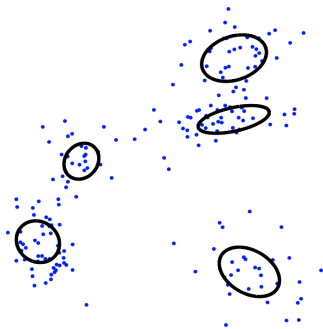
$$p(x|\phi, \pi) = \sum_{k=1}^K \pi_k p(x|\phi_k)$$

where  $\pi$  and  $\phi = (\phi_1, \dots, \phi_K)$  are the parameters, and here we assume the same parameterized family for each cluster for simplicity.

- ▶ Data  $\{x_i\}_{i=1}^N$  are assumed to be generated conditionally iid from this mixture model.



- ▶ For Gaussian mixtures,  $\phi_k = (\mu_k, \Sigma_k)$  and  $p(x|\phi_k)$  is a Gaussian density with mean  $\mu_k$  and covariance matrix  $\Sigma_k$



- ▶ Mixture models make the assumption that each data point arises from a single mixture component, i.e., the  $k$ th cluster is by definition the set of data points arising from the  $k$ th mixture component.
- ▶ Can capture this explicitly via a latent multinomial variable  $Z$ :

$$\begin{aligned} p(x|\phi, \pi) &= \sum_{k=1}^K p(Z = k|\pi)p(x|Z = k, \phi) \\ &= \sum_{k=1}^K \pi_k p(x|\phi_k) \end{aligned}$$



- ▶ Another way to express this: define an underlying measure

$$G = \sum_{k=1}^K \pi_k \delta_{\phi_k}$$

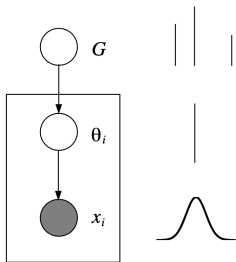
where  $\delta_{\phi_k}$  is an *atom* (Dirac delta function) at  $\phi_k$ .

- ▶ Now we can redefine the process of obtaining a sampling from a finite mixture model as follows. For  $i = 1, \dots, n$ :

$$\begin{aligned}\theta_i &\sim G \\ x_i &\sim p(\cdot | \theta_i)\end{aligned}$$

- ▶ Note that each  $\theta_i$  is equal to one of the underlying  $\phi_k$ . Indeed, the subset of  $\{\theta_i\}$  that maps to  $\phi_k$  is exactly the  $k$ th cluster





$$G = \sum_{k=1}^K \pi_k \delta_{\phi_k}$$

$$\theta_i \sim G$$

$$x_i \sim p(\cdot | \theta_i)$$

Adapted from M. I. Jordan



- ▶ Bayesian approaches allow us to integrate out model parameters
- ▶ Need to place priors on the parameters  $\phi$  and  $\pi$
- ▶ The choice of prior for  $\phi$  is model-specific; e.g., we may use conjugate normal/inverse-gamma priors for a Gaussian mixture model. Let us denote this prior as  $G_0$ .
- ▶ What to choose for the mixture weights  $\pi$ ? A common choice is a symmetric Dirichlet prior,  $\text{Dir}(\alpha_0/K, \dots, \alpha_0/K)$ 
  - ▶ the symmetry accords with the common assumption of the order-free of the labels of the mixture components
  - ▶ the concentration parameter  $\alpha_0$  controls concentration level of the labels

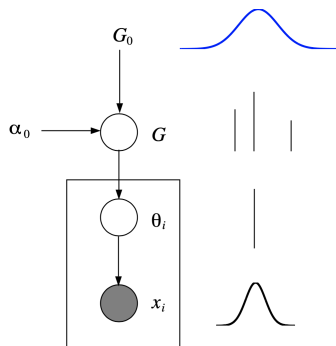
$$\phi_k \sim G_0$$

$$\pi_k \sim \text{Dir}(\alpha_0/K, \dots, \alpha_0/K)$$

$$G = \sum_{k=1}^K \pi_k \delta_{\phi_k}$$

$$\theta_i \sim G$$

$$x_i \sim p(\cdot | \theta_i)$$



- Note that  $G$  is now a random measure



- ▶ Posterior distributions can't be found analytically; nor can predictive distributions (for future observations)
- ▶ However, a variety of MCMC sampling algorithms are available
- ▶ Use the indicators  $Z$  within a Gibbs sampler. Given  $Z$ , we know which data points belong to which cluster, so:
  - ▶  $p(\pi|Z, \phi)$ : standard multinomial-Dirichlet conjugacy
  - ▶  $p(\phi|Z, \pi)$ : separate updates for each cluster; i.e., for each  $\phi_k$  (and conjugacy of  $G_0$  and  $p(\cdot|\phi)$  can make this easy)
  - ▶  $p(Z|\pi, \phi)$ : multinomial classification
- ▶ We can also use variational inference.

- ▶ How to choose  $K$ , the number of mixture components?
- ▶ Various generic model selection methods can be considered: e.g., cross-validation, bootstrap, AIC, BIC, DIC, Laplace, bridge sampling, etc
- ▶ Or we can place a parametric prior on  $K$  (e.g., Poisson) and use Bayesian methods
- ▶ The Dirichlet process provides a nonparametric Bayesian alternative.

- ▶ Make sure we always have more clusters than we need.
- ▶ How about infinite clusters a priori?

$$p(x|\phi, \pi) = \sum_{k=1}^{\infty} \pi_k p(x|\phi_k)$$

- ▶ A finite data set will always use a finite, but random, number of clusters.
- ▶ How to choose the prior?
- ▶ We need something like a Dirichlet prior, but with an infinite number of components.

- ▶ Relation to gamma distribution: If  $\eta_k \sim \text{Gamma}(\alpha_k, \beta)$  independently, then

$$S = \sum_k \eta_k \sim \text{Gamma} \left( \sum_k \alpha_k, \beta \right)$$

and

$$V = (v_1, \dots, v_k) = (\eta_1/S, \dots, \eta_k/S) \sim \text{Dir}(\alpha_1, \dots, \alpha_K)$$

- ▶ Therefore, if  $(\pi_1, \dots, \pi_K) \sim \text{Dir}(\alpha_1, \dots, \alpha_K)$  then

$$(\pi_1 + \pi_2, \pi_3, \dots, \pi_K) \sim \text{Dir}(\alpha_1 + \alpha_2, \alpha_3, \dots, \alpha_K)$$

This is known as the **collapsing** property.

- ▶ The beta distribution is a Dirichlet distribution on the 1-simplex
- ▶ Let  $(\pi_1, \dots, \pi_K) \sim \text{Dir}(\alpha_1, \dots, \alpha_K)$  and  $\theta \sim \text{Beta}(\alpha_1 b, \alpha_1(1-b))$ ,  $0 < b < 1$ .
- ▶ Then

$$(\pi_1 \theta, \pi_1(1-\theta), \pi_2, \dots, \pi_K) \sim \text{Dir}(\alpha_1 b_1, \alpha_1(1-b_1), \alpha_2, \dots, \alpha_K)$$

- ▶ More generally, if  $\theta \sim \text{Dir}(\alpha_1 b_1, \alpha_1 b_2, \dots, \alpha_1 b_N)$ ,  $\sum_i b_i = 1$ , then

$$(\pi_1 \theta_1, \dots, \pi_1 \theta_N, \pi_2, \dots, \pi_K) \sim \text{Dir}(\alpha_1 b_1, \dots, \alpha_1 b_N, \alpha_2, \dots, \alpha_K)$$

This is known as the **splitting** property.



- ▶ Renormalization. If  $(\pi_1, \dots, \pi_K) \sim \text{Dir}(\alpha_1, \dots, \alpha_K)$ , and

$$V = (V_2, V_3, \dots, V_K), \quad V_k = \frac{\pi_k}{\sum_{k \geq 2} \pi_k}$$

- ▶ What is the distribution of  $V$ ?

$$V \sim \text{Dir}(\alpha_2, \dots, \alpha_K)$$

- ▶ All these properties can be easily verified using the aforementioned gamma distribution representation.

- ▶ Let  $G_0$  be a distribution on some space  $\Omega$ , e.g. a Gaussian distribution on the real line.
- ▶ Assume that  $\pi, \phi$  have the following distributions

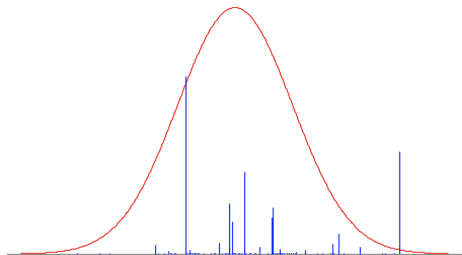
$$\phi_k \sim G_0$$

$$\pi \sim \lim_{K \rightarrow \infty} \text{Dir} \left( \frac{\alpha_0}{K}, \dots, \frac{\alpha_0}{K} \right)$$

- ▶ Then  $G := \sum_{k=1}^{\infty} \pi_k \delta_{\phi_k}$  defines an infinite distribution over  $G_0$ .
- ▶ We say (informally) that  $G$  follows a **Dirichlet Process**

$$G \sim \text{DP}(\alpha_0 G_0)$$

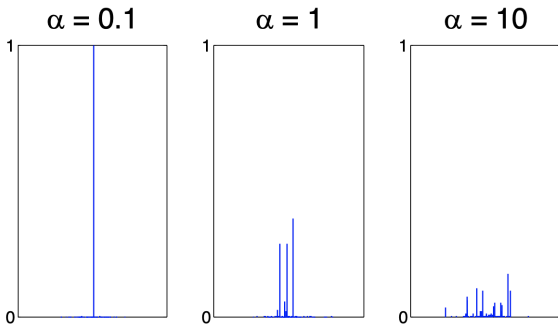
- ▶ Samples from the Dirichlet process are *discrete*.
- ▶ We call the point masses in the resulting distribution, *atoms*.



- ▶ The *base measure*  $G_0$  determines the *locations* of the atoms.



- ▶ The *concentration parameter*  $\alpha_0$  determines the distribution over atom sizes.
- ▶ Small values of  $\alpha_0$  gives *sparse* distributions.



- ▶ Let  $(\Omega, \mathcal{B})$  be a measurable space, with  $G_0$  a probability measure on the space, and let  $\alpha_0$  be a positive real number.
- ▶ A **Dirichlet process** is the distribution of a random probability measure  $G$  over  $(\Omega, \mathcal{B})$  such that, for any finite partition  $(A_1, \dots, A_r)$  of  $\Omega$ , the random vector  $(G(A_1), \dots, G(A_r))$  follows a finite-dimensional Dirichlet distribution:

$$(G(A_1), \dots, G(A_r)) \sim \text{Dir}(\alpha_0 G_0(A_1), \dots, \alpha_0 G_0(A_r))$$

- ▶ We write  $G \sim \text{DP}(\alpha_0 G_0)$ , and call  $G_0$  the **base measure**,  $\alpha_0$  the **concentration parameter**.

- ▶ Let  $A_1, \dots, A_K$  be a partition of  $\Omega$ . Let  $G(A_k)$  be the mass assigned by  $G \sim \text{DP}(\alpha_0 G_0)$  to partition  $A_k$ . Then

$$(G(A_1), \dots, G(A_K)) \sim \text{Dir}(\alpha_0 G_0(A_1), \dots, \alpha_0 G_0(A_K))$$

- ▶ If we see an observation in the  $j$ -th segment, then

$$\begin{aligned} &(G(A_1), \dots, G(A_K) | \theta_1 \in A_j) \\ &\sim \text{Dir}(\alpha_0 G_0(A_1), \dots, \alpha_0 G_0(A_j) + 1, \dots, \alpha_0 G_0(A_K)). \end{aligned}$$

- ▶ This is true for all possible partitions of  $\Omega$ .
- ▶ Therefore, the posterior distribution of  $G$ , given an observation  $\phi$ , is given by

$$G | \theta_1 = \phi \sim \text{DP}(\alpha_0 G_0 + \delta_\phi)$$

- ▶ The Dirichlet process clusters observations.
- ▶ A new data point can either join an existing cluster, or start a new cluster.
- ▶ Question: What is the predictive distribution for a new data point?
- ▶ Assume  $G_0$  is a continuous distribution on  $\Omega$ . This means for every point  $\phi$  in  $\Omega$ ,  $G_0(\phi) = 0$ .
- ▶ First data point:
  - ▶ Start a new cluster
  - ▶ Sample a parameter  $\phi_1 \sim G_0$  for that cluster.



- ▶ We have now split our parameter space in two: the singleton  $\phi_1$ , and everything else.
- ▶ Let  $\pi_1$  be the size of atom at  $\phi_1$ .
- ▶ The combined mass of all the other atoms is  $\pi_* = 1 - \pi_1$ .
- ▶ According to the DP,

$$(\pi_1, \pi_*) \sim \text{Dir}(0, \alpha_0)$$

- ▶ Given  $\theta_1 = \phi_1$ , the posterior is

$$(\pi_1, \pi_*) | \theta_1 = \phi_1 \sim \text{Dir}(1, \alpha_0)$$





- If we integrate out  $\pi_1$ , we get

$$\begin{aligned} p(\theta_2 = \phi_k | \theta_1 = \phi_1) &= \int p(\theta_2 = \phi_k | (\pi_1, \pi_*)) p((\pi_1, \pi_*) | \theta_1 = \phi_1) d\pi_1 \\ &= \int \pi_k \text{Dir}((\pi_1, 1 - \pi_1) | 1, \alpha_0) d\pi_1 \\ &= \mathbb{E}_{\text{Dir}(1, \alpha_0)} \pi_k \\ &= \begin{cases} \frac{1}{1 + \alpha_0} & \text{if } k = 1 \\ \frac{\alpha_0}{1 + \alpha_0} & \text{for new } k. \end{cases} \end{aligned}$$



- ▶ Lets say we choose to start a new cluster, and sample a new parameter  $\phi_2 \sim G_0$ . Let  $\pi_2$  be the size of the atom at  $\phi_2$ .
- ▶ Similarly, the posterior is

$$(\pi_1, \pi_2, \pi_*) | \theta_1 = \phi_1, \theta_2 = \phi_2 \sim \text{Dir}(1, 1, \alpha_0)$$

- ▶ If we integrate out  $\pi = (\pi_1, \pi_2, \pi_*)$ , we get

$$\begin{aligned} p(\theta_3 = \phi_k | \theta_1 = \phi_1, \theta_2 = \phi_2) &= \int p(\theta_3 = \phi_k | \pi) p(\pi | \theta_1 = \phi_1, \theta_2 = \phi_2) d\pi \\ &= \mathbb{E}_{\text{Dir}(1,1,\alpha_0)} \pi_k \\ &= \begin{cases} \frac{1}{2+\alpha_0} & \text{if } k = 1 \\ \frac{1}{2+\alpha_0} & \text{if } k = 2 \\ \frac{\alpha_0}{2+\alpha_0} & \text{for new } k. \end{cases} \end{aligned}$$

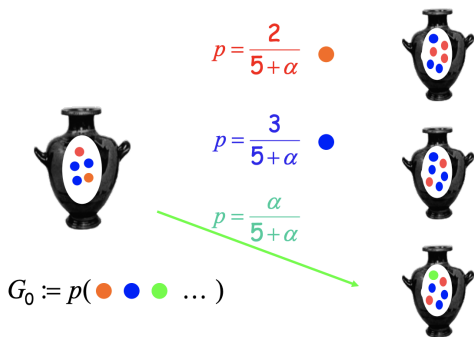


- ▶ In general, if  $m_k$  is the number of times we have seen  $X_i = k$ , and  $K$  is the total number of observed values,

$$\begin{aligned} p(\theta_{n+1} = \phi_k | \theta_1, \dots, \theta_n) &= \int p(\theta_{n+1} = \phi_k | \pi) p(\pi | \theta_1, \dots, \theta_n) d\pi \\ &= \mathbb{E}_{\text{Dir}(m_1, \dots, m_K, \alpha_0)} \pi_k \\ &= \begin{cases} \frac{m_k}{n + \alpha_0} & \text{if } k \leq K \\ \frac{\alpha_0}{n + \alpha_0} & \text{for new cluster.} \end{cases} \end{aligned}$$

- ▶ We tend to see observations that we have seen before, i.e., **rich-get-richer property**
- ▶ We can always add new features, a typical nonparametric behavior.





Adapted from Eric Xing

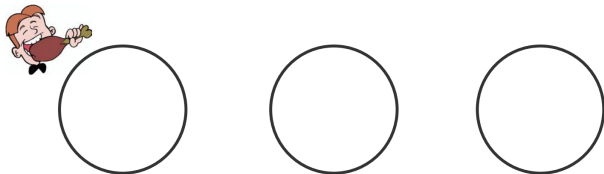
- ▶ Joint:  $G(\text{urn}) \sim \text{DP}(\alpha_0 G_0)$
- ▶ Marginal:  $\theta_{n+1} | \theta_{\leq n}, \alpha_0, G_0 \sim \sum_{k=1}^K \frac{m_k}{n+\alpha_0} \delta_{\phi_k} + \frac{\alpha_0}{n+\alpha_0} G_0$ .



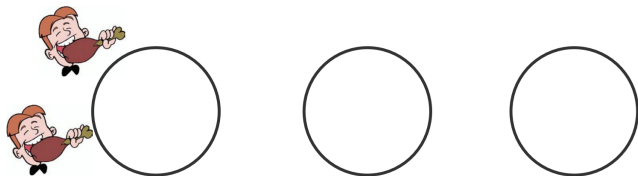
- ▶ The resulting distribution over data points can be thought of using the following urn scheme (Blackwell and MacQueen, 1973).
- ▶ An urn initially contains a black ball of mass  $\alpha_0$ .
- ▶ For  $n = 1, 2, \dots$ , sample a ball from the urn with probability proportional to its mass.
- ▶ If the ball is black, choose a previously unseen color, record that color, and return the black ball plus a unit-mass ball of the new color to the urn.
- ▶ If the ball is not black, record its color and return it, plus another unit-mass ball of the same color, to the urn.

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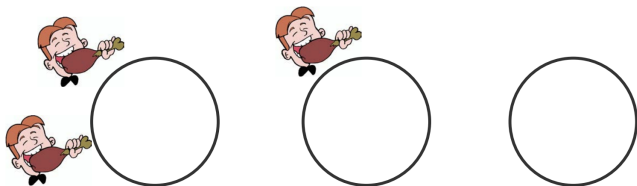


- ▶ The distribution over partitions can also be described in terms of the following restaurant metaphor:
- ▶ The first customer enters a restaurant, and picks a table.
- ▶ The  $n$ -th customer enters the restaurant. He sits at an existing table with probability  $\frac{m_k}{n-1+\alpha_0}$ , where  $m_k$  is the number of people sat the table  $k$ . He starts a new table with probability  $\frac{\alpha_0}{n-1+\alpha_0}$ .

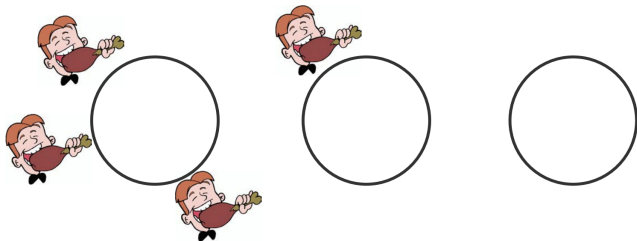




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- ▶ An interesting fact: the distribution over the clustering of the first  $N$  customers **does not depend on the order in which they arrived**.
- ▶ However, the customers are not independent. They tend to sit at popular tables.
- ▶ We say that distributions like this are *exchangeable*.

$$p(\theta_1, \dots, \theta_N) = p(\theta_{\sigma(1)}, \dots, \theta_{\sigma(n)})$$

- ▶ By **de Finetti's** theorem, there exists a random distribution  $G$  and a prior  $P(G)$  such that

$$p(\theta_1, \dots, \theta_N) = \int \prod_{i=1}^N G(\theta_i) dP(G)$$

- ▶ In our setting, the prior  $p(G)$  is just  $\text{DP}(\alpha_0 G_0)$ , thus establishing existence.



- ▶ Define an infinite sequence of Beta random variables:

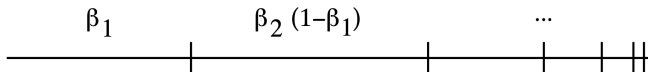
$$\beta_k \sim \text{Beta}(1, \alpha_0), \quad k = 1, 2, \dots$$

- ▶ Now define an infinite sequence of mixing proportions as:

$$\pi_1 = \beta_1$$

$$\pi_k = \beta_k \prod_{\ell=1}^{k-1} (1 - \beta_\ell), \quad k = 2, 3, \dots$$

- ▶ This can be viewed as breaking off portions of a stick:



- ▶ We now have an explicit formula for each  $\pi_k$ :

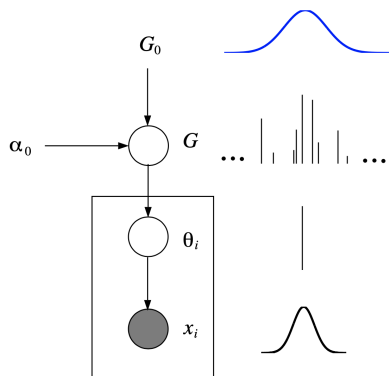
$$\pi_k = \beta_k \prod_{\ell=1}^{k-1} (1 - \beta_\ell)$$

- ▶ We can easily see that  $\sum_{k=1}^{\infty} \pi_k = 1$ :

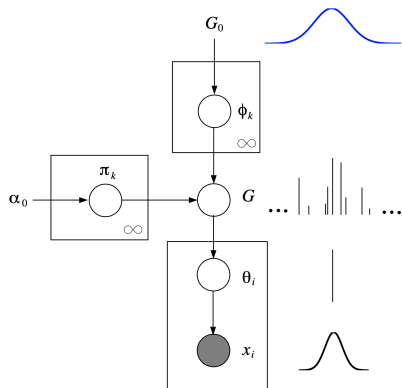
$$\begin{aligned} 1 - \sum_{k=1}^{\infty} \pi_k &= 1 - \beta_1 - \beta_2(1 - \beta_1) - \beta_3(1 - \beta_1)(1 - \beta_2) - \dots \\ &= (1 - \beta_1)(1 - \beta_2 - \beta_3(1 - \beta_2) - \dots) \\ &= \prod_{k=1}^{\infty} (1 - \beta_k) = 0 \end{aligned}$$

- ▶ Let  $\phi_k \sim G_0, \forall k$ ,  $G = \sum_{k=1}^{\infty} \pi_k \delta_{\phi_k}$  has a clean definition as a random measure. In fact,

$$G \sim \text{DP}(\alpha_0 G_0).$$



Polya urn construction



Stick breaking construction



- ▶ Now we can use a Dirichlet process as the prior for an unknown mixture distribution (with potentially infinite mixture components).
- ▶ Suppose we have  $x_1, \dots, x_n$  observations from some unknown distribution.
- ▶ We can model the unknown distribution of  $x$  as a mixture of simple distributions of the form  $f(\cdot|\theta)$ .
- ▶ We denote the mixing distribution over  $\theta$  as  $G$  and let the prior over  $G$  be a Dirichlet process

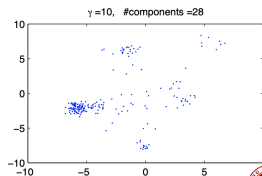
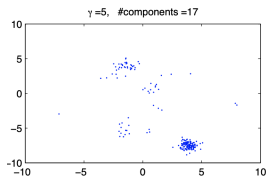
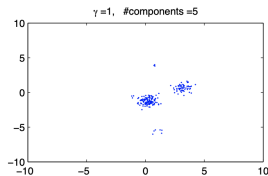
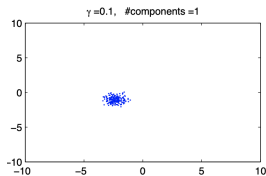
$$x_i|\theta_i \sim f(\theta_i)$$

$$\theta_i|G \sim G$$

$$G \sim \text{DP}(\alpha_0 G_0)$$



- ▶ Multiple subjects can be mapped to the same  $\phi$ . This creates a clustering of subjects.
- ▶ The following graphs shows 4 different data sets ( $n = 200$ ) randomly generated from distributions sampled from Dirichlet process mixture priors with  $\alpha_0$ .





- ▶ We can integrate out  $G$  to get the CRP. Note that the CRP is exchangeable, which induces the conditional priors

$$p(\theta_i | \theta_{-i}, \alpha_0, G_0) = \frac{\alpha_0}{n - 1 + \alpha_0} G_0(\theta_i) + \sum_{k=1}^{K^{(-i)}} \frac{m_k^{(-i)}}{n - 1 + \alpha_0} \delta_{\phi_k^{(-i)}}$$

- ▶ Let  $z_i$  be the cluster allocation of the  $i$ -th data point. The collapsed Gibbs sampler alternates between

- ▶ update  $z_i$

$$p(z_i = k | x_i, z_{-i}, \phi_{1:K}) \propto \begin{cases} m_k^{(-i)} f(x_i | \phi_k^{(-i)}) & k \leq K^{(-i)} \\ \alpha_0 \int f(x_i | \theta) dG_0(\theta) & k = K^{(-i)} + 1 \end{cases}$$

- ▶ update  $\phi_k$

$$p(\phi_k | z_{1:n}, x_{1:n}) \propto G_0(\phi_k) \prod_{i: z_i = k} f(x_i | \phi_k)$$

- ▶ If  $G_0$  is conjugate to  $f$ , the above steps can be evaluated accurately.

- ▶ For the concentration parameter  $\alpha_0$ , we have

$$p(K|\alpha_0) \propto \alpha_0^K \frac{\Gamma(\alpha_0)}{\Gamma(\alpha_0 + n)}$$

where  $K$  is the number of unique  $\phi$ 's (e.g., the number of clusters).

- ▶ Therefore, given  $K$  and the prior distribution  $P(\alpha_0)$  we can sample from the posterior distribution of  $\alpha_0$  using the MH algorithm or the Gibbs sampling method of Escobar and West (1995).

- ▶ We are only updating one data point at a time.
- ▶ Imagine two “true” clusters are merged into a single cluster, a single data point is unlikely to “break away”.
- ▶ Getting to the true distribution involves going through low probability states, i.e., mixing can be slow.
- ▶ If the likelihood is not conjugate, integrating out parameter values for new features can be difficult.
- ▶ Neal (2000) offers a variety of algorithms.

- ▶ The stick-breaking representation orders the mixture components so that the weights are stochastically decreasing. For a sufficiently large  $T$ , we will have  $\sum_{k>T} \pi_k \approx 0$ .
- ▶ Therefore, we can truncate the stick-breaking construction at a fixed value  $T$  and let  $\beta_T = 1$ .
- ▶ This implies  $\pi_k = 0, \forall k > T$ , and the distribution of

$$G_T = \sum_{k=1}^T \pi_k \delta_{\phi_k}$$

is known as a **truncated Dirichlet process**.

- ▶ Variational distance between distributions of marginals from a DP and from its truncation at  $T$  is approximately  $4n \exp(-(T-1)/\alpha_0)$ .  $T$  doesn't have to be very large to get a good approximation.



- ▶ State of the Markov chain:  $(\beta_{1:T-1}, \phi_{1:T}, z_{1:n})$ .
- ▶ Update  $z_i$  by multinomial sampling with

$$p(z_i = k | \beta, \phi, x_i) \propto \pi_k f(x_i | \phi_k)$$

- ▶ Update  $\beta_k$  by sampling from the conditional posterior

$$\beta_k \sim \text{Beta}(1 + m_k, \alpha_0 + \sum_{j>k} m_j)$$

- ▶ Update  $\phi_k$  by sampling from the conditional posterior

$$p(\phi_k | z_{1:n}, x_{1:n}) \propto G_0(\phi_k) \prod_{i:z_i=k} f(x_i | \phi_k)$$

- ▶ One can monitor  $\max_i z_i$  to verify that truncation at  $T$  is good enough, and increase  $T$  if necessary.

- ▶ We can also use truncated steak-breaking representation to form a mean field approximation of DP mixtures

$$q(\beta, \phi, z) = \prod_{i=1}^n q(z_i | w_i) \prod_{k=1}^T q(\phi_k | \tau_k) \prod_{k=1}^{T-1} q(\beta_k | \gamma_k)$$

- ▶ For a conjugate DP mixture in the exponential family

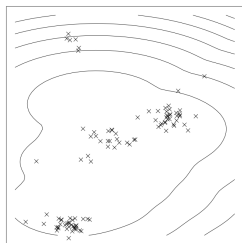
$$\begin{aligned} \gamma_{k,1} &= 1 + \sum_{i=1}^n w_{i,k}, & \gamma_{k,2} &= \alpha_0 + \sum_{i=1}^n \sum_{j>k} w_{i,j} \\ \tau_{k,1} &= \lambda_1 + \sum_{i=1}^n w_{i,k} t(x_i), & \tau_{k,2} &= \lambda_2 + \sum_{i=1}^n w_{i,k} \\ w_{i,k} &\propto \exp(S_k) \end{aligned}$$

where

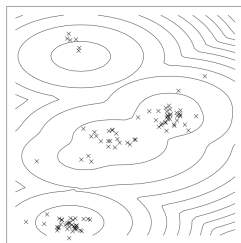
$$S_k = \mathbb{E} \log \beta_k + \sum_{j<k} \mathbb{E} \log(1 - \beta_j) + \mathbb{E} \phi_k^T t(x_i) - \mathbb{E} A(\phi_k)$$



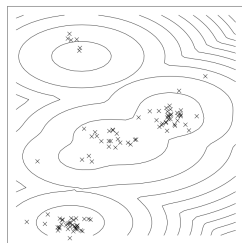
- ▶ The approximate predictive distribution given by variational inference at different stages of the algorithm. The data are 100 points generated by a Gaussian DP mixture model with fixed diagonal covariance.



Initial state



1st iteration



5th (and last) iteration



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