

Cluster and Feature Modeling from Combinatorial Stochastic Processes

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Abstract. One of the focal points of the modern literature on Bayesian nonparametrics has been the problem of *clustering*, or *partitioning*, where each data point is modeled as being associated with one and only one of some collection of groups called clusters or partition blocks. Underlying these Bayesian nonparametric models are a set of interrelated stochastic processes, most notably the Dirichlet process and the Chinese restaurant process. In this paper we provide a formal development of an analogous problem, called *feature modeling*, for associating data points with arbitrary nonnegative integer numbers of groups, now called features or topics. We review the existing combinatorial stochastic process representations for the clustering problem and develop analogous representations for the feature modeling problem. These representations include the beta process and the Indian buffet process as well as new representations that provide insight into the connections between these processes. We thereby bring the same level of completeness to the treatment of Bayesian nonparametric feature modeling that has previously been achieved for Bayesian nonparametric clustering.

Key words and phrases: Cluster, feature, Dirichlet process, beta process, Chinese restaurant process, Indian buffet process, nonparametric, Bayesian, combinatorial stochastic process.

1. INTRODUCTION

Bayesian nonparametrics is the area of Bayesian analysis in which the finite-dimensional prior distributions of classical Bayesian analysis are replaced with stochastic processes. While the rationale for allowing infinite collections of random variables into Bayesian inference is often taken to be that of diminishing the role of prior assumptions, it is also possible to view the move to nonparametrics as supplying the Bayesian paradigm with a richer collection of distributions with which to express prior belief, thus in some sense em-

phasizing the role of the prior. In practice, however, the field has been dominated by two stochastic processes—the Gaussian process and the Dirichlet process—and thus the flexibility promised by the nonparametric approach has arguably not yet been delivered. In the current paper we aim to provide a broader perspective on the kinds of stochastic processes that can provide a useful toolbox for Bayesian nonparametric analysis. Specifically, we focus on *combinatorial stochastic processes* as embodying mathematical structure that is useful for both model specification and inference.

The phrase “combinatorial stochastic process” comes from probability theory (Pitman, 2006), where it refers to connections between stochastic processes and the mathematical field of combinatorics. Indeed, the focus in this area of probability theory is on random versions of classical combinatorial objects such as partitions, trees and graphs—and on the role of combinatorial analysis in establishing properties of these processes. As we wish to argue, this connection is also fruitful in a statistical setting. Roughly speaking, in statistics it is often natural to model observed data as

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arising from a combination of underlying factors. In the Bayesian setting, such models are often embodied as latent variable models in which the latent variable has a compositional structure. Making explicit use of ideas from combinatorics in latent variable modeling cannot only suggest new modeling ideas but can also provide essential help with calculations of marginal and conditional probability distributions.

The Dirichlet process already serves as one interesting exhibit of the connections between Bayesian non-parametrics and combinatorial stochastic processes. On the one hand, the Dirichlet process is classically defined in terms of a partition of a probability space (Ferguson, 1973), and there are many well-known connections between the Dirichlet process and urn models (Blackwell and MacQueen, 1973; Hoppe, 1984). In the current paper, we will review and expand upon some of these connections, beginning our treatment (nontraditionally) with the notion of an *exchangeable partition probability function* (EPPF) and, from there, discussing related urn models, stick-breaking representations, subordinators and random measures.

On the other hand, the Dirichlet process is limited in terms of the statistical notion of a “combination of underlying factors” that we referred to above. Indeed, the Dirichlet process is generally used in a statistical setting to express the idea that each data point is associated with one and only one underlying factor. In contrast to such *clustering models*, we wish to also consider *featural models*, where each data point is associated with a set of underlying features and it is the interaction among these features that gives rise to an observed data point. Focusing on the case in which these features are binary, we develop some of the combinatorial stochastic process machinery needed to specify featural priors. Specifically, we develop a counterpart to the EPPF, which we refer to as the *exchangeable feature probability function* (EFPF), that characterizes the combinatorial structure of certain featural models. We again develop connections between this combinatorial function and suite of related stochastic processes, including urn models, stick-breaking representations, subordinators and random measures. As we will discuss, a particular underlying random measure in this case is the *beta process*, originally studied by Hjort (1990) as a model of random hazard functions in survival analysis, but adapted by Thibaux and Jordan (2007) for applications in featural modeling.

For statistical applications it is not enough to develop expressive prior specifications, but it is also essential that inferential computations involving the posterior

distribution are tractable. One of the reasons for the popularity of the Dirichlet process is that the associated urn models and stick-breaking representations yield a variety of useful inference algorithms (Neal, 2000). As we will see, analogous algorithms are available for featural models. Thus, as we discuss each of the various representations associated with both the Dirichlet process and the beta process, we will also (briefly) discuss some of the consequences of each for posterior inference.

The remainder of the paper is organized as follows. We start by reviewing partitions and introducing feature allocations in Section 2 in order to define distributions over these models (Section 3) via the EPPF in the partition case (Section 3.1) and the EFPF in the feature allocation case (Section 3.2). Illustrating these exchangeable probability functions with examples, we will see that the well-known *Chinese restaurant process* (CRP) (Aldous, 1985) corresponds to a particular EPPF choice (Example 1) and the *Indian buffet process* (IBP) (Griffiths and Ghahramani, 2006) corresponds to a particular choice of EFPF (Example 5). From here, we progressively build up richer models by first reviewing stick lengths (Section 4), which we will see represent limiting frequencies of certain clusters or features, and then subordinators (Section 5), which further associate a random label with each cluster or feature. We illustrate these progressive augmentations for both the CRP (Examples 1, 6, 10, 18 and 20) and IBP examples (Examples 5, 7, 11 and 15). We augment the model once more to obtain a random measure on a general space of cluster or feature parameters in Section 6, and discuss how marginalization of this random measure yields the CRP in the case of the Dirichlet process (Example 23) and the IBP in the case of the beta process (Example 24). Finally, in Section 7, we mention some of the other combinatorial stochastic processes, beyond the Dirichlet process and the beta process, that have begun to be studied in the Bayesian nonparametrics literature, and we provide suggestions for further developments.

2. PARTITIONS AND FEATURE ALLOCATIONS

While we have some intuitive ideas about what constitutes a cluster or feature model, we want to formalize these ideas before proceeding. We begin with the underlying combinatorial structure on the data indices. We think of $[N] := \{1, \dots, N\}$ as representing the indices of the first N data points. There are different groupings that we apply in the cluster case (*partitions*) and feature case (*feature allocations*); we describe these below.

First, we wish to describe the space of *partitions* over the indices $[N]$. In particular, a partition π_N of $[N]$ is defined to be a collection of mutually exclusive, exhaustive, nonempty subsets of $[N]$ called *blocks*; that is, $\pi_N = \{A_1, \dots, A_K\}$ for some number of partition blocks K . An example partition of $[6]$ is $\pi_6 = \{\{1, 3, 4\}, \{2\}, \{5, 6\}\}$. Similarly, a partition of $\mathbb{N} = \{1, 2, \dots\}$ is a collection of mutually exclusive, exhaustive, nonempty subsets of \mathbb{N} . In this case, the number of blocks may be infinite, and we have $\pi_N = \{A_1, A_2, \dots\}$. An example partition of \mathbb{N} into two blocks is $\{\{n : n \text{ is even}\}, \{n : n \text{ is odd}\}\}$.

We introduce a generalization of a partition called a *feature allocation* that relaxes both the mutually exclusive and exhaustive restrictions. In particular, a feature allocation f_N of $[N]$ is defined to be a multiset of nonempty subsets of $[N]$, again called *blocks*, such that each index n can belong to any finite number of blocks. Note that the constraint that no index should belong to infinitely many blocks coincides with our intuition for the meaning of these blocks as groups to which the index belongs. Consider an example where the data points are images expressed as pixel arrays, and the latent features represent animals that may or may not appear in each picture. It is impossible to display an infinite number of animals in a picture with finitely many pixels.

We write $f_N = \{A_1, \dots, A_K\}$ for some number of feature allocation blocks K . An example feature allocation of $[6]$ is $f_6 = \{\{2, 3\}, \{2, 4, 6\}, \{3\}, \{3\}, \{3\}\}$. Just as the blocks of a partition are sometimes called *clusters*, so are the blocks of a feature allocation sometimes called *features*. We note that a partition is always a feature allocation, but the converse statement does not hold in general; for instance, f_6 given above is not a partition.

In the remainder of this section we continue our development in terms of feature allocations since partitions are a special case of the former object. We note that we can extend the idea of random partitions (Aldous, 1985) to consider *random feature allocations*. If \mathcal{F}_N is the space of all feature allocations of $[N]$, then a random feature allocation F_N of $[N]$ is a random element of this space.

We next introduce a few useful assumptions on our random feature allocation. Just as exchangeability of observations is often a central assumption in statistical modeling, so will we make use of *exchangeable feature allocations*. To rigorously define such feature allocations, we introduce the following notation. Let $\sigma : \mathbb{N} \rightarrow \mathbb{N}$ be a finite permutation. That is, for some

finite value N_σ , we have $\sigma(n) = n$ for all $n > N_\sigma$. Further, for any block $A \subset \mathbb{N}$, denote the permutation applied to the block as follows: $\sigma(A) := \{\sigma(n) : n \in A\}$. For any feature allocation F_N , denote the permutation applied to the feature allocation as follows: $\sigma(F_N) := \{\sigma(A) : A \in F_N\}$. Finally, let F_N be a random feature allocation of $[N]$. Then we say that F_N is exchangeable if $F_N \stackrel{d}{=} \sigma(F_N)$ for every finite permutation σ .

Our second assumption in what follows will be that we are dealing with a *consistent* feature allocation. We often implicitly imagine the indices arriving one at a time: first 1, then 2, up to N or beyond. We will find it useful, similarly, in defining random feature allocations to suppose that the randomness at stage n somehow agrees with the randomness at stage $n + 1$. More formally, we say that a feature allocation f_M of $[M]$ is a *restriction* of a feature allocation f_N of $[N]$ for $M < N$ if

$$f_M = \{A \cap [M] : A \in f_N\}.$$

Let $\mathcal{R}_N(f_M)$ be the set of all feature allocations of $[N]$ whose restriction to $[M]$ is f_M . Then we say that the sequence of random feature allocations (F_n) is *consistent* if for all M and N such that $M < N$, we have that

$$(1) \quad F_N \in \mathcal{R}_N(F_M) \quad \text{a.s.}$$

With this consistency condition in hand, we can define a random feature allocation F_∞ of \mathbb{N} . In particular, such a feature allocation is characterized by the sequence of consistent finite restrictions F_N to $[N]$: $F_N := \{A \cap [N] : A \in F_\infty\}$. Then F_∞ is equivalent to a consistent sequence of finite feature allocations and may be thought of as a random element of the space of such sequences: $F_\infty = (F_n)_n$. We let \mathcal{F}_∞ denote the space of consistent feature allocations, of which each random feature allocation is a random element, and we see that the sigma-algebra associated with this space is generated by the finite-dimensional sigma-algebras of the restricted random feature allocations F_n .

We say that F_∞ is exchangeable if $F_\infty \stackrel{d}{=} \sigma(F_\infty)$ for every finite permutation σ . That is, when the permutation σ changes no indices above N , we require $F_N \stackrel{d}{=} \sigma(F_N)$, where F_N is the restriction of F_∞ to $[N]$. A characterization of distributions for F_∞ is provided by Broderick, Pitman and Jordan (2013), where a similar treatment of the introductory ideas of this section also appears.

In what follows, we consider particular useful ways of representing distributions for exchangeable, consistent random feature allocations with emphasis on partitions as a special case.

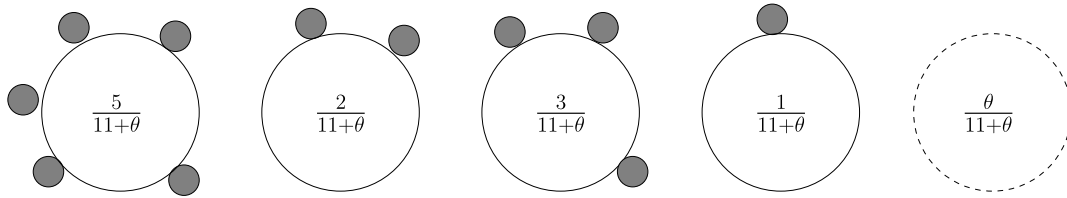


FIG. 1. The diagram represents a possible CRP seating arrangement after 11 customers have entered a restaurant with parameter θ . Each large white circle is a table, and the smaller gray circles are customers sitting at those tables. If a 12th customer enters, the expressions in the middle of each table give the probability of the new customer sitting there. In particular, the probability of the 12th customer sitting at the first table is $5/(11 + \theta)$, and the probability of the 12th customer forming a new table is $\theta/(11 + \theta)$.

3. EXCHANGEABLE PROBABILITY FUNCTIONS

Once we know that we can construct (exchangeable and consistent) random partitions and feature allocations, it remains to find useful representations of distributions over these objects.

3.1 Exchangeable Partition Probability Function

Consider first an exchangeable, consistent, random partition (Π_n) . By the exchangeability assumption, the distribution of the partition should depend only on the (unordered) sizes of the blocks. Therefore, there exists a function p that is symmetric in its arguments such that, for any specific partition assignment $\pi_n = \{A_1, \dots, A_K\}$, we have

$$(2) \quad \mathbb{P}(\Pi_n = \pi_n) = p(|A_1|, \dots, |A_K|).$$

The function p is called the *exchangeable partition probability function* (EPPF) (Pitman, 1995).

EXAMPLE 1 (Chinese restaurant process). The Chinese restaurant process (CRP) (Blackwell and MacQueen, 1973) is an iterative description of a partition via the conditional distributions of the partition blocks to which increasing data indices belong. The Chinese restaurant metaphor forms an equivalence between customers entering a Chinese restaurant and data indices; customers who share a table at the restaurant represent indices belonging to the same partition block.

To generate the label for the first index, the first customer enters the restaurant and sits down at some table, necessarily unoccupied since no one else is in the restaurant. A “dish” is set out at the new table; call the dish “1” since it is the first dish. The customer is assigned the label of the dish at her table: $Z_1 = 1$. Recursively, for a restaurant with concentration parameter θ , the n th customer sits at an occupied table with probability in proportion to the number of people at the table and at a new table with probability proportional to θ . In the former case, Z_n takes the value of the existing

dish at the table, and, in the latter case, the next available dish k (equal to the number of existing tables plus one) appears at the new table, and $Z_n = k$. By summing over all possibilities when the n th customer arrives, one obtains the normalizing constant for the distribution across potential occupied tables: $(n - 1 + \theta)^{-1}$. An example of the distribution over tables for the n th customer is shown in Figure 1. To summarize, if we let $K_n := \max\{Z_1, \dots, Z_n\}$, then the distribution of table assignments for the n th customer is

$$(3) \quad \mathbb{P}(Z_n = k | Z_1, \dots, Z_{n-1}) = (n - 1 + \theta)^{-1} \begin{cases} \#\{m : m < n, Z_m = j\}, & \text{for } j \leq K_{n-1}, \\ \theta, & \text{for } k = K_{n-1} + 1. \end{cases}$$

We note that an equivalent generative description follows a Pólya urn style in specifying that each incoming customer sits next to an existing customer with probability proportional to 1 and forms a new table with probability proportional to θ (Hoppe, 1984).

Next, we find the probability of the partition induced by considering the collection of indices sitting at each table as a block in the partition. Suppose that N_k individuals sit at table k so that the set of cardinalities of nonzero table occupancies is $\{N_1, \dots, N_K\}$ with $N := \sum_{k=1}^K N_k$. That is, we are considering the case when N customers have entered the restaurant and sat at K different tables in the specified configuration.

We can see from equation (3) that when the n th customer enters ($n > 1$), we obtain a factor of $n - 1 + \theta$ in the denominator. Using the following notation for the rising and falling factorial

$$x_{M \uparrow a} := \prod_{m=0}^{M-1} (x + ma), \quad x_{M \downarrow a} := \prod_{m=0}^{M-1} (x - ma),$$

we find a factor of $(\theta + 1)_{N-1 \uparrow 1}$ must occur in the denominator of the probability of the partition of $[N]$.

Similarly, each time a customer forms a new table except for the first table, we obtain a factor of θ in the numerator. Combining these factors, we find a factor of θ^{K-1} in the numerator. Finally, each time a customer sits at an existing table with n occupants, we obtain a factor of n in the numerator. Thus, for each table k , we have a factor of $(N_k - 1)!$ once all customers have entered the restaurant.

Having collected all terms in the process, we see that the probability of the resulting configuration is

$$(4) \quad \mathbb{P}(\Pi_N = \pi_N) = \frac{\theta^{K-1} \prod_{k=1}^K (N_k - 1)!}{(\theta + 1)_{N-1 \uparrow 1}}.$$

We first note that equation (4) depends only on the block sizes and not on the order of arrival of the customers or dishes at the tables. We conclude that the partition generated according to the CRP scheme is exchangeable. Moreover, as the partition Π_M is the restriction of Π_N to $[M]$ for any $N > M$ by construction, we have that equation (4) satisfies the consistency condition. It follows that equation (4) is, in fact, an EPPF.

3.2 Exchangeable Feature Probability Function

Just as we considered an exchangeable, consistent, random partition above, so we now turn to an exchangeable, consistent, random feature allocation (F_n) . Let $f_N = \{A_1, \dots, A_K\}$ be any particular feature allocation. In calculating $\mathbb{P}(F_N = f_N)$, we start by demonstrating in the next example that this probability in some sense undercounts features when they contain exactly the same indices: for example, $A_j = A_k$ for some $j \neq k$. For instance, consider the following example.

EXAMPLE 2 (A two-block, Bernoulli feature allocation). Let $q_A, q_B \in (0, 1)$ represent the frequencies of features A and B . Draw $Z_{A,n} \stackrel{\text{i.i.d.}}{\sim} \text{Bern}(q_A)$ and $Z_{B,n} \stackrel{\text{i.i.d.}}{\sim} \text{Bern}(q_B)$, independently. Construct the random feature allocation by collecting those indices with successful draws:

$$F_N := \{ \{n : n \leq N, Z_{A,n} = 1\}, \{n : n \leq N, Z_{B,n} = 1\} \}.$$

Then the probability of the feature allocation $F_5 = f_5 := \{\{2, 3\}, \{2, 3\}\}$ is

$$q_A^2(1 - q_A)^3 q_B^2(1 - q_B)^3,$$

but the probability of the feature allocation $F_5 = f'_5 := \{\{2, 3\}, \{2, 5\}\}$ is

$$2q_A^2(1 - q_A)^3 q_B^2(1 - q_B)^3.$$

The difference is that in the latter case the features can be distinguished, and so we must account for the two possible pairings of features to frequencies $\{q_A, q_B\}$.

Now, instead, let \tilde{F}_N be F_N with a uniform random ordering on the features. There is just a single possible ordering of f_5 , so the probability of $\tilde{F}_5 = \tilde{f}_5 := (\{2, 3\}, \{2, 3\})$ is again

$$q_A^2(1 - q_A)^3 q_B^2(1 - q_B)^3.$$

However, there are two orderings of f'_5 , so the probability of $\tilde{F}_5 = \tilde{f}'_5 := (\{2, 5\}, \{2, 3\})$ is

$$q_A^2(1 - q_A)^3 q_B^2(1 - q_B)^3,$$

and the same holds for the other ordering.

For reasons suggested by the previous example, we will find it useful to work with the random feature allocation after uniform random ordering, \tilde{F}_N . One way to achieve such an ordering and maintain consistency across different N is to associate some independent, continuous random variable with each feature; for example, assign a uniform random variable on $[0, 1]$ to each feature and order the features according to the order of the assigned random variables. When we view feature allocations constructed as marginals of a *subordinator* in Section 5, we will see that this construction is natural.

In general, given a probability of a random feature allocation, $\mathbb{P}(F_N = f_N)$, we can find the probability of a *random ordered feature allocation*, $\mathbb{P}(\tilde{F}_N = \tilde{f}_N)$ as follows. Let H be the number of unique elements of F_N , and let $(\tilde{K}_1, \dots, \tilde{K}_H)$ be the multiplicities of these unique elements in decreasing size. Then

$$(5) \quad \mathbb{P}(\tilde{F}_N = \tilde{f}_N) = \binom{K}{\tilde{K}_1, \dots, \tilde{K}_H}^{-1} \mathbb{P}(F_N = f_N),$$

where

$$\binom{K}{\tilde{K}_1, \dots, \tilde{K}_H} := \frac{K!}{\tilde{K}_1! \dots \tilde{K}_H!}.$$

We will see in Section 5 that augmentation of an exchangeable partition with a random ordering is also natural. However, the probability of an ordered random partition is not substantively different from the probability of an unordered version since the factor contributed by ordering a partition is always $1/K!$, where K here is the number of partition blocks.

With this framework in place, we can see that some ordered feature allocations have a probability function p nearly as in equation (2), that is, moreover, symmetric in its block-size arguments. Consider again the previous example.

EXAMPLE 3 (A two-block, Bernoulli feature allocation (continued)). Consider any F_N with block sizes N_1 and N_2 constructed as in Example 2. Then

$$\begin{aligned} \mathbb{P}(\tilde{F}_N = \tilde{f}_N) &= \frac{1}{2}q_A^{N_1}(1 - q_A)^{N-N_1}q_B^{N_2}(1 - q_B)^{N-N_2} \\ &\quad + \frac{1}{2}q_A^{N_2}(1 - q_A)^{N-N_2}q_B^{N_1}(1 - q_B)^{N-N_1} \\ (6) \quad &= p(N, N_1, N_2), \end{aligned}$$

where p is some function of the number of indices N and the block sizes (N_1, N_2) that we note is symmetric in all arguments after the first. In particular, we see that the order of N_1 and N_2 was immaterial.

We note that in the partition case, $\sum_{k=1}^K |A_k| = N$, so N is implicitly an argument to the EPPF. In the feature case, this summation condition no longer holds, so we make the argument N explicit in equation (6).

However, it is not necessarily the case that such a function, much less a symmetric one, exists for exchangeable feature models—in contrast to the case of exchangeable partitions and the EPPF.

EXAMPLE 4 (A general two-block feature allocation). We here describe an exchangeable, consistent random feature allocation whose (ordered) distribution does not depend only on the number of indices N and the sizes of the blocks of the feature allocation.

Let p_1, p_2, p_3, p_4 be fixed frequencies that sum to one. Let Y_n represent the collection of features to which index n belongs. For $n \in \{1, 2\}$, choose Y_n independently and identically according to

$$Y_n = \begin{cases} \{A\}, & \text{with probability } p_1, \\ \{B\}, & \text{with probability } p_2, \\ \{A, B\}, & \text{with probability } p_3, \\ \emptyset, & \text{with probability } p_4. \end{cases}$$

We form a feature allocation from these labels as follows. For each label (A or B), collect those indices n with the given label appearing in Y_n to form a feature.

Now consider two possible outcome feature allocations: $f_2 = \{\{2\}, \{2\}\}$ and $f'_2 = \{\{1\}, \{2\}\}$. The likelihood of any random ordering \tilde{f}_2 of f_2 under this model is

$$\mathbb{P}(\tilde{F}_2 = \tilde{f}_2) = p_1^0 p_2^0 p_3^1 p_4^1.$$

The likelihood of any ordering \tilde{f}'_2 of f'_2 is

$$\mathbb{P}(\tilde{F}_2 = \tilde{f}'_2) = p_1^1 p_2^1 p_3^0 p_4^0.$$

It follows from these two likelihoods that we can choose values of p_1, p_2, p_3, p_4 such that $\mathbb{P}(\tilde{F}_2 = \tilde{f}_2) \neq \mathbb{P}(\tilde{F}_2 = \tilde{f}'_2)$. But \tilde{f}_2 and \tilde{f}'_2 have the same block counts and N value ($N = 2$). So there can be no such symmetric function p , as in equation (6), for this model.

When a function p exists in the form

$$(7) \quad \mathbb{P}(\tilde{F}_N = \tilde{f}_N) = p(N, |A_1|, \dots, |A_K|)$$

for some random ordered feature allocation $\tilde{f}_N = (A_1, \dots, A_K)$ such that p is symmetric in all arguments after the first, we call it the *exchangeable feature probability function* (EFPF). Note that the EPPF is not a special case of the EFPF. The EPPF assigns zero probability to any multiset in which an index occurs in more than one element of the multiset; only the sizes of the multiset blocks are relevant in the EFPF case.

We next consider a more complex example of an EFPF.

EXAMPLE 5 (Indian buffet process). The Indian buffet process (IBP) (Griffiths and Ghahramani, 2006) is a generative model for a random feature allocation that is specified recursively like the Chinese restaurant process. Also like the CRP, this culinary metaphor forms an equivalence between customers and the indices n that will be partitioned: $n \in \mathbb{N}$. Here, “dishes” again correspond to feature labels just as they corresponded to partition labels for the CRP. But in the IBP case, a customer can sample multiple dishes.

In particular, we start with a single customer, who enters the buffet and chooses $K_1^+ \sim \text{Pois}(\gamma)$ dishes. Here, $\gamma > 0$ is called the *mass parameter*, and we will also see the *concentration parameter* $\theta > 0$ below. None of the dishes have been sampled by any other customers since no other customers have yet entered the restaurant. We label the dishes $1, \dots, K_1^+$ if $K_1^+ > 0$. Recursively, the n th customer chooses which dishes to sample in two parts. First, for each dish k that has previously been sampled by any customer in $1, \dots, n - 1$, customer n samples dish k with probability $N_{n-1,k}/(\theta + n - 1)$ for $N_{n,k}$ equal to the number of customers indexed $1, \dots, n$ who have tried dish k . As each dish represents a feature, and sampling a dish represents that the customer index n belongs to that feature, $N_{n,k}$ is the size of the block of the feature labeled k in the feature allocation of $[n]$. Next, customer n chooses $K_n^+ \sim \text{Pois}(\theta\gamma/(\theta + n - 1))$ new dishes to try. If $K_n^+ > 0$, then the dishes receive unique labels $K_{n-1} + 1, \dots, K_n$. Here, K_n represents the number of sampled dishes after n customers: $K_n = K_{n-1} + K_n^+$.

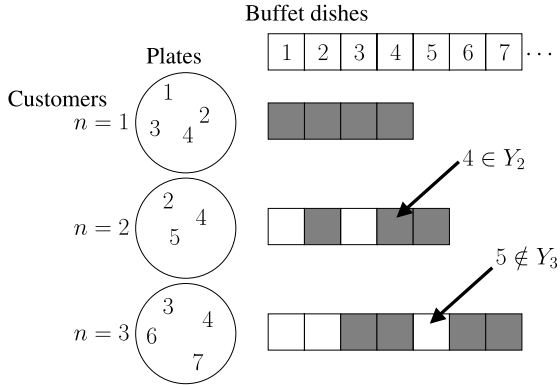


FIG. 2. Illustration of an Indian buffet process. The buffet (top) consists of a vector of dishes, corresponding to features. Each customer—corresponding to a data point—who enters first decides whether or not to eat dishes that the other customers have already sampled and then tries a random number of new dishes, not previously sampled by any customer. A gray box in position (n, k) indicates customer n has sampled dish k , and a white box indicates the customer has not sampled the dish. In the example, the second customer has sampled exactly those dishes indexed by $2, 4$ and 5 : $Y_2 = \{2, 4, 5\}$.

An example of the first few steps in the Indian buffet process is shown in Figure 2.

With this generative model in hand, we can find the probability of a particular feature allocation. We discover its form by enumeration as for the CRP EPPF in Example 1. At each round n , we have a Poisson number of new features, K_n^+ , represented. The probability factor associated with these choices is a product of Poisson densities:

$$\prod_{n=1}^N \frac{1}{K_n^+!} \left(\frac{\theta\gamma}{\theta + n - 1} \right)^{K_n^+} \exp\left(-\frac{\theta\gamma}{\theta + n - 1}\right).$$

Let M_k be the round on which the k th dish, in order of appearance, is first chosen. Then the denominators for future dish choice probabilities are the factors in the product $(\theta + M_k) \cdot (\theta + M_k + 1) \cdots (\theta + N - 1)$. The numerators for the times when the dish is chosen are the factors in the product $1 \cdot 2 \cdots (N_{N,k} - 1)$. The numerators for the times when the dish is not chosen yield $(\theta + M_k - 1) \cdots (\theta + N - 1 - N_{N,k})$. Let $A_{n,k}$ represent the collection of indices in the feature with label k after n customers have entered the restaurant. Then $N_{n,k} = |A_{n,k}|$. Finally, let $\tilde{K}_1, \dots, \tilde{K}_H$ be the multiplicities of unique feature blocks formed by this model. We note that there are

$$\left[\prod_{n=1}^N K_n^+! \right] / \left[\prod_{h=1}^H \tilde{K}_h! \right]$$

rearrangements of the features generated by this process that all yield the same feature allocation. Since they all have the same generating probability, we simply multiply by this factor to find the feature allocation probability. Multiplying all factors together and taking $f_N = \{A_{N,1}, \dots, A_{N,K_N}\}$ yields

$$\begin{aligned} \mathbb{P}(F_N = f_N) &= \frac{\prod_{n=1}^N K_n^+!}{\prod_{h=1}^H \tilde{K}_h!} \\ &\cdot \left[\prod_{n=1}^N \frac{1}{K_n^+!} \left(\frac{\theta\gamma}{\theta + n - 1} \right)^{K_n^+} \exp\left(-\frac{\theta\gamma}{\theta + n - 1}\right) \right] \\ &\cdot \left[\prod_{k=1}^{K_N} \frac{\Gamma(\theta + M_k)}{\Gamma(\theta + N)} \Gamma(N_{N,k}) \frac{\Gamma(\theta + N - N_{N,k})}{\Gamma(\theta + M_k - 1)} \right] \\ &= \left(\prod_{h=1}^H \tilde{K}_h! \right)^{-1} \left[\prod_{n=1}^N (\theta\gamma)^{K_n^+} \exp\left(-\frac{\theta\gamma}{\theta + n - 1}\right) \right] \\ &\cdot \left[\frac{\prod_{k=1}^{K_N} (\theta + M_k - 1)}{\prod_{n=1}^N (\theta + n - 1)^{K_n^+}} \right] \\ &\cdot \left[\prod_{k=1}^{K_N} \frac{\Gamma(N_{N,k}) \Gamma(\theta + N - N_{N,k})}{\Gamma(\theta + N)} \right] \\ &= \left(\prod_{h=1}^H \tilde{K}_h! \right)^{-1} (\theta\gamma)^{K_N} \\ &\cdot \exp\left(-\theta\gamma \sum_{n=1}^N (\theta + n - 1)^{-1}\right) \\ &\cdot \prod_{k=1}^{K_N} \frac{\Gamma(N_{N,k}) \Gamma(N - N_{N,k} + \theta)}{\Gamma(N + \theta)}. \end{aligned}$$

It follows from equation (5) that the probability of a uniform random ordering of the feature allocation is

$$\begin{aligned} \mathbb{P}(\tilde{F}_N = \tilde{f}_N) &= \frac{1}{K_N!} (\theta\gamma)^{K_N} \exp\left(-\theta\gamma \sum_{n=1}^N (\theta + n - 1)^{-1}\right) \\ (8) \quad &\cdot \prod_{k=1}^{K_N} \frac{\Gamma(N_{N,k}) \Gamma(N - N_{N,k} + \theta)}{\Gamma(N + \theta)}. \end{aligned}$$

The distribution of \tilde{F}_N has no dependence on the ordering of the indices in $[N]$. Hence, the distribution of F_N depends only on the same quantities—the number of indices and the feature block sizes—and the feature multiplicities. So we see that the IBP construction yields an exchangeable random feature allocation.

Consistency follows from the recursive construction and exchangeability. Therefore, equation (8) is seen to be in EPPF form [cf. equation (7)].

Above, we have seen two examples of how specifying a conditional distribution for the block membership of index n given the block membership of indices in $[n - 1]$ yields an exchangeable probability function, for example, the EPPF in the CRP case (Example 1) and the EPPF in the IBP case (Example 5). This conditional distribution is often called a *prediction rule*, and study of the prediction rule in the clustering case may be referred to as *species sampling* (Pitman, 1996; Hansen and Pitman, 1998; Lee et al., 2008). We will see next that the prediction rule can conversely be recovered from the exchangeable probability function specification and, therefore, the two are equivalent.

3.3 Induced Allocations and Block Labeling

In Examples 1 and 5 above, we formed partitions and feature allocations in the following way. For partitions, we assigned labels Z_n to each index n . Then we generated a partition of $[N]$ from the sequence $(Z_n)_{n=1}^N$ by saying that indices m and n are in the same partition block ($m \sim n$) if and only if $Z_n = Z_m$. The resulting partition is called the *induced partition* given the labels $(Z_n)_{n=1}^N$. Similarly, given labels $(Z_n)_{n=1}^\infty$, we can form an induced partition of \mathbb{N} . It is easy to check that, given a sequence $(Z_n)_{n=1}^\infty$, the induced partitions of the subsequences $(Z_n)_{n=1}^N$ will be consistent.

In the feature case, we first assigned label collections Y_n to each index n . Y_n is interpreted as a set containing the labels of the features to which n belongs. It must have finite cardinality by our definition of a feature allocation. In this case, we generate a feature allocation on $[N]$ from the sequence $(Y_n)_{n=1}^N$ by first letting $\{\phi_k\}_{k=1}^K$ be the set of unique values in $\bigcup_{n=1}^N Y_n$. Then the features are the collections of indices with shared labels: $f_N = \{\{n : \phi_k \in Y_n\} : k = 1, \dots, K\}$. The resulting feature allocation f_N is called the *induced feature allocation* given the labels $(Y_n)_{n=1}^N$. Similarly, given label collections $(Y_n)_{n=1}^\infty$, where each Y_n has finite cardinality, we can form an induced feature allocation of \mathbb{N} . As in the partition case, given a sequence $(Y_n)_{n=1}^\infty$, we can see that the induced feature allocations of the subsequences $(Y_n)_{n=1}^N$ will be consistent.

In reducing to a partition or feature allocation from a set of labels, we shed the information concerning the labels for each partition block or feature. Conversely, we introduce *order-of-appearance* labeling schemes to give partition blocks or features labels when we have, respectively, a partition or feature allocation.

In the partition case, the order-of-appearance labeling scheme assigns the label 1 to the partition block containing index 1. Recursively, suppose we have seen n indices in k different blocks with labels $\{1, \dots, k\}$. And suppose the $n + 1$ st index does not belong to an existing block. Then we assign its block the label $k + 1$.

In the feature allocation case, we note that index 1 belongs to K_1^+ features. If $K_1^+ = 0$, there are no features to label yet. If $K_1^+ > 0$, we assign these K_1^+ features labels in $\{1, \dots, K_1^+\}$. Unless otherwise specified, we suppose that the labels are chosen uniformly at random. Let $K_1 = K_1^+$. Recursively, suppose we have seen n indices and K_n different features with labels $\{1, \dots, K_n\}$. Suppose the $n + 1$ st index belongs to K_{n+1}^+ features that have not yet been labeled. Let $K_{n+1} = K_n + K_{n+1}^+$. If $K_{n+1}^+ = 0$, there are no new features to label. If $K_{n+1}^+ > 0$, assign these K_{n+1}^+ features labels in $\{K_n + 1, \dots, K_{n+1}\}$, for example, uniformly at random.

We can use these labeling schemes to find the prediction rule, which makes use of partition block and feature labels, from the EPPF or EPPF as appropriate. First, consider a partition with EPPF p . Then, given labels $(Z_n)_{n=1}^N$ with $K_N = \max\{Z_1, \dots, Z_N\}$, we wish to find the distribution of the label Z_{N+1} . Using an order-of-appearance labeling, we know that either $Z_{N+1} \in \{Z_1, \dots, Z_N\}$ or $Z_{N+1} = K_N + 1$. Let $\pi_N = \{A_{N,1}, \dots, A_{N,K_N}\}$ be the partition induced by $(Z_n)_{n=1}^N$. Let $N_{N,k} = |A_{N,k}|$. Let $\mathbb{1}(A)$ be the indicator of event A ; that is, $\mathbb{1}(A)$ equals 1 if A holds and 0 otherwise. Let $N_{N+1,k} = N_k + \mathbb{1}\{Z_{N+1} = k\}$ for $k = 1, \dots, K_{N+1}$, and set $N_{N,K_{N+1}} = 0$ for completeness. $K_{N+1} = K_N + \mathbb{1}\{Z_{N+1} > K_N\}$ is the number of partition blocks in the partition of $[N + 1]$. Then the conditional distribution satisfies

$$\begin{aligned} \mathbb{P}(Z_{N+1} = z | Z_1, \dots, Z_N) &= \frac{\mathbb{P}(Z_1, \dots, Z_N, Z_{N+1} = z)}{\mathbb{P}(Z_1, \dots, Z_N)}. \end{aligned}$$

But the probability of a certain labeling is just the probability of the underlying partition in this construction, so

$$\begin{aligned} \mathbb{P}(Z_{N+1} = z | Z_1, \dots, Z_N) &= \frac{p(N_{N+1,1}, \dots, N_{N+1,K_{N+1}})}{p(N_{N,1}, \dots, N_{N,K_N})}. \end{aligned}$$

EXAMPLE 6 (Chinese restaurant process). We continue our Chinese restaurant process example by deriving the Chinese restaurant table assignment

scheme from the EPPF in equation (4). Substituting in the EPPF for the CRP, we find

$$\begin{aligned}
& \mathbb{P}(Z_{N+1} = z | Z_1, \dots, Z_N) \\
&= \frac{p(N_{N,1}, \dots, N_{N+1, K_{N+1}})}{p(N_{N,1}, \dots, N_{N, K_N})} \\
&= \left(\theta^{K_{N+1}-1} \prod_{k=1}^{K_{N+1}} (N_{N+1,k} - 1)! \right) \\
&\quad \cdot ((\theta + 1)_{(N+1)-1 \uparrow 1})^{-1} \\
&\quad / \left(\left(\theta^{K_N-1} \prod_{k=1}^{K_N} (N_{N,k} - 1)! \right) \right. \\
&\quad \left. \cdot ((\theta + 1)_{N-1 \uparrow 1})^{-1} \right) \\
(9) \quad &= (N + \theta)^{-1} \begin{cases} N_{N,k}, & \text{for } z = k \leq K_N, \\ \theta, & \text{for } z = K_N + 1, \end{cases}
\end{aligned}$$

just as in equation (3).

To find the feature allocation prediction rule, we now imagine a feature allocation with EPPF p . Here we must be slightly more careful about counting due to feature multiplicities. Suppose that after N indices have been seen, we have label collections $(Y_n)_{n=1}^N$, containing a total of K_N features, labeled $\{1, \dots, K_N\}$. We wish to find the distribution of Y_{N+1} . Suppose $N + 1$ belongs to K_{N+1}^+ features that do not contain any index in $[N]$. Using an order-of-appearance labeling, we know that, if $K_{N+1}^+ > 0$, the K_{N+1}^+ new features have labels $K_N + 1, \dots, K_N + K_{N+1}^+$. Let $f_N = \{A_1, \dots, A_{K_N}\}$ be the feature allocation induced by $(Y_n)_{n=1}^N$. Let $N_{N,k} = |A_{N,k}|$ be the size of the k th feature. So $N_{N+1,k} = N_{N,k} + \mathbb{1}\{k \in Y_{N+1}\}$, where we let $N_{K_N+j} = 0$ for all of the features that are first exhibited by index $N + 1$: $j \in \{1, \dots, K_{N+1}^+\}$. Further, let the number of features, including new ones, be written $K_{N+1} = K_N + K_{N+1}^+$. Then the conditional distribution satisfies

$$\mathbb{P}(Y_{n+1} = y | Y_1, \dots, Y_N) = \frac{\mathbb{P}(Y_1, \dots, Y_N, Y_{N+1} = y)}{\mathbb{P}(Y_1, \dots, Y_N)}.$$

As we assume that the labels Y are consistent across N , the probability of a certain labeling is just the probability of the underlying ordered feature allocation times a combinatorial term. The combinatorial term accounts first for the uniform ordering of the new features among themselves for labeling and then for the uniform ordering of the new features among the old

features in the overall uniform random ordering:

$$\begin{aligned}
& \mathbb{P}(Y_{N+1} = y | Y_1, \dots, Y_N) \\
&= \frac{1}{K_{N+1}^+!} \cdot [(K_N + 1) \cdot (K_N + 2) \cdots K_{N+1}] \\
&\quad \cdot \frac{p(N, N_{N+1,1}, \dots, N_{N+1, K_{N+1}})}{p(N, N_{N,1}, \dots, N_{N, K_N})} \\
&= \frac{1}{K_{N+1}^+!} \cdot \frac{K_{N+1}!}{K_N!} \\
(10) \quad &\cdot \frac{p(N, N_{N+1,1}, \dots, N_{N+1, K_{N+1}})}{p(N, N_{N,1}, \dots, N_{N, K_N})}.
\end{aligned}$$

EXAMPLE 7 (Indian buffet process). Just as we derived the Chinese restaurant process prediction rule [equation (9)] from its EPPF [equation (4)] in Example 6, so can we derive the Indian buffet process prediction rule from its EPPF [equation (8)] by using equation (10). Substituting the IBP EPPF into equation (10), we find

$$\begin{aligned}
& \mathbb{P}(Y_{n+1} = y | Y_1, \dots, Y_N) \\
&= \frac{1}{K_{N+1}^+!} \cdot \frac{K_{N+1}!}{K_N!} \left(\frac{1}{K_{N+1}!} \right) (\theta \gamma)^{K_{N+1}} \\
&\quad \cdot \exp \left(-\theta \gamma \sum_{n=1}^{N+1} (\theta + n - 1)^{-1} \right) \\
&\quad \cdot \left[\prod_{k=1}^{K_{N+1}} \Gamma(N_{N+1,k}) \Gamma((N+1) - N_{N+1,k} + \theta) \right. \\
&\quad \left. / (\Gamma((N+1) + \theta)) \right] \\
&\quad / \left\{ \left(\frac{1}{K_N!} \right) (\theta \gamma)^{K_N} \right. \\
&\quad \cdot \exp \left(-\theta \gamma \sum_{n=1}^N (\theta + n - 1)^{-1} \right) \\
&\quad \cdot \left[\prod_{k=1}^{K_N} \Gamma(N_{N,k}) \Gamma(N - N_{N,k} + \theta) \right. \\
&\quad \left. / (\Gamma(N + \theta)) \right] \left. \right\} \\
&= \left[\frac{1}{K_{N+1}^+!} \exp \left(-\frac{\theta \gamma}{\theta + (N+1) - 1} \right) \right. \\
&\quad \left. \cdot \left(\frac{\theta \gamma}{\theta + (N+1) - 1} \right)^{K_{N+1}^+} \right]
\end{aligned}$$

$$\begin{aligned}
 & \cdot (\theta + (N + 1) - 1)^{K_{N+1}^+} \\
 & \cdot \left[\prod_{k=K_{N+1}}^{K_{N+1}} (\theta + (N + 1) - 1)^{-1} \right] \\
 & \cdot \prod_{k=1}^{K_N} \frac{N_k^{\mathbb{1}\{k \in z\}} (N - N_{N,k} + \theta)^{\mathbb{1}\{k \notin z\}}}{N + \theta} \\
 = & \text{Pois} \left(K_{N+1}^+ \mid \frac{\theta \gamma}{\theta + (N + 1) - 1} \right) \\
 & \cdot \prod_{k=1}^{K_N} \text{Bern} \left(\mathbb{1}\{k \in z\} \mid \frac{N_{N,k}}{N + \theta} \right).
 \end{aligned}$$

The final line is exactly the Poisson distribution for the number of new features times the Bernoulli distributions for the draws of existing features, as described in Example 5.

3.4 Inference

The prediction rule formulation of the EPPF or EFPP is particularly useful in providing a means of inferring partitions and feature allocations from a data set. In particular, we assume that we have data points X_1, \dots, X_N generated in the following manner. In the partition case, we generate an exchangeable, consistent, random partition Π_N according to the distribution specified by some EPPF p . Next, we assign each partition block a random parameter that characterizes that block. To be precise, for the k th partition block to appear according to an order-of-appearance labeling scheme, give this block a new *random* label $\phi_k \sim H$, for some continuous distribution H . For each n , let $Z_n = \phi_k$ where k is the order-of-appearance label of index n . Finally, let

$$(11) \quad X_n \stackrel{\text{indep}}{\sim} \mathcal{L}(Z_n)$$

for some distribution \mathcal{L} with parameter Z_n . The choices of both H and \mathcal{L} are specific to the problem domain.

Without attempting to survey the vast literature on clustering, we describe a stylized example to provide intuition for the preceding generative model. In this example, let n index an animal observed in the wild; $Z_n = Z_m$ indicates that animals n and m belong to the same (latent, unobserved) species; $Z_n = Z_m = \phi_k$ is a vector describing the (latent, unobserved) height and weight for that species; and X_n is the observed height and weight of the n th animal.

X_n need not even be directly observed, but equation (11) together with an EPPF might be part of a larger generative model. In a generalization of the

previous stylized example, Z_n indicates the dominant species in the n th geographical region; $Z_n = \phi_k$ indicates some overall species height and weight parameters (for the k th species); X_n indicates the height and weight parameters for species k in the n th region. That is, the height and weight for the species may vary by region. We measure and observe the height and weight $(E_{n,j})_{j=1}^J$ of some J animals in the n th region, believed to be i.i.d. draws from a distribution depending on X_n .

Note that the sequence $(Z_n)_{n=1}^N$ is sufficient to describe the partition Π_N since Π_N is the collection of blocks of $[N]$ with the same label values Z_n . The continuity of H is necessary to guarantee the a.s. uniqueness of the block values. So, if we can describe the posterior distribution of $(Z_n)_{n=1}^N$, we can in principle describe the posterior distribution of Π_N .

The posterior distribution of $(Z_n)_{n=1}^N$ conditional on $(X_n)_{n=1}^N$ cannot typically be solved for in closed form, so we turn to a method that approximates this posterior. We will see that prediction rules facilitate the design of a Markov Chain Monte Carlo (MCMC) sampler, in which we approximate the desired posterior distribution by a Markov chain of random samples proven to have the true posterior as its equilibrium distribution.

In the Gibbs sampler formulation of MCMC (Geman and Geman, 1984), we sample each parameter in turn and conditional on all other parameters in the model. In our case, we will sequentially sample each element of $(Z_n)_{n=1}^N$. The key observation here is that $(Z_n)_{n=1}^N$ is an exchangeable sequence. This observation follows by noting that the partition is exchangeable by assumption, and the sequence (ϕ_k) is exchangeable since it is i.i.d.; (Z_n) is an exchangeable sequence since it is a function of (Π_n) and (ϕ_k) . Therefore, the distribution of Z_n , given the remaining elements $\mathbf{Z}_{-n} := (Z_1, \dots, Z_{n-1}, Z_{n+1}, \dots, Z_N)$, is the same as if we thought of Z_n as the final, N th element in a sequence with $N - 1$ preceding values given by \mathbf{Z}_{-n} . And the distribution of Z_N given \mathbf{Z}_{-N} is provided by the prediction rule. The full details of the Gibbs sampler for the CRP in Examples 1 and 6 were introduced by Escobar (1994), MacEachern (1994), Escobar and West (1995) and are covered in fuller generality by Neal (2000).

It is worth noting that the sequence of order-of-appearance labels is not exchangeable; for instance, the first label is always 1. However, the prediction rule for Z_N given (Z_1, \dots, Z_{N-1}) breaks into two parts: (1) the probability of Z_N taking either a value

in $\{Z_1, \dots, Z_{N-1}\}$ or a new value and (2) the distribution of Z_N when it takes a new value. When programming such a sampler, it is often useful to simply encode the sets of unique values, which may be done by retaining any set of labels that induce the correct partition (e.g., integer labels) and separately retaining the set of unique parameter values. Indeed, updating the parameter values and partition block assignments separately can lead to improved mixing of the sampler (MacEachern, 1994).

Similarly, in the feature case, we imagine the following generative model for our data. First, let F_N be a random feature allocation generated according to the EPPF p . For the k th feature block in an order-of-appearance labeling scheme, assign a random label $\phi_k \sim H$ to this block for some continuous distribution H . For each n , let $Y_n = \{\phi_k : k \in J_n\}$, where J_n is here the set of order-of-appearance labels of the features to which n belongs. Finally, as above,

$$X_n \stackrel{\text{indep}}{\sim} \mathcal{L}(Y_n),$$

where the likelihood \mathcal{L} and parameter distribution H are again application-specific and where now \mathcal{L} depends on the variable-size collection of parameters in Y_n .

Griffiths and Ghahramani (2011) provide a review of likelihoods used in practice for feature models. To motivate some of these modeling choices, let us consider some stylized examples that provide helpful intuition. For example, let n index customers at a book-selling website; ϕ_k describes a book topic such as economics, modern art or science fiction. If ϕ_k describes science fiction books, $\phi_k \in Y_n$ indicates that the n th customer likes to buy science fiction books. But Y_n might have cardinality greater than one (the customer is interested in multiple book topics) or cardinality zero (the customer never buys books). Finally, X_n is a set of book sales for customer n on the book-selling site.

As a second example, let n index pictures in a database; ϕ_k describes a pictorial element such as a train or grass or a cow; $\phi_k \in Y_n$ indicates that picture n contains, for example, a train; finally, the observed array of pixels X_n that form the picture is generated to contain the pictorial elements in Y_n . As in the clustering case, X_n might not even be directly observed but might serve as a random effect in a deeper hierarchical model.

We observe that although the order-of-appearance label sets are not exchangeable, the sequence (Y_n) is. This fact allows the formulation of a Gibbs sampler via

the observation that the distribution of Y_n , given the remaining elements $\mathbf{Y}_{-n} := (Y_1, \dots, Y_{n-1}, Y_{n+1}, \dots, Y_N)$, is the same as if we thought of Y_n as the final, N th element in a sequence with $N - 1$ preceding values given by \mathbf{Y}_{-n} . The full details of such a sampler for the case of the IBP (Examples 5 and 7) are given by Griffiths and Ghahramani (2006).

As in the partition case, in practice, when programming the sampler, it is useful to separate the feature allocation encoding from the feature parameter values. Griffiths and Ghahramani (2006) describe how *left order form* matrices give a convenient representation of the feature allocation in this context.

4. STICK LENGTHS

Not every symmetric function defined for an arbitrary number of arguments with values in the unit interval is an EPPF (Pitman, 1995), and not every symmetric function with an additional positive integer argument is an EPPF. For instance, the consistency property in equation (1) implies certain additivity requirements for the function p .

EXAMPLE 8 (Not an EPPF). Consider the function p defined with

$$(12) \quad p(1) = 1, \quad p(1, 1) = 0.1, \quad p(2) = 0.8, \dots$$

From the information in equation (12), p may be further defined so as to be symmetric in its arguments for any number of arguments, but since it does not satisfy $p(1) = p(1, 1) + p(2)$, it cannot be an EPPF.

EXAMPLE 9 (Not an EPPF). Consider the function p defined with

$$(13) \quad \begin{aligned} p(N = 1) &= 0.9, & p(N = 1, 1) &= 0.9, \\ p(N = 1, 1, 1) &= 0.9, & \dots \end{aligned}$$

From the information in equation (13), p may be further defined so as to be symmetric in its arguments for any number of arguments after the initial N argument, but since $p(N = 1) + p(N = 1, 1) + p(N = 1, 1, 1) > 1$, it cannot be an EPPF.

It therefore requires some care to define a suitable distribution over consistent, exchangeable random feature allocations or partitions using the exchangeable probability function framework.

Since we are working with exchangeable sequences of random variables, it is natural to turn to de Finetti's theorem (De Finetti, 1931; Hewitt and Savage, 1955) for clues as to how to proceed. De Finetti's theorem

tells us that any exchangeable sequence of random variables can be expressed as an independent and identically distributed sequence when conditioned on an underlying random *mixing measure*. While this theorem may seem difficult to apply directly to, for example, exchangeable partitions, it may be applied more naturally to an exchangeable sequence of numbers derived from a sequence of partitions. The argument below is due to Aldous (1985).

Suppose that (Π_n) is an exchangeable, consistent sequence of random partitions. Consider the k th partition block to appear according to an order-of-appearance labeling scheme, and give this block a new *random label*, $\phi_k \sim \text{Unif}([0, 1])$, such that each random label is drawn independently from the rest. This construction is the same as the one used for parameter generation in Section 3.4, and (Π_n) is exchangeable by the same arguments used there. Let Z_n equal ϕ_k exactly when n belongs to the partition with this label.

If we apply de Finetti’s theorem to the sequence (Z_n) and note that (Z_n) has at most countably many different values, we see that there exists some random sequence (ρ_k) such that $\rho_k \in (0, 1]$ for all k and, conditioned on the frequencies (ρ_k) , (Z_n) has the same distribution as i.i.d. draws from (ρ_k) . In this description, we have brushed over technicalities associated with partition blocks that contain only one index even as $N \rightarrow \infty$ (which may imply $\sum_k \rho_k < 1$).

But if we assume that every partition block eventually contains at least two indices, we can achieve an exchangeable partition of $[N]$ as follows. Let (ρ_k) represent a sequence of values in $(0, 1]$ such that $\sum_{k=1}^{\infty} \rho_k \stackrel{\text{a.s.}}{=} 1$. Draw $Z_n \stackrel{\text{i.i.d.}}{\sim} \text{Discrete}((\rho_k)_k)$. Let Π_N be the induced partition given $(Z_n)_{n=1}^N$. Exchangeability follows from the i.i.d. draws, and consistency follows from the induced partition construction.

When the frequencies (ρ_k) are thought of as subintervals of the unit interval, that is, a partition of the unit interval, they are collectively called *Kingman’s paint-box* (Kingman, 1978). As another naming convention, we may think of the unit interval as a *stick* (Ishwaran and James, 2001). We partition the unit interval by breaking it into various *stick lengths*, which represent the frequencies of each partition block.

A similar construction can be seen to yield exchangeable, consistent random feature allocations. In this case, let (ξ_k) represent a sequence of values in $(0, 1]$ such that $\sum_{k=1}^{\infty} \xi_k \stackrel{\text{a.s.}}{<} \infty$. We generate feature collections independently for each index as follows. Start with $Y_n = \emptyset$. For each feature k , add k to the set Y_n , independently from all other features, with probability ξ_k . Let F_N be the induced feature allocation given $(Y_n)_{n=1}^N$. Exchangeability of F_N follows from the i.i.d. draws of Y_n , and consistency follows from the induced feature allocation construction. The finite sum constraint ensures each index belongs to a finite number of features a.s.

It remains to specify a distribution on the partition or feature frequencies. The frequencies cannot be i.i.d. due to the finite summation constraint in both cases. In the partition case, any infinite set of frequencies cannot even be independent since the summation is fixed to one. One scheme to ensure summation to unity is called *stick-breaking* (McCloskey, 1965; Patil and Taillie, 1977; Sethuraman, 1994; Ishwaran and James, 2001). In stick-breaking, the stick lengths are obtained by recursively breaking off parts of the unit interval to return as the atoms ρ_1, ρ_2, \dots (cf. Figure 3). In particular, we generate stick-breaking proportions V_1, V_2, \dots as $[0, 1]$ -valued random variables. Then ρ_1 is the first proportion V_1 times the initial stick length 1; hence, $\rho_1 = V_1$. Recursively, after k breaks, the remaining

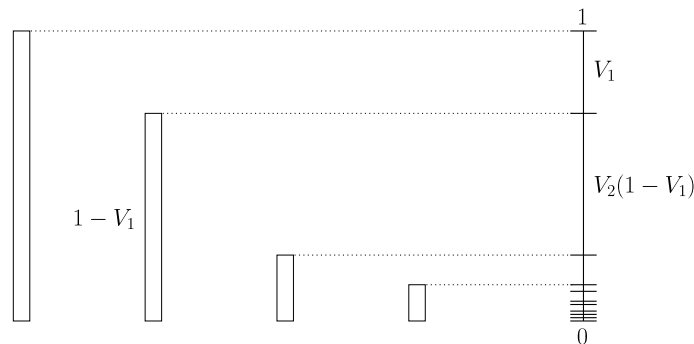


FIG. 3. An illustration of how stick-breaking divides the unit interval into a sequence of probabilities Broderick, Jordan and Pitman (2012). The stick proportions (V_1, V_2, \dots) determine what fraction of the remaining stick is appended to the probability sequence at each round.

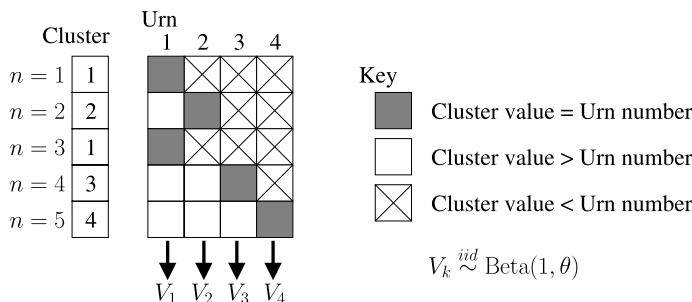


FIG. 4. An illustration of the proof based on the Pólya urn that Dirichlet process stick-breaking gives the underlying partition block frequencies for a Chinese restaurant process model. The k th column in the central matrix corresponds to a tallying of when the k th table is chosen (gray), when a table of index larger than k is chosen (white), and when an index smaller than k is chosen (\times). If we ignore the \times tallies, the gray and white tallies in each column (after the first) can be modeled as balls drawn from a Pólya urn. The limiting frequency of gray balls in each column is shown below the matrix.

length of the initial unit interval is $\prod_{j=1}^k (1 - V_j)$. And ρ_{k+1} is the proportion V_{k+1} of the remaining stick; hence, $\rho_{k+1} = V_{k+1} \prod_{j=1}^k (1 - V_j)$.

The stick-breaking construction yields ρ_1, ρ_2, \dots such that $\rho_k \in [0, 1]$ for each k and $\sum_{k=1}^{\infty} \rho_k \leq 1$. If the V_k do not decay too rapidly, we will have $\sum_{k=1}^{\infty} \rho_k \stackrel{\text{a.s.}}{=} 1$. In particular, the partition block proportions ρ_k sum to unity a.s. iff there is no remaining stick mass: $\prod_{k=1}^{\infty} (1 - V_k) \stackrel{\text{a.s.}}{=} 0$.

We often make the additional, convenient assumption that the V_k are independent. In this case, a necessary and sufficient condition for $\sum_{k=1}^{\infty} \rho_k \stackrel{\text{a.s.}}{=} 1$ is $\sum_{k=1}^{\infty} \mathbb{E}[\log(1 - V_k)] = -\infty$ (Ishwaran and James, 2001). When the V_k are independent and of a canonical distribution, they are easily simulated. Moreover, if we assume that the V_k are such that the ρ_k decay sufficiently rapidly in k , one strategy for simulating a stick-breaking model is to ignore all $k > K$ for some fixed, finite K . This approximation is known as truncation (Ishwaran and James, 2001). It is fortuitously the case that in some models of particular interest, such useful assumptions fall out naturally from the model construction (e.g., Examples 10 and 11).

EXAMPLE 10 (Chinese restaurant process). In the original exchangeability result due to de Finetti (De Finetti, 1931), the exchangeable random variables were zero/one-valued, and the mixing measure was a distribution on a single frequency so that the outcomes were conditionally Bernoulli. We will find a similar result in obtaining the stick-breaking proportions associated with the Chinese restaurant process.

We can construct a sequence of binary-valued random variables by dividing the customers in the CRP who are sitting at the first table from the rest; color

the former collection of customers gray and the latter collection of customers white. Then, we see that the first customer must be colored gray. And thus we begin with a single gray customer and no white customers. This binary valuation for the first table in the CRP is illustrated by the first column in the matrix in Figure 4.

At this point, it is useful to recall the Pólya urn construction (Pólya, 1930; Freedman, 1965), whereby an urn starts with G_0 gray balls and W_0 white balls. At each round N , we draw a ball from the urn, replace it, and add κ of the same color of ball to the urn. At the end of the round, we have G_N gray balls and W_N white balls. Despite the urn metaphor, the number of balls need not be an integer at any time. By checking equation (3), which defines the CRP, we can see that the coloring of the gray/white customer matrix assignments starting with the second customer has the same distributions as a sequence of balls from a Pólya urn as a Pólya urn with $G_{1,0} = 1$ initial gray balls, $W_{1,0} = \theta$ initial white balls and $\kappa_1 = 1$ replacement balls. Let $G_{1,N}$ and $W_{1,N}$ represent the numbers of gray and white balls, respectively, in the urn after N rounds. The important fact about the Pólya urn we use here is that there exists some $V \sim \text{Beta}(G_0/\kappa, W_0/\kappa)$ such that $\kappa^{-1}(G_{N+1} - G_N) \stackrel{i.i.d.}{\sim} \text{Bern}(V)$ for all N . In this particular case of the CRP, then, $G_{1,N+1} - G_{1,N}$ is one if a customer sits at the first table (or zero otherwise), and $G_{1,N+1} - G_{1,N} \stackrel{i.i.d.}{\sim} \text{Bern}(V_1)$ with $V_1 \sim \text{Beta}(1, \theta)$.

We now look at the sequence of customers who sit at the second and subsequent tables. That is, we condition on customers not sitting at the first table or equivalently on the sequence with $G_{1,N+1} - G_{1,N} = 0$. Again, we have that the first customer sits at the second table, by the CRP construction. Now let customers at the second table be colored gray and customers at

the third and later tables be colored white. This valuation is illustrated in the second column in Figure 4; each \times in the figure denotes a data point where the first partition block is chosen and, therefore, the current Pólya urn is not in play. As before, we begin with one gray customer and no white customers. We can check equation (3) to see that customer coloring once more proceeds according to a Pólya urn scheme with $G_{2,0} = 1$ initial gray balls, $W_{2,0} = \theta$ initial white balls and $\kappa_2 = 1$ replacement balls. Thus, contingent on a customer not sitting at the first table, the N th customer sits at the second table with i.i.d. distribution $\text{Bern}(V_2)$ with $V_2 \sim \text{Beta}(1, \theta)$. Since the sequence of individuals sitting at the second table has no other dependence on the sequence of individuals sitting at the first table, we have that V_2 is independent of V_1 .

The argument just outlined proceeds recursively to show us that the N th customer, conditional on not sitting at the first $K - 1$ tables for $K \geq 1$, sits at the K th table with i.i.d. distribution $\text{Bern}(V_K)$ and $V_K \sim \text{Beta}(1, \theta)$ with V_K independent of the previous (V_1, \dots, V_{K-1}) .

Combining these results, we see that we have the following construction for the customer seating patterns. The V_k are distributed independently and identically according to $\text{Beta}(1, \theta)$. The probability ρ_K of sitting at the K th table is the probability of not sitting at the first $K - 1$ tables, conditional on not sitting at the previous table, times the conditional probability of sitting at the K th table: $\rho_K = [\prod_{k=1}^{K-1} (1 - V_k)] \cdot V_K$. Finally, with the vector of table frequencies (ρ_k) , each customer sits independently and identically at the corresponding vector of tables according to these frequencies. This process is summarized here:

$$\begin{aligned}
 & V_k \stackrel{\text{i.i.d.}}{\sim} \text{Beta}(1, \theta), \\
 (14) \quad & \rho_K := V_K \prod_{k=1}^K (1 - V_k), \\
 & Z_n \stackrel{\text{i.i.d.}}{\sim} \text{Discrete}((\rho_k)_k).
 \end{aligned}$$

To see that this process is well-defined, first note that $\mathbb{E}[\log(1 - V_k)]$ exists, is negative and is the same for all k values. It follows that $\sum_{k=1}^{\infty} \mathbb{E}[\log(1 - V_k)] = -\infty$, so by the discussion before this example, we must have $\sum_{k=1}^K \rho_k \stackrel{\text{a.s.}}{=} 1$.

The feature case is easier. Since it does not require the frequencies to sum to one, the random frequencies can be independent so long as they have an a.s. finite sum.

EXAMPLE 11 (Indian buffet process). As in the case of the CRP, we can recover the stick lengths for the Indian buffet process using an argument based on an urn model.

Recall that on the first round of the Indian buffet process, $K_1^+ \sim \text{Pois}(\gamma)$ features are chosen to contain index 1. Consider one of the features, labeled k . By construction, each future data point N belongs to this feature with probability $N_{N-1,k}/(\theta + N - 1)$. Thus, we can model the sequence after the first data point as a Pólya urn of the sort encountered in Example 10 with initially $G_{k,0} = 1$ gray balls, $W_{k,0} = \theta$ white balls and $\kappa_k = 1$ replacement balls. As we have seen, there exists a random variable $V_k \sim \text{Beta}(1, \theta)$ such that representation of this feature by data point N is chosen, i.i.d. across all N , as $\text{Bern}(V_k)$. Since the Bernoulli draws conditional on previous draws are independent across all k , the V_k are likewise independent of each other; this fact is also true for k in future rounds. Draws according to such an urn are illustrated in each of the first four columns of the matrix in Figure 5.

Now consider any round n . According to the IBP construction, $K_n^+ \sim \text{Pois}(\gamma\theta/(\theta + n - 1))$ new features are chosen to include index n . Each future data point N (with $N > n$) represents feature k among these features with probability $N_{N-1,k}/(\theta + N - 1)$. In this case, we can model the sequence after the n th data point as a Pólya urn with $G_{k,0} = 1$ initial gray balls, $W_{k,0} = \theta + n - 1$ initial white balls and $\kappa_k = 1$ replacement balls. So there exists a random variable $V_k \sim \text{Beta}(1, \theta + n - 1)$ such that representation of feature k by data point N is chosen, i.i.d. across all N , as $\text{Bern}(V_k)$.

Finally, then, we have the following generative model for the feature allocation by iterating across $n = 1, \dots, N$ (Thibaux and Jordan, 2007):

$$\begin{aligned}
 (15) \quad & K_n^+ \stackrel{\text{indep}}{\sim} \text{Pois}\left(\frac{\gamma\theta}{\theta + n - 1}\right), \\
 & K_n = K_{n-1} + K_n^+, \\
 (16) \quad & V_k \stackrel{\text{indep}}{\sim} \text{Beta}(1, \theta + n - 1), \\
 & \qquad \qquad \qquad k = K_{n-1} + 1, \dots, K_n, \\
 & I_{n,k} \stackrel{\text{indep}}{\sim} \text{Bern}(V_k), \quad k = 1, \dots, K_n.
 \end{aligned}$$

$I_{n,k}$ is an indicator random variable for whether feature k contains index n . The collection of features to which index n belongs, Y_n , is the collection of features k with $I_{n,k} = 1$.

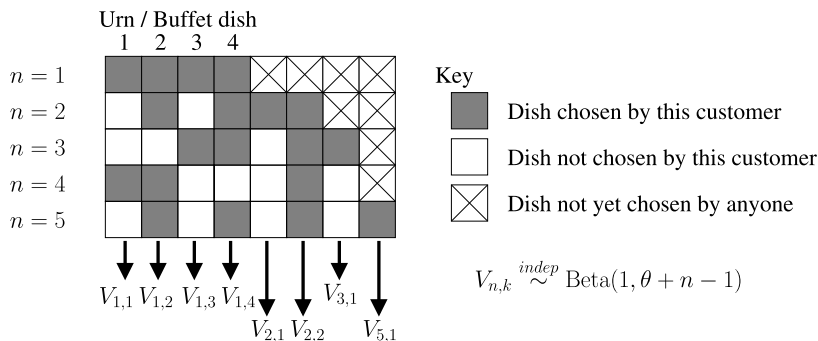


FIG. 5. Illustration of the proof that the frequencies of features in the Indian buffet process are given by beta random variables. For each feature, we can construct a sequence of zero/one variables by tallying whether (gray, one) or not (white, zero) that feature is represented by the given data point. Before the first time a feature is chosen, we mark it with an \times . Each column sequence of gray and white tallies, where we ignore the \times marks, forms a Pólya urn with limiting frequencies shown below the matrix.

4.1 Inference

As we have seen above, the exchangeable probability functions of Section 3 are the marginal distributions of the partitions or feature allocations generated according to stick-length models with the stick lengths integrated out. It has been proposed that including the stick lengths in MCMC samplers of these models will improve mixing (Ishwaran and Zarepour, 2000). While it is impossible to sample the countably infinite set of partition block or feature frequencies in these models (cf. Examples 10 and 11), a number of ways of getting around this difficulty have been investigated. Ishwaran and Zarepour (2000) examine two separate finite approximations to the full CRP stick-length model: one uses a parametric approximation to the full infinite model, and the other creates a truncation by setting the stick break at some fixed size K to be 1: $V_K = 1$. There also exist techniques that avoid any approximations and deal instead directly with the full model, in particular, retrospective sampling (Papaspiliopoulos and Roberts, 2008) and slice sampling (Walker, 2007).

While our discussion thus far has focused on MCMC sampling as a means of approximating the posterior distribution of either the block assignments or both the block assignments and stick lengths, including the stick lengths in a posterior analysis facilitates a different posterior approximation; in particular, *variational methods* can also be used to approximate the posterior. These methods minimize some notion of distance to the posterior over a family of potential approximating distributions (Jordan et al., 1999). The practicality and, indeed, speed of these methods in the case of stick-breaking for the CRP (Example 10) have been demonstrated by Blei and Jordan (2006).

A number of different models for the stick lengths corresponding to the features of an IBP (Example 11) have been discovered. The distributions described in Example 11 are covered by Thibaux and Jordan (2007), who build on work from Hjort (1990), Kim (1999). A special case of the IBP is examined by Teh, Görür and Ghahramani (2007), who detail a slice sampling algorithm for sampling from the posterior of the stick lengths and feature assignments. Yet another stick-length model for the IBP is explored by Paisley et al. (2010), who show how to apply variational methods to approximate the posterior of their model.

Stick-length modeling has the further advantage of allowing inference in cases where it is not straightforward to integrate out the underlying stick lengths to obtain a tractable exchangeable probability function.

5. SUBORDINATORS

An important point to reiterate about the labels Z_n and label collections Y_n is that when we use the order-of-appearance labeling scheme for partition or feature blocks described above, the random sequences (Z_n) and (Y_n) are not exchangeable. Often, however, we would like to make use of special properties of exchangeability when dealing with these sequences. For instance, if we use Markov Chain Monte Carlo to sample from the posterior distribution of a partition (cf. Section 3.4), we might want to Gibbs sample the cluster assignment of data point n given the assignments of the remaining data points: Z_n given $\{Z_m\}_{m=1}^N \setminus \{Z_n\}$. This sampling is particularly easy in some cases (Neal, 2000) if we can treat Z_n as the last random variable in the sequence, but this treatment requires exchangeability.

A way to get around this dilemma was suggested by Aldous (1985) and appeared above in our motivation for using stick lengths. Namely, we assign to the k th partition block a uniform random label $\phi_k \sim \text{Unif}([0, 1])$; analogously, we assign to the k th feature a uniform random label $\phi_k \sim \text{Unif}([0, 1])$. We can see that in both cases, all of the labels are a.s. distinct. Now, in the partition case, let Z_n be the uniform random label of the partition block to which n belongs. And in the feature case, let Y_n be the (finite) set of uniform random feature labels for the features to which n belongs. We can recover the partition or feature allocation as the induced partition or feature allocation by grouping indices assigned to the same label. Moreover, as discussed above, we now have that each of (Z_n) and (Y_n) is an exchangeable sequence.

If we form partitions or features according to the stick-length constructions detailed in Section 4, we know that each unique partition or feature label ϕ_k is associated with a frequency ξ_k . We can use this association to form a random measure:

$$(17) \quad \mu = \sum_{k=1}^{\infty} \xi_k \delta_{\phi_k},$$

where δ_{ϕ_k} is a unit point mass located at ϕ_k . In the partition case, $\sum_k \xi_k = 1$, so the random measure is a random probability measure, and we may draw $Z_n \stackrel{\text{i.i.d.}}{\sim} \mu$. In the feature case, the weights have a finite sum but do not necessarily sum to one. In the feature case, we draw Y_n by including each ϕ_k for which $\text{Bern}(\xi_k)$ yields a draw of 1.

Another way to codify the random measure in equation (17) is as a monotone increasing stochastic process on $[0, 1]$. Let

$$T_s = \sum_{k=1}^{\infty} \xi_k \mathbb{1}\{\phi_k \leq s\}.$$

Then the atoms of μ are in one-to-one correspondence with the jumps of the process T .

This increasing random function construction gives us another means of choosing distributions for the weights ξ_k . We have already seen that these cannot be i.i.d. due to the finite summation condition. However, we will see that if we require that the increments of a monotone, increasing stochastic process are independent and stationary, then we can use the jumps of that function as the atoms in our random measure for partitions or features.

DEFINITION 12. A subordinator (Bochner, 1955; Bertoin 1996, 1999) is a stochastic process $(T_s, s \geq 0)$ that has the following properties:

- Nonnegative, nondecreasing paths (a.s.),
- Paths that are right-continuous with left limits, and
- Stationary, independent increments.

For our purposes, wherein the subordinator values will ultimately correspond to (perhaps scaled) probabilities, we will assume the subordinator takes values in $[0, \infty)$, though alternative ranges with a sense of ordering are possible.

Subordinators are of interest to us because they not only exhibit the stationary independent increments property but they also can always be decomposed into two components: a deterministic drift component and a Poisson point process. Recall that a Poisson point process on space S with rate measure $\nu(dx)$, where $x \in S$, yields a countable subset of points of S . Let $N(A)$ be the number of points of the process in set A for $A \subseteq S$. The process is characterized by the fact that, first, $N(A) \sim \text{Pois}(\nu(A))$ for any A and, second, for any disjoint A_1, \dots, A_K , we have that $N(A_1), \dots, N(A_K)$ are independent random variables. See Kingman (1993) for a thorough treatment of these processes. An example subordinator with both drift and jump components is shown on the left-hand side of Figure 6.

The subordinator decomposition is detailed in the following result (Bertoin, 1996).

THEOREM 13. Every subordinator $(T_s, s \geq 0)$ can be written as

$$(18) \quad T_s = cs + \sum_{k=1}^{\infty} \xi_k \mathbb{1}\{\phi_k \leq s\}$$

for some constant $c \geq 0$ and where $\{(\xi_k, \phi_k)\}_k$ is the countable set of points of a Poisson point process with intensity $\Lambda(d\xi) d\phi$, where Λ is a Lévy measure; that is,

$$\int_0^{\infty} (1 \wedge \xi) \Lambda(d\xi) < \infty.$$

In particular, then, if a subordinator is finite at time t , the jumps of the subordinator up to t may be used as feature block frequencies if they have support in $[0, 1]$. Or, in general, the normalized jumps may be used as partition block frequencies. We can see from the right-hand side of Figure 6 that the jumps of a subordinator partition intervals of the form $[0, t)$, as long as the subordinator has no drift component. In either the feature or cluster case, we have substituted the condition of independent and identical distribution for the partition or feature frequencies (i.e., the jumps) with a more natural continuous-time analogue: independent, stationary intervals.

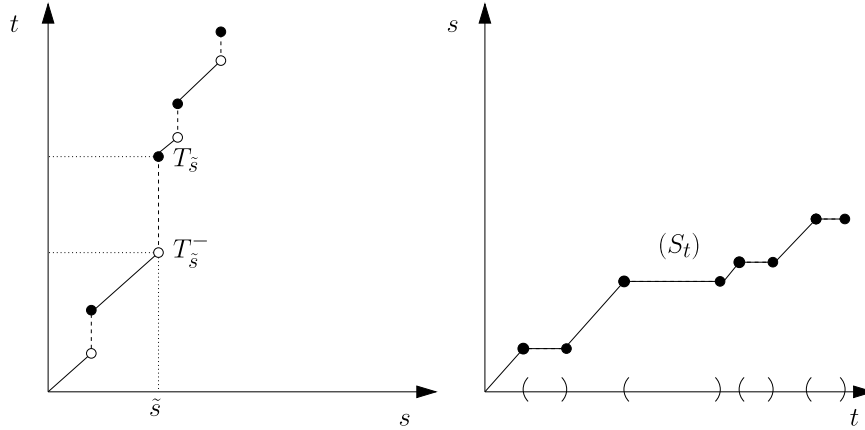


FIG. 6. Left: The sample path (T_s) of a subordinator. $T_{\tilde{s}}^-$ is the limit from the left of (T_s) at $s = \tilde{s}$. Right: The right-continuous inverse (S_t) of a subordinator: $S_t := \inf\{s : T_s > t\}$. The open intervals along the t axis correspond to the jumps of the subordinator (T_s) .

Just as the Laplace transform of a positive random variable characterizes the distribution of that random variable, so does the Laplace transform of the subordinator—which is a positive random variable at any fixed time point—describe this stochastic process (Bertoin 1996, 1999).

THEOREM 14 (Lévy–Khinchin formula for subordinators). *If $(T_s, s \geq 0)$ is a subordinator, then for $\lambda \geq 0$ we have*

$$(19) \quad \mathbb{E}(e^{-\lambda T_s}) = e^{-\Psi(\lambda)s}$$

with

$$(20) \quad \Psi(\lambda) = c\lambda + \int_0^\infty (1 - e^{-\lambda\xi})\Lambda(d\xi),$$

where $c \geq 0$ is called the drift constant and Λ is a non-negative, Lévy measure on $(0, \infty)$.

The function $\Psi(\lambda)$ is called the *Laplace exponent* in this context. We note that a subordinator is characterized by its drift constant and Lévy measure.

Using subordinators for feature allocation modeling is particularly easy; since the jumps of the subordinators are formed by a Poisson point process, we can use Poisson process methodology to find the stick lengths and EFPF. To set up this derivation, suppose we generate feature membership from a subordinator by taking Bernoulli draws at each of its jumps with success probability equal to the jump size. Since every jump has strictly positive size, the feature associated with each jump will eventually score a Bernoulli success for some index n with probability one. Therefore, we can enumerate all jumps of the process in order of appearance; that is, we first enumerate all features in which

index 1 appears, then all features in which index 2 appears but not index 1, and so on. At the n th iteration, we enumerate all features in which index n appears but not previous indices. Let K_n^+ represent the number of indices so chosen on the n th round. Let $K_0 = 0$ so that recursively $K_n := K_{n-1} + K_n^+$ is the number of subordinator jumps seen by round n , inclusive. Let ξ_k for $k = K_{n-1} + 1, \dots, K_n$ be the distribution of a particular subordinator jump seen on round n . We now turn to connecting the subordinator perspective to the earlier derivation of stick lengths in Section 4.

EXAMPLE 15 (Indian buffet process). In our earlier discussion, we found a collection of stick lengths to represent the featural frequencies for the IBP [equation (16) of Example 11 in Section 4]. To see the connection to subordinators, we start from the *beta process subordinator* (Kim, 1999) with zero drift ($c = 0$) and Lévy measure

$$(21) \quad \Lambda(d\xi) = \gamma\theta\xi^{-1}(1 - \xi)^{\theta-1} d\xi.$$

We will see that the mass parameter $\gamma > 0$ and concentration parameter $\theta > 0$ are the same as those introduced in Example 5 and continued in Example 11.

THEOREM 16. *Generate a feature allocation from a beta process subordinator with Lévy measure given by equation (21). Then the sequence of subordinator jumps (ξ_k) , indexed in order of appearance, has the same distribution as the sequence of IBP stick lengths (V_k) described by equations (15) and (16).*

PROOF. Recall the following fact about Poisson thinning (Kingman, 1993), illustrated in Figure 7. Suppose that a Poisson point process with rate measure

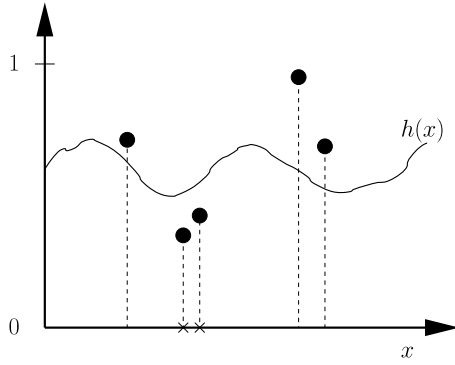


FIG. 7. An illustration of Poisson thinning. The x -axis values of the filled black circles, emphasized by dotted lines, are generated according to a Poisson process. The $[0, 1]$ -valued function $h(x)$ is arbitrary. The vertical axis values of the points are uniform draws in $[0, 1]$. The “thinned” points are the collection of x -axis values corresponding to vertical axis values below $h(x)$ and are denoted with a \times symbol.

λ generates points with values x . Then suppose that, for each such point x , we keep it with probability $h(x) \in [0, 1]$. The resulting set of points is also a Poisson point process, now with rate measure $\lambda'(A) = \int_A \lambda(dx)h(x) dx$.

We prove Theorem 16 recursively. Define the measure

$$\mu_n(d\xi) := \gamma\theta\xi^{-1}(1 - \xi)^{\theta+n-1} d\xi,$$

so that μ_0 is the beta process Lévy measure Λ in equation (21). We make the recursive assumption that μ_n is distributed as the beta process measure without atoms corresponding to features chosen on the first n iterations.

There are two parts to proving Theorem 16. First, we show that, on the n th iteration, the number of features chosen and the distribution of the corresponding atom weights agree with equations (15) and (16), respectively. Second, we check that the recursion assumption holds.

For the first part, note that on the n th round we choose features with probability equal to their atom weight. So we form a thinned Poisson process with rate measure $\xi \cdot \mu_{n-1}(d\xi)$. This rate measure has total mass

$$\int_0^1 \xi \cdot \mu_{n-1}(d\xi) = \gamma \frac{\theta}{\theta + n - 1} =: \gamma_{n-1}.$$

So the number of features chosen is Poisson-distributed with mean $\gamma\theta(\theta + n - 1)^{-1}$, as desired [cf. equation (15)]. And the atom weights have distribution equal to the normalized rate measure

$$\begin{aligned} & \gamma_{n-1}^{-1} \xi \cdot \gamma\theta\xi^{-1}(1 - \xi)^{\theta+(n-1)-1} d\xi \\ & = \text{Beta}(\xi|1, \theta + n - 1) d\xi \end{aligned}$$

as desired [cf. equation (16)].

Finally, to check the recursion assumption, we note that those sticks that remain were chosen for having Bernoulli failure draws; that is, they were chosen with probability equal to one minus their atom weight. So the thinned rate measure for the next round is

$$(1 - \xi) \cdot \gamma\theta\xi^{-1}(1 - \xi)^{\theta+(n-1)-1} d\xi,$$

which is just μ_n . \square

The form of the EPPF of the feature allocation generated from the beta process subordinator follows immediately from the stick-length distributions we have just derived by the discussion in Example 11 in Section 4.

We see from the previous example that feature allocation stick lengths and EPPFs can be obtained in a straightforward manner using the Poisson process representation of the jumps of the subordinator. Partitions, however, are not as easy to analyze, principally due to the fact that the subordinator jumps must first be normalized to obtain a probability measure on $[0, 1]$; a random measure with finite total mass is not sufficient in the partition case. Hence, we must compute the stick lengths and EPPF using partition block frequencies from these normalized jumps instead of directly from the subordinator jumps.

In the EPPF case, we make use of a result that gives us the exchangeable probability function as a function of the Laplace exponent. Though we do not derive this formula here, its derivation can be found in Pitman (2003); the proof relies on, first, calculating the joint distribution of the subordinator jumps and partition generated from the normalized jumps and, second, integrating out the subordinator jumps to find the partition marginal.

THEOREM 17. *Form a probability measure μ by normalizing jumps of the subordinator with Laplace exponent Ψ . Let (Π_n) be a consistent set of exchangeable partitions induced by i.i.d. draws from μ . For each exchangeable partition $\pi_N = \{A_1, \dots, A_K\}$ of $[N]$ with $N_k := |A_k|$ for each k ,*

$$\begin{aligned} & \mathbb{P}(\Pi_N = \pi_N) \\ & = p(N_1, \dots, N_K) \\ (22) \quad & = \frac{(-1)^{N-K}}{(N-1)!} \int_0^\infty \lambda^{N-1} e^{-\Psi(\lambda)} \prod_{k=1}^K \Psi^{(N_k)}(\lambda) d\lambda, \end{aligned}$$

where $\Psi^{(N_k)}(\lambda)$ is the N_k th derivative of the Laplace exponent Ψ evaluated at λ .

EXAMPLE 18 (Chinese restaurant process). We start by introducing the *gamma process*, a subordinator that we will see below generates the Chinese restaurant process EPPF. The gamma process has Laplace exponent $\Psi(\lambda)$ [equation (19)] characterized by

$$(23) \quad c = 0 \quad \text{and} \quad \Lambda(d\xi) = \theta \xi^{-1} e^{-b\xi} d\xi$$

for $\theta > 0$ and $b > 0$ [cf. equation (20) in Theorem 14]. We will see that θ corresponds to the CRP concentration parameter and that b is arbitrary and does not affect the partition model.

We calculate the EPPF using Theorem 17.

THEOREM 19. *The EPPF for partition block membership chosen according to the normalized jumps (ρ_k) of the gamma subordinator with parameter θ is the CRP EPPF [equation (4)].*

PROOF. By Theorem 17, if we can find all order derivatives of the Laplace exponent Ψ , we can calculate the EPPF for the partitions generated with frequencies equal to the normalized jumps of this subordinator. The derivatives of Ψ , which are known to always exist (Bertoin, 2000; Rogers and Williams, 2000), are straightforward to calculate if we begin by noting that, from equation (20) in Theorem 14, we have in general that

$$\Psi'(\lambda) = c + \int_0^\infty \xi e^{-\lambda\xi} \Lambda(d\xi).$$

Hence, for the gamma process subordinator,

$$\Psi'(\lambda) = \int_0^\infty e^{-\lambda\xi} \theta e^{-b\xi} d\xi = \frac{\theta}{\lambda + b}.$$

Then simple integration and differentiation yield

$$\Psi(\lambda) = \theta \log(\lambda + b) - \theta \log(b)$$

since $\Psi(0) = 0$ and

$$\Psi^{(n)}(\lambda) = (-1)^{n-1} \frac{(n-1)! \theta}{(\lambda + b)^n}, \quad n \geq 1.$$

We can substitute these quantities into the general EPPF formula in equation (22) of Theorem 17 to obtain

$$\begin{aligned} p(N_1, \dots, N_K) &= \frac{(-1)^{N-K}}{(N-1)!} \int_0^\infty \lambda^{N-1} (\lambda + b)^{-\theta} b^\theta \\ &\quad \cdot \prod_{k=1}^K (-1)^{N_k-1} \frac{(N_k-1)! \theta}{(\lambda + b)^{N_k}} d\lambda \end{aligned}$$

$$\begin{aligned} &= b^\theta \frac{\theta^K}{(N-1)!} \left[\prod_{k=1}^K (N_k - 1)! \right] b^{N-1-N-\theta+1} \\ &\quad \cdot \int_0^\infty x^{N-1} (x+1)^{-N-\theta} dx \quad \text{for } x = \lambda/b \\ &= \frac{\theta^K}{(N-1)!} \left[\prod_{k=1}^K (N_k - 1)! \right] \frac{\Gamma(N)\Gamma(\theta)}{\Gamma(N+\theta)} \\ &= \theta^K \left[\prod_{k=1}^K (N_k - 1)! \right] \frac{1}{\theta(\theta+1)_{N-1 \uparrow 1}}. \end{aligned}$$

The penultimate line follows from the form of the beta prime distribution. The final line is the CRP EPPF from equation (4), as desired. We note in particular that the parameter b does not appear in the final EPPF. \square

Whenever the Laplace exponent of a subordinator is known, Theorem 17 can similarly be applied to quickly find the EPPF of the partition generated by sampling from the normalized subordinator jumps.

To find the distributions of the stick lengths—that is, the partition block frequencies—from the subordinator representation for a partition, we must find the distributions of the normalized subordinator jumps.

As in the feature case, we may enumerate the jumps of a subordinator used for partitioning in the order of their appearance. That is, let ρ_1 be the normalized subordinator jump size corresponding to the cluster of the first data point. Recursively, suppose index n joins a cluster to which none of the indices in $[n-1]$ belong, and suppose there are k clusters among $[n-1]$. Then let ρ_{k+1} be the normalized subordinator jump size corresponding to the cluster containing n .

EXAMPLE 20 (Chinese restaurant process). We continue with the CRP example.

THEOREM 21. *The normalized subordinator jumps (ρ_k) in order of appearance of the gamma subordinator with concentration parameter θ (and arbitrary parameter $b > 0$) have the same distribution as the CRP stick lengths [equation (14) of Example 10 in Section 4].*

PROOF. First, we introduce some notation. Let $\tau = \sum_k \xi_k$, the sum over all of the jumps of the subordinator. Second, let $\tau_k = \tau - \sum_{j=1}^k \xi_k$, the total sum minus the first k elements (in order of appearance). Note that $\tau = \tau_0$. Finally, let $W_k = \tau_k / \tau_{k-1}$ and $V_k = 1 - W_k$. Then a simple telescoping of factors shows that $\rho_k =$

$$V_k \prod_{j=1}^{k-1} (1 - V_j):$$

$$\begin{aligned} V_k \prod_{j=1}^{k-1} (1 - V_j) &= \left(1 - \frac{\tau_k}{\tau_{k-1}}\right) \prod_{j=1}^{k-1} \frac{\tau_j}{\tau_{j-1}} \\ &= \frac{\tau_{k-1} - \tau_k}{\tau_0} = \frac{\xi_k}{\tau} = \rho_k. \end{aligned}$$

It remains to show that the V_k have the desired distribution. To that end, it is easier to work with the W_k . We will find the following lemma (Pitman, 2006) useful.

LEMMA 22. Consider a subordinator with Lévy measure Λ , and suppose τ equals the sum of all jumps of the subordinator. Let ρ be the density of Λ with respect to the Lebesgue measure. And let f be the density of the distribution of τ with respect to the Lebesgue measure. Then

$$\begin{aligned} \mathbb{P}(\tau_0 \in dt_0, \dots, \tau_k \in dt_k) \\ = f(t_k) dt_k \left(\prod_{j=0}^{k-1} \frac{(t_j - t_{j+1}) \rho(t_j - t_{j+1})}{t_j} dt_j \right). \end{aligned}$$

With this lemma in hand, the result follows from a change of variables calculation; we use a bijection between $\{W_1, \dots, W_k, \tau\}$ and $\{\tau_0, \dots, \tau_k\}$ defined by $\tau_k = \tau \prod_{j=1}^k W_j$. The determinant of the Jacobian for the transformation to the former variables from the latter is

$$J = \prod_{j=1}^k \left[\tau \prod_{i=1}^{j-1} W_i \right] = \prod_{j=0}^{k-1} \tau_j(\tau, W_1, \dots, W_k).$$

In the derivation that follows, we start by expressing results in terms of the τ_j terms with the dependence on $\{\tau, W_1, \dots, W_k\}$ suppressed to avoid notational clutter, for example, $J = \prod_{j=0}^{k-1} \tau_j$. At the end, we will evaluate the τ_j terms as functions of $\{\tau, W_1, \dots, W_k\}$.

For now, then, we have

$$\begin{aligned} \mathbb{P}(W_1 \in dw_1, \dots, W_k \in dw_k, \tau \in dt_0) \\ = \mathbb{P}(\tau_0 \in dt_0, \dots, \tau_k \in dt_k) \cdot J \\ = f(t_k) dt_k \left(\prod_{j=0}^{k-1} (t_j - t_{j+1}) \rho(t_j - t_{j+1}) \right). \end{aligned}$$

In the case of the gamma process, we can read $\rho(\xi) = \theta \xi^{-1} e^{-b\xi}$ from equation (23). The function f is determined by ρ and in this case (Pitman, 2006),

$$f(t) = \text{Ga}(t|\theta, b) = b^\theta \Gamma(\theta)^{-1} t^{\theta-1} e^{-bt}.$$

So

$$\begin{aligned} \mathbb{P}(W_1 \in dw_1, \dots, W_k \in dw_k, \tau \in dt_0) \\ \propto t_k^{\theta-1} e^{-bt_0} = t_0^{\theta-1} e^{-bt_0} \prod_{j=1}^k w_j^{\theta-1}. \end{aligned}$$

Since the distribution factorizes, the $\{W_k\}$ are independent of each other and of τ . Second, we can read off the distributional kernel of each W_k to establish $W_k \stackrel{\text{i.i.d.}}{\sim} \text{Beta}(\theta, 1)$, from whence it follows that $V_k \stackrel{\text{i.i.d.}}{\sim} \text{Beta}(1, \theta)$. \square

5.1 Inference

In some sense, we skipped ahead in describing inference in Sections 3.4 and 4.1. There, we made use of the fact that random labels for partitions and features imply exchangeability of the data partition block assignments (Z_n) and data feature assignments (Y_n) . In the discussion above, we study the object that associates random uniformly distributed labels with each partition or feature. Assuming the labels come from a uniform distribution rather than a general continuous distribution is a special case of the discussion in Section 3.4, and we defer the general case to the next section (Section 6).

We have seen above that it is particularly straightforward to obtain an EPPF or EFPPF formulation, which yields Gibbs sampling steps as described in Section 3.4, when the stick lengths are generated according to a normalized Poisson process in the partition case or a Poisson process in the feature case. Examples 15 and 18 illustrate how to find such exchangeable probability functions. Further, we have already seen the usefulness of the stick representation in inference, and Examples 15 and 20 illustrate how stick-length distributions may be recovered from the subordinator framework.

6. COMPLETELY RANDOM MEASURES

In our discussion of subordinators, the jump sizes of the subordinator corresponded to the feature frequencies or unnormalized partition frequencies and were the quantities of interest. By contrast, the locations of the jumps mainly served as convenient labels for the frequencies. These locations were chosen uniformly at random from the unit interval. This choice guaranteed the a.s. uniqueness of the labels and the exchangeability of the sequence of index assignments: (Z_n) in the clustering case or (Y_n) in the feature case.

However, a labeling retains exchangeability and a.s. uniqueness as long as the labels are chosen i.i.d. from any continuous distribution (not just the uniform distribution). Moreover, in typical applications, we wish to

associate some parameter, often referred to as a “random effect,” with each partition block or feature. In the partition case, we usually model the n th data point X_n as being generated according to some likelihood depending on the parameter corresponding to its block assignment. For example, an individual animal’s height and weight, X_n , varies randomly around the height and weight of its species, Z_n . Likewise, in the feature case, we typically model the observed data point X_n as being generated according to some likelihood depending on the collection of parameters corresponding to its collection of feature block assignments [cf. equation (11)]. For example, the book-buying pattern of an online consumer, X_n , varies with some noise based on the topics this person likes to read about: Y_n is a collection, possibly empty, of such topics.

In these cases, it can be useful to suppose that the partition block labels (or feature labels) ϕ_k are not necessarily \mathbb{R}_+ -valued but rather are generated i.i.d. according to some continuous distribution H on a general space Φ . Then, whenever k is the order-of-appearance partition block label of index n , we let $Z_n = \phi_k$. Similarly, whenever k is the order-of-appearance feature label for some feature to which index n belongs, $\phi_k \in Y_n$. Finally, then, we complete the generative model in the partition case by letting $X_n \stackrel{\text{indep}}{\sim} \mathcal{L}(Z_n)$ for some distribution function \mathcal{L} depending on parameter Z_n . And in the feature case, $X_n \stackrel{\text{indep}}{\sim} \mathcal{L}(Y_n)$, where now the distribution function \mathcal{L} depends on the collection of parameters Y_n .

When we take the jump sizes (ξ_k) of a subordinator as the weights of atoms with locations (ϕ_k) drawn i.i.d. according to H as described above, we find ourselves with a *completely random measure* μ :

$$(24) \quad \mu = \sum_{k=1}^{\infty} \xi_k \delta_{\phi_k}.$$

A completely random measure is a random measure μ such that whenever A and A' are disjoint sets, we have that $\mu(A)$ and $\mu(A')$ are independent random variables.

To see that associating these more general atom locations to the jumps of a subordinator yields a completely random measure, note that Theorem 13 tells us that the subordinator jump sizes are generated according to a Poisson point process, with some intensity measure $\nu(d\xi)$. The Marking Theorem for Poisson point processes (Kingman, 1993) in turn yields that the tuples $\{(\xi_k, \phi_k)\}_k$ are generated according to a Poisson point process with intensity measure $\nu(d\xi)H(d\phi)$. By

Kingman (1967), whenever the tuples $\{(\xi_k, \phi_k)\}_k$ are drawn according to a Poisson point process, the measure in equation (24) is completely random.

EXAMPLE 23 (Dirichlet process). We can form a completely random measure from the gamma process subordinator and a random labeling of the partition blocks. Specifically, suppose that the labels come from a continuous measure H . Then we generate a completely random measure G , called a *gamma process* (Ferguson, 1973), in the following way:

$$(25) \quad \nu(d\xi \times d\phi) = \theta \xi^{-1} e^{-b\xi} d\xi \cdot H(d\phi),$$

$$(26) \quad \{(\xi_k, \phi_k)\}_k \sim \text{PPP}(\nu),$$

$$(27) \quad G = \sum_{k=1}^{\infty} \xi_k \delta_{\phi_k}.$$

Here, $\text{PPP}(\nu)$ denotes a draw from a Poisson point process with intensity measure ν . The parameters $\theta > 0$ and $b > 0$ are the same as for the gamma process subordinator. A gamma process draw, along with its generating Poisson point process intensity measure, is illustrated in Figure 8.

The *Dirichlet process* (DP) is the random measure formed by normalizing the gamma process (Ferguson, 1973). Since the Dirichlet process atom weights sum to one, it cannot be completely random. We can write the Dirichlet process D generated from the gamma process G above as

$$\tau = \sum_{k=1}^{\infty} \xi_k,$$

$$\rho_k = \xi_k / \tau,$$

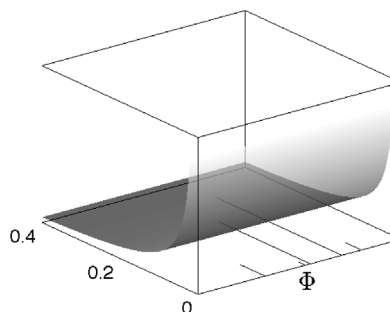


FIG. 8. The gray manifold depicts the Poisson point process intensity measure ν in equation (25) for the choice $\Phi = [0, 1]$ and H the uniform distribution on $[0, 1]$. The endpoints of the line segments are points drawn from the Poisson point process as in equation (26). Taking the positive real-valued coordinate (leftmost axis) as the atom weights, we find the random measure G (a gamma process) on Φ from equation (27) in the bottom plane.

$$D = \sum_{k=1}^{\infty} \rho_k \delta_{\phi_k}.$$

The random variables ρ_k have the same distribution as the Dirichlet process sticks [equation (14)] or normalized gamma process subordinator jump lengths, as we have seen above (Example 18).

Consider sampling points from a Dirichlet process and forming the induced partition of the data indices. Theorem 19 shows us that the distribution of the induced partition is the Chinese restaurant process EPPF.

EXAMPLE 24 (Beta process). We can form a completely random measure from the beta process subordinator and a random labeling of the feature blocks. If the labels are generated i.i.d. from a continuous measure H , then we say the completely random measure B , generated as follows, is called a *beta process*:

$$(28) \quad \nu(d\xi \times d\phi) = \gamma \theta \xi^{-1} (1 - \xi)^{\theta-1} d\xi \cdot H(d\phi),$$

$$(29) \quad \{(\xi_k, \phi_k)\}_k \sim \text{PPP}(\nu),$$

$$(30) \quad B = \sum_{k=1}^{\infty} \xi_k \delta_{\phi_k}.$$

The beta process, along with its generating intensity measure, is depicted in Figure 9. The (ξ_k) have the same distribution as the beta process sticks [equation (16)] or the beta process subordinator jump lengths (Example 15).

Now consider sampling a collection of atom locations according to Bernoulli draws from the atom weights of a beta process and forming the induced feature allocation of the data indices. Theorem 16 shows us that the distribution of the induced feature allocation is given by the Indian buffet process EPPF.

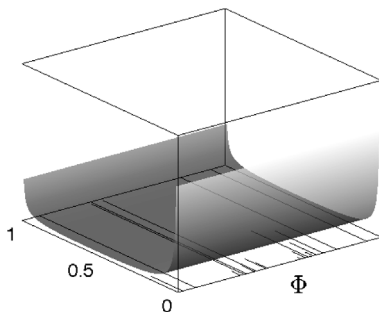


FIG. 9. The gray manifold depicts the Poisson point process intensity measure ν in equation (28) for the choice $\Phi = [0, 1]$ and H the uniform distribution on $[0, 1]$. The endpoints of the line segments are points drawn from the Poisson point process as in equation (29). Taking the $[0, 1]$ -valued coordinate (leftmost axis) as the atom weights, we find the measure B (a beta process) on Φ from equation (30) in the bottom plane.

6.1 Inference

In this section we finally study the full model first outlined in the context of inference of partition and feature structures in Section 3.4. The partition or feature labels described in this section are the same as the block-specific parameters first described in Section 3.4. Since this section focuses on a generalization of the partition or feature labeling scheme beyond the uniform distribution option encoded in subordinators, inference for the atom weights remains unchanged from Sections 3.4, 4.1 and 5.1.

However, we note that, in the course of inferring underlying partition or feature structures, we are often also interested in inferring the parameters of the generative model of the data given the partition block or the feature labels. Conditional on the partition or feature structure, such inference is handled as in a normal hierarchical model with fixed dependencies. Namely, the parameter within a particular block may be inferred from the data points that depend on this block as well as the prior distribution for the parameters. Details for the Dirichlet process example inferred via MCMC sampling are provided by MacEachern (1994), Escobar and West (1995), Neal (2000); Blei and Jordan (2006) work out details for the Dirichlet process using variational methods. In the beta process case, Griffiths and Ghahramani (2006), Teh, Görür and Ghahramani (2007), Thibaux and Jordan (2007) describe MCMC sampling, and Paisley et al. (2010) describe a variational approach.

7. CONCLUSION

In the discussion above we have pursued a progressive augmentation from (1) simple distributions over partitions and feature allocations in the form of exchangeable probability functions to (2) the representation of stick lengths encoding frequencies of the partition block and feature occurrences to (3) subordinators, which associate random \mathbb{R}_+ -valued labels with each partition block or feature, and finally to (4) completely random measures, which associate a general class of labels with the stick lengths and whose labels we generally use as parameters in likelihood models built from the partition or feature allocation representation.

Along the way, we have focused primarily on two vignettes. We have shown, via these successive augmentations, that the Chinese restaurant process specifies the marginal distribution of the induced partition formed from i.i.d. draws from a Dirichlet process, which is

in turn a normalized completely random measure. And we have shown that the Indian buffet process specifies the marginal distribution of the induced feature allocation formed by i.i.d. Bernoulli draws across the weights of a beta process.

There are many extensions of these ideas that lie beyond the scope of this paper. A number of extensions of the CRP and Dirichlet process exist—in either the EPPF form (Pitman, 1996; Blei and Frazier, 2011), the stick-length form (Dunson and Park, 2008) or the random measure form (Pitman and Yor, 1997). Likewise, extensions of the IBP and beta process have been explored (Teh, Görür and Ghahramani, 2007; Paisley et al., 2010; Broderick, Jordan and Pitman, 2012).

More generally, the framework above demonstrates how alternative partition and feature allocation models may be constructed—either by introducing different EPPFs (Pitman, 1996; Gnedin and Pitman, 2006) or EFPFs, different stick-length distributions (Ishwaran and James, 2001) or different random measures (Wolpert and Ickstadt, 2004).

Finally, we note that expanding the set of combinatorial structures with useful Bayesian priors from partitions to the superset of feature allocations suggests that further such structures might be usefully examined. For instance, the *beta negative binomial process* (Broderick et al., 2011; Zhou et al., 2012) provides a prior on a generalization of a feature allocation where we allow the features themselves to be multisets; that is, each index may have nonnegative integer multiplicities of features. Models on trees (Adams, Ghahramani and Jordan, 2010; McCullagh, Pitman and Winkel, 2008; Blei, Griffiths and Jordan, 2010), graphs (Li and McCallum, 2006) and permutations (Pitman, 1996) provide avenues for future exploration. And there likely remain further structures to be fitted out with useful Bayesian priors.

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