Marginalized Samplers for Normalized Random Measure Mixture Models

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Outline

Dirichlet Process Mixture Models

Normalized Random Measures

Posterior and Marginal Characterisations

MCMC Samplers for NRM Mixture Models

Numerical Illustrations

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Numerical Illustrations

Dirichlet Process

- ► Random probability measure $\mu \sim \mathsf{DP}(\alpha, H)$.
- For each partition (A_1, \ldots, A_m) ,

 $(\mu(A_1),\ldots,\mu(A_m)) \sim \text{Dirichlet}(\alpha H(A_1),\ldots,\alpha H(A_m))$

Prior used in Bayesian nonparametric analysis.

 $X_i | \mu \sim \mu$

for *i* = 1, . . . , *n*.

- Large support over space of probability measures.
- Analytically tractable posterior distribution.

Dirichlet Process

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Dirichlet Process Mixture Models

Draws from Dirichlet processes are discrete probability measures,

$$\mu = \sum_{k=1}^{\infty} \mathbf{w}_k \delta_{\phi_k^*}$$

where w_k, ϕ_k^* are random.

Density estimation by convolving with a smooth kernel

$$\int f(\cdot|\phi)\mu(d\phi) = \sum_{k=1}^{\infty} w_k f(\cdot|\phi_k^*)$$

A mixture model with an infinite number of components.

$$x_i | \mu \sim \sum_{k=1}^{\infty} w_k f(\cdot | \phi_k^*)$$

Bayesian Nonparametric Clustering

 $\mu \sim \mathsf{DP}(\alpha, H)$ $\phi_i | \mu \sim \mu$ $\chi_i | \phi_i \sim F(\phi_i)$

- Repeated values among \u03c6_{1:n}
 - Induces a partition π of observations $x_{1:n}$.
 - Each cluster $c \in \pi$ corresponds to a distinct value ϕ_c^* .
 - Leads to a clustering model with a varying number of clusters.
- Properties of model for cluster analysis depends on the properties of the induced random partition π .
- Generalisations of DPs allow for more flexible prior specifications.

MCMC Inference in DP Mixtures

- Conditional samplers [Ishwaran and James 2001, Walker 2007, Kalli and Walker 2006, Papaspiliopoulos and Roberts 2008]
 - Simulate from the joint posterior of μ and $\phi_{1:n}$.
 - Difficulty is in the infinite nature of µ, requiring truncations and retrospective sampling techniques.
 - Easier to understand, parallelizable.
- Marginal samplers [Escobar and West 1995, Bush and MacEachern 1996, Neal 2000]
 - Marginalize out μ , and simulate from posterior for π and $\{\phi_c^*\}$.

```
egin{aligned} &\pi \sim \mathsf{CRP}(lpha) \ &\phi^*_{\mathit{c}} \sim \mathcal{H} \ &x_i | \pi, \{\phi^*_{\mathit{c}}\} \sim \mathcal{F}(\phi^*_{\mathit{c}}) \end{aligned}
```

for $c \in \pi$ for $c \in \pi$ with $i \in c$

Generally better mixing.

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Completely Random Measures

 A completely random measure (CRM) ν is a random measure such that

$\nu(A) \perp \!\!\!\perp \nu(B)$

whenever *A* and *B* are disjoint sets.

The CRM can always be decomposed into 3 components:

$$\nu = \nu_0 + \sum_{j=1}^{\infty} \mathbf{v}_j \delta_{\eta_j^*} + \sum_{k=1}^{\infty} \mathbf{w}_k \delta_{\phi_k^*}$$

- $\nu_0, \{\eta_i^*\}$ are not random,
- $\{v_j\}$ are mutually independent and independent of $\{w_k, \phi_k^*\}$,
- {(w_k, φ^{*}_k)} is drawn from a Poisson process over ℝ₊ × Φ with rate measure ρ(w, φ)dwdφ (the Lévy measure).
- In most modelling applications, only require third component.

Completely Random Measures



The Lévy Measure

~

$$\nu = \sum_{k=1}^{\infty} w_k \delta_{\phi_k^*} \qquad \{(w_k, \phi_k^*)\} \sim \mathsf{Poisson}(\rho)$$

► Homogeneous CRMs have $\rho(w, \phi) = \rho(w)h(\phi)$, implies:

 $\{w_k\} \sim \mathsf{Poisson}(\rho) \qquad \qquad \phi_k^* \sim H$

Want v to have infinitely many atoms:

$$\Rightarrow \int_0^\infty \rho(w) dw = \infty$$

Want v to have finite total mass:

$$\Rightarrow \int_0^\infty (1 - e^{-w}) \rho(w) dw < \infty$$



Normalized Random Measures

Normalizing a CRM gives a normalized random measure (NRM):

$$\mu = \frac{\nu}{\nu(\Phi)}$$

- μ is a random discrete probability measure.
- To study random partition structure, it suffices to assume
 - No fixed measure ($\nu_0 = 0$),
 - No fixed atoms ($v_i = 0$),
 - Homogeneous NRMs $\rho(w, \phi) = \rho(w)h(\phi)$.

Normalized Gamma Process

A gamma process is a CRM with Lévy measure

$$\rho_{\alpha,\tau}(\boldsymbol{w},\phi) = \alpha \boldsymbol{w}^{-1} \boldsymbol{e}^{-\tau \boldsymbol{w}} \boldsymbol{h}(\phi)$$

> The gamma process has gamma marginals:

$$\nu(\mathbf{A}) \sim \operatorname{Gamma}\left(\alpha \int_{\mathbf{A}} \mathbf{h}(\phi) \mathbf{d}\phi, \tau\right)$$

Normalizing a gamma process gives a Dirichlet process, with mass parameter α and base distribution H with density h.

Normalized Stable Process

A stable process is a CRM with Lévy measure

$$\rho_{\sigma}(w) = \frac{\sigma}{\Gamma(1-\sigma)} w^{-\sigma-1}$$

- It has positive stable marginals with index σ .
- Normalizing a stable process, and deriving the induced exchangeable partition process, leads to the following random partition:
 - Customer 1 sits at first table.
 - Subsequent customer n + 1:
 - sits at table *c* with probability $\frac{|c|-\sigma}{p}$,
 - sits at new table with probability $\frac{|\pi|\sigma}{n}$.
- Related to the two-parameter Poisson-Dirichlet process (aka Pitman-Yor process).

Normalized Generalized Gamma Process

There is one NRM that encompasses both DP and normalized stable. It is obtained by normalizing the generalized gamma process:

$$\rho_{\sigma,\alpha,\tau}(w) = \frac{\alpha}{\Gamma(1-\sigma)} w^{-\sigma-1} e^{-\tau w}$$

- It also has power-law properties (like the Pitman-Yor).
- Specializes to DP when $\sigma = 0$.
- Specializes to normalised stable when $\tau = 0, \alpha = \sigma$.
- Specializes to normalized inverse Gaussian process when $\sigma = \frac{1}{2}$.

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NRM Hierarchical Model

for *i* = 1 . . . *n*

What is the posterior distribution?

 $\nu |\phi_{1:n}|$

What is the marginal distribution?

 $\phi_{1:n}$

NRM Hierarchical Model

 $u \sim \operatorname{CRM}(\rho, H)$ $\mu = \nu/\nu(\Phi)$ $\phi_i | \mu \sim \mu \quad \text{for } i = 1 \dots n$

- Let φ^{*}₁,..., φ^{*}_K be the K unique values among φ_{1:n}, with φ^{*}_k occurring n_k times.
- Intuitively,

$$\nu |\phi_{1:n} = \nu^* + \sum_{k=1}^{K} \nu_k \delta_{\phi_k^*}$$
$$p(\phi_{1:n} | \nu) = \frac{\prod_{k=1}^{K} \nu(\{\phi_k^*\})^{n_k}}{\nu(\Phi)^n}$$



Data Augmentation

$$p(\phi_{1:n}|\nu) = \frac{\prod_{k=1}^{K} \nu(\{\phi_k^*\})^{n_k}}{\nu(\Phi)^n}$$
$$= \prod_{k=1}^{K} \nu(\{\phi_k^*\})^{n_k} \int_0^\infty \frac{u^{n-1} e^{-u\nu(\Phi)}}{\Gamma(n)} du$$

Introducing an auxiliary variable u:

$$p(\phi_{1:n}, u|\nu) = \frac{u^{n-1}}{\Gamma(n)} e^{-u(\nu^*(\Phi) + \sum_{k=1}^{K} v_k)} \prod_{k=1}^{K} v_k^{n_k}$$



[James et al. 2009] 19/43

NRM Marginal Characterisation

$$p(\phi_{1:n}, u|\nu) = \frac{u^{n-1}}{\Gamma(n)} e^{-u\nu^*(\Phi)} \prod_{k=1}^{K} v_k^{n_k} e^{-uv_k}$$
$$p(\phi_{1:n}, u) = \mathbb{E}[p(\phi_{1:n}, u|\nu)] = \frac{u^{n-1}}{\Gamma(n)} e^{-\psi(u)} \prod_{k=1}^{K} \kappa(u, n_k) h(\phi_k^*)$$

The Laplace transform of ρ is

$$\psi(u) = -\log \mathbb{E}[e^{-u\nu(\Phi)}] = \int_0^\infty (1 - e^{-uw})\rho(w)dw$$

The *m*th moment of the *u*-exponentially tilted Lévy measure:

$$\kappa(u,m) = \int_0^\infty w^m e^{-uw} \rho(w) dw$$

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NRM Posterior Characterisation

$$\nu |\phi_{1:n}, u = \nu^{*} + \sum_{k=1}^{K} v_{k} \delta_{\phi_{k}^{*}}$$

$$p(\phi_{1:n}, u | \nu) = \frac{u^{n-1} e^{-u\nu^{*}(\Phi)}}{\Gamma(n)} \prod_{k=1}^{K} v_{k}^{n_{k}} e^{-uv_{k}}$$

$$\nu^{*} |\phi_{1:n} \sim \operatorname{CRM}(\rho^{*}, H)$$

$$\rho^{*}(w, \phi) = e^{-uw} \rho(w)$$

$$p(v_{k} | \phi_{1:n}, u) \propto v_{k}^{n_{k}} e^{-uv_{k}} \rho(v_{k})$$

$$\nu | \phi_{1:n}$$

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NRM Mixture Model

 $u \sim \mathsf{CRM}(
ho, H)$ $\mu =
u /
u(\Phi)$

For *i* = 1, . . . , *n*:

 $\phi_i | \mu \sim \mu$ $x_i | \phi_i \sim F(\phi_i)$

- The distinct values among $\phi_{1:n}$ induces a partition π .
- Let ϕ_c^* be the distinct value associated with cluster $c \in \pi$.

Conjugate case

$$p(\pi, u, x_{1:n}) = \frac{u^{n-1}e^{-\psi(u)}}{\Gamma(n)} \prod_{c \in \pi} \left(\kappa(u, |c|) \int_{\Phi} h(\phi_c^*) \prod_{i \in c} f(x_i | \phi_c^*) d\phi_c^* \right)$$

- Gibbs sampling:
 - Marginalize out cluster parameters $\{\phi_c^*\}$.
 - Update u given π using slice sampling.
 - ► Gibbs sample cluster assignment of each item *x_i* in turn:

$$p(i \in c | \pi_{\backslash i}, x_{1:n}) \\ \propto \begin{cases} \frac{\kappa(u, |c|+1)}{\kappa(u, |c|)} \int f(x_i | \phi_c^*) p(\phi_c^* | (x_j)_{j \in c}) d\phi_c^* & \text{for } c \in \pi_{\backslash i}, \\ \kappa(u, 1) \int f(x_i | \phi_c^*) p(\phi_c^*) d\phi_c^* & \text{for } c = \text{new cluster.} \end{cases}$$

Conjugate case

Normalised generalised gamma processes:

$$p(i \in c | \pi_{\backslash i}, x_{1:n}) \\ \propto \begin{cases} (|c| - \sigma) \int f(x_i | \phi_c^*) p(\phi_c^* | (x_j)_{j \in c}) d\phi_c^* & \text{for } c \in \pi_{\backslash i}, \\ \alpha(u + \tau)^{\sigma} \int f(x_i | \phi_c^*) p(\phi_c^*) d\phi_c^* & \text{for } c = \text{new cluster.} \end{cases}$$

Non-conjugate case

• Cannot marginalize out cluster parameters $\{\phi_c^*\}$.

$$p(\pi, u, \{\phi_c^*\}, x_{1:n}) \propto \frac{u^{n-1}e^{-\psi(u)}}{\Gamma(n)} \prod_{c \in \pi} \left(\kappa(u, |c|)h(\phi_c^*) \prod_{i \in c} f(x_i | \phi_c^*)\right)$$

Gibbs sample cluster assignment of item x_i:

$$p(i \in c | \pi_{\setminus i}, x_{1:n}) \propto egin{cases} rac{\kappa(u, |c|+1)}{\kappa(u, |c|)} f(x_i | \phi_c^*) & ext{for } c \in \pi_{\setminus i}, \ \kappa(u, 1) \int f(x_i | \phi_c^*) p(\phi_c^*) d\phi_c^* & ext{for } c = ext{new cluster.} \end{cases}$$

- Integral for new clusters expensive to evaluate.
- ► When singleton cluster is emptied, parameter is discarded.

Neal's Algorithm 8

- Framed as a data augmentation scheme:
 - Introduce *M* "new" clusters, with parameters ψ_k for k = 1, ..., M.

 $\psi_{\mathbf{k}} \sim \mathbf{H}$

- Drawn before each Gibbs update.
- Exists only during the Gibbs update, and discarded afterwards.
- The parameter of an emptied cluster is used as the parameter for one of the new cluster.

$$p(i \in \boldsymbol{c} | \pi_{\backslash i}, x_{1:n}) \propto \begin{cases} \frac{\kappa(\boldsymbol{u}, |\boldsymbol{c}| + 1)}{\kappa(\boldsymbol{u}, |\boldsymbol{c}|)} f(x_i | \phi_{\boldsymbol{c}}^*) & \text{for } \boldsymbol{c} \in \pi_{\backslash i}, \\ \frac{\kappa(\boldsymbol{u}, 1)}{M} f(x_i | \psi_k) & \text{for } \boldsymbol{c} = k \in \{1, \dots, M\}. \end{cases}$$







Xi



Xi







Reuse Algorithm

- Computationally expensive to generate many parameters from base distribution *h*.
- Would like to somehow reuse unused parameters.
- A transdimensional algorithm:
 - Augment state space permanently with *M* new clusters.
 - Reversible jump Metropolis-Hastings updates.

Reuse Algorithm



- Augment state space with *M* new clusters.
- Unassign x_i; if current cluster is a singleton,
 - Replace the parameter of a randomly chosen new cluster with its parameter.
- Reassign cluster assignment of x_i .
- If x_i is assigned to a new cluster,
 - Create a cluster with the parameter,
 - Generate a new parameter from base distribution.
- Acceptance probability always one.

Reuse Algorithm



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NRM Mixture of Normals

- One-dimensional examples:
 - ► Galaxy (*n* = 82)
 - Acidity (*n* = 155)
- Multi-dimensional examples:
 - ► Old Faithful, (*p* = 2, *n* = 272)
 - ▶ Neural spike sorting, (*p* = 6, *n* = 1000, 2000)
- ► Non-conjugate prior over mean and covariance of normals: $m \sim \mathcal{N}(m_0, S_0)$ $\Sigma \sim \mathcal{IW}(\alpha_0, \Sigma_0)$
- Hierarchical prior for $\Sigma_0 \sim \mathcal{IW}(\beta_0, \gamma_0 S_0)$.
- Weakly informative, using prior knowledge of data range.
- In 1D case reduces to prior used in [Richardson and Green 1997].

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Efficiency Evaluation

- 10000 iterations burn-in, 10000 samples collected from 200000 iterations.
- ► Effective sample size of number of clusters *K* using Coda.
- Reports mean ESS and standard error over 10 repeats.
- Compared:
 - Conjugate marginalized sampler
 - Neal's Algorithm 8 marginalized sampler
 - Reuse Algorithm marginalized sampler
 - Slice sampler based on posterior representation
 - Variation on [Griffin and Walker 2011]
 - Truncation required.

Galaxy Dataset



Acidity Dataset



Comparative Results

Galaxy and Acidity datasets (conjugate model)

Sampler	Galaxy		Acidity	
	Runtime (s)	ESS	Runtime (s)	ESS
Cond Slice	239.1 ± 4.2	2004 ± 178	196.5 ± 1.0	910 ± 142
Marg (<i>C</i> = 1)	215.7 ± 1.4	$\textbf{7809} \pm \textbf{87}$	395.5 ± 1.7	$\textbf{5236} \pm \textbf{181}$
Cond Slice	133.0 ± 3.2	1594 ± 117	77.4 ± 0.7	1099 ± 49
Marg Neal 8 (C=1)	74.4 ± 0.6	5815 ± 145	133.3 ± 1.8	4175 ± 85
Marg Neal 8 (C=2)	87.9 ± 0.6	$\textbf{6292} \pm \textbf{94}$	163.8 ± 1.5	4052 ± 158
Marg Neal 8 (C=3)	101.9 ± 0.7	$\textbf{6320} \pm \textbf{137}$	188.2 ± 1.1	$\textbf{4241} \pm \textbf{99}$
Marg Neal 8 (C=4)	115.9 ± 0.6	$\textbf{6283} \pm \textbf{86}$	216.6 ± 1.7	4266 ± 122
Marg Neal 8 (C=5)	130.0 ± 0.6	6491 ± 203	243.8 ± 2.0	4453 ± 123
Marg Reuse (C=1)	64.3 ± 0.3	4451 ± 79	114.6 ± 2.0	3751 ± 65
Marg Reuse (C=2)	67.6 ± 0.5	5554 ± 112	123.1 ± 1.9	4475 ± 110
Marg Reuse (C=3)	71.3 ± 0.5	5922 ± 157	128.2 ± 2.2	4439 ± 158
Marg Reuse (C=4)	74.9 ± 0.5	6001 ± 101	140.1 ± 1.6	4543 ± 108
Marg Reuse (C=5)	78.7 ± 0.6	6131 ± 124	147.7 ± 1.5	$\textbf{4585} \pm \textbf{116}$

Comparative Results

Galaxy and Acidity datasets (non-conjugate model)

Sampler	Galaxy		Acidity	
campion	Runtime (s)	ESS	Runtime (s)	ESS
Cond Slice	75.5 ± 1.2	939 ± 92	50.9 ± 0.5	949 ± 70
Marg Neal 8 (C=1)	65.0 ± 0.5	4313 ± 172	110.9 ± 0.8	4144 ± 64
Marg Neal 8 (C=2)	78.6 ± 0.4	4831 ± 168	139.2 ± 1.8	4290 ± 125
Marg Neal 8 (C=3)	92.5 ± 0.5	4785 ± 97	162.7 ± 0.9	4368 ± 72
Marg Neal 8 (C=4)	106.3 ± 0.5	4849 ± 120	187.6 ± 1.1	4234 ± 142
Marg Neal 8 (C=5)	119.7 ± 0.6	$\textbf{5029} \pm \textbf{89}$	215.4 ± 1.3	4144 ± 213
Marg Reuse (C=1)	55.2 ± 0.5	$\textbf{3830} \pm \textbf{103}$	91.3 ± 0.9	4007 ± 122
Marg Reuse (C=2)	58.7 ± 0.5	$\textbf{4286} \pm \textbf{101}$	98.1 ± 0.9	4192 ± 138
Marg Reuse (C=3)	$\textbf{62.4} \pm \textbf{0.6}$	4478 ± 124	105.1 ± 0.9	4260 ± 136
Marg Reuse (C=4)	$\textbf{66.1} \pm \textbf{0.5}$	4825 ± 63	112.3 ± 1.0	4191 ± 139
Marg Reuse (C=5)	69.8 ± 0.6	4755 ± 141	121.0 ± 1.8	4186 ± 121

Comparative Results

Old Faithful and spike sorting datasets (non-conjugate model)

Sampler	Old Faithful		Spike Sorting	
	Runtime (s)	ESS	Runtime (s)	ESS
Cond Slice	142.6 ± 1.1	574 ± 36	732.6 ± 8.1	17.1 ± 2.3
Marg Reuse (C=1)	$\textbf{208.0} \pm \textbf{1.3}$	2770 ± 209	1120.3 ± 8.8	35.7 ± 2.4
Marg Reuse (C=2)	$\textbf{225.3} \pm \textbf{1.4}$	$\textbf{3236} \pm \textbf{73}$	1164.5 ± 5.4	$\textbf{46.9} \pm \textbf{2.9}$
Marg Reuse (C=3)	241.5 ± 1.3	3148 ± 71	1204.1 ± 7.3	$\textbf{57.0} \pm \textbf{3.9}$
Marg Reuse (C=4)	257.7 ± 1.7	$\textbf{3291} \pm \textbf{145}$	1238.5 ± 7.8	61.4 ± 3.3
Marg Reuse (C=5)	$\textbf{274.8} \pm \textbf{1.7}$	3144 ± 70	1291.8 ± 7.9	69.8 ± 4.9
Marg Reuse (C=10)	356.3 ± 2.5	$\textbf{3080} \pm \textbf{135}$	1513.8 ± 11.9	90.8 ± 5.6
Marg Reuse (C=15)	446.6 ± 4.9	3312 ± 154	1746.3 ± 10.7	95.9 ± 4.2
Marg Reuse (C=20)	550.4 ± 3.5	$\textbf{3336} \pm \textbf{109}$	1944.0 ± 14.7	114.5 ± 8.4

Spike Sorting Dataset



Spike Sorting Dataset



Discussion

- Marginalised samplers for NRMs more efficient than conditional slice samplers.
- Simple algorithms, introducing an additional auxiliary variable *u*.
- Pitman-Yor processes are not normalised random measures.
- Marginalised samplers for all σ-stable Poisson-Kingman mixture models (including Pitman-Yor) (Lomeli et al).
- Motivation for normalised random measures?
 - Power-law properties
 - Dependent normalised random measures

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