# Summary of Short-term Research Objectives

Detian Deng

May 14, 2015

# 1 Model Specification

Let L be a K-dimensional Bernoulli random variable denoting the true state. Consider the general log linear model:

$$f(l;\Theta) = \exp\{\Theta_1^T l + \Theta_2^T u_2 + \ldots + \Theta_K^T u_K - A^*(\Theta)\}\$$

where  $U_k$  is a  $\binom{K}{k} \times 1$  vector of k-way cross-products, k = 1, ..., K, and  $\Theta = (\Theta_1, ..., \Theta_K)$  contains the the natural parameters, which is a  $(2^K - 1) \times 1$  vector.

Model restrictions, let  $\tilde{l} = (l, u_2, \dots, u_K)^T$ , and  $S = \sum_{j=1}^K L_j = s$  has some fixed pmf

$$\pi(s) := P(S = s) = \frac{1}{A(\Theta)} \sum_{\tilde{l}: S = s} \exp\{\Theta^T \tilde{l}\}, \ s = 0, 1, \dots, K$$
 (1)

$$A(\Theta) = \sum_{\tilde{l}: l \in \{0,1\}^K} \exp\{\Theta^T \tilde{l}\}$$
 (2)

### 1.1 Additional Definition of Parameters

Using previous notations, with some properly defined  $\gamma$ , we have:

$$P(L_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) = P(L_i, S_i; \boldsymbol{\pi}, \boldsymbol{\gamma})$$
  
=  $P(L_i | S_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) P(S_i; \boldsymbol{\pi}, \boldsymbol{\gamma})$ 

where  $P(L_i|S_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) = P(L_i|S_i; \boldsymbol{\gamma})$  because  $1(S_i = s)$  is the sufficient statistic for  $\pi_s$ , furthermore  $S_i$  is the sufficient statistic for  $\boldsymbol{\pi}$ . Also  $P(S_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) = \pi_{S_i}$  by definition.

Therefore, we have

$$P(L_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) = P(L_i|S_i; \boldsymbol{\gamma})\pi_{S_i}$$

Then we define the following parameters:

$$\gamma_{j_1,...,j_s} = P(L_{ij_1} = \dots = L_{ij_s} = 1 | S_i = s)$$

$$\gamma = (\gamma_1, \gamma_2, \dots, \gamma_{12}, \dots, \gamma_{1...K})^T, \text{ where } \sum_j \gamma_j = \sum_{j \neq j'} \gamma_{jj'} = \sum_{j \neq j' \neq j''} \gamma_{jj'j''} = \dots = \gamma_{1...K} = 1$$

Therefore  $(\boldsymbol{\pi}, \boldsymbol{\gamma})^T$  is a vector of length  $2^K + K$  with degrees of freedom  $2^K - 1$ .

Let  $J_i = \{j : L_{ij} = 1\}$ . We have,

$$P(L_i; \boldsymbol{\pi}, \boldsymbol{\gamma}) = \gamma_{J_i} \pi_{S_i} \tag{4}$$

The relation between  $(\pi, \gamma)$  and  $\Theta$  is defined by equation (2) together with:

$$\gamma_{J_i} = \frac{\exp(\Theta^T \tilde{l}_i)}{\sum_{l:l^T 1 = S_i} \exp(\Theta^T \tilde{l})}$$
(5)

with  $2^K - 1 - K$  degrees of freedom.

Therefore (1), (3) and (5) together define  $2^K-1$  non-linear equations for  $2^K-1$  unknowns. If there exists a unique root for the above non-linear system, then there is a one-to-one mapping between  $(\pi, \gamma)$  and  $\Theta$ , which provides the re-parameterization.

## 1.2 Find the Re-parameterization

### 1.2.1 Quasi-Newton Method

Numerically solve the system defined by (1), (3) and (5). As the dimension of L grows (K > 6), multiple sets of starting values are needed to reach the solution. Also, solutions to high order  $\Theta$  are subject to larger error.

See code GammaToTheta() in the appendix.

### 1.2.2 Restrict $\Theta$ to QE model

Setting all high order interaction parameter to 0, using only the equations defined by  $pi_0, pi_1, \gamma_1, \ldots, \gamma_{K-1}$  and  $\gamma_{11}, \ldots, \gamma_{K-2,K}$ , which are in total  $\frac{K(K+1)}{2}$  equations, we can solve for  $\Theta$  for larger value of K.

See code GammaToTheta.QE() in the appendix.

# 2 Posterior Distribution

$$\begin{split} P(\mu, \theta^{(2)} | L) \propto & P(L, \mu, \theta^{(2)}) \\ \propto & P(L, \mu, \theta^{(1)}, \theta^{(2)}, \pi) \\ \propto & P(L | \mu, \theta^{(1)}, \theta^{(2)}, \pi) P(\mu, \theta^{(1)}, \theta^{(2)}, \pi) \\ \propto & P(L | \theta^{(1)}, \theta^{(2)}) P(\theta^{(1)}, \theta^{(2)} | \mu, \pi) P(\mu) P(\pi) \\ \propto & QE(L; \theta^{(1)}, \theta^{(2)}) UFR(\theta^{(1)}, \theta^{(2)} | \mu, \pi) N(\text{logit}(\mu), \Sigma) tPois(\pi) \end{split}$$

where QE is the second-order log linear model.

UFR is a Multivariate distribution of  $[\theta^{(1)}, \theta^{(2)}|\mu, \pi]$  subject to non-linear constrains:  $M(\theta^{(1)}, \theta^{(2)}) = \mu$  and  $\Pi(\theta^{(1)}, \theta^{(2)}) = \pi$ , which can be sampled by a two-step procedure. tPois is a truncated conjugate Poisson distribution defined as:

$$\pi \sim \text{Dirichilet(hist(\vec{s}))}$$

$$s \sim \frac{\lambda^s}{s!} e^{-\lambda} / [1 - \sum_{s > K} \frac{\lambda^s}{s!} e^{-\lambda}]$$

# **2.1** On sampling $[\theta^{(1)}, \theta^{(2)}|\mu, \pi]$

### 2.1.1 QE model

For QE model, we have a  $J_1 \times J_2$  design matrix  $\tilde{L}$ , where  $J_1 = 2^K - 1$ ,  $J_2 = \frac{1}{2}K(K+1)$ . Recall that

$$A(\Theta) = \frac{1}{\pi(0)}$$

$$\pi(s) = \frac{1}{A(\Theta)} \sum_{\tilde{l}:S=s} \exp\{\Theta^T \tilde{l}\}, s = 1, \dots, K$$

$$\mu_k = \frac{1}{A(\Theta)} \sum_{\tilde{l}:J=1} \exp\{\Theta^T \tilde{l}\}, k = 1, \dots, K$$
(6)

Define intermediate parameter  $\phi_j = \exp(\theta^T \tilde{l}_j) > 0$ ,  $\theta = (\theta^{(1)}, \theta^{(2)}), j = 1, \dots, J_1$  and two  $K \times J_1$  sub-design matrices B, C, where  $B[k, j] = 1(\tilde{l}_j^T 1 = k), C[k, j] = 1(\tilde{L}[j, k] = 1)$ . Thus (6) and (7) become

$$\vec{\phi} > 0$$

$$B\vec{\phi} = \vec{\pi}/\pi(0)$$

$$C\vec{\phi} = \vec{\mu}/\pi(0)$$

Based on [1][2], we can sample  $\vec{\phi}$  from Uniform distribution subject to the above linear constraints efficiently and robustly. Now we have a over-determined linear system: ( $J_1$  equations with  $J_2$  unknowns)

$$\tilde{L}\theta = \log \vec{\phi}$$

Then we can use Least Square method to solve for  $\theta$ .

### 2.1.2 General model

For General model,  $\tilde{L}$  is  $J_1 \times J_1$ , the intermediate parameters  $\vec{\phi}$  and  $\pi(0)$  fully specify all cell probabilities, thus the posterior distribution becomes

$$P(\mu, \vec{\phi}|L) \propto \mathrm{LL}(L; \vec{\phi}) \mathrm{UFR}(\vec{\phi}|\mu, \pi) \mathrm{N}(\mathrm{logit}(\mu), \Sigma) \mathrm{tPois}(\pi)$$

# Reference

- 1. Smith RL . Efficient Monte-Carlo Procedures for Generating Points Uniformly Distributed over Bounded Regions. Operations Research, 32(6), 12961308 (1984).
- 2. Van den Meersche, Karel, K. E. R. Soetaert, and D. J. Van Oevelen. "xsample (): an R function for sampling linear inverse problems." Journal of Statistical Software 30 (2009).