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Abstract For square contingency tables, Iki, Tahata and Tomizawa (2011) considered the measure to represent the degree of departure from the marginal homogeneity model. Using the first-order term in the Taylor series expansion, the estimated measure with the cell probabilities replaced by the corresponding sample proportions is an approximately unbiased estimator when the sample size is large. The present paper proposes the improved approximate unbiased estimator of the measure which is obtained by using the second-order term in the Taylor series expansion. Also, it shows that the improved estimator approaches to the true measure faster than the original estimator as the sample size becomes larger by the simulation studies.

Keywords: Estimation, marginal homogeneity, marginal odds, square contingency table, taylor series expansion, unbiased estimator.

1 Introduction

Consider an $R \times R$ square contingency table with the same row and column ordinal classifications. Let p_{ij} denote the probability that an observation will fall in the *i*th row and *j*th column of the table (i = 1, ..., R; j = 1, ..., R), and let X and Y denote the row and column variables, respectively. The marginal homogeneity model is defined by

$$p_{i\cdot} = p_{\cdot i} \quad (i = 1, \dots, R),$$

where $p_{i.} = \sum_{t=1}^{R} p_{it}$ and $p_{\cdot i} = \sum_{s=1}^{R} p_{si}$; see Stuart (1955). This indicates that the row marginal distribution is identical to the column marginal distribution. This model is also expressed as

$$F_i^X = F_i^Y$$
 $(i = 1, \dots, R-1),$

where $F_i^X = \sum_{k=1}^i p_k$ and $F_i^Y = \sum_{k=1}^i p_k$. Using the marginal logit, this model can be expressed as $L_i^X = L_i^Y$ (i = 1, ..., R - 1),

where

$$L_i^X = \log\left(\frac{F_i^X}{1 - F_i^X}\right), \quad L_i^Y = \log\left(\frac{F_i^Y}{1 - F_i^Y}\right).$$

This states that the log odds that X is i or below instead of i + 1 or above is equal to the log odds that Y is i or below instead of i + 1 or above for i = 1, ..., R - 1. Further, the marginal homogeneity model is expressed as

$$H_{1(i)} = H_{2(i)}$$
 $(i = 1, \dots, R-1)$

where

$$H_{1(i)} = \sum_{s=1}^{i} \sum_{t=i+1}^{R} p_{s.} p_{.t} = F_i^X (1 - F_i^Y),$$

$$H_{2(i)} = \sum_{s=i+1}^{R} \sum_{t=1}^{i} p_{s.} p_{.t} = (1 - F_i^X) F_i^Y.$$

This indicates that the probability that the row variable X selected at random from the row marginal distribution is in category i or below and the column variable Y selected independently at random from the column marginal distribution is in category i + 1 or above is equal to the probability that such X is in category i + 1 or above and such Y is in category i or below.

Since the marginal homogeneity model indicates that $\{H_{1(i)}\}\$ are equal to corresponding $\{H_{2(i)}\}\$, when the marginal homogeneity model does not hold, we are interested in a measure for seeing how far the probabilities $\{H_{1(i)}\}\$ and $\{H_{2(i)}\}\$ are distant from marginal homogeneity. Iki et al. (2011) considered the measure $\Phi^{(\lambda)}$ to represent the degree of departure from marginal homogeneity for the ordinal data, which is expressed by using the power-divergence (Read and Cressie, 1988, p. 15) or the Patil and Taillie's (1982) diversity index, and as a function of $\{H_{1(i)}\}\$ and $\{H_{2(i)}\}\$. Assuming that $\{H_{1(i)} + H_{2(i)} > 0\}$, let

$$\Delta = \sum_{m=1}^{R-1} \left(H_{1(m)} + H_{2(m)} \right),$$

and let

$$H_{1(i)}^{*} = \frac{H_{1(i)}}{\Delta}, \quad H_{2(i)}^{*} = \frac{H_{2(i)}}{\Delta}, \quad Q_{i}^{*} = \frac{1}{2}(H_{1(i)}^{*} + H_{2(i)}^{*}),$$
$$H_{1(i)}^{c} = \frac{H_{1(i)}}{H_{1(i)} + H_{2(i)}}, \quad H_{2(i)}^{c} = \frac{H_{2(i)}}{H_{1(i)} + H_{2(i)}} \quad (i = 1, \dots, R-1).$$

For $\lambda > -1$, the measure of departure from the marginal homogeneity model considered by Iki et al. (2011), is defined by

$$\begin{split} \Phi^{(\lambda)} &= \frac{1}{2^{\lambda} - 1} \sum_{i=1}^{R-1} \left[H_{1(i)}^{*} \left\{ \left(\frac{H_{1(i)}^{*}}{Q_{i}^{*}} \right)^{\lambda} - 1 \right\} + H_{2(i)}^{*} \left\{ \left(\frac{H_{2(i)}^{*}}{Q_{i}^{*}} \right)^{\lambda} - 1 \right\} \right] \\ &= 1 - \frac{2^{\lambda}}{2^{\lambda} - 1} \sum_{i=1}^{R-1} (H_{1(i)}^{*} + H_{2(i)}^{*}) \left[1 - (H_{1(i)}^{c})^{\lambda+1} - (H_{2(i)}^{c})^{\lambda+1} \right], \end{split}$$

and the value at $\lambda = 0$ is taken to be the limit as $\lambda \to 0$. The measure $\Phi^{(\lambda)}$ must lie between 0 and 1, and it would be useful for comparing the degrees of departure from marginal homogeneity toward the maximum departure in several tables.

Using the first-order term in the Taylor series expansion, the estimated measure with the cell probabilities replaced by the corresponding sample proportions is an approximately unbiased estimator when the sample size is large. Using the second-order term, Tahata et al. (2014) proposed the refined estimators of measures for marginal homogeneity proposed by Tomizawa and Makii (2001) and Tomizawa et al. (2003). So we are now interested in proposing the improved approximate unbiased estimator of $\Phi^{(\lambda)}$.

The purpose of the present paper is to propose the improved approximate unbiased estimator of $\Phi^{(\lambda)}$. Section 2 gives such a estimator. Section 3 shows that the proposed estimator works well in many cases by the simulation studies.

2 Improved Approximate Unbiased Estimator

Assume that the observed frequencies $\{n_{ij}\}$ have a multinomial distribution. Let p be the $R^2 \times 1$ probabilities vector

$$p = (p_{11}, p_{12}, \dots, p_{1R}, p_{21}, p_{22}, \dots, p_{2R}, \dots, p_{R1}, p_{R2}, \dots, p_{RR})^t,$$

where AgtAhmeans transpose. Also let $\{\hat{p}_{ij}\}$ be the sample proportion, where $\hat{p}_{ij} = n_{ij}/n$ with $n = \sum \sum n_{ij}$ and let \hat{p} be the $R^2 \times 1$ vector in the similar way. We assume that g has a nonzero differential at p, i.e., that g has the following expansion as $\hat{p} \to p$:

$$g(\hat{p}) = g(p) + \left[\frac{\partial g(p)}{\partial p^t}\right](\hat{p} - p) + o(\|\hat{p} - p\|).$$

where $[\partial g(p)/\partial p^t]$ denotes $[\partial g(\hat{p})/\partial \hat{p}^t]$ evaluated at $\hat{p} = p$. For the details, see e.g., Agresti (2013, p. 589) and Bishop et al. (1975, p. 486). For large n, we can see from above equation that $g(\hat{p})$ is an approximate unbiased estimator of g(p) because mean of \hat{p} equals p. Similarly, the sample version of $\Phi^{(\lambda)}$, i.e., $\hat{\Phi}^{(\lambda)}$ is given by $\Phi^{(\lambda)}$ with $\{p_{ij}\}$ replaced by $\{\hat{p}_{ij}\}$, is an asymptotically unbiased estimator of $\Phi^{(\lambda)}$ when the sample size n is large.

Assuming that g has a second differential at p, $g(\hat{p})$ has the following expansion as $\hat{p} \to p$:

$$g(\hat{p}) = g(p) + \left[\frac{\partial g(p)}{\partial p^t}\right](\hat{p} - p) + \frac{1}{2}(\hat{p} - p)^t \left[\frac{\partial^2 g(p)}{\partial p \partial p^t}\right](\hat{p} - p) + o(\|\hat{p} - p\|^2),$$

where $[\partial^2 g(p)/\partial p \partial p^t]$ denotes $[\partial^2 g(\hat{p})/\partial \hat{p} \partial \hat{p}^t]$ evaluated at $\hat{p} = p$. Therefore when the sample size n is large, the mean of $g(\hat{p})$, i.e., $E(g(\hat{p}))$, is approximately equal to

$$g(p) + \frac{1}{2n} tr\left(\left[\frac{\partial^2 g(p)}{\partial p \partial p^t}\right] (D(p) - pp^t)\right),$$

where D(p) denotes the $R^2 \times R^2$ diagonal matrix with the *i*th element of *p* as the *i*th diagonal element, because $\operatorname{Var}(\hat{p}) = \frac{1}{n}(D(p) - pp^t)$. Thus the mean of

$$g(\hat{p}) - \frac{1}{2n} tr\left(\left[\frac{\partial^2 g(p)}{\partial p \partial p^t}\right] (D(p) - pp^t)\right)$$

is approximately equal to g(p), and it would approach g(p) faster than $g(\hat{p})$ as the sample size n becomes larger. However, since the second term is unknown, the improved estimator of g(p) is given as follows:

$$g(\hat{p}) - \frac{1}{2n} tr\left(\left[\frac{\partial^2 g(\hat{p})}{\partial p \partial p^t}\right] (D(\hat{p}) - \hat{p}\hat{p}^t)\right),\,$$

where $[\partial^2 g(\hat{p})/\partial p \partial p^t]$ is given by $[\partial^2 g(p)/\partial p \partial p^t]$ with $\{p_{ij}\}$ replaced by $\{\hat{p}_{ij}\}$ and $D(\hat{p})$ denotes D(p) with $\{p_{ij}\}$ replaced by $\{\hat{p}_{ij}\}$.

We now propose the improved estimator of the true measure $\Phi^{(\lambda)}$ as follows:

$$\hat{\varPhi}^{(\lambda)*} = \hat{\varPhi}^{(\lambda)} - \frac{1}{2n} tr\left(\left[\frac{\partial^2 \hat{\varPhi}^{(\lambda)}}{\partial \hat{p} \partial \hat{p}^t}\right] (D(\hat{p}) - \hat{p}\hat{p}^t)\right),$$

where $[\partial^2 \hat{\Phi}^{(\lambda)} / \partial \hat{p} \partial \hat{p}^t]$ is given by $[\partial^2 \Phi^{(\lambda)} / \partial p \partial p^t]$ with $\{p_{ij}\}$ replaced by $\{\hat{p}_{ij}\}$. Then, since $tr[\partial^2 \Phi^{(\lambda)} / \partial p \partial p^t]pp^t = 0$, we note that

$$tr\left(\left[\frac{\partial^2 \Phi^{(\lambda)}}{\partial p \partial p^t}\right] (D(p) - pp^t)\right) = \sum_{k=1}^R \sum_{l=1}^R \frac{\partial^2 \Phi^{(\lambda)}}{\partial p_{kl}^2} p_{kl}$$

where

$$\frac{\partial^2 \Phi^{(\lambda)}}{\partial p_{kl}^2} = \frac{1 - \Phi^{(\lambda)}}{\Delta} \left(K_{2(kl)} - \frac{2}{\Delta} \left(K_{1(kl)} \right)^2 \right) + \frac{1}{\Delta^2} \left\{ 2K_{1(kl)} L_{1(kl)}^{(\lambda)} - \Delta L_{2(kl)}^{(\lambda)} \right\},$$

$$K_{1(kl)} = \sum_{m=1}^{R-1} \left(W_{1(kl)}^{(m)} + W_{2(kl)}^{(m)} \right),$$

$$K_{2(kl)} = \sum_{m=1}^{R-1} \left(W_{3(kl)}^{(m)} + W_{4(kl)}^{(m)} \right),$$

for $\lambda \neq 0$,

$$\begin{split} L_{1(kl)}^{(\lambda)} &= \frac{2^{\lambda}}{2^{\lambda} - 1} \sum_{i=1}^{R-1} \bigg[W_{1(kl)}^{(i)} \left\{ 1 - \left(H_{1(i)}^{c}\right)^{\lambda} - \lambda H_{2(i)}^{c} \left(\left(H_{1(i)}^{c}\right)^{\lambda} - \left(H_{2(i)}^{c}\right)^{\lambda} \right) \right\} \\ &\quad + W_{2(kl)}^{(i)} \left\{ 1 - \left(H_{2(i)}^{c}\right)^{\lambda} - \lambda H_{1(i)}^{c} \left(\left(H_{2(i)}^{c}\right)^{\lambda} - \left(H_{1(i)}^{c}\right)^{\lambda} \right) \right\} \bigg], \\ L_{2(kl)}^{(\lambda)} &= \frac{2^{\lambda}}{2^{\lambda} - 1} \sum_{i=1}^{R-1} \bigg[W_{3(kl)}^{(i)} \left\{ 1 - \left(H_{1(i)}^{c}\right)^{\lambda} - \lambda H_{2(i)}^{c} \left(\left(H_{1(i)}^{c}\right)^{\lambda} - \left(H_{2(i)}^{c}\right)^{\lambda} \right) \right\} \right] \\ &\quad + W_{4(kl)}^{(i)} \left\{ 1 - \left(H_{2(i)}^{c}\right)^{\lambda} - \lambda H_{1(i)}^{c} \left(\left(H_{2(i)}^{c}\right)^{\lambda} - \left(H_{1(i)}^{c}\right)^{\lambda} \right) \right\} \\ &\quad - \frac{\lambda(1 + \lambda)}{H_{1(i)} + H_{2(i)}} \left(\left(H_{1(i)}^{c}\right)^{\lambda-1} + \left(H_{2(i)}^{c}\right)^{\lambda-1} \right) \left(W_{1(kl)}^{(i)} H_{2(i)}^{c} - W_{2(kl)}^{(i)} H_{1(i)}^{c} \right)^{2} \bigg], \end{split}$$

and for $\lambda = 0$,

$$\begin{split} L_{1(kl)}^{(0)} &= \frac{1}{\log 2} \sum_{i=1}^{R-1} \left(-W_{1(kl)}^{(i)} \log H_{1(i)}^{c} - W_{2(kl)}^{(i)} \log H_{2(i)}^{c} \right), \\ L_{2(kl)}^{(0)} &= \frac{1}{\log 2} \sum_{i=1}^{R-1} \bigg\{ -W_{3(kl)}^{(i)} \log H_{1(i)}^{c} - W_{4(kl)}^{(i)} \log H_{2(i)}^{c} \\ &- \frac{1}{H_{1(i)}} W_{1(kl)}^{(i)} \left(W_{1(kl)}^{(i)} H_{2(i)}^{c} - W_{2(kl)}^{(i)} H_{1(i)}^{c} \right) \\ &- \frac{1}{H_{2(i)}} W_{2(kl)}^{(i)} \left(W_{2(kl)}^{(i)} H_{1(i)}^{c} - W_{1(kl)}^{(i)} H_{2(i)}^{c} \right) \bigg\}, \end{split}$$

with

$$W_{1(kl)}^{(m)} = \sum_{a=1}^{m} \sum_{b=m+1}^{R} \left(I_{(a=k)} p_{\cdot b} + p_{a \cdot} I_{(b=l)} \right),$$

$$W_{2(kl)}^{(m)} = \sum_{a=m+1}^{R} \sum_{b=1}^{m} \left(I_{(a=k)} p_{\cdot b} + p_{a \cdot} I_{(b=l)} \right),$$

$$W_{3(kl)}^{(m)} = \sum_{a=1}^{m} \sum_{b=m+1}^{R} \left(I_{(a=k)} I_{(b=l)} + I_{(a=k)} I_{(b=l)} \right),$$

$$W_{4(kl)}^{(m)} = \sum_{a=m+1}^{R} \sum_{b=1}^{m} \left(I_{(a=k)} I_{(b=l)} + I_{(a=k)} I_{(b=l)} \right),$$

and where $I_{(\cdot)}$ is the indicator function, $I_{(\cdot)} = 1$ if true, 0 if not. Therefore, the improved estimator $\hat{\Phi}^{(\lambda)*}$ is also expressed as follows:

$$\hat{\varPhi}^{(\lambda)*} = \hat{\varPhi}^{(\lambda)} - \frac{1}{2n} \sum_{k=1}^{R} \sum_{l=1}^{R} \frac{\partial^2 \hat{\varPhi}^{(\lambda)}}{\partial \hat{p}_{kl}^2} \hat{p}_{kl}.$$

3 Simulation Studies

By the simulation studies, we calculate the values of estimated measures $\hat{\Phi}^{(\lambda)}$ and $\hat{\Phi}^{(\lambda)*}$ from the observed frequencies of sample size n = 30, 40, 50, 100, 500 and 1000, which are obtained from the true probability distribution (see Tables 1a to 6a). We shall compare the mean of the values of $\hat{\Phi}^{(\lambda)}$ and $\hat{\Phi}^{(\lambda)*}$ obtained by 1000 times simulations, for each sample size. The results of simulations are given in Tables 1c to 6c.

Tables 1a, 3a and 5a have a characteristic that the sum of the probabilities of main-diagonal cells is very small ($p_{ii} = 0.020$ for i = 1, 2, 3, 4) and Tables 2a, 4a and 6a have a characteristic that the sum of the probabilities of main-diagonal cells is large ($p_{ii} = 0.100$ for i = 1, 2, 3, 4). Also the true values of measures for Tables 1a and 2a are small, while those for Tables 3a and 4a are medium, and those for Tables 5a and 6a are large, respectively.

We can see that the improved estimator $\hat{\Phi}^{(\lambda)*}$ approaches the true value $\Phi^{(\lambda)}$ faster than the original estimator $\hat{\Phi}^{(\lambda)}$ when $\lambda \geq 1$ from Tables 1c to 6c. Especially, we can see great improvement when sample size is small.

Table 1. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 1a.

(a)					(c)			
	(1)	(2)	(3)	(4)	λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
(1)	0.020	0.101	0.085	0.038	1.0	30	0.1940	0.1286
(2)	0.066	0.020	0.123	0.140		40	0.1660	0.1144
(3)	0.042	0.063	0.020	0.110		50	0.1581	0.1162
(4)	0.040	0.051	0.061	0.020		100	0.1330	0.1112
						500	0.1146	0.1101
	(b)				1000	0.1138	0.1115
) <u></u>	(λ)		3.0	30	0.1677	0.1028
			118			40	0.1533	0.1041
			978			50	0.1415	0.1013
		0.0 0.0	910			100	0.1185	0.0978
						500	0.1001	0.0959
						1000	0.1005	0.0984

Table 2. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 2a.

(a)					(c)			
	(1)	(2)	(3)	(4)	λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
(1)	0.100	0.052	0.068	0.110	1.0	30	0.1614	0.1153
(2)	0.044	0.100	0.054	0.058		40	0.1492	0.1140
(3)	0.038	0.042	0.100	0.052		50	0.1431	0.1149
(4)	0.020	0.020	0.042	0.100		100	0.1318	0.1175
						500	0.1154	0.1125
	(b)				1000	0.1141	0.1127
		$\frac{\lambda}{\lambda} = \Phi^{(1)}$	(λ)		3.0	30	0.1504	0.1062
			129			40	0.1302	0.0960
			129 986			50	0.1234	0.0961
		5.0 0.0	980			100	0.1118	0.0979
						500	0.1016	0.0988
						1000	0.1014	0.1000

4 Concluding Remarks

The present paper has proposed the improved approximate unbiased estimator $\hat{\Phi}^{(\lambda)*}$ of the true measure $\Phi^{(\lambda)}$ proposed by Iki et al. (2011), however, Tahata et al. (2014) proposed the refined estimators of measures proposed by Tomizawa and Makii (2001) and Tomizawa et al. (2003).

76	

(a)					(c)			
	(1)	(2)	(3)	(4)	λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
(1)	0.020	0.096	0.100	0.068	1.0	30	0.4669	0.4371
(2)	0.041	0.020	0.163	0.170		40	0.4501	0.4272
(3)	0.027	0.023	0.020	0.160		50	0.4556	0.4381
(4)	0.040	0.011	0.021	0.020		100	0.4515	0.4431
						500	0.4410	0.4393
	(b)				1000	0.4398	0.4390
		$\frac{1}{\lambda} \Phi^{(1)}$	(λ)		3.0	30	0.4375	0.4035
		$\frac{1}{1.0}$ $\frac{\Psi}{0.4}$				40	0.4319	0.4065
						50	0.4250	0.4044
		3.0 0.4	066			100	0.4193	0.4092
						500	0.4091	0.4070
						1000	0.4049	0.4039

Table 3. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 3a.

Table 4. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 4a.

(a)					(c)			
	(1)	(2)	(3)	(4)	λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
(1)	0.100	0.052	0.128	0.135	1.0	30	0.4670	0.4465
(2)	0.014	0.100	0.079	0.088		40	0.4571	0.4416
(3)	0.008	0.012	0.100	0.062		50	0.4562	0.4439
(4)	0.005	0.005	0.012	0.100		100	0.4490	0.4428
						500	0.4442	0.4429
	(b)				1000	0.4438	0.4432
		$\frac{\partial}{\partial t} \Phi^{(i)}$	(λ)		3.0	30	0.4386	0.4141
		-	430			40	0.4186	0.3997
		0.4				50	0.4194	0.4041
		0.0 0.4	079			100	0.4122	0.4045
						500	0.4102	0.4086
						1000	0.4089	0.4081

Table 5. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 5a.

(a)				
	(1)	(2)	(3)	(4)
(1)	0.020	0.181	0.105	0.202
(2)	0.006	0.020	0.133	0.110
(3)	0.002	0.013	0.020	0.143
(4)	0.003	0.011	0.011	0.020

(b)	
λ	$arPhi^{(\lambda)}$
1.0	0.7797
3.0	0.7564

(c)			
λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
1.0	30	0.7830	0.7787
	40	0.7814	0.7786
	50	0.7814	0.7792
	100	0.7817	0.7808
	500	0.7799	0.7798
	1000	0.7803	0.7802
3.0	30	0.7693	0.7623
	40	0.7590	0.7538
	50	0.7580	0.7539
	100	0.7573	0.7554
	500	0.7562	0.7559
	1000	0.7563	0.7561

(a)						(c)			
	(1)	(2)	(3)	(4)	_	λ	n	$\hat{\varPhi}^{(\lambda)}$	$\hat{\varPhi}^{(\lambda)*}$
(1)	0.100	0.002	0.008	0.559	_	1.0	30	0.7647	0.7638
(2)	0.004	0.100	0.004	0.006			40	0.7664	0.7660
(3)	0.008	0.002	0.100	0.002			50	0.7675	0.7673
(4)	0.001	0.002	0.002	0.100			100	0.7664	0.7664
							500	0.7663	0.7766
	(b)					1000	0.7663	0.7663
			(λ)			3.0	30	0.7473	0.7445
			662				40	0.7405	0.7383
							50	0.7417	0.7401
		3.0 0.7	407				100	0.7402	0.7395
							500	0.7411	0.7410
					_		1000	0.7399	0.7399

Table 6. (a) The artificial probabilities $\{p_{ij}\}$, (b) the value of $\Phi^{(\lambda)}$ and (c) the mean of the values of estimated measures obtained by generating 1000 times simulations, with each sample size n, for Table 6a.

From the simulation studies, we conclude that the improved estimator $\hat{\Phi}^{(\lambda)*}$ tends to approach to the true value $\Phi^{(\lambda)}$ faster than the estimator $\hat{\Phi}^{(\lambda)}$ as the sample size *n* becomes larger, when $\lambda \geq 1$.

When $\lambda < 1$, we can calculate the improved estimator $\hat{\Phi}^{(\lambda)*}$ for only the case of $H_{1(i)} > 0$ and $H_{2(i)} > 0$ for i = 1, ..., R - 1, i.e., $p_{1.} > 0$, $p_{R.} > 0$, $p_{.1} > 0$ and $p_{.R} > 0$. On the other hand, the original estimator $\hat{\Phi}^{(\lambda)}$ can be calculated for the case of $H_{1(i)} + H_{2(i)} > 0$ for $i = 1, \ldots, R-1$. In other words, the calculable conditions are different between the improved estimator and the original estimator. Thus, it seems difficult to evaluate whether the improved estimator tends to approach the true value faster than the original estimator by simulation study when $\lambda < 1$. Therefore, we recommend that the proposed estimator should be used for the case of $\lambda \geq 1$. Then this estimator works very well.

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