## COMPARISON OF EXPERIMENTS AND INFORMATION MEASURES<sup>1</sup>

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Let  $\mathcal{E}_X = \{X, S_X; P_\theta, \theta \in \Theta\}$  and  $\mathcal{E}_Y = \{Y, S_Y; Q_\theta, \theta \in \Theta\}$  be two statistical experiments with the same parameter space  $\Theta$ . Some implications of the sufficiency of  $\mathcal{E}_X$  for  $\mathcal{E}_Y$ , according to Blackwell's definition, are given in terms of Kullback-Leibler information and Fisher information matrices. For a scale parameter  $\theta$ , and  $k_1 > k_2 > 0$ , the experiment with parameter  $\theta^{k_1}$  is proved to be sufficient for the experiment with parameter  $\theta^{k_2}$  for a class of distributions including the gamma distribution and the normal distribution with known mean. Some results of Stone are generalized to the class of experiments with both location and scale parameters. A concept of sufficiency is proposed in which  $\mathcal{E}_X$  is more informative than  $\mathcal{E}_Y$  for a fixed prior distribution of  $\theta$  if the expected Bayes risk from  $\mathcal{E}_X$  is not greater than that from  $\mathcal{E}_Y$  for every decision problem involving  $\theta$ . This concept is then used to develop a definition of marginal Bayesian sufficiency in the presence of nuisance parameters.

1. Introduction and summary. Let  $\mathcal{E}_{X} = \{X, S_{X}; P_{\theta}, \theta \in \Theta\}$  denote a statistical experiment in which a random variable or random vector X defined on some sample space  $S_X$  is to be observed, and the distribution  $P_{\theta}$  of X depends on a parameter  $\theta$  whose value is unknown and lies in some parameter space  $\Theta$ . Also, let  $\mathcal{E}_Y = \{Y, S_Y; Q_\theta, \theta \in \Theta\}$  denote another statistical experiment with the same parameter space Θ. Blackwell's (1951) method for comparing two experiments states that the experiment  $\mathcal{E}_X$  is sufficient for the experiment  $\mathcal{E}_Y$  (denoted  $\mathcal{E}_X \geqslant \mathcal{E}_Y$ ) if there exists a stochastic transformation of X to a random variable Z(X) such that, for each  $\theta \in \Theta$ , the random variables Z(X) and Y have identical distributions. It was proved by Blackwell (1953) and Boll (1955) that this method of comparison is equivalent to Bohnenblust, Shapley and Sherman's method for comparing two experiments [see Blackwell (1951)] which states that  $\mathcal{E}_X$  is more informative than  $\mathcal{E}_{\mathbf{r}}$  if for every decision problem involving  $\boldsymbol{\theta}$  and every prior distribution on  $\Theta$ , the expected Bayes risk from  $\mathcal{E}_X$  is not greater than that from &<sub>Y</sub>. LeCam (1964) generalized this notion to a concept of approximate sufficiency or  $\varepsilon$ -deficiency of  $\mathcal{E}_X$  relative to  $\mathcal{E}_Y$ . Torgersen (1970, 1972, 1976) and Hansen and Torgersen (1974) have extended these results and applied them to interesting special classes of experiments. Some other papers on this topic are DeGroot (1962,

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1966), Torgersen (1977), Feldman (1972), and Gerber (1977). We shall now give a summary of the results presented in this paper.

In Section 2, we summarize various implications of the relation  $\mathcal{E}_X \gg \mathcal{E}_Y$  in terms of Kullback-Leibler (K-L) information and Fisher information.

In Section 3, it is assumed that  $\theta$  is a scale parameter in the distribution of Y and that X has this same distribution except that  $\theta$  is replaced by  $\theta^k(k > 1)$ . Let W denote a random variable with distribution identical to that of X with  $\theta = 1$ , and let  $\varphi(t)$  denote the characteristic function of log W. As stated in Lemma 2, if  $\varphi(t)/\varphi(t/k)$  is a characteristic function, then  $\mathcal{E}_X \geq \mathcal{E}_Y$ . It is noted that  $\varphi(t)$  satisfies this condition if and only if  $\varphi(t)$  is a self-decomposable characteristic function [see Lukacs (1970), page 161]. Let  $G_n(a, b)$  denote an experiment in which a random sample of size n is taken from a gamma distribution G(a, b) with parameters a > 0 and b > 0, for which the density function is

(1.1) 
$$g(w) = \frac{w^{a-1}}{b^a \Gamma(a)} \exp(-w/b) \text{ for } w > 0.$$

Also, let  $N_n(\mu, \sigma^2)$  denote an experiment in which a random sample of size n is taken from a normal distribution with mean  $\mu$  and variance  $\sigma^2$ . The result in Lemma 2 is used to prove that for any known numbers a > 0 and b > 0,  $G_n(a, b\theta^k) \ge G_n(a, b\theta)$  for all k > 1, and that  $N_n(0, \sigma^{2k}) \ge N_n(0, \sigma^2)$  for all k > 1.

In Section 4, we extend some of the results obtained by Stone (1961) for a class of experiments with location parameter  $\theta$  to the class of experiments with both a location parameter  $\mu$  and a scale parameter  $\sigma$ . Various results for normal distributions are summarized. We consider an experiment  $\mathcal{E}^*(c)$  involving a random variable having a pdf of the form  $(c/\sigma)f[c(x-\mu)/\sigma]$ . In Theorem 5 we study the relation between the Fisher information matrices for two experiments  $\mathcal{E}^*(c_1)$  and  $\mathcal{E}^*(c_2)$  when f is a symmetric function.

Since two experiments  $\mathcal{E}_X$  and  $\mathcal{E}_Y$  may not be comparable in Blackwell's sense, Feldman (1972) introduced another definition in which  $\mathcal{E}_X$  is more informative than  $\mathcal{E}_Y$  for a fixed decision problem involving  $\theta$  if, for every prior distribution on  $\Theta$ , the expected Bayes risk from  $\mathcal{E}_X$  is not greater than that from  $\mathcal{E}_Y$ . In Section 5, we propose an alternative definition in which  $\mathcal{E}_X$  is more informative than  $\mathcal{E}_Y$  with respect to a fixed prior distribution on  $\Theta$  if, for every decision problem involving  $\theta$ , the expected Bayes risk from  $\mathcal{E}_X$  is not greater than that from  $\mathcal{E}_Y$ . We then apply this concept to problems in which  $\theta$  is a vector with a given prior distribution, and we are interested in decision problems involving only some of the components of  $\theta$ . We present a definition of the marginal Bayesian sufficiency of  $\mathcal{E}_X$  for  $\mathcal{E}_Y$  in this context. Some examples are given to illustrate the usefulness of this concept.

2. Relationships between sufficiency and information. Consider again two arbitrary experiments  $\mathcal{E}_X$  and  $\mathcal{E}_Y$  with the same parameter space  $\Theta$  as defined at the beginning of Section 1. We shall assume that there exist generalized probability density functions (gpdf's)  $p(x|\theta)$  and  $q(y|\theta)$  for the distributions  $P_{\theta}$  and  $Q_{\theta}$ , with

respect to some  $\sigma$ -finite measures  $\mu$  and  $\nu$  respectively. We shall now investigate the implications of the relation  $\mathcal{E}_X \geq \mathcal{E}_Y$  in terms of some well-known information measures. Let  $\Xi$  denote the class of all prior distributions on the parameter space  $\Theta$ . Given two prior distributions  $\xi_1, \xi_2 \in \Xi$ , let  $p_i(x)$  denote the marginal gpdf  $\int_{\Theta} p(x|\theta) d\xi_i(\theta)$ , for i=1,2, and let  $I_X(\xi_1,\xi_2)$  denote the K-L information contained in  $\mathcal{E}_X$  for discriminating between  $p_1(x)$  and  $p_2(x)$ , defined by

(2.1) 
$$I_X(\xi_1, \xi_2) = \int_{S_X} p_1(x) \log \frac{p_1(x)}{p_2(x)} d\mu(x).$$

If  $\xi_1$  assigns probability 1 to a point  $\theta = \theta_0$ , we shall denote  $I_X(\xi_1, \xi_2)$  by  $I_X(\theta_0, \xi_2)$ . The K-L information  $I_Y(\xi_1, \xi_2)$  contained in  $\mathcal{E}_Y$  is defined analogously.

Lindley (1956) has shown that if  $\mathcal{E}_X \geq \mathcal{E}_Y$ , then the Shannon information contained in  $\mathcal{E}_X$  is at least as large as that contained in  $\mathcal{E}_Y$ . That is, if  $\mathcal{E}_X \geq \mathcal{E}_Y$  then

If (2.2) holds for  $\mathcal{E}_X$  and  $\mathcal{E}_Y$ , we shall denote it by  $\mathcal{E}_X \geq_L \mathcal{E}_Y$ . The following stronger version of Lindley's result was proved by Sakaguchi (1964).

LEMMA 1. If  $\mathcal{E}_{x} \geq \mathcal{E}_{y}$ , then

(2.3) 
$$I_{\nu}(\theta, \xi) \ge I_{\nu}(\theta, \xi)$$
 for all  $\theta \in \Theta$  and  $\xi \in \Xi$ .

In fact, this result can be easily extended as follows:

THEOREM 1. If  $\mathcal{E}_X \geq \mathcal{E}_Y$ , then

(2.4) 
$$I_X(\xi_1, \xi_2) \ge I_Y(\xi_1, \xi_2)$$
 for all  $\xi_1, \xi_2 \in \Xi$ .

PROOF. The relation  $\mathcal{E}_X \geqslant \mathcal{E}_Y$  is preserved when the parameter space  $\Theta$  is enlarged to include all convex combinations of the basic distributions  $P_{\theta}$  of X and convex combinations of the distributions  $Q_{\theta}$  of Y. Therefore, the distribution  $\xi$  in (2.4) can be assumed, without loss of generality, to be degenerate. Hence, the theorem follows from Lemma 1.  $\square$ 

The following example shows that Lemma 1 is stronger than Lindley's result.

EXAMPLE 1. Let  $\mathcal{E}(\theta_1, \theta_2, \theta_3)$  denote an experiment in which a coin with unknown probability of heads  $\theta$  is flipped n times and the parameter space  $\Theta$  contains only three points  $0 \le \theta_1 \le \theta_2 \le \theta_3 \le 1$ . Blackwell (1951) remarks that the experiment  $\mathcal{E}_X \equiv \mathcal{E}(0, \frac{1}{2}, 1)$  is not sufficient for the experiment  $\mathcal{E}_Y \equiv \mathcal{E}(0, \frac{1}{2}, \frac{1}{2})$  even though our intuition suggests the contrary. In other words, suppose that there are only three possible states of nature and we have a choice of either (i) observing n flips of a coin  $\mathcal{E}_X$  for which the probability of heads is  $0, \frac{1}{2}$ , or 1 according as  $\theta_1$ ,  $\theta_2$  or  $\theta_3$  is correct, or (ii) observing n flips of a coin  $\mathcal{E}_Y$  for which the probability of heads is 0 if  $\theta_1$  is correct, but is  $\frac{1}{2}$  if either  $\theta_2$  or  $\theta_3$  is correct. It would seem at first glance that  $\mathcal{E}_X$  must always be at least as useful as  $\mathcal{E}_Y$ , but Blackwell pointed out that this conclusion is not correct.

Lindley (1956) showed that  $\mathcal{E}_X \geq_L \mathcal{E}_Y$  for these experiments. However, it can be shown that  $I_X(\theta_2, \xi) \geq I_Y(\theta_2, \xi)$  if and only if  $3\xi_1 + \xi_2 \leq 1$ , where  $\xi_i = \xi(\theta_i)$ . Hence (2.3) does not hold for  $\theta = \theta_2$  and a prior distribution  $\xi$  for which  $3\xi_1 + \xi_2 > 1$ .

Since  $\mathcal{E}_X$  is not sufficient for  $\mathcal{E}_Y$  in this example, there must be a decision problem in which the expected Bayes risk from  $\mathcal{E}_Y$  is less than that from  $\mathcal{E}_X$ . The following simple decision problem has this property. Suppose that the hypothesis  $H_0: \theta = \theta_2$  is to be tested against the alternative  $H_1: \theta \neq \theta_2$  with the usual 0-1 loss function and the prior distribution  $\xi$  defined by  $\xi(\theta_2) = \lambda$  and  $\xi(\theta_1) = \xi(\theta_3) = (1-\lambda)/2$ . If  $\lambda$  satisfies  $2^{n-1} < \lambda/(1-\lambda) < 2^{n-1} + \frac{1}{2}$ , then it can be shown that the Bayes rule for the experiment  $\mathcal{E}_Y$  is to reject  $H_0$  if the number of heads is 0 and the Bayes rule for the experiment  $\mathcal{E}_X$  is to accept  $H_0$  regardless of the outcome. Since the outcome of  $\mathcal{E}_X$  is of no value to the experimenter in this decision problem, it can be shown that the expected Bayes risk from  $\mathcal{E}_X$  is larger than that from  $\mathcal{E}_Y$ .

The converse of Theorem 1 does not necessarily hold. In fact, it follows from Torgersen (1970, Proposition 18) that if  $\Theta$  is finite and, in his terminology,  $\mathcal{E}_X$  is more informative than  $\mathcal{E}_Y$  for testing problems, then (2.4) holds, even if  $\mathcal{E}_X$  is not sufficient for  $\mathcal{E}_Y$ .

The relation  $\mathcal{E}_X \geqslant \mathcal{E}_Y$  implies a similar ordering in terms of Fisher information. Let  $\boldsymbol{\theta} = (\theta_1, \dots, \theta_k)$  and suppose that  $\Theta$  is an open subset of  $R^k$ . Let  $\mathbf{i}_X(\boldsymbol{\theta})$  and  $\mathbf{i}_Y(\boldsymbol{\theta})$  denote the  $k \times k$  Fisher information matrices for the experiments  $\mathcal{E}_X$  and  $\mathcal{E}_Y$  respectively, under the standard regularity conditions such as those given in Kullback ((1968), pages 26-27). We shall use the notation  $\mathcal{E}_X \geqslant_F \mathcal{E}_Y$  whenever  $\mathbf{i}_X(\boldsymbol{\theta}) - \mathbf{i}_Y(\boldsymbol{\theta})$  is nonnegative definite. The next theorem essentially follows from the development given in Kullback ((1968), pages 26-28). A sketch of the proof of this result is also given by Barndorff-Nielsen in his discussion of Torgersen (1976). For k = 1, the result had been proved by Stone (1961).

THEOREM 2. If 
$$\mathcal{E}_X \geq \mathcal{E}_Y$$
, then  $\mathcal{E}_X \geq_F \mathcal{E}_Y$ .

REMARK 1. It should be noted that if  $\mathbf{i}_X(\theta) - \mathbf{i}_Y(\theta)$  is nonnegative definite and  $\mathbf{i}_Y(\theta)$  is nonnegative definite, then  $|\mathbf{i}_X(\theta)| \ge |\mathbf{i}_Y(\theta)|$  (see Rao (1973), page 70). In other words, the generalized Fisher information in  $\mathcal{E}_X$  is at least as large as that in  $\mathcal{E}_Y$ .

REMARK 2. A counterexample which shows that the converse of Theorem 2 does not necessarily hold is given by Hansen and Torgersen (1974).

3. Comparison of normal experiments with known mean and unknown variance. In this section we shall consider experiments in which a random sample can be taken from a normal distribution for which the mean is known and the variance is unknown. Without loss of generality we shall assume that the known value of the mean is 0.

To begin we note that  $N_n(0, \sigma^2) \ge N_n(0, \sigma^2 + k^2)$  where k is a given constant. To see this, suppose that the random variable X is distributed as  $N(0, \sigma^2)$ , the random variable Y is distributed as  $N(0, \sigma^2 + k^2)$ , and the random variable W is independent of X and has the distribution  $N(0, k^2)$ . Then X + W has the same distribution as Y for every possible value of  $\sigma^2$ . Hence,  $N_1(0, \sigma^2) \ge N_1(0, \sigma^2 + k^2)$ . However, it is well known that if  $\mathcal{E}_X \ge \mathcal{E}_Y$  when only one observation is taken in each experiment, then this same relation holds when a random sample of n observations is taken from each experiment [See Blackwell (1951)]. It now follows that  $N_n(0, \sigma^2) \ge N_n(0, \sigma^2 + k^2)$ .

Next, we note that for any given constant  $k \neq 0$ , the experiments  $N_n(0, \sigma^2)$  and  $N_n(0, k^2\sigma^2)$  are equivalent in the sense that each is sufficient for the other. This follows from the fact that multiplying each observation in the first experiment by k or multiplying each observation in the second experiment by 1/k maps each experiment into the other one.

We turn now to the much more difficult problem of determining whether either of the experiments  $N_n(0, \sigma^2)$  and  $N_n(0, \sigma^{2k})$  is sufficient for the other, where k is a given positive constant. We shall prove that for  $k_1 > k_2 > 0$ ,  $N_n(0, \sigma^{2k_1}) \ge N_n(0, \sigma^{2k_2})$ . First we present some related results.

LEMMA 2. Let W be a nonnegative random variable with pdf g(w) and let  $\varphi(t)$  denote the characteristic function of  $\log W$ . Let  $\theta > 0$  be an unknown parameter, let k > 0 be a given constant, and let  $G_n(\theta^k)$  denote the experiment in which a random sample of n observations is taken from the distribution with pdf  $(1/\theta^k)g(w/\theta^k)$ . For any given constant c > 0, define

(3.1) 
$$\psi_c(t) = \frac{\varphi(t)}{\varphi(ct)}, -\infty < t < \infty.$$

If  $\psi_{k_2/k_1}(t)$  is a characteristic function, then  $G_n(\theta^{k_1}) \ge G_n(\theta^{k_2})$ . Moreover,  $G_1(\theta^{k_1}) \ge G_1(\theta^{k_2})$  if and only if  $\psi_{k_2/k_1}(t)$  is a characteristic function.

The proof of this lemma follows from results in Boll (1955) or Torgersen (1972) by regarding  $\log \theta$  as a translation parameter in the distribution of  $\log W$ .

It should be noted that  $\psi_c(t)$ , defined in (3.1), is a characteristic function for all  $c \in (0, 1)$  if and only if  $\varphi(t)$  belongs to the class of self-decomposable characteristic functions, introduced by P. Lévy and A. Ya. Khinchine (See Lukacs (1970), subsection 5.11). Some interesting properties of this class (also called L-class by Gnedenko and Kolmogorov (1954)) are as follows:

- (i) All self-decomposable characteristic functions are infinitely divisible.
- (ii) If  $\varphi(t)$  is a self-decomposable characteristic function, then  $\psi_c(t)$  is infinitely divisible.
  - (iii) All stable characteristic functions are self-decomposable.
- (iv) The necessary and sufficient conditions for  $\varphi(t)$  to be self-decomposable in terms of Lévy's and Kolmogorov's canonical representations of an infinitely divisible characteristic functions are given in Theorems 1 and 2 of Chapter 6 in Gnedenko and Kolmogorov (1954).

We shall now assume that g is the density function of a gamma distribution G(a, b), defined in (1.1), with known values of a and b, and prove that the assumptions in Lemma 2 hold for this pdf.

THEOREM 3. Let  $G_n(a, b)$  denote the experiment in which a random sample of n observations is taken from the gamma distribution G(a, b) with pdf (1.1). Then  $G_n(a, b\theta^{k_1}) \ge G_n(a, b\theta^{k_2})$ , where  $\theta > 0$  is an unknown parameter, and  $a, b, k_1$  and  $k_2$  are given positive constants with  $k_1 > k_2$ .

PROOF. Let  $\varphi(t)$  denote the characteristic function of  $\log W$ , where W is a random variable with pdf (1.1), and let  $\alpha = k_2/k_1$ . Then

(3.2) 
$$\varphi(t) = \int_0^\infty \frac{1}{b^a \Gamma(a)} w^{it} w^{a-1} e^{-w/b} dw,$$
$$= b^{it} \Gamma(a+it) / \Gamma(a).$$

For  $\alpha < 1$ , consider

(3.3) 
$$\psi_{\alpha}(t) = \frac{\varphi(t)}{\varphi(\alpha t)} = \frac{b^{it(1-\alpha)}\Gamma(a+it)}{\Gamma(a+i\alpha t)}.$$

The Weerstrass expansion of  $1/\Gamma(z)$  (see Whittaker and Watson (1935), page 236) is

(3.4) 
$$\frac{1}{\Gamma(z)} = ze^{\gamma z} \prod_{j=1}^{\infty} \left\{ \left( 1 + \frac{z}{j} \right) \exp(-z/j) \right\}$$

where  $\gamma$  is the Euler's constant. Therefore, after some algebraic manipulation,  $\psi_{\alpha}(t)$  can be written as

$$(3.5) \quad \psi_{\alpha}(t) = b^{it(1-\alpha)} \exp\{-\gamma it(1-\alpha)\} \left(\frac{a+it\alpha}{a+it}\right) \prod_{j=1}^{\infty} \left(\frac{j+a+it\alpha}{j+a+it}\right) e^{it(1-\alpha)/j}$$

or, equivalently, as

(3.6) 
$$\psi_{\alpha}(t) = \exp\{-it(1-\alpha)(\gamma - \log b)\} \left[\alpha + (1-\alpha)\left(1 + \frac{it}{a}\right)^{-1}\right]$$

$$\times \prod_{j=1}^{\infty} \left[ \alpha + (1-\alpha) \left( 1 + \frac{it}{j+a} \right)^{-1} \right] e^{it(1-\alpha)/j}.$$

The first factor in (3.6) is the characteristic function of a degenerate random variable  $T_0$  with probability one at the point  $[-(1-\alpha)(\gamma-\log b)]$ , and the factor  $e^{it(1-\alpha)/j}$  is the characteristic function of a degenerate random variable  $T_j$  with probability one at the point  $(1-\alpha)/j$ ,  $j=1,2,\cdots$ . Furthermore,  $[\alpha+(1-\alpha)(1+(it/(j+a))^{-1}]$  is the characteristic function of a random variable  $Z_j$ ,  $j=0,1,2,\cdots$ , which takes the value 0 with probability  $\alpha$  and, with probability  $(1-\alpha)$ , has the pdf

(3.7) 
$$f_j(z) = (j+a)\exp[(j+a)z] \quad \text{for } z \le 0$$
$$= 0 \quad \text{for } z > 0.$$

Let  $\{T_i, i = 0, 1, 2, \dots\}$  and  $\{Z_i, i = 0, 1, 2, \dots\}$  be independent sequences of independent random variables with the distributions defined above. Define  $S_j = \sum_{l=0}^{j} (T_l + Z_l)$ , and let  $\psi_j(t)$  denote the characteristic function of  $S_j$ . It follows from (3.6) and the above discussion that

$$\psi_{\alpha}(t) = \lim_{i \to \infty} \psi_i(t).$$

It is obvious that  $\psi_{\alpha}(0) = 0$ . Furthermore, since the gamma function  $\Gamma(z)$  is analytic for complex arguments except at the points  $z = 0, -1, -2, \cdots$ , where it has simple poles,  $\psi_{\alpha}(t)$  is continuous at t = 0. Hence by the continuity theorem (Theorem 3.6.1, Lukacs (1970)),  $\psi_{\alpha}(t)$  is a characteristic function. It now follows from Lemma 2 that  $G_n(a, b\theta^{k_1}) \geq G_n(a, b\theta^{k_2})$  for all  $k_1 > k_2 > 0$ .  $\square$ 

- REMARK. (i) An alternative way to prove that  $\sum_{i=0}^{j} (Z_i + T_i)$  converges in distribution to a random variable Z, is to use Theorem 3.7.3, Lukacs (1970). Since  $\sum_{i=0}^{\infty} \operatorname{Var}(Z_i + T_i) = \alpha(1-\alpha)\sum_{i=0}^{\infty} 1/(i+a)^2 < \infty$ ,  $\psi_j(t)$  converges to a characteristic function  $\psi_{\alpha}(t)$  as  $j \to \infty$ .
- (ii) Another proof of the fact that  $\psi_{\alpha}(t)$  is a characteristic function could be given by proving that  $\varphi(t)$  is self-decomposable by using either Theorem 1 or Theorem 2 in Chapter 6 of Gnedenko and Kolmogorov (1954). However, we prefer the proof given above because it gives the specific form of the random variable Z in the proof of Lemma 2.
- (iii) Using Theorem 3.7.6 of Lukacs (1970), it can be shown that the distribution function of the random variable Z is continuous.

We shall now prove the main result of this section.

THEOREM 4. Let  $k_1$  and  $k_2$  be given constants satisfying  $k_1 > k_2 > 0$ . Then  $N_n(0, \sigma^{2k_1}) \ge N_n(0, \sigma^{2k_2})$ .

PROOF. As explained earlier, we can assume, without loss of generality, that  $k_2 = 1$  and  $k_1 = k$ . Let X and Y denote the observations in the experiments  $N_1(0, \sigma^{2k})$  and  $N_1(0, \sigma^2)$ , respectively. Since  $X^2/\sigma^{2k}$  and  $Y^2/\sigma^2$  have the same  $\chi^2$  distribution, it follows from Theorem 3 that the experiments in which  $X^2$  is observed is sufficient for that in which  $Y^2$  is observed. Furthermore, since  $X^2$  is a sufficient statistic for the experiment  $N_1(0, \sigma^{2k})$  and  $Y^2$  is a sufficient statistic for the experiment  $N_1(0, \sigma^2)$ , it follows that  $N_1(0, \sigma^{2k}) \ge N_1(0, \sigma^2)$ . Hence,  $N_n(0, \sigma^{2k}) \ge N_n(0, \sigma^2)$ .

If X and Y have the distributions specified in the proof of Theorem 4, we now know how to generate a random variable equivalent to an observation on Y from an observation on X.

Let the random variable Z be as defined in the proof of Theorem 3, independently of X, with  $a = \frac{1}{2}$ , b = 2, and  $\alpha = k_2/k_1$ , and let Y' be defined as follows

$$Y' = |X|^{\alpha} e^{Z/2}$$
 with probability  $\frac{1}{2}$ ,  
=  $-|X|^{\alpha} e^{Z/2}$  with probability  $\frac{1}{2}$ .

Then Y' has the same distribution as Y for every possible value of  $\sigma^2$ .

4. Comparison of experiments with location and scale parameters. Stone (1961) considers the class of experiments  $\{\mathcal{E}(c); c>0\}$  where  $\mathcal{E}(c)$  is the experiment in which an observation is taken from the pdf  $cf[c(x-\theta)]$ , for a fixed pdf f and  $\Theta=R$ . For given values of  $c_1$  and  $c_2$ , he obtains conditions under which  $\mathcal{E}(c_1) \geq \mathcal{E}(c_2)$ ,  $\mathcal{E}(c_1) \geq_L \mathcal{E}(c_2)$ , or  $\mathcal{E}(c_1) \geq_F \mathcal{E}(c_2)$ . Let  $\varphi(t)$  denote the characteristic function of the pdf f. Stone shows that if  $f(\cdot)$  is bounded and  $c_1 > c_2 > 0$ , then a sufficient condition for  $\mathcal{E}(c_1) \geq_F \mathcal{E}(c_2)$  is that

(4.1) 
$$\psi(t) = \frac{\varphi(t/c_2)}{\varphi(t/c_1)}$$

be a characteristic function. However, it follows from the references mentioned after Lemma 2 that the boundedness of  $f(\cdot)$  is not needed in this result. Furthermore, it follows that if  $\varphi(t)$  is a self-decomposable characteristic function, then  $\mathcal{E}(c_1) \geq \mathcal{E}(c_2)$  for all  $c_1 > c_2 > 0$ .

Stone also established that if  $f(\cdot)$  is bounded and the family of pdf's  $\{f(u-\theta); \theta \in R\}$  is boundedly complete, then a necessary condition that  $\mathcal{E}(c_1) \geq \mathcal{E}(c_2)$  whenever  $c_1 > c_2 > 0$  is that  $\psi(t)$  be a characteristic function. In addition, if all the cumulants of  $f(\cdot)$  exist, Stone proves that  $\psi(t)$  is a characteristic function, only if (i)  $f(\cdot)$  is a normal density or (ii) the even-order cumulants of  $f(\cdot)$  are positive. After proving this result, he states that "it is possible that condition (ii) is inconsistent with  $\mathcal{E}(c_1) \geq \mathcal{E}(c_2)$  whenever  $c_1 > c_2$ , in which event, yet another characterization of the normal distribution would be provided". However, for  $f(u) = \exp(-u)$ , u > 0, and  $c_1 > c_2 > 0$ , it can be shown that  $\psi(t)$  is a characteristic function and therefore  $\mathcal{E}(c_1) \geq \mathcal{E}(c_2)$ . Furthermore, all the cumulants of  $f(\cdot)$  exist, all the even order cumulants of  $f(\cdot)$  are positive,  $f(\cdot)$  is bounded, and the family of distributions  $\{f(u-\theta); \theta \in R\}$  is boundedly complete. Hence, this result does not provide yet another characterization of the normal distribution, as suggested by Stone.

A natural extension of the above results is to consider the class of experiments  $\{\mathcal{E}^*(c); c>0\}$  such that  $\mathcal{E}^*(c)$  is the experiment in which an observation is taken from the pdf  $(c/\sigma)f[c(x-\mu)/\sigma]$ , where f is a given pdf and the parameter space is  $\Theta = \{(\mu, \sigma) : \mu \in R, \sigma > 0\}$ . One may ask whether  $\mathcal{E}^*(c_1) \geq \mathcal{E}^*(c_2)$  for  $c_1 > c_2 > 0$ . In particular, one may ask whether  $N_n(\mu, \sigma^2)$  is sufficient for  $N_n(\mu, \sigma^2/c^2)$ , where c < 1 is a known constant. The answer is negative, as shown by Boll (1955). In fact, it follows from Theorem 3.1 of Hansen and Torgersen (1974) that  $N_{n_1}(\mu, \sigma^2/c_1^2) \geq N_{n_2}(\mu, \sigma^2/c_2^2)$  if and only if either  $n_1c_1^2 = n_2c_2^2$  and  $n_1 \geq n_2$  or  $n_1c_1^2 > n_2c_2^2$  and  $n_1 \geq n_2 + 1$ .

Stone (1961) also proved that in the location parameter case,  $\mathcal{E}(c_1) \geq_F \mathcal{E}(c_2)$  for  $c_1 > c_2 > 0$ , whenever the Fisher information exists. We shall now extend this result to the family of experiments  $\mathcal{E}*(c)$  with both location and scale parameters, defined at the beginning of this section.

THEOREM 5. Suppose that the pdf f(x) is symmetric around the point x = 0. For i = 1, 2, let  $\mathfrak{S}_i$  denote the experiment in which  $n_i$  independent replications of the experiment  $\mathfrak{S}^*(c_i)$  are performed, where  $c_1$  and  $c_2$  are given positive constants. Assume

that the Fisher information matrices  $\mathbf{i}_1(\mu, \sigma)$  and  $\mathbf{i}_2(\mu, \sigma)$  exist for these two experiments. Then  $\mathcal{E}_1 \geq_F \mathcal{E}_2$  if and only if  $n_1 \geq n_2$  and  $n_1 c_1^2 \geq n_2 c_2^2$ .

PROOF. It can be shown that if  $f(\cdot)$  is symmetric, then the matrix  $\mathbf{i}_1(\mu, \sigma) - \mathbf{i}_2(\mu, \sigma)$  is diagonal with diagonal elements  $(n_1c_1^2 - n_2c_2^2)A$  and  $(n_1^2 - n_2^2)B$ , where A and B are the positive diagonal elements of the information matrix for the experiment  $\mathfrak{S}^*(1)$ . Hence,  $\mathbf{i}_1(\mu, \sigma) - \mathbf{i}_2(\mu, \sigma)$  will be nonnegative definite if and only if  $n_1c_1^2 \geqslant n_2c_2^2$  and  $n_1^2 \geqslant n_2^2$ .  $\square$ 

It has been pointed out by a referee that for any given pdf f and any positive constants  $c_1 \neq c_2$ , the experiments  $\mathcal{E}^*(c_1)$  and  $\mathcal{E}^*(c_2)$  are never comparable with respect to the sufficiency ordering  $\geq$ .

5. Marginally sufficient experiments. In general, the relation  $\mathcal{E}_X \geqslant \mathcal{E}_Y$  is equivalent to the requirement that  $\mathcal{E}_X$  is at least as preferred as  $\mathcal{E}_Y$  for every decision problem involving the parameter  $\theta$  and every prior distribution on  $\Theta$ . Therefore, it is a very restrictive relation and induces only a partial ordering on the class  $E(\Theta)$  of all possible experiments with parameter space  $\Theta$ . Feldman (1972) studied certain properties of orderings of  $E(\Theta)$  induced by the weakened requirement that in a fixed decision problem, the expected Bayes risk from  $\mathcal{E}_X$  be not greater than that from  $\mathcal{E}_Y$  for every prior distribution  $\xi \in \Xi$ . Following DeGroot (1962), he identified the decision problem with an uncertainty function  $\mathfrak{A}(\xi)$  defined on  $\Xi$  and considered the experiment  $\mathcal{E}_X$  to be at least as informative as the experiment  $\mathcal{E}_Y$  with respect to  $\mathfrak{A}$  if  $\mathfrak{A}(\xi|X) \leq \mathfrak{A}(\xi|Y)$  for all  $\xi \in \Xi$ , where  $\mathfrak{A}(\xi|X)$  is the expected posterior uncertainty if X is observed and the prior distribution is  $\xi$  and  $\mathfrak{A}(\xi|Y)$  is the corresponding value for the observation Y.

An alternative possibility for comparing experiments is to consider a fixed prior distribution  $\xi$  and study the ordering on  $E(\Theta)$  induced by the requirement that the expected Bayes risk from  $\mathcal{E}_X$  be not greater than that from  $\mathcal{E}_Y$  for every decision problem involving  $\theta$ . In this case, we will say that  $\mathcal{E}_X$  is at least as informative as  $\mathcal{E}_Y$  with respect to the prior distribution  $\xi$ .

If  $\mathcal{E}_X$  is at least as informative as  $\mathcal{E}_Y$  for a fixed decision problem, then every experimenter interested in that decision problem will prefer  $\mathcal{E}_X$  to  $\mathcal{E}_Y$  since any risk function that can be attained from  $\mathcal{E}_Y$  can be matched or dominated by one from  $\mathcal{E}_X$ . On the other hand, if  $\mathcal{E}_X$  is at least as informative as  $\mathcal{E}_Y$  with respect to a prior distribution  $\xi$ , then an experimenter with prior distribution  $\xi$  on  $\Theta$  will prefer  $\mathcal{E}_X$  to  $\mathcal{E}_Y$  regardless of his decision problem. We shall now give an example to illustrate this concept.

EXAMPLE 2. Let  $c_1 > c_2 > 0$  be given constants and for i = 1, 2, let  $X_i$  denote a random variable with the normal distribution  $N(\mu, \sigma^2/c_i^2)$ . Suppose that the prior distribution of  $(\mu, \sigma)$  is concentrated on just two points such that  $\Pr[(\mu, \sigma) = (0, 1)] = \xi$  and  $\Pr[(\mu, \sigma) = (\mu_0, \sigma_0)] = 1 - \xi$  where  $0 < \xi < 1$ ,  $\sigma_0$  and  $\mu_0$  are known and arbitrary. It follows from Bradt and Karlin (1956) that when the parameter space is regarded as containing just these two points,  $N_1(\mu, \sigma^2/c_1^2)$  is sufficient for

 $N_1(\mu, \sigma^2/c_2^2)$ , and therefore  $N_n(\mu, \sigma^2/c_1^2) \ge N_n(\mu, \sigma^2/c_2^2)$ . Hence  $N_n(\mu, \sigma^2/c_1^2)$  is more informative than  $N_n(\mu, \sigma^2/c_2^2)$  with respect to this prior distribution.

Torgersen (1976, Theorem 1) shows that if  $\Theta$  is countable, then the partial ordering of experiments with respect to a given prior distribution  $\xi$  such that  $\xi(\theta) > 0$  for each  $\theta \in \Theta$  is equivalent to the usual sufficiency partial ordering  $\geq$ . However, the concept of relative informativeness with respect to a prior distribution  $\xi$  is especially useful when the parameter  $\theta$  is vector valued,  $\theta = (\theta_1, \theta_2)$ , and the experimenter is interested only in  $\theta_1$ ; i.e.,  $\theta_2$  is a nuisance parameter. For example, in the experiment  $\mathcal{E}^*(c)$  defined in Section 4, corresponding to the pdf  $(c/\sigma)f[(c/\sigma)(x-\mu)]$ , the decision problems of interest may involve only  $\mu$  or only  $\sigma$ . A detailed discussion on the elimination of nuisance parameters in the framework of classical statistical inference is given by Basu (1977). For a given prior distribution,  $\xi(\theta_1, \theta_2) = \xi_1(\theta_1)\xi_2(\theta_2|\theta_1)$ , a Bayesian statistician who is interested only in  $\theta_1$ , because the loss depends only on  $\theta_1$ , will eliminate  $\theta_2$  from the analysis and use the prior pdf  $\xi_1(\theta_1)$  together with the conditional pdf

(5.1) 
$$g(x|\boldsymbol{\theta}_1) = \int_{\boldsymbol{\Theta}_2} p(x|\boldsymbol{\theta}_1, \,\boldsymbol{\theta}_2) d\xi_2(\boldsymbol{\theta}_2|\boldsymbol{\theta}_1).$$

Consider a particular decision problem with  $\Theta = \{\theta = (\theta_1, \theta_2) | \theta_1 \in \Theta_1, \theta_2 \in \Theta_2\}$  and a given class D of all possible decisions d, and let  $l(\theta, d)$  denote the loss incurred from any decision  $d \in D$  when  $\theta \in \Theta$  is true. We shall say that the decision problem involves only  $\theta_1$  if, for every pair  $(\theta_1, d)$ ,

$$(5.2) l[(\boldsymbol{\theta}_1, \boldsymbol{\theta}_2), d] = l[(\boldsymbol{\theta}_1, \boldsymbol{\theta}_2^*), d] \text{for all } \boldsymbol{\theta}_2^* \in \Theta_2,$$

i.e., l depends only on the value of  $\theta_1$  and the value of d, and not on the value of  $\theta_2$ . For such decision problems, we now present a natural and useful concept of marginal Bayesian sufficiency with respect to a given prior distribution  $\xi(\theta_1, \theta_2)$ .

**DEFINITION.** The experiment  $\mathcal{E}_X$  is marginally sufficient for  $\mathcal{E}_Y$ , denoted by  $\mathcal{E}_X \geqslant \mathcal{E}_Y(\boldsymbol{\theta}_1)$ , with respect to the prior distribution  $\xi(\boldsymbol{\theta}_1, \boldsymbol{\theta}_2)$  if the expected Bayes risk from  $\mathcal{E}_X$  is not greater than that from  $\mathcal{E}_Y$  for every decision problem involving only  $\boldsymbol{\theta}_1$ , when the prior distribution is  $\xi(\boldsymbol{\theta}_1, \boldsymbol{\theta}_2)$ .

For each  $\theta_1 \in \Theta_1$ , let  $G_{\theta_1}$  denote the distribution on  $S_X$  represented by the conditional pdf given by (5.1), and let  $H_{\theta_1}$  denote the analogous distribution on  $S_Y$ . Also, let  $\mathcal{E}_X' = \{X, S_X; G_{\theta_1}, \theta_1 \in \Theta_1\}$  and  $\mathcal{E}_Y' = \{Y, S_Y; H_{\theta_1}, \theta_1 \in \Theta_1\}$ . If  $\mathcal{E}_X' \geq \mathcal{E}_Y'$  then it will be true that  $\mathcal{E}_X \geq \mathcal{E}_Y(\theta_1)$  with respect to any prior distribution  $\xi^*$  that yields the same conditional distribution  $\xi_2(\theta_2|\theta_1)$  as  $\xi$ . In this case, we shall say that  $\mathcal{E}_X \geq \mathcal{E}_Y(\theta_1)$  with respect to the conditional prior distribution  $\xi_2(\theta_2|\theta_1)$ .

We shall now give some examples of marginal Bayesian sufficiency.

EXAMPLE 3. For a given pdf f, let  $\mathcal{E}^*(c)$  denote the experiment defined in Section 4, and let  $\varphi(t) = \int_R e^{itu} f(u) du$ . For any joint prior distribution of  $\mu$  and  $\sigma$ , let  $\xi_2(\sigma)$  denote the marginal prior distribution of  $\sigma$  and let  $\varphi_1(t) = \int_0^\infty \varphi(t\sigma) d\xi_2(\sigma)$ . It follows from (5.1) that if  $\mu$  and  $\sigma$  are independent under their joint prior distribution, then  $g(x|\mu)$  is of the form  $cg^*[c(x-\mu)]$ . Therefore, if  $\varphi_1(t)$  is a self-decomposable characteristic function, then it follows from the result of Stone,

presented at the beginning of Section 4, that  $\mathscr{E}^*(c_1) \geq \mathscr{E}^*(c_2)(\mu)$  with respect to the conditional prior distribution  $\xi_2(\sigma)$  for  $c_1 > c_2 > 0$ .

In particular, let f(u) be the standard normal pdf and let either (i) the prior density of  $\sigma^2$  be a gamma distribution of the form  $G(\alpha, \beta)$ , or (ii) the prior density of  $(1/\sigma^2)$  be a gamma distribution of the form  $G(\frac{1}{2}, \beta)$ . By carrying out the analysis indicated in this example, it can be shown that  $N_1(\mu, \sigma^2/c_1) \ge N_1(\mu, \sigma^2/c_2)(\mu)$  with respect to both of these conditional prior distributions of  $\sigma^2$  given  $\mu$ .

EXAMPLE 4. Let  $c_1 > c_2 > 0$  be given constants and, for i = 1, 2, let  $X_i$  denote a random variable with the normal distribution  $N[\mu, \sigma^2/c_i]$ . Suppose that our interest lies in decision problems involving only  $\sigma^2$ . If  $\mu$  and  $\sigma$  are independent under their joint prior distribution, and if the marginal distribution of  $\mu$  is a normal distribution  $N(m, \tau^2)$ , then it follows that, given  $\sigma^2$ ,  $X_i$  is distributed as  $N[m, \sigma^2/c_i^2 + \tau^2]$ . Let W be distributed as  $N[(1 - c_2/c_1)m, (1 - (c_2^2/c_1^2))\tau^2]$  independently of  $X_2$ . Then it can be verified that  $(c_2/c_1)X_2 + W$  has the same distribution as  $X_1$  for every possible value of  $\sigma^2$ . Hence,  $N_1(\mu, \sigma^2/c_2^2) \ge N_1(\mu, \sigma^2/c_1^2)(\sigma^2)$  with respect to this conditional prior distribution of  $\mu$ . In fact, using the joint distribution of  $\overline{X}$  and  $S^2$  from a random sample of n observations given  $\sigma^2$ , it can be shown that  $N_n(\mu, \sigma^2/c_2^2) \ge N_n(\mu, \sigma^2/c_1^2)(\sigma^2)$  with respect to this conditional prior distribution of  $\mu$ .

However, if the conditional prior distribution of  $\mu$  given  $\sigma^2$  is  $N(m, \sigma^2/\tau^2)$ , then it follows that  $X_i$  is distributed as  $N[m, \sigma^2(1/c_i^2 + 1/\tau^2)]$ , given  $\sigma^2$ . Therefore, the experiments  $N_1(\mu, \sigma^2/c_1^2)$  and  $N_1(\mu, \sigma^2/c_2^2)$  are sufficient for each other with respect to this conditional prior distribution of  $\mu$ . Furthermore, it can be shown that the experiments  $N_n(\mu, \sigma^2/c_1^2)$  and  $N_n(\mu, \sigma^2/c_2^2)$  are sufficient for each other with respect to this conditional prior distribution of  $\mu$ .

EXAMPLE 5. For given constants  $c_1 > c_2 > 0$ , consider again the normal experiments  $N_n(\mu, \sigma^2/c_2^2)$ , i = 1, 2, and suppose that the joint prior distribution of  $\mu$  and  $\sigma^2$  is a conjugate normal-gamma distribution such that the conditional distribution of  $\mu$  given  $\sigma^2$  is  $N[m, \sigma^2/\tau^2]$  and the distribution of  $(1/\sigma^2)$  is  $G(\alpha, \beta)$ . It follows from Example 3 that for decision problems involving only  $\sigma^2$ , the experiments  $N_n(\mu, \sigma^2/c_1^2)$  and  $N_n(\mu, \sigma^2/c_2^2)$  are marginally equivalent with respect to this joint prior distribution. However, for decision problems involving only  $\mu$ , it is not known whether one of these experiments is marginally sufficient for the other with respect to this conjugate joint prior distribution. We can prove, however, that for estimating any of the functions  $\mu$ ,  $\mu/\sigma$ ,  $\mu/\sigma^2$ ,  $\mu\sigma$  and  $\mu\sigma^2$  with squared-error loss, the experiment  $N_n(\mu, \sigma^2/c_1^2)$  has a smaller expected Bayes risk than the experiment  $N_n(\mu, \sigma^2/c_1^2)$  for this conjugate prior distribution.

EXAMPLE 6. If a statistic T(Y) is partially sufficient for the parameter  $\theta_1$ , according to Fraser's definition (1956), then it can be proved that the experiment  $\mathcal{E}_T$ , in which only T is observed, satisfies  $\mathcal{E}_T \geq \mathcal{E}_Y(\theta_1)$  with respect to any prior distribution  $\xi(\theta_1, \theta_2)$  under which  $\theta_1$  and  $\theta_2$  are independent. For example, if

 $Y_1, \dots, Y_n$  are independent and identically distributed with a gamma distribution  $G(\alpha, \beta)$ , then  $T = \sum_{i=1}^{n} Y_i$  is partially sufficient for  $\beta$  in Fraser's sense and, therefore,  $\mathcal{E}_T \geq \mathcal{E}_Y(\beta)$  with respect to any prior distribution for which  $\alpha$  and  $\beta$  are independent.

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