## STK3405 – Week 43 (2)

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## Chapter 7

Computing reliability using conditional Monte Carlo methods

#### Section 7.1

Monte Carlo simulation and conditioning

Let  $\boldsymbol{X}=(X_n,\ldots,X_n)$  be a random vector with a known distribution, and let  $\phi=\phi(\boldsymbol{X})$ . The distribution of  $\phi$  cannot be derived analytically in polynomial time with respect to n. We want to estimate:

$$h = E[\phi(\mathbf{X})]$$

We run a Monte Carlo simulation generating N independent vectors  $X_1, \ldots, X_N$ , all having the same distribution as X, and estimate h by the simple *Monte Carlo* estimate:

$$\hat{h}_{MC} = \frac{1}{N} \sum_{r=1}^{N} \phi(\boldsymbol{X}_r).$$

Then  $E[\hat{h}_{MC}] = h$ , and we have:

$$\operatorname{Var}(\hat{h}_{MC}) = \frac{1}{N^2} \sum_{r=1}^{N} \operatorname{Var}(\phi) = \frac{1}{N} \operatorname{Var}(\phi)$$

Let S = S(X) is a *discrete* variable with values in the set  $\{s_1, \ldots, s_k\}$ , and with a distribution that can be calculated analytically in polynomial time with respect to n.

We also introduce:

$$\theta_j = E[\phi \mid S = s_j], \quad j = 1, \ldots, k.$$

We then have:

$$h = E[\phi] = \sum_{j=1}^{k} E[\phi \mid S = s_j] P(S = s_j) = \sum_{j=1}^{k} \theta_j P(S = s_j)$$

Instead of generating N samples from the distribution of X, we divide the sample set into k groups, one for each possible value of S.

The *j*-th group has size  $N_j$ , j = 1, ..., k, and  $N_1 + \cdots + N_k = N$ .

The samples in the *j*-th group,  $X_{1,j}, \ldots, X_{N_j,j}$ , are sampled from the conditional distribution of X given that  $S = s_j$ ,  $j = 1, \ldots, k$ .

Then  $\theta_i$  is estimated by:

$$\hat{\theta}_j = \frac{1}{N_j} \sum_{r=1}^{N_j} \phi(\boldsymbol{X}_{r,j}), \quad j = 1, \dots, k.$$

We then have:

$$E[\hat{\theta}_j] = \frac{1}{N_j} \sum_{r=1}^{N_j} E[\phi(\boldsymbol{X}_{r,j})] = \frac{1}{N_j} \sum_{r=1}^{N_j} E[\phi|S = s_j]$$
$$= \frac{1}{N_j} \sum_{r=1}^{N_j} \theta_j = \theta_j, \qquad j = 1, \dots, k.$$

Moreover, the variances of these estimates are:

$$\operatorname{Var}(\hat{\theta}_{j}) = \frac{1}{N_{j}^{2}} \sum_{r=1}^{N_{j}} \operatorname{Var}(\phi(\boldsymbol{X}_{r,j})) = \frac{1}{N_{j}^{2}} \sum_{r=1}^{N_{j}} \operatorname{Var}(\phi|S = s_{j})$$
$$= \frac{1}{N_{j}} \operatorname{Var}(\phi|S = s_{j}), \qquad j = 1, \dots, k.$$

By combining  $\hat{\theta}_1, \dots, \hat{\theta}_k$ , we get the *conditional Monte Carlo estimate*:

$$\hat{h}_{CMC} = \sum_{j=1}^{k} \hat{\theta}_j P(S = s_j).$$

Since  $\hat{\theta}_1, \dots, \hat{\theta}_k$  are independent, the variance of the conditional Monte Carlo estimate is:

$$\operatorname{Var}(\hat{h}_{CMC}) = \operatorname{Var}[\sum_{j=1}^{k} \hat{\theta}_{j} P(S = s_{j})] = \sum_{j=1}^{k} \operatorname{Var}(\hat{\theta}_{j}) [P(S = s_{j})]^{2}$$
$$= \sum_{j=1}^{k} \frac{1}{N_{j}} \operatorname{Var}(\phi | S = s_{j}) [P(S = s_{j})]^{2}.$$

NOTE: The variance of  $\hat{h}_{CMC}$  depends on the choices of the  $N_j$ 's.

In order to compare the result with the variance of  $\hat{h}_{MC}$ , we let:

$$N_j \approx N \cdot P(S = s_j), \quad j = 1, \ldots, k.$$

Hence:

$$\sum_{j=1}^k N_j \approx \sum_{j=1}^k N \cdot P(S=s_j) = N \cdot \sum_{j=1}^k P(S=s_j) = N.$$

With this choice we get:

$$\begin{aligned} Var(\hat{h}_{CMC}) &\approx \sum_{j=1}^{k} \frac{1}{N \cdot P(S = s_j)} \operatorname{Var}(\phi | S = s_j) [P(S = s_j)]^2 \\ &= \frac{1}{N} \sum_{j=1}^{k} \operatorname{Var}(\phi | S = s_j) P(S = s_j) = \frac{1}{N} E[\operatorname{Var}(\phi | S)] \\ &= \frac{1}{N} (\operatorname{Var}(\phi) - \operatorname{Var}[E(\phi | S)]) \leq \frac{1}{N} \operatorname{Var}(\phi) = \operatorname{Var}(\hat{h}_{MC}). \end{aligned}$$

Here we have used the well-known relation:

$$Var(\phi) = Var[E(\phi|S)] + E[Var(\phi|S)]$$



NOTE:  $Var(\hat{h}_{CMC}) < Var(\hat{h}_{MC})$  whenever  $Var[E(\phi \mid S)] > 0$ .

 $Var[E(\phi \mid S)]$  can be interpreted as a measure of how much information S contains relative to  $\phi$ .

Thus, we should look for S containing as much information about  $\phi$  as possible. At the same time S must be such that:

- ullet The distribution of S can be derived analytically in polynomial time
- The number of possible values of S, i.e., k, must be bounded by a polynomial function of n
- It must be possible to sample efficiently from the distribution of X given S

#### Section 7.2

Conditioning on the sum of the component state variables

## Conditioning on the sum

Consider a binary monotone system  $(C, \phi)$  where  $C = \{1, ..., n\}$ , and let  $\mathbf{X} = (X_1, ..., X_n)$  be the vector of the component state variables.

Moreover, let:

$$S = S(X) = \sum_{i=1}^{n} X_i$$

Thus, the set of possible values for S is  $\{0, 1, ..., n\}$ , and we let:

$$\theta_s = E[\phi|S=s], \qquad s=0,1,\ldots,n.$$

We must find an efficient way of sampling from the conditional distribution of  $\boldsymbol{X}$  given S = s, s = 0, 1, ..., n.

This can be done as follows:

STEP 1. Sample  $X_1$  from the conditional distribution of  $X_1|S=s$ , and let  $x_1$  be the result.

STEP 2. Sample  $X_2$  from the conditional distribution of  $X_2|S=s, X_1=x_1$ , and let  $x_2$  be the result.

. . .

STEP n. Sample  $X_n$  from the conditional distribution of  $X_n | S = s, X_1 = x_1, \dots X_{n-1} = x_{n-1}$ , and let  $x_n$  be the result.

To compute the necessary conditional distributions we introduce the partial sums,  $S_1, \ldots, S_n$ :

$$S_m = \sum_{i=m}^n X_i, \quad m = 1, \ldots, n.$$

We then have:

$$P(X_{m} = x_{m} \mid X_{1} = x_{1}, \dots, X_{m-1} = x_{m-1}, S = s)$$

$$= \frac{P(X_{m} = x_{m}, S = s \mid X_{1} = x_{1}, \dots, X_{m-1} = x_{m-1})}{P(S = s \mid X_{1} = x_{1}, \dots, X_{m-1} = x_{m-1})}$$

$$= \frac{P(X_{m} = x_{m}, S_{m+1} = s - \sum_{j=1}^{m} x_{j})}{P(S_{m} = s - \sum_{j=1}^{m-1} x_{j})}$$

$$= \frac{p_{m}^{x_{m}} (1 - p_{m})^{1 - x_{m}} P(S_{m+1} = s - \sum_{j=1}^{m} x_{j})}{P(S_{m} = s - \sum_{j=1}^{m-1} x_{j})},$$

The distributions of the partial sums  $S_1, \ldots, S_n$  can be calculated before running the simulations using *generating functions*. See Exercise 2.10.



NOTE: We only calculate the conditional probabilities we actually need along the way during the simulations, not the entire set of all possible conditional probabilities corresponding to all possible combinations of values of the  $X_i$ 's.

Thus, in each simulation we calculate n probabilities, one for each  $X_j$ .

Each probability is calculated using a fixed number of operations (independent of n).

Hence, sampling from the conditional distribution of X given S = s, can be done in O(n) time, i.e., just as fast as sampling from the unconditional distribution.

When choosing the sample sizes  $N_0, \ldots, N_n$ , we may let  $N_s \approx N \cdot P(S = s)$ ,  $s = 0, 1, \ldots, n$ .

However, it is possible to improve the results slightly. By examining the system, it is often easy to determine the size of the smallest minimal path and cut sets, and we introduce:

d = The size of the smallest minimal path set

c = The size of the smallest minimal cut set

We then have that  $\theta_s = 0$  for s < d, and  $\theta_s = 1$  for s > n - c, implying that  $Var(\phi|S = s) = 0$  for s < d, or s > n - c.

Hence, there is no point in spending simulations on estimating  $\theta_s$  for s < d, or s > n - c, so we let  $N_s = 0$  for s < d, or s > n - c. As a result, we have more simulations to spend on the remaining quantities.

An extreme situation occurs when the system is a k-out-of-n-system, i.e., when  $\phi(\mathbf{X}) = I(S \ge k)$ . For such systems d = k, and c = (n - k + 1).

This implies that n-c=n-(n-k+1)=k-1.

Hence,  $\theta_s = 0$  for s < k and  $\theta_s = 1$  for  $s \ge k$ .

Thus, the CMC-estimate is equal to the true value of the reliability, and can be calculated without doing *any simulations at all*.

For all other nontrivial systems, however, it is easy to see that we always have that  $d \le n - c$ . In order to ensure that we get improved results, we assume that  $P(d \le S \le n - c) > 0$ , and let:

$$N_s \approx N \cdot P(S=s)/P(d \leq S \leq n-c), \quad s=d,\ldots,n-c$$
.

We can now show that:

$$\operatorname{Var}(\hat{h}_{CMC}) \leq P(d \leq S \leq n-c) \operatorname{Var}(\hat{h}_{MC})$$

Hence, if d and n - c are close, the variance is reduced considerably.

#### Section 7.3

System reliability when all the component state variables have identical reliabilities

## Identical component reliabilities

If  $p_1 = \cdots = p_n = p$ , it is possible to improve things even further. Then S has a binomial distribution, and the conditional distribution of X given S is given by:

$$P(X = x \mid S = s) = \frac{p^{\sum_{i=1}^{n} x_i} (1-p)^{n-\sum_{i=1}^{n} x_i}}{\binom{n}{s} p^s (1-p)^{n-s}} = \frac{1}{\binom{n}{s}},$$

for all  $\mathbf{x}$  such that  $\sum_{i=1}^{n} x_i = \mathbf{s}$ .

From this it follows that:

$$egin{aligned} heta_{m{s}} = E[\phi \mid m{S} = m{s}] = \sum_{\left\{m{X} \mid \sum_{i=1}^{n} x_i = m{s}
ight\}} \phi(m{x}) P(m{X} = m{x} \mid m{S} = m{s}) = rac{m{b_s}}{\left(rac{n}{m{s}}
ight)} \;, \end{aligned}$$

where  $b_s$  is the number of path sets with s components,  $s = 0, \dots, n$ .

#### Identical component reliabilities (cont.)

The system reliability, h, expressed as a function of p, is given by:

$$h(p) = \sum_{s=0}^{n} \theta_{s} P(S = s)$$

$$= \sum_{s=0}^{n} \frac{b_{s}}{\binom{n}{s}} \binom{n}{s} p^{s} (1-p)^{n-s} = \sum_{s=0}^{n} b_{s} p^{s} (1-p)^{n-s}.$$

NOTE 1:  $\theta_0, \dots, \theta_n$  do not depend on p. Thus, by estimating these quantities, we get an estimate of the entire h(p)-function for all  $p \in [0, 1]$ .

NOTE 2:  $\theta_s$  can be interpreted as the fraction of path sets of size s among all sets of size s. Thus,  $\theta_s$  can be estimated by sampling random sets of size s and calculating the frequency of path sets among the sampled sets.

#### Identical component reliabilities (cont.)

ALGORITHM: Let  $\mathbf{T} = (T_1, \dots, T_n)$  be a vector for storing results from the simulations. This vector is initialized as  $(0, \dots, 0)$ .

Then for i = 1, ..., N do:

- 1. Sample a component from the set C, say component  $i_1$ , and define  $A_1 = \{i_1\}$ . If  $A_1$  is a path set,  $T_1$  is incremented with 1.
- 2. Sample a component from the set  $C \setminus A_1$ , say component  $i_2$ , and define  $A_2 = \{i_1, i_2\}$ . If  $A_2$  is a path set,  $T_2$  is incremented with 1. . . .
- *n*. Sample the last remaining component, say component  $i_n$ , and define  $A_n = C$ . If  $A_n$  is a path set,  $T_n$  is incremented with 1.

When all the simulations are carried out, the vector T contains the number of observed path sets of sizes  $1, \ldots, n$ . From this we get the resulting estimates of  $\theta_1, \ldots, \theta_n$  simply as:

$$\theta_s = T_s/N, \quad s = 1, \ldots, n.$$