

Stochastic flow networks: How component criticality changes with component reliability

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ABSTRACT

Consider a directed flow network whose arcs have random two-state capacities. The paper describes how a single Monte Carlo experiment allows one to estimate the sensitivity of the probability that a particular cutset is critical, given that the maximum flow between a pair of nodes takes values in an interval, in response to varying arc reliabilities. The technique improves considerably on alternative potential methods.

THE PROBLEM

Flow networks model several distribution systems, electricity and water supply systems being examples. In practice, several of the parameters of a flow network are random variables and the network is called stochastic. The model under study is a directed network $\mathcal{G} = (\mathcal{N}, \mathcal{A})$ with node set $\mathcal{N} = \{1, \dots, n\}$ and arc set $\mathcal{A} = \{1, \dots, a\}$. Assume that the nodes do not restrict the transmission of flow through the network and that the arcs have finite capacities which are independent two-state random variables B_i with states 0 and b_i for $i \in \mathcal{A}$. Let $X_i = 1$ if $B_i = b_i$ and $X_i = 0$ otherwise. Also, let s and t denote a pair of selected nodes in \mathcal{N} . For given capacity state-vector $x \in \mathcal{X}$ where $\mathcal{X} = \{0, 1\}^a$, let $\Lambda(x)$ denote the value of a maximum s - t flow. Every directed minimal s - t cutset (or s - t cutset hereafter) \mathcal{C} separates \mathcal{N} into two sets X and \bar{X} such that $X \cup \bar{X} = \mathcal{N}$, $X \cap \bar{X} = \emptyset$, $s \in X$ and $t \in \bar{X}$. The capacity of \mathcal{C} , denoted by $Z(\mathcal{C}, x)$, is defined as $\sum_{i \in (X, \bar{X})} b_i x_i$ and satisfies $Z(\mathcal{C}, x) \geq \Lambda(x)$. The cutset \mathcal{C} is called critical when $Z(\mathcal{C}, x) = \Lambda(x)$. For a review on maximum flows see Ford and Fulkerson (1962).

Fix two maximum flow values $l < u \leq \Lambda(b_1, \dots, b_a)$ and a cutset \mathcal{C} . This paper proposes an efficient Monte Carlo sampling plan for estimating the variation of the probability $h(\mathcal{C}, q)$ that the cutset \mathcal{C} is critical given that the value of a maximum s - t flow $\Lambda(X)$ lies in the interval $(l, u]$ in response to varying reliability vectors q in a set $\mathcal{S} \subset (0, 1)^a$, where

$q_i = \text{pr}(B_i = b_i)$ for $i \in \mathcal{A}$. We call the vectors in $(0, 1)^a$ points. The evaluation of $h(\mathcal{C}, q)$ at a single point q is an intractable (NP-hard) problem.

Identifying s - t cutsets \mathcal{C}^* with high $h(\mathcal{C}^*, q)$, for fixed q , as well as points q for which the probabilities $h(\mathcal{C}^*, q)$, for a fixed s - t cutset \mathcal{C}^* , become large or small are very important issues in repairing and designing stochastic networks.

For a fixed point q , let $g(q)$ denote the probability that $l < \Lambda(X) \leq u$ and let $f(\mathcal{C}, q)$ denote the probability that \mathcal{C} is a critical s - t cutset and $l < \Lambda(X) \leq u$. We then define

$$h(\mathcal{C}, q) = \begin{cases} f(\mathcal{C}, q)/g(q) & \text{if } g(q) > 0 \\ 0 & \text{otherwise.} \end{cases} \quad (1)$$

Hereafter, assume $g(q) > 0 \forall q \in \mathcal{S}$.

We first describe briefly an efficient Monte Carlo importance sampling plan in Alexopoulos and Fishman (1988) for estimating $h(\mathcal{C}, q)$ at a single point q and then we show how this plan can be extended for estimating this flow performance measure at all points in \mathcal{S} . The sampling plans use an upper bound on the flow probability $g(q)$, $q \in [0, 1]^a$ to gain their advantage over a crude Monte Carlo sampling plan.

Let $P(x, q) = \prod_{i=1}^a q_i^{x_i} (1-q_i)^{1-x_i}$ denote the p.m.f. of the random vector $X = (X_1, \dots, X_a)$. Also, let $\mathcal{P}_1, \dots, \mathcal{P}_I$ denote arc-disjoint directed s - t paths and $\mathcal{C}_1, \dots, \mathcal{C}_J$ denote arc-disjoint s - t cutsets. Define $\Lambda_1(X) = \sum_{i=1}^I [\min_{j \in \mathcal{P}_i} b_j] \prod_{j \in \mathcal{P}_i} X_j$, that is the maximum amount of flow that can be transmitted from s to t through the paths $\mathcal{P}_1, \dots, \mathcal{P}_I$ and $\Lambda_2(X) = \min_{j=1, \dots, J} \sum_{i \in \mathcal{C}_j} b_i X_i$.

One can readily show that

$$\begin{aligned}
g(\mathbf{q}) \leq g_U(\mathbf{q}) &= \text{pr}[\Lambda_2(X) > l, \Lambda_1(X) \leq u] \\
&= \text{pr}[\Lambda_2(X) > l] - \text{pr}[\Lambda_1(X) \leq u].
\end{aligned} \tag{2}$$

For $x \in \mathcal{X}$ let $\phi(x) = I_{(l,u]}(\Lambda(x))$, $\phi_U(x) = I_{(l,\infty)}(\Lambda_2(x))I_{(-\infty,u]}(\Lambda_1(x))$ and $\psi(\mathcal{E}, x) = I_{\{0\}}(Z(\mathcal{E}, x) - \Lambda(x))$ where $I_{(c,d]}(x)$ is the indicator function on the interval $(c, d]$.

To take advantage of this upper bound, one constructs a conditional distribution

$$Q(x, \mathbf{q}) = \phi_U(x)P(x, \mathbf{q})/g_U(\mathbf{q}) \tag{3}$$

(see ref. [AF] for details) and draws K independent samples $X^{(1)}, \dots, X^{(K)}$ from it. An estimate for $h(\mathcal{E}, \mathbf{q})$ is given by

$$\begin{aligned}
\tilde{h}_K(\mathcal{E}, \mathbf{q}) &= \tilde{f}_K(\mathcal{E}, \mathbf{q})/\tilde{g}_K(\mathbf{q}) && \text{if } \tilde{g}_K(\mathbf{q}) > 0 \\
&= 0 && \text{otherwise}
\end{aligned} \tag{4}$$

where

$$\tilde{f}_K(\mathcal{E}, \mathbf{q}) = g_U(\mathbf{q}) \frac{1}{K} \sum_{k=1}^K \phi(X^{(k)})\psi(\mathcal{E}, X^{(k)}) \tag{5}$$

and

$$\tilde{g}_K(\mathbf{q}) = g_U(\mathbf{q}) \frac{1}{K} \sum_{k=1}^K \phi(X^{(k)})$$

are unbiased estimates of $f(\mathcal{E}, \mathbf{q})$ and $g(\mathbf{q})$, respectively, with variances

$$\text{var } \tilde{f}_K(\mathcal{E}, \mathbf{q}) = f(\mathcal{E}, \mathbf{q})[g_U(\mathbf{q}) - f(\mathcal{E}, \mathbf{q})]/K \leq g_U(\mathbf{q})^2/K$$

and

$$\text{var } \tilde{g}_K(\mathcal{E}, \mathbf{q}) = g(\mathbf{q})[g_U(\mathbf{q}) - g(\mathbf{q})]/K \leq g_U(\mathbf{q})^2/K.$$

From Alexopoulos (1988) one has

$$E\tilde{h}_K(\mathcal{E}, \mathbf{q}) = h(\mathcal{E}, \mathbf{q})[1 - (1 - g(\mathbf{q}))^K]$$

and

$$\text{var } \tilde{h}_K(\mathcal{E}, \mathbf{q}) = [g(\mathbf{q})/g_U(\mathbf{q})]h(\mathcal{E}, \mathbf{q})[1 - h(\mathcal{E}, \mathbf{q})]/K + o(K^{-1})$$

where $o(y)$ denotes a function of y such that $o(y) \rightarrow 0$ as $y \rightarrow 0$. To order $1/K$ this variance improves on the variance of a crude estimate by a factor of $g_U(\mathbf{q})$ demonstrating the efficiency of the importance sampling plan based on the bound information.

One can compute the bound $g_U(\mathbf{q})$ in time polynomial in b_i , $i \in \{\bigcup_{k=1}^I \mathcal{E}_i\} \cup \{\bigcup_{j=1}^J \mathcal{E}_j\}$ and sample from the distribution $\{Q(x, \mathbf{q})\}$ in (3) in time $O(|\mathcal{A}|)$.

The sampling plan described above is designed to provide an estimate for $h(\mathcal{E}, \mathbf{q})$ at a single reliability vector \mathbf{q} . Therefore, the estimation of the function $\{h(\mathcal{E}, \mathbf{q}), \mathbf{q} \in \mathcal{E}\}$ requires, in principle, $|\mathcal{E}|$ experiments. We now show how sampling data collected for estimating $h(\mathcal{E}, \mathbf{q})$ at a single point can be used to provide estimates for $h(\mathcal{E}, \mathbf{q})$ at all points $\mathbf{q} \in \mathcal{E}$.

Let \mathbf{p} be a point and define the *importance functions*

$$R(x, \mathbf{q}, \mathbf{p}) = P(x, \mathbf{q})/P(x, \mathbf{p})$$

$$x \in \mathcal{X}; \mathbf{q} \in \mathcal{E}$$

$$= \prod_{i=1}^a \left[\frac{q_i}{p_i} \right]^{x_i} \left[\frac{1 - q_i}{1 - p_i} \right]^{1 - x_i}.$$

Suppose one draws K independent samples $X^{(1)}, \dots, X^{(K)}$ from $\{Q(x, \mathbf{p})\}$. Then

$$\tilde{g}_{aK}(\mathbf{q}, \mathbf{p}) = g_U(\mathbf{p}) \frac{1}{K} \sum_{k=1}^K \phi(X^{(k)})R(X^{(k)}, \mathbf{q}, \mathbf{p})$$

and

$$\mathbf{q} \in \mathcal{E} \tag{8}$$

$$\tilde{g}_{bK}(\mathbf{q}, \mathbf{p}) = g_U(\mathbf{q}) - g_U(\mathbf{p}) \frac{1}{K} \sum_{k=1}^K [1 - \phi(X^{(k)})]R(X^{(k)}, \mathbf{q}, \mathbf{p})$$

are unbiased estimates of $g(\mathbf{q})$, $\mathbf{q} \in \mathcal{E}$ with

$$\begin{aligned}
K \text{ var } \tilde{g}_{aK}(\mathbf{q}, \mathbf{p}) &= K \text{ var } \tilde{g}_K(\mathbf{q}) \\
&\quad + \{c(\mathbf{q}, \mathbf{p})g_U(\mathbf{p})g_U(\mathbf{q}^*) - g_U(\mathbf{q})g(\mathbf{q})\}
\end{aligned}$$

and

$$\tag{9}$$

case bound $\max_{q \in \mathcal{S}} c(q, p) g_{ij}(p) g_{ij}(q^*)$. Other approaches, related directly to the accuracy of the estimates $\tilde{h}_{ijK}(\mathcal{G}, q)$, $i, j = a, b$ in (11), are currently under investigation.

Confidence intervals (individual and simultaneous) for $h(\mathcal{G}, q)$, $q \in \mathcal{S}$ are necessary for evaluating the accuracy of the estimates the proposed method produces. Since convergence to normality is not uniform in general, confidence intervals which are derived with exclusive use of statistical inequalities and are, therefore, valid for each finite sample size K deserve special attention. Such non normal confidence intervals are currently under development.

Finally, the general case of multiple capacity levels and set \mathcal{S} consisting of joint p.m.f.s. with independent marginals and common support is also under development by the authors of this paper.

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