

Limitations of Realistic Monte-Carlo Techniques

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Abstract

Data processing means applying some algorithm $f(x_1, \dots, x_n)$ to the values of the quantities x_1, \dots, x_n , resulting in a value $y = f(x_1, \dots, x_n)$. Values x_i usually come from measurements. Measurements are never absolutely accurate; the measurement result \tilde{x}_i is, in general, different from the actual (unknown) value x_i of the corresponding quantity: $\Delta x_i \stackrel{\text{def}}{=} \tilde{x}_i - x_i \neq 0$. Because of this, the computed value $\tilde{y} = f(\tilde{x}_1, \dots, \tilde{x}_n)$ is, in general, different from the ideal value $y = f(x_1, \dots, x_n)$. It is therefore desirable to estimate the accuracy $\Delta y \stackrel{\text{def}}{=} \tilde{y} - y$.

In many practical situations, the measurement errors are relatively small. In such cases, we can safely ignore terms which are quadratic or higher order in Δx_i , and conclude that $\Delta y = \sum_{i=1}^n c_i \cdot \Delta x_i$, where $c_i \stackrel{\text{def}}{=} \frac{\partial f}{\partial x_i}$.

When we know the probability distributions for all Δx_i (and we know that they are independent), then we can use Monte-Carlo techniques: several times $k = 1, \dots, N$, we simulate $\Delta x_i^{(k)}$, then the differences $\tilde{y} - f(\tilde{x}_1 - \Delta x_1^{(k)}, \dots, x_n - \Delta x_n^{(k)})$ have the same distribution as Δy .

Alternatively, we can use numerical differentiation to estimate all the derivatives c_i , and then use the above formula, but this would require $n + 1$ calls to the algorithm f , which for large n can be too long. In contrast, the Monte-Carlo method needs $N + 1$ calls, where N is determined only by the accuracy with which we want Δ (and does not depend on n).

In many practical situations, we only know the upper bound Δ_i on each measurement error Δx_i : $|\Delta x_i| \leq \Delta_i$. In this case, the only information that we have about the unknown (actual) value x_i is that it is in the interval $[\tilde{x}_i - \Delta_i, \tilde{x}_i + \Delta_i]$. In this case, the value Δy is bounded by $\Delta = \sum_{i=1}^n |c_i| \cdot \Delta_i$. A straightforward computation of Δ requires $n + 1$ calls to f , but there is a faster method based on using Cauchy distribution: we simulate $\Delta x_i^{(k)}$ based on Cauchy with parameter Δ_i , then the differences $\tilde{y} - f(\tilde{x}_1 - \Delta x_1^{(k)}, \dots, x_n - \Delta x_n^{(k)})$ are Cauchy distributed with the desired parameter Δ .

This method works, but its simulation is not realistic: we know that $|\Delta x_i| \leq \Delta_i$, but a Cauchy distribution goes beyond this bound. It has been known that if we consider simulations in which all $\Delta x_i^{(k)}$ are independent, then no realistic Monte-Carlo technique can always compute Δ . In this paper, we prove that this result holds even without the independence assumption, i.e., that the simulated values $\Delta x_i^{(k)}$ have to go beyond $[-\Delta_i, \Delta_i]$.