

On Completion Times of Networks of Concurrent and Sequential Tasks

Daniel Berleant, Jianzhong Zhang, and Gerald Sheblé
Department of Electrical and Computer Engineering
2215 Coover Hall
Iowa State University, Ames, Iowa 50011

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Contact:
Daniel Berleant
Department of Electrical and Computer Engineering
2215 Coover Hall
Iowa State University
Ames, Iowa 50011

Email: berleant@iastate.edu
Phone: (515) 294-3959
Fax: (515) 294-8432

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Iowa State University, Ames, Iowa 50011

Abstract

The problem of determining the time to complete multiple tasks that may proceed concurrently, sequentially, or both is considered. In the solution offered, each individual task completion time may be described with a number, interval, or distribution function. In the case of distribution functions, two task completion times might be independent random variables, as when the tasks are performed in different environments and proceed independently. Alternatively, completion times might be positively correlated, as when both depend on the quality of management and proceed within the same managerial environment, or they could be negatively correlated, as when resource sharing means that faster completion of one implies slower completion of the other. Finally, various factors might interact to make completion times dependent in a way that is difficult to characterize accurately. The solution offered avoids assuming that individual task completion times are independent or have any other dependency relationship. One application of the results is in project management, as in the context of PERT (Program Evaluation and Review Technique) diagrams.

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Iowa State University, Ames, Iowa 50011
berleant@iastate.edu

1 Introduction

Determining the time to complete all tasks in a network of tasks is easy when the time to complete each individual task in the network has a numerical value, harder when individual completion times are described using probability distributions, and still more challenging when these distributions are neither assumed independent nor assumed to have any other dependency relationship. A method is described here for determining completion times of task networks in the last case. We begin by describing each task completion time with a probability distribution function, noting that this includes as a special case a completion time described with a precise number since a number may be represented as a step distribution function (Figure 1, *left*). We later generalize to the case of left and right envelopes enclosing a family of cumulative distribution functions (CDFs) which, as a special case, allows a completion time to be represented as an interval describing a range of plausible values with high and low bounds but no information about the probability distribution within those bounds (Figure 1, *right*).

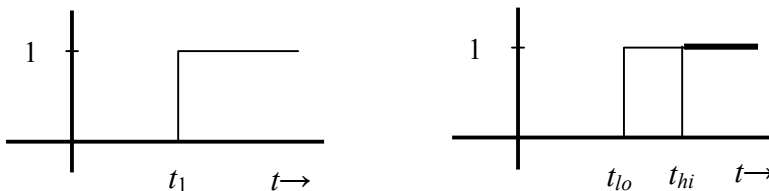


Figure 1. (*Left*) the numerical value of time t_1 is a special case of a cumulative distribution function (CDF) which is 0 below t_1 , and 1 at t_1 or above. (*Right*) an interval $[t_{lo}, t_{hi}]$ is a special case of a family of distributions containing any CDF which is 0 below t_{lo} and 1 (at or) above t_{hi} .

In real situations, two task completion times might be independent random variables, as when each is done in a different environment and they proceed independently. Alternatively, completion times might be highly positively correlated, as could occur if the tasks depend on the quality of management and proceed within the same managerial environment. As a third possibility, completion times might be quite negatively correlated, as could occur if the tasks proceed concurrently with shared personnel or other resources and faster completion of one entails slower completion of the other. A final and quite likely possibility is that various factors interact to make completion times dependent in a way that is difficult to characterize accurately. Therefore in the general case we wish to avoid assuming that individual task completion times are independent or have any other particular dependency relationship. A solution to this general case is offered.

The results have application to project management, where task completion time analyses can be useful as illustrated by the well-known PERT (Program Evaluation and Review Technique) method.

2 Solution for the case of two concurrent tasks

This section discusses the case of two concurrent tasks. Generalization to larger networks of tasks is discussed in Section 3.

Consider concurrent tasks X and Y , each beginning when the task environment is in a start state S and whose joint completion brings about desired finish state F (Figure 2). Let F_x be the CDF of random variable t_x , the completion time of task X , and let F_y be the CDF of random variable t_y , the completion time of task Y . We begin by reviewing solution strategies when t_x and t_y are independent, and then generalize by removing the independence assumption.

One solution strategy is the analytical one. The analytical approach to arithmetic on random variables is limited in the forms of the distributions it can handle and usually relies on the assumption of independence (e.g. Springer 1979). The Monte Carlo approach is a numerical

strategy that does not produce definite bounds, does not handle cases where one operand is a CDF and the other an interval except under severe restrictions, does not handle the case of unknown dependency between random variables, and has other limitations (Ferson 1996). Numerical convolution (Ingram et. al 1968; Colombo and Jaarsma 1980; Kaplan 1981) is an alternative numerical strategy that allows arithmetic operations to be applied to random variables with a wide variety of CDFs, and has been extended to capture discretization error via error bounds that propagate through the calculations and lead to left and right envelopes around the true solution (Williamson and Downs 1990; Berleant 1993; Cooper et al. 1996). See Figure 3.

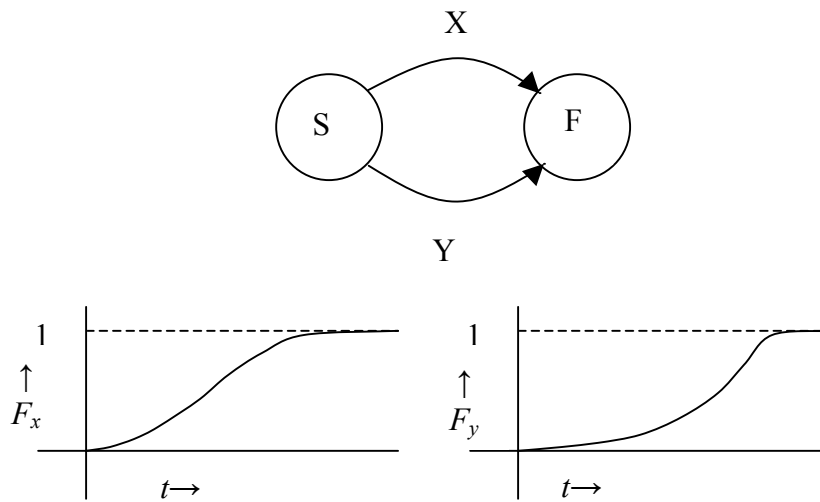


Figure 2. PERT diagram showing a starting state S, a finish state F, and two tasks X and Y that must be completed to reach state F. Two different distribution functions F_x and F_y describe random variables t_x and t_y , which represent the completion times of tasks X and Y.

Envelopes consist of non-crossing CDFs that enclose the paths of all CDFs consistent with the problem. These envelopes are often called *probability bounds* (Ferson et al. 1998) and, because they do not cross, the right envelope has *first order stochastic dominance* over the left (Levy 1992). Coarse discretizations for random variables t_x and t_y (e.g. Figure 3) lead to correspondingly large discretization error and therefore more widely spaced left and right envelopes. Finer discretizations would result in left and right envelopes that have more and

smaller steps and are closer together. The CDF for result t , the time to complete both tasks, must be some CDF enclosed by the left and right envelopes.

$$t = \max(t_x, t_y), \quad t_x \text{ and } t_y \text{ independent}$$

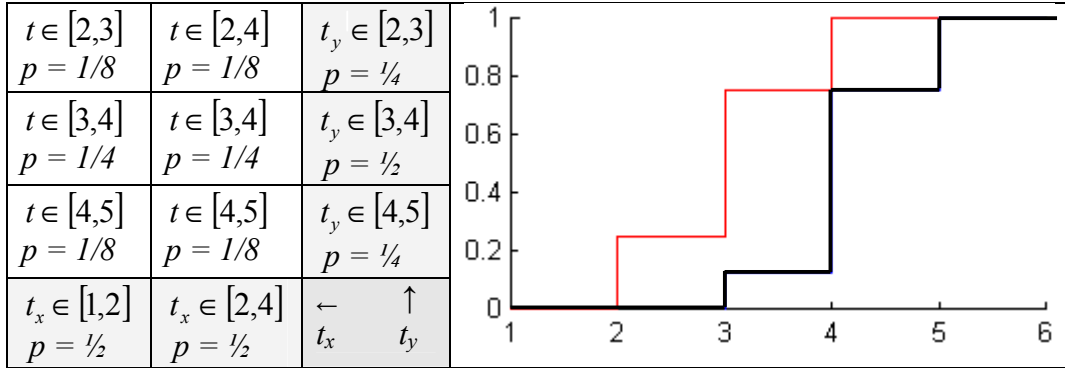


Figure 3. Random variable t_x is coarsely discretized (bottom row), and similarly for t_y (right hand column). The binary operation appropriate to the task completion problem is $\max(t_x, t_y)$ because, for any samples of t_x and t_y , both tasks are complete when the one that finishes last is complete. The distribution of joint completion times is implicit in the set of interior cells (*unshaded*) of the joint distribution table, each of which is calculated from its corresponding marginal cells. For example, the upper left cell contains probability mass $1/8$, which is the product of the probabilities of its marginal cells in the right hand column and bottom row, $1/4$ and $1/2$ respectively. The product is used because t_x and t_y are assumed independent (this assumption will be relaxed later). The upper left cell contains the interval $[2,3]$ because its marginal cells have task X complete in time $[1,2]$ and Y in time $[2,3]$, so the time to complete both could potentially be anywhere within that interval. The cumulation over t of the interior cells is bounded by the left and right envelopes shown, with the separation between the envelopes due to the undetermined distribution of each cell's mass across its interval which could, in extreme cases, be concentrated at the interval low or high bound (Berleant 1993).

Left and right envelopes are each derived from a joint distribution table such as that shown in Figure 3. The probability mass shown associated with each interior cell of a joint distribution

table is the product of the probability masses in its corresponding marginal cells if the operands are independent, but relaxing the assumption of independence leaves them undetermined. Therefore when the dependency relationship between the operands is unknown, the process illustrated in Figure 3 requires significant modification (Berleant and Goodman-Strauss 1998). Regardless of the dependency relationship between the marginals, the masses of the interior cells are constrained to some extent by the marginals, which require the masses of all the interior cells in a row to sum to the mass of the marginal cell at the right of that row, and the masses of the interior cells in a column to sum to the mass of the corresponding marginal cell at the bottom of that column. Consequently the summed mass of any particular subset of interior cells will typically have a range of possible values, and for a properly chosen subset the maximum or minimum of this range yields a point on the left or right envelope. More specifically, obtaining the height of the left envelope at time t requires maximizing the collective probability mass of the interior cells whose intervals have low bounds below (or equal to) t subject to the row and column constraints, because the mass of each of those cells either may (if the interval's high bound is above t) or must (if the interval's high bound is not above t) be in the cumulation at t . The process is analogous for finding the height of the right envelope: minimize the sum of the probability masses of the interior cells whose intervals have high bounds below or equal to t (Berleant and Goodman-Strauss 1998). Figure 4 explains the process, which can be done by hand for a very small table although in the general case linear programming (LP) is more practical. The left and right envelopes have staircase-like forms. In Figure 4, for example, the heights of the left and right envelopes at $t=3.5$ hold for all other values of t between 3 and 4. Because for staircase-shaped curves the heights for only a limited number of values of t need to be found to fully characterize the envelopes, the number of LP problems is correspondingly limited. Figure 4 also shows the full envelopes.

Figure 4. An example.

Each interior cell interval in the following joint distribution table has bounds defined by $\max(t_x, t_y)$ for its associated (shaded) marginal cell intervals. While interior cell probabilities are constrained by the marginal cell probabilities, they are not fully determined because no assumptions are made about the dependency relationship between t_x and t_y .

$t \in [2,3]$ p_{11}	$t \in [2,4]$ p_{12}	$t_y \in [2,3]$ $p = 1/4$
$t \in [3,4]$ p_{21}	$t \in [3,4]$ p_{22}	$t_y \in [3,4]$ $p = 1/2$
$t \in [4,5]$ p_{31}	$t \in [4,5]$ p_{32}	$t_y \in [4,5]$ $p = 1/4$
$t_x \in [1,2]$ $p = 1/2$	$t_x \in [2,4]$ $p = 1/2$	$\leftarrow \uparrow$ $t_x \quad t_y$

Consider for example the cumulative probability of t at 3.5. Bolded probabilities masses p_{11} , p_{12} , p_{21} , and p_{22} can contribute to the left envelope of t at 3.5, because the low bounds of the intervals in those cells are ≤ 3.5 . Therefore those probabilities could all be in the cumulation at $t=3.5$, and in the extreme case that p_{12} , p_{21} , & p_{22} happen to be concentrated at the low bounds of their intervals, will be (and to find points on the envelopes, we are interested in extreme cases). To maximize this cumulation of p_{11} , p_{12} , p_{21} , and p_{22} , their sum must be maximized (at the expense of non-bolded probabilities p_{31} and p_{32}), yielding $p_{11}+p_{12}+p_{21}+p_{22}=3/4$ as shown in the following solution:

$t \in [2,3]$ $p_{11}=1/4$	$t \in [2,4]$ $p_{12}=0$	$t_y \in [2,3]$ $p = 1/4$
$t \in [3,4]$ $p_{21}=0$	$t \in [3,4]$ $p_{22}=1/2$	$t_y \in [3,4]$ $p = 1/2$
$t \in [4,5]$ $p_{31}=1/4$	$t \in [4,5]$ $p_{32}=0$	$t_y \in [4,5]$ $p = 1/4$
$t_x \in [1,2]$ $p = 1/2$	$t_x \in [2,4]$ $p = 1/2$	$\leftarrow \uparrow$ $t_x \quad t_y$

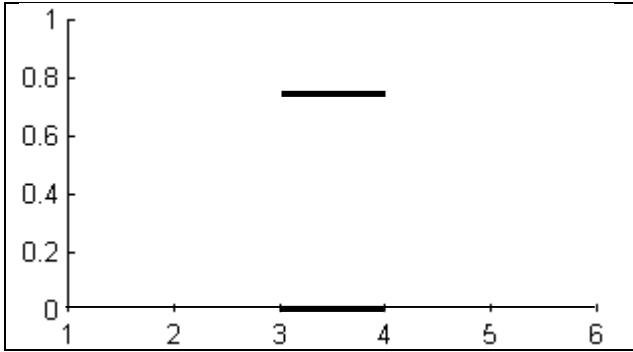
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(Figure 4 continued)

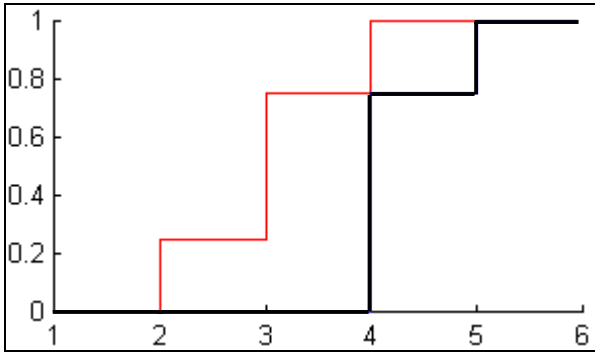
For the other envelope, the (unary “sum” of) italicized probability mass p_{1l} is minimized, yielding 0 as shown in the following solution:

$t \in [2,3]$ $p_{11}=\mathbf{0}$	$t \in [2,4]$ $p_{12}=\mathbf{1/4}$	$t_y \in [2,3]$ $p = \mathbf{1/4}$
$t \in [3,4]$ $p_{21}=\mathbf{1/2}$	$t \in [3,4]$ $p_{22}=\mathbf{0}$	$t_y \in [3,4]$ $p = \mathbf{1/2}$
$t \in [4,5]$ $p_{31}=\mathbf{0}$	$t \in [4,5]$ $p_{32}=\mathbf{1/4}$	$t_y \in [4,5]$ $p = \mathbf{1/4}$
$t_x \in [1,2]$ $p = \mathbf{1/2}$	$t_x \in [2,4]$ $p = \mathbf{1/2}$	$\leftarrow \uparrow$ $t_x \quad t_y$

These maximum and minimum cumulations of $3/4$ and 0 hold not only for $t=3.5$ but also for all other t from 3 to 4, because no interior cell has an interval with an endpoint in that range, as graphed next.



Repeating this process for appropriate values of t yields the following full envelopes around $t=\max(t_x, t_y)$.



Although the marginals used here are the same as in Figure 3, the envelopes are farther apart because the dependency between the random variables is unspecified, so the inferential power of the independence assumption is absent. The discretization, coarse in this example, also affects the degree of separation of

the envelopes. Finer discretization would yield smaller steps in the envelopes and hence envelopes that are, on average, closer together.

Figure 4 (end).

When linear programming is applied to minimization and maximization problems of this type the objective function is the sum of the probabilities of the subset of interior cells to be maximized or minimized, and the constraint set consists of one for each row and one for each column. A general-purpose linear programming algorithm such as the simplex method can be used, but a faster choice is the transportation simplex method, which applies to certain problems such as these containing only row and column constraints.

To apply the transportation simplex method to optimize the distribution of probability masses across interior cells, the cost coefficients of the cells in the subset whose probability mass is to be maximized or minimized are set to one, the cost coefficients of the remaining cells are set to zero, and the allocations of the cells are their probability masses. In our software implementation, problems involving generating envelopes from a 16x16 joint distribution table require approximately 14 seconds using the simplex algorithm but only 1 second using the transportation simplex algorithm, on a Pentium III PC running Windows NT.

3 Generalizing the solution to networks of concurrent and sequential tasks

Extending the approach from two concurrent tasks to larger networks of tasks requires solving three problems: (1) determining the completion time of two tasks that run not concurrently but sequentially, (2) determining the completion time of three or more concurrent tasks, and (3) using results as inputs to obtain further downstream results. These problems may be solved as follows.

- (1) To determine the completion time of two sequential tasks, their individual completion times are added, because one completes and then the next begins. To add them, the same procedure that was described earlier for concurrent tasks is applied except that the intervals in the interior cells of the joint distribution table are obtained by performing

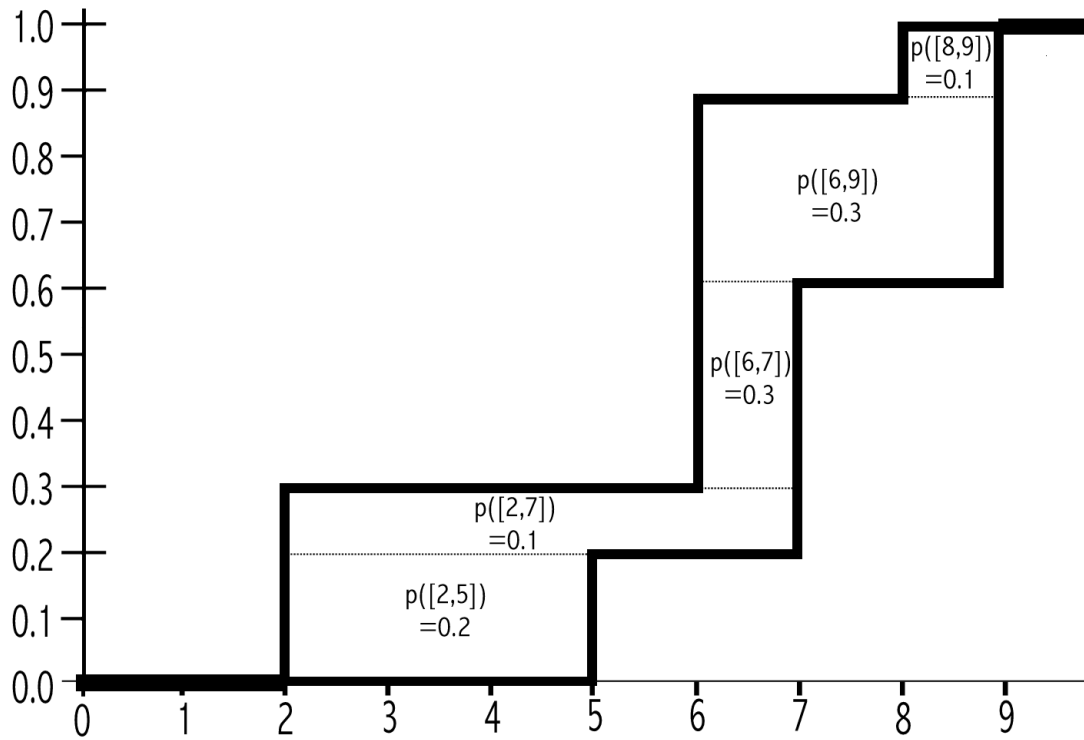
t_x+t_y instead of $\max(t_x,t_y)$. Thus for each joint distribution tables in Figure 4, the top left cell would contain the interval $[3,5]=[1,2]+[2,3]$ instead of $[2,3]=\max([1,2],[2,3])$.

- (2) To handle three concurrent tasks, the result for two of them is calculated, and that result used as the completion time for a composite task that proceeds concurrently with the third task. In other words, for concurrent tasks X, Y, & Z, we wish to calculate $\max(\max(x,y), z)$. This is a case of using intermediate results as inputs, discussed next.
- (3) To use a result as an input to another calculation, we must convert a pair of left and right envelopes, which is what a result looks like, into a set of intervals and associated probability masses, which is what a marginal in a joint distribution table looks like. To convert, first note that the envelopes consist of horizontal and vertical line segments. This allows the space they enclose to be partitioned into a stack of rectangles (Figure 5, *top*). Each rectangle defines an interval whose low bound is a value on the horizontal axis at which there is a vertical segment of the left envelope (forming the left side of the rectangle), and whose high bound is a value on the horizontal axis at which there is a vertical segment of the right envelope (forming the right side of the rectangle). The mass of the interval is the increment in the cumulative probability represented by the (bottom-to-top) height of the rectangle. The result of this partition process is a set of intervals and their associated probabilities, usable as a marginal in a joint distribution table for another arithmetic operation (Figure 5, *bottom*).

4 Using inferences from result envelopes

Consider three types of inference that may be drawn from a pair of left and right envelopes.

- 1) The probability of finishing all the tasks by some time T_0 is at least P_0 in Figure 6. Similarly, the probability of not finishing by time T_0 is at least $(1-P_1)$.



$$t = \max(t_z, t_w)$$

$t \in [2,5]$ $p =$	$t \in [2,7]$ $p =$	$t \in [6,7]$ $p =$	$t \in [6,9]$ $p =$	$t \in [8,9]$ $p =$	$t_w \in [2,3]$ $p = 0.25$
$t \in [3,5]$ $p =$	$t \in [3,7]$ $p =$	$t \in [6,7]$ $p =$	$t \in [6,9]$ $p =$	$t \in [8,9]$ $p =$	$t_w \in [3,4]$ $p = 0.5$
$t \in [4,5]$ $p =$	$t \in [4,7]$ $p =$	$t \in [6,7]$ $p =$	$t \in [6,9]$ $p =$	$t \in [8,9]$ $p =$	$t_w \in [4,5]$ $p = 0.25$
$t_z \in [2,5]$ $p = 0.2$	$t_z \in [2,7]$ $p = 0.1$	$t_z \in [6,7]$ $p = 0.3$	$t_z \in [6,9]$ $p = 0.3$	$t_z \in [8,9]$ $p = 0.1$	$\leftarrow \begin{matrix} \uparrow \\ t_z \quad t_w \end{matrix}$

Figure 5. Staircase shaped envelopes partitioned into a set of intervals and masses (*top*).

These might represent a random variable $t_z = \max(t_x, t_y)$, used as a marginal in the last row of a joint distribution table (*bottom*), and combined with the concurrent completion time t_w of some other task W. The interior cell probabilities of the table are undetermined since no dependency relationship was defined between the marginals, and so cannot be given values.

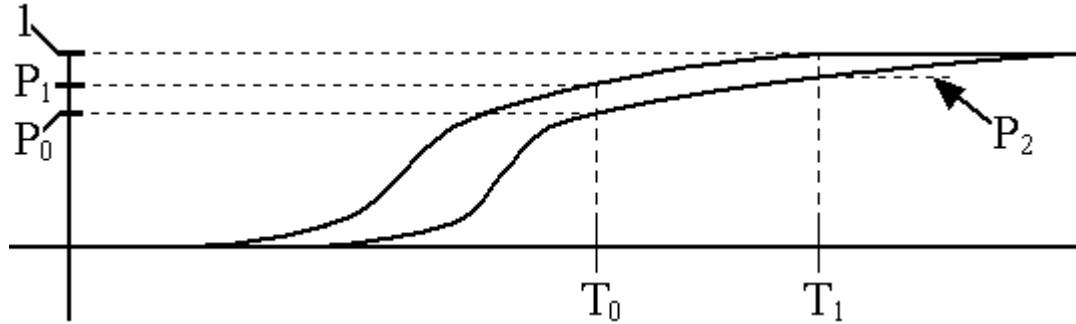


Figure 6. Left and right envelopes associate probability intervals with time points. If the envelopes describe cumulative probability of task completion over time, then the probability of completion by time T_0 is within the interval $[P_0, P_1]$, and by time T_1 , $[P_2, 1]$.

- 2) Suppose that $p(\text{some outcome}) \in [P, 1]$. For example in Figure 6, $p(\text{task completion by time } T_1) \in [P_2, 1]$. The interval $[P_2, 1]$ is qualitatively different from a point estimate somewhere between P_2 and 1 that would derive from an analysis that produced a single distribution function instead of left and right envelopes. This is because, unlike a point estimate, $p \in [P_2, 1]$ indicates the plausibility of two distinct scenarios with different implications, (1) certain completion (within the model limits), and (2) uncertain completion. Decisions about resource allocation on the overall project or about deadlines to contract for could depend on which scenario is correct, yet the implied opportunity to seek further information to enable discriminating, or at least to reduce the second order uncertainty in completion time would not be available from an analysis that produced a point probability estimate.
- 3) Consider the problem of determining the probability that one task will finish later than another, $p(t_y > t_x)$. The probability of one task or path taking longer than another is relevant in such applications as project management where task networks represent PERT diagrams describing the prerequisite structure of tasks in a project. A simple example is

two tasks that begin at the same time, as in Figure 1. A generalization is two tasks embedded in a network of tasks, such as in Figure 7 for final tasks CF and EF. In the generalization the tasks need not start at the same time, and the times at which they complete depend on both the task itself and any prerequisite tasks in the network. These prerequisite tasks may form a simple sequence as in the case of task EF with prerequisite partial path SDE, or contain concurrency as in the case of task CF with prerequisite, concurrent, partial paths SAC and SBC.

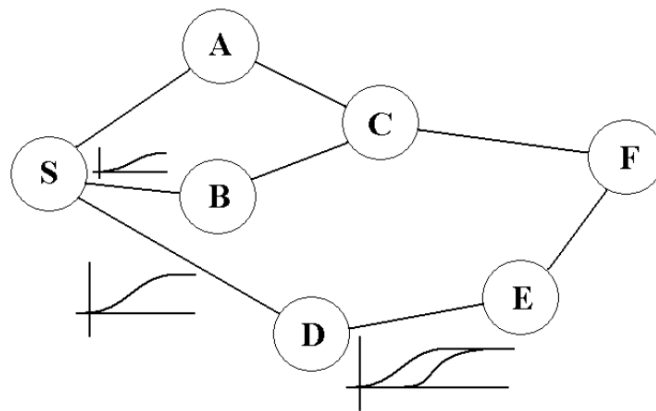


Figure 7. A network of tasks. The times to complete tasks SB and SD are shown as cumulative distributions. The time to reach state E is the sum of the times to complete SD and DE, and if the dependency relationship between the completion times for SD and DE are unknown the sum is a pair of envelopes rather than a single cumulative distribution.

Solving this type of problem requires determining $p(t_y > t_x)$, where t_x and t_y are sample values of random variables for the time points at which two tasks X and Y, or CF and EF, etc., complete. To do this, and relate it to standard techniques, we first provide a continuous solution for the case of independent distributions, then give the discrete form of the solution, then an intervalized discrete form, and finally remove the independence assumption.

In the case of a continuous solution for independent distributions, if the density functions of the task completion times are $f_x(t)$ and $f_y(t)$ and sample completion times are t_x and t_y , then

$$(1) \quad p(t_y > t_x) = \int_{t=-\infty}^{\infty} \left(f_y(t) dt \int_{-\infty < t_0 < t} f_x(t_0) dt \right).$$

Intuitively, $\int_{-\infty < t_0 < t} f_x(t_0) dt$ is the area under f_x over all times earlier than some given time t ,

which is $p(t > t_x)$, or the probability that t is later than the completion time t_x of task X. The probability that the completion time of task Y is within a time period centered at t with width dt is $p(t_y \in t \pm \frac{1}{2} dt) = f_y(t) dt$. The probability of both $(t > t_x)$ and $(t_y \in t \pm \frac{1}{2} dt)$ is

therefore the product of their individual probabilities, $f_y(t) dt \int_{-\infty < t_0 < t} f_x(t_0) dt$, and

integrating this expression over all possibilities for t gives equation (1).

Discretizing (1) gives $p(t_y > t_x) = \sum_{t=-\infty}^{\infty} \left(f_y(t) \Delta t \sum_{-\infty < t_0 < t} f_x(t_0) \Delta t \right)$ for values of t and

t_0 spaced Δt apart. This can be intervalized, bounding the discretization error and giving

$$(2) \quad p(t_y > t_x) = \left[\sum_{T_y} \left(p(T_y) \sum_{T_x, \underline{T}_y > \underline{T}_x} p(T_x) \right), \sum_{T_y} \left(p(T_y) \sum_{T_x, \overline{T}_y > \overline{T}_x} p(T_x) \right) \right]$$

where the T_x and T_y are intervals over t_x and t_y , such as might appear in the marginals of a joint distribution table, $p(T_x)$ and $p(T_y)$ are their associated probability masses, and $\underline{T}_x, \underline{T}_y, \overline{T}_x$ and \overline{T}_y are their low and high bounds.

As an example of equation (2) consider the joint distribution table in Figure 8. The low bound of $p(t_y > t_x)$ is the sum of the probability masses of cells labeled True, which is 0.789. The high bound is the sum of the masses of cells labeled True or Uncertain, which is 0.939, yielding $p(t_y > t_x) \in [0.789, 0.939]$.

To remove the independence assumption, the masses of the interior cells are reapportioned among the interior cells within the limits imposed by the row and column constraints using linear programming to minimize the summed masses of the cells labeled True, giving a low bound of 0.61, and then reapportioned again to maximize the summed masses of the cells labeled True or Uncertain, giving a high bound of 1. The new result, $p(t_y > t_x) \in [0.61, 1]$, as expected is wider than the earlier result of $p(t_y > t_x) \in [0.789, 0.939]$, which benefited from assuming independence.

$t_y > t_x$, t_x and t_y independent

True p=.005	True p=.006	True p=.008	True p=.01	True p=.021	Uncertain p=.021	Uncertain p=.01	False p=.008	False p=.006	False p=.005	[10.1,11.1] p=.1
True p=.01	True p=.012	True p=.016	True p=.02	True p=.042	True p=.042	Uncertain p=.02	Uncertain p=.016	False p=.012	False p=.01	[11.1,12.1] p=.2
True p=.02	True p=.024	True p=.032	True p=.04	True p=.084	True p=.084	True p=.04	Uncertain p=.032	Uncertain p=.024	False p=.02	[12.1,13.1] p=.4
True p=.01	True p=.012	True p=.016	True p=.02	True p=.042	True p=.042	True p=.02	True p=.016	Uncertain p=.012	Uncertain p=.01	[13.1,14.1] p=.2
True p=.005	True p=.006	True p=.008	True p=.01	True p=.021	True p=.021	True p=.01	True p=.008	True p=.006	Uncertain p=.005	[14.1,15.1] p=.1
[5,6] p=.05	[6,7] p=.06	[7,8] p=.08	[8,9] p=.1	[9,10] p=.21	[10,11] p=.21	[11,12] p=.1	[12,13] p=.08	[13,14] p=.06	[14,15] p=.05	← ↑ t_x t_y

Figure 8. Joint distribution table representing $t_y > t_x$, for independent t_x and t_y . Each interior cell is labeled True if $t_y > t_x$ for t_y and t_x in the intervals of the marginal cells of that interior cell, False if instead $t_y < t_x$, and Uncertain if the marginal cell intervals overlap (indicating that the unspecified details of the distributions of the marginal cell masses over their intervals determine whether $t_y > t_x$ for all, some, or none of the interior cell mass).

To restate an example, this process could be used to bound the probability that the completion time of task X will be later than that of task Y in a PERT diagram conforming to Figure 1. The process could also be used in a more complex example such as bounding the probability that task CF will complete later than task EF in Figure 7. The completion time of each of these tasks will be in the form of envelopes, which when

converted to marginals will have overlapping intervals as in Figure 5. However any overlap is irrelevant to equation (2), which justifies Figure 8. Ultimately such results could support management decisions about resource allocation intended to optimize the overall completion time of the entire project.

5 Software

Crystal Ball (www.decisioneering.com) and @risk (www.palisade.com) are well-known commercial products that rely on Monte Carlo simulation, thereby inheriting the shortcomings of Monte Carlo simulation noted earlier in Section 2.

RiskCalc (Ferson et al. 1998) is a commercially available package that can do the operations on random variables used here, although its algorithm (Williamson and Downs, 1990) is different and more complicated than the one used here, some further details of which have been described by Berleant and Goodman-Strauss (1998). Our software, Statool, is downloadable from <http://www.public.iastate.edu/~berleant/statool.html>. Both systems run under Windows on a PC. Statool requires Visual Basic to be installed.

6 Conclusion

We have shown how to solve a simply stated problem with significant implications: determining completion times of networks of tasks in the absence of assumptions about both the forms of distribution functions and their independence or other dependency relationships. Results are left and right envelopes bounding the space of plausible CDFs. Completion times of individual tasks may be expressed as numbers, intervals, distribution functions, or left and right envelopes.

Real problems frequently pose a variety of uncertainties. Therefore methods for obtaining results with minimal assumptions and while accounting for uncertainty remain an important area of investigation.

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8 References

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