### Runtime distributions

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#### Overview of talk

- Time-to-target plots
- Runtime distribution of GRASP
- Comparing algorithms with exponential runtime distributions
- Comparing algorithms with general runtime distributions
- Numerical applications to sequential algorithms
  - ▶ DM-D5 and GRASP algorithms for server replication
  - Multistart and tabu search algorithms for routing and wavelength assignment
  - GRASP algorithms for 2-path network design
- Comparing and evaluating parallel algorithms

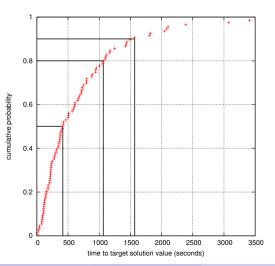
#### Introduction

- Runtime distributions or time-to-target plots display on the ordinate axis the probability that an
  algorithm will find a solution at least as good as a given target value within a given running time,
  shown on the abscissa axis.
- They provide a very useful tool to characterize the running times of stochastic algorithms for combinatorial optimization problems and to compare different algorithms or strategies for solving a given problem.
- They have been widely used as a tool for algorithm design and comparison.

### Time-to-target plots

- ullet Let  ${\mathcal P}$  be an optimization problem and  ${\mathcal H}$  a randomized heuristic for this problem.
- Furthermore, let  $\mathcal{I}$  be a specific instance of  $\mathcal{P}$  and let *look4* be a solution cost *target value* for this instance.
- Heuristic  $\mathcal{H}$  is run N times on the fixed instance  $\mathcal{I}$  and the algorithm is made to stop as soon as a solution whose objective function is at least as good as the given target value *look4* is found.
- For each of the N runs, the random number generator used in the implementation of the heuristic is initialized with a distinct seed and, therefore, the runs are assumed to be independent.
- The solution time of each run is recorded and saved.
- To compare their empirical and theoretical distributions, we follow a standard graphical methodology for data analysis.
- This methodology is used to produce the time-to-target plots (TTT-plots).

### Time-to-target plots



- After concluding the N independent runs, solution times are sorted in increasing order.
- The *i*-th sorted solution time  $t_i$  is associated with a probability  $p_i = (i 1/2)/N$ , and the points  $z_i = (t_i, p_i)$ , for i = 1, ..., N, are plotted.
- The figure illustrates this estimated cumulative probability distribution plot for problem P, a GRASP heuristic H, instance I, and target look4.
- Observation: The probability that the heuristic finds a solution at least as good as the target value in at most 416 seconds is about 50%, in at most 1064 seconds is about 80%, and in at most 1569 seconds is about 90%.

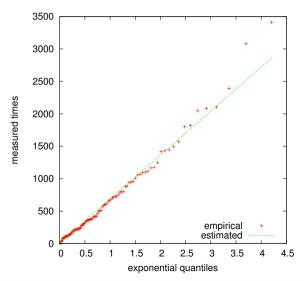
- The previous plot appears to fit an exponential distribution, or more generally, a shifted exponential distribution.
- To estimate the parameters of this two-parameter exponential distribution, we first draw the theoretical quantile-quantile plot (or Q-Q plot) for the data.
- To describe Q-Q plots, we recall that the cumulative distribution function for the two-parameter exponential distribution is given by:

$$F(t) = 1 - e^{-(t-\mu)/\lambda}$$

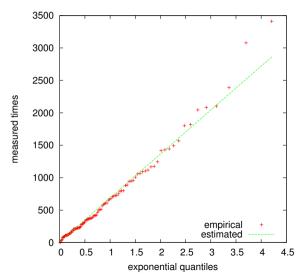
- $\triangleright$   $\lambda$  is the mean of the distribution data (and also is the standard deviation of the data).
- $\blacktriangleright$   $\mu$  is the shift of the distribution with respect to the ordinate axis.

- The *quantiles* of the data of an empirical distribution are derived from the (sorted) raw data, which in our case are N measured (sorted) running times.
- Quantiles are cutpoints that group a set of sorted observations into classes of equal (or approximately equal) size.
- For each value  $p_i$ , i = 1, ..., N, we associate a  $p_i$ -quantile  $q(p_i)$  of the theoretical distribution.
- For each  $p_i$ -quantile we have, by definition, that  $F((q(p_i)) = p_i$ .
- Hence,  $q(p_i) = F^{-1}(p_i)$  and therefore, for the two-parameter exponential distribution, we have  $q(p_i) = -\lambda \cdot \ln(1-p_i) + \mu$ .
- Note that if we were to use  $p_i = 1/N$ , for i = 1, ..., N, then  $q(p_N)$  would be undefined.

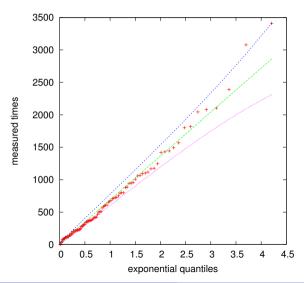
- A theoretical quantile-quantile plot (or theoretical Q-Q plot) is obtained by plotting the quantiles of the data of an empirical distribution against the quantiles of a theoretical distribution. This involves three steps:
  - First, the data (in this case, the measured solution times) are sorted in ascending order.
  - Second, the quantiles of the theoretical exponential distribution are obtained.
  - Finally, a plot of the data against the theoretical quantiles is made.
- In a situation where the theoretical distribution is a close approximation of the empirical distribution, the points in the Q-Q plot will have a nearly straight configuration.
- In a plot of the data against a two-parameter exponential distribution with  $\lambda=1$  and  $\mu=0$ , the points would tend to follow the line  $y=\hat{\lambda}\cdot x+\hat{\mu}$ .
- Consequently, parameters  $\lambda$  and  $\mu$  of the two-parameter exponential distribution can be estimated, respectively, by the slope  $\hat{\lambda}$  and the intercept  $\hat{\mu}$  of the line depicted in the Q-Q plot.



- The Q-Q plot shown is obtained by plotting the measured times in the ordinate against the quantiles of a two-parameter exponential distribution with  $\lambda=1$  and  $\mu=0$  in the abscissa, given by  $q(p_i)=-\ln(1-p_i)$ , for  $i=1,\ldots,n$ .
- To avoid possible distortions caused by outliers, we do not estimate the distribution mean with the data mean or by linear regression on the points of the Q-Q plot.
- Instead, we estimate the slope  $\hat{\lambda}$  of the line  $y = \lambda \cdot x + \mu$  using the upper quartile  $q_u$  and lower quartile  $q_l$  of the data.



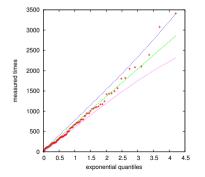
- The upper quartile  $q_u$  and lower quartile  $q_l$  are, respectively, the q(1/4) and q(3/4) quantiles.
- We take  $\hat{\lambda} = (z_u z_l)/(q_u q_l)$  as an estimate of the slope, where  $z_u$  and  $z_l$  are the u-th and l-th points of the ordered measured times, respectively.
- This informal estimation of the distribution of the measured data mean is robust since it will not be distorted by a few outliers
- Consequently, the estimate for the shift is  $\hat{\mu} = z_l \hat{\lambda} q_l.$



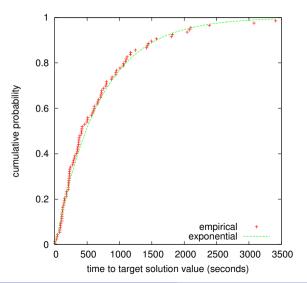
- To analyze the straightness of the Q-Q plots, we superimpose them with variability information.
- For each plotted point, we show plus and minus one standard deviation in the vertical direction from the line fitted to the plot.
- An estimate of the standard deviation for point  $z_i$ ,  $i=1,\ldots,n$ , of the Q-Q plot is  $\hat{\sigma}=\hat{\lambda}[p_i/(1-p_i)n]^{\frac{1}{2}}$ .
- The figure shows an example of a Q-Q plot with superimposed variability information.

- When observing a theoretical quantile-quantile plot with superimposed standard deviation information, one should avoid turning such information into a formal test.
- One important fact that must be kept in mind is that the natural variability of the data generates departures from the straightness, even if the model of the distribution is valid.
- The most important reason for portraying standard deviation is that it gives us a sense of the relative variability of the points in the different regions of the plot.
- However, since one is trying to make simultaneous inferences from many individual inferences, it is difficult to use standard deviations to judge departures from the reference distribution.
- For example, the probability that a particular point deviates from the reference line by more than two standard deviations is small.
- However, the probability that any of the points deviates from the line by two standard deviations is probably much greater.

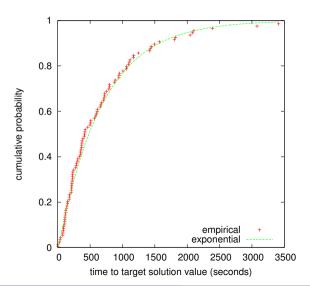
- In order statistics, this is made more difficult by the high correlation that exists between neighboring points.
- If one plotted point deviates by more than one standard deviation, there is a good chance that a whole bunch of them will too.



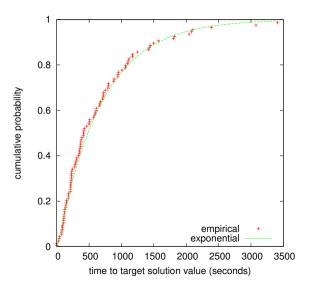
- Another point to keep in mind is that standard deviations vary substantially in the Q-Q plot.
- As one can observe in the previous Q-Q plot, the standard deviation of the points near the high end is substantially larger than the standard deviation of the points near the other end.



- Once the two parameters of the distribution have been estimated, a superimposed plot of the empirical and theoretical distributions can be made.
- The figure depicts the superimposed empirical and theoretical distributions corresponding to the Q-Q plot in the previous figure.
- The runtime distribution of a pure GRASP heuristic has been shown experimentally to behave as a random variable that fits an exponential distribution.



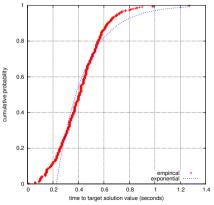
- However, in the case of more elaborate heuristics where setup times are not negligible, the runtimes fit a two-parameter or shifted exponential distribution.
- Therefore, the probability density function of the time-to-target random variable is given by  $f(t) = (1/\lambda) \cdot e^{-t/\lambda}$  in the first case (exponential distribution) and by  $f(t) = (1/\lambda) \cdot e^{-(t-\mu)/\lambda}$  in the second (shifted exponential distribution), with the parameters  $\lambda \in \mathbb{R}^+$  and  $\mu \in \mathbb{R}^+$  being associated with the shape and the shift of the exponential function, respectively.



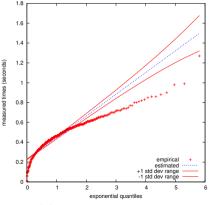
- The figure illustrates this result, depicting the superimposed empirical and theoretical distributions observed for an instance of the maximum covering problem where one wants to choose 500 out of 1000 facility locations such that, of the 10,000 customers, the sum of the weights of those that are covered is maximized.
- The best known solution for this instance is 33,343,542 and the target solution value used was 33,339,175 (about 0.01% off of the best known solution).

- However, if path-relinking is applied as an intensification step at the end of each GRASP iteration, then the iterations are no longer independent and the memoryless characteristic of GRASP is destroyed.
- This also happens in the case of cooperative parallel implementations of GRASP.
- Consequently, the time-to-target random variable may not fit an exponential distribution in such situations.
- This result is illustrated by two implementations of GRASP with bidirectional path-relinking:
  - ▶ The first is an application to the 2-path network design problem.
  - ▶ The second is an application to the three-index assignment problem.

The figure depicts the runtime distribution and the corresponding quantile-quantile plot for GRASP with bidirectional path-relinking of an instance of the 2-path network design problem with 80 nodes and 800 origin-destination pairs, with target set to 588.



(a) Runtime distribution

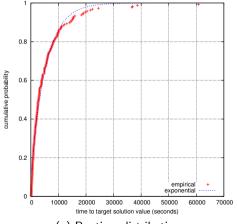


(b) Quantile-quantile plot

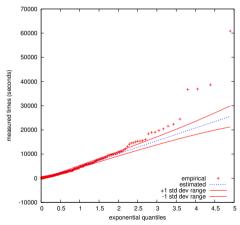
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The figure shows the runtime distribution and the corresponding quantile-quantile plot for GRASP with bidirectional path-relinking on Balas and Saltzman problem 22.1, with the target value set to 8.

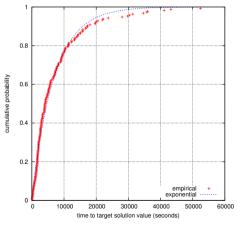


(a) Runtime distribution

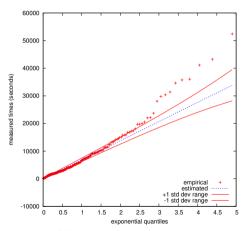


(b) Quantile-quantile plot

The figure shows the runtime distribution and the corresponding quantile-quantile plot for GRASP with bidirectional path-relinking on Balas and Saltzman problem 24.1, with the target value set to 7.



(a) Runtime distribution



(b) Quantile-quantile plot

#### Observations:

- For both heuristics and these three example instances, we observe that points steadily deviate by
  more than one standard deviation from the estimate for the upper quantiles in the quantile-quantile
  plots.
  - That is, many points associated with large computation times fall outside the plus or minus one standard deviation bounds.
- Therefore, we cannot say that these runtime distributions are exponentially distributed.

- We assume the existence of two randomized algorithms  $A_1$  and  $A_2$  for the approximate solution of some optimization problem.
- Furthermore, we assume that their solution times fit exponential (or shifted exponential) distributions.
- We denote by  $X_1$  (resp.  $X_2$ ) the continuous random variable representing the time needed by algorithm  $A_1$  (resp.  $A_2$ ) to find a solution as good as a given target value:

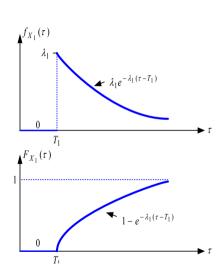
$$X_1 \mapsto \left\{ egin{array}{ll} 0, & au < T_1 \ \lambda_1 \cdot e^{-\lambda_1( au - T_1)}, & au \geq T_1 \end{array} 
ight.$$

and

$$X_2 \mapsto \left\{ egin{array}{ll} 0, & au < T_2 \ \lambda_2 \cdot e^{-\lambda_2( au - T_2)}, & au \geq T_2 \end{array} 
ight.$$

where  $T_1$ ,  $\lambda_1$ ,  $T_2$ , and  $\lambda_2$  are parameters ( $\lambda_1$  and  $\lambda_2$  define the shape of each shifted exponential distribution, whereas  $T_1$  and  $T_2$  denote by how much each of them is shifted).

The figure depicts the cumulative probability distribution and the probability density function of the random variable  $X_1$ .



- Since both algorithms stop when they find a solution at least as good as the target, we can say that algorithm  $A_1$  performs better than  $A_2$  if the former stops before the latter.
- Therefore, we must evaluate the probability  $Pr(X_1 \leq X_2)$  that the random variable  $X_1$  takes a value smaller than or equal to  $X_2$ .
- Conditioning on the value of  $X_2$  and applying the total probability theorem, we obtain

$$Pr(X_{1} \leq X_{2}) = \int_{-\infty}^{\infty} Pr(X_{1} \leq X_{2} | X_{2} = \tau) \cdot f_{X_{2}}(\tau) \cdot d\tau =$$

$$= \int_{T_{2}}^{\infty} Pr(X_{1} \leq X_{2} | X_{2} = \tau) \cdot \lambda_{2} \cdot e^{-\lambda_{2}(\tau - T_{2})} \cdot d\tau = \int_{T_{2}}^{\infty} Pr(X_{1} \leq \tau) \cdot \lambda_{2} \cdot e^{-\lambda_{2}(\tau - T_{2})} \cdot d\tau.$$

Let  $\nu = \tau - T_2$ . Then,  $d\nu = d\tau$  and

$$Pr(X_1 \le X_2) = \int_0^\infty Pr(X_1 \le (\nu + T_2)) \cdot \lambda_2 \cdot e^{-\lambda_2 \nu} \cdot d\nu. \tag{1}$$

Using the formula of cumulative probability function of the random variable  $X_1$  (see previous figure ), we obtain

$$Pr(X_1 \le (v + T_2)) = 1 - e^{-\lambda_1(v + T_2 - T_1)}.$$
 (2)

Replacing (2) in (1) and solving the integral, we conclude that

$$Pr(X_1 \le X_2) = 1 - e^{-\lambda_1(T_2 - T_1)} \cdot \frac{\lambda_2}{\lambda_1 + \lambda_2}.$$
 (3)

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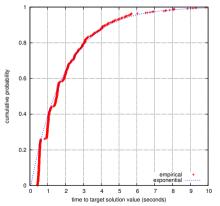
This result can be better interpreted by rewriting expression (3) as

$$Pr(X_1 \le X_2) = (1 - e^{-\lambda_1(T_2 - T_1)}) + e^{-\lambda_1(T_2 - T_1)} \cdot \frac{\lambda_1}{\lambda_1 + \lambda_2}.$$
 (4)

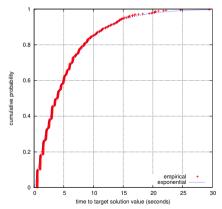
- The first term of the right-hand side of equation (4) is the probability that  $0 \le X_1 \le T_2$ , in which case  $X_1$  is clearly less than or equal to  $X_2$ .
- The second term is given by the product of the factors  $e^{-\lambda_1(T_2-T_1)}$  and  $\lambda_1/(\lambda_1+\lambda_2)$ , in which the former corresponds to the probability that  $X_1 \geq T_2$  and the latter to the probability that  $X_1$  be less than or equal to  $X_2$ , given that  $X_1 > T_2$ .

- To illustrate the above result, we consider two algorithms for solving the server replication for reliable multicast problem.
  - ▶ Algorithm  $A_1$  is an implementation of pure GRASP with  $\alpha = 0.2$ .
  - ▶ Algorithm  $A_2$  is a pure GRASP heuristic with  $\alpha = 0.9$ .
- The runs were performed on an Intel Core2 Quad with 2.40 GHz of clock speed and 4 GB of RAM memory.

The figure depicts the runtime distributions of each algorithm, obtained after 500 runs with different seeds of an instance of the server replication for reliable multicast problem with m=25 and the target value set at 2830.



(a) Pure GRASP with  $\alpha = 0.2$ 

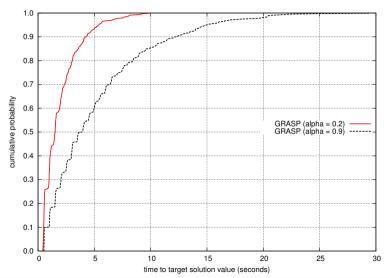


(b) Pure GRASP with  $\alpha = 0.9$ 

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- The parameters of the two distributions are  $\lambda_1=0.524422349,\ T_1=0.36,\ \lambda_2=0.190533895,$  and  $T_2=0.51.$
- Applying expression (3), we get  $Pr(X_1 \le X_2) = 0.684125$ .
- This probability is consistent with the following figure, in which we superimposed the runtime distributions of the two pure GRASP heuristics for the same instance.



The plots in this figure show that the pure GRASP with  $\alpha = 0.2$  outperforms one with  $\alpha = 0.9$ , since the runtime distribution of the former is to the left of the runtime distribution of the latter

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#### Observations:

- If the solution times do not fit exponential (or two-parameter shifted exponential) distributions, as for the case of GRASP with path-relinking heuristics, the closed form result established in expression (3) does not hold.
- Algorithms in this situation cannot be compared by this approach.
- The next topic extends this approach to general runtime distributions.

Let  $X_1$  and  $X_2$  be two continuous random variables, with cumulative probability distributions  $F_{X_1}(\tau)$  and  $F_{X_2}(\tau)$  and probability density functions  $f_{X_1}(\tau)$  and  $f_{X_2}(\tau)$ , respectively. Then,

$$Pr(X_1 \leq X_2) = \int_{-\infty}^{\infty} Pr(X_1 \leq \tau) \cdot f_{X_2}(\tau) \cdot d\tau = \int_{0}^{\infty} Pr(X_1 \leq \tau) \cdot f_{X_2}(\tau) \cdot d\tau,$$

since  $f_{X_1}(\tau) = f_{X_2}(\tau) = 0$  for any  $\tau < 0$ . For an arbitrary small real number  $\varepsilon$ , the above expression can be rewritten as

$$Pr(X_1 \le X_2) = \sum_{i=0}^{\infty} \int_{i \cdot \varepsilon}^{(i+1) \cdot \varepsilon} Pr(X_1 \le \tau) \cdot f_{X_2}(\tau) \cdot d\tau. \tag{5}$$

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Since  $Pr(X_1 \leq i \cdot \varepsilon) \leq Pr(X_1 \leq \tau) \leq Pr(X_1 \leq (i+1) \cdot \varepsilon)$  for  $i \cdot \varepsilon \leq \tau \leq (i+1) \cdot \varepsilon$ , then replacing  $Pr(X_1 \leq \tau)$  by  $Pr(X_1 \leq i \cdot \varepsilon)$  and by  $Pr(X_1 \leq (i+1) \cdot \varepsilon)$  in (5) leads to

$$\sum_{i=0}^{\infty} F_{X_1}(i \cdot \varepsilon) \int_{i \cdot \varepsilon}^{(i+1) \cdot \varepsilon} f_{X_2}(\tau) \cdot d\tau \leq Pr(X_1 \leq X_2) \leq \sum_{i=0}^{\infty} F_{X_1}((i+1) \cdot \varepsilon) \int_{i \cdot \varepsilon}^{(i+1) \cdot \varepsilon} f_{X_2}(\tau) \cdot d\tau.$$

Let  $L(\varepsilon)$  and  $R(\varepsilon)$  be the value of the left and right hand sides of the above expression, respectively, with  $\Delta(\varepsilon) = R(\varepsilon) - L(\varepsilon)$  being the difference between the upper and lower bounds to  $Pr(X_1 \leq X_2)$ .

Then, we have that

$$\Delta(\varepsilon) = \sum_{i=0}^{\infty} \left[ F_{X_1}((i+1) \cdot \varepsilon) - F_{X_1}(i \cdot \varepsilon) \right] \int_{i \cdot \varepsilon}^{(i+1) \cdot \varepsilon} f_{X_2}(\tau) \cdot d\tau. \tag{6}$$

Let  $\delta = \max_{\tau \geq 0} \{f_{X_1}(\tau)\}$ . Since  $|F_{X_1}((i+1) \cdot \varepsilon) - F_{X_1}(i \cdot \varepsilon)| \leq \delta \cdot \varepsilon$  for  $i \geq 0$ , expression (6) turns out to be

$$\Delta(\varepsilon) \leq \sum_{i=0}^{\infty} \delta \cdot \varepsilon \int_{i \cdot \varepsilon}^{(i+1) \cdot \varepsilon} f_{X_2}(\tau) \cdot d\tau = \delta \cdot \varepsilon \int_0^{\infty} f_{X_2}(\tau) \cdot d\tau = \delta \cdot \varepsilon.$$

Consequently,

$$\Delta(\varepsilon) \le \delta \cdot \varepsilon,\tag{7}$$

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i.e., the difference  $\Delta(\varepsilon)$  between the upper and lower bounds to  $Pr(X_1 \leq X_2)$  (or the absolute error in the integration) is smaller than or equal to  $\delta \varepsilon$ .

Therefore, this difference can be made as small as desired by choosing a sufficiently small value for  $\varepsilon$ .

In order to numerically evaluate a good approximation to  $Pr(X_1 \le X_2)$ , we select the appropriate value of  $\varepsilon$  such that the resulting approximation error  $\Delta(\varepsilon)$  is sufficiently small.

Next, we compute  $L(\varepsilon)$  and  $R(\varepsilon)$  to obtain the approximation

$$Pr(X_1 \le X_2) \approx \frac{L(\varepsilon) + R(\varepsilon)}{2}.$$
 (8)

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In practice, the above probability distributions are unknown.

Instead of the distributions, the information available is limited to a sufficiently large number  $N_1$  (resp.  $N_2$ ) of observations of the random variable  $X_1$  (resp.  $X_2$ ).

Since the value of  $\delta = \max_{\tau \geq 0} \{f_{X_1}(\tau)\}$  is also unknown beforehand, the appropriate value of  $\varepsilon$  cannot be estimated.

Then, we proceed iteratively as follows.

Let  $t_1(j)$  (resp.  $t_2(j)$ ) be the value of the j-th smallest observation of the random variable  $X_1$  (resp.  $X_2$ ), for  $j = 1, \ldots, N_1$  (resp.  $N_2$ ).

We set the bounds  $a = \min\{t_1(1), t_2(1)\}$  and  $b = \max\{t_1(N_1), t_2(N_2)\}$  and choose an arbitrary number h of integration intervals to compute an initial value  $\varepsilon = (b-a)/h$  for each integration interval.

For sufficiently small values of the integration interval  $\varepsilon$ , the probability density function  $f_{X_1}(\tau)$  in the interval  $[i \cdot \varepsilon, (i+1) \cdot \varepsilon]$  can be approximated by  $\hat{f}_{X_1}(\tau) = (\hat{F}_{X_1}((i+1) \cdot \varepsilon) - \hat{F}_{X_1}(i \cdot \varepsilon))/\varepsilon$ , where

$$\hat{F}_{X_1}(i\cdot\varepsilon)=|\{t_1(j),j=1,\ldots,N_1:t_1(j)\leq i\cdot\varepsilon\}|. \tag{9}$$

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The same approximations hold for random variable  $X_2$ .

Finally, the value of  $Pr(X_1 \leq X_2)$  can be computed as in expression (8), using the estimates  $\hat{f}_{X_1}(\tau)$  and  $\hat{f}_{X_2}(\tau)$  in the computation of  $L(\varepsilon)$  and  $R(\varepsilon)$ .

If the approximation error  $\Delta(\varepsilon) = R(\varepsilon) - L(\varepsilon)$  becomes sufficiently small, then the procedure stops.

Otherwise, the value of  $\varepsilon$  is halved and the above steps are repeated until convergence.

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### Numerical applications to sequential algorithms

We illustrate next an application of the procedure described in the previous topic for the comparison of randomized algorithms (running on the same instance) on three problems:

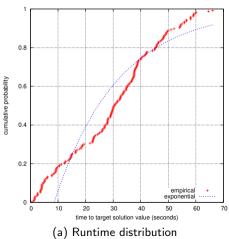
- server replication for reliable multicast,
- routing and wavelength assignment, and
- 2-path network design.

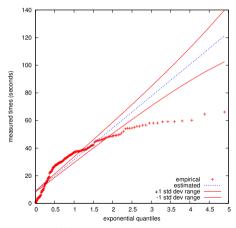
# DM-D5 and GRASP algorithms for server replication

- Multicast communication consists of simultaneously delivering the same information to many receivers, from single or multiple sources.
- Network services specially designed for multicast are needed.
- The scheme used in current multicast services create a delivery tree, whose root represents the sender, whose leaves represent the receivers, and whose internal nodes represent network routers or relaying servers.
- Transmission is performed by creating copies of the data at split points of the tree.
- An important issue regarding multicast communication is how to provide reliable service, ensuring the delivery of all packets from the sender to receivers.
- A successful technique to provide reliable multicast service is the server replication approach, in which
  data is replicated at some of the multicast-capable relaying hosts (also called replicated or repair
  servers) and each of them is responsible for the retransmission of packets to receivers in its group.
- The problem consists in selecting the best subset of the multicast-capable relaying hosts to act as replicated servers in a multicast scenario.
- It is a special case of the p-median problem.

- DM-GRASP is a hybrid version of GRASP which incorporates a data-mining process.
- We compare two heuristics for the server replication problem:
  - algorithm A<sub>1</sub> is an implementation of the DM-D5 version of DM-GRASP, in which the mining algorithm is periodically applied,
  - $\blacktriangleright$  while  $A_2$  is a pure GRASP heuristic.
- We present results for two instances using the same network scenario, with m=25 and m=50 replication servers.
- Each algorithm was run 200 times with different seeds.
- The target was set at 2,818.925 (the best known solution value is 2,805.89) for the instance with m = 25 and at 2,299.07 (the best known solution value is 2,279.84) for the instance with m = 50.

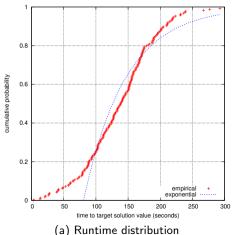
The figure depicts the runtime distribution and quantile-quantile plot for algorithm DM-D5 on the instance with m=25 and the target value set at 2,818.925.

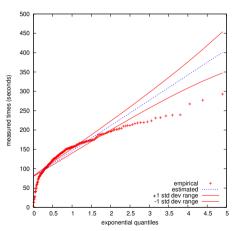




(b) Quantile-quantile plot

The figure depicts the runtime distribution and quantile-quantile plot for algorithm DM-D5 on the instance with m = 50 and the target value set at 2,299.07.



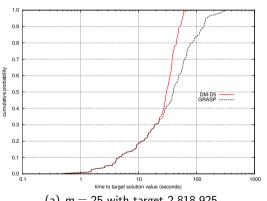


(b) Quantile-quantile plot

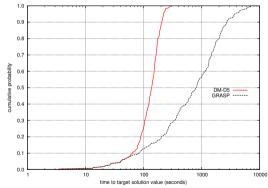
#### Observations:

- Running times of DM-D5 did not fit exponential distributions for any of the instances.
- GRASP solution times were exponential for both.

The figure shows the superimposed empirical runtime distributions of DM-D5 and GRASP.



(a) m = 25 with target 2,818.925  $Pr(X_1 \le X_2) = 0.619763$ 



(b) m = 50 with target 2,299.07  $Pr(X_1 < X_2) = 0.854113$ 

#### Observations:

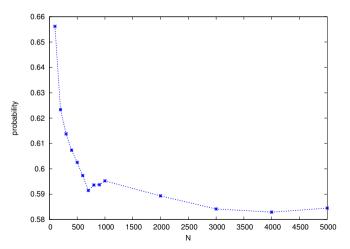
- Algorithm DM-D5 outperformed GRASP, since the runtime distribution of the DM-D5 is to the left of the distribution for GRASP on the both instances, with m = 25 and m = 50.
- Consistently, the computations show that:
  - ▶  $Pr(X_1 \le X_2) = 0.619763$  (with  $L(\varepsilon) = 0.619450$ ,  $R(\varepsilon) = 0.620075$ ,  $\Delta(\varepsilon) = 0.000620$ , and  $\varepsilon = 0.009552$ ) for the instance with m = 25.
  - ▶  $Pr(X_1 \le X_2) = 0.854113$  (with  $L(\varepsilon) = 0.853800$ ,  $R(\varepsilon) = 0.854425$ ,  $\Delta(\varepsilon) = 0.000625$ , and  $\varepsilon = 0.427722$ ) for the instance with m = 50.

- We also investigate the convergence of the proposed measure with the sample size (i.e., with the number of independent runs of each algorithm).
- Convergence with the sample size is illustrated next for the same m=25 instance of the server replication problem, with the same target 2,818.925 already used in the previous experiment.
- Once again, algorithm A<sub>1</sub> is the DM-D5 version of DM-GRASP and algorithm A<sub>2</sub> is the pure GRASP heuristic.
- The estimation of  $Pr(X_1 \le X_2)$  is computed for N = 100, 200, 300, 400, 500, 600, 700, 800, 900, 1000, 2000, 3000, 4000, and 5000 independent runs of each algorithm.

Convergence of the estimation of  $Pr(X_1 \le X_2)$  with the sample size for the m=25 instance of the server replication problem.

N	L(arepsilon)	$Pr(X_1 \leq X_2)$	R(arepsilon)	$\Delta(arepsilon)$	$\varepsilon$
100	0.655900	0.656200	0.656500	0.000600	0.032379
200	0.622950	0.623350	0.623750	0.000800	0.038558
300	0.613344	0.613783	0.614222	0.000878	0.038558
400	0.606919	0.607347	0.607775	0.000856	0.038558
500	0.602144	0.602548	0.602952	0.000808	0.038558
600	0.596964	0.597368	0.597772	0.000808	0.038558
700	0.591041	0.591440	0.591839	0.000798	0.038558
800	0.593197	0.593603	0.594009	0.000812	0.042070
900	0.593326	0.593719	0.594113	0.000788	0.042070
1000	0.594849	0.595242	0.595634	0.000785	0.042070
2000	0.588913	0.589317	0.589720	0.000807	0.047694
3000	0.583720	0.584158	0.584596	0.000875	0.047694
4000	0.582479	0.582912	0.583345	0.000866	0.047694
5000	0.584070	0.584511	0.584953	0.000882	0.050604

Convergence of the estimation of  $Pr(X_1 \leq X_2)$  with the sample size for the m=25 instance of the server replication problem.

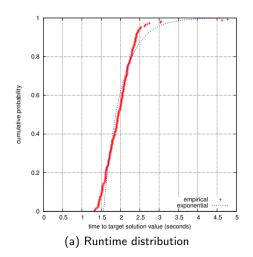


#### Observation:

The estimation of  $Pr(X_1 \le X_2)$  stabilizes as the sample size N increases.

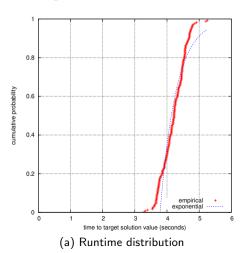
- A point-to-point connection between two endnodes of an optical network is called a lightpath.
- Two lightpaths may use the same wavelength, provided they do not share any common link.
- The routing and wavelength assignment problem is that of routing a set of lightpaths and assigning a wavelength to each of them, minimizing the number of wavelengths needed.
- A decomposition strategy is compared with a multistart greedy heuristic.
- Two networks are used for benchmarking:
  - ▶ The first has 27 nodes representing the capitals of the 27 states of Brazil, with 70 links connecting them. There are 702 lightpaths to be routed.
  - ▶ Instance Finland is formed by 31 nodes and 51 links, with 930 lightpaths to be routed.
- Each algorithm was run 200 times with different seeds.
- The target was set at 24 (the best known solution value) for instance Brazil and at 50 for instance Finland (the best known solution value is 47).
- Algorithm  $A_1$  is the multistart heuristic, while  $A_2$  is the tabu search decomposition scheme.
- The multistart solution times fit exponential distributions for both instances.

The figure displays the runtime distribution and quantile-quantile plot for tabu search on Brazil instance with the target value set at 24.



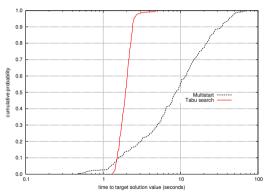
4.5 neasured times (seconds) 3.5 2.5 1.5 0 exponential quantiles

The figure displays the runtime distribution and quantile-quantile plot for tabu search on Finland instance with the target value set at 50.

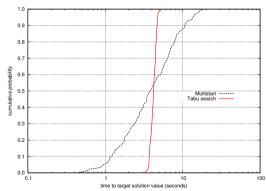


6.5 neasured times (seconds) 5.5 4.5 3.5 0 0.5 exponential quantiles

The figure shows the superimposed empirical runtime distributions of multistart and tabu search algorithms.



(a) Brazil instance with target 24  $Pr(X_1 \le X_2) = 0.13$ 



(b) Finland instance with target 50  $Pr(X_1 \le X_2) = 0.536787$ 

#### Observations:

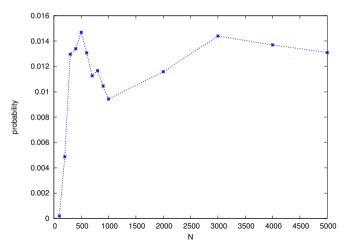
- The direct comparison of the two approaches shows that decomposition clearly outperformed the multistart strategy for instance Brazil, since  $Pr(X_1 \le X_2) = 0.13$  in this case (with  $L(\varepsilon) = 0.129650$ ,  $R(\varepsilon) = 0.130350$ ,  $\Delta(\varepsilon) = 0.000700$ , and  $\varepsilon = 0.008163$ ).
- However, the situation changes for instance Finland.
- Although both algorithms have similar performances, multistart is slightly better with respect to the measure proposed in this presentation, since  $Pr(X_1 \le X_2) = 0.536787$  (with  $L(\varepsilon) = 0.536525$ ,  $R(\varepsilon) = 0.537050$ ,  $\Delta(\varepsilon) = 0.000525$ , and  $\varepsilon = 0.008804$ ).

- As done for the server replication problem, we also investigate the convergence of the proposed measure with the sample size (i.e., with the number of independent runs of each algorithm).
- Convergence with the sample size is illustrated next for the Finland instance of the routing and wavelength assignment problem, with the target set at 49.
- Once again, algorithm A<sub>1</sub> is the multistart heuristic and algorithm A<sub>2</sub> is the tabu search decomposition scheme.
- The estimation of  $Pr(X_1 \le X_2)$  is computed for N = 100, 200, 300, 400, 500, 600, 700, 800, 900, 1000, 2000, 3000, 4000, and 5000 independent runs of each algorithm.

Convergence of the estimation of  $Pr(X_1 \leq X_2)$  with the sample size for the Finland instance of the routing and wavelength assignment problem.

N	$L(\varepsilon)$	$Pr(X_1 \leq X_2)$	$R(\varepsilon)$	$\Delta(arepsilon)$	$\varepsilon$
100	0.000001	0.000200	0.000400	0.000400	1.964844
200	0.000100	0.004875	0.009650	0.009550	0.000480
300	0.006556	0.012961	0.019367	0.012811	0.000959
400	0.007363	0.013390	0.019425	0.012063	0.000959
500	0.007928	0.014694	0.021460	0.013532	0.000610
600	0.006622	0.013069	0.019517	0.012894	0.000610
700	0.005722	0.011261	0.016800	0.011078	0.000610
800	0.005033	0.011667	0.018302	0.013269	0.000610
900	0.004556	0.010461	0.016367	0.011811	0.000610
1000	0.004100	0.009425	0.014750	0.010650	0.000610
2000	0.006049	0.011580	0.017112	0.011063	0.000610
3000	0.007802	0.014395	0.020987	0.013185	0.000610
4000	0.007408	0.013698	0.019988	0.012580	0.000610
5000	0.006791	0.013090	0.019389	0.012598	0.000623

Convergence of the estimation of  $Pr(X_1 \leq X_2)$  with the sample size for the Finland instance of the routing and wavelength assignment problem.



#### Observation:

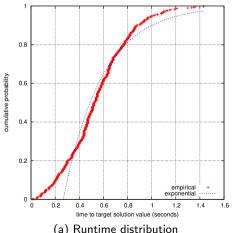
Once again, the estimation of  $Pr(X_1 \leq X_2)$  stabilizes as the sample size N increases.

## GRASP algorithms for 2-path network design

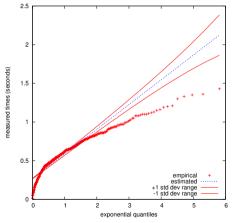
- Given a connected undirected graph with non-negative weights associated with its edges, together
  with a set of origin-destination nodes, the 2-path network design problem consists in finding a
  minimum weighted subset of edges containing a path formed by at most two edges between every
  origin-destination pair.
- Applications can be found in the design of communication networks, in which paths with few edges
  are sought to enforce high reliability and small delays.

- We first compare four GRASP heuristics for solving an instance of the 2-path network design problem with 90 nodes.
  - ▶ The first heuristic is a pure GRASP (algorithm  $A_1$ ).
  - ► The others integrate different path-relinking strategies for search intensification at the end of each GRASP iteration:
    - \* forward path-relinking (algorithm  $A_2$ ),
    - $\star$  bidirectional path-relinking (algorithm  $A_3$ ), and
    - ★ backward path-relinking (algorithm  $A_4$ ).
- Each algorithm was run 500 independent times on the benchmark instance with 90 nodes and 900 origin-destination pairs, with the solution target value set at 673 (the best known solution value is 639).
- The runtime distributions and quantile-quantile plots for the different versions of GRASP with path-relinking are shown in the following figures.

The figure shows the runtime distribution and quantile-quantile plot for GRASP with forward path-relinking on 90-node instance with the target value set at 673.

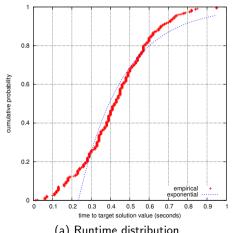


(a) Runtime distribution

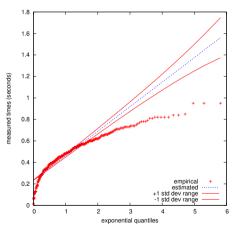


(b) Quantile-quantile plot

The figure shows the runtime distribution and quantile-quantile plot for GRASP with bidirectional path-relinking on 90-node instance with the target value set at 673.

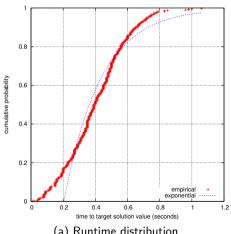


(a) Runtime distribution

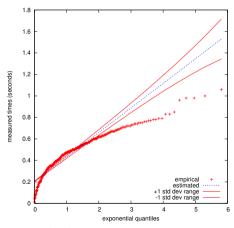


(b) Quantile-quantile plot

The figure shows the runtime distribution and quantile-quantile plot for GRASP with backward path-relinking on 90-node instance with the target value set at 673.

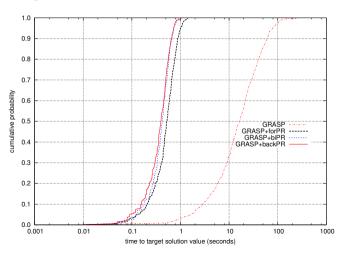


(a) Runtime distribution



(b) Quantile-quantile plot

The figure shows the superimposed empirical runtime distributions of pure GRASP and three versions of GRASP with path-relinking.



#### Observations:

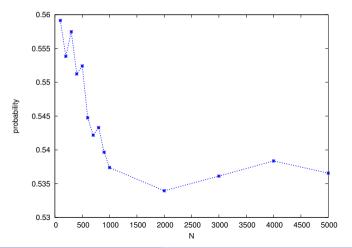
- Algorithm  $A_2$  (as well as  $A_3$  and  $A_4$ ) performs much better than  $A_1$ , as indicated by  $Pr(X_2 \le X_1) = 0.986604$  (with  $L(\varepsilon) = 0.986212$ ,  $R(\varepsilon) = 0.986996$ ,  $\Delta(\varepsilon) = 0.000784$ , and  $\varepsilon = 0.029528$ ).
- Algorithm  $A_3$  outperforms  $A_2$ , as illustrated by the fact that  $Pr(X_3 \le X_2) = 0.636000$  (with  $L(\varepsilon) = 0.630024$ ,  $R(\varepsilon) = 0.641976$ ,  $\Delta(\varepsilon) = 0.011952$ , and  $\varepsilon = 1.354218 \times 10^{-6}$ ).
- Finally, we observe that algorithms  $A_3$  and  $A_4$  behave very similarly, although  $A_4$  performs slightly better for this instance, since  $Pr(X_4 \le X_3) = 0.536014$  (with  $L(\varepsilon) = 0.528560$ ,  $R(\varepsilon) = 0.543468$ ,  $\Delta(\varepsilon) = 0.014908$ , and  $\varepsilon = 1.001358 \times 10^{-6}$ ).

- As for the problems considered in previous topics, we also investigate the convergence of the proposed measure as a function of sample size (i.e., with the number of independent runs of each algorithm).
- Convergence with the sample size is illustrated next for the 90-node instance of the 2-path network design problem, with the same target 673 previously used.
- We recall that algorithm  $A_1$  is the GRASP with backward path-relinking heuristic, while algorithm  $A_2$  is the GRASP with bidirectional path-relinking heuristic.
- The estimation of  $Pr(X_1 \le X_2)$  is computed for N = 100, 200, 300, 400, 500, 600, 700, 800, 900, 1000, 2000, 3000, 4000, and 5000 independent runs of each algorithm.

Convergence of the estimation of  $Pr(X_1 \leq X_2)$  with the sample size for the 90-node instance of the 2-path network design problem.

N	$L(\varepsilon)$	$Pr(X_1 \leq X_2)$	$R(\varepsilon)$	$\Delta(\varepsilon)$	$\varepsilon$
100	0.553300	0.559150	0.565000	0.011700	$4.387188 \times 10^{-7}$
200	0.553250	0.553850	0.554450	0.001199	$4.501629 \times 10^{-7}$
300	0.551578	0.557483	0.563389	0.011811	$4.501629 \times 10^{-7}$
400	0.545244	0.551241	0.557238	0.011994	$4.730511 \times 10^{-7}$
500	0.546604	0.552420	0.558236	0.011632	$5.035686 \times 10^{-7}$
600	0.538867	0.544749	0.550631	0.011764	$5.073833 \times 10^{-7}$
700	0.536320	0.542181	0.548041	0.011720	$5.073833 \times 10^{-7}$
800	0.537533	0.543298	0.549064	0.011531	$5.073833 \times 10^{-7}$
900	0.533912	0.539671	0.545430	0.011517	$5.073833 \times 10^{-7}$
1000	0.531595	0.537388	0.543180	0.011585	$5.073833 \times 10^{-7}$
2000	0.528224	0.533959	0.539698	0.011469	$5.722427 \times 10^{-7}$
3000	0.530421	0.536128	0.541835	0.011414	$6.027603 \times 10^{-7}$
4000	0.532695	0.538364	0.544033	0.011338	$6.027603 \times 10^{-7}$
5000	0.530954	0.536566	0.542178	0.011225	$6.027603 \times 10^{-7}$

Convergence of the estimation of  $Pr(X_1 \leq X_2)$  with the sample size for the 90-node instance of the 2-path network design problem.

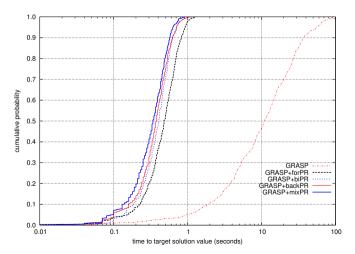


#### Observation:

Once again, the estimation of  $Pr(X_1 \leq X_2)$  stabilizes as the sample size N increases.

- We next compare five GRASP heuristics for the 2-path network design problem, with and without path-relinking, for solving an instance with 80 nodes and 800 origin-destination pairs, with target value set at 588 (the best known solution value is 577).
- In this example, the first algorithm is a pure GRASP (algorithm  $A_1$ ).
- The other heuristics integrate different path-relinking strategies at the end of each GRASP iteration:
  - ▶ forward path-relinking (algorithm A₂),
  - ▶ bidirectional path-relinking (algorithm A<sub>3</sub>),
  - **b** backward path-relinking (algorithm  $A_4$ ), and
  - ightharpoonup mixed path-relinking (algorithm  $A_5$ ).
- As before, each heuristic was run independently 500 times.

The figure shows the superimposed empirical runtime distributions of pure GRASP and four GRASP with path-relinking heuristics.

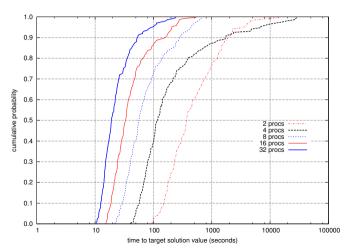


#### Observations:

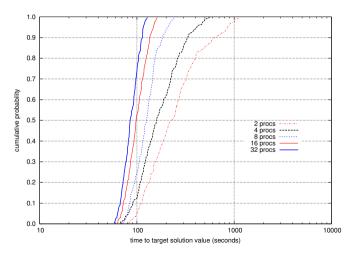
- Algorithm  $A_2$  (as well as  $A_3$ ,  $A_4$ , and  $A_5$ ) performs much better than  $A_1$ , as indicated by  $Pr(X_2 \le X_1) = 0.970652$  (with  $L(\varepsilon) = 0.970288$ ,  $R(\varepsilon) = 0.971016$ ,  $\Delta(\varepsilon) = 0.000728$ , and  $\varepsilon = 0.014257$ ).
- Algorithm  $A_3$  outperforms  $A_2$ , as shown by the fact that  $Pr(X_3 \le X_2) = 0.617278$  (with  $L(\varepsilon) = 0.610808$ ,  $R(\varepsilon) = 0.623748$ ,  $\Delta(\varepsilon) = 0.012940$ , and  $\varepsilon = 1.220703 \times 10^{-6}$ ).
- Algorithm  $A_4$  performs slightly better than  $A_3$  for this instance, since  $Pr(X_4 \le X_3) = 0.537578$  (with  $L(\varepsilon) = 0.529404$ ,  $R(\varepsilon) = 0.545752$ ,  $\Delta(\varepsilon) = 0.016348$ , and  $\Delta(\varepsilon) = 1.201630 \times 10^{-6}$ ).
- Algorithms  $A_5$  and  $A_4$  also behave very similarly, but  $A_5$  is slightly better for this instance since  $Pr(X_5 \le X_4) = 0.556352$  (with  $L(\varepsilon) = 0.547912$ ,  $R(\varepsilon) = 0.564792$ ,  $\Delta(\varepsilon) = 0.016880$ , and  $\varepsilon = 1.001358 \times 10^{-6}$ ).

- We conclude this presentation by describing the use of the runtime distribution methodology to evaluate and compare parallel implementations of stochastic local search algorithms.
- Once again, the 2-path network design problem is used to illustrate this application.
- The next two figures superimpose the runtime distributions of, respectively, cooperative and independent parallel implementations of GRASP with bidirectional path-relinking for the same problem on 2, 4, 8, 16, and 32 processors, on an instance with 100 nodes and 1000 origin-destination pairs, using 683 as target value.
- Each algorithm was run independently 200 times.
- We denote by  $A_1^k$  (resp.  $A_2^k$ ) the cooperative (resp. independent) parallel implementation running on k processors, for k = 2, 4, 8, 16, 32.

The figure shows the superimposed empirical runtime distributions of cooperative parallel GRASP with bidirectional path-relinking running on 2, 4, 8, 16, and 32 processors.



The figure shows the superimposed empirical runtime distributions of independent parallel GRASP with bidirectional path-relinking running on 2, 4, 8, 16, and 32 processors.



Comparing cooperative (algorithm  $A_1$ ) and independent (algorithm  $A_2$ ) parallel implementations.

Processors (k)	$Pr(X_1^k \leq X_2^k)$		
2	0.309784		
4	0.597253		
8	0.766806		
16	0.860864		
32	0.944938		

#### Observations:

- The table shows the probability that the cooperative parallel implementation performs better than the independent implementation on 2, 4, 8, 16, and 32 processors.
- The independent implementation performs better than the cooperative implementation on two processors.
- In that case, the cooperative implementation does not benefit from the availability of two processors, since only one of them performs iterations, while the other acts as the master.
- However, as the number of processors increases from two to 32, the cooperative implementation
  performs progressively better than the independent implementation, since more processors are
  devoted to perform GRASP iterations.

- The proposed methodology is clearly consistent with the relative behavior of the two parallel versions for any number of processors.
- Furthermore, it illustrates that the cooperative implementation becomes progressively better than the independent implementation when the number of processors increases.

Comparing the parallel implementations on  $2^{j+1}$  (algorithm  $A_1$ ) and  $2^j$  (algorithm  $A_2$ ) processors, for j = 1, 2, 3, 4.

Processors (a)	Processors (b)	$Pr(X_1^a \leq X_1^b)$	$Pr(X_2^a \leq X_2^b)$
4	2	0.766235	0.651790
8	4	0.753904	0.685108
16	8	0.724398	0.715556
32	16	0.747531	0.669660

#### Observations:

- Both implementations scale appropriately as the number of processors grows.
- Once again, we can see that the performance measure appropriately describes the relative behavior
  of the two parallel strategies and provides insight on how parallel algorithms scale with the number
  of processors.
- The table shows numerical evidence to evaluate the trade-offs between computation times and the number of processors in parallel implementations.

#### Concluding remarks

The material in this talk is taken from

• Chapter 6 - Runtime distributions

of our book, *Optimization by GRASP: Greedy Randomized Adaptive Search Procedures* (Resende & Ribeiro, Springer. 2016).

