Flexible selection of heterogeneous and unreliable services in large-scale grids

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As grids become larger and more interconnected in nature, scientists can benefit from a growing number of distributed services that may be invoked on demand to complete complex computational workflows. However, it also means that these scientists become dependent on the cooperation of third-party service providers, whose behaviour may be uncertain, failure prone and highly heterogeneous. To address this, we have developed a novel decision-theoretic algorithm that automatically selects appropriate services for the tasks of an abstract workflow and deals with failures through redundancy and dynamic re-invocation of functionally equivalent services. In this paper, we summarize our approach, describe in detail how it can be applied to a real-world bioinformatics workflow and show that it offers a significant improvement over current service selection techniques.

**Keywords:** service selection; grids; reliability; service-oriented computing

1. Introduction

Grid computing is a key enabling technology for the field of e-Science. As scientists rely increasingly on computational support for their work, grids allow research organizations to pool their resources, spread the cost of expensive hardware and use idle machines (Coveney 2005). However, as grids become larger, more open and distributed, new challenges need to be addressed (Foster et al. 2004). In particular, the behaviour of services may be highly uncertain, as manifested by frequent failures and execution delays. This uncertainty may be due to network delays, power cuts, competition for resources or machine failures. Moreover, services will increasingly be owned by agents that act autonomously and cannot be assumed to always behave cooperatively or deterministically (Jennings 2001). Rather, such agents will follow their own decision-making procedures and may even defect when this is beneficial. For example, the owner of a computational cluster may decide to temporarily suspend all jobs in its queue in order to run an important internal job.

Clearly, this uncertainty must be addressed, and this is particularly important when consumers execute large workflows of interdependent tasks, when service providers demand remuneration and when workflows have deadlines. To this end, we concentrate on the process of deciding which service instances to select for the workflow tasks.
for the constituent tasks of an abstract workflow in an automated fashion. This allows the consumer to consider non-functional quality-of-service (QoS) parameters (such as the cost or reliability) of functionally equivalent services and to react to failures by dynamically selecting alternative services when previous ones have failed.

Current work does not address the selection problem satisfactorily. Typically, QoS parameters are considered only locally for each task, either to meet user-imposed quality constraints (Thain & Livny 2003), or to optimize some of the parameters, e.g. to select the cheapest or most reliable service (Vazhkudai et al. 2001). Other work optimizes a weighted sum of global QoS parameters, such as the workflow cost and reliability (Zeng et al. 2004). However, it is often unclear how weights are set and both the local and global approaches select only a single service for each task, thus producing brittle workflows that are vulnerable to failures.

A number of existing grid technologies do consider service failures explicitly. In particular, failures are sometimes handled by submitting the failed task to a different service provider (Thain & Livny 2003; Oinn et al. 2004). In other systems, fixed redundancy is employed to increase the probability of success (Anderson et al. 2002). However, neither of these approaches explicitly models the impact of failures on the overall workflow (e.g. whether failures will result in missing the overall deadline and whether the cost of redundancy outweighs its benefit).

To address these shortcomings, we have previously developed a decision-theoretic service selection algorithm (Stein et al. 2007). This reasons explicitly about the value of choosing particular service instances and balances the overall workflow cost with its expected reward. Furthermore, the algorithm uses both service redundancy and dynamic reselection, but applies them in a principled manner based on their impact on the expected performance of the workflow. While our previous work concentrated on the abstract description of the algorithm, in this paper, we examine its applicability to e-Science by considering a bioinformatics workflow. Furthermore, we present novel experimental results that show an average 46.8 per cent improvement in utility over existing selection approaches.

The remainder of this paper is organized as follows. In §2, we provide a high-level description of our algorithm. Then, in §3, we show how it can be applied to a real grid problem. In §4, we evaluate our work experimentally and conclude in §5.

2. Flexible service selection

In this section, we summarize our assumptions about how services are provided and consumed in a grid system and we provide a high-level description of our selection algorithm. The full technical details of our approach and associated assumptions can be found in our previous work (Stein et al. 2007).

Throughout our work, we consider a generic grid system that is based on service-oriented technologies (Foster et al. 2002). Here, providers advertise their services with a registry (which could be potentially federated), which can then be discovered and invoked by consumers when required. Specifically, we assume that there is no centralized resource or workload manager, but rather that each consumer is responsible for selecting and communicating with the service providers directly. We believe that such a view is more appropriate for open, large-scale, highly heterogeneous systems, where a centralized solution would
represent a potentially unacceptable communication and performance bottleneck, create a single point of failure and significantly decrease the autonomy of both consumers and providers.

Figure 1 shows the typical activities of a service consumer in our model, which consists of four steps. First, during workflow selection, a consumer selects an appropriate workflow to meet its current high-level goal. Such a workflow describes the types of services required by the consumer and related ordering constraints (for an example, see §3). In practice, it may be chosen from a repository of workflows, automatically synthesized by a planner or submitted by a human user. In more detail, we assume that a workflow is a directed acyclic graph, where nodes represent tasks, each mapped to a specific service type, and edges represent ordering constraints. Furthermore, we assume that a consumer derives a time-dependent utility value when completing the workflow (or none if a fixed deadline is reached).

Given such an abstract workflow, the consumer next engages in a service discovery phase. Here, it discovers candidate services by searching a public registry or by contacting an appropriate service broker. This discovery step uses the service types provided by the workflow, mapping each type to a set of suitable services.

This is followed by service selection, which constitutes the main focus of this paper. Here, the consumer makes predictions about the performance of the workflow and explores the space of service allocations. In doing so, we assume that the consumer has access to some QoS information about the matching services as follows.

— Cost. The invocation cost, expressed in the same units as the workflow utility. This is paid on invocation and regardless of the eventual outcome.
— Failure probability. The probability that it will fail to complete a task. In this case, a service is not generally assumed to send explicit failure messages.
— Duration function. A probability distribution for the time between invoking the service and receiving notification of a successful outcome (as observed by the consumer and including any data transmission or queueing delays).

As described in §1, we take a decision-theoretic approach and are therefore interested in selecting services that maximize the consumer’s expected utility, as given by the difference between the reward of completing the workflow and the overall cost incurred (we assume that the consumer is risk neutral). Formally, the aim of our algorithm is to find \( \alpha^* \),

\[
\alpha^* = \arg\max_\alpha E(R(\alpha) - C(\alpha)),
\]

(2.1)

where \( \alpha \) is a service allocation that maps tasks to selected services; \( R(\alpha) \) and \( C(\alpha) \) are random variables representing, respectively, the reward when following allocation \( \alpha \) and the associated cost; and \( E(\cdot) \) is the expectation operator.
Now, while existing selection approaches typically map each task to a single service, we are interested in introducing redundancy where necessary and in planning for potential failures by selecting and invoking multiple services in series. Hence, we represent the service allocation $\alpha$ as a mapping that associates each task with the selected services for that task and their respective invocation time steps (conditional on the task not having been completed successfully). An example of such an allocation for a single task is shown in figure 2. Here, three cheap and unreliable services are invoked immediately when the task first becomes executable. If they have not been successful after 15 time steps, more services are gradually invoked to ensure that the task is completed eventually.

Unfortunately, the optimization problem shown in equation (2.1) is non-deterministic polynomial-time-hard (NP-hard) due to its combinatorial nature. This means that it is generally unfeasible to compute a globally optimal solution for larger workflows with many candidate services (as is common in large grid systems). For this reason, we carry out a local search, as summarized in figure 3. This starts with a random candidate solution and then iteratively improves this until no more improvements can be made. As calculating the expected utility of a given allocation is intractable, we use a heuristic approximation for this ($\tilde{u}(\alpha)$). This first calculates a number of performance parameters for each workflow task (i.e. a success probability, expected cost, mean duration and variance), then aggregates these over the entire workflow and finally uses them to obtain an estimated expected utility. Such a heuristic typically converges quickly to a good solution, but is not guaranteed to find the global optimum $\alpha^*$.

In the following, we illustrate our approach using a real-world grid workflow.

3. Illustrative example

To demonstrate how our work can be employed in practice, we use a simple workflow from the bioinformatics domain—an area that relies heavily on computationally intensive services and that has increasingly seen the establishment of large distributed grid systems for sharing resources (Oinn et al. 2004).
In our example, a scientist has just sequenced a previously unknown gene of a bacterium, and is now interested in visualizing the shape of the associated protein. For this, she has to carry out a number of tasks, which are shown as a workflow in Figure 4a.

To summarize this briefly, the scientist first uses a base-calling service (BaseCall) to annotate DNA chromatograms and attach quality values to each base. These are then assembled into a single continuous DNA sequence by a sequence-assembling service (GeneAssemble), which also identifies the coding region of the gene. Next, a translation service is used to derive a corresponding amino acid sequence (Translate), and a particularly computationally extensive folding service (Fold) predicts the three-dimensional shape of the coded protein. This is then rendered in high resolution using a graphics service (Render). In parallel with the folding service, the scientist is also interested in comparing the gene with known sequences. To this end, she searches through public collections of proteins to find the closest match using a specialized service (Blast), and then accesses database services to retrieve structural information about this protein (LookUp). This is rendered again, and both images are printed as part of a report on a local printer (Print).

Now, some service types in this example require a significant computational effort and may take a considerable amount of time to complete. In this context, the workflow utility function allows the scientist to succinctly encode the overall value of the workflow and how this relates to the time taken. For example, if the scientist needs the results later in the day, but is not overly concerned about waiting a bit longer, a utility function that decreases slowly over time, such as $u_1(x)$ in Figure 4b, is appropriate. If, on the other hand, the results are critical for a presentation she is giving to a funding committee just hours later, a utility function such as $u_2(x)$ expresses the urgency and high value of the workflow more suitably.

To illustrate how our approach deals with many heterogeneous services, we assume that each workflow task has associated with it a number of services, as shown in Table 1. Here, each row of the table describes a group of services that share the same performance characteristics. For example, the population $P_0$ in the first row represents 20 services that each cost £1, fail 20 per cent of the time and follow a gamma distribution with shape $k=1.5$ and scale $\theta=2$. Generally, we have chosen these service populations to offer certain trade-offs, e.g. services in $P_1$ are not only more reliable and faster than those in $P_0$, but also three times as expensive, while services in $P_{29}$ are slower, but more reliable than those in $P_{21}$.

To demonstrate the types of decisions our algorithm makes, Figure 5 shows the final service allocation that is found by a local search when the less urgent utility function $u_1$ is used (this rewards a constant utility of 150 if the workflow is completed within

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1 Full details can be found in Stein (2008).

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4 hours, then deducts 1 from this for every minute that the workflow is late, until no more utility is gained after 390 min). Here, services are selected redundantly for each task, such that their respective success probabilities are between 0.99 and 1. However, the algorithm generally chooses cheaper and slower services where possible, resulting in low expected costs and large mean durations and variances for each workflow task. Two example task allocations are shown here for the Fold and the lower Render tasks. For the former, the algorithm initially chooses two services from $P_1$ and one from $P_2$. Although each of these has a non-negligible failure probability, their parallel invocation, followed some time later by another service from $P_1$, results in an overall success probability of 0.99. However, despite the redundancy, the services are slow and the overall completion time has a high variance (as shown by the cumulative success probability of the task over time, which is plotted along with the allocation). Owing to the lower failure probability for $P_1$, the algorithm proceeds more cautiously for the Render task, invoking a single service only, followed almost 2 hours later by a second one.

Given the performance characteristics for each task, these are aggregated to yield an overall success probability for the workflow (0.98) and an expected cost (65.26). To estimate the overall duration of the workflow, we sum all mean durations and variances along the critical path of the workflow (i.e. the longest path when considering the mean duration of each task), and then use a normal

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Table 1. Service types used in the example workflow.

<table>
<thead>
<tr>
<th>service type</th>
<th>ID</th>
<th>failure probability</th>
<th>cost (£)</th>
<th>number</th>
<th>duration</th>
<th>mean (min)</th>
<th>variance (min$^2$)</th>
</tr>
</thead>
<tbody>
<tr>
<td>BaseCall</td>
<td>$P_0$</td>
<td>0.2</td>
<td>1</td>
<td>20</td>
<td>gamma (1.5, 2)</td>
<td>3</td>
<td>6</td>
</tr>
<tr>
<td></td>
<td>$P_1$</td>
<td>0.1</td>
<td>3</td>
<td>10</td>
<td>gamma (1, 2)</td>
<td>2</td>
<td>4</td>
</tr>
<tr>
<td></td>
<td>$P_2$</td>
<td>0.1</td>
<td>1</td>
<td>1</td>
<td>gamma (1, 2)</td>
<td>2</td>
<td>4</td>
</tr>
<tr>
<td>GeneAssemble</td>
<td>$P_3$</td>
<td>0.1</td>
<td>5</td>
<td>25</td>
<td>gamma (5, 2)</td>
<td>10</td>
<td>20</td>
</tr>
<tr>
<td></td>
<td>$P_4$</td>
<td>0</td>
<td>10</td>
<td>1</td>
<td>gamma (5, 2)</td>
<td>10</td>
<td>20</td>
</tr>
<tr>
<td></td>
<td>$P_5$</td>
<td>0.3</td>
<td>1</td>
<td>10</td>
<td>gamma (10, 2)</td>
<td>20</td>
<td>40</td>
</tr>
<tr>
<td>Blast</td>
<td>$P_6$</td>
<td>0.3</td>
<td>2</td>
<td>50</td>
<td>gamma (5, 3)</td>
<td>15</td>
<td>45</td>
</tr>
<tr>
<td></td>
<td>$P_7$</td>
<td>0.8</td>
<td>0.1</td>
<td>50</td>
<td>gamma (10, 10)</td>
<td>100</td>
<td>1000</td>
</tr>
<tr>
<td></td>
<td>$P_8$</td>
<td>0.05</td>
<td>10</td>
<td>5</td>
<td>gamma (2, 1)</td>
<td>2</td>
<td>2</td>
</tr>
<tr>
<td>LookUp</td>
<td>$P_9$</td>
<td>0.5</td>
<td>5</td>
<td>10</td>
<td>gamma (1.5, 1.5)</td>
<td>2.25</td>
<td>3.375</td>
</tr>
<tr>
<td></td>
<td>$P_{10}$</td>
<td>0.5</td>
<td>4</td>
<td>2</td>
<td>gamma (1.5, 1.5)</td>
<td>2.25</td>
<td>3.375</td>
</tr>
<tr>
<td></td>
<td>$P_{11}$</td>
<td>0.75</td>
<td>5</td>
<td>10</td>
<td>gamma (0.5, 0.5)</td>
<td>0.25</td>
<td>0.125</td>
</tr>
<tr>
<td>Render</td>
<td>$P_{12}$</td>
<td>0.1</td>
<td>10</td>
<td>25</td>
<td>gamma (30, 3)</td>
<td>90</td>
<td>270</td>
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<tr>
<td></td>
<td>$P_{13}$</td>
<td>0.01</td>
<td>100</td>
<td>5</td>
<td>gamma (20, 2)</td>
<td>40</td>
<td>80</td>
</tr>
<tr>
<td></td>
<td>$P_{14}$</td>
<td>0.9</td>
<td>1</td>
<td>25</td>
<td>gamma (30, 3)</td>
<td>90</td>
<td>270</td>
</tr>
<tr>
<td>Translate</td>
<td>$P_{15}$</td>
<td>0.7</td>
<td>0.5</td>
<td>50</td>
<td>gamma (1, 1)</td>
<td>1</td>
<td>1</td>
</tr>
<tr>
<td></td>
<td>$P_{16}$</td>
<td>0.7</td>
<td>0.1</td>
<td>50</td>
<td>gamma (10, 2)</td>
<td>20</td>
<td>40</td>
</tr>
<tr>
<td></td>
<td>$P_{17}$</td>
<td>0</td>
<td>25</td>
<td>10</td>
<td>gamma (1, 1)</td>
<td>1</td>
<td>1</td>
</tr>
<tr>
<td>Fold</td>
<td>$P_{18}$</td>
<td>0.2</td>
<td>10</td>
<td>5</td>
<td>gamma (3, 30)</td>
<td>90</td>
<td>2700</td>
</tr>
<tr>
<td></td>
<td>$P_{19}$</td>
<td>0.05</td>
<td>50</td>
<td>1</td>
<td>gamma (3, 5)</td>
<td>15</td>
<td>75</td>
</tr>
<tr>
<td></td>
<td>$P_{20}$</td>
<td>0.75</td>
<td>1</td>
<td>1</td>
<td>gamma (50, 2)</td>
<td>100</td>
<td>200</td>
</tr>
<tr>
<td>Print</td>
<td>$P_{21}$</td>
<td>0.2</td>
<td>2</td>
<td>20</td>
<td>gamma (2, 3)</td>
<td>6</td>
<td>18</td>
</tr>
<tr>
<td></td>
<td>$P_{22}$</td>
<td>0.05</td>
<td>2</td>
<td>10</td>
<td>gamma (5, 5)</td>
<td>25</td>
<td>125</td>
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<tr>
<td></td>
<td>$P_{23}$</td>
<td>0.9</td>
<td>0.1</td>
<td>30</td>
<td>gamma (2, 3)</td>
<td>6</td>
<td>18</td>
</tr>
</tbody>
</table>

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distribution \( d_W \) with the resulting summed mean and variance to approximate the completion time. This allows us to estimate the overall expected utility as 
\[
\tilde{u} = 0.98 \int_0^\infty d_W(x) u_1(x) \, dx - 65.26 = 73.25.
\]

Next, figure 6 shows the outcome of the algorithm when the more urgent utility function \( u_2 \) is used. Here, it is clear that the strategy adapts appropriately to the changed deadline and higher workflow reward. Generally, the algorithm relies on higher redundancy, as exemplified by the lower Render task, where the consumer now starts with two services from \( P_{12} \), followed just minutes later by a third. In other cases, the consumer also switches to entirely different populations. For the Fold task, it now selects a fast and reliable, but expensive, service from \( P_{19} \), followed by more services from \( P_{18} \) and \( P_{20} \). This indicates that the algorithm now proceeds more aggressively, not only incurring higher service costs, but also significantly reducing the mean workflow duration, so that the overall deadline can be met.

In §4, we evaluate our approach experimentally.

### 4. Empirical evaluation

In order to evaluate the performance of our strategy, we implemented it using the Java programming language and applied it to a range of realistic problems, such as the one described in §3. To aid the scientist and visualize the workflow before and during execution, we also developed an interactive user interface, as shown in figure 7. This is divided into three main sections: the top part shows the workflow and its current progress; the middle panel shows global workflow statistics, including the predicted success probability and an estimated completion time distribution; while the lower part gives detailed information about a particular task and its service allocation over time. Moreover, to...
generalize our results over a range of environments with varying degrees of uncertainty, we decided to conduct an empirical analysis using a simulated grid system.

In more detail, we first generate 12 random service types and attach an average cost, duration shape and scale (of a gamma distribution) to each, which are drawn from the continuous uniform distributions $U_c(1, 80)$, $U_c(1, 40)$ and $U_c(1, 10)$, respectively. Next, we generate a random number of service populations for each type (drawn from the discrete uniform distribution $U_d(3, 10)$), and add to each a random number of service (drawn from $U_d(1, 200)$). These populations represent groups of services with the same performance characteristics. Then, we vary the characteristics between the populations of each type, based on the type-specific values above. More specifically, we generate the failure probability of a population by sampling from $B(F, 0.005)$, where $F$ is the overall average failure probability of services and $B(m, v)$ is a beta distribution with mean $m$ and variance $v$. Furthermore, we determine the cost of each service as the product $2yz$, where $y$ is the type-specific average cost and $z$ is sampled from $B(0.5, 0.05)$. This is repeated for the duration parameters, resampling $z$ for each. Finally, workflows always consist of 50 tasks, with precedence constraints that are generated by randomly populating an adjacency matrix until at least 25 per cent of all possible edges are added, ignoring any that would create a cycle. We define the utility function $u$ by awarding a utility of $u_{max} = 25000$ up to time step $t_{max} = 2000$ and then deduct $\delta = 37.5$ for each further time step until the utility is 0. Finally, we map each task to a random type.

For each run, we record the net profit our strategy achieves by selecting and then executing a workflow. To obtain statistical significance, we repeat this 250 times with different services and workflows, and carry out analysis of variance (ANOVA) and two-sample $t$-tests at the 99% confidence level. We have also repeated our
experiments using a large range of controlled variables and obtained the same broad trends as described below. To evaluate our strategy, we compare it with the following benchmarks.

— *local*(\(n, w\)). This represents approaches that select services myopically for each task by optimizing a weighted sum of QoS parameters. In more detail, we use the technique described by Zeng *et al.* (2004), weighing the cost, reliability and expected service duration equally.\(^2\) To include approaches that consider redundancy and service timeouts, the strategy selects the \(n\) best services and waits \(w\) time steps until selecting more services.

— *global*(\(w\)). This strategy selects a single service for each task, maximizing a weighted sum of global performance metrics. Here, we consider the overall workflow cost, reliability and duration, as outlined in Zeng *et al.* (2004) (weighing all parameters equally). Furthermore, we impose a budget constraint \(u_{\text{max}}\) and deadline 2666. To allow the strategy to react to failures, we include a timeout \(w\), after which a new service allocation is computed.

Using this set-up, our experimental results are shown in figure 8. In particular, figure 8a contains the results of our *flexible* approach and a number of representative *local*(\(n, w\)) strategies. The first of these, *local*(1, \(\infty\)), represents selection approaches that do not consider failures and so select only the single best service that is available. Clearly, this strategy is unsuitable where there are services failures: at \(\Phi = 0.1\), it fails to complete most of its workflows, making an overall loss.

Next, *local*(2, 150) and *local*(6, 150) are shown as two of the overall best-performing local strategies. These deal with failures by including redundancy and timeouts. For this reason, they perform better than the *local*(1, \(\infty\)) in most cases.

\(^2\) We have experimented with other weight distributions, but obtained similar or worse results.
environments, and the local(6, 150) strategy is able to achieve a high positive profit even when the failure probability reaches $\Phi = 0.6$. However, two disadvantages of these strategies are immediately obvious. First, neither of them clearly dominate—at low failure probabilities, local(2, 150) performs better as it makes a smaller investment by only invoking two services at a time, but when the failure probability rises above $\Phi = 0.3$, local(6, 150) begins to perform better, as higher redundancy is needed to complete the workflow on time. Second, both of them make a significant loss when the failure probability rises to high levels. This is due to their inability to identify unfeasible workflows and stop executing when they are likely to miss their deadline or when the cost of repeatedly invoking services begins to outweigh the benefit. To give an indication of the best possible performance achievable by any local(n, w) strategy, we also plot an upper bound (a hypothetical best local strategy), which was obtained by exhaustively testing a large range of parameters for n and w.

Finally, the figure also contains the average profit of our proposed flexible strategy. Clearly, it outperforms all local(n, w) strategies. This is due to its ability to automatically vary the levels of redundancy it employs and to select different services, depending on the current environment and the characteristics of its workflow. In some cases, it makes an overall positive profit when no other local(n, w) is able to do so. Furthermore, by explicitly predicting the outcome of the workflow, it avoids an overall loss in any of the environments tested here.

Figure 8b highlights similar trends for the performance of the global(w) strategies. Again, they are not only able to cope with service uncertainty to some extent, but are also sensitive to the right choice of parameter and make an overall loss when the failure probability rises to a higher level. In this case, the loss is not as severe as that of the local(n, w) strategies, due to the overall budget and deadline constraints that eventually cause the consumer to stop executing.
Summarizing these results, the best local strategy achieves a profit of $8550.74 \pm 324.79$ (averaged over all values for $\Phi$ shown in figure 8), the best global achieves a profit of $7841.02 \pm 365.00$, while the flexible strategy achieves $12552.12 \pm 311.44$. Therefore, in the environments tested here, our approach results in an average improvement of 46.8 per cent over current techniques.

5. Conclusions

In this paper, we have described a novel decision-theoretic service selection algorithm that uses service QoS information to choose between heterogeneous, but functionally equivalent, services and to introduce redundancy when they are particularly failure prone. Using an example from the bioinformatics domain, we have demonstrated how this algorithm can be used by scientists to select uncertain service providers. Finally, by conducting a thorough empirical evaluation over a range of environments, we have shown that our approach consistently outperforms current techniques, without requiring manually supplied redundancy and timeout parameters.

Our algorithm is most applicable in large-scale grid systems where consumers are able to discover and invoke the services of a large number of heterogeneous providers. Hence, it is not directly suitable for closed systems where service selection and invocation are typically performed by centralized middleware, but rather for the open, highly distributed systems that are supported by emerging service-oriented frameworks, such as the open grid services architecture and the semantic grid (Foster et al. 2002; De Roure et al. 2005). In such systems, our algorithm can execute as an additional decision-making layer on top of existing workflow engines, requiring no additional human input (apart from an appropriate utility function $u$ to encode the value of workflow), and thereby removing the burden of detecting and recovering from most failures from the user.

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