Scalable and Efficient Resource Discovery in P2P Grids^{*}

[Extended Abstract]

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ABSTRACT

Distributed computing enables us to harness all the resources and computing power of the millions of computers connected to the Internet. Therefore, this work describes the construction of an efficient and scalable resource discovery mechanism, capable of searching not only for physical resources (e.g. CPU, Memory, etc.), but also services (e.g. facial recognition, high-resolution rendering, etc.) and applications (e.g. ffmpeg video encoder, programming language compilers, etc.) from computers connected to the same Peer-to-Peer Grid network. This is done in a novel way by combining all resource information into Attenuated Bloom Filters, which also allows us to efficiently route messages in a completely decentralized unstructured P2P network (no super-peers). The research shows that previous P2P, Grid, and Cycle Sharing systems tackled this problem by focusing on each resource type in isolation, such as (physical) resource discovery and service discovery. Methods to minimize storage and transmission costs were also researched. The discovery mechanism was evaluated with a number of different test scenarios that varied resource distribution, resource values, topologies, etc. For comparison, we also evaluated the Random Walk discovery method which served as a baseline. The results were favorable over Random Walk, having higher query success rates with less hops while requiring increase in message size and storage space at each node (for routing information), thus attaining our objectives of effectiveness, efficiency, and scalability.

1. INTRODUCTION

There are millions of computers connected to the Internet¹ with more and more going online each day due to laptops,

9¹http://www.internetworldstats.com

netbooks, PDAs, and smartphones. With so many devices connected to the same network, distributed computing on such a large scale cannot be ignored. As such, resource sharing has become immensely popular and has led to the development of Grid and Peer-to-Peer (P2P) infrastructures dedicated to that purpose. These infrastructures ease the sharing of various types of resources, that range from simple files, to software offering different services, and even hardware like CPUs and Printers.

The most popular form of resource sharing across the Internet is File Sharing via Peer-to-Peer applications, occupying roughly 50%-90% of all Internet traffic.² A lot of work has been done in this area to create robust and scalable systems, capable of efficiently supporting a large number of users in a decentralized manner. P2P Infrastructures can be divided between those that do not perform any node organization (Unstructured systems), such as Gnutella [1] and Freenet [2]; and those that structure their nodes to improve message routing (Structured systems), such as Chord [3], CAN [4], and Pastry [5].

Grid and Cycle Sharing systems are similar in nature, as their objective is to perform large-scale parallel computations in scientific and commercial communities. While Grid systems harness the power of many interconnected networks of computers, which are usually centrally or hierarchically managed by the institutions that run them; Cycle Sharing systems take advantage of the many idle computers and game consoles already connected to the Internet, volunteered by home users.

Even though Peer-to-Peer and Grid systems are different, the literature [6–9] says that they will eventually converge. In this fashion, GINGER³ [10], or simply GiGi, is a P2P Grid infrastructure that fuses three approaches (grid infrastructures, distributed cycle sharing, and decentralized P2P architectures) into one. GiGi's objective is to bring a Grid processing infrastructure to home users, i.e. a "grid-for-themasses" (e.g. achieve faster video compression, face recognition in pictures/movies, high-res rendering, molecular modeling, chemical reaction simulation, etc.).

The common theme between these different systems is that

^{9&}lt;sup>*</sup>A scientific paper describing a preliminary version of this work was published and presented at the conference *INForum 2010* under the title "Scalable and Efficient Discovery of Resources, Applications, and Services in P2P Grids."

^{9&}lt;sup>2</sup>http://torrentfreak.com/bittorrent-dominates-internet-traffic-070901

^{9&}lt;sup>3</sup>Grid Infrastructure for Non-Grid EnviRonments

users have a task that they want to accomplish: share files in P2P file sharing systems; perform scientific calculations in Grids; or perform CPU intensive tasks over a massive amount of idle home user computers in Cycle Sharing systems. Tasks require discoverable resources that satisfy certain requirements that can range from almost no requirements (file sharing), to simple requirements (idle CPU), to complex requirements (free CPU with X much RAM, with at least Y much storage space, and with application Z installed). This is where the work described in this paper comes in, where the objective is to create an effective, efficient, and scalable discovery protocol of resources, applications, and services for inclusion in the GINGER project.

The rest of this work is structured as follows. In Section 2 we discuss similar systems that also provide service or resource discovery. Section 3 describes the architecture of $SERD^4$, while in Section 5 we show some relevant performance results. Section 6 concludes this paper offering final remarks.

2. RELATED WORK

This section can be divided into three main areas: i) efficient data representation where reducing the size of data storage and transmission is the objective, ii) resource discovery which only deals with the discovery of physical (e.g. CPU, RAM, etc.) or virtual (e.g. files) resources, and iii) service discovery where the main concern is discovering the services (e.g. facial recognition, high-resolution rendering, applications, etc.) provided by computers in a network.

2.1 Efficient Data Representation

Efficient Data Representation is important in this work because nodes have to store and transmit resource information about themselves and neighbors. Compression reduces the size of highly redundant information via a dictionary based (LZW [11]) or statistic based (Huffman coding [12]) encoding process. RSync [13] and the Low-Bandwidth File System [14] use Chunks and Hashing to divide data into chunks, calculating the hash of each chunk, and only transmitting those that have changed between versions of the same file. Erasure codes take another approach, and encode a message into a few symbols which can then be used to reconstruct a partially received message. Reperasure [15] uses this technique to provide data replication without storing full-replicas. The three techniques, although important, are not directly applicable in this work. The reduced message size cancels the need to compress messages or divide them into chunks. We also do not need to perform any forward error correction nor replicate data.

The final and most useful technique is a space-efficient probabilistic data structure called **Bloom Filters**, which efficiently test whether an element is a member in a set with the possibility of a false-positive occurring. A set $S = \{x_1, x_2, ..., x_n\}$ of *n* elements is stored in an array of *m* bits all initially set to 0. It must also use *k* different hash functions, each of which maps some element to one position in the *m* bit array. Because Bloom filters are implemented as bit arrays, the union of two sets can be computed by performing the OR operation between the two, while their approximate intersections can be computed using the AND operation. Insertion is performed by passing the element through each of the kdifferent hash functions and setting the resulting position in the m bit array to one. To test whether an element is in the set or not, it has to be passed through all hash functions and if all the resulting positions in the array are set to one, then the element has a high probability of being in the set. If any position has the value zero, then we know for definite that it is not in the set (no false negatives). The small false positive rate arises from the fact that when querying for an element that is not in the set, some hash functions may result in positions that were already used (have the value one) for a previously inserted item. Therefore, the more elements are inserted into the Bloom filter, the higher the chance of a query resulting in a false positive. Another shortcoming is the inability to remove an element from the Bloom filter, as simply setting the positions given by the k hash functions to zero have the side effect of removing other elements as well.

Bloom Filter variations exist to either extend their functionality or address some limitation. Counting Bloom Filters [16] allow both insertion and removal of elements by using an array of counters, instead of bits. In [17], Mitzenmacher shows that Compressed Bloom Filters can either occupy the same space but have a lower false-positive rate, or reduce their size and maintain their false-positive rate. Almeida et al. [18] created a Scalable Bloom Filter that dynamically grows in order to support the desired false-positive rate.

Finally, **Attenuated Bloom filters** were proposed in [19] to optimize search performance w.r.t. locality of objects. It uses an array of Bloom filters with depth d, where each row i, for $1 \leq i \leq d$, corresponds to the information stored at nodes i hops away. As the depth increases, more information will be stored in that Bloom filter row, making the respective filter more attenuated and resulting in a higher probability of false positives. Therefore, information closest to the node is more accurate, and less so the further away. The major advantage of this technique is that it permits us to efficiently locate objects up to d hops away, using as little storage space as possible (due to the Bloom filters) at the cost of a certain false positive rate. The disadvantage is that it *only* lets us search information about nodes up to d hops away.

2.2 **Resource Discovery**

Resource Discovery systems do a subset of what we want to accomplish with this work: locating physical or virtual resources to perform jobs. They can be split into three categories: Peer-to-Peer, Grid, and Cycle Sharing.

Peer-to-Peer systems do not distinguish between clients and servers; all nodes are equal and have no central coordination, making them decentralized. This leads to the various types of node topology organization: unstructured, structured, and hybrid. **Unstructured** system nodes are randomly connected to a fixed number of neighbors; there is no information about where resources are located so message routing has to be performed by flooding. Searching can be uninformed or informed. Uninformed searches use no addition information to route queries, they are either flooded to all neighbors (Gnutella [1]), or are forwarded to a randomly selected neighbor (Iamnitchi et al. [20]). Informed searches are more intelligent and route messages based on

^{9&}lt;sup>4</sup>Scalable and Efficient Resource Discovery

collected information, but require more memory. Lie et al. [21] and the learning-based technique in Iamnitchi et al. forward queries to nodes that have replied to similar requests. Another strategy called best-neighbor in Iamnitchi et al. [20] just forwards queries to nodes with the highest success rate. Structured systems, such as Chord [3] and CAN [4] organize nodes into a rigid structure, called a Distributed Hash Table (DHT), which enables efficient exactmatch query routing. Each node is assigned an identifier (key) which makes him responsible for all content (values) whose hash resolves to that key. Finally, Hybrid systems try to combine the best of both worlds without their disadvantages. Some systems in this category, like Pastry [5] and Kademlia [22], tend more towards structured systems, albeit with a less "rigid" structure, where any node belonging to a defined key subspace can act as a contact for those values. Others follow a more unstructured approach and use super-peers [23] that communicate between themselves on the behalf of less capable nodes (in terms of bandwidth or CPU performance), thus increasing routing performance.

Grid and Cycle Sharing systems share the same objective: to combine many geographically dispersed computer resources in order to perform tasks that require lots of CPU processing power, or that need to process huge amounts of data. Tasks like these are common when dealing with scientific, technical, or business problems. Grid systems can run in LAN environments such as that of a university, or in a much larger network compromised of interconnected networks that belong to different institutions, corporations, or universities. Condor [24] and Legion [25] are typical examples of such systems, where information about all resources are stored in a central component, known as the Matchmaker in Condor, and in Legion is divided into 3 subcomponents: the Collection, Scheduler, and Enactor. This central component receives job requests, tries to match their requirements to available resources, and reserve those resources while notifying the requester. Cycle Sharing systems rather operate over the Internet, which can be highly unreliable with variable connection quality. Another important difference is that anyone with a computer can join a cycle sharing project of interest (e.g. SETI@Home [26] or Folding@Home) and volunteer their resources during idle times. This brings the additional problem of unreliable peer connections and possibly forged results from untrusted peers.

2.3 Service Discovery

Service Discovery systems, like Resource Discovery, do the missing subset of this work: enabling the automatic detection of services provided by computers in small LAN environments, like home networks, or in large-scale enterprise networks, like a corporation or university. SLP [27] and Jini [28] use a client/server architecture, where servers collect service information and perform lookups for clients. SLP can function without directory servers using multicast to find services, but only in small LAN environments.

The systems presented by Goering et al. [29] and Lv and Cao [30] use a Peer-to-Peer architecture instead, with the objective of being able to function in ad-hoc networks. Goering et al. propose a service discovery protocol based on the use of Attenuated Bloom Filters, which provide a method to locate objects, giving preference to objects located nearby. It is simply an array of Bloom Filters of depth d, where each row represents objects at different distances which, in this case, is in term of hops. Each node has an Attenuated Bloom Filter for each of its neighbors, which is consulted when a query is received in order to send it in a direction it will have a higher chance of success. The first level of the Attenuated Bloom Filter corresponds to the services that are one hop away, the second to services two hops away, and so forth. Therefore, the larger the distance from the node, the more services will be contained in the corresponding Attenuated Bloom Filter which will increase the chance of false positives. Relying solely on Attenuated Bloom Filters gives this system a big limitation: only the services located up to d-hops away can be easily found. Lv and Cao resolve this drawback by having nodes more than d+1 hops away cooperate among themselves. Thus, when a query is received, it follows the same process of checking the Attenuated Bloom Filters of its neighbors like Goering et al, but if no services are found, then the query is forwarded to a node d+1 hops away where the search begins again.

3. ARCHITECTURE

The objective of this work is to enhance the resource discovery mechanism in GINGER [10], also known as GiGi, by making it completely decentralized and more complete. This completeness regards the system's ability to discover, not only basic resources (e.g. CPU, Bandwidth, Memory, etc.), but also specific installed applications (e.g. video encoders, simulators, etc.) and services (e.g. face recognition, high-res rendering, etc.). Because GiGi can be used in many different ways ("grid-for-the-masses"), it has to be flexible enough to run different types of jobs normally performed by home-users.

In order to cope with a dynamic peer population and high churn rate, this system uses an unstructured peer-to-peer approach to resource discovery, even though message routing may not have optimum efficiency. If a structured system were to be used, the messages needed to keep the structure intact with an unstable population, such as home-users, could possibly result in a high overhead. Attenuated Bloom Filters are used to enhance message routing and speed up resource location. Note that this solution is different to the systems mentioned in the Related Work because it combines all types of different resources into one discovery mechanism. It is especially different to the works [29, 30] that also make use of Attenuated Bloom Filters due to usage of one aggregated Attenuated Bloom Filter (explained next), and the fact that all the different types of basic resources, services, and applications are encoded in the Bloom Filter.

Each node in the network stores a cached version of the Attenuated Bloom Filters of their neighbors. This information is then merged into one single Attenuated Bloom Filter by inserting the union (OR operation) of all neighbor Bloom Filters at a certain depth k into depth k + 1 (Figure 1). The consequence of using an Attenuated Bloom Filter of, for example, depth d = 2 is that a node will only know about the resources of nodes up to 2 hops away. A solution for this problem is discussed further in Section 3.



Figure 1: Example of a node A creating a single Attenuated Bloom Filter by merging each Level i of its neighbors' Attenuated Bloom Filters into Level i + 1.

Discovery Mechanism. The discovery of resources, applications, and services (illustrated as a flowchart in Figure 2) will be performed in the following way. When a node receives a query, it will check its own information to see if it can satisfy the requirements. If it does, a reply is sent directly to the node that originated the query. If not, it goes through its aggregated Attenuated Bloom Filter, which contains the combined information from its neighbors Attenuated Bloom Filters. This way, we can quickly determine if the query cannot be satisfied with nodes up to d hops away, in which case it will be sent directly to a node d+1 hops away to restart the search. If the query can be satisfied with nodes at most d hops away, the node then needs to determine the direction to send the query for it to be resolved. This is done by checking all the cached Attenuated Bloom Filters of its neighbors to determine which one has the requested resources. If found, it then forwards the query to that neighbor. If not, then it is because the aggregated Attenuated Bloom Filter returned a false positive, which is mitigated by simply sending the query to a node more than d+1 hops away so it can be resolved. As each message is forwarded to a node, the sender adds his own ID to the resource query's Bloom Filter which keeps track of where the message has been sent. This Bloom Filter is cleared when a query jumps to a node d + 1 hops away. If any node received a query message and its ID is in the Bloom Filter, then there must have been a false positive and therefore the query should fail.

Dynamic Resources. Some resources are mostly static and do not change often, like the Operating System, CPU and Disk speed, certain application versions, etc. But there are other resources whose values can change quite often, such as amount of RAM occupied, amount of CPU in use, etc. For those cases, if we used a classic Bloom Filter then it would need to be rebuilt periodically since it does not support the removal of elements. More, this rebuilding procedure would require resending information about resources that are not expected to change, thus wasting bandwidth.

Therefore, instead of using a classic Bloom Filter to store the information about the dynamic resources, a separate Counting Bloom Filter is used. To compensate the fact that a Counting Bloom Filter occupies more storage space than a classic one, we use a smaller Counting Bloom Filter size (less precision), as the number of static resources is greater than dynamic ones. The usage of this new Bloom Filter mirrors that described in the previous sections: queries for dynamic resources use Aggregated Counting Bloom Filters instead



Figure 2: Flowchart of resource, service, and application discovery from Section 3



Figure 3: Example showing how resource queries are forwarded with an Attenuated Bloom Filter of d = 1. When a neighbor has information about the desired resource, such as Node 3, then query is forwarded to that peer, who in turn forwards the query to Node 6 which contains the resource. In another case, when there is no information about the desired resource in Node 1's area (consisting of Nodes 1, 2 and 3), then the query is forwarded to an Outer Limit Node 4, where the search is then restarted.

and are checked after the static Aggregated Bloom Filter. The difference is that when a dynamic resource changes, it is removed from all Attenuated Bloom Filters. Only if that alteration was drastic and affected the key used for the resource in the Bloom Filter (explained next) do the neighbors need to be notified. This is done by defining a periodic interval which checks for alterations to the resources in the main Attenuated Bloom Filter, which is then sent to the node's neighbors. Each neighbor also does this periodic check for alterations, and then resends its own Attenuated Bloom Filter with the changes to its neighbors. This continues until everyone is up to date. By using this periodic interval to send updates, we avoid wasted messanges and bandwith when resource values jitter.

Node Entry/Departure. For a node to join the network, it has to contact an already participating member. When the new node establishes a connection, the already existing member returns its Aggregated Attenuated Bloom Filter information. After the new node integrates this new information, it sends its own Aggregated Attenuated Bloom Filter to the already existing node which the updates its tables, and then sends the Aggregated Attenuated Bloom Filters with dynamic and static resource information to its direct neighbors. Those direct neighbors will eventually do the same until all proper neighbors are updated.

With regards to node departure/failure, each member of the P2P network periodically sends a Ping message in order to verify if its neighbors are still alive. If there is no response, then in the next periodic check, that neighbor's information is purged from the Aggregated Attenuated Bloom Filter and its cache is deleted. The node that detected the failure then needs to rebuild its Aggregated Attenuated Bloom Filter and resend it to its neighbors. Bloom Filters are then exchanged until all neighbors are up to date. Note that this system assumes that the TCP protocol is used for Ping messages so if there is no reply, we can consider the node has left the network.

Outer Limit Peer Discovery. Using an Attenuated Bloom Filter of a certain depth d limits the amount of information a node has about its surrounding neighbors. If a query is received and cannot be satisfied using the information the node knows about its peers in the same area, then it forwards the query to another node that is d + 1 hops away (which, conceptually, is part of another area).

To find outer limit peers, a simple random walk strategy is used, where a peer discovery query is forwarded to a random neighbor until it reaches a node l hops away, in which case a reply is sent directly to the originating node with contact information (e.g. IP address). If a node is not able to forward the discovery message to a node that has not seen the message before, then it replies to the originating node letting it know that the path did not lead to an outer limit node. The originator node then restarts the discovery process, this time sending it to a different neighboring node (this information is stored along with the query message) in order to try another path that might result in an outer limit peer.

3.1 Resource Representation

Information about resources, applications, and services that each node offers are represented inside a Bloom Filter. But, because a Bloom Filter is only capable of performing membership tests given a key (in this case a string), we need to add information about the actual resource (like type, value, etc.) to that key on insertion for it to be useful in discovering resources. Therefore, keys use namespaces to differentiate between resources and their values, which also helps with performing membership tests for resources. The naming convention uses a 3-level namespace, each separated using the colon (":") as a delimiter, and follows the following rules:

- Level 1: Name of the Resource, Service, or Application (e.g. CPU or ffmpeg)
- Level 2: Type of the Resource, Service, or Application (e.g. MHz or version)
- Level 3: Actual value of the Resource, Service, or Application

For instance, if we wanted to store the fact that a node has a CPU of 3 GHz, the key we would insert into the Bloom Filter would be: "CPU:GHz:3". Or, if a node has the application ffmpeg version 2.3 installed, the key would look like: "ffmpeg:version:2.3". But, for different nodes to be able to communicate with each other and search for the same resources, the naming of resources, services, and applications need to be the same between all of them. An ontology could be used, but that is out of the scope of this work. For the time being, the system reads a configuration file that specifies the name of the resource among other things. This configuration file needs to be the same for all nodes in the network.

Insertion. However, just following a naming convention will not suffice for the discovery of resources. We also need to take into account the values used for each resource. If we do not restrict the possible values, we would need to employ a brute force strategy when querying for resources, trying each value combination and testing the Bloom Filter. For example, to find a node that at least contains a CPU of 2.6 GHz, we would need to test for values such as 2.6, 2.7, 2.8, 2.9, 3.0, etc., which is highly inefficient. To speed this up, we define a *minimum*, *maximum*, and a *quantum* for each resource value type (which are also specified in a configuration file). The *minimum* (resp. *maximum*) is the smallest (resp. largest) value that the resource will have encoded in the Bloom Filter. The quantum defines how the value space, from *minimum* to *maximum*, will be divided. When a resource is inserted into the Bloom Filter, it is first inserted with the key that corresponds to its range, and then with all the other keys that correspond to ranges smaller than the resource's value. For example, if we define minimum = 0, maximum = 4000, and quantum = 1000 for CPU values in MHz, then the range of values is divided into the following segments: [0, 1000]; [1000, 2000]; [2000, 3000]; and [3000, 4000]. Or, if a CPU of 999MHz were to be inserted into the Bloom Filter, it would need to be inserted under the value 1000: "CPU:MHz:1000"; and so on.

Querying. Now, when querying a Bloom Filter for a value, the range the value falls under needs to be determined for the specified resource and checked. For instance, if a query requires a CPU of at least 2600 MHz, we would only need to perform one exact match query using the range the value in the requirements belongs to, which in this case is 3000 $(2600 \subset]2000, 3000]$). Therefore, we only need to test the key "CPU:MHz:3000" against a Bloom Filter because processors with a faster CPU will also be registered under this key. This strategy avoids the brute-force approach and efficiently speeds up the querying process. However, one needs to take care when specifying the quantum value due to precision problems. In this example, a CPU of at least 2600 MHz is required, but testing the Bloom Filter with key "CPU:MHz:3000" can result in CPUs that belong to the interval [2000, 2599], thus not satisfying the requirements. In a real-world system, using a quantum = 200 would probably be more suitable, giving enough precision without requiring too much overhead. This, and searching for a resource with a key one *quantum* value higher than required will ensure query satisfaction.

4. IMPLEMENTATION DETAILS

This work was implemented using the PeerSim [31] simulator with its Event Driven capabilities, approximating the simulation more to real-life as opposed to a Cycle Driven simulation. Because PeerSim is implemented in Java, the SERD discovery mechanism is also implemented in Java, which also allowed us to use an open source Bloom Filter implementation from the well known Hadoop project, providing us a certain amount of confidence w.r.t. its quality.

In order to be able to evaluate this work, we had to build an infra-structure around PeerSim to allow things such as topology creation, resource distribution, and node activity specification. The Topology Manager allows the generation of random topologies, with minimum and maximum number of neighbors and network size as parameters. It also allows the loading of existing topologies in two different file formats. Node Resource Description Language (NRDL) allows us to distribute resources among nodes either in a static way (specifying each node's resources), or in a more random fashion by specifying criteria to select a certain number of nodes to distribute the resource to. Distribution criteria can be the number of hops between nodes, the density/frequency of nodes that have the resource, or even the homogeneity of resource distribution. Node Activity Specification Language (NASL) allow us to script the actions of the virtual nodes where we can select nodes using various types of specifiers (e.g. randomly, exact match, nodes with a certain resource, etc.) along with the actions that they should perform (e.g. search for some resource) and when that action should be executed (in terms of simulation cycles or periodicity). The Scenario Manager combines all the previous components into one which allows us to save and load simulation scenarios, and thus easily reproduce experiment results.

Test scenario generation was performed using the Ruby utility rake and ERB which enabled us to automate test generation and embed Ruby code into the configuration files of the aforementioned components. This way, we were able to include complex logic during test generation. During the simulation of those tests, various metrics were collected in order to be analyzed later on. This was done using Peersim **Controls** that intercepted various calls from the discovery mechanism via hooks (or callbacks) and stored the data into a key-value backend called Redis.

5. EVALUATION

We created various test scenarios in order to determine if the SERD discovery mechanism was able to achieve its goals of being effective, efficient, and scalable. We also compared our system to another, albeit simpler, discovery mechanism called Random Walk which functioned as our baseline.

5.1 Test Scenarios

The test scenarios were generated using the components from Section 4, and were executed with the Random Walk protocol (RW) and three variations of SERD: SERD1, SERD2, and SERD3 which correspond to the Attenuated Bloom Filter depths of 1, 2, and 3, respectively. The test scenarios were generated with the following parameters and values (one set focused on static resources and the other on dynamic resources):

- *Network Size*: 5000 nodes and 10000 nodes (the topology statistics can be seen in Figure 4)
- *Number of Neighbors*: 3 neighbors and 6 neighbors per node, for each Network Size
- Resource distribution: 50% (very abundant resource), 25% (abundant resource, and 5% (scarce resource), for each combination of Network Size and Number of Neighbors

For each of the three resource distribution categories (very abundant, abundant, and scarce), two types of resources were distributed accordingly: one type with uniform values (almost no variation) and another with non-uniform values (with a lot of variation). The uniform resource chosen for the static tests was the GCC v4.2 application, and the availability of a node to be used exclusively for the dynamic tests (either the systems have the resource or do not). For the non-uniform resources, the CPU speed (MHz) was chosen with a minimum, maximum and quantum of 1000, 3000, and 1000, respectively for the static tests; whereas for the non-uniform resource in the dynamic tests, the resource Hard Drive storage (GB) was chosen with a minimum, maximum, and quantum of 0, 1000, and 50, respectively.

During each of the static and dynamic tests, every 5 cycles 10% of the nodes in the network sent resource queries that could be satisfied by at least one node in the network. The difference between the two types of tests was that in the dynamic scenarios, the resource values had to change over time. This was done by defining typical times of resource consumption of 5 cycles (short task), 10 cycles (typical task), and 20 cycles (long running task). Each of these values were then picked randomly until the total was at least 5 cycles less than the maximum defined for the Peersim simulation (100 cycles). At each point in that list, one third of each resource went down 20%, the other one third of the resources maintained their value, and the rest of resources were increased by 20% of their value.

Network Size	Degree Statistics				Graph Statistics	
	Max	Min	Avg.	Variance	Avg of clustering coefficients	Avg. Of distances to all other nodes
5000	3	3	3	0	2.00E-04	10.37341428
5000	6	5	5.9996	4.00E-04	0.001	5.185918544
10000	3	3	3	0	2.00E-04	11.37325297
10000	6	5	5.9996	2.00E-04	3.20E-04	5.60869937

Figure 4: Statistics about the generated topologies

Parameter	Value	Description	
Join Protocol Halt	1	Number of cycles of inac-	
		tivity to stop the join pro-	
		tocol	
Outer Limit	$log_2(NW)$	Maximum number of	
Jumps	SIZE)	outer limit jumps	
Attenuated Bloom	2	Period that defines when	
Filter Rebuild		filter should be rebuilt af-	
		ter receiving an update	
Dynamic Update	3	Period that defines when	
Period		to send resource updates	
Bloom Filter - N	100	Number of items to store	
Bloom Filter - P	$1.0e^{-9}$	False positive probability	
Couting Bloom	50	Number of items to store	
Filter - N			
Couting Bloom	$1.0e^{-9}$	False positive probability	
Filter - P			

Table 1: The parameters used for the SERD protocol during the tests with RW

5.2 SERD Protocol Parameters

The discovery mechanism in this work has many configurable parameters. As it would be impossible to test the effect of all parameters, we executed some preliminary tests in order to figure out reasonable parameters to use in the tests against the RW protocol. These tests were executed with two topologies of 10381, and 10000 nodes with 3 and 6 maximum number of neighbors respectively (both topologies were intended to have 10000 nodes, but due to the topology manager generation process, from Section 4, one topology resulted in 10381). Every 5 cycles 10% of the nodes send queries for resources that can be satisfied by at least one in the population, for each resource. The resource included both the static and dynamic resource categories (*uniform*) and *non-uniform*) where the distribution was the worst possible: 5 percent (scarce). The parameter values used, along with their respective description, can be seen in Table 1.

5.3 Result Analysis

During the execution of the many test scenarios we collected various metrics in order to help evaluate the effectiveness, efficiency, and scalability of our system. These metrics we collected for the static and dynamic tests were:

- Resource Query Satisfaction
- Average Number of Resource Query Hops
- Total Number of Sent Messages
- Average Size of Storage at each Node, and Message Size

Static Scenario Results. With regards to the satisfaction of resource queries (Figure 5), SERD1 and SERD2 consistently got a percentage rate above 90% except for the scarce scenarios with a maximum of 3 neighbors. This can be explained by the fact that the depth of the Attenuated Bloom Filters did not allow the forwarding of queries with much hindsight, especially in a scenario where very little nodes actual contain the resource and where each node only has a maximum of 3 neighbors, thus further limiting a node's knowledge about the network. SERD3 in almost all scenarios had a satisfaction rate of 100%, and in others 99%. As the algorithm had a greater depth, it was able to direct queries in the right direction for them to be satisfied. The RW algorithm's lack of intelligence in the forwarding of queries is a great contrast, with almost all satisfaction rates below or around 80%. While it performs better in scenarios where the resources are abundant, it suffers in the scarce ones.



Figure 5: Query Satisfaction for Static Scenarios

The number of hops a query messages takes in order for it to be satisfied is another important aspect in a discovery system, which needs to be as low as possible due to network latency. As we can see in Figure 6, RW queries were consistently higher than any of the SERD protocols because of the lack of query success, which made the query reach the maximum number of hops (Outer Limit Jumps value) and fail. SERD1 to SERD3 all had an average below 3 hops in all tests except for the *scarce* ones with a maximum of 3 neighbors. In those cases, SERD1 performed the worst, while SERD3 the best. This can be explained by the fact that the lack of resource knowledge (defined by the Attenuated Bloom Filter depth) made queries take non-optimum routes while looking for the resource, or even fail.

In Figure 7, we can see the total messages sent by each protocol. It is to be expected that in this case, the RW protocol typically uses a lot less messages because it does not have to trade resource information. The cases where RW uses more messages than any of the SERD protocol is because of the low query success rate, which means that there were a lot of messages that traveled until the maximum depth. It is also to be expected that SERD3 uses more messages than SERD1 or SERD2, especially in the scenarios with 6 maximum neighbors, due to the greater Attenuated Bloom Filter depth. Its depth and the amount of neighbors each node



Figure 6: Average Query Hops for Static Scenarios

has knowledge of influences greatly the joining phase of the discovery process where Attenuated Bloom Filters have to be traded among nodes until everyone is up-to-date.



Figure 7: Total Sent Messages for Static Scenarios

Figure 8 confirms what we already expected: the greater the Attenuated Bloom Filter depth, the higher the storage costs at each node and the bigger the message size due to the trading of resource information. The RW protocol uses so little storage space that it does not appear on the graph (average of 483.38), which is normal as it has no information about neighboring nodes and shows in terms of query satisfaction. Nonetheless, in a real scenario, RW would have to store increasingly larger information regarding local resources at each node, which unoptimized would occupy much space. SERD not only keeps information about its own resources, but also caches the Attenuated Bloom Filters of its neighbors. Note that these results do not depend on the number of items actually stored in the Bloom Filters as they have a fixed size (defined in Subsection ??).

Dynamic Scenario Results. Figure 9 shows us the query satisfaction for the dynamic resource scenarios, which are expected to not be as high as the static scenarios due to the varying values of the resources. Once again SERD outperformed the RW protocol, which display a success rate of 80% and lower. In almost all tests, the SERD protocols were above 80%, except for the scarce scenario tests. In



Figure 8: Average Message Size and Storage Size for Static Scenarios

those, SERD1 struggled the most seeing as it hardly has information about the neighborhood. SERD2 and SERD3 only displayed a satisfaction rate lower than 80% when the *scarce* scenario was combined with a maximum of 3 neighbors, which limited the available options when forwarding query messages. RW in those cases was hardly able to reach 20% query satisfaction, making its lack of intelligence ever so apparent.



Figure 9: Query Satisfaction for Dynamic Scenarios

With not so stellar satisfaction results, the RW protocol in Figure 10 shows that with high hop averages, being mostly around or higher than 6 hops. The SERD protocols continued to show consistency with lower hop averages, although they had a higher increase in the scenarios where query satisfaction was lower than usual. An increase is query failures leads to a higher amount of hops as queries only fail if they reach the maximum Outer Limit Jumps.

Contrary to the static scenarios, where there were cases that the RW protocol consumed more messages than the SERD protocol, in the dynamic scenarios (Figure 11) SERD consistently used much more messages than RW. This is not at all surprising given that not only do nodes exchange resource information when new peers join the network, but also resource information when the dynamic resources change values during the simulation. The Figure display an interesting



Figure 10: Average Query Hops for Dynamic Scenarios

result: no matter the resource distribution for each topology, the number of sent messages stayed more or less the same. Another interesting result is that the topology of 5000 nodes with a maximum of 6 neighbors, and the 10000 topology with 3 maximum neighbors did not vary that much. Even though the former topology has less nodes, it sent more messages due to the bigger number of connections; whereas the latter has more nodes sending messages, but were doing so to a smaller number of connections.



Figure 11: Total Sent Messages for Dynamic Scenarios

Figure 12 does not present us with any new information and just confirms what happened in the static scenarios: the deeper the Attenuated Bloom Filter, the bigger the storage requirements are and the bigger the messages sent in the network are.

6. CONCLUSION

GiGi [10] allows home users to take advantage of Grid computing which was previously only available to scientific and corporate communities. Tasks that would usually take a lot of time, such as audio and video compression, signal processing related to multimedia content (e.g. photo, video, and audio enhancement), intensive calculus for content generation (e.g. ray-tracing, fractal generation), among others, can now be sped up by parallelizing and distributing them over



Figure 12: Average Message Size and Storage Size for Dynamic Scenarios

many computers.

However, to distribute the tasks GiGi needs to locate the resources that satisfy task prerequisites. Therefore, the architecture presented in this work is a discovery mechanism capable of locating physical resources, services, and applications from many computers connected to the same P2P Grid. This is done in a novel way by storing all resource, application, and service information in Attenuated Bloom Filters. We created a decentralized discovery mechanism that is efficient and scalable for the GiGi project and uses an unstructured P2P network in order to accommodate a highly dynamic node population. Even though this work addresses the GiGi project, it is completely independent and can be used in other types of networks, such as cycle-sharing networks.

In conclusion, the SERD discovery mechanism described in this dissertation performed well in the various test scenarios that included static and dynamic resoures, and outperformed the RW protocol which was our baseline. Our system proved to be effective in locating various types of resources, and scalable as the number of nodes in the network did not affect the mechanism's resource query satisfaction. The results obtained are encouraging towards our objective of efficiency, taking into consideration that the more resources each node has (expected in real case scenarios), the more space RW will use and thus incur a higher storage cost than SERD (with the Bloom Filters). Although message size in SERD is larger than RW, it is also able to satisfy a lot more resource queries than RW. Taking these points into consideration, we also conclude that SERD is an efficient discovery mechanism.

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