



## Basic networks: Definition and applications

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### ABSTRACT

We define basic networks as the undirected subgraphs with minimal number of units in which the distances (geodesics, minimal path lengths) among a set of selected nodes, which we call seeds, in the original graph are conserved. The additional nodes required to draw the basic network are called connectors. We describe a heuristic strategy to find the basic networks of complex graphs. We also show how the characterization of these networks may help to obtain relevant biological information from highly complex protein–protein interaction data.

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### 1. Introduction

In biology, as in other sciences, many types of complex data can be expressed as undirected graphs, in which units (nodes) are connected by bidirectional edges. Among the most significant examples are protein–protein interaction networks, networks based on cooccurrence of protein domains or graphs generated by assigning edges to genes with a given level of coexpression (Aittokallio and Schwikowski, 2006; Sharan et al., 2007). All these graphs have emerged very recently, with the advent of genomic and proteomic high-throughput technologies that have generated huge amounts of data. How to obtain useful information from this kind of data is one of the main challenges in modern biology.

A significant problem occurs when a set of units of one of those graphs is selected by any given criteria and we are interested in determining its functional or structural meaning. There are two ways of tackling this situation. First, we can exclusively use information not included in the graph. For example, if we are trying to understand the function of a set of proteins, we can simply explore the available literature or, if possible, we can use standard classifications, such as gene ontologies (GOs), in which gene products are classified according to the biological processes in which they act (Ashburner et al., 2000). These approaches are very useful if the available information is sufficient, but may fail if the set contains many proteins of unknown function. A second alternative is to explore the graph to determine the context in which the selected units are working. For example, in a

protein–protein interaction network, we can determine whether most of the selected proteins are located in a given region of that network. If this is the case, we can then use the information provided by proteins not included in our set, but closely linked to ours, to infer the functions that characterize our dataset. Conceptually, this second strategy can be schematized as follows. First, our set defines a group of *seed* nodes from which to start the analysis. Second, we characterize a second group of nodes which are sufficiently closely linked to the seed nodes, which may be called *connectors*. Third, we use the information provided by the connectors to infer information about the seeds. The problem then consists in defining the best way to determine the connectors. In the current literature, this has been generally solved by devising strategies to define modules (e.g. Del Rio et al., 2001; Bader, 2003; Asthana et al., 2004; Hashimoto et al., 2004; Krauthammer et al., 2004; Arnau et al., 2005; Scott et al., 2005; Lubovac et al., 2006; Lucas et al., 2006; Li and Horvath, 2007; see review by Sharan et al., 2007). A module is loosely defined as a group of closely linked nodes, with an internal cohesion that allows its separation from the rest of units in the network. The problem is that the characterization of modules is based on conventional, a priori criteria of unknown efficiency. The ability of any module definition to efficiently characterize connectors will depend on the features of the graph (size, connectivity) and the proximity among the seeds, and different definitions may lead to quite different results (e.g. Luo et al., 2007). In fact, to define modules may be difficult if the seeds are very distant.

The starting point of our work is the appreciation that any strategy able to determine significant connectors which is based on non-conventional criteria could be a significant advance in network exploration. We introduce here the concept of *basic*

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network and we show that it allows for a non-ambiguous definition of significant connectors. We describe a heuristic strategy to characterize the significant connectors and also, using real examples, we show the usefulness of the characterization of basic networks to explore complex biological graphs.

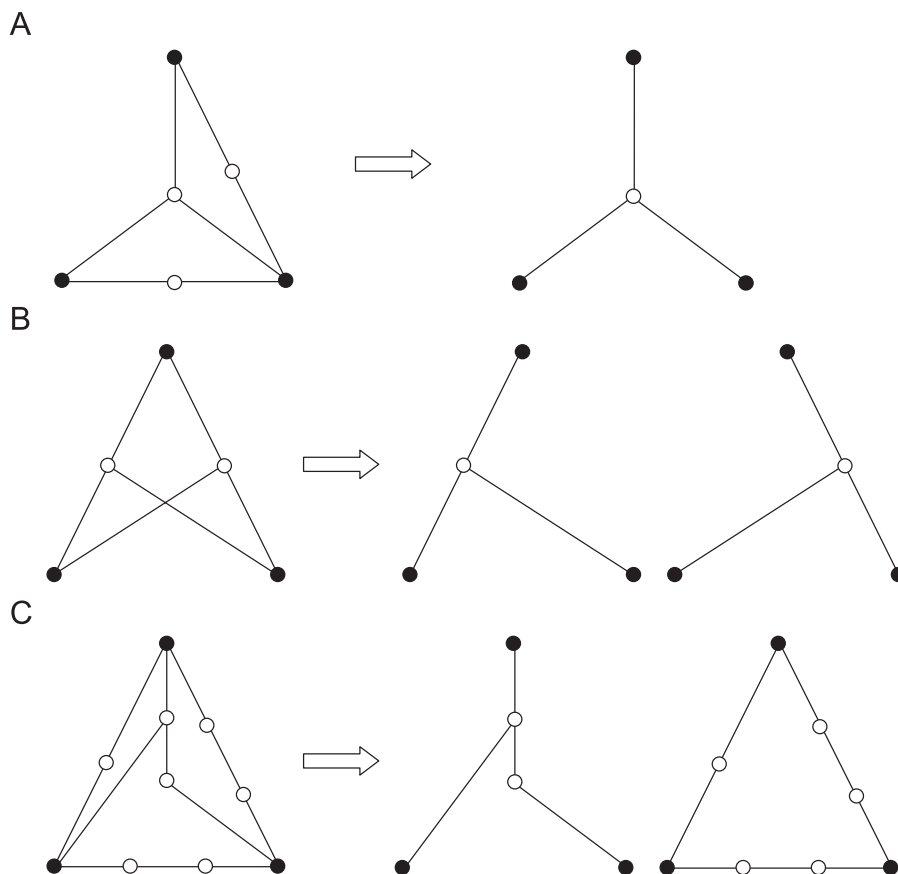
## 2. Concept of basic network and some theoretical considerations

We define *basic network* of an undirected, fully connected graph as the fully connected, undirected subgraph which includes a set of preselected units (seeds) plus the minimal number of connectors required to conserve the distances (geodesics, minimal path lengths) found among the seeds in the original graph. We define as *basic units* the set of seeds plus the set of connectors that are required to generate the basic network. This simple definition lacks any ambiguity: for any starting graph and set of seeds, there is at least one basic network (see e.g. Fig. 1A). However, it will occur often that more than one basic network, all with identical number of connectors, can be obtained. A typical example of *tied basic networks* is shown in Fig. 1B. A tied basic network can be further simplified (*pruned*; see below) to obtain several basic networks (Fig. 1B). It is very simple to demonstrate that randomly eliminating nodes from the graph will not necessarily lead to obtaining the basic network. On the contrary, random elimination of nodes will generally lead to subgraphs that cannot be further minimized but are much larger than the basic network of the graph (Fig. 1C). We call these results *local minimal networks*. Thus, any strategy to determine the basic network(s) of a graph will

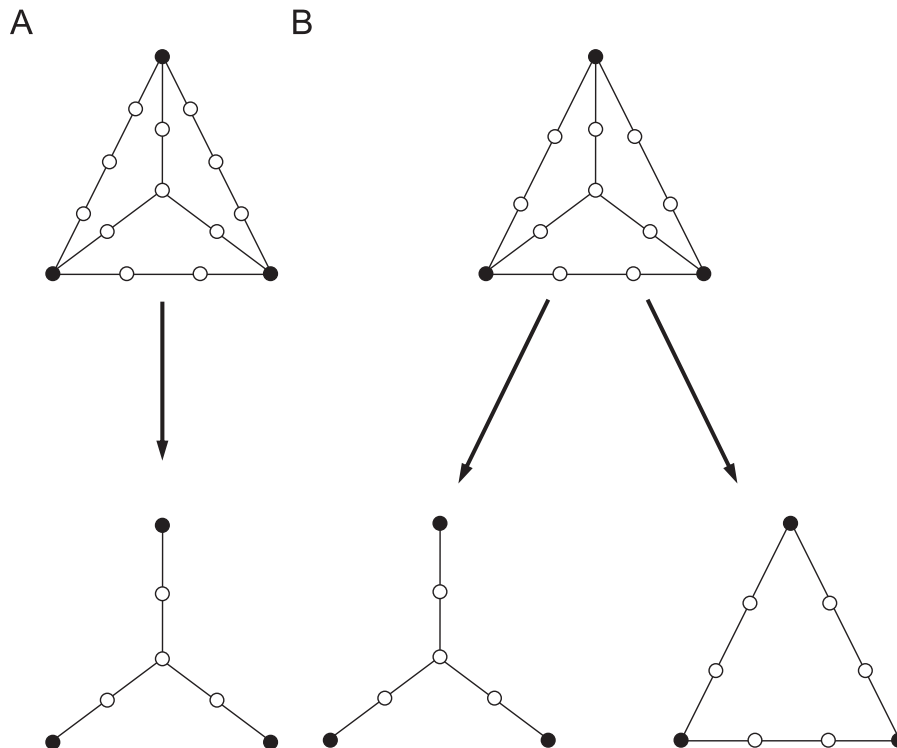
have to effectively deal with the presence of local minimal networks.

A significant theoretical consideration is the relationship among Steiner trees and basic networks. In the context of undirected graphs in which all edges are equivalent, a Steiner tree for a given set of seeds is defined as the tree that minimizes the number of edges (i.e. the total distance) to connect those seeds (Chartrand and Zhang, 2004). Although the goal of both Steiner trees and basic networks is to establish minimal subgraphs determined by the disposition of certain seeds in a starting graph, it is easy to demonstrate that they are often unrelated. The main reason is that, to obtain the basic networks, we demand the minimal distances among seeds to be conserved, while this is not the case for Steiner trees. Thus, although the basic network of a graph can coincide with the Steiner tree (Fig. 2A), it will most often occur that the basic network is larger and totally different from the Steiner tree (Fig. 2B). As can be seen in Fig. 2B, and can be easily visualized in larger graphs, Steiner trees are based on finding “internal” units (i.e. units that are quite away from the seeds and more or less equidistant from all of them) to minimize the total path length. Thus, in Steiner trees the seeds may end up at large distances. On the contrary, basic networks select “external” units, which are as closely linked to the seeds as possible.

We may now ask why the definition of basic networks may be significant in the exploration of complex graphs. There are three main reasons. First, the close proximity among seeds and connectors just mentioned is obviously a useful feature, if we wish to obtain information about the seeds. Second, as we already



**Fig. 1.** Three situations that may occur when searching for basic networks: (A) in the simplest case, a single basic network exists, and all connectors are basic units; (B) a tied basic network cannot be further simplified without pruning (left). When pruning is performed, multiple alternative basic networks are found (right). (C) Randomly eliminating nodes may lead to the basic network (left) but, most often, will lead to local minimal networks (right; this would be the optimal solution if any of the two internal nodes is eliminated).



**Fig. 2.** Relationships between Steiner trees and basic networks: (A) when the only geodesic among the seeds is through the Steiner tree, the basic network and the Steiner tree are identical; (B) a typical example in which the Steiner tree (left) is totally different from the basic network (right), for the same starting graph and starting seeds (black dots). The shortest paths do not include the nodes in the Steiner tree.

discussed above, the definition of basic network is not ambiguous, avoiding the potential pitfalls of all module definitions. This has another direct advantage: the fact that basic networks have a given, clear-cut number of seeds allows for their statistical comparison with *random basic networks*, which are those obtained by randomly taking units from the same starting graph. Thus, we can demonstrate that some seeds are unrelated if the basic network obtained starting with them is not significantly different in size from a set of random basic networks. Finally, by determining the minimal number of basic units, we focus on the nodes that best explain the connections among the seeds. These must necessarily be hubs, highly connected units, given that only hubs may contribute to minimize its size.

### 3. A heuristic strategy to characterize basic networks

It is obvious that it is impossible to test for all the possible combinations of connectors to obtain the basic network of any large graph. Thus, a heuristic strategy must be devised to obtain the basic network, or at least to get a good approximation to it, avoiding falling in local minimal networks. We have devised a strategy which is implemented in a program that we have called *Netbasic*. Although the algorithmic details of the program are complicated, and will be described elsewhere, the strategy is actually quite simple to understand. The basic pseudocode of the program is as follows:

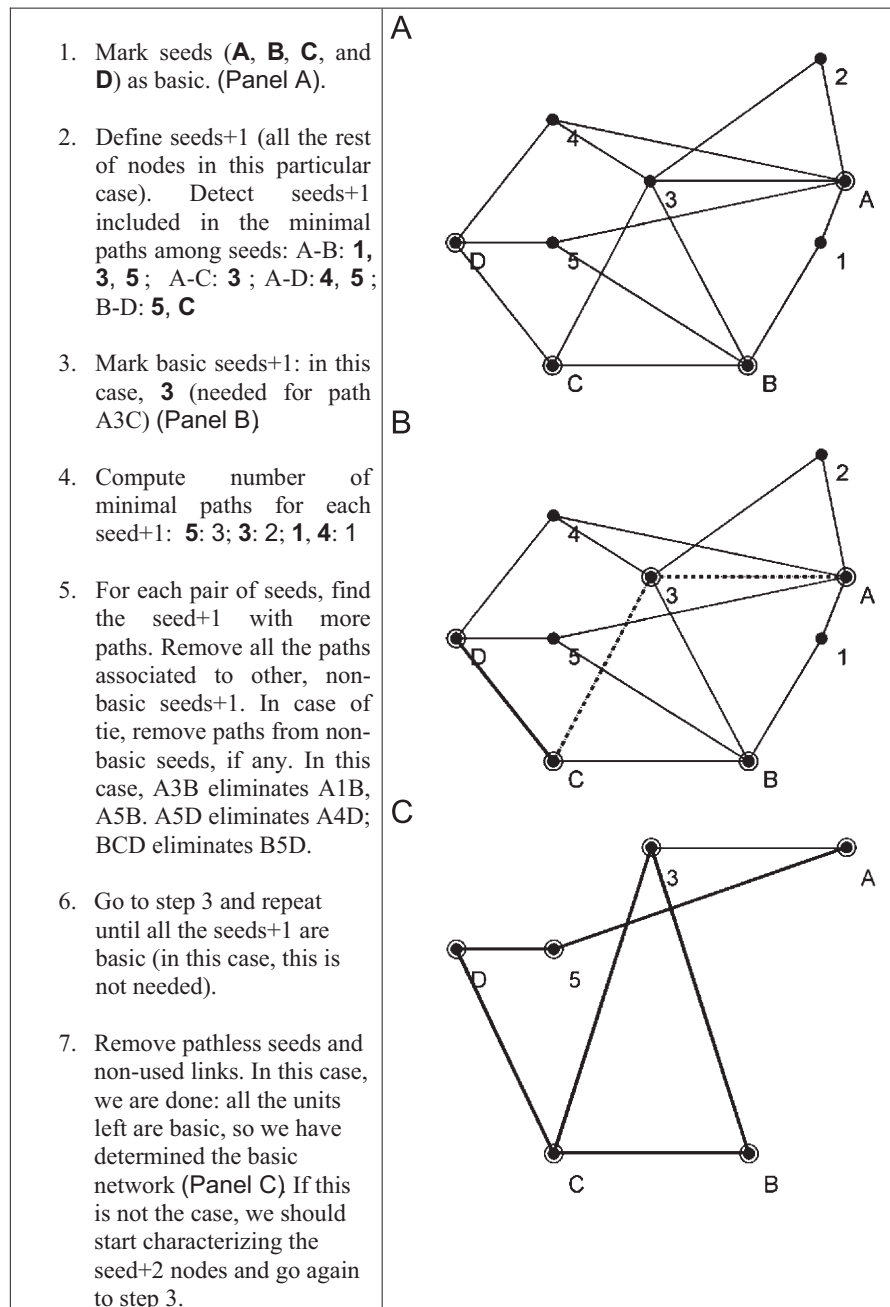
- (1) Mark seeds as basic units.
- (2) Compute the distances among the seeds.
- (3) For each seed, define seed+1 units as the second node in any geodesic that starts in that seed and ends in a different seed.
- (4) Determine all the basic seed+1 units, which are those that have no alternative seed+1 units and therefore are essential to

connect two or more seeds with a minimal number of steps. Mark them as basic units.

- (5) Remove paths among seeds, following these rules:
  - (a) If a seed+1 unit is basic, none of the paths in which it is included can be eliminated.
  - (b) Count paths among seeds for each seed+1 unit. If two or more seed+1 units have alternative paths, which connect the same two seeds, but one of them has assigned more total paths than the rest, the paths that pass through the rest of seed+1 units are eliminated.
  - (c) In case of ties in the number of paths, all paths are kept unless one of the tied seed+1 units is basic. If that occurs, alternative paths through non-basic units are eliminated.
  - (d) Determine again whether seed+1 units are basic.
  - (e) Repeat steps 5a–c until there are no more eliminations.
- (6) Remove nodes that lack paths.
- (7) If needed, define seed+2 units (and, if needed, seed+3 units, seed+4 units, ..., seed+n units) and repeat the process described in steps 4–6 until all nodes are either marked as basic or no additional node can be eliminated.

Steps 1–4 are easy to follow. Most critical is step 5. In it, seed+1 units are compared and paths that go through basic or highly connected units (i.e. those that are included in many paths among seeds) are conserved, while those that go through non-basic, poorly connected units are eliminated. The same applies to step 7, for seed+2 units, seed+3 units, etc. These steps very effectively eliminate most units that are very unlikely to be part of the basic network, given its scarce participation in connecting the seeds. Thus, this strategy is based on keeping hubs and eliminating less connected units. A simple example of how this strategy works is shown in Fig. 3.

Even though this heuristic search is logical, it is obviously not certain it will obtain the basic network. As we already described



**Fig. 3.** An example of how to obtain basic networks. Units are named in bold. **A**, **B**, **C**, **D** are the seed units. The whole graph contains only nine units. Basic units are indicated with double circles.

above, the two complications are (1) ties and (2) local minimal networks. To deal with them, we have included refinements to the basic strategy. First, after the program goes through steps 1–7, it determines whether all nodes left are basic. If not, this may be caused by tied solutions as those that we shown in Fig. 1B. If this is the case, the program can, if the user is interested, to further minimize the graph by checking in turn which connectors can be eliminated without affecting the distances among the seeds, in a process that we have called *pruning*. Once the pruning has been completed, ties are eliminated. However, the pruning process also eliminates the information about alternative solutions of the basic graph and, in many cases, it may be interesting to consider the full solution (see below). The second complication is much more difficult to cope with, given that to avoid the problem of local minimal networks with any heuristic search is

intrinsically impossible, except in trivial cases. Thus, we have chosen to minimize the problem by using a permutation method. In steps 5 and 7, it is obviously critical the order in which seed+n units are considered. The process of path elimination depends on counting paths, and that count varies each time that a seed+n unit has been processed. Thus, we can easily envisage a situation in which taking the seed+n units in a given order will lead to a local minimal network, while considering them in a different order will lead to obtaining the basic network. By randomly taking the seeds and performing many alternative runs of the program ( $10^3$ – $10^5$ ), we can avoid as much as possible this problem. The multiple runs also give us a clear idea of the variation among the solutions obtained, and which is the set of proteins that appear in all the solutions and what proteins appear less frequently.



**Table 1**  
Characterization of the basic networks obtained when the proteins annotated to several GO terms are used as seeds and comparison with random basic networks obtained starting from the same number of random seeds.

GO term	GO domain	No. of seeds	No. of connectors in the basic networks tied/pruned	No. of connectors in random basic networks tied/pruned	Z-scores tied/pruned
Mitochondrial large ribosomal subunit	Cellular component	43	22 ± 1 (20)/19 (19)	207 ± 27 (163)/206 ± 27 (163)	6.85/6.92
Mitochondrial small ribosomal subunit	Cellular component	33	11 ± 1 (10)/10 (10)	150 ± 23 (109)/149 ± 22 (109)	6.04/6.32
Mitochondrial ribosome	Cellular component	76	36 ± 2 (34)/34 ± 1 (34)	382 ± 37 (298)/381 ± 37 (298)	9.35/9.38
Spliceosome	Cellular component	79	19 ± 2 (17)/19 ± 1 (17)	402 ± 37 (339)/401 ± 36 (339)	10.35/10.61
Phosphoprotein phosphatase activity	Molecular function	46	151 ± 2 (149)/151 ± 2 (149)	220 ± 27 (171)/219 ± 26 (171)	<b>2.56/2.58</b>
Protein kinase activity	Molecular function	127	381 ± 1 (380)/381 ± 1 (380)	637 ± 36 (551)/637 ± 36 (551)	7.11/7.11
Protein folding	Biological process	72	224 ± 2 (221)/224 ± 2 (221)	362 ± 32 (301)/361 ± 31 (301)	4.31/4.42
Sexual reproduction	Biological process	110	403 ± 3 (399)/403 ± 3 (399)	560 ± 33 (499)/560 ± 33 (499)*	4.76/4.76
Sporulation	Biological process	110	525 ± 3 (522)/525 ± 3 (522)		<b>1.06/1.06</b>

In all cases, 1500 permutations of the seeds were performed. Values are expressed as mean ± standard deviation and, in parenthesis, the minimum number (i.e. the optimal solution for the essential network characterized after obtaining all replicates). Z-scores (calculated as the absolute value of the difference between the average sizes of the random and experimental networks divided by the standard deviation of the random networks) are also indicated. The asterisk refers to the fact that, the number of seeds was identical in the two last GO terms, a single characterization of random basic networks, with number of seeds equal to 110, was performed. Thus, the Z-scores obtained for the GO terms *sexual reproduction* and *sporulation* are based on the same random basic networks. Values in bold are not significant, after Bonferroni's correction.

links is a good evidence for strong functional relationships of those connectors with the seed proteins.

Very interestingly, the other 16 connectors are both unrelated to ribosome function (according to SGD) and are only linked to at most a few of the large subunit proteins (Fig. 4). These results can be explained noticing that there is an outlier among the seeds: the protein MET13, which is far apart from the rest of the proteins in this GO term (Fig. 4). The fact that MET13 is not closely linked to the rest of seeds explains the need to include those 16 additional connectors. MET13 is a methyltetrahydrofolate reductase involved in methionine biosynthesis and therefore its relationship with the mitochondrial ribosome is unclear. According to SGD, MET13 was purified together with units of the mitochondrial large ribosomal subunit by Kitakawa et al. (1997). However, a direct interaction with any of those units has not been yet described. Its purification may thus have been artifactual.

To generalize the results obtained with the large subunit of the mitochondrial ribosome, we performed related searches with other sets of proteins, defined according to diverse GO terms. Results are summarized in Table 1. Significantly, in all cases that we studied, the tied and pruned basic networks were very similar (Table 1). This is due to the fact that most units in the tied basic networks were basic units and the number of ties was therefore very low. If we now consider each GO term in detail, we can easily conclude that they can be divided into two classes. Some of them, especially those that are included in the *cellular component* GO domain (as the *mitochondrial large ribosomal subunit* term just described), are known to have a strong structural basis, that is, many of the proteins in a given term interact. They all generated basic networks which are much smaller than those generated by random seeds (Table 1; ribosome- and spliceosome-related terms). On the contrary, other GO terms, especially some of those included in the *molecular function* and *biological process* GO domains, generate basic networks that are either larger or not much smaller than those obtained starting with random seeds (see data in brackets in Table 1). In two cases, the sizes of the random basic networks were statistically not significantly different from those of the GO-based basic networks (see Z-scores in Table 1). These results demonstrate that these GO terms are not

characterized by including gene products which are closely linked in the protein-protein interaction network.

## 5. Discussion

We think that our definition of basic network is conceptually significant. First, it is an intuitively simple concept. Second, it provides a natural way to obtain groups of nodes related to a set of predefined seeds that escapes from any, more or less controversial, module definition. Finally, the results described above show that it is useful in very different contexts. On one hand, basic networks can be used to determine whether units are related or not. We have shown that this can be done by comparing the minimal graphs obtained with those units as seeds with the minimal graphs obtained with the same number of seeds, but randomly taken from the graph. On the other hand, basic networks point out proteins which are very closely linked to the seeds, for which a function related to the function of the seed proteins is likely (as we have discussed above for the connectors MHR1, MRP4, YLH47 and MDM38). Alternatively, by defining groups of nodes that are closely linked to the selected seeds, the determination of basic networks may contribute to understand what those seeds have in common. A final aspect is that basic networks may provide clues about potentially false connections in the graphs. When adding a single unit to the set of seeds involves a substantial increase in the number of connectors, we may suspect that the unit is unrelated to the rest of seeds. This can be quite easily observed by depicting the basic network: the offending unit stands out as very distant from the rest (e.g. the case of MET13 in Fig. 4).

Our solution to the problem of obtaining basic networks, the strategy described above is, given the impossibility of analyzing all possible combinations of units in large graphs, a heuristic search based on conserving the nodes with the highest number of paths traveling among seeds. As any other heuristic search, it does not guarantee finding the true basic network. However, the cases examined suggest that, in general, the solutions obtained will be very similar if the seeds are close in the graph (see the very low standard deviations in Table 1). In any case, the user may perform

a large number of trials, in order to obtain a progressively improving approximation to the basic network. Saturation, i.e. lack of further minimization of the size of the graphs after a large number of trials, may indicate that the basic network has been already obtained. Further exploration of complex graphs may determine the usefulness and limitations of this and other potentially competing heuristic searches for basic network characterization.

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I.M. developed the concept of basic networks and devised the heuristic strategy to characterize them and wrote the manuscript. S.H. wrote the Netbasic program to implement that strategy and performed the analyses. This project was supported by grant 2007201021 (Proyectos intramurales especiales, CSIC, Spain) and project BIO2008-05067 (Programa Nacional de Biotecnología; Ministerio de Ciencia e Innovación, Spain) awarded to I.M. We would like to thank Antonio Marco for his help in the development of the examples shown in this study.

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