

LABORATORY CLASSES

Implementation a simplistic version of the AGDS or AANG network for some inference reasons

AGH



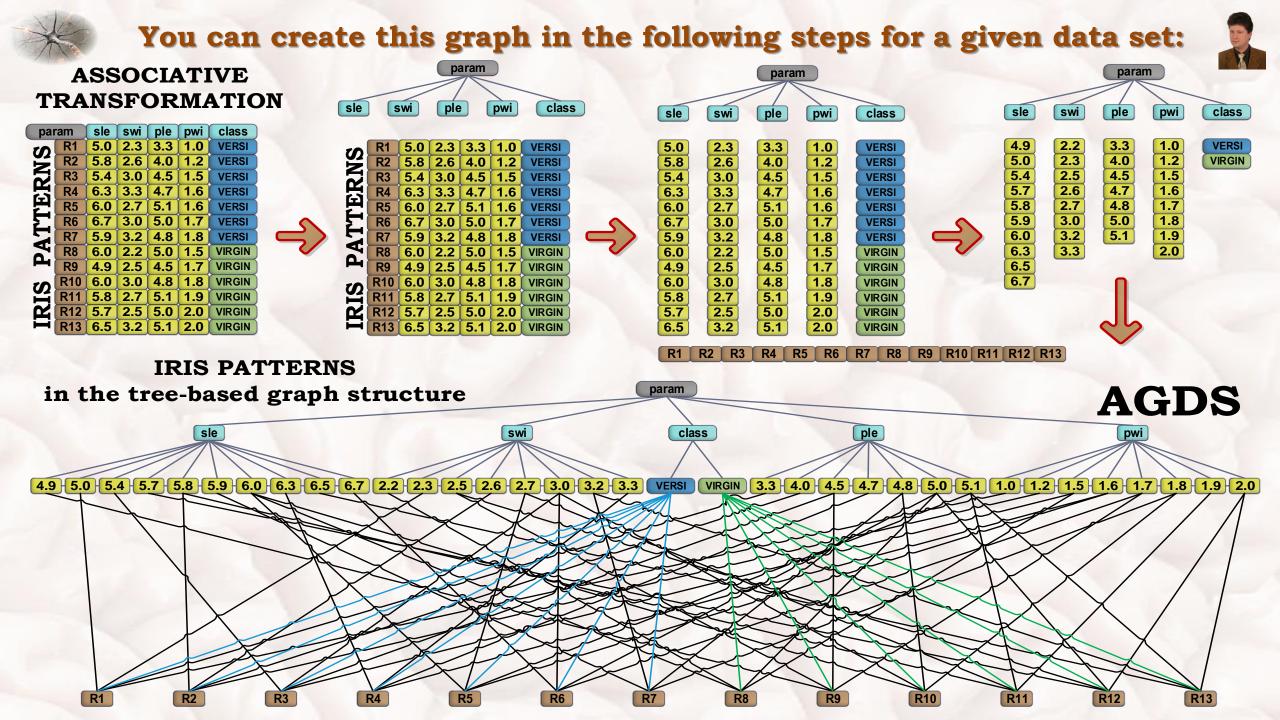




Implement the simplistic version of an AGDS structure or AANG neural graph for Iris data. You don't need to use AVB-trees, AVB+trees, or ASSORT-2 to create it. You can use hash-tables, classic sorting algorithms, sorted lists, dictionaries, but the result should be similar and correct. You don't need to present your results or a network structure graphically, Only text form of results is required. If you like to choose one of the projects implementing AANG (the proposal 1 or 2) you can include this task in your final projects and do both at one go. Independently from your choice of implementation AGDS or AANG use the selected one to:

1. Find the most similar object(s) to the given object.

- 2. Find the most similar object(s) to any combination of input data given on the input.
- 3. Using associations generate the list of all objects sorted by the similarity to the give object (Task 1) or any combination of input data (Task 2).





WEGIHTS COMPUTATION

AGDS nodes representing neighboring (subsequent) values of each attribute a_k are connected and the weight of this connection (edge) is computed after the following formula:

$$w_{v_i^{a_k}, v_j^{a_k}} = 1 - \frac{\left|v_i^{a_k} - v_j^{a_k}\right|}{r^{a_k}}$$

where

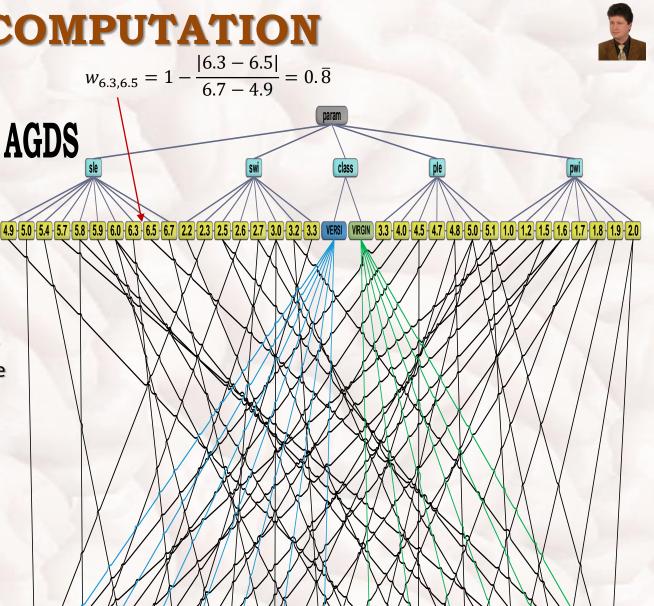
 $v_i^{a_k}$, $v_i^{a_k}$ - are values represented by the neighboring attribute nodes, which are connected by an edge in the AGDS graph,

 $r^{a_k} = v^{a_k}_{max} - v^{a_k}_{min}$ - is the current range of values of the attribute a_k . The weight of the connection from the value node $v_i^{a_k}$ of the attribute a_k to the object node R_m is determined after the number of occurrences $N_i^{a_k}$ of this value $(v_i^{a_k})$ in all objects:

$$W_{v_i^{a_k},R_m} = \frac{1}{N_i^{a_k}} = \frac{1}{\|v_i^{a_k}\|}$$

These numbers $(N_i^{a_k} = ||v_i^{a_k}||)$ are stored in the individual value nodes of each attribute. This number is equal to the number or all connections of this value node to all object nodes if there are no duplicated objects in the table used to create the AGDS structure. In the opposite direction, the weights of connections from the object nodes to the value nodes are always equal to one:

$$W_{R_m,v_i}^{a_k} = 1$$



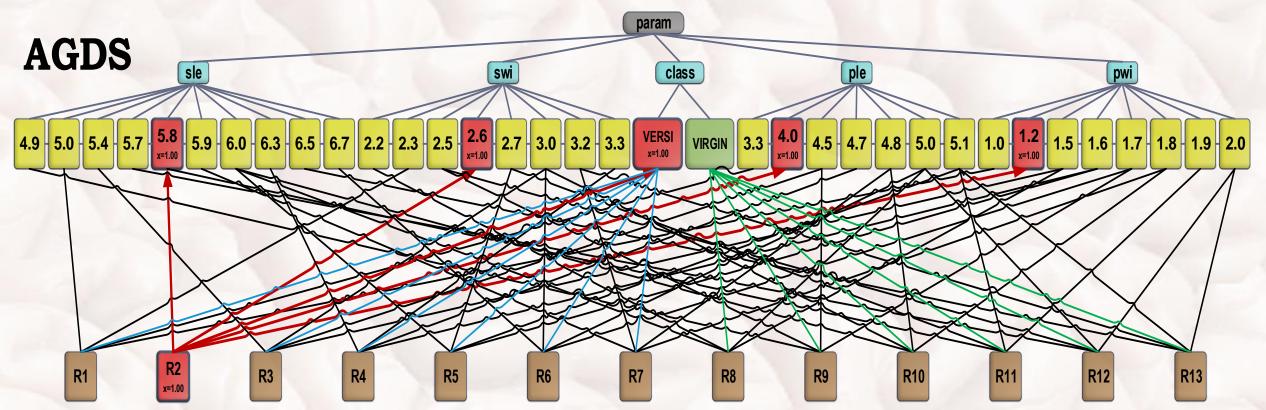


ASSOCIATIVE INFERENCE USING AGDS STRUCTURES



Associative data structures AGDS can be now used for associative inference, which is based on moving along the connections to the connected nodes and computing some values in these nodes on the basis of the send values multiplied by weights of these connections. In such a way we get the information about e.g. similarity of objects represented by other nodes of the same kind or about the objects that satisfy some given conditions defined by the represented attribute values. Let's use our AGDS graph created for 13 Irises for such inference looking for objects (Irises) Rx which are most similar to R2.

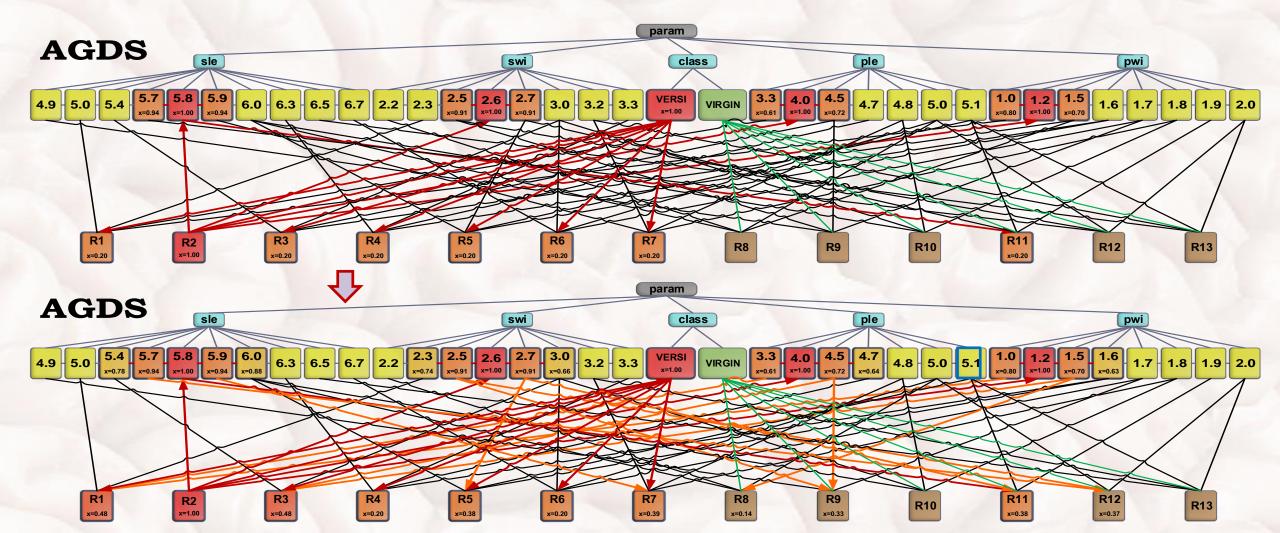
- 1. We start in the node R2 which assumes the similarity value x=1.0, because this node is 100% similar to itself.
- 2. Next, we assign values x of the connected nodes representing the following values: 5.8, 2.6, VERSI, 4.0, and 1.2 by multiplying the value coming from the node R2 with the connection weights, which are equal 1.0. So, as a result, we achieve x=1.0 for all these connected nodes.





ASSOCIATIVE INFERENCE USING AGDS STRUCTURES

- 3. Subsequently, the values computed for these nodes are multiplied by next connection weights and send to the neighbor connected value nodes, for which we also compute their similarity values x.
- 4. Similarly, we compute the similarity values x for connected object nodes with regards to the necessity to add the passed weighted values to the sums already stored in these nodes, e.g. for the node R3 we compute x = 1.0 * 0.2 + 0.72 * 0.2 + 0.7 * 0.2 = 0.48



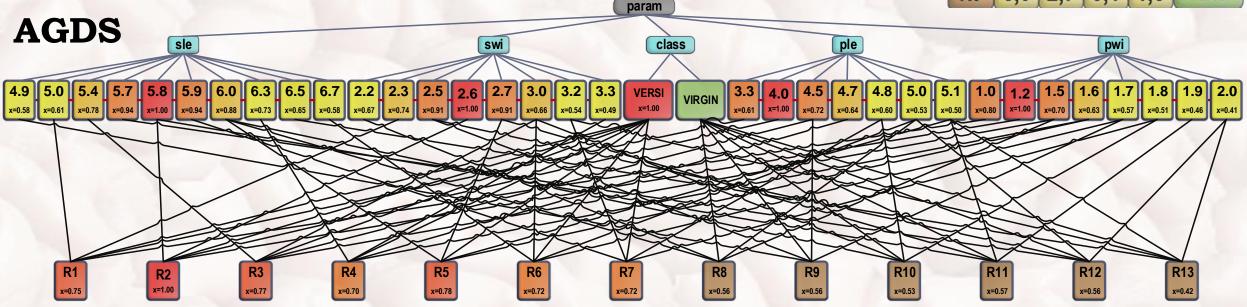


ASSOCIATIVE INFERENCE USING AGDS STRUCTURES



5. Finally, when we go through all the connected (associated) values nodes computing theirs values of similarities by multiplying the sender similarity values by connection weights. We also computed weighted sums for all object nodes, where these weights are here equal w = 1/5 = 0.2. The computed similarity values for the nodes Rx can be used to compare and designate the most similar objects to the object R2: R5 (78%), R3 (77%), R1 (75%), ...



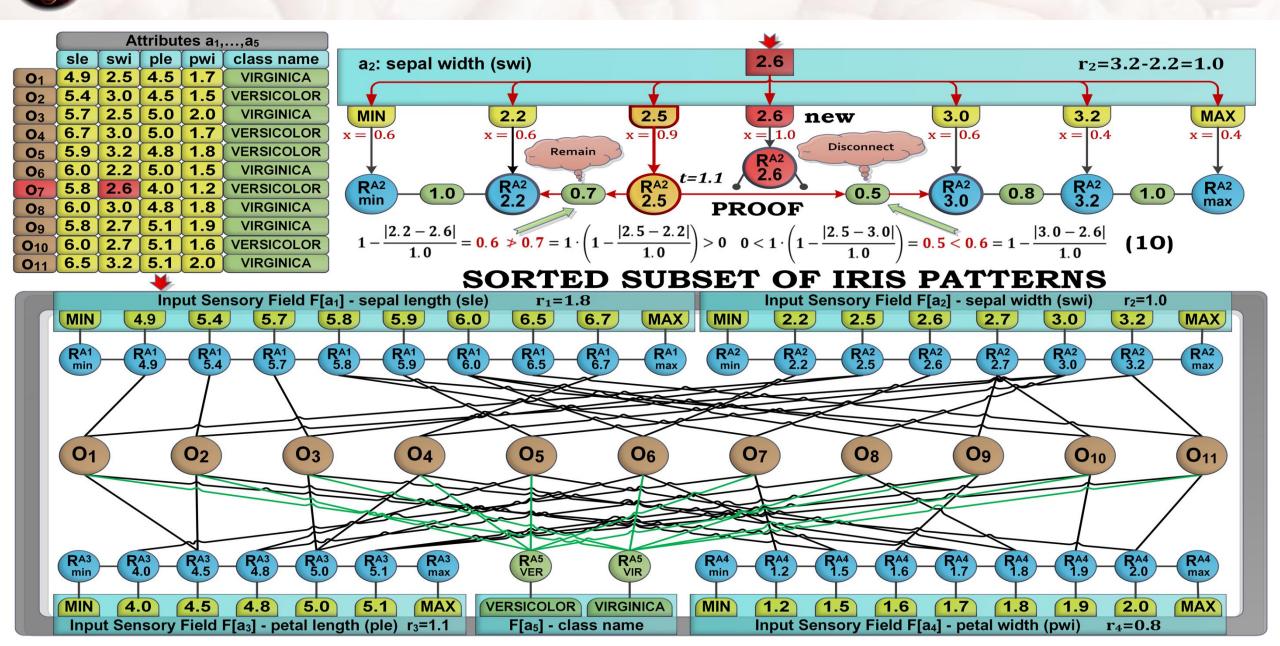


It is also worth noting that AGDS graphs are not neural structures, so we are not obligated to multiply the nodes similarity values by connection weights, but we can also use another formulas, e.g. we can subtract the complement of the connection weight value from the similarity value represented by the sender: x' = x - (1 - w).

Consequently, we get another measure of similarity represented by the value nodes and object nodes.

We can also used DASNG graph formulas to calculate weights between value nodes and object nodes to emphasize rarity of the value using the frequency of connections coming out from value nodes: w = 1 / the number of outgoing connections.

You can also choose to implement AANG neural network





Bibliography and References



http://home.agh.edu.pl/~horzyk/lectures/ahdydci.php



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LECTURES

COMPUTATIONAL INTELLIGENCE

Introduction to Artificial, Cognitive and Computational Intelligence

Introduction to Neural Networks

Models of Neurons

Construction and Learning Strategies

Overview of Neural Network Models

Multilayer Perceptron MLP

Radial Basis Function Networks RBF

Support Vector Machine SVM

Reinforcement Learning

Deep Learning DL

Unsupervised Training and Networks

Self Organizing Maps SOM

Recurrent Neural Networks

Associative and Semantic Memories

Associative Neural Systems

What is this course about? This course is intended to give students a broad overview and deep knowledge about popular solutions and efficient neural network models as well as to learn how to construct and train intelligent learning systems in order to use them in everyday life and work. During the course we will deal with the popular and most efficient models and methods of neural networks, fuzzy systems and other learning systems that enable us to find specific highly generalizing models solving difficult tasks. We will also tackle with

is already under construction and will be available at the spring semester in 2016/2017.

order to use them in everyday life and work. During the course we will deal with the popular and most efficient models and methods of neural networks, fuzzy systems and other learning systems that enable us to find specific highly generalizing models solving difficult tasks. We will also tackle with various CI and AI problems and work with various data and try to model their structures in such a way to optimize operations on them throughout making data available without necessity to search for them. This is a unique feature of associative structures and systems. These models and methods will allow us to form and represent knowledge in a modern and very efficient way which will enable us to mine it and automatically draw conclusions. You will be also able to understand solutions associated with various tasks of motivated learning and cognitive intelligence.

This course will include 28 lectures, 14 laboratory classes and 14 project classes.

