#### Dottorato Internazionale in Fisica Applicata Nonlinear Stochastic Dynamics in Complex Systems

Dipartimento di Fisica e Tecnologie Relative - Università degli Studi di Palermo Mercoledì 14 marzo 2007

### DYNAMICAL CLUSTERING: The role of Synchronization for detecting Community Structures in Complex Networks

#### **Alessandro Pluchino\***

in collaboration with A. Rapisarda\*, V. Latora\*, M. Ivanchenko\*\* and S. Boccaletti\*\*\*



\* Dipartimento di Fisica e Astronomia and INFN sezione di Catania University of Catania, Italy - Group web page: www.ct.infn.it/~cactus

\*\* Moscow University

\*\*\* Nat. Ist. Applied Optycs - Florence

## Outline

The Problem: Finding Community Structures in Complex networks The Approach: Synchronization of Dynamical Oscillators in Weighted Networks

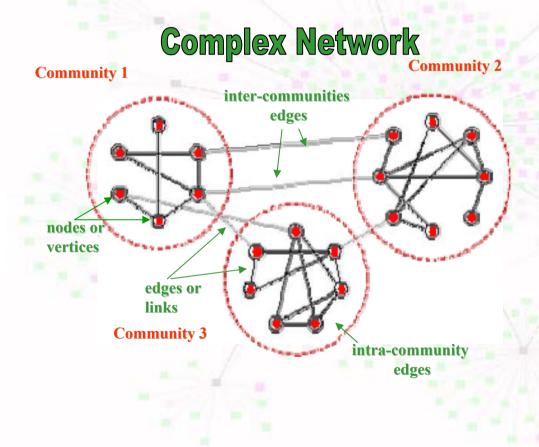
The Solution:

Dynamical Clustering algorithm for the identification of Community Structures in Trial and Real Networks

**Discussion and Numerical Results** 

## **The problem: Finding Community Structures in Complex Networks**

An important open problem in complex networks analysis is the identification of modular structures.



Distinct modular structures, usually called Communities, can loosely be defined as subset of nodes (vertices) which are more densely linked, when compared to the rest of the network.

# **The problem:** Finding Community Structures in Complex Networks

In a limiting case, communities can be also defined as nonconnected clusters of interconnected nodes:

# **Complex Network Community 1 Community 2 Community 3**

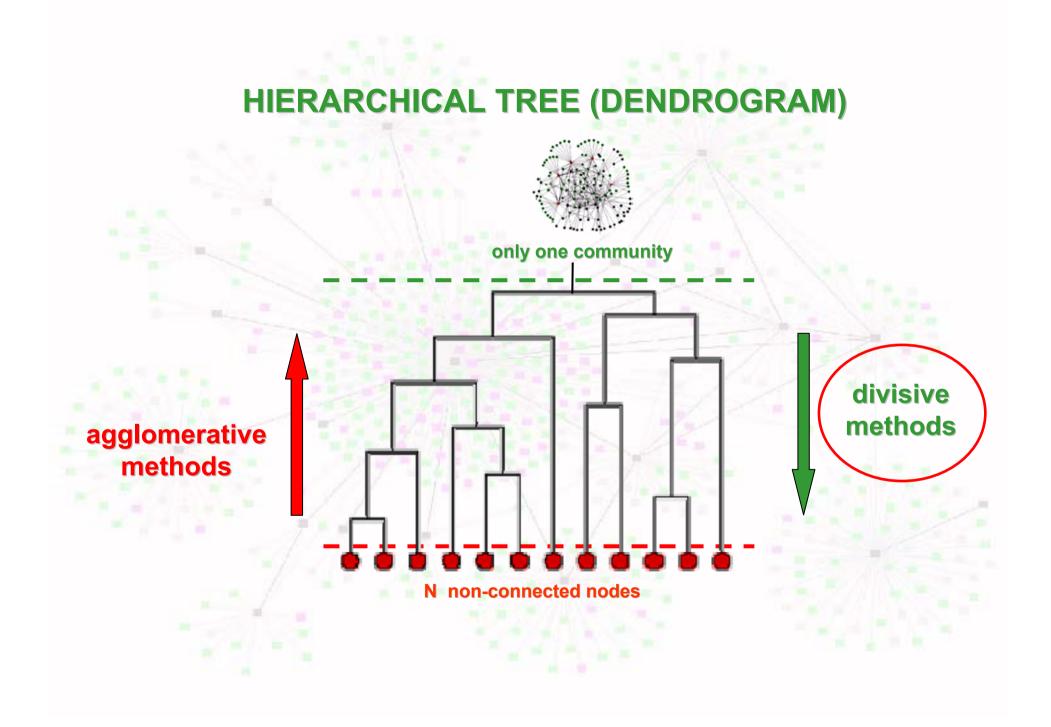
Communities, of course, are fundamental in social networks (parties, cultures, elites), but are also very important in biochemical, metabolic or neural networks (functional groups), in the world wide web (thematic pages), in economic networks, food webs, computer clusters and so on...

A useful set of techniques for the detection of community structures was firstly developed in social network analysis and is known as

#### HIERARCHICAL CLUSTERING METHODS.

These techniques are aimed at discovering natural divisions of networks into groups, based on various metric of similarity or strength of connection between vertices.

They fall into two broad classes: **agglomerative** and **divisive** methods, depending on whether they focus on the addition or the removal of edges to or from the network...

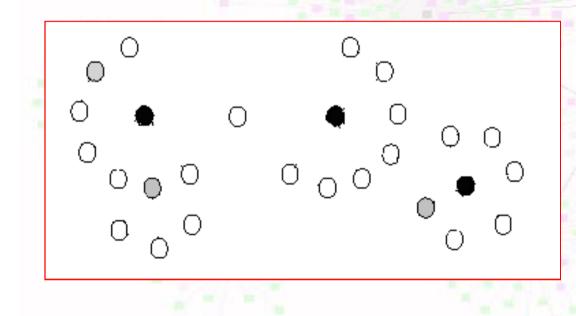


Divisive methods progressively remove the edges of the networks in terms of their importance:

- in propagating some information over the network (*information centrality*) S.Fortunato, V.Latora, M.Marchiori, 2004 *Phys. Rev. E* **70** 056104
- in connecting many pairs of nodes (*shortest-path edge betweenness*, i.e. the number of shortest paths which are making use of a given edge)

M.E.J.Newman and M.Girvan, 2004 Phys. Rev. E 69 026113

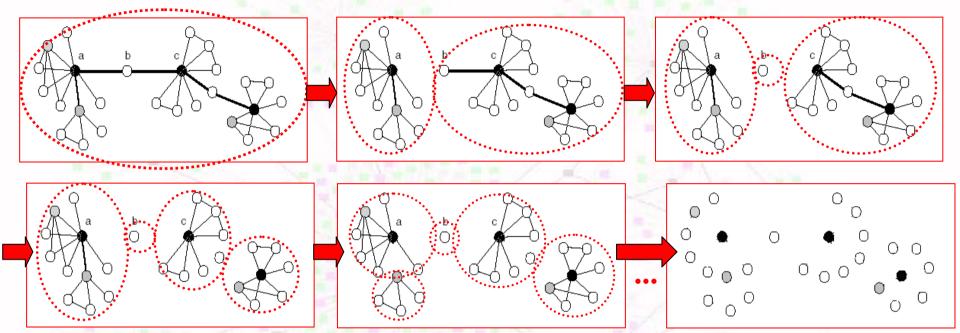
- other quantities...



By doing this repeatedly, and recalculating the betweenness at each step, the network breaks iteratively into smaller and smaller groups of nodes...

...until it breaks into a collection of single non-connected nodes...

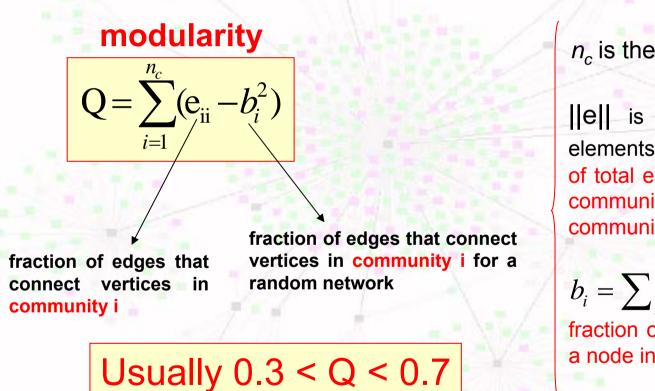
The divisive algorithm produces a hierarchy of subdivisions of the network in isolated groups of interconnected nodes (communities)...



## But which subdivision level does give the best communities configuration for a given network?

Clearly we need some parameter to quantify the reliability of each communities configuration...

This parameter is the "modularity" Q \*, a quantity that, at each step, compares the fraction of intra-community edges with the expected value of the same quantity in an equivalent network with random connections, and allows us to test which communities configuration found by the divisive algorithm is the best one:



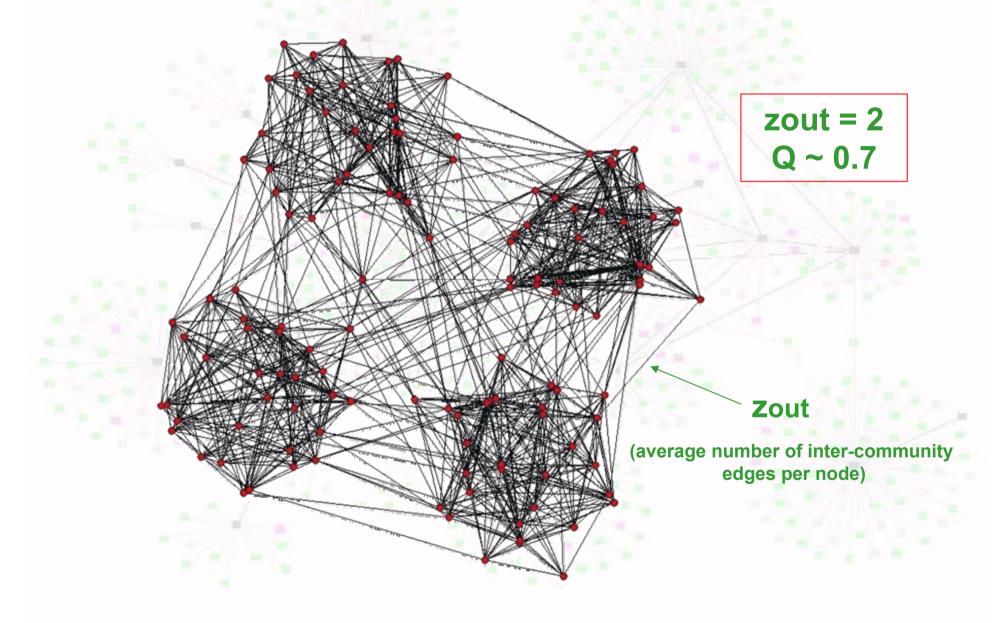
 $n_c$  is the number of communities

 $||\mathbf{e}||$  is a  $n_c \times n_c$  matrix whose elements  $\mathbf{e}_{ij}$  represent the fraction of total edges connecting a node in community-i with a node in community-j

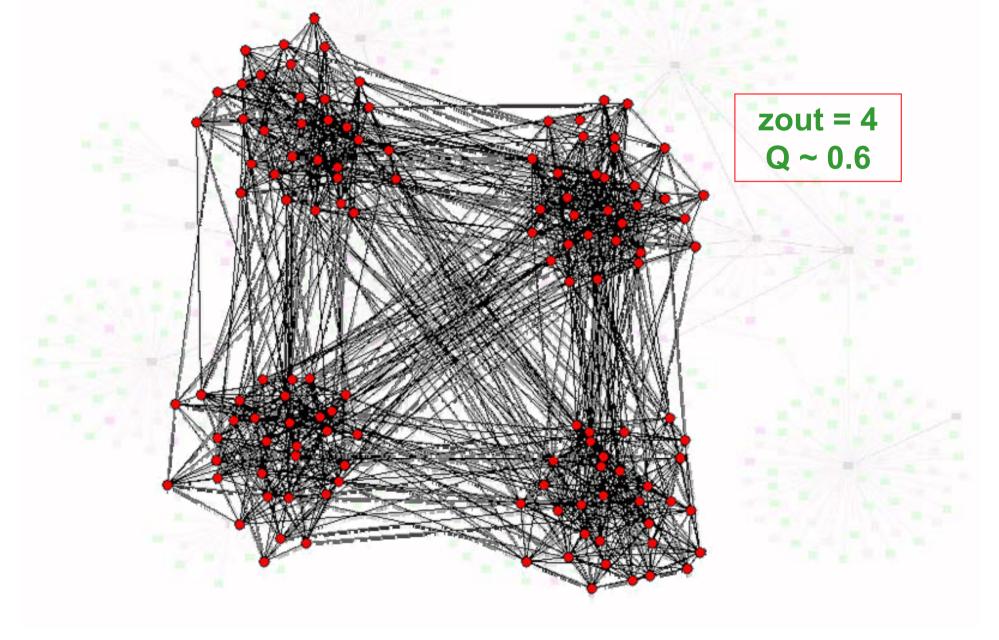
 $b_i = \sum_j e_{ij}$  represents the fraction of total edges connected to a node in community-i

\*M.E.J.Newman and M.Girvan, 2004 Phys. Rev. E 69 026113

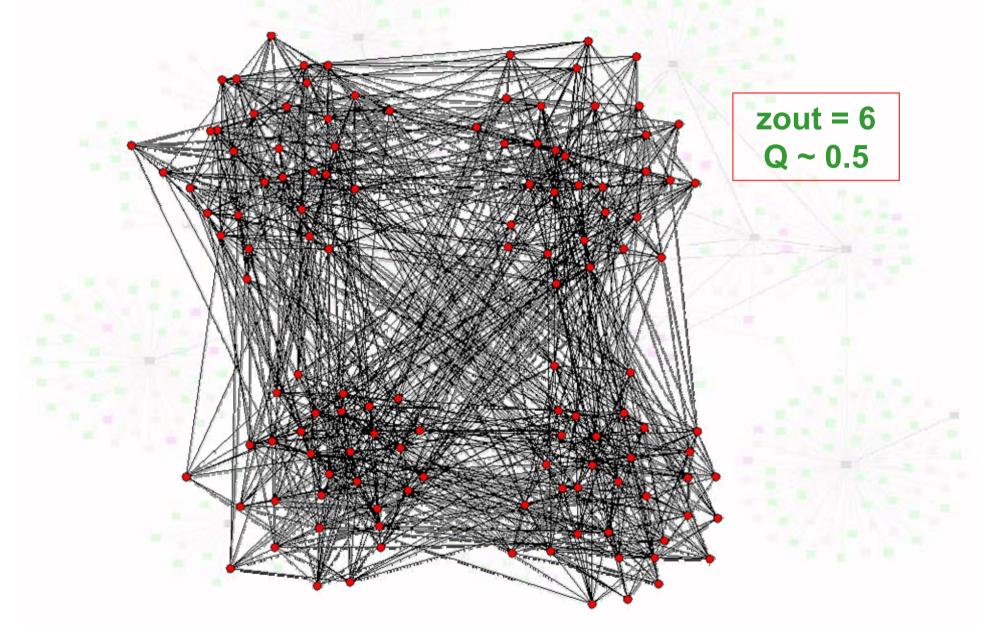
#### Modularity in "ad hoc" random trial networks (N=128, <k>=16, 4 communities)

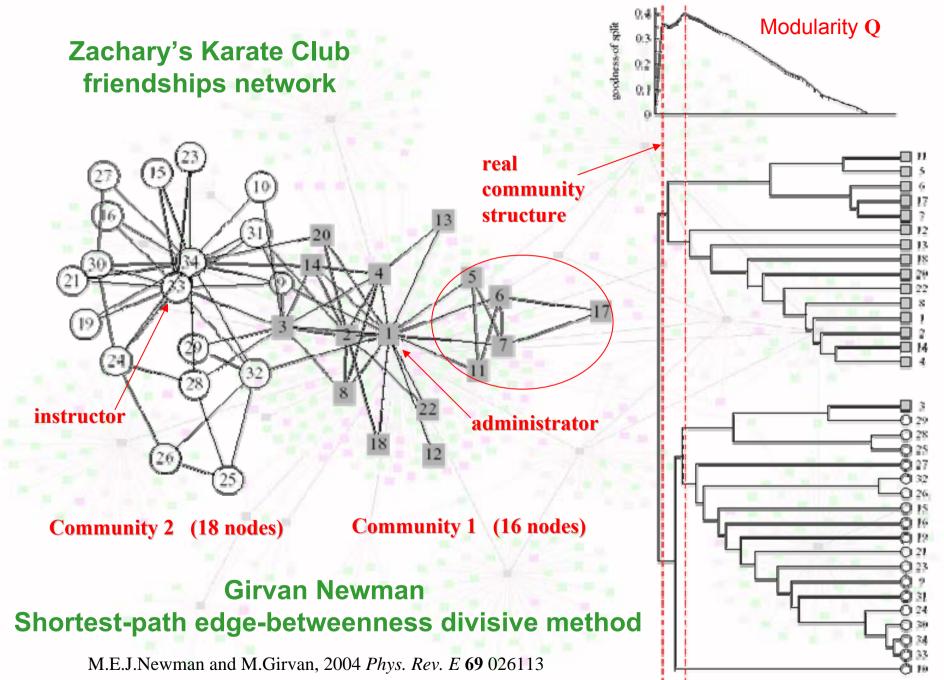


#### Modularity in "ad hoc" random trial networks (N=128, <k>=16, 4 communities)



#### Modularity in "ad hoc" random trial networks (N=128, <k>=16, 4 communities)





W.Zachary (1977) J.Anthropol.Res. 33 452-473

Topological Divisive Algorithms like GN have the problem of recalculating betweennesses at each step.

Since a single-step calculus of all the edge-betweennesses takes O(N<sup>2</sup>) operations, and the whole process takes N steps, these algorithms are quite slow – O(N<sup>3</sup>) –

#### DIFFERENT APPROACH: Synchronization of Dynamical Oscillators in Weighted Networks

## **The Kuramoto model\***

The simplest models for synchronization available on the market is the celebrate Kuramoto model and consists of N fully connected phase oscillators with natural frequencies  $\omega_i$  and coupling parameter K:

 $\begin{aligned}
\underbrace{\frac{d\vartheta_i(t)}{dt} = \omega_i + \frac{K}{N} \sum_{j=1}^{N} \sin(\vartheta_j - \vartheta_i), \quad i = 1, \dots, N}_{\text{natural frequencies}} \\
\theta_i(t) \in \begin{bmatrix} 0, 2\pi \end{bmatrix} \xrightarrow{\text{natural frequencies}} \text{phases of oscillators} \\
The coherence of the system is measured by the mean field order parameter r (<math>0 \le r(t) \le 1$ ):  $re^{i\psi} = \frac{1}{N} \sum_{j=1}^{N} e^{i\vartheta_j}$ Through r the rate equations can be rewritten as:

$$\frac{\vartheta_i(t)}{dt} = \omega_i + Kr\sin(\psi - \vartheta_i), \quad i = 1, \dots N$$

\*proposed by Y.Kuramoto in 1975

## The Kuramoto model (2)

As Kuramoto showed analitically in a beautiful analysis, one observes synchronization above a certain *critical* value of the control parameter  $K_c \dots$ 

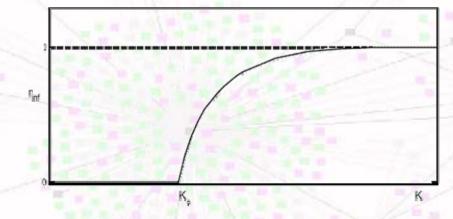
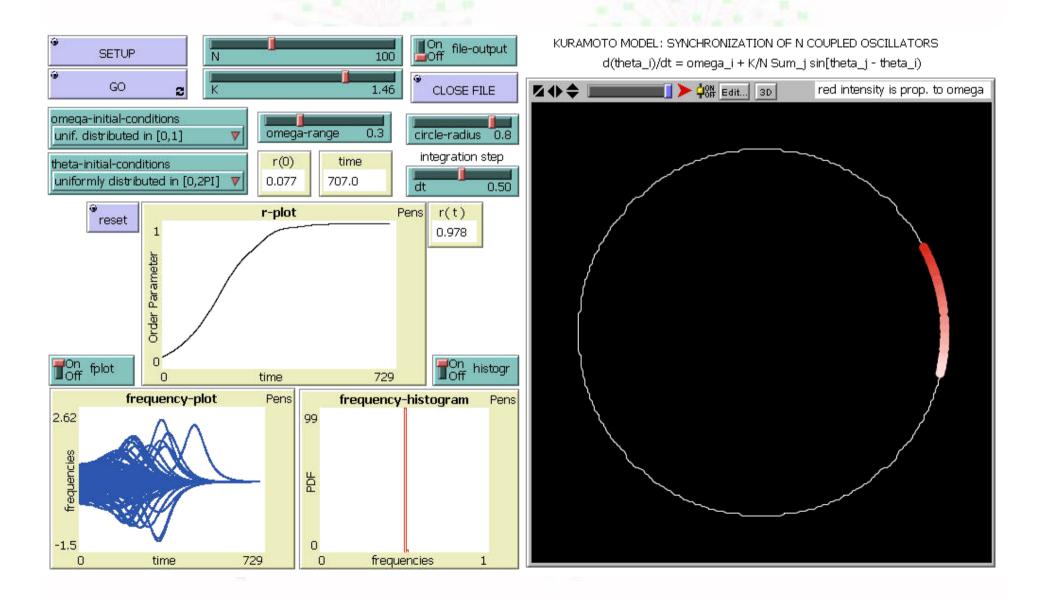


Fig. 1. Asymptotic order parameter  $r_\infty$  as a function of the coupling in the Kuramoto model

 $K \to 0$  $\vartheta_i(t) \approx \omega_i t + \vartheta_i(0)$  $r \to 0$ Incoherent phase $K \to \infty$  $\vartheta_i(t) \approx \psi(t)$  $r \to 1$ Global synchronization

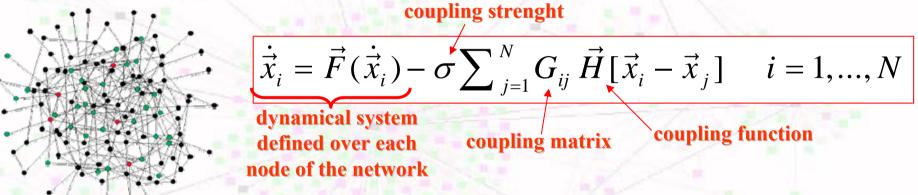
## The Kuramoto model (3)



#### THE MASTER STABILITY FUNCTION APPROACH TO ENHANCE SYNCHRONIZATION IN COMPLEX NETWORKS

Suppose to have a (unweighted, undirected) network of N linearly coupled identical oscillators\*. The equation of motion reads:

**Network with N nodes** 



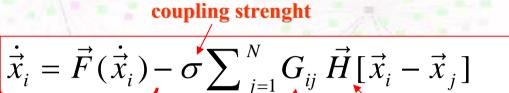
If G has a real spectrum of eigenvalues  $\lambda_i$  (for symmetric coupling) and if we associate  $\lambda_1$  to the state  $\mathbf{x}_s(t)$ , the stability of the synchronous manifold  $(\mathbf{x}_i(t)=\mathbf{x}_s(t), \forall i)$  requires that all the conditional Lyapunov exponents  $\Lambda$ associated with  $\lambda_2 \leq \ldots \leq \lambda_i \leq \ldots \leq \lambda_N$  would be negative...

\*M.Chavez, D.U.Hwang, A.Amann, H.G.E.Hentschel and S.Boccaletti, Phys. Rev. Lett. 94 218701 (2005)

#### THE MASTER STABILITY FUNCTION APPROACH TO ENHANCE SYNCHRONIZATION IN COMPLEX NETWORKS (2)

**Network with N nodes** 





$$i = 1, ..., N$$

dynamical system defined over each node of the network

coupling matrix `co

**`coupling function** 

Defining the Master Stability Function (MSF) as the largest Lyapunov exponent  $\Lambda_{max}$  versus a parameter  $\nu = \sigma \lambda$ , it can be shown\* that, for a large class of dynamical systems, the MSF is negative in a finite parameter interval ( $\nu_1$ ,  $\nu_2$ ).

Thus the condition for synchronization stability is governed by the ratio  $\lambda_N/\lambda_2$ : the more packed the eigenvalues of G are, the higher is the chance of having all Lyapunov exponents into the stability range for some  $\sigma$ .

\*M.Chavez, D.U.Hwang, A.Amann, H.G.E.Hentschel and S.Boccaletti, Phys. Rev. Lett. 94 218701 (2005)

One can use the master stability function approach:

- 1) to find the best synchronization condition of a given network\*
- 2) to tune the synchronization of a network in order to identify community structures\*\* (DYNAMICAL CLUSTERING)

Both the results can be realized with an opportune choice of the coupling matrix  $G_{ij}$  in the network equation, by means of a weighting procedure that assignes to each edge a 'load'  $I_{ij}$  equal to its betweenness (i.e. the number of shortest paths that are making use of that edge):

 $coupling matrix G=G(\alpha)$ 

$$\dot{\vec{x}}_{i} = \vec{F}(\dot{\vec{x}}_{i}) - \sigma \sum_{j \in K_{i}} \left( \frac{l_{ij}}{\sum_{j \in K_{i}} l_{ij}} \vec{H}[\vec{x}_{i} - \vec{x}_{j}] \right) \qquad i = 1, ..., N$$

where  $\alpha$  is a real tunable parameter and K<sub>i</sub> is the set of neighbors of the i<sup>th</sup> node.

#### master stability function arguments apply

\*M.Chavez, D.U.Hwang, A.Amann, H.G.E.Hentschel and S.Boccaletti, *Phys. Rev. Lett.* **94** 218701 (2005) \*\* A.Pluchino., M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti, *in preparation* 

#### Most common dynamical systems defined over the network

$$\dot{\vec{x}}_{i} = \vec{F}(\dot{\vec{x}}_{i}) + \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha}} \sum_{j \in K_{i}} l_{ij}^{\alpha} \vec{H}[\vec{x}_{i} - \vec{x}_{j}] \qquad i = 1, ..., N$$

**Chaotic Rössler** identical 3D oscillators

$$\begin{cases} \dot{x}_{i} = -\omega y_{i} - z_{i} - \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha(t)}} \sum_{j \in K_{i}} l_{ij}^{\alpha(t)} (x_{i} - x_{j}) \\ \dot{y}_{i} = \omega x_{i} + 0.165 y_{i} \end{cases}$$

$$i = 1, ..., N$$

 $\dot{z}_i = 0.2 + z_i (x_i - 10)$ 

Kuramoto's non identical 1D oscillators

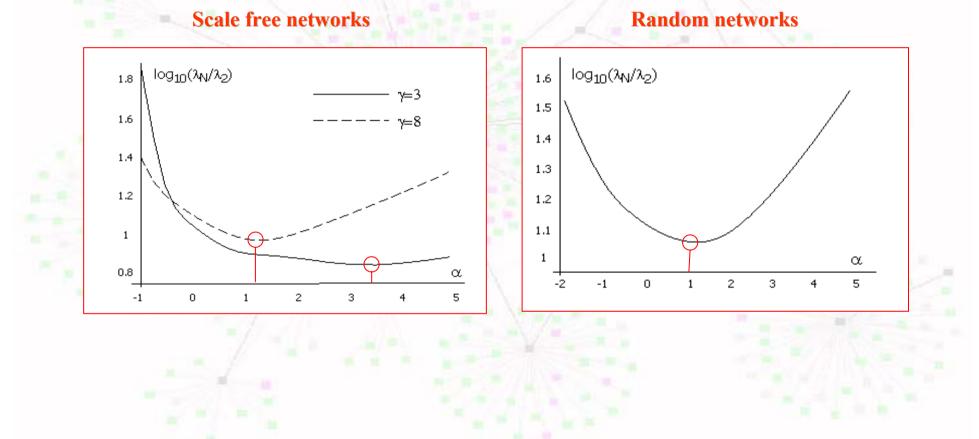
$$\dot{\vartheta}_{i} = \omega_{i} + \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha(t)}} \sum_{j \in K_{i}} l_{ij}^{\alpha(t)} \sin(\theta_{j} - \theta_{i}) \qquad i = 1, \dots, N$$

#### Sine-Circle Map: non identical 1D oscillators

$$x_{i}(n+1) = x_{i}(n) + \omega_{i} + \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha(t)}} \sum_{j \in K_{i}} l_{ij}^{\alpha(t)} \sin(x_{j} - x_{i}) \qquad i = 1, ..., N$$

## 1. Finding the best synchronization condition for a network of Rössler oscillators

$$\dot{\vec{x}}_{i} = \vec{F}(\dot{\vec{x}}_{i}) - \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha}} \sum_{j \in K_{i}} l_{ij}^{\alpha} \vec{H}[\vec{x}_{i} - \vec{x}_{j}] \qquad i = 1, ..., N$$



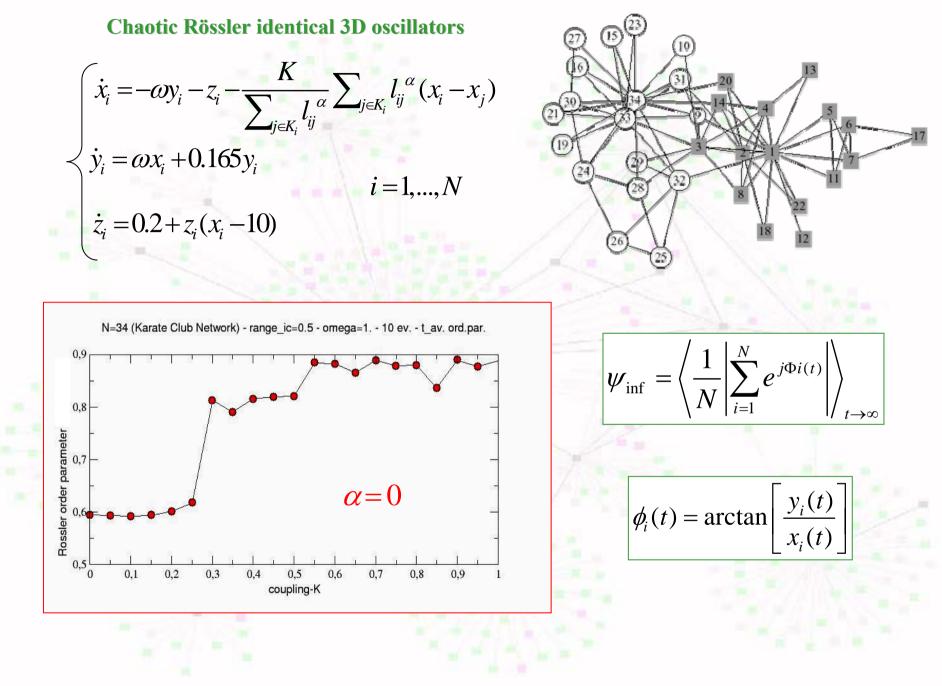
M.Chavez, D.U.Hwang, A.Amann, H.G.E.Hentschel and S.Boccaletti, Phys. Rev. Lett. 94 218701 (2005)

## 2. Tuning the synchronization of a network of oscillators for finding community structures

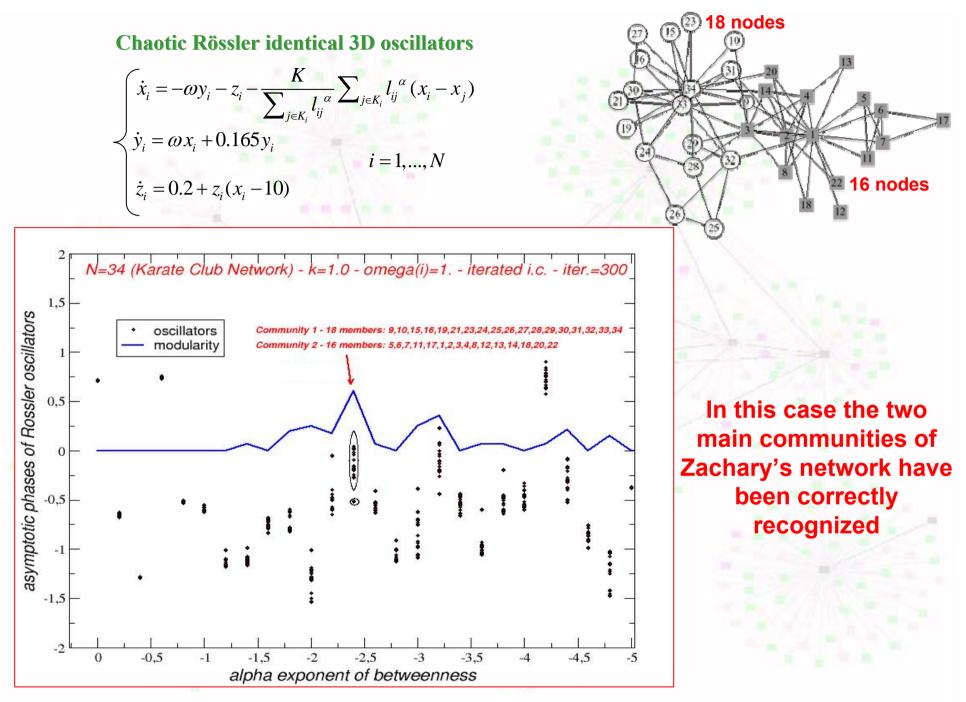
#### **DYNAMICAL CLUSTERING ALGORITHMS**

$$\dot{\vec{x}}_{i} = \vec{F}(\dot{\vec{x}}_{i}) - \frac{\sigma}{\sum_{j \in K_{i}} l_{ij}^{\alpha(t)}} \sum_{j \in K_{i}} l_{ij}^{\alpha(t)} \vec{H}[\vec{x}_{i} - \vec{x}_{j}] \qquad i = 1, \dots, N$$

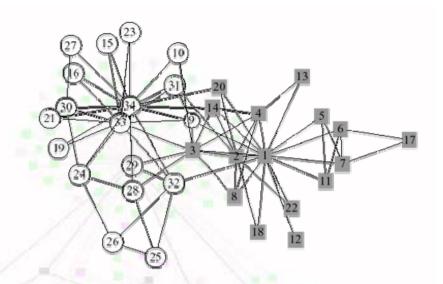
- 1. At variance with the topological methods we calculate the edge betweennesses (i.e. the edge's loads  $l_{ii}$ ) of the network only one time before starting the simulation;
- 2.  $t=0: \alpha(0) \sim 0$  The system starts from a state with perfectly synchronized frequencies for a given value of the coupling strenght.;
- 3. t > 0:  $\alpha(t) \rightarrow -\infty$  Decreasing  $\alpha$ , the edges with the greatest betweenness are weighted less and less and the oscillators progressively desynchronize;
- 4. We look to clusters of nodes oscillating with a common frequency (communities) and we select the clusters configuration with the highest modularity Q.



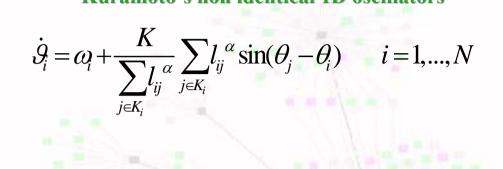
A.Pluchino, V.Latora, A.Rapisarda and S.Boccaletti, in preparation

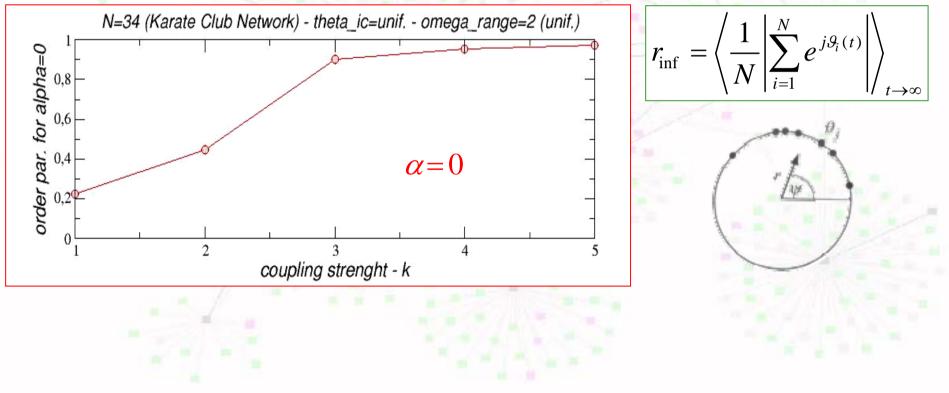


A.Pluchino, V.Latora, A.Rapisarda and S.Boccaletti, in preparation

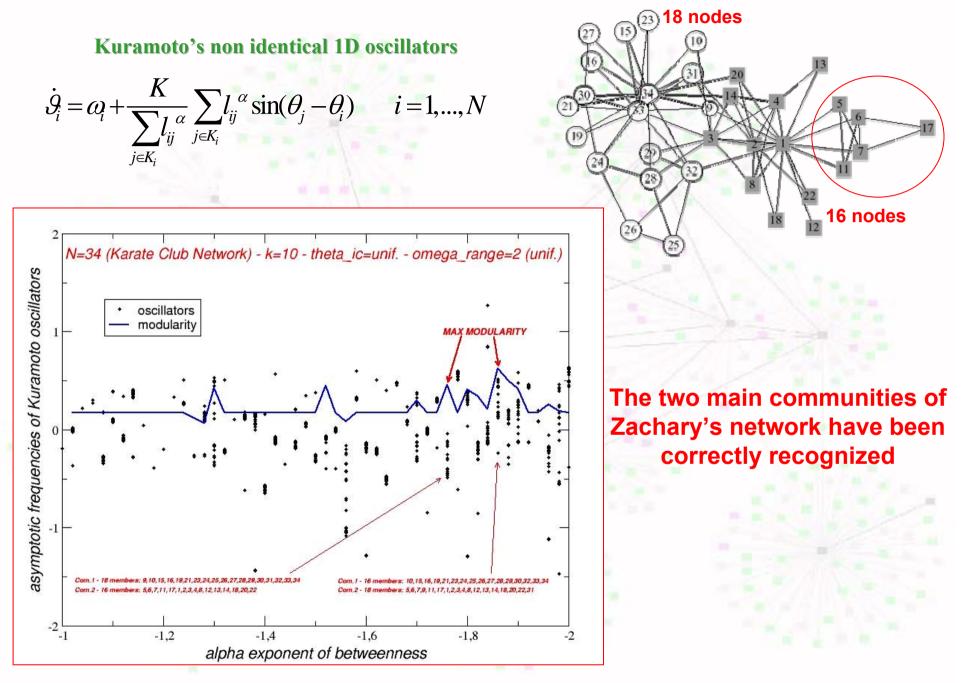


Kuramoto's non identical 1D oscillators





A.Pluchino, V.Latora, A.Rapisarda and S.Boccaletti, in preparation

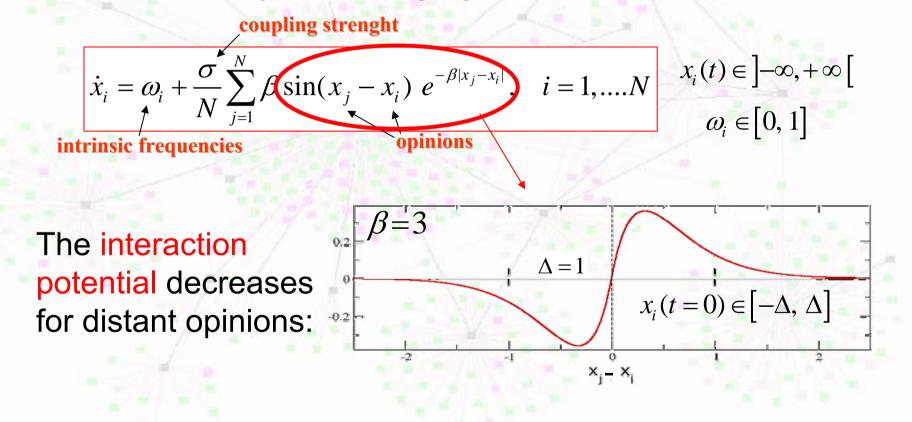


A.Pluchino, V.Latora, A.Rapisarda and S.Boccaletti, in preparation

But the results of sensitivity tests on trial network with a known community structure were not so excellent.....

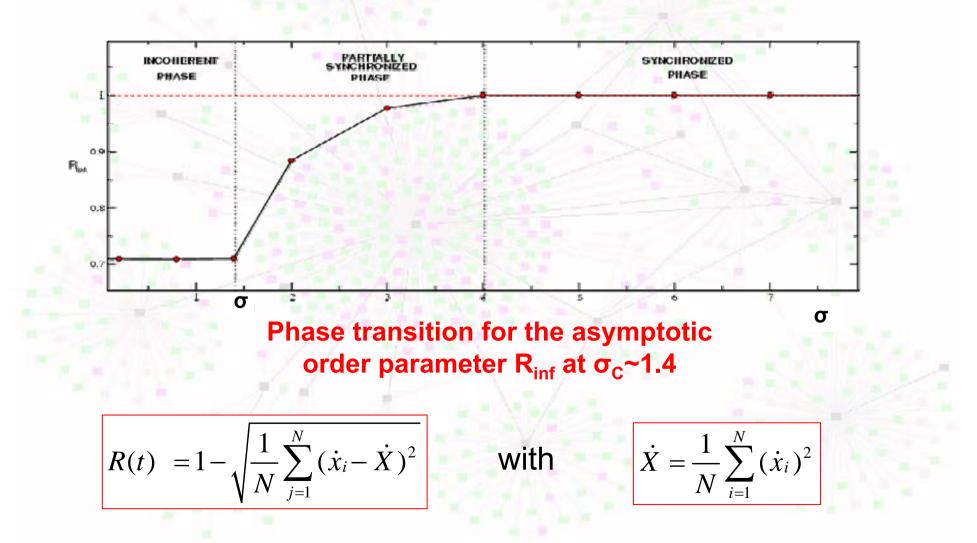
#### The Opinion Changing Rate (OCR) model\*

It is a modification of the Kuramoto model and consists of the following rate equations describing the opinions evolution of N fully interacting agents:



\*A.Pluchino, V.Latora, A.Rapisarda, Int.Journ.of Mod.Phys. C 16 515 (2005)

#### The Opinion Changing Rate (OCR) model\*

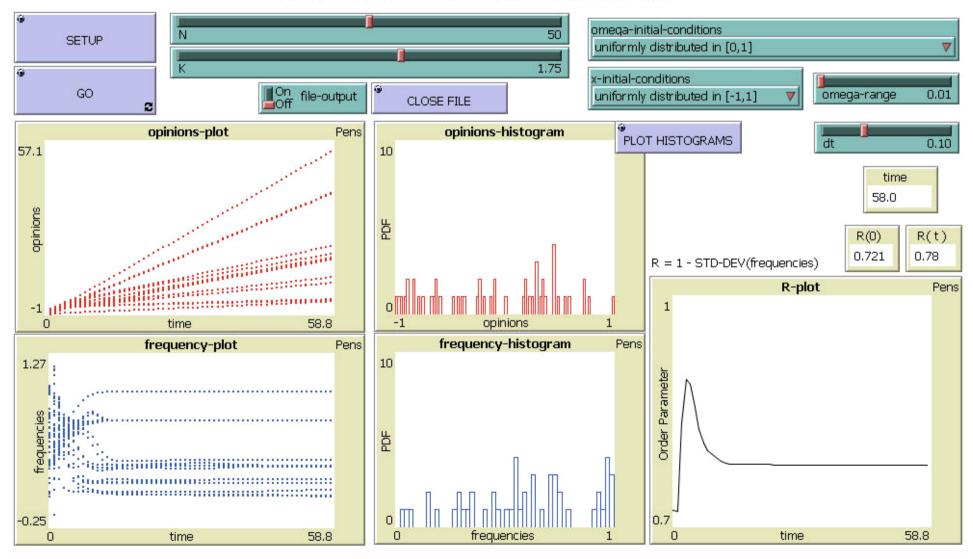


\*A.Pluchino, V.Latora, A.Rapisarda, Int.Journ.of Mod.Phys. C 16 515 (2005)

#### The Opinion Changing Rate (OCR) model\*

OCR MODEL: SYNCHRONIZATION OF CHANGING OPINIONS

d(x\_i)/dt = omega\_i + K/N Sum\_j (alfa sin[x\_j - x\_i) exp(-alfa|xj - xi|))



Here we further modified the standard OCR model forcing the oscillators natural frequencies to follow the so called Heigselmann-Krause dynamics...

$$\dot{x}_{i} = \omega_{i} + \frac{\sigma}{N} \sum_{j=1}^{N} \beta \sin(x_{j} - x_{i}) e^{-\beta |x_{j} - x_{i}|}, \quad i = 1, ....N$$

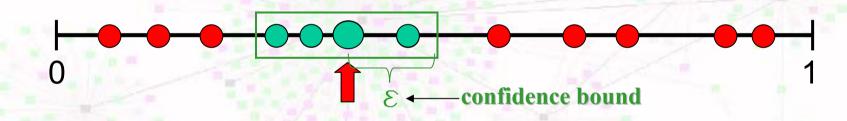
intrinsic frequencies

 $\omega_i = \omega_i(t)$  time dependent

#### **Heigselmann-Krause Dynamics**

The Hegselmann-Krause (HK) opinion dynamics<sup>\*</sup> is based on the presence of a parameter  $\varepsilon$ , called "confidence bound", which expresses the range of compatibility of opinions of agents put on a network (real space).

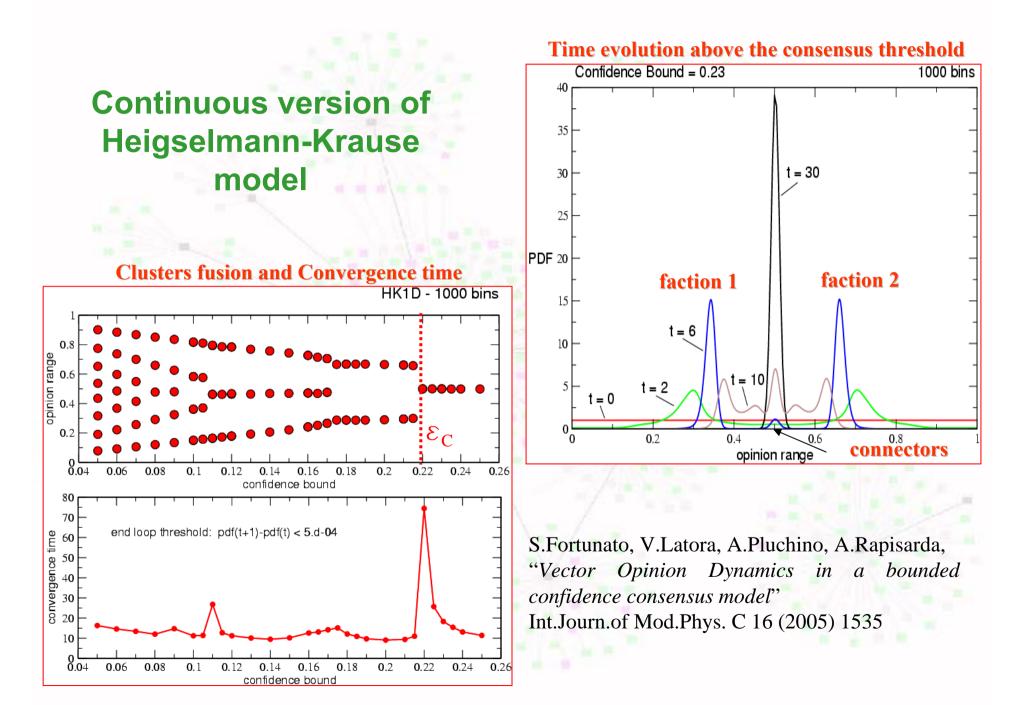
The 1-D opinion space is represented by the points of a [0,1] line, where the opinions are randomly distributed:



At each step, one chooses at random one opinion and checks how many opinions, belonging to first neighbours agents on the network, are compatible with him, i.e. are inside the confidence bound...

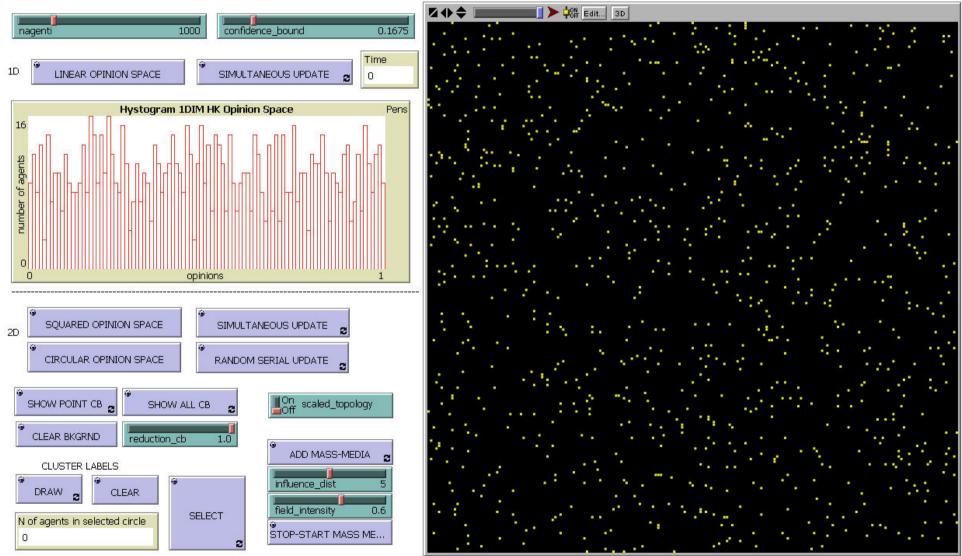
...at the next step, the agent takes the **average opinion** of its compatible neighbours...

\*R. Hegselmann and U. Krause, *Journal of Articial Societies and Social Simulation 5*, issue 3, paper 2 (jasss.soc.surrey.ac.uk) (2002);



#### **Heigselmann-Krause Dynamics**

HEGSELMANN-KRAUSE MODEL



## OCR-HK Dynamical Clustering Algorithm

istantaneous frequencies (opinion changing rates)

ben changing rates)  $\begin{aligned}
& \text{loads (betweennesses)} \\
& \text{tuning parameter} \\$ 

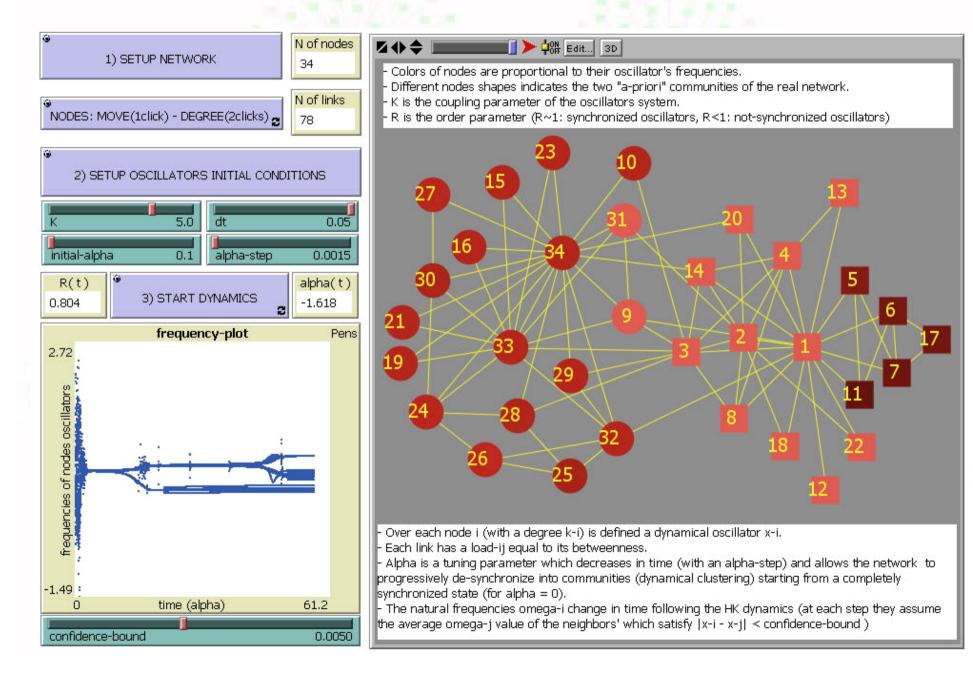
intrinsic frequencies, no more constant but updated with HK dynamics

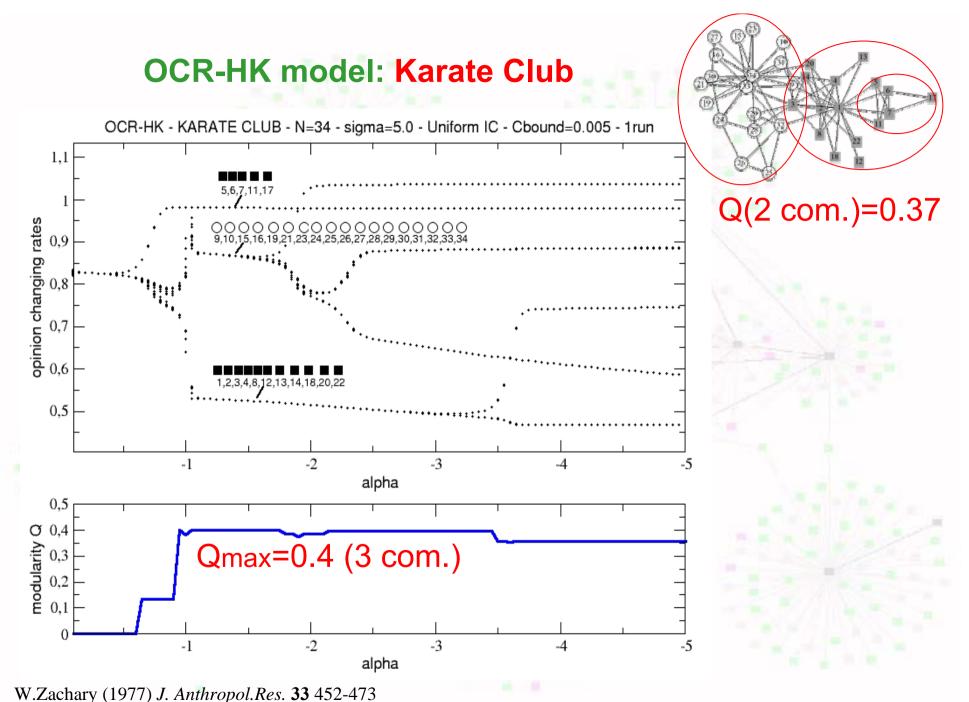
Starting from  $\alpha=0$  (synchronized state) we let  $\alpha$  to decrease in time and we look at the evolution of clusters in frequency during a single run (frequency desynchronization).

We repeat the procedure for several runs, with different initial conditions, then we select the configuration with the highest modularity Q

A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - physics/0607179

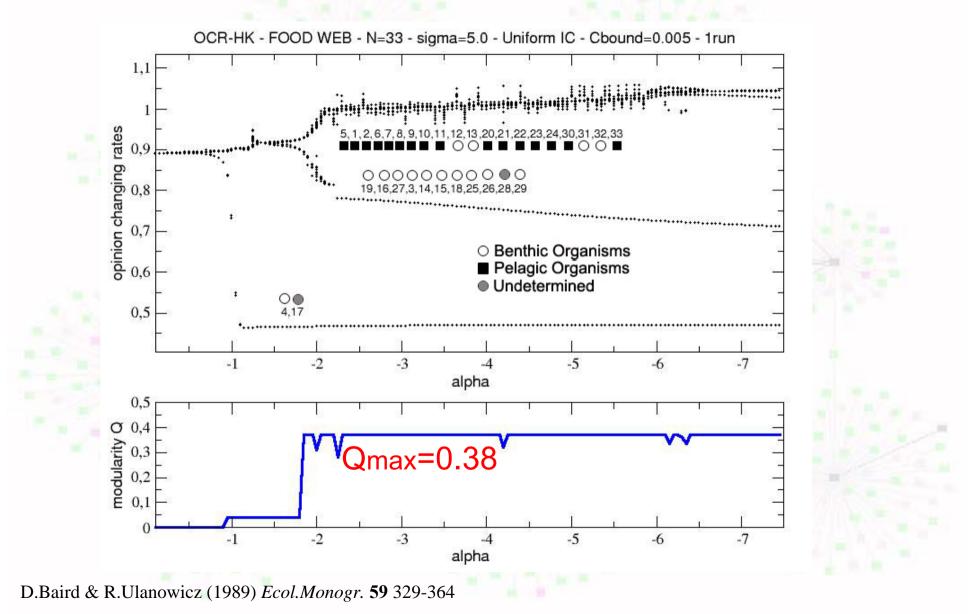
#### **OCR-HK model: Karate Club**





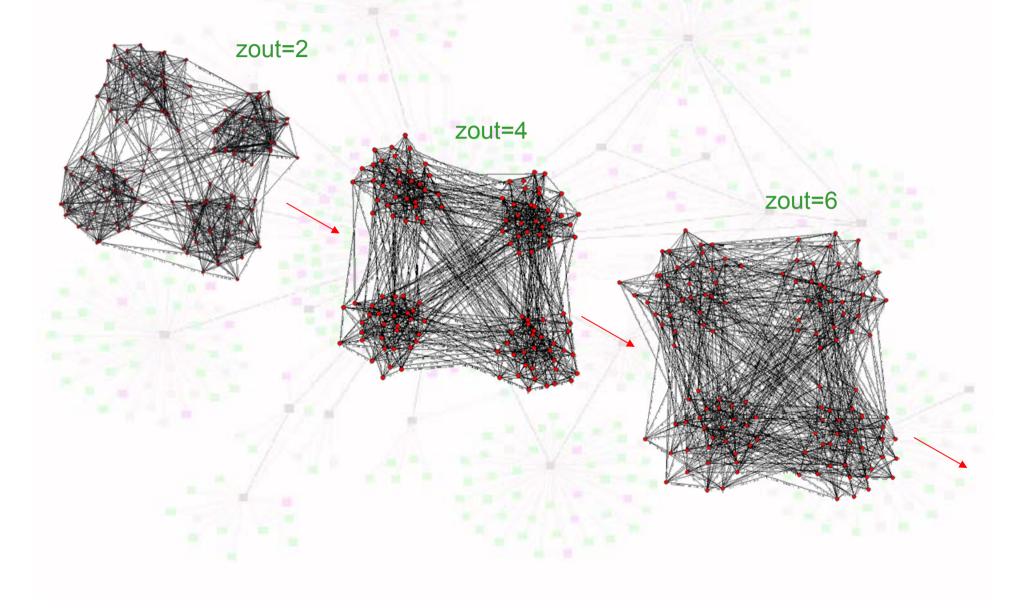
A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - *physics/0607179* 

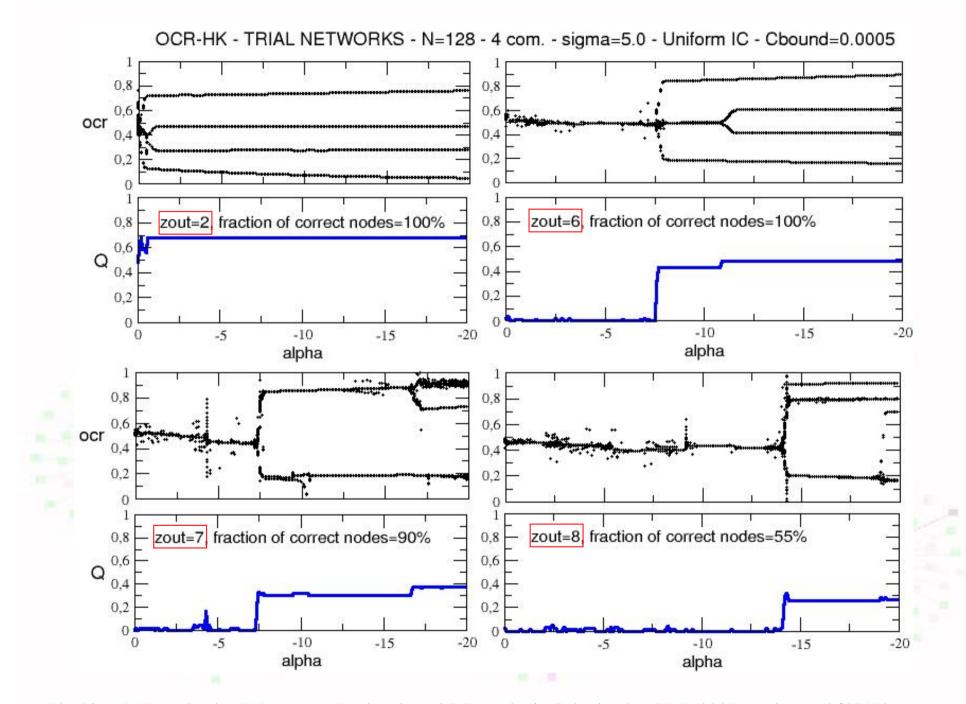
#### **OCR-HK model: Chesapeake Bay food web (USA)**



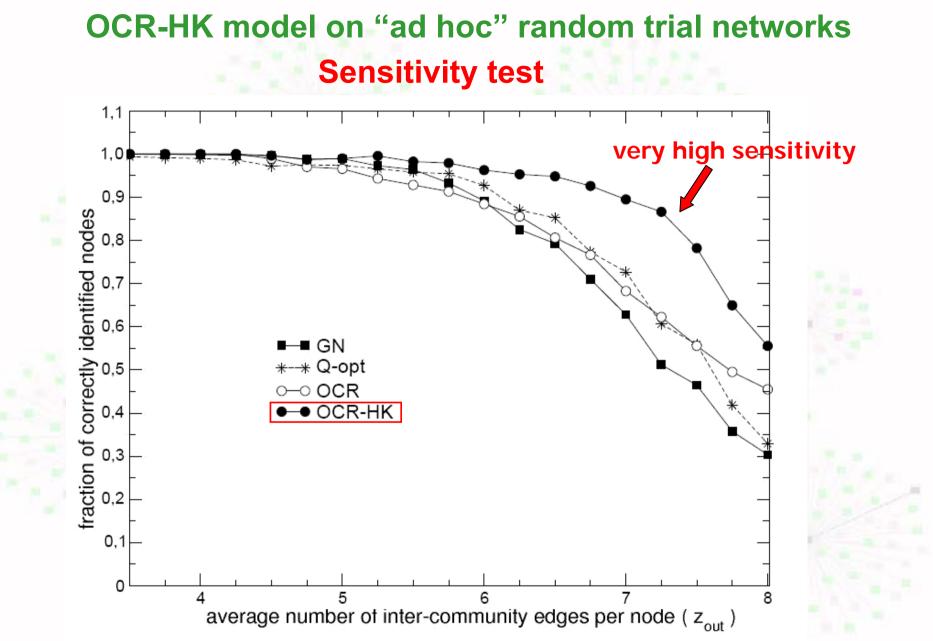
A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - physics/0607179

#### OCR-HK model: "ad hoc" random trial networks (N=128, <k>=16, 4 communities) with increasing zout





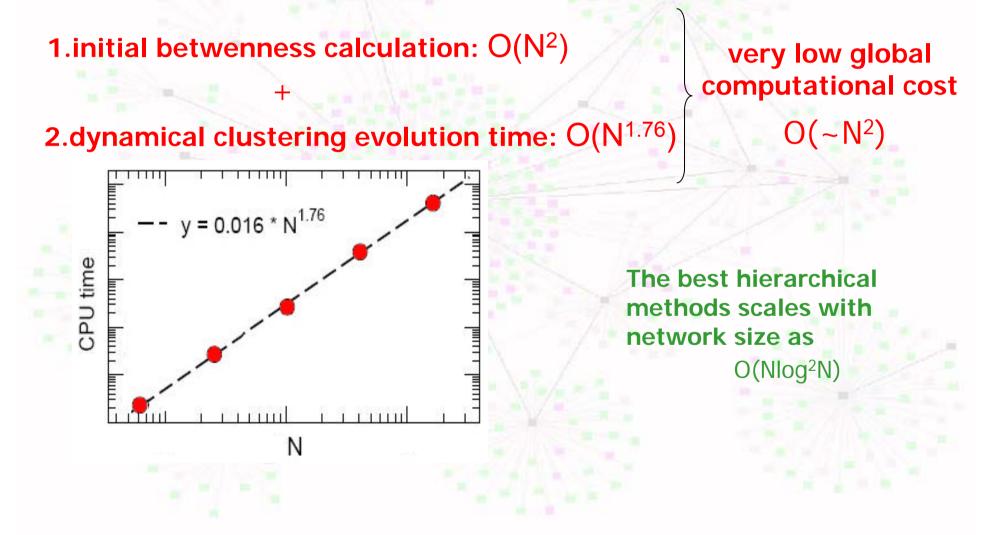
A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - physics/0607179



L.Danon, A.Diaz-Guilera, J.Duch and A.Arenas J.of Stat.Mech.: Theory and Exp. (2005)

A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - physics/0607179

#### OCR-HK model on "ad hoc" random trial networks Computational cost



A.Pluchino, M.Ivanchenko, V.Latora, A.Rapisarda and S.Boccaletti - Submitted to PRE (2007) - physics/0607179

#### Conclusions

• The problem of finding the best community structure subdivision of a network is very important

• Divisive topological methods have a good sensitivity but have also an high computational cost

• We developed a new algorithm based on a dynamical clustering tecnique that shows a very high sensitivity and at the same time is very fast

 It makes also an interesting bridge between researches in complex network and those in synchronization of dynamical systems

• Future investigations regard the application of our algorithm to large real networks (genetic networks, social networks, etc...) – work in preparation...

## **Thanks for the attention!**



http://www.ct.infn.it/~cactus