

Data visualization

Networks

David Hoksza

siret.ms.mff.cuni.cz/hoksza

bioinformatika.mff.cuni.cz/cusbg

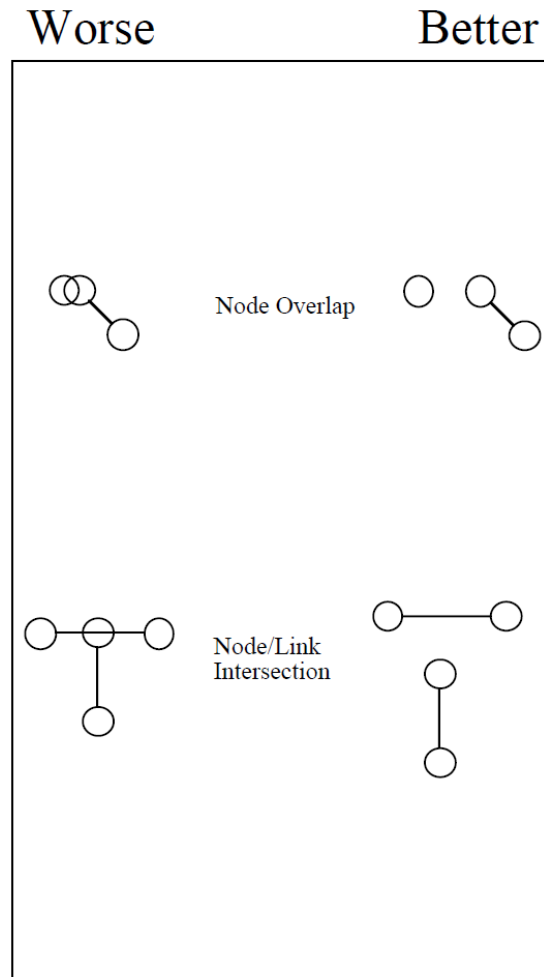
Application areas

- Physical networks
 - Power grid, train routes
- Community modelling
 - Users interaction
- Information (knowledge) networks
 - Citations, web pages, peer-to-peer networks, preference networks
- Biology
 - Metabolic pathways, genetic regulatory networks, protein-protein interaction networks

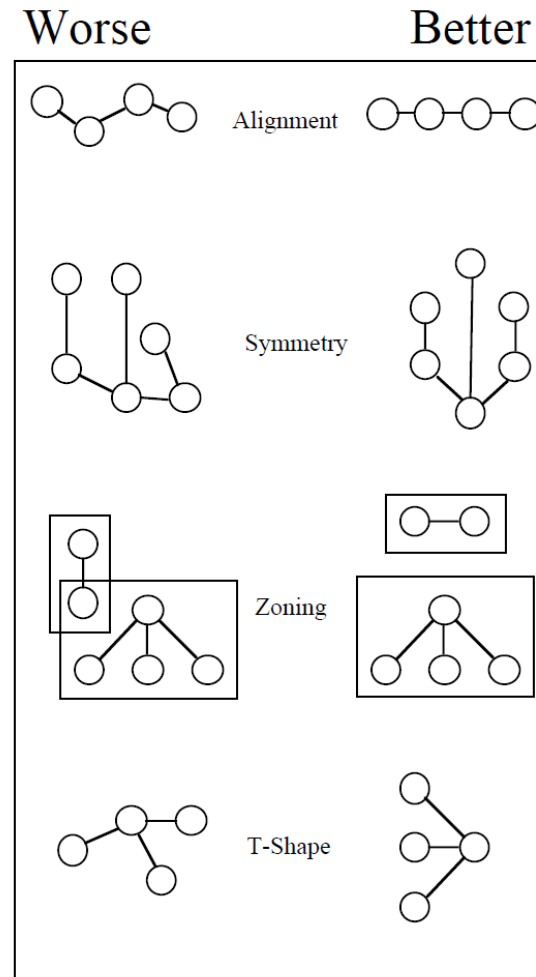
Graph drawing

- Mapping graph attributes (derived from the underlying data) to the aesthetics geometric objects, which represent the graph nodes
 - We will focus on general graphs (there exist separate solutions for trees)
- **Graph drawing** problem
 - **Input:** set of nodes and edges
 - **Output:** positions of nodes and curved shapes representing edges → graph layout

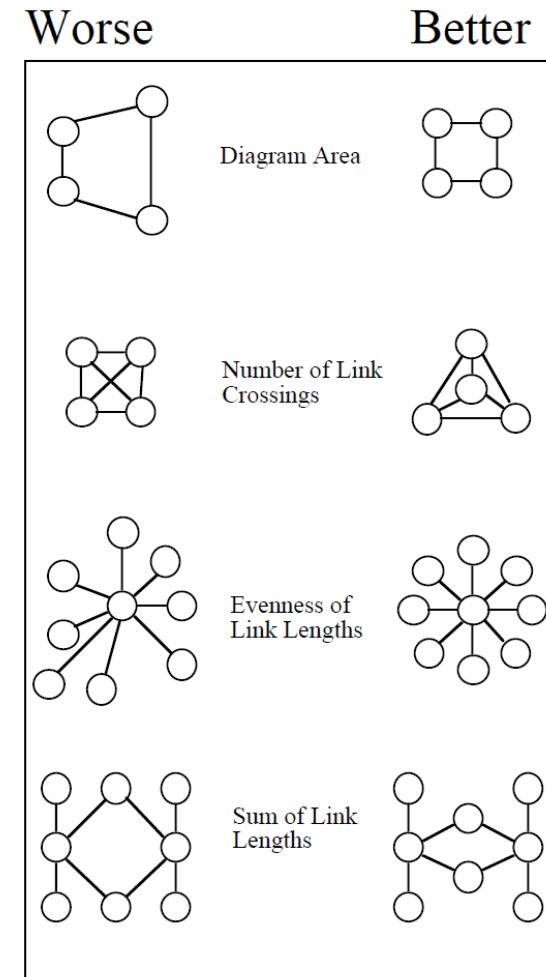
Graph drawing criteria (1)



Syntactic Validity



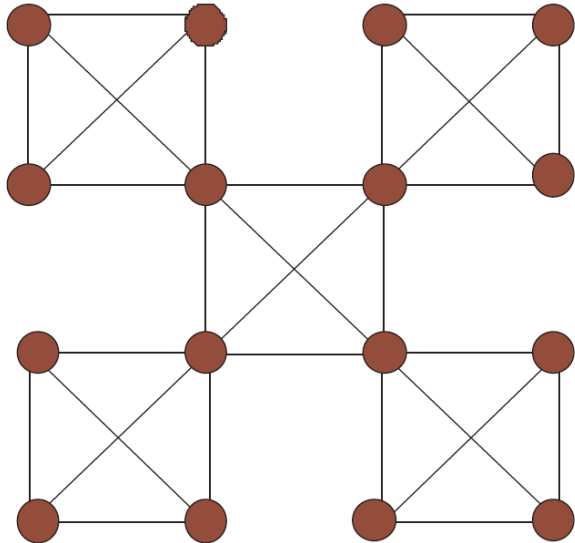
Perceptual Organization



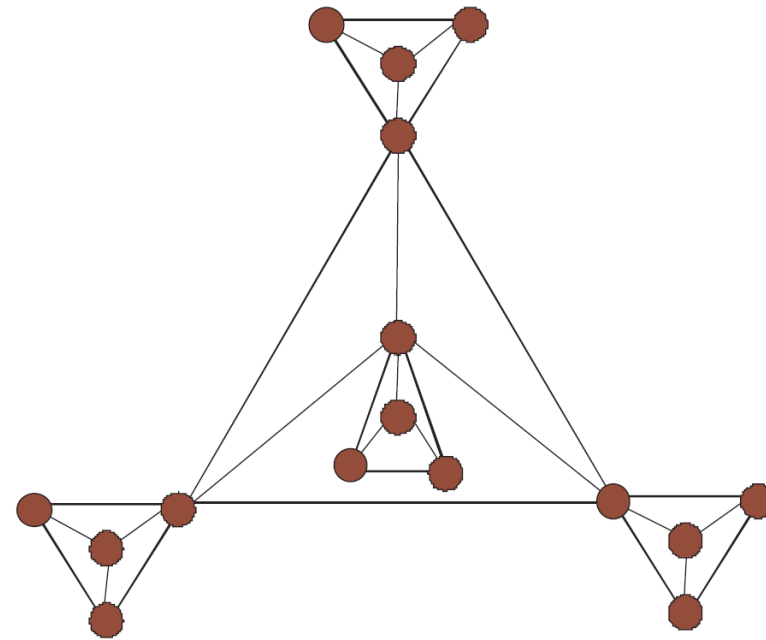
Aesthetic Optimality

Graph drawing criteria (2)

- However, some of aesthetics criteria can be mutually exclusive



Symmetry optimization



edge crossings (planarity) optimization

Basic graph layout strategies

- **Tree** layouts (not covered here)
- **Energy-based** layouts
 - Often called **force-directed placement** algorithms or **spring embedders**
- **Circular** layouts

Force-directed layouts

Force-directed layout

- **Force-directed placement (FDP)**

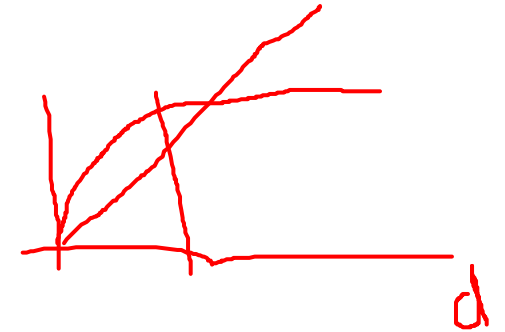
- Classical approaches
- Eades (84)
 - Kamada and Kawai (89)
 - Fruchterman and Reingold (90)
 - Davidson and Harel (96)
 - **Optimization**
 - Multi-scale approaches
 - ...

- **Clustering**

- Linlog (Noack (07))
- ...



FDP – Eades (1)



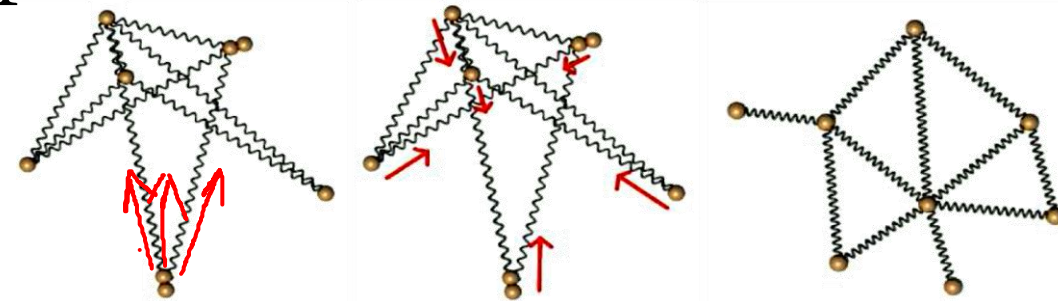
- **Spring-based** mechanical model that embeds a graph in 2D
 - **Vertices** = charged steel **rings**, **edges** = **springs**
 - Let the **spring forces** on the rings move the **system to a minimal energy state**
 - **Spring** exerted force $f_a = c_1 \times \log(d/c_2)$
 - **Nonadjacent vertices repel** each other with force $f_r = c_3/\sqrt{d}$

No force for $d = c_2$

Linear springs (Hook's Law) turn out to be too strong when vertices are far apart.

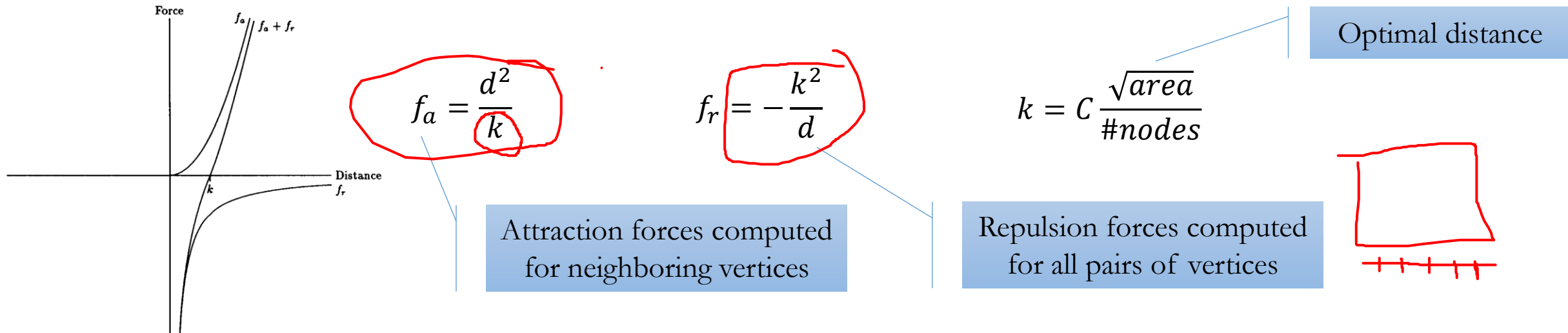
Algorithm

1. Place vertices in random locations. Set $i = 1$
2. Stop if $i = M, i = i + 1$
3. Calculate forces acting on each vertex
4. Move each vertex $c_4 \times \text{force_on_vertex}$



FDP - Fruchterman and Reingold

- Includes (even) **vertex distribution** in the layout into the set of criteria

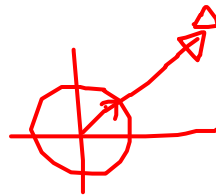


- Includes **temperature cooling**
 - Cooling bounds vertex movement similarly to simulated annealing

```

area ← W * L ;
initialize G = (V, E) ;
k ← √(area/|V|) ;
function fr(x) = k2/x ;
for i = 1 to iterations do
    foreach v ∈ V do
        v.disp := 0 ;
        for u ∈ V do
            if (u ≠ v) then
                Δ ← v.pos - u.pos ;
                v.disp ← v.disp + (Δ/|Δ|) * fr(|Δ|) ;
        function fa(x) = x2/k ;
        foreach e ∈ E do
            Δ ← e.v.pos - e.u.pos ;
            e.v.disp ← e.v.disp - (Δ/|Δ|) * fa(|Δ|);
            e.u.disp ← e.u.disp + (Δ/|Δ|) * fa(|Δ|);
        foreach v ∈ V do
            /* limit max displacement to frame; use temp. t to scale */
            v.pos ← v.pos + (v.disp/|v.disp|) * min(v.disp, t);
            v.pos.x ← min(W/2, max(-W/2, v.pos.x));
            v.pos.y ← min(L/2, max(-L/2, v.pos.y));
            t ← cool(t) ;

```



/* initialize displacement vector */

/* distance between u and v */

/* displacement */

/* compute attractive force */

/* e is ordered vertex pair .v and .u */

/* limit max displacement to frame; use temp. t to scale */

/* reduce temperature for next iteration */

FDP – Davidson and Harel (1)

- Adds **edge crossings minimization** criterion and uses **simulated annealing** for optimization
- Simulated annealing
 1. Set initial configuration σ
 2. Set initial temperature T
 3. Choose a new configuration σ' from a close neighborhood of σ
 4. Evaluate energy functions E and E' of configurations σ and σ'
 5. If $(E' < E \text{ or } r < e^{-(E-E')/T})$ then $\sigma \leftarrow \sigma'$
 6. Reduce temperature and go to $\textcircled{3}$

FDP – Davidson and Harel (2)

- **Energy function**

1. **Repulsive component**
2. **Placement component**

$$\sum_{i,j} \frac{\lambda_1}{d_{ij}^2}$$
$$\lambda_2 \left(\frac{1}{r_i^2} + \frac{1}{l_i^2} + \frac{1}{t_i^2} + \frac{1}{b_i^2} \right)$$

Distance to the right, left, top and bottom edges. High λ_2 value favors centered layouts.

- 3. **Edge length component**

$$\lambda_3 d_k^2$$

λ_3 is a normalization constant causing shortening edges to the necessary minimum.

4. **Crossings component**

$$\lambda_4 (\# \text{crossings})$$

Increasing λ_4 more heavily penalizes crossings.

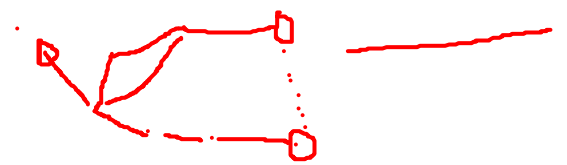
- **Cooling**

- Displacement of a vertex limited by circle of decreasing radius centered in current position



FDP - Kamada and Kawai (1)

- **Replacement of the ideal spring lengths by the shortest path distance** → graph-theoretic approach
 - In a connected graph, all pairs of vertices are connected → **no need for repulsive forces**
- **Correlates graph distances with the Euclidean distances** → minimizing the energy of the system corresponds to minimizing the difference between the geometric and graph distances



Energy function to be minimized

$$E = \sum_{i < j} \frac{1}{2} k_{ij} (|pos_i - pos_j| - l_{ij})^2$$

“Spring” strength

$$l_{ij} = L \times d_{ij}$$

$$L = \frac{\text{display_width}}{\max_{i < j} d_{ij}}$$

$$k_{ij} = \frac{K}{d_{ij}^2}$$

Shortest path between i and j

FDP - Kamada and Kawai (2)

$$E = \sum_{i < j} \frac{1}{2} k_{ij} (|pos_i - pos_j| - l_{ij})^2 = \sum_{i < j} \frac{1}{2} k_{ij} \left((x_i - x_j)^2 + (y_i - y_j)^2 + l_{ij}^2 - 2l_{ij} \sqrt{(x_i - x_j)^2 + (y_i - y_j)^2} \right)$$

- At a **local minimum** all the partial derivatives of $E(x_1, y_1, \dots, x_n, y_n)$ equal zero ($\forall j \frac{\partial E}{\partial x_j} = \frac{\partial E}{\partial y_j} = 0$)
- In Kamada and Kawai, **one node at a time is optimized** using the Newton method. At each step, the node m with the highest distance from the minimum (maximum gradient value) is chosen (the others are frozen):

$$\Delta_m = \sqrt{\left(\frac{\partial E}{\partial x_m}\right)^2 + \left(\frac{\partial E}{\partial y_m}\right)^2}$$

$O(n^3)$ using Floyd-Warshall algorithm

compute pairwise distances $d_{i,j}$ for $1 \leq i \neq j \leq n$;
compute pairwise ideal lengths $l_{i,j}$ for $1 \leq i \neq j \leq n$;
compute pairwise spring strength $k_{i,j}$ for $1 \leq i \neq j \leq n$;
initialize particle positions p_1, p_2, \dots, p_n ;

while ($\max_i \Delta_i > \epsilon$) **do**

let p_m be the particle satisfying $\Delta_m = \max_i \Delta_i$;

while ($\Delta_m > \epsilon$) **do**

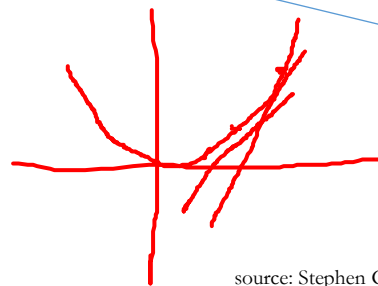
compute δx and δy by solving the following system of equations:

$$\frac{\partial^2 E}{\partial x_m^2}(x_m^{(t)}, y_m^{(t)})\delta x + \frac{\partial^2 E}{\partial x_m \partial y_m}(x_m^{(t)}, y_m^{(t)})\delta y = -\frac{\partial E}{\partial x_m}(x_m^{(t)}, y_m^{(t)});$$

$$\frac{\partial^2 E}{\partial y_m \partial x_m}(x_m^{(t)}, y_m^{(t)})\delta x + \frac{\partial^2 E}{\partial y_m^2}(x_m^{(t)}, y_m^{(t)})\delta y = -\frac{\partial E}{\partial y_m}(x_m^{(t)}, y_m^{(t)})$$

$$x_m \leftarrow x_m + \delta x;$$

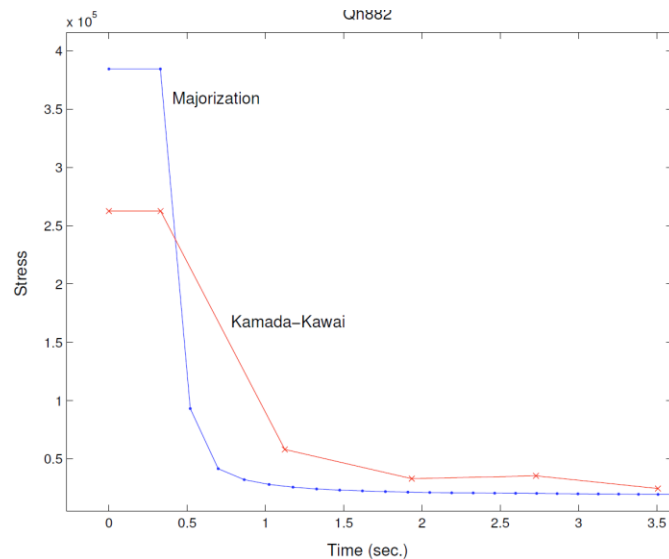
$$y_m \leftarrow y_m + \delta y;$$



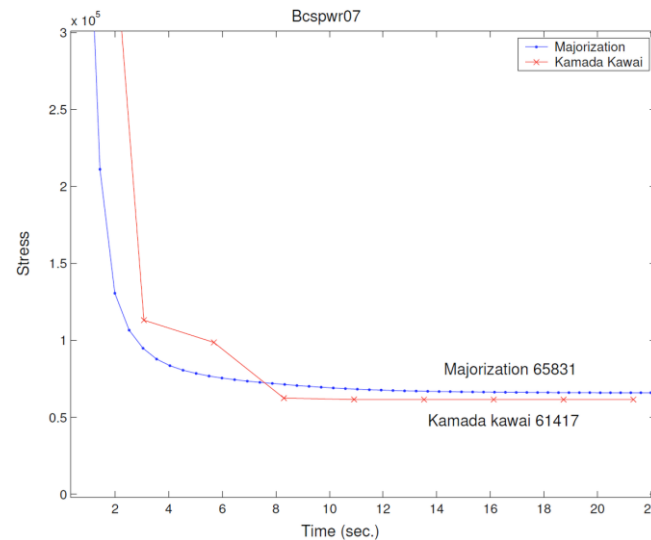
Two-dimensional
Newton-Raphson method

Relation of KK algorithm to MDS

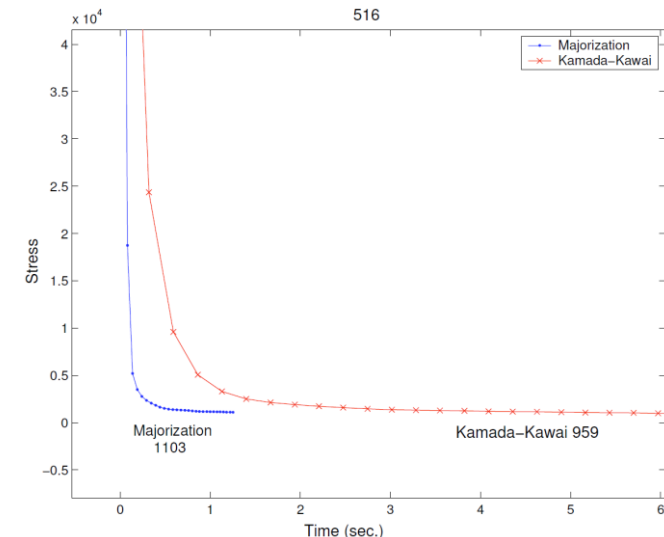
- The layout **energy function** $E = \sum_{i < j} \frac{1}{2} k_{ij} (|pos_i - pos_j| - l_{ij})^2$ is basically the same as the **stress** known in MDS \rightarrow instead of the Newton method **stress majorization** can be used, which is guaranteed to converge



$|V| = 882, |E| = 1533$



$|V| = 882, |E| = 1533$



$|V| = 516, |E| = 729$

Aesthetic criteria of FDP layouts

Criteria	Eades (1984)	Kamada and Kawai (1989)	Fruchterman and Reingold (1991)	Dauids and Harel (1996)
Symmetric	YES	YES		
Evenly distributed nodes		YES	YES	YES
Uniform edge length	YES	YES	YES	YES
Minimized edge crossings		YES	YES	YES

Source: Chen, Information visualization: Beyond the Horizon. Springer (2006)

Pros and cons of FDP

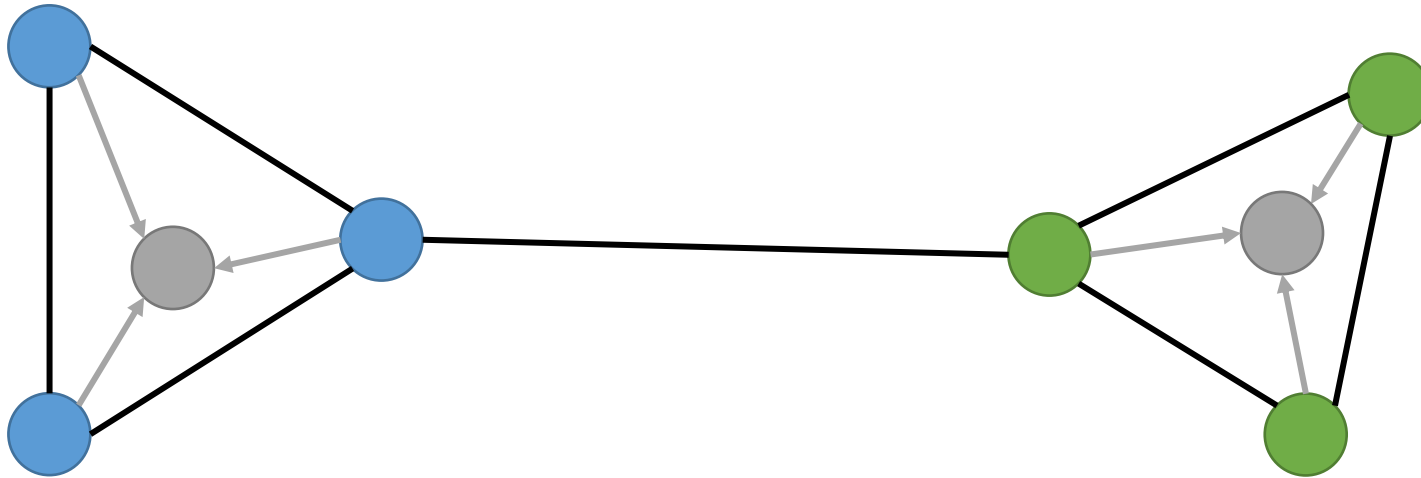
- Easy **interpretation**
- Easy **implementation**
- Easy to add **new heuristics**

- **Run time**
- Can end up in **local minima**
- Does **not reflect** the inherent graph **cluster structure** (if required) – see the following slides

Cluster separation of FDP layouts

Dummy attractors

- One can achieve better **separation of clusters** based on **underlying information** by adding **dummy attractors**
 - Requires prior knowledge about the structure of the underlying data

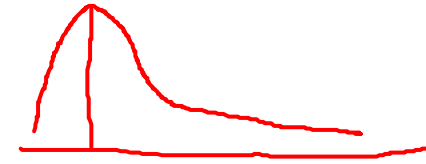


Graph-clustering focused energy modelling

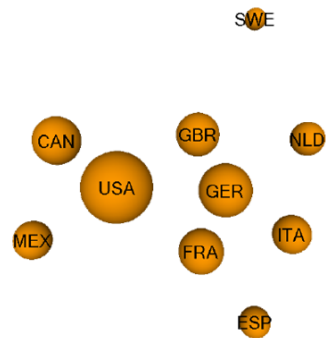
- Layout algorithms can be viewed as having two components – *energy model* and *energy minimization algorithm*
- **Up to now**, the focus of the energy models has been on producing generally readable layouts enforced by small and uniform edge lengths
→ tendency to group large-degree nodes in the center of the layout
- There exist energy models focusing rather on **good separation of clusters** → LinLog model

LinLog energy models (1)

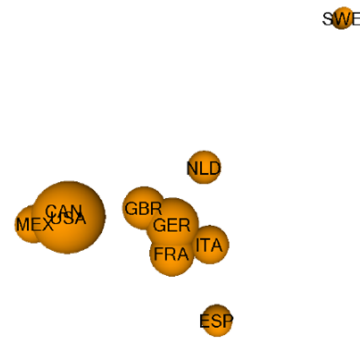
- Two energy models
 - **Node-repulsion**
 - **Edge-repulsion**



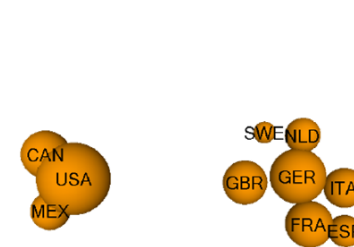
- **Not biased** towards grouping together **nodes with high degree** → appropriate for many real-world graphs with right-skewed degree distributions



(a) Fruchterman-Reingold

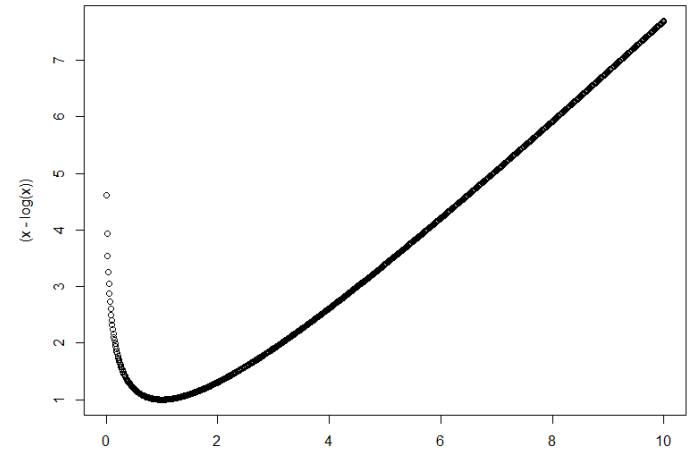


(b) Node-repulsion LinLog



(c) Edge-repulsion LinLog

LinLog energy models (2)



- **Node-repulsion** model of a layout p

$$U_{Node}(p) = \sum_{\{u,v\} \in E} \|p(u) - p(v)\| - \sum_{\{u,v\} \in V^2} \ln \|p(u) - p(v)\|$$

Attraction forces

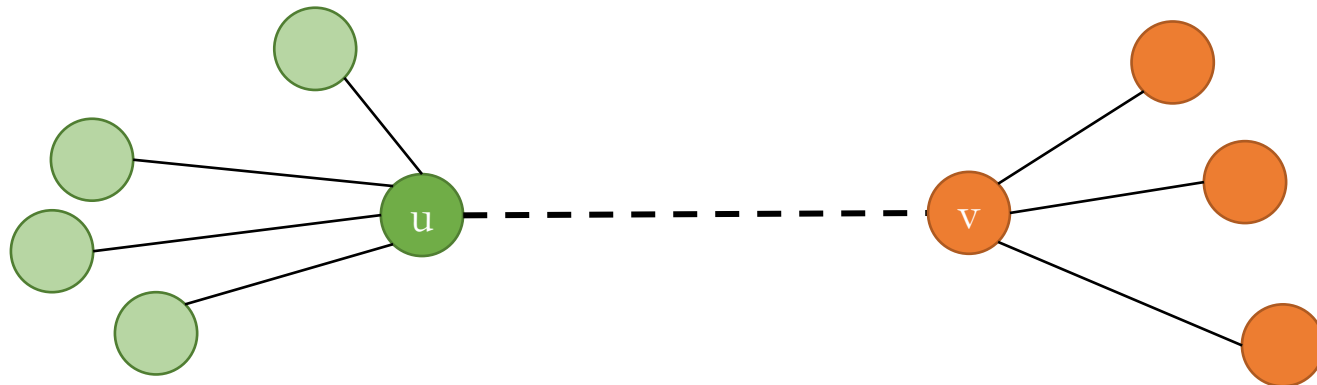
Repulsion forces

- **Edge-repulsion** model of a layout p

$$U_{Edge}(p) = \sum_{\{u,v\} \in E} \|p(u) - p(v)\| - \sum_{\{u,v\} \in V^2} \deg(u) \deg(v) \ln \|p(u) - p(v)\|$$

Nodes represented in terms of their neighboring nodes \rightarrow edges

Each node's influence on the layout is proportional to its degree.



Clustering criteria

- **Good clustering** consists of subgraphs with **many internal** and **few external edges** - **clustering criteria** formalize this notion

- $V_1, V_2 \subset V$: $\mathbf{cut}(V_1, V_2) = |\{ \{v_1, v_2\} \in E \mid v_1 \in V_1, v_2 \in V_2 \}|$

- Expected cut is

Probability of an edge between two arbitrary nodes

$$\frac{|E|}{(|V|^2 - |V|)/2} (|V_1| |V_2|) = \frac{2|E|(|V_1| |V_2|)}{|V|^2 - |V|}$$

#possible pairs between two node sets

- which is **biased** towards **uneven cluster sizes** → layout strategies driven with this criterion will favor uneven clusters

LinLog clustering criteria

- **Node-normalized cut**

$$\mathbf{ncut}(V_1, V_2) = \frac{\text{cut}(V_1, V_2)}{|V_1||V_2|}$$

- Still **biased** towards uneven partitions if the **number of edges used as a measure** of subgraph size

- **Edge-normalized cut**

$$\mathbf{ecut}(V_1, V_2) = \frac{\text{cut}(V_1, V_2)}{\text{deg}(V_1) \text{deg}(V_2)}$$

LinLog energy and clustering

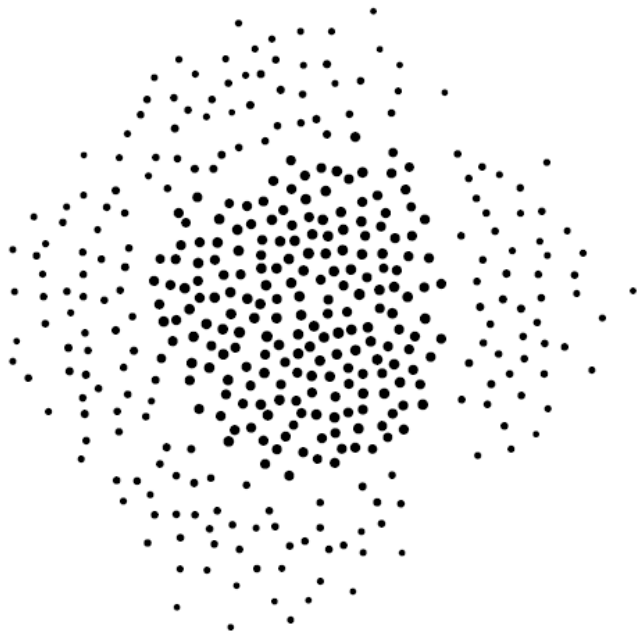
Each node v can be imagined as an end point of the $\deg(v)$ connected edges

- Minimal-energy **node-repulsion** layouts **minimize the ratio** of the **mean distance** between **connected nodes** to the mean distance **between all nodes**
- The distance of two dense and sparsely connected clusters V_1, V_2 approximates $1/\text{ncut}(V_1, V_2)$ in minimal-energy node-repulsion LinLog layouts

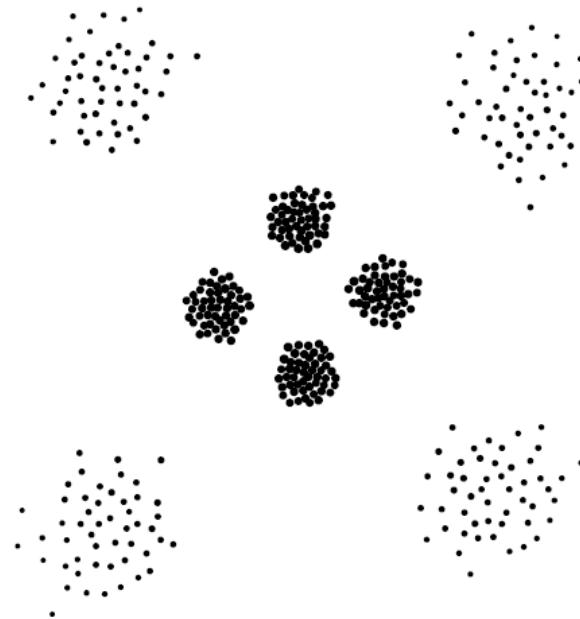
- Minimal-energy **edge-repulsion** layouts **minimize the ratio** of the **mean distance** between **connected *end* nodes** to the mean distance between **all *end* nodes**
- The distance of two dense and sparsely connected clusters V_1, V_2 approximates $1/\text{ecut}(V_1, V_2)$ in minimal-energy edge-repulsion LinLog layouts

Pseudorandom graph with 400 nodes consisting of 8 equal-sized clusters:

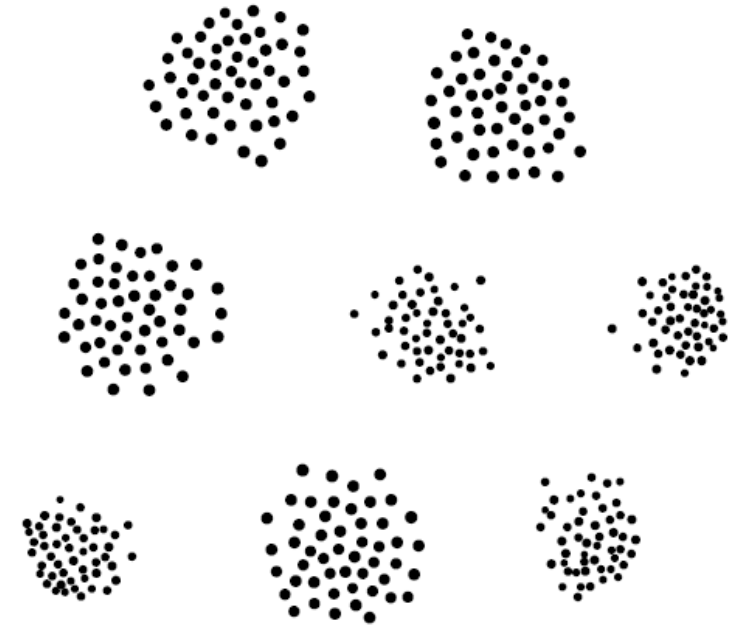
- Cluster 1-4 : intra-cluster edge probability = 1
- Cluster 5-8 : intra-cluster edge probability = 0.5
- Cluster 1-4 : inter-cluster edge probability = 0.2
- Cluster 5-8 : inter-cluster edge probability = 0.05
- Cluster 1-4 vs 5-8 : inter-cluster edge probability = 0.1



Fruchterman-Reingold



Node-repulsion LinLog

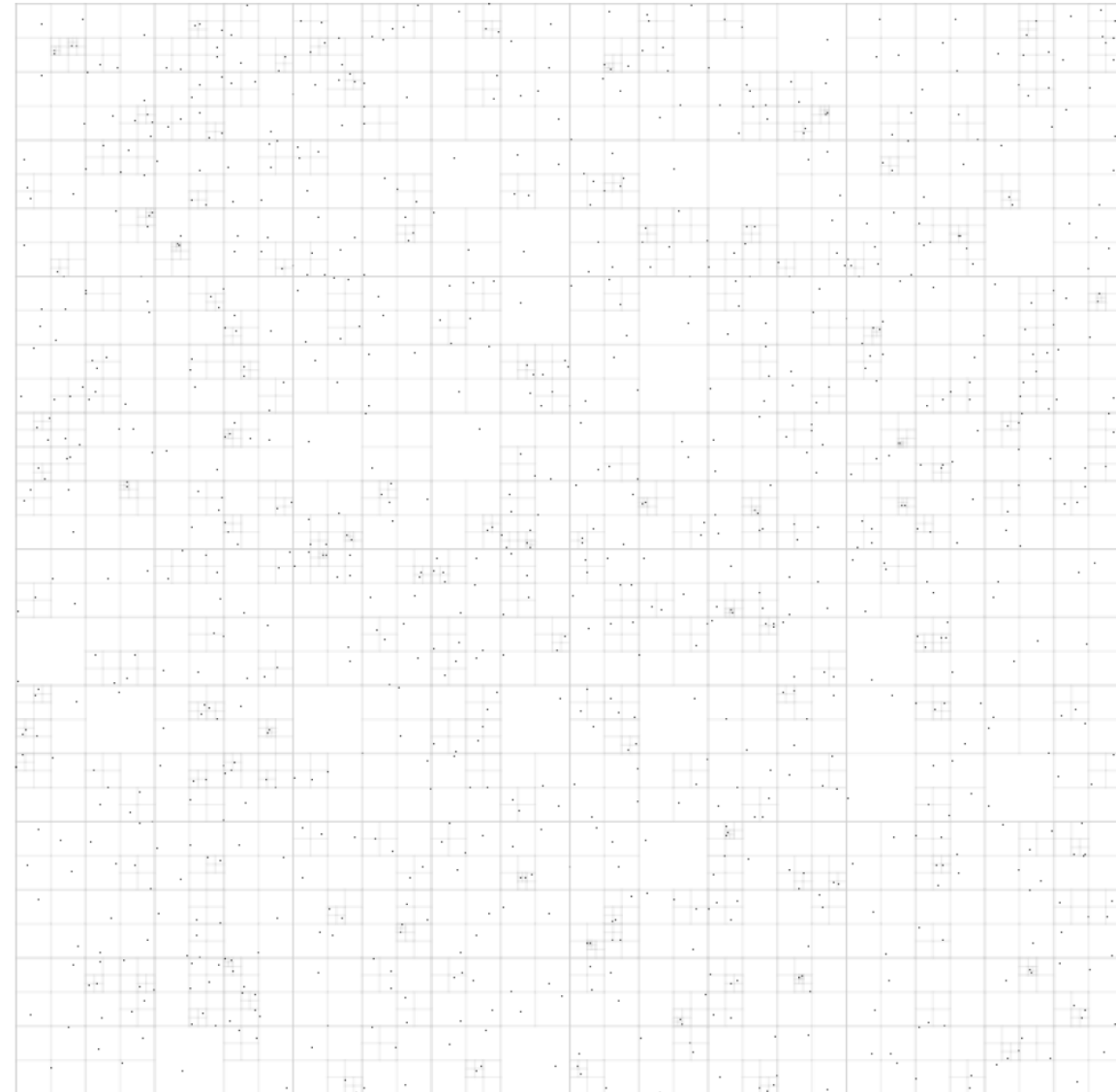
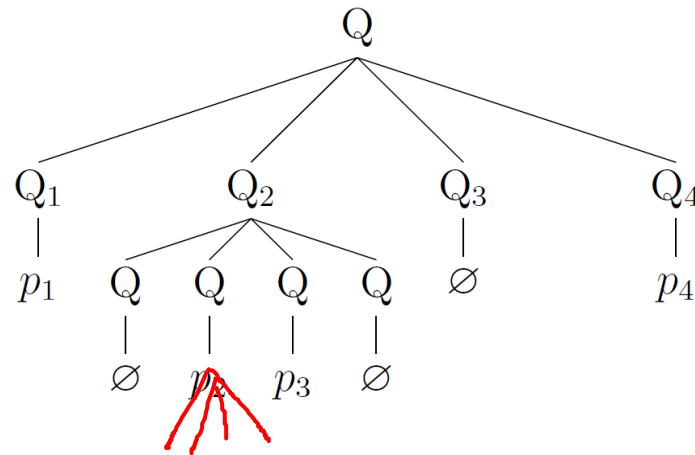
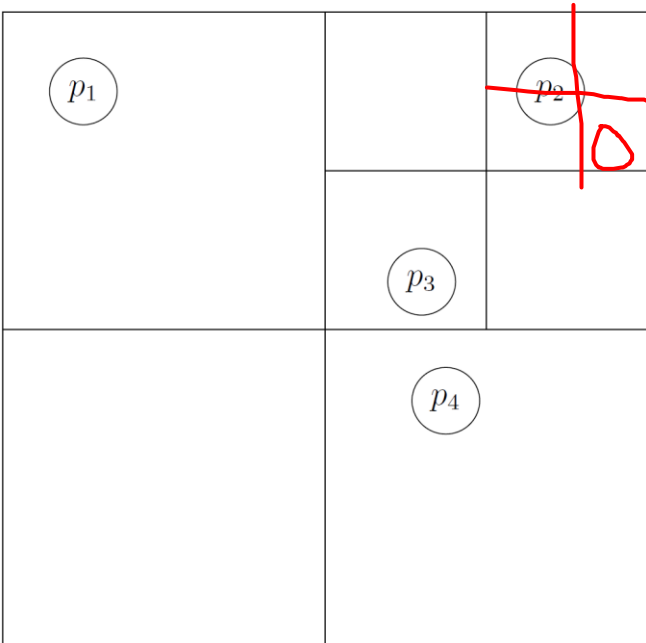


Edge-repulsion LinLog

Speed optimization of the FDP layouts

Barnes-Hut approximation (1)

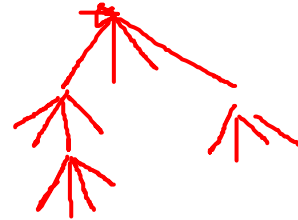
- Division of space using quad-tree \rightarrow inner nodes form clusters for approximation



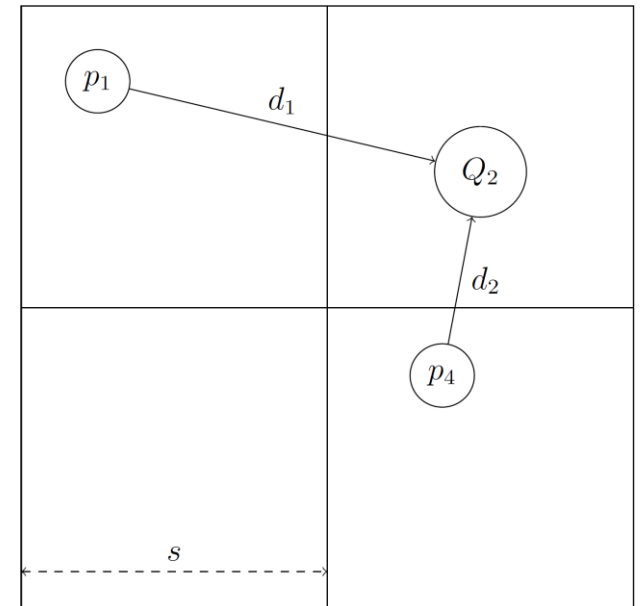
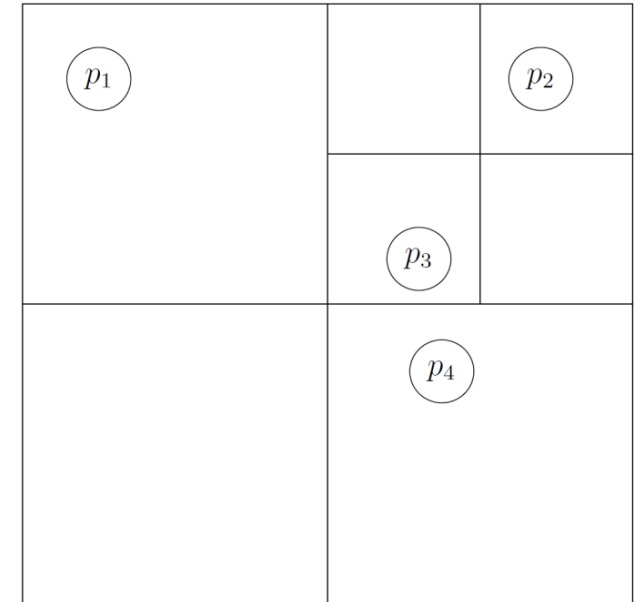
Barnes-Hut approximation (2)

- An object is compared either to other objects or cluster centers based on **Barnes-Hut criterion**

$$\frac{s}{d} \leq \theta$$



- If for given level of tree the **condition holds**, the **repulsive force** is computed only **with respect to the cluster**
- θ affects how often the criterion is applied \rightarrow
 $O(n^2)$ vs $O(n)$



Barnes-Hut approximation (3)

Center of gravitation

- **Repulsive forces** need to be **modified**

$$f_r(x_i, x_j) = -\frac{k^2}{d(x_i, x_j)} \rightarrow f_r(x_i, x_Q) = -\frac{|Q|k^2}{d(x_i, x_Q)} \dots x_Q = \frac{\sum_{j \in Q} x_j}{|Q|}$$

- The algorithm is modified so that **quad tree** is **constructed** at the **beginning of each iteration**
 - Complexity of construction is $O(dn)$, where d is depth of the tree
- **Repulsive forces** are computed using the modified f_r to **clusters centers of gravitation** only

Small world networks

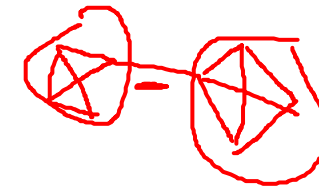
- Many of the real-world networks have so-called **small world networks (SWN)** property
 - Average path length in a SWN compares to a path length in a random graph
 - Clustering indexes of SWN nodes are on average orders of magnitudes larger

$$c(v) = \frac{|edges(neighbors(v))|}{(k(k-1))/2}$$

Number of vertices in the neighborhood of v

How close neighbors of v resemble a clique.

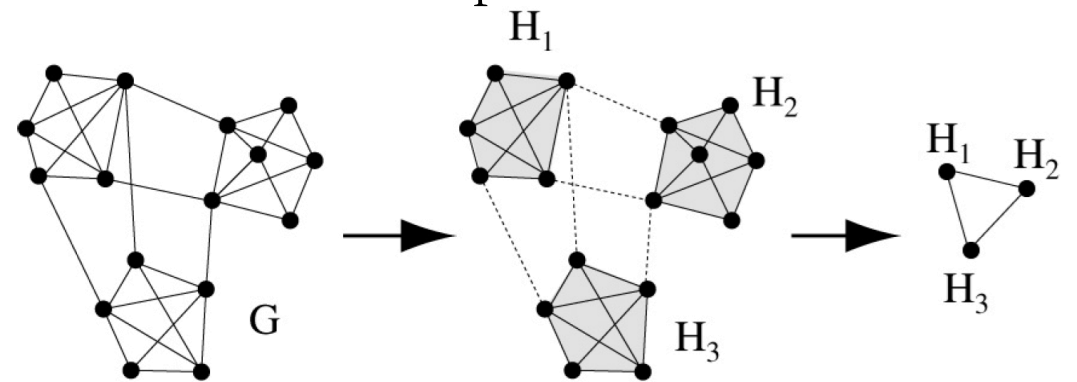
Graph	Clustering index	Ave Path Length	Random graph (same number of nodes and edges)	
IMDB	0.9666	3.2043	0.0243	2.6694
“Resyn assistant”	0.9518	3.2847	0.1942	1.8195
Mac OS 9	0.3875	2.8608	0.0179	3.3196
Web ²	0.1078	3.1	2.3e-4	-
.edu sites ²	0.156	4.062	0.0012	4.048



Multi-scale algorithms

- **Most networks** are not only SWN, but their **components form SWN** as well → hierarchy of SWNs → **multi-scale networks**
 - Leaf level of the hierarchy consists of cliques
- Series of graph representations with **different levels of details** → **optimization of the layout** with respect to these coarse abstractions of the original graph
 - One has to define the notion of coarse-scale representations of a graph, in which the combinatorial structure is significantly simplified but features important for visualization are well preserved.

- The multi-scale algorithms differ in
 - Bottom-up vs top-down
 - Coarsening
 - Underlying layout methods



We will show here
just two examples

Multi-scale algorithms – Harel & Koren

- **Coarsening** process based on a **k -center approximation** –
 - Find k nodes, so that the maximum distance of any node to arbitrary of the k nodes is minimized
 - NP-hard, but can be 2-approximated
- The algorithm in **each iteration**
 - Finds **centers** of clusters
 - Finds **layout for the cluster centers**
 - **Moves nodes close** to their respective cluster **centers**

Input: graph $G = (V, E)$

Output: Find L , a nice layout of G

$Iterations = 4$; /* number of iterations in local beautification */

$Ratio = 3$; /* ratio of vertex sets in consecutive levels */

$Rad = 7$; /* radius of local neighborhood */

$MinSize = 10$; /* size of coarsest graph */

Compute all-pairs shortest path lengths $d_{i,j}$;

Initialize layout L by placing vertices at random;

$k \leftarrow MinSize$;

while $k \leq |V|$ **do**

$C \leftarrow \mathbf{K-Centers}(G(V, E), k)$;

$radius = \max_{v \in C} \min_{u \in C} \{d_{vu}\} * Rad$;

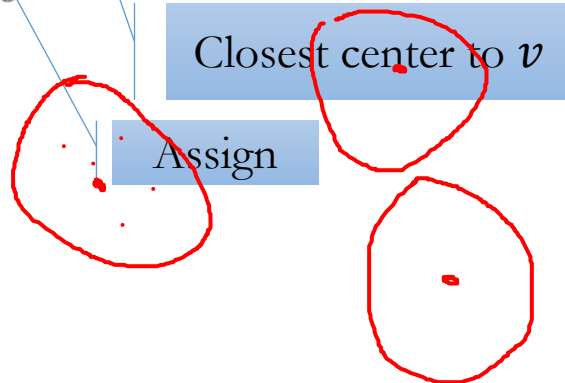
$\mathbf{LocalLayout}(d_{C \times C}, L(C), radius, Iterations)$;

foreach $v \in V$ **do**

$L(v) \in L(\mathit{center}(v)) + \mathit{rand}$;

$k \leftarrow kRatio$

return L ;



$\mathbf{K-Centers}(G(V, E), k)$;

Input: Graph $G = (V, E)$ and constant k

Output: Compute set $S \subseteq V$ of size k , s.t. $\max_{v \in V} \min_{s \in S} \{d_{sv}\}$ is minimized

$S \leftarrow \{v\}$ for some arbitrary $v \in V$;

for $i = 2$ **to** k **do**

 find vertex u farthest away from S ;

 (i.e., such that $\min_{s \in S} \{d_{us}\} \geq \min_{s \in S} \{d_{ws}\}, \forall w \in V$);

$S \leftarrow S \cup \{u\}$;

return S ;

$\mathbf{LocalLayout}(d_{V \times V}, L, k, Iterations)$;

Input: APSP matrix $d_{V \times V}$, layout L , constants k and $Iterations$

Output: Compute locally nice layout L by beautifying k -neighborhoods

for $i = 1$ **to** $Iterations * |V|$ **do**

 Choose the vertex v with the maximal Δ_v^k ;

 Compute δ_v^k as in Kamada-Kawai;

$L(v) \leftarrow L(v) + (\delta_v^k(x), \delta_v^k(y))$;

Multi-scale algorithms - Auber

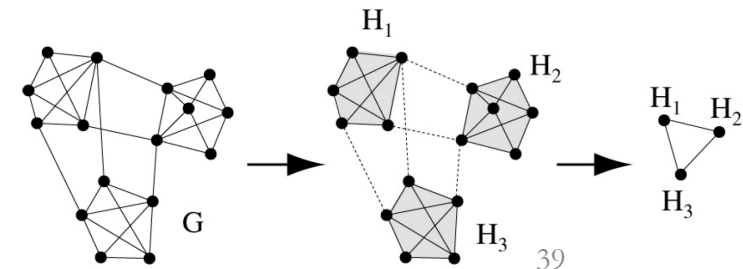
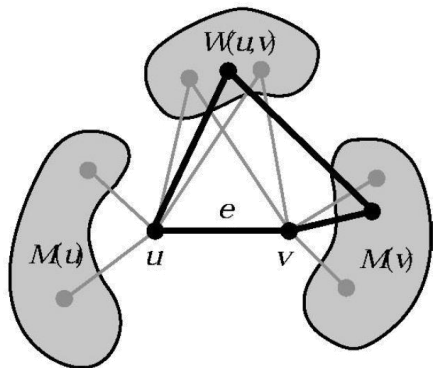
- Iteratively **filter out edges** with low edge clustering index \rightarrow remaining **connected components** form nodes of the **quotient graph**

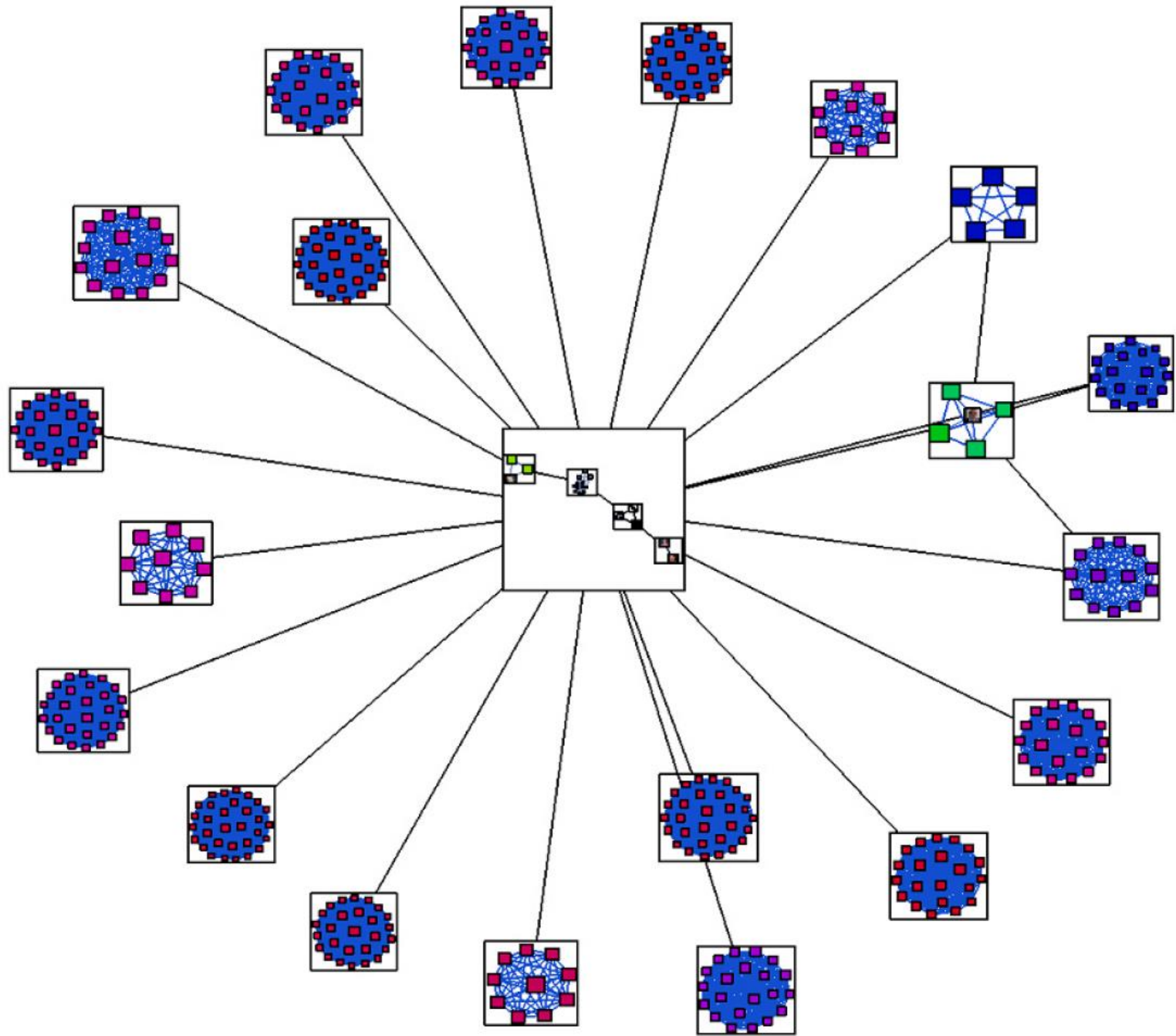
$$s(A, B) = \mathit{cut}(A, B) / |A||B|$$

Any edge connecting two of $M(u), M(v), W(u, v)$ is a part of a 4-cycle going through (u, v)

$$\begin{aligned} & \mathit{strength}(u, v) \\ &= s(M(u), W(u, v)) + s(W(u, v), M(v)) + s(M(u), M(v)) \\ &+ s(W(u, v)) + \underbrace{|W(u, v)| / |M(u) + W(u, v) + M(v)|}_{\text{Ratio of 3-cycles using } \{u, v\}} \end{aligned}$$

Ratio of 3-cycles using $\{u, v\}$



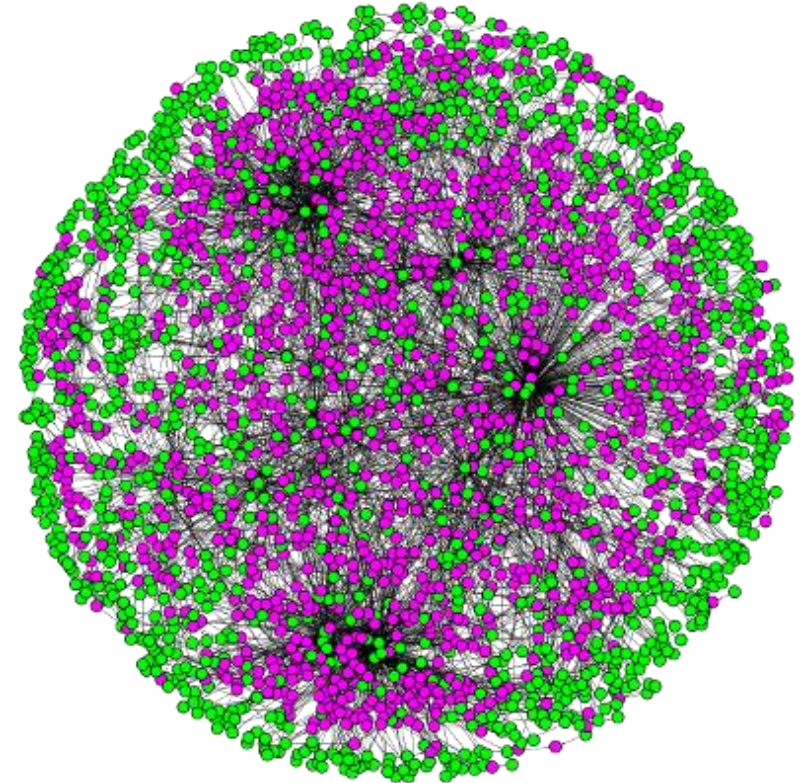


Very complex or specific networks

Hairballs

In case of complex data, node-link diagrams are notoriously difficult to visualize and interpret → *hairballs*

- *Junk food of network visualization* – very low nutritional value, leaving the user hungry
- A complex network does not necessarily communicate complex information

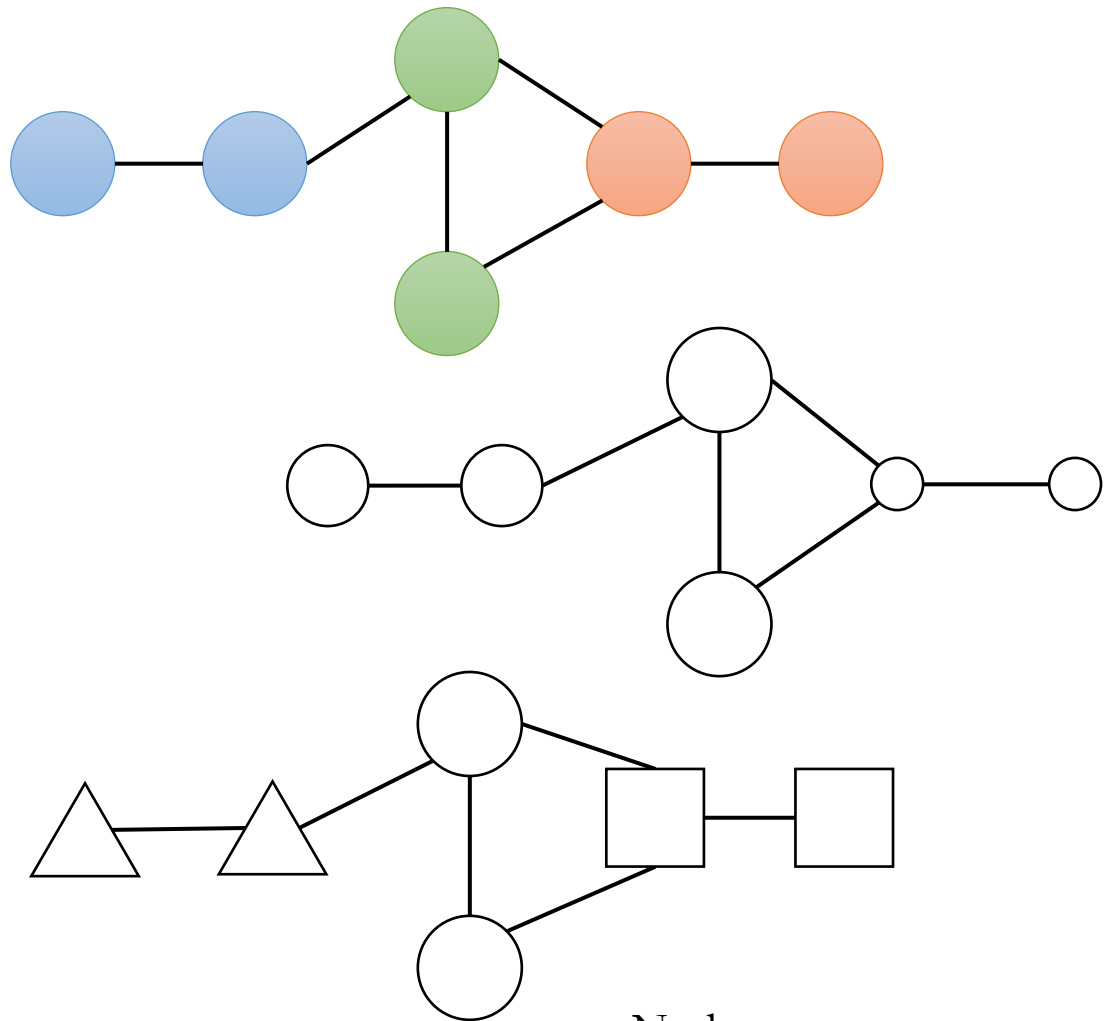


*E. coli metabolic network
visualized using Cytoscape*

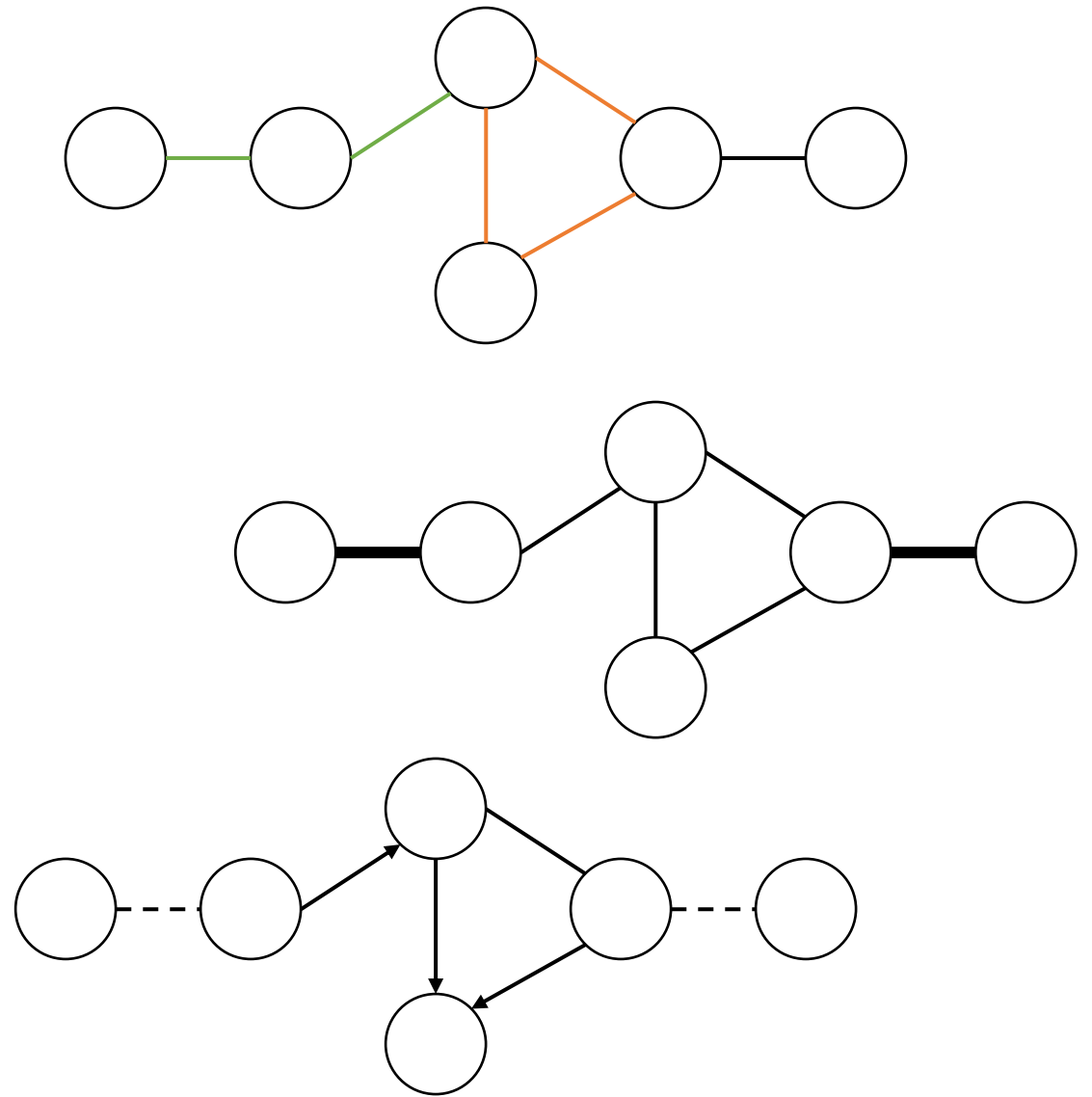
Interpretability of graph layouts

- **Interface problem**
 - Graph **sizes** tend to **grow** while the **size of the medium** we use to visualize them tends to stay **the same**
 - **Solutions**
 - **Visual encodings**
 - **Smart layouts**
 - **Flow diagrams, Circular layouts, matrix methods, hive plots, hierarchies, ...**
 - Often aim at specific use cases
 - **Interactivity** (navigation)
 - **Zooming, panning, fish-eye, ...**

Visual encodings



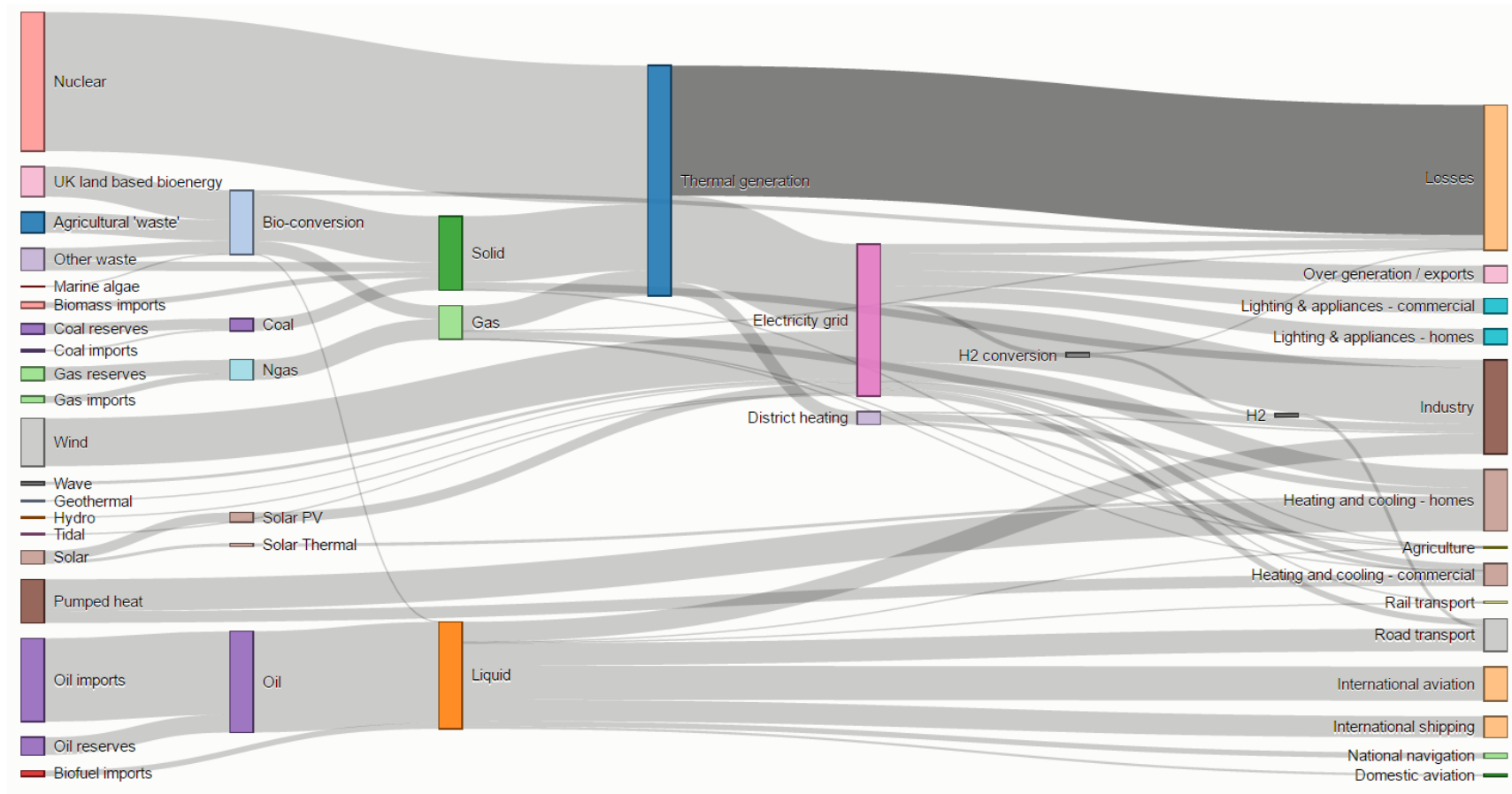
Nodes



Edges

Flow diagrams

- **Sankey diagrams** – flow magnitude proportional to connection width

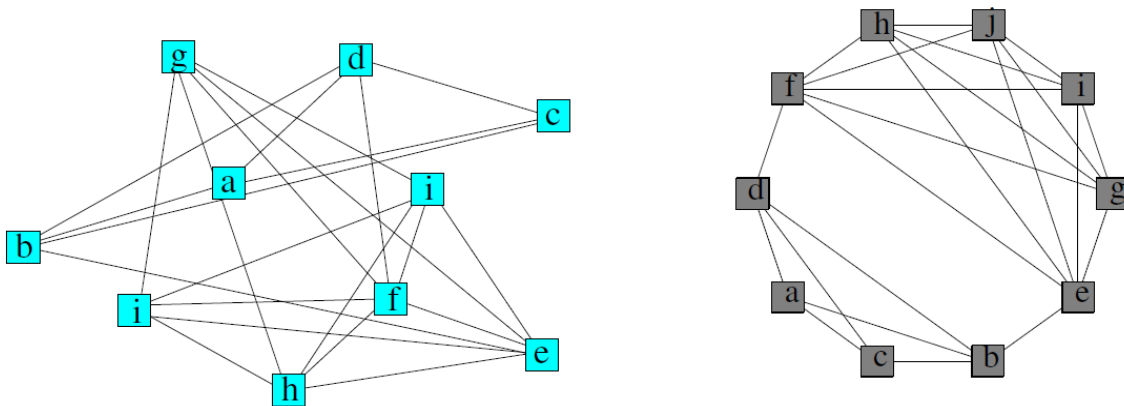


Circular layout

- Vertices placed on the circumference of an embedding circle
 - Often the vertices are first grouped into clusters

Basic layout

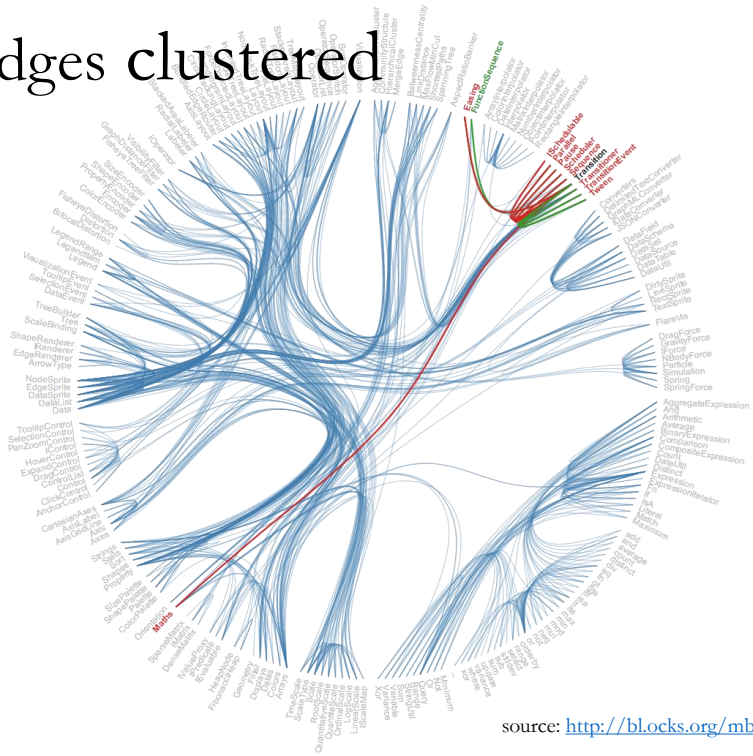
- Edges drawn as straight lines



source: Six and Tollis: Circular Drawing Algorithms. Handbook of Graph Drawing and Visualization (2013)

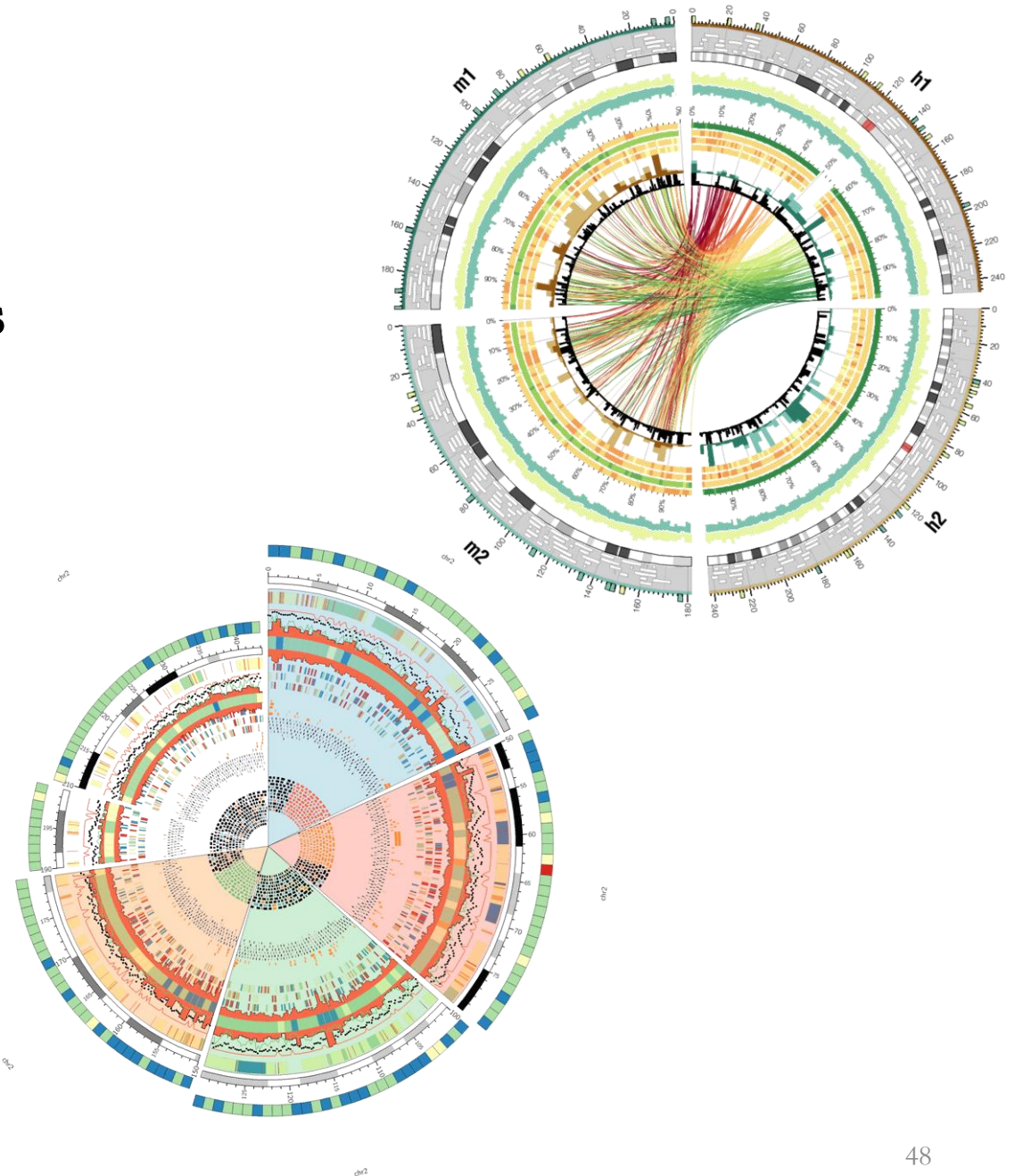
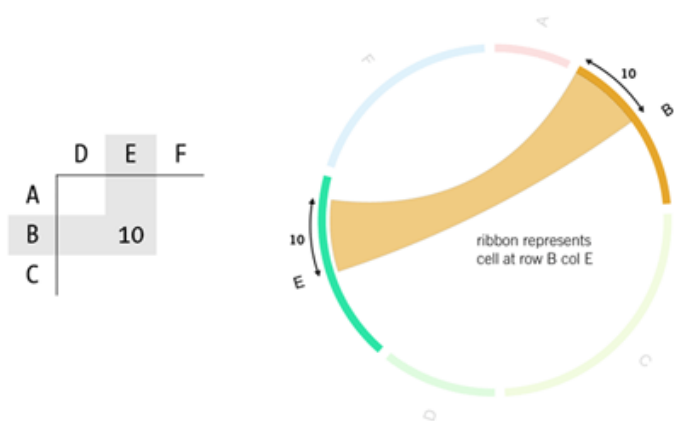
Edge binding

- Edges clustered



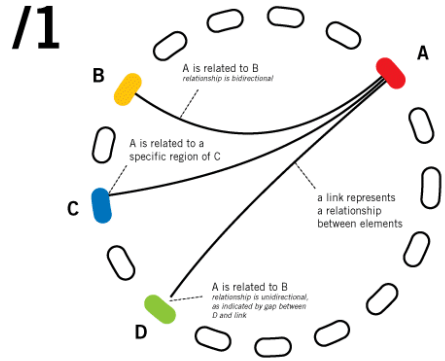
Circos

- [Software](#) for laying out **relationships** (graphs)
- Basically **converts tabular data to a circular layout** (any graph can be expressed as a table)

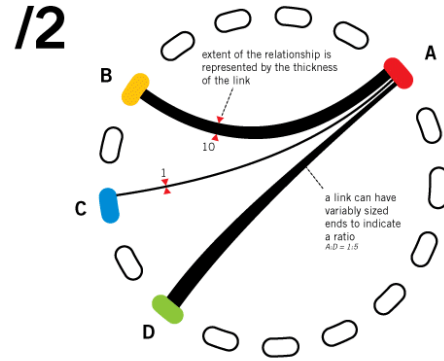




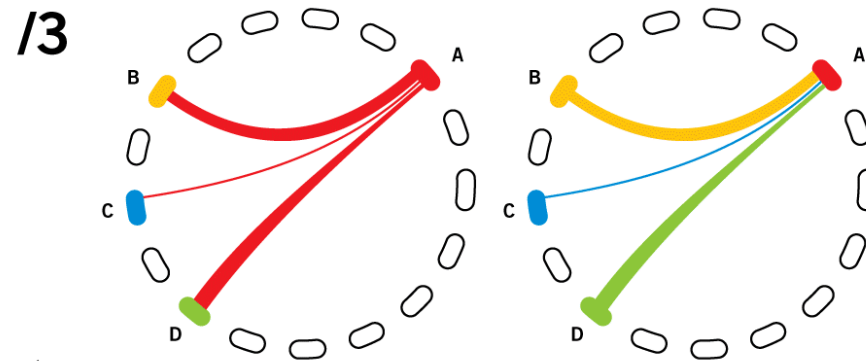
Applying Circos to Visualizing Tables using a Circular Layout



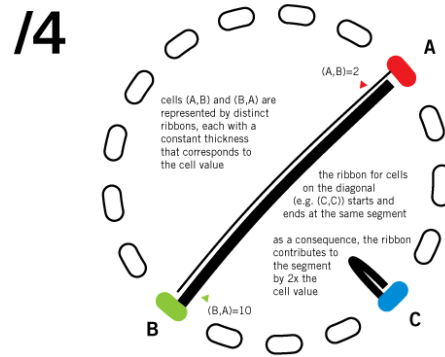
Relationships between elements in your data set are indicated by links. Links can indicate a simple relationship (A-B), a relationship that has positional information (A-C), or a unidirectional relationship (A-D). In each case, the link is formatted differently.



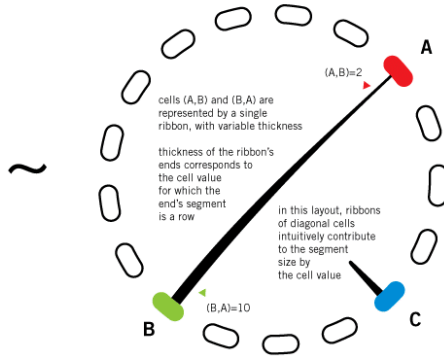
Links with variable thickness can represent the extent or magnitude of the relationship between elements.



When links are colored based on the elements that they relate, spotting patterns is easier. In particular, when relationships have a direction, links can be colored by source or target element.

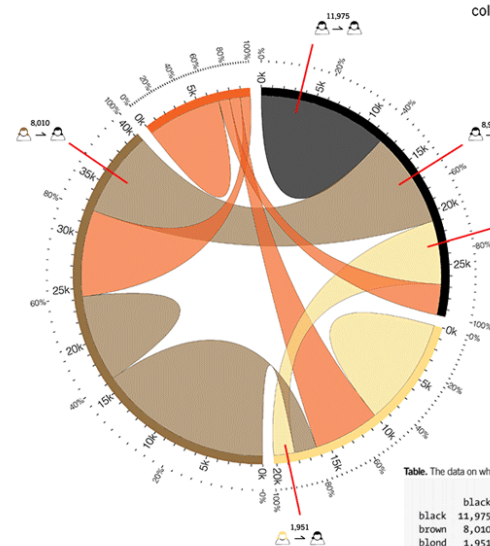
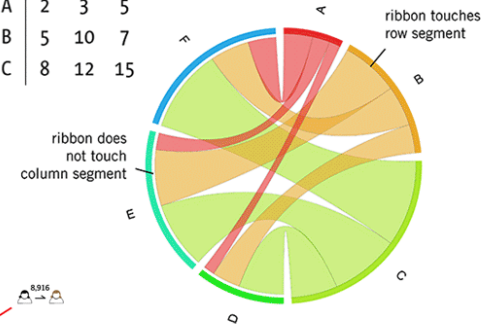


Transpositive cells (e.g. (A,B) and (B,A)) are here shown by a single ribbon (right). The ribbon now ends directly at both row and column segments and its ends are of variable thickness, which is the cell value for which the end's segment is a row. For example, if (A,B)=2 and (B,A)=10 then the ribbon's end touching A is thickness 2 and the end touching



/5 Ribbons touch the row segments, but terminate a short distance before reaching the column segment.

	D	E	F
A	2	3	5
B	5	10	7
C	8	12	15



/6

In the ratio layout, a ribbon represents two cells, except for cells on the diagonal which are represented by a single ribbon. By following the size of the ribbon from one segment to another, you can visually estimate the ratio of (A,B):(B,A).

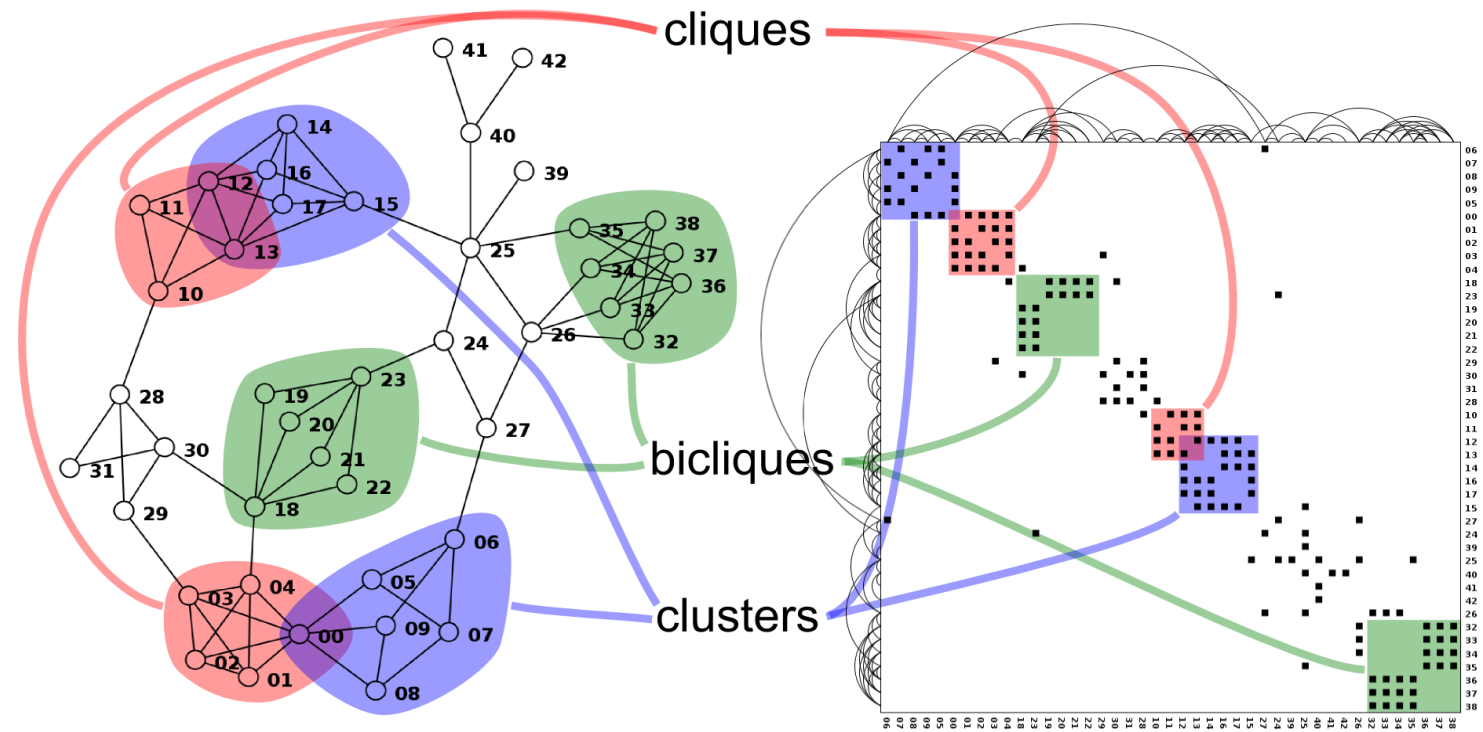
Table visualization is implemented in the tableviewer utility, bundled with Circos. The parse-table script parses a tabular data file into an intermediate format, which is then used by make-conf to create Circos configuration and data files. Circos is then used to generate the final image.

Table. The data on which this figure is based.

	black	brown	blond	red
black	11,975	8,916	5,871	2,868
brown	8,010	8,090	16,145	8,045
blond	1,951	2,060	10,048	6,171
red	1,013	940	990	6,907

Matrix methods

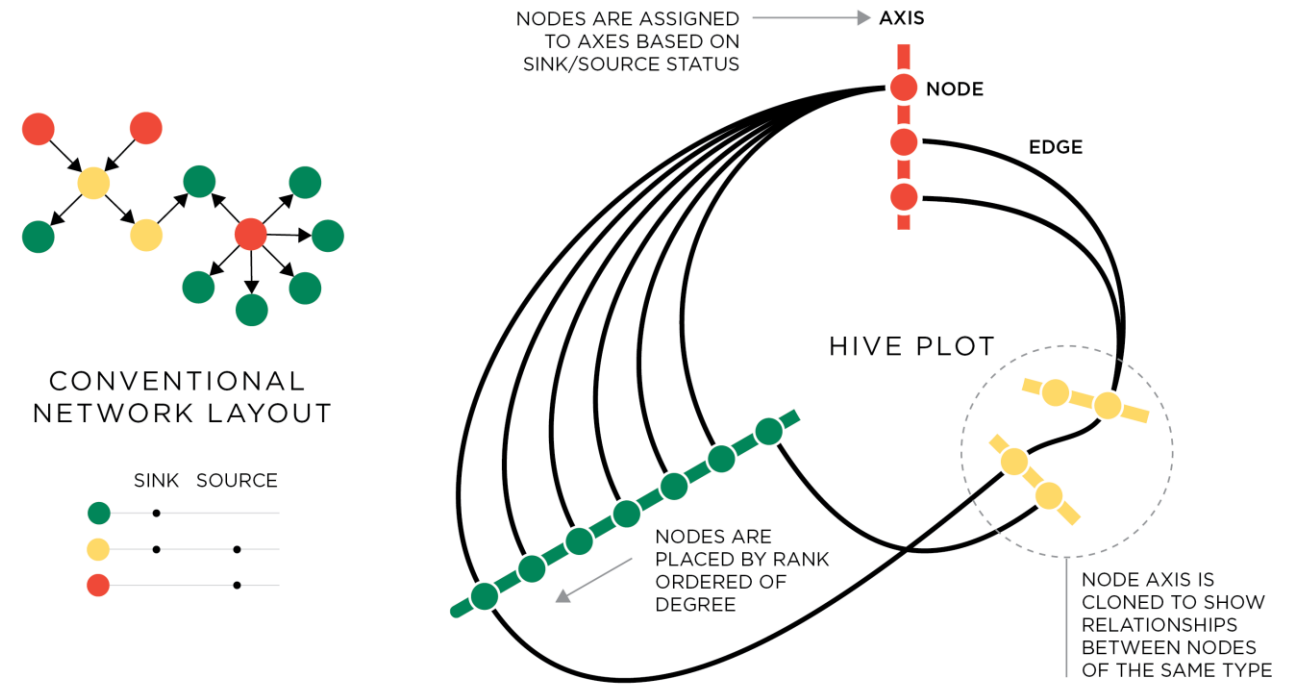
- Methods based on **analysis of the adjacency matrix**
- Requires some **support from the visualization tool**
 - Rows and columns of the matrix can be rearranged → patterns



source: McGuffin. Simple Algorithms for Network Visualization: A Tutorial (2012)

Hive plots

- **Nodes** mapped to and positioned on **radially distributed linear axes** → linear layout of nodes
 - Can be divided into **segments**
- **Edges** drawn as **curved links**
- Graph structure can be mapped to
 - Axis, Position, Color

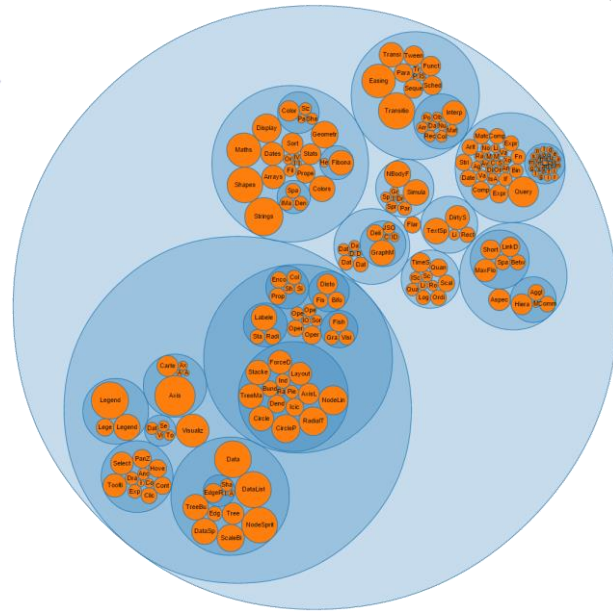


source: <http://www.hiveplot.net/conference/vizbi2011/poster/krzywinski-hiveplot-poster.png>

Hierarchical layouts

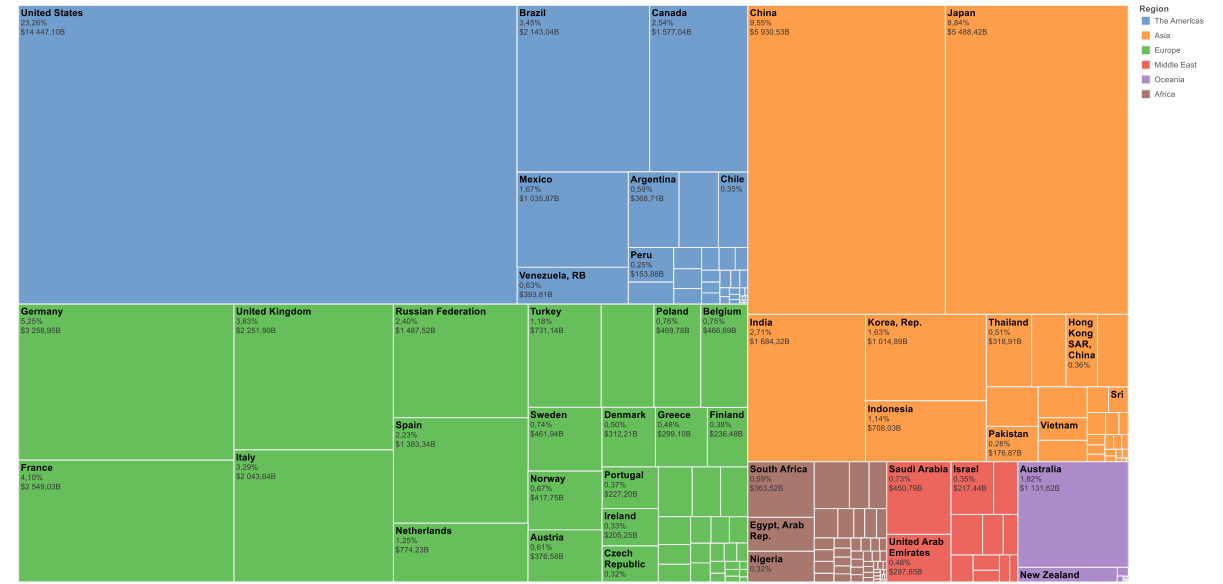


Sunburst



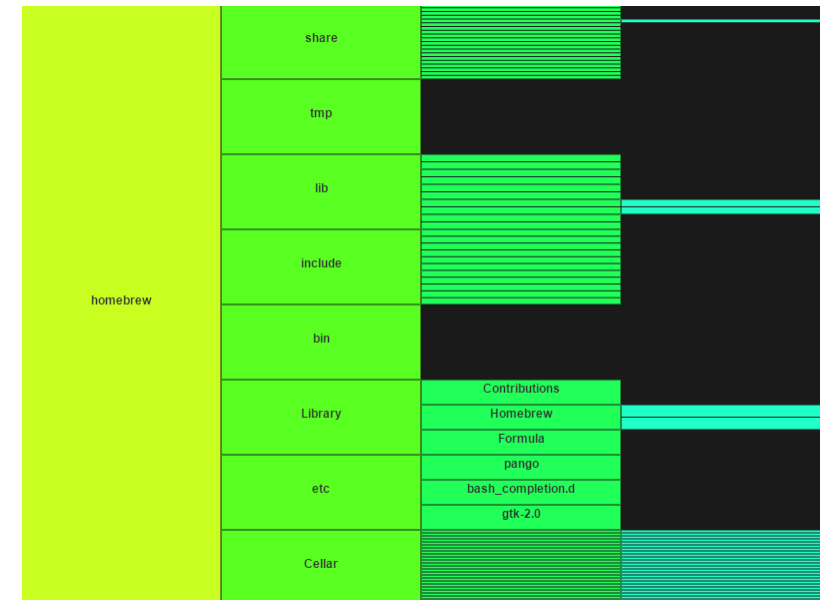
Circle packing

2010 Country Share of World GDP



Country / Region, % of Total Sum of GDP (curr \$) and sum of GDP (curr \$). Color shows details about Region. Size shows sum of GDP (curr \$). The marks are labeled by Country / Region, % of Total Sum of GDP (curr \$) and sum of GDP (curr \$). The data is filtered on Date (Year, which keeps 2010). The view is filtered on sum of GDP (curr \$), which keeps non-NaN values only.

Tree map



Icicle

Interactivity

Zooming & panning

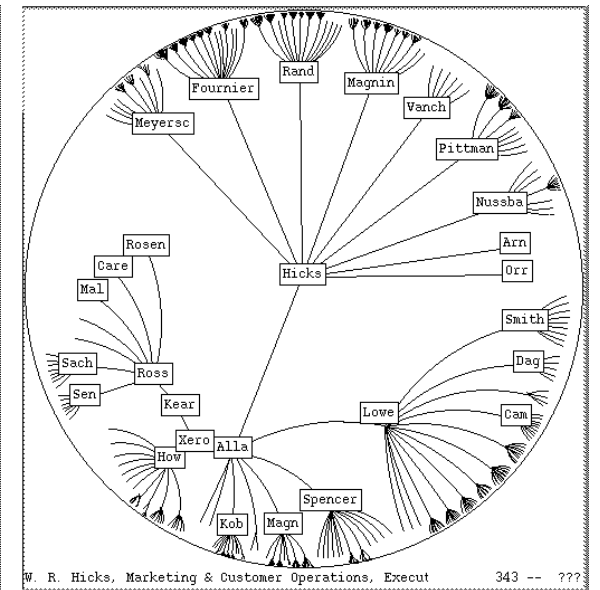
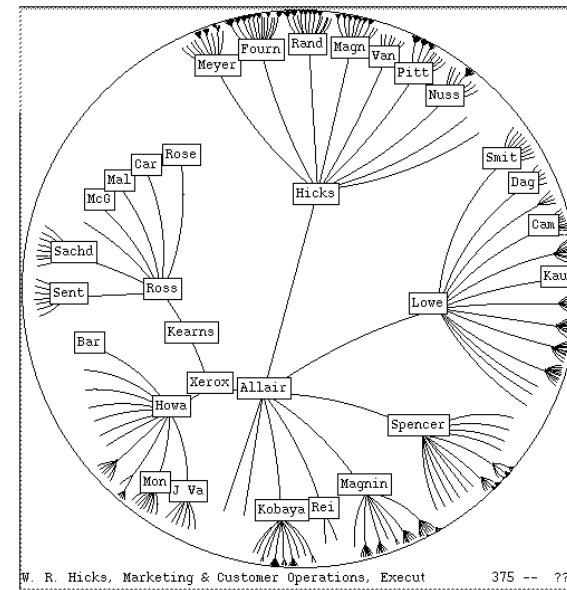
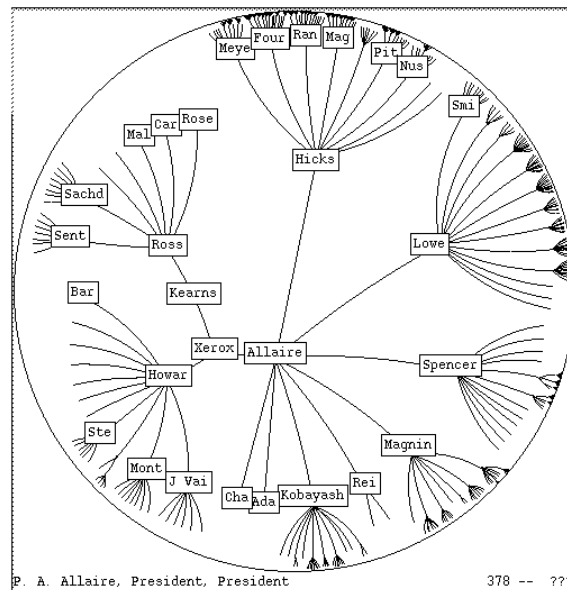
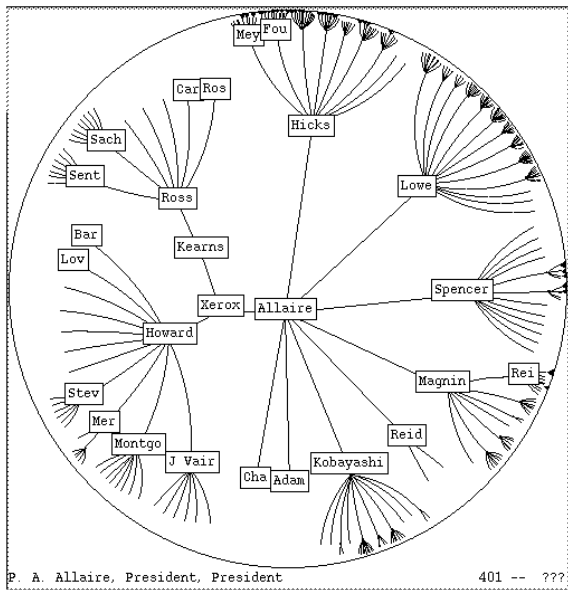
- Not always is the interface problem solvable with the layout algorithm → zoom & pan, i.e. navigation in visualization
- **Zooming is simple for graphs**, which contain only simple geometrical objects (nodes and edges)
 - **Geometric** (standard) – only changes the scale of magnification
 - **Semantic** – modifies what is being shown; either more details, different representation of the data or even different data

Focus and context

- Zooming suffers from **loosing context** – when zoomed in, all context is lost → difficult to pan → **decreased usability**
- A technique **combining** both **overview** (context) and **detail** information (focus) in one view would be desirable → **focus+context** techniques

Fisheye (1)

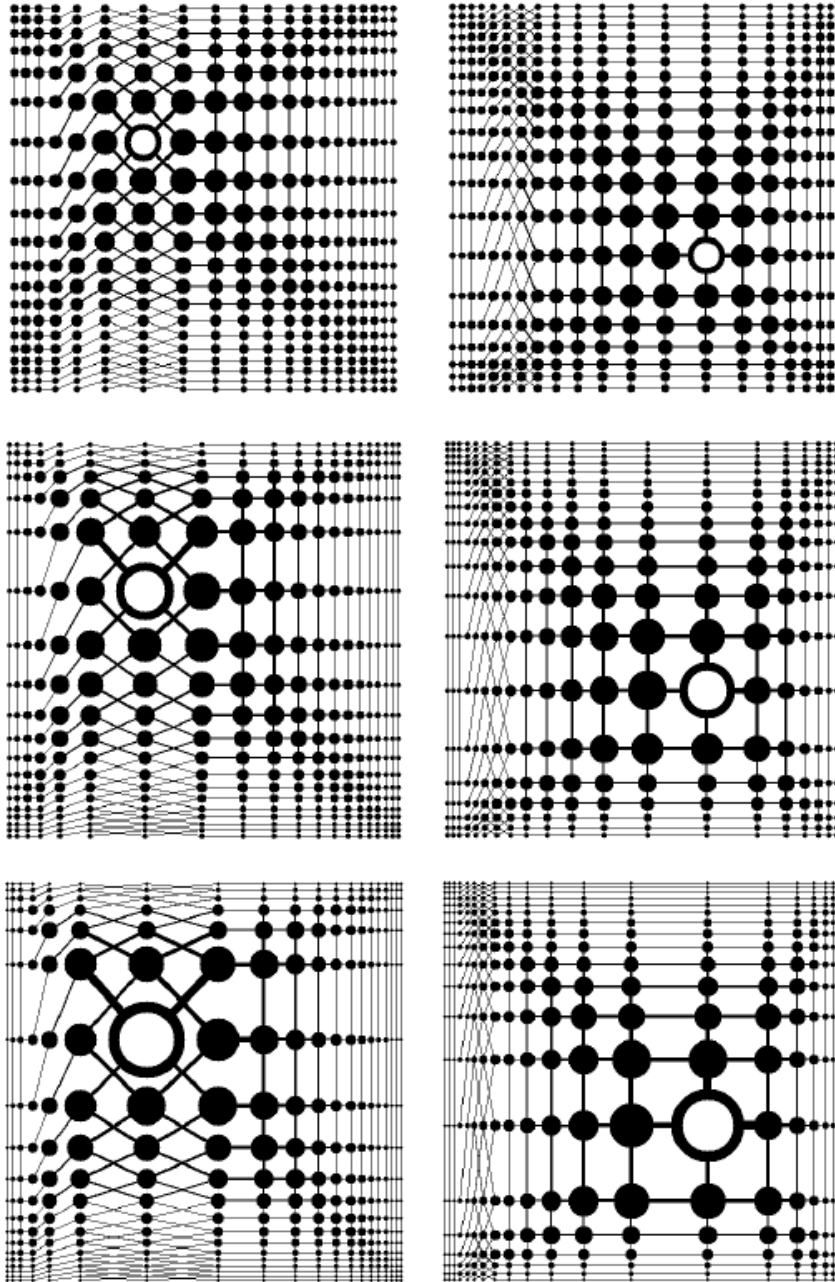
- Popular **distortion** viewing technique that magnifies nearby objects while shrinking distant objects
- With **graphs**, we want to see **details of the specific subgraph** while still seeing the **whole structure**



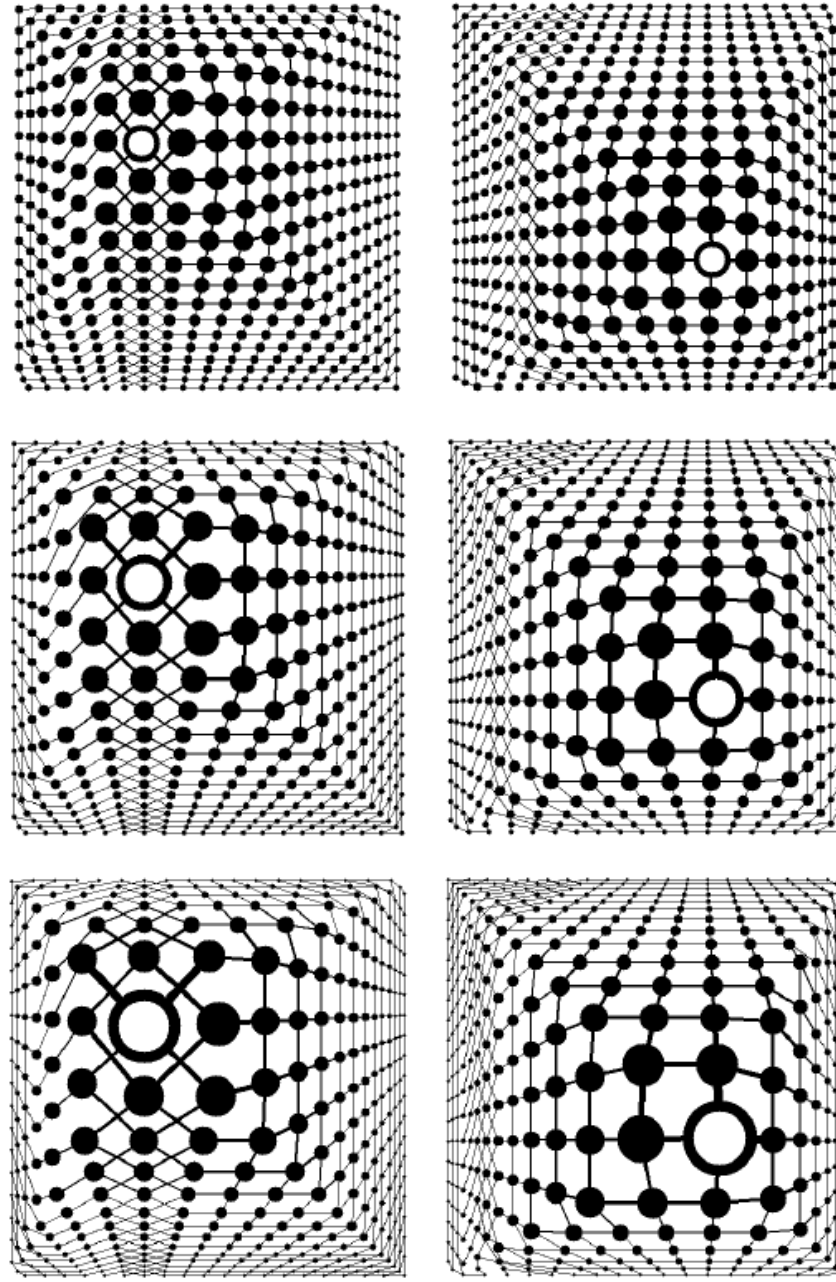
Fisheye (2)

- Distortion technique – user selects a **focus point** and the **layout is distorted**
 - **Polar** transformation – transformation on a polar coordinate system (r, θ) - distortion applied equally in all the directions from the focus points (θ kept unchanged) up to some point
 - suitable for objects such as maps <https://bost.ocks.org/mike/fisheye/>
 - **Cartesian** transformation – scales x and y positions individually, does not require curving the lines
 - better for, e.g., scatter plots

Cartesian distortion



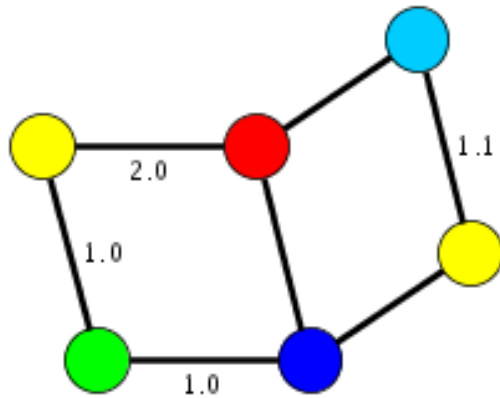
Polar distortion



Tools and data

Graph languages - GraphML

- [XML-based format](#) for exchanging graph structure data
- Stores structural information as well as the graphical information

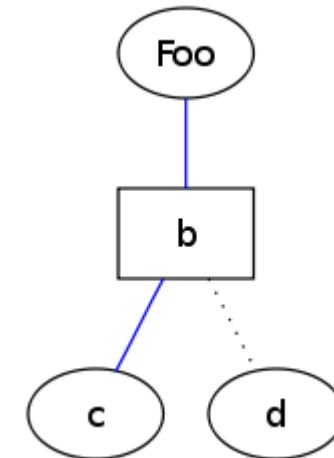


```
<?xml version="1.0" encoding="UTF-8"?>
<graphml xmlns="http://graphml.graphdrawing.org/xmlns"
  xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance"
  xsi:schemaLocation="http://graphml.graphdrawing.org/xmlns
  http://graphml.graphdrawing.org/xmlns/1.0/graphml.xsd">
  <key id="d0" for="node" attr.name="color" attr.type="string">
    <default>yellow</default>
  </key>
  <key id="d1" for="edge" attr.name="weight" attr.type="double"/>
  <graph id="G" edgedefault="undirected">
    <node id="n0">
      <data key="d0">green</data>
    </node>
    <node id="n1"/>
    <node id="n2">
      <data key="d0">blue</data>
    </node>
    <node id="n3">
      <data key="d0">red</data>
    </node>
    <node id="n4"/>
    <node id="n5">
      <data key="d0">turquoise</data>
    </node>
    <edge id="e0" source="n0" target="n2">
      <data key="d1">1.0</data>
    </edge>
    <edge id="e1" source="n0" target="n1">
      <data key="d1">1.0</data>
    </edge>
    <edge id="e2" source="n1" target="n3">
      <data key="d1">2.0</data>
    </edge>
    <edge id="e3" source="n3" target="n2"/>
    <edge id="e4" source="n2" target="n4"/>
    <edge id="e5" source="n3" target="n5"/>
    <edge id="e6" source="n5" target="n4">
      <data key="d1">1.1</data>
    </edge>
  </graph>
</graphml>
```

Graph languages - DOT

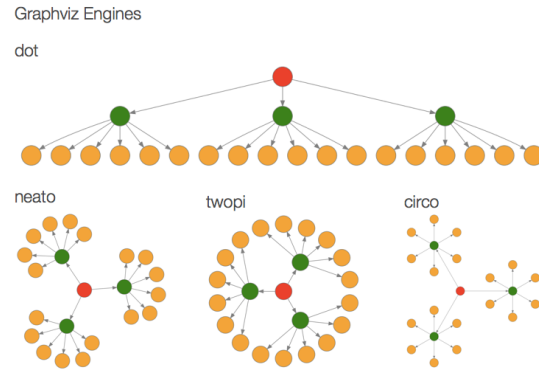
- Plain text graph description language
- Supported by GraphViz, Gephi, ..

```
graph graphname {  
    // This attribute applies to the graph itself  
    size="1,1";  
    // The label attribute can be used to change the  
label of a node  
    a [label="Foo"];  
    // Here, the node shape is changed.  
    b [shape=box];  
    // These edges both have different line properties  
    a -- b -- c [color=blue];  
    b -- d [style=dotted];  
}
```



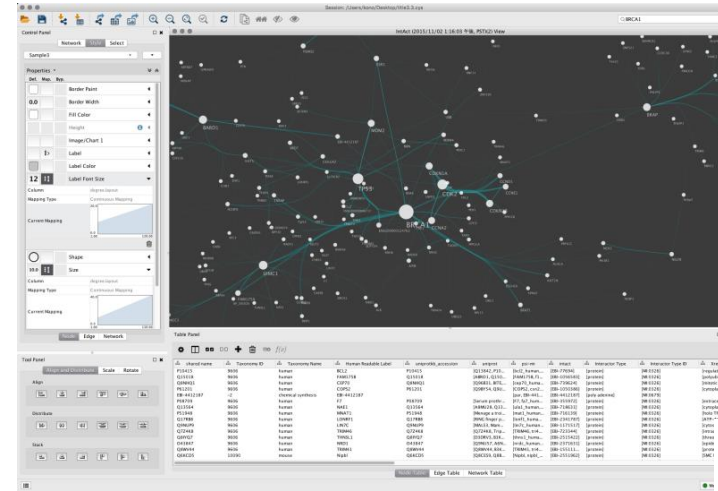
Software

- GraphViz

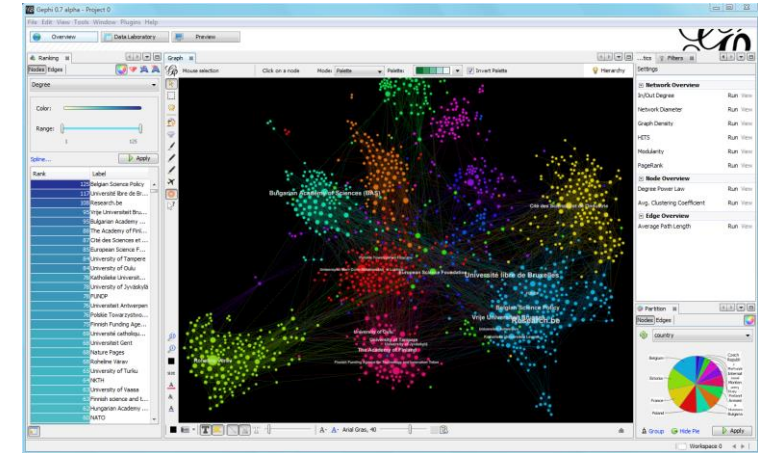


source: http://rich-iannone.github.io/DiagrammeR/graphviz_and_mermaid.html

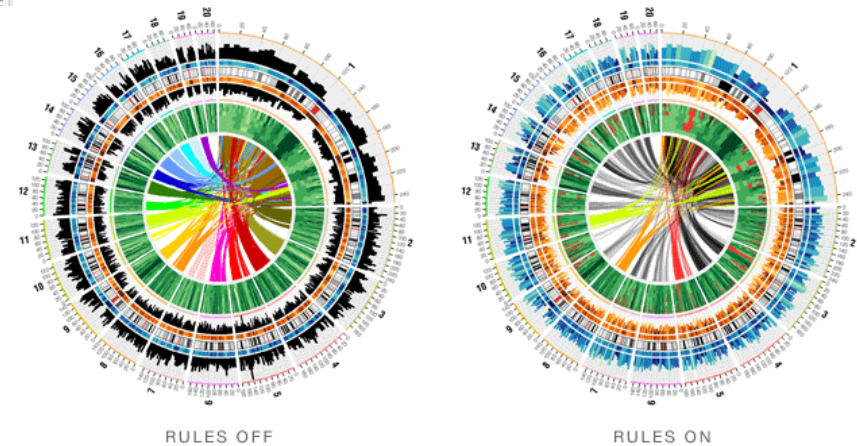
- Gephi



- Cytoscape
 - Cytoscape.js



- Circos



Network data sets collections

- [Stanford Large Network Dataset Collection](#)
- [Pajek data sets](#)
- R package [igraphdata](#)

Sources

- Roberto Tamassia (2013) Handbook of Graph Drawing and Visualization. Chapman and Hall/CRC
- Chaomei Chen (2006) Information Visualization: Beyond the Horizon. Springer