Assoc.Prof. Dr. Andreas Holzinger

185.A83 Machine Learning for Health Informatics
2020S, VU, 2.0 h, 3.0 ECTS
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Lecture 04 - Week 17

From Decision Making under Uncertainty to Probabilistic Graphical Models

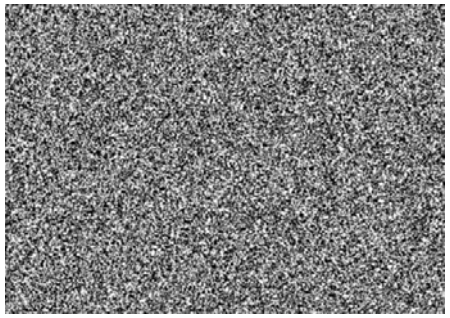
Contact: andreas.holzinger AT tuwien.ac.at

https://human-centered.ai/machine-learning-for-health-informatics-class-2020

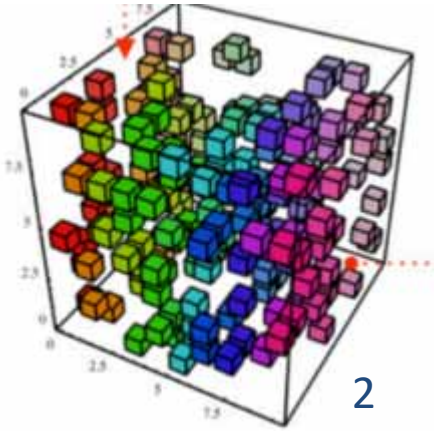
- 00 Reflection from last lecture
- 01 Decision Making under uncertainty
- 02 Some Basics of Graphs/Networks
- 03 Bayesian Networks (BN)
- 04 Markov Chain Monte Carlo (MCMC)
- 05 Metropolis Hastings Algorithm (MH)
- 06 Probabilistic Programming (PP)

00 Reflection

Warm-up Quiz



1



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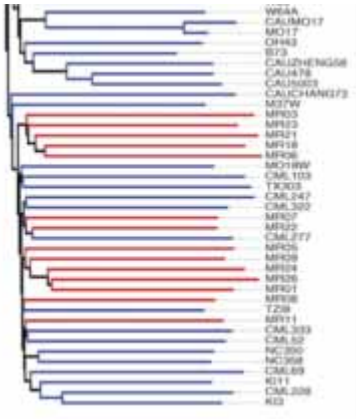

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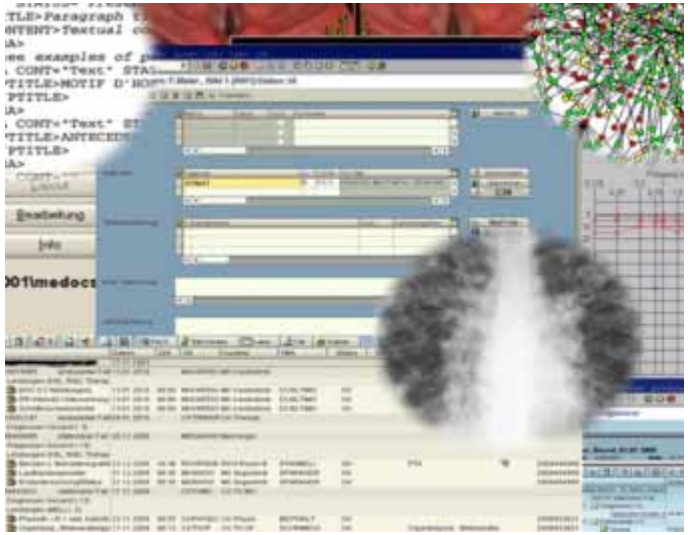

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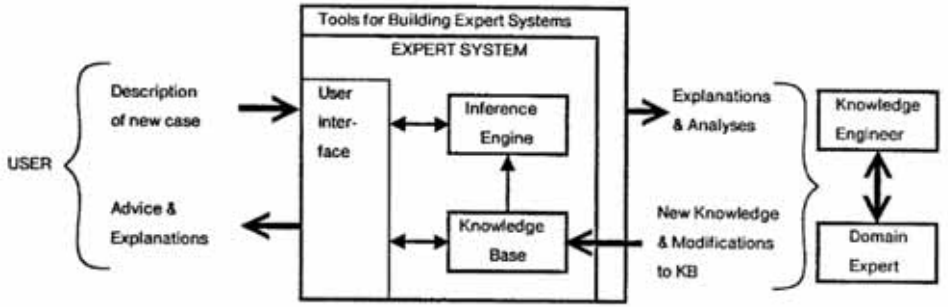


5

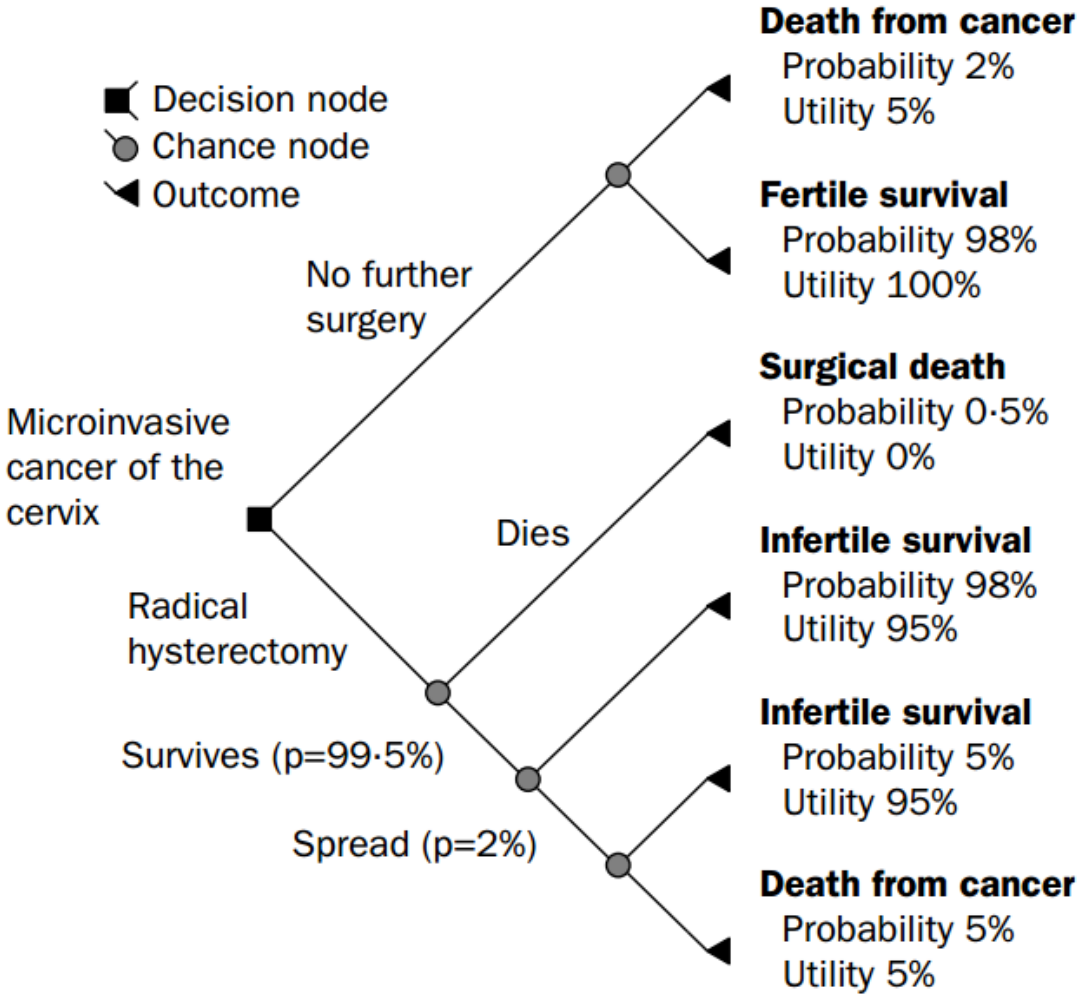
Which ML approaches do we know ?

- Symbolic ML
  - First order logic, inverse deduction, knowledge composition
  - Tom Mitchell, Steve Muggleton, Ross Quinlan, …
- Bayesian ML
  - Statistical learning, probabilistic inference
  - Judea Pearl, Michael Jordan, David Heckermann, …
- Cognitive ML
  - Analogisms from Psychology, Kernel machines
  - Vladimir Vapnik, Peter Hart, Douglas Hofstaedter, …
- Connectionist ML
  - Neuroscience, Backpropagation
  - Geoffrey Hinton, Yoshua Bengio, Yann LeCun, …
- Evolutionary ML
  - Nature-inspired concepts, genetic programming
  - John Holland (1929-2015), John Koza, Hod Lipson, …

Pedro Domingos 2015. *The Master Algorithm: How the Quest for the Ultimate Learning Machine Will Remake Our World*, Penguin UK.

https://learning.acm.org/techtalks/machinelearning

Why is this important for us?



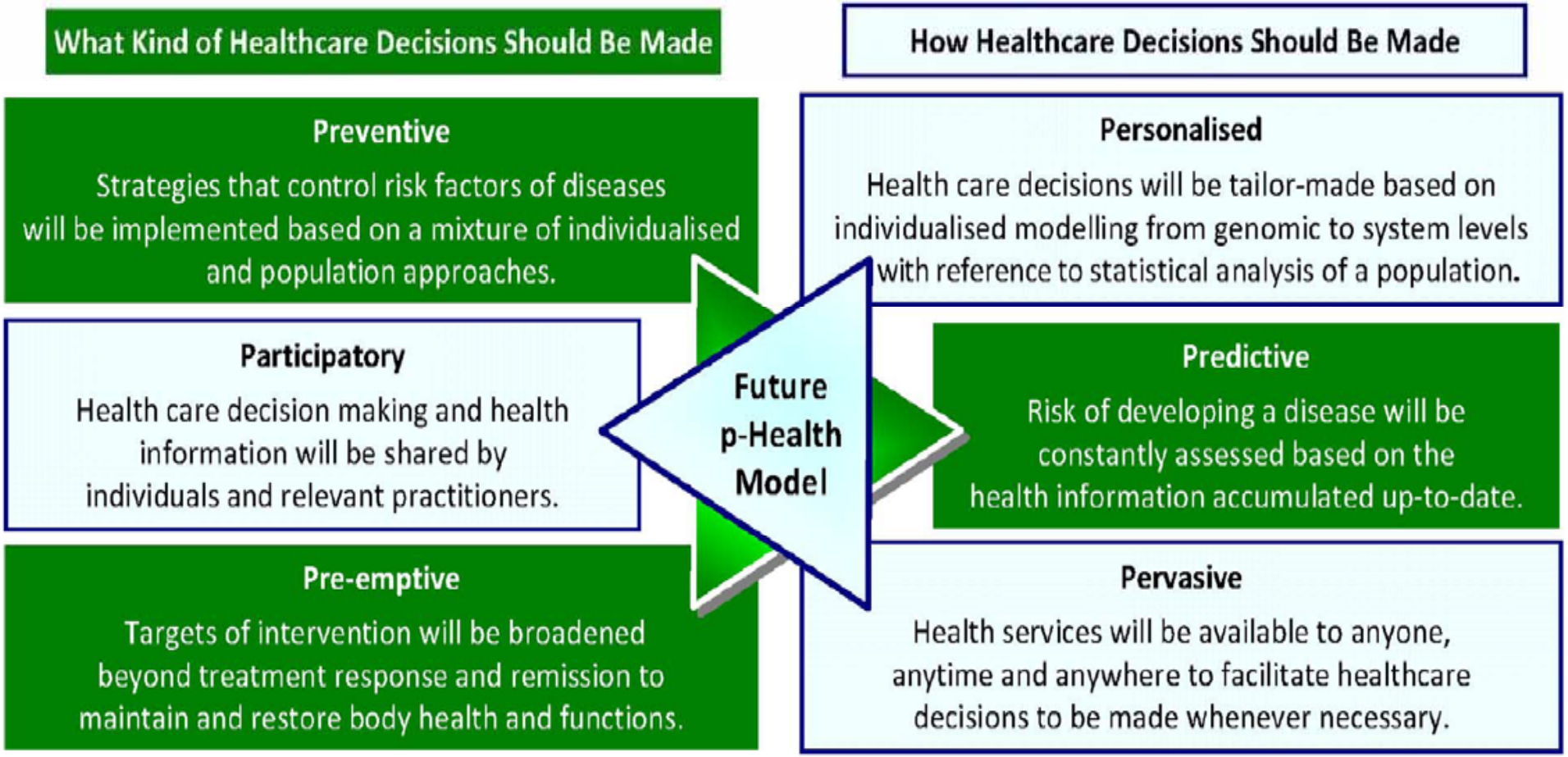
Physician treating a patient
approx. 480 B.C.
Beazley (1963), Attic Red-figured Vase-Painters, 813, 96.
Department of Greek, Etruscan and Roman Antiquities, Sully, 1st floor, Campana Gallery, room 43
Louvre, Paris

Elwyn, G., Edwards, A., Eccles, M. & Rovner, D. 2001. Decision analysis in patient care. The Lancet, 358, (9281), 571-574.

A short historical digression: whom do you see here?

Both Images are in the public domain

Remember: What is “personalized medicine” ?



Zhang, Y. T. & Poon, C. C. Y. (2010) Editorial Note on Bio, Medical, and Health Informatics. *Information Technology in Biomedicine, IEEE Transactions on, 14, 3, 543-545.*

01 Decision Making under uncertainty

Pierre-Simon Laplace 1781. Mémoire sur les probabilités. *Mémoires de l'Académie Royale des sciences de Paris*, 1778, 227-332.

How to go from System 1 deep learning to System 2 deep learning?

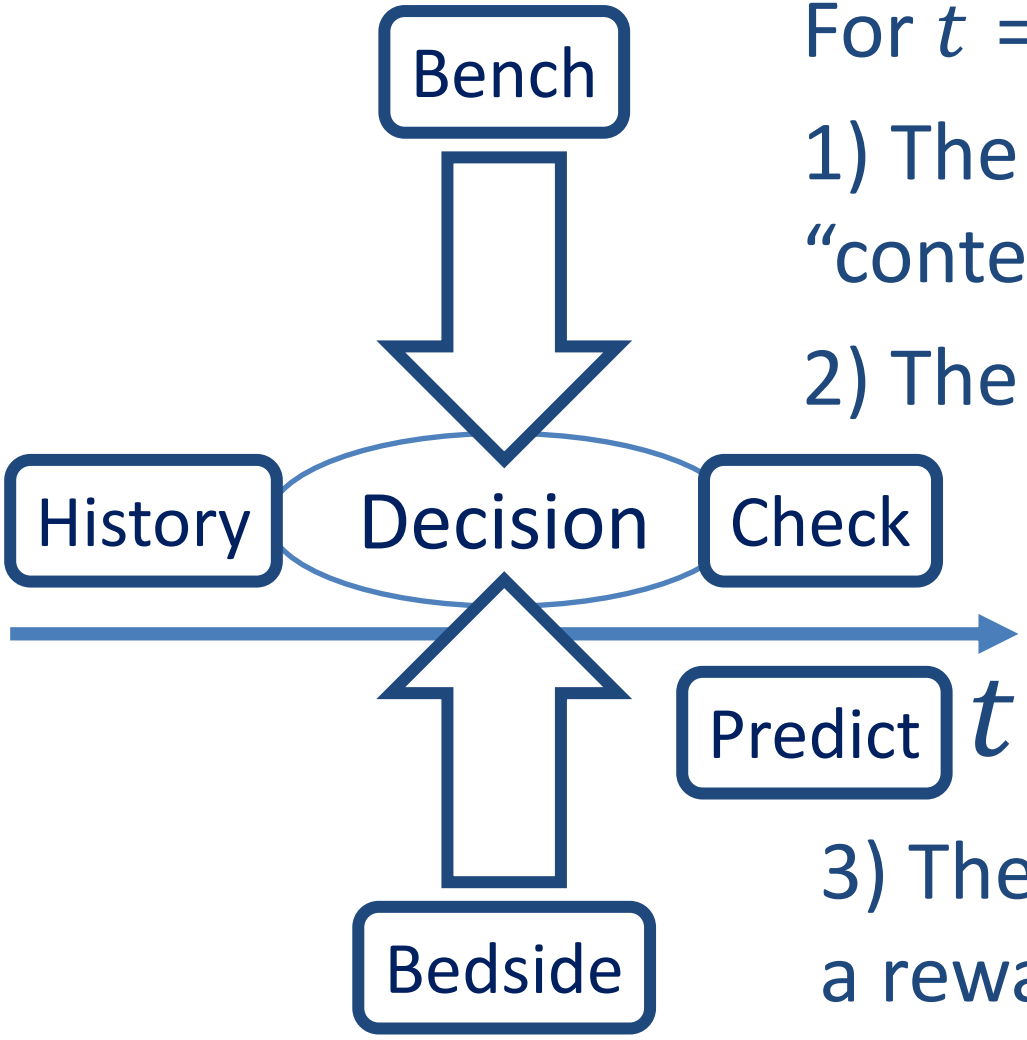
https://www.youtube.com/watch?v=T3sxeTgT4qc

Daniel Kahneman 2011. Thinking, fast and slow, New York, Macmillan.

Amos Tversky & Daniel Kahneman 1974. Judgment under uncertainty: Heuristics and biases. Science, 185, (4157), 1124-1131, doi:10.1126/science.185.4157.1124.

... permanent decision making
under uncertainty!

Goal: Learn an **optimal policy** for selecting best actions within a given **context**



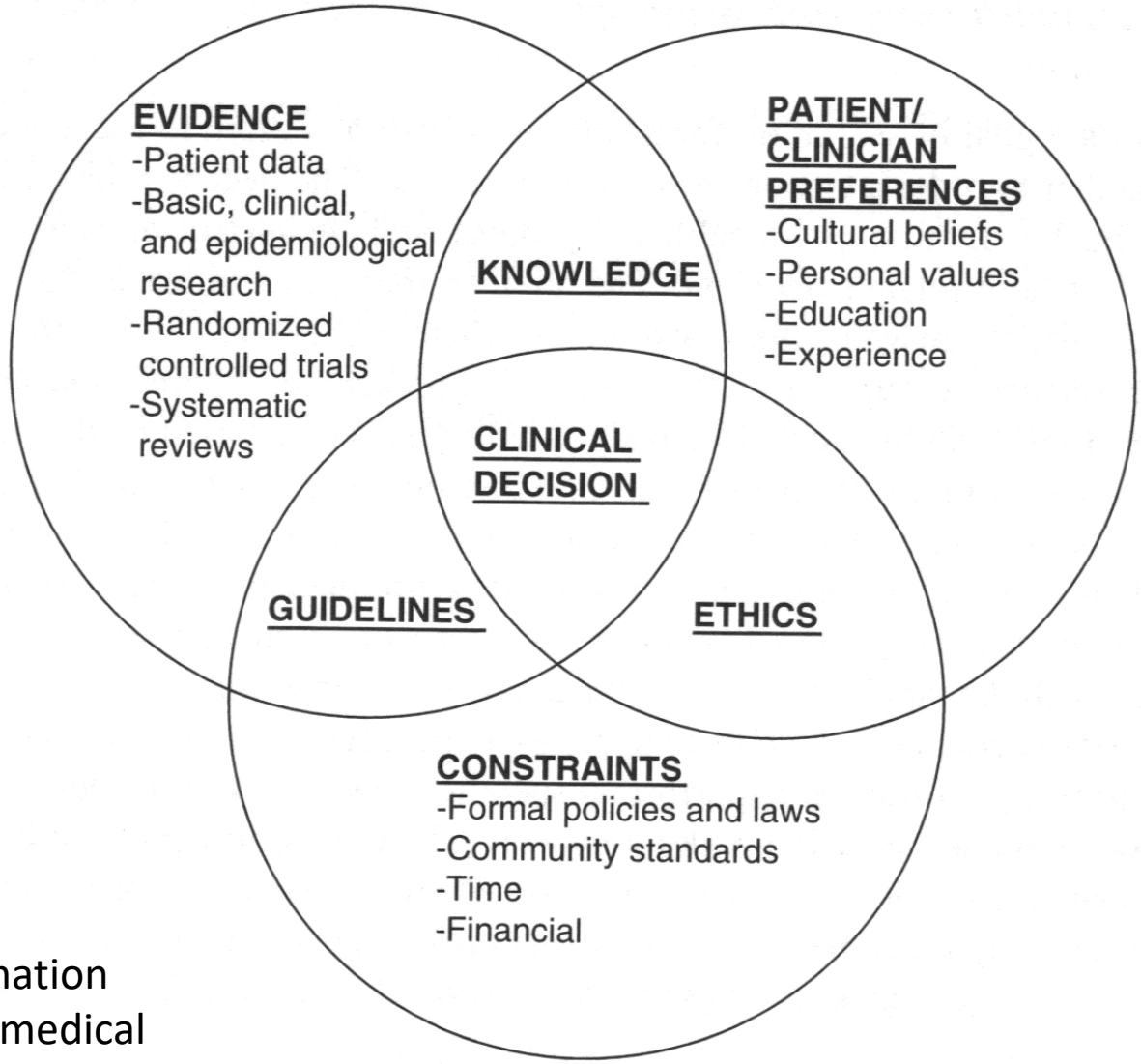
For $t = 1, \ldots, T$

1) The world produces a "context" $x_t \in X$

2) The learner selects an action $a_t \in \{1, \ldots, K\}$

3) The world reacts with a reward $r_t(a_t) \in [0,1]$

Why is decision making so difficult ?



William Hersh 2010. Information Retrieval: A Health and Biomedical Perspective, New York, Springer.

How does a medical doctor make a decision ?

3 July 1959, Volume 130, Number 3366

SCIENCE

Reasoning Foundations of Medical Diagnosis

Symbolic logic, probability, and value theory aid our understanding of how physicians reason.

Robert S. Ledley and Lee B. Lusted

The purpose of this article is to analyze the complicated reasoning processes inherent in medical diagnosis. The importance of this problem has received recent emphasis by the increasing interest in the use of electronic computers as an aid to medical diagnostic processes

fitted into a definite disease category, or that it may be one of several possible diseases, or else that its exact nature cannot be determined." This, obviously, is a greatly simplified explanation of the process of diagnosis, for the physician might also comment that after seeing a

ance are the ones who do remember and consider the most possibilities."

Computers are especially suited to help the physician collect and process clinical information and remind him of diagnoses which he may have overlooked. In many cases computers may be as simple as a set of hand-sorted cards, whereas in other cases the use of a large-scale digital electronic computer may be indicated. There are other ways in which computers may serve the physician, and some of these are suggested in this paper. For example, medical students might find the computer an important aid in learning the methods of differential diagnosis. But to use the computer thus we must understand how the physician makes a medical diagnosis. This, then, brings us to the subject of our investigation: the reasoning foundations of medical diagnosis and treatment.

Medical diagnosis involves processes that can be systematically analyzed, as well as those characterized as "intangible." For instance, the reasoning foundations of medical diagnostic procedures

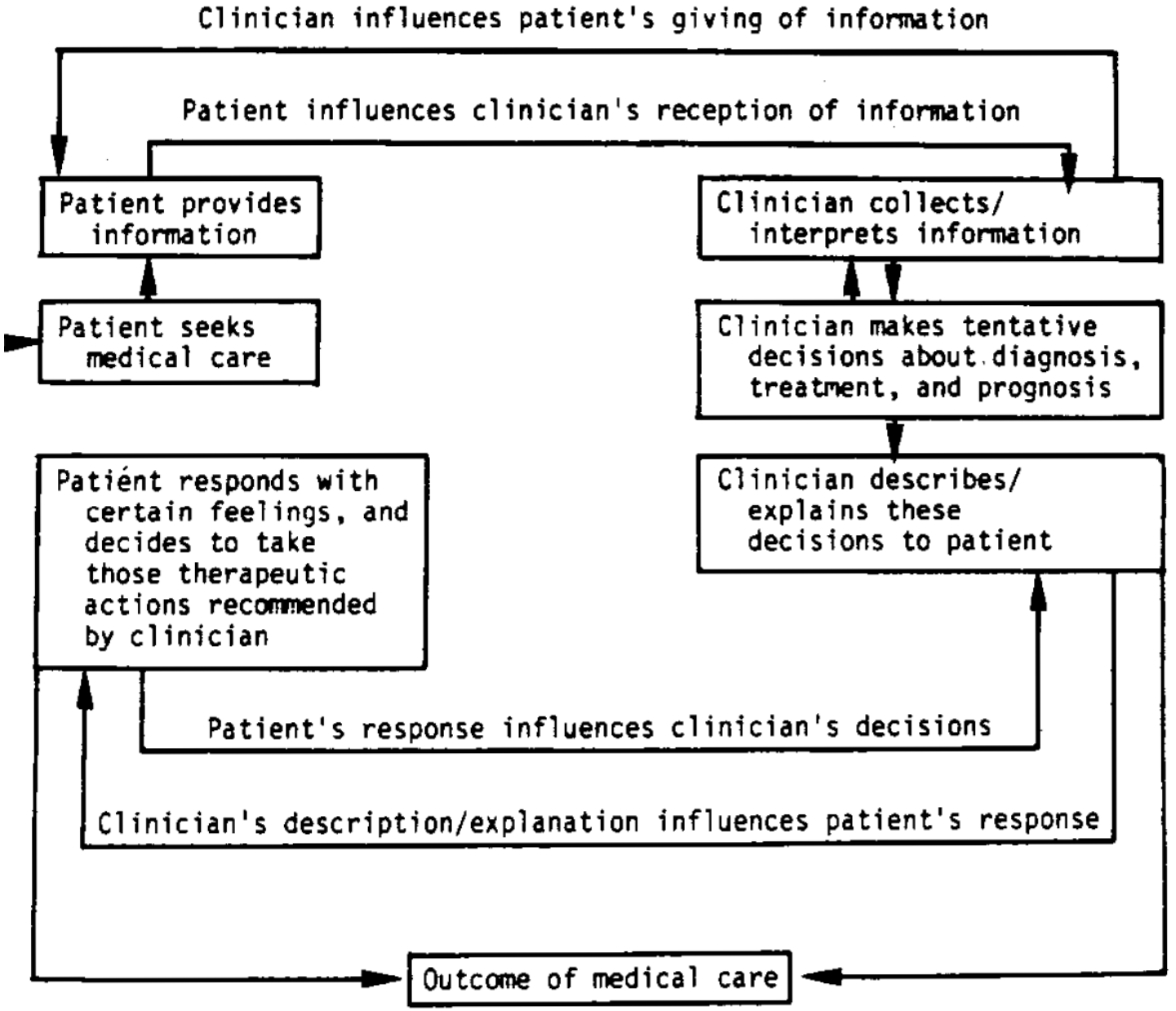
What about the accuracy and uncertainty in medicine ?

- Medical (clinical) data are defined and detected disturbingly “soft” …
- … having an obvious degree of **variability** and **inaccuracy**.
- Taking a medical history, the performance of a physical examination, the interpretation of laboratory tests, even the definition of diseases … are surprisingly **inexact**.
- Data is defined, collected, and interpreted with a degree of variability and inaccuracy which falls far short of the standards **which engineers do expect from most data**.
- Moreover, standards might be **interpreted variably** by different medical doctors, different hospitals, different medical schools, different medical cultures, …

Anthony L. Komaroff 1979. The variability and inaccuracy of medical data. Proceedings of the IEEE, 67, (9), 1196-1207.

Why is the patient-doctor dialogue so important ?

Anthony L. Komaroff 1979. The variability and inaccuracy of medical data. Proceedings of the IEEE, 67, (9), 1196-1207.



d … *data*

h … *hypotheses*

$\mathcal{H}$ … $\{h_1, h_2, \ldots, h_n\}$

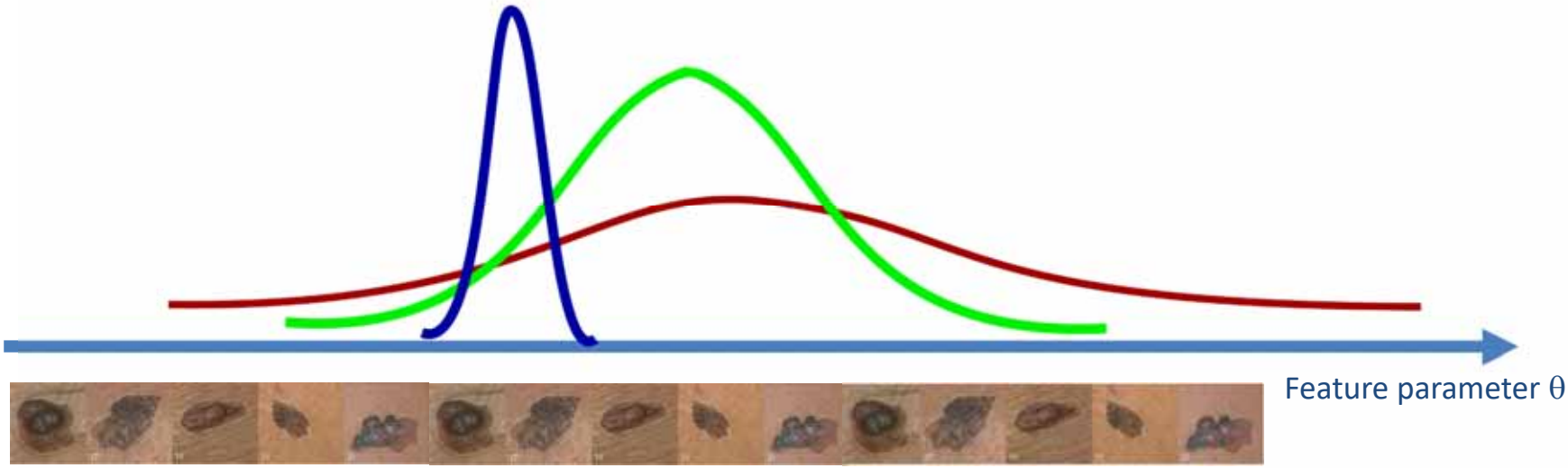
Likelihood

Prior Probability

$$p(h|d) = \frac{p(d|h) * p(h)}{\sum_{h \in H} p(d|h')\, p(h')}$$

Posterior Probability

Problem in $\mathbb{R}^n \rightarrow$ complex



How do humans make decisions under uncertainty ?

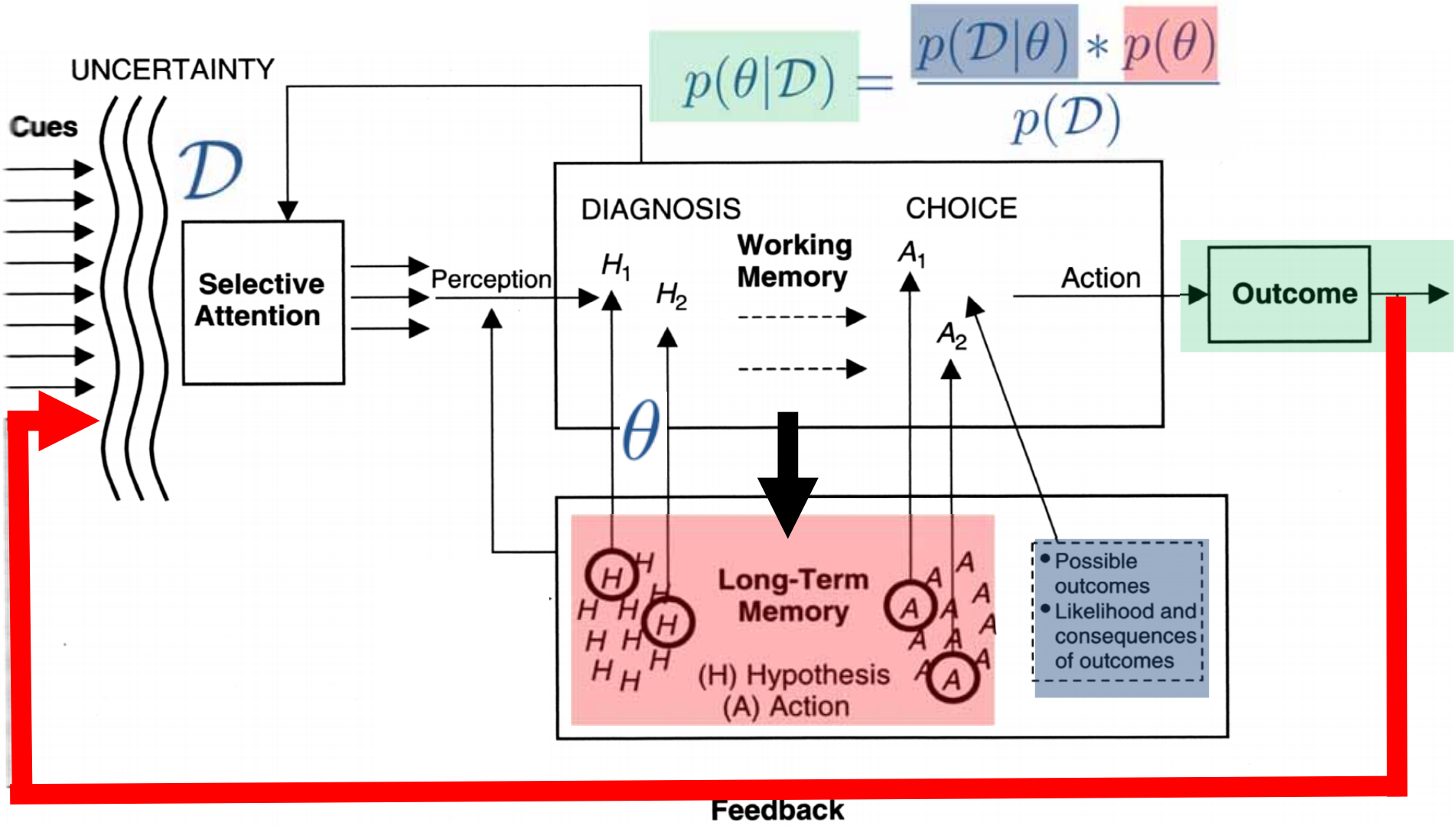
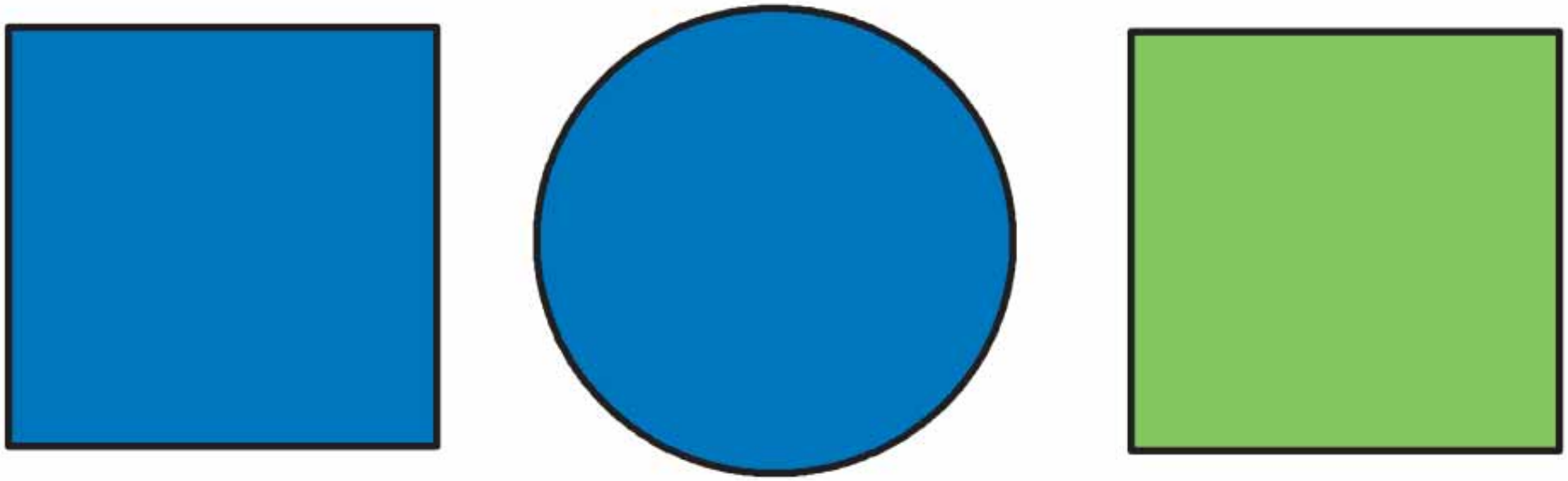


Image by Christopher D. Wickens 1984, modified by Andreas Holzinger 2004

You are talking to you colleague and want to refer to the middle object – which wording would you prefer: circle or blue?



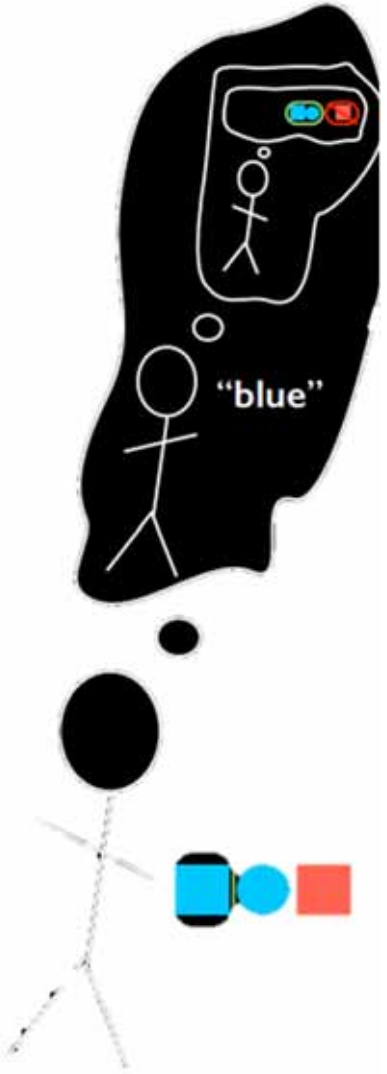
Michael C. Frank & Noah D. Goodman 2012. Predicting pragmatic reasoning in language games. *Science*, 336, (6084), 998-998, doi:10.1126/science.1218633.

Recursive reasoning: a case for probabilistic programming

```
var literalListener = function(property){
 Infer(function(){
    var object = refPrior(context)
    condition(object[property])
    return object
```

```
var speaker = function(object) {
  Infer(function(){
    var property = propPrior()
    condition(
       object ==
```

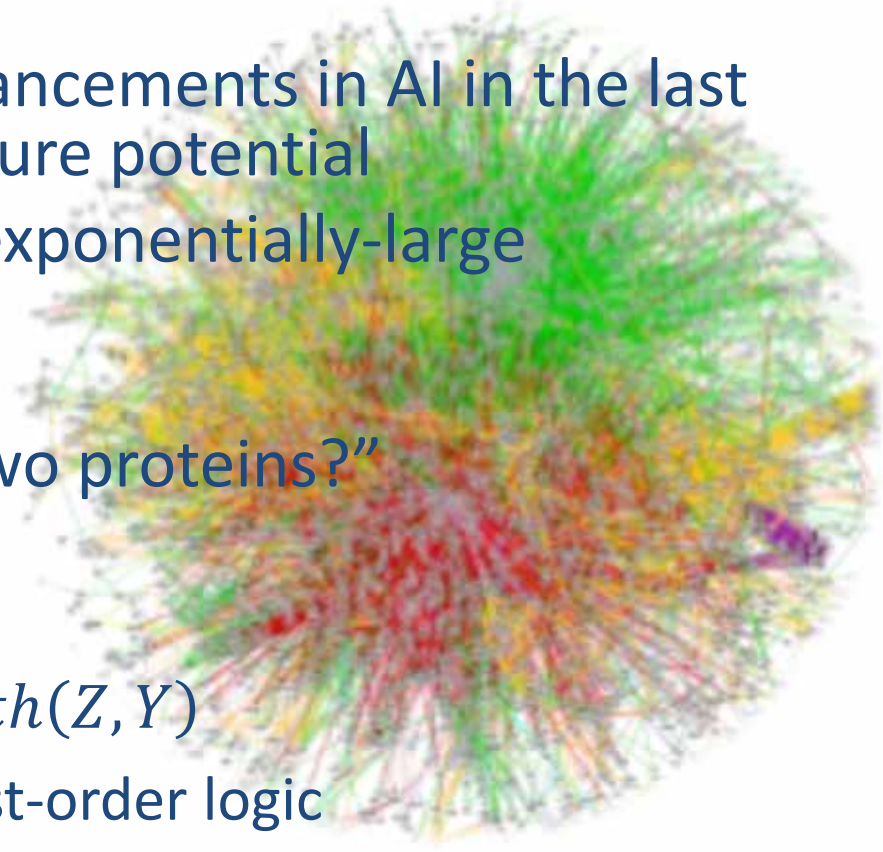
```
var listener = function(property) {
  Infer(function(){
    var object = refPrior(context)
    condition(utterance ==
                sample(speaker(object)))
    return object
  })}
```



Noah D. Goodman & Michael C. Frank 2016. Pragmatic language interpretation as probabilistic inference. *Trends in Cognitive Sciences*, 20, (11), 818-829, doi:10.1016/j.tics.2016.08.005.

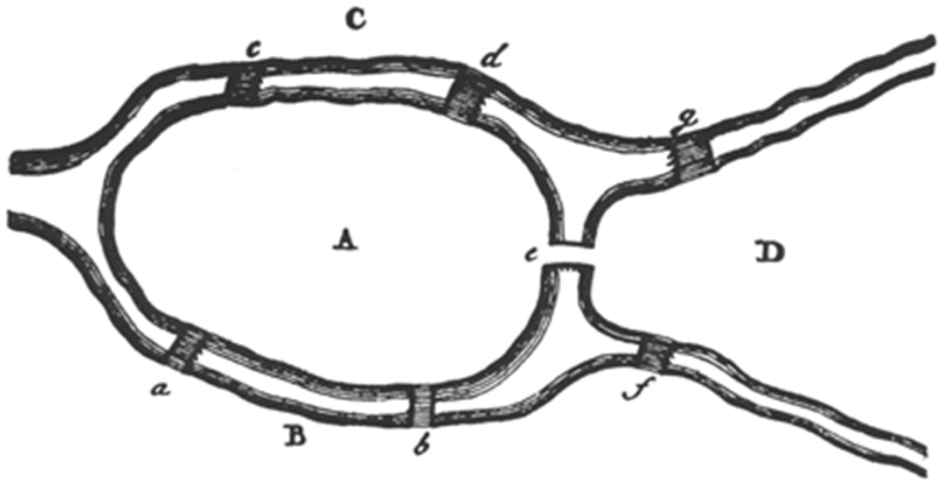
02 Graphs = Networks

- PGM can be seen as a combination between
- **Graph Theory + Probability Theory + Machine Learning**
- One of the most exciting advancements in AI in the last decades – with enormous future potential
- Compact representation for exponentially-large probability distributions
- Example Question:
  "Is there a path connecting two proteins?"



- $Path\ (X,Y) :=\ edge\ (X,Y)$
- $Path\ (X,Y) :=\ edge\ (X,Y), path(Z,Y)$
- This can NOT be expressed in first-order logic

We start in 1736



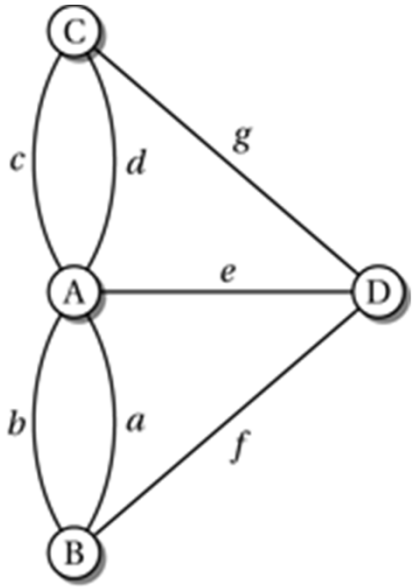
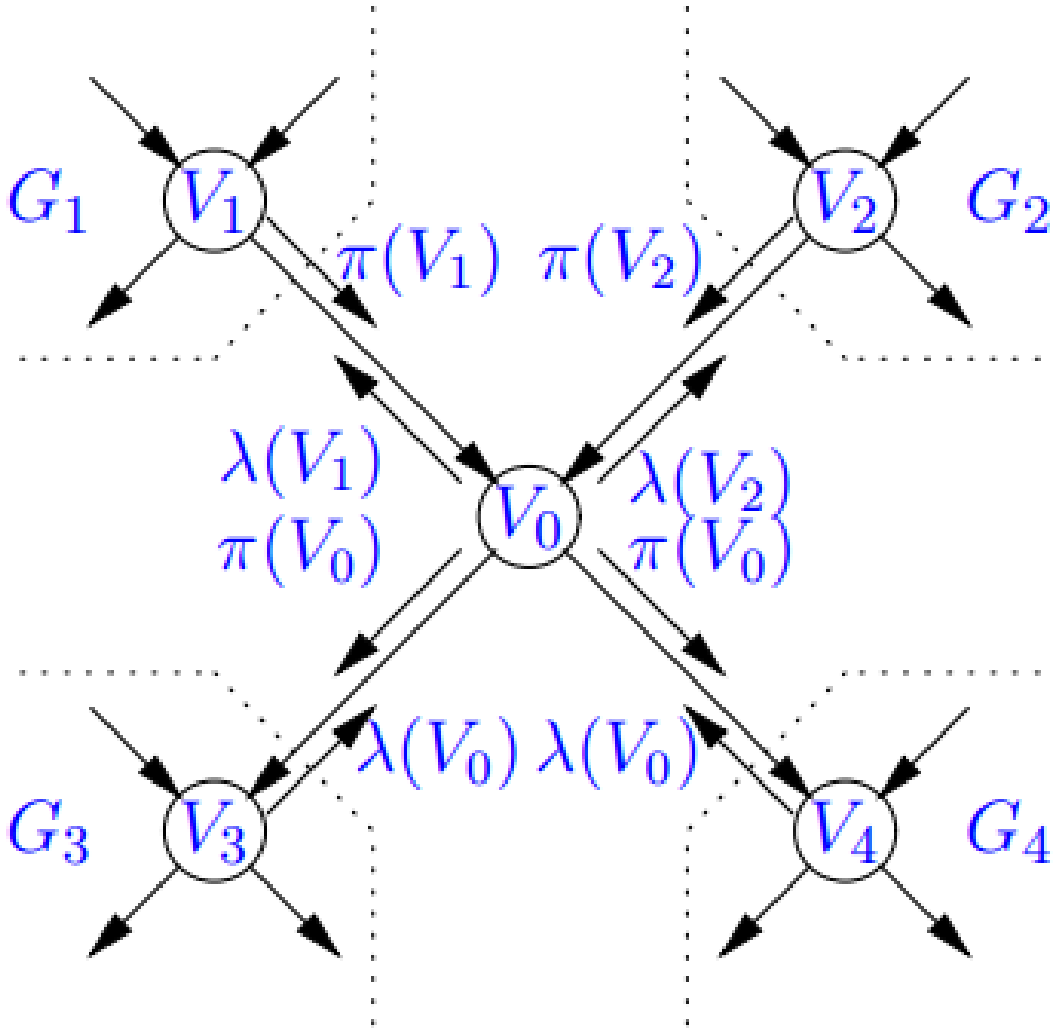


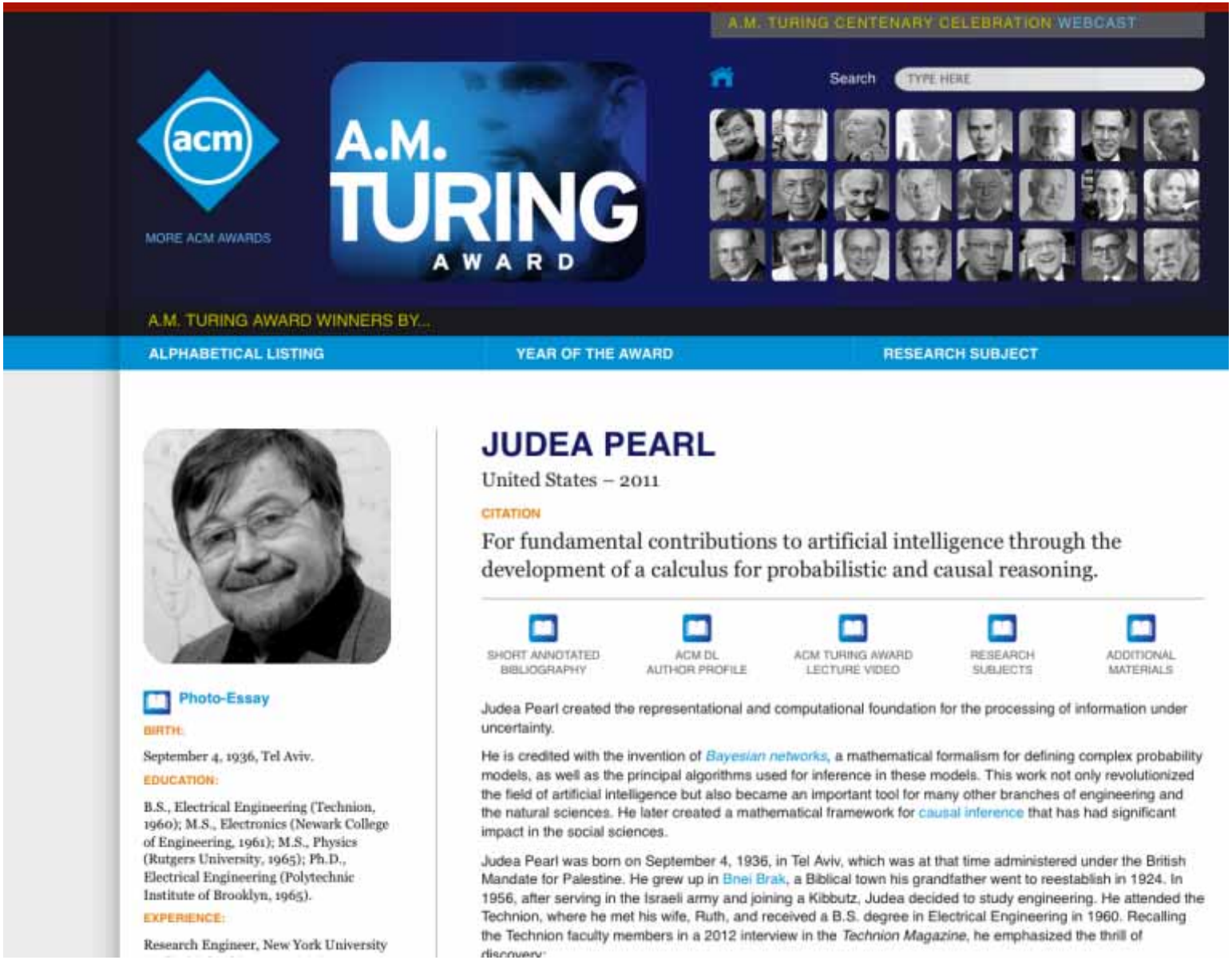
Image from https://people.kth.se/~carlofi/teaching/FEL3250-2013/courseinfo.html

Leonhard Euler 1741. Solutio problematis ad geometriam situs pertinentis. Commentarii academiae scientiarum Petropolitanae, 8, 128-140.



Pearl, J. 1988. Embracing causality in default reasoning. Artificial Intelligence, 35, (2), 259-271.

275 years later … the “Nobel-prize in Computer Science”



http://amturing.acm.org/vp/pearl_2658896.cfm

Nobel Prize in Chemistry 2013

Scientific Background on the Nobel Prize in Chemistry 2013

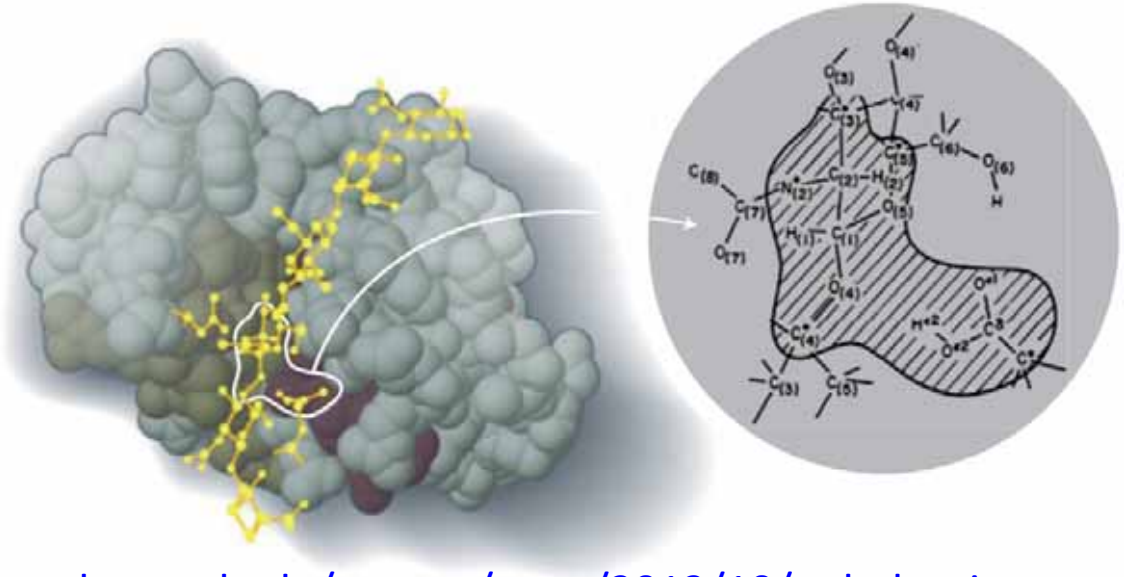
DEVELOPMENT OF MULTISCALE MODELS FOR COMPLEX CHEMICAL SYSTEMS

Photo: A. Mahmoud
Martin Karplus
Prize share: 1/3

Photo: A. Mahmoud
Michael Levitt
Prize share: 1/3

Photo: A. Mahmoud
Arieh Warshel
Prize share: 1/3

http://www.nobelprize.org/nobel_prizes/chemistry/laureates/2013

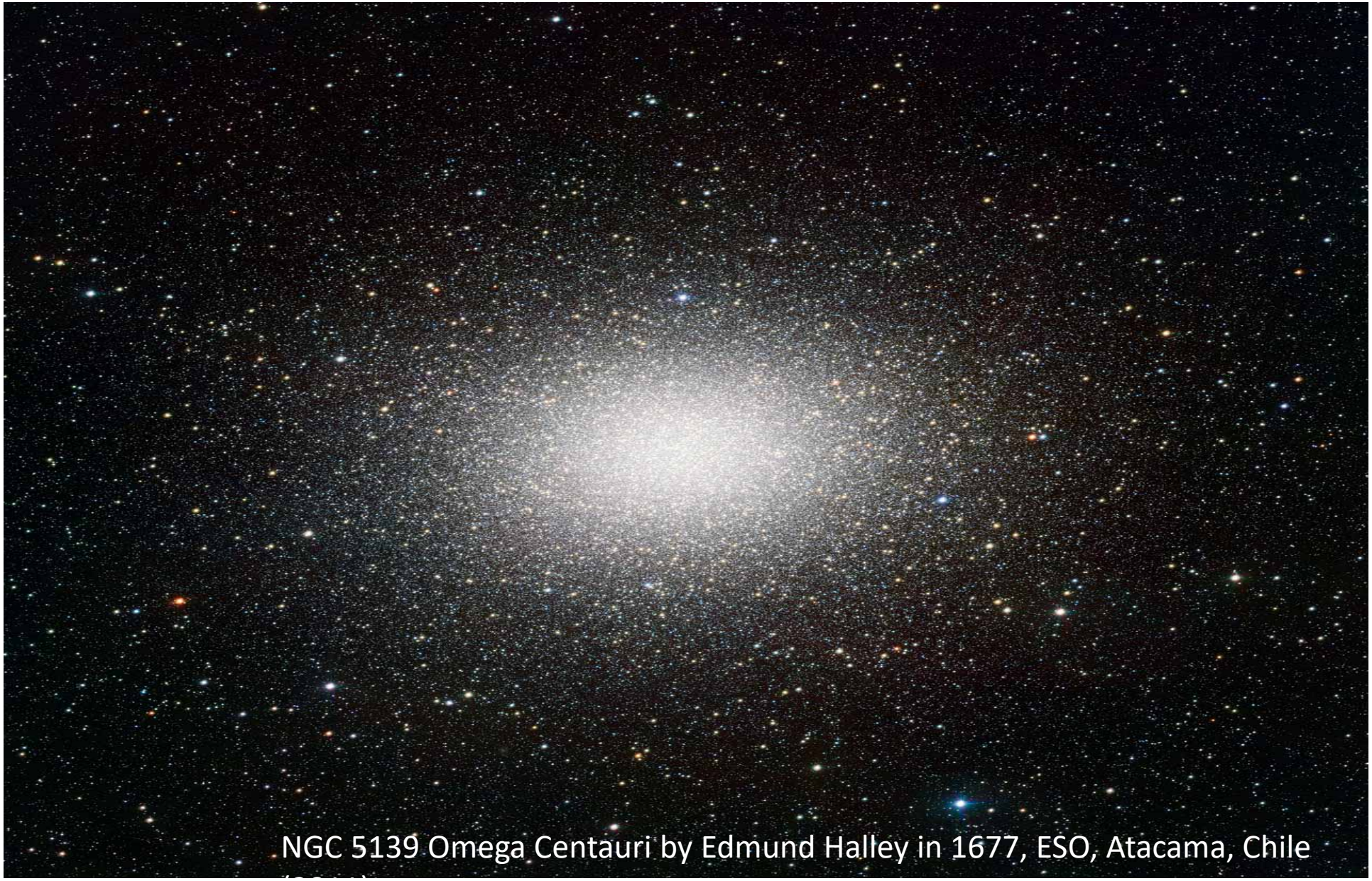


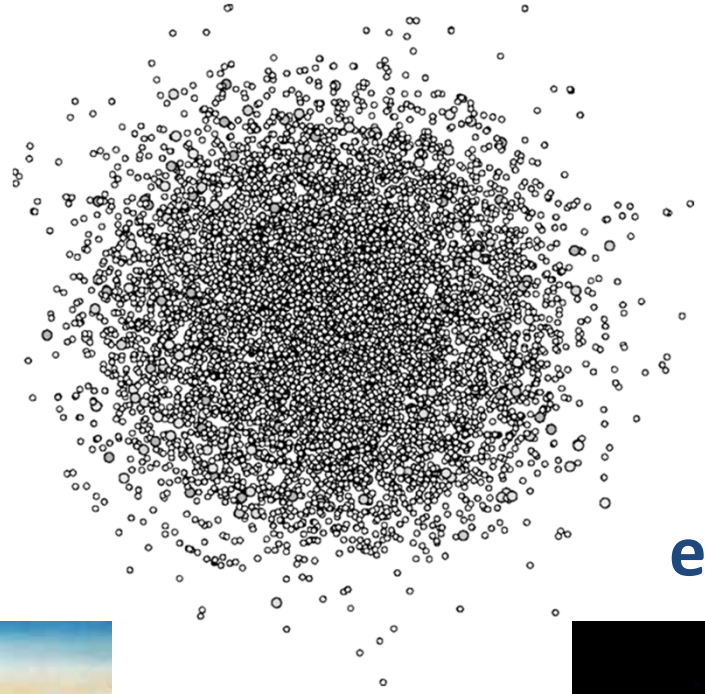
http://news.harvard.edu/gazette/story/2013/10/nobel_prize_awarded_2013

- **Graphs as models for networks**
- given as direct input (point cloud data sets)
- Given as properties of a structure
- Given as a representation of information (e.g. Facebook data, viral marketing, etc., …)

- **Graphs as nonparametric basis**
- we learn the structure from samples and infer
- flat vector data, e.g. similarity graphs
- encoding structural properties (e.g. smoothness, independence, …)

What do you see here ?





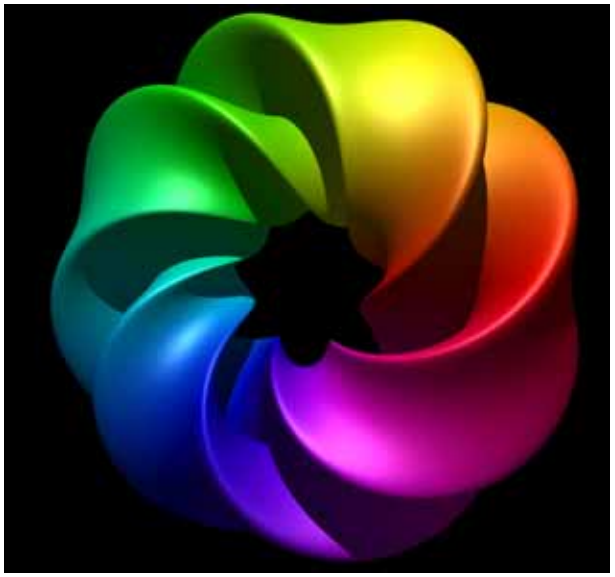
Time

e.g. Entropy

Dali, S. (1931) The persistence of memory

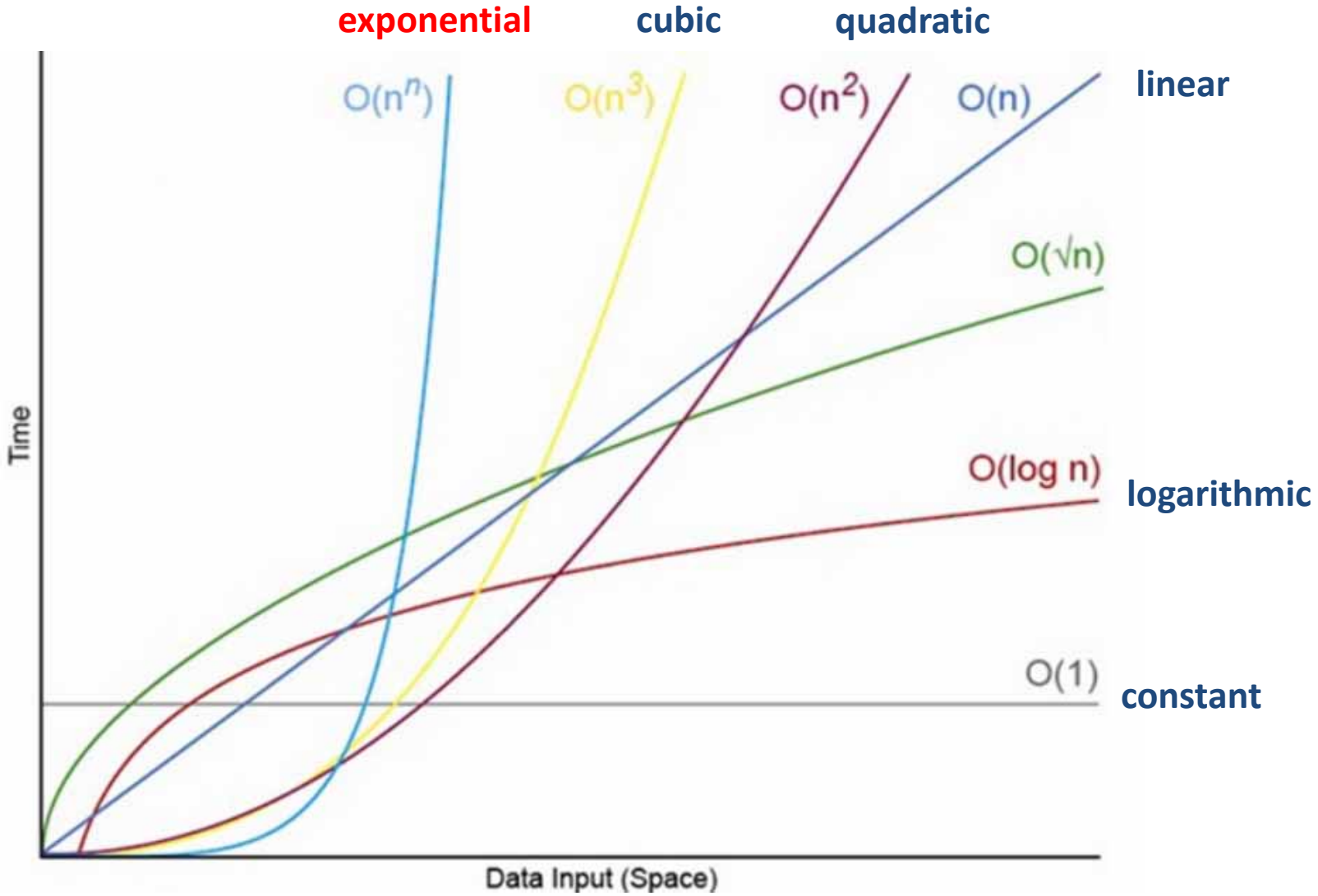
Space

e.g. Topology



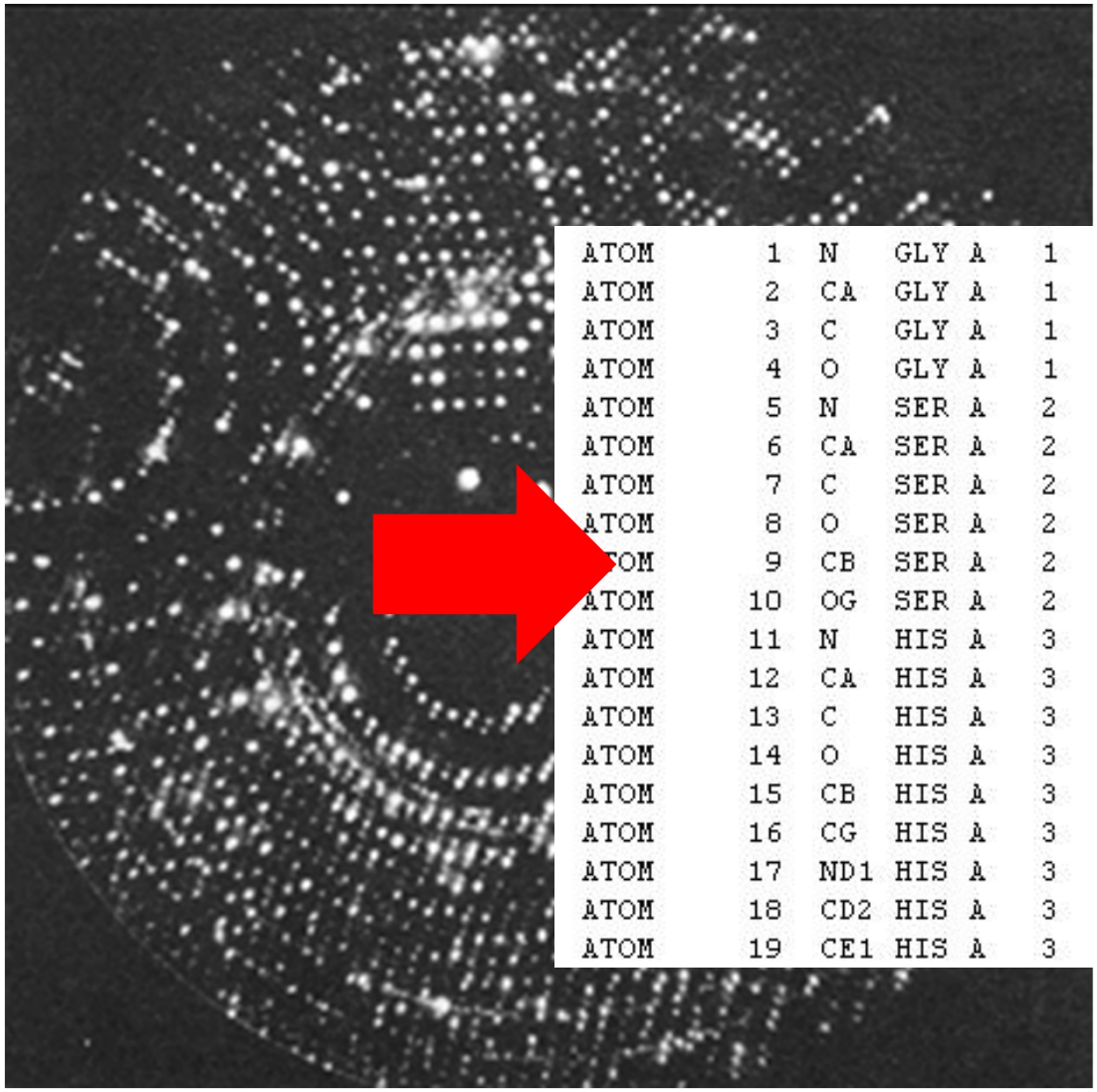
Bagula & Bourke (2012) Klein-Bottle

Complexity Problem: Time versus Space



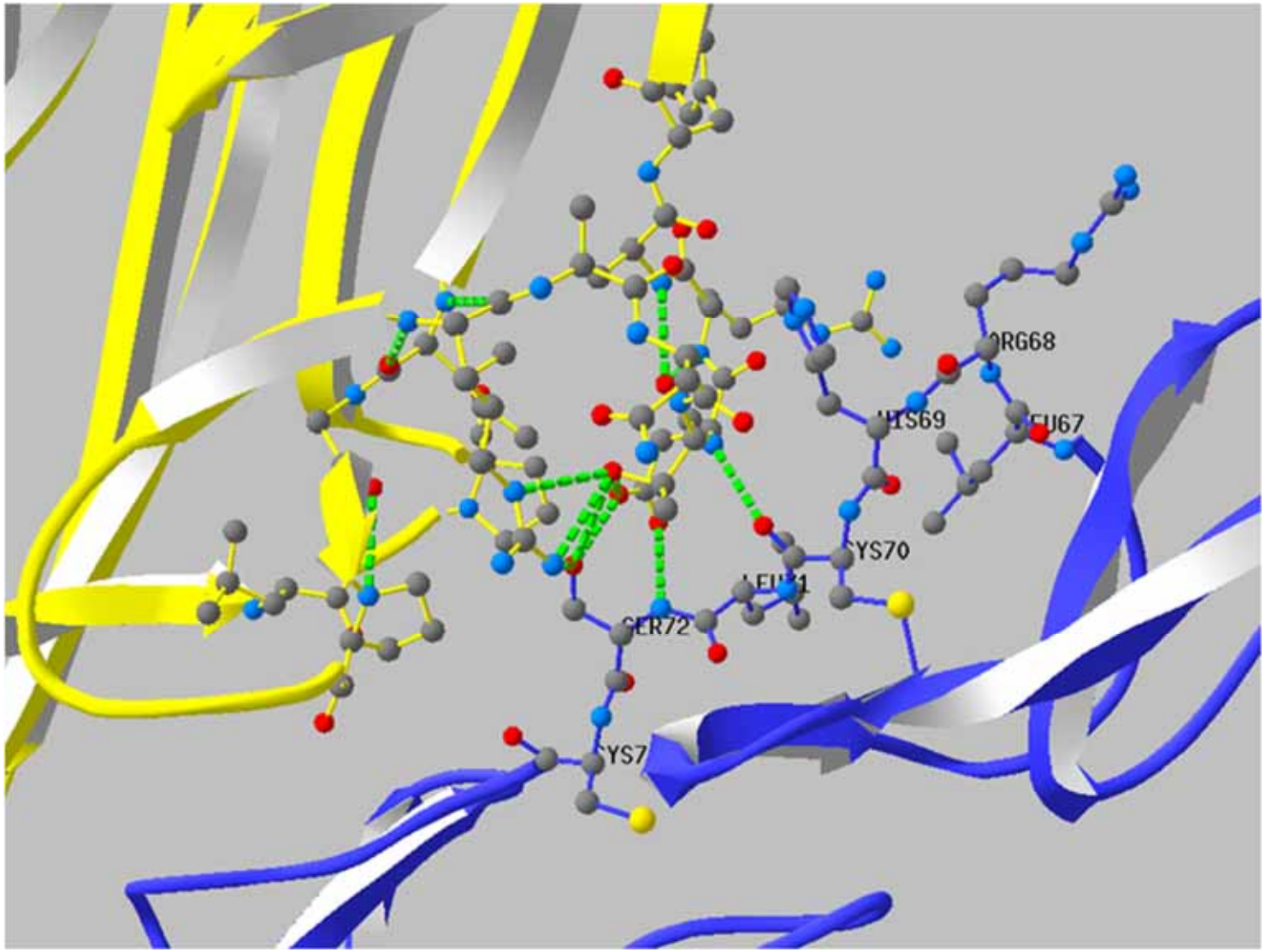
P versus NP and the Computational Complexity Zoo, please have a look at https://www.youtube.com/watch?v=YX40hbAHx3s

Why are protein structures so important ?



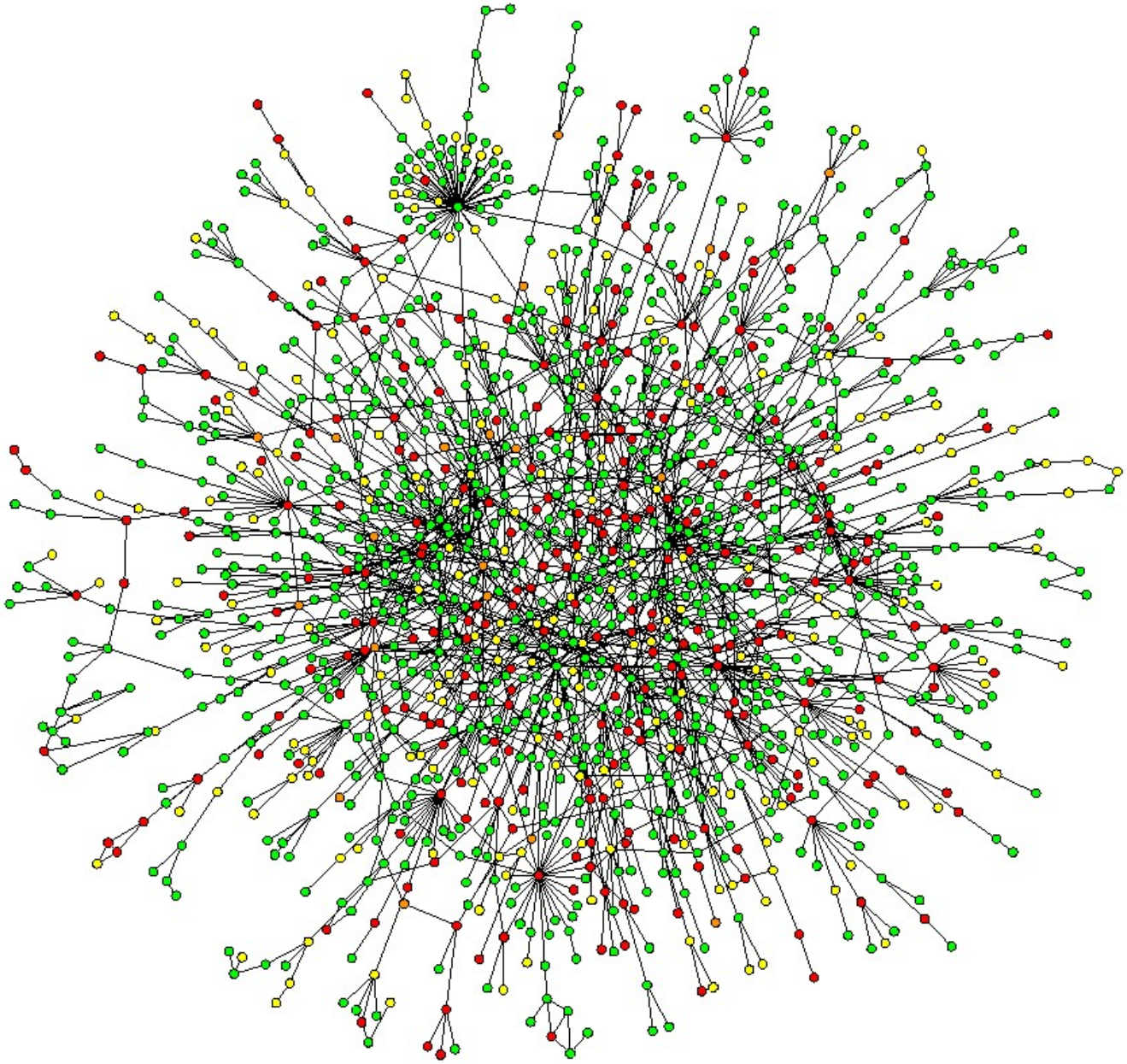
```
ATOM      1  N   GLY A   1      44.842  51.034 101.284  0.01 27.20
ATOM      2  CA  GLY A   1      45.640  50.230 100.389  0.01 26.99
ATOM      3  C   GLY A   1      46.692  49.648 101.308  0.01 26.80
ATOM      4  O   GLY A   1      46.895  50.222 102.381  0.01 26.91
ATOM      5  N   SER A   2      47.283  48.516 100.951  1.00 26.26
ATOM      6  CA  SER A   2      48.277  47.866 101.761  1.00 26.17
ATOM      7  C   SER A   2      49.212  47.031 100.845  1.00 24.21
ATOM      8  O   SER A   2      49.060  47.195  99.630  1.00 19.77
ATOM      9  CB  SER A   2      47.438  47.091 102.800  1.00 26.31
ATOM     10  OG  SER A   2      46.276  46.356 102.404  1.00 27.99
ATOM     11  N   HIS A   3      50.147  46.186 101.370  1.00 23.93
ATOM     12  CA  HIS A   3      51.129  45.389 100.609  1.00 21.44
ATOM     13  C   HIS A   3      50.953  43.905 100.849  1.00 20.32
ATOM     14  O   HIS A   3      50.530  43.595 101.950  1.00 22.00
ATOM     15  CB  HIS A   3      52.555  45.674 100.990  1.00 19.69
ATOM     16  CG  HIS A   3      52.940  47.090 100.611  1.00 21.44
ATOM     17  ND1 HIS A   3      53.371  47.470  99.422  1.00 20.87
ATOM     18  CD2 HIS A   3      52.956  48.175 101.433  1.00 21.69
ATOM     19  CE1 HIS A   3      53.676  48.730  99.476  1.00 20.57
```

Wiltgen, M. & Holzinger, A. (2005) Visualization in Bioinformatics: Protein Structures with Physicochemical and Biological Annotations. In: *Central European Multimedia and Virtual Reality Conference. Prague, Czech Technical University (CTU), 69-74*



Wiltgen, M., Holzinger, A. & Tilz, G. P. (2007) Interactive Analysis and Visualization of Macromolecular Interfaces Between Proteins. In: *Lecture Notes in Computer Science (LNCS 4799). Berlin, Heidelberg, New York, Springer, 199-212.*

First yeast protein-protein interaction network

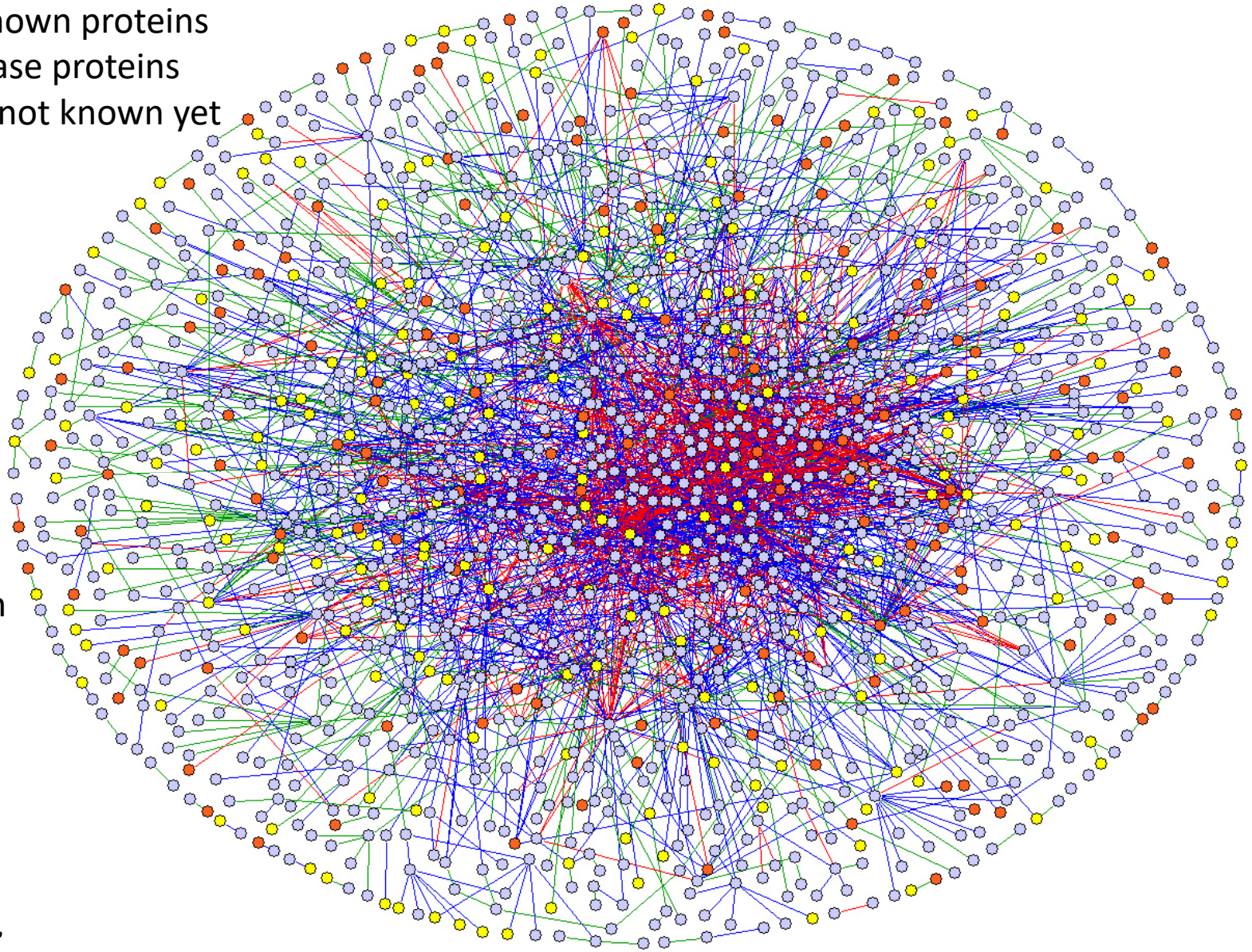


Nodes = proteins
Links = physical interactions (bindings)
Red Nodes = lethal
Green Nodes = non-lethal
Orange = slow growth
Yellow = not known

Jeong, H., Mason, S. P., Barabasi, A. L. & Oltvai, Z. N. (2001) Lethality and centrality in protein networks. *Nature, 411, 6833, 41-42.*

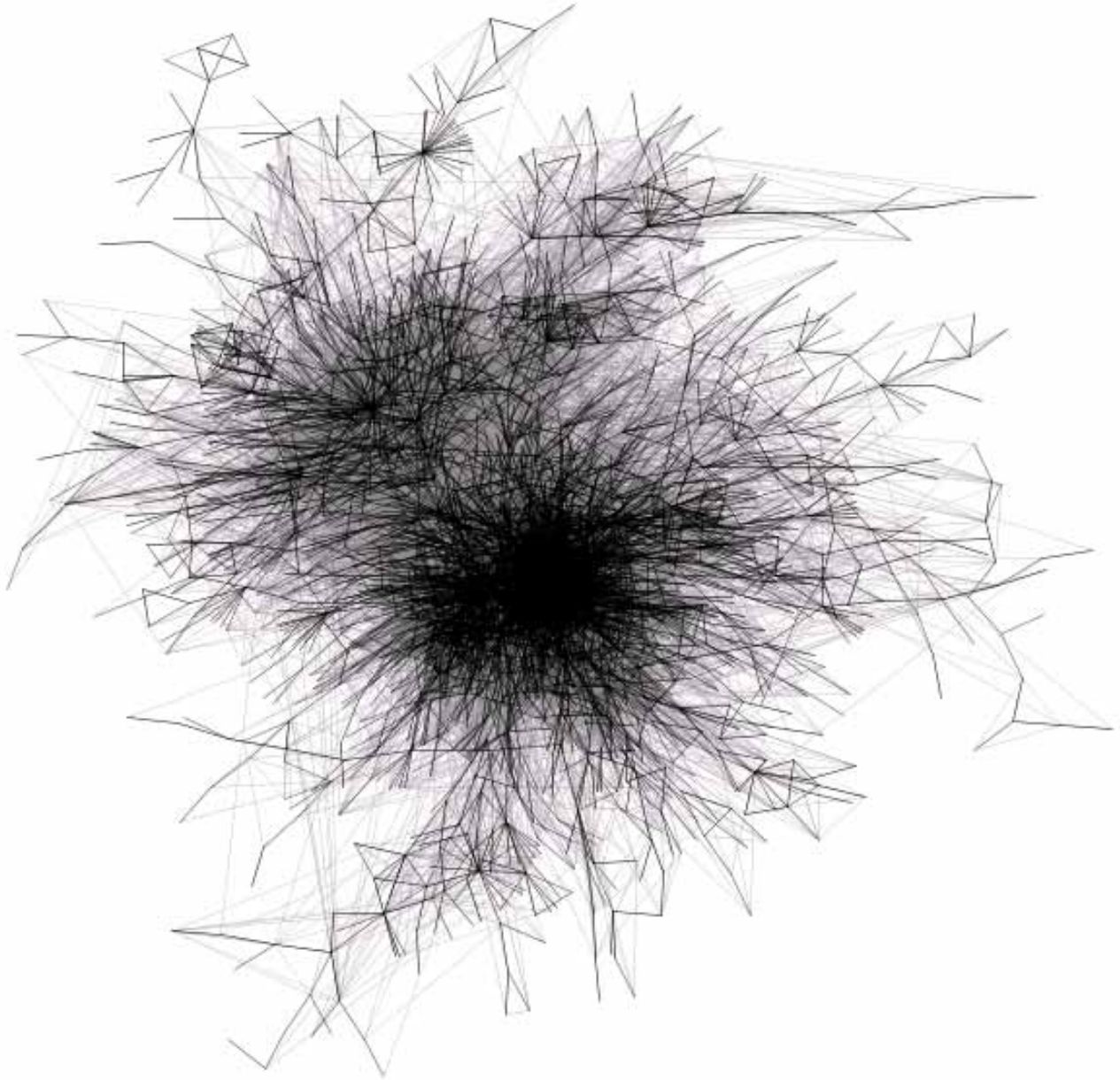
First human protein-protein interaction network

Light blue = known proteins
Orange = disease proteins
Yellow ones = not known yet



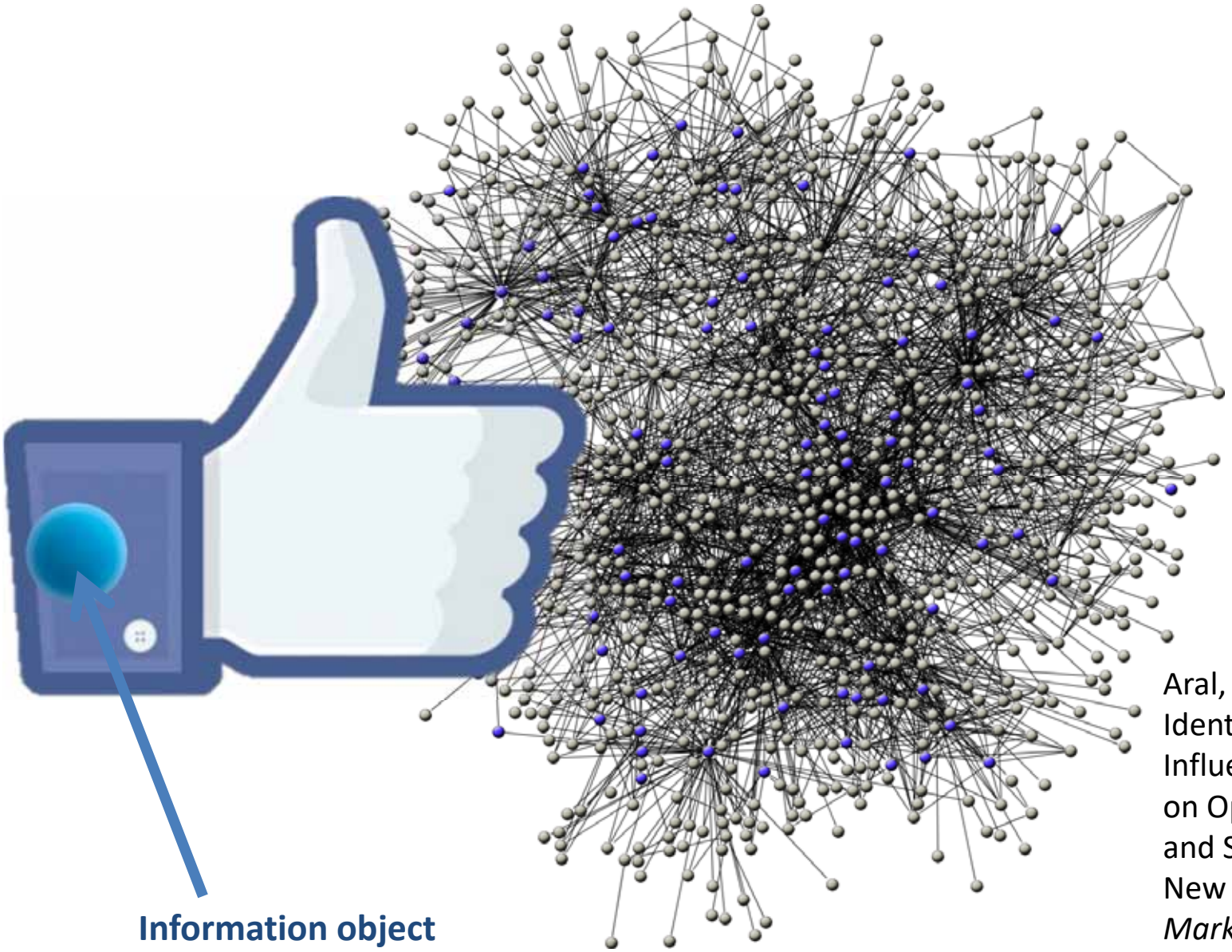
Stelzl, U. et al. (2005) A Human Protein-Protein Interaction Network: A Resource for Annotating the Proteome. *Cell, 122, 6, 957-968*.

Non-Natural Network Example: Blogosphere

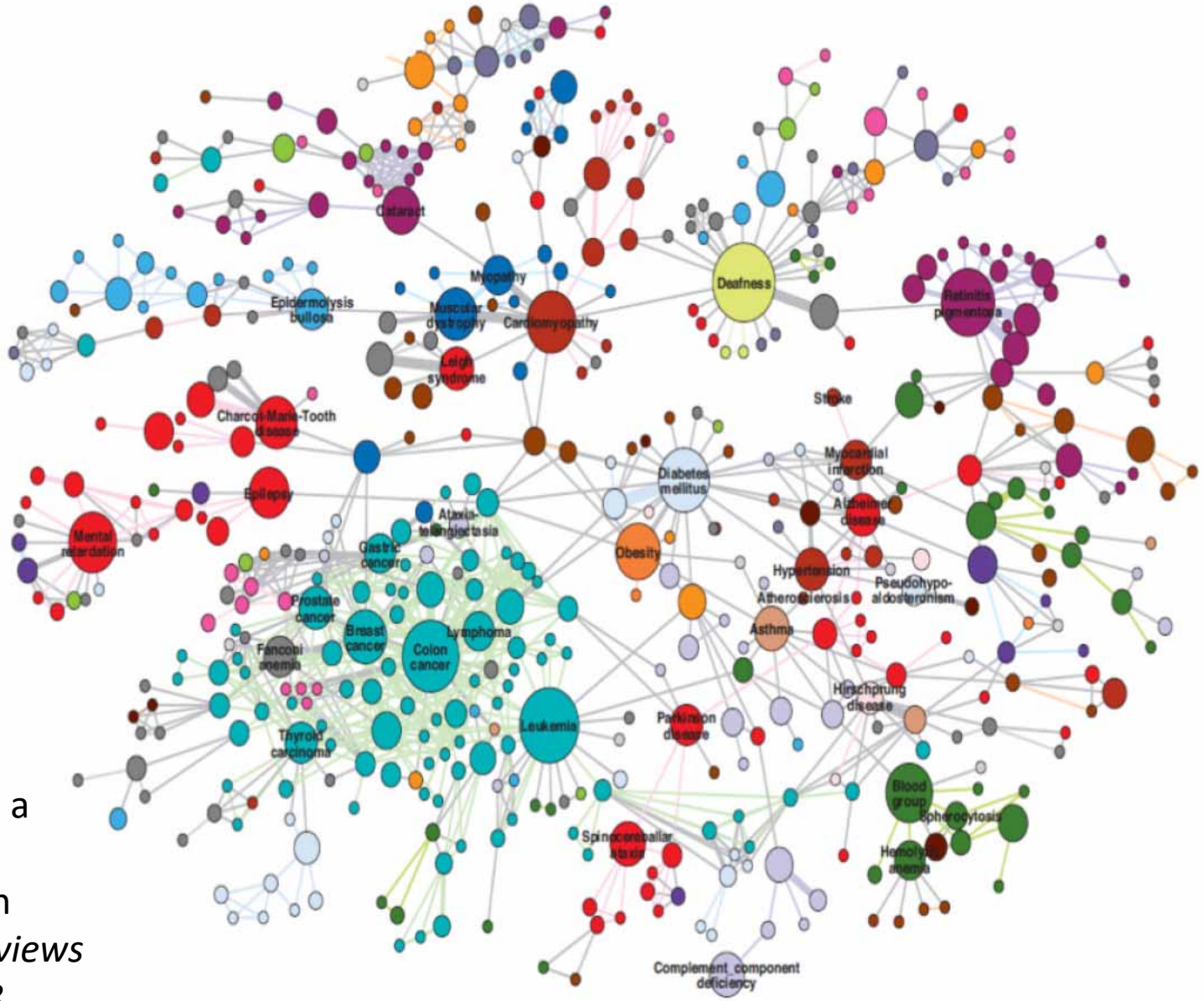


Hurst, M. (2007), Data Mining: Text Mining, Visualization and Social Media. Online available: http://datamining.typepad.com/data_mining/2007/01/the_blogosphere.html, last access: 2011-09-24

Social Behavior Contagion Network



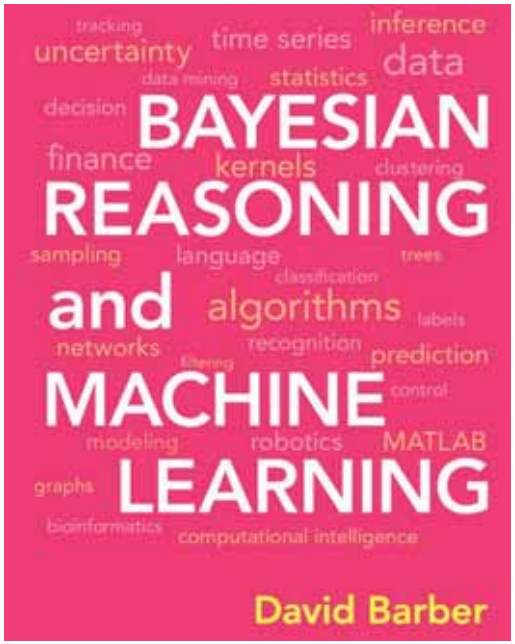
Aral, S. (2011) Identifying Social Influence: A Comment on Opinion Leadership and Social Contagion in New Product Diffusion. *Marketing Science, 30, 2, 217-223.*



Barabási, A. L., Gulbahce, N. & Loscalzo, J. 2011. Network medicine: a network-based approach to human disease. *Nature Reviews Genetics, 12*, ***56-68.***

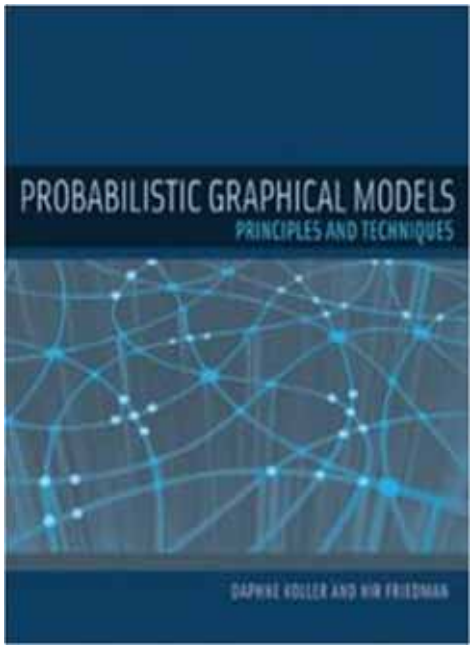
03 Bayesian Networks "Bayes' Nets"

Book recommendations 1



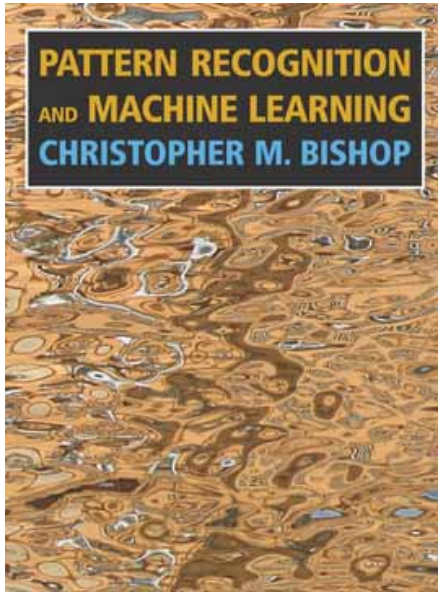
David Barber 2012. Bayesian reasoning and machine learning, Cambridge, Cambridge University Press.

http://www.cs.ucl.ac.uk/staff/d.barber/brml/



Daphne Koller & Nir Friedman 2009. Probabilistic graphical models: principles and techniques, MIT press.

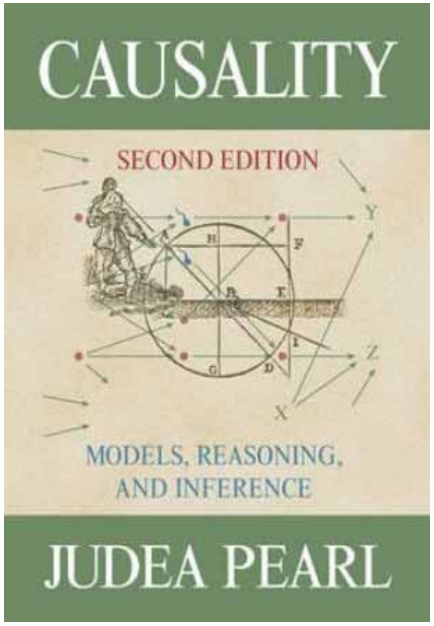
Book recommendations 2



https://goo.gl/6a7rOC

Chapter 8 Graphical Models is as sample chapter fully downloadable for free

Chris Bishop 2006. Pattern Recognition and Machine Learning, Heidelberg, Springer.



http://bayes.cs.ucla.edu/BOOK-2K/

Judea Pearl 2009. Causality: Models, Reasoning, and Inference (2nd Edition), Cambridge, Cambridge University Press.

$$P(x) = \sum_y P(x, y)$$

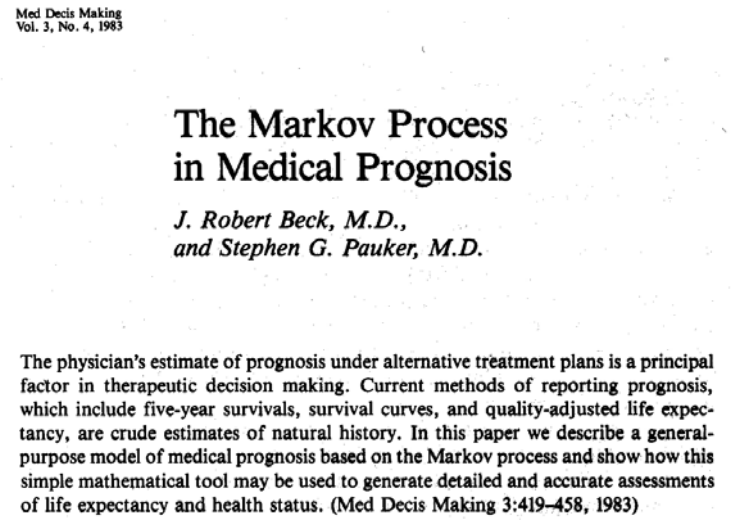
$$P(x, y) = P(y|x)P(x)$$

$$P(y|x) = \frac{P(x|y)P(y)}{P(x)}$$

$$P(x) = \sum_y P(x|y)P(y)$$

Digression: Markov Processes in Machine Learning

Why are Markov decision processes so important ?



Med Decis Making
Vol. 3, No. 4, 1983

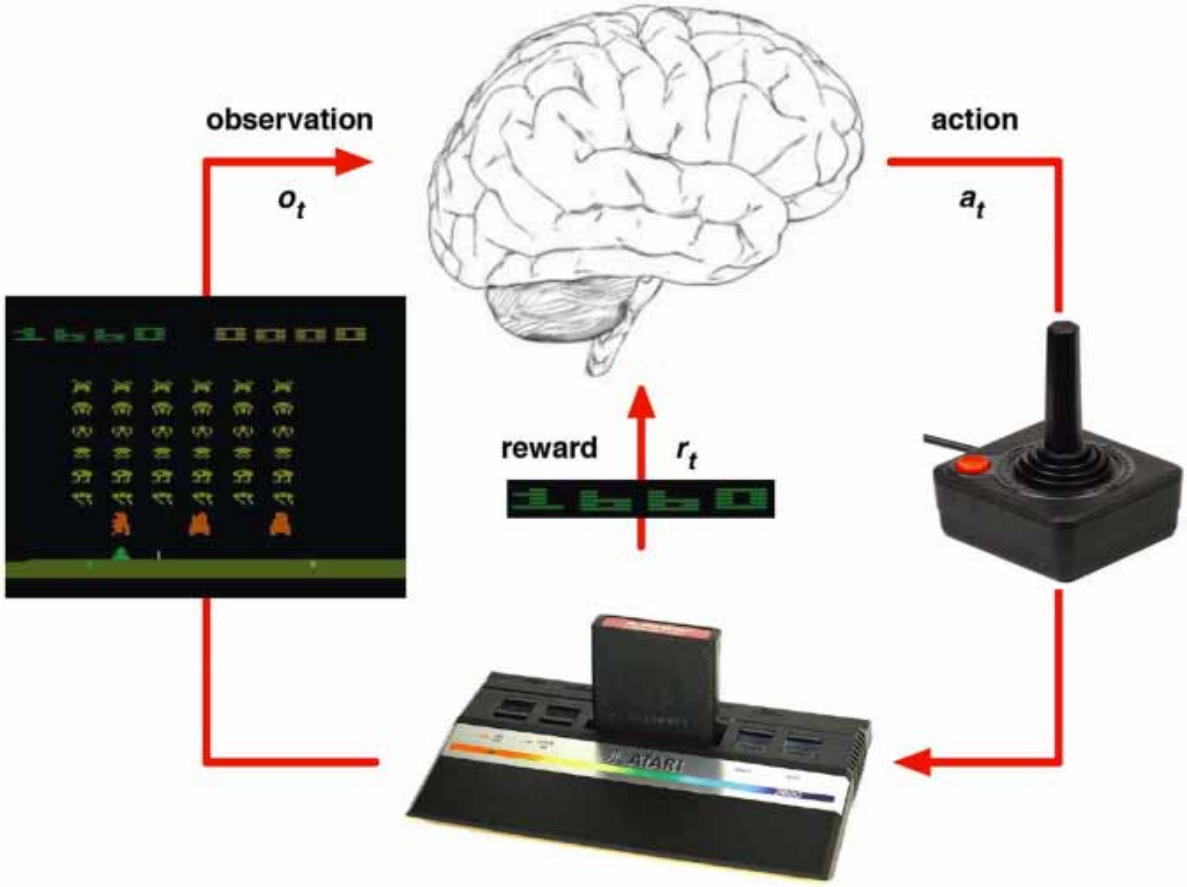
The Markov Process in Medical Prognosis

J. Robert Beck, M.D., and Stephen G. Pauker, M.D.

The physician's estimate of prognosis under alternative treatment plans is a principal factor in therapeutic decision making. Current methods of reporting prognosis, which include five-year survivals, survival curves, and quality-adjusted life expectancy, are crude estimates of natural history. In this paper we describe a general-purpose model of medical prognosis based on the Markov process and show how this simple mathematical tool may be used to generate detailed and accurate assessments of life expectancy and health status. (Med Decis Making 3:419–458, 1983)

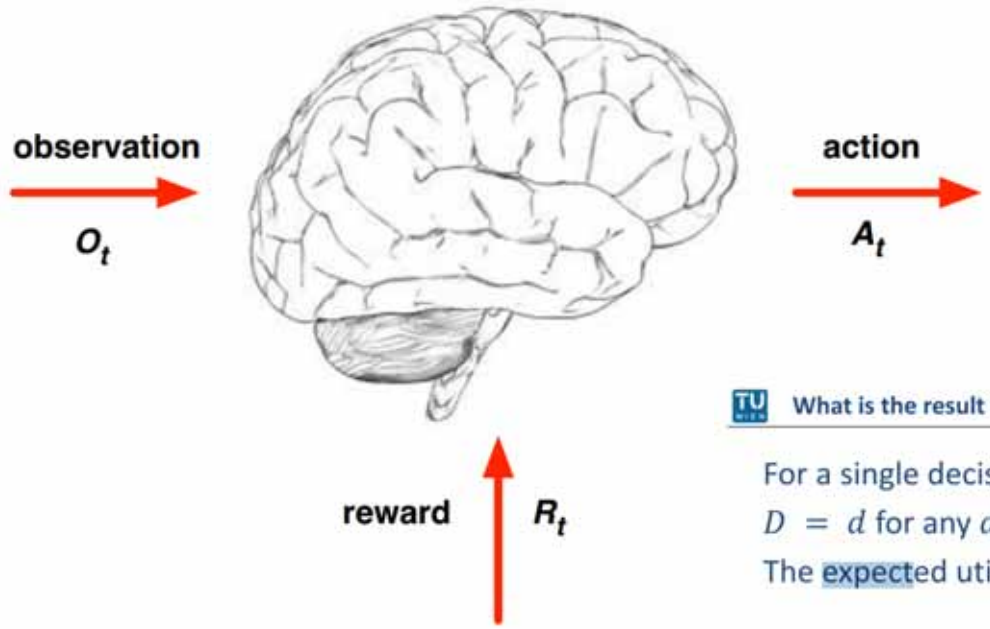
- Markov decision processes (MDP) are …
- random processes in which the future, given the present, is independent of the past!
- one of the most important classes of random processes!

How can MDP be useful for machine learning ?



Mnih, V., Kavukcuoglu, K., Silver, D., Rusu, A. A., Veness, J., Bellemare, M. G., Graves, A., Riedmiller, M., Fidjeland, A. K., Ostrovski, G., Petersen, S., Beattie, C., Sadik, A., Antonoglou, I., King, H., Kumaran, D., Wierstra, D., Legg, S. & Hassabis, D. 2015. Human-level control through deep reinforcement learning. Nature, 518, (7540), 529-533, doi:10.1038/nature14236

From where do we know such behaviour ?



What is the result of the Expected Utility Theory $E\,(U|d)$?

For a single decision variable an agent can select $D = d$ for any $d \in dom(D)$.

The expected utility of decision $D = d$ is

http://www.eoht.info/page/Oskar+Morgenstern

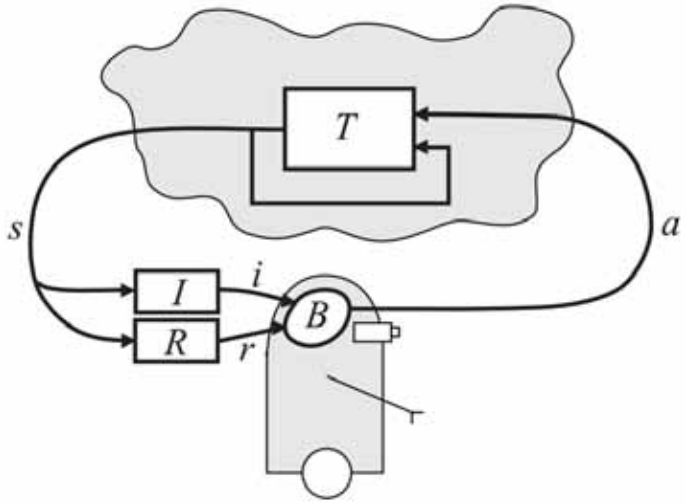
$$E(U \mid d) = \sum_{x_1,\ldots,x_n} P(x_1,\ldots,x_n \mid d)U(x_1,\ldots,x_n,d)$$

An optimal single decision is the decision $D = dmax$ whose expected utility is maximal:

$$d_{\max} = \arg \max_{d \in \mathrm{dom}(D)} E(U \mid d)$$

John Von Neumann & Oskar Morgenstern 1944. Theory of games and economic behavior, Princeton university press.

human-centered.ai (Holzinger Group) 47 2020 health AI 01

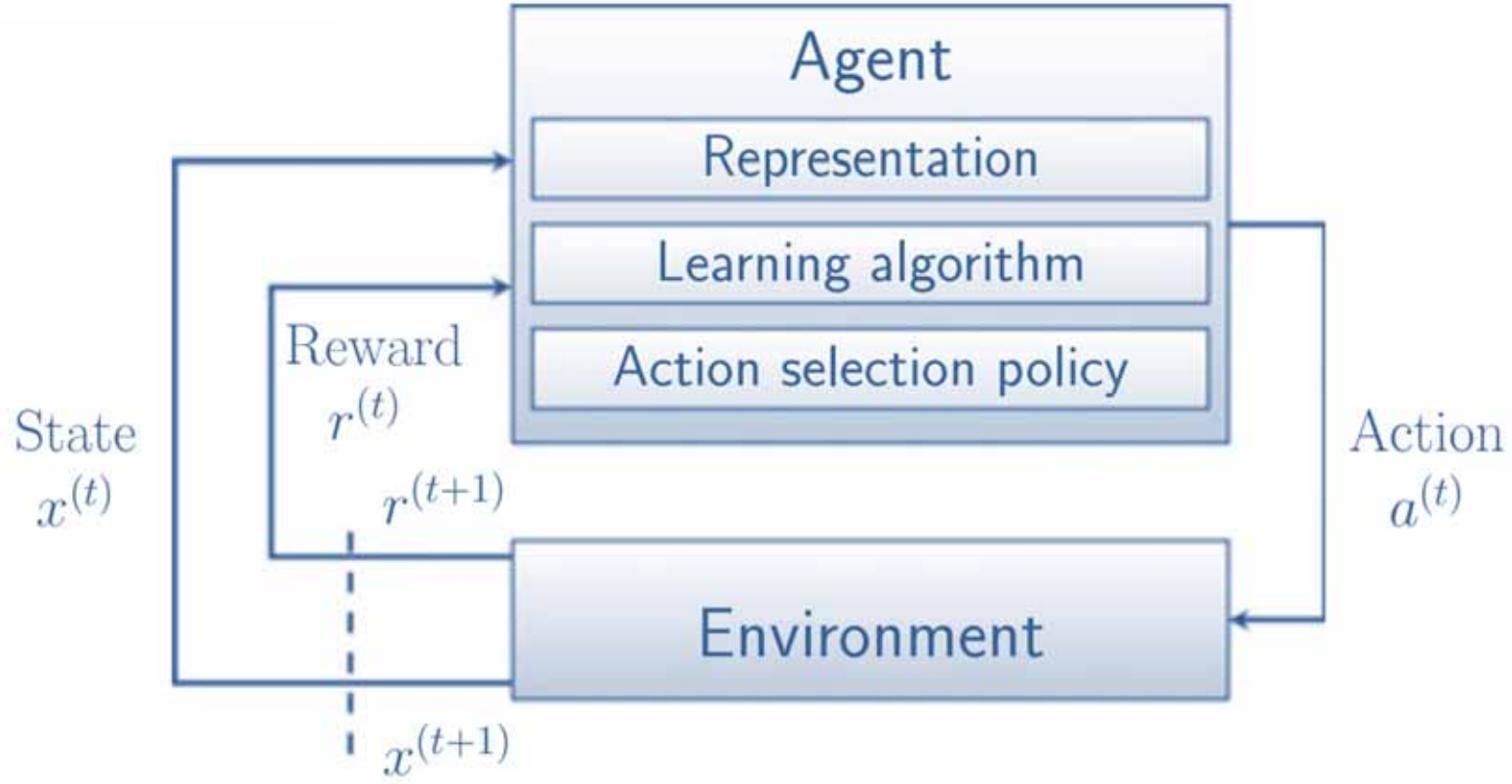



```
initialize V(s) arbitrarily
loop until policy good enough
    loop for s ∈ S
        loop for a ∈ A
            Q(s,a) := R(s,a) + γ Σ_{s'∈S} T(s,a,s') V(s')
        V(s) := max_a Q(s,a)
    end loop
end loop
```

Kaelbling, L. P., Littman, M. L. & Moore, A. W. 1996. Reinforcement learning: A survey. Journal of Artificial Intelligence Research, 4, 237-285.

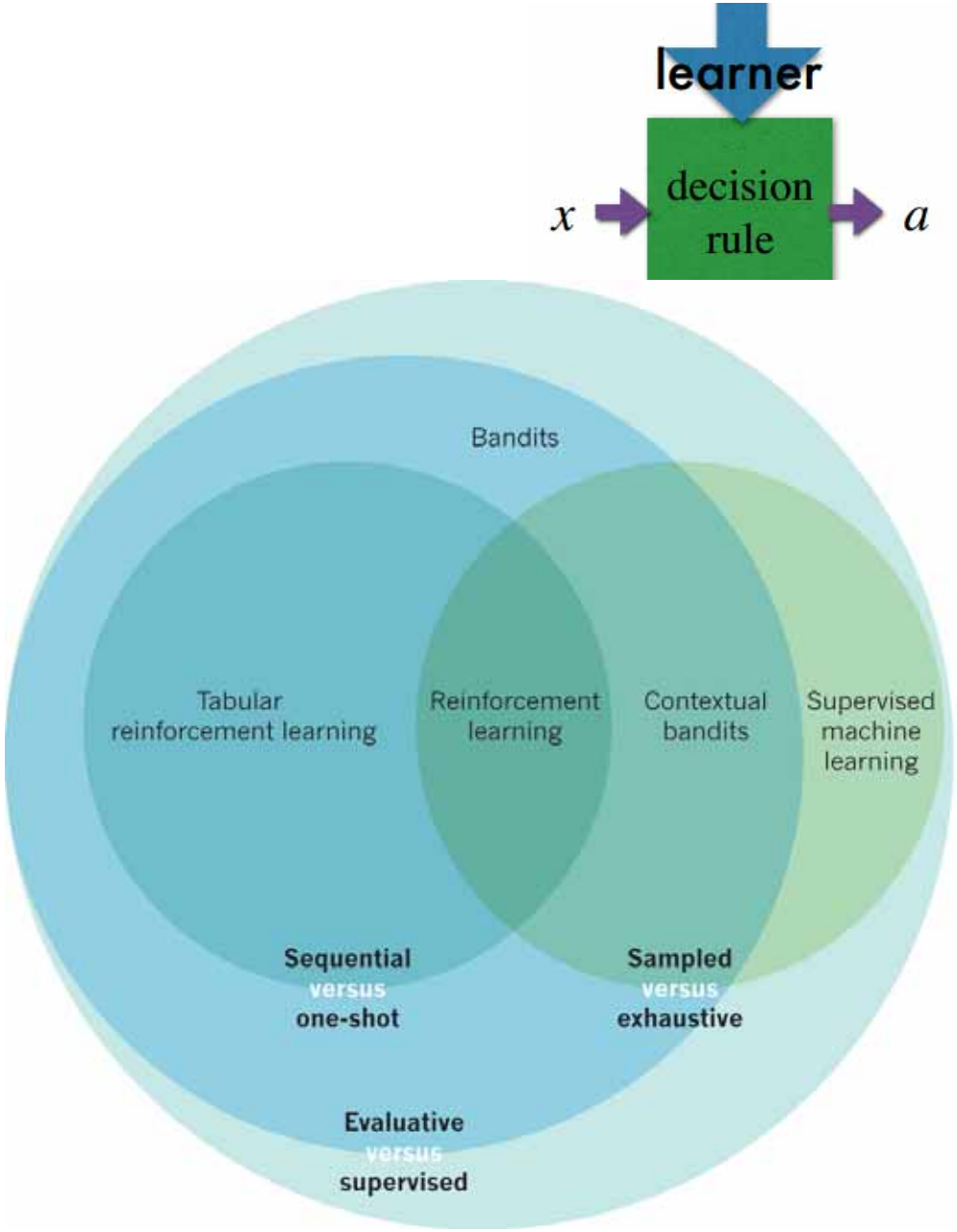
```
for t = 1, ..., n do
    The agent perceives state s_t
    The agent performs action a_t
    The environment evolves to s_{t+1}
    The agent receives reward r_t
end for
```

Intelligent behavior arises from the actions of an individual seeking to **maximize its received reward** signals in a **complex and changing world**



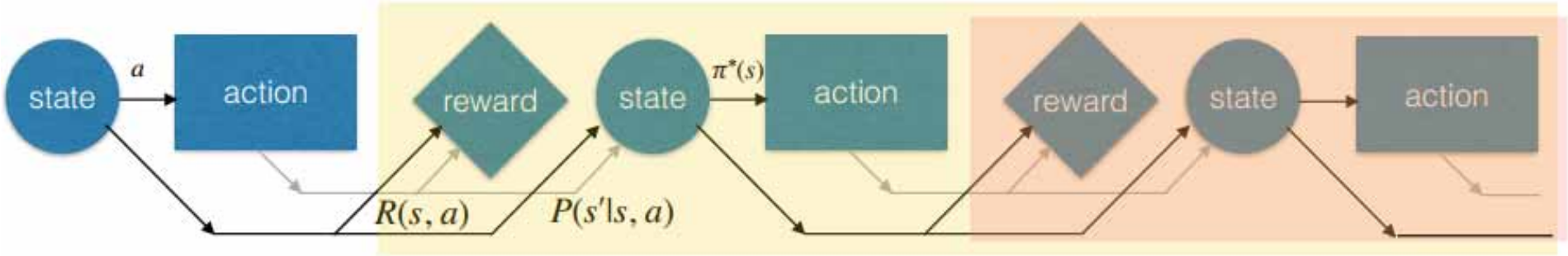
Sutton, R. S. & Barto, A. G. 1998. Reinforcement learning: An introduction, Cambridge MIT press

- Supervised:
  Learner told best a
- Exhaustive:
  Learner shown every possible x
- One-shot: Current x independent of past a



Littman, M. L. 2015. Reinforcement learning improves behaviour from evaluative feedback. Nature, 521, (7553), 445-451.

- Markov decision processes specify setting and tasks
- Planning methods use knowledge of P and R to compute a good policy π
- Markov decision process model captures both sequential feedback and the more specific one-shot feedback (when $P(s'|s,a)$ is independent of both s and a



$$Q^*(s,a) = R(s,a) + \gamma \Sigma P(s'|s,a) \max_{a'} Q^*(s',a')$$

Littman, M. L. 2015. Reinforcement learning improves behaviour from evaluative feedback. Nature, 521, (7553), 445-451.

- 1) Overserves
- 2) Executes
- 3) Receives Reward
- Executes action A_t:
- $O_t = sa_t = se_t$
- Agent state = environment state = information state
- Markov decision process (MDP)

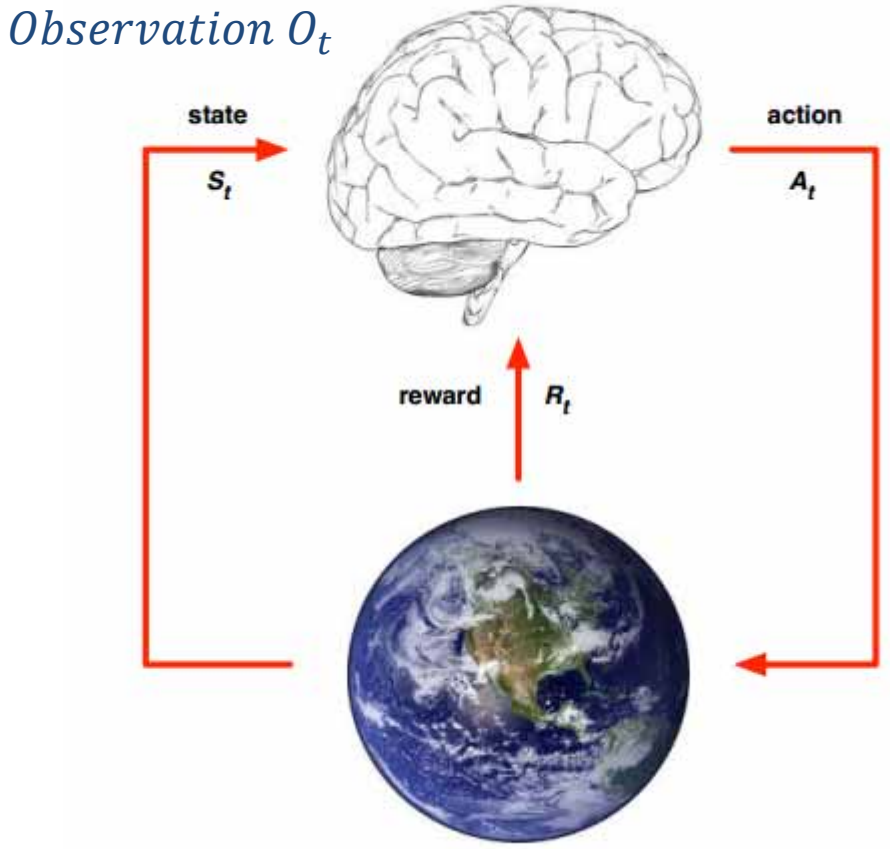
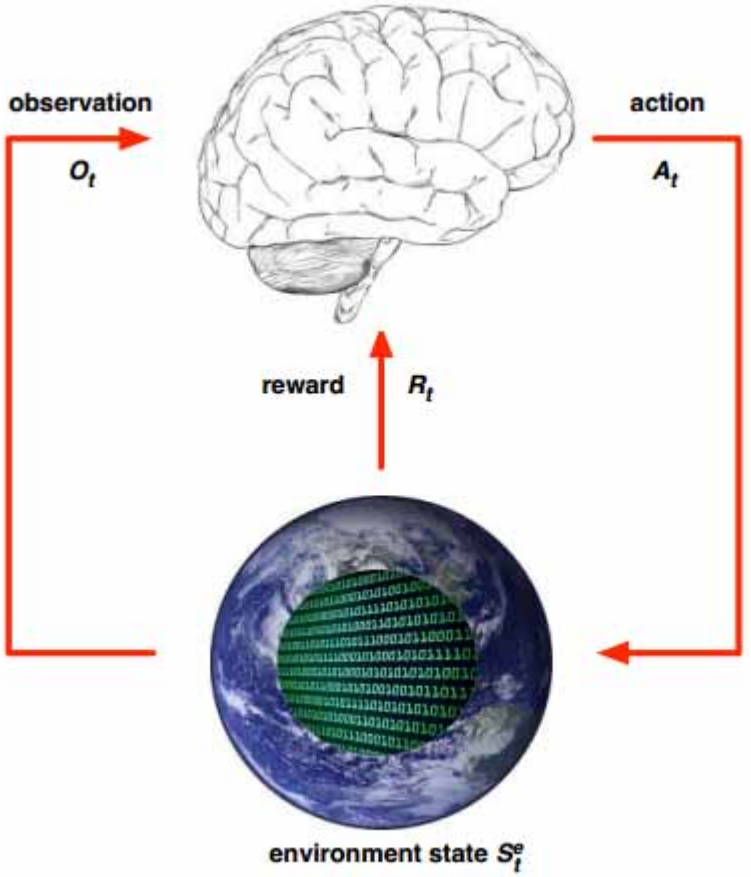


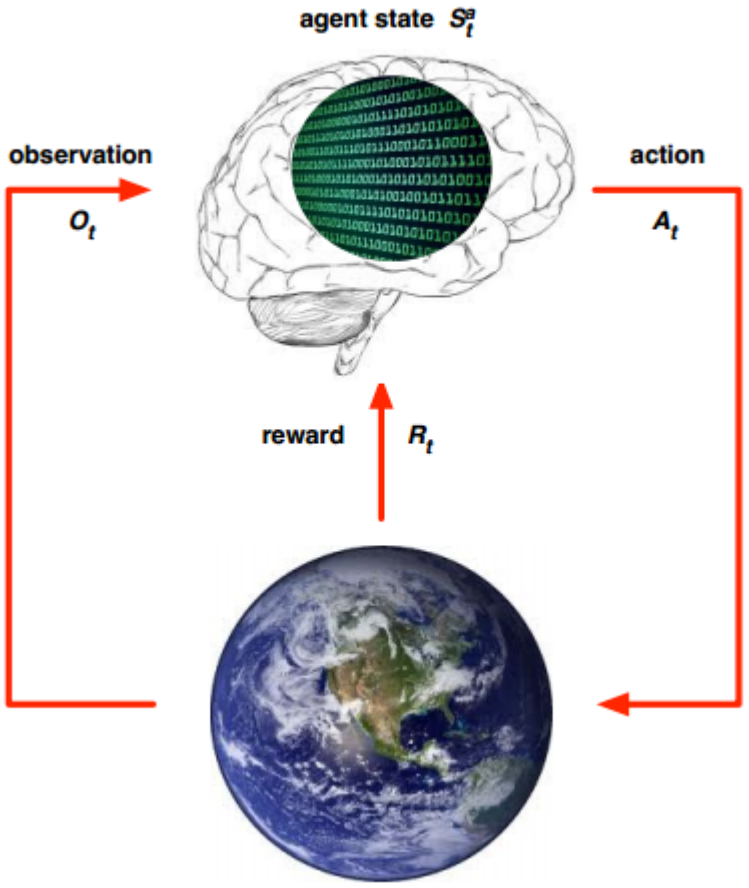
Image credit to David Silver, UCL

- i.e. whatever data the environment uses to pick the next observation/reward
- The environment state is not usually visible to the agent
- Even if S is visible, it may contain irrelevant information
- A State S_t is Markov iff:



$$\mathbb{P}[S_{t+1}|S_t] = \mathbb{P}[S_{t+1}|S_1, \dots, S_t]$$

- i.e. whatever information the agent uses to pick the next action
- it is the information used by reinforcement learning algorithms
- It can be any function of history:
- $S = f(H)$



$$H_t = O_1, R_1, A_1, \ldots, A_{t-1}, O_t, R_t$$

- RL agent components:
  - Policy: agent's behaviour function
  - Value function: how good is each state and/or action
  - Model: agent's representation of the environment
- Policy as the agent's behaviour
  - is a map from state to action, e.g.
  - Deterministic policy: a = (s)
  - Stochastic policy: (ajs) = P[At = ajS t = s
- Value function is prediction of future reward:

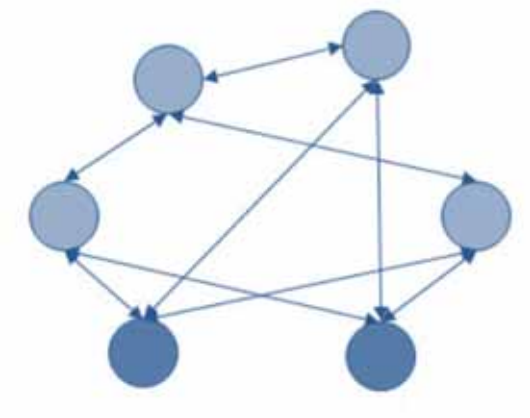
$$v_\pi(s) = \mathbb{E}_\pi \left[R_{t+1} + \gamma R_{t+2} + \gamma^2 R_{t+3} + \dots \mid S_t = s \right]$$

- Partial observability: when agent only indirectly observes environment
- Formally this is a Partially Observable Markov Decision Process (POMDP):
  - Agent must construct its own state representation S,

for example:

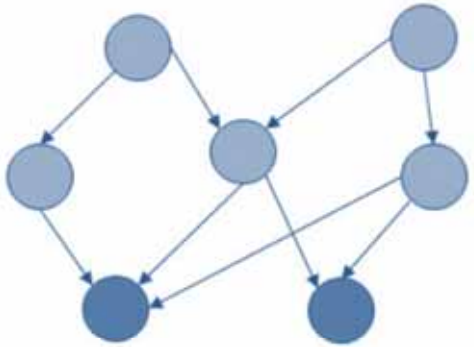
- Complete history: $S_t^a = H_t$
- Beliefs of environment state: $S_t^a = (\mathbb{P}[S_t^e = s^1], \ldots, \mathbb{P}[S_t^e = s^n])$
- Recurrent neural network: $S_t^a = \sigma(S_{t-1}^a W_s + O_t W_o)$

Back to Bayesian Networks



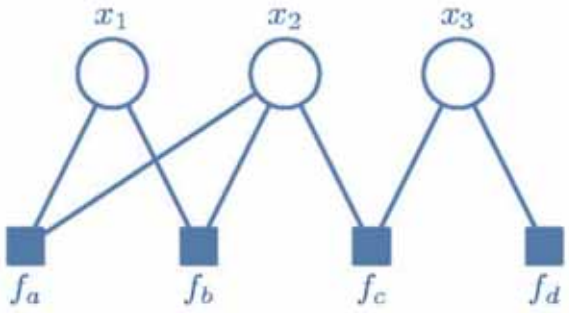
Undirected: Markov random fields, useful e.g. for computer vision (Details: Murphy 19)

$$P(\boldsymbol{X}) = \frac{1}{Z}\exp\left(\sum_{ij} W_{ij}\, x_i x_j + \sum_i x_i b_i\right)$$



Directed: Bayes Nets, useful for designing models (Details: Murphy 10)

$$p(\mathbf{x}) = \prod_{k=1}^{K} p(x_k|\mathrm{pa}_k)$$



Factored: useful for inference/learning

$$p(\mathbf{x}) = \prod_s f_s(\mathbf{x}_s)$$

- is a **probabilistic model**, consisting of two parts:
- 1) a dependency structure and
- 2) local probability models.

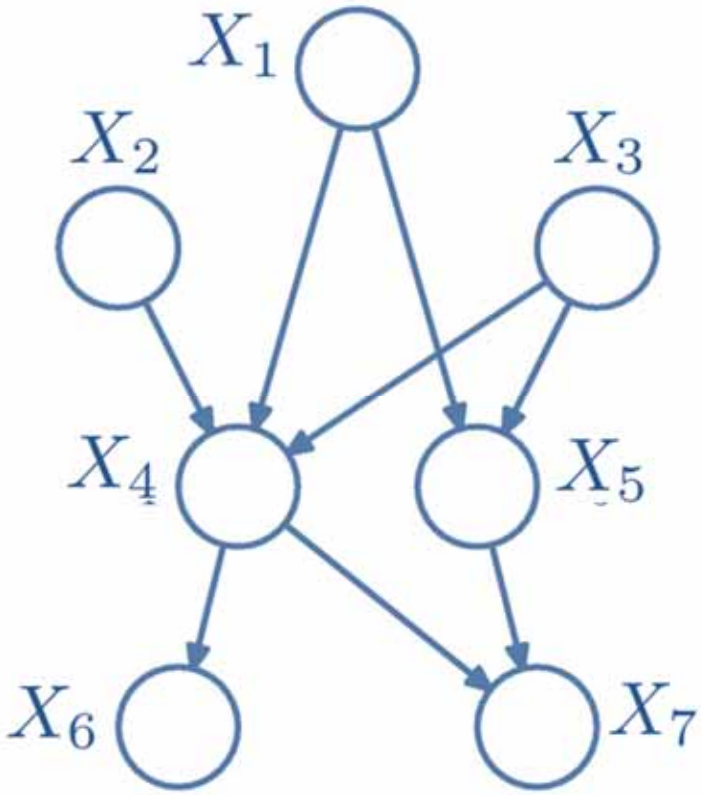
$$p(x_1, \ldots, x_n) = \prod_{i=1}^{n} p(x_i \mid Pa(x_i))$$

Where $Pa(x_i)$ are the parents of x_i

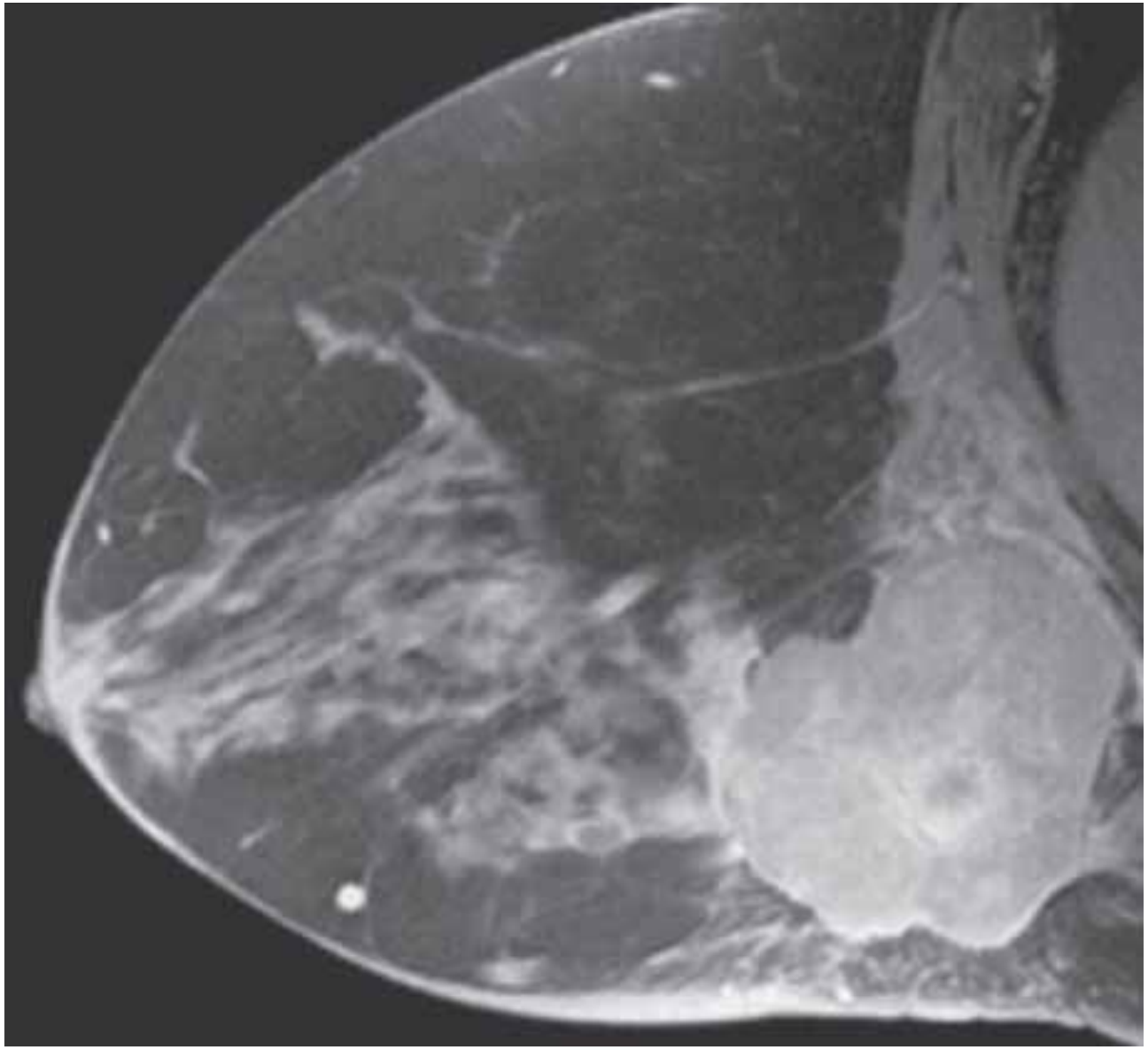
BN inherently model the **uncertainty in the data.** They are a successful marriage between probability theory and graph theory; allow to model a multidimensional probability distribution in a sparse way by searching independency relations in the data. Furthermore this model allows different strategies to integrate two data sources.

Pearl, J. (1988) *Probabilistic reasoning in intelligent systems: networks of plausible inference. San Francisco, Morgan Kaufmann.*

$$p(X_1, \ldots, X_7) =$$
$$p(X_1)p(X_2)p(X_3)p(X_4|X_1, X_2, X_3) \cdot$$
$$p(X_5|X_1, X_3)p(X_6|X_4)p(X_7|X_4, X_5)$$

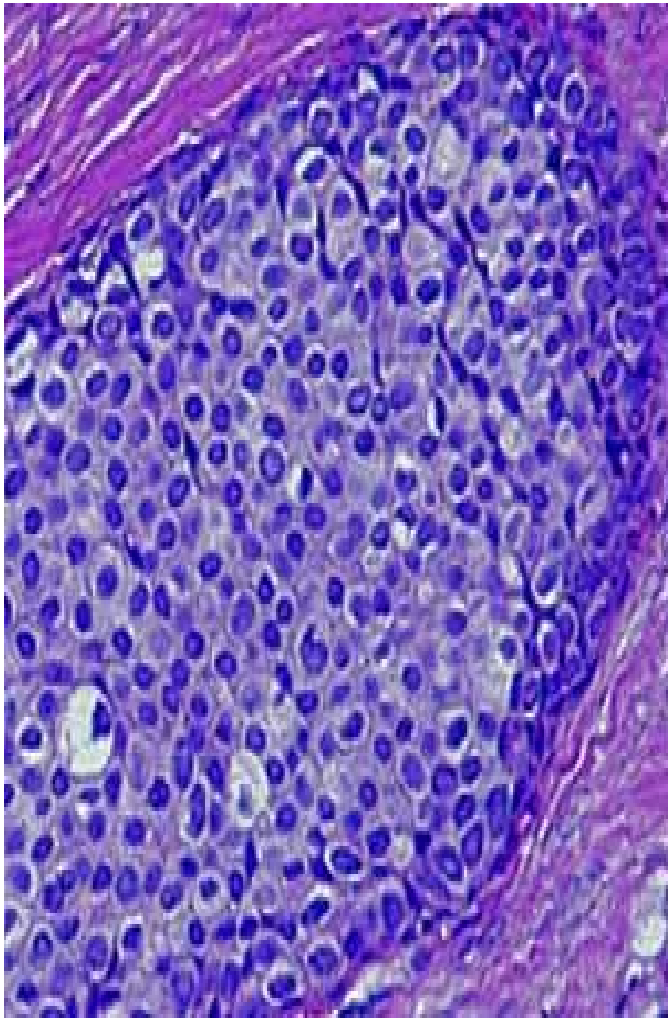


Clinical Case Example

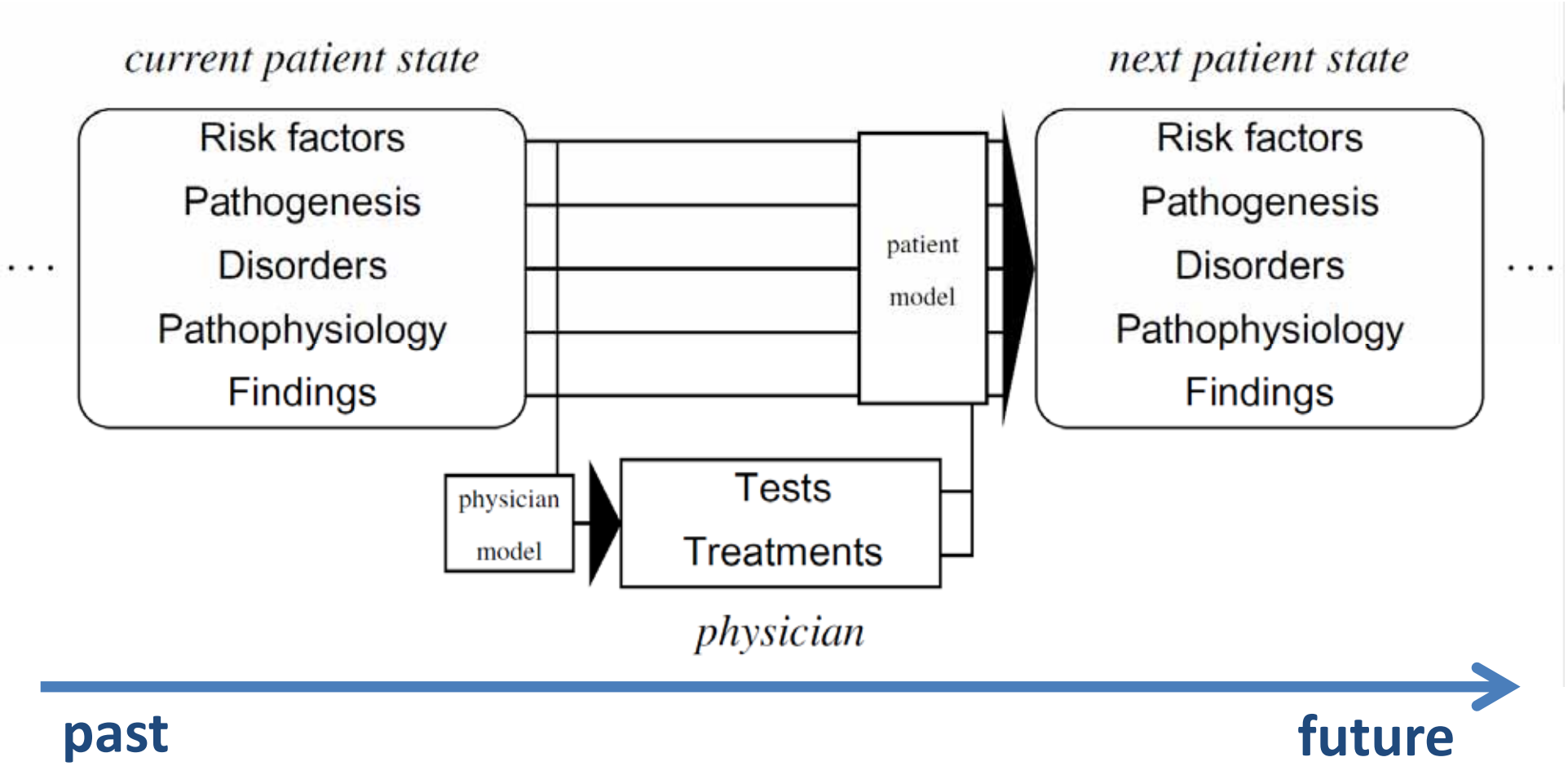


Overmoyer, B. A., Lee, J. M. & Lerwill, M. F. (2011) Case 17-2011 A 49-Year-Old Woman with a Mass in the Breast and Overlying Skin Changes. *New England Journal of Medicine, 364, 23, 2246-2254.*

- = the prediction of the future course of a disease conditional on the patient's history and a projected treatment strategy
- Danger: probable Information !
- Therefore valid prognostic models can be of great benefit for clinical decision making and of great value to the patient, e.g., for notification and quality of-life decisions



Knaus, W. A., Wagner, D. P. & Lynn, J. (1991) Short-term mortality predictions for critically ill hospitalized adults: science and ethics. *Science, 254, 5030, 389.*

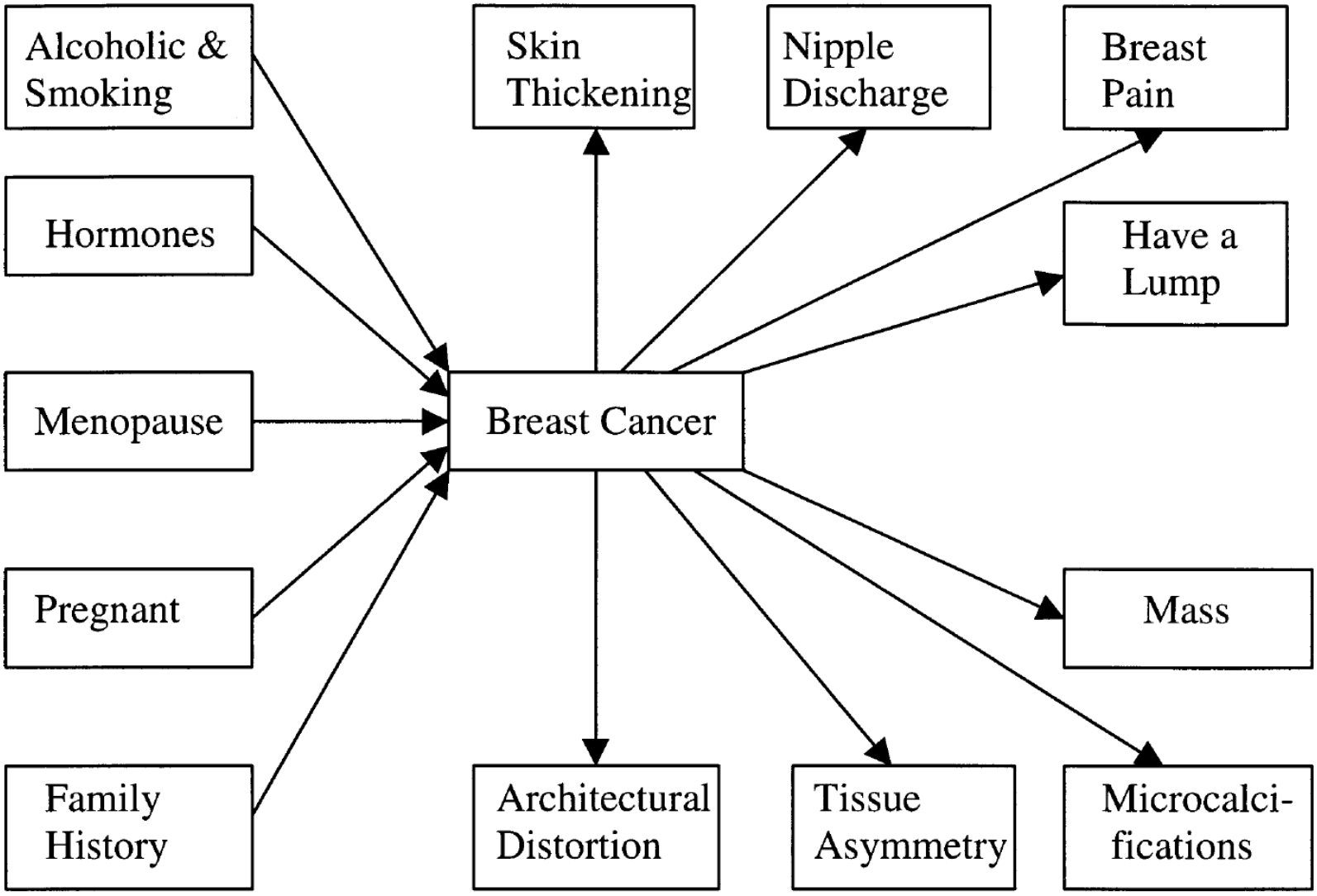


van Gerven, M. A. J., Taal, B. G. & Lucas, P. J. F. (2008) Dynamic Bayesian networks as prognostic models for clinical patient management. *Journal of Biomedical Informatics, 41, 4, 515-529.*

Example: Breast cancer - Probability Table

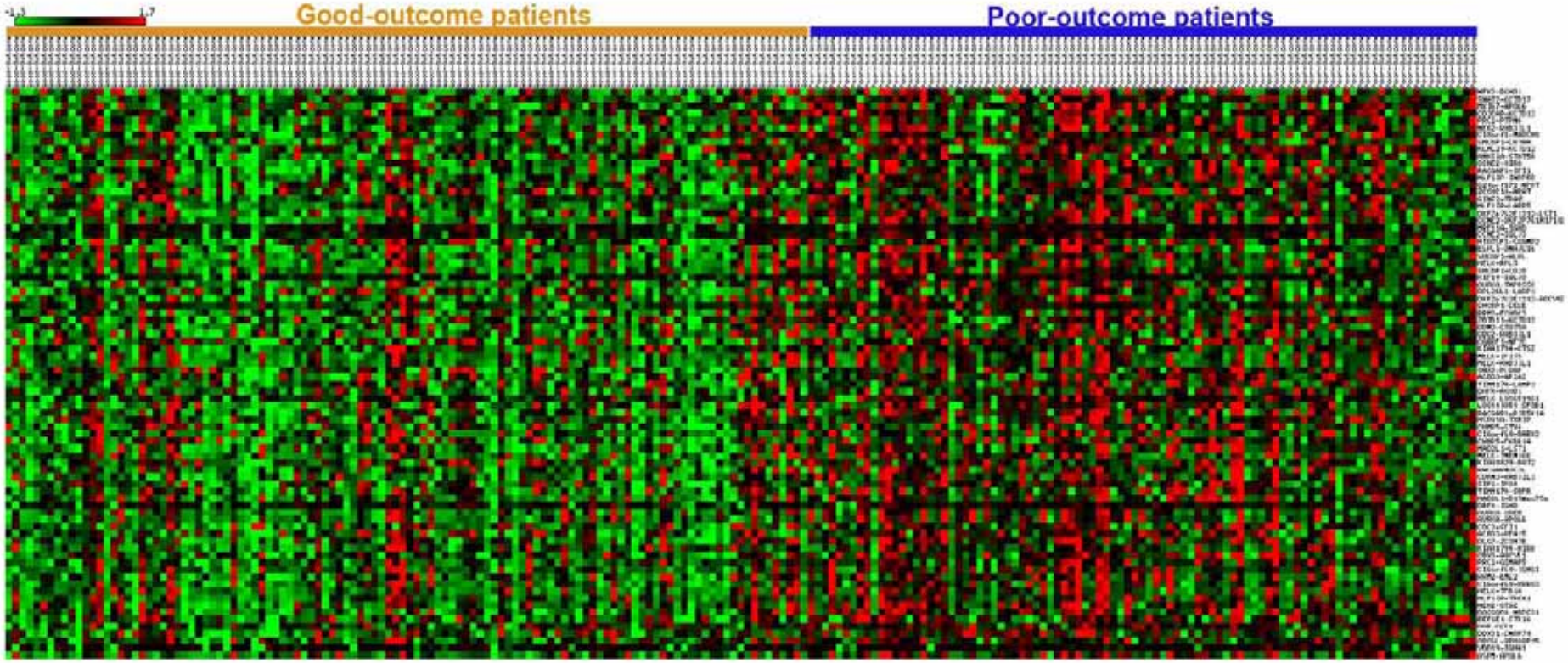
Category	Node description	State description
Diagnosis	Breast cancer	Present, absent.
Clinical history	Habit of drinking alcoholic beverages and smoking	Yes, no.
	Taking female hormones	Yes, no.
	Have gone through menopause	Yes, no.
	Have ever been pregnant	Yes, no.
	Family member has breast cancer	Yes, no.
Physical findings	Nipple discharge	Yes, no.
	Skin thickening	Yes, no.
	Breast pain	Yes, no.
	Have a lump(s)	Yes, no.
Mammographic findings	Architectural distortion	Present, absent.
	Mass	Score from one to three, score from four to five, absent
	Microcalcification cluster	Score from one to three, score from four to five, absent
	Asymmetry	Present, absent.

Wang, X. H., et al. (1999) Computer-assisted diagnosis of breast cancer using a data-driven Bayesian belief network. *International Journal of Medical Informatics, 54, 2, 115-126.*



Wang, X. H., et al. (1999) Computer-assisted diagnosis of breast cancer using a data-driven Bayesian belief network. *International Journal of Medical Informatics, 54, 2, 115-126.*

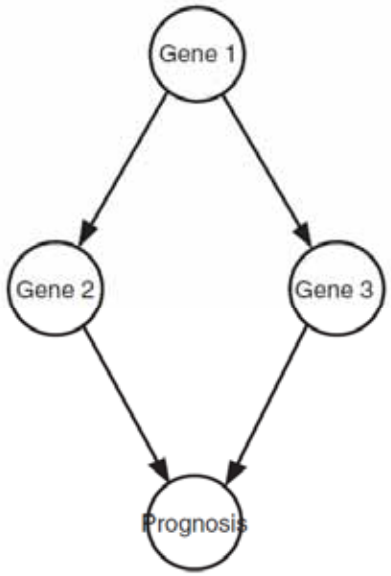
- Integrating microarray data from multiple studies to increase sample size;
- = approach to the development of more robust prognostic tests



Xu, L., Tan, A., Winslow, R. & Geman, D. (2008) Merging microarray data from separate breast cancer studies provides a robust prognostic test. *BMC Bioinformatics, 9, 1, 125-139.*

Example: Bayes Net with four binary variables

Gene 1	
P(on)	0.8
P(off)	0.2

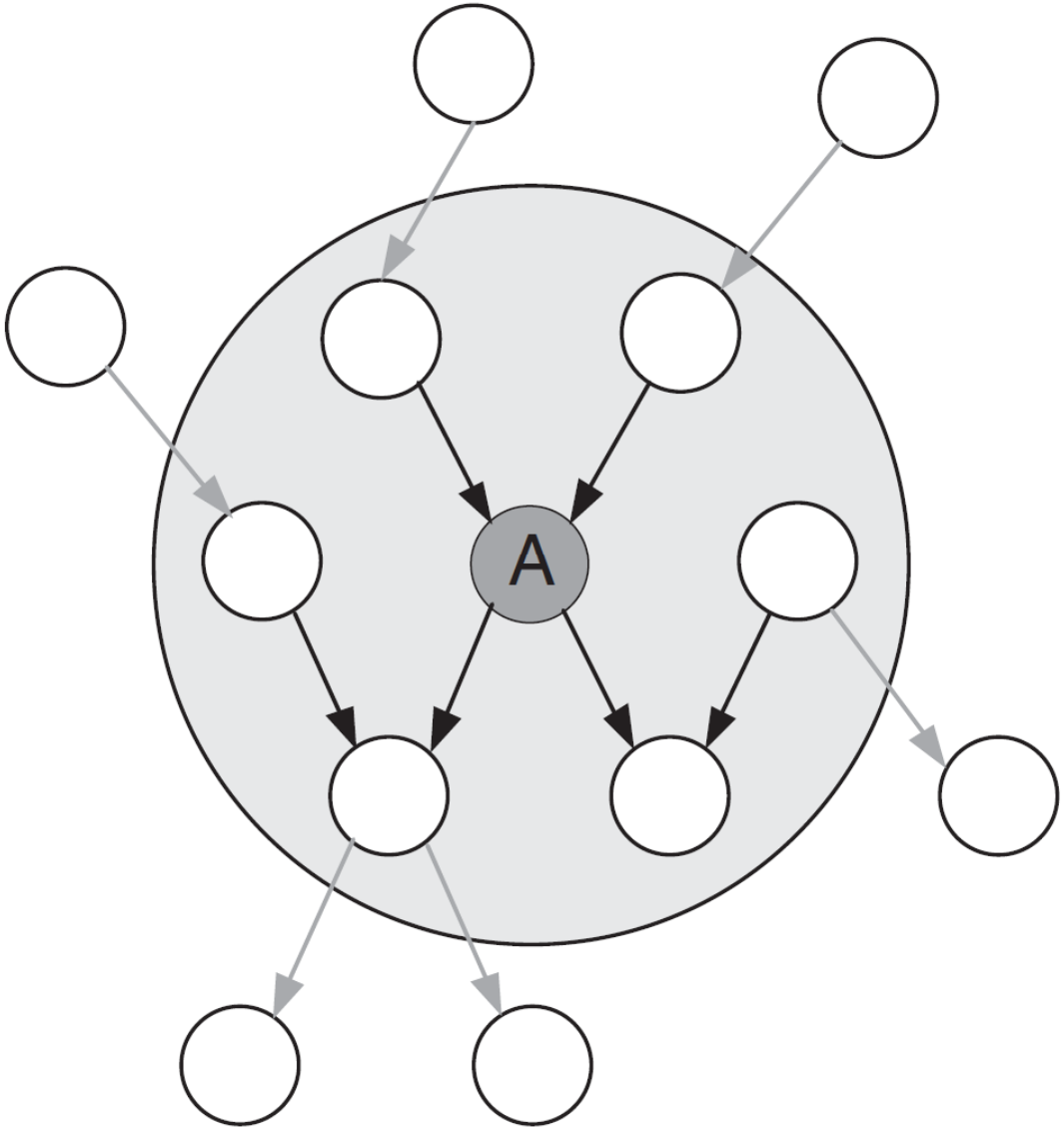


Gene 2	Gene 1 on	Gene 1 off
P(on)	0.3	0.6
P(off)	0.7	0.4

Gene 2	Gene 1 on	Gene 1 off
P(on)	0.3	0.6
P(off)	0.7	0.4

Prognosis	Gene 2 on Gene 3 on	Gene 2 on Gene 3 off	Gene 2 off Gene 2 on	Gene 2 off Gene 3 off
P(good)	0.6	0.1	0.9	0.5
P(poor)	0.4	0.9	0.1	0.5

Gevaert, O., Smet, F. D., Timmerman, D., Moreau, Y. & Moor, B. D. (2006) Predicting the prognosis of breast cancer by integrating clinical and microarray data with Bayesian networks. *Bioinformatics, 22, 14, 184-190.*



Gevaert, O., Smet, F. D., Timmerman, D., Moreau, Y. & Moor, B. D. (2006) Predicting the prognosis of breast cancer by integrating clinical and microarray data with Bayesian networks. *Bioinformatics, 22, 14, 184-190.*

- First the structure is learned using a search strategy.
- Since the number of possible structures increases super exponentially with the number of variables,
- the well-known greedy search algorithm K2 can be used in combination with the Bayesian Dirichlet (BD) scoring metric:

$$p(S|D) \propto p(S) \prod_{i=1}^{n} \prod_{j=1}^{q_i} \left[\frac{\Gamma(N'_{ij})}{\Gamma(N'_{ij} + N_{ij})} \prod_{k=1}^{r_i} \frac{\Gamma(N'_{ijk} + N_{ijk})}{\Gamma(N'_{ijk})} \right]$$

N_{ijk} … number of cases in the data set D
having variable i in state k associated with the j-th instantiation of its parents in current structure S.
n is the total number of variables.

Dependency Structure – first step (2/2)

- Next, N_{ij} is calculated by summing over all states of a variable:
- $N_{ij} = \sum_{k=1}^{r_i} N_{ijk} \cdot N'_{ijk}$ and N'_{ij} have similar meanings but refer to prior knowledge for the parameters.
- When no knowledge is available they are estimated using $N_{ijk} = N/(r_i q_i)$
- with N the equivalent sample size,
- r_i the number of states of variable i and
- q_i the number of instantiations of the parents of variable i.
- $\Gamma(.)$ corresponds to the gamma distribution.
- Finally $p(S)$ is the prior probability of the structure.
- $p(S)$ is calculated by:
- $p(S) = \prod_{i=1}^{n} \prod_{l_i=1}^{p_i} p(l_i \longrightarrow x_i) \prod_{m_i=1}^{o_i} p(m_i x_i)$
- with p_i the number of parents of variable x_i and o_i all the variables that are not a parent of x_i.
- Next, $p(a \longrightarrow b)$ is the probability that there is an edge from a to b while $p(ab)$ is the inverse, i.e. the probability that there is no edge from a to b

- Estimating the parameters of the local probability models corresponding with the dependency structure.
- CPTs are used to model these local probability models.
- For each variable and instantiation of its parents there exists a CPT that consists of a set of parameters.
- Each set of parameters was given a uniform Dirichlet prior:

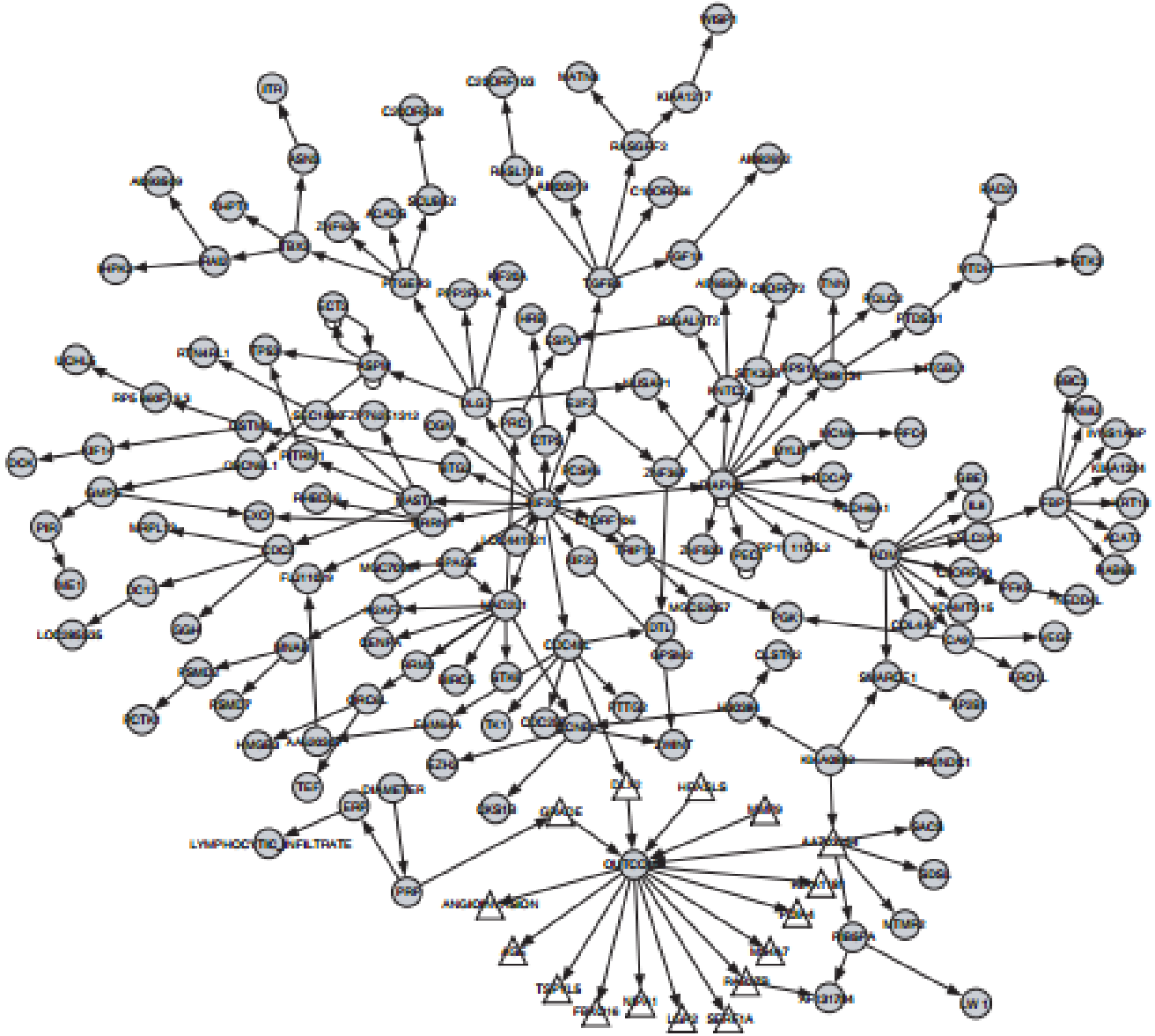
$$\mathrm{p}(\theta_{ij}|S) = Dir(\theta_{ij}|N'_{ij1}, \ldots, N'_{ijk}, \ldots, N'_{ijr_i})$$

Note: With θ_{ij} a parameter set where i refers to the variable and j to the j-th instantiation of the parents in the current structure. θ_{ij} contains a probability for every value of the variable x_i given the current instantiation of the parents. Dir corresponds to the Dirichlet distribution with $(N'_{ij1}, \ldots, N'_{ijr_i})$ as parameters of this Dirichlet distribution. Parameter learning then consists of updating these Dirichlet priors with data. This is straightforward because the multinomial distribution that is used to model the data, and the Dirichlet distribution that models the prior, are conjugate distributions. This results in a Dirichlet posterior over the parameter set:

$$\mathrm{p}(\theta_{ij}|D,S) = Dir(\theta_{ij}|N'_{ij1} + N_{ij1}, \ldots, N'_{ijk} + N_{ijk}, \ldots, N'_{ijr_i} + N_{ijr_i})$$

with N_{ijk} defined as before.

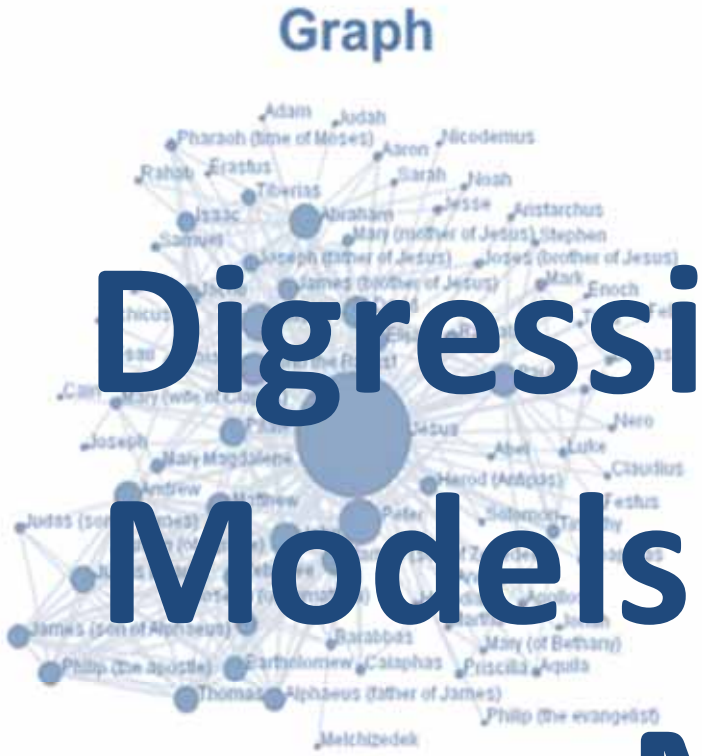
Predicting the prognosis of breast cancer (integrated a.)



Gevaert, O., Smet, F. D., Timmerman, D., Moreau, Y. & Moor, B. D. (2006) Predicting the prognosis of breast cancer by integrating clinical and microarray data with Bayesian networks. *Bioinformatics, 22, 14, 184-190.*

- For certain cases it is tractable if:
  - Just one variable is unobserved
  - We have singly connected graphs (no undirected loops -> belief propagation)
  - Assigning probability to fully observed set of variables
- Possibility: Monte Carlo Methods (generate many samples according to the Bayes Net distribution and then count the results)
- Otherwise: approximate solutions …

Often it is better to have a good solution within time – than an perfect solution too late …

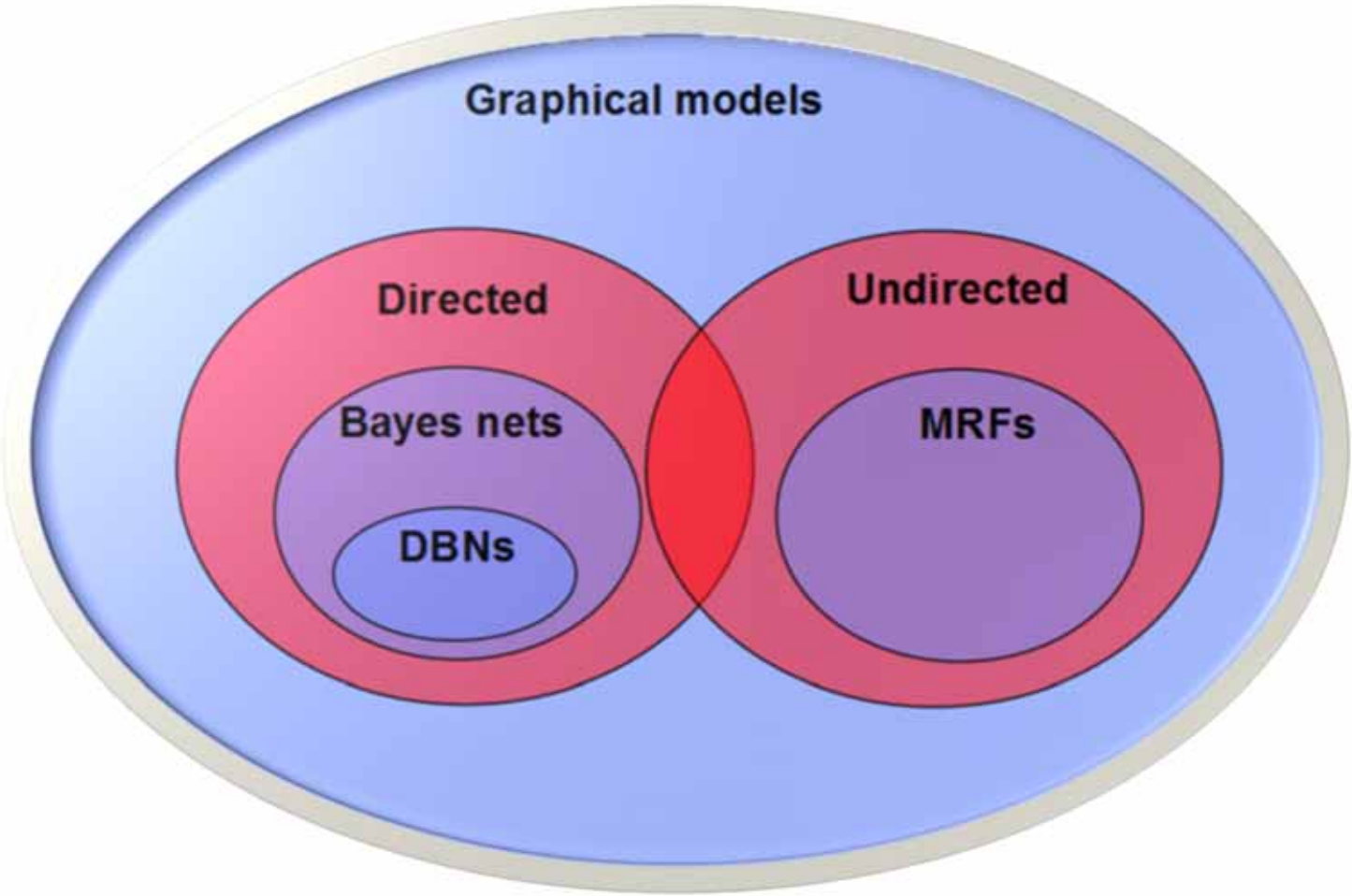


Digression: Graphical Models and Decision Making

$\mathcal{M}$

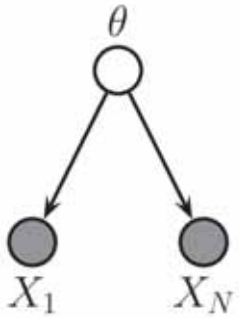
$$\mathcal{D} \equiv \{X_1^{(i)}, X_2^{(i)}, \ldots, X_m^{(i)}\}_{i=1}^{N}$$

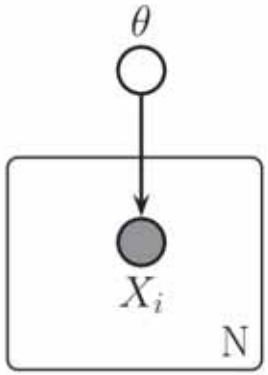
What Classes of Graphical Models do we know ?



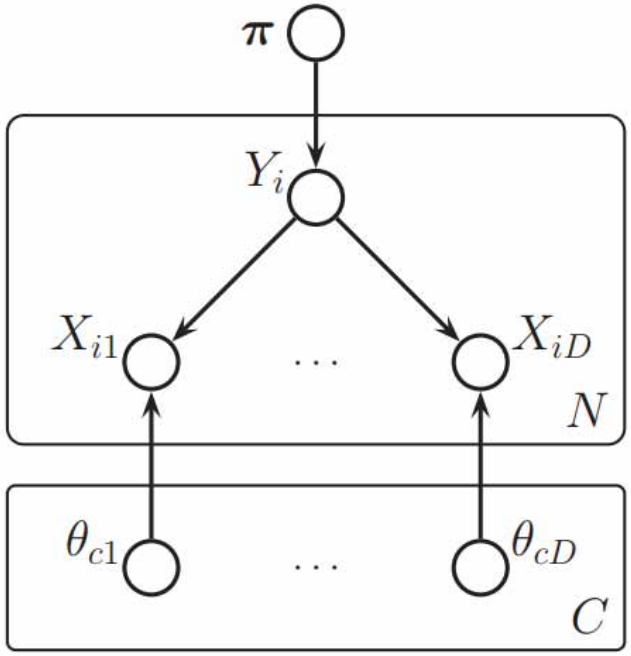
Murphy, K. P. 2012. Machine learning: a probabilistic perspective, Cambridge (MA), MIT press.

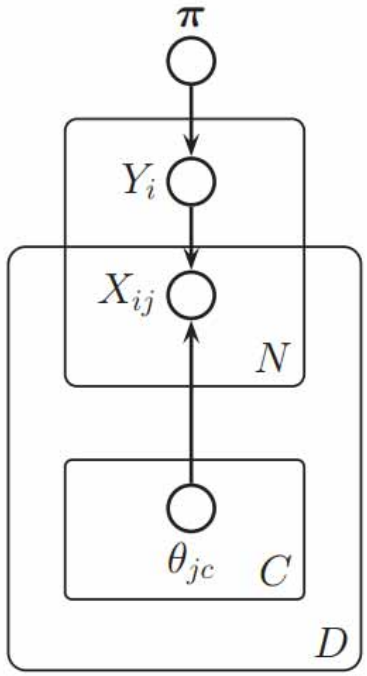
Naïve Bayes classifier as DGM (single/nested plates)





$$p(\boldsymbol{\theta}, \mathcal{D}) = p(\boldsymbol{\theta}) \left[\prod_{i=1}^{N} p(\mathbf{x}_i|\boldsymbol{\theta}) \right]$$





Murphy, K. P. 2012. Machine learning: a probabilistic perspective, Cambridge (MA), MIT press.

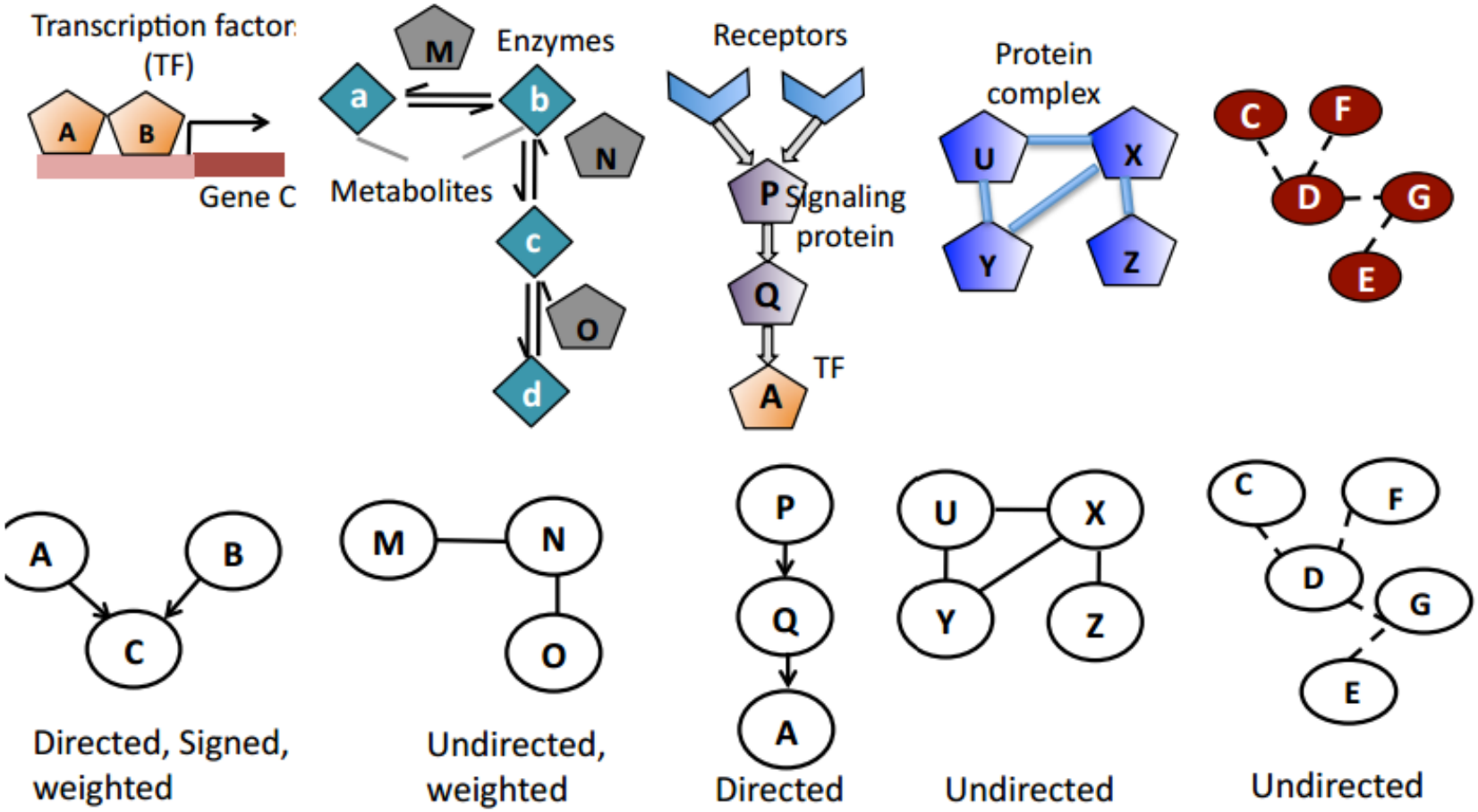
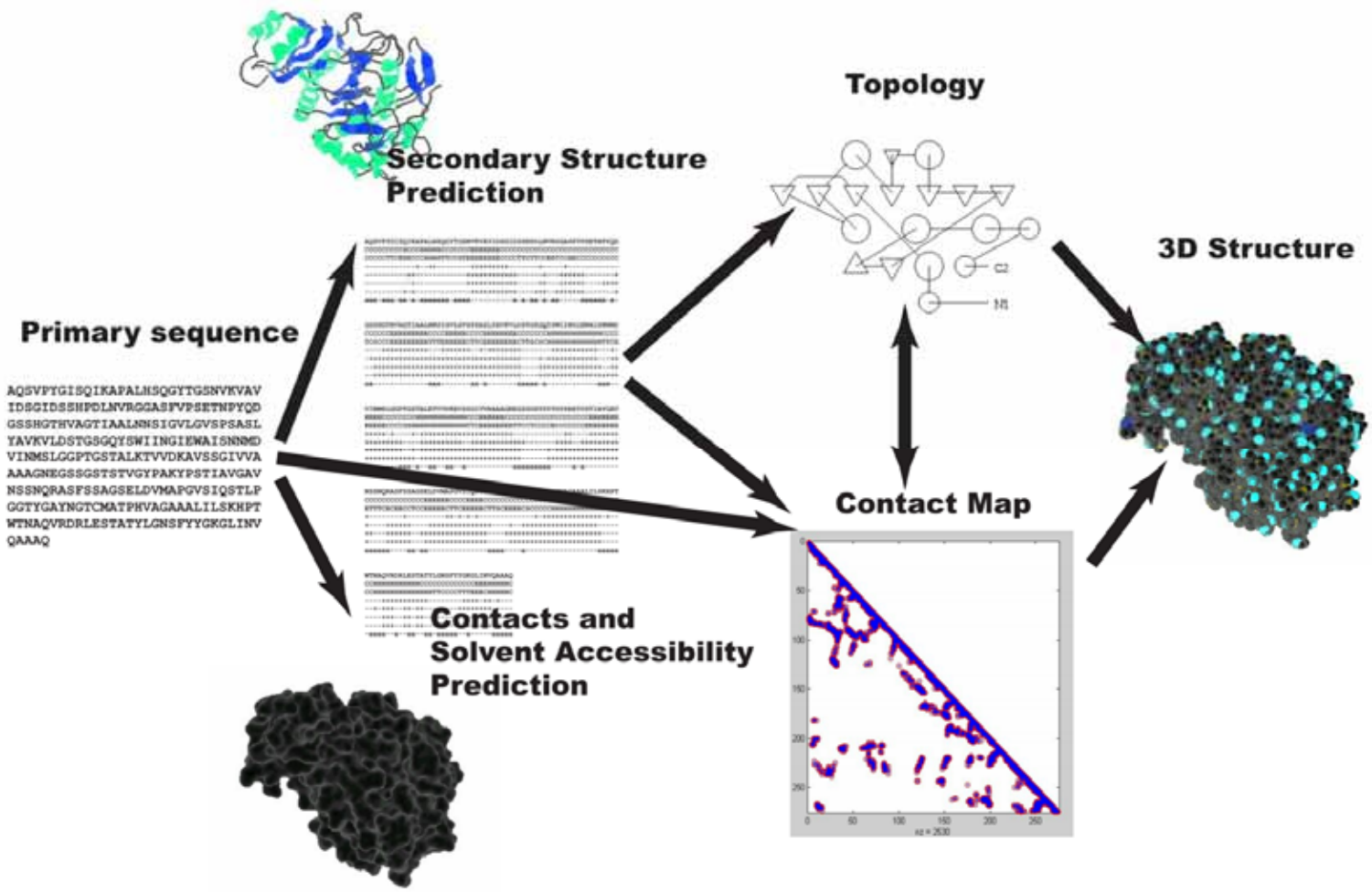


Image credit to Anna Goldenberg, Toronto

- Medicine is an extremely complex application domain – dealing most of the time with uncertainties -> **probable information!**
- When we have big data but little knowledge automatic ML can help to gain insight:
- **Structure learning and prediction in large-scale biomedical networks with probabilistic graphical models**
- If we have little data and deal with NP-hard problems we still need the human-in-the-loop

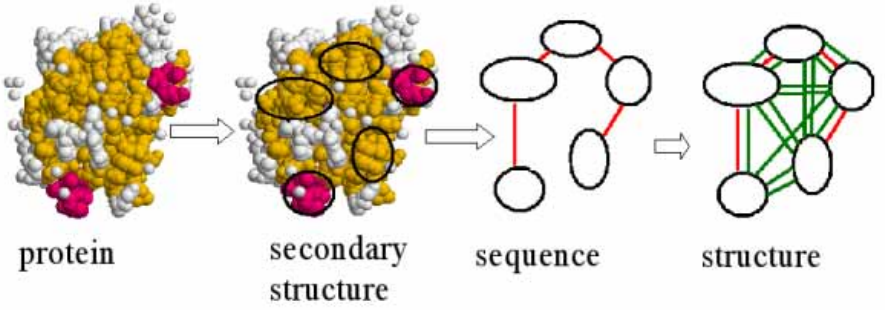


Baldi, P. & Pollastri, G. 2003. The principled design of large-scale recursive neural network architectures--dag-rnns and the protein structure prediction problem. The Journal of Machine Learning Research, 4, 575-602.

- Hypothesis: most biological functions involve the interactions between many proteins, and the complexity of living systems arises as a result of such interactions.
- In this context, the problem of inferring a global protein network for a given organism,
- - using all (genomic) data of the organism,
- is one of the main challenges in computational biology

Yamanishi, Y., Vert, J.-P. & Kanehisa, M. 2004. Protein network inference from multiple genomic data: a supervised approach. Bioinformatics, 20, (suppl 1), i363-i370.

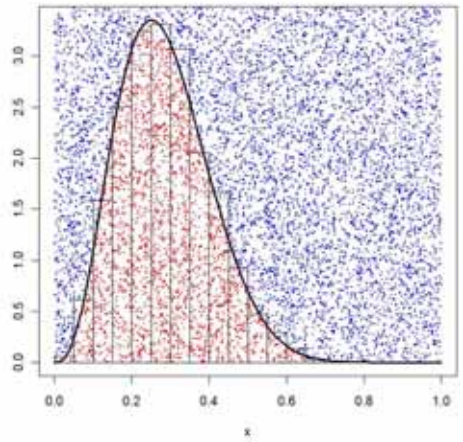
Borgwardt, K. M., Ong, C. S., Schönauer, S., Vishwanathan, S., Smola, A. J. & Kriegel, H.-P. 2005. Protein function prediction via graph kernels. Bioinformatics, 21, (suppl 1), i47-i56.

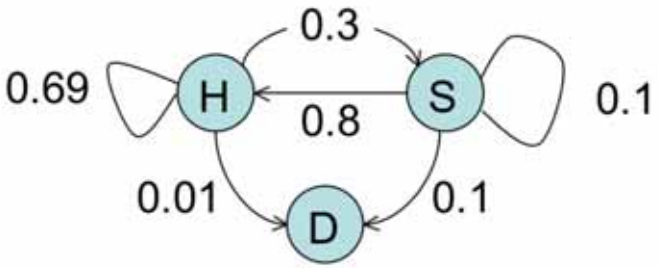


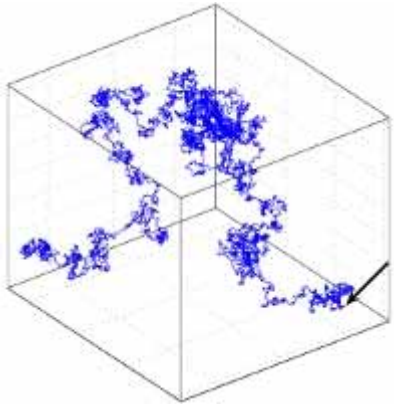
- Important for health informatics: Discovering relationships between biological components
- Unsolved problem in computer science:
- Can the graph isomorphism problem be solved in polynomial time?
  - So far, no polynomial time algorithm is known.
  - It is also not known if it is NP-complete
  - We know that subgraph-isomorphism is NP-complete

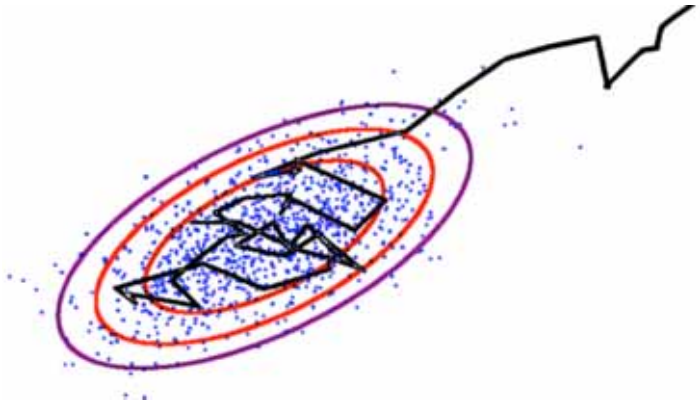
04 Markov Chain Monte Carlo (MCMC)

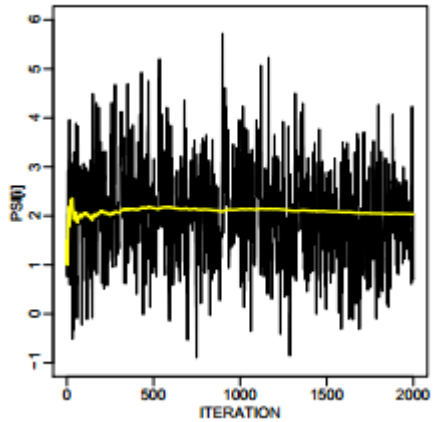
Monte Carlo Method (MC)
Monte Carlo Sampling
Markov Chains (MC)
MCMC
Metropolis-Hastings











- Often we want to calculate characteristics of a **high-dimensional** probability distribution … $p(\mathcal{D}|\theta)$

$$p(h|d) \propto p(\mathcal{D}|\theta) * p(h)$$

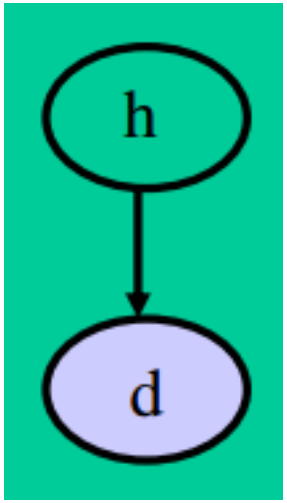
Posterior integration problem: (almost) all statistical inference can be deduced from the posterior distribution by calculating the appropriate sums, which involves an integration:

$$J = \int f(\theta) * p(\theta|\mathcal{D}) d\theta$$

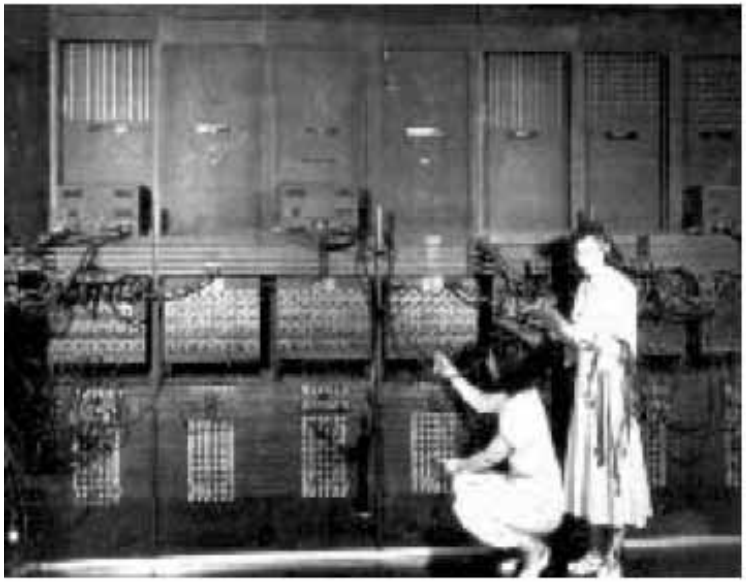
- **Statistical physics:** computing the partition function – this is evaluating the posterior probability of a hypothesis and this requires summing over all hypotheses … remember:

$$\mathcal{H} = \{H_1, H_2, ..., H_n\} \quad \forall(h, d)$$

$$P(h|d) = \frac{P(d|h) * P(h)}{\sum_{h' \in \mathcal{H}} P(d|h')P(h')}$$



What was the origin of MCMC ?

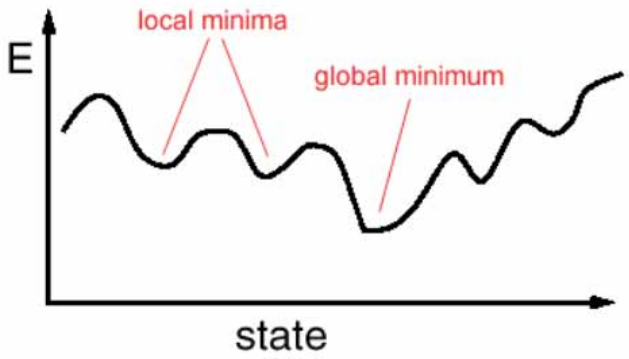


- Class of algorithms that rely on **repeated random sampling**
- Basic idea: using **randomness** to solve problems with high uncertainty (Laplace, 1781)
- For solving **multidimensional integrals** which would otherwise intractable
- For simulation of systems with **many dof**
- e.g. fluids, gases, particle collectives, **cellular structures** - see our last tutorial on Tumor growth simulation!

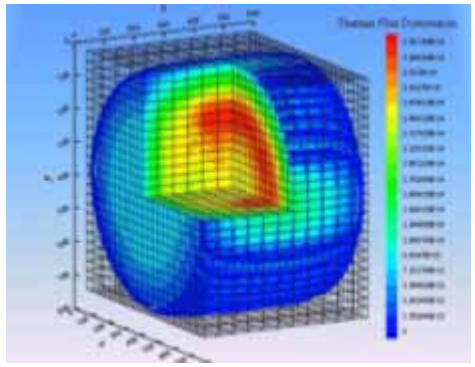
- for solving problems of probabilistic inference involved in developing computational models
- as a source of hypotheses about how the human mind might solve problems of inference
- For a function $f(x)$ and distribution $P(x)$, the expectation of f with respect to P is generally the average of f, when x is drawn from the probability distribution $P(x)$

$$\mathbb{E}_{p(x)}(f(x)) = \sum_{X} f(x)P(x)dx$$

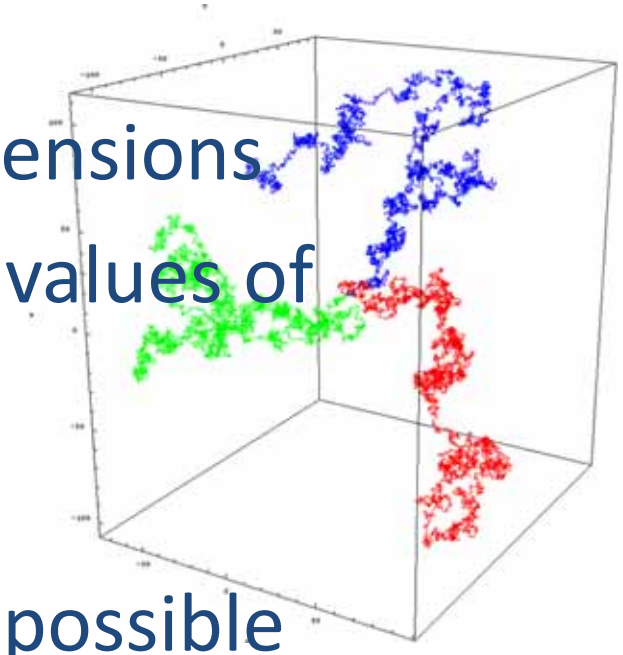
- Solving intractable integrals
- Bayesian statistics: **normalizing** constants, expectations, marginalization
- Stochastic Optimization



- Generalization of simulated annealing
- Monte Carlo expectation maximization (EM)



- Physical simulation
- estimating neutron diffusion time
- Computing expected utilities and best responses toward Nash equilibria
- Computing volumes in high-dimensions
- Computing eigen-functions and values of operators (e.g. Schrödinger)
- Statistical physics
- Counting many things as fast as possible



- Expectation of a function $f(x, y)$ with respect to a random variable x is denoted by $\mathbb{E}_x[f(x, y)]$
- In situations where there is no ambiguity as to which variable is being averaged over, this will be simplified by omitting the suffix, for instance $\mathbb{E}x$.
- If the distribution of x is conditioned on another variable z, then the corresponding conditional expectation will be written $\mathbb{E}_x[f(x)|z]$
- Similarly, the variance is denoted $var[f(x)]$, and for vector variables the covariance is written $cov[x, y]$

$$\underset{x}{\operatorname{argmax}}\, f(x)$$

Normalization: $$p(x|y) = \frac{p(y|x) * p(x)}{\int_X p(y|x) * p(x) dx}$$

Marginalization: $$p(x) = \int_Z p(x, z) dz$$

Expectation: $$\mathbb{E}_{p(x)}(f(x)) = \int_X f(x) p(x) dx$$

05 Metropolis-Hastings Algorithm

JOURNAL OF THE AMERICAN STATISTICAL ASSOCIATION

Number 247 SEPTEMBER 1949 *Volume 44*

THE MONTE CARLO METHOD

NICHOLAS METROPOLIS AND S. ULAM
Los Alamos Laboratory

We shall present here the motivation and a general description of a method dealing with a class of problems in mathematical physics. The method is, essentially, a statistical approach to the study of differential equations, or more generally, of integro-differential equations that occur in various branches of the natural sciences.

ALREADY in the nineteenth century a sharp distinction began to appear between two different mathematical methods of treating physical phenomena. Problems involving only a few particles were studied in classical mechanics, through the study of systems of ordinary differential equations. For the description of systems with very many particles, an entirely different technique was used, namely, the method of statistical mechanics. In this latter approach, one does not concentrate on the individual particles but studies the properties of *sets of particles*. In pure mathematics an intensive study of the properties of sets of points was the subject of a new field. This is the so-called theory of sets, the basic theory of integration, and the twentieth century development of the theory of probabilities prepared the formal apparatus for the use of such models in theoretical physics, i.e., description of properties of aggregates of points rather than of individual points and

Image Source:
http://www.manhattanprojectvoices.org/oral-histories/nicholas-metropolis-interview

THE JOURNAL OF CHEMICAL PHYSICS VOLUME 21, NUMBER 6 JUNE, 1953

Equation of State Calculations by Fast Computing Machines

NICHOLAS METROPOLIS, ARIANNA W. ROSENBLUTH, MARSHALL N. ROSENBLUTH, AND AUGUSTA H. TELLER,
Los Alamos Scientific Laboratory, Los Alamos, New Mexico

AND

EDWARD TELLER,* *Department of Physics, University of Chicago, Chicago, Illinois*

(Received March 6, 1953)

A general method, suitable for fast computing machines, for investigating such properties as equations of state for substances consisting of interacting individual molecules is described. The method consists of a modified Monte Carlo integration over configuration space. Results for the two-dimensional rigid-sphere system have been obtained on the Los Alamos MANIAC and are presented here. These results are compared to the free volume equation of state and to a four-term virial coefficient expansion.

I. INTRODUCTION

THE purpose of this paper is to describe a general method, suitable for fast electronic computing machines, of calculating the properties of any substance which may be considered as composed of interacting individual molecules. Classical statistics is assumed, only two-body forces are considered, and the potential field of a molecule is assumed spherically symmetric. These are the usual assumptions made in theories of liquids. Subject to the above assumptions, the method is not restricted to any range of temperature or density. This paper will also present results of a preliminary two-dimensional calculation for the rigid-sphere system. Work on the two-dimensional case with a Lennard-Jones potential is in progress and will be reported in a later paper. Also, the problem in three dimensions is being investigated.

* Now at the Radiation Laboratory of the University of California, Livermore, California.

II. THE GENERAL METHOD FOR AN ARBITRARY POTENTIAL BETWEEN THE PARTICLES

In order to reduce the problem to a feasible size for numerical work, we can, of course, consider only a finite number of particles. This number N may be as high as several hundred. Our system consists of a square† containing N particles. In order to minimize the surface effects we suppose the complete substance to be periodic, consisting of many such squares, each square containing N particles in the same configuration. Thus we define d_{AB}, the minimum distance between particles A and B, as the shortest distance between A and any of the particles B, of which there is one in each of the squares which comprise the complete substance. If we have a potential which falls off rapidly with distance, there will be at most one of the distances AB which can make a substantial contribution; hence we need consider only the minimum distance d_{AB}.

† We will use the two-dimensional nomenclature here since it is easier to visualize. The extension to three dimensions is obvious.

Metropolis, N., Rosenbluth, A. W., Rosenbluth, M. N., Teller, A. H. & Teller, E. 1953. Equation of State Calculations by Fast Computing Machines. The Journal of Chemical Physics, 21, (6), 1087-1092, doi:10.1063/1.1699114.

Hastings, W. K. 1970. Monte Carlo sampling methods using Markov chains and their applications. Biometrika, 57, (1), 97-109.

Biometrika (1970), **57**, 1, *p.* 97
Printed in Great Britain

Monte Carlo sampling methods using Markov chains and their applications

BY W. K. HASTINGS
University of Toronto

SUMMARY

A generalization of the sampling method introduced by Metropolis *et al.* (1953) is presented along with an exposition of the relevant theory, techniques of application and methods and difficulties of assessing the error in Monte Carlo estimates. Examples of the methods, including the generation of random orthogonal matrices and potential applications of the methods to numerical problems arising in statistics, are discussed.

1. INTRODUCTION

For numerical problems in a large number of dimensions, Monte Carlo methods are often more efficient than conventional numerical methods. However, implementation of the Monte Carlo methods requires sampling from high dimensional probability distributions and this may be very difficult and expensive in analysis and computer time. General methods for sampling from, or estimating expectations with respect to, such distributions are as follows.

(i) If possible, factorize the distribution into the product of one-dimensional conditional distributions from which samples may be obtained.

(ii) Use importance sampling, which may also be used for variance reduction. That is, in order to evaluate the integral

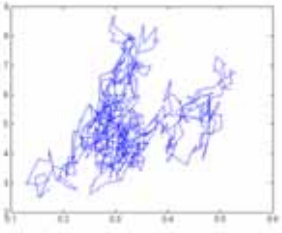
$$J = \int f(x)p(x)dx = E_p(f),$$

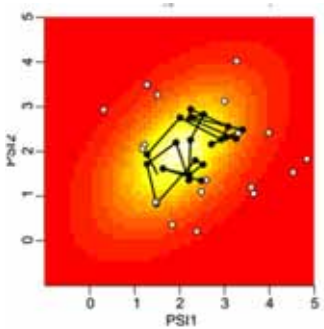
where $p(x)$ is a probability density function, instead of obtaining independent samples $x_1, \ldots, x_N$ from $p(x)$ and using the estimate $\hat{J}_1 = \Sigma f(x_i)/N$, we instead obtain the sample from

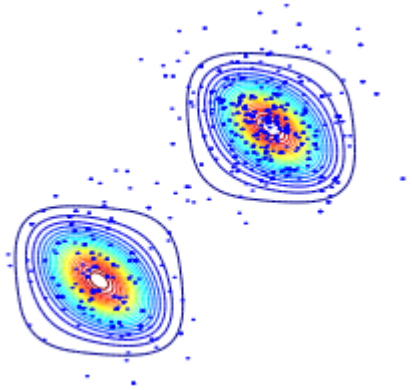
So what is the MH-algorithm doing ?

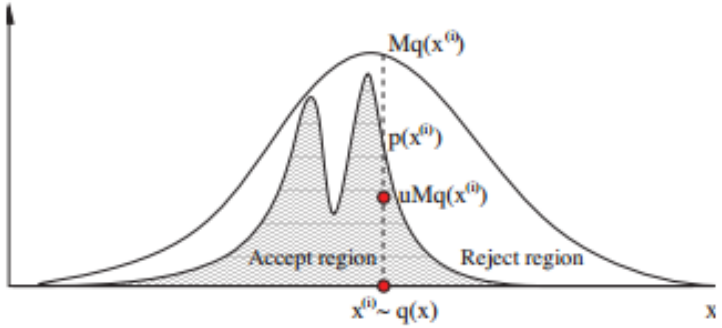
Barber, D. 2012. Bayesian reasoning and machine learning, Cambridge, Cambridge University

1: Choose a starting point x^1.
2: **for** $i = 2$ to L **do**
3: Draw a candidate sample x^{cand} from the proposal $\tilde{q}(x'|x^{l-1})$.
4: Let $a = \frac{\tilde{q}(x^{l-1}|x^{cand})p(x^{cand})}{\tilde{q}(x^{cand}|x^{l-1})p(x^{l-1})}$
5: **if** $a \geq 1$ **then** $x^l = x^{cand}$
6: **else**
7: draw a random value u uniformly from the unit interval $[0, 1]$.
8: **if** $u < a$ **then** $x^l = x^{cand}$
9: **else**
10: $x^l = x^{l-1}$
11: **end if**
12: **end if**
13: **end for**









- Importance sampling is a technique to approximate averages with respect to an intractable distribution $p(x)$.
- The term 'sampling' is arguably a misnomer since the method does not attempt to draw samples from $p(x)$.
- Rather the method draws samples from a simpler importance distribution $q(x)$ and then reweights them
- such that averages with respect to $p(x)$ can be approximated using the samples from $q(x)$.

- The Gibbs Sampler is an interesting special case of MH:

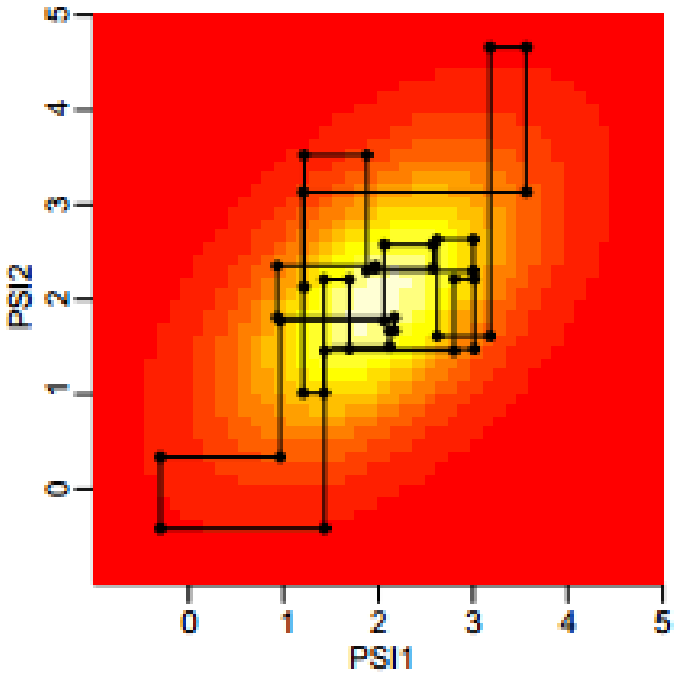
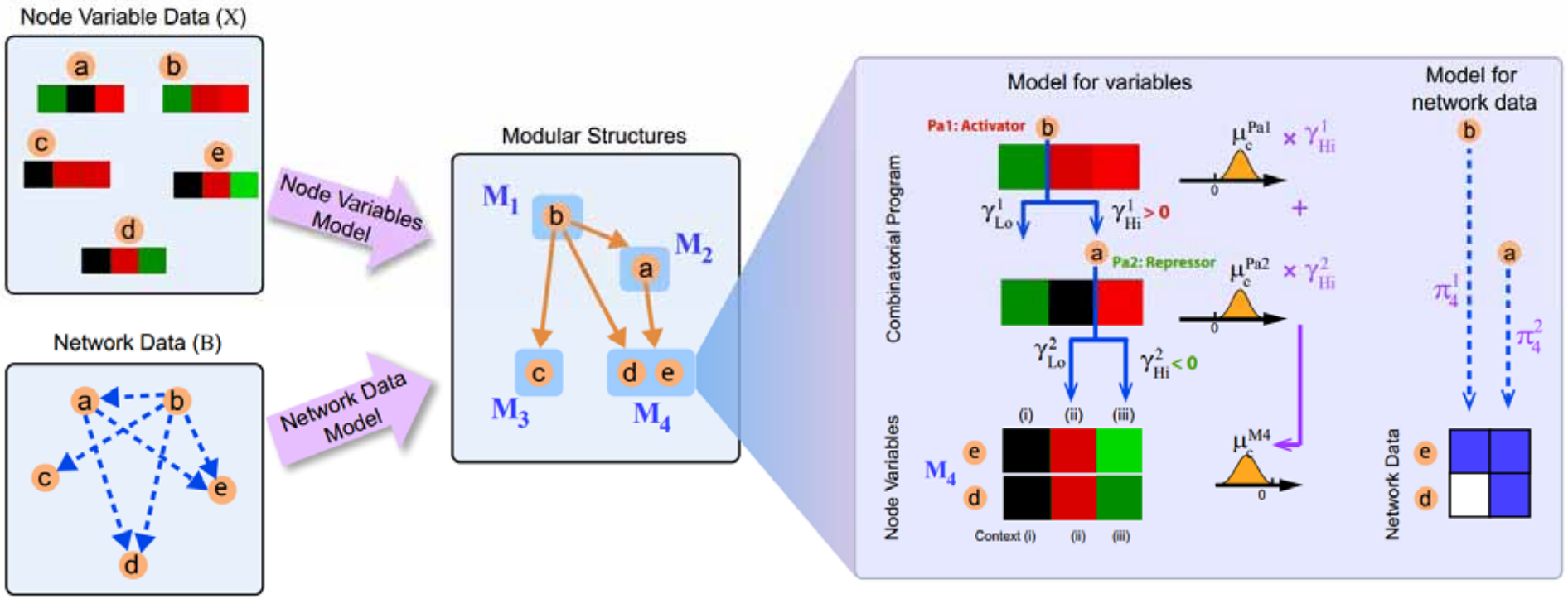
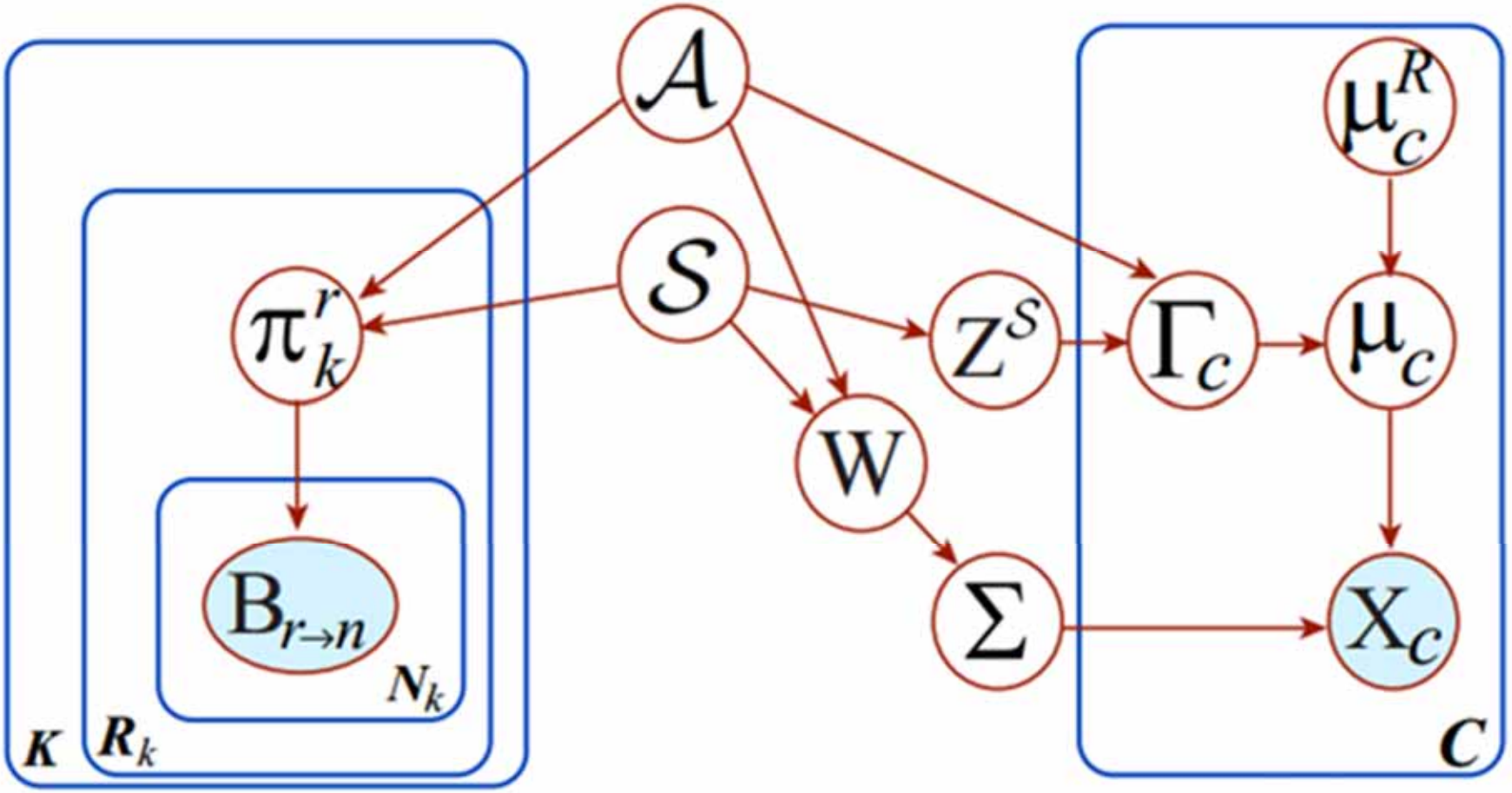


Image Source: Peter Mueller, Anderson Cancer Center

How to learn modular structures from Network Data ?



Elham Azizi, Edoardo M. Airoldi & James E. Galagan. Learning Modular Structures from Network Data and Node Variables. In: Xing, Eric P. & Jebara, Tony, eds. Proceedings of the 31st International Conference on Machine Learning (ICML), 2014 Beijing. JMLR, 1440-1448.

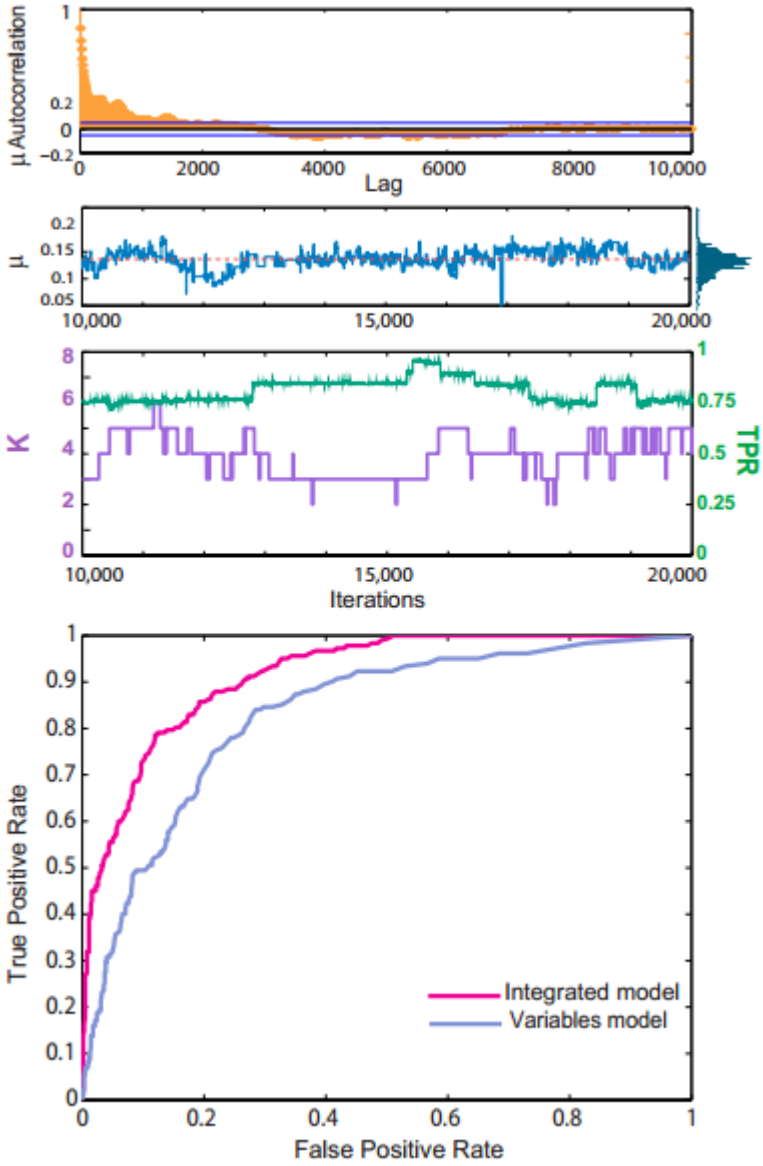


Elham Azizi, Edoardo M. Airoldi & James E. Galagan. Learning Modular Structures from Network Data and Node Variables. In: Xing, Eric P. & Jebara, Tony, eds. Proceedings of the 31st International Conference on Machine Learning (ICML), 2014 Beijing. JMLR, 1440-1448.

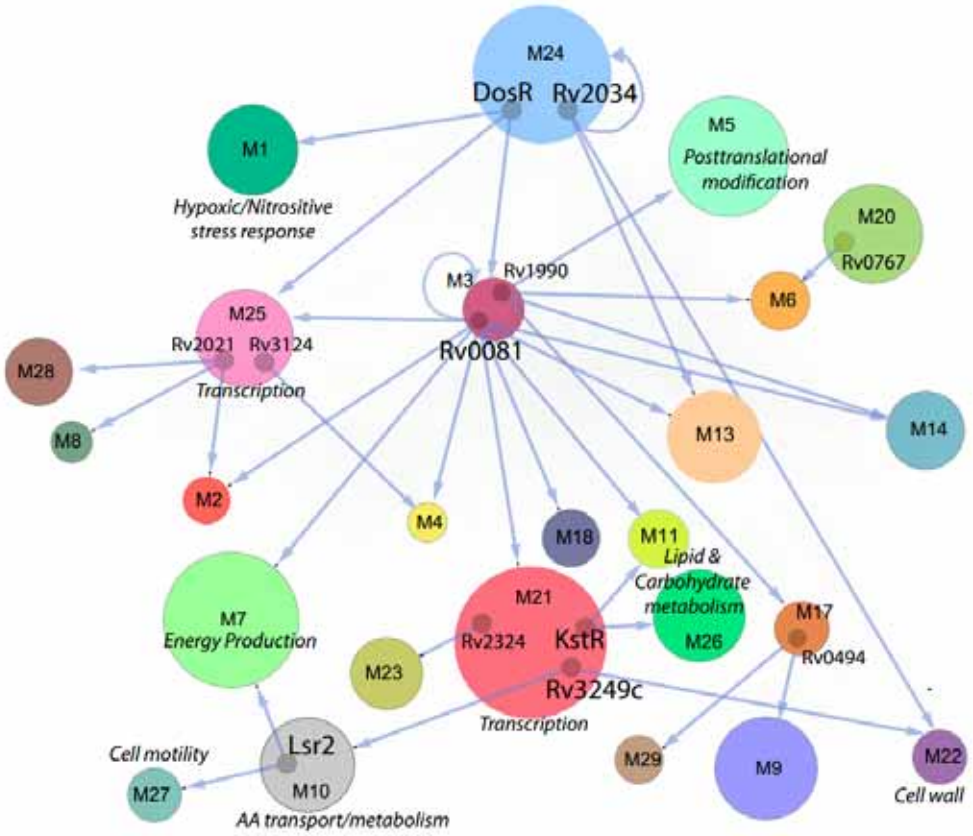
How can you estimate the parameters using RJMCMC ?

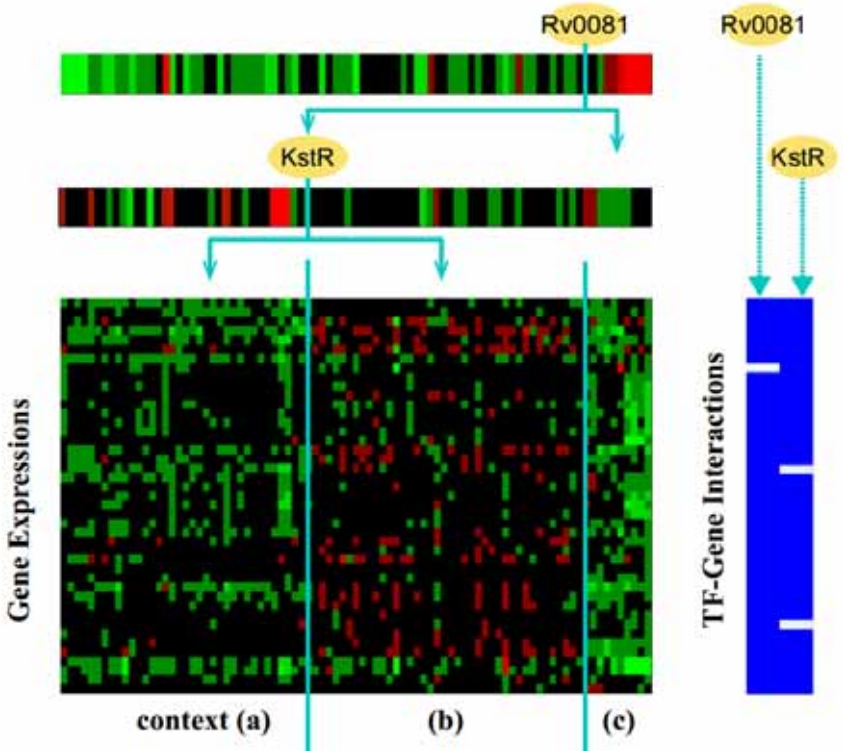
Algorithm 1 RJMCMC for sampling parameters

Inputs:
Node Variables Data $\mathbf{X}$
Network Data $\mathbf{B}$
for iterations $j = 1$ **to** J **do**
  Sample $\mathcal{A}^{(j+1)}$ given $\mathcal{A}^{(j)}$ using Alg 2 in (Azizi et al., 2014)
  Sample $\mathcal{S}^{(j+1)}$ given $\mathcal{S}^{(j)}$ using Alg 3 in (Azizi et al., 2014)
  for modules $k = 1$ **to** $K^{(j)}$ **do**
    Propose $w_k^{(j+1)} \sim \mathcal{N}(w_k^{(j)}, I)$
    Accept with probability P_{mh}; update $\Sigma^{(j+1)}$
    for parents $r = 1$ **to** R_k **do**
      Propose $z_k^{r(j+1)} \sim \mathcal{N}(z_k^{r(j)}, I)$; accept with P_{mh}
      Propose $\pi_k^{r(j+1)} \sim \mathcal{N}(\pi_k^{r(j)}, I)$; accept with P_{mh}
    end for
  end for
  for condition $c = 1$ **to** C **do**
    Propose $\boldsymbol{\mu}_c^{\mathbf{R}(j+1)} \sim \mathcal{N}(\boldsymbol{\mu}_c^{\mathbf{R}(j)}, I)$; accept with P_{mh}
    Propose $\boldsymbol{\gamma}_c^{\mathbf{R}(j+1)} \sim \mathcal{N}(\boldsymbol{\gamma}_c^{\mathbf{R}(j)}, I)$; accept with P_{mh}
  end for
end for

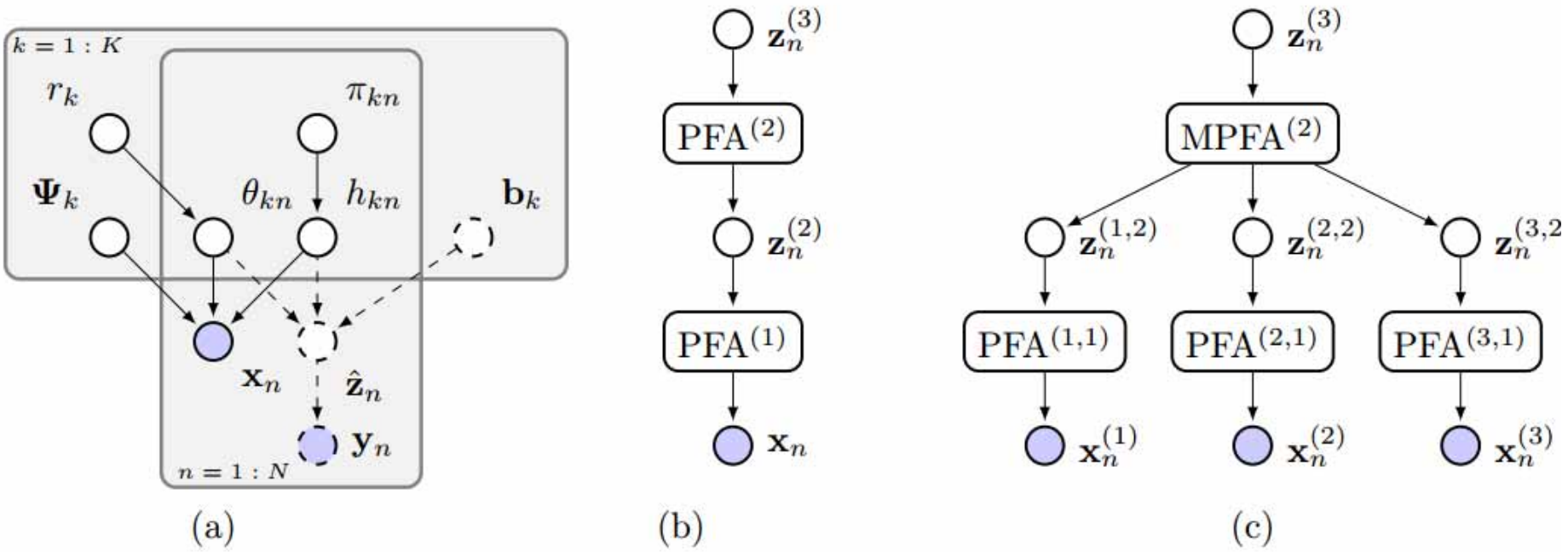


Azizi, E., Airoldi, E. M. & Galagan, J. E. 2014. Learning Modular Structures from Network Data and Node Variables. Proceedings of the 31st International Conference on Machine Learning (ICML). Beijing: JMLR. 1440-1448.



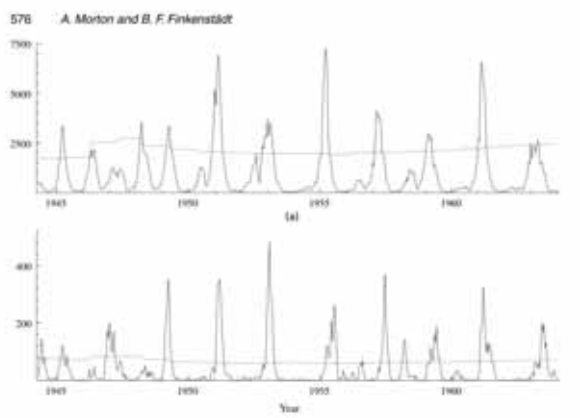


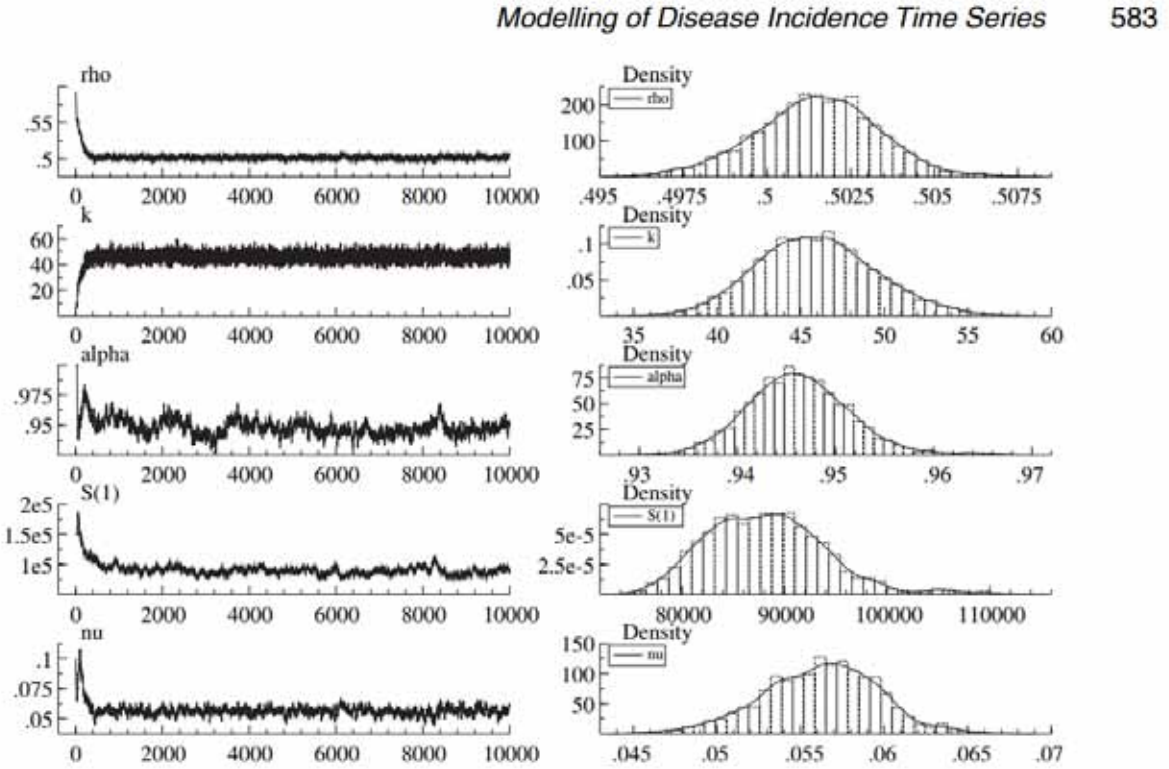
Azizi, E., Airoldi, E. M. & Galagan, J. E. 2014. Learning Modular Structures from Network Data and Node Variables. Proceedings of the 31st International Conference on Machine Learning (ICML). Beijing: JMLR. 1440-1448.



Henao, R., Lu, J. T., Lucas, J. E., Ferranti, J. & Carin, L. 2016. Electronic health record analysis via deep poisson factor models. Journal of Machine Learning Research JMLR, 17, 1-32.

Discrete time modelling of disease incidence time series by MCMC



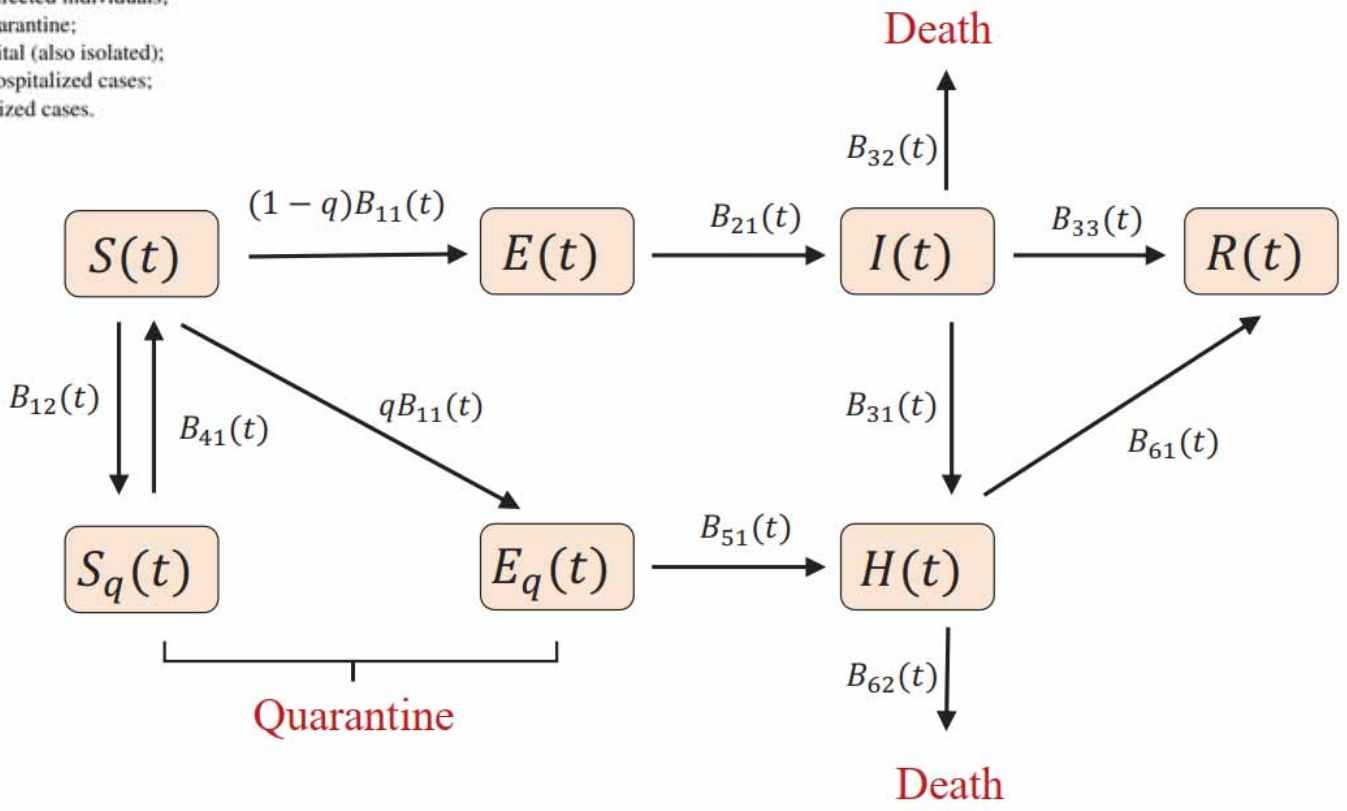


Alexander Morton & Bärbel F. Finkenstädt 2005. Discrete time modelling of disease incidence time series by using Markov chain Monte Carlo methods. Journal of the Royal Statistical Society: Series C (Applied Statistics), 54, (3), 575-594, doi:10.1111/j.1467-9876.2005.05366.x

Discrete-time stochastic epidemic model of COVID-19

- $B_{11}(t)$ is the number of susceptible individuals who become newly infected;
- $B_{12}(t)$ is the number of quarantined susceptible individuals who have contact with infected individuals but are not infected;
- $B_{21}(t)$ is the number of new cases with symptom onset;
- $B_{31}(t)$ is the number of new confirmed and admitted patients;
- $B_{32}(t)$ is the number of new death from infected individuals;
- $B_{33}(t)$ is the number of newly recovered from infected individuals;
- $B_{41}(t)$ is the number of people released from quarantine;
- $B_{51}(t)$ is the number of people admitted to hospital (also isolated);
- $B_{61}(t)$ is the number of newly recovered from hospitalized cases;
- $B_{62}(t)$ is the number of new death from hospitalized cases.

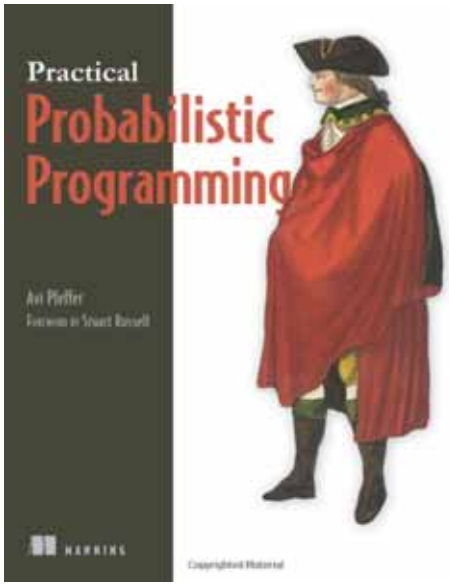
Sha He, Sanyi Tang & Libin Rong 2020. A discrete stochastic model of the COVID-19 outbreak: Forecast and control. Journal of Mathematical Biosciences & Engineering, 17, (4), 2792-2804, doi:10.3934/mbe.2020153 https://www.aimspress.com/MBE/2020/4/2792 (Online open available)

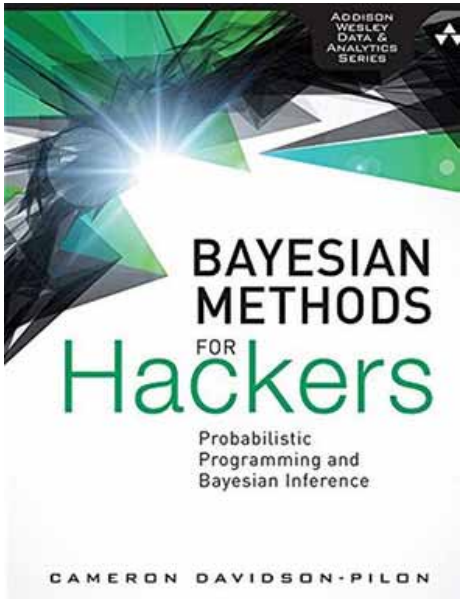


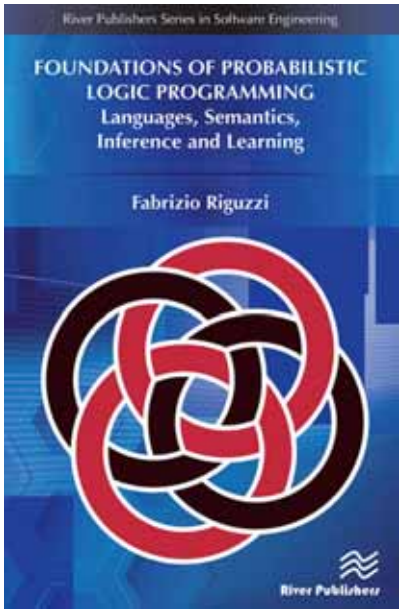
$$L(B_{11}(t), B_{12}(t), B_{21}(t), B_{31}(t), B_{32}(t), B_{33}(t), B_{41}(t), B_{51}(t), B_{61}(t), B_{62}(t)|\Theta) = \prod_{t=0}^{T^n} g_{i,j}(B_{ij}(t)|\cdot)$$

06 Probabilistic Programming

Book recommendations







Avi Pfeffer 2016. Practical probabilistic programming, Shelter Island (NY), Manning.

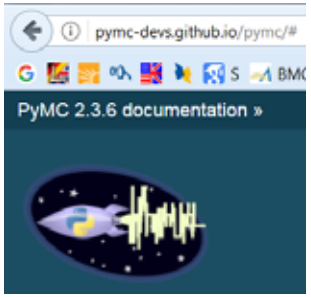
Cameron Davidson-Pilon 2015. Bayesian methods for hackers: probabilistic programming and Bayesian inference, Addison-Wesley Professional.

Fabrizio Riguzzi 2018. Foundations of Probabilistic Logic Programming, River Publishers.

Arnaud N. Fadja & Fabrizio Riguzzi 2017. Probabilistic Logic Programming in Action. In: Holzinger, Andreas, Goebel, Randy, Ferri, Massimo & Palade, Vasile (eds.) Towards Integrative Machine Learning and Knowledge Extraction: BIRS Workshop, Banff, AB, Canada, July 24-26, 2015, Revised Selected Papers. Cham: Springer, pp. 89-116, doi:10.1007/978-3-319-69775-8_5.

- Probabilistic thinking is a valuable tool for decision making
- Overcoming uncertainties is the huge success currently in machine learning (and for AI ;-)
- Probabilistic reasoning is a versatile tool
- PPLs are domain specific languages that use probabilistic models and the methods to make inferences in those models
- The “magic” is in combining “probability methods” with “representational power”

- C → Probabilistic-C
- Scala → Figaro
- Scheme → Church
- Excel → Tabular
- Prolog → Problog
- Javascript → webPP
- → Venture
- **Python → PyMC**



PyMC3

PyMC Pythonic Markov chain Monte Carlo

Medical Example

Sequence	Outcome
CGTCGGAGGTACATGATTGGAAGAAAACCT	Y
GCGCCTTTGCACATCTCTTAATCTCAGTCA	X
TTAAAATAGCAGAGACACTTCTACTGATAC	Y
CCAAGAGCCTCGTAATTAAGTATTGCAATA	Y
TTATGACGTCGTTTCGAGTGGATTTGTCTT	X
...	...

1

- Simple example: Nucleotide "A" may follow nucleotide "T" in the sequences more frequently for outcome X than for outcome Y,

$$P(A \mid T, X) > P(A \mid T, Y)$$

2

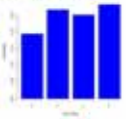
- Compute maximum *a posteriori* estimates of the probabilities:

```
from pymc import MAP, Model
model = Model({'f_x' : f_x, 'prob_dist' : prob_dist})
M = MAP(model)
M.fit()    # Nelder-Mead Optimization
```

- The MAP estimates are now contained in th M.prob_dist value:

```
>>> print M.prob_dist.value
[ 0.19472259  0.24842748  0.25265728]
```

Posterior Distribution of the Nucleotides



$$P(\theta \mid D) = \frac{P(D \mid \theta) \cdot P(\theta)}{P(D)}$$

6

- Specify the prior distribution:

```
import numpy as np
from pymc import Dirichlet  # conjugate prior
alpha = np.array([30.0,25.0,20.0,25.0])
prob_dist = Dirichlet('prob_dist', alpha)
```

Prior Distribution of the Nucleotides

3

$$P(\theta \mid D) = \frac{P(D \mid \theta) \cdot P(\theta)}{P(D)}$$

- Specify the value to maximize using numerical simulation, as well as the expected form of the posterior distribution:

```
from pymc import Categorical
f_x = Categorical('cat', prob_dist, value=exp_data, observed=True)
```

5

$$P(\theta \mid D) = \frac{P(D \mid \theta) \cdot P(\theta)}{P(D)}$$

- Specify the experimental data:

```
exp_data = np.array([1, 1, 3, 2, 2, 1, 0, ...])
```

Experimental Data

Observation #	Nucleotide
1	1
2	1
3	3
4	2
5	2
6	1
7	0

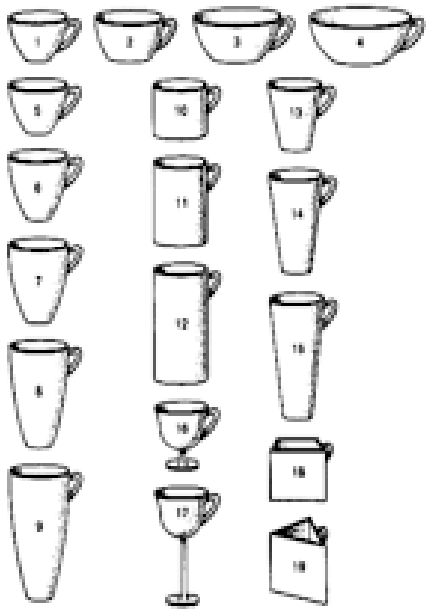
$$P(\theta \mid D) = \frac{P(D \mid \theta) \cdot P(\theta)}{P(D)}$$

4

Image Source: Dan Williams, Life Technologies, Austin TX

Digression on Concept Learning

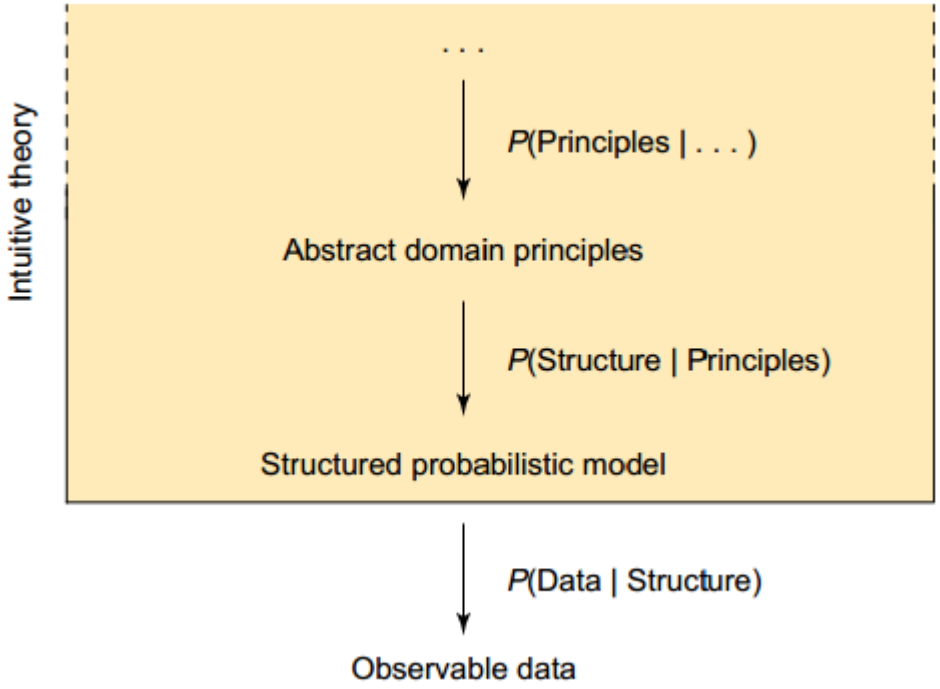
- Bruner, Goodnow, and Austin (1956) published “A Study of Thinking”, which became a landmark in cognitive science and has much influence on machine learning.
  - Rule-Based Categories
  - A concept specifies conditions for membership



Jerome S. Bruner, Jacqueline J. Goodnow & George A. Austin 1986. A Study of Thinking, Transaction Books.

- which is highly relevant for ML research, concerns the factors that determine the subjective difficulty of concepts:
- Why are some concepts psychologically extremely simple and easy to learn,
- while others seem to be extremely difficult, complex, or even incoherent?
- These questions have been studied since the 1960s but are still unanswered …

Feldman, J. 2000. Minimization of Boolean complexity in human concept learning. Nature, 407, (6804), 630-633, doi:10.1038/35036586.



$$P(h|x,T) = \frac{P(x|h,T)P(h|T)}{\sum_{h' \in H_T} P(x|h',T)P(h'|T)}$$

Joshua B. Tenenbaum, Thomas L. Griffiths & Charles Kemp 2006. Theory-based Bayesian models of inductive learning and reasoning. Trends in cognitive sciences, 10, (7), 309-318, doi:10.1016/j.tics.2006.05.009.

How does our mind get so much out of it ?

Salakhutdinov, R., Tenenbaum, J. & Torralba, A. 2012. One-shot learning with a hierarchical nonparametric Bayesian model. Journal of Machine Learning Research, 27, 195-207.

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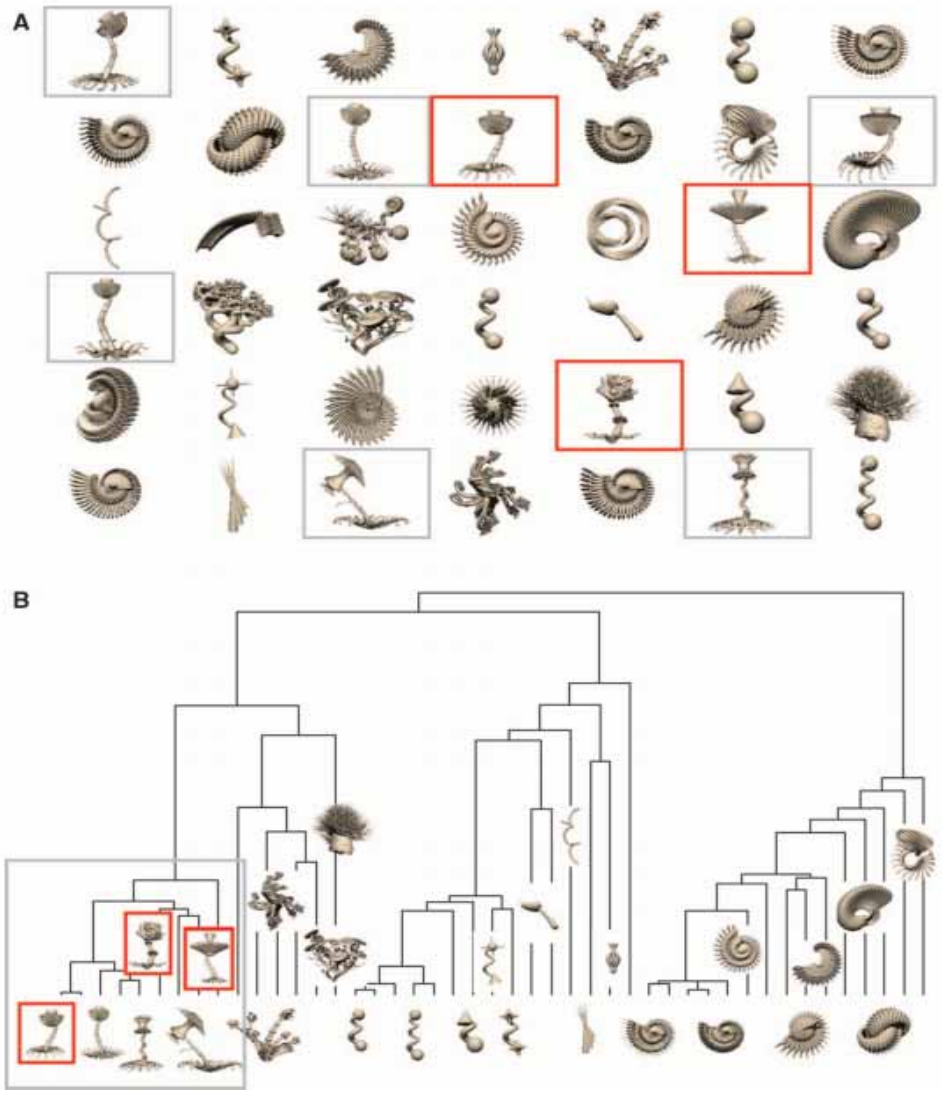
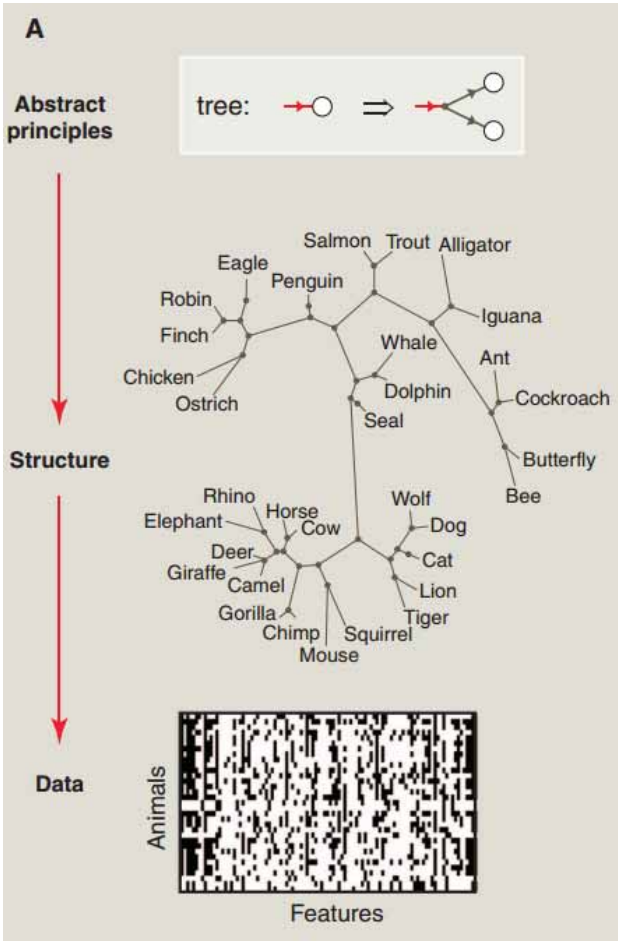
Salakhutdinov, R., Tenenbaum, J. & Torralba, A. 2012. One-shot learning with a hierarchical nonparametric Bayesian model. Journal of Machine Learning Research, 27, 195-207.

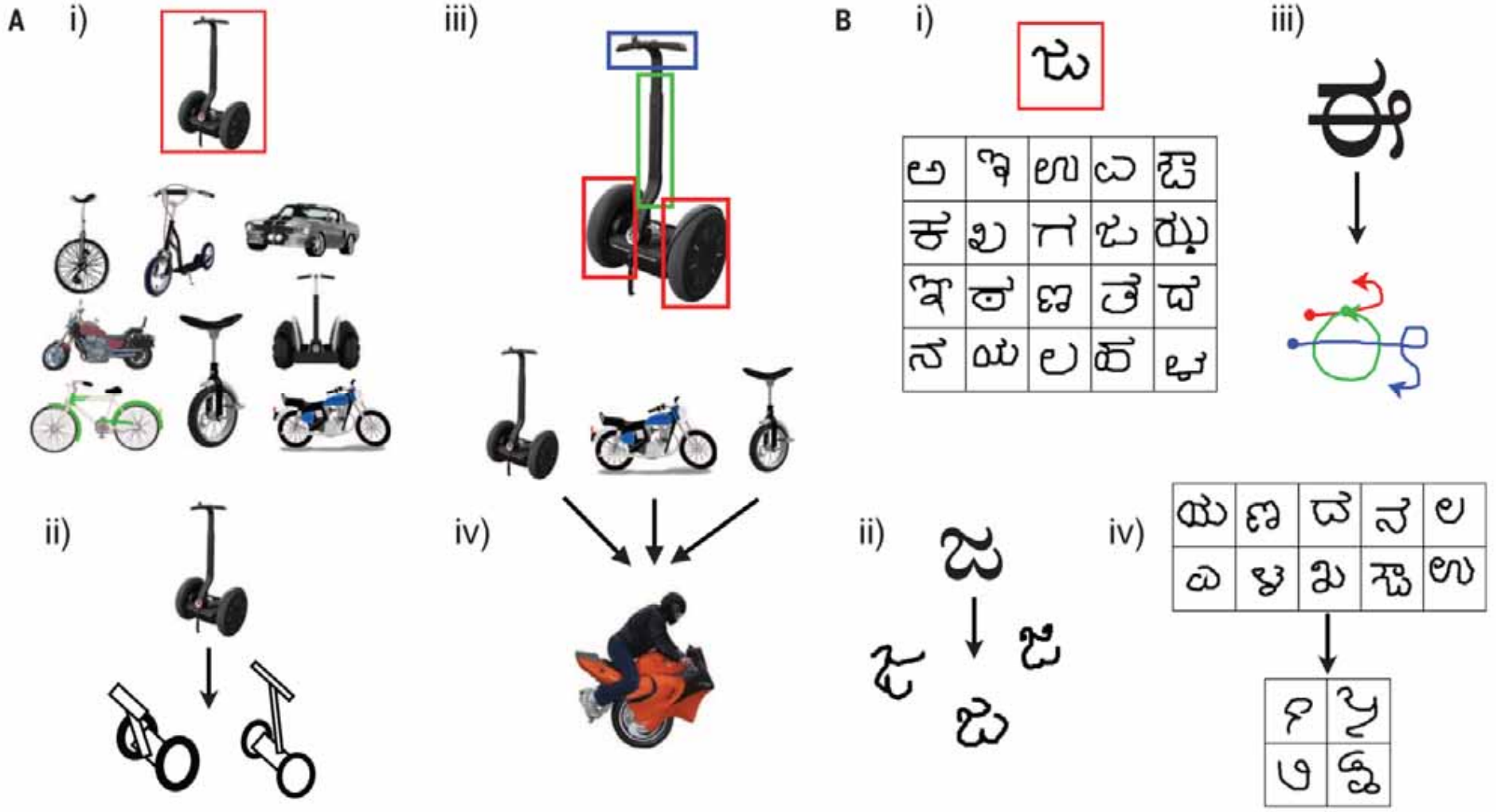


$$P(h|d) = \frac{P(d|h)P(h)}{\sum_{h' \in H} P(d|h')P(h')} \propto P(d|h)P(h)$$



Tenenbaum, J. B., Kemp, C., Griffiths, T. L. & Goodman, N. D. 2011. How to grow a mind: Statistics, structure, and abstraction. Science, 331, (6022), 1279-1285.

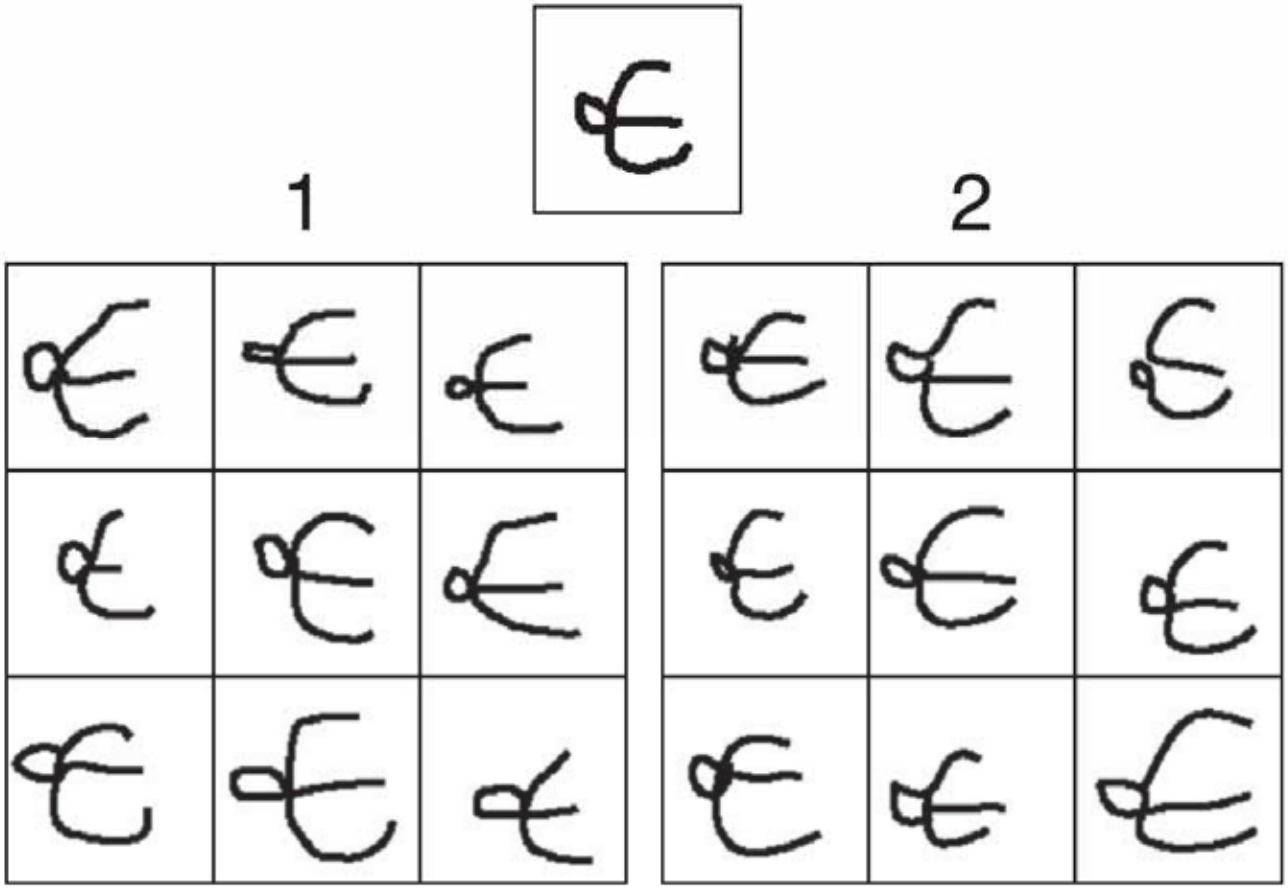
What is probabilistic program induction ?



Brenden M. Lake, Ruslan Salakhutdinov & Joshua B. Tenenbaum 2015. Human-level concept learning through probabilistic program induction. Science, 350, (6266), 1332-1338, doi:10.1126/science.aab3050.

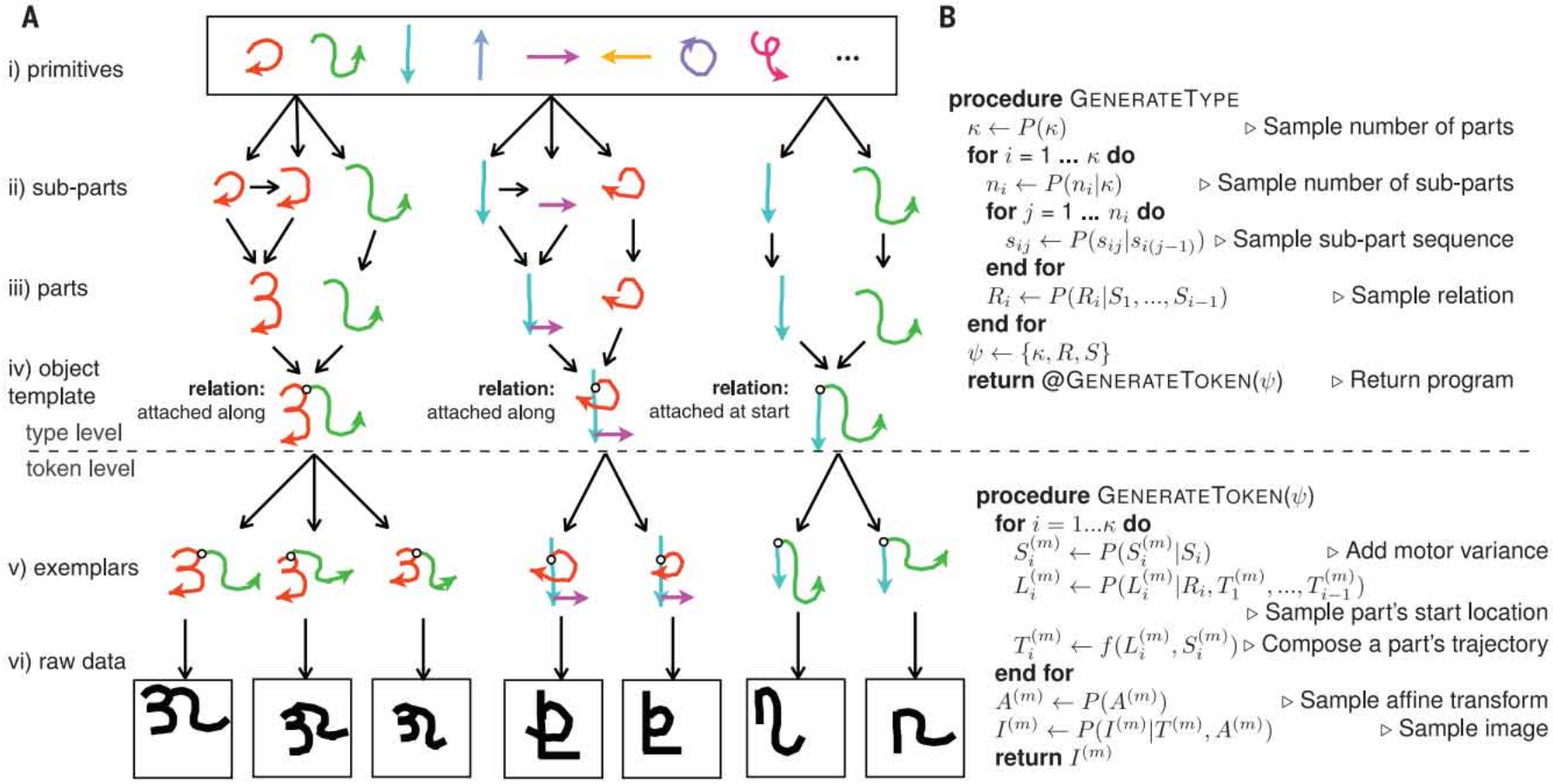
- **Deductive Reasoning** = Hypothesis > Observations > Logical Conclusions (general → specific – proven correctness)
  - DANGER: Hypothesis must be correct! DR defines whether the truth of a conclusion can be determined for that rule, based on the truth of premises: A=B, B=C, therefore A=C
- **Inductive reasoning** = makes broad generalizations from specific observations (specific → general – not proven correctness)
  - DANGER: allows a conclusion to be false if the premises are true
  - generate hypotheses and use DR for answering specific questions
- **Abductive reasoning** = inference = to get the best explanation from an incomplete set of preconditions.
  - Given a true conclusion and a rule, it attempts to select some possible premises that, if true also, may support the conclusion, though not uniquely.
  - Example: "When it rains, the grass gets wet. The grass is wet. Therefore, it might have rained." This kind of reasoning can be used to develop a hypothesis, which in turn can be tested by additional reasoning or data.

Drawn by Human or Machine Learning Algorithm ?



Brenden M. Lake, Ruslan Salakhutdinov & Joshua B. Tenenbaum 2015. Human-level concept learning through probabilistic program induction. Science, 350, (6266), 1332-1338, doi:10.1126/science.aab3050.

What can a Bayesian program learning (BPL) framework do ?



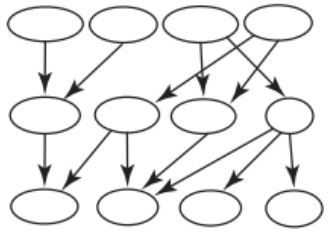
Brenden M. Lake, Ruslan Salakhutdinov & Joshua B. Tenenbaum 2015. Human-level concept learning through probabilistic program induction. Science, 350, (6266), 1332-1338, doi:10.1126/science.aab3050.

What can we do with graphical models ?

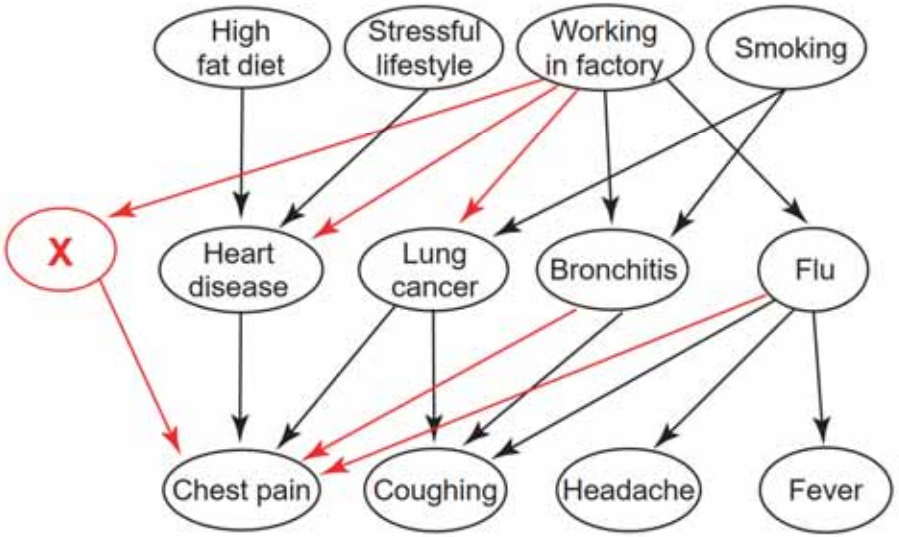
Principles

Classes: {R, D, S} (Risks, Diseases, Symptoms)
Causal laws: R → D, D → S

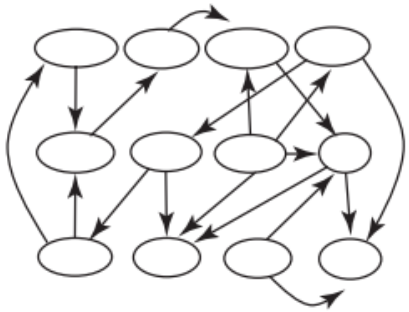
Classes: {R, D, S}
Causal laws: R → D, D → S



Structure



Classes: {C}
Causal laws: C → C

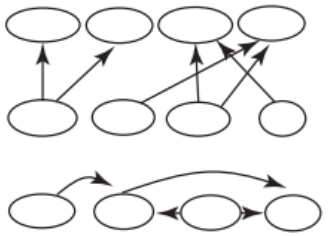


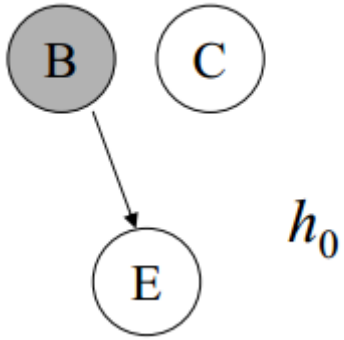
Joshua B. Tenenbaum, Thomas L. Griffiths & Charles Kemp 2006. Theory-based Bayesian models of inductive learning and reasoning. Trends in cognitive sciences, 10, (7), 309-318, doi:10.1016/j.tics.2006.05.009.

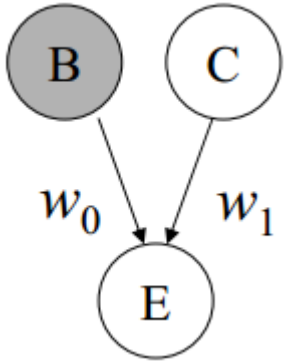
Data

Patient 1: Stressful lifestyle
Chest Pain
Patient 2: Smoking
Coughing
Patient 3: Working in factory
Chest Pain
. . .

Classes: {R, D, S}
Causal laws: D → R, S → S







- Cognition as probabilistic inference
  - Visual perception, language acquisition, motor learning, associative learning, memory, attention, categorization, reasoning, causal inference, decision making, theory of mind
- Learning concepts from examples
- Learning causation from correlation
- Learning and applying intuitive theories (balancing complexity vs. fit)

Thank you!