

Causal discovery: additional approaches

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Learning models from data: a Bayesian approach

- Fundamental concepts

- Equivalence between DAGs

- The GES algorithm

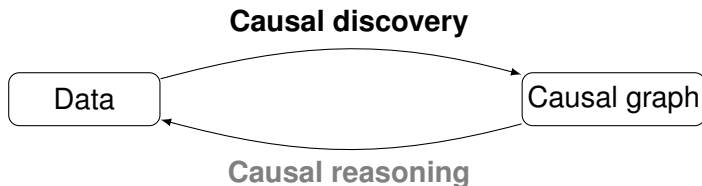
Granger causality

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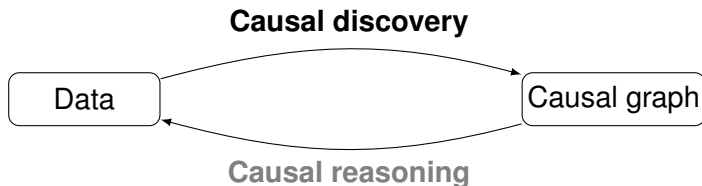
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What is our concern?



Infer a causal graph from observed data following a Bayesian approach

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Infer a causal graph from observed data following a Bayesian approach

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Bayesian network models and DAG models

Parametrized Bayesian-network model A pair (\mathcal{G}, θ) where $\mathcal{G} = (\mathbf{V}, \mathbf{E})$ is a DAG in which nodes correspond to variables and θ is a set of parameter values that specify all conditional probability distributions ($\theta_i \subset \theta$ subset of parameter values that define the conditional probability if X_i given its parents in \mathcal{G})

$$P(X_1 = x_1, \dots, X_n = x_n) = \prod_{i=1}^n P(X_i = x_i | \mathbf{Pa}_i^{\mathcal{G}} = \mathbf{pa}_i^{\mathcal{G}}, \theta_i) \quad (1)$$

- ▶ The structure \mathcal{G} is a DAG model that represents the independence constraints that must hold in any distribution represented by the network
- ▶ The set of independence constraints imposed by \mathcal{G} are represented by the Markov conditions (independence of non-descendants given parents)

Complete network

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Complete network

Remark: Markov conditions vs faithfulness

absent edge \Rightarrow conditional independence



see conditional dependence \Rightarrow infer edge

edge \Rightarrow conditional dependence



see conditional independence \Rightarrow absent edge

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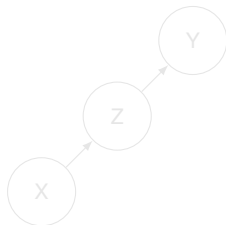
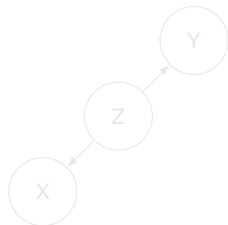


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Bayesian-network learning problem (1)

Learning one or more DAG models that fit a set of observed data \mathbf{D} well according to some scoring criterion $S(\mathcal{G}, \mathbf{D})$

Hypothesis \mathcal{G}^h for \mathcal{G} The observed data is a set of iid samples from a distribution that contains exactly the independence constraints implied by \mathcal{G}

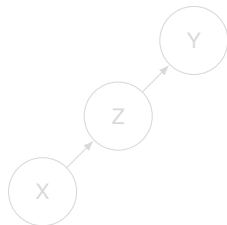
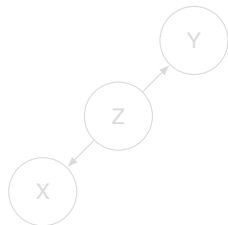


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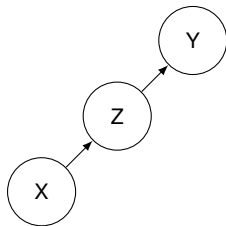
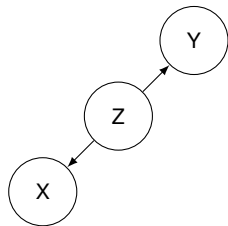


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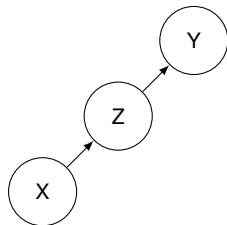
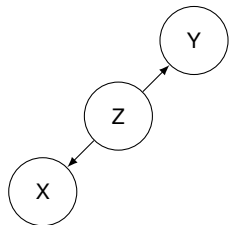


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*Perfect map We say that \mathcal{G} is a *perfect map* of P if every independence constraint in P is implied by \mathcal{G} and every independence implied by \mathcal{G} holds in P . In this case, P is *DAG-perfect**

Assumption Each record in \mathbf{D} is an iid sample from a DAG-perfect probability distribution

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Bayesian scoring criterion

Bayesian scoring criterion: $S_B(\mathcal{G}, \mathbf{D}) = \log P(\mathcal{G}^h) + \log P(\mathbf{D} | \mathcal{G}^h)$

- ▶ $P(\mathcal{G}^h)$: prior probability of \mathcal{G}^h
- ▶ $P(\mathbf{D} | \mathcal{G}^h)$: marginal likelihood obtained by integrating over the unknown parameters the likelihood function (Eq. 1) applied to each record in \mathbf{D} (illustration on board)

Bayesian information criterion (BIC - Schwarz, 1978) Under some assumptions:

$$S_B(\mathcal{G}, \mathbf{D}) = \underbrace{\log P(\mathbf{D} | \hat{\theta}, \mathcal{G}^h)}_{BIC} - \frac{d}{2} \log m + O(1)$$

$\hat{\theta}$: maximum-likelihood values of θ ; d : number of free parameters; m : number of records in \mathbf{D} ; $O(1)$: constant

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Decomposability, local consistency

Decomposability A scoring $S(\mathcal{G}, \mathbf{D})$ is decomposable if
$$S(\mathcal{G}, \mathbf{D}) = \sum_{i=1}^n s(X_i, \mathbf{Pa}_i^{\mathcal{G}})$$

Is the Bayesian scoring criterion decomposable?

Local consistency Let \mathbf{D} be m iid samples from distribution P , \mathcal{G} be any DAG and \mathcal{G}' the DAG obtained from \mathcal{G} by adding the edge $X_i \rightarrow X_j$. A scoring $S(\mathcal{G}, \mathbf{D})$ is *locally consistent* if the following properties hold:

1. If $X_j \not\perp\!\!\!\perp_P X_i \mid \mathbf{Pa}_j^{\mathcal{G}}$, then $S(\mathcal{G}', \mathbf{D}) > S(\mathcal{G}, \mathbf{D})$
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During the construction of graph inferred from data:

- ▶ Bayesian scoring criterion favours addition of edges that eliminate independence constraints not contained in the generative distribution
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The GES algorithm

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Markov equivalence

Theorem (Markov equivalence) Two DAGs are equivalent *iff* they have the same skeleton and the same v-structures

- ▶ Markov equivalence defines an equivalence relation (reflexive, symmetric, transitive)
- ▶ Equivalence class of \mathcal{G} : $\mathcal{E}(\mathcal{G})$

Covered edges An $X \rightarrow Y$ is covered in \mathcal{G} if $\mathbf{Pa}(Y) = \mathbf{Pa}(X) \cup X$

Lemma (Chickering, 1995) Let \mathcal{G} be a DAG and let \mathcal{G}' the result of reversing the edge $X \rightarrow Y$ in \mathcal{G} . \mathcal{G} and \mathcal{G}' are equivalent *iff* $X \rightarrow Y$ is covered in \mathcal{G}

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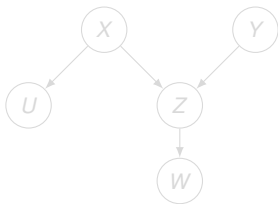
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Markov equivalence (cont'd)

CPDAG: completed PDAG; PDAG: partially DAG

CPDAG of an equivalence class The CPDAG of an equivalence class consists of a directed edge for every *compelled* edge, and an undirected edge for every *reversible* edge (compelled: exists in all graphs of the equivalence class; reversible: not compelled)

What's the CPDAG of the equivalence class of the following DAG?



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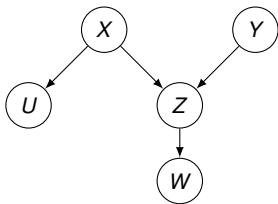


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Fundamental result and neighbour classes

Remark If \mathcal{G} and \mathcal{H} are in the same equivalence class, then $\mathcal{G}^h = \mathcal{H}^h$ and $S_B(\mathcal{G}, \mathbf{D}) = S_B(\mathcal{H}, \mathbf{D}) := S_B(\mathcal{E}(\mathcal{G}), \mathbf{D})$

Proposition Let \mathcal{E}^* denote the equivalence class that is a perfect map of distribution P , and let m be the number of records in \mathbf{D} . Then in the limit of large m , $S_B(\mathcal{E}^*, \mathbf{D}) > S_B(\mathcal{E}, \mathbf{D})$ for $\mathcal{E} \neq \mathcal{E}^*$

Neighbour classes $\mathcal{E}' \in \mathcal{E}^+(\mathcal{E})$ iff one can transform any DAG \mathcal{G} in \mathcal{E} to any DAG \mathcal{G}' in \mathcal{E}' through a sequence of covered edge reversals followed by a single edge addition followed by a sequence of covered edge reversals (same definition for $\mathcal{E}^-(\mathcal{E})$ with a single edge deletion)

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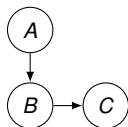
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Example

What are the equivalence class $\mathcal{E} = \mathcal{E}(\mathcal{G})$, $\mathcal{E}^+(\mathcal{E})$ and $\mathcal{E}^-(\mathcal{E})$ of the following graph \mathcal{G} ?



GES: greedy equivalence search

GES algorithm

1. Initialisation: set \mathcal{E} to the equivalence class corresponding to the DAG with no edge
2. Repeatedly replace \mathcal{E} with the member of $\mathcal{E}^+(\mathcal{E})$ that has the highest score, until no such replacement increases the score
3. Repeatedly replace \mathcal{E} with the member of $\mathcal{E}^-(\mathcal{E})$ that has the highest score, until no such replacement increases the score
4. Output the current class \mathcal{E}

Consistency of GES Let \mathcal{E} denote the equivalence class that results from GES, let P denote the DAG-perfect distribution associated with \mathbf{D} , and let m denote the number of records in \mathbf{D} . Then in the limit of large m , \mathcal{E} is a perfect map of P

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Remarks

1. Well-founded algorithm with consistency proof first established by Meek (Meek, 1997) based on a conjecture proven by Chickering (Chickering, 2002)
2. Main disadvantage: computational complexity
 - ▶ Learning optimal structure with Bayesian scoring criterion is NP-hard (Chickering, 1996)
 - ▶ Fast implementations exist when the underlying graph is sparse (Chickering, 2020)
3. Another (faster) approach exists based on the EM (expectation-maximisation) algorithm called MS-EM for *model selection EM* described in (Friedman, 1997)
4. Several other extensions for different data types, *e.g.* for time series (Assaad *et al.*, 2022)

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Learning models from data: a Bayesian approach

Granger causality

Causality according to Granger

Granger causality A time series X^p Granger-causes X^q if past values of X^p provide unique, statistically significant information about future values of X^q

Standard pairwise version Under the assumption of stationary linear systems, one considers the following autoregression models:

$$X_t^q = a_{q,0} + \sum_{i=1}^{\tau} a_{q,i} X_{t-i}^q + \zeta_t^q \quad (\text{Mres})$$

$$X_t^q = a_{q,0} + \sum_{i=1}^{\tau} a_{q,i} X_{t-i}^q + \sum_{i=1}^{\tau} a_{p,i} X_{t-i}^p + \zeta_t^q \quad (\text{Mfull})$$

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If the full model is significantly more accurate than the restricted model, one concludes that X^p Granger-causes X^q

Remarks

- ▶ Statistical test such as the F -test can be used to determine whether the full model is significantly better than the restricted one (null hypothesis: X^p does not Granger-cause X^q)
- ▶ Optimal lag τ estimated using an information criterion, as AIC (Akaike information criterion) or BIC

Causality according to Granger (cont'd)

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$$X_t^q = a_{q,0} + \sum_{i=1}^{\tau} a_{q,i} X_{t-i}^q + \sum_{i=1}^{\tau} a_{p,i} X_{t-i}^p + \zeta_t^q \quad (\text{Mfull})$$

If the full model is significantly more accurate than the restricted model, one concludes that X^p Granger-causes X^q

Remarks

- ▶ Statistical test such as the F -test can be used to determine whether the full model is significantly better than the restricted one (null hypothesis: X^p does not Granger-cause X^q)
- ▶ Optimal lag τ estimated using an information criterion, as AIC (Akaike information criterion) or BIC

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Associated algorithm

input X a d -dimensional time series, $\tau_{\max} \in \mathbb{N}$ optimal lag
initialisation Form an empty graph \mathcal{G} with d nodes V
Standardize data and check if it is covariance stationary
for $X^q \in V$ **do**
 Fit Mres and compute its residuals
 for $X^p \in V \setminus \{X^q\}$ **do**
 Fit Mfull and compute its residuals
 Compare Mres and Mfull
 if null hypothesis rejected **then** add $X^p \rightarrow X^q$ to \mathcal{G}
return \mathcal{G}

Multivariate extension

$$\mathcal{X}_t^q = \mathbf{a}_{q,0} + \sum_{\substack{r=1 \\ r \neq p}}^d \sum_{i=1}^{\tau} \mathbf{a}_{r,i} \mathcal{X}_{t-i}^p + \zeta_t^q \quad (\text{mvMres})$$

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Remarks

- ▶ Yields better results than previous version
- ▶ Computationally costly so that people mostly rely on pairwise version

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Other extensions

Several other extensions have been proposed (Assaad *et al.*, 2022), including

- ▶ Dealing with non-stationary processes (Luo *et al.*, 2015)
- ▶ Using deep learning to learn complex, non linear relations (Nauta *et al.*, 2019)

Granger causality is not causality: no explicit way to distinguish causal relations from spurious correlations

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Conclusion

We have reviewed the major methods for **causal discovery**

- ▶ Constraint-based methods
- ▶ Noise-based methods
- ▶ Score-based methods
- ▶ Granger causality

Other methods exist but are less used: logic-based approaches, topology-based approaches (not true causality), difference-based approaches (Assaad *et al.*, 2022)

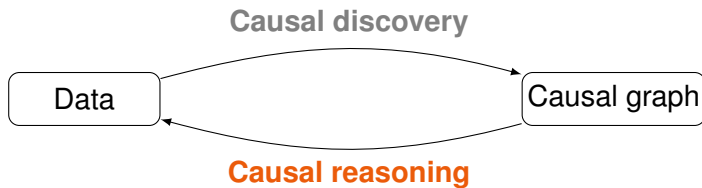
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Next courses



References (1)

1. *Estimating the dimension of a model*, G. E. Schwarz, 1978
2. *A transformational characterization of Bayesian network structures*, D. M. Chickering, 1995
3. *Learning Bayesian networks is NP-complete*, D. M. Chickering, 1996
4. *Learning belief networks in the presence of missing values and hidden variables*, N. Friedman, 1997
5. *Optimal structure identification with greedy search*, D. M. Chickering, 2002
6. *Discovering causal structures from time series data via enhanced Granger causality*, L. Luo, W. Liu, I. Koprinska, F. Chen, 2015

References (2)

7. *Causal discovery with attention-based convolutional neural networks*, M. Nauta, D. Bucur, C. Seifert, 2019
8. *Statistically Efficient Greedy Equivalence Search*, D. M. Chickering, 2020
9. *Survey and evaluation of causal discovery methods for time series*, K. C. Assaad, E. Devijver, E. Gaussier, 2022