Let's talk about Thurstone & Co.: An information-theoretical model

for comparative judgments, and its statistical translation

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Abstract

This study revisits Thurstone's law of comparative judgments (CJ) by addressing two key limitations

in traditional approaches. Firstly, it addresses the overreliance on the assumptions of Thurstone's

Case V in the statistical analysis of CJ data. Secondly, it addresses the apparent disconnect

between CJ's approach to trait measurement and hypothesis testing. We put forward a systematic

approach based on causal analysis and Bayesian statistical methods, which results in a model that

facilitates a more comprehensive understanding of the factors influencing CJ experiments while

offering a robust statistical translation. The new model accommodates unequal dispersions and

correlations between stimuli, enhancing the reliability and validity of CJ's trait estimation, thereby

ensuring the accurate measurement and interpretation of comparative data. The paper highlights

the relevance of this updated framework for modern empirical research, particularly in education

and social sciences. This contribution advances current research methodologies, providing a robust

foundation for future applications in diverse fields.

Keywords: causal inference, directed acyclic graphs, structural causal models, bayesian statistical

methods, thurstonian model, comparative judgement, probability, statistical modeling

1. Introduction

In comparative judgment (CJ) studies, judges assess a specific trait or attribute across different

stimuli by performing pairwise comparisons (Thurstone, 1927b,a). Each comparison produces a

dichotomous outcome, indicating which stimulus is perceived to have a higher trait level. For

example, when assessing writing quality, judges compare pairs of written texts (the stimuli) to

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determine the relative writing quality each text exhibit (the trait) (Laming, 2004; Pollitt, 2012b; Whitehouse, 2012; van Daal et al., 2016; Lesterhuis, 2018a; Coertjens et al., 2017; Goossens and De Maeyer, 2018; Bouwer et al., 2023).

Numerous studies have documented the effectiveness of CJ in assessing traits and competencies over the past decade. These studies have highlighted three aspects of the method's effectiveness: its reliability, validity, and practical applicability. Research on reliability suggests that CJ requires a relatively modest number of pairwise comparisons (Verhavert et al., 2019; Crompvoets et al., 2022) to generate trait scores that are as precise and consistent as those generated by other assessment methods (Coertjens et al., 2017; Goossens and De Maeyer, 2018; Bouwer et al., 2023). In addition, the evidence suggests that the reliability and time efficiency of CJ are comparable, if not superior, to those of other assessment methods when employing adaptive comparison algorithms (Pollitt, 2012b; Verhavert et al., 2022; Mikhailiuk et al., 2021). Meanwhile, research on the validity of CJ scores indicates their capacity to represent accurately the traits under measurement (Whitehouse, 2012; van Daal et al., 2016; Lesterhuis, 2018a; Bartholomew et al., 2018; Bouwer et al., 2023). Moreover, research on its practical applicability highlights CJ's versatility across both educational and non-educational contexts (Kimbell, 2012; Jones and Inglis, 2015; Bartholomew et al., 2018; Jones et al., 2019; Marshall et al., 2020; Bartholomew and Williams, 2020; Boonen et al., 2020).

Nevertheless, despite the increasing number of CJ studies, research in this domain remains unsystematic and fragmented, leaving several critical issues unresolved. This study identifies and discusses two prominent issues that can undermine the reliability and validity of CJ's trait estimates. The first issue arises from the excessive reliance on Thurstone's Case V assumptions in the statistical analysis of CJ data. The second stems from the apparent disconnect between CJ's approach to trait measurement and hypothesis testing. This study addresses these issues by extending Thurstone's general form through a systematic and integrated approach based on Causal and Bayesian inference methods.

As a result, the study divides its content into six main sections. Section 2 provides an overview of Thurstone's theory. Section 3 discusses the identified issues in detail. Section 4 extends Thurstone's general form to address these challenges. The extension integrates core theoretical principles alongside key assessment design features relevant to CJ experiments, such as the selection of judges, stimuli, and comparisons. Section 5 translates these theoretical and practical elements into a probabilistic statistical model for the analysis of dichotomous pairwise comparison data. Section 6

discusses the findings, limitations, challenges and explores avenues for future research. Finally, Section 7 summarizes the study's key insights.

2. Thurstone's theory

In its most general form, Thurstone's theory addresses pairwise comparisons wherein a single judge evaluates multiple stimuli (Thurstone, 1927a, pp. 267). The theory posits that two key factors determine the dichotomous outcome of these comparisons: the discriminal process of each stimulus and their discriminal difference. The discriminal process captures the psychological impact each stimulus exerts on the judge or, more simply, his perception of the stimulus trait. The theory assumes that the discriminal process for any given stimulus forms a Normal distribution along the trait continuum (Thurstone, 1927a, pp. 266). The mode (mean) of this distribution, known as the modal discriminal process, indicates the stimulus position on this continuum, while its dispersion, referred to as the discriminal dispersion, reflects variability in the perceived trait of the stimulus.

Figure 1a illustrates the hypothetical discriminal processes along a quality trait continuum for two written texts. The figure indicates that the modal discriminal process for Text B is positioned further along the continuum than that of Text A $(T_B > T_A)$, suggesting that Text B exhibits higher quality. Additionally, the figure highlights that Text B has a broader distribution compared to Text A, which arises from its larger discriminal dispersion $(\sigma_B > \sigma_A)$.

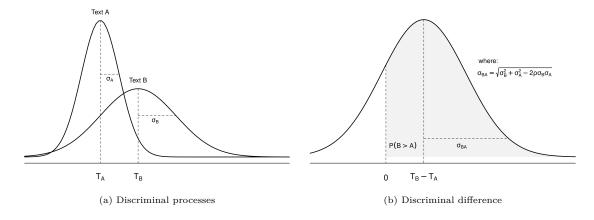


Figure 1: Hypothetical discriminal processes and discriminant difference along a quality trait continuum for two written texts.

However, since the individual discriminal processes of the stimuli are not directly observable, the theory introduces the *law of comparative judgment*. This law posits that in pairwise comparisons, a judge perceives the stimulus with a discriminal process positioned further along the trait continuum

as possessing more of the trait (Bramley, 2008, pp. 251). This suggests that the relative distance between stimuli, rather than their absolute positions on the continuum, likely defines the outcome of pairwise comparisons. Indeed, the theory assumes that the difference between the underlying discriminal processes of the stimuli, referred to as the discriminal difference, determines the observed dichotomous outcome. Furthermore, the theory assumes that because the individual discriminal processes form a Normal distribution on the continuum, the discriminal difference will also conform to a Normal distribution (Andrich, 1978, pp. 452). In this distribution, the mode (mean) represents the relative separation between the stimuli, and its dispersion indicates the variability of that separation.

Figure 1b illustrates the distribution of the discriminal difference for the hypothetical texts depicted in Figure 1a. The figure indicates that the judge perceives Text B as having significantly higher quality than Text A. This conclusion is supported by two key observations: the positive difference between their modal discriminal processes $(T_B - T_A > 0)$ and the probability area where the discriminal difference distinctly favors Text B over Text A, represented by the shaded gray area denoted as P(B > A). As a result, the dichotomous outcome of this comparison is more likely to favor Text B over Text A.

3. Two prominent issues in CJ practice

Thurstone noted from the outset that his general form, described in Section 2, led to a trait scaling problem. Specifically, the model required estimating more "unknown" parameters than the number of available pairwise comparisons (Thurstone, 1927a, pp. 267). For instance, in a CJ experiment with five texts, the general form would require estimating 20 parameters: five modal discriminal processes, five discriminal dispersions, and 10 correlations—one per comparison (see Table 1). However, a single judge could provide only $\binom{5}{2} = 10$ unique comparisons, an insufficient data set to estimate the required parameters.

To address this issue and facilitate the practical implementation of the theory, Thurstone developed five cases derived from this general form, each progressively incorporating additional simplifying assumptions into the model. In Case I, Thurstone postulated that pairs of stimuli would maintain a constant correlation across all comparisons. In Case II, he allowed multiple judges to undertake comparisons instead of confining evaluations to a single judge. In Case III, he posited that there was no correlation between stimuli. In Case IV, he assumed that the stimuli exhibited similar dispersions. Finally, in Case V, he replaced this assumption with the condition that stimuli had

Table 1: Thurstones cases and their asumptions

	General	Thurstone's					\mathbf{BTL}
Assumption	\mathbf{form}	Case I	Case II	Case III	Case IV	Case V	model
Discriminal process (distribution)	Normal	Normal	Normal	Normal	Normal	Normal	Logistic
Discriminal dispersion (between stimuli)	Different	Different	Different	Different	Similar	Equal	Equal
Correlation (between stimuli)	One per pair	Constant	Constant	Zero	Zero	Zero	Zero
How many judges compare?	Single	Single	Multiple	Multiple	Multiple	Multiple	Multiple

equal discriminal dispersions. Table 1 summarizes the assumptions of the general form and the five cases. For a detailed discussion of these cases and their progression, refer to Thurstone (1927a) and Bramley (2008, pp. 248–253).

It is evident that although Thurstone developed Case V to provide a "rather coarse scaling" of traits (Thurstone, 1927a, pp. 269) and cautioned that its use "should not be made without experimental test" (Thurstone, 1927a, pp. 270), the CJ literature has nevertheless relied the most on it in practice. This preference is particularly notable given that Case V incorporates the most extensive set of simplifying assumptions among all the cases (Bramley, 2008, pp. 253; Kelly et al., 2022, pp. 677) (see Table 1).

Furthermore, since Thurstone's primary goal was to define a psychological scale and "allocate the compared stimuli on this continuum" (Thurstone, 1927a, pp. 269), he did not provide guidance on how to use these measurements when, for instance, researchers aim to test hypotheses related to the measured traits. Although the CJ practice has circumvented this issue by using scores' point estimates or their transformations for such analyses, a fundamental question remains: Is this method appropriate for addressing these inquiries?

Thus, this section discusses these two prominent issues, as they could undermine the reliability and validity of CJ's trait estimates (Perron and Gillespie, 2015, pp. 2). Specifically, Section 3.1 examines the the excessive reliance on Thurstone's Case V assumptions in the statistical analysis of CJ data. Moreover, Section 3.2 focuses on the apparent disconnect between the approaches to trait measurement and hypothesis testing in CJ.

3.1. The overreliance on Case V in the statistical analysis of CJ data

As previously discussed, despite relying on the most extensive set of simplifying assumptions (Bramley, 2008, pp. 253; Kelly et al., 2022, pp. 677), Case V remains the most widely used case

in the CJ literature. This popularity stems mainly from its simplified statistical representation in the Bradley-Terry-Luce (BTL) model (Bradley and Terry, 1952; Luce, 1959). The BTL model mirrors the assumptions of Case V, with one notable distinction: whereas Case V assumes a Normal distribution for the stimuli discriminal processes, the BTL model uses the more mathematically tractable Logistic distribution (Andrich, 1978; Bramley, 2008, pp. 254) (see Table 1). As noted by van der Linden (2017a, pp. 16), this substitution has little impact on the model's estimation or interpretation because the discriminal process scale is arbitrary up to a non-monotonic transformation. Furthermore, this minimal impact is reinforced by the fact that the Normal and Logistic distributions exhibit analogous statistical properties, differing only by a scaling factor of approximately 1.7.

However, as mentioned earlier, Thurstone originally developed Case V to provide a "rather coarse scaling" of traits (Thurstone, 1927a, pp. 269), prioritizing statistical simplicity over precision in trait measurement (Kelly et al., 2022, pp. 677). He explicitly warned against its untested application, stating that its use "should not be made without (an) experimental test" (Thurstone, 1927a, pp. 270). Furthermore, he acknowledged that some assumptions could prove problematic when researchers assess complex traits or heterogeneous stimuli (Thurstone, 1927b, pp. 376). Consequently, given that modern CJ applications frequently involve such traits and stimuli, two main assumptions of Case V and, by extension, of the BTL model may not consistently hold in theory or practice, namely the assumption of equal dispersion and zero correlation between stimuli.

3.1.1. The assumption of equal dispersions between stimuli

Thurstone argued that the assumption of equal dispersions may not apply when researchers assess complex traits or heterogeneous stimuli (Thurstone, 1927b, pp. 376), as these traits and stimuli can introduce judgment discrepancies due to their unique characteristics (van Daal et al., 2016; Lesterhuis, 2018b; Chambers and Cunningham, 2022). Indeed, the CJ literature may already provide evidence of this violation through "misfit" statistics. For example, labeling texts as misfits indicates that comparisons involving these texts result in more judgment discrepancies than those involving other texts (Pollitt, 2004, pp. 11; Pollitt, 2012b, pp. 290). These discrepancies imply that the discriminal differences for misfit texts have broader distributions, further suggesting that their discriminal processes exhibit more dispersion than those of other texts. This idea can be illustrated using Figure 1a, where text B shows a wider discriminal dispersion than text A. As a result, the texts' discriminal processes overlap more, making it difficult for a judge to determine which text possesses more of the trait.

A similar line of reasoning applies to the concept of "misfit" judges, whose evaluations deviate substantially from the shared consensus due to the unique characteristics of the stimuli or the judges themselves (Pollitt, 2012a, pp. 164-165; Pollitt, 2012b, pp. 289-290; van Daal et al., 2016, pp. 4; Goossens and De Maeyer, 2018, pp. 20). These "misfit" judges and their associated deviations can give rise to additional statistical and measurement issues, which we discuss in more detail in Section 3.1.2.

According to the theory, discrepancies in the discriminal dispersions of stimuli shape the distribution of the discriminal difference, exerting a direct influence on the outcome of pairwise comparisons. A thought experiment can illustrate this idea. In this experiment, researchers can observe the discriminal processes for the texts depicted in Figure 1a. Furthermore, the experiment assumes that the discriminal dispersion for Text A remains constant and that the texts are uncorrelated ($\rho = 0$). Figure 2a reveals that an increase in the uncertainty associated with the perception of Text B compared to Text A ($\sigma_B - \sigma_A$), broadens the distribution of their discriminal difference. This broadening affects the probability area where the discriminal difference distinctly favors Text B over Text A, expressed as P(B > A), ultimately influencing the comparison outcome. Additionally, the figure reveals that when the discriminal dispersions of the texts are equal ($\sigma_B - \sigma_A = 0$), the discriminal difference is more narrow compared to situations where their dispersions differ. Consequently, the discriminal difference is more likely to favor Text B over Text A (shaded gray area).

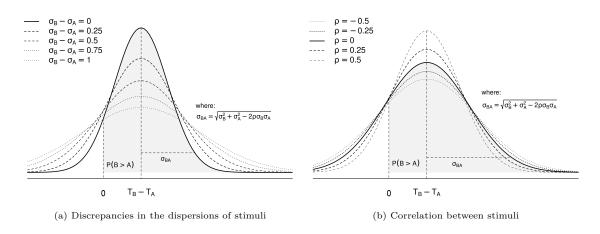


Figure 2: The effect of dispersion discrepancies and stimuli correlation on the distribution of the discriminal difference.

In experimental practice, however, the thought experiment occurs in reverse. Researchers first observe the comparison outcome and then use the BTL model to infer the discriminal difference between stimuli and their respective discriminal processes (Thurstone, 1927b, pp. 373). Consequently, the ability of the outcome to reflect *true* differences between stimuli largely depends on the validity

of the model's assumptions (Kohler et al., 2019, pp. 150), particularly the assumption of equal dispersions. For example, when researchers correctly assume equal dispersions between texts, the BTL model estimates a discriminal difference distribution that accurately represents the *true* discriminal difference between them. This scenario is illustrated in Figure 2a, when the model's discriminal difference distribution aligns with the *true* discriminal difference distribution, represented by the thick continuous line corresponding to $\sigma_B - \sigma_A = 0$. The accuracy of this discriminal difference then ensures reliable estimates for the texts' discriminal processes.

However, erroneously assuming equal dispersions between stimuli can give rise to significant statistical and measurement issues. Specifically, the model may overestimate the trait's reliability, that is, the degree to which the outcome accurately reflects the *true* discriminal differences between stimuli. This overestimation, in turn, results in spurious conclusions about these differences (McElreath, 2020, pp. 370) and, by extension, about the underlying discriminal processes of stimuli. Figure 2a also illustrates this scenario when the model's discriminal difference distribution aligns with the thick continuous line for $\sigma_B - \sigma_A = 0$, while the *true* discriminal difference follows any discontinuous line where $\sigma_B - \sigma_A \neq 0$. Additionally, if researchers recognize that misfit statistics highlight these critical differences in dispersions, the conventional CJ practice of excluding stimuli based on these statistics (Pollitt, 2012a,b; van Daal et al., 2016; Goossens and De Maeyer, 2018) can unintentionally discard valuable information, introducing bias into trait estimates (Zimmerman, 1994; McElreath, 2020, chap. 12). The direction and magnitude of these biases are often unpredictable, as they depend on which stimuli are excluded from the analysis.

3.1.2. The assumption of zero correlation between stimuli

In contrast, Thurstone assumed that stimuli were uncorrelated because judges' biases, arising from two opposing and equally weighted effects occurring during the pairwise comparisons, canceled each other out (Thurstone, 1927a, pp. 268). Andrich (1978) provided a mathematical demonstration of this cancellation using the BTL model under the assumption of discriminal processes with additive biases. However, the assumption of zero correlation is not met in practice in at least two scenarios. First, when intricate aspects of multidimensional, complex traits or heterogeneous stimuli influence judges' perceptions, generating biases that resist cancellation. Second, when an additional hierarchical structure is relevant to the stimuli.

In the first scenario, research on text quality suggests that when judges evaluate multidimensional, complex traits or heterogeneous stimuli, they often rely on various intricate characteristics of the stimuli to form their judgments (van Daal et al., 2016; Lesterhuis, 2018b; Chambers and Cunningham, 2022). These additional relevant characteristics are unlikely to be equally weighted or opposing. As a result, they may exert an uneven influence on judges' perceptions, creating biases in their judgments that resist cancellation. For example, this could occur when a judge assessing the argumentative quality of a text places more weight on its grammatical accuracy than other judges, thereby favoring texts with fewer errors but weaker arguments. Furthermore, since the discriminal difference of the stimuli becomes an observable outcome only through the judges' perceptions, these biases could ultimately introduce dependencies between the stimuli (van der Linden, 2017b, pp. 346). While direct evidence for this particular scenario is lacking, studies such as Pollitt and Elliott (2003) demonstrate the presence of such biases, supporting the notion that the factors influencing pairwise comparisons may not always cancel out.

In the second scenario, the shared context or inherent connections introduced by additional hierarchical structures may create dependencies between stimuli, a statistical phenomenon known as clustering (Everitt and Skrondal, 2010). Although the CJ literature acknowledges the existence of such hierarchical structures, the statistical approaches to account for this additional source of dependence have been insufficient. For instance, when CJ data incorporates multiple samples of stimuli from the same individuals, researchers frequently rely on (averaged) point estimates of the BTL scores to conduct subsequent analyses and tests at the individual hierarchical level (Bramley and Vitello, 2019; Boonen et al., 2020; Bouwer et al., 2023; van Daal et al., 2017; Jones et al., 2019; Gijsen et al., 2021). However, this approach can introduce additional statistical and measurement issues, which we discuss in greater detail in Section 3.2.

According to the theory, the correlation ρ measures how much the judges' perception of a specific trait in one stimulus depends on their perception of the same trait in another. As with the discriminal dispersions, this correlation shapes the distribution of the discriminal difference, directly impacting the outcomes of pairwise comparisons. A similar thought experiment, as the one depicted in Section 3.1.1, can illustrate this idea. The experiment now assumes that the discriminal dispersions for both texts remain constant. Figure 2b reveals that as the correlation between the texts increases, the distribution of their discriminal difference becomes narrower. This narrowing affects the area under the curve where the discriminal difference distinctly favors Text B over Text A, denoted as P(B > A), thus influencing the comparison outcome. Furthermore, the figure shows that when two texts are independent or uncorrelated ($\rho = 0$), their discriminal difference is less narrow compared to scenarios where the texts are highly correlated. Consequently, the discriminal

difference is less likely to favor Text B over Text A (shaded gray area).

Again, in experimental practice, researchers approach this process in reverse. They begin by observing the outcomes and use the BTL model to estimate the stimuli' discriminal difference and discriminal processes. Given that the BTL model assumes independent discriminal processes across comparisons, if this assumption holds, the model estimates a discriminal difference distribution that accurately reflects the *true* discriminal difference between the texts. This scenario is illustrated in Figure 2b when the discriminal difference distribution of the model aligns with the *true* distribution, represented by the thick continuous line corresponding to $\rho = 0$. Once more, the accuracy of the discriminal difference ensures reliable estimates for the discriminal processes of the texts.

However, erroneously assuming zero correlation between stimuli can lead to significant statistical and measurement issues. Specifically, neglecting judges' biases or hierarchical structures can create dimensional mismatches in the model, leading to the over- or underestimation of trait reliability (Hoyle, 2023, pp. 341, 482). In other words, the outcome may fail to accurately reflect the *true* discriminal differences between stimuli. These inaccuracies can result in spurious conclusions about these differences (McElreath, 2020, pp. 370) and, by extension, the underlying discriminal processes (Ackerman, 1989). This issue is illustrated in Figure 2b when the model discriminal difference distribution follows the thick continuous line ($\rho = 0$), while the *true* discriminal difference follows any discontinuous line where ($\rho \neq 0$). Conversely, excluding judges based on misfit statistics risks discarding valuable information, potentially biasing the trait estimates (Zimmerman, 1994; McElreath, 2020, chap. 12). Again, the direction and magnitude of these biases are often unpredictable, as they depend on which judges are excluded from the analysis.

3.2. The disconnect between trait measurement and hypothesis testing

Building on the previous section, it is clear that, despite its limitations, the BTL model commonly serves as the measurement model for CJ assessments. Specifically, the model provides a method for transforming manifest into latent variables (psychological scales) and allocating the compared stimuli on this continuum (Thurstone, 1927a, pp. 269; Everitt and Skrondal, 2010). For example, when evaluating writing quality, researchers use the BTL model to process the dichotomous outcomes resulting from the pairwise comparisons (the manifest variables) to estimate scores that reflect the underlying level of writing quality (the latent variable) (Laming, 2004; Pollitt, 2012b; Whitehouse, 2012; van Daal et al., 2016; Lesterhuis, 2018a; Coertjens et al., 2017; Goossens and De Maeyer, 2018; Bouwer et al., 2023).

Researchers then typically use the BTL scores' point estimates, or their transformations, to conduct additional analyses or hypothesis tests. For example, these scores have been used to identify 'misfit' judges and stimuli (Pollitt, 2012b; van Daal et al., 2016; Goossens and De Maeyer, 2018), detect biases in judges' ratings (Pollitt and Elliott, 2003; Pollitt, 2012b), calculate correlations with other assessment methods (Goossens and De Maeyer, 2018; Bouwer et al., 2023), or test hypotheses related to the underlying trait of interest (Casalicchio et al., 2015; Bramley and Vitello, 2019; Boonen et al., 2020; Bouwer et al., 2023; van Daal et al., 2017; Jones et al., 2019; Gijsen et al., 2021).

However, the statistical literature advises caution when using estimated scores for additional analyses and tests. A key consideration is that BTL scores are parameter estimates that inherently carry uncertainty. Ignoring this uncertainty can bias the analysis and reduce the precision of hypothesis tests. Notably, the direction and magnitude of such biases are often unpredictable. Results may be attenuated, exaggerated, or remain unaffected depending on the degree of uncertainty in the scores and the actual effects being tested (Kline, 2023, pp. 25; Hoyle, 2023, pp. 137). Finally, the reduced precision in hypothesis tests diminishes their statistical power, increasing the likelihood of committing type-I or type-II errors (McElreath, 2020).

In aggregate, excessive reliance on Thurstone's Case V assumptions in the statistical analysis of comparative data can compromise the reliability of trait estimates. This overreliance may also undermine their validity (Perron and Gillespie, 2015, pp. 2), particularly when coupled with the disconnect between CJ's approach to trait measurement and hypothesis testing. However, the structural approach to causal inference can address these issues by offering a systematic and integrated framework that strengthens measurement reliability and validity while enhancing the statistical accuracy of hypothesis tests.

4. Extending Thurstone's general form

The structural approach to causal inference provides a formal framework for identifying causes and estimating their effects using data. The approach uses structural causal models (SCMs) and directed acyclic graphs (DAGs) (Pearl, 2009; Pearl et al., 2016; Gross et al., 2018; Neal, 2020) to formally and graphically represent the assumed causal structure of a system, such as the one found in CJ experiments. Essentially, SCMs and DAGs function as conceptual models on which identification analysis rests (Schuessler and Selb, 2023, pp. 4). Identification analysis helps researchers to determine whether an estimator can accurately compute an estimand based solely on its (causal)

assumptions, regardless of random variability (Schuessler and Selb, 2023, pp. 4). Here, estimands represent the specific quantities researchers aim to determine (Everitt and Skrondal, 2010). Estimators denote the methods or functions that transform data into an estimate, while estimates are the numerical values approximating the estimand (Neal, 2020; Everitt and Skrondal, 2010).

A motivating example that will appear throughout the document clarifies these concepts. In this example, researchers aim to determine: "To what extent do different teaching methods influence students' ability to produce high-quality written texts?" To investigate this, researcher design a CJ experiment by randomly assigning students (individuals) to two groups, each receiving a different teaching method. Judges then compare pairs of students' written texts (stimuli), to produce a dichotomous outcome that reflects the relative quality of each text (trait). Based on this setup, researchers reformulate the research question as the estimand: "On average, is there a difference in the ability to produce high-quality written texts between the two groups of students?". Finally, following current CJ practices, researchers then rely on estimates from the BTL model, or its transformations, to approximate this estimand.

However, Section 3 presents compelling evidence that Thurstone's Case V, and by extension the BTL model, suffers from several statistical and measurement limitations. These limitations hinder the model's ability to identify various estimands relevant to CJ inquiries, including the one described in the example. Identification is crucial because it is a necessary condition for ensuring consistent estimators. *Consistency* refers to the property of an estimator whose estimates converge to the "true" value of the estimand as the data size approaches infinity (Everitt and Skrondal, 2010). Without identification, consistency cannot be achieved, even with "infinite" and error-free data. Thus, deriving meaningful insights from finite data becomes impossible (Schuessler and Selb, 2023, pp. 5).

Luckily, SCMs and DAGs support identification analysis through two key advantages. First, regardless of complexity, they can represent various causal structures using only five fundamental building blocks (Neal, 2020; McElreath, 2020). This feature allows researchers to decompose complex structures into manageable components, facilitating their analysis (McElreath, 2020). Second, they depict causal relationships in an interactive, non-parametric way. This flexibility enables feasible identification strategies without requiring specification of the types of variables, the functional forms relating them, or the parameters of those functional forms (Pearl et al., 2016, pp. 35).

Thus, this section addresses the issues identified in Section 3 by extending Thurstone's general

form using the structural approach to causal inference. Specifically, it formalizes the combination of the core theoretical principles outlined in Section 2 with key assessment design features relevant to CJ experiments, such as the selection of judges, stimuli, and comparisons. In addition to enhancing statistical accuracy and strengthening measurement reliability and validity, the approach offers two key advantages. First, it clarifies the interactions among all actors and processes involved in CJ experiments. Second, it shifts the current comparative data analysis paradigm from passively accepting the model assumptions to actively testing whether those assumptions fit the data under analysis.

Accordingly, Section 4.1 incorporates the theoretical principles into what we call the *conceptual-population model*. This model assumes that researchers have access to a *census* of comparative data, that is, data representing all repeated judgments made by every available judge for each pair of stimuli produced by each pair of individuals in the population, hence its name. Conversely, Section 4.2 integrates the assessment design features into what we call the *sample-comparison model*. This model assumes a more realistic scenario where researchers only have access to a sample of judges, individuals, stimuli, and comparisons from the conceptual population.

4.1. The conceptual-population model

4.1.1. Integrating initial theoretical principles

Before incorporating the theoretical principles into the conceptual-population model, it is essential to further define SCMs. SCMs are formal mathematical models characterized by a set of endogenous variables V, a set of exogenous variables E, and a set of functions F (Pearl, 2009; Cinelli et al., 2020). Endogenous variables are those whose causal mechanisms a researcher chooses to model (Neal, 2020). In contrast, exogenous variables represent errors or disturbances arising from omitted factors that the investigator chooses not to model explicitly (Pearl, 2009, pp. 27,68). Lastly, the functions, referred to as structural equations, express the endogenous variables as non-parametric functions of other endogenous and exogenous variables. These functions use the symbol ':=' to denote the asymmetrical causal dependence between variables and the symbol ' \bot ' to represent d-separation, a concept akin to (conditional) independence.

SCM 3a presents the initial integration of theoretical principles into the conceptual-population model for the example, applied to examine the effect of different teaching methods on students' writing ability. This SCM maps the relationship between the conceptual-population outcome (O^{cp}_{hiabjk}) and several related variables. The subscripts h and i label the students who authored the texts

(individuals), while the indices a and b represent the compared texts (stimuli). Moreover, the index j refers to the judge, and k specifies the judgment number, accounting for experimental conditions where judges perform the same judgment multiple times, i.e., a repeated measures design (Lawson, 2015, pp. 366-376). Notably, the indexing system allows comparisons between texts written by the same student $(h = i, a \neq b)$ and between texts written by distinct students $(h \neq i, \text{ where } a = b \text{ is possible})$. However, it excludes cases where judges compare a student's text to itself (h = i, a = b), as such comparison lacks practical meaning within the CJ framework.

In line with Thurstone's theory, SCM 3a depicts the texts' discriminal processes (T_{ha}, T_{ib}) and their discriminal difference (D_{hiabjk}) (see Section 2). Additionally, based on the arguments developed under the first scenario of Section 3.1.2, and the recommendations of Andrich (1978) and Wainer et al. (1978), the SCM incorporates the judges' biases (B_{kj}) . Together with the outcome, these variables form the preliminary set of endogenous variables, $V = \{O_{hiabjk}, D_{hiabjk}, T_{ha}, T_{ib}, B_{kj}\}$. Finally, the SCM shows the preliminary set of structural equations, $F = \{f_O, f_D\}$, which define the non-parametric dependencies among these variables.

$$O_{hiabjk}^{cp} := f_O(D_{hiabjk})$$

$$D_{hiabjk} := f_D(T_{ha}, T_{ib}, B_{jk})$$
 (a) SCM
$$T_{ha}$$

$$O_{hiabjk}^{cp}$$

$$O_{hiabjk}^{cp}$$

$$O_{hiabjk}^{cp}$$

$$O_{hiabjk}^{cp}$$
 (b) DAG

Figure 3: Conceptual-population model, scalar form.

Notably, every SCM has an associated DAG (Pearl et al., 2016; Cinelli et al., 2020). A DAG is a graph consisting of nodes connected by edges, where nodes represent random variables. The term directed indicates that edges or arrows extend from one node to another, indicating the direction of causal influence, while the absence of an edge implies no relationship between the nodes. The term acyclic means that the causal influences do not form loops, ensuring the influences do not cycle back on themselves (McElreath, 2020). DAGs conventionally represent observed variables with solid

black circles and unobserved (latent) variables with open circles (Morgan and Winship, 2014). They also display conditioned variables by surrounding them with a square. Here, conditioning refers to restricting the statistical analysis to a subset of the population defined by the values of the given variable (Neal, 2020, pp. 32). While DAGs often omit exogenous variables for simplicity, following a standard representation, this section adopts the magnified representation of DAGs, as including exogenous variables can improve clarity and help reveal potential issues related to conditioning and confounding (Cinelli et al., 2020).

Figure 3b displays the DAG corresponding to SCM 3a, illustrating the causal relationships among variables based on Thurstone's theory. Particularly, the graph shows how the texts' discriminal processes (T_{ha}, T_{ib}) influence their discriminal difference (D_{hiabjk}) , which in turn determines the outcome (O_{hiabjk}^{cp}) . It also depicts the judges' biases (B_{kj}) as an additional factor influencing the texts' discriminal difference. Furthermore, the DAG distinguish between observed endogenous variables, such as the outcome, and latent endogenous variables, which include the texts' discriminal processes, their discriminal difference, and the judges' biases.

4.1.2. The comparative census data structure

Although specifying a system's data structure is not mandatory when using SCMs and DAGs, defining one can improve clarity and facilitate the description of the system. Thus, to re-express the scalar form of the CJ system, outlined in the previous section, into an simpler equivalent vectorized form, we first define the vectors I and J, along with the matrices IA and JK, as follows:

where an element of I corresponds to a unique individual h or i, while each row of IA corresponds to a unique pairing of individuals h, i with stimuli a, b. Here, n_I represents the number of individuals, while n_A denotes the number of stimuli available per individual. Thus, vector I has a length of n_I , and matrix IA has a dimension of $(n_I \cdot n_A)$ rows and 2 columns. The vector and the matrix consist of scalars from the set of natural numbers \mathbb{N} .

Similarly, an element of the vector J corresponds to a unique judge j. Moreover, each row of JK specifies a unique pairing of a judge j with a judgment number k. In this case, n_J indicates the total number of judges, and n_K represents the number of judgments each judge performs. As a result, vector J has a length of n_J , and matrix JK has a dimension of $(n_J \cdot n_K)$ rows and 2 columns. Like the previous structures, the vector and the matrix consist of scalars from the set of natural numbers \mathbb{N} .

Additionally, we define the matrix RPUV, where each row uniquely pairs the corresponding rows from the matrices IA and JK. This matrix reflects the fundamental idea that the discriminal difference of the stimuli becomes an observable outcome only through the judges' perceptions. Thus, the matrix is defined as follows:

$$RPUV_{n\times 6} = \begin{bmatrix} 1 & 1 & 1 & 2 & 1 & 1 \\ 1 & 1 & 1 & 2 & 1 & 2 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ 1 & 1 & 1 & 2 & 1 & n_K - 1 \\ 1 & 1 & 1 & 2 & 1 & n_K \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ h & i & a & b & j & k \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ n_I & n_I & n_A - 1 & n_A & n_J & 1 \\ n_I & n_I & n_A - 1 & n_A & n_J & 2 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ n_I & n_I & n_A - 1 & n_A & n_J & n_K - 1 \\ n_I & n_I & n_A - 1 & n_A & n_J & n_K \end{bmatrix}$$

$$(2)$$

This matrix also consists of scalars from the set of natural numbers \mathbb{N} and has a dimension of $n=\binom{n_I\cdot n_A}{2}\cdot n_J\cdot n_K$ rows and 6 columns. The matrix's dimension results from the conceptual-population model's assumption that researchers have access to a *census* of comparative data, capturing all repeated judgments made by every available judge for each pair of stimuli produced by each pair of individuals in the population. Here, the term $\binom{n_I\cdot n_A}{2}$ represents the Binomial coefficient, which quantifies the total number of individual-stimuli comparison pairs within the conceptual population. The coefficient is defined as follows:

$${n_I \cdot n_A \choose 2} = \frac{(n_I \cdot n_A)!}{2!(n_I \cdot n_A - 2)!}$$

We visualize the structure of these vectors and matrices by considering small values for the parameters corresponding to number of individuals, the number of available stimuli per individual, the total number of judges, and the number of judgments each judge performs. For instance, assuming $n_I = 5$, $n_A = 2$, $n_J = 3$, and $n_K = 3$, the matrices described in equations (1) and (2) take the following form:

$$I_{5} = \begin{bmatrix} 1 \\ 2 \\ 3 \\ 4 \\ 5 \end{bmatrix}; IA_{10 \times 2} = \begin{bmatrix} 1 & 1 \\ 1 & 2 \\ 2 & 1 \\ 3 & 2 \\ 4 & 1 \\ 4 & 2 \\ 5 & 1 \\ 5 & 2 \end{bmatrix}; J_{3} = \begin{bmatrix} 1 \\ 2 \\ 2 \\ 3 \\ 3 \end{bmatrix}; JK_{9 \times 2} = \begin{bmatrix} 1 & 1 \\ 1 & 2 \\ 1 & 3 \\ 2 & 1 \\ 2 & 2 \\ 3 & 3 \end{bmatrix}; RPUV_{405 \times 6} = \begin{bmatrix} 1 & 1 & 1 & 2 & 1 & 1 \\ 1 & 1 & 2 & 1 & 3 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ 1 & 1 & 5 & 1 & 1 & 1 \\ 1 & 1 & 5 & 1 & 1 & 2 \\ 1 & 1 & 5 & 1 & 1 & 3 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ 1 & 1 & 5 & 1 & 3 & 1 \\ 1 & 1 & 5 & 1 & 3 & 1 \\ 1 & 1 & 5 & 1 & 3 & 2 \\ 1 & 1 & 5 & 1 & 3 & 3 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ 5 & 1 & 5 & 2 & 3 & 1 \\ 5 & 1 & 5 & 2 & 3 & 2 \\ 5 & 1 & 5 & 2 & 3 & 3 \end{bmatrix}$$

Thus, using the data structure defined by the matrices in equations (1) and (2), we can reexpress SCM 3a and DAG 3b in a simpler equivalent vectorized form, as shown in Figure 4. In this depiction, the outcome O_{RPUV}^{ep} , the texts' discriminal difference D_{RPUV} , their discriminal processes T_{IA} , and the judges' biases B_{JK} are represented as vectors rather than scalar values. These vectors encompass all values from the conceptual population. Specifically, O_{RPUV}^{ep} is an observed vector of length $\binom{nI \cdot n_A}{2} \cdot n_J \cdot n_K$, while D_{RPUV} is a latent vector of the same dimension. Additionally, T_{IA} and B_{JK} are latent vectors of lengths $(n_I \cdot n_A)$ and $(n_J \cdot n_K)$, respectively.

$$O_{RPUV}^{cp} := f_O(D_{RPUV})$$

$$D_{RPUV} := f_D(T_{IA}, B_{JK})$$
 (a) SCM
$$T_{IA} \bigcirc O_{RPUV}^{cp} \bigcirc O_{RPUV}^{cp}$$

$$O_{RPUV}^{cp} \bigcirc O_{RPUV}^{cp}$$
 (b) DAG

Figure 4: Conceptual-population model, vectorized form.

4.1.3. Integrating hierarchical structural components

Building on the principles of Structural Equation Modeling (SEM) (Hoyle, 2023, pp. 138) and Item Response Theory (IRT) (Fox, 2010, chap. 6; van der Linden, 2017a, chap. 24), the conceptual-population model incorporates two hierarchical structural components. Each component describes how different observed or latent variables influence the primary latent variable of interest (Everitt and Skrondal, 2010). This addition addresses the problems arising when an additional hierarchical structure is relevant to the stimuli, as outlined in the second scenario of Section 3.1.2. Furthermore, it facilitates analyses that integrate both trait measurement and hypothesis testing processes, thereby accounting for both traits and their associated uncertainties, reducing bias and improving measurement precision (see Section 3.2).

Figure 5 displays the two hierarchical structural components for the example, applied to examine the effect of different teaching methods on students' writing ability. Specifically, the top branch of DAG 5b depicts the first component. This component describes how the students' writing-quality trait (T_I) and a set of relevant text-related variables (X_{IA}) , such as text length, causally influence the texts' written-quality trait (T_{IA}) . The DAG also highlights the idiosyncratic errors (e_{IA}) , which capture variations in the texts' trait that neither the students' trait nor other relevant text-related variables can explain. Here, T_I and e_{IA} are latent vectors of length n_I and $(n_I \cdot n_A)$, respectively. Additionally, X_{IA} is an observed matrix with a dimension of $(n_I \cdot n_A)$ rows and q_{IA} columns.

Furthermore, the top branch of DAG 5b describes how a set of relevant student-related variables (X_I) , including teaching methods, causally influence the students' writing-quality trait (T_I) . The DAG also includes the idiosyncratic errors (e_I) , which capture variations in the students' traits that the relevant student-related variables cannot explain. Here, e_I is a latent vector of length n_I , and X_I is an observed matrix with n_I rows and q_I columns.

$$\begin{split} O_{RPUV}^{cp} &:= f_O(D_{RPUV}) \\ D_{RPUV} &:= f_D(T_{IA}, B_{JK}) \\ T_{IA} &:= f_T(T_I, X_{IA}, e_{IA}) \\ T_I &:= f_T(X_I, e_I) \\ B_{JK} &:= f_B(B_J, Z_{JK}, e_{JK}) \\ B_J &:= f_B(Z_J, e_J) \\ e_I &\perp \{e_J, e_{IA}, e_{JK}\} \\ e_J &\perp \{e_{IA}, e_{JK}\} \\ e_{IA} &\perp e_{JK} \end{split}$$

(a) SCM

Units X_I T_I e_I Sub-units T_{IA} e_{IA} O_{RPUV}^{cp} D_{RPUV} Judgments Z_{JK}^{\bullet} B_{JK} e_{JK} Judges \widetilde{B}_J e_J Z_J

Figure 5: Conceptual-population model, final vectorized form.

(b) DAG

Similarly, the bottom branch of DAG 5b depicts the second component. This component describes how the judges' bias (B_J) and a set of relevant judgment-related variables (Z_{JK}) , such as number of judgments or time for judgment, causally influence the biases associated with each text B_{JK} . The DAG also highlights the idiosyncratic errors (e_{JK}) , which capture variations in the texts' bias that neither the judges' bias nor other relevant judgment-related variables can explain. In this description, B_J and e_{JK} are latent vectors of length n_J and $(n_J \cdot n_K)$, respectively. Additionally, Z_{JK} is an observed matrix with a dimension of $(n_J \cdot n_K)$ rows and q_{JK} columns.

Moreover, the bottom branch of DAG 5b describes how a set of relevant judge-related variables (Z_J) , including judgment expertise, causally influence the judges' bias (B_J) . The DAG also accounts for idiosyncratic errors (e_J) , capturing the variations in the judges' bias that relevant judge-related variables cannot explain. Here, e_J is a latent vector of length n_J , and Z_J is an observed matrix with n_J rows and q_J columns.

Notably, all variables and functions shown in SCM 5a and DAG 5b define the set of endogenous variables V, structural equations F, and exogenous variables E for the conceptual-population model. Additionally, the figures demonstrate that all exogenous variables are independent of one another, as indicated by the relationships $e_{IA} \perp \{e_I, e_{JK}, e_J\}$, $e_I \perp \{e_{JK}, e_J\}$ and $e_{JK} \perp e_J$.

Overall, the conceptual-population model extends Thurstone's general form by introducing key innovations to address the limitations discussed in Section 3. These enhancements include accounting for judges' biases and integrating hierarchical structural components. Nevertheless, despite its promise of enhancing measurement accuracy and precision, the model still depends on the unrealistic assumption that researchers have access to a *census* of comparative data, encompassing all repeated judgments from every judge for each pair of stimuli produced by every individual in the population. Since researchers rarely meet this assumption in practice, they must consider a more realistic scenario. The extensions to the conceptual-population model, derived from this realistic scenario, are concisely described in the sample-comparison model.

4.2. The sample-comparison model

4.2.1. The sample mechanism

In the sample-comparison model, we assume a more realistic scenario compared to the conceptualpopulation model. Specifically, we assume that researchers have access to all comparisons from a limited number of judgments (n_K^s) from a sample of judges (n_J^s) , along with a specific number of texts (n_A^s) from a sample of students (n_I^s) drawn from the conceptual population. Figure 6 illustrates this model, incorporating only this sample mechanism.

Following the convention of McElreath (2020, pp. 499-516), the DAG 6b depicts the conceptual-population outcome O_{RPUV}^{cp} as an unobserved variable, emphasizing the researchers' inability to access such an outcome due to the sampling mechanism. Moreover, the DAG shows the conditioned sample design variable S, which causally influence the sample-comparison outcome O_{RPUV}^{sc} . This variable reflects how this sample-comparison outcome is a sample drawn from the conceptual-population outcome O_{RPUV}^{cp} . Specifically, the variable is a vector $S = \{s_1, \dots, s_n\}$, where n corresponds to the number of rows in the matrix RPUV (refer to Section 4.1.2), and each s_i is a binary variable indicating the presence (1) or absence (0) of data in the corresponding positions of the conceptual-population outcome O_{RPUV}^{cp} . Specifically,

$$s_i = \begin{cases} 1 & \text{if the data for the } i^{th} \text{ element of } O_{RPUV}^{cp} \text{ is observed} \\ 0 & \text{if the data for the } i^{th} \text{ element of } O_{RPUV}^{cp} \text{ is missing} \end{cases}$$

$$\begin{split} O^{sc}_{RPUV} &:= f_S(O^{cp}_{RPUV}, S) \\ O^{cp}_{RPUV} &:= f_O(D_{RPUV}) \\ D_{RPUV} &:= f_D(T_{IA}, B_{JK}) \\ T_{IA} &:= f_T(T_I, X_{IA}, e_{IA}) \\ T_I &:= f_T(X_I, e_I) \\ B_{JK} &:= f_B(B_J, Z_{JK}, e_{JK}) \\ B_J &:= f_B(Z_J, e_J) \\ e_I &\perp \{e_J, e_{IA}, e_{JK}\} \\ e_J &\perp \{e_{IA}, e_{JK}\} \\ e_{IA} &\perp e_{JK} \end{split}$$

(a) SCM

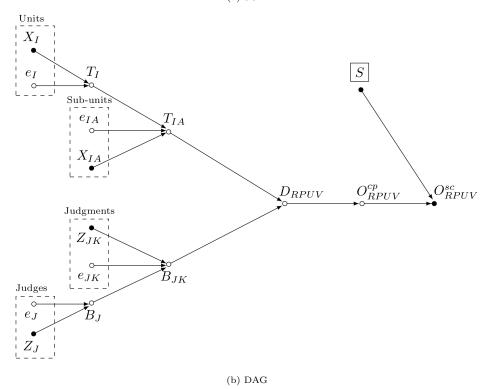


Figure 6: Sample-comparison model including sampling mechanism

Notice,

4.2.2. The comparison mechanism

Conversely, we make the scenario depicted in Section 4.2.1 more realistic by integrating the comparison mechanism, that is, we assume that researchers do not have access to all comparisons from the sample, but rather a sufficient number of comparisons (n_C) so that a proportion, P(B > A), may be determined for each pair of stimuli (Thurstone, 1927a, pp. 267). Figure 7 illustrates this model, incorporating both the sample and comparison mechanisms.

Following the convention of McElreath (2020, pp. 499-516), the DAG 7b depicts the sample-comparison outcome O_{RPUV}^{sc} as an unobserved variable, emphasizing the researchers' inability to access such an outcome due to the comparison mechanism. Moreover, the DAG shows the conditioned comparison design variable C, which causally influence the observed outcome O_{RPUV} . This variable reflects how this observed outcome is a sample drawn from the sample-comparison outcome O_{RPUV}^{sp} . Specifically, the variable is a vector $C = \{c_1, \dots, c_n\}$, where n corresponds to the number of rows in the matrix RPUV (refer to Section 4.1.2), and each c_i is a binary variable indicating the presence (1) or absence (0) of data in the corresponding positions of the sample-comparison outcome O_{RPUV}^{sp} . Specifically,

$$c_i = \begin{cases} 1 & \text{if the data for the } i^{th} \text{ element of } O_{RPUV}^{sp} \text{ is observed} \\ 0 & \text{if the data for the } i^{th} \text{ element of } O_{RPUV}^{sp} \text{ is missing} \end{cases}$$

$$\begin{split} O_{RPUV} &:= f_C(O^{sc}_{RPUV}, C) \\ O^{sc}_{RPUV} &:= f_S(O^{cp}_{RPUV}, S) \\ O^{cp}_{RPUV} &:= f_O(D_{RPUV}) \\ D_{RPUV} &:= f_D(T_{IA}, B_{JK}) \\ T_{IA} &:= f_T(T_I, X_{IA}, e_{IA}) \\ T_I &:= f_T(X_I, e_I) \\ B_{JK} &:= f_B(B_J, Z_{JK}, e_{JK}) \\ B_J &:= f_B(Z_J, e_J) \\ e_I &\perp \{e_J, e_{IA}, e_{JK}\} \\ e_J &\perp \{e_{IA}, e_{JK}\} \\ e_{IA} &\perp e_{JK} \end{split}$$

(a) SCM

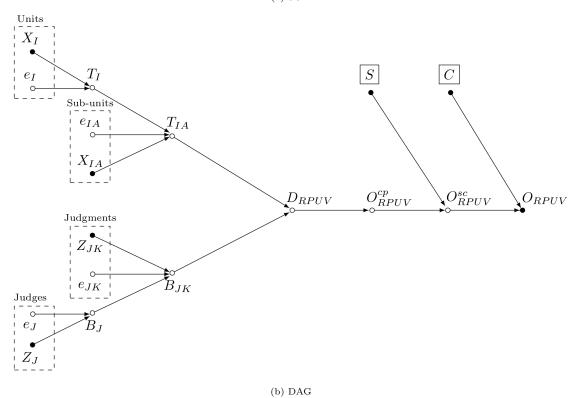


Figure 7: Sample-comparison model including sample and comparison mechanisms

Notice,

5. Abandoning the BTL model

$$\begin{aligned} O_{RPUV} &= O_{RPUV}^{sc} \odot C \\ O_{RPUV}^{sc} &= O_{RPUV}^{sc} \odot S \\ O_{RPUV}^{cp} &\sim \operatorname{Bernoulli}\left[\operatorname{inv_logit}(D_{RPUV})\right] \\ O_{RPUV} &= (T_{IA}[R_1, P_1] - T_{IA}[R_2, P_2]) + B_{JK}[U, V] \\ T_{IA} &= T_I + \beta_{XA}X_{IA} + e_{IA} \\ T_I &= \beta_{XI}X_I + e_I \\ B_{JK} &= B_J + \beta_{ZK}Z_{JK} + e_{JK} \\ B_J &= \beta_{ZJ}Z_J + e_J \\ e_I &\sim \operatorname{Normal}(0, 1) \\ e_{IA} &\sim \operatorname{Normal}(0, p_{IA}) \\ e_{JK} &\sim \operatorname{Normal}(0, p_{JK}) \end{aligned} \qquad \begin{aligned} O_{RPUV}^{sc} &= f_C(O_{RPUV}^{sc}, C) \\ O_{RPUV}^{sc} &= f_S(O_{RPUV}^{cp}, C) \\ O_{RPUV}^{sc} &= f_S(O_{RPUV}^{sc}, C) \\ O_{RPUV}^{sc} &= f_S(O_{$$

Figure 8: Sample-comparison model, including sample and comparison mechanisms and assuming equal discriminal dispersions for the student's traits

$$\begin{split} O_{RPUV} &= O_{RPUV}^{sc} \odot C \\ O_{RPUV}^{sc} &= O_{RPUV}^{cp} \odot S \\ O_{RPUV}^{cp} \sim \text{Bernoulli} \left[\text{inv_logit}(D_{RPUV}) \right] \\ O_{RPUV} &\sim \text{Bernoulli} \left[\text{inv_logit}(D_{RPUV}) \right] \\ O_{RPUV} &= (T_{IA}[R_1, P_1] - T_{IA}[R_2, P_2]) + B_{JK}[U, V] \\ T_{IA} &= T_I + \beta_{XA} X_{IA} + e_{IA} \\ T_I &= \beta_{XI} X_I + e_I \\ B_{JK} &= B_J + \beta_{ZK} Z_{JK} + e_{JK} \\ B_J &= \beta_{ZJ} Z_J + e_J \\ e_I &\sim \text{Normal}(0, s_{XI}) \\ e_J &\sim \text{Normal}(0, s_{ZJ}) \\ e_{IA} &\sim \text{Normal}(0, p_{IA}) \\ e_{JK} &\sim \text{Normal}(0, p_{JK}) \\ \\ &\leftarrow Contraints: \\ \sum_{g=1}^{G_{XI}} s_{XI}/G_{XI} = 1; \quad \sum_{g=1}^{G_{ZJ}} s_{ZJ}/G_{ZJ} = 1 \\ \end{split} \label{eq:optimizero} O_{RPUV} &:= f_C(O_{RPUV}^{sc}, C) \\ O_{RPUV}^{sc} &:= f_S(O_{RPUV}^{cp}, C) \\ O_{RPUV}^{sc} &:= f_S(O_{RPUV}^$$

Figure 9: Sample-comparison model, including sample and comparison mechanisms and assuming different discriminal dispersions for the student's traits

6. Discussion

- 6.1. Findings
- 6.2. Limitations and further research

(a) Statistical model

7. Conclusion

Declarations

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8. Appendix

8.1. Statistical and Causal inference

This section introduces fundamental statistical and causal inference concepts necessary for understanding the core theoretical principles described in this document. It does not, however, offer a comprehensive overview of causal inference methods. Readers seeking more in-depth understanding may wish to explore introductory papers such as Pearl (2010), Rohrer (2018), Pearl (2019), and Cinelli et al. (2020). They may also find it helpful to consult introductory books like Pearl and Mackenzie (2018), Neal (2020), and McElreath (2020). For more advanced study, readers may refer to seminal intermediate papers such as Neyman (1923), Rubin (1974), Spirtes et al. (1991), and Sekhon (2009), as well as books such as Pearl (2009), Morgan and Winship (2014), and Hernán and Robins (2020).

8.1.1. Empirical research and randomized experiments

Empirical research uses evidence from observation and experimentation to address real-world challenges. In this context, researchers typically formulate their research questions as estimands or targets of inference, i.e., the specific quantities they seek to determine (Everitt and Skrondal, 2010). For instance, researchers might be interested in answering the following question: "To what extent do different teaching methods (T) influence students' ability to produce high-quality written texts (Y)?" To investigate this, researchers could randomly assign students to two groups, each exposed to a different teaching method $(T_i = \{1,2\})$. Then, they would perform pairwise comparisons, generating a dichotomous outcome $(Y_i = \{0,1\})$ showing which student exhibits more of the ability. In this scenario, the research question can be rephrased as the estimand, "On average, is there a difference in the ability to produce high-quality written texts between the two groups of students?" and this estimand can be mathematically represented by the random associational quantity in Equation 3, where $E[\cdot]$ denotes the expected value.

$$E[Y_i \mid T_i = 1] - E[Y_i \mid T_i = 2] \tag{3}$$

Researchers then proceed to identify the estimands. *Identification* determines whether an estimator can accurately compute the estimand based solely on its assumptions, regardless of random variability (Schuessler and Selb, 2023, pp. 4). An *estimator* refers to a method or function that transforms data into an estimate (Neal, 2020). *Estimates* are numerical values that approximate

the estimand derived through the process of *estimation*, which integrates data with an estimator (Everitt and Skrondal, 2010). The Identification-Estimation flowchart (McElreath, 2020; Neal, 2020), shown in Figure 10, visually represents the transition from estimands to estimates.



Figure 10: Identification-Estimation flowchart. Extracted and slightly modified from Neal (2020, pp. 32)

Identification is a necessary condition to ensure *consistent* estimators. An estimator achieves *consistency* when it converges to the "true" value of an estimand as the data size approaches infinity (Everitt and Skrondal, 2010). Without identification, researchers cannot achieve consistency, even with "infinite" and error-free data. As a result, deriving meaningful insights about an estimand from finite data becomes impossible (Schuessler and Selb, 2023, pp. 5). Therefore, to ensure accurate and reliable estimates, researchers prioritize estimators with desirable identification properties. For instance, the Z-test is a widely used estimator for comparing group proportions, yielding accurate estimates when its underlying assumptions are satisfied (Kanji, 2006). Furthermore, researchers can interpret estimates from the Z-test as causal, provided the data is collected through a randomized experiment.

Randomized experiments are widely recognized as the gold standard in evidence-based science (Hariton and Locascio, 2018; Hansson, 2014). This recognition stems from their ability to enable researchers interpret associational estimates as causal. They achieve this by ensuring data, and by extension an estimator, satisfies several key identification properties, such as common support, no interference, and consistency (Morgan and Winship, 2014; Neal, 2020). The most critical property, however, is the elimination of confounding. Confounding occurs when an external variable X simultaneously influences the outcome Y and the variable of interest T, resulting in spurious associations (Everitt and Skrondal, 2010). Randomization addresses this issue by decoupling the association between the intervention allocation T from any other variable X (Morgan and Winship, 2014; Neal, 2020).

Nevertheless, researchers often face constraints that limit their ability to conduct randomized experiments. These constraints include ethical concerns, such as the assignment of individuals to potentially harmful interventions, and practical limitations, such as the infeasibility of, for example, assigning individuals to genetic modifications or physical impairments (Neal, 2020). In these cases, causal inference offers a valuable alternative for generating causal estimates and understanding the mechanisms underlying specific data. In addition, the framework can provide significant theoretical insights that can enhance the design of experimental and observational studies (McElreath, 2020).

8.1.2. Identification under causal inference

Unlike classical statistical modeling, which focuses primarily on summarizing data and inferring associations, the *causal inference* framework is designed to identify causes and estimate their effects using data (Shaughnessy et al., 2010; Neal, 2020). The framework uses rigorous mathematical techniques to address the *fundamental problem of causality* (Pearl, 2009; Pearl et al., 2016; Morgan and Winship, 2014). This problem revolves around the question, "What would have happened 'in the world' under different circumstances?" This question introduces the concept of counterfactuals, which are instrumental in defining and identifying causal effects.

Counterfactuals are hypothetical scenarios that are contrary to fact, where alternative outcomes resulting from a given cause are neither observed nor observable (Neal, 2020; Counterfactual, 2024). The structural approach to causal inference (Pearl, 2009; Pearl et al., 2016) provides a formal framework for defining counterfactuals. For instance, in the scenario described in Section 8.1.1, the approach begins by defining the *individual causal effect* (ICE) as the difference between each student's potential outcomes, as in Equation 4.

$$\tau_i = Y_i \mid do(T_i = 1) - Y_i \mid do(T_i = 2) \tag{4}$$

where $do(T_i = t)$ represents the intervention operator, $Y_i \mid do(T_i = 1)$ represents the potential outcome under intervention $T_i = 1$, and $Y_i \mid do(T_i = 1)$ represents the potential outcome under intervention $T_i = 2$. Here, an intervention involves assigning a constant value to the treatment variable for each student's potential outcomes. Note that if a student is assigned to intervention $T_i = 1$, the potential outcome under $T_i = 2$ becomes a counterfactual, as it is no longer observed nor observable. To address this challenge, the structural approach extends the ICE to the average causal effect (ACE, Equation 5), representing the average difference between the students' observed potential outcomes and their counterfactual counterparts.

$$\begin{split} \tau &= E[\tau_i] \\ &= E[Y_i \mid do(T_i = 1)] - E[Y_i \mid do(T_i = 2)] \end{split} \tag{5}$$

Even though counterfactuals are unobservable, researchers can still identify the ACE from associational estimates by leveraging the structural approach. The approach identifies the ACE by statistically conditioning data on a *sufficient adjustment set* of variables X (Pearl, 2009; Pearl et al., 2016; Morgan and Winship, 2014). This *sufficient* set (potentially empty) must block all non-causal paths between T to Y without opening new ones. When such a set exists, then T and Y are d-separated by X ($T \perp Y \mid X$) (Pearl, 2009), and X satisfies the backdoor criterion (Neal, 2020, pp 37). Here, conditioning describes the process of restricting the focus to the subset of the population defined by the conditioning variable (Neal, 2020, pp. 32) (see Equation 6).

Conditioning on a sufficient adjustment set enables researchers to estimate the ACE, even when the data comes from an observational study. This process is feasible because such conditioning ensures that the ACE estimator satisfies several critical properties, including confounding elimination (Morgan and Winship, 2014). Naturally, the validity of claims about the causal effects of T on Y now hinges on the assumption that X serves as a sufficient adjustment set. However, as Kohler et al. (2019, pp. 150) noted, drawing conclusions about the real world from observed data inevitably requires assumptions. This requirement holds true for both observational and experimental data.

For instance, if researchers cannot conduct the randomized experiments described in Section 8.1.1 and must instead rely on observational data, they can still identify the ACE as long as an observed variable X, such as the socio-economic status of the school, satisfies the backdoor criterion. Under these circumstances, researchers first identify the *conditional average causal effect* (CACE, Equation 6)

$$CACE_t = E[Y_i \mid T_i = t, X] \tag{6}$$

From the CACE, researchers can identify the ACE from associational quantities as in Equation 7. This identification process is commonly known as the *backdoor adjustment*. Here, $E_X[\cdot]$ represents the marginal expected value over X (Morgan and Winship, 2014).

$$\begin{split} \tau &= E[Y_i \mid do(T_i = 1)] - E[Y_i \mid do(T_i = 2)] \\ &= E_X[CACE_1 - CACE_2] \\ &= E_X\left[E[Y_i \mid T_i = 1, X] - E[Y_i \mid T_i = 2, X]\right] \end{split} \tag{7}$$

Notably, the approach extends the ACE identification for a continuous variable T as in Equation 8, ensuring broad applicability across different causal scenarios (Neal, 2020, pp. 45)

$$\begin{split} \tau &= E[Y_i \mid do(T_i = t)] \\ &= dE_X \left[E[Y_i \mid T_i = t, X] \right] / dt \end{split} \tag{8}$$

8.1.3. Diving into the specifics

The structural approach to causal inference uses SCMs and DAGs to formally and graphically represent the presumed causal structure underlying the ACE (Pearl, 2009; Pearl et al., 2016; Gross et al., 2018; Neal, 2020). Essentially, these tools serve as conceptual (theoretical) models on which identification analysis rests (Schuessler and Selb, 2023, pp. 4). Thus, using these tools, researchers can determine which statistical models can identify (ACE, CACE, or other), assuming the depicted causal structure is correct (McElreath, 2020), thus enabling valid causal inference. Figure 10 shows the role of theoretical models in the inference process.

SCMs and DAGs support identification analysis through two key advantages. First, regardless of complexity, they can represent various causal structures using only five fundamental building blocks (Neal, 2020; McElreath, 2020). This feature allows researchers to decompose complex structures into manageable components, facilitating their analysis (McElreath, 2020). Second, they depict causal relationships in a non-parametric and fully interactive way. This flexibility enables feasible ACE identification strategies without defining the variables' data types, the functional form between them, or their parameters (Pearl et al., 2016, pp. 35).

Thus, Section 8.1.3.1 and Section 8.1.3.2 elaborate on the first advantage, while Section 8.1.3.2 and Section 8.1.3.3 do so for the second. Finally, Section 8.1.3.4 explains how researchers use SCMs and DAGs alongside Bayesian inference methods in the estimation process.

8.1.3.1. The five fundamental block for SCMs and DAGs.

Figures 11, 12, 13, 14, and 15 display the five fundamental building blocks for SCMs and DAGs. The left panels of the figures show the formal mathematical models, represented by the SCMs, defined in terms of a set of endogenous variables $V = \{X_1, X_2, X_3\}$, a set of exogenous variables $E = \{e_{X1}, e_{X2}, e_{X3}\}$, and a set of functions $F = \{f_{X1}, f_{X2}, f_{X3}\}$ (Pearl, 2009; Cinelli et al., 2020). Endogenous variables are those whose causal mechanisms a researcher chooses to model (Neal, 2020). In contrast, exogenous variables represent errors or disturbances arising from omitted factors that the investigator chooses not to model explicitly (Pearl, 2009, pp. 27,68). Lastly, the functions, referred to as structural equations, express the endogenous variables as non-parametric functions of other variables. These functions use the symbol ':=' to denote the asymmetrical causal dependence of the variables and the symbol ' \bot ' to represent d-separation, a concept akin to (conditional) independence.

Notably, every SCM has an associated DAG (Pearl et al., 2016; Cinelli et al., 2020). The right panels of the figures display these DAGs. A DAG is a graph consisting of nodes connected by edges, where the nodes represent random variables. The term directed means that the edges extend from one node to another, with arrows indicating the direction of causal influence. The term acyclic implies that the causal influences do not form loops, ensuring the influences do not cycle back on themselves (McElreath, 2020). DAGs represent observed variables as solid black circles, while they use open circles for unobserved (latent) variables (Morgan and Winship, 2014). Although the standard representation of DAGs typically omits exogenous variables for simplicity, the magnified representation depicted in the figures offers one key advantage: including exogenous variables can help researchers highlight potential issues related to conditioning and confounding (Cinelli et al., 2020).

$$X_1:=f_{X1}(e_{X1})$$

$$X_3:=f_{X3}(e_{X3})$$

$$e_{X1}\perp e_{X3}$$
 (b) DAG

Figure 11: Two unconnected nodes

$$X_1 := f_{X1}(e_{X1})$$

$$X_3 := f_{X3}(X_1, e_{X3})$$

$$e_{X1} \perp e_{X3}$$
 (b) DAG

(a) SCM

Figure 12: Two connected nodes or descendant

$$X_1 := f_{X1}(e_{X1})$$

$$X_2 := f_{X2}(X_1, e_{X2})$$

$$X_3 := f_{X3}(X_2, e_{X3})$$

$$e_{X1} \perp e_{X2}$$

$$e_{X1} \perp e_{X3}$$
 (b) DAG
$$e_{X2} \perp e_{X3}$$

Figure 13: Chain or mediator

$$X_1 := f_{X1}(X_2, e_{X1})$$

$$X_2 := f_{X2}(e_{X2})$$

$$X_3 := f_{X3}(X_2, e_{X3})$$

$$e_{X1} \perp e_{X2}$$

$$e_{X1} \perp e_{X3}$$
 (b) DAG
$$e_{X2} \perp e_{X3}$$
 (a) SCM

Figure 14: Fork or confounder

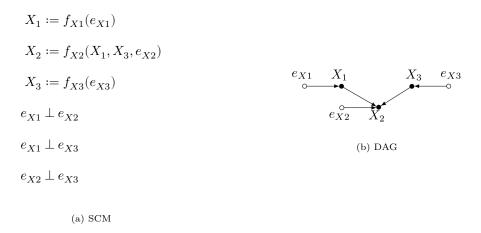


Figure 15: Collider or inmorality

A careful examination of these building blocks highlights the theoretical assumptions underlying their observed variables. SCM 11a and DAG 11b depict two unconnected nodes, representing a scenario where variables X_1 and X_3 are independent or not causally related. SCM 12a and DAG 12b illustrate two connected nodes, representing a scenario where a parent node X_1 exerts a causal influence on a child node X_3 . In this setup, X_3 is considered a descendant of X_1 . Additionally, X_1 and X_3 are described as adjacent because there is a direct path connecting them. SCM 13a and DAG 13b depict a chain, where X_1 influences X_2 , and X_2 influences X_3 . In this configuration, X_1 is a parent node of X_2 , which is a parent node of X_3 . This structure creates a directed path between X_1 and X_3 . Consequently, X_1 is an ancestor of X_3 , and X_2 fully mediates the relationship between the two. SCM 14a and DAG 14b illustrate a fork, where variables X_1 and X_3 are both influenced by X_2 . Here, X_2 is a parent node that *confounds* the relationship between X_1 and X_3 . Finally, SCM 15a and DAG 15b show a *collider*, where variables X_1 and X_3 are concurrent causes of X_2 . In this configuration, X_1 and X_3 are not causally related to each other but both influence X_2 (an inmorality). Notably, all building blocks assume the errors are independent of each other and from all other variables in the graph, as evidenced by the pairwise relations $e_{X1} \perp e_{X2}$, $e_{X1} \perp e_{X3}$, and $e_{X2} \perp e_{X3}$.

Researchers can then use these building blocks to represent the scenario outlined in Section 8.1.2. SCM 16a and DAG 16b depict the plausible causal structure for this example. In this context, the variable X (socio-economic status of the school) is thought to be a confounder in the relationship between the teaching method T and the outcome Y. The figures display multiple descendant relationships such as $X \to T$, $X \to Y$, and $T \to Y$. They also highlight unconnected node pairs,

evident from the relationships $e_T \perp e_X$, $e_T \perp e_Y$, and $e_X \perp e_Y$. Additional, the figures show one fork, $X \to \{T,Y\}$, and two colliders: $\{X,e_T\} \to T$ and $\{X,T,e_Y\} \to Y$.

$$X := f_X(e_X)$$

$$T := f_T(X, e_T)$$

$$Y := f_Y(T, X, e_Y)$$

$$e_T \perp e_X$$

$$e_T \perp e_Y$$
 (b) DAG
$$e_X \perp e_Y$$
 (a) SCM

Figure 16: Plausible causal structure the scenario outlined in Section 8.1.2.

8.1.3.2. The probabilistic implications of these blocks.

Beyond their graphical capabilities, SCMs and DAGs can encode the probabilistic information embedded within a causal structure. They achieve this encoding by relying on three fundamental assumptions: the local Markov, the minimality, the causal edges assumption. The local Markov assumption encodes probabilistic independencies between variables by declaring that nodes in a graph are independent of all its non-descendants, given its parents (Neal, 2020, pp. 20). Meanwhile, the minimality assumption encodes probabilistic dependencies among variables by stating that every pair of adjacent nodes exhibits a dependency (Neal, 2020, pp. 21). Finally, the causal edges assumption encodes causal relationships between variables by declaring that each parent node acts as a direct cause of its children (Neal, 2020, pp. 22). Figure 17 illustrates how these assumptions influence the statistical and causal interpretations of graphs.

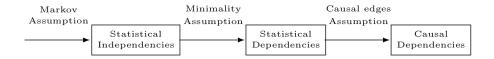


Figure 17: The flow of association and causation in graphs. Extracted and slightly modified from Neal (2020, pp. 31)

A notable implication of the assumptions underlying the probabilistic encoding is that any conceptual model described by an SCM and DAG can represent the joint distribution of variables more efficiently (Pearl et al., 2016, pp. 29). This expression takes the form of a product of conditional probability distributions (CPDs) of the type $P(child \mid parents)$. This property is formally known

as the Bayesian Network factorization (BNF, Equation 9) (Pearl et al., 2016, pp. 29; Neal, 2020, pp. 21). In this expression, $pa(X_i)$ denotes the set of variables that are the parents of X_i .

$$\begin{split} P(X_1,X_2,\dots,X_P) &= X_1 \cdot \prod_{p=2}^P P(X_i \mid X_{i-1},\dots,X_1) \quad \text{(by chain rule)} \\ &= X_1 \cdot \prod_{p=2}^P P(X_i \mid pa(X_i)) \qquad \qquad \text{(by BNF)} \end{split}$$

This encoding enables researchers with conceptual (theoretical) knowledge in the form of an SCM and DAG to predict patterns of (in)dependencies in the data. As highlighted by Pearl et al. (2016, pp. 35), these predictions depend solely on the structure of these conceptual models without requiring the quantitative details of the equations or the distributions of the errors. Moreover, once researchers observe empirical data, the patterns of (in)dependencies in the data can provide significant insights into the validity of the proposed conceptual model.

The five fundamental building blocks described in Section 8.1.3.1 clearly illustrate which (in)dependencies can SMCs and DAGs predict. For instance, applying the BNF to the causal structure shown in the SCM 11a and DAG 11b enables researchers to express the joint probability distribution of the observed variables as $P(X_1, X_3) = P(X_1)P(X_3)$, supporting the theoretical assumption that the observed variables X_1 and X_3 are unconditionally independent $(X_1 \perp X_3)$ (Neal, 2020, pp. 24). Conversely, when X_3 is unconditionally dependent on X_1 ($X_1 \not\perp X_3$), as depicted in the SCM 12a and DAG 12b, the BNF express their joint probability distribution as $P(X_1, X_3) = P(X_3 \mid X_1)P(X_1)$. Notably, these descriptions demonstrate the clear correspondence between the structural equations illustrated in Section 8.1.3.1 and the CPDs.

Beyond the insights gained from two-node structures, researchers can uncover more nuanced patterns of (in)dependencies from chains, forks, and colliders. These (in)dependencies apply to any data set generated by a causal model with those structures, regardless of the specific functions attached to the SCM (Pearl et al., 2016, pp. 36). For instance, applying the BNF to the chain structure depicted in the SCM 13a and DAG 13b allow researchers to represent the joint distribution for the observed variables as $P(X_1, X_2, X_3) = P(X_1)P(X_2 \mid X_1)P(X_3 \mid X_2)$. This expression implies that X_1 and X_3 are unconditionally dependent $(X_1 \not\perp X_3)$, but conditionally independent when controlling for X_2 ($X_1 \perp X_3 \mid X_2$). Moreover, in the fork structure shown in the SCM 14a and DAG 14b, researchers can express the joint distribution of the observed variables as $P(X_1, X_2, X_3) = P(X_1 \mid X_2)P(X_2)P(X_3 \mid X_2)$. Similar to the chain structure, this expression allows researchers to further infer that X_1 and X_3 are unconditionally dependent $(X_1 \not\perp X_3)$, but conditionally independent when controlling for X_2 $(X_1 \perp X_3 \mid X_2)$. Finally, researchers analyzing the collider structure illustrated in the SCM 15a and DAG 15b can express the joint distribution of the observed variables as $P(X_1, X_2, X_3) = P(X_1)P(X_2 \mid X_1, X_3)P(X_3)$. This representation allows researchers to infer that X_1 and X_3 are unconditionally independent $(X_1 \perp X_3)$, but conditionally dependent when controlling for X_2 $(X_1 \not\perp X_3 \mid X_2)$. The authors Pearl et al. (2016, pp. 37, 40, 41) and Neal (2020, pp. 25–26) provide the mathematical proofs for these conclusions.

Using these additional probabilistic insights, researchers can revisit the scenario in Section 8.1.2. In this context, applying the BNF to the SCM 18a structure, enables the representation of the joint probability distribution of the observed variables as $P(Y, T, X) = P(Y \mid T, X)P(T \mid X)P(X)$. From this expression, researchers can infer that the outcome Y is unconditionally dependent on the teaching method $T(Y \not\perp T)$. This dependency arises from two key structures: a direct causal path from the teaching method T to the outcome Y, represented by the two-connected-nodes structure $T \to Y$ (black path in DAG 18b), and a confounding non-causal path from the teaching method T to the outcome Y through the socio-economic status of the school X, represented by the fork structure $T \leftarrow X \to Y$ (gray path in DAG 18b).

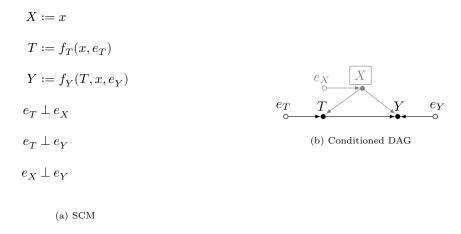


Figure 18: Plausible causal structure the scenario outlined in Section 8.1.2.

8.1.3.3. From probability to causality.

The structural approach to causal inference translates probabilistic insights into actionable strategies seeking to identify the ACE from associational quantities. The approach achieves this by relying on the *modularity assumption*, which posits that intervening on a node alters only the causal mechanism of that node, leaving others unchanged (Neal, 2020, pp. 34).

The modularity assumption underpins the concepts of manipulated graphs and Truncated Factorization, which are essential for representing interventions $P(Y_i \mid do(T_i = t))$ within SCMs and DAGs. Manipulated graphs simulate physical interventions by removing specific edges from a DAG, while preserving the remaining structure unchanged (Neal, 2020, pp. 34). In parallel, Truncated Factorization (TF) achieves a similar simulation by removing specific functions from the conceptual model and replacing them with constants, while keeping the rest of the structure unchanged (Pearl, 2010). The probabilistic implications of this factorization are formalized in Equation 10, where S represents the subset of variables X_p directly influenced by the intervention, while an example illustrating these concepts follows below.

$$P(X_1, X_2, \dots, X_P \mid do(S)) = \begin{cases} \prod P(X_p \mid pa(X_p)) & \text{if } p \notin S \\ 1 & \text{otherwise} \end{cases}$$
 (10)

Using the TF, researchers can define the backdoor adjustment to identify the ACE. This adjustment states that if a variable $X_p \in S$ serves as a sufficient adjustment set for the effect of X_a on X_b , then the ACE can be identified using Equation 11. The sufficient adjustment set (potentially empty) must block all non-causal paths between X_a and X_b without introducing new paths. If such a set exists, then X_a and X_b are d-separated by X_p ($X_a \perp X_b \mid X_p$) (Pearl, 2009), and X_p satisfies the backdoor criterion (Neal, 2020, pp. 37).

$$P(X_a \mid do(X_b = x)) = \sum_{X_p} P(X_a \mid X_b = x, X_p) P(X_p) \tag{11} \label{eq:11}$$

Ultimately, the backdoor adjustment enables researchers to express the ACE as:

$$\begin{split} \tau &= E[X_a \mid do(X_b = 1)] - E[X_a \mid do(X_b = 2)] \\ &= E_{Xp} \left[E[X_a \mid do(X_b = 1), X_p] - E[X_a \mid do(X_b = 2), X_p] \right] \\ &= \sum_{Xp} X_a \cdot P(X_a \mid X_b = 1, X_p) \cdot P(X_p) - \sum_{Xp} X_a \cdot P(X_a \mid X_b = 2, X_p) \cdot P(X_p) \end{split} \tag{12}$$

With these new insights, researchers revisiting the scenario in Section 8.1.3.2 can infer that the socio-economic status of the school, X, satisfies the backdoor criterion, assuming the causal structure depicted by the SCM 18a and DAG 18b is correct. This means that X serves as a sufficient adjustment set, as it effectively blocks all confounding non-causal paths introduced by the fork

structure. Nevertheless, since Y remains dependent on T even after conditioning $(Y \not\perp T \mid X)$, this dependency can only be attributed to the direct causal effect $T \to Y$. Notably, for the purpose of identification, the conditioned DAG 18b is equivalent to the manipulated DAG 19b, because X satisfies the backdoor criterion.

$$X:=f_X(e_X)$$

$$T:=t$$

$$Y:=f_Y(t,X,e_Y)$$

$$e_T\perp e_X$$

$$t \qquad T \qquad Y \qquad e_Y$$

$$e_T\perp e_Y$$
 (b) Manipulated DAG
$$e_X\perp e_Y$$

Figure 19: Plausible causal structure the scenario outlined in Section 8.1.3.2.

Researchers can then apply the backdoor adjustment to identify the ACE of T on Y. They achieve this by first identifying the CACE of T on Y by conditioning on X, and then marginalizing this effect over X to obtain the ACE. This process is expressed in Equation 13 (see Section 8.1.2).

$$\begin{split} \tau &= E[Y_i \mid do(T_i = 1)] - E[Y_i \mid do(T_i = 2)] \\ &= E_X \left[E[Y_i \mid T_i = 1, X] - E[Y_i \mid T_i = 2, X] \right] \\ &= \sum_X Y_i \cdot P(Y_i \mid T_i = 1, X) \cdot P(X) - \sum_X Y_i \cdot P(Y_i \mid T_i = 2, X) \cdot P(X) \end{split} \tag{13}$$

8.1.3.4. The estimation process.

Ultimately, researchers can use Bayesian inference methods to estimate the ACE. The approach begins by defining two probability distributions: the likelihood of the data, $P(X_1, X_2, ..., X_P | \theta)$, and the prior distribution, $P(\theta)$ (Everitt and Skrondal, 2010), where X_P represents a random variable, and θ represents a one-dimensional parameter space for simplicity. After observing empirical data, researchers can update the priors to posterior distributions using Bayes' rule in Equation 14:

$$P(\theta \mid X_1, X_2, \dots, X_P) = \frac{P(X_1, X_2, \dots, X_P \mid \theta) \cdot P(\theta)}{P(X_1, X_2, \dots, X_P)} \tag{14}$$

Given that the denominator on the right-hand side of Equation 14 serves as a normalizing constant independent of the parameter θ , researchers can simplify the posterior updating process into three steps. First, they integrate new empirical data through the likelihood. Second, they update the parameters' priors to a posterior distribution according to Equation 15. Ultimately, they normalize these results to obtain a valid probability distribution.

$$P(\theta \mid X_1, X_2, \dots, X_P) \propto P(X_1, X_2, \dots, X_P \mid \theta) \cdot P(\theta) \tag{15}$$

Temporarily setting aside the definition of prior distributions $P(\theta)$, note that the posterior updating process depends heavily on the assumptions underlying the likelihood of the data. However, as the number of random variables, P, increases, this joint distribution quickly becomes intractable (Neal, 2020). This intractability is evident from Equation 16, where the likelihood distribution is expressed by multiple chained CPDs.

$$P(X_1, X_2, \dots, X_P \mid \theta) = P(X_1 \mid \theta) \prod_{p=2}^{P} P(X_i \mid X_{i-1}, \dots, X_1, \theta)$$
 (16)

Nevertheless, researchers can manage the complexity of the likelihood by assuming specific local (in)dependencies among variables. SCMs and DAGs provide a formal framework to represent these assumptions, as detailed in Section 8.1.3.2. These assumptions improve model tractability and simplify the estimation process by enabling the derivation of the BNF of the likelihood (Equation 17). With this simplified structure, any probabilistic programming language can model the system and compute the parameter's posterior distribution using Equation 14.

$$P(X_1, X_2, \dots, X_P \mid \theta) = P(X_1 \mid \theta) \prod_{p=2}^{P} P(X_i \mid pa(X_i), \theta) \tag{17}$$

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