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# Causal Feature Selection via Orthogonal Search

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## Abstract

The problem of inferring the direct causal parents of a response variable among a large set of explanatory variables is of high practical importance in many disciplines. Recent work in the field of causal discovery exploits invariance properties of models across different experimental conditions for detecting direct causal links. However, these approaches generally do not scale well with the number of explanatory variables, are difficult to extend to nonlinear relationships, and require data across different experiments. Inspired by *Debiased* machine learning methods, we study a one-vs.-the-rest feature selection approach to discover the direct causal parent of the response. We propose an algorithm that works for purely observational data, while also offering theoretical guarantees, including the case of partially nonlinear relationships. Requiring only one estimation for each variable, we can apply our approach even to large graphs, demonstrating significant improvements compared to established approaches.

## 1 Introduction

Identifying causal relationships is a profound and hard problem pervading the sciences. While randomized controlled intervention studies are considered the gold standard, they are in many cases ruled out by financial or ethical concerns [26, 36]. In order to improve understanding of systems and help design relevant interventions, the subset of causes which have a direct effect (*direct causes* / *direct causal parents*) need to be identified.

Let us consider the setup described in Figure 1, corresponding to a linear system only for the response  $Y$ ,

$$Y = \langle \theta, X \rangle + U. \quad (1)$$

where  $U$  is an independent noise variable with zero mean,  $\theta, X \in \mathbb{R}^d$ ,  $Y \in \mathbb{R}$  and  $\langle \cdot, \cdot \rangle$  denotes the inner product. We investigate how to find the direct causes of  $Y$  among a high-dimensional vector of covariates  $X$ , where the covariates can have arbitrary non-linear relationships between

them. From our formulation, a given entry of  $\theta$  should be non-zero if and only if the variable corresponding to that particular coefficient is a direct causal parent [28], e.g.,  $X_1$  and  $X_2$  in Figure 1. We restrict ourselves to the setting of *linear direct causes* of  $Y$  (LDC) and *no feature descending from  $Y$*  (NFD). LDC is justified as an approximation when the effects of each causal feature are weak such the possibly non-linear effects can be linearized; NFD is justified in some applications where we can exclude any influence of  $Y$  on a covariate. This is for example the case when  $X$  are genetic factors, and  $Y$  is a particular trait/phenotype.

While applicable to full graph discovery rather than the simplified problem of finding causal parents, state of the art methods for causal discovery often rely on strong assumptions or the availability of interventional data, and have prohibitive computational cost as explained in section 1.1 in more detail. In addition to and despite their strong assumptions, causal discovery methods may perform worse than simple regression baselines [16, 18, 42].

While plain regression techniques have appealing computational cost, they come without guarantees. When using unregularized least-square regression to estimate  $\theta$ , there can be infinitely many possible choices for  $\theta$  recovered with equivalent prediction accuracy for regressing  $Y$ , in case of over-parametrized models. However, none of these choices provide any information about the features which, when intervened upon, directly cause the output variable  $Y$ . On the other hand, when using a regularized method such as Lasso, a critical issue is the bias induced by regularization [20].

When knowing the distinction between covariates and direct causes, Double ML approaches [7] have shown promising bias compensation results in the context of high dimensional observed confounding of a single variable. In the present paper, we generalize them to the problem of finding direct causes. Our key contributions are:

- We show that under the assumption that no feature of  $X$  is a child of  $Y$ , the Double ML [8] principle can be applied in an iterative and parallel way to find the subset of direct causes, even in the challenging case of only having observational data.
- Our approach has low computational complexity requirement ( $\mathcal{O}(d)$  in terms of number of variables), outperforming the usual  $\mathcal{O}(2^d)$ .
- Our method provides asymptotic guarantees that the set can be recovered exactly. Importantly, this result is valid without additional assumptions, and especially neither requires linear interactions among the covariates, nor faithfulness.
- Extensive experimental results demonstrate state-of-the-art performance of our method. Our approach significantly outperforms all other methods, especially in the case of non-linear interactions between covariates, despite relying only on linear projection.

## 1.1 Related work

The question of finding direct causal parents is also addressed in the literature as mediation analysis [3, 15, 34]. Several principled approaches have been proposed (relying for instance on Instrumental Variables (IVs)) [1, 2, 4] to test for a single direct effect in the context of specific causal graphs. Extensions of the IV based approach to generalized IVs based approaches [5, 39] are the closest known result to discover direct causal parents. However, no algorithm is provided in [5] to identify the instrumental set. Subsequently, an algorithm is provided in [39] for discovering the instrumental set in the simple setting where all the interactions are linear and the graph is acyclic. In contrast, our method allows non-linear interaction amongst the variables.

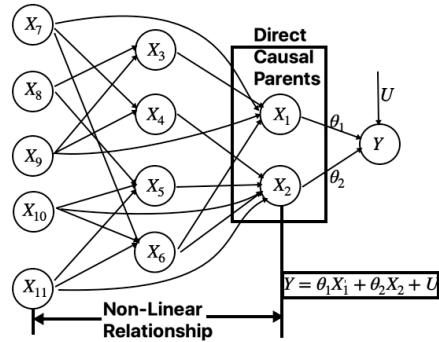


Figure 1: Graphical representation of Causal Feature Selection in our setting. Out of variables  $\{X_1, \dots, X_{11}\}$ , only  $X_1$  and  $X_2$  are direct causal parents of  $Y$  such that  $Y = \theta_1 X_1 + \theta_2 X_2 + U$ ,  $U$  being an independent zero-mean noise. We propose an approach to find  $X_1$  and  $X_2$  in this paper under assumptions discussed in the text. An example of this setup in the real-world is finding genes which directly cause a phenotype.

Several other works have also tried to address the problem of discovering causal features. The authors review work on causal feature selection in [14]. More recent papers on causal feature selection have appeared since [6, 25, 40], but none of those claims to recover all the direct causal parents asymptotically or non-asymptotically as we do in our case.

There has been another line of works on inferring causal relationships from observational data most of which require strong assumptions, such as faithfulness [21, 26, 36]. Classical approaches along these lines include the PC-algorithm [36], which can only reconstruct the network up to a Markov equivalence class. Another approach is to restrict the class of interactions among the covariates and the functional form of the signal-noise mixing (typically considered additive) or the distribution (e.g., non-Gaussianity) to achieve identifiability (see [17, 30]); this includes linear approaches like LiNGAM [33] and nonlinear generalizations with additive noise [29]. For a recent review of the empirical performance of structure learning algorithms and a detailed description of causal discovery methods, we refer to [16]. Recently, there have been several attempts at solving the problem of causal inference by exploiting the invariance of a prediction under a causal model given different experimental settings [13, 27]. The computational cost to run both of the algorithms is exponential in the number of variables, when aiming to discover the full causal graph.

Our method mainly takes inspiration from Debiased/Double ML method [7] which utilizes the concept of orthogonalization to overcome the bias introduced due to regularization. We will discuss this in detail in the next section. Considering specific example, the Lasso suffers from the fact that the estimated coefficients are shrunk towards zero which is undesirable [37]. To overcome this limitation, an approach of debiasing the Lasso was proposed in several papers [19, 20, 41]. However, unlike our approach, Debiased Lasso methods don't recover all the non-zero coefficients of the parameter vector  $\theta$  under generic assumptions we have in the paper.

## 2 Methodology

Before describing the proposed method, we quickly discuss Double ML and Neyman orthogonality in the next section which will be helpful in building the theoretical framework for our method.

### 2.1 Double Machine Learning (Double ML)

Given a fixed set of policy variables  $D$  and control variables  $X$ , an unbiased estimator of the parameter  $\theta_0$  for the partial regression model in Equation (2) can be obtained via the orthogonalization approach as in [7].

$$\begin{aligned} Y &= D\theta_0 + g_0(X) + U, \quad \mathbb{E}[U|X, D] = 0 \\ D &= m_0(X) + V, \quad \mathbb{E}[V|X] = 0, \end{aligned} \tag{2}$$

where  $Y$  is the outcome variable,  $U, V$  are disturbances and  $g_0, m_0 : \mathbb{R}^d \rightarrow \mathbb{R}$  are non-linear functions. Orthogonalization is obtained via the use of "Neyman Orthogonality Condition" which we describe below.

**Neyman Orthogonality Condition:** the traditional estimator of  $\theta_0$  in Equation (2) can be simply obtained by finding the zero of the empirical average of a score function  $\psi$  such that  $\psi(W; \theta, g) = D^\top(Y - D\theta - g(X))$ . However, the estimation of  $\theta_0$  is sensitive to the bias in the estimation of the function  $g$ . Neyman [24] proposed an orthogonalization approach to get the estimate for  $\theta$  which is more robust to the bias in the estimation of nuisance parameter ( $m_0, g_0$ ). For a moment, if we assume that the true nuisance parameter is  $\eta_0$  (which represents  $m_0$  and  $g_0$  in Equation (2)) then the orthogonalized "score" function  $\psi$  should satisfy the property that the Gateaux derivative operator with respect to  $\eta$  vanishes when evaluated at the true parameter values:

$$\partial_\eta \mathbb{E}\psi(W; \theta_0, \eta_0)[\eta - \eta_0] = 0 \tag{3}$$

Orthogonalized or Double/Debiased ML estimator  $\check{\theta}_0$  solves

$$\frac{1}{n} \sum_{i=1}^n \psi(W; \check{\theta}_0, \hat{\eta}_0) = 0,$$

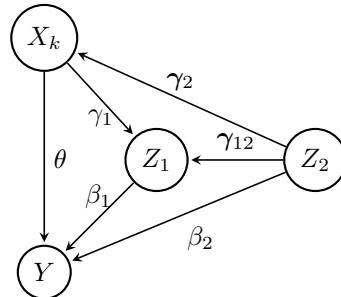


Figure 2: A generic example of identification of a causal effect  $\theta$  in the presence of causal and anti-causal interactions between the causal predictor and other putative parents, and possibly arbitrary non-linear and non-invertible assignments for all nodes except  $Y$  (see Proposition 2). We have  $X_{-k} = Z_1 \cup Z_2$ .

where  $\hat{\eta}_0$  is the estimator of  $\eta_0$  and  $\psi$  satisfies condition in Equation (3). For the partially linear model discussed in Equation (2), the score function  $\psi$  is,

$$\psi(W; \theta, \eta) = (Y - D\theta - g(X))(D - m(X)) \quad (4)$$

with  $\eta = (m, g)$ .

**From Double ML to Causal Discovery:** The distinction between policy variables and confounding variables is not always known in advance, which motivates us to consider the more general setting of causal discovery. To this end, we consider a set of variables  $X = \{X_1, X_2, \dots, X_d\}$  which includes direct causal parents of the outcome variable  $Y$  as well as other variables. We also reiterate our assumption that the relationship between outcome variable and direct causal parents of the outcome variable is linear. The relationship among other variables can be nonlinear. We now provide a general approach to scanning putative direct causes scaling “linearly” with their number, based on the application of a statistical test and a parameter estimation based on Debiased machine learning method. We describe first the algorithm and then provide theoretical support for its performance.

## 2.2 Informal Search Algorithm Description

We provide pseudo-code for our proposed method in Algorithm 1. Intuitively, the idea is to do a one-vs-rest split for each variable in turn, and try to estimate the link between that particular variable and the outcome variable using Debiased approach. To do so, we decompose eq. 1 to single out a variable  $D = X_k$  as policy variable and take the remaining variables  $Z = X_{-k} = X \setminus X_k$  as multidimensional control variables, and run Double ML estimation assuming the partial regression model presented in Sec. 2.1, which now takes the form

$$\begin{aligned} Y &= D\theta_k + g_k(Z) + U, \quad \mathbb{E}[U|X, D] = 0 \\ D &= m_k(Z) + V, \quad \mathbb{E}[V|X] = 0. \end{aligned} \quad (5)$$

The step-wise description of our estimation algorithm goes as follows:

- (a) Select one of the variables  $X_i$  to estimate its (hypothetical) linear causal effect  $\theta$  on  $Y$ .
- (b) Set all of the other variables  $X_{-i}$  as the set of possible confounders.
- (c) Use the Double ML approach to estimate the parameter  $\theta$  i.e. the causal effect of  $X_i$  on  $Y$ .
- (d) If the variable  $X_i$  is not a causal parent, the distribution of the conditional covariance  $\chi_i$  (Proposition 3) is a Gaussian centered around zero. We use a simple normality test for  $\chi_i$  to select or discard  $X_i$  as one of the direct causal parents of  $Y$ .

We iteratively repeat the procedure on each of the variables until completion. Pseudo-code for the entire procedure is given below in Algorithm 1.

Note that only for (5) this corresponds to the Double ML approach by estimating  $\theta_k$ . For all other cases (for instance  $D$  not corresponding to direct causes) it is a none trivial result to show that one obtains a non-zero parameter.

**Remarks on Algorithm 1:**  $X_i^{[k]}$  is a vector which corresponds to the samples chosen in the  $k^{\text{th}}$  subsampling procedure,  $X_{\setminus i}^{[k]} = (X_1^{[k]}, \dots, X_{i-1}^{[k]}, X_{i+1}^{[k]}, \dots, X_d^{[k]})$  for any  $i \in [d]$ . In general the subscript  $i$  represents the estimation for the  $i^{\text{th}}$  variable and super-script  $k$  represents the  $k^{\text{th}}$  subsampling procedure.  $K$  represents the set obtained after sample splitting.  $m_i^{[\setminus k]}$  are (possibly nonlinear) parametric functions fitted using  $(1^{\text{st}}, \dots, k-1^{\text{th}}, k+1^{\text{th}}, \dots, K^{\text{th}})$  subsamples.

## 2.3 Orthogonal Scores

Now we describe the execution of our algorithm for a simple graph with 3 nodes. Let us consider the following linear structural equation model as an example of our general formulation:

$$Y := \theta_1 X_1 + \theta_2 X_2 + \varepsilon_3, \quad X_2 := a_{12} X_1 + \varepsilon_2, \quad \text{and } X_1 := \varepsilon_1. \quad (6)$$

**Example 1.** Let us consider the system whose structural equation model is given in (6). If  $\varepsilon_1, \varepsilon_2$  and  $\varepsilon_3$  are independent uncorrelated noise terms with zero mean, then Algorithm 1 will recover the coefficients  $\theta_1$  and  $\theta_2$ .

A detailed proof is given in the Appendix. While the estimation of the parameter  $\theta_1$  is in line with the assumed partial regression model of eq. (5), the estimation of  $\theta_2$  does not follow the same. However, it can be seen from the proof that  $\theta_2$  can also be estimated from the orthogonal score (Equation (4)).

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**Algorithm 1** Efficient Causal Structure Search

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1: Input: response  $Y \in \mathbb{R}^N$ , covariates
    $\mathbb{X} \in \mathbb{R}^{N \times d}$ , significance level  $\alpha$ , number
   of partitions  $K$ .
2: Split  $N$  observations into K-fold random
   partitions,  $I_k$  for  $k = 1, 2, \dots, K$ , each
   having  $n = N/K$  observations.
3: for  $i = 1, \dots, d$  do
4:   for Subsample  $k \in [K]$  do
5:      $D_k \leftarrow X_i^{[k]}$  and  $Z_k \leftarrow X_{\setminus i}^{[k]}$ 
6:     Fit  $m_i^{[\setminus k]}(Z_{\setminus k})$  to  $D_{\setminus k}$ 
7:     Fit  $g_i^{[\setminus k]}(Z_{\setminus k})$  to  $Y^{[\setminus k]}$ 
8:      $\hat{V}_i^{[k]} \leftarrow D_k - m_i^{[\setminus k]}(Z_k)$ 
9:      $\check{\theta}_i^{[k]} \leftarrow (\frac{1}{n} \sum_{j \in I_k} \hat{V}_{ij}^{[k]} D_{kj})^{-1}$ 
10:     $\hat{\chi}_i^{[k]} \leftarrow \frac{1}{n} \sum_{j \in I_k} (-Y_j^{[k]} m_{ij}^{[\setminus k]}(Z_{kj})$ 
       $- D_{kj} g_{ij}^{[\setminus k]}(Z_{kj}) + m_{ij}^{[\setminus k]}(Z_{kj})$ 
       $g_{ij}^{[\setminus k]}(Z_{kj}) + \hat{V}_{ij}^{[k]} D_{kj})$ 
11:     $(\hat{\sigma}_i^{[k]})^2 \leftarrow \frac{1}{n} \sum_{j \in I_k} (-Y_j^{[k]} m_{ij}^{[\setminus k]}(Z_{kj})$ 
       $- D_{kj} g_{ij}^{[\setminus k]}(Z_{kj}) + m_{ij}^{[\setminus k]}(Z_{kj}) g_{ij}^{[\setminus k]}(Z_{kj})$ 
       $+ \hat{V}_{ij}^{[k]} D_{kj} - \hat{\chi}_i^{[k]})^2$ 
12:  end for
13:   $\hat{\theta}_i \leftarrow \frac{1}{K} \sum_{k \in K} \check{\theta}_i^{[k]}$ 
14:   $\hat{\chi}_i \leftarrow \frac{1}{K} \sum_{k \in K} \hat{\chi}_i^{[k]}$ 
15:   $\hat{\sigma}_i^2 \leftarrow \frac{1}{K} \sum_{k \in K} (\hat{\sigma}_i^{[k]})^2$ 
16: end for
17: DecVec:=[]
18: for  $i \in [d]$  do
19:   Gaussian normality test for  $\hat{\chi}_i \approx N\left(0, \frac{\hat{\sigma}_i^2}{N}\right)$ 
   with  $\alpha$  significance level.
20:   if rejected then
21:     DecVec[i] = 1
22:   else
23:     DecVec[i] = 0
24:   end if
25: end for
26: Return DecVec

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We now show that this result holds for a more general graph structure given in figure 2, allowing also for non-linear interactions among features.

**Proposition 2.** Assume the partially linear Gaussian model of Fig. 2, denote  $X_{-k} = [Z_1^\top, Z_2^\top]^\top$  the control variables,  $\gamma = (\gamma_1, \gamma_1, \gamma_{12})$  the parameter vector of the (possibly non-linear) assignments between putative parents of  $Y$ , and  $\beta = (\beta_1, \beta_2)$ , the vector of causal coefficients for encoding linear effects of  $X_{-k}$  on outcome  $Y$ . Then, independently of the  $\gamma$  parameters and of the functional form of the associated assignments between parents of  $Y$ , the score

$$\psi(W; \theta, \beta) = (Y - X_k \theta - X_{-k}^\top \beta)(X_k - r_{XX_{-k}} X_{-k}), \quad (7)$$

with  $r_{XX_{-k}} = \mathbb{E}[X_k X_{-k}]^\top \mathbb{E}[X_{-k} X_{-k}]^{-1}$ , follows the Neyman orthogonality condition for the estimation of  $\theta$  with nuisance parameters  $\eta = (\beta, \gamma)$  which reads

$$\mathbb{E} \left[ (Y - X_k \theta - X_{-k}^\top \beta)(X_k - r_{XX_{-k}} X_{-k}) \right] = 0.$$

Comparing the score in Equation (7) with the score in Equation (4), there are two takeaways from Proposition 2: (i) the orthogonality condition remains invariant irrespective of the causal direction between  $X_k$  and  $Z$ , and (ii) the second term in Equation (15) replaces function  $m$  by the (unbiased) linear regression estimator for modelling all the relations; given that the relation between  $Z$  and  $Y$  is linear, even if relationships between  $Z$  and  $X_k$  are non-linear. Combining with the Double ML theoretical results [7], this suggests that regularized predictors based on Lasso or ridge regression are tools of choice for fitting functions  $(m, g)$ .

## 2.4 Statistical Test

Now that we have illustrated and justified the fitting procedure of our algorithm, we provide a theoretically justified statistical decision criterion for the direct causes after the model has been fitted. Consider  $(Y, X)$ ,  $Y \in \mathbb{R}$ ,  $X \in \mathbb{R}^d$ , satisfying

$$Y = \langle \theta, X \rangle + U, \quad (8)$$

$$E(Y^2) < \infty, E(U^2) < \infty, E(U) = 0, E(U | X_j) = 0, \forall j, \text{ and } E(\|X\|_2^2) < \infty, \quad (9)$$

$$E \left[ (X_j - E(X_j | X_{-j}))^2 \right] \neq 0, \quad \text{for all } j \in \{1, \dots, p\}. \quad (10)$$

where  $U$  is noise variable and  $X_{-j}$  represents all the variables except  $X_j$ . The assumptions made with the above formulation are standard in the orthogonal machine learning literature [31, 35, 8]. Let us define the quantity  $\chi_j = E[(Y - E(Y | X_{-j}))(X_j - E(X_j | X_{-j}))]$  for  $j \in \{1, \dots, d\}$ , which is the expected conditional covariance of  $X_j$  given  $X_{-j}$ .

**Proposition 3.** *Let  $PA_Y = \{j \in \{1, \dots, p\} : \theta_j \neq 0\}$ . For each  $j \in \{1, \dots, p\}$  let  $X_{-j}$  be the vector equals to  $X$  but excluding coordinate  $j$  and define  $\theta_{-j}$  similarly. Define for  $j \in \{1, \dots, p\}$*

$$\chi_j = \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j}))(X_j - \mathbb{E}(X_j | X_{-j}))],$$

which also has the double robustness property ([8, 31]) then under the conditions given in Equations (8)-(10)

- a) If  $j \in PA_Y$  then  $\chi_j = \theta_j \mathbb{E}[(X_j - \mathbb{E}(X_j | X_{-j}))^2]$ .
- b) If  $j \notin PA_Y$  then  $\chi_j = 0$ .
- c) We also have (with notations of Prop. 2)  $\chi_j = \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j}))(X_j - r_{XX_{-k}} X_{-j})]$ .

There are two main implications of the results provided in Proposition 3. (i)  $\chi_j$  is non-zero only for direct causal parents of the outcome variable and  $\chi_j$  has double robustness property as shown in [31, 35, 8], hence one can use regularized methods like ridge regression or Lasso to estimate the function  $m$ . Afterwards, one can perform statistical tests on top of it to decide between zero or non-zero test. (ii) In line with the above orthogonal score results, we see that this quantity can be estimated using linear (unbiased) regression to fit the function  $m$ , although interactions between features may be non-linear. Next, we discuss the variance of our estimator so that later a statistical test can be used to differentiate between zero and non-zero test.

**Variance of Empirical Estimates of  $\chi_j$ :** Suppose we have  $n$  i.i.d. observations  $\mathcal{D}_n = \{(X_i, Y_i), i = 1, \dots, n\}$ . Randomly split the data in two halves, say  $\mathcal{D}_{n1}$  and  $\mathcal{D}_{n2}$ . Take  $j \in \{1, \dots, p\}$ . For  $k = 1$  let  $\bar{k} = 2$ , for  $k = 2$  let  $\bar{k} = 1$ . For  $k = 1, 2$ , compute estimates of  $\widehat{\mathbb{E}}^{\bar{k}}(Y | X_{-j})$  and  $\widehat{\mathbb{E}}^{\bar{k}}(X_j | X_{-j})$  using the data in sample  $\bar{k}$ . Computing  $\widehat{\mathbb{E}}^{\bar{k}}(Y | X_{-j})$  and  $\widehat{\mathbb{E}}^{\bar{k}}(X_j | X_{-j})$  can be considered as regularized regression problems. We use Lasso as the estimator for conditional expectation (eq. (17)) in the experiments. Now, we compute the empirical estimates of  $\chi_j$ . Theorem 1 of [35] provides conditions under which (see also [8]), when the estimators

$$\widehat{\mathbb{E}}^{\bar{k}}(Y | X_{-j}) \quad \text{and} \quad \widehat{\mathbb{E}}^{\bar{k}}(X_j | X_{-j}) \tag{11}$$

are Lasso-type regularized linear regressions, it holds that asymptotically  $\widehat{\chi}_j \approx N\left(\chi_j, \frac{\widehat{\sigma}_j^2}{n}\right)$ .

In this case, the test that rejects  $\chi_j = 0$  when  $|\widehat{\chi}_j| \geq 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}$  will have approximately 5% level. The probability of rejecting the null when it is false is

$$P\left(|\widehat{\chi}_j| \geq 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}\right) \geq P\left(|\widehat{\chi}_j - \chi_j| \leq |\chi_j| - 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}\right) \rightarrow 1.$$

Expressions for  $\widehat{\chi}_j$  and  $\widehat{\sigma}_j$  are provided in the Appendix A.2. In order to account for multiple testing, we use Bonferroni correction.

**Comments about Estimator:** In this paper, we use Lasso for the nuisance parameter estimation as the variance of the conditional covariance is known [35]. One can also use other estimators instead, assuming one obtains a good enough estimate of the nuisance parameter (upto  $N^{-1/4}$ -neighbourhood [7]) with the right variance term, which is beyond the scope of this paper.

### 3 Experiments

In this section, we perform extensive empirical evaluation for our method.

#### 3.1 Experimental Setup

For every combination of number of nodes (#nodes), sparsity ( $p_s$ ), noise level ( $\sigma^2$ ), number of observations ( $z$ ), and non-linear probability ( $p_n$ ) (see Table C.1), 100 examples (DAGs) are generated and stored as csv files (altogether 72.000 DAGs are simulated, comprising a dataset of overall  $>10\text{GB}$ ).

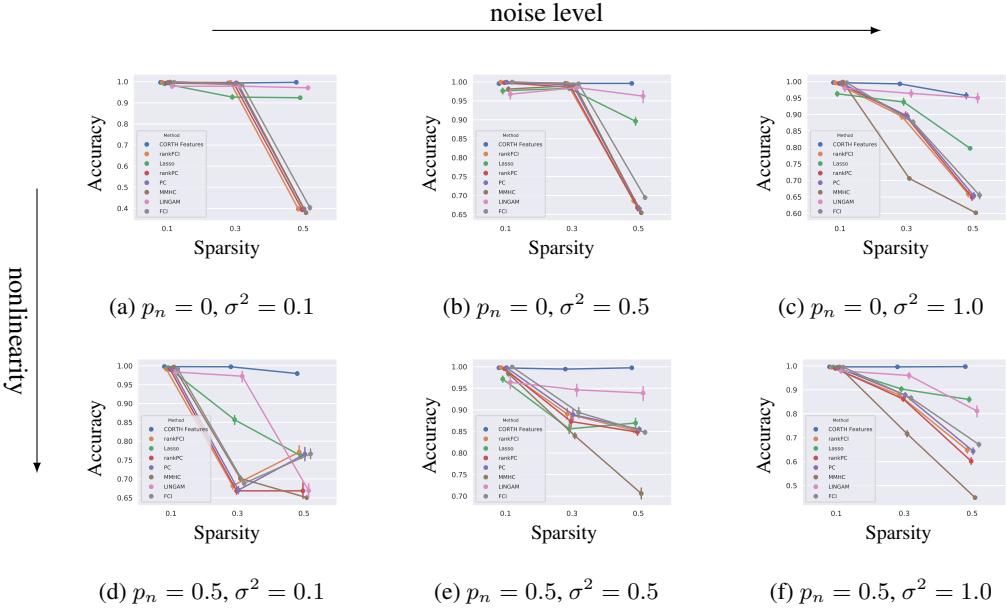


Figure 3: Overall performance for a single random DAG with 100 simulations for each setting, having 20 nodes and 500 observations.

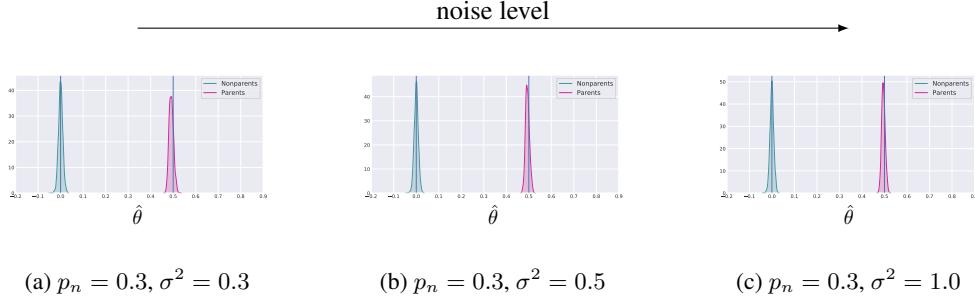


Figure 4: Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations of the graph with 20 nodes, 20000 observations and 0.3 as sparsity. The vertical lines indicate the ground truth values for the linear coefficients corresponding to causal parents. Plots for all other estimation settings are in the Appendix F.

For each DAG,  $z$  number of samples are generated. We provide more details about the parameters (#nodes,  $p_s$ ,  $p_n$  and  $z$ ) and data generation process in Appendix C. For future benchmarking, the generated files with the code will be made available later.

The baselines we compare our method against are: LINGAM [33], order-independent PC [11], rankPC, MMHC [38], GES [9], rankGES, ARGES (adaptively restricted GES [23]), rankARGES, FCI+ [10] and Lasso which are suitable for observational data. CompareCausalNetworks Package<sup>1</sup> is used to run the baselines methods. We use 10-fold cross validation to choose the parameters of all approaches. Recall, Fall-out, Critical Success Index, Accuracy, F1 Score, and Matthews correlation coefficient [22] are considered as metric for the evaluation. These metrics are described in Appendix C.

**Regression Technique and Hyper-parameters:** We use Lasso as the estimator of conditional expectation for our method because the variance bound for  $\chi_j$  with Lasso type estimator of conditional expectation (Equation (17)) is provided in Equation (16). Further, using more splits than 2 splits in the experiment increases the performance of parameter estimation. Plots for parameter estimation are provided in Appendix F.

<sup>1</sup> <https://cran.r-project.org/web/packages/CompareCausalNetworks/index.html>

### 3.2 Results

Results aggregated by number of nodes over all simulations (corresponding to 18000 simulations per entry in the table), are illustrated in Table 1. It shows our method performs better than the competing baselines in terms of accuracy and F1 score, especially for a higher number of nodes. To provide a visual comparison, we plot the accuracy of all the methods w.r.t. the sparsity parameter ( $p_s$ ) in Figure 3 for different values of  $p_n$  and  $\sigma^2$  on 1800 samples.

It can be observed that the accuracies of the competing baselines significantly drop with increasing noise level, while our method is more robust to it. More plots are given in Appendix G and Appendix H for several other combinations of varying parameters in the simulation. We also extensively compare all the metrics (Recall, Fall-out, Critical Success Index, Accuracy, F1 Score and Matthews correlation coefficient) for all the methods in Appendix E. According to these metrics, our approach performs better than baselines in most of the cases regardless of the set of parameters used for generating data. Our method shows in particular stability in performance w.r.t. the number of nodes (Table E.1), partially non-linear relationships (Table E.3), sparsity (Table E.2), number of observations (Table E.4), and noise level (Table E.5). We also show the plot of parameter estimation for direct causal parents vs. non-causal parents in Figure 4. In the plots and tables, we denote our approach as **CORTH Features**.

### 3.3 Scaling Causal Inference to Large Graphs

Figure 5 shows the runtime of the method in seconds as a function of the graph’s size. Notice that the runtime of our algorithm in the log-log plot is roughly linear. Since we used 5000 observations, any additional overhead is coming from cross-validation.

### 3.4 Real-World Data

We also apply our algorithm on a real world data sets and evaluate the performance of our algorithm on a recent COVID-19 Dataset [12] where the task is to predict COVID-19 cases (confirmed using RT-PCR) amongst suspected ones. For an existing and extensive analysis of the dataset with predictive methods we refer to [32]. We apply our algorithm to discover the features which directly cause the diagnosed infection. We found that the following were the most common causes across different runs of our approach: Patient age quantile, Arterial Lactic Acid, Promyelocytes, and Base excess venous blood gas analysis. Lacking medical ground truth, we report these not as a corroboration of our approach, but rather as a potential contribution to causal discovery in this challenging problem. It is encouraging that some of these variables are consistent with other studies [32]. Details on data preprocessing and more results are available in Appendix D.

Method	Number of Nodes					
	10		20		50	
ACC	F1	ACC	F1	ACC	F1	
GES	0.85	0.78	0.74	0.53	0.70	0.32
rankGES	0.85	0.75	0.74	0.51	0.70	0.32
ARGES	0.80	0.58	0.75	0.52	0.71	0.22
rankARGES	0.79	0.57	0.75	0.51	0.71	0.22
FCI+	0.87	0.81	0.83	0.70	0.77	0.49
LINGAM	<b>0.95</b>	0.89	0.89	0.78	0.75	0.39
PC	0.86	0.79	0.82	0.66	0.76	0.46
rankPC	0.85	0.77	0.81	0.64	0.75	0.43
MMHC	0.84	0.74	0.77	0.51	0.73	0.28
Lasso	0.91	0.90	0.90	0.87	0.77	0.63
<b>CORTH Features</b>	<b>0.95</b>	<b>0.93</b>	<b>0.95</b>	<b>0.91</b>	<b>0.80</b>	<b>0.66</b>

Table 1: Performance across all the settings for different number of nodes (10,20 and 50). Each entry in the table is averaged over 18000 simulations.

Table 1: Performance across all the settings for different number of nodes (10,20 and 50). Each entry in the table is averaged over 18000 simulations.

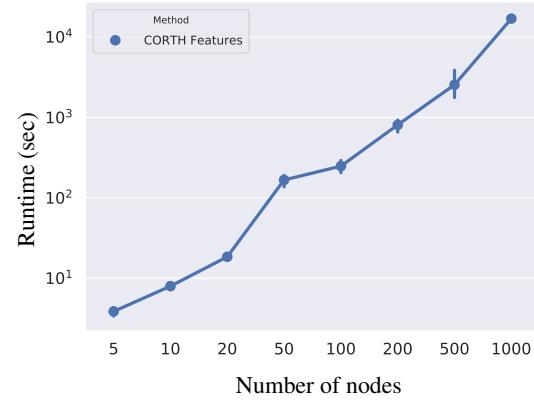


Figure 5: Runtime as a function of the number of variables for 10 simulations per number of nodes. In these simulations sparsity, number of observations, nonlinaer prob., and noise level are set to 0.3, 5000, 0, and 1 respectively.

## 4 Discussion

A recent empirical evaluation of different causal discovery methods highlighted the desirability of more efficient algorithms [16]. In the present work, we provide identifiability results for the set of direct causal parents including the case of partially nonlinear models, as well as a highly efficient algorithm that scales linearly in the number of variables and exhibits state-of-the-art performance across extensive experiments. Our approach builds on the Double ML method of [7], however, we stress that it goes beyond this since the distinction between policy variables and confounding variables is not always known in advance. Whilst not amounting to full causal graph discovery, identification of causal parents is of major interest in real-world applications, e.g., when assaying the causal influence of genes on the phenotype. A natural direction worth exploring is to extend this approach for discovering direct causal parents in the case when non-linear relation exists between output variable and its direct causal parents.

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# Appendix

## A Causal Discovery via Orthogonalization

*Proof of Example 1.* Let us start from the easier case first. Let us first try to estimate the coefficient of interaction between  $X_2$  and  $Y$  but it is also very clear that the estimation of  $\theta_2$  will be unbiased as the given setting precisely match with the double machine learning setting. However, we will see in this example that given the population,  $\theta_1$  can be approximated as well. Let us write down the structural equation model first:

$$\begin{aligned} Y &:= \theta_1 X_1 + \theta_2 X_2 + \varepsilon_3 \\ X_2 &:= a_{12} X_1 + \varepsilon_2 \\ X_1 &:= \varepsilon_1 \end{aligned} \tag{12}$$

From the set of equations we have:

$$X_1 = a_{12}^{-1} X_2 - a_{12}^{-1} \varepsilon_2$$

Let also denote  $\mathbb{E}[\varepsilon_1^2] = \sigma_1^2$  and  $\mathbb{E}[\varepsilon_2^2] = \sigma_2^2$ . Hence,  $\mathbb{E}[X_1^2] = \sigma_1^2$ ,  $\mathbb{E}[X_1 X_2] = a_{12} \sigma_1^2$  and  $\mathbb{E}[X_2^2] = a_{12} \mathbb{E}[X_1 X_2] + \mathbb{E}[\varepsilon_2 X_2] = a_{12}^2 \sigma_1^2 + \sigma_2^2$ . Let us first try to find the regression co-efficient of fitting  $X_2$  on  $Y$ .

$$Y = \hat{\theta}_2 X_2 + \eta_1$$

Hence,  $\hat{\theta}_2 = \frac{\mathbb{E}[X_2 Y]}{\mathbb{E}[X_2^2]}$  if  $\eta$  is independent of  $X_2$ .

$$\hat{\theta}_2 = \frac{\mathbb{E}[X_2 Y]}{\mathbb{E}[X_2^2]} = \frac{\mathbb{E}[X_2(\theta_1 X_1 + \theta_2 X_2 + \varepsilon_3)]}{\mathbb{E}[X_2^2]} = \theta_2 + \theta_1 a_{12} \frac{\sigma_1^2}{\sigma_2^2 + a_{12}^2 \sigma_1^2} \tag{13}$$

Similarly, if we fit  $X_2$  on  $X_1$  then

$$X_1 = \hat{a}_{12}^{-1} X_2 + \eta_2$$

then  $\hat{a}_{12}^{-1} = \frac{\mathbb{E}[X_1 X_2]}{\mathbb{E}[X_2^2]}$ . However  $\mathbb{E}[X_1 X_2]$  can also be written as following:

$$\mathbb{E}[X_1 X_2] = a_{12}^{-1} \mathbb{E}[X_2^2] - a_{12}^{-1} \mathbb{E}[\varepsilon_2 X_2]$$

Hence,

$$\hat{a}_{12}^{-1} = a_{12}^{-1} \left( 1 - \frac{\sigma_2^2}{\sigma_2^2 + a_{12}^2 \sigma_1^2} \right) = a_{12}^{-1} \left( \frac{a_{12}^2 \sigma_1^2}{\sigma_2^2 + a_{12}^2 \sigma_1^2} \right)$$

Residual  $\hat{V} = X_1 - \hat{a}_{12}^{-1} X_2$ . Hence we can have

$$\mathbb{E}(\hat{V} X_1) = \mathbb{E}[X_1^2] - \hat{a}_{12}^{-1} \mathbb{E}[X_1 X_2] = \mathbb{E}[\varepsilon_1^2] - \hat{a}_{12}^{-1} a_{12} \mathbb{E}[\varepsilon_1^2] = \frac{\sigma_1^2 \sigma_2^2}{\sigma_2^2 + a_{12}^2 \sigma_1^2}$$

We now calculate,

$$\begin{aligned} \mathbb{E}[\hat{V}(Y - \hat{\theta}_2 X_2)] &= \mathbb{E}[(X_1 - \hat{a}_{12}^{-1} X_2)(Y - \hat{\theta}_2 X_2)] \\ &= \mathbb{E}[(X_1 - \hat{a}_{12}^{-1} X_2)((\theta_2 - \hat{\theta}_2)X_2 + \theta_1 X_1 + \varepsilon_3)] \\ &= (\theta_2 - \hat{\theta}_2)a_{12} \sigma_1^2 + \theta_1 \sigma_1^2 - \hat{a}_{12}^{-1}(\theta_2 - \hat{\theta}_2)(\sigma_2^2 + a_{12}^2 \sigma_1^2) - \hat{a}_{12}^{-1} \theta_1 a_{12} \sigma_1^2 \\ &= \frac{\theta_1 \sigma_1^2 \sigma_2^2}{\sigma_2^2 + a_{12}^2 \sigma_1^2} \end{aligned}$$

Last equation was written after step of minor calculation. Since the estimator is

$$\hat{\theta}_1 = [\mathbb{E}(\hat{V} X_1)]^{-1} \mathbb{E}[\hat{V}(Y - \hat{\theta}_2 X_2)] = \theta_1$$

□

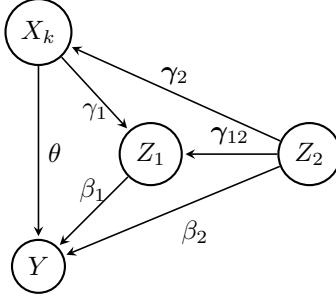


Figure A.1: Generic example of identification of a causal effect  $\theta$  in the presence causal and anticausal interactions between the causal predictor and other putative parents, and possibly arbitrary nonlinear and non-invertible assignments for all nodes except  $Y$  (see Proposition 2) We have  $X_{-k} = Z_1 \cup Z_2$ .

### A.1 Influence of the interactions between parents

In this section, we use a generic example shown in Figure 2 which we show again in Figure A.1 to illustrate the role of interactions between the covariates on the proposed causal discovery algorithm.

The estimator discussed can simply be derived from the Neyman orthogonality condition. We now provide the below the proof for Proposition 2. For the sake of completeness, we also rewrite the statement of the proposition again.

**Proposition (Restatement of Proposition 2).** *Assume the partially linear Gaussian model of Fig. A.1, denote  $X_{-k} = [Z_1^\top, Z_2^\top]^\top$  the control variables,  $\boldsymbol{\gamma} = (\gamma_1, \gamma_1, \gamma_{12})$  the parameter vector of the (possibly non-linear) assignments between putative parents of  $Y$ , and  $\boldsymbol{\beta} = (\beta_1, \beta_2)$  the vector of causal coefficients for encoding linear effects of  $X_{-k}$  on outcome  $Y$ . Then, independently from the  $\boldsymbol{\gamma}$  parameters and of the functional form of the associated assignments between parents of  $Y$ , the score*

$$\psi(W; \theta, \boldsymbol{\beta}) = (Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta})(X_k - r_{XX_{-k}} X_{-k}), \quad (14)$$

with  $r_{XX_{-k}} = \mathbb{E}[X_k X_{-k}^\top] \mathbb{E}[X_{-k} X_{-k}^\top]^{-1}$ , follows the Neyman orthogonality condition for the estimation of  $\theta$  with nuisance parameters  $\boldsymbol{\eta} = (\beta, \boldsymbol{\gamma})$  which reads

$$\mathbb{E} \left[ (Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta})(X_k - \mathbb{E}[X_k X_{-k}^\top] \mathbb{E}[X_{-k} X_{-k}^\top]^{-1} X_{-k}) \right] = 0. \quad (15)$$

*Proof of Proposition 2.* Using the Markov factorization

$$P(W; \theta, \boldsymbol{\eta}) = P(Y|X_{-k}, X_k; \theta, \boldsymbol{\beta})P(X_{-k}, X_k; \boldsymbol{\gamma})$$

due to linearity and gaussianity of the assignement of  $Y$ , we obtain a negative log likelihood of the form (up to additive constants)

$$\ell(W; \theta, \boldsymbol{\eta}) = \frac{1}{2\sigma_Y^2} (Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta})(Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta}) + f(X_k, X_{-k}; \boldsymbol{\gamma})$$

where  $f$  stands for the negative log likelihood of the second factor. Following [7][eq. (2.7)], this leads to the Neyman orthogonal score

$$\begin{aligned} \psi(W; \theta, \boldsymbol{\eta}) &= \partial_\theta \ell(W; (\theta, \boldsymbol{\eta})) - \boldsymbol{\mu} \partial_\eta \ell(W; (\theta, \boldsymbol{\eta})) = -\frac{1}{\sigma_Y^2} (Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta}) X_k \\ &\quad - \boldsymbol{\mu} \left( -\frac{1}{\sigma_Y^2} (Y - X_k \theta - X_{-k}^\top \boldsymbol{\beta}) X_{-k} + \partial_\gamma f(X_k, X_{-k}; \boldsymbol{\gamma}) \right) \end{aligned}$$

Following eq. (2.8) of the same paper, we derive the expression of  $\boldsymbol{\mu}$  as

$$\boldsymbol{\mu} = J_{\theta, \eta} J_{\eta, \eta}^{-1}$$

with

$$J_{\eta, \eta} = \partial_{\eta^\top} \mathbb{E} [\partial_\eta \ell(W, \theta, \boldsymbol{\eta})] = \begin{bmatrix} \sigma_Y^{-2} \mathbb{E} [X_{-k}^\top X_{-k}] & 0 \\ 0 & \partial_\gamma \mathbb{E} [\partial_\gamma f(X_k, X_{-k}; \boldsymbol{\gamma})] \end{bmatrix},$$

and

$$J_{\theta,\eta} = \partial_{\eta^\top} \mathbb{E} [\partial_\theta \ell(W, \theta, \eta)] = \sigma_Y^{-2} \begin{bmatrix} \mathbb{E}[X_k^\top X_{-k}] & 0 \end{bmatrix},$$

resulting in

$$\boldsymbol{\mu} = \mathbb{E}[X_k^\top X_{-k}] \mathbb{E}[X_{-k}^\top X_{-k}]^{-1}$$

Reintroducing  $\boldsymbol{\mu}$  in the expression of  $\psi$  leads to the result.  $\square$

**Proposition (Restatement of Proposition 3).** Let  $PA_Y = \{j \in \{1, \dots, p\} : \theta_j \neq 0\}$ . For each  $j \in \{1, \dots, p\}$  let  $X_{-j}$  be the vector equals to  $X$  but excluding coordinate  $j$  and define  $\theta_{-j}$  similarly. Define for  $j \in \{1, \dots, p\}$

$$\chi_j = \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j})) (X_j - \mathbb{E}(X_j | X_{-j}))],$$

then under the conditions given in Equations (8)-(10)

- a) If  $j \in PA_Y$  then  $\chi_j = \theta_j \mathbb{E}[(X_j - \mathbb{E}(X_j | X_{-j}))^2]$ .
- b) If  $j \notin PA_Y$  then  $\chi_j = 0$ .
- c) We also have (with notations of Prop. 2)  $\chi_j = \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j})) (X_j - r_{XX_{-k}} X_{-j})]$ .

*Proof.* Take  $j \in PA_Y$ . Then, from (8)

$$\begin{aligned} \mathbb{E}(Y | X_{-j}) &= \mathbb{E}(\langle \theta, X \rangle | X_{-j}) + \mathbb{E}(U | X_{-j}) \\ &= \mathbb{E}(\langle \theta_{-j}, X_{-j} \rangle | X_{-j}) + \theta_j \mathbb{E}(X_j | X_{-j}) \\ &= \langle \theta_{-j}, X_{-j} \rangle + \theta_j \mathbb{E}(X_j | X_{-j}) \\ &= \langle \theta, X \rangle - \theta_j X_j + \theta_j \mathbb{E}(X_j | X_{-j}) \\ &= Y - U - \theta_j (X_j - \mathbb{E}(X_j | X_{-j})). \end{aligned}$$

Thus

$$\begin{aligned} \chi_j &= \mathbb{E}[(U + \theta_j (X_j - \mathbb{E}(X_j | X_{-j}))) (X_j - \mathbb{E}(X_j | X_{-j}))] \\ &= \mathbb{E}[U(X_j - \mathbb{E}(X_j | X_{-j}))] + \theta_j \mathbb{E}[(X_j - \mathbb{E}(X_j | X_{-j}))^2] \\ &= \theta_j \mathbb{E}[(X_j - \mathbb{E}(X_j | X_{-j}))^2]. \end{aligned}$$

Now take  $j \notin PA_Y$ . Then

$$\begin{aligned} \mathbb{E}(Y | X_{-j}) &= \mathbb{E}(\langle \theta, X \rangle | X_{-j}) + \mathbb{E}(U | X_{-j}) \\ &= \langle \theta, X \rangle. \end{aligned}$$

Thus,  $\chi_j = \mathbb{E}[U(X_j - \mathbb{E}(X_j | X_{-j}))] = 0$ .

For c), we rewrite

$$\begin{aligned} \chi_j &= \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j})) (X_j - r_{XX_{-k}} X_{-j})] \\ &\quad + \mathbb{E}[(Y - \mathbb{E}(Y | X_{-j})) (r_{XX_{-k}} X_{-j} - \mathbb{E}(X_j | X_{-j}))]. \end{aligned}$$

Let  $\mathcal{G}$  the sub-sigma algebra generated by  $X_{-j}$ , under our assumptions,  $\mathbb{E}(Y | X_{-j})$  is the orthogonal projection of  $Y$  on the subspace of  $\mathcal{G}$ -measurable square integrable RV's  $L^2(\Omega, \mathcal{G})$ , so  $Y - \mathbb{E}(Y | X_{-j})$  is orthogonal to any elements of  $L^2(\Omega, \mathcal{G})$ . Noticing that  $(r_{XX_{-k}} X_{-j} - \mathbb{E}(X_j | X_{-j}))$  is an element of  $L^2(\Omega, \mathcal{G})$ , the second right-hand side term of the above equation vanishes and we get the result.  $\square$

## A.2 Variance of Empirical Estimates of $\chi_j$

Suppose we have  $n$  i.i.d. observations  $\mathcal{D}_n = \{(X_i, Y_i), i = 1, \dots, n\}$ . Randomly split the data in two halves, say  $\mathcal{D}_{n1}$  and  $\mathcal{D}_{n2}$ . Take  $j \in \{1, \dots, p\}$ . For  $k = 1$  let  $\bar{k} = 2$ , for  $k = 2$  let  $\bar{k} = 1$ . For  $k = 1, 2$ , compute estimates of  $\widehat{\mathbb{E}}^{\bar{k}}(Y | X_{-j})$  and  $\widehat{\mathbb{E}}^{\bar{k}}(X_j | X_{-j})$  using the data in sample  $\bar{k}$ . Computing  $\widehat{\mathbb{E}}^{\bar{k}}(Y | X_{-j})$  and  $\widehat{\mathbb{E}}^{\bar{k}}(X_j | X_{-j})$  can be considered as regularized regression problems.

We use Lasso as the estimator for conditional expectation (eq. (17)) in the experiments. Now, we compute the empirical estimates of  $\chi_j$ . Let,

$$\widehat{\chi}_j^k = \mathbb{P}_{nk} \left[ -Y\widehat{\mathbb{E}}^k(X_j | X_{-j}) - X_j\widehat{\mathbb{E}}^k(Y | X_{-j}) + \widehat{\mathbb{E}}^k(Y | X_{-j})\widehat{\mathbb{E}}^k(X_j | X_{-j}) + YX_j \right].$$

and

$$\begin{aligned} (\widehat{\sigma}_j^k)^2 &= \mathbb{P}_{nk} \left[ \left( -Y\widehat{\mathbb{E}}^k(X_j | X_{-j}) - X_j\widehat{\mathbb{E}}^k(Y | X_{-j}) \right. \right. \\ &\quad \left. \left. + \widehat{\mathbb{E}}^k(Y | X_{-j})\widehat{\mathbb{E}}^k(X_j | X_{-j}) + YX_j - \widehat{\chi}_j^k \right)^2 \right]. \end{aligned} \quad (16)$$

$\mathbb{P}_{nk}$  here denotes empirical average and  $\widehat{\sigma}_j^k$  denotes empirical variance of  $\chi_j$ . Finally, let

$$\widehat{\chi}_j = \frac{\widehat{\chi}_1 + \widehat{\chi}_2}{2}, \quad \widehat{\sigma}_j^2 = \frac{(\widehat{\sigma}_j^1)^2 + (\widehat{\sigma}_j^2)^2}{2}.$$

Theorem 1 of [?] provides conditions under which (see also [8]), when the estimators

$$\widehat{\mathbb{E}}^k(Y | X_{-j}) \quad \text{and} \quad \widehat{\mathbb{E}}^k(X_j | X_{-j}) \quad (17)$$

are Lasso-type regularized linear regressions, it holds that asymptotically  $\widehat{\chi}_j \approx N\left(\chi_j, \frac{\widehat{\sigma}_j^2}{n}\right)$ .

In this case, the test that rejects  $\chi_j = 0$  when  $|\widehat{\chi}_j| \geq 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}$  will have approximately 5% level. The probability of rejecting the null when it is false is

$$P\left(|\widehat{\chi}_j| \geq 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}\right) \geq P\left(|\widehat{\chi}_j - \chi_j| \leq |\chi_j| - 1.96 \frac{\widehat{\sigma}_j}{\sqrt{n}}\right) \rightarrow 1.$$

## B Examples

The result discussed in Proposition 2 is not directly intuitive. In simple words, there are two takeaways from Proposition 2: (i) the orthogonality condition remains invariant irrespective of the causal direction between  $X_k$  and  $Z$ , and (ii) the second term in equation 15 suggests to use a linear estimator for modeling all the relations, given that the relation between  $Z$  and  $Y$  is linear.

To generate more intuition, we provide a few examples. Let us go back again to the three variable interaction assuming the following structural equation model:

$$\begin{aligned} Y &:= \theta_1 X_1 + \theta_2 X_2 + \varepsilon_3 \\ X_2 &:= f(X_1) + \varepsilon_2 \\ X_1 &:= \varepsilon_1, \end{aligned} \quad (18)$$

where  $f$  is a nonlinear function and  $\varepsilon_1, \varepsilon_2$  and  $\varepsilon_3$  are zero mean Gaussian noises.

- Consider the case when  $f(x) = x^2$ . The goal is to estimate the parameter  $\theta_1$  which we call  $\hat{\theta}_1$ . We follow the standard double ML procedure assuming policy variable  $X_1$  and control  $X_2$ , although the ground truth causal dependency  $X_1 \rightarrow X_2$  in contradiction with such setting (see equation 2). The estimate of  $\theta_2$  following the double ML procedure, which we call  $\hat{\theta}_2 = \frac{\mathbb{E}[X_2 Y]}{\mathbb{E}[X_2^2]} = \theta_2 + \theta_1 \frac{\mathbb{E}[X_1 X_2]}{\mathbb{E}[X_2^2]}$ . Similarly, we want to estimate  $X_1 = \alpha X_2 + \eta$  from which we get,  $\alpha = \frac{\mathbb{E}[X_1 X_2]}{\mathbb{E}[X_2^2]}$ . It is easy to see that  $\mathbb{E}[X_1 X_2] = \mathbb{E}[X_1^3] = 0$ . Hence,  $\alpha = 0$  and it is easy to see  $\hat{\theta}_1 = \theta_1$ .
- Consider now the more general case where  $f$  is any nonlinear function. As in the previously discussed example, the goal is to estimate  $\theta_1$ . We have  $\hat{\theta}_2 = \frac{\mathbb{E}[X_2 Y]}{\mathbb{E}[X_2^2]} = \theta_2 + \theta_1 \frac{\mathbb{E}[X_1 X_2]}{\mathbb{E}[X_2^2]}$ . Similarly,

sparsity ( $p_s$ )	# nodes	nonlinear probability ( $p_n$ )	# observ. z	noise level ( $\sigma^2$ )
0.1	5	0	100	0.01
0.3	10	0.3	500	0.1
0.5	20	0.5	1.000	0.3
	50	1		0.5
				1

Table C.1: Experimental Setup: In the experiments we vary a sparsity parameter, the number of nodes in the graph, the non-linear probability, the number of observations and the noise level and generate 100 graphs for each setting.

$\alpha = \frac{\mathbb{E}[X_1 X_2]}{E[X_2^2]}$ . We substitute these estimates into the orthogonality condition (15):

$$\begin{aligned}
& \mathbb{E} \left[ (Y - X_1 \hat{\theta}_1 - X_2 \hat{\theta}_2)(X_1 - \alpha X_2) \right] = 0. \\
\Rightarrow & \mathbb{E} \left[ \left( Y - X_1 \hat{\theta}_1 - X_2 \hat{\theta}_2 \right) \left( X_1 - \frac{\mathbb{E}[X_1 X_2]}{E[X_2^2]} X_2 \right) \right] = 0. \\
\Rightarrow & \mathbb{E} \left[ \left( X_1(\theta_1 - \hat{\theta}_1) + (a_2 - \hat{\theta}_2)X_2 + \varepsilon_3 \right) \right. \\
& \quad \left. \left( X_1 - \frac{\mathbb{E}[X_1 X_2]}{E[X_2^2]} X_2 \right) \right] = 0. \\
\Rightarrow & \hat{\theta}_1 = \theta_1
\end{aligned}$$

From the above two examples it is clear that even though the internal relations between the variables are nonlinear, all we need is an unbiased linear estimate to estimate the causal parameter in a directed acyclic graph.

## C Data Generation and Evaluation Metric

### C.1 Data Generation

For every combination of number of nodes (#nodes), sparsity ( $p_s$ ), noise level ( $\sigma^2$ ), number of observation ( $z$ ), and non-linear probability ( $p_n$ ) (look at Table C.1), 100 examples (DAGs) are generated and stored as csv files (altogether 72.000 DAGs are simulated, comprising a dataset of overall >10GB). For each DAG,  $z$  number of samples are generated by sampling noise ( $\epsilon$  in Equation (19)) with variance  $\sigma^2$  starting from root of the DAG. For future benchmarking, the generated files will be made available with the code later on.

We generate DAGs (Direct Acyclic Graphs) in multiple steps: i) a random permutation of nodes is chosen as a topological order of a DAG. ii) Based on this order, directed edges are added to this DAG from each node to its followers with a certain probability  $p_s$  (sparsity). iii) For each observation, values are assigned to nodes according to the topological order of the DAG in such a way that each node's value is determined by summing over transformations (linear or nonlinear with a certain nonlinear probability  $p_n$ ) of values of its direct causes with the addition of Gaussian distributed noise. The non-linear transformation used is  $\alpha \tanh(\beta x)$ , with  $\alpha = 0.5$  and  $\beta = 1.5$ . If the set of parents for the node  $X'$  is denoted as  $PA_{X'}$  as before then value assignment for a node  $X'$  is as follow:

$$X' = \varepsilon + \sum_{X \in PA_{X'}} \iota_\ell(p_n) \theta X + (1 - \iota_\ell(p_n)) \alpha \tanh(\beta X), \quad (19)$$

where  $\varepsilon \sim N(0, \sigma^2)$  in which  $\sigma^2$  represents noise level.  $\iota_\ell(X)$  is an indicator functions which decides between linear or non-linear contribution of  $X$  in  $X'$ . We decide the value of  $\iota_\ell(p_n)$  by generating a binary random number which is 1 with probability  $p_n$  and 0 with probability  $1 - p_n$ . The value of  $\theta$  is set to 2 for the small DAGs (number of nodes equal to 5 or 10) and 0.5 for large DAGs (number of nodes equal to 20 or 50) due to the value exploitation that might happen in large graphs.

We vary and investigate the effect of non-linear relationships, the number of nodes, number of observations, effect of sparsity and noise level while simulating the data. We summarize the factors in the data generation in table C.1.

## C.2 Evaluation Metric

Let the total number of true positives, false positives, true negatives ,and false negatives denoted by TP, FP, TN, and FN, we evaluate our method using following metrics:

- Recall (true positive rate):

$$TPR = \frac{TP}{TP + FN}$$

- Fall-out (false positive rate):

$$FPR = \frac{FP}{FP + TN}$$

- Critical Success Index (CSI): also known as Threat Score.

$$CSI = \frac{TP}{TP + FN + FP}$$

- Accuracy:

$$ACC = \frac{TP + TN}{P + N}$$

- F1 Score: harmonic mean of precision and sensitivity.

$$F_1 = \frac{2TP}{2TP + FP + FN}$$

- Matthews correlation coefficient (MCC): a metric for evaluating quality of binary classification introduced in [22].

$$MCC = \frac{TP \times TN - FP \times FN}{\sqrt{(TP + FP)(TP + FN)(TN + FP)(TN + FN)}}$$

In some rare cases, we encountered zero-divided-by-zero and divided-by-zero cases for some of these metrics. In these situations, scores are reported 1 and 0 respectively while Fall-out is reported 0 and 1.

## D Real-World Data Experiment-Covid19

### D.1 Preprocessing

The preprocessing stage for this dataset is the same as [32] except that, for each target variable upsampling is used to resolve data imbalance.

### D.2 Results

Table D.1: Ranks of the features based on the times being predicted as direct causes of **SARS-CoV-2 exam result** out of 1000 different runs of our proposal approach. Not mentioned features were not predicted even once, note that preprocessed dataset has 331 features.

Rank	Feature	Rate of being Predicted as a Direct Cause
1	Patient age quantile Arterial Lactic Acid Promyelocytes Base excess venous blood gas analysis	1
5	pH venous blood gas analysis	0.999
6	MISSING Mean platelet volume	0.992
7	MISSING Lactic Dehydrogenase	0.966
8	Segmented	0.934
9	Myelocytes	0.904
10	Eosinophils	0.794
11	Leukocytes	0.784
12	Total CO <sub>2</sub> arterial blood gas analysis	0.450
13	Potassium	0.340
14	MISSING International normalized ratio INR	0.289
15	Metapneumovirus not detected	0.234
16	Arterial Fio2	0.092
17	HCO <sub>3</sub> arterial blood gas analysis.	0.046
18	Creatinine	0.035
19	MISSING Magnesium	0.034
20	pO <sub>2</sub> arterial blood gas analysis	0.031
21	MISSING Arterial Fio2	0.024
22	Direct Bilirubin	0.016
23	MISSING Ferritin Respiratory Syncytial Virus detected	0.014
25	MISSING Albumin Creatine phosphokinase CPK	0.010
27	Strepto A positive	0.008
28	Neutrophils Red blood cell distribution width RDW Coronavirus HKU1 detected Influenza A rapid test positive	0.004
32	Hb saturation venous blood gas analysis	0.002
33	Urine pH Inf A H1N1 2009 detected MISSING Serum Glucose Aspartate transaminase Urine Esterase nan	0.001

Table D.2: Ranks of the features based on the times being predicted as direct causes of **Patient addmited to regular ward** out of 1000 different runs of our propsal approach. Not mentiond features were not predicted even once, note that preprocessed dataset has 331 features.

Rank	Feature	Rate of being Predicted as a Direct Cause
1	Patient age quantile HCO3 venous blood gas analysis Total CO2 venous blood gas analysis Gamma glutamyltransferase	1
5	MISSING Lactic Dehydrogenase	0.987
6	Alanine transaminase	0.845
7	MISSING International normalized ratio INR	0.804
8	Serum Glucose	0.652
9	pH venous blood gas analysis	0.631
10	Base.excess venous blood gas analysis	0.341
11	MISSING Arteiral Fio2	0.334
12	Urine Density	0.334
13	Magnesium	0.323
14	Metapneumovirus not detected	0.261
15	MISSING Mean platelet volume	0.118
16	Creatine phosphokinase CPK	0.086
17	Creatinine	0.058
18	International normalized ratio INR	0.049
19	MISSING Ferritin	0.046
20	Urea	0.044
21	Respiratory Syncytial Virus detected	0.032
22	MISSING Magnesium	0.021
23	MISSING Albumin	0.018
24	MISSING Potassium	0.016
25	Inf A H1N1 2009 detected	0.014
26	Coronavirus HKU1 detected	0.010
27	Strepto A positive	0.008
28	Influenza A rapid test positive	0.007
29	MISSING Sodium Urine Protein nan	0.002
31	ctO2 arterial blood gas analysis Influenza A detected Influenza B detected	0.001

Table D.3: Ranks of the features based on the times being predicted as direct causes of **Patient addmited to semi-intensive unit** out of 1000 different runs of our proposal approach. Not mentioned features were not predicted even once, note that preprocessed dataset has 331 features.

Rank	Feature	Rate of being Predicted as a Direct Cause
1	Patient age quantile Creatinine MISSING Lactic Dehydrogenase Total CO2 venous blood gas analysis Magnesium Gamma glutamyltransferase Alanine transaminase	1
8	ctO2 arterial blood gas analysis HCO3 venous blood gas analysis	0.999
10	Relationship Patient Normal	0.786
11	MISSING Arterial Fio2	0.595
12	Base excess venous blood gas analysis	0.578
13	pO2 venous blood gas analysis	0.449
14	MISSING International normalized ratio INR	0.435
15	Mean platelet volume	0.366
16	Metapneumovirus not detected	0.308
17	Proteina C reactiva mg dL	0.235
18	Sodium	0.212
19	Phosphor	0.164
20	Urine Density	0.085
21	Respiratory Syncytial Virus detected	0.068
22	MISSING Mean platelet volume	0.056
23	MISSING Ferritin	0.054
24	pH venous blood gas analysis	0.021
25	Strepto A positive	0.018
26	Inf A H1N1 2009 detected	0.016
27	Influenza A rapid test positive	0.014
28	MISSING Albumin Coronavirus HKU1 detected	0.012
30	MISSING Magnesium	0.008
31	Aspartate transaminase	0.004
32	Urine Ketone Bodies absent Red blood cell distribution width RDW Influenza A detected Urine Esterase absent Urine Protein nan	0.001

Table D.4: Ranks of the features based on the times being predicted as direct causes of **Patient addmited to intensive care unit** out of 1000 different runs of our propsal approach. Not mentiond features were not predicted even once, note that preprocessed dataset has 331 features.

Rank	Feature	Rate of being Predicted as a Direct Cause
1	Patient age quantile MISSING Mean platelet volume Total CO2 venous blood gas analysis HCO3 venous blood gas analysis Alanine transaminase Gamma glutamyltransferase Magnesium MISSING Lactic Dehydrogenase Creatinine	1
10	pO2 venous blood gas analysis	0.982
11	ctO2 arterial blood gas analysis	0.962
12	pH venous blood gas analysis	0.938
13	MISSING Arteiral Fio2	0.667
14	MISSING International normalized ratio INR	0.586
15	Red blood cell distribution width RDW	0.503
16	Urine Density	0.414
17	Creatine phosphokinase CPK	0.380
18	Base excess venous blood gas analysis	0.352
19	Potassium	0.234
20	Promyelocytes	0.221
21	MISSING Ferritin	0.174
22	Metapneumovirus not detected	0.132
23	Phosphor	0.082
24	Sodium	0.036
25	MISSING Magnesium	0.032
26	Proteina C reativa mg dL	0.016
27	Aspartate transaminase	0.015
28	Respiratory Syncytial Virus detected	0.010
29	Relationship Patient Normal	0.007
30	MISSING Albumin Arterial Lactic Acid	0.006
32	Coronavirus HKU1 detected Eosinophils	0.005
34	Inf A H1N1 2009 detected	0.004
35	Influenza A rapid test positive International normalized ratio INR	0.002
37	Urine Crystals Ausentes Leukocytes Strepto A positive	0.001

## E Supplementary Tables for Performance in Inferring Direct Causes

Table E.1: Performance across all the settings for different number of nodes. Each single entry in the table is averaged over 18000 simulations. Our method is almost state of the art in every case.

Method	Number of Nodes											
	5			10			20			50		
ACC	CSI	F1	ACC	CSI	F1	ACC	CSI	F1	ACC	CSI	F1	
GES	0.935	0.890	0.911	0.854	0.730	0.779	0.743	0.442	0.526	0.698	0.245	0.323
rankGES	0.923	0.857	0.883	0.846	0.700	0.753	0.740	0.428	0.514	0.697	0.237	0.316
ARGES	0.922	0.864	0.885	0.797	0.551	0.584	0.752	0.447	0.524	0.705	0.186	0.221
rankARGES	0.914	0.838	0.861	0.793	0.537	0.572	0.750	0.435	0.514	0.705	0.181	0.216
FCI+	0.963	0.918	0.932	0.873	0.744	0.808	0.830	0.602	0.703	0.766	0.368	0.486
LINGAM	<b>0.991</b>	<b>0.978</b>	<b>0.982</b>	<b>0.953</b>	0.865	0.889	0.891	0.712	0.778	0.750	0.318	0.385
PC	0.957	0.913	0.929	0.864	0.723	0.786	0.823	0.569	0.664	0.763	0.348	0.457
rankPC	0.946	0.891	0.912	0.854	0.701	0.768	0.813	0.541	0.638	0.754	0.324	0.431
MMHC	0.929	0.878	0.905	0.841	0.675	0.739	0.767	0.432	0.507	0.725	0.218	0.281
Lasso	0.965	0.948	0.968	0.905	0.834	0.892	0.894	0.786	0.866	0.773	0.489	0.627
<b>CORTH Features (Ours)</b>	0.988	0.968	0.973	0.949	<b>0.908</b>	<b>0.934</b>	<b>0.949</b>	<b>0.865</b>	<b>0.905</b>	<b>0.795</b>	<b>0.559</b>	<b>0.663</b>
Method	Number of Nodes											
	5			10			20			50		
TPR	FPR	MCC	TPR	FPR	MCC	TPR	FPR	MCC	TPR	FPR	MCC	
GES	0.934	0.056	0.891	0.790	0.090	0.711	0.502	0.088	0.436	0.304	0.083	0.221
rankGES	0.924	0.068	0.877	0.780	0.098	0.695	0.493	0.089	0.425	0.297	0.083	0.215
ARGES	0.903	0.046	0.906	0.590	0.041	0.841	0.500	0.073	0.557	0.220	0.020	0.794
rankARGES	0.897	0.054	0.896	0.584	0.044	0.832	0.495	0.075	0.549	0.216	0.020	0.789
FCI+	0.969	0.029	0.948	0.797	0.054	0.759	0.642	0.042	0.645	0.389	0.030	0.454
LINGAM	0.991	0.007	<b>0.988</b>	0.886	0.008	<b>0.934</b>	0.770	0.055	0.759	0.391	0.072	0.471
PC	0.950	0.024	0.941	0.759	0.041	0.759	0.600	0.032	0.650	0.363	0.021	0.468
rankPC	0.944	0.039	0.925	0.750	0.053	0.734	0.580	0.034	0.629	0.341	0.024	0.427
MMHC	0.895	0.011	0.903	0.691	0.015	0.724	0.444	0.009	0.523	0.219	0.005	0.330
Lasso	0.999	0.074	0.949	0.944	0.119	0.817	0.954	0.147	0.794	0.681	0.148	0.488
<b>CORTH Features (Ours)</b>	0.999	0.016	0.986	0.952	0.044	0.906	0.884	0.011	<b>0.894</b>	0.609	0.101	<b>0.567</b>

Table E.2: Performance across all the settings for different sparsities. Each single entry in the table is averaged over 24000 simulations. Our method is state of the art in every case.

Method	Sparsity											
	0.1				0.3				0.5			
ACC	CSI	F1	MCC	ACC	CSI	F1	MCC	ACC	CSI	F1	MCC	
GES	0.961	0.786	0.825	0.857	0.815	0.539	0.598	0.522	0.646	0.405	0.482	0.315
rankGES	0.954	0.746	0.790	0.840	0.809	0.522	0.584	0.511	0.642	0.398	0.475	0.308
ARGES	0.965	0.794	0.828	0.876	0.805	0.456	0.501	0.726	0.612	0.286	0.330	0.720
rankARGES	0.959	0.763	0.801	0.863	0.802	0.447	0.494	0.721	0.611	0.282	0.328	0.716
FCI+	0.974	0.819	0.853	0.910	0.866	0.631	0.714	0.674	0.734	0.524	0.629	0.521
LINGAM	0.966	0.763	0.796	0.889	0.896	0.710	0.753	0.761	0.827	0.682	0.727	0.715
PC	0.975	0.819	0.849	0.921	0.861	0.609	0.689	0.676	0.718	0.486	0.588	0.516
rankPC	0.971	0.797	0.831	0.912	0.852	0.587	0.670	0.653	0.701	0.458	0.560	0.470
MMHC	0.978	0.834	0.867	0.901	0.830	0.497	0.561	0.574	0.639	0.321	0.397	0.385
Lasso	0.976	0.886	0.925	0.926	0.876	0.725	0.811	0.737	0.800	0.682	0.778	0.622
<b>CORTH Features (Ours)</b>	<b>0.988</b>	<b>0.915</b>	<b>0.934</b>	<b>0.959</b>	<b>0.926</b>	<b>0.813</b>	<b>0.858</b>	<b>0.833</b>	<b>0.847</b>	<b>0.747</b>	<b>0.814</b>	<b>0.724</b>

Table E.3: Performance across all the settings for different number of nonlinear probabilities. Each single entry in the table is averaged over 18000 simulations. Our method is almost state of the art in every case.

Method	Nonlinear Probability											
	0			0.3			0.5			1		
ACC	CSI	F1	ACC	CSI	F1	ACC	CSI	F1	ACC	CSI	F1	
GES	0.803	0.583	0.646	0.806	0.566	0.622	0.811	0.577	0.632	0.810	0.581	0.641
rankGES	0.796	0.559	0.625	0.801	0.546	0.605	0.805	0.556	0.613	0.805	0.561	0.623
ARGES	0.781	0.476	0.515	0.786	0.486	0.525	0.792	0.506	0.546	0.818	0.581	0.628
rankARGES	0.778	0.461	0.503	0.782	0.474	0.515	0.788	0.490	0.531	0.814	0.564	0.615
FCI+	0.827	0.599	0.674	0.860	0.663	0.745	0.872	0.685	0.764	0.873	0.685	0.746
LINGAM	<b>0.907</b>	0.738	0.778	0.886	0.689	0.725	0.880	0.684	0.724	0.911	0.762	0.808
PC	0.818	0.574	0.641	0.854	0.641	0.720	0.864	0.665	0.7430	0.869	0.672	0.731
rankPC	0.813	0.560	0.630	0.841	0.614	0.694	0.848	0.627	0.704	0.864	0.656	0.720
MMHC	0.797	0.516	0.578	0.815	0.549	0.610	0.823	0.566	0.625	0.826	0.571	0.620
Lasso	0.847	0.694	0.773	0.891	0.776	0.853	0.902	0.797	0.869	0.896	0.790	0.857
<b>CORTH Features (Ours)</b>	0.871	<b>0.768</b>	<b>0.824</b>	<b>0.934</b>	<b>0.830</b>	<b>0.873</b>	<b>0.943</b>	<b>0.851</b>	<b>0.891</b>	<b>0.933</b>	<b>0.852</b>	<b>0.887</b>
Method	Nonlinear Probability											
	0			0.3			0.5			1		
TPR	FPR	MCC	TPR	FPR	MCC	TPR	FPR	MCC	TPR	FPR	MCC	
GES	0.643	0.093	0.564	0.620	0.074	0.557	0.629	0.071	0.568	0.637	0.079	0.570
rankGES	0.633	0.100	0.550	0.612	0.080	0.546	0.620	0.076	0.557	0.628	0.083	0.559
ARGES	0.514	0.041	0.789	0.526	0.041	0.793	0.547	0.043	0.791	0.626	0.055	0.725
rankARGES	0.509	0.044	0.780	0.522	0.044	0.788	0.540	0.046	0.783	0.620	0.059	0.715
FCI+	0.638	0.045	0.637	0.704	0.037	0.708	0.728	0.035	0.731	0.728	0.037	0.730
LINGAM	0.775	0.025	<b>0.832</b>	0.723	0.028	0.759	0.722	0.034	0.741	0.819	0.053	0.822
PC	0.605	0.037	0.649	0.672	0.027	0.707	0.695	0.025	0.728	0.702	0.029	0.734
rankPC	0.597	0.043	0.626	0.656	0.040	0.680	0.668	0.036	0.695	0.692	0.031	0.714
MMHC	0.528	0.017	0.581	0.561	0.008	0.623	0.578	0.007	0.636	0.582	0.008	0.639
Lasso	0.823	0.130	0.684	0.907	0.120	0.778	0.926	0.116	0.800	0.921	0.122	0.787
<b>CORTH Features (Ours)</b>	0.840	0.119	0.730	0.849	0.007	<b>0.872</b>	0.870	0.008	<b>0.888</b>	0.885	0.038	<b>0.863</b>

Table E.4: Performance across all the settings for different number of observations. Each single entry in the table is averaged over 24000 simulations. Our method is almost state of the art in every case.

Method	Number of Observations											
	100				500				1000			
ACC	CSI	F1	MCC	ACC	CSI	F1	MCC	ACC	CSI	F1	MCC	
GES	0.797	0.524	0.588	0.539	0.811	0.593	0.650	0.572	0.815	0.612	0.666	0.583
rankGES	0.788	0.495	0.561	0.522	0.806	0.576	0.636	0.564	0.810	0.595	0.652	0.573
ARGES	0.780	0.446	0.489	0.786	0.799	0.535	0.576	0.773	0.803	0.555	0.595	0.764
rankARGES	0.776	0.428	0.473	0.778	0.795	0.523	0.566	0.766	0.800	0.542	0.584	0.757
FCI+	0.837	0.589	0.671	0.652	0.865	0.684	0.755	0.720	0.871	0.702	0.771	0.732
LINGAM	0.840	0.578	0.650	0.678	0.908	0.719	0.743	0.825	0.941	0.858	0.883	0.862
PC	0.830	0.568	0.642	0.661	0.858	0.662	0.732	0.719	0.866	0.684	0.752	0.733
rankPC	0.821	0.544	0.617	0.632	0.849	0.639	0.711	0.696	0.855	0.660	0.733	0.707
MMHC	0.800	0.495	0.557	0.579	0.820	0.570	0.625	0.633	0.826	0.587	0.642	0.647
Lasso	0.870	<b>0.729</b>	<b>0.812</b>	0.732	0.889	0.778	0.848	0.773	0.893	0.786	0.854	0.780
<b>CORTH Features (Ours)</b>	<b>0.883</b>	0.710	0.780	<b>0.754</b>	<b>0.935</b>	<b>0.874</b>	<b>0.906</b>	<b>0.874</b>	<b>0.942</b>	<b>0.891</b>	<b>0.920</b>	<b>0.887</b>

Table E.5: Performance across all the settings for different noise levels. Each single entry in the table is averaged over 14400 simulations. Our method is state of the art in every case.

Method	Noise Level											
	0.01				0.5				1			
	ACC	CSI	F1	MCC	ACC	CSI	F1	MCC	ACC	CSI	F1	MCC
GES	0.804	0.579	0.639	0.559	0.808	0.571	0.629	0.562	0.818	0.586	0.644	0.589
rankGES	0.797	0.557	0.619	0.548	0.802	0.552	0.613	0.551	0.812	0.565	0.625	0.577
ARGES	0.810	0.572	0.625	0.653	0.789	0.496	0.534	0.814	0.774	0.434	0.460	0.897
rankARGES	0.804	0.549	0.605	0.643	0.786	0.483	0.523	0.806	0.774	0.433	0.459	0.895
FCI+	0.843	0.617	0.691	0.674	0.865	0.678	0.753	0.717	0.874	0.697	0.766	0.740
LINGAM	0.888	0.703	0.744	0.763	0.899	0.723	0.763	0.797	0.903	0.732	0.773	0.803
PC	0.837	0.595	0.664	0.683	0.859	0.659	0.731	0.716	0.870	0.686	0.752	0.745
rankPC	0.831	0.584	0.657	0.653	0.845	0.626	0.699	0.688	0.856	0.655	0.724	0.714
MMHC	0.806	0.526	0.585	0.605	0.818	0.557	0.615	0.626	0.829	0.586	0.639	0.652
Lasso	0.868	0.728	0.807	0.725	0.891	0.780	0.852	0.779	0.898	0.794	0.861	0.793
<b>CORTH Features (Ours)</b>	<b>0.899</b>	<b>0.789</b>	<b>0.839</b>	<b>0.795</b>	<b>0.929</b>	<b>0.842</b>	<b>0.883</b>	<b>0.858</b>	<b>0.934</b>	<b>0.854</b>	<b>0.891</b>	<b>0.866</b>

## F Supplementary Figures for Parameter Estimation

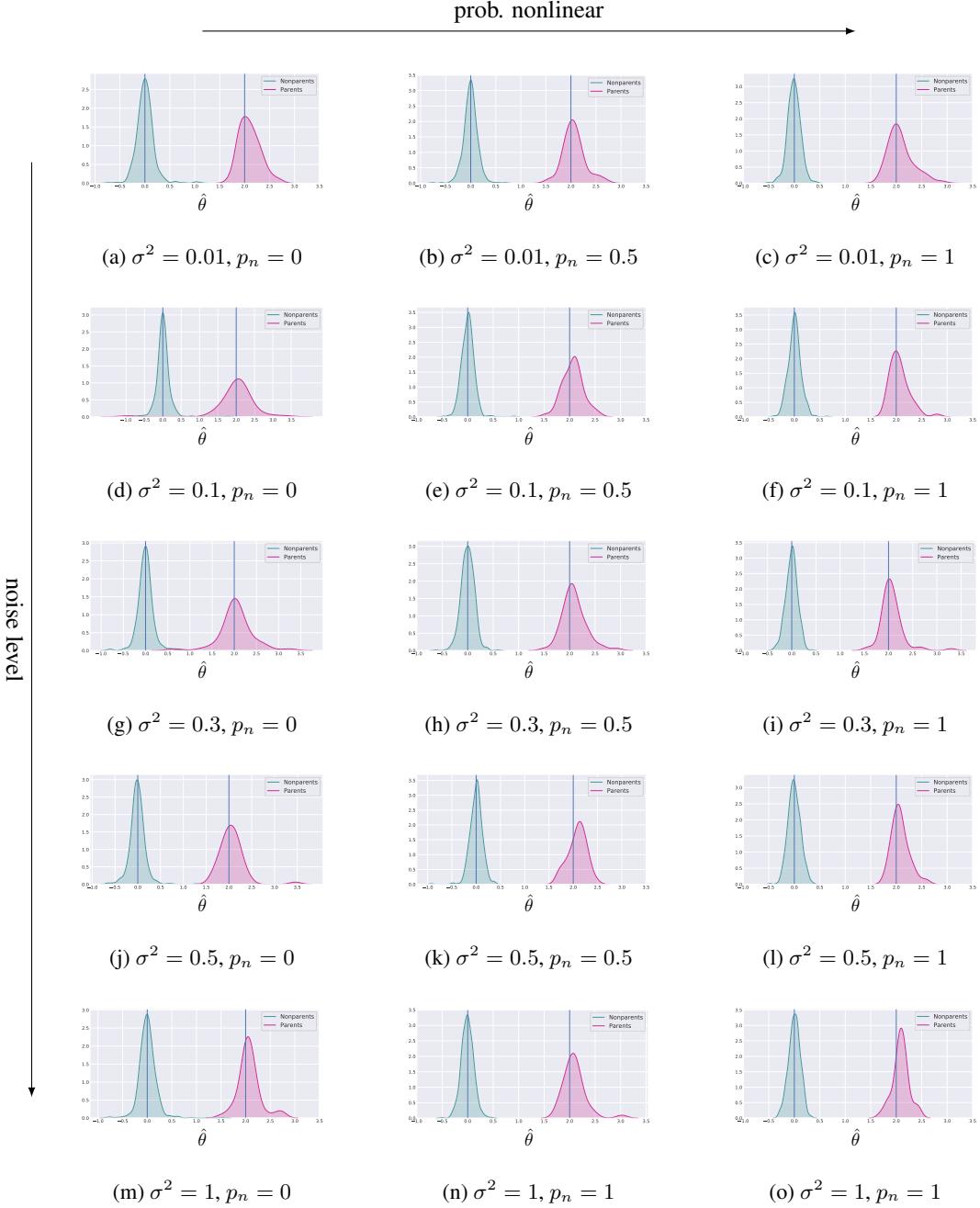


Figure F.1: 0.1 sparsity, 5 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

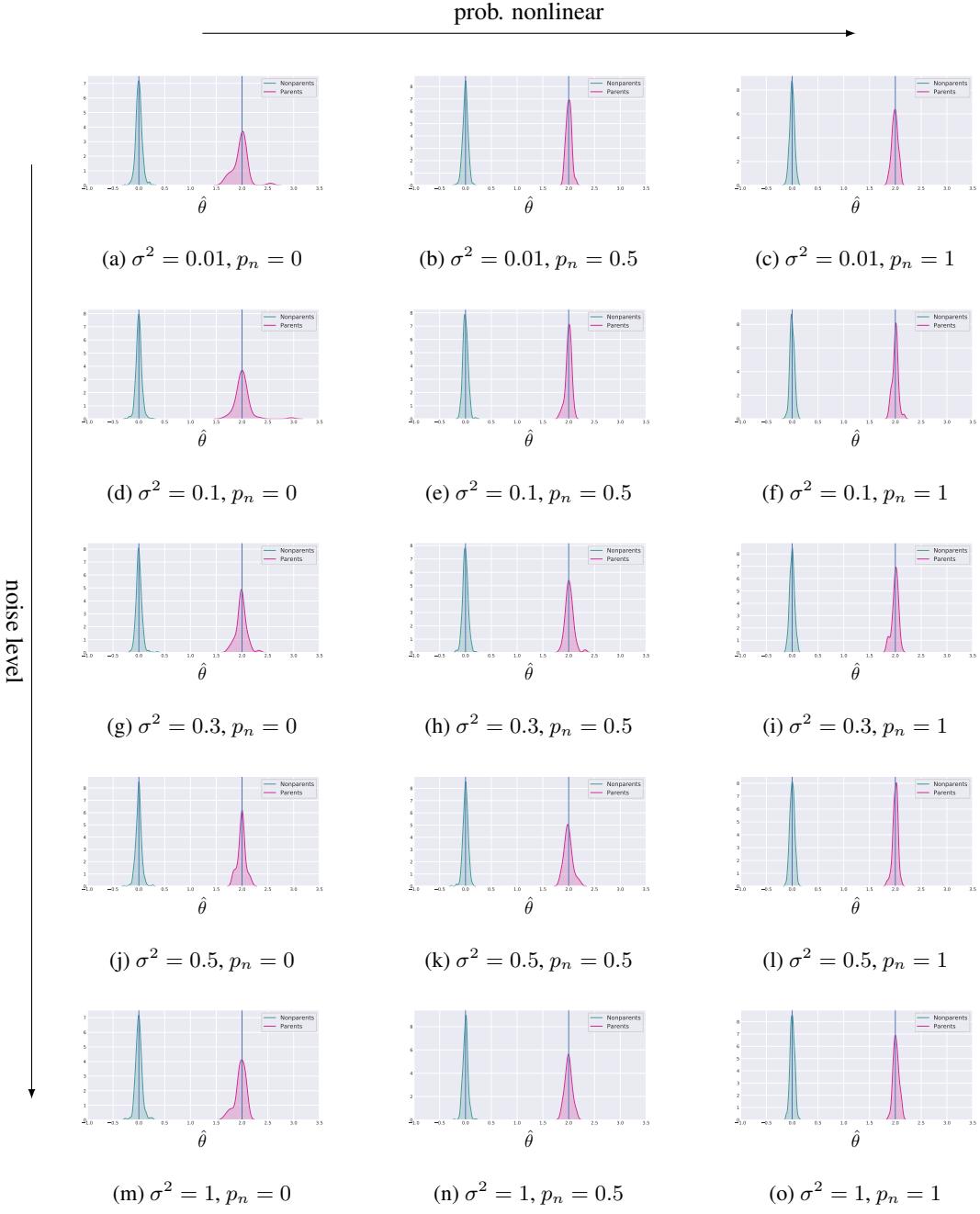


Figure F.2: 0.1 sparsity, 5 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

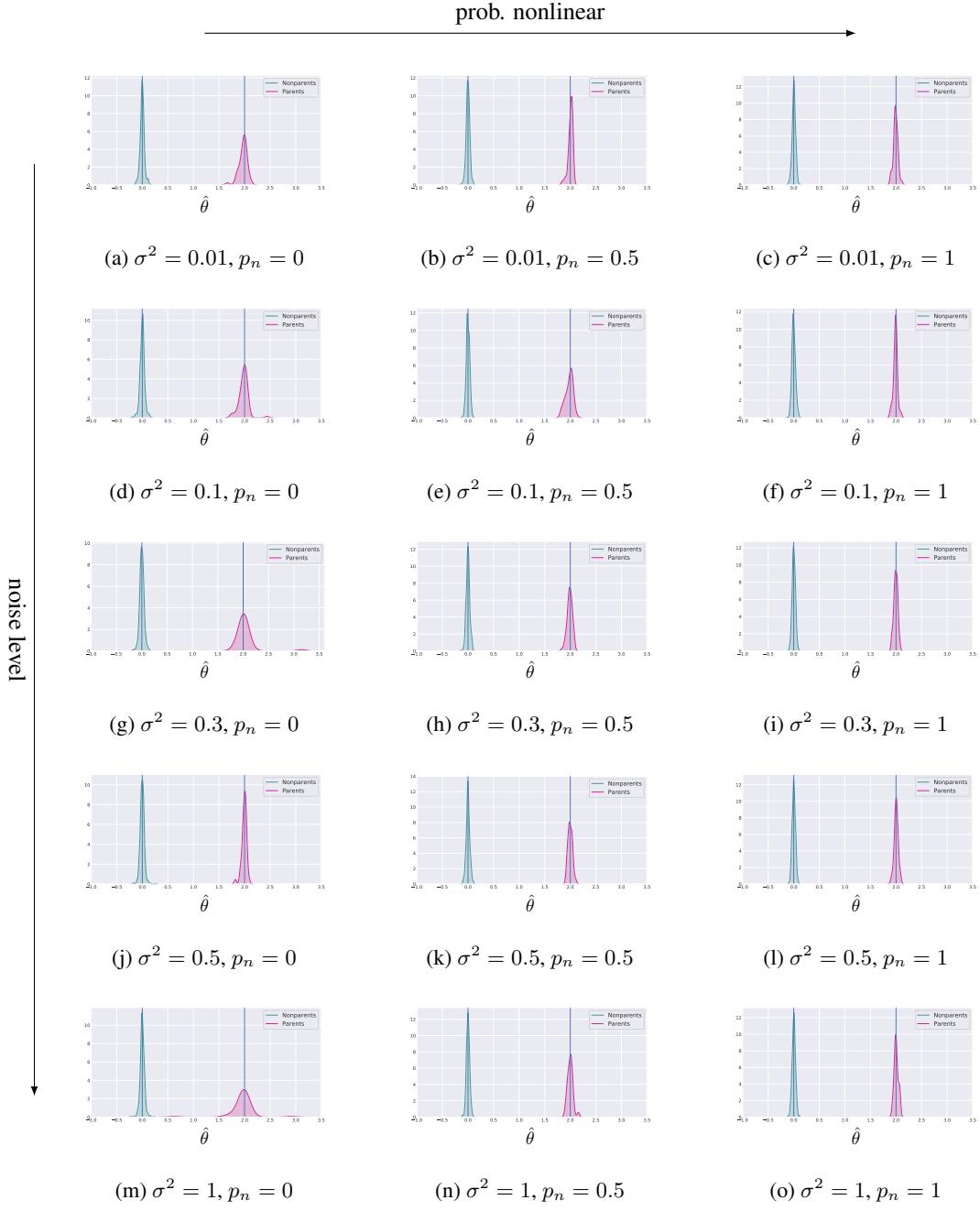


Figure F.3: 0.1 sparsity, 5 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

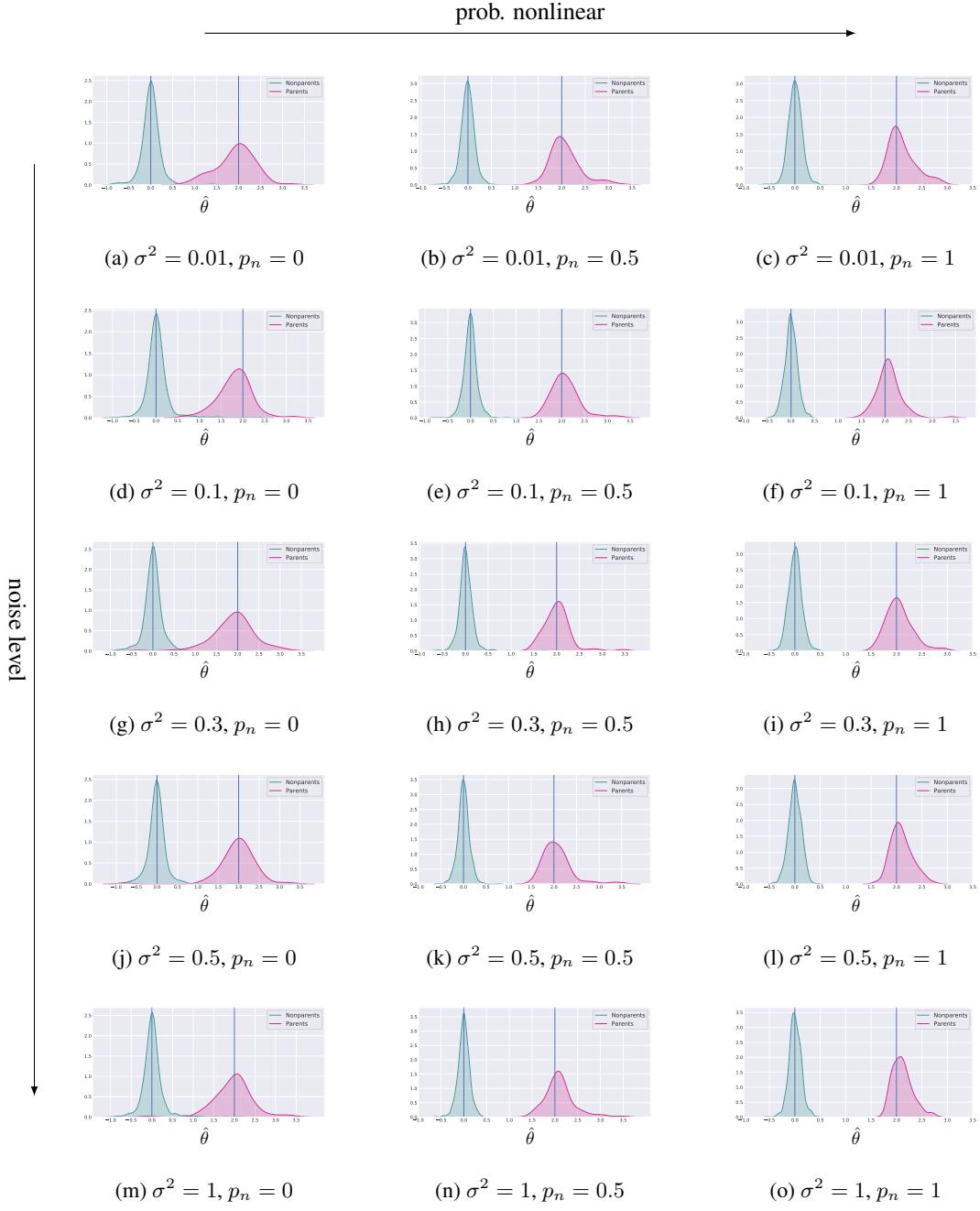


Figure F.4: 0.1 sparsity, 10 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

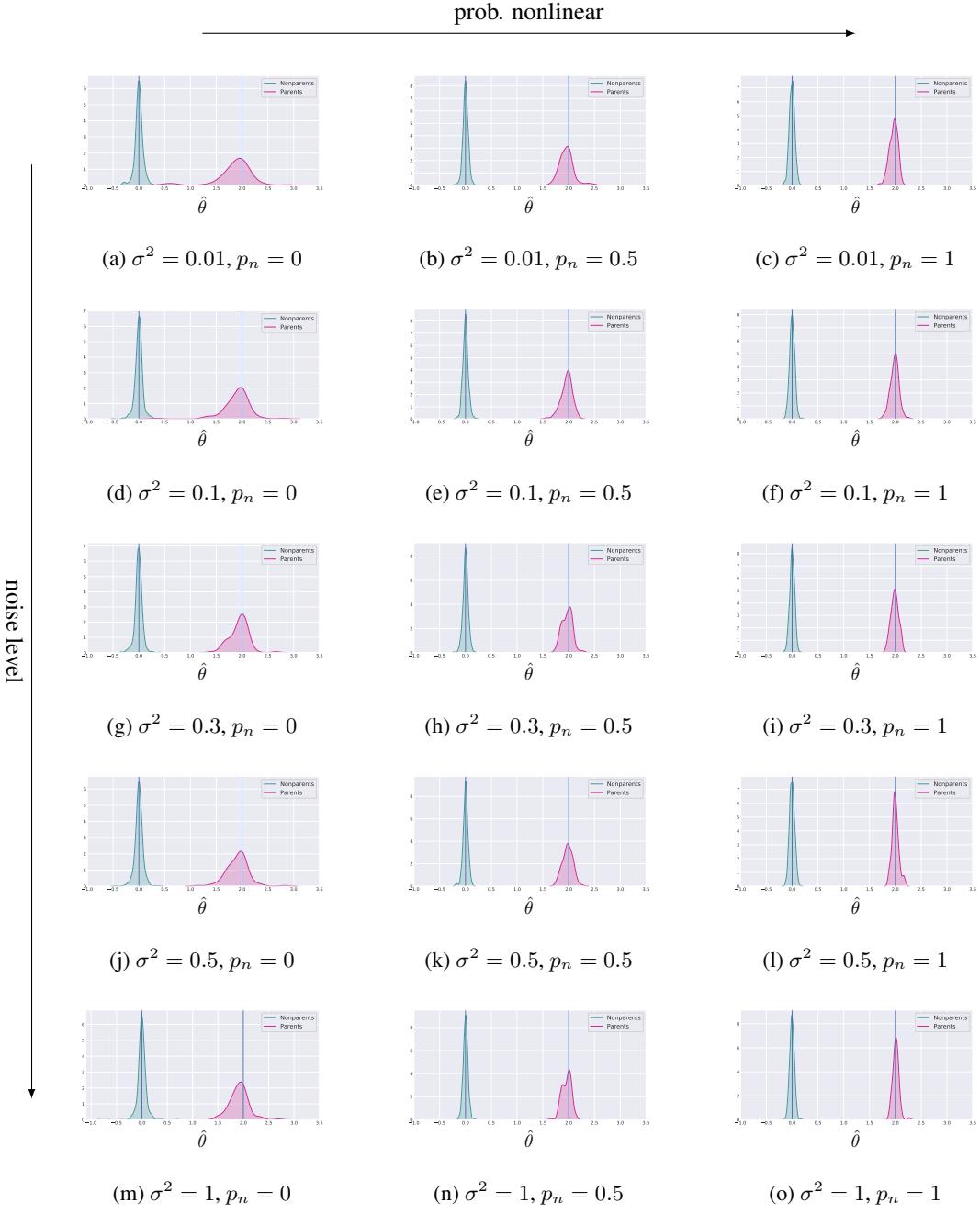


Figure F.5: sparsity 0.1, 10 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

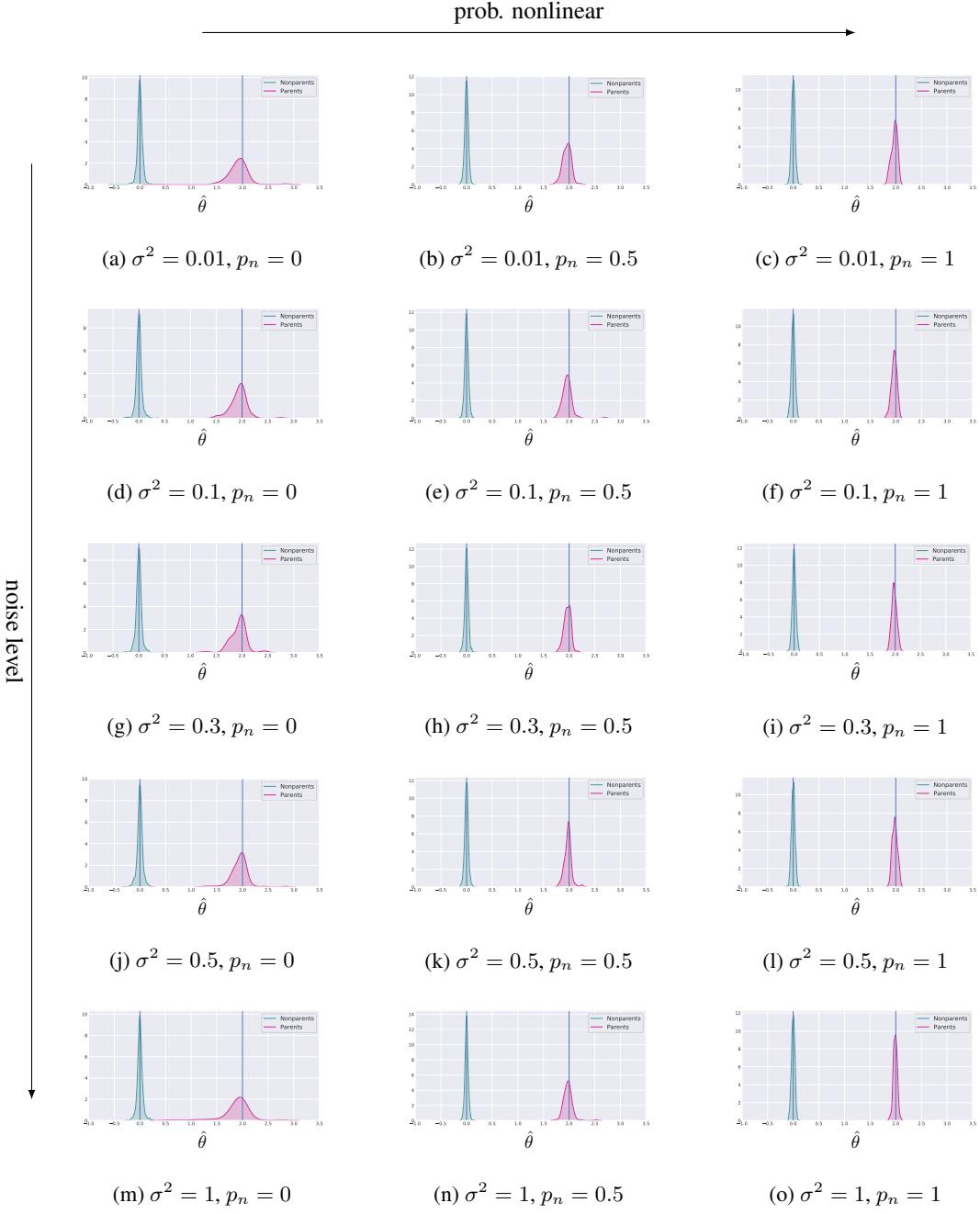


Figure F.6: sparsity 0.1, 10 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

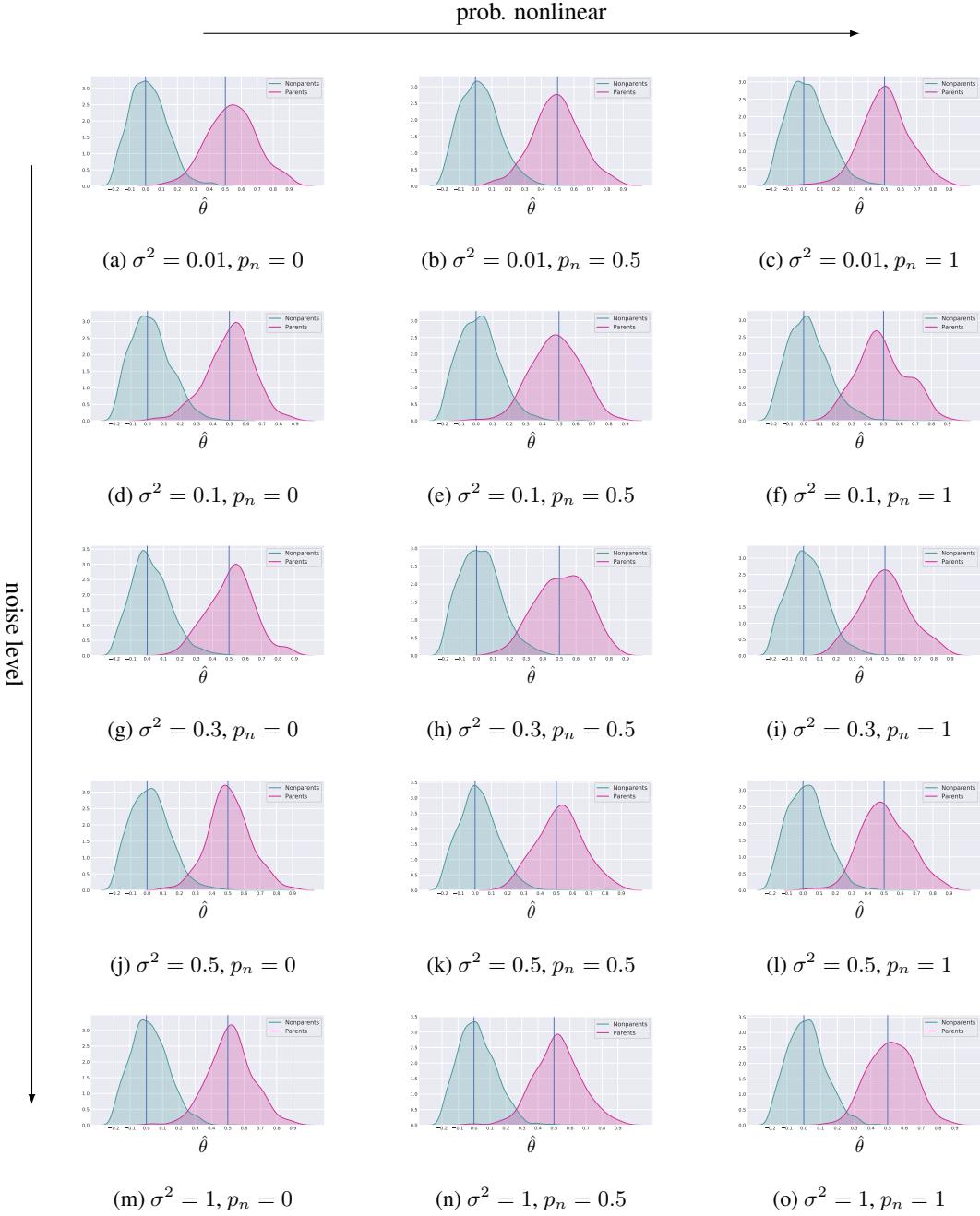


Figure F.7: sparsity 0.1, 20 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

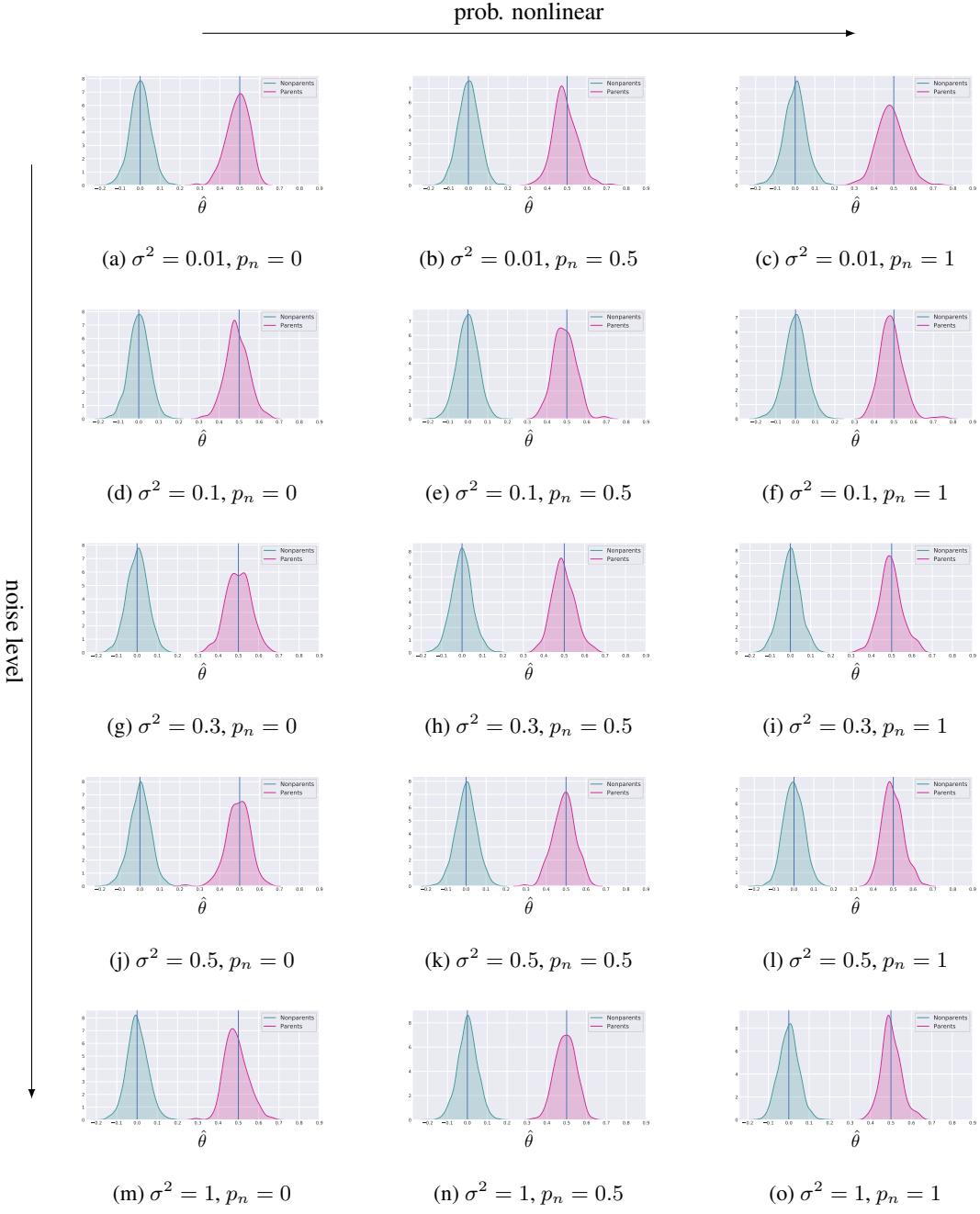


Figure F.8: sparsity 0.1, 20 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

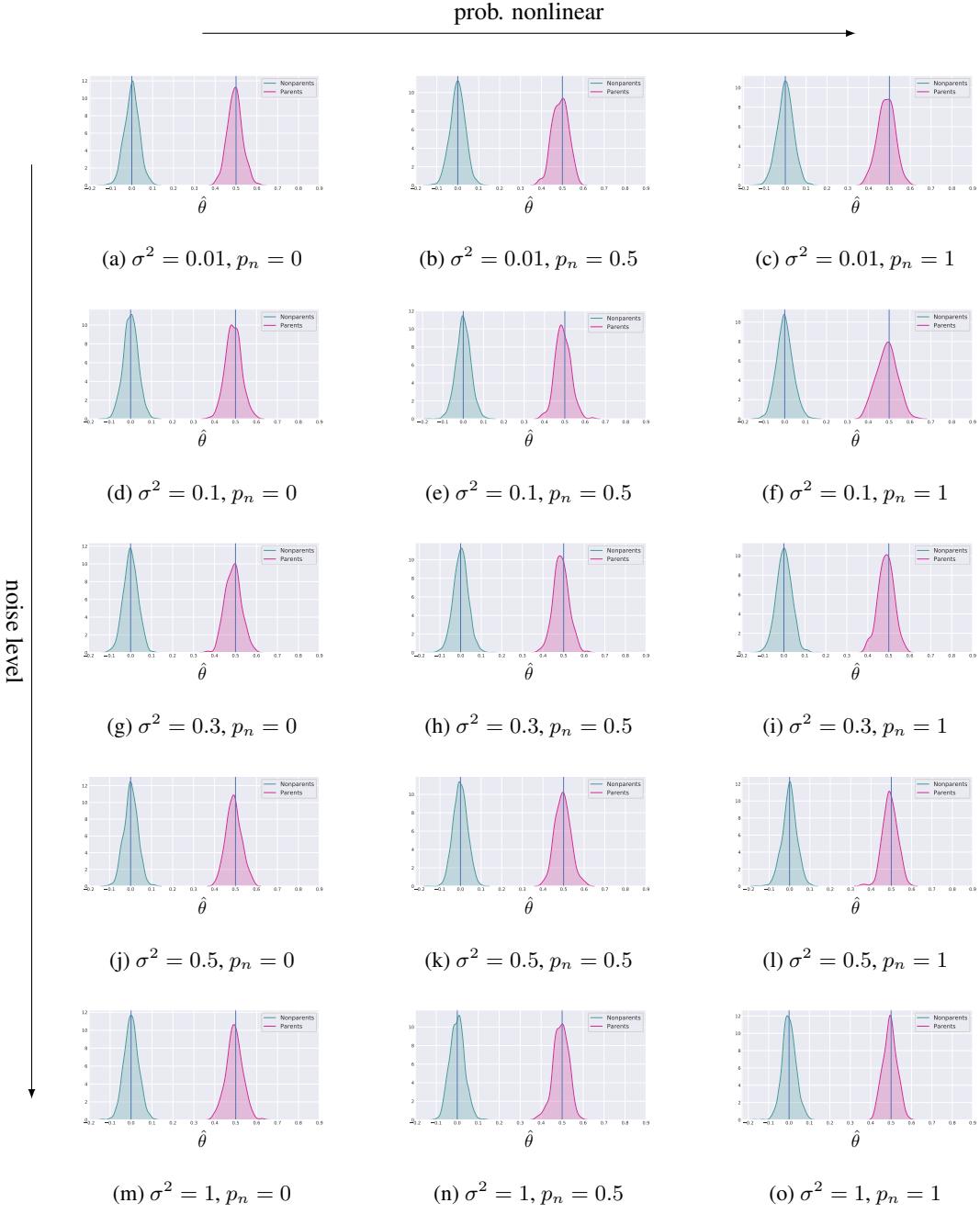


Figure F.9: sparsity 0.1, 20 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

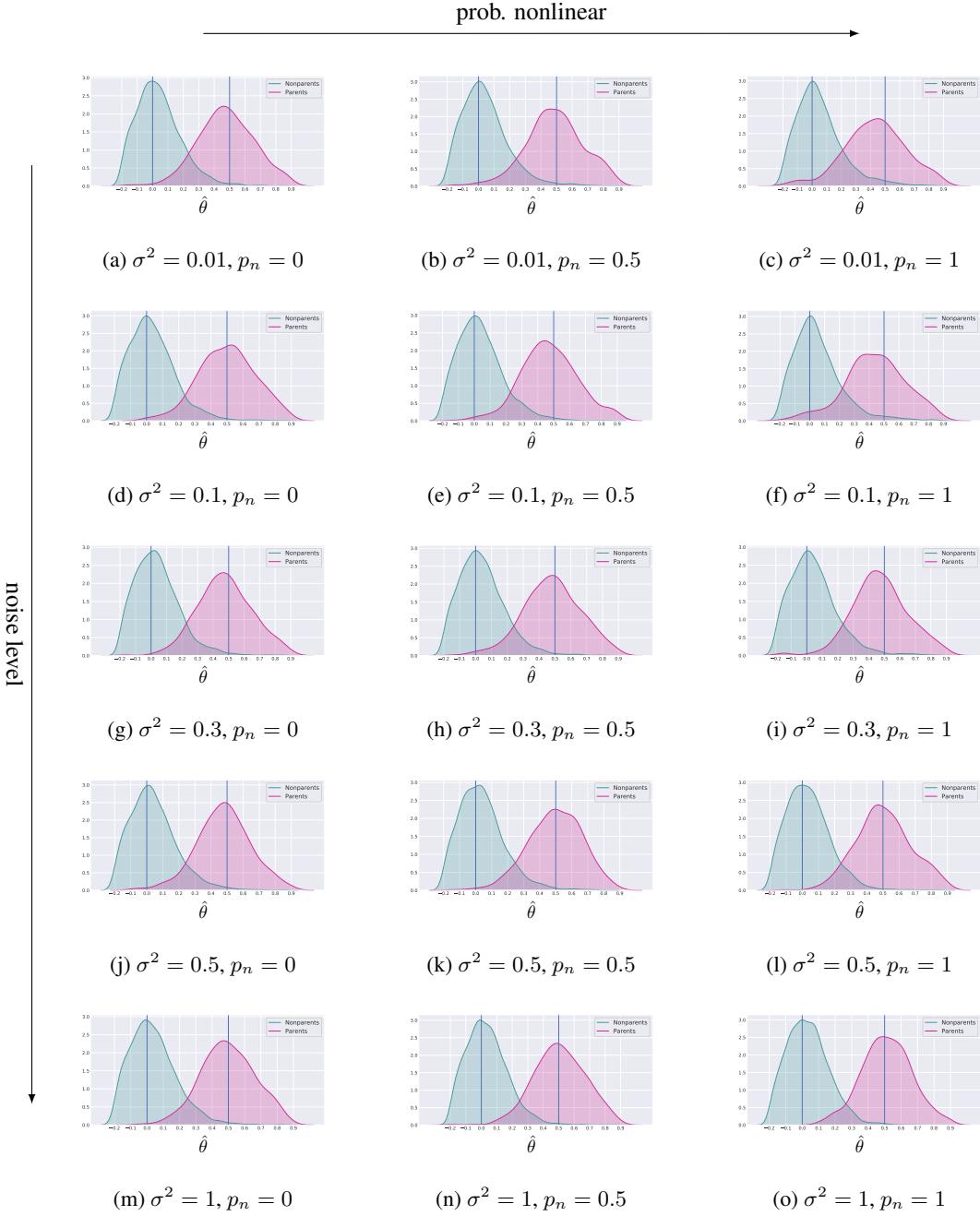


Figure F.10: sparsity 0.1, 50 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

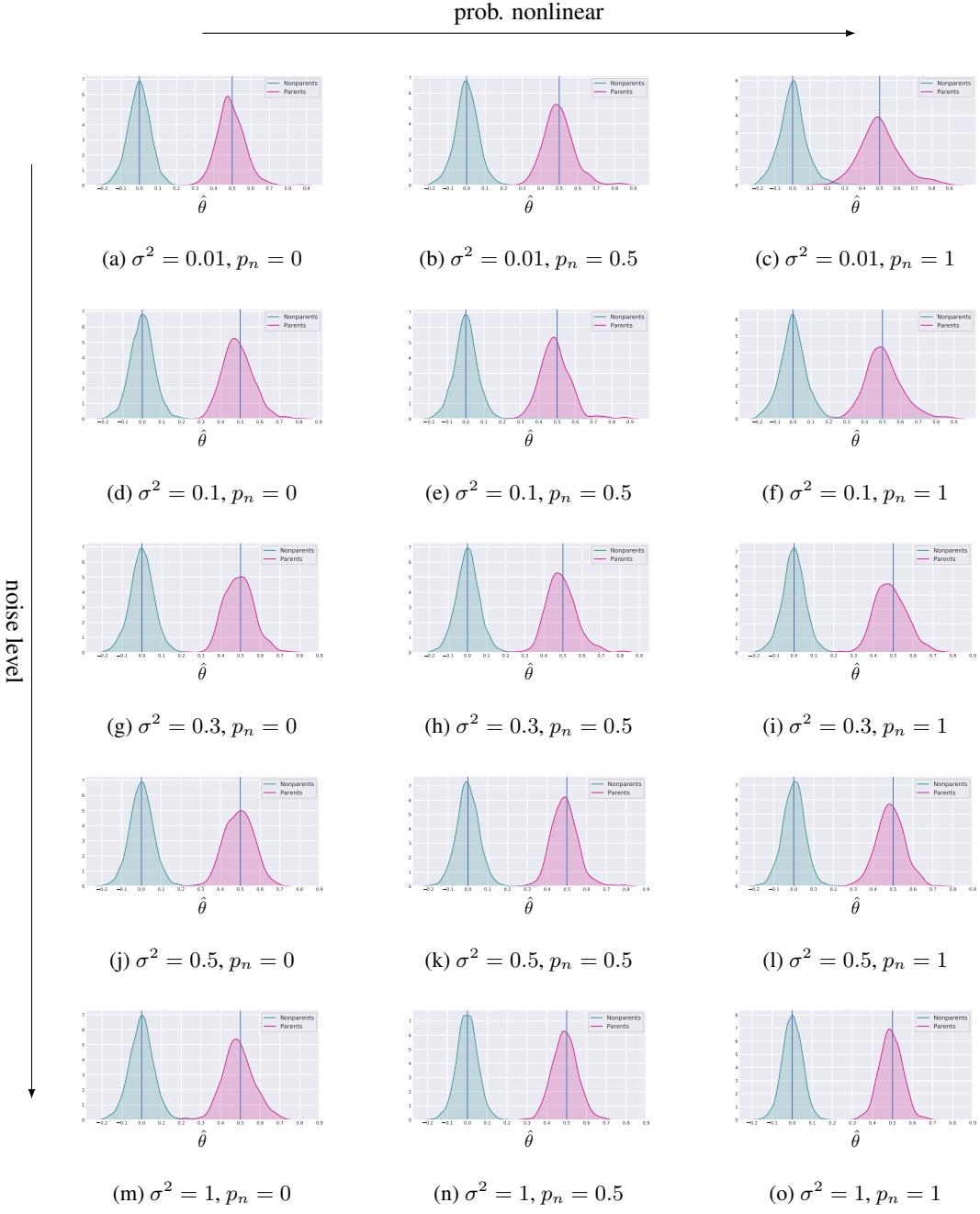


Figure F.11: sparsity 0.1, 50 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

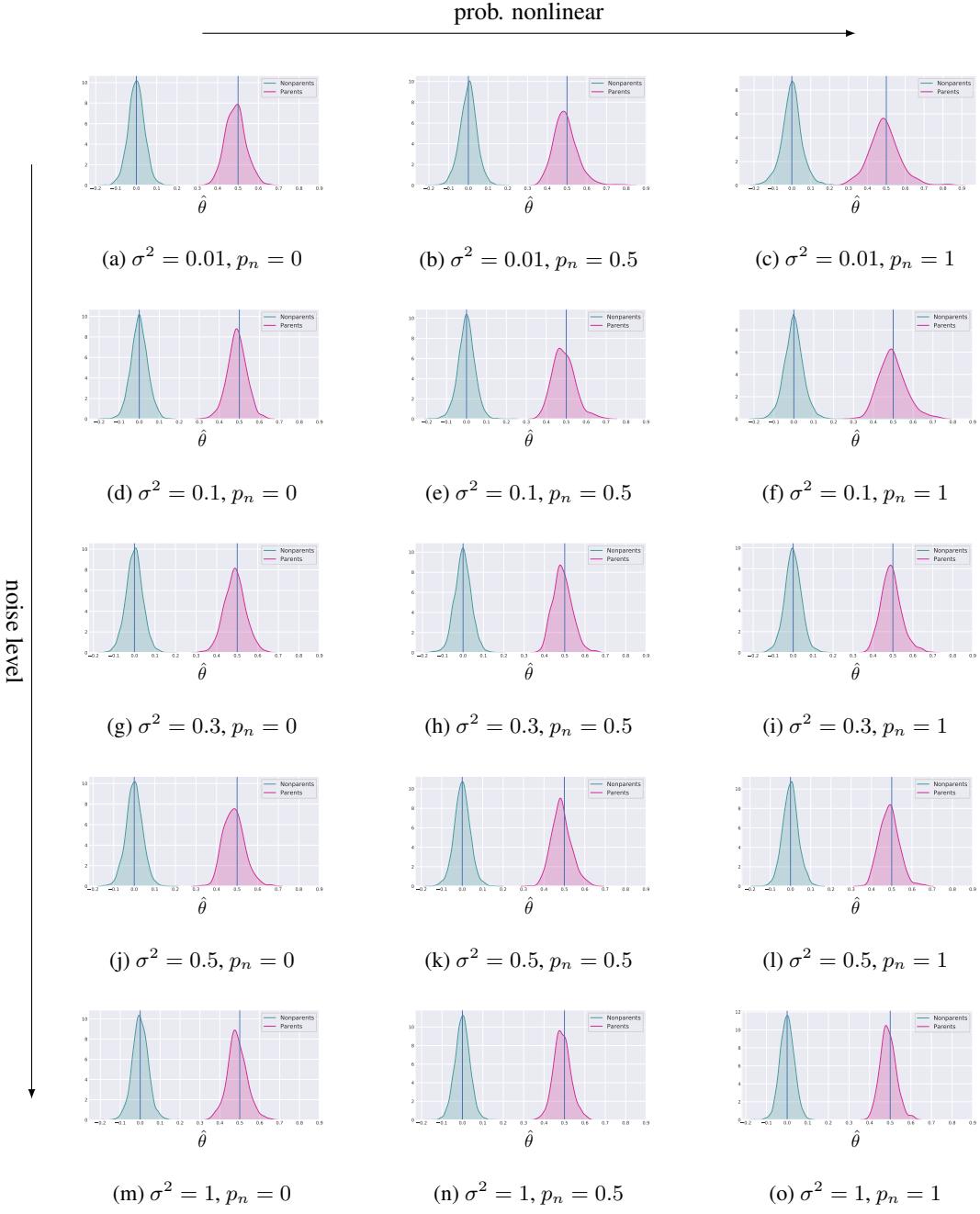


Figure F.12: sparsity 0.1, 50 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

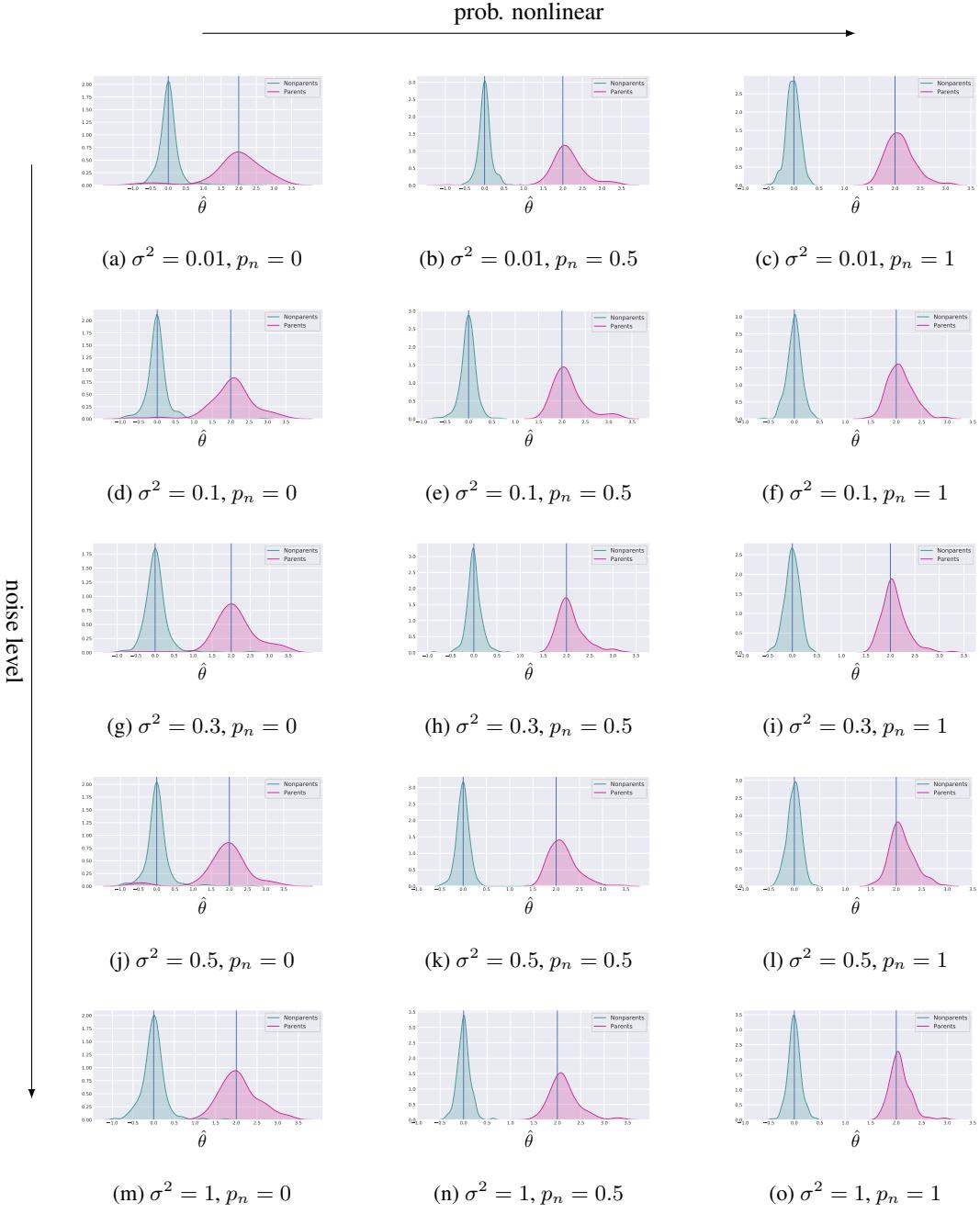


Figure F.13: 0.3 sparsity, 5 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

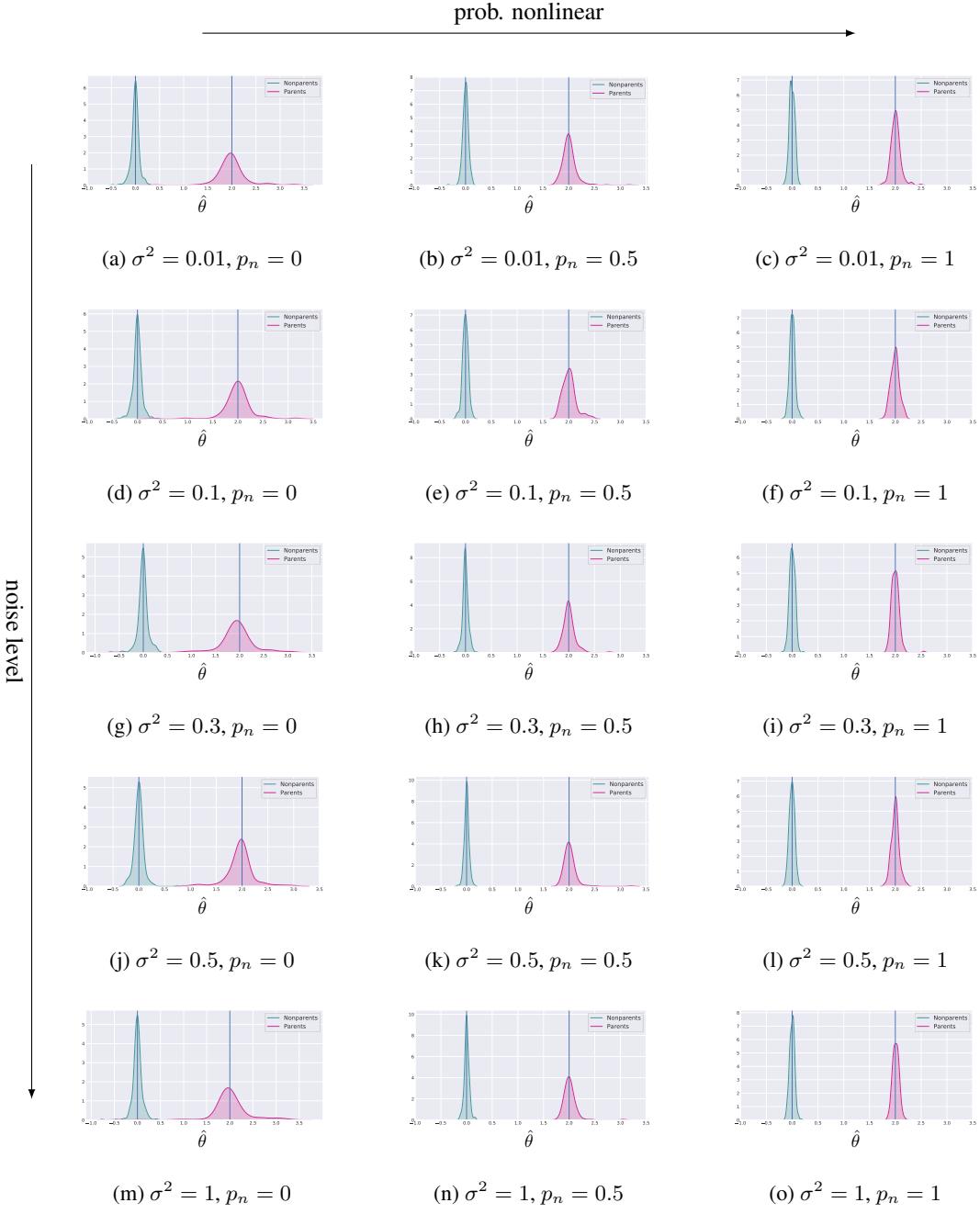


Figure F.14: 0.3 sparsity, 5 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

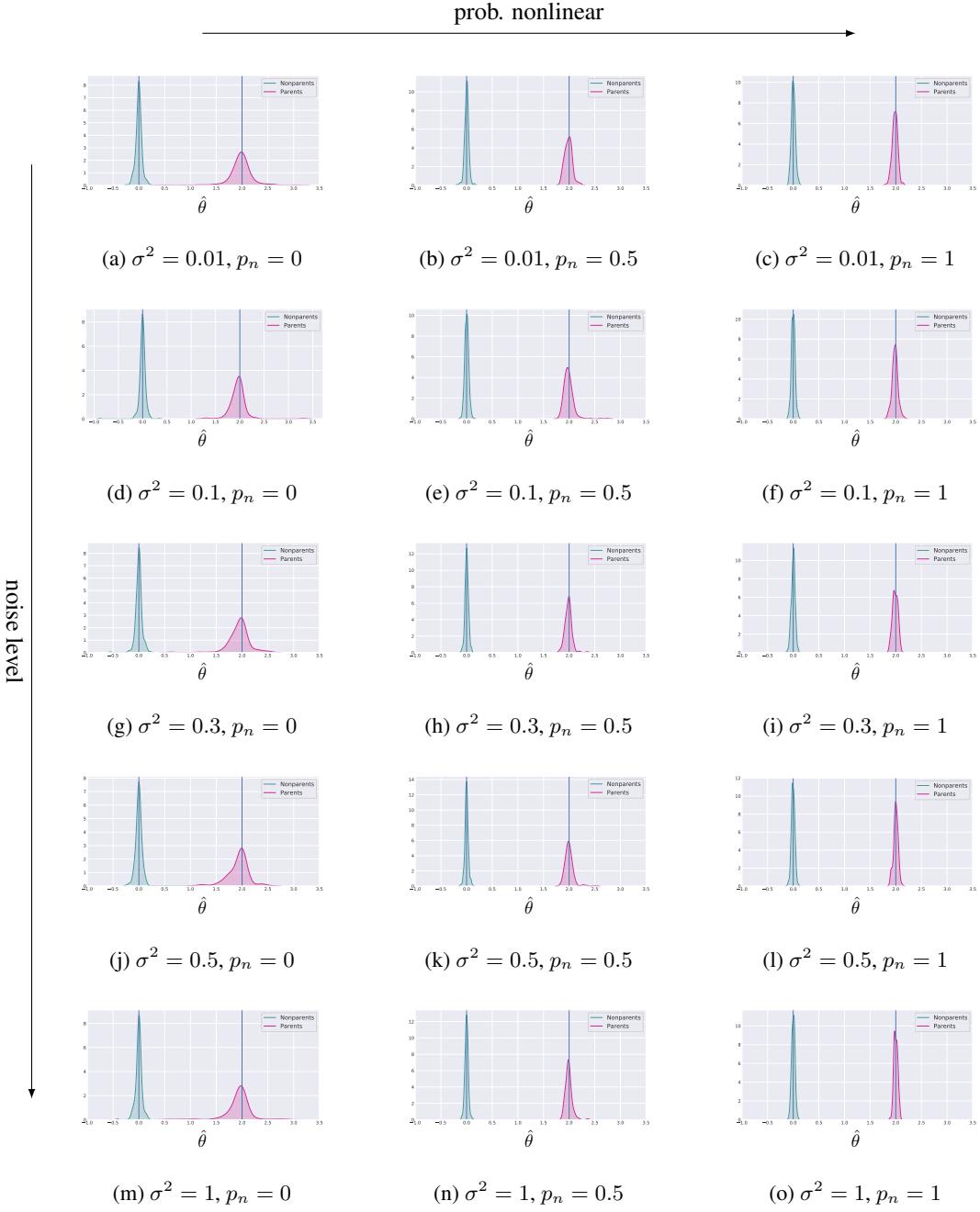


Figure F.15: 0.3 sparsity, 5 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

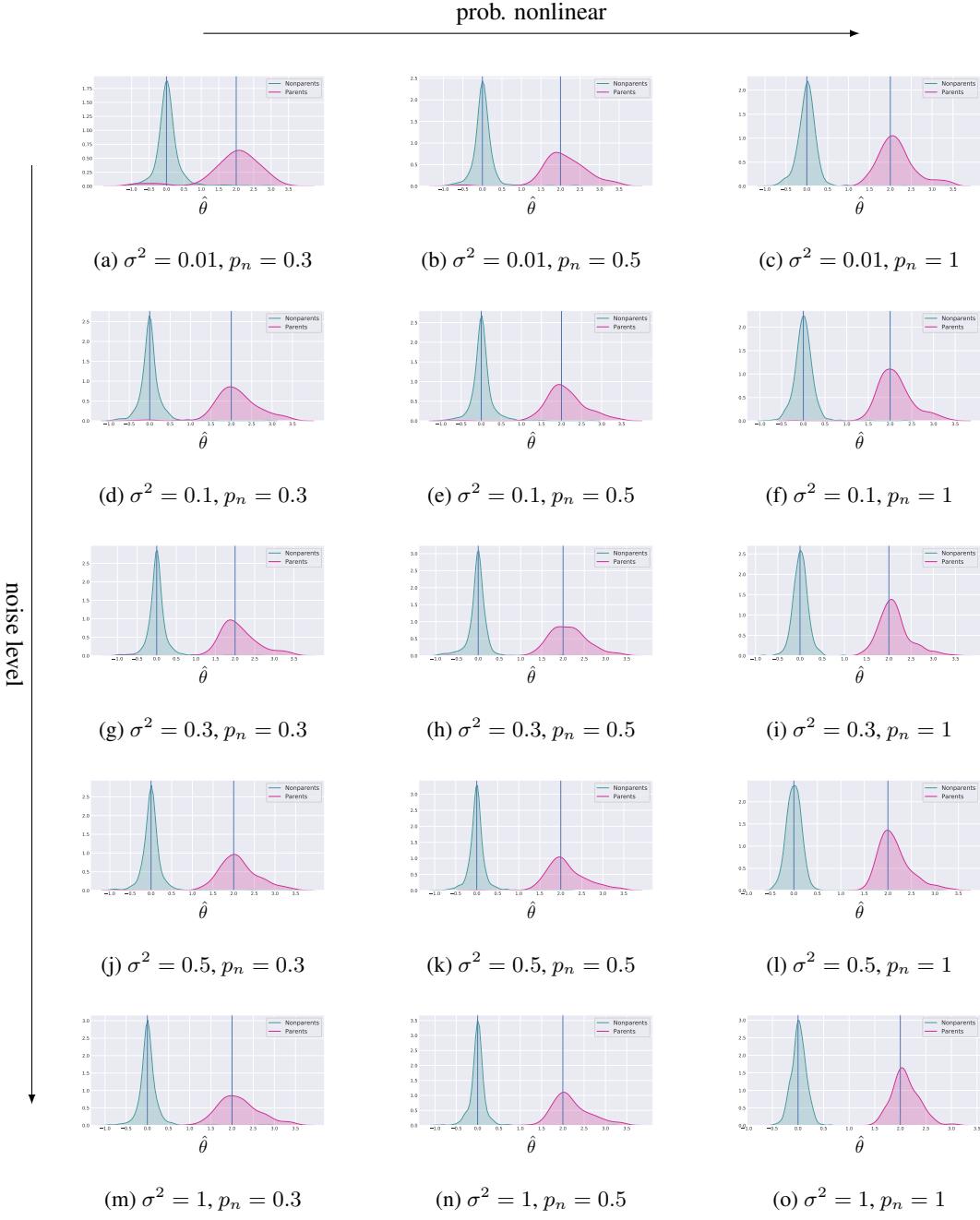


Figure F.16: 0.3 sparsity, 10 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

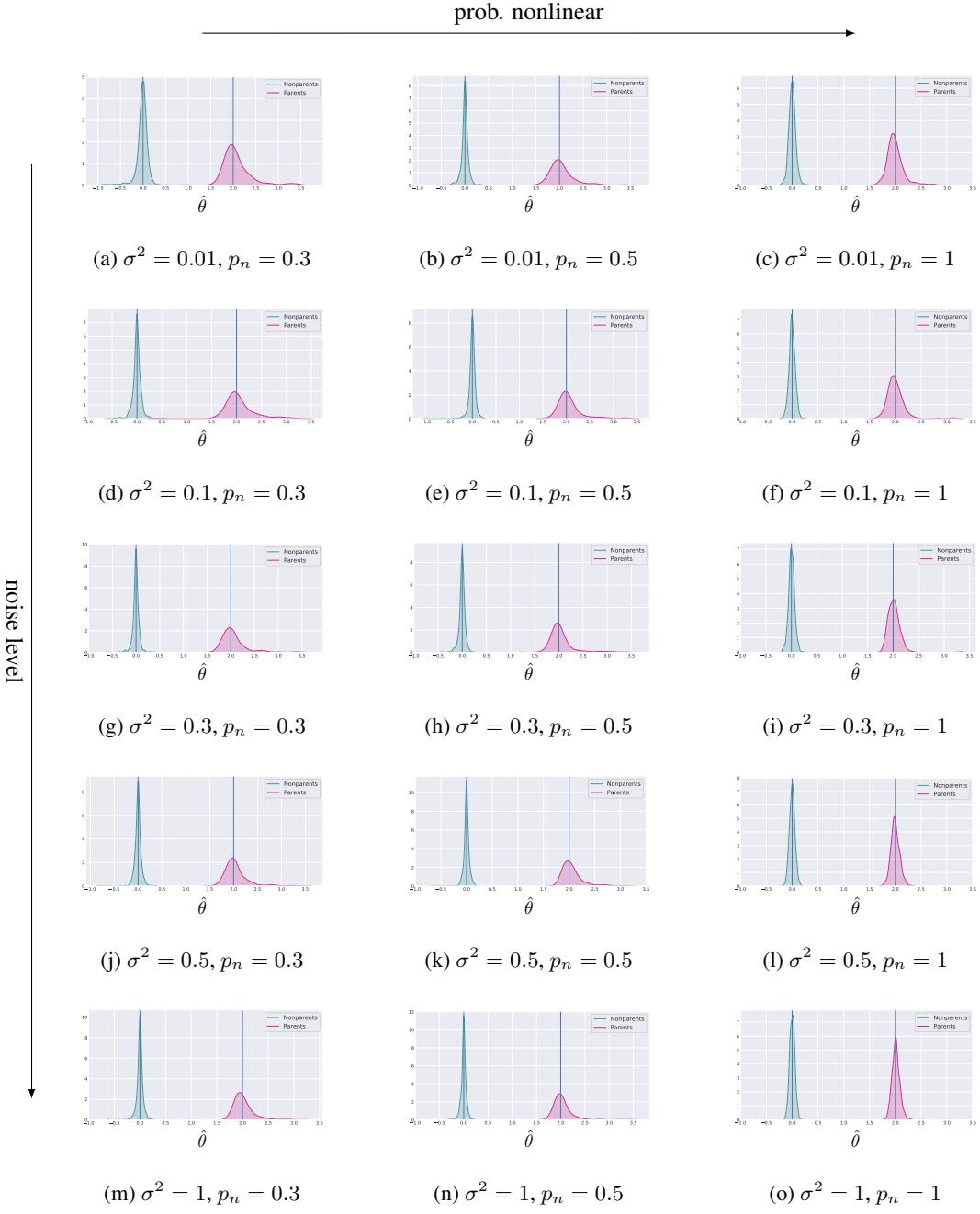


Figure F.17: 0.3 sparsity, 10 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

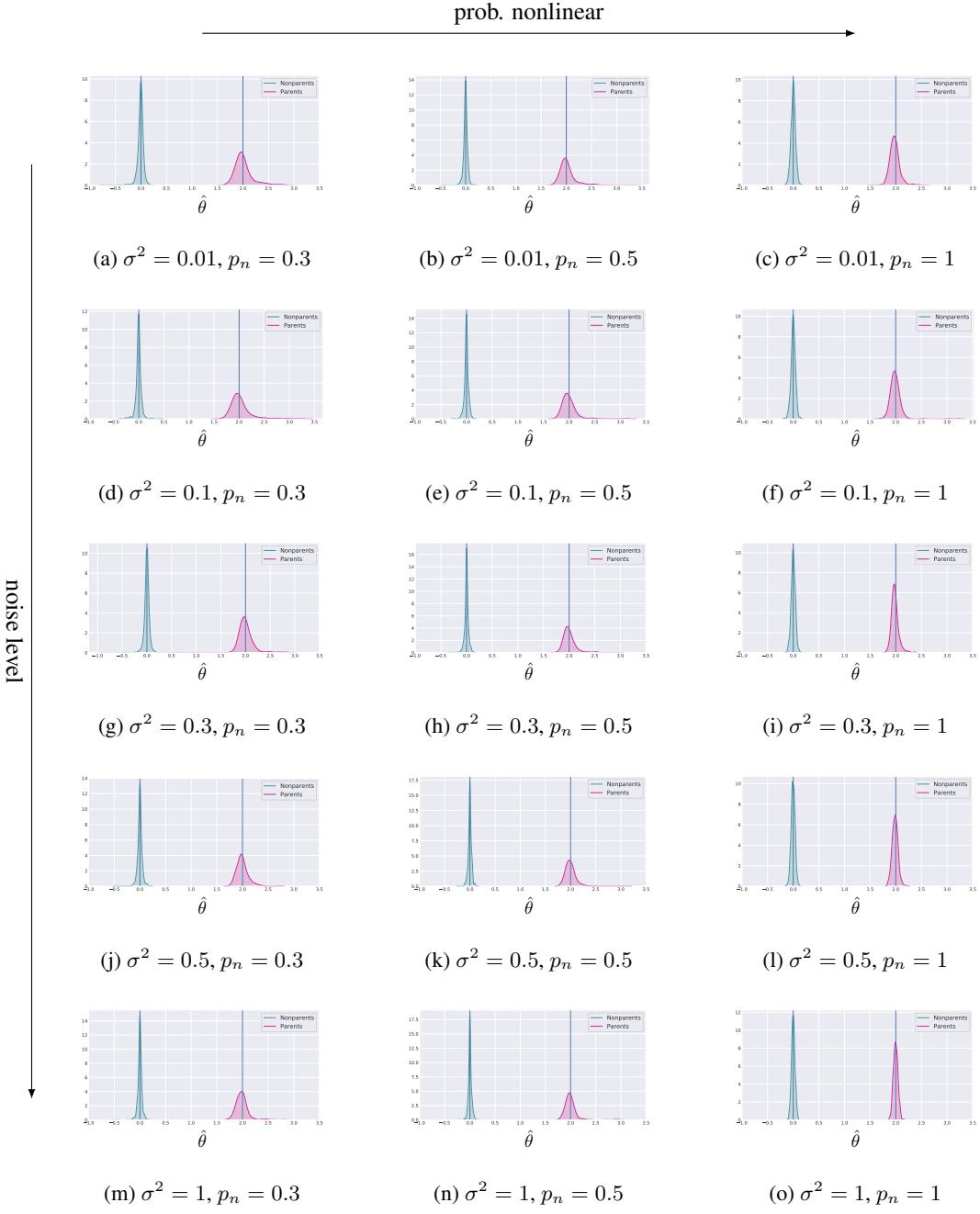


Figure F.18: 0.3 sparsity, 10 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

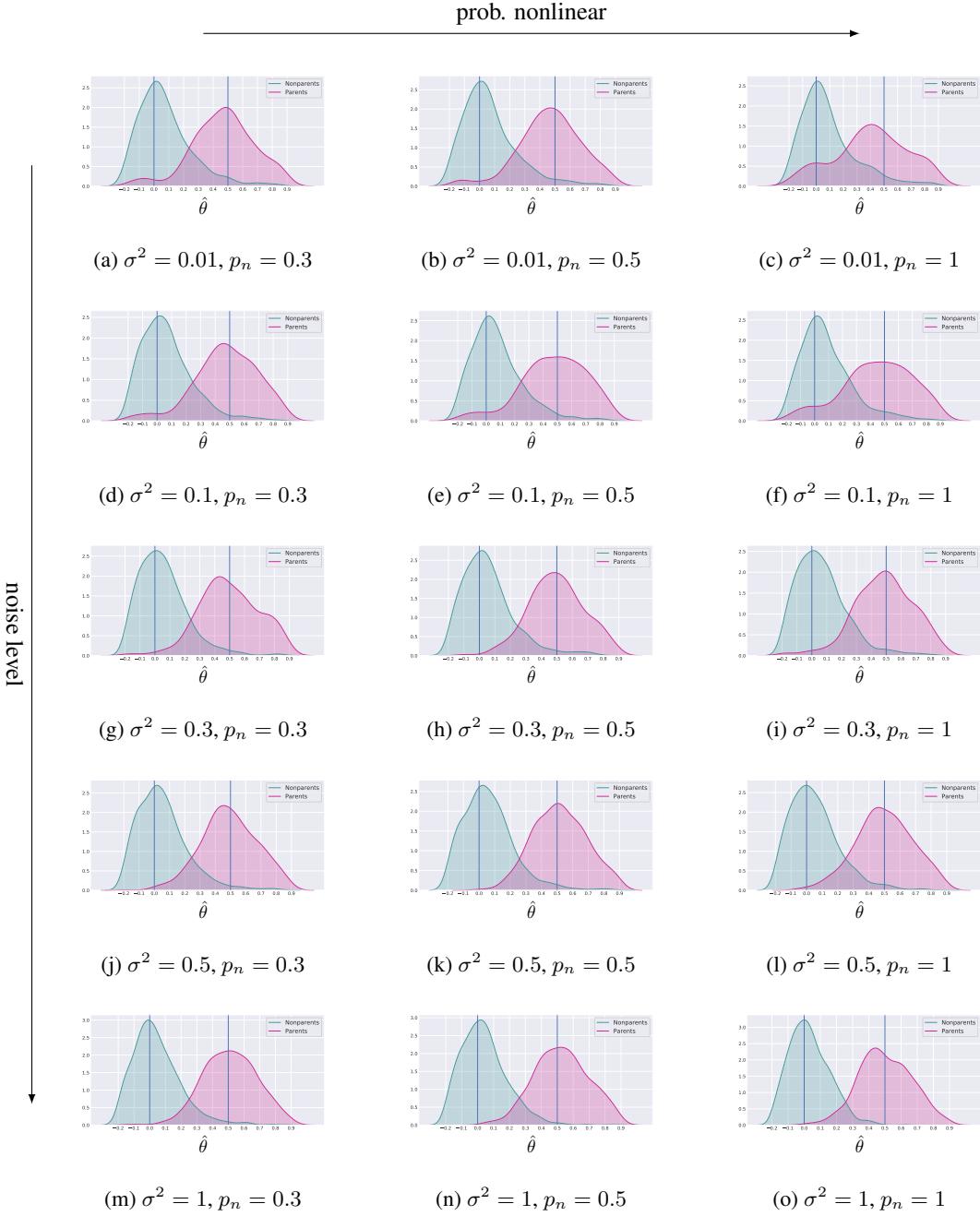


Figure F.19: 0.3 sparsity, 20 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

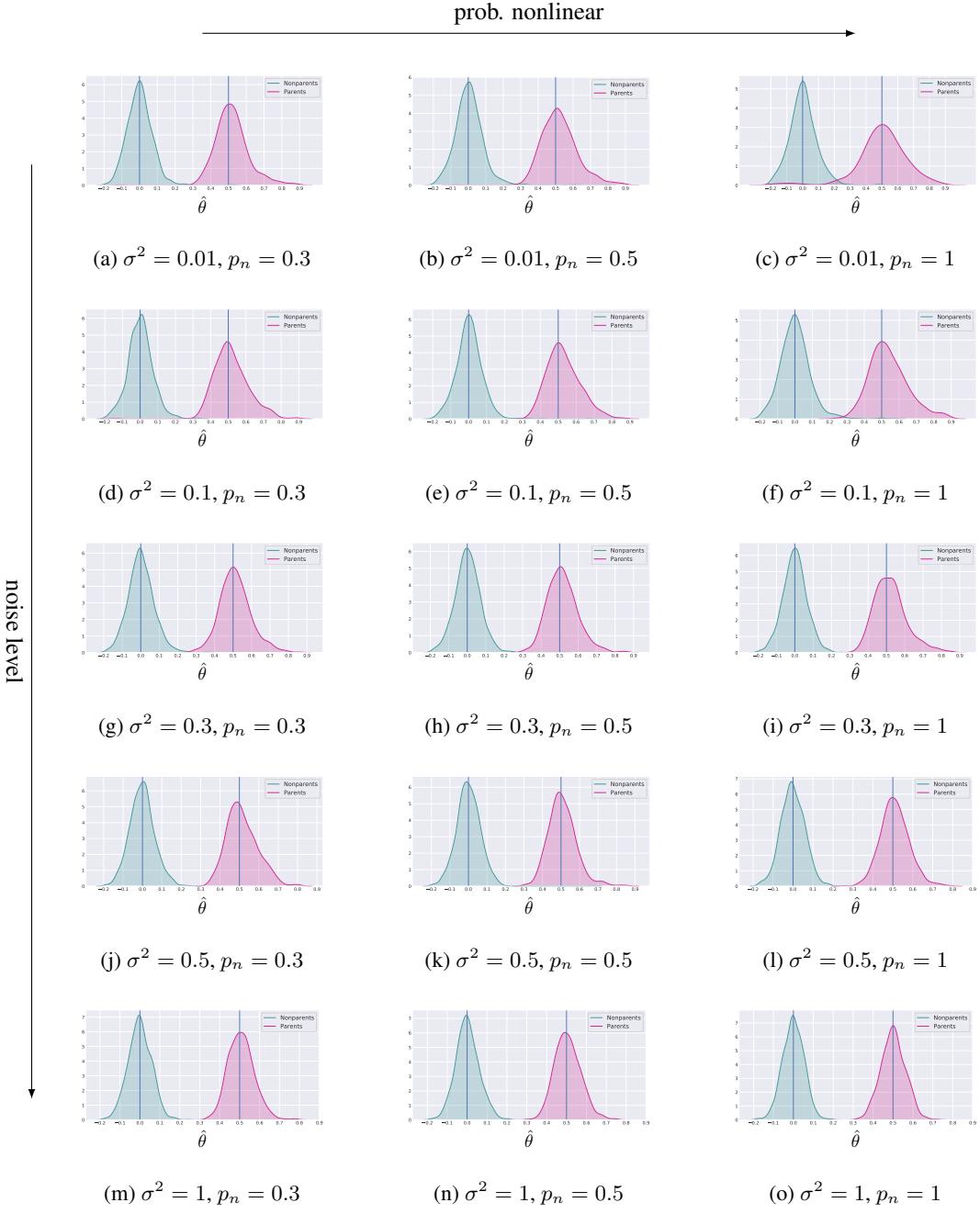


Figure F.20: 0.3 sparsity, 20 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

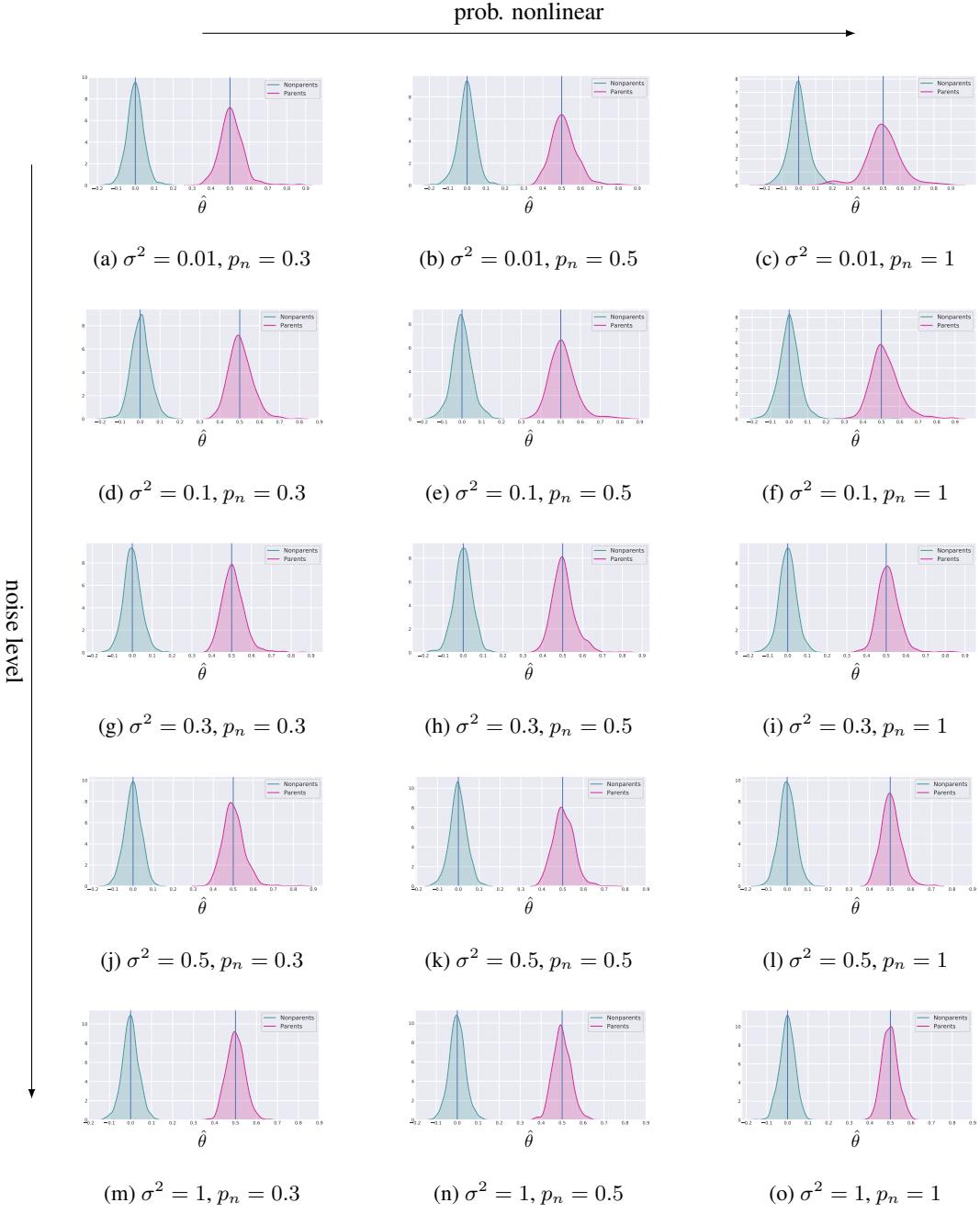


Figure F.21: 0.3 sparsity, 20 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

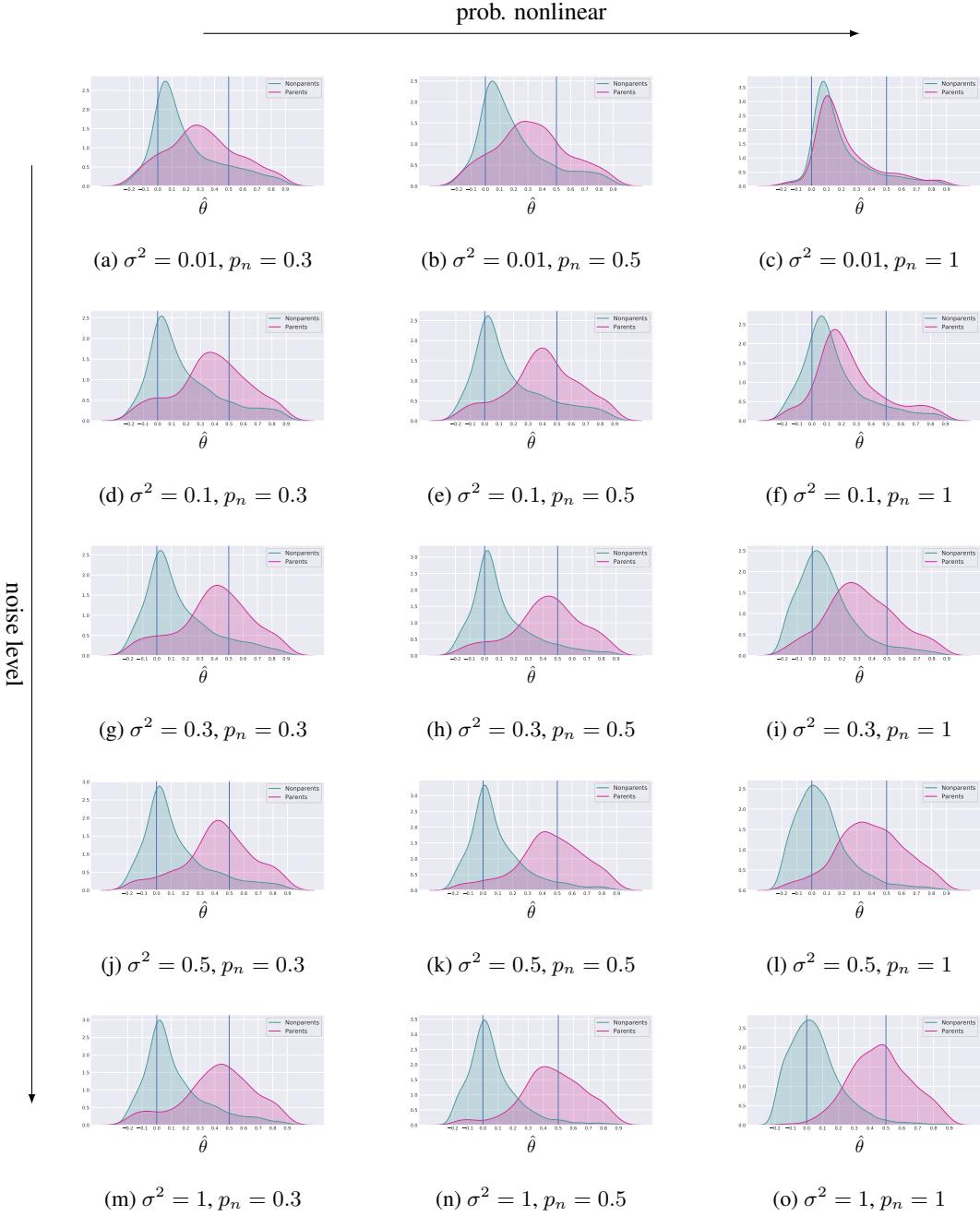


Figure F.22: 0.3 sparsity, 50 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

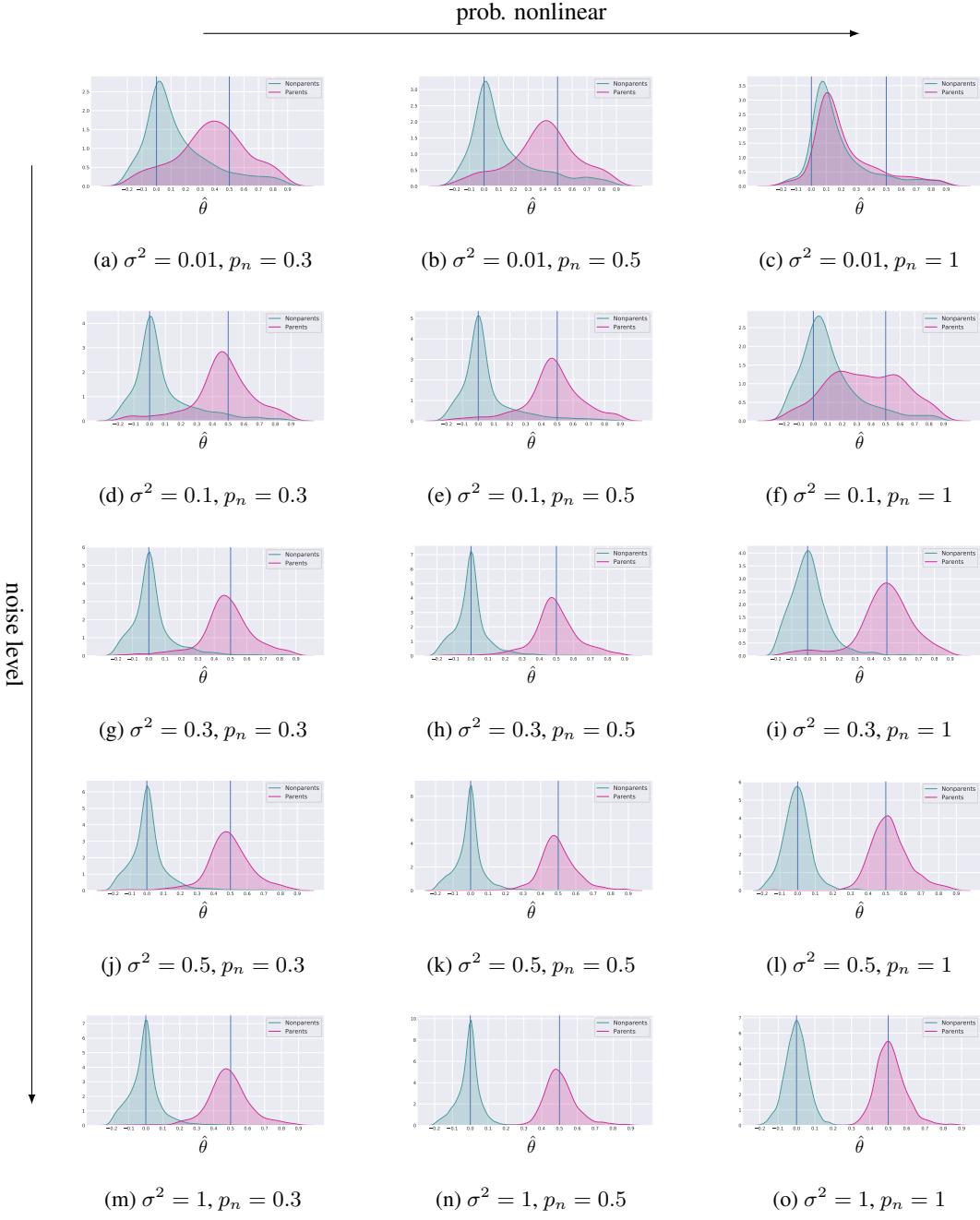


Figure F.23: 0.3 sparsity, 50 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

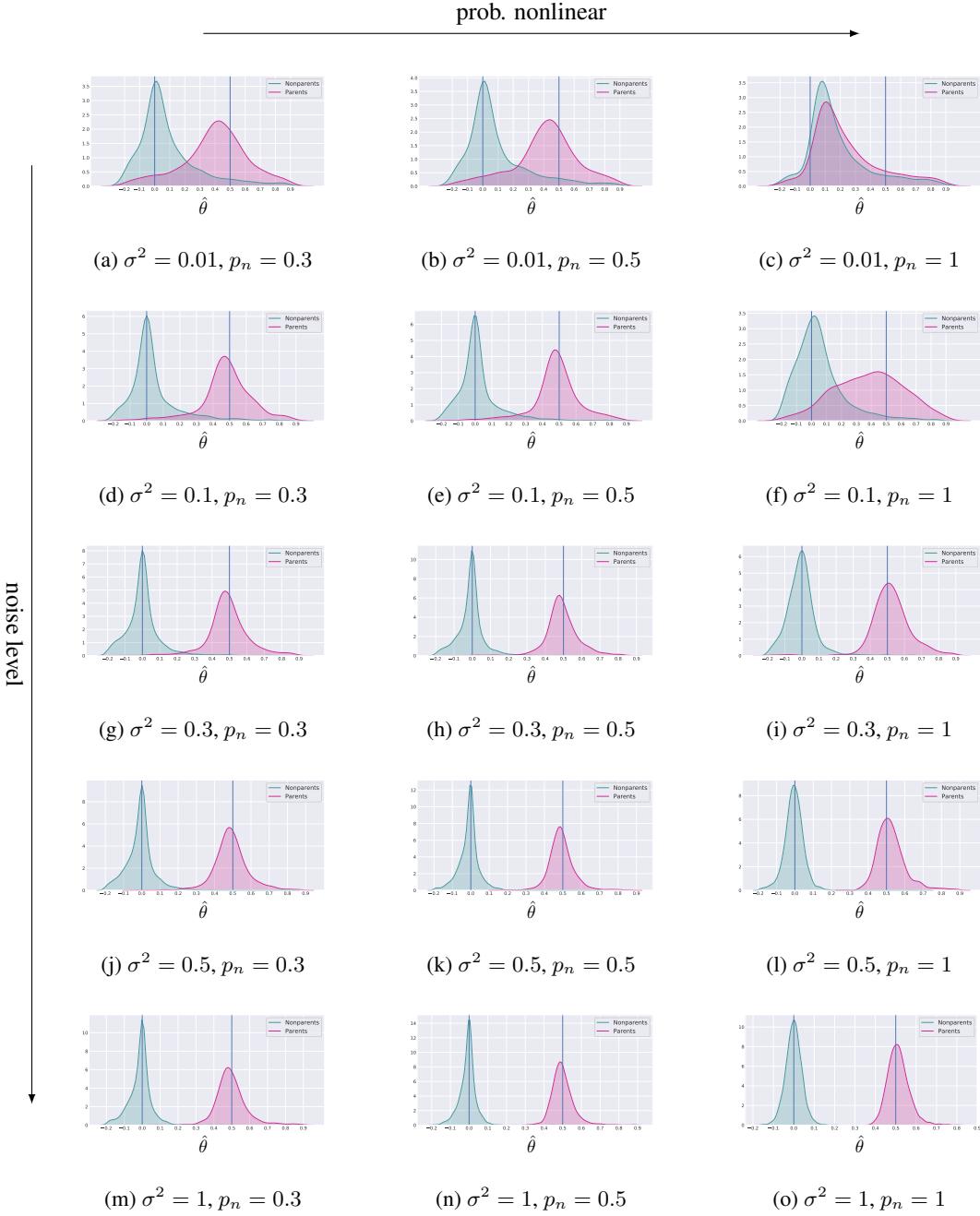


Figure F.24: 0.3 sparsity, 50 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

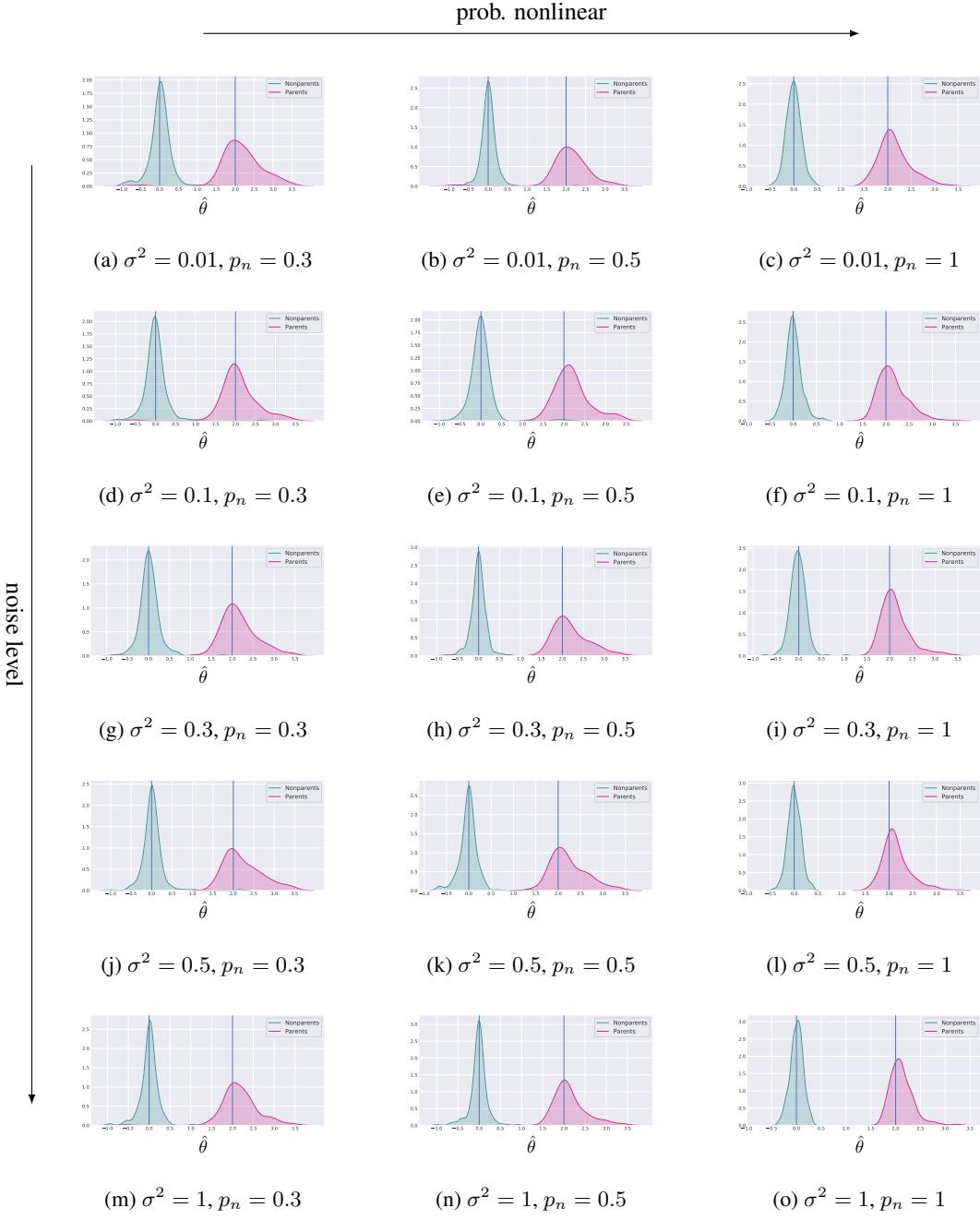


Figure F.25: 0.5 sparsity, 5 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

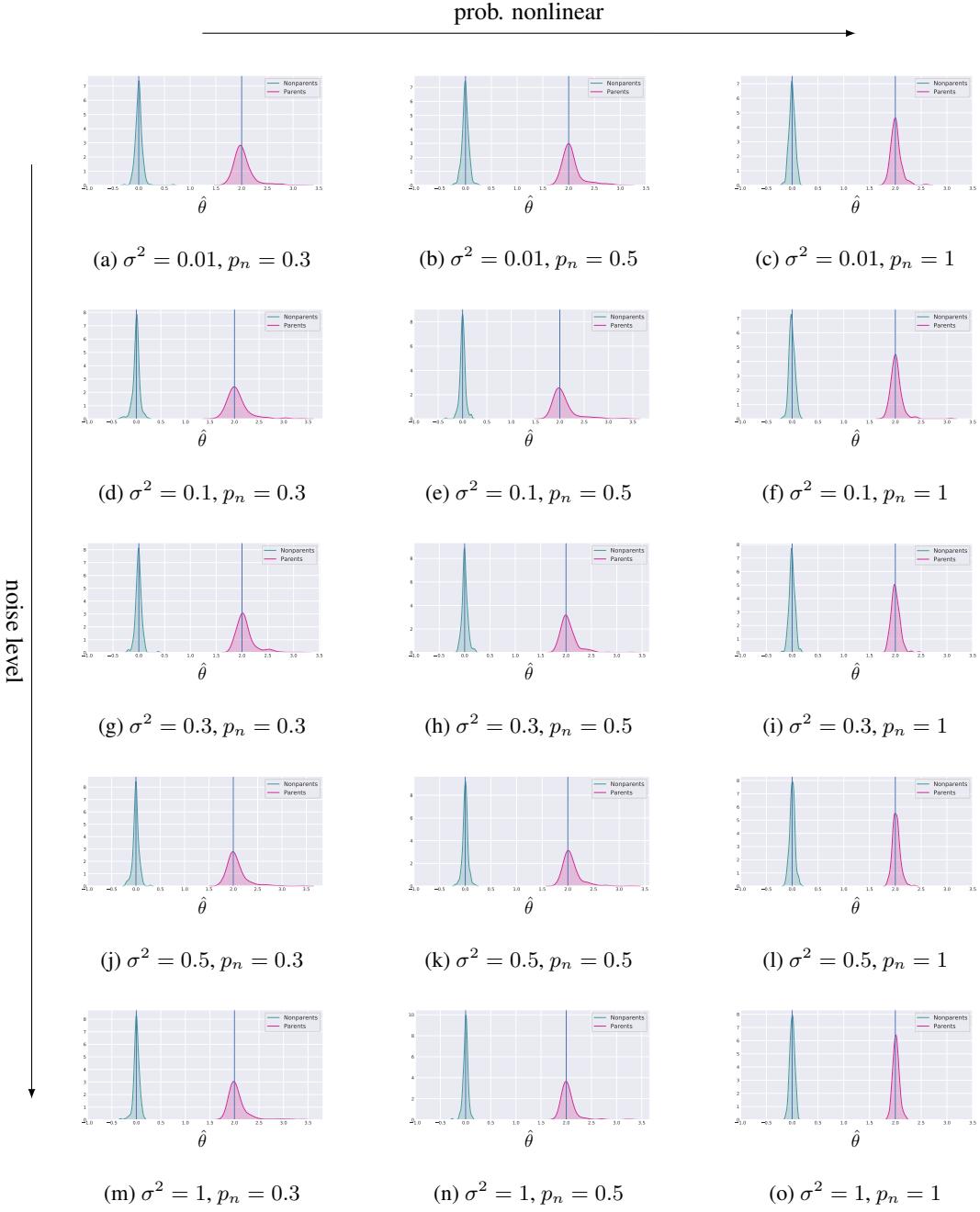


Figure F.26: 0.5 sparsity, 5 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

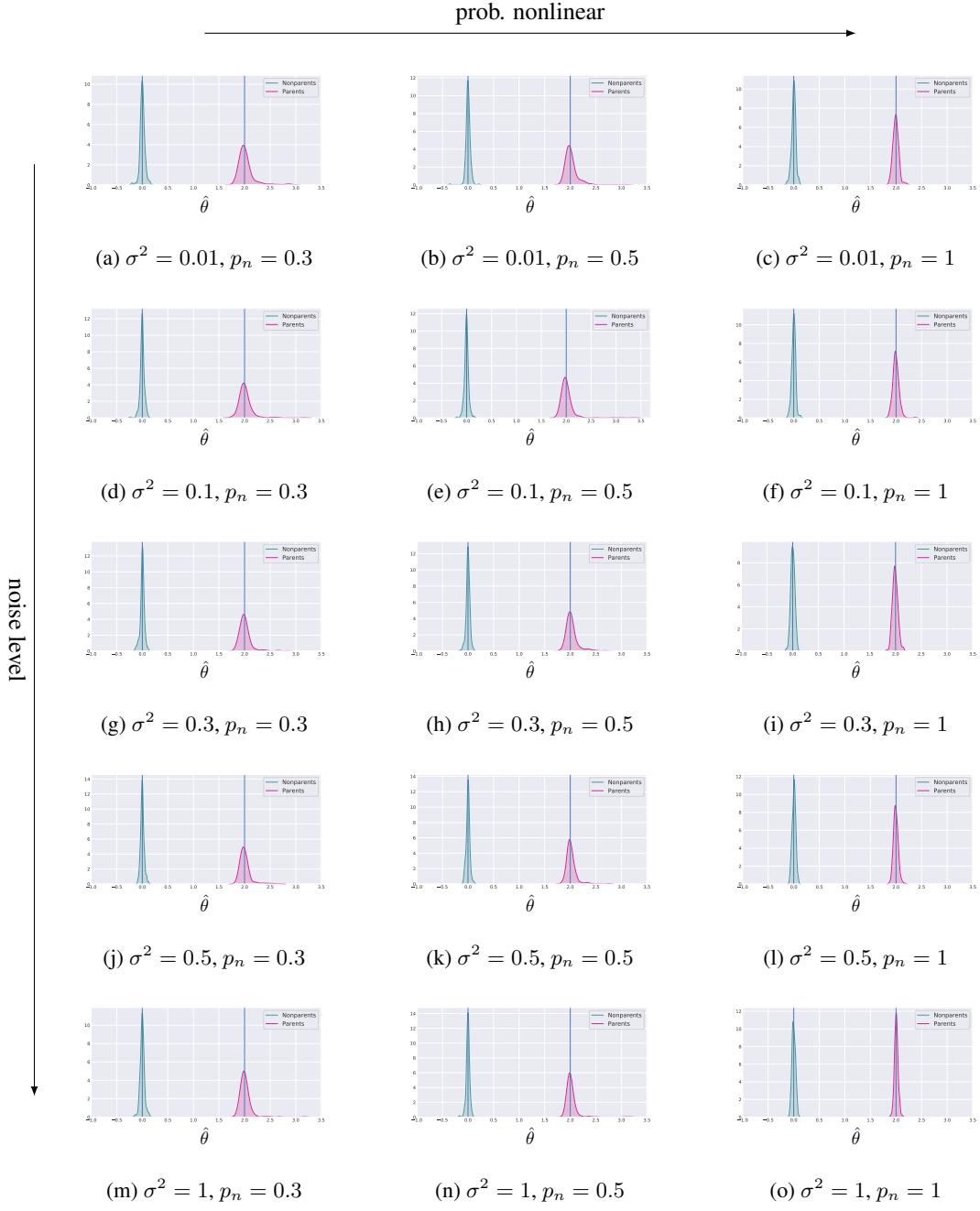


Figure F.27: 0.5 sparsity, 5 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

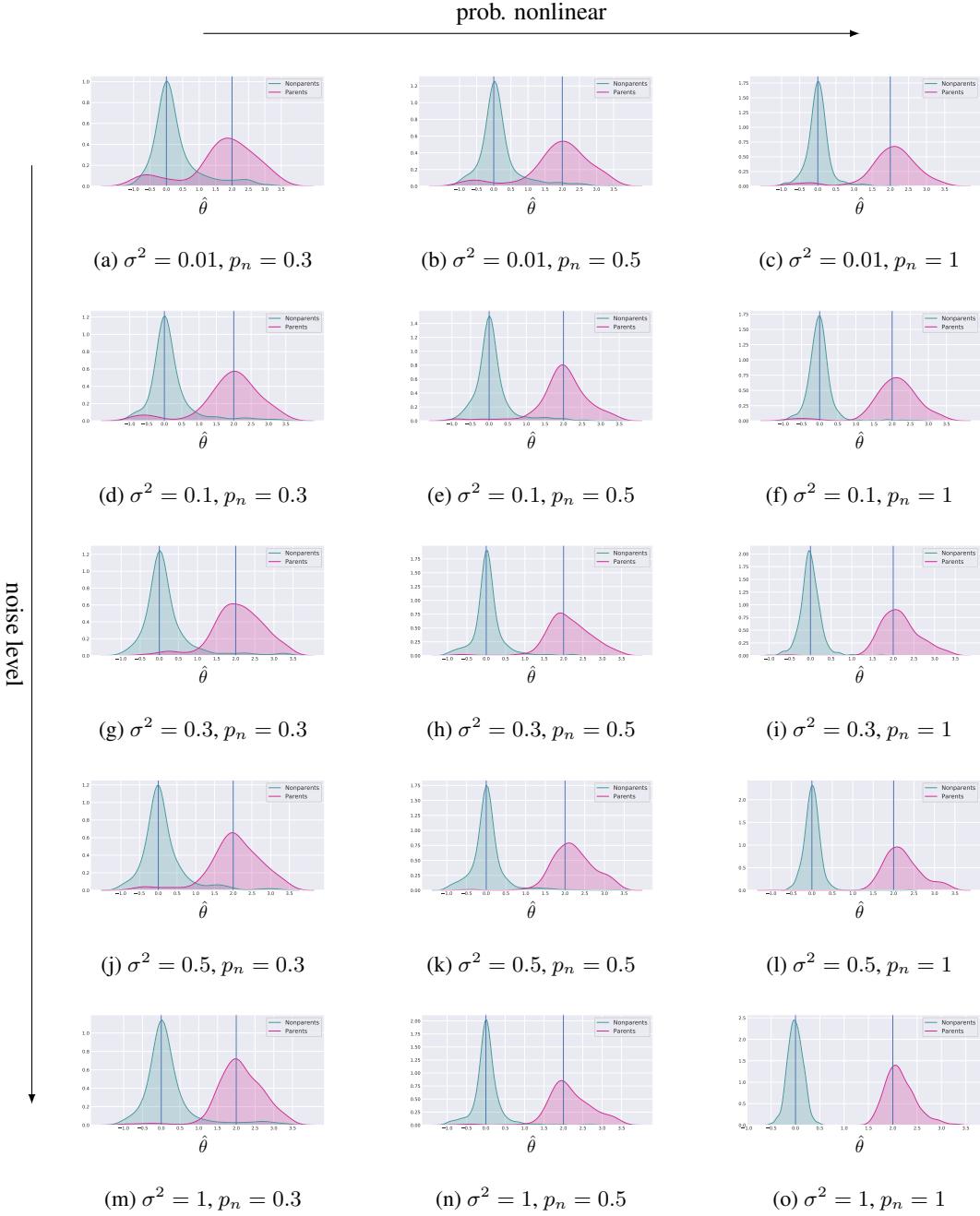


Figure F.28: 0.5 sparsity, 10 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

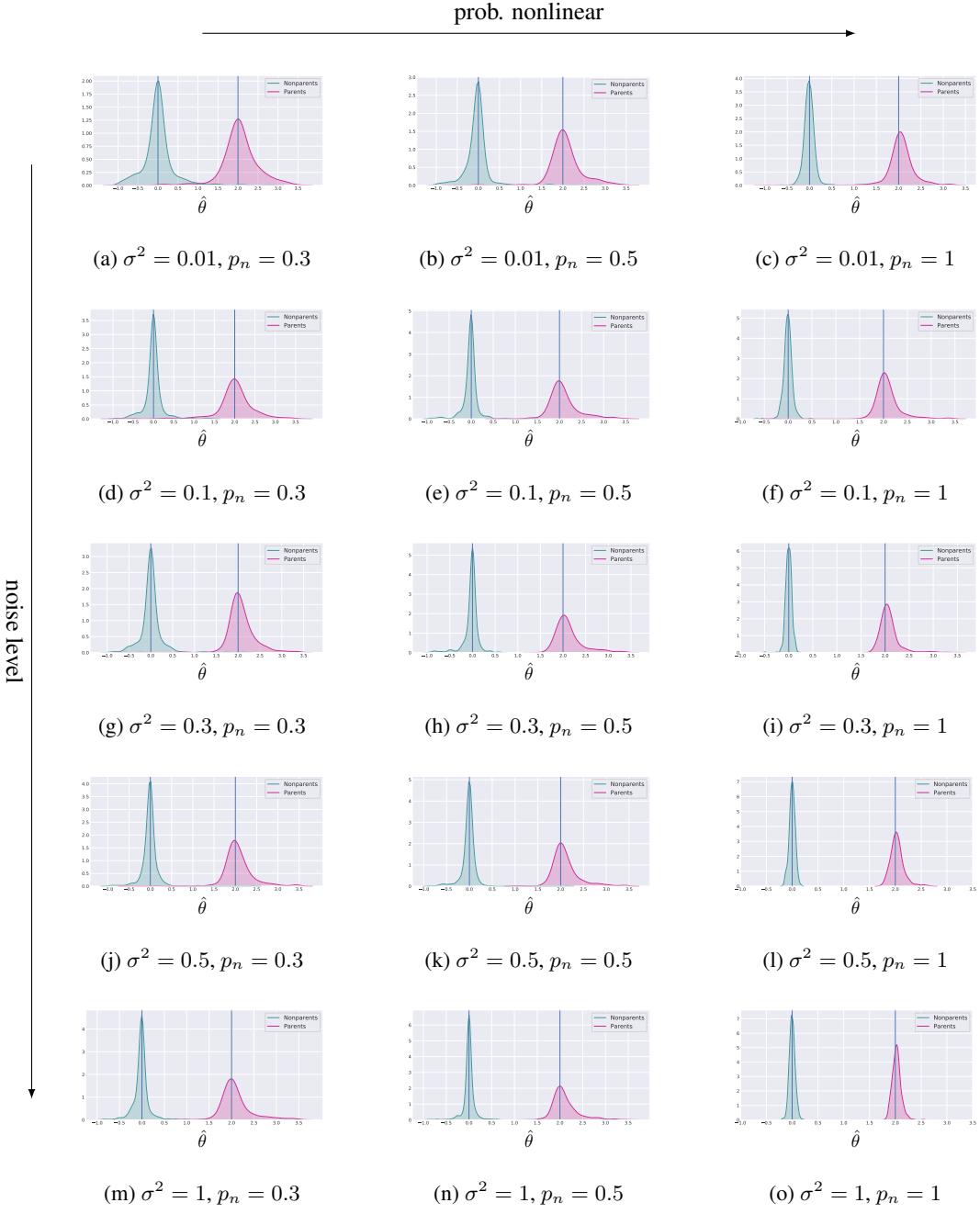


Figure F.29: 0.5 sparsity, 10 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

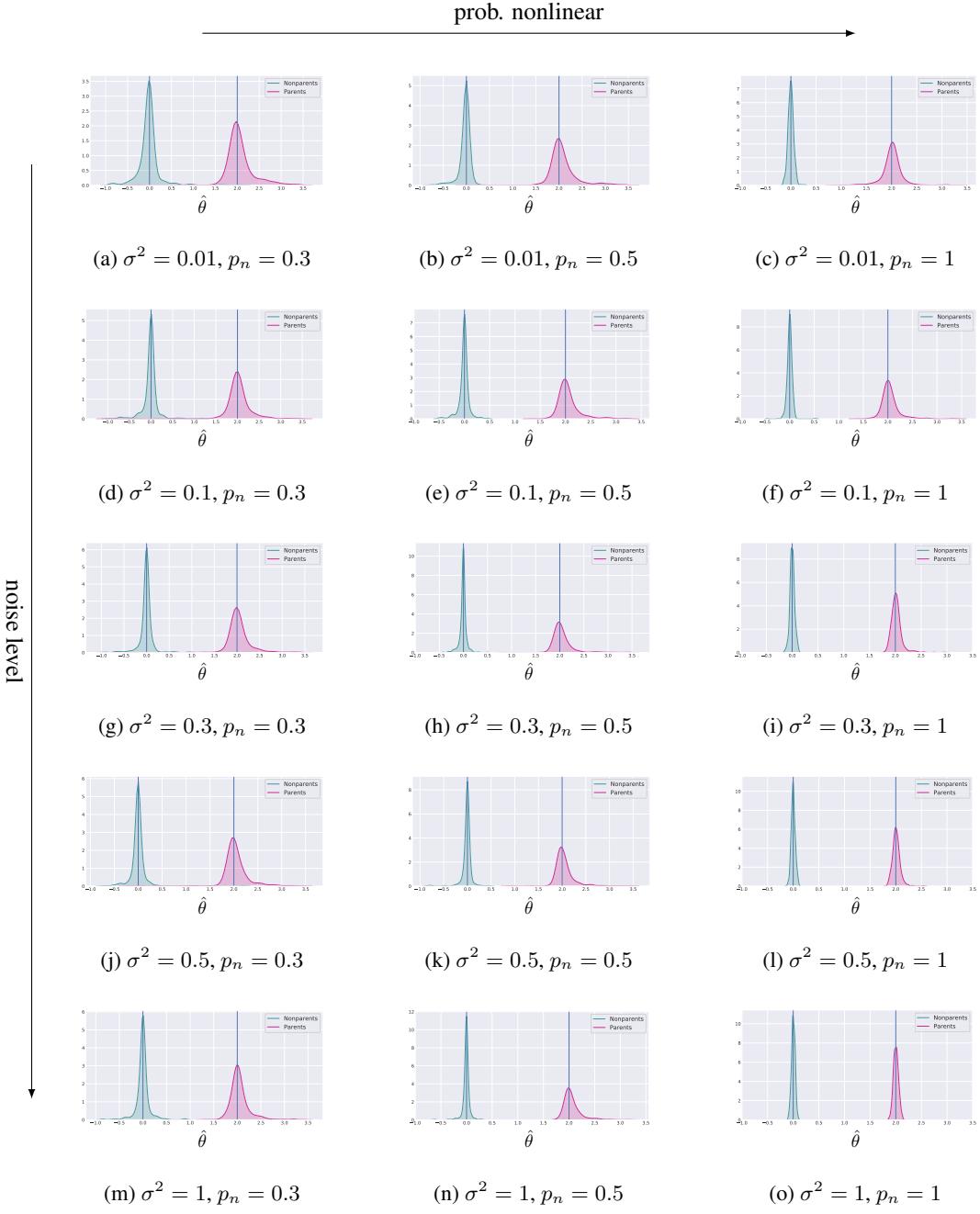


Figure F.30: 0.5 sparsity, 10 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

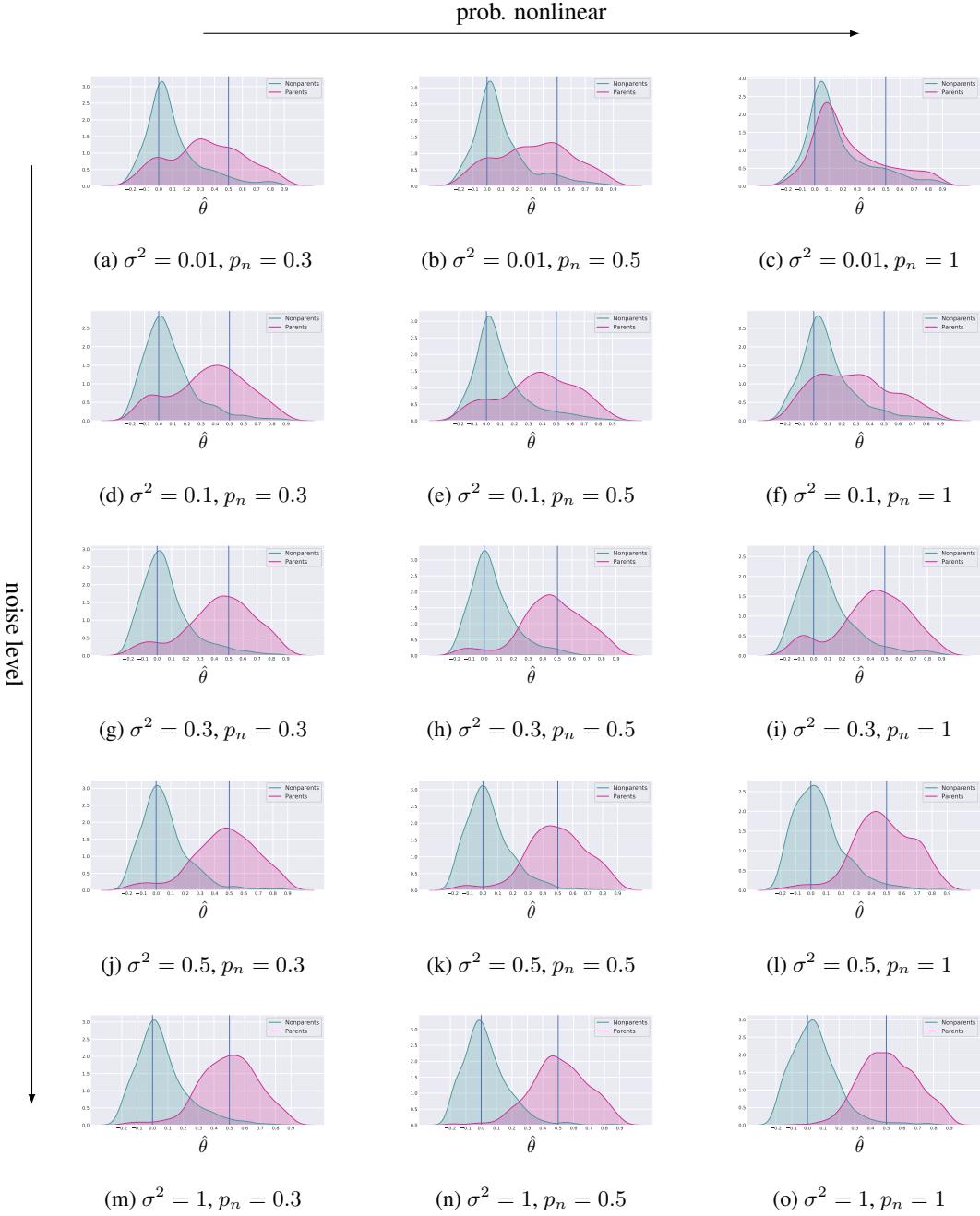


Figure F.31: 0.5 sparsity, 20 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

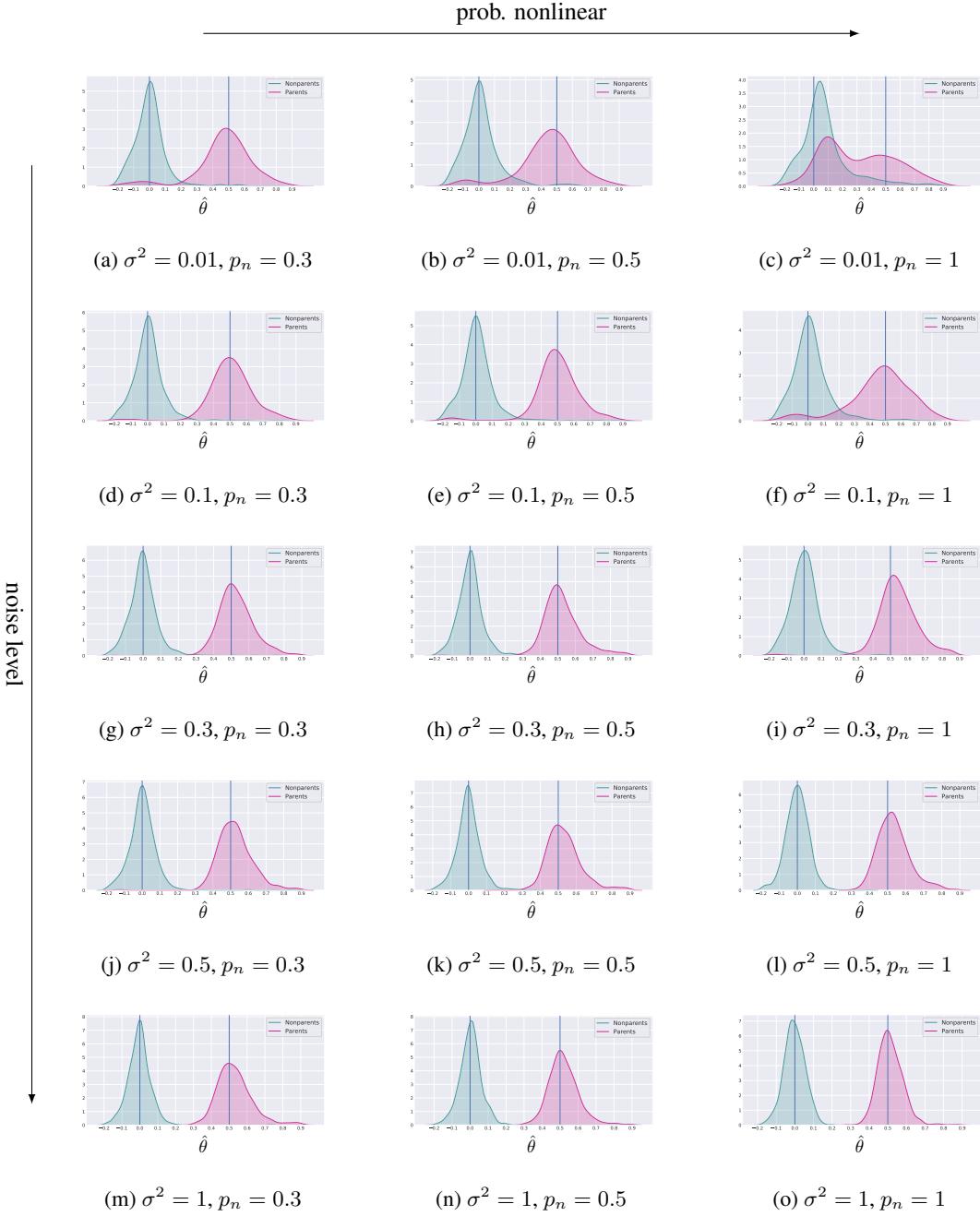


Figure F.32: 0.5 sparsity, 20 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\hat{\theta}$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

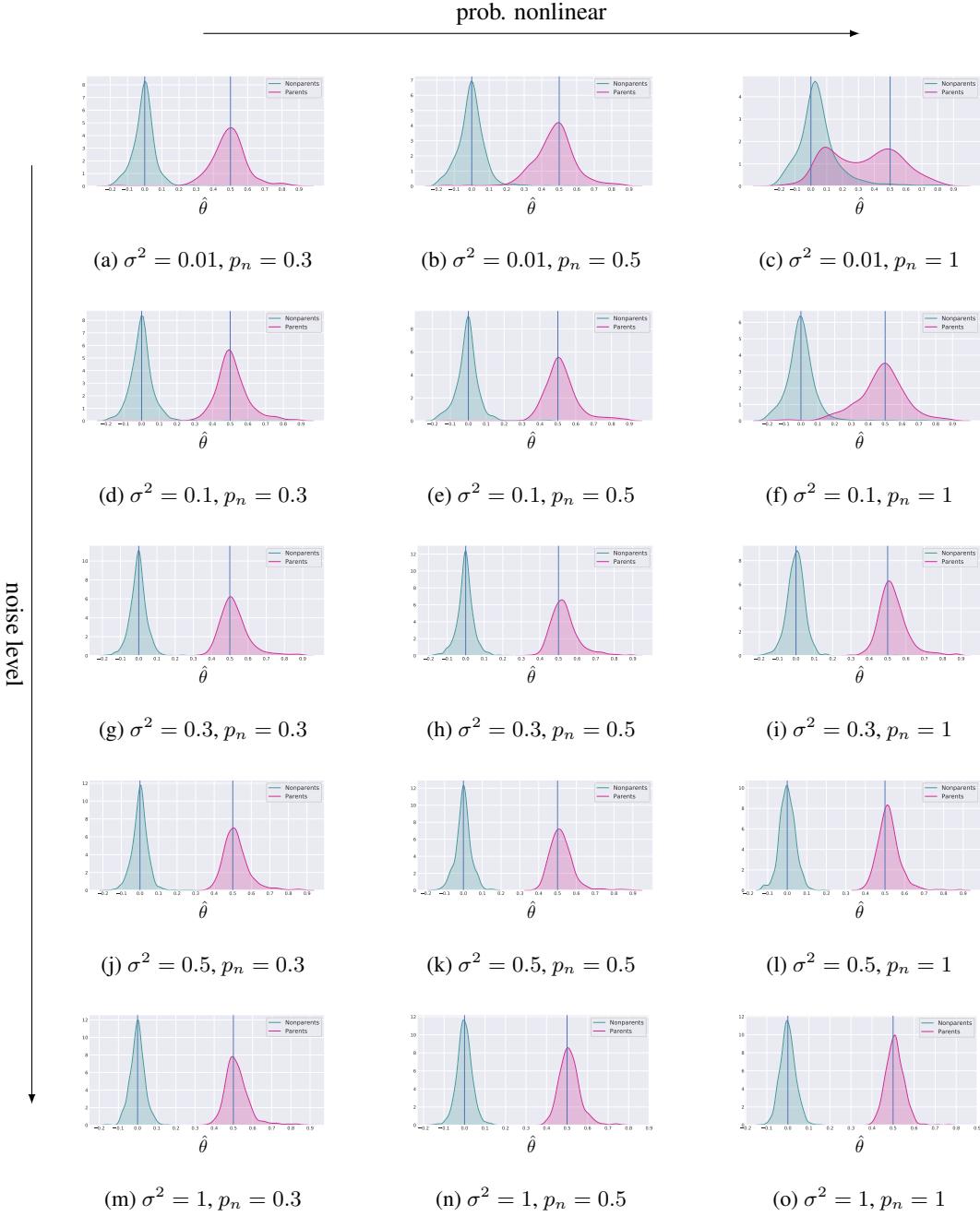


Figure F.33: 0.5 sparsity, 20 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

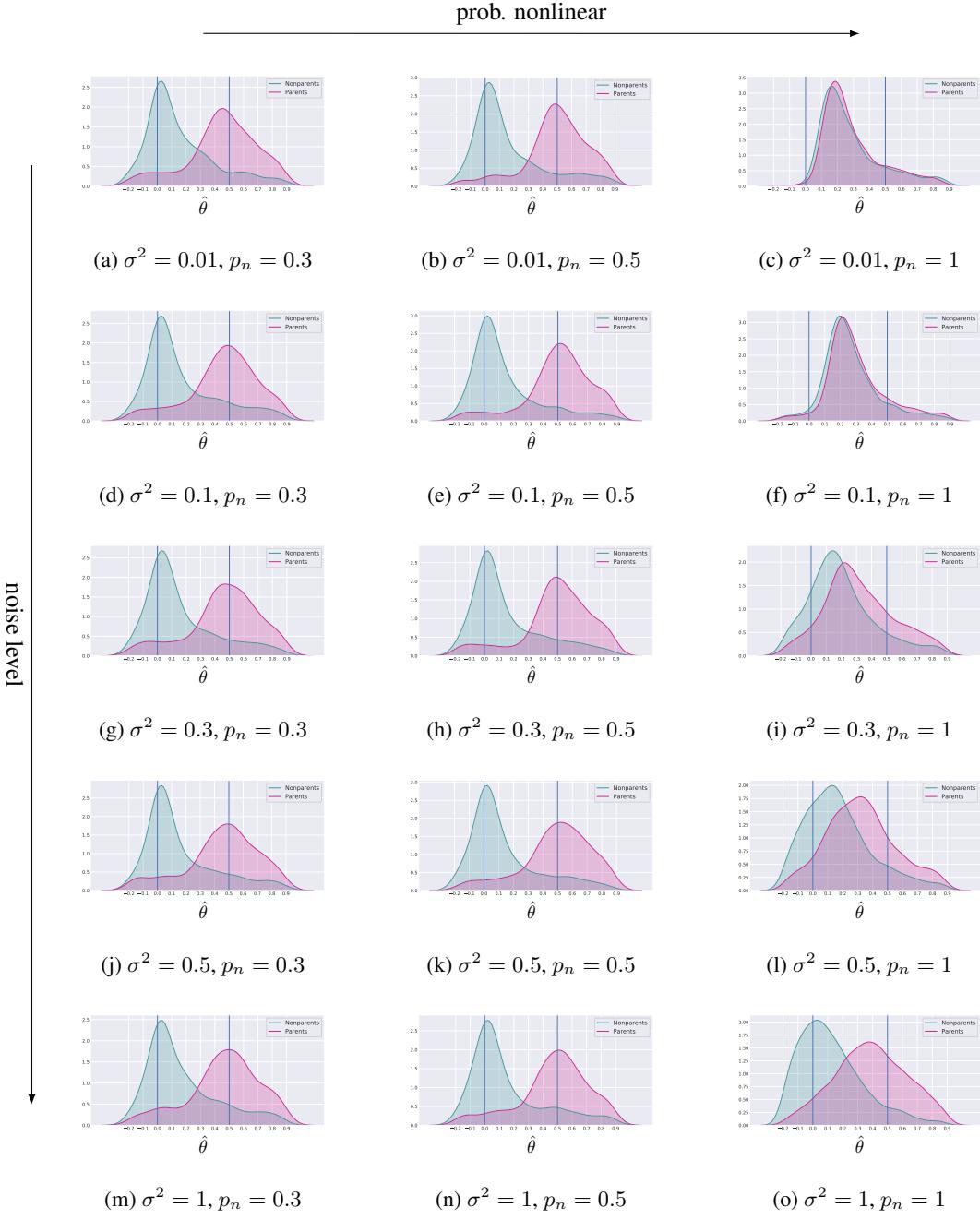


Figure F.34: 0.5 sparsity, 50 nodes, 100 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the paremater estimation works reliably.

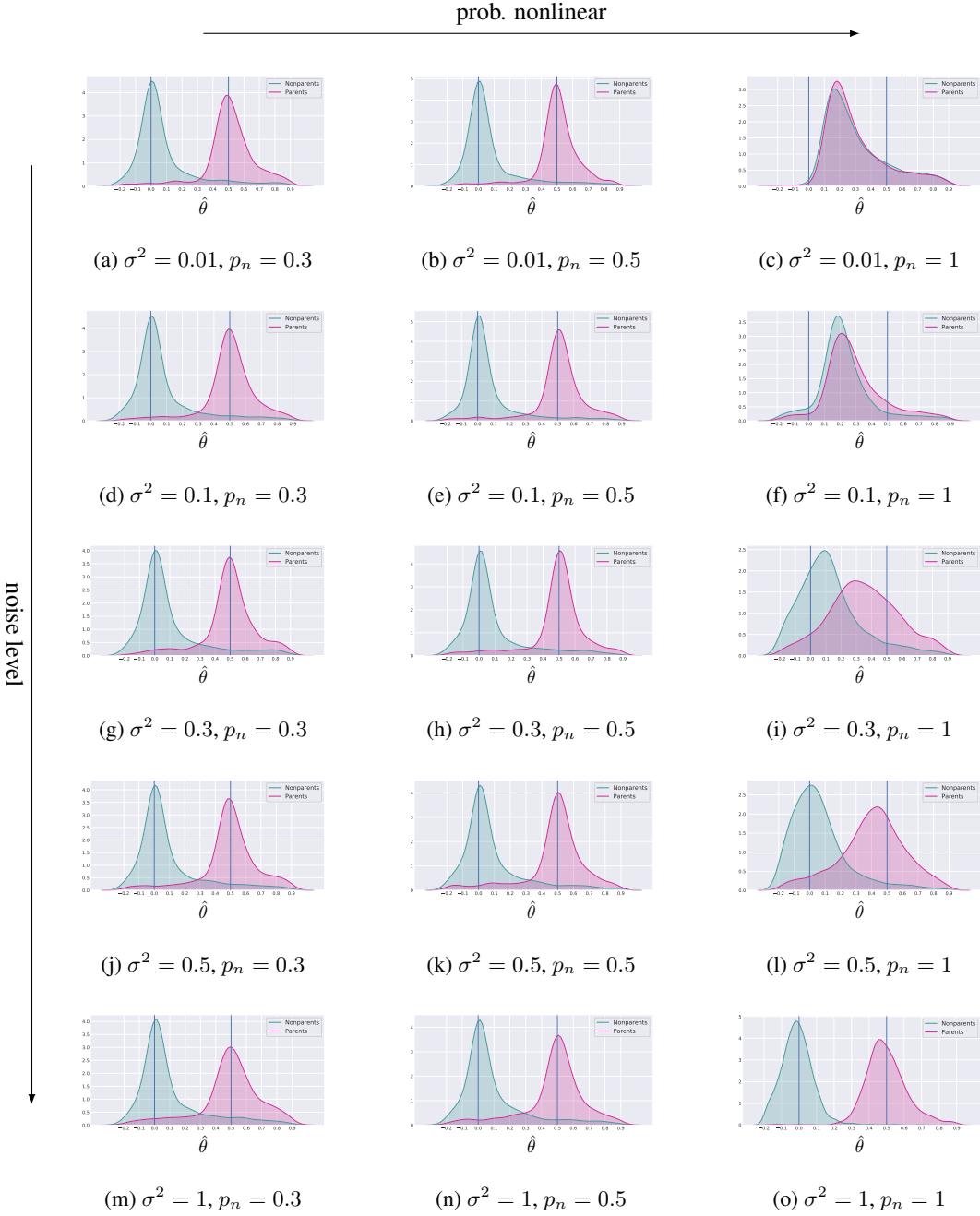


Figure F.35: 0.5 sparsity, 50 nodes, 500 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

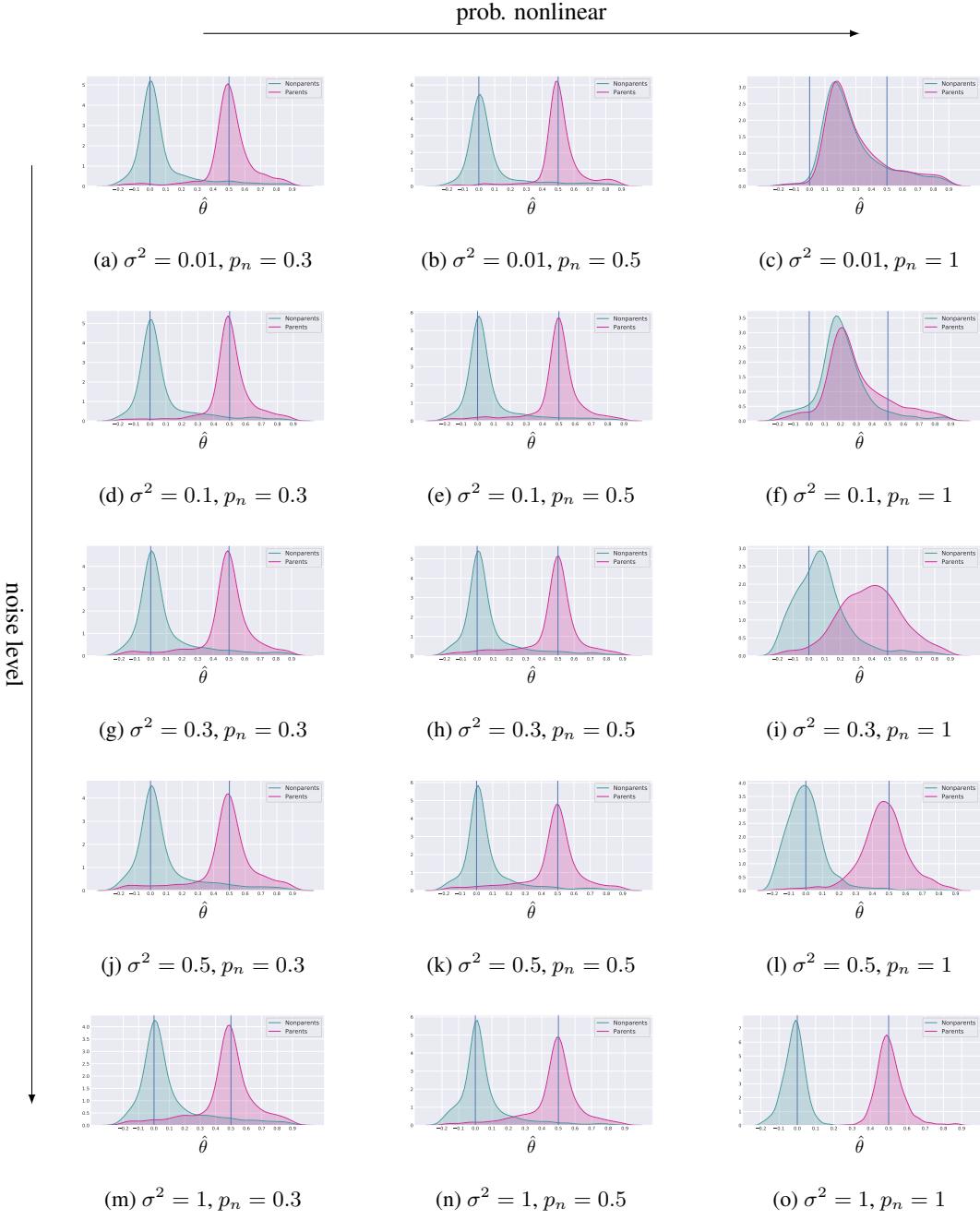


Figure F.36: 0.5 sparsity, 50 nodes, 1000 observations, 100 simulations. Distribution of the estimated  $\theta$  values for the true and false causal parents in 100 simulations. The vertical lines indicate the ground truth values for the causal parents linear coefficients. In general we observe that in all settings with enough observations the parameter estimation works reliably.

## G Supplementary Figures for Performance in Inferring Direct Causes, Accuracy as Metric

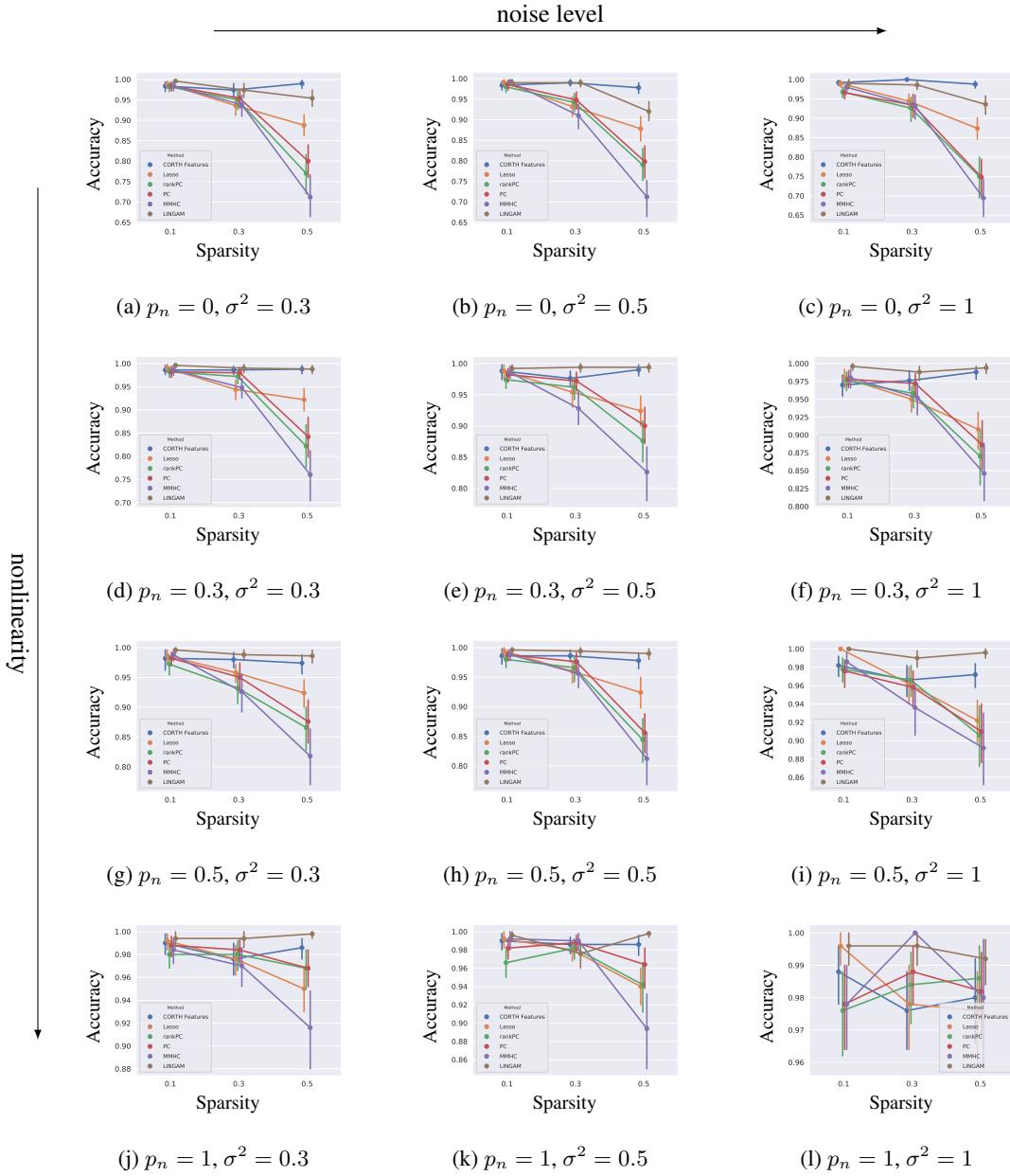


Figure G.1: 5 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

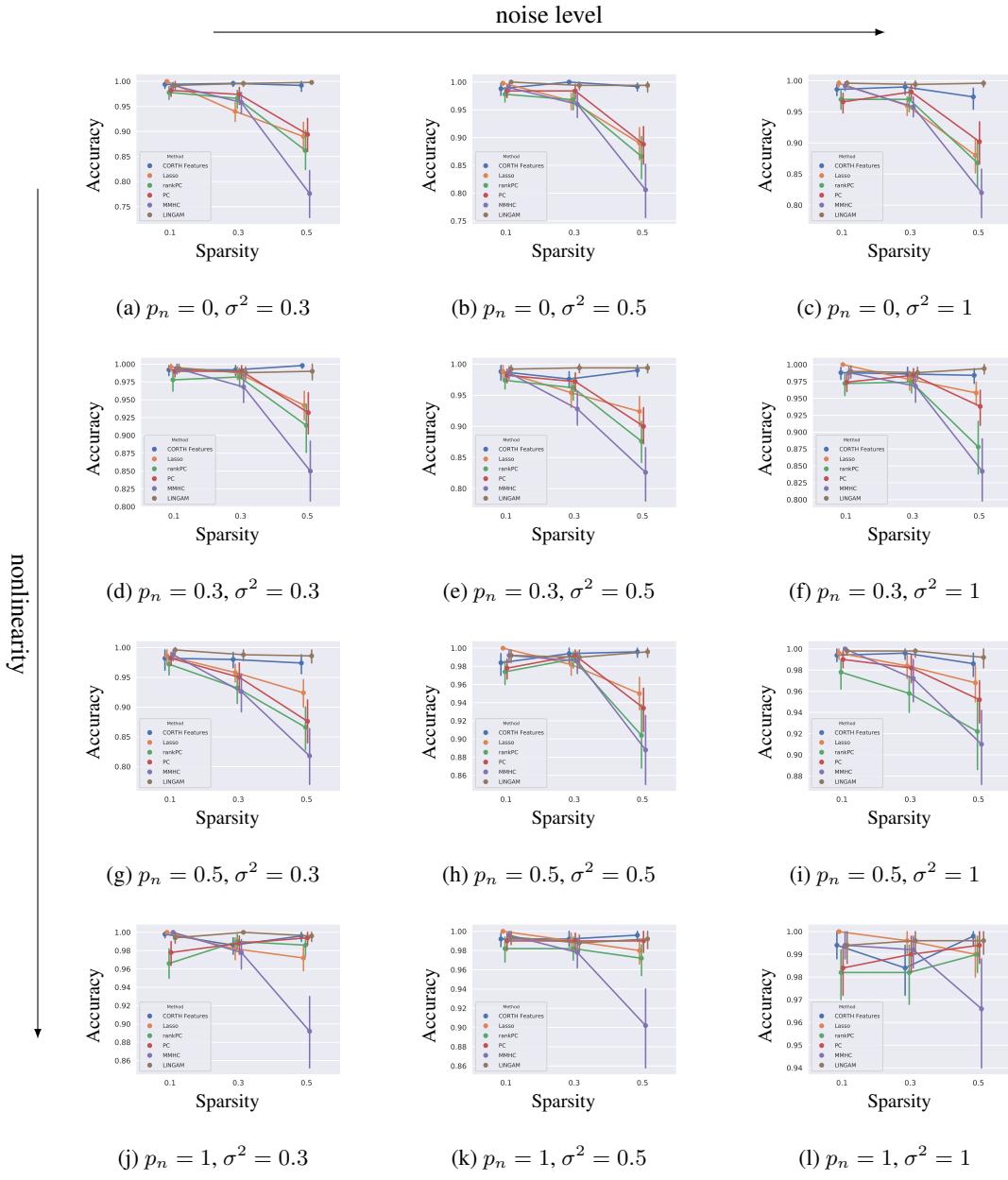


Figure G.2: 5 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

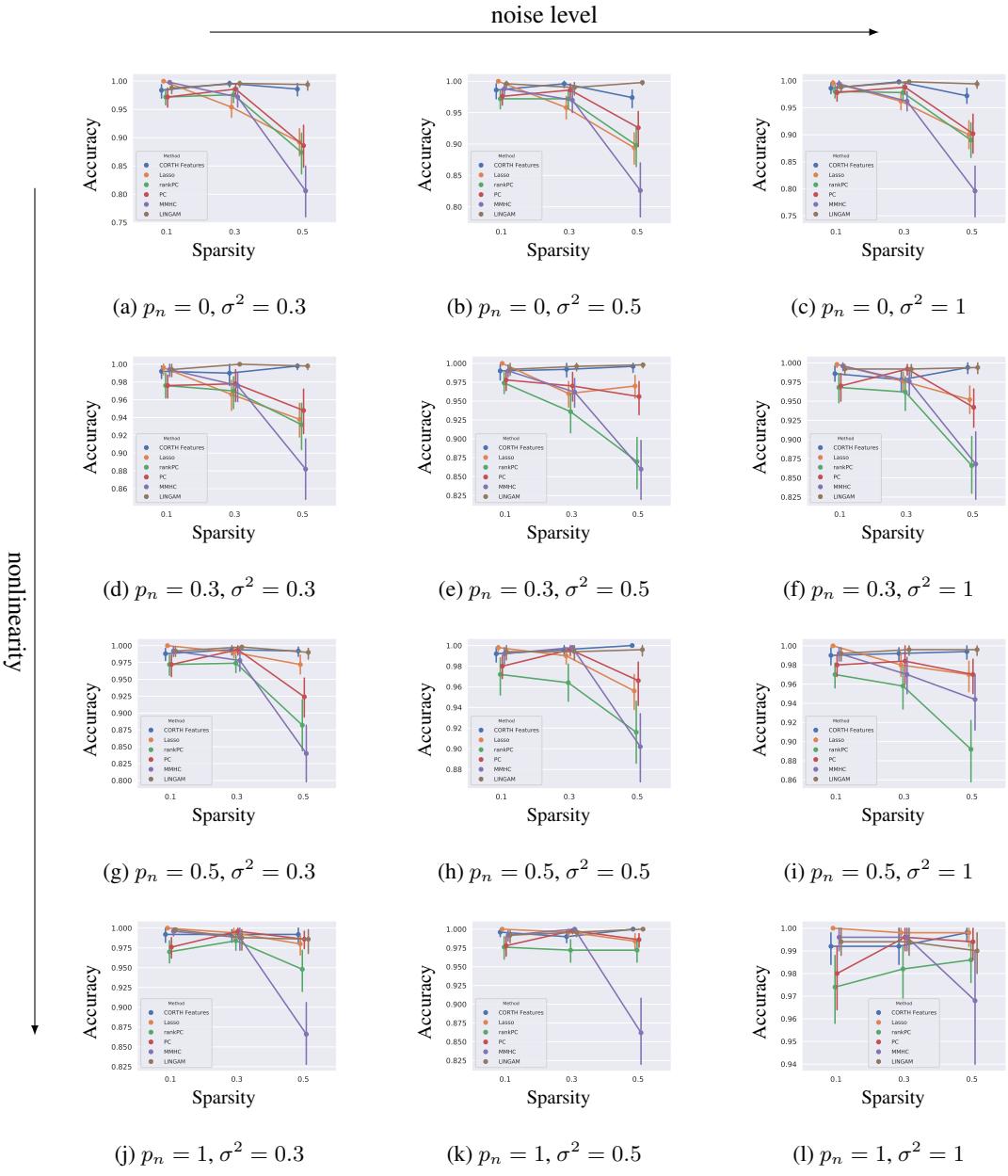


Figure G.3: 5 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

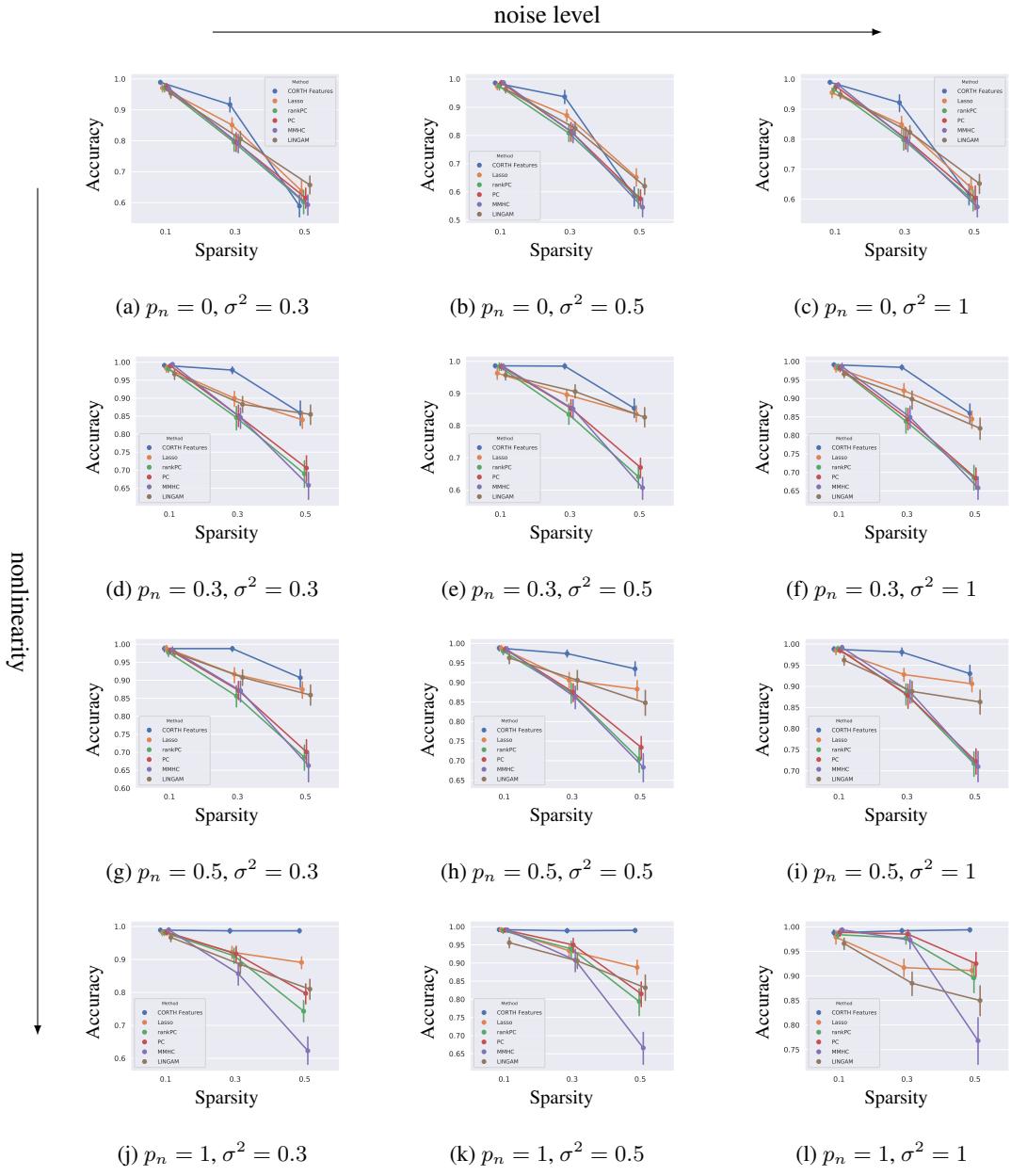


Figure G.4: 10 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

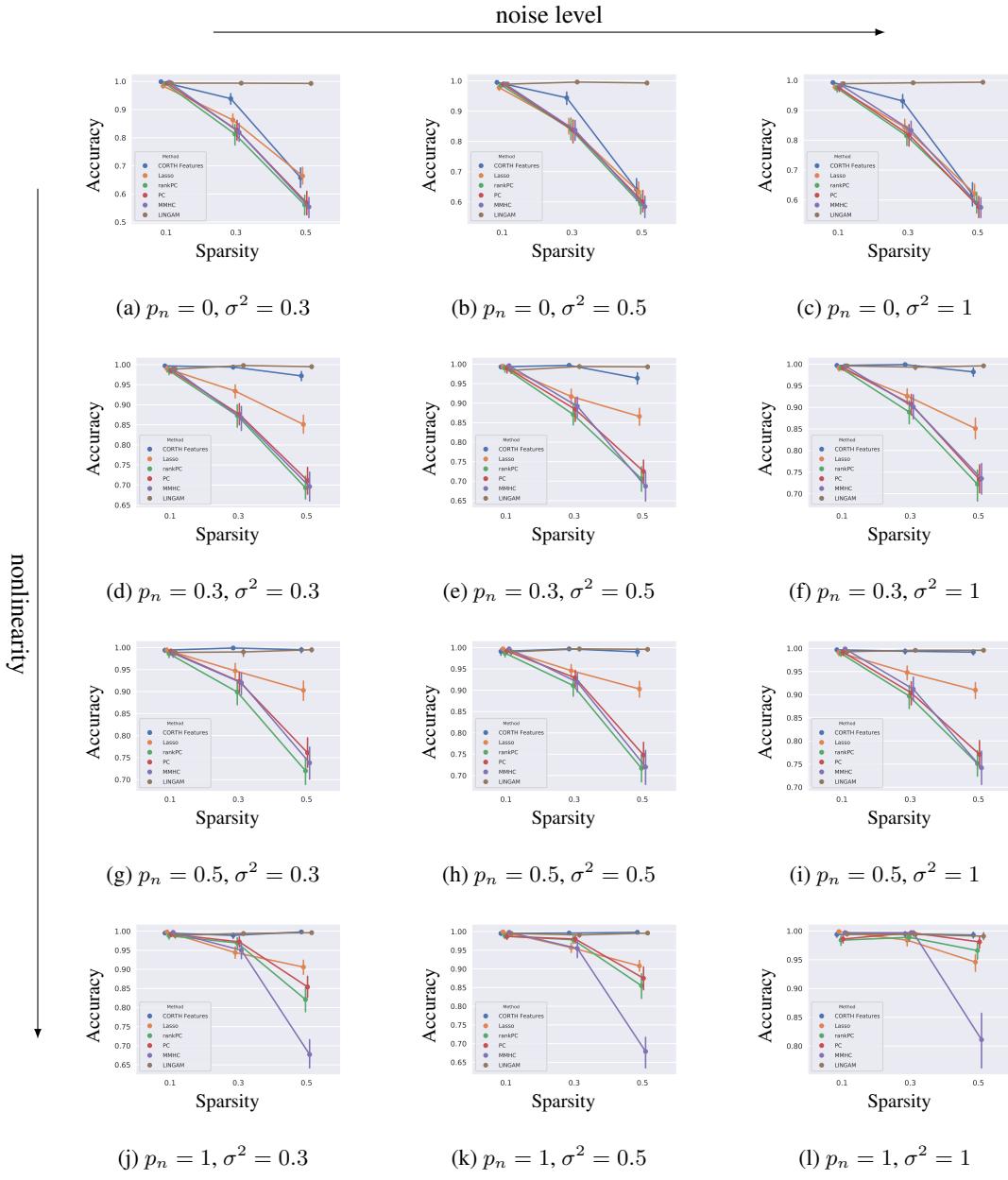


Figure G.5: 10 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

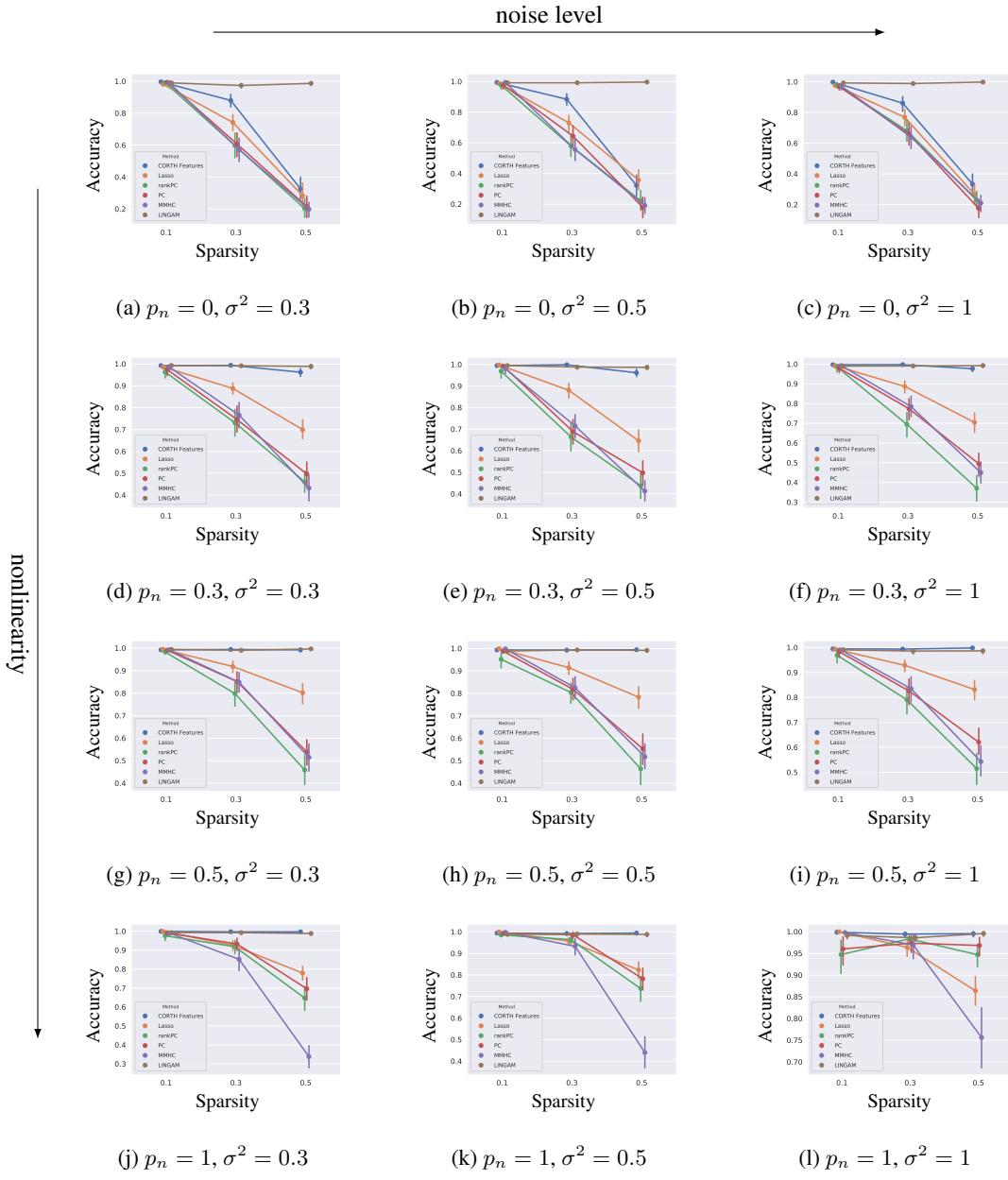


Figure G.6: 10 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

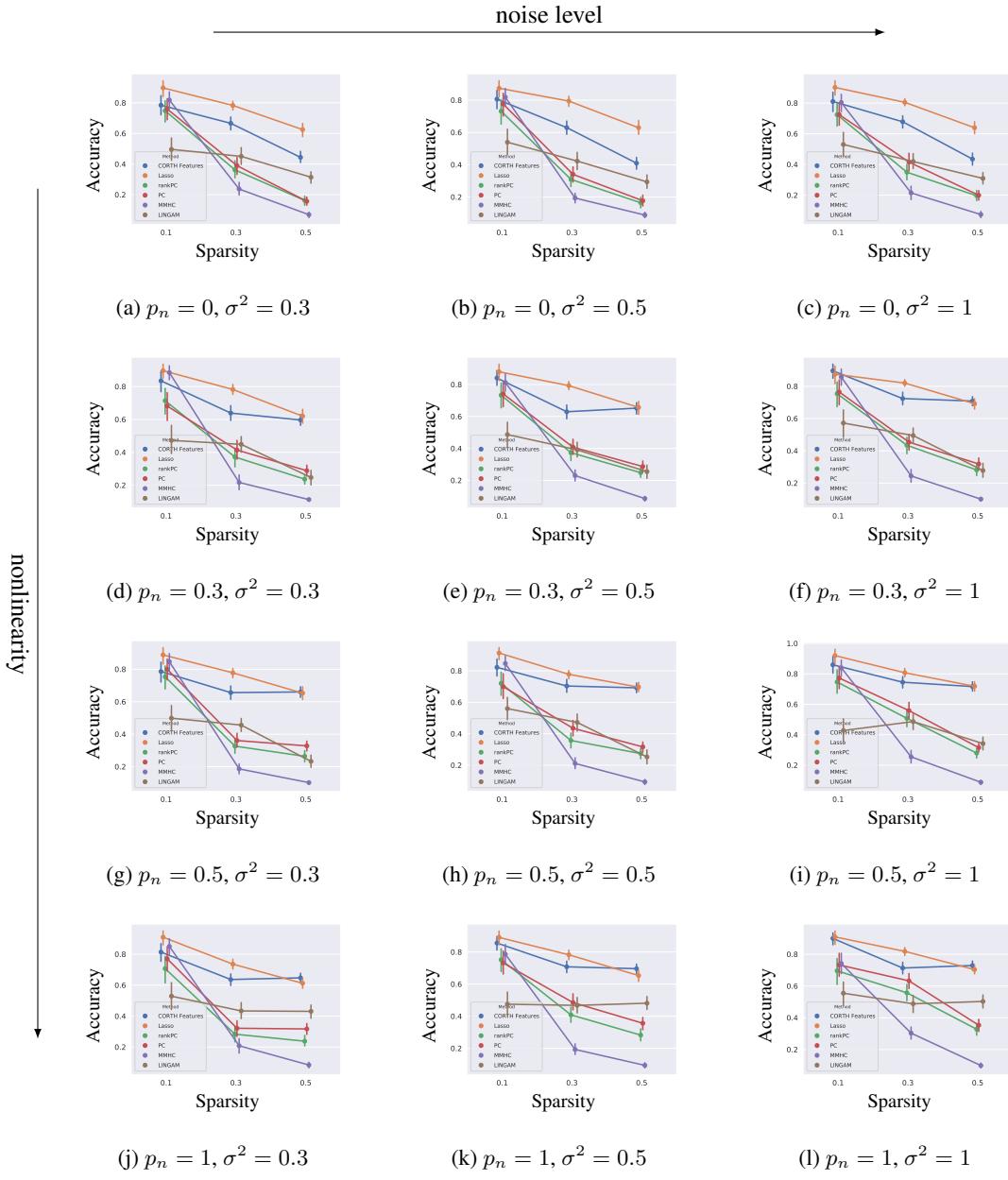


Figure G.7: 20 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

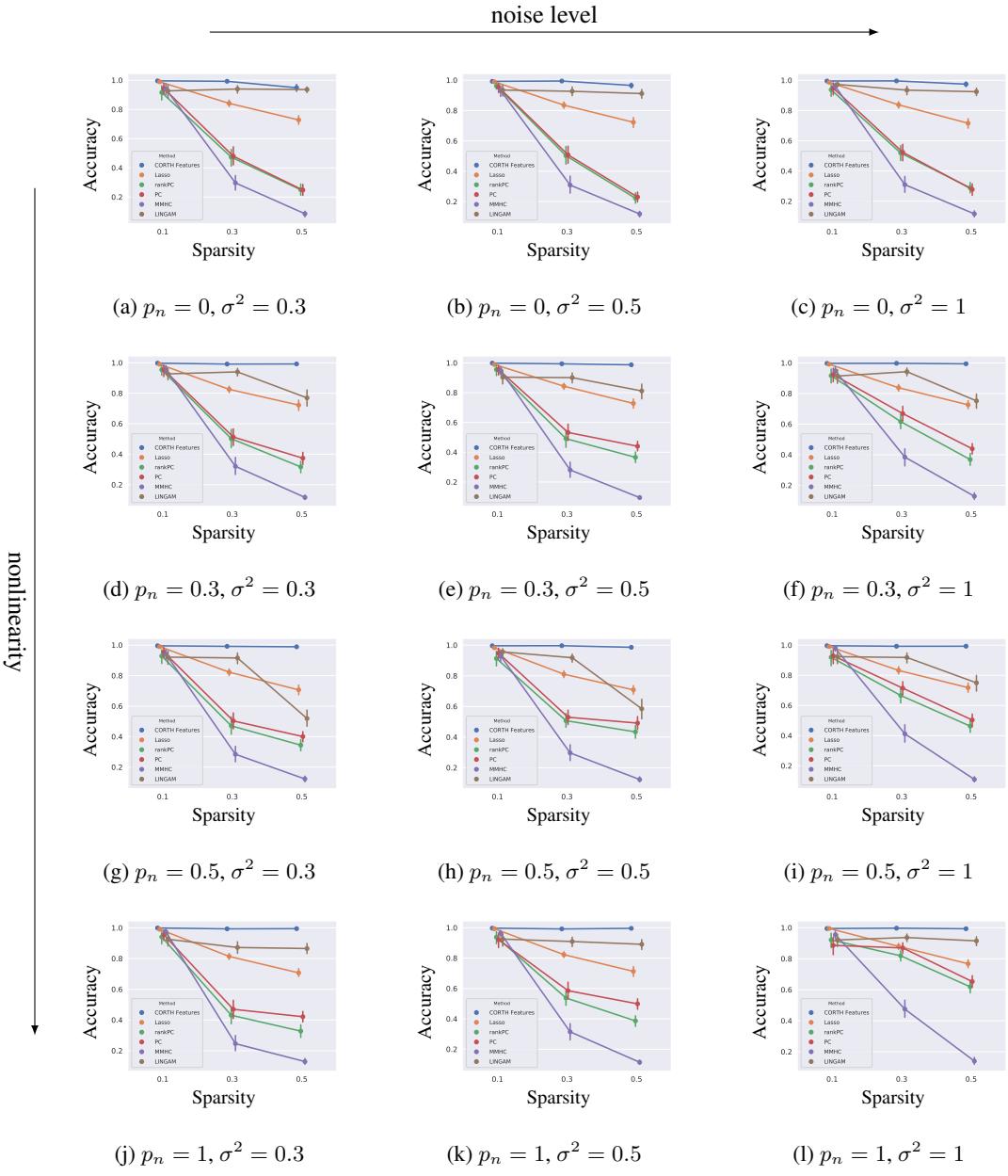


Figure G.8: 20 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

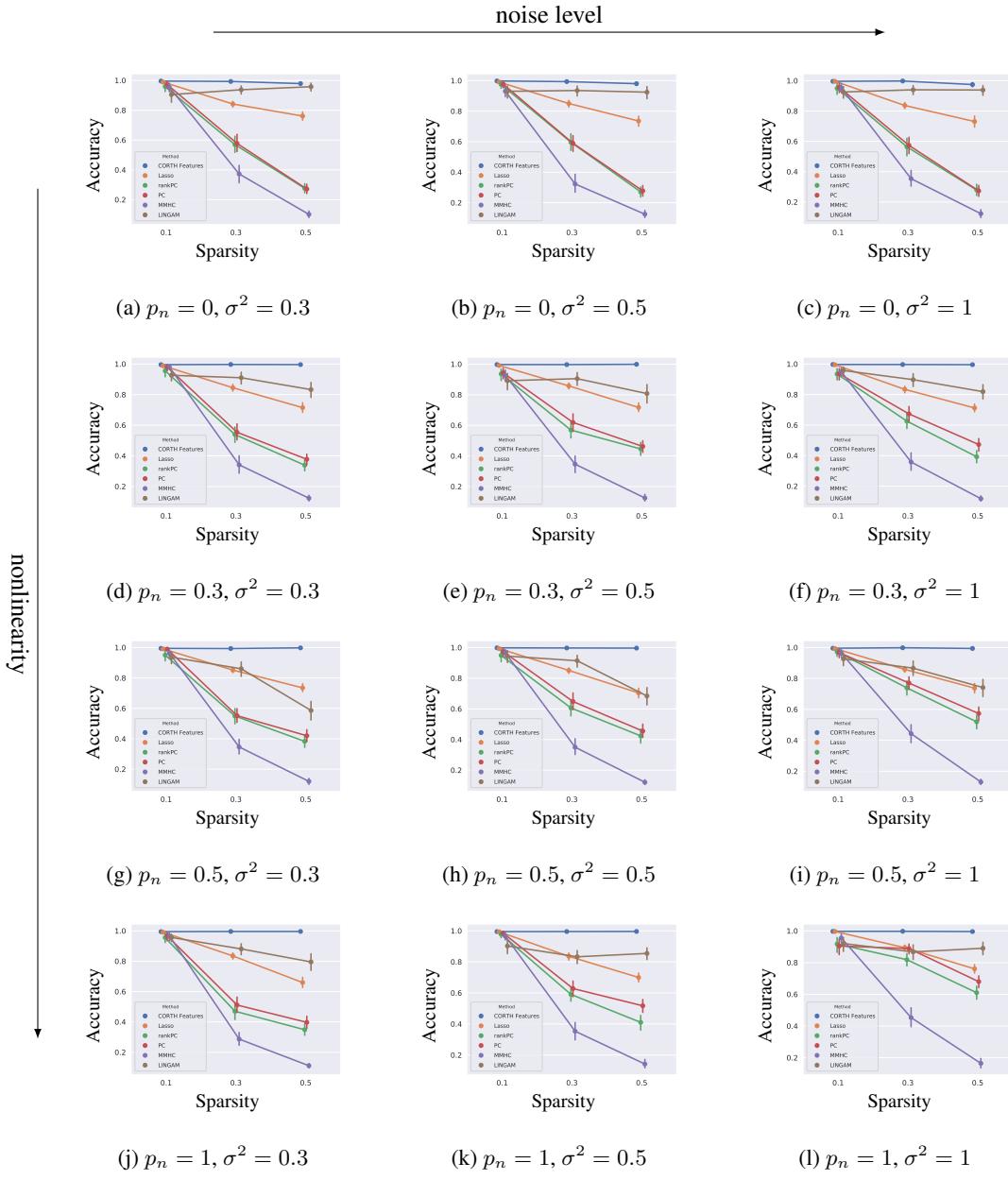


Figure G.9: 20 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

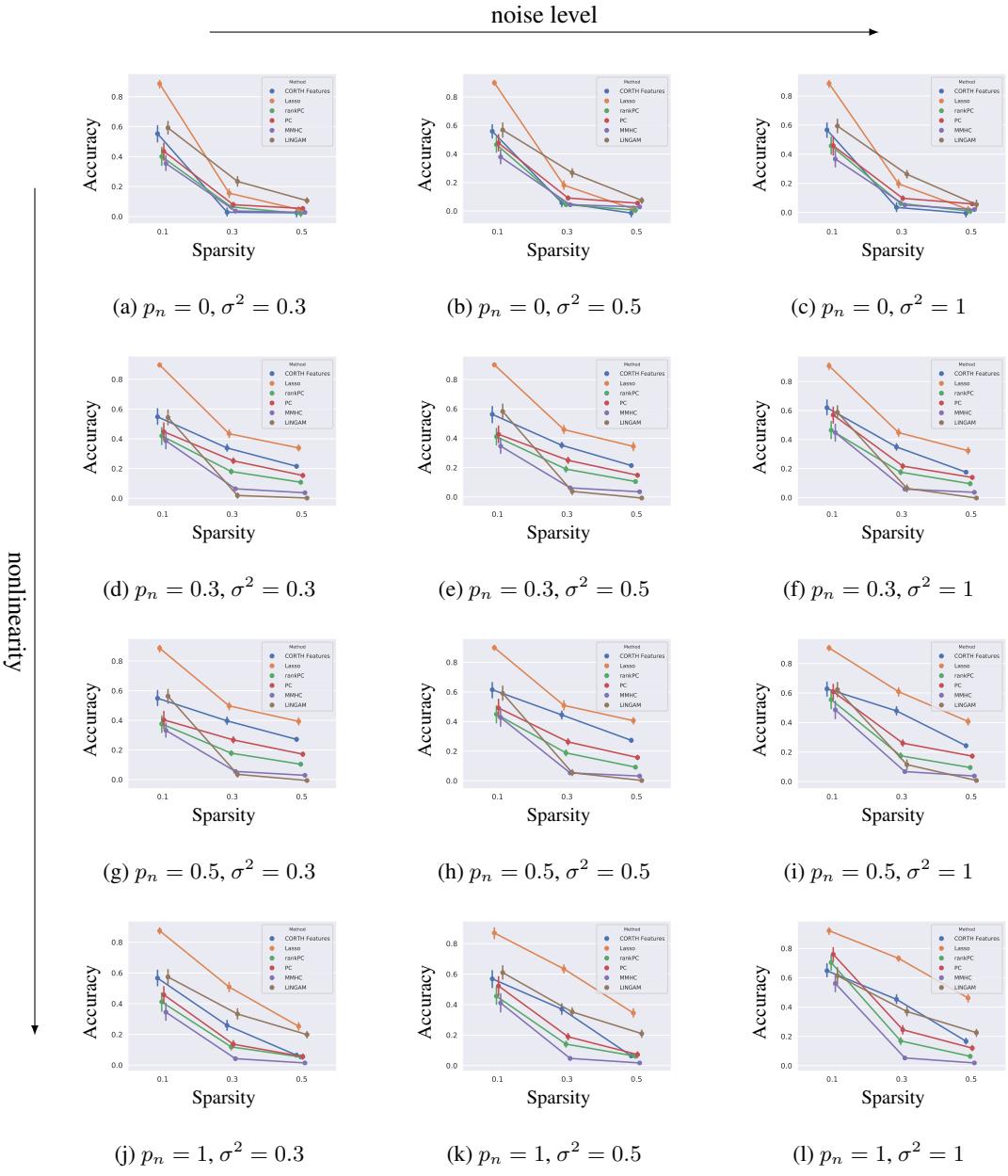


Figure G.10: 50 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

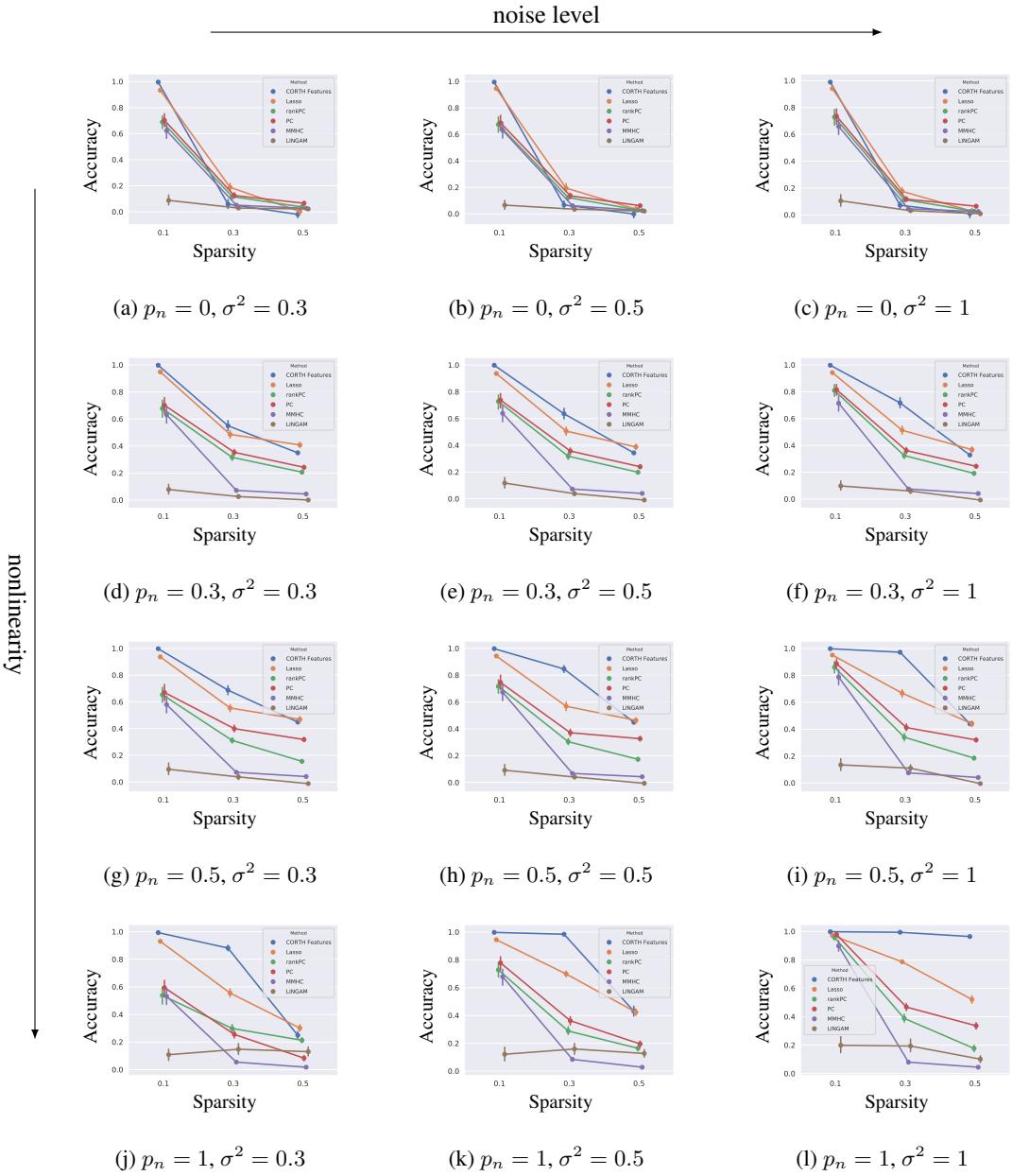


Figure G.11: 50 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

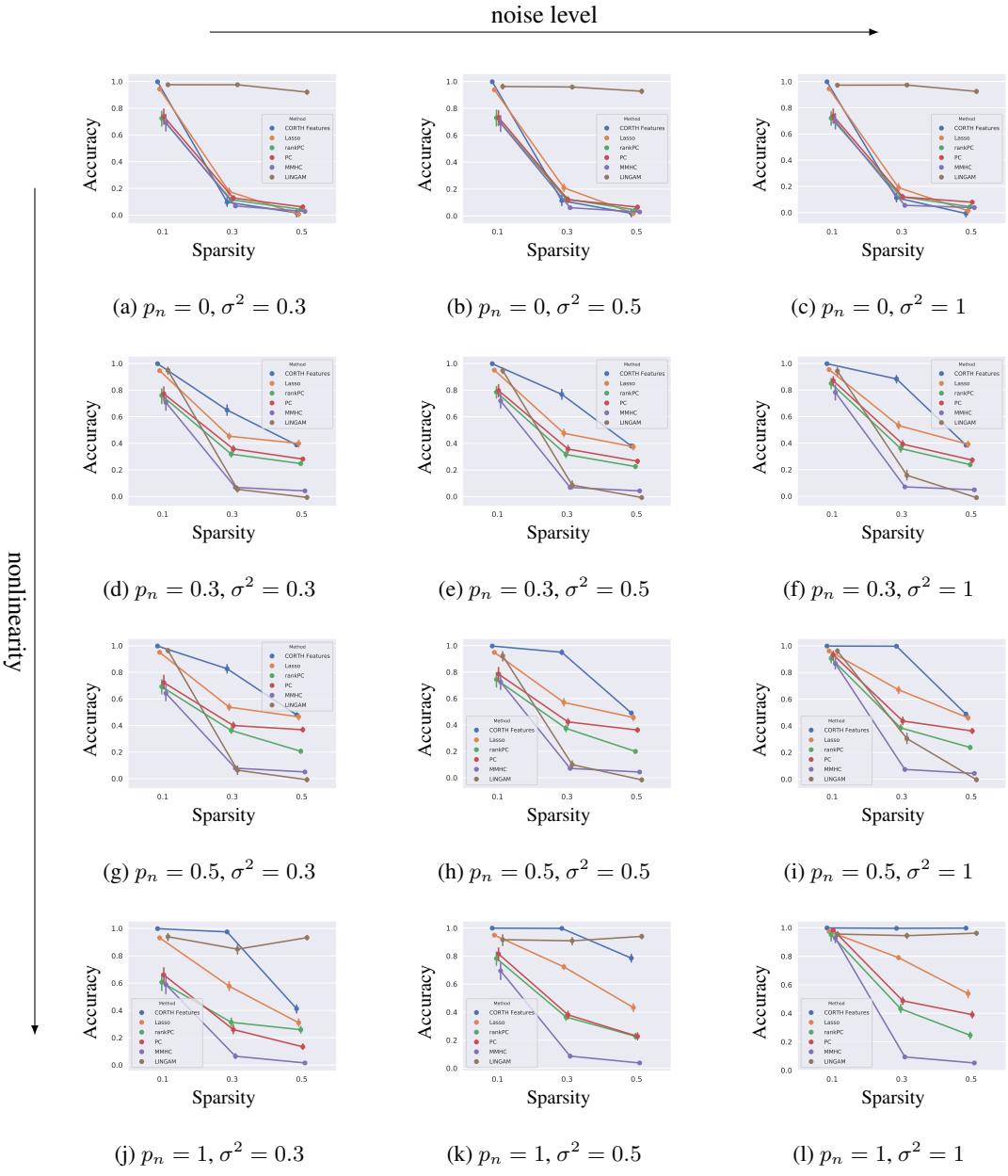


Figure G.12: 50 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

## H Supplementary Figures for Performance in Inferring Direct Causes, F1 Score as Metric

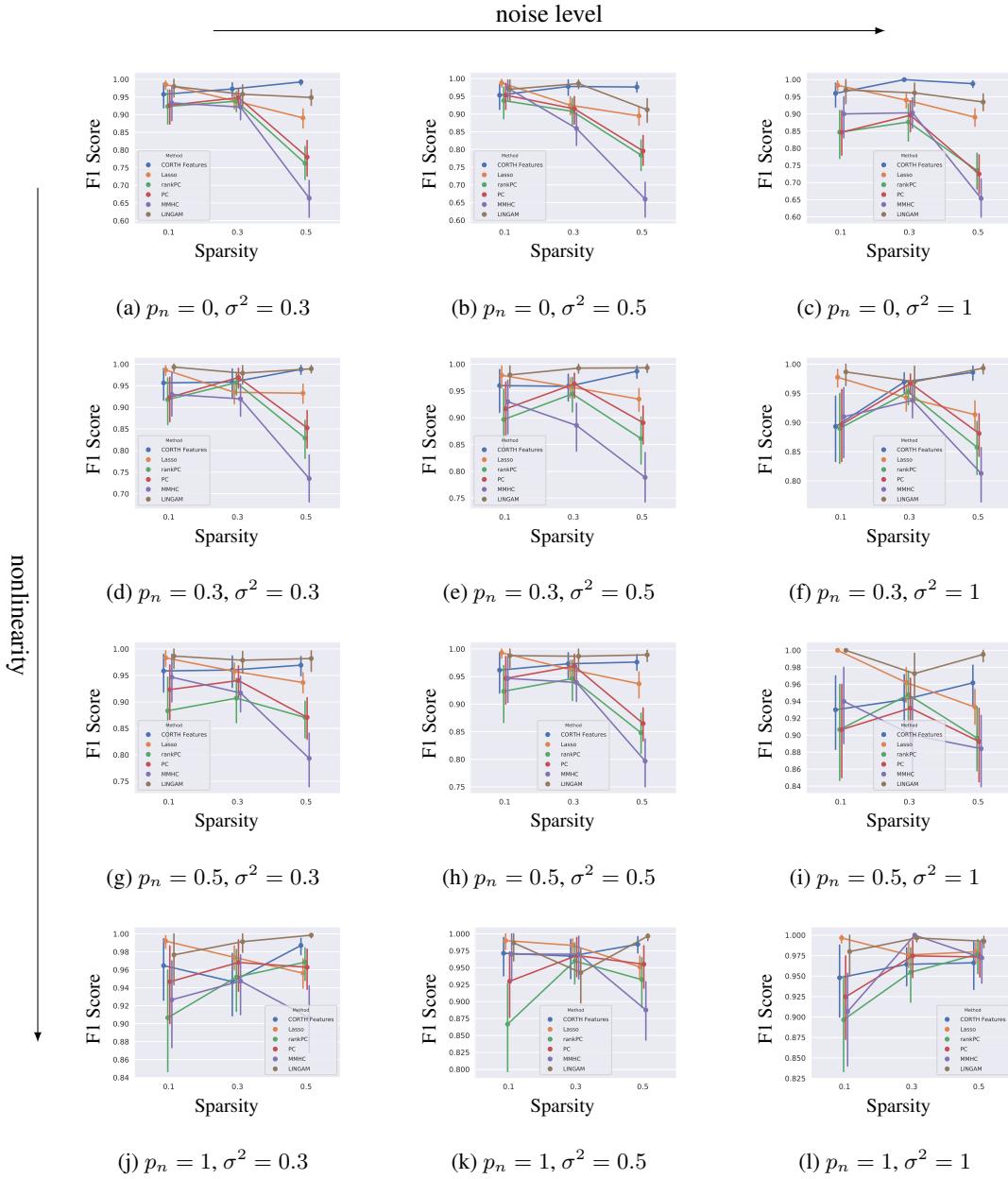


Figure H.1: 5 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

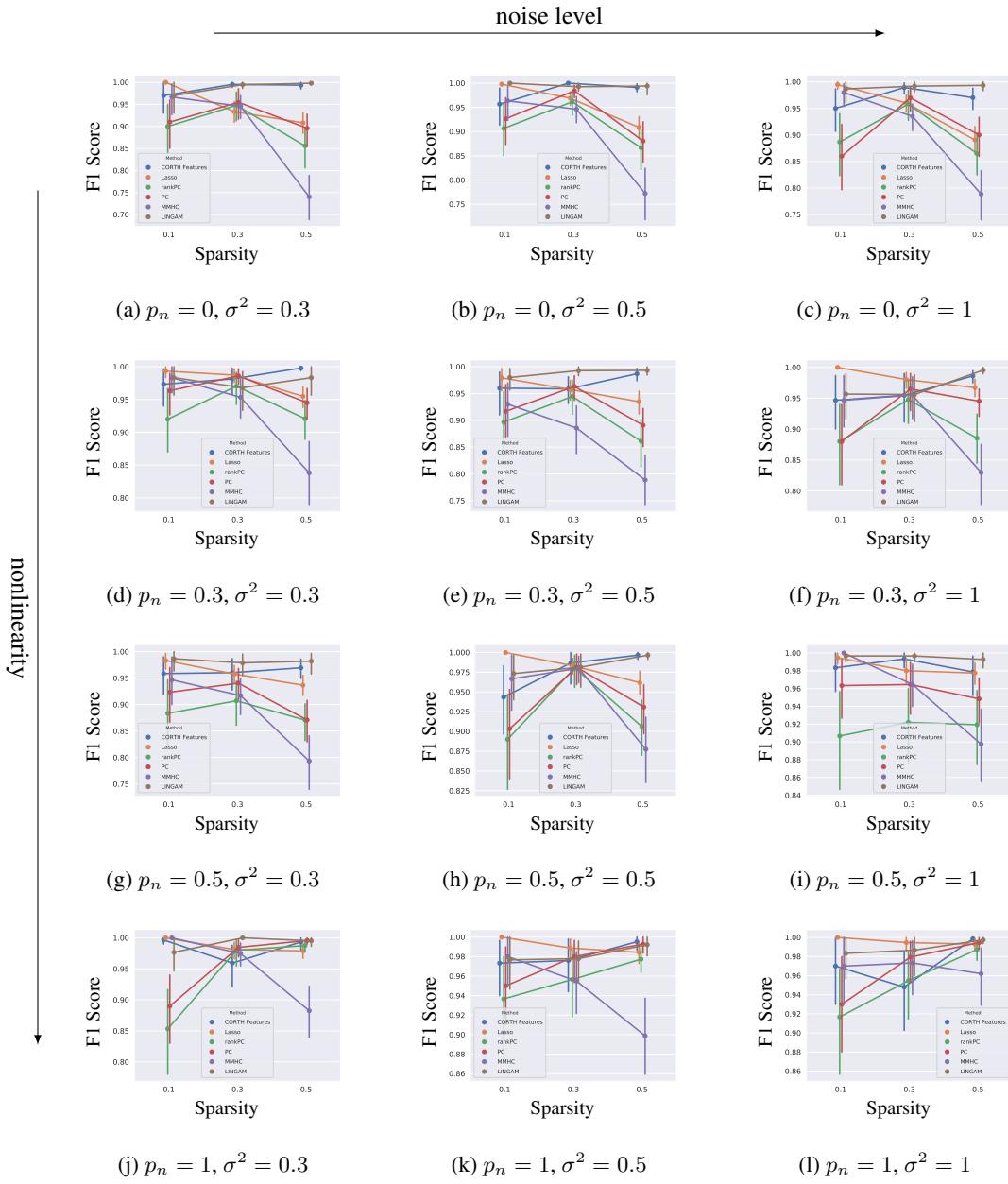


Figure H.2: 5 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

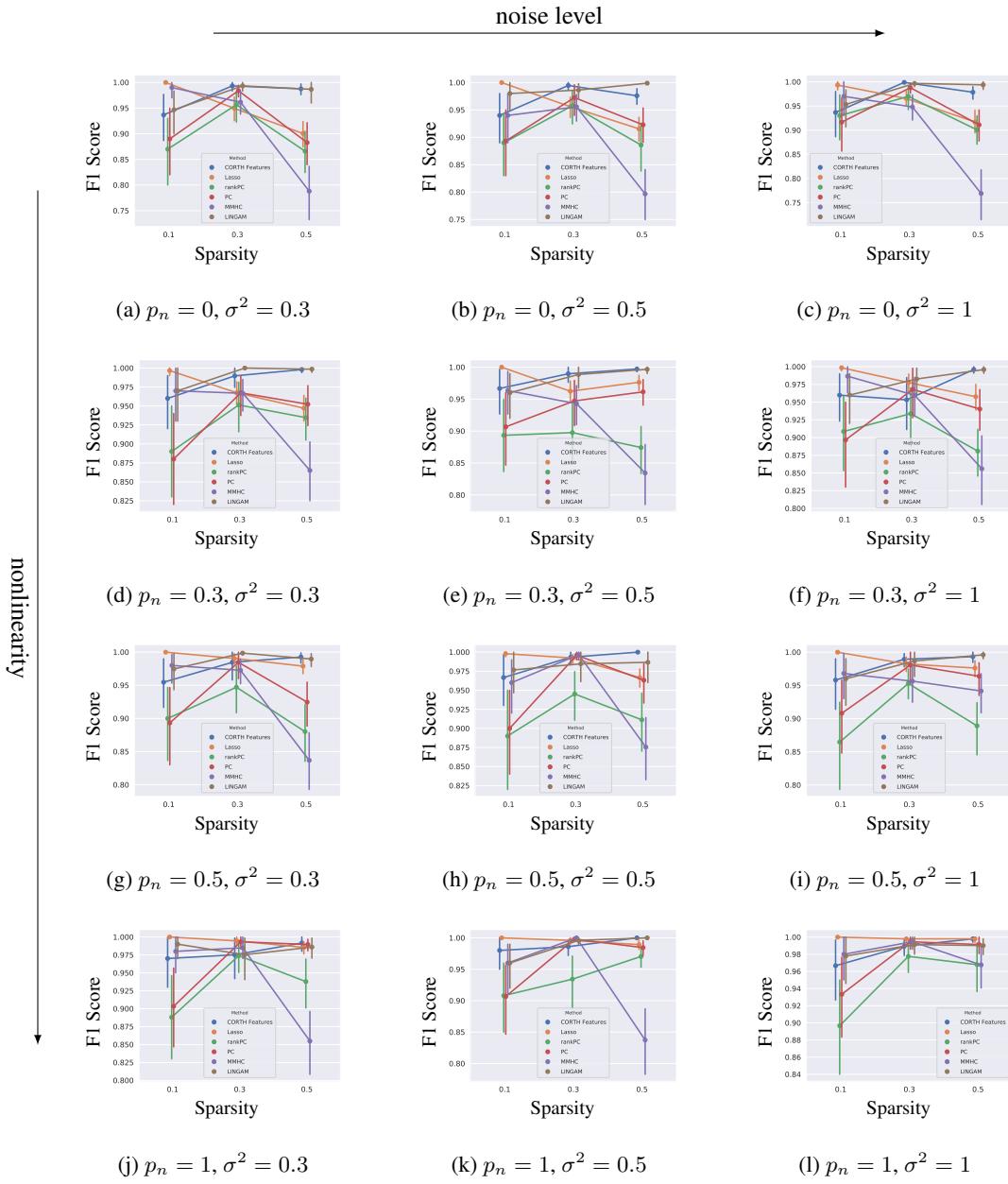


Figure H.3: 5 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

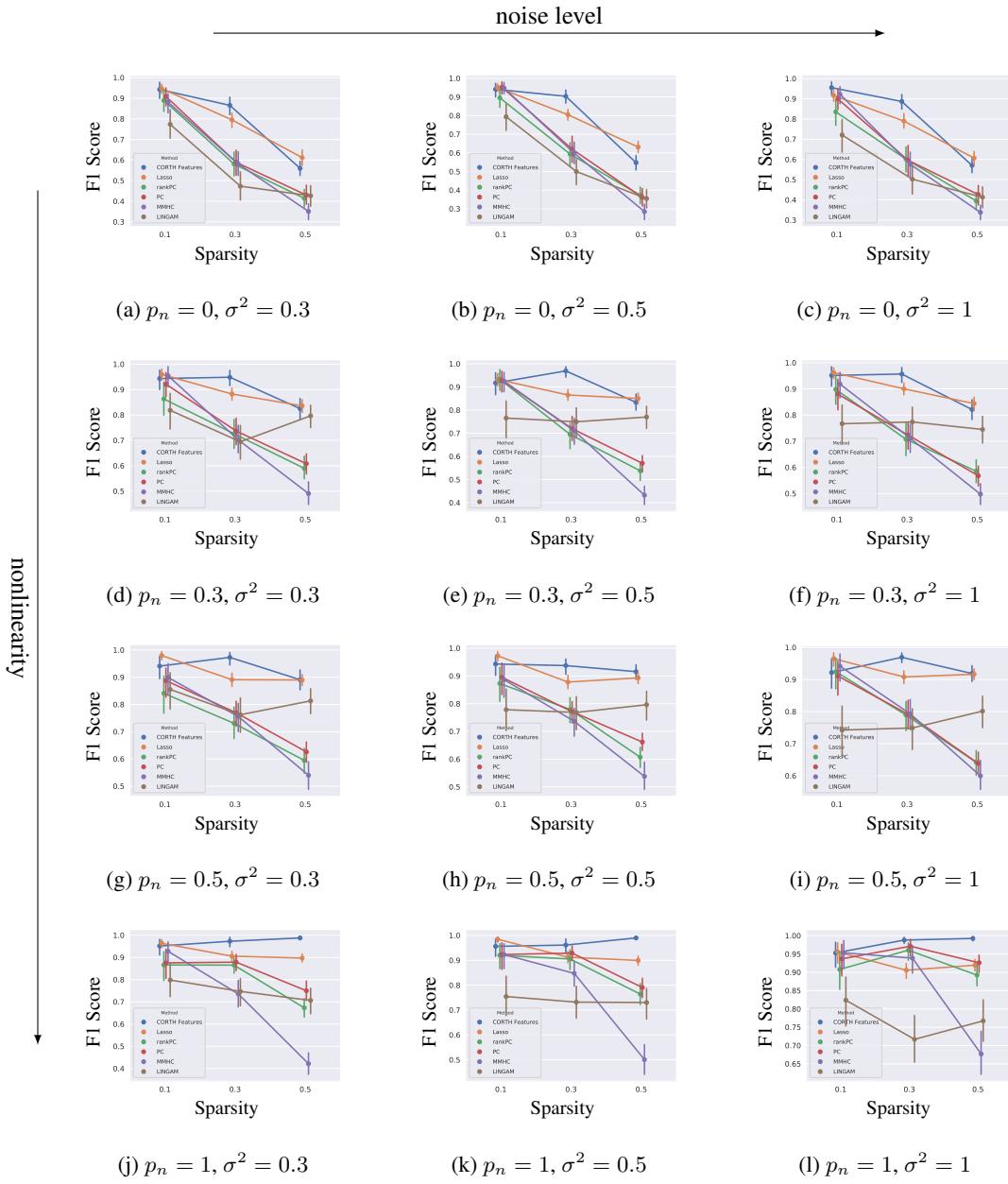


Figure H.4: 10 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

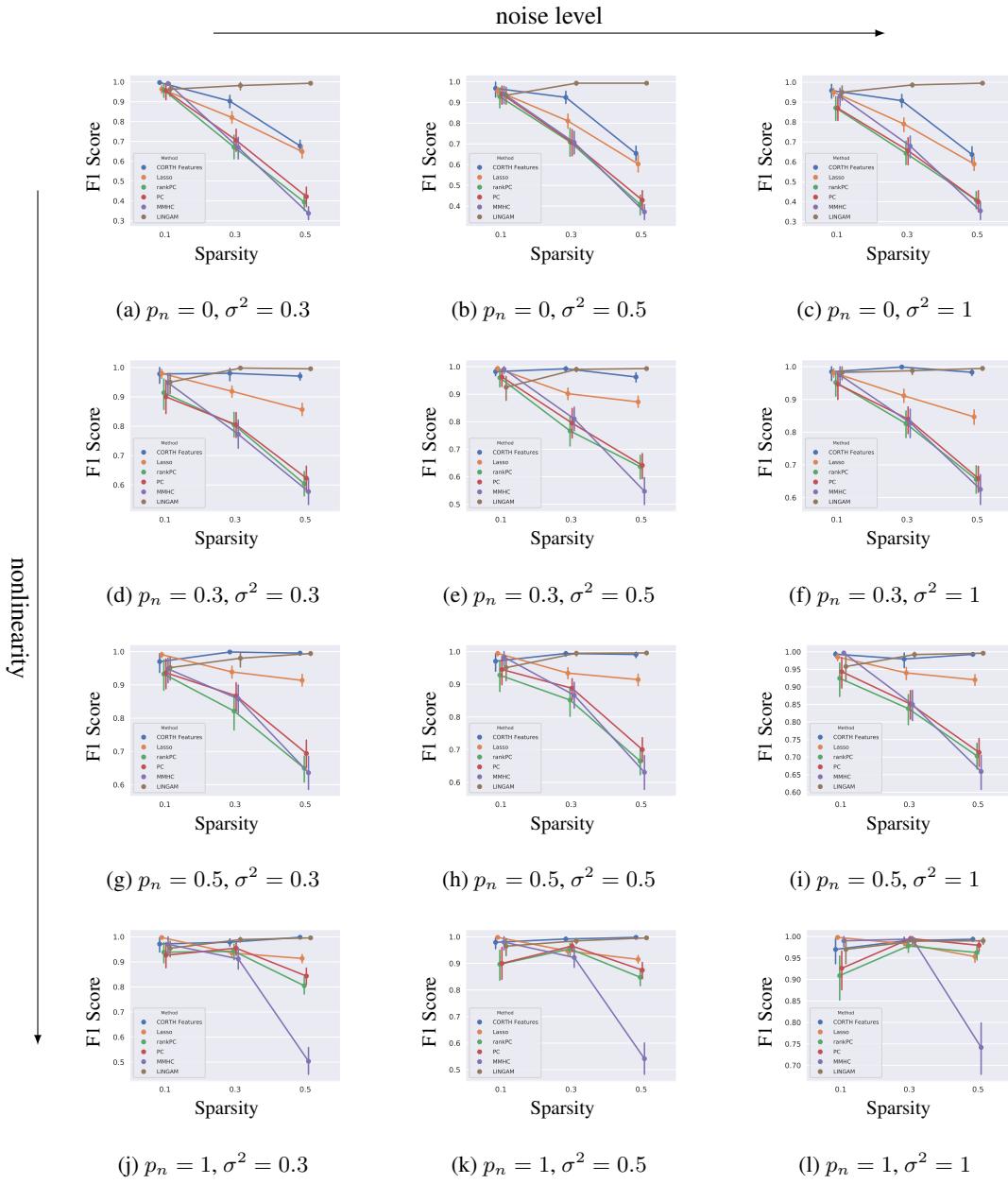


Figure H.5: 10 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

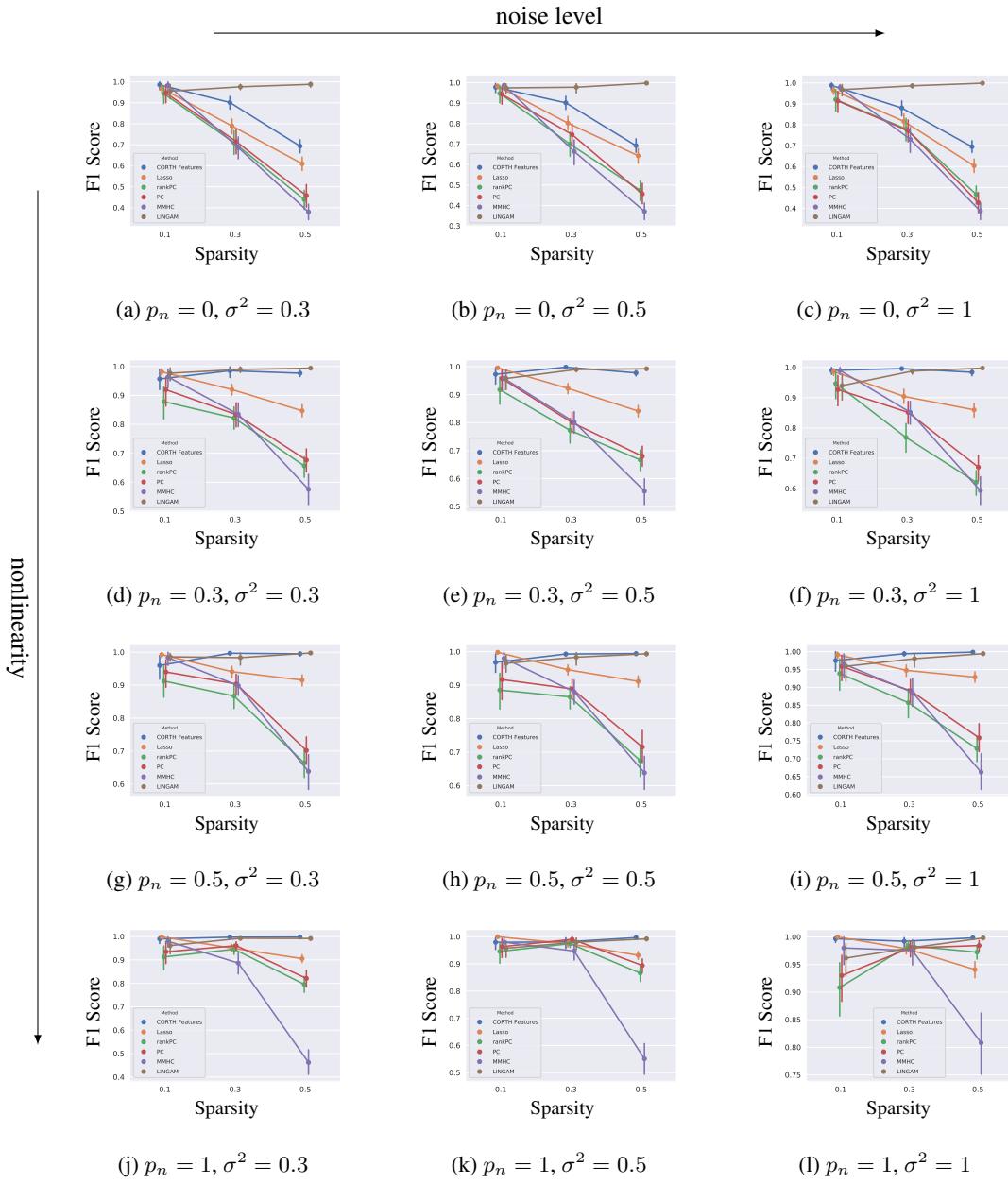


Figure H.6: 10 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

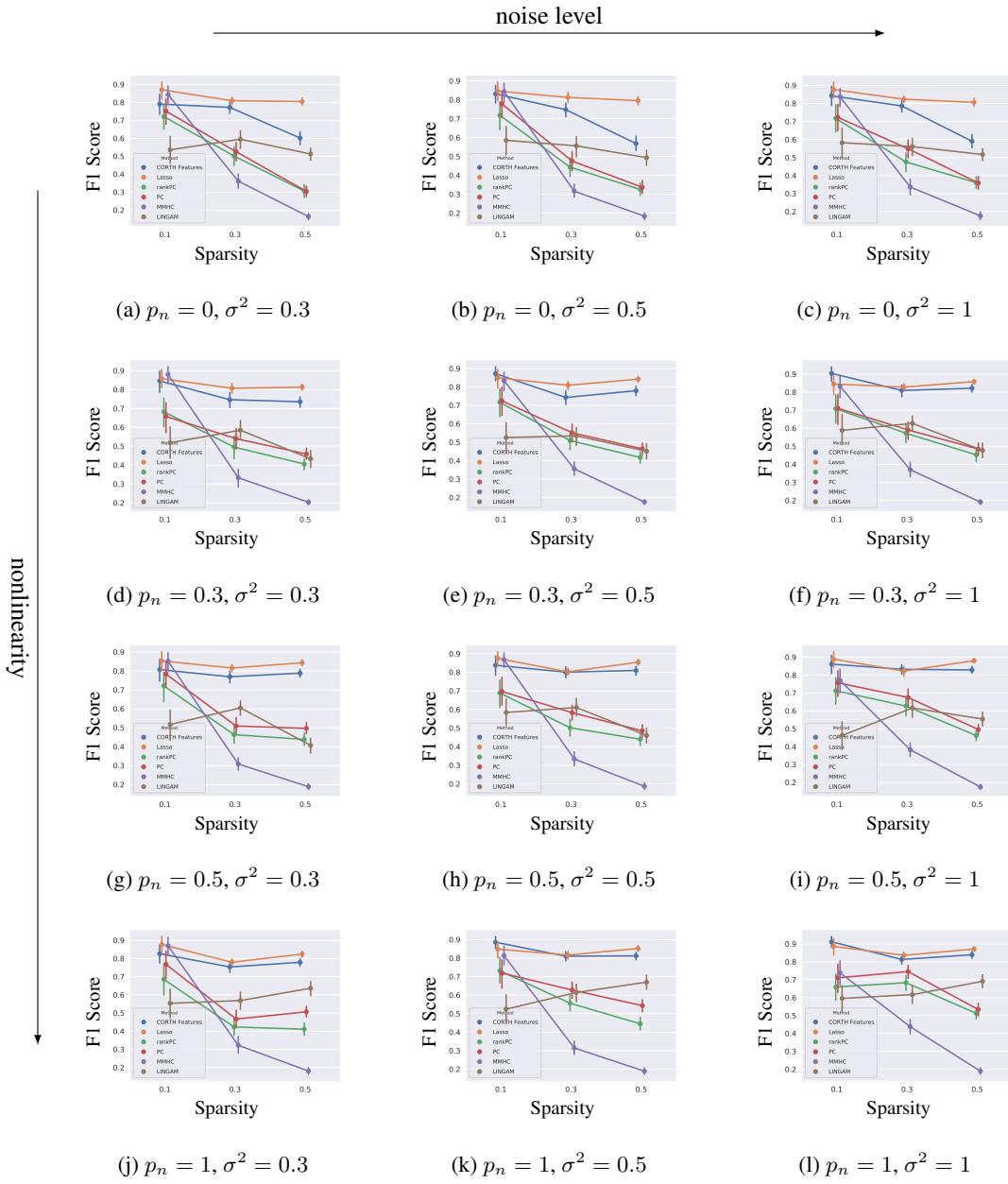


Figure H.7: 20 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

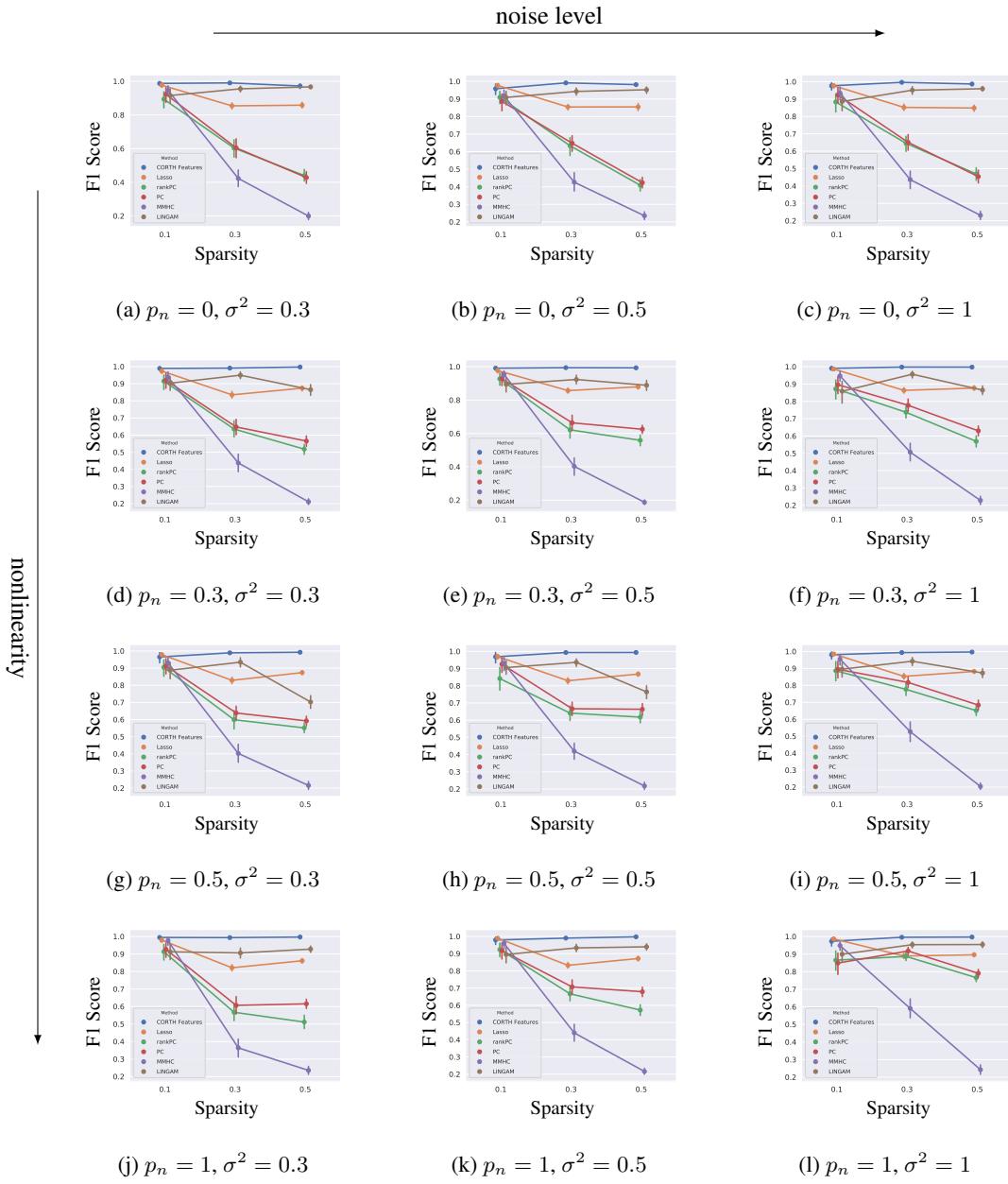


Figure H.8: 20 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

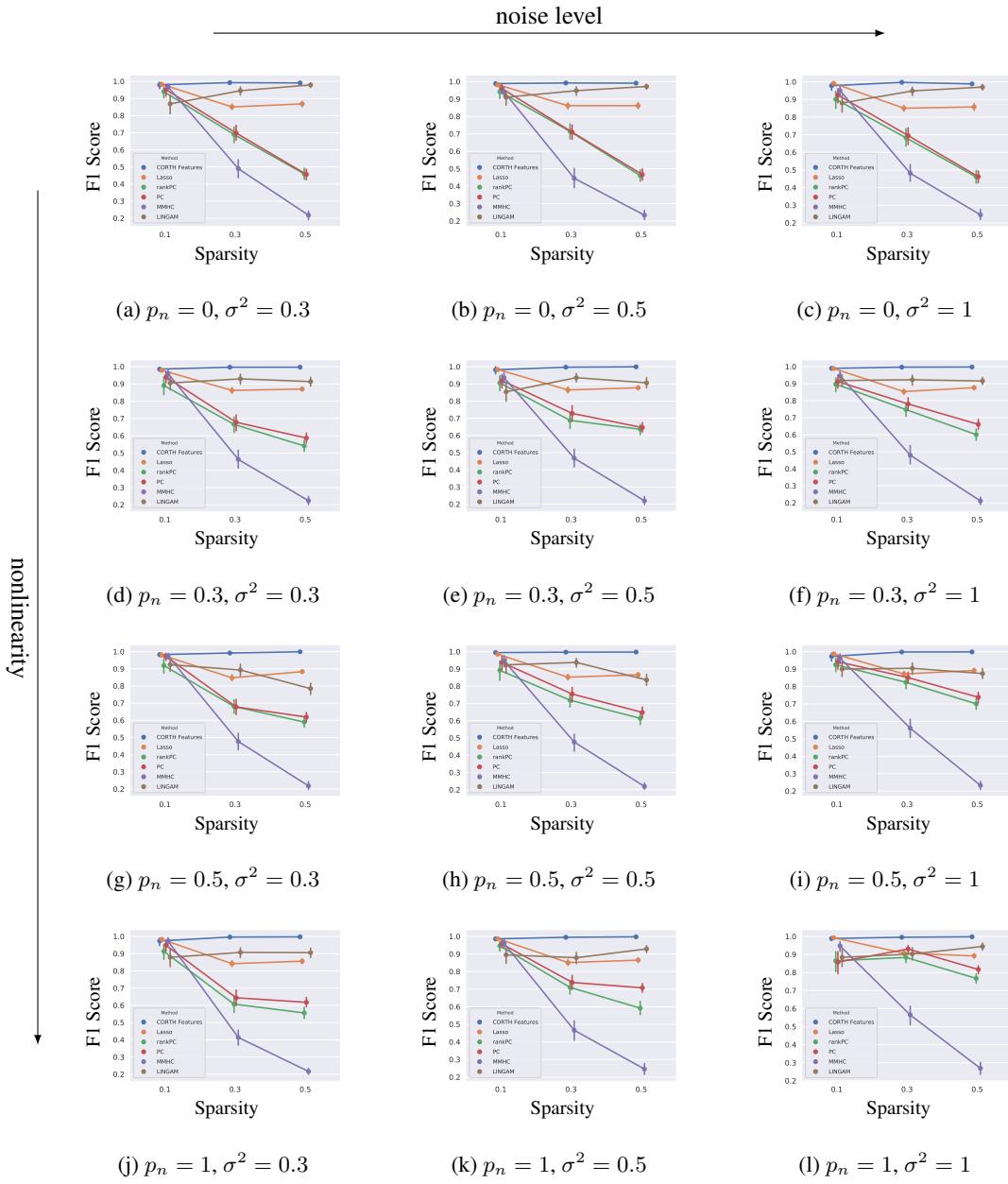


Figure H.9: 20 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

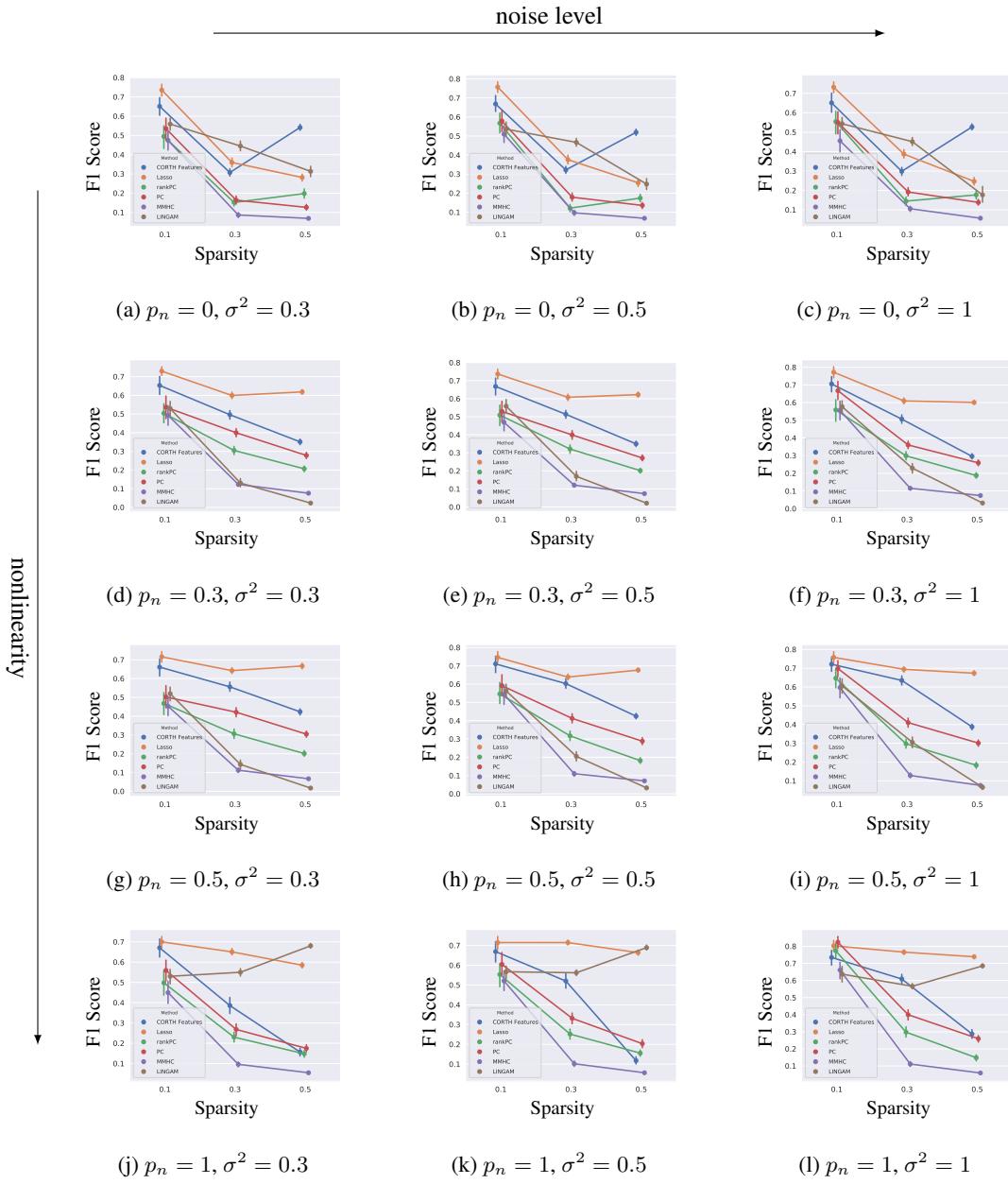


Figure H.10: 50 nodes, 100 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

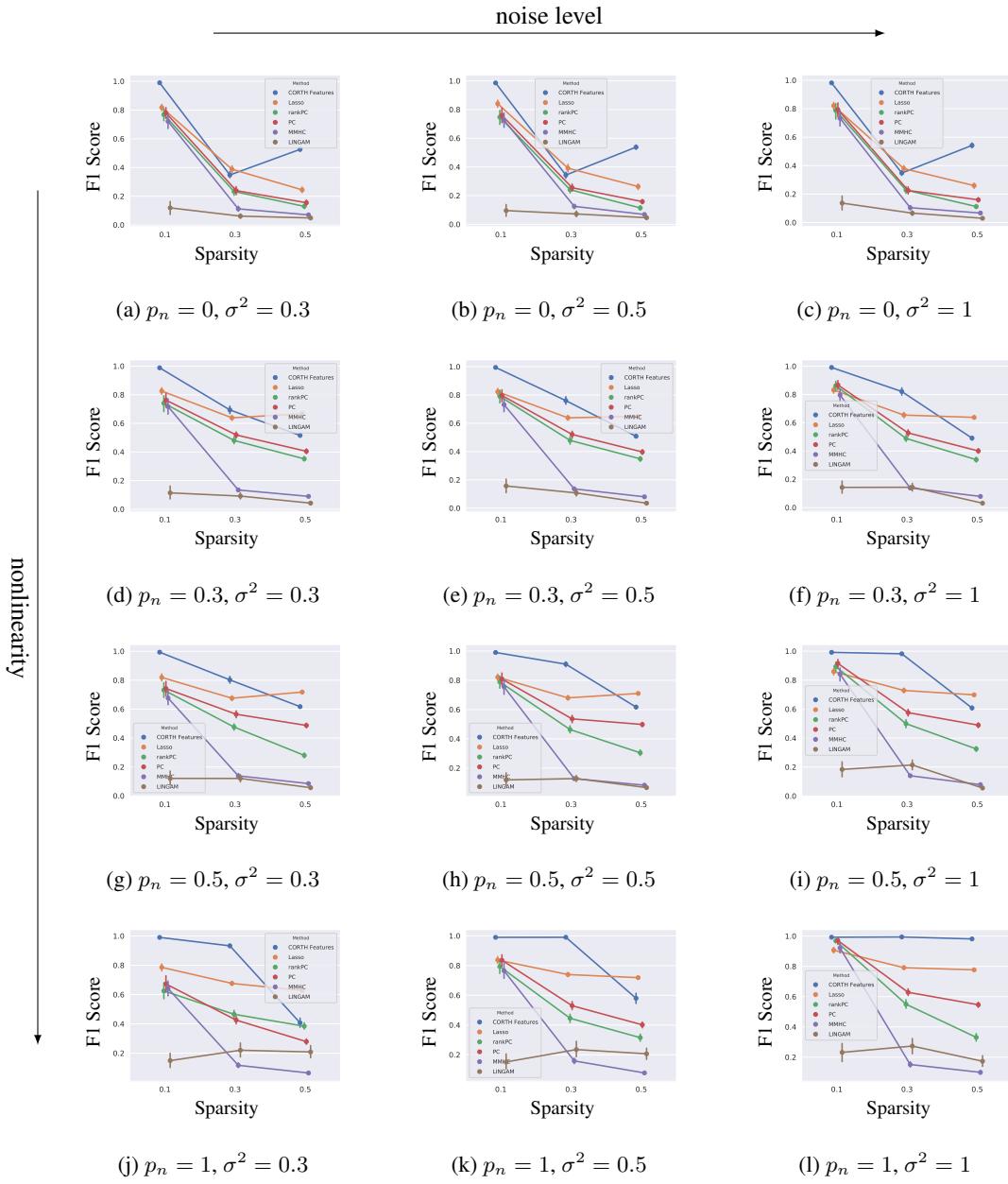


Figure H.11: 50 nodes, 500 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.

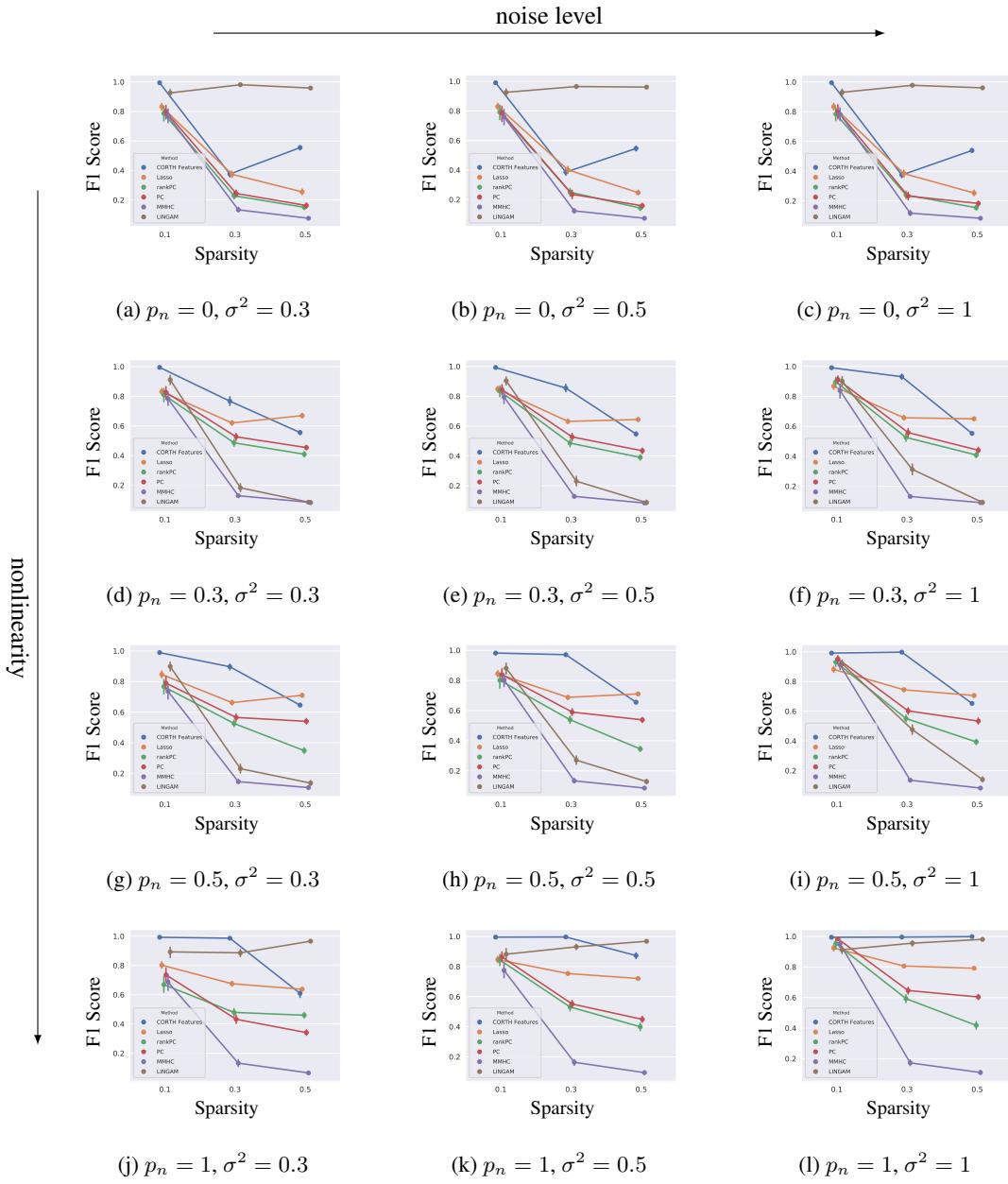


Figure H.12: 50 nodes, 1000 observations, 100 simulations. Compared to other methods, our approach is stable wrt. noise level, sparsity and even partially non-linear relationships, significantly outperforming the majority of baselines.