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#### Abstract

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# Locally- but not globally-identified SVARs 

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cemmap working paper CWP40/20

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# Locally- but not Globally-identified SVARs* 

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#### Abstract

This paper analyzes Structural Vector Autoregressions (SVARs) where identification of structural parameters holds locally but not globally. In this case there exists a set of isolated structural parameter points that are observationally equivalent under the imposed restrictions. Although the data do not inform us which observationally equivalent point should be selected, the common frequentist practice is to obtain one as a maximum likelihood estimate and perform impulse response analysis accordingly. For Bayesians, the lack of global identification translates to nonvanishing sensitivity of the posterior to the prior, and the multi-modal likelihood gives rise to computational challenges as posterior sampling algorithms can fail to explore all the modes. This paper overcomes these challenges by proposing novel estimation and inference procedures. We characterize a class of identifying restrictions that deliver local but non-global identification, and the resulting number of observationally equivalent parameter values. We propose algorithms to exhaustively compute all admissible structural parameter given reduced-form parameters and utilize them to sampling from the multi-modal posterior. In addition, viewing the set of observationally equivalent parameter points as the identified set, we develop Bayesian and frequentist procedures for inference on the corresponding set of impulse responses. An empirical example illustrates our proposal.


Keywords: local identification, Bayesian inference, Markov Chain Monte Carlo, robust Bayesian inference, frequentist inference. multi-modal posterior
JEL codes: C01,C13,C30,C51.

[^1]
## I Introduction

Macroeconomic policy analysis makes extensive use of impulse response analysis based on Structural Vector Autoregressions (SVARs). Various types of identifying assumptions have been proposed, including equality and sign restrictions, and analytical investigation of whether they point- or setidentify the objects of interest is an active area of research. The seminal work of Rubio-Ramirez et al. (2010) (henceforth RWZ) shows a necessary and sufficient condition for zero restrictions to achieve global identification. This class of zero restrictions, however, does not exhaust the universe of zero and non-zero equality restrictions that are relevant in practice. Questions regarding identification, estimation, and inference when identification is not global remain largely open.

This paper focuses on a class of SVARs where the imposed identifying restrictions guarantee local identification but do not attain global identification. The set of observationally equivalent structural parameters then consists of multiple isolated points, which implies that the likelihood can have multiple peaks of the same height. Such locally- but non-globally identified SVARs appear in various settings of practical relevance. Examples include non-zero restrictions which set the structural parameters to calibrated values, non-recursive zero restrictions, equality restrictions across shocks and/or equations, and heteroskedastic SVARs with across-regime restrictions on the structural coefficients. Although the data do not inform us which observationally equivalent point should be selected, the common frequentist practice is to obtain one as a maximum likelihood estimator and perform impulse response analysis as if it were the only maximizer. In our view, this practice is prevalent due to the lack of an efficient algorithm that can uncover all the local maxima. Standard Bayesian analysis also faces challenges when the likelihood has multiple modes. First, the lack of global identification leads the posterior to remain sensitive to the choice of prior even asymptotically. Second, posterior sampling algorithms may fail to explore all the modes, resulting in an inaccurate approximation of the posterior.

This paper proposes methods for estimation and inference that overcome these challenges. We first characterize a class of equality and sign restrictions that delivers local but non-global identification. Second, we show a necessary and sufficient condition for local identification that can be easily checked under a general class of equality constraints imposed on the structural parameters or functions of them. Third, we investigate how many observationally equivalent parameter values exist under such identifying restrictions, and propose algorithms to exhaustively compute them given reduced-form parameter values. Specifically, we exploit the orthogonal matrix parametrization of Uhlig (2005) and Rubio-Ramirez et al. (2010) and pin down the observationally equivalent parameter points (conditional on the reduced-form parameters) by sequentially exhausting the admissible orthogonal vectors satisfying the imposed restrictions, or in some case solving a system of polynomial equations. We provide an intuitive geometric exposition that illustrates the mechan-
ism driving the lack of global identification and the number of observational equivalent parameter values. As a byproduct, we also characterize the set of reduced-form parameter values that yield no admissible structural parameters (i.e, the empty identified set) despite the condition for local identification being met.

Our proposal for computing the identified set contributes to standard Bayesian inference by simplifying and stabilising sampling from the multi-modal posterior. In addition, Bayesian inference requires specifying a prior over the observationally equivalent parameter values. For the case where the user cannot form this prior or is not confident about the choice, we consider (multiple prior) robust Bayesian procedures that draw posterior inference for the impulse responses under local but non-global identification. Viewing the set of observationally equivalent parameter points as the identified set (a set-valued map from the reduced-form parameters to the set of observationally equivalent structural parameters), we extend the approach of Kline and Tamer (2016) and Giacomini and Kitagawa (2020), designed primarily for models with interval identified sets, to cases where the identified set consists of a finite number of points. Specifically, we consider projecting the posterior credible region for the reduced-form parameters to the impulse responses through the discrete identified set mapping. This approach obtains an asymptotically frequentist valid confidence intervals in the presence of local identification.

To illustrate our proposal, we apply the method to a locally identified New-Keynesian monetary policy SVAR. We show that, when a single element is selected from the identified set, the choice of element can lead to significantly different and arguably contradictory results. Our proposal for robust Bayesian frequentist-valid inference, in contrast, explores the outcome from every admissible impulse response, and provides their summary.

## I. 1 Related literature

The theory of identification for linear simultaneous equation models has a long history in econometrics. See Dhrymes (1978), Fisher (1966) and Hausman (1983), among others. Rothenberg (1971) analyses identification in parametric models. Building on this, Giannini (1992) proposes a criterion for verifying local identification for SVAR models. This criterion takes the form of rank conditions for the Hessian matrix of the average likelihood. It is much weaker than the necessary and sufficient condition for global identification shown in Rubio-Ramirez et al. (2010). The focus of this paper is the class of identifying restrictions that satisfies the former but not the latter. Once local identification is guaranteed, Giannini (1992) proposes estimating the parameters of the SVAR by numerically maximizing the likelihood. This approach is also recommended by the textbooks Amisano and Giannini (1997), Lütkepohl (2006) and Hamilton (1994). For the locally identified models considered in this paper, however, the maximum likelihood estimate is not necessarily unique, and
a typical numerical maximization routine will select only one point in a non-systematic manner (e.g. depending on a choice of initial value). Sims and Zha (1999) and Hamilton et al. (2007) include discussions of the existence of multiple likelihood peaks due to local identification.

Following Uhlig (2005), Rubio-Ramirez et al. (2010), and Granziera et al. (2018), we parameterize an SVAR by its reduced-form VAR-parameters and the orthogonal matrix relating its reduced-form error covariance matrix and structural parameters. Fixing the reduced-form parameters, finding all the observationally equivalent structural parameters reduces to finding all the admissible orthogonal matrices that satisfy the imposed identifying restrictions. Compared to expressing the non-linear equation system by the reduced form and structural parameters, this formulation is advantageous in terms of geometric interpretability and analytical tractability. In addition, it simplifies not only assessing local identification (e.g., Magnus and Neudecker, 2007), but also obtaining all the solutions given the reduced-form parameters.

Our paper is related to the growing literature on SVARs that are set-identified through sign and zero restrictions (Faust 1998; Canova and de Nicoló 2002; Uhlig 2005; Mountford and Uhlig 2009, Arias et al. 2018a, Gafarov et al. 2018, Giacomini and Kitagawa 2020, Granziera et al. 2018, among others). The identified set of impulse responses in this class of models is a set with a positive measure if nonempty, whereas the identified set here consists of a finite number of isolated points, each corresponding to a solution of a non-linear system of equations. This difference in the topological features of the identified set distinguishes our inferential procedure from these works.

Our inference proposals include an application of the multiple-prior robust Bayesian approach of Giacomini and Kitagawa (2020) to the discrete-point identified set. We believe that the robust Bayes approach is attractive as the prior knowledge of the researcher is typically exhausted in constructing the prior distribution for the reduced form parameters (e.g., the Minnesota prior) and the identifying restrictions. It is then difficult to construct a credible prior over the parameter values in the identified set. The use of multiple priors is a way of reflecting this lack of prior knowledge while delivering posterior conclusions that are robust to the choice of a prior within a given class. Depending on the application, the class of priors considered in Giacomini and Kitagawa (2020) could be too large. In such cases, refining the set of priors would be sensible. This can be done, for instance, by applying the approaches considered in Giacomini et al. (2018) and Giacomini et al. (2019b), although we do not present them in this paper. Chen et al. (2018), Kline and Tamer (2016), Liao and Simoni (2019), Moon and Schorfheide (2012), and Norets and Tang (2014) propose Bayesian approaches for drawing posterior inference for the identified set. To our knowledge, none of these proposals have been applied to the case where the identified set consists of isolated points.

The results and proposals of this paper, from identification to estimation and inference, can also contribute to the literature that bridges Dynamic Stochastic General Equilibrium (DSGE) and

VAR models. The solution of a linearized DSGE model can be summarized by a state-space representation that implies, under appropriate invertibility conditions, an (infinite order) SVAR subject to specific identifying restrictions (see, Christiano et al. 2006, Fernandez-Villaverde et al. 2007, and Ravenna, 2007 for example). As stressed by Canova (2005, chapter 4) among others, popular identification schemes that lead to global identification, such as the Cholesky decomposition, cannot be justified in a large class of DSGE models. Hence, if the mapping between the DSGE and the SVAR is unique as in Christiano et al. (2006, Proposition 1), DSGE-based identifying restrictions can result in local (but not global) identification. This is due to the non-recursive nature of the identification scheme, and the possible multiplicity of solutions characterizing the DSGE model. See Iskrev (2010), Komunjer and Ng (2011) and Qu and Tkachenko (2012) for DSGE models, and Al-Sadoon and Zwiernik (2019) for local identification in linear rational expectation models. This paper offers estimation and inference methods to handle local identification in these models.

The remainder of the paper is organized as follows. Section II introduces notation and a general analytical framework for SVARs whose identifying restrictions take the form of equality and sign restrictions. It also presents a new necessary and sufficient condition for local identification in SVARs. Section III discusses a battery of examples of locally- but not globally-identified SVARs. Section IV presents algorithms for computing observationally equivalent parameter values, and Section V proposes inference methods that accommodate frequentist, Bayesian, and robust Bayesian perspectives. Section VI presents an empirical example and Section VII concludes. Further results on local identification are reported in Appendices A and B, and the proofs omitted from the main text are presented in Appendix C.

## II Econometric framework

Let $y_{t}$ be a $n \times 1$ vector of variables observed over $t=1 \ldots T$. The SVAR model is specified as

$$
\begin{equation*}
A_{0} y_{t}=a+\sum_{j=1}^{p} A_{j} y_{t-j}+\varepsilon_{t} \tag{1}
\end{equation*}
$$

where $\varepsilon_{t}$ is a $n \times 1$ multivariate normal white noise process with null expected value and covariance matrix equal to the identity matrix $I_{n}$. The quantities $A_{0}, A_{1}, \ldots, A_{p}$ are $n \times n$ matrices of parameters, and $a$ is a $n \times 1$ vector of constant terms. The set of structural parameters is denoted by $A=\left(A_{0}, A_{+}\right) \in \mathcal{A} \subset \mathbb{R}^{(n+m) n}$, with $m \equiv n p+1$ and $A_{+} \equiv\left(a, A_{1}, \ldots, A_{p}\right)$ being a $n \times m$ matrix. We also assume that the initial conditions $y_{1}, \ldots, y_{p}$ are given.

The reduced-form representation of the SVAR, obtained by pre-multiplying by the inverse of
$A_{0}$, is the standard VAR model

$$
\begin{equation*}
y_{t}=b+\sum_{j=1}^{p} B_{j} y_{t-j}+u_{t} \tag{2}
\end{equation*}
$$

where $B_{j}=A_{0}^{-1} A_{j}, j=1, \ldots, p, b=A_{0}^{-1} a, u_{t}=A_{0}^{-1} \varepsilon_{t}$ and $E\left(u_{t} u_{t}^{\prime}\right) \equiv \Sigma=A_{0}^{-1} A_{0}^{-1 \prime}$. The set of reduced-form parameters is $\phi=(B, \Sigma) \in \Phi \subset \mathbb{R}^{n+n^{2} p} \times \Omega$, where $B=\left(b, B_{1}, \ldots, B_{p}\right)$ and $\Omega$ is the space of positive semi-definite matrices.

Assuming further that the VAR Eq. (2) is invertible, it has the VMA $(\infty)$ representation:

$$
y_{t}=c+\sum_{j=0}^{\infty} C_{j}(B) u_{t-j}=c+\sum_{j=0}^{\infty} C_{j}(B) A_{0}^{-1} \varepsilon_{t-j}
$$

where $C_{j}(B)$ is the $j$-th coefficient matrix of the inverted lag polynomial $\left(I_{n}-\sum_{j=1}^{p} B_{j} L^{j}\right)^{-1}$. We define the impulse response matrix at horizon $h\left(I R^{h}\right)$, the long-run impulse response matrix $\left(I R^{\infty}\right)$ and the long-run cumulative impulse response matrix $\left(C I R^{\infty}\right)$ to be

$$
\begin{align*}
I R^{h} & =C_{h}(B) A_{0}^{-1},  \tag{3}\\
I R^{\infty} & =\lim _{h \rightarrow \infty} I R^{h}=\left(I_{n}-\sum_{j=1}^{p} B_{j}\right)^{-1} A_{0}^{-1},  \tag{4}\\
C I R^{\infty} & =\sum_{j=0}^{\infty} I R^{h}=\left(\sum_{j=0}^{\infty} C_{j}(B)\right) A_{0}^{-1}, \tag{5}
\end{align*}
$$

In what follows throughout, we denote the Cholesky decomposition of $\Sigma$ by $\Sigma=\Sigma_{t r} \Sigma_{t r}^{\prime}$, where $\Sigma_{t r}$ is the unique lower-triangular Cholesky factor with non-negative diagonal elements. The column vectors of $\Sigma_{t r}^{-1}$ and $\Sigma_{t r}^{\prime}$ are denoted by $\Sigma_{t r}^{-1} \equiv\left(\tilde{\sigma}_{1}, \tilde{\sigma}_{2}, \ldots, \tilde{\sigma}_{n}\right)$ and $\Sigma_{t r}^{\prime} \equiv\left(\sigma_{1}, \sigma_{2}, \ldots, \sigma_{n}\right)$. The $i$-th entry of $\tilde{\sigma}_{j}$ and $\sigma_{j}$ are denoted by $\tilde{\sigma}_{j, i}$ and $\sigma_{j, i}$, respectively.

## II. 1 Identification of SVAR models

Identification analysis of SVAR models concerns solving $\Sigma=A_{0}^{-1} A_{0}^{-1 \prime}$ to decompose the reduced form error variance-covariance matrix $\Sigma$ into the matrix of structural coefficients $A_{0}$. Following Uhlig (2005), any structural matrix $A_{0}$ defined by a rotation of the Cholesky factor $A_{0}=Q^{\prime} \Sigma_{t r}^{-1}$ admits the decomposition $\Sigma=A_{0}^{-1} A_{0}^{-1 \prime}$ and, given the reduced-form parameters $\phi$, the set of admissible $A_{0}$ matrices can be represented by $\mathcal{A}_{0}(\phi) \equiv\left\{A_{0}=Q^{\prime} \Sigma_{t r}^{-1}: Q \in \mathcal{O}(n)\right\}$, where $\mathcal{O}(n)$ is the set of $n \times n$ orthogonal matrices. Let $R$ be generic notation denoting identifying restrictions. The identifying restrictions constrain the admissible values of $A$ to a subset of $\mathcal{A}$. We denote this subset by $\mathcal{A}_{R}$, and its projection for $A_{0}$ by $\mathcal{A}_{R, 0}$. Accordingly, let $\Phi_{R} \subset \Phi$ be the space of
reduced-form parameters formed by projecting $A \in \mathcal{A}_{R}$, and let

$$
\begin{equation*}
\mathcal{A}_{R, 0}(\phi) \equiv \mathcal{A}_{0}(\phi) \cap \mathcal{A}_{R, 0} \tag{6}
\end{equation*}
$$

which is nonempty for $\phi \in \Phi_{R}$.
We define global and local identification for an SVAR as follows.
Definition 1 (Global identification). An SVAR model is globally identified under identifying restrictions $R$ if for almost every $A \in \mathcal{A}_{R}$ there is no other observationally equivalent $A$ in $\mathcal{A}_{R}$.

Definition 2 (Local identification). An SVAR model is locally identified under identifying restrictions $R$ if for almost every $A \in \mathcal{A}_{R}$, there exists an open neighborhood $G$ such that $G \cap \mathcal{A}_{R}$ contains no other observationally equivalent $A$.

Some remarks on these two notions of identification are in order. An equivalent definition of global identification would be that, for almost every $\phi \in \Phi_{R}$, there exists a unique corresponding structural parameter point. In other words, $\mathcal{A}_{R, 0}(\phi)$ is singleton-valued at almost every $\phi \in \Phi_{R}$. In addition, the case where $\Phi_{R}=\Phi$, i.e. the imposed identifying assumptions are not observationally restrictive, is what RWZ refer to as exact identification. In contrast, the definition of local identification says that, if there are multiple observationally equivalent structural parameter points, they must be far apart. This implies that for almost every $\phi \in \Phi_{R}$, if $\mathcal{A}_{R, 0}(\phi)$ is not singleton, it consists of isolated points. In Proposition 2 below, we characterize a class of locally identified SVARs. For this class of SVARs, the space of reduced-form parameters $\Phi$ can be partitioned into three subsets. The first, of positive measure, contains parameters for which the model is locally- but not globally-identified; the second, of positive measure, on which there is no structural parameter satisfying the identifying assumption (i.e., $\mathcal{A}_{R, 0}(\phi)$ is empty); and the third, of measure zero, on which the model is globally identified. This feature of locally identified SVARs stands in contrast to exactly identified SVARs and globally and over-identified SVARs, where the mapping from the reduced-form parameter space $\Phi$ to structural parameters that satisfy the identifying restrictions is guaranteed to be either singleton-valued or empty at almost every $\phi \in \Phi$.

## II. 2 Normalization, sign, zero and non-zero identifying restrictions

This section introduces the types of identifying restriction considered in this paper. We begin with sign normalization restrictions, and then move to zero, non-zero, and sign restrictions.

## Sign Normalization restrictions

Following Waggoner and Zha (2003) and Hamilton et al. (2007), and in line with RWZ and Giacomini and Kitagawa (2020), we impose sign normalization restrictions on the structural shocks. Specifically, we restrict the diagonal elements of $A_{0}$ to be non-negative.

$$
\begin{equation*}
\operatorname{diag}\left(Q^{\prime} \Sigma_{t r}^{-1}\right) \geq 0 \tag{7}
\end{equation*}
$$

Under these assumptions, a unit positive change in a structural shock can be interpreted as a one standard-deviation ceteris paribus positive shock in the corresponding endogenous variable.

## Zero and non-zero equality restrictions

While sign normalization restrictions on the diagonal elements of $A_{0}$ restrict the set of admissible structural matrices, they are not enough to obtain identification. The standard approach in the literature is to impose equality restrictions either on the structural parameters or particular linear and non-linear functions of them. ${ }^{1}$

Following RWZ, we represent identifying restrictions as restrictions on the reduced-form parameters $\phi$ and the column vectors $\left(q_{1}, q_{2}, \ldots, q_{n}\right)$ of the orthogonal matrix $Q$.

$$
\begin{align*}
\left((i, j) \text {-th element of } A_{0}^{-1}\right)=0 & \Longleftrightarrow\left(e_{i}^{\prime} \Sigma_{t r}\right) q_{j}=0,  \tag{8}\\
\left((i, j) \text {-th element of } A_{0}\right)=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} e_{j}\right)^{\prime} q_{i}=0,  \tag{9}\\
\left((i, j) \text {-th element of } A_{l}\right)=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} B_{l} e_{j}\right)^{\prime} q_{i}=0,  \tag{10}\\
\left((i, j) \text {-th element of } C I R^{\infty}\right)=0 & \Longleftrightarrow\left[e_{i}^{\prime} \sum_{h=0}^{\infty} C_{h}(B) \Sigma_{t r}\right] q_{j}=0,  \tag{11}\\
\left((i, j) \text {-th element of } A_{0}^{-1}\right)=c & \Longleftrightarrow\left(e_{i}^{\prime} \Sigma_{t r}\right) q_{j}=c, \tag{12}
\end{align*}
$$

(linear restriction between $(i, j)$-th

$$
\begin{equation*}
\text { and } \left.(h, k) \text {-th elements of } A_{0}^{-1}\right) \Longleftrightarrow\left(e_{i}^{\prime} \Sigma_{t r}\right) q_{j}-d\left(e_{h}^{\prime} \Sigma_{t r}\right) q_{k}=c, \tag{13}
\end{equation*}
$$

where $e_{i}$ is the $i$-th column of the identity matrix $I_{n}$, and $c$ and $d$ are known non-zero scalars. Eq. (8) and Eq. (9) cover short-run identifying restrictions including the causal ordering restrictions of Sims (1980) and Bernanke (1986). Eq. (10) corresponds to restrictions that exclude some of the right-hand side variables in the structural equations. Eq. (11) corresponds to long-run identifying restriction as considered in Blanchard and Quah (1989). These first four equality re-

[^2]strictions were considered in RWZ and Giacomini and Kitagawa (2020), but the remaining two were not. The additional restrictions we allow are non-zero equality restrictions (also referred to as nonhomogeneous restrictions) and cross-equation restrictions on the structural parameters and impulse responses. As we clarify in Section III, these last two types of restriction drive a departure from global identification to local identification.

We represent these equality restrictions by

$$
\begin{align*}
\mathbf{F}(\phi, Q) & \equiv\left(\begin{array}{cccc}
F_{11}(\phi) & F_{12}(\phi) & \cdots & F_{1 n}(\phi) \\
F_{21}(\phi) & F_{22}(\phi) & \cdots & F_{2 n}(\phi) \\
\vdots & \vdots & \ddots & \vdots \\
F_{n 1}(\phi) & F_{n 2}(\phi) & \cdots & F_{n n}(\phi)
\end{array}\right)\left(\begin{array}{c}
q_{1} \\
q_{2} \\
\vdots \\
q_{n}
\end{array}\right)-\left(\begin{array}{c}
c_{1} \\
c_{2} \\
\vdots \\
c_{n}
\end{array}\right)=\mathbf{0} \\
& \equiv \mathbf{F}(\phi) \operatorname{vec} Q-\mathbf{c}=\mathbf{0} \tag{14}
\end{align*}
$$

where $F_{i j}(\phi), 1 \leq i \leq j \leq n$, is a matrix of dimension $f_{i} \times n$, which depends only on the reducedform parameters $\phi=(B, \Sigma)$. The dimension of $\mathbf{F}(\phi)$ is $f \times n^{2}$, where $f=f_{1}+\cdots+f_{n}$ denotes the total number of restrictions imposed. We allow $f_{i}=0$ for some $i$, in which case the $i$-th block row in $\mathbf{F}(\phi)$ is null. Finally, vec $Q \equiv\left(q_{1}^{\prime}, \ldots q_{n}^{\prime}\right)^{\prime}$ is the vectorization of $Q$, and $\mathbf{c} \equiv\left(c_{1}^{\prime} \ldots, c_{n}^{\prime}\right)^{\prime}$ is a vector of known constants with length $f$, where each $c_{i}$ is a $f_{i} \times 1$ vector.

If $F_{i j}(\phi)=0$ for any $i \neq j$, there are no cross equation restrictions. If $c_{i}=0$ for all $i$, then only zero restrictions are imposed. This represention of the identifying restrictions is in line with Lütkepohl (2006) and Bacchiocchi and Lucchetti (2018), both of which allow non-homogeneous and across-shock restrictions. We provide the following formal definitions.

Definition 3 (Recursive restrictions). The restrictions are said to be recursive if $F_{i j}(\phi)=0$ for $j>i$, and $f_{i}=n-i$, for $i=1, \ldots, n$.

Definition 4 (Homogeneous and non-homogeneous restrictions). The restrictions are said to be homogeneous if $\mathbf{c}=\mathbf{0}$ and non-homogeneous if $\mathbf{c} \neq \mathbf{0}$.

Defined in this way, recursive restrictions pin down a unique ordering of the variables with Eq. (14) becoming a lower-triangular block matrix. Otherwise, our framework allows for the ordering of variables to be non-unique. Even with the order of variables fixed, if the restrictions include across-shock restrictions, then Eq. (14) allows for a multiple block-matrix representation. The general identification results of this section are valid independent of how the variables are ordered or how the imposed restrictions are represented within Eq. (14), unless the recursive structure is assumed explicitly.

## Sign restrictions

In addition to equality restrictions, sign restrictions can be imposed on impulse responses or structural parameters. These sign restrictions can be seen as additional constraints on the columns of the $Q$ matrix. Suppose we impose $s_{h, i} \leq n$ number of sign restrictions on the impulse responses to $i$-th shock at $h$-th horizon. They can be expressed as

$$
\begin{equation*}
S_{h, i}(\phi) q_{i} \geq \mathbf{0}, \tag{15}
\end{equation*}
$$

where $S_{h, i} \equiv D_{h, i} C_{h}(B) \Sigma_{t r}$ is a $s_{h, i} \times n$ matrix, $D_{h, i}$ is the $s_{h, i} \times n$ signed selection matrix, which indicates by $1(-1)$ the impulse responses whose signs are retricted to being positive (negative), and $C_{h}(B)$ is from the definition of an impulse response Eq. (3). The inequality in Eq. (15) is component-wise. Sign restrictions on structural parameters are linear inequality constraints on the columns of the matrix $Q$, so can also be accommodated. Stacking all the $S_{h, i}$ matrices involving sign restrictions on $q_{i}$ at different horizons into a matrix $S_{i}$, we have

$$
\begin{equation*}
S_{i}(\phi) q_{i} \geq \mathbf{0} \tag{16}
\end{equation*}
$$

We represent the set of all sign restrictions by

$$
\begin{equation*}
\mathbf{S}(\phi, Q) \geq \mathbf{0} \tag{17}
\end{equation*}
$$

## Admissible structural parameters and identified set

Given identifying restrictions of the form introduced above, we hereafter let $R$ be the collection of restrictions $\left\{\mathbf{F}(\phi, Q)=0, \mathbf{S}(\phi, Q) \geq 0, \operatorname{diag}\left(Q^{\prime} \Sigma_{t r}^{-1}\right) \geq 0\right\}$, or $R=(F, S)$ for short. We call $A=\left(A_{0}, A_{+}\right)$admissible if it satisfies $R$. The set of all these admissible structural parameters can be represented by
$\mathcal{A}_{R}(\phi) \equiv\left\{\left(A_{0}, A_{+}\right)=\left(Q^{\prime} \Sigma_{t r}^{-1}, Q^{\prime} \Sigma_{t r}^{-1} B\right): Q \in \mathcal{O}(n), \mathbf{F}(\phi, Q)=\mathbf{0}, \mathbf{S}(\phi, Q) \geq \mathbf{0}, \operatorname{diag}\left(Q^{\prime} \Sigma_{t r}^{-1}\right) \geq 0\right\}$.

The projection of $\mathcal{A}_{R}(\phi)$ for $A_{0}$ gives $\mathcal{A}_{R, 0}(\phi)$ as defined in Eq. (6). The identified set for $Q$ is defined as the set of admissible orthogonal matrices given the reduced-form parameters:

$$
\mathcal{Q}_{R}(\phi) \equiv\left\{Q \in \mathcal{O}(n): \mathbf{F}(\phi, Q)=\mathbf{0}, \mathbf{S}(\phi, Q) \geq \mathbf{0}, \quad \operatorname{diag}\left(Q^{\prime} \Sigma_{t r}^{-1}\right) \geq 0\right\}
$$

The objects of interest may also include transformations of structural parameters such as impulse
response functions. We denote a scalar parameter of interest by $\eta=\eta(\phi, Q)$ and define its identified set as

$$
I S_{\eta}(\phi) \equiv\left\{\eta(\phi, Q): Q \in \mathcal{Q}_{R}(\phi)\right\}
$$

When $\eta(\phi, Q)$ is a restriction on an impulse response

$$
\eta(\phi, Q)=I R_{i j}^{h}=e_{i}^{\prime} C_{h}(B) \Sigma_{t r} Q e_{j} \equiv c_{i h}^{\prime}(\phi) q_{j},
$$

where $I R_{i j}^{h}$ is the $(i, j)$-th element of $I R^{h}$ and $c_{i h}^{\prime}(\phi)$ is the $i$-th row of $C_{h}(B) \Sigma_{t r}$. Similar definitions can be obtained for restrictions on long-run impulse responses and cumulative impulse responses. If, instead, the object of interest is the $(i, j)$-th element of $A_{l}$, then $\eta(\phi, Q)=e_{j}^{\prime}\left(\Sigma_{t r}^{-1} B_{l}\right)^{\prime} q_{i}$, with $B_{0}=I_{n}$.

When $A$ is globally identified, $I S_{\eta}(\phi)$ is a singleton for almost every $\phi \in \Phi_{R}$. If $A$ is only locally identified, $I S_{\eta}(\phi)$ can be a set of multiple isolated points generated by observationally equivalent structural parameters. Local identification can be certainly viewed as a special case of set identification, although it is not covered by standard set identification analysis where the identified set is typically an interval or a set with positive Lebesgue measure.

## II. 3 Conditions for local identification

This section presents conditions for global and local identification when the identifying restrictions are equality restrictions in the form Eq. (14). In the case of local identification, we present an analytical characterization of the number of observationally equivalent structural parameter values.

We begin with a modified version of the well known condition for global identification developed in Theorem 7 of RWZ, as recently discussed in Bacchiocchi and Kitagawa (2020). ${ }^{2}$ This condition for global identification acts as a reference point in our discussion of local identification.

Proposition 1 (Necessary and Sufficient condition for global identification, RWZ and Bacchiocchi and Kitagawa (2020)). Consider an SVAR with identifying restrictions of the form Eq. (8) - Eq. (13) collected in $\boldsymbol{F}(\phi, Q)$. Assume $F_{i j}(\phi)=0$ for $i \neq j$, and $\boldsymbol{c}=\boldsymbol{0}$.

The SVAR is globally identified at $A=\left(A_{0}, A_{+}\right) \in \mathcal{A}_{R}$ if and only if the following conditions hold at $\phi$ implied by $A$ :

## 1. It holds

$$
\begin{equation*}
\operatorname{rank}\left(\left(F_{11}(\phi)^{\prime}, \tilde{\sigma}_{1}\right)\right)=n \tag{18}
\end{equation*}
$$

[^3]2. Let $q_{1}$ be a unit length vector satisfying $F_{11}(\phi) q_{1}=0$ and the sign normalization restriction, which is unique under Eq. (18). For $i=2, \ldots, n$
\[

$$
\begin{equation*}
\operatorname{rank}\left(\left(F_{i i}(\phi)^{\prime}, q_{1}, \ldots, q_{i-1}, \tilde{\sigma}_{i}\right)\right)=n \tag{19}
\end{equation*}
$$

\]

hold, where the orthonormal vectors $q_{2}, \ldots, q_{n}$ solve

$$
\begin{equation*}
\left(F_{i i}(\phi)^{\prime}, q_{1}, \ldots, q_{i-1}\right)^{\prime} q_{i}=\mathbf{0} \tag{20}
\end{equation*}
$$

sequentially, and satisfy the sign normalization restrictions.
This proposition characterizes a boundary separating cases where an SVAR is globally identified and cases where it is not guaranteed to be globally identified. In what follows, we consider departures from this proposition's conditions for global identification, and show implications for local identification and the failure of global identification. In particular, we allow $F_{i j}(\phi)$ to be nonzero for some $i \neq j$ and/or $\mathbf{c} \neq \mathbf{0}$ by including restrictions of the form Eq. (12)- Eq. (13). With this expanded set of identifying restrictions, Proposition 2 derives a rank condition that is necessary and sufficient for local identification. Lütkepohl (2006) and Bacchiocchi and Lucchetti (2018) provide similar conditions for local identification in a setting that is less general in terms of the kind of restrictions that can be imposed. Their rank condition is expressed in terms of the structural parameter matrices $A$, while our Proposition 2 presents the rank condition in terms of the coefficient matrix of the equality restrictions $\boldsymbol{F}(\phi)$ and the orthogonal matrix $Q$. We define $\operatorname{Chol}(\cdot)$ to be the Cholesky factor of $(\cdot)$ and $g: \mathbb{R}^{(n+m) n} \rightarrow \mathbb{R}^{n+n^{2} p} \times \Omega \times \mathcal{O}(n)$, to be the function mapping structural to reduced-form parameters and the admissible orthogonal matrix.

Proposition 2 (Rank condition - necessary and sufficient condition for local identification). Consider an SVAR with equality restrictions of the form Eq. (8) - Eq. (13) collected in $\boldsymbol{F}(\phi, Q)$. Let $\tilde{D}_{n}$ be the $n^{2} \times n(n-1) / 2$ full-column rank matrix such that for any $n(n-1) / 2$-dimensional vector $v, \tilde{D}_{n} v \equiv$ vec $(H)$ holds, where $H$ is an $n \times n$ skew-symmetric matrix satisfying $H=-H^{\prime}$ (see Appendix $D$ for the specific construction of $\tilde{D}_{n}$ for $n=2,3,4$ ).
(i) The SVAR is locally identified at $A=\left(A_{0}, A_{+}\right) \in \mathcal{A}_{R}$, i.e., there exists an open neighborhood about $A$ containing no other observationally equivalent structural parameter point, if and only if

$$
\begin{equation*}
\operatorname{rank}\left[\boldsymbol{F}(\phi)\left(I_{n} \otimes Q\right) \tilde{D}_{n}\right]=n(n-1) / 2 \tag{21}
\end{equation*}
$$

holds, where the reduced-form parameters $\phi=(B, \Sigma) \in \Phi$ and the orthogonal matrix $Q \in \mathcal{O}(n)$ are such that $(B, \Sigma, Q)=g\left(A_{0}, A_{+}\right)=\left(A_{0}^{-1} A_{+}, A_{0}^{-1} A_{0}^{-1 \prime}, \operatorname{Chol}\left(A_{0}^{-1} A_{0}^{-1 \prime}\right)^{\prime} A_{0}^{\prime}\right)$. Hence, a necessary condition for the rank condition Eq. (21) is $f=\sum_{i=1}^{n} f_{i} \geq n(n-1) / 2$.
(ii) Let $\mathcal{K}$ be the set of structural parameters in $\mathcal{A}_{R}$ satisfying the rank condition of Eq. (21),

$$
\mathcal{K} \equiv\left\{A \in \mathcal{A}_{R}: \operatorname{rank}\left[\boldsymbol{F}(\phi)\left(I_{n} \otimes Q\right) \tilde{D}_{n}\right]=n(n-1) / 2\right\}
$$

Either $\mathcal{K}$ is empty or the complement of $\mathcal{K}$ in $\mathcal{A}_{R}$ is of measure zero.
Proof. See Appendix C.
Statement (i) of this proposition provides a necessary and sufficient condition for local identification at a given $A \in \mathcal{A}_{R}$ in the form of a rank condition for a matrix that is a function of $A$, i.e., $(\phi, Q)$ is a function of $A$. Eq. (21) as stated is of limited practical use since the true $A$ is generally unknown, which means that verifying Eq. (21) is infeasible.

Statement (ii) of this proposition makes the rank condition Eq. (21) useful by showing that it holds either nowhere or almost everywhere in the parameter space $\mathcal{A}_{R}$. This means that, similar to the proposals following Theorem 3 in RWZ and Theorem 1 in Bacchiocchi and Lucchetti (2018), one can assess local identification by randomly generating structural parameters $A \in \mathcal{A}_{R}$ and checking whether the rank condition holds or not. Specifically, we can consider drawing reducedform parameters $\phi \in \Phi_{R}$ from its prior or posterior and solving a constrained non-linear optimization problem of the form ${ }^{3}$

$$
\begin{align*}
& \arg \min _{Q \in \mathbb{R}^{n^{2}}}(\mathbf{F}(\phi) \operatorname{vec} Q-\mathbf{c})^{\prime}(\mathbf{F}(\phi) \operatorname{vec} Q-\mathbf{c})  \tag{22}\\
& \text { s.t. } \operatorname{diag}\left(Q^{\prime} \Sigma_{t r}^{-1}\right) \geq 0, \mathbf{S}(\phi, Q) \geq \mathbf{0} \text { and } Q^{\prime} Q=I_{n}
\end{align*}
$$

If the value of the optimization is zero, then the obtained $Q$ is an admissible orthogonal matrix at the given $\phi$. If such an admissible $Q$ satisfies the rank condition in Eq. (21), then the SVAR is locally identified at $(\phi, Q)$. If the rank condition is not met, the SVAR is not locally identified at $(\phi, Q)$. Proposition 2 (ii) says that only one of the two possibilities occurs with positive measure, while the other has zero measure. Hence, by checking the rank condition at a few parameter values drawn from a probability distribution supporting $\mathcal{A}_{R}$ or $\Phi_{R}$, we can learn whether the rank condition holds nowhere or almost everywhere on the space of structural parameters. Confirming the latter can be seen as a strong support for local identification holding at the true $A$, unless the true structural parameter value is believed to belong to the null set in the parameter space.

Allowing only recursive identifying restrictions, the next proposition provides a simple necessary and sufficient condition for the rank condition of Proposition 2 (i). It extends to local identification

[^4]the condition for global identification presented in Proposition 1.
Proposition 3 (Necessary and sufficient condition for local identification in recursive SVARs). Consider an SVAR with recursive identifying restrictions of the form Eq. (14). Let $\tilde{F}_{i i}(\phi)=F_{11}(\phi)$ for $i=1$, and
\[

$$
\begin{equation*}
\tilde{F}_{i i}(\phi)=\left(F_{i i}^{\prime}(\phi), q_{1}, \ldots, q_{i-1}\right)^{\prime} \tag{23}
\end{equation*}
$$

\]

for $i=2, \ldots, n$, where $q_{1}, \ldots, q_{i}$ are the first $i$ column vectors of $Q \in \mathcal{O}(n)$ satisfying the equality restrictions $\boldsymbol{F}(\phi)$ vec $Q-\mathbf{c}=\mathbf{0}$ given $\phi \in \Phi_{R}$. The rank condition of Eq. (21) holds at $(\phi, Q)$ if and only if $\operatorname{rank}\left(\tilde{F}_{i i}(\phi)\right)=n-1$ holds for all $i=1, \ldots, n$.

Proof. See Appendix C.
Since the rank condition of Proposition 2 (i) is necessary and sufficient for local identification, the condition shown in Proposition 3 is also a necessary and sufficient for local identification for SVARs under recursive identifying restrictions. Moreover, the claim of Proposition 2 (ii) carries over to the setting of Proposition 3, so knowing that the condition shown in Proposition 3 holds at a few $\phi \in \Phi_{R}$ drawn from its prior or posterior allows us to conclude local identification holds almost everywhere in the parameter space. The condition in Proposition 3 exploits sequential determination of $q_{i}, i=1, \ldots, n$, given $\phi$, so checking it does not require nonlinear optimization for $Q$.

The proof of Proposition 3 leads to the following corollary showing a necessary and sufficient condition for the local identification of impulse response to a particular shock.

Corollary 1 (Sufficient condition for local identification of the $j$-th shock). Under the assumptions of Proposition 3, the impulse responses for the $j$-th structural shock, $1 \leq j \leq n$, are locally identified at the parameter point $A=\left(A_{0}, A_{+}\right) \in \mathcal{A}_{R}$ if and only if $\operatorname{rank}\left(\tilde{F}_{i i}(\phi)\right)=n-1$ holds for all $i=1, \ldots, j$.

## II. 4 The number of observationally equivalent parameter points

The results presented so far are silent about how many observationally equivalent structural parameter point there are. As the next proposition shows, our constructive identification argument through the orthogonal matrix $Q$ allows us to characterize the number of observationally equivalent parameter points.

Proposition 4 (Number of locally identified points). Consider an SVAR with equality restrictions of the form Eq. (8)-Eq. (13) collected in $\boldsymbol{F}(\phi, Q)=\mathbf{0}$. Given $\phi \in \Phi$ and provided that the rank condition in Eq. (21) is met, the number of admissible $Q$ matrices ( $Q$ matrices solving $\boldsymbol{F}(\phi, Q)=\mathbf{0}$ )
is zero or finite. In particular, if the equality identifying restrictions are recursive, the number of admissible $Q$ matrices is at most $2^{n}$. If the equality identifying restrictions are non-recursive, the number of admissible $Q$ matrices is at most $2^{n(n+1) / 2}$.

Proof. See Appendix C.
The proposition provides an upper bound for the number of locally identified observationally equivalent parameter points. It corresponds to the maximal number of modes that the likelihood of the structural parameters can have. The maximum number of observationally equivalent structural parameters is considerably lower when the SVAR is identified through recursive equality restrictions rather than non-recursive restrictions. The intuition for this result is that, if the identification of the columns of $Q$ can be performed recursively, the equations concerning the orthogonality conditions among the columns of $Q$ are linear, rather than quadratic.

In comparison to the exact (global) identification case of RWZ and Proposition 1, Proposition 4 highlights that non-homogenous restrictions $(\mathbf{c} \neq \mathbf{0})$ lead to the possibility that, given $\phi \in \Phi$, (i) an admissible $Q$ does not exist, or (ii) the admissible $Q$ is no longer unique. Adding sign restrictions to the sign normalization restrictions can reduce the number of admissible $Q$ 's, but cannot generally guarantee uniqueness of the admissible $Q$ 's. Section II. 5 below illustrates the transition from exact global identification to local identification through a simple example.

## II. 5 The geometry of identification

We present an intuitive geometric exposition for why the introduction of nonhomogeneous restrictions Eq. (12) and/or across-shock restrictions Eq. (13) can lead to local identification. This exposition also provides intuition for the number of local identified parameter points shown in Proposition 4. Appendix A provides the algebraic analysis behind our geometric discussion.

To make exposition as simple as possible, consider a bivariate VAR with a single non-homogeneous identifying restriction imposed on the structural parameter matrix:

$$
\begin{equation*}
\left(A_{0}\right)_{[1,1]}^{-1}=c \Longleftrightarrow\left(e_{1}^{\prime} \Sigma_{t r}\right) q_{1}=c \tag{24}
\end{equation*}
$$

where $c>0$ is a known (positive) scalar and $e_{1}$ is the first column of $I_{2}$. Denoting the first column of $\Sigma_{t r}^{\prime}=\left(\begin{array}{cc}\sigma_{1,1} & \sigma_{2,1} \\ 0 & \sigma_{2,2}\end{array}\right)$ by $\sigma_{1}=\left(\sigma_{1,1}, 0\right)^{\prime}$, this identifying restriction can be written as $\sigma_{1}^{\prime} q_{1}=c$. Hence, given $\phi, q_{1}$ must satisfy the two equations,

$$
\left\{\begin{array}{l}
\sigma_{1}^{\prime} q_{1}=c \\
q_{1}^{\prime} q_{1}=1 .
\end{array}\right.
$$

Figure 1: Identification of $q_{1}$ in the bivariate SVAR with non-zero restriction.


Notes: The vertical red line represents the non-zero restriction $\left(A_{0}\right)_{[1,1]}^{-1}=c$. The two black arrows represent the identified vectors $q_{1}^{(1)}$ and $q_{1}^{(2)}$.

Figure 1 depicts these two constraints. Letting the x -axis correspond to the vector $\sigma_{1}$, the set of $q_{1}$ vectors satisfying the first constraint is a vertical line whose location is determined by $\sigma_{1}$ and c. The second constraint imposes that $q_{1}$ lies on the unit circle. Points at the intersection of the vertical line and the unit circle, if any exist, are solutions to this system of equations.

When the imposed restriction is a zero restriction $(c=0)$, the vertical line passes through the origin and intersects the circle at two points. The two solutions for $q_{1}, q_{1}^{(1)}$ and $q_{1}^{(2)}$ are symmetric across the origin, and the sign normalization restriction Eq. (7) is guaranteed to rule one of them out (see Appendix A for details). Thus, the first column of $Q$ is globally identified.

The vertical line in Figure 1 corresponds to a non-zero restriction ( $c>0$ ). If the vertical line is perfectly tangent to the unit circle, we continue to have global identification. Otherwise, there are two distinct solutions for $q_{1}$, as shown in Figure 1. Compared to the case where $c=0$, a crucial difference is that there are some values of $\phi$ and $c$ where the sign normalization restriction cannot rule out one solution. In this case, they are both admissible and the first column of $Q$ is locallybut not globally-identified. ${ }^{4}$

The second column of $Q$, i.e. the unit-length vector $q_{2}$, can be pinned down through its ortho-

[^5]gonality with $q_{1}$
\[

\left\{$$
\begin{align*}
q_{2}^{\prime} q_{1} & =0  \tag{25}\\
q_{2}^{\prime} q_{2} & =1
\end{align*}
$$\right.
\]

If $q_{1}$ is only locally identified with two admissible vectors $q_{1}^{(1)}$ and $q_{1}^{(2)}$, Eq. (25) needs to be solved given both. Solving the system when $q_{1}=q_{1}^{(1)}$ provides two solutions for $q_{2}$ that are depicted in the left panel of Figure 2. As the two solutions mirror each other across the origin, only one will satisfy the sign normalization restriction for the second shock. A similar picture is obtained when $q_{1}=q_{1}^{(2)}$ (the right panel of Figure 2), and here too one of the solutions for $q_{2}$ can be ruled out by the sign normalization restriction.

To summarize, an equality restriction with $c>0$ leads to local but non-global identification for $q_{1}$, and there are then two admissible $Q$ matrices, $Q_{1}=\left[q_{1}^{(1)}, q_{2}^{(1)}\right]$ and $Q_{2}=\left[q_{1}^{(2)}, q_{2}^{(2)}\right]$ given $\phi$. This implies that both $A_{0}=Q_{1}^{\prime} \Sigma_{t r}^{-1}$ and $A_{0}=Q_{2}^{\prime} \Sigma_{t r}^{-1}$ are admissible. In this example, we obtain two observationally equivalent $Q$ matrices, which is consistent with the upper bound on the number of observationally equivalent $Q$ matrices in Proposition 4.

For a specific numerical illustration, let the bivariate VAR be characterized by constants that are zero and a single lag with reduced-form parameters

$$
B_{1}=\left(\begin{array}{cc}
0.8 & -0.2 \\
0.1 & 0.6
\end{array}\right), \quad \Sigma=\left(\begin{array}{cc}
0.49 & -0.14 \\
-0.14 & 0.13
\end{array}\right), \quad \Sigma_{t r}=\left(\begin{array}{cc}
0.7 & 0 \\
-0.2 & 0.3
\end{array}\right)
$$

and consider imposing restriction $\left(A_{0}\right)_{[1,1]}^{-1}=0.5 \Longleftrightarrow\left(e_{1}^{\prime} \Sigma_{t r}\right) q_{1}=0.5$. Following Eq. (56) and Eq. (61) - Eq. (62) in Appendix A, we calculate the two admissible matrices $Q_{1}$ and $Q_{2}$

$$
Q_{1}=\left(\begin{array}{cc}
0.714 & -0.700 \\
0.700 & 0.714
\end{array}\right) \quad \text { and } \quad Q_{2}=\left(\begin{array}{cc}
0.714 & 0.700 \\
-0.700 & 0.714
\end{array}\right)
$$

with associated admissible $A_{0}$ matrices

$$
A_{0}^{(1)}=\left(\begin{array}{cc}
1.687 & 2.333 \\
-0.320 & 2.381
\end{array}\right) \quad \text { and } \quad A_{0}^{(2)}=\left(\begin{array}{cc}
0.354 & -2.333 \\
1.680 & 2.381
\end{array}\right)
$$

Based on these structural parameter values, Figure 3 shows the impulse response of $y_{t}=\left(y_{1 t}, y_{2 t}\right)^{\prime}$ to the structural shocks $\varepsilon_{1 t}$ and $\varepsilon_{2 t}$. Despite the simplicity of this example, it clearly illustrates the extent to which conclusions depend on the choice of observationally equivalent $Q$ matrices.

Figure 2: Identification of $q_{2}$ in the bivariate SVAR with non-zero restriction.


Case 1: $q_{1}^{(1)}$


Case 2: $q_{1}^{(2)}$

Notes: The left panel shows the identification of the $q_{2}^{(1)}$ and $q_{2}^{(2)}$ vectors (in blue), conditional on the identified $q_{1}^{(1)}$ (in black). Similarly, the right panel shows the identification of the $q_{2}^{(1)}$ and $q_{2}^{(2)}$ vectors (in blue), conditional on the identified $q_{1}^{(2)}$ (in black).

## III Locally identified SVARs: some examples

Hamilton et al. (2007) discuss local identification as a normalization problem. As shown in the previous section, in the presence of non-homogeneous equality restrictions Eq. (12) and/or acrossshock restrictions Eq. (13), proper sign normalization restrictions are not enough to resolve the issue of local identification in SVARs. The examples below illustrate that this issue is of practical relevance.

## III. 1 Calibrated identifying restrictions

One strategy employed in the literature is to calibrate some parameters instead of estimating them. Calibration can be viewed as imposing non-homogeneous restrictions. Depending on whether or not the calibrated parameters are normalized by the structural error variance, these non-homogeneous restrictions can deliver either global identification or local identification. The following two examples clarify this.

Following Bacchiocchi et al. (2005), consider a trivariate SVAR analyzing the impact of privatization policies on the real economy. Let $Y_{t}=\left(p_{t}, g_{t}, x_{t}\right)^{\prime}$ denote privatization proceeds, public spending and real output, respectively. In line with their results, suppose that the government in the model adopts an automatic rule to redirect privatization proceeds to public spending.

Figure 3: Impulse response functions related to the locally identified SVAR discussed in Section II. 5.


Notes: The bivariate SVAR is characterized by a non-zero restriction. In solid lines we report the IRFs obtained through $Q_{1}=\left(q_{1}^{(1)}, q_{2}^{(1)}\right)$ while in dashed lines those obtained through $Q_{2}=\left(q_{1}^{(2)}, q_{2}^{(2)}\right)$.

With the dynamics concentrated out, the model is specified as

$$
\begin{align*}
a_{11} p_{t}+a_{12} g_{t} & =\varepsilon_{t}^{p} \\
a_{21} p_{t}+a_{22} g_{t} & =\varepsilon_{t}^{g}  \tag{26}\\
a_{31} p_{t}+a_{23} g_{t}+a_{33} x_{t} & =\varepsilon_{t}^{x},
\end{align*}
$$

where the vector $\varepsilon_{t}=\left(\varepsilon_{t}^{p}, \varepsilon_{t}^{g}, \varepsilon_{t}^{x}\right)^{\prime}$ denotes the three structural shocks with normalized variances $E\left(\varepsilon_{t} \varepsilon_{t}^{\prime}\right)=I_{3}$. In addition, we assume that privatization shocks are transmitted to public spending through a fixed mechanism governed by the nonzero parameter $c$, i.e. $\left(A_{0}\right)_{[2,1]}^{-1}=c \Longleftrightarrow\left(e_{2}^{\prime} \Sigma_{t r}\right) q_{1}=$
c. This set of restrictions can be written as

$$
\begin{align*}
a_{13}=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime} q_{1}=0 \\
\left(A_{0}\right)_{[2,1]}^{-1}=c & \Longleftrightarrow\left(e_{2}^{\prime} \Sigma_{t r}\right) q_{1}=c  \tag{27}\\
a_{23}=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime} q_{2}=0 .
\end{align*}
$$

We have $f_{1}=2, f_{2}=1$ and $f_{3}=0$, and Proposition 3 guarantees that the model is locally identified. The next proposition shows a condition that leads to failure of global identification and provides a concrete characterization of the number of locally identified parameter points.

Proposition 5. Let the $j$-th column vector of $\Sigma_{t r}^{\prime}$ be denoted by $\sigma_{j}=\left(\sigma_{j, 1}, \sigma_{j, 2}, \sigma_{j, 2}\right)^{\prime}, j=1,2,3$. In the privatization policy SVAR Eq. (26) subject to the restrictions Eq. (27) with $c \neq 0$, if $\sigma_{2,1}^{2}+\sigma_{2,2}^{2}>c^{2}$ holds and the half-space $\left\{q_{1}: \sigma_{1}^{\prime} q_{1} \geq 0\right\}$ representing the sign normalization restriction for $q_{1}$ contains the slice of sphere $\left\{q_{1}:\left\|q_{1}\right\|=1, \sigma_{2}^{\prime} q_{1}=c\right\}$, then global identification fails and we have two distinct locally identified structural parameter points.

Proof. See Appendix C.
The non-homogeneous identifying restriction in Eq. (26) constrains the value of the structural parameters in the SVAR representation where the structural error variance is normalized to unity. It is worth noting a contrasting case such that imposing non-homogeneous restrictions without the normalization of the structural error variance yields global identification.

To illustrate this point, consider Blanchard and Perotti's (2002) seminal work on the dynamic effect of fiscal policy shocks on the real economy. They consider a three variable SVAR which, using the terminology of Amisano and Giannini (1997) and Lütkepohl (2006), has an AB-SVAR representation ${ }^{5}$ where $Y_{t}=\left(T_{t}, G_{t}, X_{t}\right)^{\prime}$ consists of the logarithms of quarterly taxes, spending, and GDP, all in the real per capita terms. The equations of the model are

$$
\begin{aligned}
t_{t} & =a_{1} x_{t}+a_{2} \epsilon_{t}^{g}+\epsilon_{t}^{t} \\
g_{t} & =b_{1} x_{t}+b_{2} \epsilon_{t}^{t}+\epsilon_{t}^{g} \\
x_{t} & =c_{1} t_{t}+c_{2} g_{t}+\epsilon_{t}^{x}
\end{aligned}
$$

where $u_{t}=\left(t_{t}, g_{t}, x_{t}\right)^{\prime}$ is the vector of reduced-form residuals and $\epsilon_{t}^{t}, \epsilon_{t}^{g}$ and $\epsilon_{t}^{x}$ are mutually uncorrelated structural shocks with unknown variances $\sigma_{t}^{2}, \sigma_{g}^{2}$ and $\sigma_{y}^{2}$, respectively.

In addition to the zero restrictions already incorporated above, the authors propose three further restrictions: (1) $b_{1}=0$, (2) $a_{1}=2.08$ and, (3) either $a_{2}=0$ or $b_{2}=0$. Combining these restrictions

[^6]with $b_{2}=0$ (the situation is very similar for the alternative $a_{2}=0$ ), we can write the model as
\[

\left($$
\begin{array}{ccc}
\alpha_{11} & 0 & 0 \\
\alpha_{21} & \alpha_{22} & \alpha_{23} \\
\alpha_{31} & \alpha_{32} & \alpha_{33}
\end{array}
$$\right)\left($$
\begin{array}{c}
t_{t} \\
g_{t} \\
x_{t}
\end{array}
$$\right)=\left($$
\begin{array}{c}
\varepsilon_{t}^{t} \\
\varepsilon_{t}^{g} \\
\varepsilon_{t}^{x}
\end{array}
$$\right) .
\]

where the new structural shocks $\tilde{\varepsilon}_{t}=\left(\tilde{\varepsilon}_{t}^{t}, \tilde{\varepsilon}_{t}^{g}, \tilde{\varepsilon}_{t}^{x}\right)^{\prime}$ have normalized variance $E\left(\tilde{\varepsilon}_{t} \tilde{\varepsilon}_{t}^{\prime}\right)=I_{3}$, and the $\alpha_{i j}$ parameters are defined as

$$
\begin{array}{llll}
\alpha_{11}=1 / \sigma_{g} & & \\
\alpha_{21}=-2.08 / \sigma_{t} & , & \alpha_{22}=1 / \sigma_{t} & , \\
\alpha_{31}=-a_{1} / \sigma_{x} & , & \alpha_{32}=-c_{2} / \sigma_{x} & ,
\end{array} \alpha_{33}=1 / \sigma_{x} .
$$

The model imposes two restrictions on the first equation, one restriction on the second and no restriction in the third. The two zero restrictions in the first equation identify the orthogonal vector $q_{1}$, while the restriction on the second row parameters take the form $\alpha_{21}=-2.08 \alpha_{22}$. As shown in the proof of the next proposition, this equality restriction can be written as a zero restriction on $q_{2}$. In Combination with the sign normalization restrictions, we obtain global identification.

Proposition 6. The Blanchard-Perotti fiscal policy AB-SVAR is globally identified.
Proof. See Appendix C.
Although the literature has proposed only criteria for local-identification of AB-SVARs, this proposition shows that these restrictions are sufficient for global identification. The comparison of Propositions 5 and 6 highlights that the implications of identifying restrictions with calibrated coefficients for identification depend on whether or not the variances of structural shocks are normalized to unity.

## III. 2 Heteroskedastic SVAR

Bacchiocchi and Fanelli (2015) consider SVARs with a break in the structural error variances and potentially regime-dependent structural coefficients. They consider identifying assumptions that restrict some structural parameters to being invariant across the regimes.

Suppose that the two regimes are characterized by two different reduced-form error covariance matrices $\Sigma_{1}$ and $\Sigma_{2}$, which are related to the regime-dependent structural parameters through

$$
\begin{equation*}
\Sigma_{1}=A_{01}^{-1} A_{01}^{-1 \prime} \quad \text { and } \quad \Sigma_{2}=A_{02}^{-1} A_{02}^{-1 \prime}, \tag{28}
\end{equation*}
$$

where $A_{01}$ and $A_{02}$ are the matrices of regime-specific structural parameters. Let $Q_{1}$ and $Q_{2}$ be the regime specific orthogonal matrices mapping the reduced-form error variances to the structural coefficients,

$$
\begin{equation*}
A_{01}=Q_{1}^{\prime} \Sigma_{1, t r}^{-1} \quad \text { and } \quad A_{02}=Q_{2}^{\prime} \Sigma_{2, t r}^{-1} \tag{29}
\end{equation*}
$$

with $Q_{i}=\left[q_{1(i)}, \ldots, q_{n(i)}\right], \forall i, \in\{1,2\}$. We denote the $j$-th column vector of $\Sigma_{i, t r}^{\prime}$ by $\sigma_{j(i)}$ for $j=1,2$ and $i=1,2$.

For simplicity, consider a bivariate SVAR with two regimes. Impose the following identifying restrictions:

$$
\begin{array}{ll}
\left(A_{01}\right)_{[1,2]}^{-1}=0 & \Longleftrightarrow\left(e_{1}^{\prime} \Sigma_{1, t r}\right) q_{2(1)}=0  \tag{30}\\
\left(A_{01}\right)_{[2,1]}^{-1}=\left(A_{02}\right)_{[2,1]}^{-1} & \Longleftrightarrow\left(e_{2}^{\prime} \Sigma_{1, t r}\right) q_{1(1)}=\left(e_{2}^{\prime} \Sigma_{2, t r}\right) q_{1(2)} .
\end{array}
$$

The first zero restriction combined with the sign normalization pins down the orthogonal matrix $Q_{1}$ in the first regime. The second restriction in Eq. (30) gives rise to the following system of equations:

$$
\left\{\begin{align*}
\sigma_{1(2)}^{\prime} q_{1(2)} & =c  \tag{31}\\
q_{1(2)}^{\prime} q_{1(2)} & =1
\end{align*}\right.
$$

where $c=\sigma_{2(1)}^{\prime} q_{1(1)}$, which is a known constant once the $Q_{1}$ for the first regime is identified. Hence, the problem of identification for structural parameters in the second regime is reduced to the example discussed in Section II.5, in which local identification holds with two distinct solutions.

## III. 3 Restrictions across shocks or across equations

Cross-equation restrictions have been investigated in the classical literature of simultaneous equation systems (Fisher, 1966, and Kelly, 1975). ${ }^{6}$ Analogous restrictions arise in SVARs when crossrestrictions are imposed on impulse responses to different structural shocks, as illustrated by the following example.

Since the seminal work of Sims (1980), many studies of the transmission of monetary policy have adopted SVAR model with a triangular structure where real variables do not immediately respond to monetary policy shocks (see, among many others, Christiano et al., 2005). However, recent theoretical and empirical contributions have provided evidence that monetary policy SVARs are

[^7]not recursive (see Bacchiocchi et al., 2018, and the references therein). Furthermore, Bacchiocchi and Fanelli (2015) provide evidence of Taylor-rule type behaviour by the Federal Reserve during the Great Moderation period, with coefficients in line with the original values proposed by Taylor (1993), i.e. the central bank reacts to unexpected inflation shocks with a coefficient equal to its reaction to excess demand shocks.

This evidence suggests the following simple trivariate model for inflation $\pi_{t}$, output $y_{t}$ and the short term interest rate $i_{t}$

$$
\begin{align*}
a_{11} \pi_{t} & =\varepsilon_{t}^{\pi} \\
a_{21} \pi_{t}+a_{22} y_{t}+a_{23} i_{t} & =\varepsilon_{t}^{y}  \tag{32}\\
a_{31} \pi_{t}+a_{23} y_{t}+a_{33} i_{t} & =\varepsilon_{t}^{m p}
\end{align*}
$$

where $\varepsilon_{t}=\left(\varepsilon_{t}^{\pi}, \varepsilon_{t}^{y}, \varepsilon_{t}^{m p}\right)^{\prime}$ are, respectively the shocks to inflation, demand, and monetary policy. Two restrictions, $a_{12}=a_{13}=0$, have already been imposed. In addition, consider the restriction that the on-impact response of $i_{t}$ to $\varepsilon_{t}^{\pi}$ is equal to that to $\varepsilon_{t}^{y}$. These restrictions can be expressed as

$$
\begin{align*}
a_{12}=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} e_{2}\right)^{\prime} q_{1}=0 \\
a_{13}=0 & \Longleftrightarrow\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime} q_{1}=0  \tag{33}\\
A_{0[3,1]}^{-1}=A_{0[3,2]}^{-1} & \Longleftrightarrow\left(e_{3}^{\prime} \Sigma_{t r}\right) q_{1}=\left(e_{3}^{\prime} \Sigma_{t r}\right) q_{2},
\end{align*}
$$

with $f_{1}=2$ and one restriction connecting the first and second columns of the orthogonal matrix $Q$, which can be viewed as a restriction on $q_{2}$ so $f_{2}=1$. By Proposition 3 the current model is locally identified. However, we can show that global identification is not guaranteed.

Given $\Sigma$, the first two zero restrictions in Eq. (33) and the sign normalization restriction pin down a unique $q_{1}$, so the impulse response for the inflation shock is globally identified. The third restriction in Eq. (33) and orthogonality to $q_{1}$, however, cannot generally pin down a unique $q_{2}$, since they involve a non-homogeneous restriction. Taking $q_{1}$ to be determined in the previous step, we note the similarly to the first two restrictions in Eq. (27) of Section III.1. An analogous argument to the proof of Proposition 5 can be applied to conclude that $q_{2}$ cannot be uniquely determined in a set of reduced-form parameter values that has a positive measure. Thus, $A_{0}$ is locally- but not globally-identified.

## III. 4 Non-recursive SVAR models

RWZ provide an example of a locally- but not globally-identified SVAR where the sufficient condition for local identification of Proposition 3 is not met. This example involves non-recursive causal
ordering restrictions and has practical importance, as we illustrate below.
Cochrane (2006) considers the following New-Keynesian model for inflation $\pi_{t}$, output gap $x_{t}$, and the nominal interest rate $i_{t}$ :

$$
\begin{align*}
\pi_{t} & =\beta E_{t} \pi_{t+1}+\kappa x_{t}+u_{t}^{s} \\
x_{t} & =E_{t} x_{t+1}-\tau\left(i_{t}-E_{t} \pi_{t+1}\right)+u_{t}^{d}  \tag{34}\\
i_{t} & =\phi_{\pi} \phi_{t}+u_{t}^{m p}
\end{align*}
$$

with $u_{t}^{s}$, $t_{t}^{d}$ and $u_{t}^{m p}$ being, respectively, the independent supply, demand, and monetary policy shocks with variances $\sigma_{s}^{2}, \sigma_{d}^{2}$ and $\sigma_{m p}^{2}$. Fukac et al. (2007) show that, once the discount factor $\beta$ has been fixed, this model can be written as an SVAR of the form

$$
A_{0} y_{t}=\varepsilon_{t}
$$

where $y_{t}=\left(\pi_{t}, x_{t}, i_{t}\right)^{\prime}$ is the vector of observable variables, $\varepsilon_{t}=\left(\varepsilon_{t}^{s}, \varepsilon_{t}^{d}, \varepsilon_{t}^{m p}\right)$ collects the unitvariance uncorrelated structural shocks and

$$
A_{0}=\left(\begin{array}{ccc}
a_{11} & a_{12} & 0 \\
0 & a_{22} & a_{23} \\
a_{31} & 0 & a_{33}
\end{array}\right)
$$

Note that there is a well-defined mapping between the parameters in $A_{0}$ and those in the DSGE representation Eq. (34). ${ }^{7}$

This model includes the following restrictions:

$$
\begin{align*}
& a_{13}=0 \quad \Longleftrightarrow \quad\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime} q_{1}=0 \\
& a_{21}=0 \quad \Longleftrightarrow \quad\left(\Sigma_{t r}^{-1} e_{1}\right)^{\prime} q_{2}=0  \tag{35}\\
& a_{32}=0 \quad \Longleftrightarrow \quad\left(\Sigma_{t r}^{-1} e_{2}\right)^{\prime} q_{3}=0,
\end{align*}
$$

with $f_{1}=f_{2}=f_{3}=1$. The sufficient condition for local identification in Proposition 3 is clearly not satisfied, but it can be shown that the rank condition of Proposition 2 is satisfied. Specifically, Rubio-Ramirez et al. (2008) show the existence of two orthogonal matrices $Q_{1}$ and $Q_{2}$ transforming the reduced-form parameters into admissible structural parameters. The empirical application shown in Section VI performs estimation and inference for this model.

[^8]
## III. 5 Proxy-SVAR

A set of identifying restrictions similar to the non-recursive zero restrictions discussed above can appear when the identification strategy exploits proxy variables for the structural shocks.

Consider again a three-variable SVAR. Instead of imposing zero restrictions directly on any element of $A_{0}$, we consider observable variables that proxy some of the underlying structural shocks. The idea of using proxy variables to identify the structural impulse responses has been considered in Stock and Watson (2012) and Mertens and Ravn (2013), amongst others. We restrict our analysis to SVARs and focus on identification of the full system of SVARs rather than subset identification of the impulse responses. ${ }^{8}$ To be specific, consider introducing the external variables $m_{t}=\left(m_{1 t}, m_{2 t}, m_{3 t}\right)^{\prime}$, each of which acts as a proxy for some contemporaneous structural shocks. Following Angelini and Fanelli (2019), Arias et al. (2018b), and Giacomini et al. (2019a), we augment $m_{t}$ into the original SVAR,

$$
\left(\begin{array}{cc}
A_{0} & O  \tag{36}\\
\Gamma_{1} & \Gamma_{2}
\end{array}\right)\binom{y_{t}}{m_{t}}=\binom{\epsilon_{t}}{\nu_{t}}, \quad\left(\epsilon_{t}, \nu_{t}\right)^{\prime} \sim \mathcal{N}\left(0, I_{6 \times 6}\right),
$$

where $O$ is $3 \times 3$ matrix of zeros, $\Gamma_{1}$ and $\Gamma_{2}$ are $3 \times 3$ coefficient matrices in the augmented equations, and the shocks $\nu_{t}$ in the second block component of the augmented system are interpreted as measurement errors in the proxy variables. Inverting eq. (36) leads to

$$
\begin{equation*}
m_{t}=-\Gamma_{2}^{-1} \Gamma_{1} A_{0}^{-1} \epsilon_{t}+\Gamma_{2}^{-1} \nu_{t} \tag{37}
\end{equation*}
$$

In the Proxy-SVAR approach, the identifying restrictions are zero restrictions on the covariance matrix of $m_{t}$ and $\epsilon_{t}$. Consider imposing the following restrictions:

$$
E\left(m_{t} \epsilon_{t}^{\prime}\right)=\left(\begin{array}{ccc}
0 & \rho_{12} & \rho_{13}  \tag{38}\\
\rho_{21} & 0 & \rho_{23} \\
\rho_{31} & \rho_{32} & 0
\end{array}\right)
$$

where $\rho_{i j}$ is the (unconstrained) covariance of $m_{i t}$ and $\epsilon_{j t}$. The zero-covariance restrictions represented in (38) imply that variable $m_{i t}, i=1,2,3$, proxies a combination of the structural shocks excluding $\epsilon_{i t}$. Combining eq. (37) with eq. (38) and substituting $A_{0}^{-1}=\Sigma_{t r} Q, Q=\left[q_{1}, q_{2}, q_{3}\right]$, the

[^9]exogeneity restrictions of eq. (38) can be expressed as
\[

$$
\begin{align*}
& \left(e_{1}^{\prime} \Gamma_{2}^{-1} \Gamma_{1} \Sigma_{t r}\right) q_{1}=0, \\
& \left(e_{2}^{\prime} \Gamma_{2}^{-1} \Gamma_{1} \Sigma_{t r}\right) q_{2}=0,  \tag{39}\\
& \left(e_{3}^{\prime} \Gamma_{2}^{-1} \Gamma_{1} \Sigma_{t r}\right) q_{3}=0 .
\end{align*}
$$
\]

Since $\Gamma_{2}^{-1} \Gamma_{1}$ can be identified by the covariance matrix of the reduced-form VAR errors in the augmented system Eq. (36), the zero restrictions of Eq. (39) have the same form as Eq. (35). Hence, Proxy-SVAR identification under the exogeneity restrictions Eq. (38) delivers local but non-global identification of $A_{0}$ matrix.

## IV Computing identified sets of locally identified SVARs

A common approach to estimating SVAR structural parameters is constrained maximum likelihood (Amisano and Giannini, 1997), with the maximization performed numerically given some initial values. The standard gradient-based algorithm stops once it reaches a local maximum, and does not check for the existence of other observationally equivalent parameter values. Hence, the conventional maximum likelihood procedure applied to an SVAR that is locally but not globally identified will select one of the observationally equivalent structural parameters in a nonsystematic way, limiting the credibility of the resulting estimates and inference.

This section proposes computational methods that produce estimates of all the observationally equivalent $A$ matrices given the identifying restrictions. Our approach is first to obtain $\hat{\phi}=(\hat{B}, \hat{\Sigma})$ an estimate of the reduced-form parameters $\phi$, and then compute the identified set for $A_{0}$ given $\hat{\phi}$, $\mathcal{A}_{0}(\hat{\phi} \mid F, S)$ by solving a system of equations for $Q$ matrix given $\hat{\phi}$. For estimators of $\phi$, we consider (i) the unconstrained reduced-form VAR estimator for $\phi$ denoted by $\hat{\phi}_{u}$ and (ii) the estimator for $\phi$ induced by a constrained maximum likelihood estimate of $A$ under the identifying restrictions (i.e., one of the locally identified structural parameter points maximizing the likelihood), denoted by $\hat{\phi}_{r}$. In the Bayesian inference methods considered in Section V, we view $\hat{\phi}$ as a draw from the posterior of $\phi$.

In what follows, we propose two procedures to compute $\mathcal{A}_{0}(\hat{\phi} \mid F, S)$. The first procedure is general and invokes a non-linear solver. The second procedure is more constructive and involves only elementary calculus, but the allowed type of identifying restrictions is more limited. Both algorithms deal with just identified SVARs, and we presume the rank condition in Proposition 2 or the sufficient condition in Proposition 3 are ensured or have been checked empirically prior to implementation.

## IV. 1 A general computation procedure for locally identified SVARs

Given $\hat{\phi}$, this method computes the orthogonal matrices subject to the identifying restrictions by solving a non-linear system of equations. ${ }^{9}$ If the model is locally identified, then it yields at most $2^{n(n+1) / 2}$ solutions for $Q$. Some of these will be discarded by normalization and sign restrictions. The remained solutions for $Q$ are then used to span the identified set for $A_{0}, \mathcal{A}_{0}(\hat{\phi} \mid F, S)$ and its projection leads to the identified set of an impulse response $I S_{\eta}(\hat{\phi})$. All these steps are stated formally in the next algorithm.

Algorithm 1. Consider a SVAR with equality restrictions Eq. (14) and sign restrictions Eq. (17), and assume $f=n(n-1) / 2$ equality restrictions are imposed. Let $\hat{\phi}$ be a given estimator for $\phi$ such as $\hat{\phi}_{u}$ or $\hat{\phi}_{r}$.

1. Solve the system of equations for $Q$ :

$$
\left\{\begin{align*}
\boldsymbol{F}(\hat{\phi}) \operatorname{vec} Q-\mathbf{c} & =\boldsymbol{o}  \tag{40}\\
Q^{\prime} Q & =I_{n} ;
\end{align*}\right.
$$

2. If the set of real solutions for $Q$ is non-empty (which is guaranteed if $\hat{\phi}=\hat{\phi}_{r}$ ), then retain only those satisfying the normalization and sign restrictions to obtain $\mathcal{Q}_{R}(\hat{\phi}) . \mathcal{A}_{0}(\hat{\phi} \mid F, S)$ is constructed accordingly by $\left\{A_{0}=Q^{\prime} \hat{\Sigma}_{t r}^{-1}: Q \in \mathcal{Q}_{R}(\hat{\phi})\right\}$.
3. When $\hat{\phi}=\hat{\phi}_{u}$, it is possible that no real solution for $Q$ exists in Step 1. If so, we return $\mathcal{Q}(\hat{\phi} \mid F, S)=\emptyset$, i.e., $\hat{\phi}$ is not compatible with the imposed identifying restrictions.

The crucial step in this algorithm is obtaining all the solutions to the equation system (40). This is a system of polynomial equations consisting of linear and quadratic equations. ${ }^{10}$ Closedform solutions do not seem available, but numerical algorithms to compute all the roots of the polynomial equations are. Matlab, for example, has the function vpasolve, an algorithm to find all the solutions of a system of non-linear equations. According to the Matlab documentation, ${ }^{11}$ vpasolve returns the complete set of solutions in the case of polynomial equations. The strength of this algorithm is its generality, but it is a black-box function. ${ }^{12}$

[^10]When non-homogeneous restrictions or cross-shock restrictions are imposed, the model becomes observationally restrictive, as seen in Sections III.1-III.3. Hence, when $\hat{\phi}$ is obtained from the unconstrained reduced-form VAR estimator $\hat{\phi}_{r}$, if $\hat{\phi}_{r}$ happens to be outside of $\Phi_{R}$, then Step 3 of Algorithm 1 becomes relevant. When the algorithm returns $\mathcal{Q}_{R}(\hat{\phi})=\emptyset$, the maximum likelihood reduced-form model suggests that some of the imposed identifying restrictions are misspecfied. One can hence consider relaxing some of the imposed sign restrictions, or modify the value of $\mathbf{c} \neq 0$ if non-homogeneous restrictions are present. Alternatively, if we want to maintain the imposed restrictions, we can employ the constrained reduced-form estimate $\hat{\phi}=\hat{\phi}_{r}$ instead, so that $\mathcal{Q}_{R}(\hat{\phi})$ is guaranteed to be nonempty.

## IV. 2 Computational procedure for locally identified SVARs with recursive nonhomogeneous restrictions

If the identifying restrictions imposed allow the sequential determination of the column vectors of $Q$ as exploited in the identification arguments in the previous sections, we can modify Algorithm 1. In this section, we consider recursive SVARs with non-homogeneous and cross-shock restrictions, as covered in Proposition 3.

Let $Q_{1: i}, 1 \leq i \leq n$, be a $n \times i$ matrix whose column vectors are orthonormal (i.e. it consists of the first $i$ column vectors of $Q$ ). Given $\phi$, define $\tilde{F}_{11}(\phi)=F_{11}(\phi)$ and the following matrices sequentially for $i=2, \ldots, n$,

$$
\begin{equation*}
\tilde{F}_{i i}(\phi)=\binom{F_{i i}(\phi)}{Q_{1:(i-1)}(\phi)^{\prime}}, \text { where }\left(F_{j 1}(\phi), \ldots F_{j j}(\phi)\right) \operatorname{vec} Q_{1:(i-1)}(\phi)=c_{j} \tag{41}
\end{equation*}
$$

where $Q_{1:(i-1)}(\phi)$ satisfies the identifying restrictions for the first $(i-1)$ orthogonal vectors, i.e., $\left(F_{j 1}(\phi), \ldots F_{j j}(\phi)\right) \operatorname{vec} Q_{1:(i-1)}(\phi)=c_{j}$ holds for $j=1, \ldots,(i-1)$. For $i=1, \ldots, n$, we define a $(n-1) \times 1$ vector,

$$
\tilde{c}_{i}(\phi)=\left(\begin{array}{c}
c_{i}-\left(F_{i 1}(\phi), \ldots, F_{i(i-1)}(\phi)\right) \operatorname{vec} Q_{1:(i-1)}(\phi)  \tag{42}\\
0 \\
\vdots \\
0
\end{array}\right)
$$

Then, for $i=1, \ldots, n$, define

$$
\begin{align*}
d_{i}(\phi) & =\tilde{F}_{i i}(\phi)^{\prime}\left(\tilde{F}_{i i}(\phi) \tilde{F}_{i i}(\phi)^{\prime}\right)^{-1} \tilde{c}_{i}(\phi)  \tag{43}\\
B_{i}(\phi) & =\left(I_{n}-\tilde{F}_{i i}(\phi)^{\prime}\left(\tilde{F}_{i i}(\phi) \tilde{F}_{i i}(\phi)^{\prime}\right)^{-1} \tilde{F}_{i i}(\phi)\right), \tag{44}
\end{align*}
$$

and let $\alpha_{i}(\phi)$ be a $n \times 1$ basis vector of the linear space spanned by the vectors in $B_{i}(\phi)$. Note that $B_{i}(\phi)$ is the $n \times n$ matrix projecting onto the linear space orthogonal to the row vectors of $\tilde{F}_{i i}(\phi)$. Hence, given the rank of $\tilde{F}_{i i}(\phi)$ is $n-1, B_{i}(\phi)$ has a rank of 1 , so $\alpha_{i}(\phi)$ is unique up to sign, and $\tilde{F}_{i i}(\phi) \alpha_{i}(\phi)=0$ holds.

Consider the $n \times 1$ vector, $x=d_{i}(\phi)+z \alpha_{i}(\phi), z \in \mathbb{R}$. Due to the way $d_{i}(\phi)$ and $\alpha_{i}(\phi)$ are constructed, $\tilde{F}_{i i}(\phi) x=\tilde{c}_{i}$ holds. That is, by choosing $z$ so that $x$ is a unit-length vector, we can obtain $q_{i}$ vectors satisfying $\tilde{F}_{i i}(\phi) q_{i}=\tilde{c}_{i}$. Solving for $x$ is simple as it requires only finding the roots of a quadratic equation (see Eq. (45) and Eq. (46) in Algorithm 2 below). Given $\phi$, we repeat this process for every $i=1, \ldots, n$ to determine the $q_{i}$ vectors sequentially, and compute all the $Q$ matrices satisfying the equality restrictions $\mathbf{F}(\phi, Q)=\mathbf{0} . \mathcal{A}_{0}(\phi \mid F, S)$ and $I S_{\eta}(\phi)$ can then be obtained by retaining the $Q$ that satisfy the normalization and sign restrictions. We summarize this computational procedure in the next algorithm.

Algorithm 2. Consider a SVAR satisfying the normalization restrictions Eq. (7), the equality restrictions Eq. (14), and the sign restrictions Eq. (17), where the imposed equality restrictions satisfy the sufficient condition for local identification given in Proposition 3. Let $\hat{\phi}$ be a given estimator for $\phi$ such as $\hat{\phi}_{u}$ or $\hat{\phi}_{r}$. In the description of the algorithm below, we omit the argument $\hat{\phi}$ as far as it does not give rise confusion.

Let $\boldsymbol{b}=\left(b_{1}, \ldots, b_{n}\right) \in\{0,1\}^{n}$ be a bit vector which will be used to index each of the at most $2^{n}$ possible solutions for the $Q$ matrices. Beginning with $\boldsymbol{B}=\{0,1\}^{n}$, we will map each $\boldsymbol{b} \in \boldsymbol{B}$ to a possible solution of $Q$, check if it is feasible or not, and refine $\boldsymbol{B}$ accordingly. The following algorithm describes this process in detail:

1. Solve for $z \in \mathbb{R}$ in

$$
\begin{equation*}
d_{1}^{\prime} d_{1}+2 d_{1}^{\prime} \alpha_{1} z+\alpha_{1}^{\prime} \alpha_{1} z^{2}=1, \tag{45}
\end{equation*}
$$

and denote the two solutions by $z_{1}^{b_{1}}, b_{1} \in\{0,1\}$.
(a) If they are real, then define $q_{1}^{b_{1}}=d_{1}+\alpha_{1} z_{1}^{b_{1}}, b_{1} \in\{0,1\}$. Let $\boldsymbol{B}_{1} \subset\{0,1\}$ be the set of $b_{1}$ such that $q_{1}^{b_{1}}$ satisfies the sign normalization and sign restrictions for $q_{1}$. If $\boldsymbol{B}_{1}$ is empty (i.e., no $q_{1}^{b_{1}}$ satisfies the sign normalization and sign restrictions for $q_{1}$ ), then stop and conclude $\mathcal{Q}_{R}(\hat{\phi})=\emptyset$
(b) If the roots of Eq. (45) are not real, then stop and return $\mathcal{Q}(\hat{\phi} \mid F, S)=\emptyset$.
2. This step iterates sequentially for $i=2, \ldots, n$, given $\boldsymbol{B}_{i-1} \subset\{0,1\}^{i-1}$.
(a) For each $\left(b_{1}, \ldots, b_{i-1}\right) \in \boldsymbol{B}_{i-1}$, construct $\boldsymbol{B}_{i}\left(b_{1} b_{2} \cdots b_{i-1}\right) \subset\{0,1\}^{i}$ by performing the following subroutines:
i. Construct $\tilde{F}_{i i}$ from Eq. (41) by setting $Q_{1: i-1}=\left[q_{1}^{b_{1}}, q_{2}^{b_{1} b_{2}}, \ldots, q_{i-1}^{b_{1} \cdots b_{i-1}}\right]$, and obtain $d_{i}$ and $\alpha_{i}$ accordingly. Then, solve for $z \in \mathbb{R}$ in

$$
\begin{equation*}
d_{i}^{\prime} d_{i}+2 d_{i}^{\prime} \alpha_{i} z+\alpha_{i}^{\prime} \alpha_{i} z^{2}=1, \tag{46}
\end{equation*}
$$

and denote the two solutions by $z_{i}^{b_{1} b_{2} \cdots b_{i}}, b_{i} \in\{0,1\}$.
ii. If they are real, define $q_{1}^{b_{1} b_{2} \cdots b_{i}}=d_{i}+\alpha_{i} z_{1}^{b_{1} b_{2} \cdots b_{i}}, b_{i} \in\{0,1\}$. Let $\boldsymbol{B}_{i}\left(b_{1} b_{2} \cdots b_{i}\right)$ be the set of $\left(b_{1}, b_{2}, \ldots, b_{i}\right) \in\{0,1\}^{i}$ such that $q_{i}^{b_{1} b_{2}, \cdots b_{i}}$ satisfies the sign normalization and sign restrictions for the $i$-th column vector of $Q$. This can be empty if no $q_{i}^{b_{1} b_{2}, \cdots b_{i}}$ satisfies them.
iii. If the roots of Eq. (46) are not real, return $\boldsymbol{B}_{i}\left(b_{1} b_{2} \cdots b_{i-1}\right)=\emptyset$.
(b) Construct $\boldsymbol{B}_{i}=\bigcup_{\left(b_{1}, \ldots, b_{i-1}\right) \in \boldsymbol{B}_{i-1}} \boldsymbol{B}_{i}\left(b_{1} \cdots b_{i-1}\right)$. If $\boldsymbol{B}_{i} \neq \emptyset$, go back to the beginning of Step 2.
(c) If $\boldsymbol{B}_{i}=\emptyset$, then stop and return $\mathcal{Q}_{R}(\hat{\phi})=\emptyset$.
3. We obtain

$$
\mathcal{Q}_{R}(\hat{\phi})=\left\{\left(q_{1}^{b_{1}}, q_{2}^{b_{1} b_{2}}, \ldots, q_{n}^{b_{1} b_{2} \cdots b_{n}}\right): \boldsymbol{b} \in \boldsymbol{B}_{n}\right\} .
$$

Algorithm 2 computes the set of all admissible $Q \in \mathcal{Q}_{R}(\hat{\phi})$. In the description of the algorithm, they are indexed by the bit vectors $\mathbf{b} \in \mathbf{B}_{n}$. The algorithm is constructive and guaranteed to compute all the admissible $Q$ matrices. Projecting this set of admissible matrices onto the impulse response of interest, we obtain a plug-in estimate of the identified set $I S_{\eta}(\hat{\phi})$.

Algorithm 2 is more constructive than Algorithm 1, but it restricts the set of equality restrictions to be recursive. Algorithm 2 can be extended to a class of models involving non-recursive identifying restrictions (e.g., examples in Sections III. 4 and III.5) by incorporating steps that solve a certain system of quadratic equations. Such an algorithm is rather involved to present, so we do not include it in this paper. Algortihm 1 can be certainly applied to a general class of models with nonrecursive identifying restrictions.

## V Inference for locally identified SVARs

## V. 1 Bayesian inference

Standard Bayesian inference specifies a prior distribution for either the structural parameters $A$ (e.g., Baumeister and Hamilton, 2015), or the reduced-form parameters and rotation matrix ( $\phi, Q$ ) as a reparametrization of $A$ (e.g., Uhlig, 2005). When identification is local, the likelihood for the joint parameter vector $A$ can have multiple modes, which means that the posterior for the structural parameters and impulse responses may also have mutliple modes. This leads to computational challenges as commonly used Markov Chain Monte Carlo (MCMC) methods can fail to adequately explore the posterior when it is multi-modal. For instance, in the standard Metropolis-Hastings algorithm, the presence of multiple modes complicates the choice of proposal distribution. If the proposal distribution in the Metropolis-Hastings algorithm does not support some some modes well, a lack of irreducibility of the Markov chain can lead it to fail to converge to the posterior. Similarly, in the standard Gibbs sampler, the presence of multiple modes in the posterior for $A$ leads its support to be almost disconnected, which can then lead a break down of irreducibility and the Gibbs sampler to fail to converge (see Example 10.7 in Robert and Casella (2004)). By combining our constructive algorithm for computing $I S_{\eta}(\phi)$ with the posterior sampling algorithm for $\phi$, we can overcome such computational challenges.

We consider approximating the posterior for a scalar impulse response $\eta(\phi)$. Assume that the reduced-form parameters yield nonempty $\mathcal{Q}_{R}(\phi)$. Let $I S_{\eta}(\phi)$ consist of $M(\phi) \geq 1$ distinct points,

$$
\begin{equation*}
I S_{\eta}(\phi)=\left\{\eta_{1}(\phi), \eta_{2}(\phi), \ldots, \eta_{M(\phi)}(\phi)\right\}, \tag{47}
\end{equation*}
$$

where we index the observationally equivalent impulse responses to satisfy $\eta_{1}(\phi)<\eta_{2}(\phi)<\cdots<$ $\eta_{M(\phi)}(\phi)$.

We follow the "agnostic" Bayesian approach of Uhlig (2005). The posterior for $\eta$ is induced by the posterior for $\phi, \pi_{\phi \mid Y}$, which is supported on $\Phi_{R} \equiv\left\{\phi: \mathcal{Q}_{R}(\phi) \neq \emptyset\right\}$, and $Q$ has a uniform prior supported only on the admissible set of rotation matrices $\mathcal{Q}_{R}(\phi)$ given $\phi \in \Phi_{R}$. Local identification with the $M(\phi)$-point identified set as in (47) can be obtained by projecting the $M(\phi)$ admissible rotation matrices if each of them leads to distinct values of impulse response. Hence, the uniform weights assigned over these rotation matrices imply that equal weights are assigned to the points in $I S_{\eta}(\phi)$. As a result, for $G \subset \mathbb{R}$, the posterior for $\eta$ can be expressed as, :

$$
\begin{equation*}
\pi_{\eta \mid Y}(\eta \in G) \propto E_{\phi \mid Y}\left[\sum_{m=1}^{M(\phi)} 1\left\{\eta_{m}(\phi) \in G\right\}\right] . \tag{48}
\end{equation*}
$$

Since the reduced-form VAR likelihood is unimodal and concentrated around the maximum likelihood estimate, MCMC algorithms will perform well when sampling from $\pi_{\phi \mid Y}$. Hence, the posterior (48) can be approximated by combining a posterior sampler for $\phi$ with the algorithm for computing $\left\{\eta_{m}(\phi): m=1, \ldots, M(\phi)\right\}$.

## V. 2 Frequentist-valid inference

Bayesian inference as considered above can be sensitive to the choice of prior even in large samples due to the lack of global identification. The standard Bayesian procedure (assuming a unique prior for the structural parameters) specifies an allocation of the prior belief over observationally equivalent impulse responses, $I S_{\eta}(\phi)$, conditional on $\phi$. This conditional belief given $\phi$ is not updated by the data and, as a result, the shape and heights of the posterior around the modes remain sensitive to its specification. In this section, we propose an asymptotically valid frequentist inference procedure for the impulse response identified set that can draw inferential statements which are robust to the choice of prior weights over the set of locally identified parameter values.

Our approach is to project asymptotically valid frequentist confidence sets for the reduced-form parameters $\phi$ through the identified set mapping $I_{\eta}(\phi)$. In standard set-identified models where the identified set is a connected interval with positive width, the projection approach to constructing the confidence set has appeared in the 2011 working paper version of Moon and Schorfheide (2012), Norets and Tang (2014), Kline and Tamer (2016), among others. This approach generally yields asymptotically valid (but conservative) confidence sets even when the identified set consists of discrete points. However, a challenge unique to the discrete identified set case is the computation of projection confidence sets for the impulse responses based on a finite number of grid points or draws of $\phi$ from their confidence set. In what follows, we propose methods to tackle this computational challenge.

Let $C S_{\phi, \alpha}$ be an asymptotically valid confidence set for $\phi$ with coverage probability $\alpha \in(0,1)$. If the maximum likelihood estimator $\hat{\phi}$ is $\sqrt{T}$-asymptotically normal, the likelihood contour set $C S_{\phi, \alpha}$ is determined by the $\alpha$-th quantiles of the $\chi^{2}$ distribution with the degree of freedom $\operatorname{dim}(B)+$ $n(n-1) / 2$. If the posterior for $\phi$ satisfies the Bernstein-von Mises property, that is the posterior for $\sqrt{T}(\phi-\hat{\phi})$ asymptotically coincides with the sampling distribution of the maximum likelihood estimator, the Bayesian highest density posterior region with credibility $\alpha$ can be used for $C S_{\phi, \alpha}$. The MCMC confidence set procedure developed by Chen et al. (2018) can then be used to obtain draws of $\phi$ from the highest density posterior region with credibility $\alpha$. We follow this procedure in our empirical application below. The inference procedures below allows for any $C S_{\phi, \alpha}$ with asymptotically valid coverage, and takes draws or grids of $\phi$ from $C S_{\phi, \alpha}$ as given.

The projection confidence set is defined as

$$
\begin{equation*}
C S_{\eta, \alpha}^{p}=\bigcup_{\phi \in C S_{\phi, \alpha}} I S_{\eta}(\phi) \tag{49}
\end{equation*}
$$

We assume that $C S_{\phi, \alpha}$ is an asymptotically valid confidence set for $\phi$ in the sense that

$$
\lim _{T \rightarrow \infty} p_{Y^{T} \mid \phi_{0}}\left(\phi_{0} \in C S_{\phi, \alpha}\right)=\alpha
$$

where $p_{Y^{T} \mid \phi_{0}}$ is the sampling distribution of the data with sample size $T$ and $\phi_{0}$ is the true value of $\phi$. Since $\left\{\phi_{0} \in C S_{\phi, \alpha}\right\}$ implies $\left\{I S_{\eta}\left(\phi_{0}\right) \subset C S_{\eta, \alpha}^{p}\right\}, C S_{\eta, \alpha}^{p}$ (and any set including $C S_{\eta, \alpha}^{p}$ ) is an asymptotically-valid but potentially conservative confidence set for $I S_{\eta}\left(\phi_{0}\right)$,

$$
\lim _{T \rightarrow \infty} p_{Y^{T} \mid \phi_{0}}\left(I S_{\eta}\left(\phi_{0}\right) \subset C S_{\eta, \alpha}^{p}\right) \geq \alpha
$$

Let $\left\{\phi_{k}: k=1, \ldots, K\right\}$ be a finite number of Monte Carlo draws or grid points from $C S_{\phi, \alpha}$. A sample analogue of the projection confidence set, $\bigcup_{k=1, \ldots, K} I S_{\eta}\left(\phi_{k}\right)$, is less useful in approximating $C S_{\eta, \alpha}^{p}$, because each $I S_{\eta}\left(\phi_{k}\right)$ is a discrete set, whereas the underlying $C S_{\eta, \alpha}^{p}$ we want to approximate can be a union of disconnected intervals with positive widths. In addition, it is difficult to judge how many disconnected intervals $I S_{\eta, \alpha}^{p}$ has and where the possible gaps lie within $C S_{\eta, \alpha}^{p}$ from a finite number of draws of $I S_{\eta}\left(\phi_{k}\right), k=1, \ldots, K$. Reporting the convex hull of $\bigcup_{k=1, \ldots, K} I S_{\eta}\left(\phi_{k}\right)$ is simple, but it can lead to a connected confidence set that obscures the discrete feature of the identified set.

In what follows, we propose two different approaches for computing the projection confidence set for an impulse response given a set of Monte Carlo draws for $\phi$. We refer to the first as switching-label projection confidence sets. It allows the labels indexing observationally equivalent impulse responses to vary across the horizons, and produces confidence sets that can capture multimodality of the posterior distribution or the integrated likelihood for each impulse response at each horizon. We refer to the second approach as constant label projection confidence set. It maintains unique labels for observationally equivalent structural parameters across the impulse responses and over horizons, i.e., the labels for observationally equivalent structural parameters are defined in terms of the modes of the posterior for $Q$. This approach may produce confidence sets that are wider than the switching label projection confidence set, but it can better capture and visualize dependence of the impulse responses over the horizons.

## V.2.1 Switching-label projection confidence sets

The switching-label approach draws inference for each impulse response at each horizon one-by-one. We hence set $\eta(\phi)$ to a particular scalar impulse response.

Maintaining the notation of the previous subsection, let $I S_{\eta}\left(\phi_{k}\right)=\left\{\eta_{1}\left(\phi_{k}\right), \ldots, \eta_{M\left(\phi_{k}\right)}\left(\phi_{k}\right)\right\}$, where $M\left(\phi_{k}\right)$ is the number of distinct points in the identified set at $\phi=\phi_{k}$. We label these points in increasing order, $\eta_{1}\left(\phi_{k}\right)<\cdots<\eta_{M\left(\phi_{k}\right)}\left(\phi_{k}\right)$. Let $\bar{M}=\max _{k} M\left(\phi_{k}\right)$ be the largest cardinality of $I S_{\eta}\left(\phi_{k}\right)$ among the draws of $\phi_{k}, k=1, \ldots, K . \bar{M}$ indicates the largest possible number of disconnected intervals of $C S_{\eta, \alpha}^{p}$. We view them as clusters, each of which is indexed by $\tilde{m} \in\{1, \ldots, \bar{M}\}$. Let $\tilde{K}=\left|\left\{\phi_{k}: M\left(\phi_{k}\right)=\bar{M}\right\}\right|$ be the number of $\phi$ draws that has the maximal number of observationally equivalent impulse responses and define estimates of the cluster-specific mean and variance by

$$
\begin{align*}
\mu_{\tilde{m}} & =\frac{1}{\tilde{K}} \sum_{\phi_{k}: M\left(\phi_{k}\right)=\bar{M}} \eta_{\tilde{m}}\left(\phi_{k}\right), \\
\sigma_{\tilde{m}}^{2} & =\frac{1}{\tilde{K}-1} \sum_{\phi_{k}: M\left(\phi_{k}\right)=\bar{M}}\left(\eta_{\tilde{m}}\left(\phi_{k}\right)-\mu_{\tilde{m}}\right)^{2}, \tag{50}
\end{align*}
$$

for each $\tilde{m}=1, \ldots, \bar{M}$.
For each $\phi_{k}, k=1, \ldots, K$, we augment a binary vector of length $\bar{M}, D\left(\phi_{k}\right)=\left(D_{\tilde{m}}\left(\phi_{k}\right) \in\{0,1\}:\right.$ $\tilde{m}=1, \ldots, \bar{M})$, which indicates whether or not any one point of $I S_{\eta}\left(\phi_{k}\right)$ can be associated with $\tilde{m}$-th cluster. The true $D\left(\phi_{k}\right)$ is not observed, so must be imputed by, for instance, maximizing the Gaussian $\log$-likelihood criterion in the following manner. Let $\rho_{\phi_{k}}$ be an increasing injective map from $\left\{1, \ldots, M\left(\phi_{k}\right)\right\}$ to $\{1, \ldots, \bar{M}\}$, characterizing which cluster each $\eta_{m}\left(\phi_{k}\right), m=1, \ldots, M\left(\phi_{k}\right)$, belongs to. Define

$$
\begin{equation*}
\hat{\rho}_{\phi_{k}} \in \arg \min _{\rho_{\phi_{k}}} \sum_{m=1}^{M\left(\phi_{k}\right)} \frac{\left(\eta_{m}\left(\phi_{k}\right)-\mu_{\rho_{\phi_{k}}(m)}\right)^{2}}{\sigma_{\rho_{\phi_{k}}(m)}^{2}} \tag{51}
\end{equation*}
$$

which minimizes the sum of variance-weighted squared distances to the cluster-specific means. We then construct $D\left(\phi_{k}\right)=\left(D_{\tilde{m}}\left(\phi_{k}\right): \tilde{m}=1, \ldots, \bar{M}\right) \in\{0,1\}^{\bar{M}}$ from the indicators for whether $\hat{\rho}_{\phi_{k}}$ maps any $m \in\left\{1, \ldots, M\left(\phi_{k}\right)\right\}$ to $\tilde{m}$, i.e., $D_{\tilde{m}}\left(\phi_{k}\right)=1\left\{\exists m\right.$ s.t. $\left.\rho_{\phi_{k}}(m)=\tilde{m}\right\}$. We then construct an interval for each cluster $\tilde{m} \in\{1, \ldots, \bar{M}\}$ by

$$
\begin{equation*}
C_{\tilde{m}}=\left[\min _{\left.{\phi_{k}: D_{\tilde{m}}\left(\phi_{k}\right)=1} \eta_{\hat{\rho}_{\phi_{k}}^{-1}(\tilde{m})}\left(\phi_{k}\right), \max _{\phi_{k}: D_{\tilde{m}}\left(\phi_{k}\right)=1} \eta_{\hat{\rho}_{\phi_{k}}^{-1}(\tilde{m})}\left(\phi_{k}\right)\right] . . ~ . ~ . ~}^{\text {. }}\right. \tag{52}
\end{equation*}
$$

An approximation of the projection confidence set is then formed by taking the union of $C_{\tilde{m}}$ :

$$
\begin{equation*}
\widehat{C S}_{\eta, \alpha}^{p} \equiv \bigcup_{\tilde{m}=1}^{\bar{M}} C_{\tilde{m}} \tag{53}
\end{equation*}
$$

Note $\widehat{C S}_{\eta, \alpha}^{p}$ obtained in this way includes all the $I S_{\eta}\left(\phi_{k}\right), k=1, \ldots, K$, and at the same time, can yield a collection of disconnected intervals. Moreover, if the maximum likelihood estimator for $\phi$ is consistent for $\phi_{0}, I S_{\eta}(\phi)$ is a continuous correspondence at $\phi_{0}$ and $M(\phi)$ is constant in an open neighborhood of $\phi_{0}$, it can be shown that $\widehat{C S}_{\eta, \alpha}^{p}$ converges to $I S_{\eta}\left(\phi_{0}\right)$ in the Hausdorff metric. Hence, $\widehat{C S}_{\eta, \alpha}^{p}$ can consistently uncover the true identified set consisting of potentially multiple points.

We construct $\widehat{C S}_{\eta, \alpha}^{p}$ separately for each impulse response at each horizon. Hence, the labeling of the clusters $\tilde{m}=1, \ldots, \bar{M}$ defined for one impulse response does not correspond to the labeling of the clusters defined for other impulse responses or horizons. For example, a particular impulse response function labeled as $\tilde{m}=1$ in one horizon can be labeled as $\tilde{m}=2$ in another horizon. We expect that switching-label projection confidence sets can visualize well the multi-modality of the marginal posterior for each impulse response.

## V.2.2 Constant label projection confidence sets

In contrast to label-switching projection confidence sets, constant label projection confidence sets maintain fixed labeling across impulse responses and over time horizons. For example, an impulse response function labeled as $\tilde{m}=1$ at one horizon is labeled as $\tilde{m}=1$ at other horizons.

To implement this procedure, we need to anchor the labels to a particular impulse response, say, the impulse response of $i^{*}$-th variable to $j^{*}$-th structural shock at a particular horizon $h=h^{*}$, denoted hereafter by $\eta^{*}\left(\phi, q_{j^{*}}\right) \equiv e_{i^{*}}^{\prime} C_{h^{*}}(\phi) q_{j^{*}}$. Given a Monte Carlo draw of the reduced-form parameters, $\phi_{k}, k=1, \ldots, K$, from $C S_{\phi, \alpha}$, let $q_{j^{*}, m}\left(\phi_{k}\right), m=1, \ldots, M\left(\phi_{k}\right)$ be observationally equivalent $q_{j^{*}}$ vectors labeled according to the ordering of $\eta^{*}\left(\phi, q_{j^{*}}\right)$, i.e., $\eta^{*}\left(\phi_{k}, q_{j^{*}, 1}\left(\phi_{k}\right)\right)<\eta^{*}\left(\phi_{k}, q_{j^{*}, 2}\left(\phi_{k}\right)\right)<$ $\cdots<\eta^{*}\left(\phi_{k}, q_{j^{*}, M\left(\phi_{k}\right)}\left(\phi_{k}\right)\right)$. Similarly to the labeling procedure shown in Eq. (51), we assign cluster identifier $\tilde{m}=1, \ldots, \bar{M}$ to $q_{j^{*}, m}\left(\phi_{k}\right)$ by constructing $\hat{\rho}_{\phi_{k}}$ an increasing injective map from $\left\{1, \ldots, M\left(\phi_{k}\right)\right\}$ to $\{1, \ldots, \bar{M}\}$,

$$
\hat{\rho}_{\phi_{k}} \in \arg \min _{\rho_{\phi_{k}}} \sum_{m=1}^{M\left(\phi_{k}\right)} \frac{\left(\eta^{*}\left(\phi_{k}, q_{j^{*}, m}\right)-\mu_{\rho_{\phi_{k}}(m)}\right)^{2}}{\sigma_{\rho_{\phi_{k}}(m)}^{2}}
$$

where $\mu_{\tilde{m}}=\frac{1}{\tilde{K}} \sum_{\phi_{k}: M\left(\phi_{k}\right)=\bar{M}} \eta^{*}\left(\phi_{k}, q_{j^{*}, \tilde{m}}\right)$ and $\sigma_{\tilde{m}}^{2}=\frac{1}{\tilde{K}-1} \sum_{\phi_{k}: M\left(\phi_{k}\right)=\bar{M}}\left(\eta^{*}\left(\phi_{k}, q_{j^{*}, \tilde{m}}\right)-\mu_{\tilde{m}}\right)^{2}$. We then construct $D\left(\phi_{k}\right)$ in the same way as the label switching projection confidence set.

For each impulse response $\eta\left(\phi, q_{j^{*}}\right)=e_{i}^{\prime} C_{h}(\phi) q_{j^{*}}, i=1, \ldots, n$, and $h=0,1, \ldots$, we construct

$$
C_{\tilde{m}}=\left[\min _{\phi_{k}: D_{\tilde{m}}\left(\phi_{k}\right)=1} \eta\left(\phi_{k}, q_{j^{*}, \hat{\rho}_{\phi_{k}}^{-1}(\tilde{m})}\right), \max _{\phi_{k}: D_{\tilde{m}}\left(\phi_{k}\right)=1} \eta\left(\phi_{k}, q_{j^{*}, \hat{\rho}_{\phi_{k}}^{-1}(\tilde{m})}\right)\right]
$$

and form confidence sets by taking the union over $\tilde{m}$ as in Eq. (53).
In contrast to the switching label procedure, the constant label projection confidence sets keep the labeling of the observationally equivalent impulse responses $\hat{\rho}_{\phi_{k}}(m)$ fixed over variables $i=1, \ldots, n$ and different horizons $h=0,1, \ldots$. If the impulse response $\eta^{*}\left(\phi, q_{j^{*}}\right)$ chosen to anchor the labels can tie the observationally equivalent impulse responses to different economic models or hypotheses, the labels can be interpreted as indexing the underlying economic model or hypothesis and kept invariant throughout the impulse response analysis. The constant label projection confidence set approach is suitable in such a case, and allows us to track and compare the observationally equivalent impulse response functions across different models.

## V.2.3 Robust Bayesian interpretation

If we obtain $\left\{\phi_{k}: k=1, \ldots, K\right\}$ as draws from the credible region of the posterior distribution for $\phi, \widehat{C S}_{\eta, \alpha}^{p}$ can be seen as an approximation of the set $C_{\eta, \alpha}$ satisfying

$$
\pi_{\phi \mid Y}\left(I S_{\eta}(\phi) \subset C_{\eta, \alpha}\right) \geq \alpha .
$$

In terms of the robust Bayesian procedure proposed in Giacomini and Kitagawa (2020), $C_{\eta, \alpha}$ can be interpreted as a robust credible region with credibility $\alpha$; a set of $\eta$ on which a posterior distribution for $\eta$ assigns probability at least $\alpha$ irrespective of the choice of the unrevisable part of the prior $\pi_{Q \mid \phi}$. Our construction of the robust credible region can be conservative and is not guaranteed to provide the shortest one. We leave the construction of the shortest robust credible region for future research.

This link to robust Bayes inference also suggests that the range of posterior probabilities (lower and upper probabilities) spanned by arbitrary conditional priors for $Q$ given $\phi$ can be computed straightforwardly based on the draws $\left\{\phi_{\ell}: \ell=1, \ldots, L\right\}$ from the posterior $\pi_{\phi \mid Y}$. Let $H_{0} \subset \mathbb{R}$ and consider the range of posterior probabilities for a hypothesis of interest $\left\{\eta(\phi, Q) \in H_{0}\right\}$. By applying Theorem 1 of Giacomini and Kitagawa (2020), the range of posterior probabilities for $\left\{\eta \in H_{0}\right\}$ is given by the convex interval:

$$
\begin{equation*}
\pi_{\eta \mid Y}\left(H_{0}\right) \in\left[\pi_{\eta \mid Y *}\left(H_{0}\right), \pi_{\eta \mid Y}^{*}\left(H_{0}\right)\right] \equiv\left[\pi_{\phi \mid Y}\left(I S_{\eta}(\phi) \subset H_{0}\right), \pi_{\phi \mid Y}\left(I S_{\eta}(\phi) \cap H_{0} \neq \emptyset\right)\right] . \tag{54}
\end{equation*}
$$

Since the algorithms given in Section IV exhaust all the locally identified parameter values in
$I S_{\eta}(\phi)$, we can approximate the lower and upper bounds of the posterior probabilities in (54) for each hypothesis of interest by the Monte Carlo frequencies for $\left\{I S_{\eta}(\phi) \subset A\right\}$ and $\left\{I S_{\eta}(\phi) \cap A \neq \emptyset\right\}$, respectively,

$$
\left[\hat{\pi}_{\eta \mid Y *}\left(H_{0}\right), \hat{\pi}_{\eta \mid Y}^{*}\left(H_{0}\right)\right] \equiv\left[\frac{1}{L} \sum_{\ell=1}^{L} 1\left\{I S_{\eta}\left(\phi_{\ell}\right) \subset H_{0}\right\}, \frac{1}{L} \sum_{\ell=1}^{L} 1\left\{I S_{\eta}\left(\phi_{\ell}\right) \cap H_{0} \neq \emptyset\right\}\right] .
$$

For a scalar impulse response, it is also straightforward to compute the range of posterior means. Let $\underline{\eta}(\phi)=\min \left\{\eta \in I S_{\eta}(\phi)\right\}$ and $\bar{\eta}(\phi)=\max \left\{\eta \in I S_{\eta}(\phi)\right\}$. Theorem 2 in Giacomini and Kitagawa (2020) shows that the range of posterior means is given by the connected interval $\left[E_{\phi \mid Y}(\underline{\eta}(\phi)), E(\bar{\eta}(\phi))\right]$, which can be approximated by Monte Carlo analogues based on draws $\left\{\phi_{\ell}\right.$ : $\ell=1, \ldots, L\}$ from $\pi_{\phi \mid Y}$.

## VI Empirical application

We illustrate how our approach works with an empirical application to the non-recursive NewKeynesian SVAR shown in Section III.4. We consider the small scale DSGE model presented in Eq. (34), which has the SVAR representation with sign normalizations and the zero restrictions Eq. (35). The vector of observables is inflation as measured by the GDP deflator $\left(\pi_{t}\right)$, real GDP per capita as a deviation from a linear trend $\left(x_{t}\right)$ and the federal funds rate $\left(i_{t}\right) .{ }^{13}$ The data are quarterly from 1965:1 to 2006:1.

As discussed in Section III.4, the imposed restrictions deliver local identification and, given the reduced form parameters, they can yield up to two admissible matrices, $Q_{1}$ and $Q_{2}$. To compute them, we apply Algorithm 1 at every draw of $\phi$ from its posterior, using the Matlab command vpasolve to solve the system of quadratic equations.

We specify the Jeffreys' prior for the reduced-form parameters. Its density function is proportional to $|\Sigma|^{-\frac{3+1}{2}}$. We draw from the posterior 2,000 times and, considering uniquely the zero restrictions in Eq. (35), obtain 2,000 realizations of $\mathcal{Q}_{R}(\hat{\phi})$, each of which is nonempty and consists of two orthogonal matrices $Q_{1}$ and $Q_{2}$. We label them by $Q_{1}$ and $Q_{2}$ according to the ordering of the contemporaneous inflation impulse response.

Figure 4 reports the impulse response to a contractionary monetary policy shock for the output gap (left panel) and inflation (right panel). It shows the posterior means and the highest posterior density regions with credibility $90 \%$ that would be obtained if the conditional prior for $Q$ given $\phi$ assigned all probability mass to either $Q_{1}$ or $Q_{2}$. That is, reporting one of the inference outputs

[^11]corresponds to the Bayesian approach that focuses only on one of the posterior modes, ignoring the other. ${ }^{14}$

The inference result based on $Q_{1}$ shows evidence for both price and output puzzles in the short run. In the medium term, on the other hand, a contractionary monetary policy shock triggers a contraction of the output gap, leaving the price dynamics mostly unaffected. Inference based on the other orthogonal matrix $Q_{2}$, however, leads to a contrasting conclusion. The reaction of prices to the monetary policy shock is significantly negative, while the output gap responds positively and significantly, particularly in the medium-long run.

This example illustrates that different locally-identified observationally equivalent parameter values can lead to strikingly different conclusions, and ignoring this distorts the information contained in the data. A standard off-the-shelf econometric package could uncover just one of the two results. For instance, Gretl and Eviews both return results in line with those obtained through $Q_{1}$. These packages rely on algorithms that maximize the likelihood starting from some initial value without checking other local maxima. Ignoring the other observationally equivalent solutions can distort the information contained in the data. Thus, we recommend checking for the existence of other local maxima and, if any exist, addressing how the conclusions change among them by applying the methods proposed in this paper.

The inference approaches proposed in Section V produce the results reported in Figure 5. The left panels plot results for the output gap while the right panels plot those for inflation. The top panels show the draws of the impulse responses obtained based on the draws of $\phi$ from its posterior. For each draw, we highlight the two observationally equivalent impulse responses corresponding to admissible $Q_{1}$ (blue) and $Q_{2}$ (red) that are coherent with the zero restrictions in Eq. (35). The labels of $Q_{1}$ and $Q_{2}$ in the plots of impulse responses for inflation are maintained in the plots for output.

The middle and bottom panels present interval estimates based on the Bayesian and frequentist inference procedures of Sections V. 1 and V.2. The Bayesian posterior (whose highest density regions are reported both in the middle and bottom panels) is obtained by specifying the uniform conditional prior for $Q$ given $\phi$, i.e., equal weights are assigned to $Q_{1}$ and $Q_{2}$ conditional on $\phi$. The highest posterior density regions are plotted with gradation in gray scale, where the credibility levels vary over $90 \%, 75 \%, 50 \%, 25 \%$, and $10 \%$, from the lightest to darkest. ${ }^{15}$

The middle left panel of 5 shows the marginal posterior distributions for the output gap impulse response. These are unimodal up to $h=4$, but become bimodal for longer horizons. While there is

[^12]evidence for output gap puzzle at the shortest horizons, the probability density is tighter and higher for the negative impulse responses in the medium-long run: the darkest gray region (highest $25 \%$ and $10 \%$ of the distribution) appears mostly for the negative part of the responses. The middle right panel of Figure 5 shows the marginal posterior distribution for the inflation impulse response. This is bimodal up to $h=10$ and becomes unimodal at longer horizons. Similarly to the output gap, for the horizons with bimodal distributions, the negative impulse responses have tighter and higher densities than the positive ones.

For both the output gap and inflation, we also present the frequentist-valid confidence intervals (in dotted-circle lines) proposed in Section V.2. These are obtained by retaining $90 \%$ of the draws of the reduced-form parameters with the highest value of the posterior density function. The middle panels show the constant-label projection confidence sets of Section V.2.2, while the bottom panels show the switching-label projection confidence sets of Section V.2.1. In addition, for both the output gap and inflation, we show the range of posterior means obtained by the robust Bayesian approach (dotted lines). It is evident that the Bayesian approach gives the narrowest interval estimates, and the highest posterior density regions well visualize the bi-modal nature of the posterior distributions at some horizons. The wider confidence intervals of the frequentist approach reflect a couple of their features. First, they are agnostic over the observationally equivalent parameters in the sense that they do not assign any weights over the observationally equivalent impulse responses. Second, our proposed frequentist procedures rely on projecting the joint confidence intervals for the reducedform parameters and do not optimize the width of the interval estimates for impulse responses. Concerning the results of the robust Bayesian approach, the bounds of the set of posterior means are in line with the two modes of the posterior distributions.

A useful strategy for reducing the number of locally-identified admissible solutions is to introduce sign restrictions. To refine the results reported in Figure 5, consider assuming no price puzzle by restricting the inflation responses to be non-positive for (a) the contemporaneous period, or (b) for the four quarters following a contractionary monetary policy shock. The results are reported in Figures 6 and 7, respectively.

The results with the contemporaneous non-positivity restriction (Figure 6) appear similar to those in Figure 5. A notable difference is in the upper bound of the frequentist confidence intervals for the inflation response, which now excludes the positive responses shown in Figure 5 (top-right panel). For the first few quarters, both switching- and constant-label projection confidence sets exclude a region between the two modes of the distribution.

Imposing the four sign restrictions (Figure 7), allows us to eliminate one of the admissible $Q$ matrices for most of the draws of $\phi$. In the top panels of Figure 7, the impulse responses plotted in black have a unique admissible $Q$ under the imposed sign restrictions. A comparison of Figure

5 and Figure 7 shows that the sign restrictions rule out the impulse responses corresponding to $Q_{1}$ matrices. The constant-label and switching-label confidence intervals produce similar results. The only notable difference appears in the response of output gap, where the switching-label confidence intervals in the bottom-left panel have narrow "gaps" from $h=11$ to $h=18$, while the constantlabel confidence intervals in the middle-left panel do not.

Figure 4: Impulse response functions for the New-Keynesian non-recursive SVAR.

(a) Output gap to $\varepsilon_{t}^{m p}$ : Single admissible $Q_{i}$.
(b) Inflation to $\varepsilon_{t}^{m p}$ : Single admissible $Q_{i}$.


Notes: In the left column we report the impulse responses for the output gap obtained as the Bayesian posterior means with the upper and lower bounds of the highest posterior density regions with credibility $90 \%$ obtained through the admissible $Q_{1}$ and $Q_{2}$ matrices, considered separately. Similarly, in the right column we report the impulse responses for inflation.

## VII Conclusion

This paper analyzes SVARs that attain local identification but may fail to attain global identification. We identify the class of identifying restrictions that delivers local but non-global identification. This is characterized by non-homogeneous, non-recursive, and/or across-shock equality restrictions. Exploiting the geometric structure of the identification problem, we propose a novel way to analyze and exhaustively compute the observationally equivalent impulse responses. The novel analytical and computational insights also contribute to the development of a posterior sampling algorithm for Bayesian inference and projection-based frequentist-valid inference in the presence of locally identified parameters.

Since locally- but not globally-identified structural models appear in other macroeconometric models, including heteroskedastic SVARs and the dynamic stochastic general equilibrium models, extending our computational and inference approaches to these models is a promising avenue for future research.

Figure 5: New-Keynesian non-recursive SVAR with zero restrictions only


Notes: The left column reports the output gap impulse responses and the right column reports the inflation impulse responses, both to a contractionary monetary policy shock. The middle and bottom panels report the posterior highest density regions at $90 \%, 75 \%, 50 \%, 25 \%$ and $10 \%$ in gray scale. The upper and loser bounds of the frequentist confidence sets are plotted by the dotted-circle lines. The dotted lines in the middle panels plot the set of posterior means.

Figure 6: New-Keynesian non-recursive SVAR with zero restrictions and one sign restriction


Notes: The left column reports the output gap impulse responses and the right column reports the inflation impulse responses, both to a contractionary monetary policy shock. The middle and bottom panels report the posterior highest density regions at $90 \%, 75 \%, 50 \%, 25 \%$ and $10 \%$ in gray scale. The upper and loser bounds of the frequentist confidence sets are plotted by the dotted-circle lines. The dotted lines in the middle panels plot the set of posterior means.

Figure 7: New-Keynesian non-recursive SVAR with zero and additional four sign restrictions


Notes: The left column reports the output gap impulse responses and the right column reports the inflation impulse responses, both to a contractionary monetary policy shock. The middle and bottom panels report the posterior highest density regions at $90 \%, 75 \%, 50 \%, 25 \%$ and $10 \%$ in gray scale. The upper and loser bounds of the frequentist confidence sets are plotted by the dotted-circle lines. The dotted lines in the middle panels plot the set of posterior means.

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## A Appendix: Some analytical results on the geometry of identification

Let the data generating process be the bivariate VAR defined in Section II. 5 with the identifying restriction

$$
\begin{equation*}
\left(A_{0}\right)_{[1,1]}^{-1}=c \Longleftrightarrow\left(e_{1}^{\prime} \Sigma_{t r}\right) q_{1}=c \tag{55}
\end{equation*}
$$

where $c>0$ is a known (positive) scalar and $e_{1}$ is the first column of the $(2 \times 2)$ identity matrix. The non-homogeneous restriction in Eq. (55) affects the orthogonal matrix $Q$ as $\sigma_{1}^{\prime} q_{1}=c$, with $\sigma_{1}$ denoting the first column of $\Sigma_{t r}^{\prime}=\left(\begin{array}{cc}\sigma_{1,1} & \sigma_{2,1} \\ 0 & \sigma_{2,2}\end{array}\right)$.

The vector $q_{1}$ must satisfy the two equations

$$
\left\{\begin{array}{l}
\sigma_{1}^{\prime} q_{1}=c \\
q_{1}^{\prime} q_{1}=1
\end{array}\right.
$$

By simple algebra, the two solutions are

$$
\begin{equation*}
q_{1}^{(1)}=\binom{c / \sigma_{1,1}}{+\sqrt{\frac{\sigma_{1,1}^{2}-c^{2}}{\sigma_{1,1}^{2}}}} \quad \text { and } \quad q_{1}^{(2)}=\binom{c / \sigma_{1,1}}{-\sqrt{\frac{\sigma_{1,1}^{2}-c^{2}}{\sigma_{1,1}^{2}}}} \tag{56}
\end{equation*}
$$

These two possible solutions are represented in Figure 1. If ( $\sigma_{1,1}^{2}<c^{2}$ ), the straight (vertical) line does not intersect the unit circle, and no real solution is admissible. The SVAR, although identified, does not admit any real solution given the reduced-form parameters $\phi$. If, instead, $\left(\sigma_{1,1}^{2}=c^{2}\right)$, i.e. the vertical red line is tangent to the unit circle, we continue to have global identification, although the imposed restriction is not coherent with those derived by RWZ. In all other situations, there will be two solutions that, a priori, can be admissible despite the sign normalization restriction. This is the case depicted in Figure 1.

Concerning the sign normalization restriction, for the first equation, the definition in Eq. (7) reduces to $q_{1}^{\prime} \tilde{\sigma}_{1} \geq 0$, where $\tilde{\sigma}_{1}$ is the first column of $\Sigma_{t r}^{-1}=1 /\left(\sigma_{1,1} \sigma_{2,2}\right)\left(\begin{array}{cc}\sigma_{2,2} & 0 \\ -\sigma_{2,1} & \sigma_{1,1}\end{array}\right)$. Through elementary algebra, we obtain that

$$
\begin{equation*}
q_{1}^{\prime} \tilde{\sigma}_{1} \geq 0 \quad \Longleftrightarrow \quad \frac{q_{1,1}}{\sigma_{1,1}} \geq \frac{q_{1,2} \sigma_{2,1}}{\sigma_{1,1} \sigma_{2,2}} \tag{57}
\end{equation*}
$$

where $q_{1,1}$ and $q_{1,2}$ are the two generic elements of $q_{1}$, i.e. $q_{1}=\left(q_{1,1}, q_{1,2}\right)^{\prime}$. Suppose, first, that from the data we have $\sigma_{2,1}<0$. In this case, if we substitute in the values of $q_{1,1}$ and $q_{1,2}$ obtained for $q_{1}^{(1)}$ in the left-hand side of Eq. (56), the sign normalization condition for the first equation becomes

$$
\begin{equation*}
\frac{c}{\sigma_{1,1}^{2}} \geq \frac{\sigma_{2,1}}{\sigma_{1,1} \sigma_{2,2}} \sqrt{\frac{\sigma_{1,1}^{2}-c^{2}}{\sigma_{1,1}^{2}}} \tag{58}
\end{equation*}
$$

As the left-hand side is always positive and the right-hand side always negative, this is always satisfied. If, instead, we substitute the values $q_{1,1}$ and $q_{1,2}$ obtained for $q_{1}^{(2)}$ in the right-hand side of Eq. (56), the sign normalization condition for the first equation becomes

$$
\begin{equation*}
\frac{c}{\sigma_{1,1}^{2}} \geq-\frac{\sigma_{2,1}}{\sigma_{1,1} \sigma_{2,2}} \sqrt{\frac{\sigma_{1,1}^{2}-c^{2}}{\sigma_{1,1}^{2}}} \tag{59}
\end{equation*}
$$

that is also satisfied when $c^{2} \geq \frac{1}{2} \frac{\sigma_{1,1}^{2} \sigma_{2,1}^{2}}{\sigma_{2,2}^{2}}$. If this is the case, both solutions $q_{1}^{(1)}$ and $q_{1}^{(2)}$ are admissible, leading to local identification. The situation is very similar when $\sigma_{2,1}>0$.

If, instead, $c=0$ as in the standard RWZ setup, the two $q_{1}$ vectors in Eq. (56) become $q_{1}^{(1)}=(0,1)^{\prime}$ and $q_{1}^{(2)}=(0,-1)^{\prime}$. If, as before, we suppose $\sigma_{2,1}<0$, the sign normalization for $q_{1}^{(1)}$ in Eq. (58) reduces to $0 \geq \sigma_{2,1} /\left(\sigma_{1,1} \sigma_{2,2}\right)$, which is always true. The sign normalization for $q_{1}^{(2)}$ in Eq. (59) is $0 \geq-\sigma_{2,1} /\left(\sigma_{1,1} \sigma_{2,2}\right)$ which, in contrast, is never true. The case where $\sigma_{2,1}>0$, is exactly the same, but with inverted results. One of the two solutions, thus, will be always ruled out by the sign normalization, and global identification is guaranteed.

The second column of $Q$, the unit-length vector $q_{2}$, although not restricted, can be pinned down through the its orthogonality to $q_{1}$

$$
\left\{\begin{array}{l}
q_{2}^{\prime} q_{1}=0  \tag{60}\\
q_{2}^{\prime} q_{2}=1
\end{array}\right.
$$

However, given that there are two admissible vectors $q_{1}^{(1)}$ and $q_{1}^{(2)}$, the system Eq. (60) must be solved for both. This can be done with simple algebra, yielding the two solutions

$$
\begin{equation*}
q_{2}^{(1)}=\binom{+\sqrt{\frac{\left(\sigma_{1,1}^{2}-c^{2}\right)\left(2 c^{2}-\sigma_{1,1}^{2}\right)}{c^{4}}}}{-\sqrt{\frac{2 c^{2}-\sigma_{1,1}^{2}}{c^{2}}}} \quad \text { and } \quad q_{2}^{(1)}=\binom{-\sqrt{\frac{\left(\sigma_{1,1}^{2}-c^{2}\right)\left(2 c^{2}-\sigma_{1,1}^{2}\right)}{c^{4}}}}{+\sqrt{\frac{2 c^{2}-\sigma_{1,1}^{2}}{c^{2}}}} \tag{61}
\end{equation*}
$$

One of the two, precisely which depends on the reduced-form parameters, will be eliminated by the sign normalization restriction. This case is represented in the left panel of Figure 2, together with $q_{1}^{(1)}$.

The other possibility, represented in the right panel of Figure 2, is when we solve the system conditional on $q_{1}^{(2)}$, obtaining

$$
\begin{equation*}
q_{2}^{(2)}=\binom{+\sqrt{\frac{\left(\sigma_{1,1}^{2}-c^{2}\right)\left(2 c^{2}-\sigma_{1,1}^{2}\right)}{c^{4}}}}{+\sqrt{\frac{2 c^{2}-\sigma_{1,1}^{2}}{c^{2}}}} \quad \text { and } \quad q_{2}^{(2)}=\binom{-\sqrt{\frac{\left(\sigma_{1,1}^{2}-c^{2}\right)\left(2 c^{2}-\sigma_{1,1}^{2}\right)}{c^{4}}}}{-\sqrt{\frac{2 c^{2}-\sigma_{1,1}^{2}}{c^{2}}}} \tag{62}
\end{equation*}
$$

where, as before, one of the two solutions is ruled out by the sign normalization restriction.

## B Appendix: Further results on local identification

In this appendix we provide a new result on local identification for SVAR models. We consider a set of equality restrictions $\mathbf{F}(\phi, Q)$ satisfying the recursive identification scheme in Definition 3.

Proposition 7 (RWZ sufficient condition for checking local identification). Consider an SVAR with recursive identifying restrictions of the form Eq. (14). The SVAR is locally identified at
$A=\left(A_{0}, A_{+}\right) \in \mathcal{A}_{R}$ if, for $i=1, \ldots, n$,

$$
M_{i}(Q) \equiv\left(\begin{array}{c}
F_{i i}(\phi)  \tag{63}\\
\underset{(n-i) \times n}{ } \underset{n \times n}{Q} \\
\left(\begin{array}{cc}
I_{i} & 0 \\
i \times i & i \times(n-i)
\end{array}\right)
\end{array}\right)
$$

is of rank n.
Proof. See Appendix C.
Proposition 7 reconciles our condition for local identification of recursive SVARs with the general rank condition for global identification provided by RWZ (their Theorem 1). In particular, under a recursive identification scheme, the RWZ condition for global identification developed for the case of homogeneous restrictions implies local identification, even though we allow non-homogeneous and across shock restrictions.

## C Appendix: Proofs

This appendix collects proofs for all propositions reported in this paper. We make use of the following matrices. $K_{n}$ is the $n^{2} \times n^{2}$ commutation matrix as defined in Magnus and Neudecker (2007) and $N_{n}=1 / 2\left(I_{n^{2}}+K_{n}\right)$. Let $\tilde{D}_{n}$ be the $n^{2} \times n(n-1) / 2$ full-column rank matrix $\tilde{D}_{n}$ defined in Magnus (1988) such that for any $n(n-1) / 2$-dimensional vector $v, \tilde{D}_{n} v \equiv \operatorname{vec}(H)$ holds, where $H$ is an $n \times n$ skew-symmetric matrix $\left(H=-H^{\prime}\right)$. See Appendix D for explicit constructions of $\tilde{D}_{n}$ for $n=2,3,4$.

## Proof of Proposition 2: necessary and sufficient condition for local identification

Fixing $\phi$, a matrix $Q$ satisfies the identifying restrictions if:

$$
\begin{align*}
\mathbf{F}(\phi) \operatorname{vec} Q & =\mathbf{c}  \tag{64}\\
Q^{\prime} Q & =I_{n} \tag{65}
\end{align*}
$$

which is a system of quadratic equations. Eq. (64) consists of $f=f_{1}+\cdots+f_{n}$ linear and nonhomogeneous equations. Eq. (65) is a set of quadratic equations stating that the columns of $Q$, the vectors $\left(q_{1}, \ldots, q_{n}\right)$, must be orthogonal and of unit length.

The system can be solved locally as:

$$
\begin{aligned}
\mathbf{F}(\phi) \operatorname{vec} d Q & =0 \\
d Q^{\prime} Q+Q^{\prime} d Q & =0
\end{aligned}
$$

which, using the Kronecker product and its properties, becomes

$$
\begin{aligned}
\mathbf{F}(\phi) \operatorname{vec} d Q & =0 \\
{\left[\left(Q^{\prime} \otimes I_{n}\right)+\left(I_{n} \otimes Q^{\prime}\right)\right] \operatorname{vec} d Q } & =0
\end{aligned}
$$

Moreover, using the commutation matrix $K_{n}$ we have

$$
\begin{aligned}
\mathbf{F}(\phi) \operatorname{vec} d Q & =0 \\
{\left[K_{n}\left(I_{n} \otimes Q^{\prime}\right)+\left(I_{n} \otimes Q^{\prime}\right)\right] \operatorname{vec} d Q } & =0
\end{aligned}
$$

and recalling $N_{n}=1 / 2\left(I_{n^{2}}+K_{n}\right)$, we obtain

$$
\begin{aligned}
\mathbf{F}(\phi) \operatorname{vec} d Q & =0 \\
2 N_{n}\left(I_{n} \otimes Q^{\prime}\right) \operatorname{vec} d Q & =0 .
\end{aligned}
$$

The Jacobian matrix can, thus, be defined as

$$
\begin{equation*}
J(Q)=\binom{\mathbf{F}(\phi)}{2 N_{n}\left(I_{n} \otimes Q^{\prime}\right)} \tag{66}
\end{equation*}
$$

Following Magnus and Neudecker (2007), a sufficient condition for local identification of $Q$ at the point $Q=Q_{0}$ is that $J\left(Q_{0}\right)$ has full column rank. If there exists an admissible neighborhood of $Q_{0}$ such that $J\left(Q_{0}\right)$ is of full column rank, this condition becomes necessary too.

The condition regarding the rank of Eq. (66) can be further simplified. Given that $Q$ is invertible (it is orthogonal), the rank of $J(Q)$ is unchanged if we post-multiply Eq. (66) by ( $\left.I_{n} \otimes Q^{-1 \prime}\right)=I_{n} \otimes Q$. Checking whether $J(Q)$ is of full column rank, thus, corresponds to checking whether the system of equations

$$
\begin{aligned}
\mathbf{F}(\phi)\left(I_{n} \otimes Q\right) x & =0 \\
2 N_{n} x & =0
\end{aligned}
$$

admits the null vector $x$ as the unique solution. However, as in Magnus (1988), the second equation, can be solved as $x=\tilde{D}_{n} z$, with $z$ a $n(n-1) / 2 \times 1$ vector. Substituting this solution into the first equation leads to the rank condition Eq. (21) of Proposition 3. Since $\tilde{D}_{n}$ is a matrix of full column rank rank $n(n-1) / 2$, a necessary condition for the rank condition Eq. (21) is that the number of rows of $F(\phi), f$, is greater than or equal to $n(n-1) / 2$. This completes the proof of (i).

To show claim (ii), let $\overline{\mathcal{F}}$ be the set of matrices of dimension $f \times n(n-1) / 2$ and denote by $X$ a generic element of $\overline{\mathcal{F}}$. Viewing the space spanning the $j$-th column of $X$ as $V_{j}$ in Lemma 3 of RWZ, and defining the set $S$ in Lemma 3 of RWZ to be the set of matrices with deficient rank $S=\{X \in \overline{\mathcal{F}}: \operatorname{rank}(X)<n(n-1) / 2\}$, Lemma 3 in RWZ shows that either $S=\overline{\mathcal{F}}$, or $S$ is a set of measure zero in $\overline{\mathcal{F}}$.

Define

$$
\begin{equation*}
\mathcal{F} \equiv\left\{\mathbf{F}(\phi)\left(I_{n} \otimes Q\right) \tilde{D}_{n}: \mathbf{F}(\phi) \operatorname{vec}(Q)=c, \quad(\phi, Q) \in \Phi \times \mathcal{O}(n)\right\} \tag{67}
\end{equation*}
$$

Since $\mathcal{F} \subset \overline{\mathcal{F}}, S \cap \mathcal{F}$ is either equal to $\mathcal{F}$ or is a set of measure zero in $\mathcal{F}$. Let $g: \mathcal{A} \rightarrow \Phi \times \mathcal{O}(n)$ be the function that reparametrizes the structural parameters $A$ to $(\phi, Q)$, and $h:(\Phi \times \mathcal{O}(n)) \rightarrow \mathcal{F}$ be the function that maps $(\phi, Q)$ to $\left[\mathbf{F}(\phi)\left(I_{n} \otimes Q\right) \tilde{D}_{n}\right] \in \mathcal{F}$. By applying Lemma 2 in RWZ (proved in Spivak, 1965) to the chain of inverse maps $h^{-1}$ and $g^{-1}$, we conclude that either $\left(g^{-1} \circ h^{-1}\right)(\mathcal{F})=\mathcal{A}_{R}$ or it is of measure zero in $\mathcal{A}_{R}$. The conclusion then follows by noting $\left(g^{-1} \circ h^{-1}\right)(\mathcal{F})=\mathcal{K}^{c}$.

## Proof of Proposition 3: local identification in recursive SVARs

Assume that the rank condition of Proposition 2 holds at parameter point $A=\left(A_{0}, A_{+}\right) \in \mathcal{A}_{R}$, and let $\phi$ be the corresponding reduced-form parameter. Since local identification holds at $A$, there is no observationally equivalent parameter point in a neighborhood of $A$. In other words, no infinitesimal rotation of the orthogonal matrix $Q$ generates observationally equivalent and admissible structural parameters in the neighborhood of $A$. Any infinitesimal rotation can be represented by $\left(I_{n}+H\right)$, where $H$ is an $n \times n$ skew-symmetric matrix (see Lucchetti 2006) whose $i$-th column we denote by $h_{i}$.

Projecting on $q_{1}$, an admissible structural parameter lying in a local neighborhood of $A$ has to satisfy

$$
F_{11}(\phi)\left[Q\left(I_{n}+H\right)\right] e_{1}=c_{1} \quad \Longrightarrow \quad F_{11}(\phi) q_{1}+F_{11} Q H e_{1}=c_{1} \quad \Longrightarrow \quad F_{11}(\phi) Q h_{1}=0
$$

where $e_{i}$ is the $i$-th column of the identity matrix $I_{n}$, and the last equation follows from the fact that $F_{11}(\phi) q_{1}=c_{1}$. The system $F_{11}(\phi) Q h_{1}=c_{1}$ is characterized by $n-1$ equations and an $n$ dimensional $h_{1}$. The first element of $h_{1}$ is zero by definition (the elements on the main diagonal of a skew-symmetric matrix are equal to zero). Hence, we have

$$
\begin{equation*}
\binom{F_{11}(\phi) Q}{e_{1}^{\prime}} h_{1}=0 \tag{68}
\end{equation*}
$$

This linear equation system has $h_{1}=0$ as its unique solution if and only if $\binom{F_{11}(\phi) Q}{e_{1}^{\prime}}$ is of rank $n$, or, equivalently, $F_{11}(\phi)\left(q_{2} \ldots q_{n}\right)$ is of full rank (equal to $n-1$ ). Since the model is locally identified by assumption, $h_{1}=0$ has to be the only solution of Eq. (68). Hence, $\operatorname{rank}\left(F_{11}(\phi)\left(q_{2} \ldots q_{n}\right)\right)=n-1$ must hold, implying that $\operatorname{rank}\left(F_{11}(\phi)\right)=n-1$.

For $q_{2}$, given a $q_{1}$ vector solving $F_{11}(\phi) q_{1}=0$, we have the following system:

$$
\left\{\begin{aligned}
F_{21}(\phi) q_{1}+F_{22}(\phi) q_{2} & =c_{2} \\
q_{1}^{\prime} q_{2} & =0,
\end{aligned}\right.
$$

Considering again an infinitesimal rotation
$\left\{\begin{aligned} & F_{21}(\phi) q_{1}+F_{22}(\phi) Q\left(I_{n}+H\right) e_{2}= \\ & c_{2} \\ & q_{1}^{\prime} Q\left(I_{n}+H\right) e_{2}=\end{aligned} \quad 0 \quad \Longrightarrow\left\{\begin{array}{rl}F_{21}(\phi) q_{1}+F_{22}(\phi) Q e_{2}+F_{22}(\phi) Q h_{2} & =c_{2} \\ q_{1}^{\prime} Q h_{2} & =\end{array} \quad 0\right.\right.$,
but, given the restrictions, $F_{21}(\phi) q_{1}+F_{22}(\phi) q_{2}=c_{2} \quad \Longrightarrow \quad F_{22}(\phi) Q h_{2}=0$, which allows the system to be written as

$$
\left\{\begin{array}{rl}
F_{22}(\phi) Q h_{2} & =0  \tag{69}\\
q_{1}^{\prime} Q h_{2} & =0
\end{array} \quad \Longrightarrow \quad\binom{F_{22}(\phi) Q}{q_{1}^{\prime} Q} h_{2}=0\right.
$$

Similarly to the argument for $h_{1}$ above, and noting that the first two entries of $h_{2}$ are zero, we can represent the linear equations as

$$
\left(\begin{array}{c}
F_{22}(\phi) Q \\
q_{1}^{\prime} Q \\
e_{2}^{\prime} \\
e_{1}^{\prime}
\end{array}\right) h_{2}=0
$$

Since $q_{1}^{\prime} Q=e_{1}^{\prime}$, the last equation in this system is redundant. Thus, in order for $h_{2}=0$ to be the unique solution, $\binom{F_{22}(\phi)}{q_{1}^{\prime}}$ must have full raw rank (equal to $n-1$ ).

To obtain the sequential rank conditions of Proposition 3, we repeat this argument further for $i=3,4$

Next, we show the reverse implication. For each column of $Q$, we consider a system of equations of the form,

$$
\left\{\begin{aligned}
\tilde{F}_{i i}(\phi) q_{i} & =\left(c_{i}^{\prime}, 0, \ldots, 0\right)^{\prime} \\
q_{i}^{\prime} q_{i} & =1
\end{aligned}\right.
$$

sequentially for $i=1, \ldots, n$, where $\tilde{F}_{i i}(\phi)$ is as defined in the statement of Proposition 3. If $\operatorname{rank}\left(\tilde{F}_{i i}(\phi)\right)=n-1$, the system of equations represents the intersection between a straight line and the unit circle in $\mathbb{R}^{n}$, which has at most two distinct solutions. Hence, any admissible $Q$ matrices are isolated points, so the SVAR is locally identified. The rank condition of Eq. (21) follows by Proposition 2.

## Proof of Proposition 4: number of admissible $Q$ 's

We split the proof into five cases based on the type of equality restrictions. The first three cases are recursive identification schemes. The remaining two are non-recursive.

We first consider cases with recursive restrictions. That is, the variables are ordered to satisfy

$$
\begin{equation*}
f_{1} \geq f_{2} \geq \ldots \geq f_{n} . \tag{70}
\end{equation*}
$$

Case 1: Recursive and homogeneous restrictions but no restrictions across shocks
Under recursive restrictions, we have shown in Proposition 3 that the rank condition of Eq. (21) is equivalent to the sequential rank conditions of Eq. (23). If the sign normalization restrictions select either the admissible $q_{i}$ or $-q_{i}$ at every $i=1, \ldots, n$, the sequential determination procedure of RWZ pins down an admissible $Q$ matrix. The sequential rank conditions do not guarantee that the sign normalizations select a unique $Q$ matrix, but the number of solutions for each $q_{i}$ is at most two. Hence, the number of admissible $Q$ matrices is at most equal to the number of distinct selections of two vectors $\left\{q_{i},-q_{i}\right\}$ over $i=1, \ldots, n$, which amounts to $2^{n}$.

Case 2: Recursive non-homogeneous restrictions but no restrictions across shocks
Under recursive and non-homogeneous restrictions, consider solving for the admissible $Q$ matrices column by column by exploiting the sequential rank conditions Eq. (23). For the first column $q_{1}$,
we have

$$
\left\{\begin{align*}
F_{11}(\phi) q_{1} & =c_{1}  \tag{71}\\
q_{1}^{\prime} q_{1} & =1
\end{align*}\right.
$$

Given that $F_{11}(\phi)$ has full row rank, the set of solutions of $q_{1}$ for the first equations can be spanned by any $n \times 1$ vector $t_{1} \in \mathbb{R}$,

$$
\begin{align*}
q_{1} & =F_{11}(\phi)^{\prime}\left(F_{11}(\phi) F_{11}(\phi)^{\prime}\right)^{-1} c_{1}+\left(I_{n}-F_{11}(\phi)^{\prime}\left(F_{11}(\phi) F_{11}(\phi)^{\prime}\right)^{-1} F_{11}(\phi)\right) t_{1} \\
& \equiv d_{1}+B_{1} t_{1} \tag{72}
\end{align*}
$$

Since the $(n \times n)$ matrix $B_{1}$ has rank $n-f_{1}=1$, it can be decomposed as $B_{1}=\alpha_{1} \beta_{1}^{\prime}$, where $\alpha_{1}$ is a basis for span $\left(B_{1}\right)$, i.e. the column space of $B_{1}$, and both $\alpha_{1}$ and $\beta_{1}$ are non-zero $n \times 1$ vectors. We can hence write

$$
\begin{equation*}
q_{1}=d_{1}+\alpha_{1} z_{1} \tag{73}
\end{equation*}
$$

with $z_{1}=\beta_{1}^{\prime} t_{1}$, being any scalar. The second (quadratic) equation in system Eq. (71) becomes

$$
\begin{aligned}
q_{1}^{\prime} q_{1} & =\left(d_{1}+\alpha_{1} z_{1}\right)^{\prime}\left(d_{1}+\alpha_{1} z_{1}\right) \\
& =d_{1}^{\prime} d_{1}+2 d_{1}^{\prime} \alpha_{1} z_{1}+\alpha_{1}^{\prime} \alpha_{1} z_{1}^{2}=1 \\
& \Rightarrow \lambda_{1}+2 \xi_{1} z_{1}+\omega_{1} z_{1}^{2}=0
\end{aligned}
$$

where $\lambda_{1}=d_{1}^{\prime} d_{1}-1, \xi_{1}=d_{1}^{\prime} \alpha_{1}$ and $\omega_{1}=\alpha_{1}^{\prime} \alpha_{1}$ are all functions of the reduced form parameters. There are hence three possibilities:

1. If $\xi_{1}^{2}-\lambda_{1} \omega_{1}>0$, we have two real solutions. It may be that none, one, or both satisfy the sign normalization restriction for $q_{1}$.
2. If $\xi_{1}^{2}-\lambda_{1} \omega_{1}=0$, we have a single real solution. It may or may not satisfy the sign normalization restriciton.
3. If $\xi_{1}^{2}-\lambda_{1} \omega_{1}<0$, we have no real solution, implying that $\phi$ is not compatible with the imposed restrictions.

In summary, at most there are two admissible $q_{1}$ 's. Denote them by $q_{1}^{(1)}$ and $q_{1}^{(2)}$ (allowing $q_{1}^{(1)}=$ $\left.q_{1}^{(2)}\right)$.

Given an admissible $q_{1} \in\left\{q_{1}^{(1)}, q_{1}^{(2)}\right\}$, consider obtaining an admissible second column vector $q_{2}$ by solving

$$
\left\{\begin{align*}
F_{22}(\phi) q_{2} & =c_{2}  \tag{74}\\
q_{1}^{\prime} q_{2} & =0 \\
q_{2}^{\prime} q_{2} & =1
\end{align*}\right.
$$

with $\operatorname{rank}\left(\left(F_{22}(\phi)^{\prime}, q_{1}\right)\right)=n-1$. This system can be transformed as

$$
\left\{\begin{array} { r l } 
{ F _ { 2 2 } ( \phi ) q _ { 2 } } & { = c _ { 2 } }  \tag{75}\\
{ q _ { 1 } ^ { \prime } q _ { 2 } } & { = 0 } \\
{ q _ { 2 } ^ { \prime } q _ { 2 } } & { = 1 }
\end{array} \Longrightarrow \left\{\begin{array}{rl}
\binom{F_{22}(\phi)}{q_{1}^{\prime}} q_{2}=\binom{c_{2}}{0} \Longrightarrow\left\{\begin{aligned}
\tilde{F}_{22}(\phi) q_{2} & =\tilde{c}_{2} \\
q_{2}^{\prime} q_{2} & =1
\end{aligned}\right\}=1
\end{array}\right.\right.
$$

where $\tilde{F}_{22}(\phi)=\left(F_{22}^{\prime}(\phi), q_{1}\right)^{\prime}$ and $\tilde{c}_{2}=\left(c_{2}^{\prime}, 0\right)^{\prime}$. Given the assumption rank $\left(\tilde{F}_{22}(\phi)\right)=n-1$, Eq. (75) can be solved in the same way as the system for $q_{1}$. We can hence obtain at most two admissible $q_{2}$ vectors for each of $q_{1}=q_{1}^{(1)}$ and $q_{1}=q_{1}^{(2)}$. So far there are at most four admissible vectors for the first two columns of $Q$.

We repeat this argument for $i=3, \ldots, n$. Given that there are at most $2^{i-1}$ admissible constructions of $\left(q_{1}, \ldots, q_{i-1}\right)$, and at each admissible $\left(q_{1}, \ldots, q_{i-1}\right)$, we solve for $q_{i}$

$$
\left\{\begin{align*}
\tilde{F}_{i i}(\phi) q_{i} & =\tilde{c}_{i}  \tag{76}\\
q_{i}^{\prime} q_{i} & =1
\end{align*}\right.
$$

where

$$
\tilde{F}_{i i}(\phi)=\left(F_{i i}(\phi)^{\prime}, q_{1}, \ldots, q_{i-1}\right)^{\prime} \quad \text { and } \quad \tilde{c}_{i}=\left(c_{i}^{\prime}, 0, \ldots, 0\right)
$$

Again, finding an admissible $q_{i}$ given $\left(q_{1}, \ldots, q_{i-1}\right)$ boils down to solving a quadratic equation, so there are at most two solutions for $q_{i}$, implying that there are at most $2^{i}$ admissible constructions of $\left(q_{1}, \ldots, q_{i-1}, q_{i}\right)$. At $i=n$, we obtain at most $2^{n}$ admissible $Q$ matrices.

Case 3: Recursive non-homogeneous restrictions and restrictions across shocks
The recursive restrictions imply that $\mathbf{F}(\phi)$ is lower block-triangular, i.e. $F_{i j}=0$ for $j>i$, and $f_{i}=n-i$ for all $i=1, \ldots, n$. The case where $i=1$ is identical to the initial step in Case 2 above, so we have at most two admissible $q_{1}$ vectors. For $i>1$ we exploit the sequential structure of the restrictions and obtain each admissible $q_{i}$ sequentially given $\left(q_{1}, \ldots, q_{i-1}\right)$ obtained in the preceding steps. The only difference with respect to case 2 is that, once $\left(q_{1}, \ldots, q_{i-1}\right)$ is given, the system of equations in Eq. (76), will be characterized by

$$
\begin{equation*}
\tilde{F}_{i i}(\phi)=\left(F_{i i}^{\prime}(\phi), q_{1}, \ldots, q_{i-1}\right)^{\prime}, \quad \text { and } \quad \tilde{c}_{i}=\left(\left(c_{i}-F_{i 1}(\phi) q_{1}-\cdots-F_{i, i-1}(\phi) q_{i-1}\right)^{\prime}, 0, \ldots, 0\right)^{\prime} \tag{77}
\end{equation*}
$$

Repeating the argument of Case 2 , we conclude there are at most $2^{n}$ admissible $Q \in \mathcal{O}(n)$.
We now move to the cases with non-recursive identifying restrictions.
Case 4: Non-recursive restrictions and no restrictions across shocks
If $f_{1}=n-1$, we can proceed as in Case 2 and globally or locally identify $q_{1}$, depending on the restrictions at hand. If, instead, $f_{1}<n-1$, we can only identify the basis spanning a subspace in $\mathbb{R}^{n}$ of dimension $n-f_{1}$ containing $q_{1}$. The system of equations characterizing $q_{1}$ is given by

$$
\left\{\begin{align*}
F_{11}(\phi) q_{1} & =c_{1}  \tag{78}\\
q_{1}^{\prime} q_{1} & =1
\end{align*}\right.
$$

Following the analysis of Case 2 , we can represent an admissible $q_{1}$ by $q_{1}=d_{1}+\alpha_{1} z_{1}$, where $z_{1}=\beta_{1}^{\prime} t_{1} \in \mathbb{R}^{n-f_{1}}, \alpha_{1}$ is a nonzero $n \times\left(n-f_{1}\right)$ matrix, $\beta_{1}$ is a nonzero $\left(n-f_{1}\right) \times n$ matrix, and $t_{1} \in \mathbb{R}^{n}$. Given this representation of $q_{1}$, the second (quadratic) equation in system Eq. (78) becomes

$$
\begin{aligned}
q_{1}^{\prime} q_{1} & =d_{1}^{\prime} d_{1}+2 d_{1}^{\prime} \alpha_{1} z_{1}+z_{1}^{\prime} \alpha_{1}^{\prime} \alpha_{1} z_{1}=1 \\
& \Rightarrow \lambda_{1}+2 \xi_{1}^{\prime} z_{1}+z_{1}^{\prime} \omega_{1} z_{1}=0
\end{aligned}
$$

where $\lambda_{1}=d_{1}^{\prime} d_{1}-1, \xi_{1}=\alpha_{1}^{\prime} d_{1}$, and $\omega_{1}=\alpha_{1}^{\prime} \alpha_{1}$. The set of real roots of this quadratic equation in $z_{1}$, if nonempty, is a singleton or a hyper-ellipsoid in $\mathbb{R}^{n-f_{1}}$ with its radius given by the constant term in the completion of squares (if nonnegative).

Assuming an admissible $q_{1}$ exists, consider the equation system for $q_{2}$,

$$
\left\{\begin{align*}
q_{2} & =d_{2}+\alpha_{2} z_{2}  \tag{79}\\
q_{2}^{\prime} q_{2} & =1,
\end{align*}\right.
$$

whose set of roots, if nonempty, is again a singleton or a $n-f_{2}$-dimensional hyper-ellipsoid. In addition, we have the following orthogonality restriction between $q_{1}$ and $q_{2}$,

$$
\begin{aligned}
q_{1}^{\prime} q_{2} & =d_{1}^{\prime} d_{2}+d_{1}^{\prime} \alpha_{2} z_{2}+z_{1}^{\prime} \alpha_{1}^{\prime} d_{2}+z_{1}^{\prime} \alpha_{1}^{\prime} \alpha_{2} z_{2} \\
& \equiv \lambda_{1,2}+\xi_{1,2}^{\prime} z_{2}+z_{1}^{\prime} \xi_{2,1}+z_{1}^{\prime} \omega_{1,2} z_{2}=0 .
\end{aligned}
$$

Enumerating these equations for all $i=1, \ldots, n$, we obtain the following system of equations:

$$
\left\{\begin{align*}
z_{1}^{\prime} \omega_{1} z_{1}+2 \xi_{1}^{\prime} z_{1}+\lambda_{1} & =0  \tag{80}\\
z_{2}^{\prime} \omega_{2} z_{2}+2 \xi_{2}^{\prime} z_{2}+\lambda_{2} & =0 \\
\vdots & \\
z_{n}^{\prime} \omega_{n} z_{n}+2 \xi_{n}^{\prime} z_{n}+\lambda_{n} & =0 \\
z_{1}^{\prime} \omega_{1,2} z_{2}+\xi_{1,2}^{\prime} z_{2}+z_{1}^{\prime} \xi_{2,1}+\lambda_{1,2} & =0 \\
z_{1}^{\prime} \omega_{1,3} z_{3}+\xi_{1,3}^{\prime} z_{3}+z_{1}^{\prime} \xi_{3,1}+\lambda_{1,3} & =0 \\
\vdots & \\
z_{n-1}^{\prime} \omega_{n-1, n} z_{n}+\xi_{n-1, n}^{\prime} z_{n}+z_{n-1}^{\prime} \xi_{n, n-1}+\lambda_{n-1, n} & =0
\end{align*}\right.
$$

The number of equations is $n+n(n-1) / 2=n(n+1) / 2$. The number of unknowns, contained in $z_{1}, z_{2}, \ldots, z_{n}$, is

$$
\begin{equation*}
\left(n-f_{1}\right)+\left(n-f_{2}\right)+\ldots+\left(n-f_{n}\right) \leq n^{2}-n(n-1) / 2=n(n+1) / 2 \tag{81}
\end{equation*}
$$

where the inquality follows by the order condition stated in Proposition $2, \sum_{i=1}^{n} f_{i} \geq n(n-1) / 2$. Hence, we have a system of $n(n+1) / 2$ equations with at most $n(n+1) / 2$ unknowns. Moreover, each one is a quadratic equation and, importantly, given the rank condition for local identification is satisfied, each of the solutions has to be an isolated point. Bézout's theorem gives that the maximum number of solutions is the product of the polynomial degree of all the equations, so the number of solutions is at most $2^{n(n+1) / 2}$.

Case 5: Non-recursive and across-shocks restrictions
In this case analysis of identification requires considering all equations jointly. We will have a
system of equations of the form

$$
\left\{\begin{align*}
\mathbf{F}(\phi) \operatorname{vec} Q & =\mathbf{c}  \tag{82}\\
q_{1}^{\prime} q_{1} & =1 \\
q_{2}^{\prime} q_{2} & =1 \\
\vdots & \\
q_{n}^{\prime} q_{n} & =1 \\
q_{1}^{\prime} q_{2} & =0 \\
\vdots & \\
q_{n-1}^{\prime} q_{n} & =0
\end{align*}\right.
$$

This system consists of $n^{2}$ equations with $n^{2}$ unknowns (the elements in $Q$ ). The first $n(n-1) / 2$ equations are linear and the latter $n(n+1) / 2$ equations are all quadratic. By Bézout's theorem, the maximum number of solutions is at most $2^{n(n+1) / 2}$.

## Proof of Proposition 7: RWZ sufficient condition for checking local identification

The result is a by-product of Proposition 3. As observed in Eq. (68), the first column $q_{1}$ is locally identified if and only if $\binom{F_{11}(\phi) Q}{e_{1}^{\prime}}$ has full column rank equal to $n$. When moving to the identification of $q_{2}$, from the system Eq. (69), and recalling that the first two elements of $h_{2}$ are zero, we have no admissible infinitesimal rotation (i.e. $h_{2}=0$ ) if

$$
\left.\operatorname{rank}\left(\begin{array}{cc}
F_{22}(\phi) \cdot \underset{ }{(n-2) \times n} & \begin{array}{c}
Q \times n \\
I_{2}
\end{array} \\
\left(\begin{array}{ll}
0 \times(n-2)
\end{array}\right.
\end{array}\right)\right)=n
$$

Repeating this argument for the remaining of columns of $Q$, we obtain the proposition.

## Proof of Proposition 5: privatization policy SVAR

Denote the column vectors of $\Sigma_{t r}^{-1}$ and $\Sigma_{t r}^{\prime}$ by $\Sigma_{t r}^{-1}=\left(\tilde{\sigma}_{1}, \tilde{\sigma}_{2}, \tilde{\sigma}_{3}\right)$ and $\Sigma_{t r}^{\prime}=\left(\sigma_{1}, \sigma_{2}, \sigma_{3}\right)$. We denote the $i$-th entry of $\tilde{\sigma}_{j}$ and $\sigma_{j}$ by $\tilde{\sigma}_{j, i}$ and $\sigma_{j, i}$, respectively. We analyze the identification of the structural parameters through sequential determination of the column vectors of $Q$ given reduced-form parameters $\Sigma$.

The first vector $q_{1}$ of the orthogonal matrix $Q$ can be determined by the first two restrictions in Eq. (27). The system of quadratic equations to be solved is

$$
\left\{\begin{array}{l}
\tilde{\sigma}_{3,1} q_{11}+\tilde{\sigma}_{3,2} q_{12}+\tilde{\sigma}_{3,3} q_{13}=0  \tag{83}\\
\sigma_{2,1} q_{11}+\sigma_{2,2} q_{12}+\sigma_{2,3} q_{13}=c \\
q_{11}^{2}+q_{12}^{2}+q_{13}^{2}=1
\end{array}\right.
$$

Since $\tilde{\sigma}_{3,1}=0$ and $\tilde{\sigma}_{3,2}=0$, the first equation reduces to $q_{13}=0$. The second equation then
becomes

$$
\sigma_{2,1} q_{11}+\sigma_{2,2} q_{12}=c \quad \Longrightarrow \quad q_{11}=-\frac{\sigma_{2,2}}{\sigma_{2,1}} q_{12}+\frac{c}{\sigma_{2,1}} \equiv \gamma_{1} q_{12}+\gamma_{2},
$$

where $\gamma_{1}=-\frac{\sigma_{2,2}}{\sigma_{2,1}}$ and $\gamma_{2}=\frac{c}{\sigma_{2,1}}$. Substituting into the third equation, we obtain

$$
\begin{equation*}
\left(\gamma_{1}^{2}+1\right) q_{12}^{2}+2 \gamma_{1} \gamma_{2} q_{12}+\left(\gamma_{2}^{2}-1\right)=0 . \tag{84}
\end{equation*}
$$

which can be solved for $q_{12}$.
Given that $\gamma_{2} \neq 0$, the existence and number of the solutions depends on the sign of $\gamma_{1}^{2} \gamma_{2}^{2}-$ $\left(\gamma_{1}^{2}+1\right)\left(\gamma_{2}^{2}-1\right)=\gamma_{1}^{2}-\gamma_{2}^{2}+1$.

1. If $\gamma_{1}^{2}-\gamma_{2}^{2}+1=0$, there is a unique solution.
2. If $\gamma_{1}^{2}-\gamma_{2}^{2}+1<0$, there is no real solution.
3. If $\gamma_{1}^{2}-\gamma_{2}^{2}+1>0$, there are two real solutions.

The condition in the current proposition can be stated equivalently to the condition for case 3 . Hence, if this condition holds, there a two distinct solutions $q_{1}^{(1)}$ and $q_{1}^{(2)}$, neither of which can be eliminated by the sign normalization restriction for $q_{1}$.

For each $q_{1}^{(1)}$ and $q_{2}^{(2)}$, the third restriction in Eq. (27) $\tilde{\sigma}_{3}^{\prime} q_{2}=0, q_{2}$ 's orthogonality to $q_{1}$, and the sign normalization restriction for $q_{2}$ pin down a unique $q_{2}$. We denote these by $q_{2}^{(1)}$ and $q_{2}^{(2)}$, respectively. Finally, orthogonality to $\left(q_{1}^{(1)}, q_{2}^{(1)}\right)$ or $\left(q_{1}^{(2)}, q_{2}^{(2)}\right)$ in $\mathbb{R}^{3}$ and the sign normalization for $q_{3}$ pin down the vectors $q_{3}^{(1)}$ and $q_{3}^{(2)}$. Thus, we obtain two distinct admissible orthogonal matrices $Q^{(1)}=\left(q_{1}^{(1)}, q_{2}^{(1)}, q_{3}^{(1)}\right)$ and $Q^{(2)}=\left(q_{1}^{(2)}, q_{2}^{(2)}, q_{3}^{(2)}\right)$. This SVAR is locally but not globally identified when the variance-covariance matrix of the reduced-form errors satisfies the stated conditions.

## Proof of Proposition 6: Blanchard-Perotti fiscal policy SVAR

The AB-SVAR representation of the Blanchard-Perotti model is

$$
\left(\begin{array}{ccc}
1 & 0 & -a_{1}  \tag{85}\\
0 & 1 & -b_{1} \\
-c_{1} & -c_{2} & 1
\end{array}\right)\left(\begin{array}{c}
t_{t} \\
g_{t} \\
x_{t}
\end{array}\right)=\left(\begin{array}{ccc}
1 & a_{2} & 0 \\
b_{2} & 1 & 0 \\
0 & 0 & 1
\end{array}\right)\left(\begin{array}{c}
\varepsilon_{t}^{t} \\
\varepsilon_{t}^{g} \\
\varepsilon_{t}^{x}
\end{array}\right) .
$$

The authors impose three further restrictions to solve the identification issue: 1) $b_{1}=0, a_{1}=2.08$ and, 3) either $a_{2}=0$ or $b_{2}=0$. Under these restrictions with $b_{2}=0$ (the analysis is similar for the alternative $a_{2}=0$ ), the model becomes

$$
\begin{aligned}
t_{t} & =2.08 x_{t}+a_{2} \epsilon_{t}^{g}+\epsilon_{t}^{t} \\
g_{t} & =\epsilon_{t}^{g} \\
x_{t} & =c_{1} t_{t}+c_{2} g_{t}+\epsilon_{t}^{x}
\end{aligned}
$$

whose SVAR representation is

$$
\begin{aligned}
& \left(\begin{array}{ccc}
1 & a_{2} & 0 \\
0 & 1 & 0 \\
0 & 0 & 1
\end{array}\right)^{-1}\left(\begin{array}{ccc}
1 & 0 & -2.08 \\
0 & 1 & 0 \\
-c_{1} & -c_{2} & 1
\end{array}\right)\left(\begin{array}{c}
t_{t} \\
g_{t} \\
x_{t}
\end{array}\right)=\left(\begin{array}{c}
\epsilon_{t}^{t} \\
\epsilon_{t}^{g} \\
\epsilon_{t}^{x}
\end{array}\right) \\
& \Leftrightarrow\left(\begin{array}{ccc}
1 & -a_{2} & -2.08 \\
0 & 1 & 0 \\
-c_{1} & -c_{2} & 1
\end{array}\right)\left(\begin{array}{c}
t_{t} \\
g_{t} \\
x_{t}
\end{array}\right)=\left(\begin{array}{c}
\epsilon_{t}^{t} \\
\epsilon_{t}^{g} \\
\epsilon_{t}^{x}
\end{array}\right)
\end{aligned}
$$

As is standard for simultaneous equation models, the normalization here has been obtained by imposing unit coefficients on the diagonal of the left-had side matrix, leaving the variances of the structural shocks unconstrained.

Let $\sigma_{t}^{2}, \sigma_{g}^{2}$ and $\sigma_{y}^{2}$, be the variances of $\epsilon_{t}^{t}, \epsilon_{t}^{g}$, and $\epsilon_{t}^{x}$, respectively. An alternative normalization to have the unit-variance structural shocks and ordering the variables to conform to Proposition 1 lead to

$$
\left(\begin{array}{ccc}
\alpha_{11} & 0 & 0 \\
\alpha_{21} & \alpha_{22} & \alpha_{23} \\
\alpha_{31} & \alpha_{32} & \alpha_{33}
\end{array}\right)\left(\begin{array}{c}
g_{t} \\
t_{t} \\
x_{t}
\end{array}\right)=\left(\begin{array}{c}
\varepsilon_{t}^{g} / \sigma_{g} \\
\varepsilon_{t}^{t} / \sigma_{t} \\
\varepsilon_{t}^{x} / \sigma_{x}
\end{array}\right)
$$

where $\alpha_{i j}$ parameters satisfy

$$
\begin{array}{lll}
\alpha_{11}=1 / \sigma_{g} & & \\
\alpha_{21}=-a_{2} / \sigma_{t} & , \quad \alpha_{22}=1 / \sigma_{t} & , \quad \alpha_{23}=-2.08 / \sigma_{t} \\
\alpha_{31}=-c_{2} / \sigma_{x} & , \quad \alpha_{32}=-c_{1} / \sigma_{x} & , \quad \alpha_{33}=1 / \sigma_{x}
\end{array}
$$

The two zero restrictions in the first equation provids two linear restrictions for $q_{1}$, which, combined with the sign normalization, pin down a unique $q_{1}$. The restriction on the second equation can be expressed as $\alpha_{22}=-2.08 \alpha_{23}$, which can be viewed as a zero restriction for $q_{2}$,

$$
\begin{aligned}
\alpha_{22}=-2.08 \alpha_{23} & \Leftrightarrow\left(\Sigma_{t r}^{-1} e_{2}\right)^{\prime} q_{2}=-2.08\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime} q_{2} \\
& \Leftrightarrow\left[\left(\Sigma_{t r}^{-1} e_{2}\right)^{\prime}+2.08\left(\Sigma_{t r}^{-1} e_{3}\right)^{\prime}\right] q_{2}=0 .
\end{aligned}
$$

As this boils down to a standard zero restriction, orthogonality to $q_{1}$ and the sign normalization uniquely pin down the second vector $q_{2}$. The third equation is unconstrained and orthogonality to ( $q_{1}, q_{2}$ ) combined with the sign normalization pins down $q_{3}$. This proves that there exists a unique orthogonal matrix $Q=\left(q_{1}, q_{2}, q_{3}\right)$ mapping the unrestricted reduced-form parameters to the structural ones satisfying the imposed restrictions. Hence, the Blanchard-Perotti AB-SVAR model is globally identified.

## D Appendix: The $\tilde{D}_{n}$ matrix

A skew-symmetric (square) matrix $A$ satisfies $A^{\prime}=-A$. Let $\tilde{v}(A)$ be a vector containing the $n(n-1) / 2$ essential elements of $A$. When $A$ is skew-symmetric, it is possible to expand the elements of $\tilde{v}(A)$ to obtain vec $A$. $\tilde{D}_{n}$, thus, can be defined to be the $n^{2} \times n(n-1 / 2)$ matrix with the property that

$$
\tilde{D}_{n} \tilde{v}(A)=\operatorname{vec} A
$$

for any skew symmetric $n \times n$ matrix $A$. For a formal definition and properties of $\tilde{D}_{n}$, see Magnus (1988). Here, we present $\tilde{D}_{n}$ for $n=2, n=3$ and $n=4$ :
where we have circled the elements selecting the last $n-i$ columns, $i=1, \ldots, n$, of the $F_{i i}(\phi) Q$ matrix in the proof of Proposition 7 .

Finally, as can be seen from $\tilde{D}_{2}, \tilde{D}_{3}$ and $\tilde{D}_{4}$, the matrix $\tilde{D}_{n}$ is always of full column rank $n(n-1) / 2$.


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[^2]:    ${ }^{1}$ Other proposals of identification strategies include the use of external instruments as in Mertens and Ravn (2013) and Stock and Watson (2018), heteroskedasticity of the structural shocks as in Rigobon (2003), Bacchiocchi and Fanelli (2015) and Bacchiocchi (2017), and the presence of non-normality as in Lanne and Lütkepohl (2010) and Lanne et al. (2017).

[^3]:    ${ }^{2}$ Theorem 7 of Rubio-Ramirez et al. (2010) claims that the exact identification of an SVAR holds if and only if $f_{i}=n-i$ for all $i=1, \ldots, n$. Bacchiocchi and Kitagawa (2020) shows by a counterexample that their condition is not sufficient and needs to be augmented by the rank conditions of Eq. (18) and Eq. (19). See Bacchiocchi and Kitagawa (2020) for further detail.

[^4]:    ${ }^{3}$ This minimization problem is constrained by the orthogonality constraints $Q^{\prime} Q=I_{n}$, which is the known as Stiefel manifold following Stiefel (1935-1936). Edelman et al. (1998) develop algorithms for optimization in the Stiefel manifold, while Boumal et al. (2014) propose a Matlab toolbox for optimization on manifolds including the Stiefel one. A Matlab code for this optimization is available from the authors upon request.

[^5]:    ${ }^{4}$ For $\phi \notin \Phi_{F}$, the vertical line does not intersect the unit circle, and no real solution for $q_{1}$ exists. If $c \neq 0$, the identifying restriction becomes observationally restrictive, and the identifying restriction can be refuted by the reduced-form models.

[^6]:    ${ }^{5}$ Identification criteria for AB-SVAR without normalizing the structural variances is known for local identification only. See Amisano and Giannini (1997), Bacchiocchi and Lucchetti (2018), Lütkepohl (2006), and Hamilton (1994).

[^7]:    ${ }^{6}$ In particular, Kelly (1975) presents cases in which economic theory might suggest imposing such restrictions. However, constraining parameters across equations is conditional on the kind of normalization considered. In simultaneous equation systems, normalization rules were generally based on imposing a unit coefficient for the variable playing the role of endogenous variable in that specific equation. In the parametrization proposed by RWZ for SVAR models the normalization rule instead consists of imposing unit variance on the uncorrelated structural shock. In this case, imposing restrictions on elasticities across equations would involve non-linear restrictions on the estimated coefficients. In fact, to obtain the elasticities, we need to normalize the coefficient for the endogenous variable in the each equation. See Hamilton et al. (2007) and Waggoner and Zha (2003) for specific details on the normalization issue in SVAR models.

[^8]:    ${ }^{7}$ As pointed out by Canova (2005), zero restrictions implied by DSGE models do not match the recursive identification schemes common in SVAR analyses.

[^9]:    ${ }^{8}$ The proxy-variable identification strategy has been shown to be useful for non-invertible structural MA models. See Stock and Watson (2018).

[^10]:    ${ }^{9}$ Kociecki and Kolasa (2018) similarly check global identification of DSGE models by examining the solutions of a non-linear system of equations.
    ${ }^{10}$ Sturmfels (2002) provides a good overview of systems of polynomial equations with potential applications in statistics and economics. As we saw in Section II.3, this system can be also seen as a minimization problem of the quadratic objective function subject to the orthogonality constraints $Q^{\prime} Q=I_{n}$. Noting that the orthogonality constraints generate the Stiefel manifold, we can consider applying algorithms for optimization on the Stiefel manifold. See Edelman et al. (1998) and Boumal et al. (2014).
    ${ }^{11}$ https://uk.mathworks.com/help/symbolic/vpasolve.html
    ${ }^{12}$ Matlab solvers are not open source, and we fail to uncover the precise numerical algorithm vpasolve uses to find roots of nonlinear equation systems.

[^11]:    ${ }^{13}$ The data are used in Aruoba and Schorfheide (2011) and downloaded from Frank Schorfheide's website: https://web.sas.upenn.edu/schorf/. For details on the construction of the series, see Appendix D from Granziera et al. (2018) and Footnote 5 of Aruoba and Schorfheide (2011).

[^12]:    ${ }^{14}$ Although not reported for saving space, the Bayesian credible intervals of Figure 4 are nearly identical to those obtained by the frequentist bootstrap-after-bootstrap approach of Kilian (1998)
    ${ }^{15}$ The highest posterior density regions are computed by slicing the posterior density approximated through kernel smoothing of the posterior draws of the impulse responses.

