EQUILIBRIA AND THEIR STABILITY IN NETWORKS WITH STEEP SIGMOIDAL NONLINEARITIES

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Abstract. In this paper we investigate equilibria of continuous differential equation models of network dynamics. The motivation comes from gene regulatory networks where each directed edge represents either down- or up-regulation, and is modeled by a sigmoidal nonlinear function. We show that the existence and stability of equilibria of a sigmoidal system is determined by a combinatorial analysis of the limiting switching system with piece-wise constant non-linearities. In addition, we describe a local decomposition of a switching system into a product of simpler cyclic feedback systems, where the cycles in each decomposition correspond to a particular subset of network loops.

1. Introduction. Analysis of large systems of ordinary differential equations is difficult, especially when we seek to understand changes in dynamics when parameters change. To set the stage, we are interested in systems of $O(10)$ differential equations with the same order of magnitude of number of parameters; big enough to be complicated but not so large that statistical approaches may be applicable. Systems of this size are important in systems biology, in particular, in models of gene regulation. Here variables usually represent concentrations of chemical species (mRNA, proteins) in the cell, and the interactions between variables are represented in a form of a network with signed directed edges. Nodes represent concentrations and directed edges monotone interactions; positive edges indicate activation and negative edges repression. The interactions are nonlinear; both on the level of the pairwise effect of $x_i$ on $x_j$ which is usually modeled by a saturating function, but also on the level of how effects from different inputs combine together to influence $x_j$. The choice of the nonlinearity that models the effect of $x_i$ on $x_j$ is not given by any fundamental law of physics; the usual choices are Hill functions, but other sigmoid functions and threshold (switching) functions are used as well. This ambiguity, coupled with technical challenges related to simultaneous measurement of time evolution of multiple chemical species in a single cell, necessarily limits the expectation of fidelity of the model simulations with the experimental data. The model should not be expected to reproduce measurements in fine detail, but still answer qualitative questions on number and types stable equilibria are present, or capacity of the network to admit oscillations. Taking into consideration ever present noise in molecular systems, there is always a need to address a question of how robust these qualitative features are under parameter changes in the model.

In this paper we concentrate on the existence and stability of equilibria in networks where the pairwise interaction is modeled by sigmoidal nonlinearities. Extending results [22, 28, 19], we show that the equilibria in a network modeled by sufficiently steep sigmoidal functions are in one-to-one correspondence with a collection of so called regular and singular equilibria of a model of the same network using switching functions. Switching functions are piece-wise constant functions with a single threshold and range with two values $\{L, U\}$, which can be interpreted as two rates of expression (L for Lower, and U for Upper) of the target gene based on whether the controlling gene is below, or above the threshold $\theta$. These models have been used for gene regulatory networks since the 70’s [15, 16, 6, 25, 7, 5, 19, 11]. However, using these functions as the right hand side of an ODE system presents several technical challenges, especially how to deal with the fact that the vector field is not defined at thresholds $\theta$. One approach to extend switching systems so that they
are defined at thresholds is to view them as differential inclusions rather than ordinary differential equations. Stability of singular equilibria of the switching systems from this point of view were studied by [2]. Convergence of switching systems dynamics was studied also by [21] who formulated conditions that are needed to construct a global Lyapunov function for a switching system by piecing together local Lyapunov functions. The idea of the new DSGRN (Dynamic Signatures Generated by Regulatory Networks) approach [5, 12, 11], supported by a suite of corresponding software [4], is to capture information about the network dynamics given by switching system models in a form of combinatorial (finite) data, and then use this data to rigorously establish results about well-defined dynamics of ODE’s with continuous right hand side that are a small perturbation of the switching functions. We emphasize that the distinction in the two approaches is that the Filippov extension approach is concerned with the dynamics of switching systems whereas the DSGRN approach only uses switching systems as a computational tool for the study of the dynamics of sigmoidal systems.

While we describe the combinatorial data in greater detail below, for the purpose of this introduction it is sufficient to note that the switching system ODE only contains stable equilibria (which we call regular), because the unstable equilibria of sigmoidal systems limit to intersections of thresholds of switching systems. In this paper we show that the intersections, the intersections that appear as such limits of unstable equilibria, which we call loop characteristic equilibrium cells, can be precisely characterized using combinatorics of the switching system. In particular, we add to the DSGRN approach by showing how to use the combinatorial data from a switching system to predict existence and stability of all equilibria for all sigmoidal functions that arise as perturbations to the switching functions. The combinatorial data only uses the type (i.e. positive feedback or negative feedback) and the number of feedback loops in the network. To obtain these results we show that switching systems can be locally (in phase space) decomposed into a product of simpler cyclic feedback systems. Each such cyclic system can be associated to a unique oriented loop in the gene regulatory network.

A potential application of description of equilibria and their stability in sigmoidal systems is in the recurrent artificial neural networks (rANN). These models were introduced by Hopfield [18] and Grossberg [17] almost 40 years ago, but they found their newest incarnation as Echo state networks [20]. There is a great variety of implementations but at the core there are network nodes (i.e. neurons) that are connected by weighted directed edges. Each node processes the input through a nonlinear function (binary, sigmoidal, or a ramp). Our work provides a characterization of the number of stable equilibria for steep sigmoidal functions based on the combinatorics of the switching system, which is ultimately tied to the structure of the connections in the network.

1.1. Organization of the paper. In Section 2, we define sigmoidal functions and switching functions. In Section 3, we define the combinatorial data associated to switching functions and show how this data can be used to identify all equilibria of steep sigmoidal systems. The proofs for these results are in Section 6. In Section 4, we use the results of Section 3 to characterize the equilibria of cyclic feedback networks and then analyze their stability. The proofs can be found in Section 7. In Section 5, we show that all switching systems can be locally (in phase space) decomposed as a product of cyclic feedback systems and use this decomposition to generalize the stability results of Section 4 from cyclic feedback systems to general networks. Finally, in Section 8, we conclude with a discussion of our results.

2. The Regulatory Network and Switching Systems.
\textbf{Definition 2.1} ([5]). A regulatory network \( \text{RN} = (V,E) \) is an annotated finite directed graph with vertices \( V = \{1, \ldots, N\} \) called network nodes and directed edges \( E \subseteq V \times V \times \{1, -1\} \). An annotated edge \((j, i, +1)\) represents an activation of node \( i \) by node \( j \) and is denoted \( j \to i \); annotated edge \((j, i, -1)\) represents repression of node \( i \) by node \( j \) and is denoted \( j \leftarrow i \). We write \( s_{ij} = 1 \) if \( j \to i \) and \( s_{ij} = -1 \) if \( j \leftarrow i \). We indicate either \( j \to i \) or \( j \leftarrow i \) without specifying which by writing \((j, i) \in E \). We allow self edges, but admit at most one edge between any two nodes. The set of sources and targets of a node are denoted by

\[
\mathbf{S}(k) = \{ j \mid (j, k) \in E \} \quad \text{and} \quad \mathbf{T}(k) = \{ j \mid (k, j) \in E \}
\]

and we require every node has a target.

We remark that the assumption that every node has a target is not a serious constraint. If a node \( j \) does not have a target, then the dynamics of the remaining nodes are independent of \( j \). Once the dynamics of the remaining nodes are understood, the dynamics of node \( j \) can be treated as a non-autonomous system driven by the remaining nodes.

To an \( \text{RN} \) we associate a switching system of the form

\[
\dot{x} = -\Gamma x + \Lambda(x)
\]

where \( \Gamma \) is a diagonal matrix with entries \( \Gamma_{jj} = \gamma_j \) and \( \Lambda \) is a nonlinear function of the form

\[
\Lambda_i(x) := \prod_{j \in I_{ii}} \sigma_{ij}(x_j)
\]

with \( I_1, \ldots, I_p \) a partition of \( S(i) \). Each \( \sigma_{ij} \) is a switching function of the form

\[
\sigma_{ij}(x_j) := \begin{cases} L_{ij}, & s_{ij} = 1 \text{ and } x_j < \theta_{ij} \text{ or } s_{ij} = -1 \text{ and } x_j > \theta_{ij} \\ U_{ij}, & s_{ij} = 1 \text{ and } x_j > \theta_{ij} \text{ or } s_{ij} = -1 \text{ and } x_j < \theta_{ij} \\ \text{undefined}, & \text{if } x_j = \theta_{ij}. \end{cases}
\]

The parameter \( Z = (L, U, \theta, \Gamma) \), where \( L := (L_{ij}), U := (U_{ij}), \theta := (\theta_{ij}) \) are vectors indexed by \((ij)\), is the switching parameter. We denote a switching system parameterized by \( Z \) by \( \text{SWITCH}(Z) \).

To an \( \text{RN} \) we also associate a sigmoidal system, \( \mathcal{S}(Z, \varepsilon) \), where \( Z \) is a switching parameter and \( \varepsilon \in \mathbb{R}^{N \times N} \) is a perturbation parameter. We say \( \varepsilon' \leq \varepsilon \) or \( \varepsilon' < \varepsilon \) when the component-wise comparisons \( \varepsilon'_{ij} \leq \varepsilon_{ij} \) or \( \varepsilon'_{ij} < \varepsilon_{ij} \) hold for each \((j, i) \in E\), respectively. The pair \((Z, \varepsilon)\) is the sigmoidal parameter. The dynamics of \( \mathcal{S}(Z, \varepsilon) \) are given by

\[
\dot{x} = -\Gamma x + \Lambda(x; \varepsilon)
\]

where \( \Lambda(x; \varepsilon) \) is obtained from \( \Lambda \) by replacing the switching functions \( \sigma_{ij} \) with sigmoidal perturbations \( \sigma_{ij}(\cdot; \varepsilon_{ij}) \), which we define below.

\textbf{Definition 2.2}. \( \sigma_{ij}(\cdot; \varepsilon_{ij}) \) is a family of sigmoidal perturbations of \( \sigma_{ij} \) at a parameter \( Z \) if for each \( \varepsilon_{ij} \in \mathbb{R}_+ \),

1. \( \sigma_{ij}(\cdot; \varepsilon_{ij}) \) is continuously differentiable and monotone non-increasing or monotone non-decreasing.
2. \( \sup_x \sigma_{ij}(x; \varepsilon_{ij}) = U_{ij} \) and \( \inf_x \sigma_{ij}(x; \varepsilon_{ij}) = L_{ij} \).

3. There is a neighborhood \( \mathcal{U}_1(\varepsilon_{ij}) \subset \mathbb{R} \) of \( \theta_{ij} \) such that \( \mathcal{U}_1(\varepsilon_{ij}) \to \{\theta_{ij}\} \) as \( \varepsilon_{ij} \to 0 \) and a constant \( C_1 > 0 \), such that if \( x \in \mathbb{R} \setminus \mathcal{U}_1(\varepsilon_{ij}) \), then \( |\sigma'_{ij}(x; \varepsilon_{ij})| \leq C_1 \varepsilon_{ij} \).

4. There is a neighborhood \( \mathcal{U}_2(\varepsilon_{ij}) \subset \mathbb{R} \) of \( \theta_{ij} \) such that \( \mathcal{U}_2(\varepsilon_{ij}) \to \{\theta_{ij}\} \) and \( \sigma_{ij}(\mathcal{U}_2(\varepsilon_{ij}); \varepsilon_{ij}) \to (L_{ij}, U_{ij}) \) as \( \varepsilon_{ij} \to 0 \) and a constant \( C_2 > 0 \) such that if \( x \in U_2(\varepsilon_{ij}) \) then \( |\sigma'_{ij}(x; \varepsilon_{ij})| \geq C_2 \varepsilon_{ij}^{-1} \).

Given a perturbation parameter \( \varepsilon \), we will write \( \sigma_{ij}(\cdot; \varepsilon) \) instead of \( \sigma_{ij}(\cdot; \varepsilon_{ij}) \) to simplify notation. Note that as \( \varepsilon \to 0 \), the sigmoidal perturbation \( \sigma_{ij}(\cdot; \varepsilon) \) converges pointwise to the step function \( \sigma_{ij} \). Given a switching parameter \( Z \) and perturbation parameter \( \varepsilon \geq 0 \), we denote the Jacobian of (2.4) at \( x \) by \( J(x; \varepsilon) \) or \( J(\varepsilon) \) when \( x \) is implied from context.

**Example 2.3.** Throughout the paper we will illustrate the concepts on a simple example of a two node network we call the positive toggle plus, where two nodes activate themselves and mutually activate each other, i.e.

\[ \text{RN} = (V, E) = \{(1, 2), (1 \to 1), (2 \to 2), (1 \to 2), (2 \to 1)\}. \]

The name "positive toggle" refers to the network without self loops and was chosen for its resemblance to the toggle switch introduced in [10], in which the nodes mutually repress each other rather than activate. The "plus" modifier refers to the addition of the self loops. The associated switching system has the form

\[
\begin{align*}
\dot{x}_1 &= -\gamma_1 x_1 + \sigma_{11}(x_1)\sigma_{12}(x_2) \\
\dot{x}_2 &= -\gamma_2 x_2 + \sigma_{22}(x_2)\sigma_{21}(x_1)
\end{align*}
\]

and we note \( s_{11} = s_{22} = s_{12} = s_{21} = 1 \). We will consider this system with a switching parameter satisfying

\[
\begin{align*}
L_{11}L_{12} &< L_{11}U_{12} < \gamma_1 \theta_{21} < U_{11}L_{12} < \gamma_1 \theta_{11} < U_{11}U_{12}, \text{ and} \\
L_{22}L_{21} &< \gamma_2 \theta_{12} < U_{22}U_{21} < L_{22}L_{21} < U_{22}U_{21} < \gamma_2 \theta_{22}.
\end{align*}
\]

**3. Equilibria of Regulatory Networks.** As observed in [5], the thresholds \( \theta_{ij} \) of a switching system impose a grid-like structure on phase space \( \mathbb{R}^N_+ \). In this section we characterize where the equilibria lie in phase space relative to this structure. We begin by defining the structure.

**Definition 3.1.**

1. For each \( j \in V \), we define \( \theta_{-\infty j} := 0, \theta_{\infty j} := \infty \), and

\[
\Theta_j(Z) := \{\theta_{ij} > 0 \mid i \in T(j)\} \cup \{\theta_{\infty j}, \theta_{-\infty j}\}.
\]

The threshold set is the collection \( \Theta(Z) := (\Theta_1(Z), \ldots, \Theta_N(Z)) \). We say \( \theta_{i_{1j}}, \theta_{i_{2j}} \in \Theta_j(Z) \) are consecutive thresholds if \( \theta_{i_{1j}} < \theta_{i_{2j}} \) and there does not exist \( \theta_{i_{3j}} \in \Theta_j(Z) \) such that \( \theta_{i_{1j}} < \theta_{i_{3j}} < \theta_{i_{2j}} \).

2. A cell \( \tau \) associated to a threshold set \( \Theta \), is a product of \( k \leq N \) thresholds and \( N-k \) open intervals whose endpoints are consecutive thresholds. By renumbering the variables we write

\[
\tau = \prod_{j=1}^{k} (\theta_{i_{1j}}, \theta_{i_{2j}}) \times \prod_{j=k+1}^{N} (\theta_{a_{ij}}, \theta_{b_{ij}}).
\]
We write $\pi_j(\tau)$ for the projection of $\tau$ onto the $j$th direction. A cell is regular if $k = 0$ and singular otherwise. The cell complex, $\chi(\Theta)$, is the collection of all cells associated to the threshold set $\Theta$. The cell complex associated to the switching system $\text{SWITCH}(Z)$ at parameter $Z$ is $\chi(\Theta(Z))$. When the switching parameter $Z$ and the threshold set $\Theta(Z)$ are clear from context we drop the argument $\Theta(Z)$ and write $\chi$.

Figure 1(a) depicts the cell complex $\chi$ for the positive toggle plus. The concept of neighboring cells is described below.

Fig. 1: The cell complex $\chi$, neighbors, and the labeling map in the positive toggle plus network. The network and switching system are defined in Example 2.3. (a): Each box, line, and point is a cell in the cell complex $\chi$. The cell $\kappa = \{\theta_{21}\} \times (\theta_{12}, \theta_{22})$ is indicated by the dashed line. For $\kappa$, the direction 1 is singular and the direction 2 is regular. The 2-neighbors of $\kappa$ (see Definition 3.5), $\kappa_2^+$ and $\kappa_2^-$, are indicated by the gray circles while the 1-neighbors, $\kappa_1^+$ and $\kappa_1^-$, are the labeled two dimensional cells. (b): The loop characteristic cells $\tau$ and $\tilde{\tau}$ are indicated by the gray circles and the regular cells $\tau'$ and $\tau''$ are labeled. The dotted arrow represents $L(\tau, 1, -)$ for which 1 is a singular direction and, at the same time, it represents $L(\tau', 2, +)$ and $L(\tau'', 2, -)$ for which 2 is a regular direction. See Example 3.10 for details. These values are equal to $-1$ so the arrow points down. The arrows on the outside cell complex boundary are not drawn because they point inwards for all choices of parameters. The arrows of the labeling map imply that $\Phi_1(\tau) = \Phi_2(\tau') = 0$. The arrows on the top and bottom boundaries of $\tau''$ indicate that $\Phi_2(\tau'') = -1$.

Observe, that in the switching system, the function $\Lambda$ is only defined on regular cells and not on singular cells. Therefore, equilibria of $\text{SWITCH}(Z)$ can only be contained in regular cells. Equilibria in singular cells can be understood if (2.1) is replaced with its Filippov extension wherein the differential equation is replaced by a differential inclusion. This was done in [2] where existence and stability of these singular equilibria was studied. However, our goal is to understand equilibria of the sigmoidal systems $S(Z, \varepsilon)$ which are perturbations of the switching system $\text{SWITCH}(Z)$, and not the equilibria of $\text{SWITCH}(Z)$. For this reason, in the next definition we define an equilibrium
cell to be a cell which a sigmoidal equilibrium converges to as \( \varepsilon \to 0 \). It is straightforward to see that regular equilibrium cells contain a unique stable equilibrium of \( \text{SWITCH}(Z) \).

**Definition 3.2.** Let \( Z \) be a switching parameter and \( \tau \in \chi \). If there is an \( A \in \mathbb{R}^{N \times N}_+ \) so that for all \( \varepsilon < A \), \( S(Z, \varepsilon) \) has a fixed point \( x^\varepsilon \) satisfying \( x^\varepsilon \to \tau \) as \( \varepsilon \to 0 \), then \( \tau \) is an equilibrium cell. If \( \tau \) is a singular cell, then \( x^\varepsilon \) is a singular stationary point (SSP).

Theorem 3.13 characterizes equilibrium cells using combinatorial information about \( \text{SWITCH}(Z) \) only. That is, information about the switching system is necessary and sufficient to characterize the equilibria of the sigmoidal system \( S(Z, \varepsilon) \) when \( \varepsilon \) is small enough. To state the theorem precisely, we need some additional definitions to describe the cell complex and a notion of a flow direction map on the cell complex.

### 3.1. Cell Complex.

**Definition 3.3.**
1. Given \( \tau \in \chi \), the singular directions of \( \tau \), denoted \( \text{sd}(\tau) \), correspond to the set of indices, \( s \), such that \( \pi_s(\tau) = \{ \theta_{i,s} \} \). An index is a regular direction if it is not singular. We define \( \rho^\tau : V \to V \) by

\[
\rho^\tau(j) := \begin{cases} 
  i_j, & j \in \text{sd}(\tau) \\
  j, & \text{otherwise}.
\end{cases}
\]

The set of cells with \( k \) singular directions is denoted \( \chi^{(N-k)} \). The map \( \rho^\tau \) depends on the cell \( \tau \), but when \( \tau \) is fixed and clear from the context, we will use \( \rho \) instead of \( \rho^\tau \).

2. If \( s \) is a singular direction of \( \tau \in \chi \) we denote the neighboring thresholds by \( \theta_{\rho_-(s)} \) and \( \theta_{\rho_+(s)} \) where \( \theta_{\rho_-(s)} < \theta_{\rho(s)} < \theta_{\rho_+(s)} \) are consecutive thresholds in \( \Theta_s(Z) \). If \( r \) is a regular direction of \( \tau \) we write \( \pi_r(\tau) = (\theta_{a,r}, \theta_{b,r}) \).

**Definition 3.4.** A cell, \( \tau \in \chi \), is a loop characteristic cell if \( \rho^\tau \) is a permutation on \( \text{sd}(\tau) \). We denote the set of loop characteristic cells by \( \text{LCC} \). Note that all \( N \)-dimensional cells \( \kappa \) are automatically loop characteristic cells, since \( \text{sd}(\kappa) = \emptyset \). Therefore \( \chi^{(N)} \subset \text{LCC} \).

A loop characteristic cell is a cell in which some number of disjoint loops of the network are active. For example, the loop characteristic cells of the positive toggle plus network (see Example 2.3) include \( \{ \theta_{21} \} \times \{ \theta_{12} \} \), where the loop 1 \( \to \) 2 \( \to \) 1 is active, \( \{ \theta_{11} \} \times (\theta_{12}, \theta_{22}) \), where the loop 1 \( \to \) 1 is active, and \( \theta_{11} \times \theta_{22} \), where the loops 1 \( \to \) 1 and 2 \( \to \) 2 are both active. The concept of loop characteristics was introduced in [22] where it was shown singular equilibria of switching systems are contained in loop characteristic cells. This was later extended to sigmoidal systems by [28] who showed that equilibrium cells are a subset of loop characteristic cells when the sigmoidal perturbations are taken to be Hill functions. Theorem 3.13 extends this work by considering a more general class of sigmoids and providing necessary and sufficient conditions for a loop characteristic cell to be an equilibrium cell.

In the following definition we introduce notation to describe the neighbor of a cell. By a neighbor to \( \tau \) we mean a cell which is directly adjacent to \( \tau \).

**Definition 3.5.** Let \( \tau \in \chi \) and \( j \in V \). The left \( j \)-neighbor (see Figure 1) of the cell \( \tau \) is a cell \( \tau^-_j \),
defined by

\[ \pi_k(\tau_j^-) := \begin{cases} \pi_j(\tau), & j \neq k \\ \{\theta_{p^-_{(j)j}}, \theta_{p^+_{(j)j}}\}, & j = k, k \in sd(\tau) \\ \{\theta_{\alpha_j}\}, & j = k, k \notin sd(\tau). \end{cases} \]

Similarly, the right \( j \)-neighbor, \( \tau_j^+ \), is defined by

\[ \pi_k(\tau_j^+) := \begin{cases} \pi_k(\tau), & j \neq k \\ \{\theta_{p^+_{(j)j}}, \theta_{p^-_{(j)j}}\}, & j = k, k \in sd(\tau) \\ \{\theta_{\alpha_j}\}, & j = k, k \notin sd(\tau). \end{cases} \]

An \( j \)-neighbor of \( \tau \) is either a left or right \( j \)-neighbor of \( \tau \). A neighbor of \( \tau \) is any \( j \)-neighbor.

On a diagram of the cell complex, the left \( j \)-neighbor of \( \tau \), \( \tau_j^- \), is the cell directly below \( \tau \) in the \( j \)th direction and the right \( j \)-neighbor of \( \tau \), \( \tau_j^+ \), is the cell directly above \( \tau \) in the \( j \)th direction. If \( j \) is a singular direction of \( \tau \) then \( j \) is a regular direction of \( \tau_j^\pm \). If \( j \) is a regular direction of \( \tau \) then \( j \) is a singular direction of \( \tau_j^\pm \). See Figure 1(a) for an example.

In the remainder of the paper we assume that the thresholds \( \theta_{ij} \) are positive. This reflects that our motivation comes from biological networks in which activities or concentrations of a reactant are always non-negative. We also assume that thresholds corresponding to the same node are not equal. This holds generically, greatly simplifies our analysis, and is typical in the literature on switching systems [22, 2, 8, 28, 19, 5, 21]. These assumptions are captured in the following definition.

**Definition 3.6.** The switching parameter \( Z \) is threshold regular if

- For all \( (j, i) \in E \), \( \theta_{ij} > 0 \), and
- for all \( j \in V \), \( i_1, i_2 \in T(j) \), \( \theta_{i_1j} \neq \theta_{i_2j} \).

**Definition 3.7 (Definition 4.6 of [5]).** Consider a threshold regular switching parameter \( Z \). For \( j \in V \), denote the ordering of the thresholds \( \{\theta_{ij} \mid i \in T(j)\} \) by \( O_j(Z) \). The order parameter of \( Z \) is the collection of these orders, \( O(Z) = (O_1(Z), \ldots, O_N(Z)) \).

### 3.2. Flow Direction Map

The goal of this section is to define a flow direction map on the cell complex \( \chi(\Theta(Z)) \), which is induced by the right hand side of the switching system (2.1). We start by introducing some notation. Observe that if \( j \) is a singular direction of \( \tau \in \chi \) and thus \( \pi_j(\tau) = \{\theta_{ij}\} \) for some \( i \), then by equation (2.3) the function \( \sigma_{ij}(x_j) \) is not defined on \( \tau \). However, if \( j \) is a regular direction of \( \tau \) and thus \( \pi_j(\tau) = (\theta_{i_1j}, \theta_{i_2j}) \), then \( \sigma_{ij}(x_j) \) is constant on \( \tau \). We denote its value by \( \sigma_{ij}(\tau) \). It follows that for \( \tau \in \chi(\chi) \) which has no singular directions \( sd(\tau) = \emptyset \), the value of \( \Lambda_i(\tau) \) is well defined and constant for very \( i \). The vector \( \Gamma^{-1}(\Lambda(\tau)) \) is sometimes called the focal point of \( \tau \) because any trajectory of the switching system with initial condition in \( \tau \) will converge to this value until the trajectory reaches the boundary of \( \tau \) [2]. In the following definition we give non-degeneracy conditions for the switching parameter \( Z \) which we will assume throughout.

**Definition 3.8 (Definition 2.7 of [5]).** The switching parameter \( Z \) is regular if

- \( Z \) is threshold regular,
• for all \((j, i) \in E, 0 < L_{ij} < U_{ij}\),
• for all \(k \in V, \gamma_k > 0\), and
• for all \(\kappa \in \chi^{(N)}\) and \((j, i) \in E, \gamma_j \theta_{ij} \neq \Lambda_j(\kappa)\) for each threshold \(\theta_{ij}\) which defines \(\kappa\).

The last condition expresses the requirement that the focal point of each regular cell \(\kappa\) does not lie in the boundary of \(\kappa\), which holds generically. This is a typical assumption for switching systems because it implies that the right hand side of the switching system (2.1) has a well defined crossing direction on all the boundaries of cells \(\kappa \in \chi^{(N)}\) [2, 8, 5, 21]. We will use this to define a labeling map that collects information about these crossing directions. We then use the labeling map to define the flow direction map which indicates the direction in which the flow of the corresponding system crosses the the threshold. The flow direction map can be viewed as a multi-valued map and represented as a state transition graph that is a combinatorial summary of the flow information given by the switching system. The labeling map, defined below, generalizes the concept of wall-labeling (Definition 3.1 of [5]) in switching systems from regular cells to loop characteristic cells.

**Definition 3.9.** Let \(Z\) be a regular switching parameter.

1. The labeling map \(LCC \times V \times \{-, +\} \rightarrow \{-1, 1\}\) describes the sign of the right hand side of the switching system on the cells that are neighbors of \(\tau \in LCC\) in a particular direction. Letting \(\rho = \rho^\tau\), we first consider regular directions \(j \notin sd(\tau)\). Here we look at the sign of the \(j\)-th equation of the switching system (2.1) on the boundary in the \(j\)-direction:

\[L(\tau, j, \beta) := \begin{cases} 
\text{sgn}(-\gamma_j \theta_{a^2, j} + \Lambda_j(\tau)), & j \notin sd(\tau), \beta = - \\
\text{sgn}(-\gamma_j \theta_{b^2, j} + \Lambda_j(\tau)), & j \notin sd(\tau), \beta = + 
\end{cases}\]

For singular direction \(j \in sd(\tau)\), we look at a \(j\)-neighbor of \(\tau\) and ask for the sign of the \(\rho(j)\)-th equation of the switching system because \(\Lambda_{\rho(j)}\) is guaranteed to be well defined on a \(j\)-neighbor (see Lemma 6.1):

\[L(\tau, j, \beta) := \begin{cases} 
\text{sgn}(-\gamma_{\rho(j)} \theta_{a^2, \rho(j)} + \Lambda_{\rho(j)}(\tau_j^-)), & j \in sd(\tau), \beta = - \\
\text{sgn}(-\gamma_{\rho(j)} \theta_{b^2, \rho(j)} + \Lambda_{\rho(j)}(\tau_j^+)), & j \in sd(\tau), \beta = + 
\end{cases}\]

2. The flow direction map, \(\Phi : LCC \rightarrow \{-1, 0, 1\}^N\) summarizes the degree of agreement in the labeling map between the neighbors of \(\tau\) in a given direction. It is defined component-wise by

\[\Phi_j(\tau) := \begin{cases} 
1, & L(\tau, j, -) = 1 = L(\tau, j, +) \\
-1, & L(\tau, j, -) = -1 = L(\tau, j, +) \\
0, & L(\tau, j, -) = -L(\tau, j, +). 
\end{cases}\]

**Example 3.10.** Consider the positive toggle plus system of Example 2.3. Let \(\tau' = (0, \theta_{21}) \times (0, \theta_{12})\) be the lower left regular cell of the cell complex as in Figure 1(b). Then \(\Lambda_2(\tau') = L_{22}L_{21}\) and \(\theta_{22} = \theta_{12}\) so that

\[L(\tau', 2, +) = \text{sgn}(-\gamma_2 \theta_{a^2, 2} + \Lambda_2(\tau')) = \text{sgn}(-\gamma_2 \theta_{12} + L_{22}L_{21}) = -1\]
which is represented by the dotted down arrow originating from the upper boundary of \( \tau' \) Figure 1(b). This indicates the flow of the switching system in the \( x_2 \) direction is downward when \( x \in \tau' \) is close to the upper boundary of \( \tau' \). We also have \( \theta_{a2} = 0 \) so that

\[
L(\tau', 2, -) = \text{sgn}(-\gamma_2 \theta_{a2} + \Lambda(\tau')) = \text{sgn}(0 + L_{22}L_{21}) = 1
\]

and \( \Phi_2(\tau') = 0 \) since \( L(\tau', 2, -) = -L(\tau', 2, +) \). This indicates that \( \dot{x}_2 = 0 \) for some \( x \in \tau' \). The cell \( \tau = \{ \theta_{21} \} \times \{ \theta_{12} \} \) is a loop characteristic cell with \( \rho(1) = 2 \) and \( \rho(2) = 1 \). We compute

\[
L(\tau, 1, -) = \text{sgn}(-\gamma_{\rho(1)} \theta_{\rho(1)} + \Lambda_{\rho(1)}(\tau^-)) = \text{sgn}(-\gamma_2 \theta_{12} + L_{22}L_{21}) = -1.
\]

which is the direction of the flow on the left neighbor of \( \tau \). This is also represented by the dotted down arrow in Figure 1(b).

In general, \( L(\tau, j, \pm) \) is represented on a diagram of the cell complex by an arrow originating from \( \tau \pm j \) pointing in direction \( \rho(j) \) either positively or negatively according to the sign of \( L(\tau, j, \pm) \). As suggested by Figure 1, the flow direction map gives rise to a state transition graph which represents admissible transitions between the states that are represented by \( \kappa \in \chi^{(N)} \). The state transition graph is explicitly constructed in [5].

Note that the flow direction map depends on the choice of parameter \( Z \). A key observation is that it only depends on inequalities between parameters. We define an equivalence relation on all parameters \( Z \) that satisfy the same inequalities and therefore produce the same flow direction map. These equivalence classes, which we now proceed to define, will be called combinatorial parameters.

**Definition 3.11 (Definitions 4.5 and 4.6 of [5]).** Consider a regular switching parameter \( Z \).

1. The input combinations of the \( i \)th node is the Cartesian product

\[
\mathbb{I}_i := \prod_{j \in S(i)} \{ \text{off}, \text{on} \}.
\]

The indicator function, \( \mathbb{I}_i : \mathbb{R}^{S(i)}_+ \to \mathbb{I}_i \), is defined component-wise by

\[
\mathbb{I}_{i,j}(x) := \begin{cases} 
\text{off}, & s_{ij} = 1 \text{ and } x_j < \theta_{ij} \text{ or } s_{ij} = -1 \text{ and } x_j > \theta_{ij} \\
\text{on}, & s_{ij} = 1 \text{ and } x_j > \theta_{ij} \text{ or } s_{ij} = -1 \text{ and } x_j < \theta_{ij} \\
\text{undefined}, & \text{otherwise}.
\end{cases}
\]

The \( \sigma \)-valuation function, \( v_{ij} : \mathbb{I}_i \to \mathbb{R} \) is defined by

\[
v_{ij}(A) := \begin{cases} 
L_{ij}, & A = \text{off} \\
U_{ij}, & A = \text{on} \\
\text{undefined}, & \text{otherwise}.
\end{cases}
\]

Note that \( \sigma_{ij} = v_{ij} \circ \mathbb{I}_{ij} \). The \( \Lambda \)-valuation function, \( \omega_i : \mathbb{I}_i \to \mathbb{R} \), is defined by

\[
\omega_i(A) := \prod_{\ell=1}^{p_i} \sum_{j \in I_i} v_{ij}(A_j).
\]
Note that $\Lambda_i = \omega_i \circ \mathbb{1}_i$.

Define $L_j : \mathbb{I}_j \times T(j) \to \{ -1, 1 \}$ by

$$L_j(A, i) := \text{sgn}(-\gamma_j \theta_{ij} + \omega_j(A)).$$

The logic parameter is the collection $L(Z) := (L_1(\cdot, \cdot), \ldots, L_N(\cdot, \cdot))$.

2. We define an equivalence relation $Z \sim Z'$ whenever $(L(Z'), O(Z')) = (L(Z), O(Z))$. The combinatorial parameter is an equivalence class of this relationship. In other words, $Z' \in \mathcal{P}(Z)$ whenever $(L(Z'), O(Z')) = (L(Z), O(Z))$.

The notion of combinatorial parameter $\mathcal{P}(Z)$ was introduced in [5]. Each combinatorial parameter is defined in terms of inequalities between real valued parameters of $Z$. Therefore each combinatorial parameter corresponds to an open domain in the real-valued parameter space of parameters $Z$. The key observation from [5] is that any two parameters $Z_1, Z_2 \in \mathcal{P}(Z)$ define identical labeling maps and therefore identical flow direction maps. This is because the logic parameters $L_j$ that enter the definition of combinatorial parameters represent the same signs of the switching differential equations that define the labeling map.

Example 3.12. Consider the positive toggle plus system of Example 2.3. We will reference $\tau'$ and $\tau''$ from Figure 1(b). The input combinations for the first node is

$$\mathbb{I}_1 = \{ \text{off}, \text{on} \} \times \{ \text{off}, \text{on} \}$$

because the first node has two inputs. The indicator function depends only on the order parameter $O(Z)$. The second component of the indicator function $\mathbb{1}_1$ satisfies

$$\mathbb{1}_{12}(x) = \begin{cases} \text{off}, & x \in \tau' \\ \text{on}, & x \in \tau'' \end{cases}.$$

The $\sigma$-valuation function $v_{12}$ satisfies $v_{12}(\text{off}) = L_{12}$ and $v_{12}(\text{on}) = U_{12}$ so that

$$\sigma_{12}(x) = v_{12}(\mathbb{1}_{12}(x)) = \begin{cases} L_{12}, & x \in \tau' \\ U_{12}, & x \in \tau'' \end{cases}.$$

The $\Lambda$-valuation function $\omega_1$ satisfies

$$\omega_1((\text{off}, \text{off})) = v_{11}(\text{off})v_{12}(\text{off}) = L_{11}L_{12}, \text{ and}$$

$$\omega_1((\text{off}, \text{on})) = v_{11}(\text{off})v_{12}(\text{on}) = L_{11}U_{12}$$

so that

$$\Lambda_1(x) = \omega_1(\mathbb{1}_1(x)) = \begin{cases} \omega_1((\text{off}, \text{off})) = L_{11}L_{12}, & x \in \tau' \\ \omega_1((\text{off}, \text{on})) = L_{11}U_{12}, & x \in \tau'' \end{cases}.$$

The first component of the logic parameter $L(Z)$ satisfies

$$L_1((\text{off}, \text{off}), 2) = \text{sgn}(-\gamma_1 \theta_{21} + \omega_1((\text{off}, \text{off}))) = \text{sgn}(-\gamma_1 \theta_{21} + L_{11}L_{12}) = -1, \text{ and}$$

$$L_1((\text{off}, \text{on}), 2) = \text{sgn}(-\gamma_1 \theta_{21} + \omega_1((\text{off}, \text{on}))) = \text{sgn}(-\gamma_1 \theta_{21} + L_{11}U_{12}) = 1.$$

These values are related to the labeling map via $\mathcal{L}(\tau', 1, +) = L_1((\text{off}, \text{off}), 2)$ and $\mathcal{L}(\tau'', 1, +) = L_1((\text{off}, \text{on}), 2)$. 
3.3. Characterization of Equilibrium Cells. We now provide a theorem which characterizes the equilibrium cells of the switching system $\text{SWITCH}(Z)$ and shows that there is a unique equilibrium of $S(Z,\varepsilon)$ which converges to each equilibrium cell. The proof for the theorem can be found in Section 6.

**Theorem 3.13.** Let $Z$ be a regular switching parameter.

(a) $\tau \in \chi$ is an equilibrium cell if and only if

1. $\tau$ is a loop characteristic cell, and
2. $\Phi_j(\tau) = 0$ for each $j$.

(b) Furthermore, there is an $0 < A \in \mathbb{R}^{N \times N}$ so that for $\varepsilon < A$ any sigmoidal system $S(Z,\varepsilon)$ has a unique equilibrium $x^\varepsilon$ such that $x^\varepsilon \to \tau$ as $\varepsilon \to 0$.

The theorem in the case of regular cells is implied by Proposition 3.6 of [5]. It was shown in [28] that loop characteristic cells are a subset of equilibrium cells in the case that the sigmoidal perturbations $\sigma_{ij}(\cdot,\varepsilon)$ are Hill functions. This theorem extends these results by providing a necessary and sufficient condition for the identification of all equilibrium cells, enlarging the class of functions for which it applies, and giving uniqueness of the equilibrium $x^\varepsilon$.

By Theorem 3.13, an equilibrium cell $\kappa$ has a unique equilibrium $x^\varepsilon$ of $S(Z,\varepsilon)$ associated to it. We associate the stability of this equilibrium with the cell through the following definition.

**Definition 3.14.** An equilibrium cell $\kappa$ is stable if the associate equilibrium $x^\varepsilon$ of $S(Z,\varepsilon)$ is stable for all $\varepsilon > 0$ small enough and unstable otherwise.

The equilibrium cells of a switching system can be computed using the DSGRN software [4]. We show in Section 5 that the analysis of their stability can be reduced to the problem of stability of multiple cyclic feedback systems that are associated to each singular equilibrium cell. We therefore first discuss the stability of equilibrium cells in cyclic feedback systems.

4. Equilibrium Cells and their Stability in Cyclic Feedback Networks. This section concentrates on a particular type of a network, a cyclic feedback network, and characterizes the equilibrium cells and the stability of the equilibria they contain. In the following section, we generalize these results to arbitrary networks.

**Definition 4.1.** A cyclic feedback network (CFN) is a regulatory network $\mathbb{R}N = (V, E)$ such that $E = \{(1,2), (2,3), \ldots, (N-1,N), (N,1)\}$. A cyclic feedback system (CFS) is a switching or sigmoidal system associated to a CFN.

Throughout this section we will assume that $\mathbb{R}N$ is a cyclic feedback network. Since each node $j$ has exactly one target, $j+1$, and one source, $j-1$, the node $j$ is associated to exactly one threshold, $\theta_{j(j+1)}$, and $A_j = \sigma_{j(j-1)}$. This implies the combinatorial parameter, and thus the flow direction map, is determined by the ordering of numbers within the sets $\{\gamma_j \theta_{j(j+1)}, L_{j(j-1)}, U_{j(j-1)}\}$, $j = 1, \ldots, N \mod N$. This observation informs the following definition.

**Definition 4.2.** Given a regular switching parameter $Z = (L,U,\theta,\Gamma)$ for a cyclic feedback system, a node $j$ is essential if $L_{j(j-1)} < \gamma_j \theta_{j(j+1)} < U_{j(j-1)}$ and inessential otherwise.
Another consequence of having exactly one threshold for each node is that there is only one singular loop characteristic cell, \( \tau \), for which all the directions are singular, i.e. \( \text{sd}(\tau) = V \). The permutation \( \rho \) for this cell is defined by \( \rho^*(j) = j + 1 \). Throughout this section, \( \tau \) always denotes this cell and \( \rho = \rho^* \) will denote the associated permutation. We associate a sign to \( \rho \) which describes whether the net effect of the cycle is positive or negative:

\[
\text{sgn}(\rho) := \prod_{i=1}^{N} s_{i+1|i}.
\]

We say \( RN \) is a positive CFN if \( \text{sgn}(\rho) = 1 \) and a negative CFN if \( \text{sgn}(\rho) = -1 \).

4.1. Equilibrium Cells. This section identifies equilibrium cells of a CFN. To simplify notation, we observe as in [14] that by changing variables we may assume without loss of generality that if \( RN \) is a positive CFN, then every edge is activating, and if \( RN \) is a negative CFN, then every edge is activating except the edge \((N, 1)\). This change of variables is of the form

\[
x_j \rightarrow \alpha_j (x_j - \theta_{(j+1)j}) + \theta_{(j+1)j},
\]

where \( \alpha_j = \pm 1 \).

The following lemma specifies the equilibrium cell for a CFS at a parameter \( Z \) for which it has an inessential node.

**Lemma 4.3.** If at a switching parameter \( Z = (L, U, \theta, \Gamma) \) the CFS has at least one inessential node, then \( \text{SWITCH}(Z) \) has a unique equilibrium cell \( \kappa \) and this cell is regular. The cell \( \kappa \) is defined as follows. If \( j \) is an inessential node, the \( j \)-th projection is

\[
\pi_j(\kappa) = \begin{cases} 
(0, \theta_{(j+1)j}), & \text{if } U_{j(j-1)} < \gamma_j \theta_{(j+1)j} \\
(\theta_{(j+1)j}, \infty), & \text{if } \gamma_j \theta_{(j+1)j} < L_{j(j-1)}. 
\end{cases}
\]

If \( j \) is essential, let \( k \) be the inessential node which forms the shortest path of the form \( k \rightarrow k+1 \rightarrow \cdots \rightarrow j \), where nodes \( k+1, \ldots, j-1 \) are essential. We have two cases:

If \( \text{sgn}(\rho) = 1 \), then

\[
\pi_j(\kappa) = \begin{cases} 
(0, \theta_{(j+1)j}), & \text{if } U_{k(k-1)} < \gamma_k \theta_{(k+1)k}, \\
L_{(j+1)j}, \infty), & \text{if } \gamma_k \theta_{(k+1)k} < L_{k(k-1)}. 
\end{cases}
\]

If \( \text{sgn}(\rho) = -1 \), then

\[
\pi_j(\kappa) = \begin{cases} 
(0, \theta_{(j+1)j}), & \text{if } U_{k(k-1)} < \gamma_k \theta_{(k+1)k} \text{ and } 1 \leq k < j, \text{ or } \gamma_k \theta_{(k+1)k} < L_{k(k-1)} \text{ and } j < k \leq N, \\
(\theta_{(j+1)j}, \infty), & \text{if } \gamma_k \theta_{(k+1)k} < L_{k(k-1)} \text{ and } 1 \leq k < j, \text{ or } U_{k(k-1)} < \gamma_k \theta_{(k+1)k} \text{ and } j < k \leq N. 
\end{cases}
\]

When \( N = 2 \), and for a parameter \( Z \) where all nodes are essential, the value of the labeling map on the neighbors of \( \tau \) for a positive CFS is depicted in Figure 2(a), and for a negative CFS in Figure 2 (b). It is apparent that the positive CFS has two regular equilibrium cells and the negative CFS has no regular equilibrium cells. In either case \( \tau \) is a singular equilibrium cell. The next lemma shows that this is true for all \( N \).
Fig. 2: **Labeling map for the two node CFSs with no inessential nodes.** The labeling map on the boundary of $\mathbb{R}_+^2$ points inward (not shown). (a): $\text{RN} = (V, E) = \{(1, 2), (1 \to 2), (2 \to 1)\}$. The arrows indicate $\Phi_i(\kappa^L) = \Phi_i(\kappa^H) = \Phi_i(\tau) = 0$, $i = 1, 2$ so that $\kappa^L, \kappa^H,$ and $\tau$ are equilibrium cells. (b): $\text{RN} = (V, E) = \{(1, 2), (1 \to 2), (2 \to 1)\}$. The arrows indicate that $\Phi_2(\kappa) = 0$ but $\Phi_1(\kappa) = 1$ so that $\kappa$ is not an equilibrium cell. However, $\Phi_1(\tau) = \Phi_2(\tau) = 0$ so $\tau$ is an equilibrium cell. There are no regular equilibrium cells.

**Lemma 4.4.** If at a switching parameter $Z = (L, U, \theta, \Gamma)$ the CFS has no inessential nodes, then $\tau$ is an equilibrium cell of $\text{SWITCH}(Z)$. Furthermore,

1. If $\text{RN}$ is a positive CFN then $\text{SWITCH}(Z)$ has exactly two regular equilibrium cells defined by

$$
\kappa^L = \prod_{j=1}^N (0, \theta_{(j+1)j}) \quad \text{and} \quad \kappa^H = \prod_{j=1}^N (\theta_{(j+1)j}, \infty).
$$

2. If $\text{RN}$ is a negative CFN then $\tau$ is the unique equilibrium cell.

The proofs for these lemmas can be found in Section 7.1.

**4.2. Stability of Equilibria in Sigmoidal CFSs.** To determine stability of equilibria of $\mathcal{S}(Z, \varepsilon)$, we compute the characteristic polynomial of the Jacobian $J(\varepsilon)$. The structure of a cyclic feedback system imposes structure on $J$. In particular, we have

$$
J = \begin{pmatrix}
-\gamma_1 & \sigma'_{1N} \\
\sigma'_{21} & -\gamma_2 \\
\vdots & \ddots \\
\sigma'_{N-1N} & -\gamma_N \\
\sigma'_{N(N-1)} & -\gamma_N
\end{pmatrix}.
$$

(4.1)
Before computing the characteristic polynomial we first note that we can obtain stability of any regular equilibrium cell \( \kappa \) from \( J \).

**Proposition 4.5.** If \( \kappa \) is a regular equilibrium cell then it is stable.

**Proof.** If \( x^\circ \) is an equilibrium of \( S(Z, \varepsilon) \) which converges to \( \kappa \), then \( \sigma'_{j(j-1)}(x^\circ; \varepsilon) \) converges to 0. Therefore, for \( \varepsilon \) small enough \( J \) is strictly diagonally dominant with negative diagonal entries and thus all eigenvalues have negative real part. \( \square \)

We now give the characteristic polynomial of \( J \), which can be computed using the Leibniz Formula for the determinant. For the proof of the following lemma and remaining results of this subsection, see Section 7.2.

**Lemma 4.6.** Let \( S(Z, \varepsilon) \) be a cyclic feedback system. The characteristic polynomial of the Jacobian \( J(x; \varepsilon) \) is given by

\[
\det(J(x; \varepsilon) - \lambda I) = (-1)^N \left( \prod_{i=1}^{N} (\gamma_i + \lambda) - \text{sgn}(\rho) M(x, \varepsilon) \right)
\]

where \( M(x, \varepsilon) := \prod_{i=1}^{N} |\sigma'_{i(i-1)}(x; \varepsilon)| \).

The sign of the CFS plays a significant role in stability of equilibria. First we address positive cyclic feedback systems.

**Proposition 4.7.** Let \( \text{RN} \) be a positive CFN and \( x \) be an equilibrium of \( S(Z, \varepsilon) \). Stability of \( x \) can be determined as follows.

1. If \( M(x, \varepsilon) < \prod_j \gamma_j \) then \( x \) is asymptotically stable.
2. If \( M(x, \varepsilon) > \prod_j \gamma_j \) then \( x \) is unstable.
3. If \( M(x, \varepsilon) = \prod_j \gamma_j \) then \( S(Z, \varepsilon) \) has a steady state bifurcation at \( x \).

If an equilibrium \( x^\circ \) of \( S(Z, \varepsilon) \) converges to the singular loop characteristic cell \( \tau \), then condition (2) of Proposition 4.7 is satisfied if \( \varepsilon \) is small enough. Therefore, we have the following.

**Proposition 4.8.** Let \( \text{RN} \) be a positive CFN. If \( \tau \) is a singular equilibrium cell, then it is unstable.

Now we discuss negative cycles. When there are \( N \leq 2 \) nodes, we can compute eigenvalues at an equilibrium and show that \( \tau \) is stable.

**Proposition 4.9.** Let \( \text{RN} \) be a negative CFN with \( N \leq 2 \). If \( \tau \) is a singular equilibrium cell, then it is stable.

To address negative cyclic feedback systems with \( N > 2 \) we make an additional assumption that the \( \gamma \)'s are identical, that is \( \Gamma = I \). This allows us to explicitly compute all the eigenvalues of \( J \).

**Lemma 4.10.** Let \( \text{RN} \) be a CFN. Consider switching parameter \( Z \) with \( \Gamma = I \). Let \( \lambda_k(x, \varepsilon) \) for \( k = 0, \ldots, N-1 \) be the eigenvalues of the Jacobian \( J(x; \varepsilon) \) evaluated at \( x \). Then

\[
\lambda_k(x, \varepsilon) = \begin{cases} 
-1 + (e^{2\pi i k} M(x, \varepsilon))^\frac{1}{N}, & \text{sgn}(\rho) = 1 \\
-1 + (e^{\pi i + 2\pi i k} M(x, \varepsilon))^\frac{1}{N}, & \text{sgn}(\rho) = -1.
\end{cases}
\]
An eigenvalue with largest real part is

\[
\lambda_0(x, \varepsilon) = \begin{cases} 
-1 + M(x, \varepsilon)^\frac{1}{N}, & \text{sgn}(\rho) = 1 \\
-1 + (e^{\pi i} M(x, \varepsilon))^\frac{1}{N}, & \text{sgn}(\rho) = -1. 
\end{cases}
\]

(4.3)

The next proposition, often referred to as the secant condition for negative feedback systems [27, 26, 23, 24, 9], follows immediately from the computation of the eigenvalues.

**Proposition 4.11.** Let \( \text{RN} \) be a negative CFN, \( Z = (L, U, \theta, \Gamma) \) be a switching parameter with \( \Gamma = I \), and \( x \) be an equilibrium of \( S(Z, \varepsilon) \). If \( N > 2 \), then stability can be determined as follows.

1. If \( M(x, \varepsilon) < \sec\left(\frac{\pi}{N}\right)^N \) then \( x \) is asymptotically stable.
2. If \( M(x, \varepsilon) > \sec\left(\frac{\pi}{N}\right)^N \) then \( x \) is unstable.
3. If \( M(x, \varepsilon) = \sec\left(\frac{\pi}{N}\right)^N \) then \( S(Z, \varepsilon) \) has a Hopf bifurcation at \( x \).

If \( N > 2 \) and an equilibrium \( x^\varepsilon \) converges to the singular loop characteristic cell \( \tau \), the second condition of Proposition 4.11 holds when \( \varepsilon \) is small enough and we have the following.

**Proposition 4.12.** Let \( \text{RN} \) be a negative CFN with \( N > 2 \) and and \( Z = (L, U, \theta, \Gamma) \) be a switching parameter with \( \Gamma = I \). If \( \tau \) is a singular equilibrium cell it is unstable.

We remark that while Proposition 4.12 is a statement about sigmoidal CFS, [8] proves a stronger statement for switching CFS. In [8] it was shown that a negative switching CFS with \( N > 2 \) and no inessential nodes has a stable periodic orbit and that the singular loop characteristic cell \( \tau \) is an unstable source. This was proven even for \( \Gamma \neq I \). We suspect Proposition 4.12 holds for \( \Gamma \neq I \) as well but do not pursue it here. We also suspect that for small \( \varepsilon \) the sigmoidal CFS with \( N > 2 \) has a stable periodic orbit. This is consistent with Proposition 4.11 which suggests the existence of a supercritical Hopf bifurcation as \( \varepsilon \) increases.

5. Equilibria, Stability, and Bifurcations in General Networks. To characterize the equilibrium cells, stability and bifurcations of a network \( \text{RN} \) which is not a cyclic feedback network, we decompose \( \text{SWITCH}(Z) \) locally on a neighborhood of an arbitrary loop characteristic cell into cyclic feedback systems and then apply the results of Section 4. Before proceeding to describe the decomposition, we define the neighborhood of a cell on which the local decomposition is valid.

**Definition 5.1.** For \( \tau \in \chi \), the cell neighborhood of \( \tau \), denoted \( \mathcal{N}(\tau) \) is defined by

\[
\mathcal{N}(\tau) := \{ \kappa \in \chi \mid \tau \subset \kappa \},
\]

where \( \overline{\kappa} \) is the closure of \( \kappa \).

We will write \( x \in \mathcal{N}(\kappa) \) to denote \( x \in \tau \) for some \( \tau \in \mathcal{N}(\kappa) \). Note that the cell neighborhood of a regular cell \( \kappa \) consists only of \( \kappa \). In Figure 1(a), the cell neighborhood of the singular cell \( \kappa \) is given by \( \mathcal{N}(\kappa) = \{ \kappa, \kappa_1, \kappa^*_1 \} \). The cell neighborhood of the singular cell \( \tilde{\tau} \) of the positive toggle plus system consists of all cells contained in the interior of the gray shaded region of Figure 3(a).
5.1. Local Decomposition of SWITCH(Z) into Cyclic Feedback Systems. The idea behind the decomposition is to examine the values of $\Lambda$ on the cell neighborhood of a loop characteristic cell $\tau$. Using the next lemma we will show that for any regular direction $r$ of $\tau$, $\Lambda_r$ is constant on $\mathcal{N}(\tau)$. This will be used to show that regular directions enter trivially into the decomposition. On the other hand, for any singular direction $s$ of $\tau$, the lemma will be used to show that $\Lambda_{\rho(s)}$ takes one of two possible values. Furthermore, the value of $\Lambda_{\rho(s)}$ can only change in the $s$ direction. This will ultimately lead to the decomposition into cyclic feedback systems along the cycles in the permutation $\rho^\tau$. See Section 6 for the proof of the lemma.

**Lemma 5.2.** Let $\tau \in \chi$ and $(j,i) \in E$. If $j$ is a regular direction of $\tau$ or $j \in sd(\tau)$ and $i \neq \rho^\tau(j)$ then

- $\sigma_{ij}(\tau)$ is well defined,
- for all $\kappa \in \mathcal{N}(\tau)$, we have $\sigma_{ij}(\kappa) = \sigma_{ij}(\tau)$ is independent of $\kappa$.

Consequently if $i \notin \{\rho^\tau(j) \mid j \in sd(\tau)\}$, then

- $\Lambda_i(\tau)$ is well defined,
- for all $\kappa \in \mathcal{N}(\tau)$ we have $\Lambda_i(\kappa) = \Lambda_i(\tau)$ is independent of $\kappa$,

Let $\tau \in \chi$ be a loop characteristic cell, $\rho = \rho^\tau$ be the corresponding permutation. By Lemma 5.2, for each $x$ in a cell-neighborhood of $\tau$, $x \in \mathcal{N}(\tau)$, and each regular direction $r$, the value of $\dot{x}_r$ in the switching system (2.1) is independent of every other variable.

Also by Lemma 5.2, $\sigma_{sj}(\kappa) = \sigma_{sj}(\tau)$ is well defined for each singular direction $s$ and $j \in S(s) \setminus \{\rho^{-1}(s)\}$ on any cell $\kappa \in \mathcal{N}(\tau)$. Therefore, the value of $\dot{x}_s$ in the switching system (2.1) for any singular direction $s$ is independent of the regular directions. We conclude that on $\mathcal{N}(\tau)$, the switching system SWITCH(Z) decomposes into two independent systems: one corresponding to the singular directions and one corresponding to the regular directions. Since for each regular direction $r$, $\dot{x}_r$ in (2.1) depends only on $x_r$ itself, the regular directions system consists of a collection of uncoupled one dimensional systems.

The dynamics of the singular directions decomposes according to the cycles that generate the permutation $\rho$. Let $\rho_{sd(\tau)} = (c_1, \ldots, c_n)$ be the cycle decomposition of $\rho$ restricted to the singular directions. Let $\ell_d = \text{length}(c_d)$ and $s_d = \sum_{j<d} \ell_j$. We reorder the variables so that $c_d$ acts on $\{s_d + 1, s_d + 2, \ldots, s_d + \ell_d\}$ and $c_d(s_d + i) = s_d + i + 1$ for $i < \ell_d$ and $c_d(s_d + \ell_d) = s_d + 1$. On $\mathcal{N}(\tau)$, the dynamics of the variables $x^d = (x_{s_d+1}, \ldots, x_{s_d+\ell_d})$ are independent of the value of all other variables so that the dynamics of the singular directions can be written as $n$ independent systems. Within each system, $\dot{x}_s$ depends only on $\rho^{-1}(s)$ so each of these $n$ systems are cyclic feedback systems.

We conclude that the switching system SWITCH(Z) restricted to $\mathcal{N}(\tau)$ decomposes into $d$ cyclic feedback systems and a diagonal system corresponding to regular directions. The sign of the cycle

$$\text{sgn}(c_d) := \prod_{j=s_d+1}^{s_d+\ell_d} s_{c_d(j)}$$

determines whether the $d$th system is a positive or negative CFS.
To explicitly write the decomposition we define $\ell_{n+1} := N - s_{n+1}$ to be the number of regular directions. For $d = 1, \ldots, n + 1$ we define the projections of the cell neighborhood $N(\tau)$ and cells $\kappa \in N(\tau)$

$$
\Lambda^d(\tau) := \prod_{j=s_d+1}^{s_d+\ell_d} \pi_j(N(\tau)) \quad \text{and} \quad \kappa^d := \prod_{j=s_d+1}^{s_d+\ell_d} \pi_j(\kappa).
$$

We set $\Lambda(\cdot; \tau) := \Lambda|_{N(\tau)}$ to be the restriction of $\Lambda$ onto $N(\tau)$. We then define

$$
\Lambda^d(\cdot; \tau) := (\Lambda_{s_d+1}(\cdot; \tau), \ldots, \Lambda_{s_d+\ell_d}(\cdot; \tau))
$$

to be the projection of the resulting function onto the directions in the $d$-th subsystem.

Further, let $\Gamma^d$ be the $\ell_d \times \ell_d$ diagonal matrix with entries $\Gamma_{ii} = \gamma_{s_d+i}$ for $i = 1, \ldots, \ell_d$. The dynamics for the $d$-th subsystem is given explicitly by

$$
\dot{x}^d = -\Gamma^d x^d + \Lambda^d(x^d; \varepsilon, \tau), \quad x^d \in N^d(\tau; \varepsilon)
$$

which we denote by $\text{SWITCH}^d(Z; \tau)$. On the collection of cells $N(\tau)$, the system is

$$
\text{SWITCH}(Z; \tau) := (\text{SWITCH}^1(Z; \tau), \ldots, \text{SWITCH}^n(Z; \tau), \text{SWITCH}^{n+1}(Z; \tau)).
$$

We denote the cyclic feedback network associated with $\text{SWITCH}^d(Z; \tau)$ by $\text{RN}^d(\tau)$.

**Example 5.3.** Consider the positive toggle plus system of Example 2.3. For the loop characteristic cell $\vec{\tau} = \{\theta_{11}\} \times \{\theta_{22}\}$ pictured in Figure 1(b), the singular directions of $\vec{\tau}$ are both directions 1 and 2 and the cycle decomposition is given by $\rho = (c_1, c_2)$ where $c_1(1) = 1$ and $c_2(2) = 2$. The cycle decomposition corresponds to the disjoint loops $1 \to 1$ and $2 \to 2$ in the network. The cell neighborhood of $\vec{\tau}$, $N(\vec{\tau})$ is the gray shaded region in Figure 3(a). For $x \in N(\vec{\tau})$, $x_1 > \theta_{21}$ and $x_2 > \theta_{12}$ so that

$$
\dot{x}_1 = -\gamma_1 x_1 + \Lambda_1(x) = -\gamma_1 x_1 + U_{12} \sigma_{11}(x_1)
$$

$$
\dot{x}_2 = -\gamma_1 x_2 + \Lambda_2(x) = -\gamma_2 x_2 + U_{21} \sigma_{22}(x_2).
$$

The decomposition at $\vec{\tau}$, $\text{SWITCH}(Z) = (\text{SWITCH}^1(Z; \vec{\tau}), \text{SWITCH}^2(Z; \vec{\tau}))$ is defined by $x^1 = x_1$, $x^2 = x_2$, $\Gamma^1 = \gamma_1$, $\Gamma^2 = \gamma_2$, $\Lambda^1(x_1; \vec{\tau}) = U_{12} \sigma_{11}(x_1)$, and $\Lambda^2(x_2; \vec{\tau}) = U_{21} \sigma_{22}(x_2)$. Note that the superscripts index the cycles, not the directions.

Next we show every regular cell in $\chi$ lies in a neighborhood of at least one singular loop characteristic cell. This implies that every regular cell will be in at least one neighborhood $N(\tau)$ on which decomposition into cyclic feedback systems is applicable. Therefore, in spite of of being local, the decomposition into cyclic feedback systems affects all the cells in $\chi$. In particular, all equilibrium cells of $\text{SWITCH}(Z)$ are contained in at least one neighborhood $N(\tau)$.

**Lemma 5.4.** Suppose that every node of $\text{RN}$ has an out-edge. Then for every regular cell $\kappa \in \chi^{(N)}$, there is a singular loop characteristic cell $\tau \in \chi$ so that $\kappa \in N(\tau)$.

**Proof.** Let $\kappa$ be a regular cell. Recall that for every $j$, $\pi_j(\kappa) = (\theta_{a_j^i, \theta b_j^i})$ where $a_j^i, b_j^i \in V \cup \{-\infty, \infty\}$. Since every $j$ has an out-edge and hence a target node $i_j$, there is a threshold $\theta_{i_j}$ so that $\theta_{i,j}$ is a boundary of $\pi_j(\kappa)$. In particular, since $i_j$ is a network node, $i_j \notin \{\theta_{-\infty}, \theta_{\infty}\}$. The
selection of such a threshold for every \( j \in V \) defines a map \( \sigma : V \to V \) by \( \sigma(j) = i_j \). Note that since \( \sigma(V) \subset V \) and \( V \) is finite, the set

\[
U := \bigcup_{k=1}^{\infty} \sigma^k(V)
\]

is non-empty and satisfies \( U = \sigma(U) \). Then a cell \( \tau \) defined by

\[
\pi_j(\tau) = \begin{cases} \pi_j(\kappa), & j \notin U \\ \{\theta_{\sigma(j)}\}, & j \in U \end{cases}
\]

is a loop characteristic cell with \( \kappa \in \mathcal{N}(\tau) \).

5.2. Equilibrium Cells. First, we generalize Section 4.1 and characterize the equilibrium cells of \( \text{SWITCH}(\mathbb{Z}) \) in \( \mathcal{N}(\tau) \). We will take advantage of the decomposition

\[
\text{SWITCH}(\mathbb{Z}) = (\text{SWITCH}^1(\mathbb{Z}; \tau), \ldots, \text{SWITCH}^n(\mathbb{Z}; \tau), \text{SWITCH}^{n+1}(\mathbb{Z}; \tau))
\]

which is valid on \( \mathcal{N}(\tau) = \mathcal{N}^1(\tau) \times \cdots \times \mathcal{N}^n(\tau) \times \mathcal{N}^{n+1}(\tau) \) (equation 5.1). However, the lemmas of Section 4.1 are only valid for cyclic feedback systems which are defined on the whole positive orthant. We therefore extend \( \text{SWITCH}^d(\mathbb{Z}; \tau) \) to all of \( \mathbb{R}^{\ell_d} \) for \( d \leq n \). The regular directions do not form a cyclic feedback system on \( \mathcal{N}(\tau) \) so we do not need to extend \( \text{SWITCH}^{n+1}(\mathbb{Z}; \tau) \).

The dynamics of \( \text{SWITCH}^d(\mathbb{Z}; \tau) \) are given by

\[
x^d = -\Gamma^d x^d + \Lambda^d(x^d; \tau), \quad x^d \in \mathcal{N}^d(\tau).
\]

To extend the domain of definition of \( \text{SWITCH}^d(\mathbb{Z}; \tau) \) on \( \mathbb{R}^{\ell_d} \), we need to define \( \Lambda^d(\cdot; \tau) \) on \( \mathbb{R}^{\ell_d} \), while ensuring \( \text{SWITCH}^d(\mathbb{Z}; \tau) \) remains a cyclic feedback system. To do so, we make the following definition.

**Definition 5.5.** Given a loop characteristic cell \( \tau \subset \mathbb{R}_+^N \) with permutation \( \rho = \rho^\tau \), the cone \( C(\kappa; \tau) \) rooted in \( \tau \) and induced by a cell \( \kappa \in \mathcal{N}(\tau) \) is defined by its \( N \) projections. For a regular direction, \( r \), of \( \tau \)

\[
\pi_r(C(\kappa; \tau)) := \pi_r(\tau).
\]

For a singular direction, \( s \in \text{sd}(\tau) \),

\[
\pi_s(C(\kappa; \tau)) := \begin{cases} \{\theta_{(s)}\}, & \text{if } \pi_s(\kappa) = \{\theta_{(s)}\} \\ \{\theta_{(s)}, \infty\}, & \text{if } \pi_s(\kappa) = (\theta_{(s)}, \theta_{(s)}^{\beta}) \\ (0, \theta_{(s)}), & \text{if } \pi_s(\kappa) = (\theta_{(s)}^{\beta}, \theta_{(s)}^{\beta+1}) \\ \{\theta_{(s)}^{\beta+1}\}, & \text{if } \pi_s(\kappa) = (\theta_{(s)}, \theta_{(s)}^{\beta}) \end{cases}
\]

For \( \sigma \in \mathcal{N}^d(\tau) \) define a \( d \)-cone in \( \mathbb{R}^{\ell_d}_+ \) by

\[
C^d(\sigma; \tau) := \prod_{j=s_d+1}^{s_d+\ell_d} \pi_j(C(\sigma; \tau)).
\]
Fig. 3: The decomposition of the positive toggle plus system at a loop characteristic cell $\tau$. The switching parameter is chosen as in Example 2.3. (a): The cell complex $\chi$ and the labeling map. The cell neighborhood $N(\tau)$ consists of all cells in the shaded gray region. The regular cells in the cell neighborhood are labeled. (b): The cell complex and labeling map formed by the product of switching systems $(\text{SWITCH}_1(Z; \tau), \text{SWITCH}_2(Z; \tau))$ after extending their domains. The cell complex is formed by the cones rooted in $\tau$. The cones induced by the regular cells are labeled. (c): The cell complex $\chi^2(\tau)$ and labeling map $L^2$ for $\text{SWITCH}_2(Z; \tau)$. The cell $\tau^-_2$ is a 2-candidate equilibrium cell because $C^2(\tau^-_2; \tau)$ is an equilibrium cell of $\text{SWITCH}_2(Z; \tau)$. It is also 2-consistent. (d): The cell complex $\chi^1(\tau)$ and labeling map $L^1$ for $\text{SWITCH}_1(Z; \tau)$. Each of $\tau$, $\tau^-_1$, and $\tau^+_1$ are candidate equilibrium cells. The cells $\tau$ and $\tau^+_1$ are 1-consistent, while $\tau^-_1$ is not. See Example 5.12 for more details on (a) and (b).

The cones for a loop characteristic cell $\tau$ in a two node network are depicted in Figure 3.

We now proceed to extend the function $\Lambda^d$ from $\mathcal{N}^d(\tau)$ to $\mathbb{R}^d_{\tau^+}$. Take $x^d \in \mathbb{R}^d_{\tau^+}$. If $x^d \notin \mathcal{N}^d(\tau)$, then $x^d \in C^d(\sigma; \tau)$ for some $\sigma \in \mathcal{N}^d(\tau)$. We define the value of $\Lambda^d$ on $C^d(\sigma; \tau)$ to be the value of $\Lambda^d$ on $\sigma$. Explicitly,

$$\Lambda^d(x^d; \tau) := \Lambda^d(\sigma), \quad x^d \in C^d(\sigma; \tau).$$

Given this extension, $\text{SWITCH}^d(Z; \tau)$ is a switching system defined on $\mathbb{R}^d_{\tau^+}$ with an associated cell complex, labeling map, and flow direction map which we denote $\chi^d(\tau)$, $\mathcal{L}^d(\cdot, \cdot; \tau)$, and $\Phi^d(\cdot; \tau)$, respectively. Here, the threshold set which generates $\chi^d(\tau)$, $\Theta^d := (\Theta_{s,+1}^d, \ldots, \Theta_{s,+d}^d)$, is defined
Lemmas 4.3 and 4.4 can be used to determine the equilibrium cells of $\text{SWITCH}^d(Z; \tau)$. Since the cell complex is formed by the cones (see Figure 3(c) and (d))
\[
\chi^d(\tau) = \{ C^d(\sigma; \tau) | \sigma \in \mathcal{N}^d(\tau) \},
\]
there is a straightforward identification between cells of $\text{SWITCH}^d(Z; \tau)$ and cells $\sigma \in \mathcal{N}^d(\tau)$.

**Definition 5.6.** The $d$-candidate equilibrium cells of $\tau$ are defined by
\[
\text{Eq}^d(\tau) := \{ \sigma^d | \sigma \in \mathcal{N}(\tau) \text{ and } C^d(\sigma; \tau) \text{ is an equilibrium cell of } \text{SWITCH}^d(Z; \tau) \}.
\]

The candidate equilibrium cells of $\tau$ are $\text{Eq}(\tau) := \prod_{d=1}^{n+1} \text{Eq}^d(\tau)$.

See Figure 3 and Example 5.12 for examples of candidate equilibrium cells in the positive toggle plus system. The next lemma justifies the name candidate equilibrium cell.

**Lemma 5.7.** If $\kappa \in \mathcal{N}(\tau)$ is an equilibrium cell, then $\kappa \in \text{Eq}(\tau)$.

Before proving the lemma, we isolate part of the proof that will be useful later.

**Lemma 5.8.** Let $\kappa \in \mathcal{N}(\tau)$. If $r$ is a regular direction of $\tau$ then $\Phi_r(\kappa) = \Phi^*_{n+1}(\kappa; \tau)$. For any singular direction $s$ of $\kappa$ that belongs to $\{ s_d + 1, \ldots, s_d + \ell_d \}$ we have $\Phi_s(\kappa) = \Phi_s^d(\kappa; \tau)$.

**Proof.** First consider the case that $r$ is a regular direction of $\tau$. Then $r$ is also a regular direction of $\kappa$ and $\pi_r(\kappa) = \pi_r(\tau) = (\theta_{a^r \tau}, \theta_{b^r \tau})$. Further, by Lemma 5.2, $\Lambda_r(\tau) = \Lambda_r(\kappa)$. Finally we have $\Lambda^*_{n+1}(C(\kappa; \tau); \tau) = \Lambda_r(\kappa) = \Lambda_r(\tau)$. These observations allow us to compute the labeling map of $\kappa$ in the $r$-th direction. We have
\[
\mathcal{L}^{n+1}(C(\kappa; \tau), r, -) = \text{sgn}(\gamma_r a^r + \Lambda^*_n(C(\kappa; \tau); \tau)) = \text{sgn}(\gamma_r a^r + \Lambda_r(\tau)) = \mathcal{L}(\tau, r, -)
\]
and
\[
\mathcal{L}^{n+1}(C(\kappa; \tau), r, +) = \text{sgn}(\gamma_r a^r + \Lambda^*_n(C(\kappa; \tau); \tau)) = \text{sgn}(\gamma_r a^r + \Lambda_r(\tau)) = \mathcal{L}(\tau, r, +).
\]

Since the labeling maps agree and the flow direction maps are defined via the labeling maps (see Definition 3.9), we have $\Phi_r(\tau) = \Phi^*_{n+1}(C(\kappa; \tau); \tau)$.

Now suppose $s \in \{ s_d + 1, \ldots, s_d + \ell_d \}$ is a singular direction of $\kappa$ for some $d = 1, \ldots, n$. Then
\[
\mathcal{L}(\kappa, s, -) = \text{sgn}(\gamma_r a^r + \Lambda^*_n(\kappa_s^-)) \quad \text{and} \quad \mathcal{L}(\kappa, s, +) = \text{sgn}(\gamma_r a^r + \Lambda^*_n(\kappa_s^+))
\]
Since $\Lambda^*_n(\kappa_s^\pm) = \Lambda^*_n(\kappa^\pm; \tau)$, we have $\mathcal{L}^d(C(\kappa; \tau), s, -) = \mathcal{L}^d(C(\kappa; \tau), s, -)$ and $\mathcal{L}(\kappa, s, +) = -\mathcal{L}(\kappa, s, +)$, or $\Phi_s(\kappa) = \Phi^d_s(C(\kappa; \tau); \tau)$.
Proof of Lemma 5.7. Let \( \kappa \in \mathcal{N}(\tau) \) be an equilibrium cell. By Theorem 3.13 we have \( \Phi_j(\kappa) = 0 \) for every \( j \in V \). By Lemma 5.8, we have
\[
\Phi^{n+1}_r(C(\kappa; \tau); \tau) = \Phi_r(\kappa) = 0 \quad \text{and} \quad \Phi^d(\kappa; \tau) = \Phi_d(\kappa) = 0
\]
both for regular directions \( r \) of \( \tau \) and singular directions \( s \) of \( \kappa \) for some \( d = 1, \ldots, n \). This implies \( \kappa^{n+1} \in \mathcal{E}^{n+1}(\tau) \) and \( \kappa^d \in \mathcal{E}^d(\tau) \) where \( s \in \{s_d + 1, \ldots, s_d + \ell_d\} \).

Suppose \( j \in \{s_d + 1, \ldots, s_d + \ell_d\} \) is a singular direction of \( \tau \) but a regular direction of \( \kappa \). Assume without loss of generality that \( \pi_{\rho(j)}(\kappa) = (\theta_{\rho(j)j}, \theta_{\rho(j)j}) \). Then we have
\[
1 = L(\kappa, j, -) = \text{sgn}(\gamma_j \theta_{\rho(j)j} + \Lambda_j(\kappa)) = - \text{sgn}(\gamma_j \theta_{\rho(j)j} + \Lambda_j(\kappa)) = -L(\kappa, j, +).
\]
Now, \( L^d(C(\kappa; \tau), j, +) = L(\kappa, j, +) \) so we only need to check \( L(\kappa, j, -) = 1 = L^d(\kappa^d, j, -) \). Since \( \Theta_j = \{0, \theta_{\rho(j)j}, \infty\} \), we have
\[
L^d(C(\kappa; \tau), j, -) = \text{sgn}(\gamma_j \theta_{\rho(j)j} \cdot 0 + \Lambda_{\rho(j)}(\kappa^d)) = 1
\]
so that \( \Phi^d(\kappa^d, \tau) = 0 \) and \( \kappa^d \in \mathcal{E}^d(\tau) \). Since \( \kappa^d \in \mathcal{E}^d(\tau) \) for each \( d, \kappa \in \mathcal{E}(\tau) \).

Once we have identified a candidate equilibrium cell \( \kappa \in \mathcal{E}(\tau) \), there is one more property to check to ascertain that it is an actual equilibrium cell. Since \( \kappa^d \in \mathcal{E}^d(\tau) \), the cell \( C^d(\kappa^d; \tau) \) must be an equilibrium cell of \( \text{SWITCH}^d(Z; \tau) \). However, \( C^d(\kappa^d; \tau) \) is a super set of \( \kappa^d \), so this equilibrium may not be contained in \( \kappa^d \).

Definition 5.9. Let \( \sigma \in \mathcal{E}^d(\tau) \) and \( d \leq n \). If the equilibrium of \( \text{SWITCH}^d(Z; \tau) \) in \( C^d(\sigma; \tau) \) is contained in \( \sigma \), then we say \( \sigma \) is a \( d \)-consistent candidate equilibrium cell.

The next proposition shows how to check if \( \sigma \) is a \( d \)-consistent candidate equilibrium cell. Notice that the condition of the proposition is vacuously satisfied when \( \sigma \) is a singular equilibrium cell of \( \text{SWITCH}^d(Z; \tau) \). In other words, if the singular loop characteristic cell \( \tau^d \) is an equilibrium cell, it is always \( d \)-consistent. It is also important to note, that while seemingly complicated, this condition is readily algorithmically checkable.

Proposition 5.10. Let \( d \leq n \) and \( \sigma \in \mathcal{E}^d(\tau) \). Then \( \sigma \) is a \( d \)-consistent candidate equilibrium cell if, and only if, for each \( j \in \{s_d + 1, \ldots, s_d + \ell_d\} \), if \( \pi_j(\sigma) \) is equal to the \( i \)-th value in first column, and \( \pi_{j-1}(\sigma) \) is equal to the \( k \)-th value in first row, then the inequality in entry \((i, k)\) of the following table is satisfied.

| \( \pi_j(\sigma) \) | \( \pi_{j-1}(\sigma) \) | \( (\theta_{\rho(j)j}, \theta_{\rho(j)j}) \) | \( (\theta_{j(j-1)(j-1)}, \theta_{j(j-1)}) \) | \( (\theta_{j(j-1)}, \theta_{j(j-1)(j-1)}) \) |
|-------------------|-------------------|---------------------|---------------------|---------------------|
| \( \gamma_j \theta_{\rho(j)j} \) | \( \Lambda_j(\tau^{j-1}_{j-1}) \) | \( \gamma_j \theta_{\rho(j)j} < \Lambda_j(\tau^{j-1}_{j-1}) \) | \( \gamma_j \theta_{\rho(j)j} < \Lambda_j(\tau^{j-1}_{j-1}) \) | \( \gamma_j \theta_{\rho(j)j} < \Lambda_j(\tau^{j-1}_{j-1}) \) |

Here \( s_d + \ell_d + 1 \) is identified with \( s_d + 1 \) and \( (s_d + 1) - 1 \) is identified with \( s_d + \ell_d \).

Proof. Let \( \Phi^d \) denote the flow direction map for the cyclic feedback system \( \text{SWITCH}^d(Z; \tau) \). We show that \( \Phi^d(\sigma) = 0 \) if and only if the conditions are satisfied. Applying Theorem 3.13 then completes the proof.
We first assume that \( \sigma \) is the singular loop characteristic cell of \( \text{SWITCH}^d(Z, \tau) \). Then \( \sigma = \mathcal{C}^d(\sigma; \tau) \). Since \( \sigma \in \text{Eq}^d(\tau) \), \( \mathcal{C}^d(\sigma; \tau) \) is an equilibrium cell of \( \text{SWITCH}^d(Z; \tau) \) so that

\[
\Phi^d(\sigma) = \Phi^d(\mathcal{C}^d(\sigma; \tau)) = 0.
\]

If \( \sigma \) is not the singular loop characteristic cell then \( \sigma \) must be a regular cell of \( \text{SWITCH}^d(Z; \tau) \) by Lemma 4.3 and 4.4. Since \( \sigma \in \mathcal{N}(\tau) \),

\[
\pi_j(\sigma) = (\theta_{(j+1)j}, \theta_{p_+(j)j}) \quad \text{or} \quad \pi_j(\sigma) = (\theta_{p_-(j)j}, \theta_{j(j+1)j})
\]

for each \( j \).

Fix \( j \) and suppose \( \sigma \) satisfies \( \pi_j(\sigma) = (\theta_{(j+1)j}, \theta_{p_+(j)j}) \). Since \( \sigma \in \text{Eq}^d(\tau) \), we have \( \Phi_j^d(\mathcal{C}(\sigma; \tau)) = 0 \). Since \( \pi_j(\sigma) \) and \( \pi_j(\mathcal{C}(\sigma; \tau)) \) have the same left boundary (by the assumption on \( \sigma \)), we have

\[
1 = \mathcal{L}^d(\mathcal{C}(\sigma; \tau), j, -; 0) = \mathcal{L}^d(\sigma, j, -; 0).
\]

Therefore, \( \Phi_j^d(\sigma) = 0 \) if and only if \( \mathcal{L}^d(\sigma, j, +; 0) = -1 \). This is equivalent to

\[
\text{sgn}(\gamma_{j}^{\theta_{p_+(j)j} + \Lambda_j(\sigma)}) = -1, \quad \text{or} \quad \Lambda_j(\sigma) < \gamma_{j}^{\theta_{p_+(j)j}}.
\]

If \( \pi_{j-1}(\sigma) = (\theta_{p_-(j-1)(j-1)}, \theta_{j(j-1)}) \) then \( \sigma \in \mathcal{N}(\tau^-_{j-1}) \) and if \( \pi_{j-1}(\sigma) = (\theta_{j(j-1)}, \theta_{p_+(j-1)(j-1)}) \) then \( \sigma \in \mathcal{N}(\tau^+_{j-1}) \). Notice that \( j - 1 \) is the only singular direction of \( \tau \) that maps to \( j \) under \( \rho \) and \( j - 1 \) is not a singular direction of \( \tau^-_{j-1} \) or \( \tau^+_{j-1} \). Therefore, by Lemma 5.2, \( \Lambda_j(\sigma) = \Lambda_j(\tau^-_{j-1}) \) or \( \Lambda_j(\sigma) = \Lambda_j(\tau^+_{j-1}) \) according to value of \( \pi_{j-1}(\sigma) \). Together with (5.2), this observation proves the proposition in the case \( \pi_j(\sigma) = (\theta_{(j+1)j}, \theta_{p_+(j)j}) \). A similar argument applies to the case \( \pi_j(\sigma) = (\theta_{p_-(j)j}, \theta_{j(j+1)j}) \).

See Example 5.12 for an example application of Proposition 5.10. The next theorem shows that any equilibrium cell in \( \mathcal{N}(\tau) \) can be written as a product over consistent candidate equilibrium cells.

**Theorem 5.11.** Let \( \kappa \in \mathcal{N}(\tau) \). Then \( \kappa \) is an equilibrium cell of \( \text{SWITCH}(Z) \) if and only if

- \( \kappa \in \text{Eq}(\tau) \), and
- for every \( d \), \( \kappa^d \) is a \( d \)-consistent candidate equilibrium cell.

**Proof.** Let \( \kappa \in \mathcal{N}(\tau) \). By Lemma 5.7, \( \kappa \in \text{Eq}(\tau) \) is a necessary condition for \( \kappa \) to be an equilibrium cell. We may therefore assume \( \kappa \in \mathcal{E}(\tau) \).

By Theorem 3.13, \( \kappa \) is an equilibrium cell if and only if \( \Phi_j(\kappa) = 0 \) for every \( j \in V \). Lemma 5.8 shows that if \( r \) is a regular direction of \( \tau \) then \( \Phi_{\rho}(\kappa) = \Phi_{\rho}^n(C(\kappa; \tau)) \) and if \( s \in \{s_d + 1, \ldots, s_d + \ell_d\} \) is a singular direction of \( \kappa \) then \( \Phi_s(\kappa) = \Phi_s^d(C(\kappa; \tau)) \). It remains to show that \( \Phi_j(\kappa) = 0 \) if and only if \( \kappa^d \) is \( d \)-consistent when \( j \in \{s_d + 1, \ldots, s_d + \ell_d\} \) is a regular direction of \( \kappa \) but singular direction of \( \tau \). There are four cases determined by the values of \( \pi_j(\kappa) \) and \( \pi_{j-1}(\kappa) \) indicated by the table in the statement of Proposition 5.10. We prove the case \( \pi_j(\kappa) = (\theta_{p_-(j)j}, \theta_{p(j)j}) \) and \( \pi_{j-1}(\kappa) = (\theta_{p_-(j-1)(j-1)}, \theta_{j(j-1)}) \). The remaining cases are similar. Since \( \pi_{j-1}(\kappa) = (\theta_{p_-(j-1)(j-1)}, \theta_{j(j-1)}) \) we have \( \Lambda_j(\kappa) = \Lambda_j(\tau^-_{j-1}) \) so that

\[
\mathcal{L}(\kappa, j, -) = \text{sgn}(\gamma_j^{\theta_{p_-(j)j} + \Lambda_j(\tau^-_{j-1}))}.
\]
Therefore $\mathcal{L}(\kappa, j, \varepsilon) = 1$ if and only if $\kappa^d$ is $d$-consistent. Since $\kappa^d \in \text{Eq}^d(\tau)$,

$$
\mathcal{L}^d(C(\kappa^d; \tau), j, \varepsilon) = \text{sgn}(-\gamma_j \cdot 0 + \Lambda_j^d(C(\kappa^d; \tau))) = 1 = -\mathcal{L}^d(C(\kappa^d; \tau), j, +).
$$

But $\mathcal{L}^d(C(\kappa; \tau), j, +) = \text{sgn}(-\gamma_j \theta(\tau) + \Lambda_j^d(C(\kappa; \tau))) = \mathcal{L}(\kappa, j, +)$ since $\Lambda_j^d(C(\kappa; \tau)) = \Lambda_j(\kappa)$. Therefore $\Phi_j(\kappa) = 0$ if and only if $\kappa^d$ is $d$-consistent, completing the proof of the theorem.

Example 5.12. Consider the decomposition of the positive toggle plus system described in Example 5.12. Consider the decomposition of the positive toggle plus system described in Example

5.3. Stability of Equilibria. The decomposition in Theorem 5.11 implies that the Jacobian $J(\varepsilon)$ for $S(Z, \varepsilon)$ has a block structure up to order $\varepsilon$. In particular, we have that there is a $B \in \mathbb{R}^{N \times N}$ such that

$$
J(\varepsilon) = \begin{pmatrix}
J^1(\varepsilon) & J^2(\varepsilon) & \cdots & J^n(\varepsilon)
\end{pmatrix} + \varepsilon B,
$$

and $J^d(\varepsilon)$ for $1 \leq d \leq n$ has the same structure as (4.1) although the entries $\sigma'_{(j+1)j}$ are replaced with $\frac{\partial}{\partial x_j} \Lambda_{j+1}(x; \varepsilon)$. The block $J^d(\varepsilon)$ is the Jacobian for the sigmoidal system $S^d(Z, \varepsilon; \tau)$ obtained from perturbing the switching system $\text{SWITCH}^d(Z; \tau)$. This implies that the stability of an equilibrium $x^\varepsilon$ associated to an equilibrium cell $\kappa \in \mathcal{N}(\tau)$ is determined by the stability of $\kappa^d$ as an equilibrium cell of $\text{SWITCH}^d(Z; \tau)$ for $d = 1, \ldots, n+1$. We formally state this observation as a theorem after the following definition.

Definition 5.13. An equilibrium cell $\kappa \in \mathcal{N}(\tau)$ is $d$-stable if $\kappa^d$ is stable as a cell of $\text{SWITCH}^d(Z; \tau)$, and $d$-unstable otherwise.

Theorem 5.14. An equilibrium cell $\kappa \in \mathcal{N}(\tau)$ is stable if and only if $\kappa$ is $d$-stable for each $d$.

As a result of Theorem 5.14, we can immediately generalize the stability results given by Propositions 4.5, 4.8, 4.9, and 4.12, which results in the following theorem.

Theorem 5.15. Let $Z = (L, U, \theta, \Gamma)$ be a switching parameter, $\tau$ be a loop characteristic cell, and $\kappa \in \mathcal{N}(\tau)$ be an equilibrium cell. Then $d$-stability of $\kappa$ can be determined as follows.

1. If $\kappa^d$ is a regular cell then $\kappa$ is $d$-stable.
2. If $\kappa^d$ is a singular cell and $c_d$ is positive, then $\kappa$ is $d$-unstable.
3. If $\kappa^d$ is a singular cell, $c_d$ is negative, and $\ell_d \leq 2$, then $\kappa$ is $d$-stable.
4. If $\Gamma = I$, $\kappa^d$ is a singular cell, $c_d$ is negative, and $\ell_d > 2$, then $\kappa$ is $d$-unstable.

We remark that an equilibrium cell $\kappa$ is always $d = n + 1$-stable since the only equilibrium cell of $\text{SWITCH}^{n+1}(Z; \tau)$ is regular.

Example 3.16. Consider the decomposition of the positive toggle plus system at $\bar{\tau}$ pictured in Figure 3. The equilibrium cell $\kappa_4$ is 1-stable because $\pi_1(\kappa_4) = C^1(\kappa_4; \bar{\tau})$ is stable in $\text{SWITCH}^1(Z; \bar{\tau})$ as can be seen in Figure 3(d). $\kappa_4$ is also 2-stable because $\pi_2(\kappa_4) = C^2(\kappa_4; \bar{\tau}) = C^2(\tilde{\tau}_2; \bar{\tau})$ is stable in $\text{SWITCH}^2(Z; \bar{\tau})$ as can be seen in Figure 3(c). The equilibrium cell $\tilde{\tau}_2$ is 1-unstable because $\pi_1(\tilde{\tau}_2) = \tau^1$ is unstable in $\text{SWITCH}^1(Z; \bar{\tau})$.

6. Proof of Theorem 3.13. Before proving the theorem, we will prove some technical lemmas. We begin by proving Lemma 5.2.

Proof of Lemma 5.2. Let $\tau \in \chi$ be a cell, $\kappa \in \mathcal{N}(\tau)$ be a cell in the cell neighborhood $\mathcal{N}(\tau)$ of $\tau$ (see Definition 5.1), and $(j, i) \in E$ be an edge. Suppose $j$ is a regular direction of $\tau$ with $\pi_j(\tau) = (\theta_{ij}, \theta_{iz})$. Then $\sigma_{ij}(\tau)$ is well defined. Since $\tau \subset \chi,$

\[
\langle \theta_{ij}, \theta_{iz} \rangle = \pi_j(\tau) \subset \pi_j(\chi) = \{\theta_{ij}, \theta_{iz}\}.
\]

Because $\theta_{ij} < \theta_{iz}$ are consecutive thresholds we must have $\theta_{ij} = \theta_{ij}$ and $\theta_{iz} = \theta_{iz}$. Therefore $\pi_j(\kappa) = \pi_j(\tau)$. This implies $\sigma_{ij}(\kappa) = \sigma_{ij}(\tau)$.

Now suppose $j \in \partial(\tau)$, and $\pi_j(\tau) = \{\theta_{0j}\}$. Assume $i_0 \neq i$. This implies that $\sigma_{ij}(\theta_{0j})$ is defined and $\sigma_{ij}(\tau) = \sigma_{ij}(\theta_{0j})$. Let $\theta_{ij} < \theta_{ij} < \theta_{ij}$ be consecutive thresholds. Then either $\pi_j(\kappa) = (\theta_{ij}, \theta_{ij}), \pi_j(\kappa) = (\theta_{ij}, \theta_{ij}),$ or $\pi_j(\kappa) = \theta_{ij}$. Since $\theta_{ij} \notin (\theta_{ij}, \theta_{ij}), \sigma_{ij}(x_j) = \sigma_{ij}(\tau)$ for all $x_j \in (\theta_{ij}, \theta_{ij})$ and $\sigma_{ij}(\kappa) = \sigma_{ij}(\tau)$.

Since $\Lambda_i$ is a product of sums of switching functions $\sigma_{ij}$, $\Lambda_i(\kappa)$ is well defined whenever $\sigma_{ij}(\tau)$ is well defined for all $j \in \mathbf{S}(i)$. That is, $\Lambda_i(\kappa)$ is well defined whenever $i \notin \rho(\partial(\tau)) = \{\rho(j) \mid j \in \partial(\tau)\}$.

Recall that $\tau^\pm_s$ denotes an $s$-neighbor of a cell $\tau \in \chi$ (see Definition 3.5). A consequence of Lemma 5.2 is that the labeling map $\mathcal{L}$ and thus the flow direction map $\Phi$ (see Definition 3.9), are well defined because $\Lambda_{\rho(s)}(\tau^\pm_s)$ is well defined for every singular direction $s$ of a loop characteristic cell $\tau$.

Lemma 6.1. Let $\tau \in \chi$ be a loop characteristic cell. If $s$ is a singular direction of $\tau$ then $\Lambda_{\rho(s)}(\tau^\pm_s)$ is well defined.

Proof. Since $s$ is the unique singular direction of $\tau$ which maps to $\rho(s)$, we have $\rho(s) \notin \rho(\partial(\tau^\pm_s))$ since $s$ is a regular direction of $\tau^\pm_s$.

Having confirmed that $\mathcal{L}$ and $\Phi$ are well defined on their domain, we now prove Theorem 3.13.

6.1. Proof of Theorem 3.13. Let $\tau \in \chi$ be an equilibrium cell and $x^\tau$ be an equilibrium of $\mathcal{S}(Z, \varepsilon)$ such that $x^\tau \rightarrow x^\tau$ as $\varepsilon \rightarrow 0^+$. Let $\rho = \rho^\tau$.

First suppose that statement (1) does not hold. Then there is a singular direction $s \in \partial(\tau)$ so that $\rho(j) \neq s$ for all $j \in \partial(\tau)$. By Lemma 5.2, $\Lambda_s(\kappa)$ is well defined and $\Lambda_s(\kappa) = \Lambda_s(\tau)$ for all $\kappa \in \mathcal{N}(\tau)$. Therefore, for all $x \in \mathcal{N}(\tau)$, $\Lambda_s(x; \varepsilon) \rightarrow \Lambda_s(\tau)$ as $\varepsilon \rightarrow 0$. Since $x^\tau \rightarrow x^\tau \in \tau$, there is an $A > 0$ so that for $\varepsilon < A$, $x^\tau \in \mathcal{N}(\tau)$. Therefore, $\Lambda_s(x^\tau, \varepsilon) \rightarrow \Lambda_s(\tau)$ as $\varepsilon \rightarrow 0$. Finally, since $\pi_s(\tau) = \{\theta_{\rho(s)s}\}$,
we have $x_r^+ \to \theta_{\rho(s)}$, so that the sth component of $x^+$ satisfies $x_s^+ = \theta_{\rho(s)}$. These arguments imply that taking the limit of the expression

$$\lim_{\varepsilon \to 0^+} -\gamma_s x_s^+ + \Lambda_s(x^+; \varepsilon) = 0$$

results in

$$-\gamma_s \theta_{\rho(s)} + \Lambda_s(\tau) = 0.$$  

But this contradicts the fact that $Z$ is a regular parameter. This proves (1).

To prove (2) we consider regular and singular directions separately. For a regular direction, $r$, let $x_r^+ = \lim_{\varepsilon \to 0} x_r^+$. We have

$$\lim_{\varepsilon \to 0^+} -\gamma_r x_r^+ + \Lambda_r(x^+; \varepsilon) = -\gamma_r x_r^+ + \Lambda_r(\tau) = 0.$$  

Since $x_r^+ \in (\theta_{a_r}, \theta_{b_r})$, we must have $-\gamma_r \theta_{a_r} + \Lambda_r(\tau) > 0$ and $-\gamma_r \theta_{b_r} + \Lambda_r(\tau) < 0$ Therefore, $\Phi_r(\tau) = 0$.

For a singular direction, $s$, let $\eta = \min \{ \frac{\theta_{s(x)} - \theta_{s(x)}}{\varepsilon}, \frac{\theta_{s(x)} - \theta_{s(x)}}{\varepsilon} \}$. Note that $\eta$ is independent of $\varepsilon$. With this $\eta$, define the sample points $x^{\pm \eta; \varepsilon}$ by

$$x^{\pm \eta; \varepsilon}_j = \begin{cases} x^+_j, & j \neq s \\ x^+_s \pm \eta, & j = s. \end{cases}$$

Note that $x^{\pm \eta; \varepsilon} \to \tau_s^\pm$ as $\varepsilon \to 0^+$. Using the fact that $x^+$ is an equilibrium of the sigmoid system $S(Z, \varepsilon)$, we have

$$0 = -\gamma_{\rho(s)} x^+_s + \Lambda_{\rho(s)}(x^+; \varepsilon) = -\gamma_{\rho(s)} x^+_s + \Lambda_{\rho(s)}(x^+; \varepsilon) - \Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon) + \Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon)$$

and

$$= -\gamma_{\rho(s)} x^+_s + \left( \prod_{x=1}^p \sum_{j \in I_x} \sigma_{\rho(s)}(x^+_j; \varepsilon) \right) (\sigma_{\rho(s)}(x^+_s; \varepsilon) - \sigma_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon) + \Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon).$$

Rearranging the resulting equality, we get

$$\text{sgn}(\gamma_{\rho(s)} x^+_s + \Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon)) = -\text{sgn} \left( (\sigma_{\rho(s)}(x^+_s; \varepsilon) - \sigma_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon)) \right) \tag{6.1}$$

The left hand side of (6.1) is $\mathcal{L}(\tau, s, \pm)$ in the limit of $\varepsilon \to 0$. To see this notice that since $\rho$ is a permutation, there is no $j \in \text{sd}(\tau^+_s)$ so that $\rho'(j) = \rho'(s)$. By Lemma 5.2, for every $x \in \mathcal{N}(\tau^+_s)$, the function values $\Lambda_{\rho(s)}(x; \varepsilon)$ converge to $\Lambda_{\rho(s)}(\tau^+_s)$ as $\varepsilon \to 0^+$. Since in the same limit $x^{\pm \eta; \varepsilon} \to \tau^+_s$, there is an $A > 0$ such that for $\varepsilon < A$, we have $x^{\pm \eta; \varepsilon} \in \mathcal{N}(\tau^+_s)$. Therefore, $\Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon) \to \Lambda_{\rho(s)}(\tau^+_s)$ as $\varepsilon \to 0^+$. Because the projection $\pi_{\rho(s)}(\tau) = \{ \theta_{\rho'(s)(\rho(s))} \}$ we conclude that $x_{\rho(s)}^+ \to \theta_{\rho'(s)(\rho(s))}$. To summarize, we have shown

$$\lim_{\varepsilon \to 0^+} \text{sgn}(\gamma_{\rho(s)} x^+_s + \Lambda_{\rho(s)}(x^{\pm \eta; \varepsilon}; \varepsilon)) = \text{sgn}(\gamma_{\rho(s)} \theta_{\rho'(s)(\rho(s))} + \Lambda_{\rho(s)}(\tau^+_s)) \mathcal{L}(\tau, s, \pm).$$  

To show that $\mathcal{L}(\tau, s, +) = -\mathcal{L}(\tau, s, -)$, we study the right hand side of (6.1). By Properties (1) and (4) of sigmoidal perturbations, there is a neighborhood $\mathcal{U} \subset \mathbb{R}_+$ of $\theta_{\rho(s)}$ such that $\sigma_{\rho(s)}(\cdot; \varepsilon)$
is monotone increasing on $\mathcal{U}$ when $\mathbf{s}_{\rho(s)} = 1$ and decreasing when $\mathbf{s}_{\rho(s)} = -1$. Since $\sigma_{\rho(s)}$ is monotone non-increasing or non-decreasing according to $\mathbf{s}_{\rho(s)}$, this implies

$$-\text{sgn}\left((\sigma_{\rho(s)}(x_s^\varepsilon;\varepsilon) - \sigma_{\rho(s)}(x_s^{\pm\eta;\varepsilon};\varepsilon)) = -\text{sgn}\left(x_s^\varepsilon - x_s^{\pm\eta;\varepsilon}\right)\mathbf{s}_{\rho(s)}\right)
$$

$$= -(\mp)\mathbf{s}_{\rho(s)} = \pm\mathbf{s}_{\rho(s)}.$$

We have shown

$$\mathcal{L}(\tau, s, \pm) = \lim_{\varepsilon \to 0^+} \text{sgn}(\gamma_{\rho(s)}x_{\rho(s)} + \Lambda_{\rho(s)}(x_s^{\pm\eta;\varepsilon};\varepsilon))
$$

$$= \lim_{\varepsilon \to 0^+} -\text{sgn}\left((\sigma_{\rho(s)}(x_s^\varepsilon;\varepsilon) - \sigma_{\rho(s)}(x_s^{\pm\eta;\varepsilon};\varepsilon))\right) = \pm\mathbf{s}_{\rho(s)}$$

where the first equality follows from (6.2) and the second equality follows from (6.1). Therefore $\mathcal{L}(\tau, s, +) = -\mathcal{L}(\tau, s, -)$ so that $\Phi_s(\tau) = 0$. This finishes the proof of (2) and thus the forward implication in statement (a). We now proceed with the backward implication in (a).

For a singular direction $s \in \text{sd}(\tau)$, let $U_{2,s}(\varepsilon)$ be the neighborhood of $\theta_{\rho(s)}$ defined for $\sigma_{\rho(s)}$ in property 4 of sigmoidal perturbations (Definition 2.2). For a regular direction $r \notin \text{sd}(\tau)$, we have $\Phi_r(\tau) = 0$ or

$$\text{sgn}(\gamma_{r}(\theta_{a_{\rho},r} + \Lambda_{r}(\tau))) = -\text{sgn}(\gamma_{r}(\theta_{b_{\rho},r} + \Lambda_{r}(\tau))).$$

Since $\Lambda_r(x;\varepsilon) \to \Lambda_r(\tau)$ for $x \in \tau$, we may choose $\eta_r > 0$ small enough so that

$$\text{sgn}(\gamma_{r}(\theta_{a_{\rho},r} + \eta_r) + \Lambda_r(\tau)) = -\text{sgn}(\gamma_{r}(\theta_{b_{\rho},r} - \eta_r) + \Lambda_r(\tau)).$$

We now define a closed neighborhood $\tau(\varepsilon)$ of $\tau$ by

$$\pi_j(\tau(\varepsilon)) := \begin{cases} U_{2,j}(\varepsilon), & j \in \text{sd}(\tau) \\ \theta_{a_{\rho},r} + \eta_r, \theta_{b_{\rho},r} - \eta_r, & j \notin \text{sd}(\tau). \end{cases}$$

Note that $\tau \subset \tau(\varepsilon)$. By properties 3 and 4 of sigmoidal perturbations we may choose $A \in \mathbf{R}^{N \times N}$ so that for $\varepsilon < A$ and $x \in \tau(\varepsilon)$

$$\left|\frac{\partial}{\partial x_s} \Lambda_{\rho(s)}(x;\varepsilon)\right| \geq 2\gamma_{\rho(s)}, \quad \text{and} \quad \left|\frac{\partial}{\partial x_r} \Lambda_r(x;\varepsilon)\right| \leq \frac{1}{2}\gamma_r.$$ (6.4)

Property 4 implies that the image of $\Lambda_{\rho(s)}(\cdot;\varepsilon)$ on $\tau(\varepsilon)$ converges, as $\varepsilon \to 0$, to the interval with endpoints $\Lambda_{\rho(s)}(\tau_s^-)$ and $\Lambda_{\rho(s)}(\tau_s^+)$, i.e. $\Lambda_{\rho(s)}(\tau(\varepsilon);\varepsilon) \to (\Lambda_{\rho(s)}(\tau_s^-), \Lambda_{\rho(s)}(\tau_s^+))$. Here we have assumed without loss of generality that $\Lambda_{\rho(s)}(\tau_s^-) < \Lambda_{\rho(s)}(\tau_s^+)$. Since $\Phi_s(\tau) = 0$, we have $\gamma_{\rho(s)}(\theta_{b_{\rho},s}(\rho(s)) \in (\Lambda_{\rho(s)}(\tau_s^-), \Lambda_{\rho(s)}(\tau_s^+))$. Therefore, we may further refine our choice of $A$ so that $\gamma_{\rho(s)}(\theta_{b_{\rho},s}(\rho(s)) \in \Lambda_{\rho(s)}(\tau(\varepsilon);\varepsilon)$ for each $\varepsilon < A$. That is, for each $s \in \text{sd}(\tau)$, there is an $x \in \tau(\varepsilon)$ so that

$$f_{\rho(s)} := -\gamma_{\rho(s)}x_{\rho(s)} + \Lambda_{\rho(s)}(x;\varepsilon) = \dot{x}_{\rho(s)}|_{x} = 0.$$

For a regular direction $r$, we also have that there is an $x \in \tau(\varepsilon)$ with

$$f_r := -\gamma_r x_r + \Lambda_r(x;\varepsilon) = \dot{x}_r|_{x} = 0$$

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by (6.3).

By (6.4), if \( x \in \tau(\varepsilon) \) and \( s \in \text{sd}(\tau) \) is a singular direction or \( r \) is a regular direction then
\[
\left| \frac{\partial f_{\rho(s)}(x)}{\partial x_s} \right| > \gamma_{\rho(s)} > 0 \quad \text{and} \quad \frac{\partial f_{r}}{\partial x_r}(x) < -\frac{1}{2}\gamma_r < 0.
\]

In either case, the derivatives are bounded away from zero so we may apply the implicit function theorem to get a function
\[
X_j : \prod_{i \neq j} \pi_i(\tau(\varepsilon)) \to \pi_j(\tau(\varepsilon))
\]
so that
\[
f_{\rho(j)}((x_1, \ldots, x_j(x_1, \ldots, x_{j-1}, x_{j+1}, \ldots, x_N), \ldots, x_N)) = 0.
\]

Define \( g : \tau(\varepsilon) \to \tau(\varepsilon) \) by
\[
g(x_1, \ldots, x_N) := (X_1(x_2, \ldots, x_N), \ldots, X_N(x_1, \ldots, x_{N-1})).
\]

By Brouwer’s fixed point theorem, \( g \) has a fixed point \( x \) in \( \tau(\varepsilon) \) so that we can simultaneously solve \( \dot{x}_j = 0 \) for all \( j \).

To prove part (b) and show that the fixed point is unique, we apply the inverse function theorem.

Let \( x^\varepsilon \) be an equilibrium of \( S(Z, \varepsilon) \). We note that except for \( O(\varepsilon) \) terms, the Jacobian \( J(\varepsilon) \) has a block diagonal structure with blocks \( J_d(\varepsilon) \) corresponding to the cyclic feedback decomposition of \( \text{SWITCH}(Z) \) at \( \tau \) (see Section 5.1). Each block corresponding to singular directions of \( \tau \) has the form of (4.1) while the block corresponding to regular directions is diagonal. Therefore the determinant is given by
\[
\det(J(x; \varepsilon)) = \prod_{d=1}^{n} (-1)^{\ell_d} \left( \prod_{\alpha = s_d + 1}^{s_d + \ell_d} \gamma_\alpha - \text{sgn}(c_d)M_d(x, \varepsilon) \right) \left( \prod_{r = s_n + 1}^{N} -\gamma_r \right) + O(\varepsilon)
\]
where \( M_d(x, \varepsilon) = \prod_{s = s_d + 1}^{s_d + \ell_d} |\sigma_{\rho(s)}'(x_s, \varepsilon)| \). Since \( M_d(x, \varepsilon) \to \infty \) for \( x \in \tau \) we have that for small enough \( \varepsilon \), \( J(x; \varepsilon) \) is non-singular on \( \tau(\varepsilon) \). The inverse function theorem then implies that the equilibrium \( x^\varepsilon \) of \( S(Z, \varepsilon) \) in \( \tau(\varepsilon) \) is unique.

7. Proof of Cyclic Feedback System Results. In this section, we prove the results of Section 4. The section is organized into subsections which correspond to the subsections of Section 4. Recall that the edges of a cyclic feedback network (CFN) are of the form \((j, j + 1)\) where \( j + 1 \) is computed modulo \( N \) (see Definition 4.1). Furthermore, by changing variables we may assume without loss of generality that all edges of a positive CFN are activating and the edges of a negative CFN are all activating except for the edge \( N + 1 \) which is repressing (see Section 4.1).

7.1. Equilibrium Cells.

Proof of Lemma 4.3. Let \( \kappa' \) be a regular equilibrium cell. We will show that \( \kappa' = \kappa \) where \( \kappa \) is the claimed unique equilibrium cell defined in the statement of the lemma. Recall that for each \( j \in V \),
\( \theta_{-\infty} = 0 \) and \( \theta_{\infty} = \infty \) (see Definition 3.1). We have First note that

\[
\text{sgn}(-\gamma_j \theta_{-\infty} + \Lambda_j(\kappa')) = \text{sgn}(0 + \Lambda_j(\kappa')) = 1, \quad \text{and}
\text{sgn}(-\gamma_j \theta_{\infty} + \Lambda_j(\kappa')) = \text{sgn}(-\infty + \Lambda_j(\kappa')) = -1.
\]

If \( j \) is inessential and \( \gamma_j \theta(j+1)_j < L_{j(j-1)} \) then

\[
\text{sgn}(-\gamma_j \theta(j+1)_j + \Lambda_j(\kappa')) = -1.
\]

If \( j \) is inessential and \( U_{j(j-1)} < \gamma_j \theta(j+1)_j \) then

\[
\text{sgn}(-\gamma_j \theta(j+1)_j + \Lambda_j(\kappa')) = 1.
\]

In either case, \( \Phi_j(\kappa') = 0 \) if and only if \( \pi_j(\kappa') = \pi_j(\kappa) \).

Now suppose that \( k \) is essential and \( k + 1 \) is inessential. If \( \pi_k(\kappa') = (0, \theta_{(k+1)}k) \), then, using the assumption on the sign of the edges,

\[
\Lambda_{k+1}(\kappa') = \begin{cases} 
  L_{(k+1)k}, & \text{sgn}(\rho) = 1 \text{ or } k < N \\
  U_{(k+1)k}, & \text{sgn}(\rho) = -1 \text{ and } k = N
\end{cases}
\]

This implies

\[
\text{sgn}(-\gamma_{k+1} \theta_{(k+2)(k+1)} + \Lambda_{k+1}(\kappa')) = \begin{cases} 
  -1, & \text{sgn}(\rho) = 1 \text{ or } k < N \\
  1, & \text{sgn}(\rho) = -1 \text{ and } k = N
\end{cases}
\]

so that \( \Phi_{k+1}(\kappa') = 0 \) if and only if \( \pi_{k+1}(\kappa') = \pi_{k+1}(\kappa) \). A similar argument shows \( \pi_{k+1}(\kappa') = \pi_{k+1}(\kappa) \) when \( \pi_k(\kappa') = (\theta_{(k+1)}k, \infty) \). An induction argument then shows that for every essential node \( j, \pi_j(\kappa') = \pi_j(\kappa) \).

We have shown that the only regular equilibrium cell is \( \kappa \). Since equilibrium cells are a subset of loop characteristic cells, to complete the proof we need only show that the only singular loop characteristic cell \( \tau = \prod \{\theta_{(j+1)}j\} \) is not an equilibrium cell. Let \( j \) be a node so that \( j + 1 \) is inessential. Assume \( U_{(j+1)j} < \gamma_{j+1} \theta(j+2)(j+1) \). Then

\[
\mathcal{L}(\tau, j, +) = \mathcal{L}(\tau, j, -) = -1
\]

since \( \text{sgn}(-\gamma_{j+1} \theta_{(j+2)(j+1)} + \Lambda_{j+1}(\tau)) = -1 \). Similarly,

\[
\mathcal{L}(\tau, j, +) = \mathcal{L}(\tau, j, -) = 1
\]

when \( \gamma_{j+1} \theta_{(j+2)(j+1)} < L_{(j+1)j} \). This shows \( \Phi_j(\tau) \neq 0 \) so that \( \tau \) is not an equilibrium cell. \( \square \)

**Proof of Lemma 4.4.** Let \( RN \) be a cyclic feedback network and \( Z \) be a switching parameter so that the corresponding cyclic feedback system \( \text{SWITCH}(Z) \) has no inessential nodes. First we show that the singular loop characteristic cell, \( \tau = \prod \{\theta_{(j+1)}j\} \), is an equilibrium cell. Using the assumption on the sign of the edges, for every \( j \in V \) we have

\[
\Lambda_{j+1}(\tau_j^+) = \begin{cases} 
  L_{(j+1)j}, & \text{sgn}(\rho) = 1 \text{ or } j < N \\
  U_{(j+1)j}, & \text{sgn}(\rho) = -1 \text{ and } j = N
\end{cases}
\]

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\[
\Lambda_{j+1}(\tau^j_j) = \begin{cases} 
U_{(j+1)j}, & \text{sgn}(\rho) = 1 \text{ or } j < N \\
L_{(j+1)j}, & \text{sgn}(\rho) = -1 \text{ and } j = N.
\end{cases}
\]

Since \(L_{(j+1)j} < \gamma_j + 1\theta_{(j+2)(j+1)} < U_{(j+1)j}\), this implies \(\Phi_j(\tau) = 0\). \(\tau\) is therefore an equilibrium cell by Theorem 3.13.

Now consider the case \(\mathbf{RN}\) is a positive CFN. Since \(\pi_j(\kappa^L) = (0, \theta_{(j+1)j})\) and all edges are activating, \(\Lambda_{j+1}(\kappa^L) = L_{(j+1)j}\). We therefore have

\[
\mathcal{L}(\kappa^L, j + 1, +) = \text{sgn}(-\gamma_j + 1\theta_{(j+2)(j+1)} + \Lambda_{j+1}(\kappa^L)) = -1.
\]

A similar argument shows \(\mathcal{L}(\kappa^H, j + 1, -) = 1\) for all \(j\). Since

\[
\text{sgn}(-\gamma_j + 1\theta_{-\infty} + \Lambda_{j+1}(\kappa)) = 1 \quad \text{and} \quad \text{sgn}(-\gamma_j + 1\theta_{\infty} + \Lambda_{j+1}(\kappa)) = -1
\]

for all cells \(\kappa\), this implies \(\Phi_{j+1}(\kappa^L) = \Phi_{j+1}(\kappa^H) = 0\) so that both cells are equilibrium cells.

Let \(\kappa\) be a regular cell different from \(\kappa^H\) and \(\kappa^L\). Then there is a \(j\) so that \(\pi_j(\kappa) = (\theta_{(j+1)j}, \infty)\) but \(\pi_{j+1}(\kappa) = (0, \theta_{(j+2)(j+1)})\). \(\pi_j(\kappa) = (\theta_{(j+1)j}, \infty)\) implies \(\Lambda_{j+1}(\kappa) = U_{(j+1)j}\). But \(\pi_{j+1}(\kappa) = (0, \theta_{(j+2)(j+1)})\) implies

\[
\mathcal{L}(\kappa, j + 1, +) = \text{sgn}(-\gamma_{j+1}\theta_{(j+2)(j+1)} + U_{(j+1)j}) = 1
\]

so that \(\Phi_j(\kappa) = 1\) and \(\kappa\) is not an equilibrium cell.

Finally, consider the case that \(\mathbf{RN}\) is a negative CFN. Let \(\kappa\) be a regular cell. Suppose there is a \(j < N\) such that

\[
\pi_j(\kappa) = (0, \theta_{(j+1)j}) \quad \text{and} \quad \pi_{j+1}(\kappa) = (\theta_{(j+2)(j+1)}, \infty).
\]

Then \(\Lambda_{j+1} = L_{(j+1)j}\) so that

\[
\mathcal{L}(\kappa, j + 1, -) = \text{sgn}(-\gamma_{j+1}\theta_{(j+2)(j+1)} + L_{(j+1)j}) = -1.
\]

This implies \(\Phi_{j+1}(\kappa) = -1\) so that \(\kappa\) is not an equilibrium cell. Similarly, if

\[
\pi_j(\kappa) = (\theta_{(j+1)j}, \infty) \quad \text{and} \quad \pi_{j+1}(\kappa) = (0, \theta_{(j+2)(j+1)}),
\]

then \(\Phi_{j+1}(\kappa) = 1\) and \(\kappa\) is not an equilibrium cell. Suppose that \(\pi_j(\kappa) = (0, \theta_{(j+1)j})\) for each \(j\). Then since \(N+1\) is repressing,

\[
\mathcal{L}(\kappa, 1, +) = \text{sgn}(-\gamma_1\theta_{21} + U_{1N}) = 1
\]

and \(\Phi_1(\kappa) = 1\). Similarly, \(\Phi_1(\kappa) = -1\) if \(\pi_j(\kappa) = (\theta_{(j+1)j}, \infty)\) for each \(j\). Therefore there are no regular equilibrium cells. Since \(\tau\) is the only singular loop characteristic cell and equilibrium cells are loop characteristic cells, \(\tau\) is the unique equilibrium cell. \(\square\)
7.2. Stability of Equilibria. We begin by providing the calculation of the characteristic polynomial of the Jacobian $J(x; \varepsilon)$. The Jacobian is given in (4.1).

**Proof of Lemma 4.6.** Let $\tilde{J} = J(\varepsilon) - \lambda I$. Let $S_N$ be the set of permutations of order $N$. For $\eta \in S_N$ let $\text{par}(\eta)$ denote the parity of $\eta$, i.e. $\text{par}(\eta) = 1$ if $\eta$ is even and $\text{par}(\eta) = -1$ if $\eta$ is odd. The Liebniz Formula gives

\[
\det(\tilde{J}) = \sum_{\eta \in S_N} \text{par}(\eta) \prod_{i=1}^{N} \tilde{J}_{\eta(i)i}.
\]

The only non-zero derivative of $\Lambda_i$ is the derivative with respect to $x_{i-1}$. Therefore, the only non-zero entries of $\tilde{J}$ are the diagonal entries $J_{ii} = -\gamma_i - \lambda$ and the entries

\[
\tilde{J}_{i(i-1)} = \frac{\partial}{\partial x_{i-1}} \Lambda_i = \sigma'_{i(i-1)}.
\]

The only non-zero entries in the sum then correspond to the identity permutation, id and the permutation $\rho$. Since $\text{par}(\text{id}) = 1$ and $\text{par}(\rho) = (-1)^{N-1}$,

\[
\det(\tilde{J}) = \prod_{i=1}^{N} (-\gamma_i - \lambda) + (-1)^{N-1} \prod_{i=1}^{N} \sigma'_{i(i-1)}
\]

\[
= (-1)^{N} \left( \prod_{i=1}^{N} (\gamma_i + \lambda) - \prod_{i=1}^{N} \sigma'_{i(i-1)} \right)
\]

\[
= (-1)^{N} \left( \prod_{i=1}^{N} (\gamma_i + \lambda) - \text{sgn}(\rho) \prod_{i=1}^{N} M \right)
\]

To determine stability of positive cyclic feedback systems, we apply Descartes’ rule of signs and the fact that $J$ is a Metzler matrix (off diagonal entries are non-negative) so that the eigenvalue with largest real part is real (see Theorem 4 of [1]).

**Proof of Proposition 4.7.** Let $p(\lambda) := (-1)^{N} \det(J - \lambda I)$ be the characteristic polynomial of $J(x, \varepsilon)$ normalized so that the leading coefficient is positive. Notice that $p(\lambda)$ has all positive coefficients except for possibly the coefficient of $\lambda^0$ which is given by $(-1)^{N} \det(J) = \prod_{j} \gamma_j - M(x, \varepsilon)$. If $M(x, \varepsilon) > \prod_{j} \gamma_j$ then by Descartes’ rule of signs, $p$ has a positive real root so that $x$ is unstable. If $M(x, \varepsilon) < \prod_{j} \gamma_j$, then by Descartes’ rule of signs, $M(x, \varepsilon)$ has no positive real roots. Theorem 4 of [1] says that $J$ has a real eigenvalue with largest real part. Since $J$ has no positive real roots, this eigenvalue must be negative, implying that $x$ is asymptotically stable. If $M(x, \varepsilon) = \prod_{j} \gamma_j$ then $\det(J) = 0$ so that $S(Z, \varepsilon)$ has a bifurcation at $x$.

The stability of equilibria of negative cyclic feedback systems when $N \leq 2$ involves only a simple computation.

**Proof of Proposition 4.9.** If $N = 1$, then $x = x_1$ and $\dot{x} = -\gamma_1 x + \sigma_{11}(x; \varepsilon)$. Since $\sigma_{11}(x; \varepsilon)$ is non-increasing we have $J(x; \varepsilon) = -\gamma_1 + \sigma'_{11}(x; \varepsilon) \leq -\gamma_1$ so that the equilibrium of $S(Z, \varepsilon)$ is stable. If $N = 2$, then $\det(J(x; \varepsilon)) = \gamma_1 \gamma_2 - \sigma'_{21}(x; \varepsilon) \sigma'_{12}(x; \varepsilon) \geq \gamma_1 \gamma_2 > 0$ since $\sigma_{12}$ is non-decreasing and $\sigma_{11}$ is non-increasing. We also have that the trace of $J(x; \varepsilon)$ is negative. Since $\det(J(x; \varepsilon)) > 0$ and $\text{Tr}(J) < 0$, the eigenvalues of $J$ have negative real part.
Proof of Propositions 4.8 and 4.12. In proving Theorem 3.13, we showed that for $\varepsilon$ small enough the equilibrium $x^\varepsilon$ which converges to $\tau$ satisfies $x^\varepsilon \in \mathcal{U}_j(\varepsilon)$ where $\mathcal{U}_j$ is defined as in property (4) of Definition 2.2. Therefore $M(x, \varepsilon) \to \infty$ as $\varepsilon \to 0$. Applying Proposition 4.7 or 4.11 as appropriate then proves the propositions.

8. Discussion. In this paper we present explicit and direct correspondence between equilibria of systems of differential equations with sigmoidal nonlinearities and equilibrium objects that are associated to a switching system. ODE models associated to switching systems are not well defined for points that lie on the family of thresholds associated to these functions. Because of the difficulties that this presents for construction of a well defined flow, we prefer to think of a switching system not as an ODE model, but as a source of combinatorial (i.e. finite) data that can be used to study sigmoidal systems. Following this philosophy we build upon work of others [28, 19] to show that all equilibria and their stability for sufficiently steep sigmoidal functions can be determined from the data associated to the corresponding switching system. Dissecting further this rigidity, the sufficient data consists of network structure and a discrete description of parameter regime in the terms of a set of monotone Boolean functions [3]. To facilitate this work, we realize that the dynamics in a neighborhood of loop characteristic cells, that contain so called singular equilibria of the switching system, can be fully understood as a product of cyclic feedback networks. These, in turn, have a simpler structure that can be fully analyzed.

There are several natural extensions of the present work. One set of questions involves asking how far the switching system can be perturbed while maintaining its predictions. Given a switching parameter, how steep must the sigmoidal functions be so that the equilibria given by the switching system data are maintained in the sigmoidal system? How can the switching parameter be chosen so that the equilibria are maintained for the shallowest possible sigmoidal functions? We are currently working on this problem in the context of ramp systems, wherein the sigmoidal functions are replaced by continuous piece-wise linear functions. In this setting, explicit analytic results to these questions can be given.

Another set of questions involves non-stationary dynamics. The switching system dynamics can be represented by a state transition graph that has information not only about the equilibria, but also about recurrent and global dynamics of sigmoidal systems. What is the relationship between recurrent dynamics, say periodic trajectories, in the state transition graph, and periodic orbits and their stability in sigmoidal systems? We have already shown that the global dynamics of the STG is closely related to global dynamics of sigmoidal perturbations in two dimensional systems [13]. We are currently working on a generalization of this result to higher dimensions. We believe that the results of this paper present only a first step in establishing a firm connection between dynamics of sigmoidal models of network dynamics and combinatorial dynamics of state transition graphs of switching systems.

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