Common Complexes of Decompositions and Complex Balanced Equilibria of Chemical Reaction Networks

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Abstract

A decomposition of a chemical reaction network (CRN) is produced by partitioning its set of reactions. The partition induces networks, called subnetworks, that are “smaller” than the given CRN which, at this point, can be called parent network. A complex is called a common complex if it occurs in at least two subnetworks in a decomposition. A decomposition is said to be incidence independent if the image of the incidence map of the parent network is the direct sum of the images of the subnetworks’ incidence maps. It has been recently discovered that the complex balanced equilibria of the parent network and its subnetworks are fundamentally connected in an incidence independent decomposition. In this paper, we utilized the set of common complexes and a developed criterion to investigate decomposition’s incidence independence properties. A framework was also developed to analyze decomposition classes with similar structure and incidence independence properties. We identified decomposition classes that can be characterized by their sets of common complexes and studied their incidence independence. Some of these decomposition classes occur in some biological and chemical models. Finally, a sufficient condition was obtained for the complex balancing of some power law kinetic (PLK) systems with incidence independent and complex balanced decompositions. This condition led to a generalization of the Deficiency Zero Theorem for some PLK systems.
1 Introduction

Decompositions of chemical reaction networks (CRNs) were first studied by Feinberg in his 1987 review [8]. A decomposition is generated by a partition of the set of reactions, with the species and complexes of a subnetwork being those occurring in the corresponding reaction subset. He also introduced the concept of an independent decomposition where the stoichiometric subspace of the parent network is the direct sum of the stoichiometric subspaces of the decomposition’s subnetworks. He showed its importance by stating the following relationship among the positive equilibria of the network and those of the decomposition’s subnetworks:

**Theorem 1** (Remark 5.4, [8]). Let $(\mathcal{N}, K)$ be a chemical kinetic system with partition $\{\mathcal{R}_1, \ldots, \mathcal{R}_k\}$. If $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ is the network decomposition generated by the partition and $E_+(\mathcal{N}_i, K_i) = \{x \in \mathbb{R}_{>0}^q \mid N_i K_i(x) = 0\}$, then

i. $\bigcap_{i=1}^k E_+(\mathcal{N}_i, K_i) \subseteq E_+(\mathcal{N}, K)$ and

ii. the equality holds if the network decomposition is independent.

Feinberg’s early result has turned out to be particularly useful in the new field of concentration robustness of kinetic systems, which he and Shinar initiated in a well-known paper in Science in 2010 [33]. A sufficient condition for deficiency one mass action systems they derived has, in the meantime, been extended to power law and Hill type kinetic systems ([16], [20]). Independent decompositions allow the detection of concentration robustness in larger and higher deficiency kinetic system via Feinberg’s theorem ([11], [12], [15], [23]).

In view of the focus of the chemical reaction network theory (CRNT) literature on mass action systems, where the subset of complex balanced equilibria is either empty or coincides with the whole set of positive equilibria, the corresponding concept and result were formulated only recently by Farinas et al. [6]: a decomposition is incidence independent if the image of the incidence map of the parent network is the direct sum of the images of the incidence maps of the subnetworks. The relationship among the complex balanced equilibria sets is fully analogous to Feinberg’s result:

**Theorem 2** (Theorem 4, [6]). Let $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ be a decomposition. Let $K$ be any kinetics, and $Z_+(\mathcal{N}, K)$ and $Z_+(\mathcal{N}_i, K_i)$ be the set of complex balanced equilibria of $\mathcal{N}$ and $\mathcal{N}_i$, respectively. Then,

$$\bigcap_{i=1}^k Z_+(\mathcal{N}_i, K_i) \subseteq Z_+(\mathcal{N}, K).$$

If the decomposition is incidence independent, then

i. $\bigcap_{i=1}^k Z_+(\mathcal{N}_i, K_i) = Z_+(\mathcal{N}, K)$ and

ii. $Z_+(\mathcal{N}, K) \neq \emptyset$ implies that $Z_+(\mathcal{N}_i, K_i) \neq \emptyset$ for $i = 1, \cdots, k$. 
The best-known example of an incidence independent decomposition is the set of
linkage classes of a network. It also has the special property that it partitions not only
the set of reactions but also the set of complexes of a network. Farinas et al. introduced
the term $C$-decomposition for a decomposition with the latter property and show that
any $C$-decomposition of a network is generated by a coarsening of the partition into
reactions of linkage classes. Inspired by the study of a model of the gene regulatory
system of *Mycobacterium tuberculosis* [7], Farinas et al. also introduced the concept of a
$C^*$-decomposition, where the subnetworks pairwise have at most the zero complex. They
also derived that any $C^*$-decomposition is incidence independent.

These results motivated the systematic study of the set of common complexes of a
decomposition, denoted by $C_\mathcal{D}$ (the subscript $\mathcal{D}$ is the symbol assigned to the decompo-
sition), and its relationships to incidence independence and complex balanced equilibria,
which is the main topic of this paper. A common complex is a complex that occurs
in at least two subnetworks in the decomposition. Clearly, $C$-decompositions are those
with $C_\mathcal{D} = \emptyset$ (or equivalently $|C_\mathcal{D}| = 0$), and $C^*$-decompositions belong to those with
$|C_\mathcal{D}| \leq 1$. For decompositions with $|C_\mathcal{D}| \geq 2$, incidence independence may not hold, and a
criterion based on counting common complexes is established. Furthermore, a framework
for describing decomposition classes based on the structure of $C_\mathcal{D}$ is developed for formu-
lating conditions for this property to hold for decomposition classes. The framework is
illustrated in detail by the class of pairwise min-max (PMM) decompositions, which form
a superset of the $C^*$-decompositions.

Although the theorems of Feinberg and Farinas et al. determine the structure of
the sets of positive equilibria and complex balanced equilibria through independent and
incidence independent decompositions of kinetic systems respectively, neither result guar-
antees that the intersections are non-empty. Specifically for an incidence independent
decomposition, given that complex balancing of the subnetworks is a necessary condition
for that of the whole system, the question is: which additional properties are needed for
it to be sufficient too?

In [6], Farinas et al. also established the surprisingly general result that the conclusion
holds for any $C$-decomposition, i.e., the additional purely structural condition of $C_\mathcal{D} = \emptyset$
and any kinetic system. We show with a simple example that this is no longer true for
$C_\mathcal{D} \neq \emptyset$. This result is singular in its purely structural character: as soon as $C_\mathcal{D} \neq \emptyset$, there
are kinetic systems with an incidence independent and complex balanced decomposition
(a decomposition where each subnetwork is complex balanced) that are not complex
balanced. In fact, mass action systems on the same network and the same incidence
independent and complex balanced decomposition, may or may not be complex balanced.
Hence, to obtain results when $C_\mathcal{D} \neq \emptyset$ for an incidence independent and complex balanced
decomposition, one needs to restrict the set of kinetics systems for which the conclusion of
complex balancing would hold. Ideally, if the latter is true for a kinetic system, it should
also hold for all kinetic systems in its semi-module (i.e., for all kinetic systems with the
same interaction function, differing only in rate constants).

Accordingly, our result, while imposing no conditions on the set of common complexes
of the incidence independent decomposition, is valid only for a class of power law kinetic
(PLK) systems. The class is defined by two properties:

- the subnetworks of the decomposition have power law reactant-determined kinetics
  (PL-RDK); and

- the induced decomposition of kinetic complexes is independent.
The additional conditions involve properties of the kinetics on the decomposition subnetworks and associated structures.

This result further underscores the importance of incidence independence for the complex balancing of the whole network. The result is analogous to a generalization of the Deficiency Zero Theorem (DZT) of Fortun et al. [14] for the set of all positive equilibria presented by Hernandez et al. [18].

The main results of the paper include:

i. a common complexes approach to study decomposition’s properties, in particular incidence independence through a developed criterion;

ii. a framework for analyzing decomposition classes with similar structure and incidence independence properties illustrated in detail by the class of PMM decompositions, which contain the linkage class decomposition and occur in biological and chemical models; and

iii. a new condition ensuring the complex balancing of a weakly reversible power law kinetic system with an incidence independent decomposition into complex balanced subnetworks.

The paper is organized as follows. Fundamental concepts on chemical reaction networks, chemical kinetic systems, and decomposition theory are given in section 2. The framework to classify network’s decompositions according to equivalent incidence independence properties is presented in section 3. In sections 4 and 5, the class of pairwise binary-sized (PBS) decompositions and its subclass of PMM decompositions are discussed. The condition ensuring complex balancing of a PLK system with an incidence independent complex balanced decomposition is derived in section 6. Section 7 provides a summary and an outlook.

2 Fundamentals of chemical reaction networks and kinetic systems

In this section, some necessary concepts and results on chemical reaction networks and chemical kinetic systems are reviewed, the details of which can be found in [1], [9], and [35].

2.1 Structure of chemical reaction networks

Consider a non-empty finite set \( \mathcal{S} \) of species. A linear combination of these species with nonnegative integer coefficients (or stoichiometric coefficients) is called a complex. A complex whose stoichiometric coefficients are all zero is called the zero complex and is denoted by 0. The set of all complexes is denoted by \( \mathcal{C} \). Here, \( m \) and \( n \) are used to denote the number of species and complexes, respectively.

The set of reactions, denoted by \( \mathcal{R} \), is a non-empty subset of \( \mathcal{C} \times \mathcal{C} \) satisfying the conditions:

i. \( (y, y) \notin \mathcal{R} \) for all \( y \in \mathcal{C} \); and

ii. for each \( y \in \mathcal{C} \), there is a \( y' \in \mathcal{C} \) such that \( (y, y') \in \mathcal{R} \) or \( (y', y) \in \mathcal{R} \).
In a reaction \((y, y')\), which can also be written as \(y \rightarrow y'\), \(y\) and \(y'\) are called reactant and product complexes, respectively. The order of \(R\) is denoted here by \(p\).

**Definition 1.** A chemical reaction network (CRN) \(N\) is a triple \((\mathcal{S}, \mathcal{C}, \mathcal{R})\) of non-empty finite sets where \(\mathcal{S}\), \(\mathcal{C}\), and \(\mathcal{R}\) are the sets of species, complexes, and reactions, respectively.

The terms “chemical reaction network”, “reaction network”, and “network” are used here interchangeably.

The reactant map \(\rho : \mathcal{R} \rightarrow \mathcal{C}\) and the product map \(\pi : \mathcal{R} \rightarrow \mathcal{C}\) map a reaction to its reactant and product complexes, respectively. The number of the reactant complexes is denoted by \(n_{\rho}\), i.e., \(|\rho(\mathcal{R})| = n_{\rho}\). A CRN is called cycle terminal if \(n_{\rho} = n\) while it is called nonbranching if \(n_{\rho} = p\).

A CRN can be viewed as a digraph with the sets of complexes and reactions serve as the digraph’s vertex and arc sets, respectively. The digraph representation of a reaction network, called reaction graph, is loopless and does not have an isolated vertex. The underlying graph of a reaction network is the graph obtained after removing the direction in each arc of its reaction graph. Paths and cycles in a CRN are the corresponding paths and cycles in the underlying graph. Directed paths and directed cycles in a reaction network are the uni-directed paths and cycles in the reaction graph.

**Example 1.** Consider a toy CRN with the following digraph representation.

\[
\begin{array}{c}
A + B \xrightarrow{R_3} A \xrightarrow{R_2} C \xleftarrow{R_4} \ 2D \\
\end{array}
\]  

This network has \(m = 4, n = 4\), and \(p = 5\) such that

\[
\mathcal{S} = \{A, B, C, D\}, \mathcal{C} = \{A + B, A, C, 2D\}, \text{ and } \mathcal{R} = \{R_1, R_2, R_3, R_4, R_5\}. \tag{2.2}
\]

The network is also cycle terminal since \(n = 4 = n_{\rho}\).

Two complexes are connected if there is a path joining the corresponding vertices in the underlying graph. A reaction network is said to be weakly connected if the underlying graph is connected, i.e., every pair of complexes are connected. Two complexes \(y\) and \(y'\) are said to be strongly connected if there is a directed path from \(y\) to \(y'\) and vice versa. The linkage classes, whose number is denoted by \(\ell\), of a network corresponds to the maximal weakly connected subgraphs of its reaction graph while the strong linkage classes are the subgraphs of the reaction graph that are maximal strongly connected. If each linkage class is also a strong linkage class, a CRN is called weakly reversible. The toy CRN in Example 1 is weakly reversible since its sole linkage class is a strong linkage class.

CRNs are often studied with the aid of finite dimensional spaces \(\mathbb{R}^{\mathcal{S}}, \mathbb{R}^{\mathcal{C}}, \text{ and } \mathbb{R}^{\mathcal{R}}\), respectively referred to as species space, complex space, and reaction space. For every reaction \(y \rightarrow y'\), a vector, called reaction vector, resulted from subtracting the reactant complex \(y\) from the product complex \(y'\) is associated. The linear subspace of \(\mathbb{R}^{\mathcal{S}}\) defined by \(S := \text{span}\{y' - y \in \mathbb{R}^{\mathcal{S}} | y \rightarrow y' \in \mathcal{R}\}\) is called the stoichiometric subspace \(S\) of a reaction network. Its dimension \(s\) also refers to the rank of the network.
CRN can be characterized by a nonnegative integer $\delta$ called deficiency which is given by $\delta = n - \ell - s$. The toy CRN in Example 1 has $S = \text{span}\{r_1, r_2, r_4\}$ where

$$r_1 = \begin{bmatrix} 0 \\ -1 \\ 0 \\ 0 \end{bmatrix}, r_2 = \begin{bmatrix} -1 \\ 0 \\ 1 \\ 0 \end{bmatrix}, \text{ and } r_4 = \begin{bmatrix} 0 \\ 0 \\ -1 \\ 2 \end{bmatrix}. \quad (2.3)$$

Hence, its rank $s = 3$ while its deficiency is $\delta = 4 - 1 - 3 = 0$, respectively.

**Definition 2.** Let $\mathcal{N} = (\mathcal{S}, \mathcal{C}, \mathcal{R})$ be a CRN. The incidence map $I_a : \mathbb{R}^\mathcal{S} \to \mathbb{R}^\mathcal{C}$ is the linear map defined by mapping for each reaction $R_i : y_i \to y'_i \in \mathcal{R}$, the basis vector $\omega_i$ to the vector $\omega_{y'_i} - \omega_{y_i} \in \mathcal{C}$.

As a linear map, the incidence map has an $n \times p$ matrix representation, called incidence matrix, whose entries are described by the following:

$$(I_a)_{(i,j)} = \begin{cases} -1 & \text{if } i \text{ is the reactant complex of reaction } j \in \mathcal{R} \\ 1 & \text{if } i \text{ is the product complex of reaction } j \in \mathcal{R} \\ 0 & \text{otherwise} \end{cases}$$

The toy CRN in Example 1 has the following incidence matrix.

$$I_a = \begin{bmatrix} R_1 & R_2 & R_3 & R_4 & R_5 \\ -1 & 0 & 1 & 0 & 0 \\ 1 & -1 & 0 & 0 & 0 \\ 0 & 1 & -1 & -1 & 1 \\ 0 & 0 & 0 & 1 & -1 \end{bmatrix} \begin{bmatrix} A + B \\ A \\ C \\ 2D \end{bmatrix} \quad (2.4)$$

Boros [2] described the rank of the incidence map through the following proposition.

**Proposition 1** (Proposition 4.2.3, [2]). $\dim \text{Im} I_a = n - \ell$.

### 2.2 Dynamics of chemical reaction networks

A kinetics is an assignment of a rate function to each reaction in a CRN. A network $\mathcal{N}$ together with a kinetics $K$ is called a chemical kinetic system (CKS) and is denoted here by $(\mathcal{N}, K)$. **Power law kinetics** (PLK) is identified by the kinetic order matrix, which is a $p \times m$ matrix $F = [F_{ij}]$, and vector $k \in \mathbb{R}^\mathcal{S}_{>0}$ called the rate vector.

**Definition 3.** A kinetics $K : \mathbb{R}^\mathcal{S}_{>0} \to \mathbb{R}^\mathcal{S}$ is a power law kinetics if

$$K_i(x) = k_i x^{F_{i,:}} \text{ for } i = 1, \ldots, r \quad (2.5)$$

where $k_i \in \mathbb{R}_{>0}$, $F_{ij} \in \mathbb{R}$, and $F_{i,:}$ is the row of $F$ associated to reaction $R_i$.

A PLK system can be classified based on the kinetic orders assigned to its branching reactions (i.e., reactions sharing a common reactant complex).

**Definition 4.** A PLK system has a reactant-determined kinetics (type PL-RDK) if for any two branching reactions $R_i, R_j \in \mathcal{R}$, the corresponding rows of kinetic orders in $F$ are identical, i.e., $F_{ih} = F_{jh}$ for $h = 1, \ldots, m$. Otherwise, a PLK system has a non-reactant-determined kinetics (type PL-NDK).
The well-known mass action kinetic (MAK) system forms a subset of PL-RDK systems. In particular, MAK is given by $K_i(x) = k_i x^{y_i}$ for all reactions $R_i : y_i \to y'_i \in R$ with $k_i \in \mathbb{R}_{>0}$ (called rate constant). The vector $Y_{ij}$ contains the stoichiometric coefficients of a reactant complex $y_i \in C$.

**Definition 5.** The species formation rate function of a chemical kinetic system is the vector field

$$f(c) = NK(c) = \sum_{y_i \to y'_i \in R} K_i(c)(y'_i - y_i)$$

where $c \in \mathbb{R}^r_{>0}$ and $N$ is the $m \times p$ matrix, called stoichiometric matrix, whose columns are the reaction vectors of the system.

The equation $dc/dt = f(c(t))$ is the ODE or dynamical system of the chemical kinetic system. An element $c^*$ of $\mathbb{R}^r_{>0}$ such that $f(c^*) = 0$ is called a positive equilibrium or steady state of the system. The set of all positive equilibria of a CKS is denoted by $E_+(N, K)$.

Analogous to the species formation rate function, we also have the complex formation rate function.

**Definition 6.** The complex formation rate function $g : \mathbb{R}^r_{>0} \to \mathbb{R}^r$ of a chemical kinetic system is the vector field

$$g(c) = I_a K(c) = \sum_{y_i \to y'_i \in R} K_i(c)(\omega_{y'_i} - \omega_{y_i})$$

where $c \in \mathbb{R}^r_{>0}$ and $I_a$ is the incidence map.

A CKS is complex balanced if it has complex balanced steady state, i.e., there is a composition $c^{**} \in \mathbb{R}^r_{>0}$ such that $g(c^{**}) = 0$. The set of all complex balanced steady states of the system is denoted by $Z_+(N, K)$.

**2.3 Decomposition theory**

Some basic concepts and results on decomposition theory are discussed in this subsection. More detailed discussions than the one contained here can be found in [6] and [15].

We state the concept of restriction on the objects of a reaction network as given in [24] and [25]. Let $N = (I, C, R)$ be a given network and consider a subset $C' \subseteq C$ and a subset $R' \subseteq R$. The restriction of $C$ to $R'$, denoted by $C|_{R'}$, is the set of (reactant and product) complexes of the reactions in $R'$. Likewise, the restriction of $I$ to $C'$, denoted by $I|_{C'}$, is the set of species that occur in the complexes in $C'$.

Let $N = (I, C, R)$ be a given network. $N' = (I', C', R')$ is said to be a subnetwork of $N$ if $R' \subseteq R$. In this notion, $C' = C|_{R'}$ and $I' = I|_{C'}$. That is, $N'$ has complex set consists of the complexes of the reactions in $R'$ and has species set consists of the species that occur in the complexes of $C|_{R'}$.

**Definition 7.** Let $N = (I, C, R)$ be a CRN. A covering of $N$ is a collection of subsets $\{R_1, \ldots, R_k\}$ whose union is $R$. A covering is called a decomposition of $N$ if $R_i$’s form a partition of $R$, i.e., $R_i$’s are pairwise disjoint.
As previously indicated, \( R_i \) defines a subnetwork \( \mathcal{N}_i \) of \( \mathcal{N} \) such that \( \mathcal{N}_i = (\mathcal{I}_i, \mathcal{C}_i, \mathcal{R}_i) \) where \( \mathcal{C}_i = \mathcal{C} \upharpoonright_{\mathcal{R}_i} \) and \( \mathcal{I}_i = \mathcal{I} \upharpoonright_{\mathcal{R}_i} \). Hence,

\[
\mathcal{R} = \bigcup \mathcal{R}_i \Rightarrow \mathcal{C} = \bigcup \mathcal{C}_i \Rightarrow \mathcal{I} = \bigcup \mathcal{I}_i.
\] (2.8)

Gross et al. [17] defined the union of networks as follows.

**Definition 8.** The **union of networks** \( \mathcal{N}_i = (\mathcal{I}_i, \mathcal{C}_i, \mathcal{R}_i), i = 1, \ldots, k \), is given by

\[
\mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k := \left( \bigcup_{i=1}^{k} \mathcal{I}_i, \bigcup_{i=1}^{k} \mathcal{C}_i, \bigcup_{i=1}^{k} \mathcal{R}_i \right)
\] (2.9)

Because of the clear connection of Definition 8 and (2.8), we adopt the notation for union of networks to indicate covering or decomposition of a network and write \( \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) to mean that \( \{\mathcal{R}_1, \ldots, \mathcal{R}_k\} \) is a covering or a decomposition of \( \mathcal{N} \).

Given a decomposition \( \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \), we may refer to \( \mathcal{N} \) as the **parent network** of its subnetworks \( \mathcal{N}_i \)'s. The number of subnetworks in a decomposition (usually denoted by \( k \)) is called the **length of the decomposition**. If the length of a decomposition is 0, it is called an **empty decomposition**. A decomposition is called **trivial** if its length is 1, i.e., it consists of only one subnetwork. Here, a decomposition is assumed to be non-empty and non-trivial unless otherwise stated. We extend the CRN notations used in this paper to fit the concept of subnetwork. The complete list of notations adapted for the subnetworks used in this paper is given in the appendix.

**Independent subnetworks**, as given in [8], are obtained from two complementary subsets of the reaction set which satisfy the condition that the stoichiometric subspace of the parent network equals the direct sum of the corresponding stoichiometric subspaces of the subnetworks. The notion of independent subnetworks was the early form of independent decomposition defined as follows.

**Definition 9.** A decomposition is **independent** if \( S \) is the direct sum of the subnetworks’ stoichiometric subspaces \( S_i \) or equivalently, \( s = s_1 + \cdots + s_k \).

**Example 2.** Let the toy CRN in Example 1 be denoted by \( \mathcal{N} \). Consider the subnetworks \( \mathcal{N}_1 \) and \( \mathcal{N}_2 \) of \( \mathcal{N} \) whose reaction graphs are given below.

\[
\begin{align*}
\mathcal{N}_1 : \quad & 2 + B \longrightarrow A \longrightarrow A + C \\
\mathcal{N}_2 : \quad & A + C \xrightarrow{\mathcal{R}} D 
\end{align*}
\] (2.10)

It is clear that \( \mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2 \) is a decomposition. Moreover, \( \mathcal{N}_1 \) and \( \mathcal{N}_2 \) have ranks 2 and 1, respectively. Since the parent network is of rank 3, the decomposition is independent.

Farinas et al. [6] introduced the concept of incidence independent decomposition that is patterned after the independent decomposition but utilizes the images of the incidence maps instead of the stoichiometric subspaces.

**Definition 10.** A decomposition \( \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) is **incidence independent** if and only if the image of the incidence map of \( \mathcal{N} \) is the direct sum of the images of the incidence maps of \( \mathcal{N}_i \)'s.

Due to Proposition 1, a decomposition is incidence independent if and only if \( n - \ell = \sum (n_i - \ell_i) \) where \( n_i \) and \( \ell_i \) are the number of complexes and linkage classes of \( \mathcal{N}_i \).
Example 3. The decomposition given in Example 2 is also incidence independent since $n - \ell = 4 - 1 = 3$ while $n_1 - \ell_1 = 3 - 1 = 2$ and $n_2 - \ell_2 = 2 - 1 = 1$.

$\mathcal{C}$- and $\mathcal{C}^*$-decompositions are decomposition classes defined in [6] which are found to be useful in several recent studies ([15], [7]).

Definition 11. A decomposition $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ with $\mathcal{N}_i = (\mathcal{I}_i, \mathcal{C}_i, \mathcal{R}_i)$ is a $\mathcal{C}$-decomposition if $\mathcal{C}_i \cap \mathcal{C}_j = \emptyset$ for $i \neq j$.

$\mathcal{C}$-decomposition is a decomposition where the set of complexes of the parent network is partitioned to become the complex sets of the subnetworks. The following structure theorem identifies $\mathcal{C}$-decomposition as a coarsening of the linkage class decomposition.

Theorem 3 (Structure Theorem for $\mathcal{C}$-decomposition, [6]). $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ is a $\mathcal{C}$-decomposition if and only if each $\mathcal{N}_i$ is the union of linkage classes and each linkage class is contained in only one $\mathcal{N}_i$.

It immediately follows from this result that, in any $\mathcal{C}$-decomposition, $\sum \ell_i = \ell$. With this result, the incidence independence of a $\mathcal{C}$-decomposition is guaranteed as stated in the following proposition.

Proposition 2 (Proposition 3, [6]). Any $\mathcal{C}$-decomposition is incidence independent.

On the other hand, $\mathcal{C}^*$-decomposition is characterized by the condition that the pair-wise intersection of the subnetworks’ complex sets can only contain the zero complex.

Definition 12. A decomposition $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ with $\mathcal{N}_i = (\mathcal{I}_i, \mathcal{C}_i, \mathcal{R}_i)$ is a $\mathcal{C}^*$-decomposition if $\mathcal{C}_i^* \cap \mathcal{C}_j^* = \emptyset$ for $i \neq j$ where $\mathcal{C}_i^*$ and $\mathcal{C}_j^*$ are the sets of non-zero complexes in $\mathcal{C}_i$ and $\mathcal{C}_j$, respectively.

Equivalently, in $\mathcal{C}^*$-decomposition, $\mathcal{C}_i \cap \mathcal{C}_j \subseteq \{0\}$. From this, it follows that $\mathcal{C}$-decompositions are contained in the class of $\mathcal{C}^*$-decompositions. An analogous structure theorem for $\mathcal{C}^*$-decomposition was also provided in [6].

Theorem 4 (Structure Theorem for $\mathcal{C}^*$-decomposition, [6]). Let $\mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ be a $\mathcal{C}^*$-decomposition and $\mathcal{L}_0$ and $\mathcal{L}_{0,i}$ be the linkage classes of $\mathcal{N}$ and $\mathcal{N}_i$ containing the zero complex (note that $\mathcal{L}_{0,i}$ is empty if $\mathcal{N}_i$ does not contain the zero complex). Then,

i. the $\mathcal{L}_{0,i}$ form a $\mathcal{C}^*$-decomposition of $\mathcal{L}_0$ and

ii. the (non-empty) $\mathcal{N}_i \setminus \mathcal{L}_{0,i}$ form a $\mathcal{C}$-decomposition of $\mathcal{N} \setminus \mathcal{L}_0$.

A criterion was formulated in [6] for the incidence independence of a $\mathcal{C}^*$-decomposition that is not $\mathcal{C}$-decomposition:

$$\sum \ell_i - \ell = k(1) - 1$$

(2.11)

where $k(1)$ is the number of subnetworks $\mathcal{N}_i$ that contain the zero complex. This criterion play an important role in proving the incidence independence of $\mathcal{C}^*$-decompositions.

Proposition 3 (Corollary 5, [6]). Any $\mathcal{C}^*$-decomposition is incidence independent.
3 The common complexes framework for decomposition classes and their incidence independence properties

In this section we introduce the concept of the set of common complexes of a decomposition, denoted by $\mathcal{C}_D$, and the associated properties which will be needed in the succeeding sections in investigating the incidence independence of decompositions.

3.1 Definition of $\mathcal{C}_D$ and some basic properties

We define the set of common complexes as follows.

**Definition 13.** Let $D$ be a decomposition. The set $\mathcal{C}_D$ of common complexes consists of all complexes contained in sets of complexes of at least two distinct subnetworks of $D$.

The order of $\mathcal{C}_D$ is denoted here by $d$, i.e., $|\mathcal{C}_D| = d$. The following proposition indicates that specifying pairwise intersections of the subnetworks' complex sets determines the set of common complexes.

**Proposition 4.** In decomposition $D$, $\mathcal{C}_D = \bigcup_{i<j}(\mathcal{C}_i \cap \mathcal{C}_j)$.

Note that for $i \neq j$, $(\mathcal{C}_i \setminus \mathcal{C}_D) \cap (\mathcal{C}_j \setminus \mathcal{C}_D) = \emptyset$. Moreover, in $\mathcal{C}$-decomposition, $\mathcal{C}_D = \emptyset$ while $\mathcal{C}_D = \{0\}$ for $\mathcal{C}^*$-decomposition that is not a $\mathcal{C}$-decomposition. Since every $\mathcal{C}$-decomposition is incidence independent, all decompositions with $d = 0$ is incidence independent. It is indicated in Proposition [3] that a decomposition with $d = 1$ is incidence independent. Here is a counterexample showing that incidence independence does not necessarily follow, in general, for $d \geq 2$.

**Example 4.** Consider the decomposition $D : \mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2$ given by the following

$\mathcal{N} : A \leftrightarrow B \quad \mathcal{N}_1 : A \rightarrow B \quad \mathcal{N}_2 : A \leftarrow B$ \hspace{1cm} (3.1)

Notice that $\mathcal{C}_D = \{A, B\}$ and $n - \ell = 2 - 1 = 1$ while $n_1 - \ell_1 = n_2 - \ell_2 = 2 - 1 = 1$. Hence, $D$ is not incidence independent.

Let $h_i$ be the number of complexes in subnetwork $\mathcal{N}_i$ that are also contained in $\mathcal{C}_D$, i.e., $h_i = |\mathcal{C}_i \cap \mathcal{C}_D|$. The following result, which generalizes (2.11), provides a criterion that determines the incidence independence of a decomposition through its $\mathcal{C}_D$. We call this connectivity-common complexes criterion (3C-criterion).

**Proposition 5.** Let $D : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ be a decomposition with $d = |\mathcal{C}_D|$. Then, the decomposition is incidence independent if and only if $\sum_{i=1}^{k} \ell_i - \ell = \sum_{i=1}^{k} h_i - d$.

**Proof.** Observe that $n = d + \sum_{i=1}^{k} (n_i - h_i)$. Then, we have
incidence independent decomposition \(\Leftrightarrow n - \ell = \sum_{i=1}^{k} (n_i - \ell_i) = \sum_{i=1}^{k} n_i - \sum_{i=1}^{k} \ell_i\)

\(\Leftrightarrow n - \ell = (d - d) + \sum_{i=1}^{k} (n_i - h_i + h_i) - \sum_{i=1}^{k} \ell_i\)

\(\Leftrightarrow n - \ell = (d - d) + \sum_{i=1}^{k} (n_i - h_i) + \sum_{i=1}^{k} h_i - \sum_{i=1}^{k} \ell_i\)

\(\Leftrightarrow \sum_{i=1}^{k} \ell_i - \ell = \sum_{i=1}^{k} h_i - d\)

Example 5. The criterion for the incidence independence of a \(\mathcal{C}^r\)-decomposition that is not \(\mathcal{C}\)-decomposition (2.11) is obtained from the 3C-criterion when \(d = 1\):

\[\sum_{i=1}^{k(1)} 1 - 1 = k(1) - 1.\]

3.2 The common complexes framework of a decomposition class

We now introduce a framework called Common Complexes Framework (CCF) that will provide a way to classify network’s decompositions into sets with equivalent incidence independence properties.

For any positive integer \(k\), an incidence independence class of decompositions of \(\mathcal{N}\) with \(k\) subnetworks is given by a set of conditions (called class conditions) defining the \(k(k-1)/2\) pairwise intersections \(\mathcal{C}_i \cap \mathcal{C}_j, i < j\), for each decomposition. The class conditions determine the range of (integer) values for \(0 \leq d < |\mathcal{C}|\). The \(k \times k\) upper triangular matrix (with zero diagonal) \(M\) given by

\[m_{ij} = \begin{cases} 0 & \text{if } i \geq j \\ |\mathcal{C}_i \cap \mathcal{C}_j| & \text{if } i < j \end{cases} \quad (3.2)\]

is called the pairwise intersection matrix (PIM).

CCF has the following three components:

i. a class condition that determines the PIM and the relevant parameters;

ii. a structure theorem for the decompositions in the class in terms of \(\mathcal{C}\)-decompositions; and

iii. a formulation of the 3C-criterion for incidence independence in terms of the class parameters.

PIM were chosen to be a matrix with elements of upper triangular nonnegative matrices which form an additive monoid and allows scalar multiplication with positive integers. These operations may have interpretable effects on the decompositions involved. The structure theorem and 3C-criterion are often useful in finding network properties that are sufficient or necessary for the incidence independence of a decomposition.
Example 6. The class of \( C \)-decompositions can be defined through the following equivalent statements

\[
C_i \cap C_j = \emptyset \text{ for } i < j \iff d = 0 \iff M = 0
\] (3.3)

where 0 is the zero matrix. The structure theorem (Theorem 3) states that every \( C \)-decomposition is a coarsening of the linkage class decomposition which led to the 3C-criterion given by \( \sum \ell_i - \ell = 0 \).

Example 7. For \( C^* \)-decomposition that is not a \( C \)-decomposition, \( d = 1 \) and the PIM is the matrix with entries 0 and 1 (for \( ij \)-entry corresponding to \( C_i \) and \( C_j \) whose intersection is the zero complex). See Theorem 4 and (2.11) for the structure theorem and 3C-criterion.

4 The class of pairwise binary-sized decompositions

In this section, we discuss a class of network decompositions characterized by the property that the intersection sets of subnetworks’ complex sets take on two sizes which are denoted by \( b \) and \( c \).

4.1 Definition and examples

Definition 14. Let \( \mathcal{D} : \mathcal{M} = \mathcal{M}_1 \cup \cdots \cup \mathcal{M}_k \) be a decomposition and \( b \) and \( c \) be integers such that \( 0 \leq b \leq c \leq d = |C_D| \). Then, \( \mathcal{D} \) is said to be pairwise binary-sized (PBS) if \( |C_i \cap C_j| = b \) or \( c \) for \( i \neq j \).

PBS decompositions for a given pair \((b, c)\) are denoted here by PBS\(_d(b, c)\). In the context of this paper, \( b \) and \( c \) are often zero and non-zero, respectively. Hence, the number \( b \) can be thought of as the “bottom number” of pairwise intersection elements while \( c \) is the common cardinality of the pairwise intersection set.

Example 8. Observe that \( C^* \)-decompositions that are not \( C \)-decompositions are PBS\(_1(0, 1)\) or PBS\(_1(1, 1)\) while \( C \)-decompositions are PBS\(_0(0, 0)\). Note that \( d = 0 \iff b = 0 \) and \( c = 0 \). Hence, \( C \)-decompositions can be denoted simply with PBS\(_0\).

We now provide several examples of PBS decompositions in models of chemical and biological systems.

Example 9. Take the following weakly reversible and deficiency zero subnetwork \( \mathcal{N} \) of the Schmitz’s pre-industrial carbon cycle model [32].

![Diagram of network](image)

(4.1)

The complexes in this network represent six carbon pools (see the table below) while the reactions describe the mass transfers. Fortun et al. [16] studied a power law representation of this model in [16].
Table 4.1. The carbon pools of the Schmitz’s pre-industrial carbon cycle model.

Now, consider the decomposition \( \mathcal{D}_1 : \mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2 \) (also given in [16]) where

\[
\begin{array}{c}
M_1 \\
M_2 \\
M_3 \\
M_4 \\
M_5 \\
M_6
\end{array}
\begin{array}{c}
\text{atmosphere} \\
\text{warm ocean surface water} \\
\text{cool ocean surface water} \\
\text{deep ocean waters} \\
\text{terrestrial biota} \\
\text{soil and detritus}
\end{array}
\]

Note that \( \mathcal{D}_1 \) is incidence independent according to Proposition 5 since \((\ell_1 + \ell_2) - \ell = (1+1) - 1 = 1\) and \((h_1 + h_2) - d = (1+1) - 1 = 1\). In addition, \( \mathcal{D}_1 \) is a PBS decomposition with \( \mathcal{C}_{\mathcal{D}_1} = \{M_1\} \) and belongs to \( \text{PBS}_1(1, 1) \).

Example 10. The \( \mathcal{C}^* \)-decomposition of a reaction network modeling the gene regulatory system of \( \text{Mycobacterium tuberculosis} \) in [7] is a PBS\(_1(1, 1) \) decomposition.

Example 11. The reaction graph below represents the CRN \( \mathcal{N} \) of \( q \)-site distributive phosphorylation/dephosphorylation \((q \geq 2) \) [4].

\[
\begin{align*}
S_0 + K & \iff S_0K \rightarrow S_1 + K \iff S_1K \iff S_2 + K \iff \cdots \rightarrow S_q + K \\
S_q + F & \iff \cdots \rightarrow S_2 + F \iff S_2F \rightarrow S_1 + F \iff S_1F \rightarrow S_0 + F
\end{align*}
\]

The \( S_i \)'s are the \( q \) substrates of a protein (the sites for both processes) while \( K \) and \( F \) are two different enzymes effecting phosphorylation and dephosphorylation respectively. Hernandez et al. [19] gave the decomposition \( \mathcal{D}_2 : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_q \) where

\[
\begin{align*}
\mathcal{N}_{i+1} : & \quad S_i + K \iff S_iK \rightarrow S_{i+1} + K \\
& \quad S_{i+1} + F \iff S_{i+1}F \rightarrow S_i + F
\end{align*}
\]

for \( i = 0, 1, \cdots, q - 1 \). They call \( \mathcal{D}_2 \) a fundamental decomposition (in the sense of Ji [22]). \( \mathcal{D}_2 \) is a PBS\(_d(0, 2) \) decomposition with \( \mathcal{C}_{\mathcal{D}_2} = \{S_{i+1} + K, S_{i+1} + F \mid i = 0, 1, \cdots, q - 2\} \) and \( d = 2(q - 1) \). Note that \( \mathcal{D}_2 \) is incidence independent (Proposition 3.38 in [19]).
We verify this using Proposition 5. Notice that \( \sum_{i=1}^{q} \ell_i - \ell = 2q - 2 \) and \( \sum_{i=1}^{q} h_i - d = h_1 + \sum_{i=2}^{q-1} h_i + h_q \) - \( d = [2 + 4(q - 2) + 2] - [2(q - 1)] = 2q - 2 \).

In a PBS decomposition, we denote by \( k(c) \) the number of subnetworks containing the “non-bottom” common pairwise intersection. As a simple illustration, a \( C \)-decomposition has \( k(c) = 0 \). For convenience, unless otherwise indicated, let the first \( k(c) \) subnetworks of a PBS decomposition (i.e., \( N_1, \ldots, N_{k(c)} \)) be the ones that contain the non-bottom common pairwise intersection. With this, the corresponding PIM \( M \) is the matrix \( [m_{ij}] \) where

\[
m_{ij} = \begin{cases} 
  c & \text{if } i < j \leq k(c) \\
  b & \text{if } j > k(c)
\end{cases} \quad (4.4)
\]

### 4.2 A structure theorem for the class of PBS decompositions

The following lemmas are vital in proving many of the succeeding results.

**Lemma 1.** Let \( \mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a PBS decomposition. If \( y' - x_1 - \cdots - x_q - y'' \) is a path in \( \mathcal{N} \) such that \( y' \) and \( y'' \) are common complexes while \( x_1, \ldots, x_q \) are not, then there is a subnetwork \( \mathcal{N}_i \) where \( y' \) and \( y'' \) are connected.

**Proof.** Suppose \( x_1 \) is a complex in subnetwork \( \mathcal{N}_i \). By definition, \( x_1 \) cannot be a complex in any other subnetwork. Hence, \( y' - x_1 \) is a subpath in \( \mathcal{N}_i \). If \( x_2 \) is not a complex in \( \mathcal{N}_i \), then the subpath \( x_1 - x_2 \) cannot occur in \( \mathcal{N}_i \) violating the assumption on \( x_1 \). So, \( x_2 \) is also a complex in \( \mathcal{N}_i \). Similarly, it can be shown that \( x_3, \ldots, x_q \) are also complexes in \( \mathcal{N}_i \). Moreover, \( y' - x_1 - \cdots - x_q - y'' \) is a path in \( \mathcal{N}_i \) indicating that \( y' \) and \( y'' \) are connected in \( \mathcal{N}_i \). \( \blacksquare \)

**Lemma 2.** Let \( \mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a PBS decomposition. If \( P : y - x_1 - \cdots - x_q \) is a path in \( \mathcal{N} \) such that \( y \) is a common complex while \( x_1, \ldots, x_q \) are not, then there is exactly one subnetwork \( \mathcal{N}_i \) that contains \( P \).

**Proof.** (Similar to the proof of Lemma 1) \( \blacksquare \)

Let a linkage class that has a common complex be called \( C \)-linkage class. We remark that Lemma 2 implies that every non-common complex of a \( C \)-linkage class of \( \mathcal{N} \) belongs to a \( C \)-linkage class of some subnetwork \( \mathcal{N}_i \).

The following result serves as the structure theorem for PBS decompositions which is a generalization of statement (ii) of the Structure Theorem for \( C \)-decomposition (Theorem 4).

**Theorem 5.** Let \( \mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a PBS decomposition such that at least one linkage class of \( \mathcal{N} \) does not contain a common complex. Let \( \mathcal{N}' \) and \( \mathcal{N}_i' \) be the networks obtained after removing from \( \mathcal{N} \) and \( \mathcal{N}_i \), respectively, the \( C \)-linkage classes. Then, \( \mathcal{D} : \mathcal{N}' = \mathcal{N}_1' \cup \cdots \cup \mathcal{N}_k' \) is a \( C \)-decomposition if \( \mathcal{D}' \) is non-trivial.

The condition that \( \mathcal{N} \) has a linkage class that has no common complex ensures that the decomposition obtained after the removal of the \( C \)-linkage classes is not going to be an empty decomposition. The worst possible scenario is that we get a trivial decomposition. We now give the proof of Theorem 5.
Proof of Theorem $5$. The length of $\mathcal{D}'$ is at least 2 since it is non-empty and non-trivial. Suppose $\mathcal{N}' = (\mathcal{D}', C', \mathcal{R}')$ and $\mathcal{N}'_j = (\mathcal{R}'_j, C'_j, \mathcal{R}'_j)$. First, we show that $\mathcal{D}'$ is a decomposition, i.e., $\{\mathcal{R}'_j\}$ is a partition of $\mathcal{D}'$. Clearly, $\mathcal{R}'_j$'s are pairwise disjoint. Let $R \in \mathcal{R}'$. Then, $R$ is not a reaction in any $C_2$-linkage class of $\mathcal{N}'$. Since $R \in \mathcal{D}$ and $\mathcal{D}$ is a decomposition, $R$ must be a reaction in exactly one subnetwork $\mathcal{N}'_j$. If $R$ belongs to a $C_2$-linkage class of $\mathcal{N}'_j$, then it also belongs to a $C_2$-linkage class of $\mathcal{N}$ which is a contradiction. Hence, $R$ occurs in $\mathcal{N}'_j'$, i.e., $R \in \mathcal{R}'_j$ which further implies that $R \in \mathcal{R}'$. Thus, $\mathcal{R} \subseteq \mathcal{R}'_j$.

Now, we show that $\bigcup \mathcal{R}'_j \subseteq \mathcal{R}'$. Let $R^* : x_1 \rightarrow x_2$ be a reaction that is not contained in $\mathcal{R}'$, i.e., $R^*$ belongs to a $C_2$-linkage class of $\mathcal{N}$. We want to show that $R^*$ is not contained in any $\mathcal{R}'_j$. If any of $x_1$ and $x_2$ are common complexes, we are done. Suppose none of them are common complexes. Let $y$ be the nearest common complex from $x_2$. Let $P : x_1 - x_2 - z_1 - z_2 - \cdots - z_p - y$, where $z_1, z_2, \ldots, z_p \notin C_2$, be the shortest path from $x_1$ to $y$. Suppose $R^*$ occurs in subnetwork $\mathcal{N}'_h$. Then, $x_1$ and $x_2$ also occur in $\mathcal{N}'_h$. By Lemma $2$, $P$ is a path in $\mathcal{N}'_h$ which means that $R^*$ belongs to a $C_2$-linkage class of $\mathcal{N}'_h$. So, $R^* \notin \mathcal{R}'_h$ or more generally, $R^*$ is not contained in any $\mathcal{R}'_j$. Hence, $\mathcal{D}'$ is a decomposition.

Finally, we show that for $i \neq j$, $C_i'$ and $C_j'$ are disjoint. Note that $C_i' \subseteq C_i$ and $C_j' \subseteq C_j$. If $C_i \cap C_j = \emptyset$, we are done. Suppose $C_i \cap C_j$ is non-empty. By definition, $C_i' \subseteq C_i \setminus C_2$ and $C_j' \subseteq C_j \setminus C_2$. So, $C_i' \cap C_j' \subseteq (C_i \setminus C_2) \cap (C_j \setminus C_2) = \emptyset$. Therefore, $\mathcal{D}'$ is a $C$-decomposition.

Example 12. Consider the decomposition $\mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2 \cup \mathcal{N}_3 \cup \mathcal{N}_4$ with the following reaction graphs.

$$\mathcal{N}:
\begin{array}{c}
x_1 \\
y_1 \\
x_2 \\
y_2 \\
x_3 \\
x_4 \\
y_1 \\
x_2 \\
x_5 \\
x_6 \\
x_7 \\
x_9 \\
x_8 \\
x_9 \\
x_{10}
\end{array}
$$

$$\begin{array}{c}
\mathcal{N}_1: \\
\mathcal{N}_2: \\
\mathcal{N}_3: \\
\mathcal{N}_4:
\end{array}
\begin{array}{c}
y_1 \\
x_1 \\
y_2 \\
x_2 \\
x_3 \\
x_4 \\
y_2 \\
x_2 \\
x_5 \\
x_6 \\
x_7 \\
x_9 \\
x_8 \\
x_{10}
\end{array}
$$

The decomposition is a PBS$2(0, 2)$ with $C_2 = \{y_1, y_2\}$ and $k(c) = 2$ (see $\mathcal{N}_1$ and $\mathcal{N}_2$). $\mathcal{N}$ has two linkage classes that do not contain any common complex. Removing the $C_2$-linkage classes from the parent network and the subnetworks yields the decomposition $\mathcal{N}' = \mathcal{N}_3' \cup \mathcal{N}_4'$ given by the following.
This decomposition is a $C$-decomposition.

### 4.3 The $C_{c,d}$ subclasses of the PBS class

**Definition 15.** For a nonnegative integer $d$, the class of pairwise binary-sized decompositions $\mathcal{PBS}_d$ is the union of all sets $\mathcal{PBS}_d(b, c)$ with $0 \leq b \leq c \leq d$.

We are interested in identifying subclasses of $\mathcal{PBS}_d$ which possess similar properties with respect to incidence independence deriving from the structure of their sets of common complexes.

**Example 13.** $C^*$-decompositions are contained in the subclass $\mathcal{PBS}_0 \cup \mathcal{PBS}_1(0, 1) \cup \mathcal{PBS}_1(1, 1)$. All decompositions in this subclass are incidence independent as will be shown in the next section. The $C^*$-decompositions consist of the special case when the single $C_y$ element $y$ is the 0 complex.

The above example motivates the following definition of subclasses with $b = 0$ or $b = c$.

**Definition 16.** The subclass $C_{c,d}$ is the union $\mathcal{PBS}_0 \cup \mathcal{PBS}_d(0, c) \cup \mathcal{PBS}_d(c, c)$.

**Example 14.** The subclass $C_{0,0}$ is the class of $C$-decompositions.

**Example 15.** The subclass $C_{1,1}$ contains the class of $C^*$-decompositions as well as the decomposition of Schmitz’s carbon cycle subnetwork given in Example 9.

### 5 The subclass $C_{d,d}$ of pairwise min-max decompositions

In this section, we we look into the subclass $C_{d,d}$ of PBS decompositions which we call class of pairwise min-max decompositions.

#### 5.1 The CCF for pairwise min-max decompositions

We define the set of pairwise min-max decompositions as follows.

**Definition 17.** The set of pairwise min-max (PMM) decompositions is defined for a set of common complexes $C_y$ with $d$ elements by either $C_i \cap C_j = \emptyset$ (minimal value) or $C_i \cap C_j = C_y$ (maximal value).
The PIM of a PMM decomposition is the matrix \([m_{ij}]\) where

\[
m_{ij} = \begin{cases} d & \text{if } i < j \leq k(c) \\ 0 & \text{if } j > k(c) \end{cases}.
\] (5.1)

A specific case of Proposition 5 gives a 3C-criterion for the incidence independence of the subclass PBS\(_d(0, d) \cup PBS_d(d, d)\):

\[
\sum \ell_i - \ell = d(k(c) - 1).
\] (5.2)

### 5.2 A fundamental property of PMM decompositions

The following proposition provides a sufficient condition for the incidence independence of a PMM decomposition where each of the CRN’s linkage classes contains at most one common complex. This result allows us to determine the incidence independence of a PMM decomposition by just observing the structure of the CRN’s reaction graph.

**Proposition 6.** Let \(D : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k\) be a PMM decomposition such that every linkage class of \(\mathcal{N}\) contains at most one common complex. Then, \(D\) is incidence independent.

*Proof.* Let \(\mathcal{N}_1, \mathcal{N}_2, \ldots, \mathcal{N}_k\) be the subnetworks containing \(\mathcal{C}_D\). Also, let \(\mathcal{N}'\) and \(\mathcal{N}'_i\) be the networks obtained after removing from \(\mathcal{N}\) and \(\mathcal{N}_i\), respectively, the \(\mathcal{C}_D\)-linkage classes. Let \(D' : \mathcal{N}' = \mathcal{N}'_1 \cup \cdots \cup \mathcal{N}'_k\). Clearly, \(\mathcal{N}'\) has \(\ell - d\) number of linkage classes.

On the other hand, Lemma 2 suggests that \(\mathcal{N}'_i\) has \(\ell_i - d\) number of linkage classes for \(i = 1, \ldots, k\).

Suppose \(D'\) is non-empty. Then, \(D'\) is either a trivial decomposition or a \(\mathcal{C}\)-decomposition (Theorem 5). The former immediately implies that

\[
\sum_{i=1}^{k} (\ell_i - d) + \sum_{i=k(c)+1}^{k} \ell_i = \ell - d.
\] (5.3)

Note that (5.3) still holds if \(D'\) is a \(\mathcal{C}\)-decomposition according to Theorem 3. Then,

\[
\sum_{i=1}^{k(c)} \ell_i - d(k(c)) + \sum_{i=k(c)+1}^{k} \ell_i = \ell - d \iff \sum_{i=1}^{k} \ell_i - \ell = d(k(c) - 1)
\] (5.4)

Now, suppose \(D'\) is empty. Then, it means that every linkage class in \(\mathcal{N}\) and \(\mathcal{N}_i\)’s is \(\mathcal{C}_D\)-linkage class. It follows that

\[
\sum_{i=1}^{k} (\ell_i - d) = 0 \text{ and } \ell - d = 0 \Rightarrow \sum_{i=1}^{k} (\ell_i - d) = \ell - d \iff \sum_{i=1}^{k} \ell_i - \ell = d(k - 1).
\] (5.5)

By Proposition 5 (5.4) and (5.5) suggest that, in any case, \(D\) is incidence independent. 

It immediately follows from Proposition 6 that PMM decompositions with \(|\mathcal{C}_D| \leq 1\) behave in the same way as \(\mathcal{C}^*\)-decompositions and are all incidence independent.
Example 16. The decomposition of the subnetwork of the of the Schmitz’s model given in Example 9 is a PMM decomposition with $\mathcal{C}_\varnothing = \{ M_1 \}$ which means that, indeed, it is incidence independent.

Example 17. Consider $\mathcal{N}$ whose reaction graph is given by

$$\mathcal{N} : \hspace{1cm} y_1 \rightarrow x_1 \quad y_2 \leftrightarrow x_4 \quad y_3 \rightarrow x_6 \quad x_9$$

$$\begin{array}{c}
\downarrow \\
 x_3 \\
 \downarrow \\
 x_2 \\
 \uparrow \\
 \downarrow \\
 x_5 \\
 \uparrow \\
 \downarrow \\
 x_8 \\
 \uparrow \\
 \downarrow \\
 x_7 \\
 \uparrow \\
 \downarrow \\
 x_10 \\
\end{array} \quad (5.6)$$

$\mathcal{N}$ can be decomposed as follows.

$$\mathcal{N}_1 : \hspace{1cm} y_1 \rightarrow x_1 \quad y_2 \leftrightarrow x_4 \quad y_3 \rightarrow x_6 \quad x_9$$

$$\begin{array}{c}
\downarrow \\
 x_2 \\
 \uparrow \\
 \downarrow \\
 \uparrow \\
 x_7 \\
 \uparrow \\
 \downarrow \\
 x_10 \\
\end{array} \quad (5.7)$$

$$\mathcal{N}_2 : \hspace{1cm} y_1 \rightarrow x_3 \quad y_2 \leftrightarrow x_5 \quad y_3 \leftrightarrow x_8$$

$\mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2$ is a PMM decomposition with $\mathcal{C}_\varnothing = \{ y_1, y_2, y_3 \}$. Notice that each linkage class of $\mathcal{N}$ contains at most one element from $\mathcal{C}_\varnothing$. Hence, by Proposition 6, this decomposition is incidence independent.

5.3 Special properties of the set PBS$_d(d, d)$ of $\mathcal{C}_\varnothing$-decompositions

We now discuss a PBS decomposition class where every subnetwork contains the whole $\mathcal{C}_\varnothing$.

Definition 18. For any subset $\mathcal{C}_\varnothing$ of $\mathcal{C}$ with $d$ elements, the $\mathcal{C}_\varnothing$-decomposition is defined by $\mathcal{C}_i \cap \mathcal{C}_j = \mathcal{C}_\varnothing$.

A special case of Proposition 5 gives $\sum \ell_i - \ell = d(k - 1)$ - the 3C-criterion for incidence independence of $\mathcal{C}_\varnothing$-decompositions.

The incidence independence of a $\mathcal{C}_\varnothing$-decomposition where every linkage class of the parent network contains at most one common complex is guaranteed by Proposition 6.

We now investigate the case where a linkage class of $\mathcal{N}$ contains more than one common complex.

Proposition 7. Let $\mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ be a $\mathcal{C}_\varnothing$-decomposition with $\mathcal{C}_\varnothing = \{ y_1, y_2 \}$ and $\ell > 2$. If there is at most one path in $\mathcal{N}$ that connects $y_1$ and $y_2$, then $\mathcal{D}$ is incidence independent.

Proof. If there is no path in $\mathcal{N}$ connecting $y_1$ and $y_2$, then $y_1$ and $y_2$ must be in separate linkage classes. So, By Proposition 6 $\mathcal{D}$ is incidence independent.

Suppose that in $\mathcal{N}$, there is exactly one path from $y_1$ to $y_2$. If such path is either of the reactions $y_1 \rightarrow y_2$ or $y_2 \rightarrow y_1$, then it will be contained in exactly one subnetwork. Now, if the path is given by $P : y_1 - a_1 - \cdots - a_q - y_2$ where $a_1, \cdots, a_q$ are not common complexes. Then, by Lemma 1 there will be a subnetwork $\mathcal{N}_j$ that contains $P$. The uniqueness of $P$
indicates that no other subnetwork will contain it, i.e., in any other subnetwork, \( y_1 \) and \( y_2 \) are in separate linkage classes.

Let \( \mathcal{N}' \) and \( \mathcal{N}'_i \) be the networks obtained after removing from \( \mathcal{N} \) and \( \mathcal{N}_i \), respectively, the linkage classes that contain \( y_1 \) or \( y_2 \). Let \( \mathcal{D}' : \mathcal{N}' = \mathcal{N}'_1 \cup \cdots \cup \mathcal{N}'_k \). From Lemma \( 2 \), \( \mathcal{N}'_j \) has \( \ell_j - 1 \) number of linkage classes while every other subnetwork has \( \ell_i - 2 \). If \( \mathcal{D}' \) is trivial, we have

\[
(\ell_j - 1) + \sum_{i \not= j} (\ell_i - 2) = \ell - 1
\]

(5.8)

If \( \mathcal{D}' \) is non-trivial, by Theorem \( 5 \) it is a \( \mathcal{G} \)-decomposition and (5.8) holds. In any case, \( \sum_{i=1}^k \ell_i - \ell = 2(k - 1) \) which means that by Proposition \( 5 \), \( \mathcal{D} \) is incidence independent.

For common complexes \( y' \) and \( y'' \), let \( P_{y'y''}^y \) denotes a path from \( y' \) to \( y'' \) whose “internal complexes” are non-common complexes. \( P_{y'y''}^y \) is the path \( y' \) – \( y'' \) if there are no internal complexes. A generalization of Proposition \( 4 \) is given in the following result.

**Theorem 6.** Let \( \mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a \( \mathcal{G} \)-decomposition such that \( \mathcal{G}_2 = \{ y_1, y_2, \ldots, y_d \} \) and \( \ell > 1 \). If there is exactly one path in \( \mathcal{N} \) that connects \( y_1, y_2, \ldots, y_d \) say \( P_{y_1}^{y_2} - \cdots - P_{y_d}^{y_{d+1}} \) such that \( P_{y_1}^{y_{d+1}} \) contains at least one non-common complex, then \( \mathcal{D} \) is incidence independent.

**Proof.** We prove this by induction on the number \( d \) of elements of \( \mathcal{G}_2 \). The case where \( d = 2 \) is given in Proposition \( 7 \). For the hypothesis of induction, we assume that every \( \mathcal{G}_2 \)-decomposition of a CRN with \( |\mathcal{G}_2| = d \) such that the common complexes are connected in the network by a unique path (having a property described in the proposition) is incidence independent.

Let \( \mathcal{D}' : \mathcal{N}' = \mathcal{N}'_1 \cup \cdots \cup \mathcal{N}'_k \) be a \( \mathcal{G} \)-decomposition with \( \mathcal{G}_2 = \{ y_1, y_2, \ldots, y_{d+1} \} \). Suppose that in \( \mathcal{N}' \), there is exactly one path that connects \( y_1, y_2, \ldots, y_{d+1} \) say \( P_{y_1}^{y_2} - \cdots - P_{y_d}^{y_{d+1}} \) with the property that \( P_{y_1}^{y_{d+1}} \) contains at least one non-common complex. Denote by \( P(y_{d+1}) \) a path with complexes consist only of \( y_{d+1} \) and non-common complexes, i.e., a path that passes through \( y_{d+1} \) whose other complexes are non-common complexes. Some of the paths \( P(y_{d+1}) \) (see the broken segments that passes through \( y_{d+1} \)) are shown in (5.9).

\[
P_{y_1}^{y_2} - \cdots - P_{y_d}^{y_{d+1}}
\]

(5.9)

Let \( \mathcal{N}^* = (\mathcal{I}^*, \mathcal{C}^*, \mathcal{R}^*) \) and \( \mathcal{N}_i^* = (\mathcal{I}_i^*, \mathcal{C}_i^*, \mathcal{R}_i^*) \) be the networks obtained after removing from \( \mathcal{N}' \) and \( \mathcal{N}_i' \) every path \( P(y_{d+1}) \). We now claim that \( \mathcal{D}^* : \mathcal{N}^* = \mathcal{N}_1^* \cup \cdots \cup \mathcal{N}_k^* \) is a \( \mathcal{G} \)-decomposition with \( \mathcal{G}_2 = \{ y_1, y_2, \ldots, y_d \} \). Let \( Q \) be the set of all paths \( P(y_{d+1}) \) and for convenience, let us call a path contained in \( Q \) a \( Q \)-path. First, we show that \( \{ \mathcal{R}_i^* \} \) is a partition of \( \mathcal{R}^* \). Clearly, \( \mathcal{R}_i^* \)'s are pairwise disjoint. Consider a reaction \( R \not\in \cup \mathcal{I}_i^* \). Then, \( R \) is “contained” in a \( Q \)-path in some subnetwork \( \mathcal{N}_i' \). It follows that \( R \) is also contained in a \( Q \)-path in \( \mathcal{N}' \). Hence, \( R \not\in \mathcal{R}^* \).
This time, let \( R : x' \to x'' \notin \mathcal{R}^* \). Then, \( R \) is in a \( Q \)-path in \( \mathcal{N} \). Consider the path \( x' - x'' - a_1 - \cdots - a_q - y_{d+1} \). From Lemma \( \Box_2 \) this path occurs in some subnetwork \( \mathcal{N}'_i \), i.e., \( R \) is in a \( Q \)-path in \( \mathcal{N}'_i \). So, \( R \notin \mathcal{R}^*_j \) and in general, \( R \notin \cup \mathcal{R}^*_i \). This means that \( \mathcal{D}^* \) is a decomposition.

Clearly, \( y_1, y_2, \ldots, y_{d-1} \) appear as complexes in every subnetwork \( \mathcal{N}'_i \). By virtue of Lemma \( \Box_2 \), \( Q \)-paths are contained in the same linkage class of every subnetwork \( \mathcal{N}'_i \). Now, due to uniqueness and Lemma \( \Box_1 \) the path \( P_{y_{d+1}} \) occurs in exactly one subnetwork say \( \mathcal{N}'_g \). This means that in \( \mathcal{N}'_g \), every \( Q \)-path is contained in the linkage class that also contains \( y_d \). On the other hand, in every other subnetwork \( \mathcal{N}'_i \), the linkage class that contains \( y_{d+1} \) does not have any other common complex. Since there is at least one non-common complex in \( P_{y_{d+1}} \), there is at least one reaction left in the linkage class of \( \mathcal{N}'_g \) that contains \( y_d \). Hence, \( y_d \) also appears as a complex in each \( \mathcal{N}'_i \). Thus, \( y_{d+1} \) is the only common complex that was removed to obtain \( \mathcal{N}^* \) and \( \mathcal{N}'_i \). So, \( \mathcal{D}^* \) is a \( \mathcal{C}_{\mathcal{D}} \)-decomposition.

Next, we show that there is exactly one path connecting \( y_1, y_2, \ldots, y_d \) in \( \mathcal{N}^* \). Suppose that such path does not exist. Then, there must be a non-common complex \( w \) in some path \( P_{w_{y_{d-1}}} \) that was removed from \( \mathcal{N}'_i \) in obtaining \( \mathcal{N}^* \), i.e., \( w \) is complex in some \( Q \)-path in \( \mathcal{N}'_i \).

\[
\begin{align*}
\text{Q-path} & \\
y_1 & \longrightarrow y_{i-1} & \longrightarrow w & \longrightarrow y_i & \longrightarrow y_d & \longrightarrow y_{d+1}
\end{align*}
\]

This implies that \( P_{w_{y_1}} - \cdots - P_{w_{y_{d-2}}} - w - \cdots - P_{w_{y_{d+1}}} \) is another path connecting the \( y_1, y_2, \ldots, y_{d+1} \) which is a contradiction. From the induction hypothesis, we conclude that \( \mathcal{N}^* = \mathcal{N}'_1 \cup \mathcal{N}'_2 \cup \cdots \cup \mathcal{N}'_3 \) is incident-independent. From Proposition \( \Box_5 \) we have

\[
\sum_{i=1}^{k'} \ell_i^* - \ell^* = d(k' - 1) \tag{5.11}
\]

where \( \ell^* \) and \( \ell_i^* \) are the number of linkage classes of \( \mathcal{N}^* \) and \( \mathcal{N}'_i \), respectively.

Let \( \ell' \) and \( \ell'_i \) be the number of linkage classes of \( \mathcal{N}' \) and \( \mathcal{N}'_i \), respectively. Then, we have \( \ell'_g = \ell_g^* \) and \( \ell'_i = \ell_i^* + 1 \) for \( i \neq g \). Also, we have \( \ell' = \ell^* \). Thus,

\[
\sum_{i=1}^{k'} \ell'_i - \ell' = \ell'_g + \sum_{i \neq g} \ell'_i - \ell' = \ell'_g + \sum_{i \neq g} (\ell_i^* + 1) - \ell^* = \sum_{i=1}^{k'} \ell_i^* - \ell^* + (k' - 1).
\]

From (5.11), we have

\[
\sum_{i=1}^{k'} \ell'_i - \ell' = d(k' - 1) + (k' - 1) = (d + 1)(k' - 1).
\]

Therefore, from Proposition \( \Box_5 \), \( \mathcal{D}' \) is incidence independent.

**Example 18.** Consider decomposition \( \mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2 \cup \mathcal{N}_3 \cup \mathcal{N}_4 \) in Example \( \Box_2 \). There is only the path \( y_1 - x_3 - y_2 \) connecting the common complexes \( y_1 \) and \( y_2 \) in \( \mathcal{N} \) such that the internal complex \( x_3 \) is a non-common complex. By Theorem \( \Box_6 \) the decomposition is incidence independent.
As shown in the next result, the incidence independence of a \( C_D \)-decomposition can still be achieved even if each linkage class of the parent network has none or has exactly two common complexes.

**Proposition 8.** Let \( \mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a \( C_D \)-decomposition with \( \frac{d}{2} < \ell \). Suppose that each linkage class of \( \mathcal{N} \) either contains zero or exactly two common complexes that are connected by a unique path. Then, \( \mathcal{D} \) is incidence independent.

**Proof.** Let \( \mathcal{L}_1, \ldots, \mathcal{L}_{d/2} \) be the \( C_D \)-linkage classes of \( \mathcal{N} \). For linkage class \( \mathcal{L}_j \), denote by \( y_1^j \) and \( y_2^j \) the common complexes it posses. From the assumption and Lemma 1, \( y_1^j \) and \( y_2^j \) belong to the same linkage class in some subnetwork say \( \mathcal{N}_h \) but are separated in every other subnetworks.

Let \( \mathcal{N}' \) and \( \mathcal{N}'_i \) be the networks obtained after removing from \( \mathcal{N} \) and \( \mathcal{N}_i \), respectively, the \( C_D \)-linkage classes. Let \( \mathcal{D}' : \mathcal{N}' = \mathcal{N}'_1 \cup \cdots \cup \mathcal{N}'_k \). If \( \mathcal{D}' \) is non-trivial, it is a \( C \)-decomposition by Theorem 5. Let \( \ell' \) and \( \ell'_i \) be the number of linkage classes of \( \mathcal{N}' \) and \( \mathcal{N}'_i \) and \( c_i \) be the number of pairs of connected \( y_1^j, y_2^j \) in \( \mathcal{N}_i \). Thus, \( \ell'_i = \ell_i - c_i - 2 \left( \frac{d}{2} - c_i \right) \).

By Theorem 3 we have
\[
\sum_{i=1}^{k} \ell'_i = \ell' \Leftrightarrow \sum_{i=1}^{k} \left[ \ell_i - c_i - 2 \left( \frac{d}{2} - c_i \right) \right] = \ell - \frac{d}{2} \Leftrightarrow \sum_{i=1}^{k} \ell_i - \ell = 2k \left( \frac{d}{2} \right) - \frac{d}{2} - \sum_{i=1}^{k} c_i. \tag{5.12}
\]

Since \( \sum_{i=1}^{k} c_i = \frac{d}{2} \), we get
\[
\sum_{i=1}^{k} \ell_i - \ell = 2k \left( \frac{d}{2} \right) - 2 \left( \frac{d}{2} \right) = d(k - 1). \tag{5.13}
\]

It can be shown that we still get (5.13) even if \( \mathcal{D}' \) is trivial. In any case, \( \mathcal{D} \) is incidence independent by Proposition 5.

## 6 A sufficient condition for complex balancing of PLK systems with incidence independent and complex balanced decompositions

In this section, we turn to the question: for an incidence independent and complex balanced decomposition (i.e., each subnetwork is complex balanced) of a kinetic system, which additional conditions must hold to ensure that the parent system is complex balanced? Recall that Farinas et al. [6] established the surprisingly general result that the conclusion holds for any \( C \)-decomposition, i.e., the additional purely structural condition of the set of common complexes \( C_D = \emptyset \) and any kinetic system (see the following theorem).

**Theorem 7** (Proposition 10, [6]). Let \((\mathcal{N}, K)\) be a CKS and \( \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \) be a weakly reversible \( C \)-decomposition (i.e., a \( C \)-decomposition where each subnetwork is weakly reversible). If \( Z_+ (\mathcal{N}_i, K_i) \neq \emptyset \) for each subnetwork, then \( Z_+ (\mathcal{N}, K) \neq \emptyset \).
In section 6.1, we show that this result is singular in its purely structural character: as soon as \( C \neq \emptyset \), there are going to be kinetic systems with an incidence independent and complex balanced decomposition that are not complex balanced. Accordingly, our new result, while imposing no conditions on the set of common complexes of the incidence independent decomposition, is valid only for a class of power law kinetic systems. After brief review of relevant results in sections 6.2 and 6.3, we state and prove the new result in section 6.4. We then briefly compare it with the generalization of the Deficiency Zero Theorem of Fortun et al. [14] and sketch some interesting research questions.

6.1 A simple counterexample for \( C \neq \emptyset \)

Consider the \( C^\ast \)-decomposition \( N = N_1 \cup N_2 \) given by the following.

\[
N : \quad A + B \overset{0}{\underset{2A + 2B}{\leq}} 0 \quad \overset{2A + 2B}{\leq} 2A
\]

\[
N_1 : \quad A + B \overset{0}{\underset{2A + 2B}{\leq}} 0
\]

\[
N_2 : \quad 0 \overset{2A + 2B}{\leq} 2A
\]

Suppose that \((N, K)\) is endowed with mass action kinetics. For the reversible and deficiency zero subsystem \((N_1, K_1)\) we have the equations

\[
dA/dt = dB/dt = k_1 - k_2 AB
\]

and \( AB = k_1/k_2 \) at equilibrium. Similarly, for the subsystem \((N_2, K_2)\), which is also reversible and deficiency zero, we get

\[
dA/dt = dB/dt = 2k_3 - 2k_4 A^2 B^2
\]

and \( AB = \sqrt{k_3/k_4} \) at equilibrium. For the parent system, we have

\[
dA/dt = dB/dt = k_1 - k_2 AB + 2k_3 - 2k_4 A^2 B^2.
\]

For the rate vector \((2, 4, 1, 1/2)\) we have

\[
Z_+(N_1, K_1) = \{(A, B) \in \mathbb{R}^2 | AB = 1/2\} \quad \text{and} \quad Z_+(N_2, K_2) = \{(A, B) \in \mathbb{R}^2 | AB = \sqrt{2}\}
\]

so that their intersection is empty, suggesting that \((N, K)\) is not complex balanced. On the other hand, \((N', K)\) is complex balanced for the rate vector \((1, 1, 1, 1)\) with

\[
E_+(N', K) = Z_+(N', K) = Z_+(N_1, K_1) = Z_+(N_2, K_2) = \{(A, B) \in \mathbb{R}^2 | AB = 1\}.
\]

Hence, even mass action systems on the same network and the same incidence independent and complex balanced decomposition (in fact a \( C^\ast \)-decomposition) may or may not be complex balanced.

6.2 Kinetic systems of LP type

In this and the next subsection, we collect some relevant concepts for the new result.
Definition 19. A kinetic system \((\mathcal{N}, K)\) is of type PLP (positive equilibria log-parametrized) if

i. \(E_+ (\mathcal{N}, K) \neq \emptyset\) and

ii. \(E_+ (\mathcal{N}, K) = \{ x \in \mathbb{R}_{>0}^S | \log x - \log x^* \in (P_E)^\perp \}\)

where \(P_E\) is a subspace of \(\mathbb{R}^S\) and \(x^*\) is a positive equilibrium.

Definition 20. A kinetic system \((\mathcal{N}, K)\) is of type CLP (complex balanced equilibria log-parametrized) if

i. \(Z_+ (\mathcal{N}, K) \neq \emptyset\) and

ii. \(Z_+ (\mathcal{N}, K) = \{ x \in \mathbb{R}_{>0}^S | \log x - \log x^* \in (P_Z)^\perp \}\)

where \(P_Z\) is a subspace of \(\mathbb{R}^S\) and \(x^*\) is a complex balanced equilibrium.

A kinetic system is bi-LP if it is of PLP and of CLP type and \(P_E = P_Z\). We will use the shorter PLP system, CLP system and bi-LP system notation as well as the collective term “LP systems”.

A key property of an LP system was in principle already derived by Feinberg in his 1979 lectures [10] as shown in [27]:

Theorem 8. Let \((\mathcal{N}, K)\) be a chemical kinetic system.

i. If \((\mathcal{N}, K)\) is a PLP system, then \(|E_+ (\mathcal{N}, K) \cap Q| = 1\) for any positive coset \(Q\) of \(P_E\) in \(\mathbb{R}^S\).

ii. If \((\mathcal{N}, K)\) is a CLP system, then \(|Z_+ (\mathcal{N}, K) \cap Q| = 1\) for any positive coset \(Q\) of \(P_Z\) in \(\mathbb{R}^S\).

iii. If \((\mathcal{N}, K)\) is a bi-LP system, then it is absolutely complex balanced, i.e., each positive equilibrium is complex balanced.

The concepts of kinetic order subspace and kinetic deficiency for cycle terminal PL-RDK systems were introduced by Müller and Regensburger in [29] where they also presented their theory of generalized mass action systems (GMAS). Hernandez and Mendoza developed these concepts in [21] by first constructing the reaction network induced by the kinetic complexes which we are adapting in this paper.

Let \(\mathcal{R}(y)\) be the set of (branching) reactions having \(y\) as reactant complex.

Definition 21. Let \((\mathcal{N}, K)\) be a cycle terminal PLK system. Given a complex \(y\), the set of kinetic complexes of \(y\) is the set \(\mathcal{C}(y) := \{ F_R \mid R \in \mathcal{R}(y) \}\) (\(F_R\) refers to the kinetic order vector that corresponds to reaction \(R\)). Moreover, for a reaction \(R : y \rightarrow y'\), \(\mathcal{R}(R) := \{ \tilde{y} \rightarrow \tilde{y}' \mid \tilde{y} \in \mathcal{C}(y) \text{ and } \tilde{y}' \in \mathcal{C}(\tilde{y}') \}\) gives the set of kinetic complex reactions of \(R\).

The reaction network of kinetic complexes of a cycle terminal PLK system is defined as follows.

Definition 22. The set of kinetic complexes induced a reaction network given by \(\tilde{\mathcal{N}} = (\tilde{\mathcal{I}}, \tilde{\mathcal{C}}, \tilde{\mathcal{R}})\) where \(\tilde{\mathcal{C}} = \bigcup_y \mathcal{C}(y)\) and \(\tilde{\mathcal{R}} = \bigcup_y \mathcal{R}(y)\).
The orders of \( \tilde{c} \) and \( \tilde{r} \) are denoted by \( \tilde{n} \) and \( \tilde{p} \), respectively. In addition, the incidence and stoichiometric maps of \( \mathcal{N} \) are respectively denoted by \( \tilde{I}_a \) and \( \tilde{N} \).

**Definition 23.** The kinetic order subspace \( \tilde{S} \) of \( (\mathcal{N}, K) \) is defined to be the image of \( \tilde{N} \) and its dimension \( \tilde{s} \) is also referred to as kinetic rank. The kinetic complex deficiency is the nonnegative number given by \( \delta_{\tilde{N}} = \tilde{n} - \tilde{\ell} - \tilde{s} \). On the other hand, the kinetic deficiency is the nonnegative number \( \tilde{\delta} = n - \ell - \tilde{s} \).

A key ingredient of the new result is the following theorem of Müller and Regensburger [29].

**Theorem 9.** Let \( (\mathcal{N}, K) \) be a weakly reversible PL-RDK system. If \( (\mathcal{N}, K) \) is complex balanced, then it is of CLP type with \( P_Z = \tilde{S} \).

With the concept of kinetic deficiency, they provide a characterization of complex balancing: if the kinetic deficiency is zero, then any weakly reversible PL-RDK system is complex balanced. If it is positive, they provide a necessary and sufficient condition for the system to be complex balanced.

### 6.3 Induced decompositions of kinetic complexes

The second part of the sufficient condition is based on the concepts of an induced subnetwork of kinetic complexes and a corresponding decomposition, which were introduced in [21]. For a PL-RDK system, a kinetic complex is the column of the \( T \) matrix assigned to a reactant complex. For a decomposition \( \tilde{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k \), we set \( \tilde{\mathcal{N}} := \text{subnetwork of } \tilde{\mathcal{N}} \text{ induced by } \mathcal{N}_i \), i.e., take the reactions defining it, then form the kinetic complexes (still in the sense of Müller-Regensburger).

**Definition 24.** The induced subnetwork \( \tilde{\mathcal{N}}_{\tilde{D}} \) is defined as the union of the \( \tilde{\mathcal{N}}_i \). If the covering \( \tilde{D} \) is a decomposition, we call it the induced decomposition.

We denote by \( \tilde{n}_{\tilde{D}} \) and \( \tilde{\ell}_{\tilde{D}} \) the number of complexes and linkage classes of \( \tilde{\mathcal{N}}_{\tilde{D}} \), respectively.

In the following, we will assume that the covering is indeed a decomposition (so that in the examples, this has to be verified). Note however, that the following proposition holds:

**Proposition 9.** If, for an induced covering \( \tilde{D} \), the flux spaces \( \tilde{S}_i \) form a direct sum, then the covering is an independent decomposition.

**Example 19.** Recall the incidence independent decomposition \( \tilde{D}_1 : \mathcal{N} = \mathcal{N}_1 \cup \mathcal{N}_2 \) given in Example 9 where \( \mathcal{N} \) is the weakly reversible and deficiency zero subnetwork of the Schmitz’s pre-industrial carbon cycle model (4.1). In [16], \( \mathcal{N} \) is assigned with the following kinetic order matrix which makes the system PL-NDK.

\[
F = \begin{bmatrix}
M_1 & M_2 & M_3 & M_4 & M_5 & M_6 \\
0 & 0 & 0 & 0 & 1 & 0 \\
0 & 0 & 0 & 0 & 1 & 0 \\
0 & 0 & 0 & 0 & 1 & 0 \\
0 & 9.4 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 1 & 0 & 0 \\
0 & 0 & 1 & 0 & 0 & 0 \\
1 & 0 & 0 & 0 & 0 & 0 
\end{bmatrix}
\] (6.2)
On the other hand, each corresponding subsystem is weakly reversible and PL-RDK. Let $\tilde{\mathcal{N}}_1$ be the network of kinetic complexes given by $\{\tilde{R}_1 : M_5 \to 0.36M_1, \tilde{R}_2 : 0.36M_1 \to M_5, \tilde{R}_3 : R_3, \tilde{R}_4 : M_6 \to 0.36M_1\}$ and $\tilde{\mathcal{N}}_2$ be given by $\{\tilde{R}_5 : 9.4M_2 \to M_1, \tilde{R}_6 : M_4 \to 9.4M_2, \tilde{R}_7 : R_7, \tilde{R}_8 : R_8\}$ (see (6.3)).

Accordingly, $\tilde{\mathcal{N}}_D = \tilde{\mathcal{N}}_1 \cup \tilde{\mathcal{N}}_2$. Observe that $\tilde{n}_D = 7$ and $\tilde{\ell}_D = 2$ implying that $\tilde{n}_D - \tilde{\ell}_D = 5$. Moreover, notice that $\tilde{s}_D = 5$, $\tilde{s}_1 = 2$, and $\tilde{s}_2 = 3$ suggesting that, by Proposition 9, the decomposition $\tilde{\mathcal{N}}_D = \tilde{\mathcal{N}}_1 \cup \tilde{\mathcal{N}}_2$ is independent.

6.4 Statement and proof of the sufficient condition

We can now formulate and derive our main result:

**Theorem 10.** Let $(\mathcal{N}, K)$ be a weakly power law system with a complex balanced PL-RDK decomposition $\mathcal{D} : \mathcal{N} = \mathcal{N}_1 \cup \cdots \cup \mathcal{N}_k$ with $P_{Z,i} = \tilde{S}_i$. If $\mathcal{D}$ is incidence independent and the induced covering $\tilde{\mathcal{D}}$ is independent, then $(\mathcal{N}, K)$ is a weakly reversible CLP system with $P_Z = \sum \tilde{S}_i$.

**Proof.** Each subnetwork of $(\mathcal{N}, K)$ in the decomposition is weakly reversible and hence, the network is weakly reversible. For the last two statements, we prove the case when $k = 2$ while the general case can be proven inductively.

Since each decomposition subnetwork is a complex balanced PL-RDK system, by Theorem 9 it is of CLP type, i.e., $Z_+ (\mathcal{N}_i, K_i) \neq \emptyset$ and for $x_i \in Z_+ (\mathcal{N}_i, K_i)$,

$$Z_+ (\mathcal{N}_i, K_i) = \{x \in \mathbb{R}^+_\times | \log x - \log x_i \in \tilde{S}_i^\perp\}.$$ 

Since the decomposition is incidence independent, we have

$$Z_+ (\mathcal{N}, K) = Z_+ (\mathcal{N}_1, K_1) \cap Z_+ (\mathcal{N}_2, K_2)$$

Hence, $x^* \in Z_+ (\mathcal{N}, K)$ if and only if

$$\log x^* \in \left( \log x_1 + \tilde{S}_1^\perp \right) \cap \left( \log x_2 + \tilde{S}_2^\perp \right).$$

From properties of cosets,

$$\left( \log x_1 + \tilde{S}_1^\perp \right) \cap \left( \log x_2 + \tilde{S}_2^\perp \right) \neq \emptyset \iff \log x_1 - \log x_2 \in \tilde{S}_1^\perp + \tilde{S}_2^\perp.$$
The independence of the induced decomposition ensures that \( \tilde{S}_1 \cap \tilde{S}_2 = \{0\} \). Then, \( \tilde{S}_1^\perp + \tilde{S}_2^\perp = (\tilde{S}_1 \cap \tilde{S}_2)^\perp = \{0\}^\perp = \mathbb{R}^\perp \). Thus,

\[
\left( \log x_1 + \tilde{S}_1^\perp \right) \cap \left( \log x_2 + \tilde{S}_2^\perp \right) \neq \emptyset.
\]

Let \( \hat{x} \in \left( \log x_1 + \tilde{S}_1^\perp \right) \cap \left( \log x_2 + \tilde{S}_2^\perp \right) \) and take \( x^* = e^{\hat{x}} \). We have \( \hat{x} \in \left( \log x_1 + \tilde{S}_1^\perp \right) \) and \( \hat{x} \in \left( \log x_2 + \tilde{S}_2^\perp \right) \), hence \( x^* \in Z_+(\mathcal{M}_1, K_1) \cap Z_+(\mathcal{M}_2, K_2) = Z_+(\mathcal{N}, K) \). To show the log parametrization, note that

\[
\left( \log x_1 + \tilde{S}_1^\perp \right) \cap \left( \log x_2 + \tilde{S}_2^\perp \right) = \log x^* + \left( \tilde{S}_1^\perp + \tilde{S}_2^\perp \right) = \log x^* + \left( \tilde{S}_1^\perp + \tilde{S}_2^\perp \right)^\perp.
\]

Hence, \( Z_+(\mathcal{N}, K) = \left\{ x \in \mathbb{R}_{>0}^\perp \mid \log x - \log x^* \in \left( \tilde{S}_1^\perp + \tilde{S}_2^\perp \right)^\perp \right\} \). \( \blacksquare \)

**Corollary 1.** Let \((\mathcal{N}, K)\) be a PL-NDK system satisfy the conditions of Theorem 10 for a decomposition into PL-NDK subnetworks with zero kinetic deficiency, i.e., \( \tilde{\delta} = 0 \). Then the induced subnetwork has zero kinetic complex deficiency and \((\mathcal{N}, K)\) is unconditionally complex balanced, i.e., it is complex balanced for any set of rate constants.

**Proof.** Since the decomposition \( \mathcal{D} \) is incidence independent, we have \( n - \ell = \sum(n_i - \ell_i) \), and due to the surjective map \( \mathcal{E} \to \mathcal{E}^* \), \( n_i - \ell_i \geq \tilde{n}_i - \tilde{\ell}_i \), we obtain the independence of the induced decomposition that the kinetic complex deficiency is less than the sum of the subnetworks’ kinetic deficiencies which is zero. The unconditional complex balancing results from the conclusion of the theorem. \( \blacksquare \)

**Example 20.** Corollary 1 holds for a PL-RDK system satisfying the conditions of the theorem for a weakly reversible PL-TIK decomposition, since Talebis et al. have shown that such system have zero kinetic deficiency \([34]\). One could in fact extend the concept of a PL-TIK system (i.e., a system with zero kinetic reactant deficiency) to such a PL-NDK systems with the additional requirement that the induced decomposition be \( \tilde{R} \)-independent, i.e., the augmented reactant subspaces of the subnetworks form a direct sum.

We remark that Corollary 1 suggests that further results on GMAS such as those of Cracium et al. on the generalized Birch’s Theorem \([5]\), and Boros et al. \([3]\) on linear stability can also be extended to such PL-NDK systems.

### 6.5 Discussion

The generalization of the main result of Fortun et al. \([14]\) is the following theorem:

**Theorem 11.** Let \((\mathcal{N}, K)\) be a power law systems with a weakly reversible PL-RDK decomposition \( \mathcal{D} : \mathcal{N} = \mathcal{M}_1 \cup \cdots \cup \mathcal{M}_k \). If \( \mathcal{D} \) is bi-level independent (both the network decomposition and the induced subnetwork decomposition are independent) and of PLP type with \( P_{E,i} = \tilde{S}_i \), then \((\mathcal{N}, K)\) is a weakly reversible PLP system with \( P_E = \sum \tilde{S}_i \).

A comparison with the main result immediately reveals that the two results are fully analogous. In fact, their proofs differ only in the use of Theorem 1 in the PLP case and
Theorem for CLP systems. On the other hand, there are only a few known examples of PLP PL-RDK systems in contrast to the very general results of Müller and Regensburger about CLP PL-RDK systems, so that currently our main result can be considered more impactful.

A further interesting point of discussion concerns research questions raised by our main result. The fact that the induced decomposition plays an essential role suggests that it may be interesting to study the set of common complexes of the induced decomposition and explore relationships between it and the common complexes of the original decomposition. These relationships could also be decomposition class specific, e.g., for PMM decompositions. Also fully unexplored are relationships between decomposition classes and conditions for independence of the induced decomposition.

Another set of research questions emerge from the observation that, so far, the results on the complex balancing of the whole networks are “binar” in the sense that Farinas et al. deal with $C_D = \emptyset$ and our main result with $C_D \neq \emptyset$ for any incidence independent decomposition. If one refined the latter set by considering only a particular decomposition class, again say PMM that are incidence independent, it might be possible to identify further classes of kinetics which enable complex balancing of the whole system.

\section{Summary and outlook}

We provide here a summary of this paper as well as some recommendations for further studies.

1. We investigated the incidence independence of decompositions using the set of common complexes of the subnetworks. A framework was created to characterize decomposition classes by their incidence independence properties.

2. Interesting decomposition classes, identified by the sets of their subnetworks’ common complexes, were introduced. Subclasses of these decomposition classes, which are incidence independent, were specified.

3. For future studies, other structural conditions sufficient or necessary for the incidence independence of other subclasses of PBS decompositions can be explored. Other decomposition classes identified by their sets of common complexes can also be considered and investigated.

4. We identified a sufficient condition that guarantees the existence of complex balanced equilibria of some PLK systems with incidence independent and complex balanced decompositions.

5. Through the result identified in 4, we obtained a generalization of the Deficiency zero Theorem for some PLK systems.

6. It is recommended to explore the capacity of other PLK systems to admit complex balanced equilibria that have incidence independent decompositions.

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References

[1] C. Arceo, E. Jose, A. Marín-Sanguino, E. Mendoza, Chemical reaction network approaches to biochemical systems theory, Math. Biosci. 269 (2015) 135–152.

[2] B. Boros, On the existence of the positive steady states of weakly reversible deficiency-one mass action systems, Math. Biosci. 245 (2013) 157–170.

[3] B. Boros, S. Müller, G. Regensburger, Complex-balanced equilibria of generalized mass-action systems: Necessary conditions for linear stability, Math. Biosci. Eng. 17 (2020) 442–459.

[4] C. Conradi, A. Shiu, Dynamics of posttranslational modification systems: recent results and future directions, Biophysical Journal 114(3) (2018) 507–515.

[5] G. Craciun, S. Müller, C. Pantea, P. Yu, A generalization of Birch’s theorem and vertex-balanced steady states for generalized mass-action systems, Math. Biosci. Eng. 16 (2019) 8243–8267.

[6] H. Farinas, E. Mendoza, A. Lao, Chemical reaction network decompositions and realizations of S-systems, Philippine Science Letters 14(1) (2021).

[7] H. Farinas, E. Mendoza, A. Lao, Structural properties of an S-system model of Mycobacterium tuberculosis gene regulation, Philippine Journal of Science 149(3) (2020) 539–555.

[8] M. Feinberg, Chemical reaction network structure and the stability of complex isothermal reactors I: The deficiency zero and deficiency one theorems, Chem. Eng. Sci. 42 (1987) 2229–2268.

[9] M. Feinberg, Foundations of chemical reaction network theory, Springer (2019).

[10] M. Feinberg, Lectures on chemical reaction networks, Notes of lectures given at the Mathematics Research Center of the University of Wisconsin (1979). Available at https://crnt.osu.edu/LecturesOnReactionNetworks.

[11] L. Fontanil, E. Mendoza, N. Fortun, A computational approach to concentration robustness in power law kinetic systems of Shinar-Feinberg type, MATCH Comm. Math. Comput. Chem. (2021, in press).
[12] N. Fortun, D. A. Talabis, E. Jose, E. Mendoza, Complex balanced equilibria of weakly reversible poly-PL systems: multiplicity, robustness and stability, arXiv: 2006.10227 (2020)

[13] N. Fortun, A. Lao, L. Razon, E. Mendoza, A deficiency-one algorithm for a power-law kinetics with reactant-determined interactions, J. Math. Chem. 56 (2018) 2929–2962.

[14] N. Fortun, E. Mendoza, L. Razon, A. Lao, A deficiency zero theorem for a class of power law kinetic systems with non-reactant determined interactions, MATCH Commun. Math. Comput. Chem. 81 (2019) 621–638.

[15] N. Fortun, E. Mendoza, Absolute concentration robustness in power law kinetic systems, MATCH Commun. Math. Comput. Chem. 85 (2021) 669–691.

[16] N. Fortun, E. Mendoza, L. Razon, A. Lao, Robustness in power-law kinetic systems with reactant-determined interactions, Lecture Notes in Computer Science (2019, in press).

[17] E. Gross, H. Harrington, N. Meshkat, and A. Shiu, Joining and decomposing reaction networks, Math. Biol. 80 (2020), 1683–1731.

[18] B. Hernandez, E. Mendoza, A. de los Reyes V, A computational approach to multistationarity of power-law kinetic systems, J. Math. Chem. 58 (2019), 56–87.

[19] B. Hernandez, E. Mendoza, A. de los Reyes V, Fundamental decompositions and multistationarity of power-law kinetic systems, MATCH Commun. Math. Comput. Chem. 83 (2020) 403–434

[20] B. Hernandez, E. Mendoza, Positive equilibria of Hill-type systems, Journal of Mathematical Chemistry (2021, in press).

[21] B. Hernandez, E. Mendoza, Weakly reversible CF-decompositions of chemical kinetic systems (2021, submitted).

[22] H. Ji, Uniqueness of equilibria for complex chemical reaction networks, Ph.D. Dissertation, Ohio State University, 2011

[23] E. Jose, E. Mendoza, D. A. Talabis, Absolutely complex balanced chemical kinetic systems (2021, submitted).

[24] B. Joshi and A. Shiu, Atoms of multistationarity in chemical reaction networks, J Math Chem 51 (2013) 153–178.

[25] B. Joshi and A. Shiu, Simplifying the Jacobian Criterion for precluding multistationarity in chemical reaction networks, SIAM J. Appl. Math. 72 (2012) 857–876.
[26] G. Magombedze and N. Mulder, Understanding TB latency using computational and
dynamic modelling procedures, *Infect Genet Evol* **13** (2013) 267–283.

[27] D. Magpantay, E. Mendoza, Concentration robustness in chemical kinetic systems of
LP type (2021, in preparation).

[28] E. Mendoza, D. A. Talabis, E. Jose, Positive equilibria of weakly reversible power
law kinetic systems with linear independent interactions, *J Math Chem*. **56** (2018)
2643–2673.

[29] S. Müller, G. Regensburger, Generalized mass action systems and positive solutions
of polynomial equations with real and symbolic exponents, *Lecture Notes in Comput.
Sci.* **8660** (2014) 302–323.

[30] A. Nazareno, R. P. Eclarin, E. Mendoza, A. Lao, Linear conjugacy of chemical kinetic
systems, *Math. Biosci. Eng*. **16** (2019) 8322–8355.

[31] I. Otero-Muras, P. Yordanov, J. Stelling, Chemical Reaction Network Theory elu-
cidates sources of multistability in interferon signaling, *PLoS Comput Biol* **13**(4)
(2017).

[32] R. Schmitz, The Earth’s carbon cycle: Chemical engineering course material, *Chem.
Engin. Edu*. **36** (2002) 296–309.

[33] G. Shinar, M. Feinberg, Structural sources of robustness in biochemical reaction
networks, *Science* **327** (2010) 1389–1391.

[34] D. A. Talabis, E. Mendoza, E. Jose, Complex balanced equilibria of weakly reversible
power law kinetic systems, *MATCH Commun. Math. Comput. Chem*. **82** (2019) 601–
624.

[35] J. Tóth, A. Nagy, D. Papp, *Reaction Kinetics: Exercises, Programs and Theorems*,
Springer, New York, 2018.
The following table list the notations used in this paper that were adapted for subnetworks.

| Notion                                              | Parent Network: $\mathcal{N}$ | Subnetwork: $\mathcal{N}_i$ |
|-----------------------------------------------------|-------------------------------|-----------------------------|
| Species, complex, and reaction sets                 | $\mathcal{S}, \mathcal{C},$ and $\mathcal{R}$ | $\mathcal{S}_i, \mathcal{C}_i, \mathcal{R}_i$ |
| Number of species, complexes, and reactions         | $m, n, \mathrm{and} \, p$     | $m_i, n_i, \mathrm{and} \, p_i$ |
| Stoichiometric subspace                             | $S$                           | $S_i$                        |
| Network rank                                        | $s$                           | $s_i$                        |
| Deficiency                                          | $\delta$                      | $\delta_i$                   |
| Number of linkage classes                           | $\ell$                        | $\ell_i$                     |
| Incidence map/matrix                                | $I_{a}$                       | $I_{a,i}$                    |
| Reaction network of kinetic complexes               | $\mathcal{N}$                 | $\mathcal{N}_i$              |
| Set of kinetic complexes                            | $\mathcal{C}$                 | $\mathcal{C}_i$              |
| Kinetic order subspace                              | $S$                           | $S_i$                        |
| Kinetic rank                                        | $s$                           | $s_i$                        |
| Kinetic deficiency                                  | $\delta$                      | $\delta_i$                   |

Notations used for the parent network and the subnetworks.