TREND TO EQUILIBRIUM FOR REACTION-DIFFUSION SYSTEMS ARISING FROM COMPLEX BALANCED CHEMICAL REACTION NETWORKS

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Abstract. The quantitative convergence to equilibrium for reaction-diffusion systems arising from complex balanced chemical reaction networks with mass action kinetics is studied by using the so-called entropy method. In the first part of the paper, by deriving explicitly the entropy dissipation, we show that for complex balanced systems without boundary equilibria, each trajectory converges exponentially fast to the unique complex balance equilibrium. Moreover, a constructive proof is proposed to explicitly estimate the rate of convergence in the special case of a cyclic reaction. In the second part of the paper, complex balanced systems with boundary equilibria are considered. We focus on a specific case involving three chemical substances for which the boundary equilibrium is shown to be unstable in some sense, so that exponential convergence to the unique strictly positive equilibrium is recovered.

1. Introduction and Main results

The foundation of the study of chemical reaction networks goes back to the pioneering works of F.J.M. Horn, R. Jackson, A.I. Volpert and M. Feinberg, (see e.g. [HJ72, Vol72, FH74, Fei79, Fei87]). The aim of the theory is to study the behaviour of chemical reaction networks regardless of specific values of the reaction rate constants, since these rates may be hard to determine in some practical situations. One of the main questions of the theory is the large time asymptotic behaviour of the dynamical system corresponding to the reaction network. While the ODE setting of chemical reaction networks has been extensively studied in the literature, the PDE setting is much less investigated. This is the main motivation of this work. The ultimate aim of this paper is to investigate the large time asymptotic behaviour of reaction-diffusion systems arising from chemical reaction networks. More precisely, by exploiting the so-called entropy method, we look for quantitative estimates of convergence to equilibrium for a class of chemical reaction networks called complex balanced reaction networks.

Before providing a general notation of chemical reaction networks, we shall consider the example of a single reaction $S_1 + S_2 \xrightarrow{k} S_3$ where $k > 0$ is a reaction rate constant. To describe this reaction, we introduce the set of chemical substances $\mathcal{S} = \{S_1, S_2, S_3\}$, the set of chemical complexes $\mathcal{C} = \{(1, 1, 0); (0, 0, 1)\}$ corresponding to $S_1 + S_2$ and $S_3$ and the set of chemical reactions $\mathcal{R} = \{(1, 1, 0) \xrightarrow{k} (0, 0, 1)\}$. Now, by following the notation of e.g. [HJ72], a general chemical reaction network consists of $N$ chemical substances $\mathcal{S} = \{S_1, \ldots, S_N\}$, a finite set $\mathcal{C}$ of chemical complexes, which appear on either sides of a chemical reaction, i.e. $\mathcal{C} = \{y\} \subset \mathbb{N}^N$ (This is a slight abuse of notation. We should indeed write $\{y^{(1)}, \ldots, y^{(|\mathcal{C}|)}\}$, but we do not because superscripts with another meaning will be used in the sequel) and the set of

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chemical reactions $\mathcal{R} = \{y_r \xrightarrow{k_r} y'_r : y_r,y'_r \in \mathcal{C}, k_r > 0\}$. We shall only consider chemical reaction networks which satisfy the following natural assumptions:

**Definition 1.1 (Chemical reaction networks).**

Let $\mathcal{S} = \{S_i\}_{i=1}^N$, $\mathcal{C} = \{y\} \subset \mathbb{N}^N$ and $\mathcal{R} = \{y_r \xrightarrow{k_r} y'_r\}_{r=1,...,|\mathcal{R}|}$ denote finite sets of species, complexes and reactions, respectively. The triple $\{\mathcal{S}, \mathcal{C}, \mathcal{R}\}$ is called a chemical reaction network as long as the following three natural requirements are met:

1. For each $S_i \in \mathcal{S}$, there exists at least one complex $y \in \mathcal{C}$ in which the stoichiometric coefficient $\nu_i$ of $S_i$ is strictly positive (subscripts for $y$ are used both to prescribe stoichiometric coefficients (usually $i$ is then used) and to prescribe the index of a chemical reaction (usually $r$ is then used); when both need to be prescribed, we use $y_{r,i}$);

2. There is no trivial reaction $y \rightarrow y \in \mathcal{R}$ for any complex $y \in \mathcal{C}$;

3. For any $y \in \mathcal{C}$, there must exist $y' \in \mathcal{C}$ such that either $y \rightarrow y' \in \mathcal{R}$ or $y' \rightarrow y \in \mathcal{R}$.

We assume that the chemical species corresponding to the reaction network $\{\mathcal{S}, \mathcal{C}, \mathcal{R}\}$ are contained in a bounded vessel (or reactor) $\Omega \subset \mathbb{R}^n$, where $\Omega$ is a bounded smooth $C^2$ domain. Moreover, we denote by $c(x,t) = (c_1(x,t), \ldots, c_N(x,t))$ the vector of concentrations, where $c_i(x,t)$ is the concentration of the specie $S_i$ at time $t > 0$ and position $x \in \Omega$. Each substance $S_i$ is assumed to diffuse in $\Omega$ with a strictly positive diffusion coefficient $d_i > 0$. The corresponding reaction-diffusion system reads as

$$\frac{\partial}{\partial t} c - \mathbb{D} \Delta c = R(c) \quad \text{for} \quad (x,t) \in \Omega \times \mathbb{R}_+,$$

where the diffusion matrix $\mathbb{D} = \text{diag}(d_1, \ldots, d_N)$ is positive definite (since $d_i > 0$ for all $i = 1, \ldots, N$). Moreover, by applying the law of mass action kinetics, the reaction vector $R(c)$ is modelled as

$$R(c) = \sum_{r=1}^{|\mathcal{R}|} k_r e^{\nu r}(y'_r - y_r), \quad \text{with} \quad e^{\nu r} := \prod_{i=1}^N e^{\nu_{r,i}},$$

where $k_r > 0$ denotes the rate constant of the $r$-th reaction. Finally, system (1.1) is associated to nonnegative initial data $c_0(x) \geq 0$ (by which we mean $c_0(x) := (c_{1,0}(x), \ldots, c_{N,0}(x))$ and $c_i(x) \geq 0$ for $i = 1, \ldots, N$), and homogeneous Neumann boundary condition

$$c(x,0) = c_0(x) \quad \text{for} \quad x \in \Omega, \quad \text{and} \quad \nabla c \cdot \nu = 0 \quad \text{for} \quad (x,t) \in \partial \Omega \times \mathbb{R}_+,$$

where $\nu := \nu(x)$ is the outward normal unit vector at point $x \in \partial \Omega$.

Concerning the set of chemical reactions $\mathcal{R}$, we denote by $W$ the Wegscheider matrix, i.e.

$$W = [(y'_r - y_r)_{r=1,...,|\mathcal{R}|}]^\top \in \mathbb{R}^{(|\mathcal{R}| \times N)}_+,\quad \text{and} \quad m = \text{dim}(\ker(W)).$$

Then, if $m > 0$, there exists a (non-unique) matrix $Q \in \mathbb{R}^{m \times N}$ such that

$$\text{rank}(Q) = m \quad \text{and} \quad Q R(c) = 0 \quad \text{for all} \quad c \in \mathbb{R}^N_{\geq 0} := (\mathbb{R}_+)^N, \quad (1.4)$$

so that the rows of $Q$ form a basis of $\ker(W)$. Thus, because of the homogeneous Neumann boundary condition, the use of the matrix $Q$ leads, at the formal level, to the following mass conservation laws for the solutions of eq. (1.1):

$$Q \overline{c}(t) = Q \overline{c}_0 =: M/|\Omega| \in \mathbb{R}^m \quad \text{for all} \quad t > 0,$$

where $\overline{c} := (\overline{c}_1, \ldots, \overline{c}_m)$ and $\overline{c}_r := \int_\Omega c_r \frac{dx}{|\Omega|}$, and $M$ is called the initial mass vector. By changing the signs of some rows of $Q$ if necessary, we can always consider (w.l.o.g.) a matrix $Q$ such that the initial mass vector $M$ is nonnegative, i.e. $M \in \mathbb{R}^m_{\geq 0}$.

To state our results, we require the following definitions concerning *equilibria* of chemical reaction networks.

**Definition 1.2 (Equilibria).**

Consider a chemical reaction network with a reaction vector $R$ defined by (1.2) and a mass vector $M \in \mathbb{R}^m_{\geq 0}$. Let $c_\infty := (c_{1,\infty}, \ldots, c_{N,\infty}) \in \mathbb{R}^N_{\geq 0}$ be such that $Q c_\infty = M$. Then,

- $c_\infty$ is called an equilibrium if $R(c_\infty) = 0$,
- $c_\infty$ is called a detailed balance equilibrium if for each forward reaction $y \xrightarrow{k_f} y'$ (with $k_f > 0$) in $\mathcal{R}$, there exists in $\mathcal{R}$ also the corresponding backward reaction $y' \xrightarrow{k_b} y$ (with $k_b > 0$) and

$$k_f c_\infty = k_b c_\infty.$$
• $c_\infty$ is called a complex balance equilibrium if for any complex $y \in C$, we have
\[
\sum_{y_i = y} k_i c^{y_i}_\infty = \sum_{y_i = y} k_i c^{y_i}_\infty.
\] (1.5)

Roughly speaking, a state $c_\infty$ is a complex balance equilibrium if and only if the total out- and inflow at $c_\infty$ are equal for every complex $y$.

• $c_\infty$ is called a boundary detailed/complex balance equilibrium (or shortly a boundary equilib-rium) if $c_\infty$ is a detailed/complex balance equilibrium and $c_\infty \in \partial \mathbb{R}_{\geq 0}$.

A chemical reaction network is called complex balanced if it possesses a strictly positive (i.e., not a boundary) complex balance equilibrium for each strictly positive mass vector $M \in \mathbb{R}^m_{>0}$.

It follows directly from the above definitions that $c_\infty$ is a detailed balance equilibrium $\Rightarrow c_\infty$ is a complex balance equilibrium $\Rightarrow c_\infty$ is an equilibrium, but the reverse is in general not true.

The principle of detailed balance goes back as far as Boltzmann for modelling collisions in kinetic gas theory and for proving the H-theorem for Boltzmann’s equation [Bol1896]. It was then applied to chemical kinetics by Wegscheider [Weg1901]. The complex balance condition was also considered by Boltzmann [Bol1887] under the name semi-detailed balance condition or cyclic balance condition, and was systematically used by Horn, Jackson and Feinberg in the seventies, see e.g. [Hor72, FH74].

It is by now a well-known fact for complex balanced chemical reaction networks that all equilibria (of such systems) are complex balanced (cf. [Hor72]). Moreover, for each strictly positive initial mass vector $M \in \mathbb{R}^m_{>0}$, there exists (for such systems) a unique strictly positive (that is, $c_\infty \in \mathbb{R}^N_{>0}$) complex balance equilibrium satisfying the mass conservation laws. One or several boundary equilibria (for such systems) can nevertheless exist (cf. [HJ72]).

There is also an extensive literature concerning the large time asymptotics of complex balanced systems in the ODE setting, i.e. by considering $c \equiv c(t)$, which satisfies the ODE system
\[
\frac{d}{dt} c = R(c),
\] (1.6)
where $R(c)$ is defined as (1.2). Indeed, it is proven that the unique strictly positive complex balance equilibrium of an ODE reaction network (of such a system) is locally stable (cf. [HJ72]). Moreover, it is conjectured that this equilibrium is in fact globally stable, i.e. that it is the unique global attractor for the dynamical system given by the ODE network (with exception of the boundary equilibria points). This statement is usually called the Global Attractor Conjecture and has remained one of the most important open problems in the theory of chemical reaction networks, see e.g. [And11, CNN13, GMS14, Pan12] and the references therein. A recently proposed proof of this conjecture in the ODE setting is currently under verification [Cra].

For the rest of the paper, we shall systematically consider only complex balance systems. We shall refer to the unique strictly positive complex balance equilibrium (of our systems) as “the” (strictly positive) complex balance equilibrium, while all the other complex balance equilibria (which are necessarily boundary equilibria) are simply named boundary equilibria.

The first part of this paper is devoted to the quantitative study of the convergence to equilibrium for complex balanced systems (1.1) of reaction diffusion PDEs without boundary equilibria, and is based on the so-called entropy method.

The idea of the entropy method consists in studying the large-time asymptotics of a dissipative PDE model by looking for a nonnegative (convex) Lyapunov functional $E(f)$ and its nonnegative dissipation
\[
\mathcal{D}(f(t)) \coloneqq -\frac{d}{dt} E(f(t)),
\] (1.7)
along the flow of the PDE model. We shall consider entropy dissipation functionals which are well-behaved in the sense that Firstly, all states satisfying $\mathcal{D}(f) = 0$ (and coherent with the conservation laws of the PDE) correspond to a unique (entropy-minimising) equilibrium $f_\infty$, i.e.
\[
\mathcal{D}(f) = 0 \quad \text{and coherence with conservation laws} \quad \iff \quad f = f_\infty,
\]
and secondly, there exists an entropy entropy-dissipation estimate of the form
\[
\mathcal{D}(f) \geq \Phi(E(f) - E(f_\infty)),
\] (1.8)
which holds for all $f$ coherent with the conservation laws of the PDE, and for some nonnegative function $\Phi$ satisfying $\Phi(x) = 0 \iff x = 0$. In this method, if $\Phi'(0) \neq 0$, one gets (at least formally) exponential
convergence toward \( f_\infty \) in relative entropy \( \mathcal{E}(f) - \mathcal{E}(f_\infty) \), with a rate given by some variant of Gronwall’s lemma.

It is well known in the ODE theory of complex balanced reaction systems (1.6) that the free energy

\[
E(c|c_\infty) = \sum_{i=1}^{N} \left( c_i \log \frac{c_i}{c_i,\infty} - c_i + c_i,\infty \right),
\]

(ODE entropy functional) (1.9)

is decreasing in time along the solutions of the ODE, i.e. \( \frac{d}{dt} E(c(t)|c_\infty) \leq 0 \) when \( c := c(t) \) satisfies eq. (1.6) (cf. [HJ72, Fei79, Gop13, And14]), and \( c_\infty \) is the complex balance equilibrium of the network. This suggests to consider as an entropy functional for the PDE system (1.1) the quantity

\[
E(c|c_\infty) = \sum_{i=1}^{N} \int_{\Omega} \left( c_i \log \frac{c_i}{c_i,\infty} - c_i + c_i,\infty \right) dx,
\]

(PDE entropy functional). (1.10)

It is remarkable, however, that the monotonicity of \( t \mapsto E(c|c_\infty) \) is in general shown indirectly, so that the explicit form of the entropy dissipation \( D(c) := -\frac{d}{dt} E(c|c_\infty) \geq 0 \) is not written down. In this paper, we therefore compute the explicit form of the entropy dissipation functionals \( D(c) = -dE/dt \) and \( \mathcal{D}(c) = -d\mathcal{E}/dt \), which is necessary for applying the entropy-entropy dissipation method. More precisely, we show that

\[
\mathcal{D}(c) := \sum_{i=1}^{N} d_i \int_{\Omega} |\nabla c_i|^2 c_i + \sum_{i=1}^{N} \int_{\Omega} \Psi \left( \frac{c_i'}{c_i^{\infty}} - \frac{c_i''}{c_i^{\infty}} \right) dx
\]

(1.11)

satisfies (1.7), where \( \mathcal{E} \) is defined by (1.10), when \( c \) is a (sufficiently integrable) solution of eq. (1.1). In (1.11), \( \Psi : (0, +\infty) \times (0, +\infty) \to \mathbb{R}_+ \) is defined by

\[
\Psi(x; y) := x \log(x/y) - x + y \quad (\geq 0),
\]

(1.12)

and it is easy to check that \( \Psi(x; y) = 0 \iff x = y \). We refer to Prop. 2.1 below for the verification that \( \mathcal{D} \) indeed satisfies (1.7).

Our first main result concerning complex balanced systems of reaction-diffusion PDEs without boundary equilibria is stated in Theorem 1.1 below. We show that all renormalised solutions, which satisfy a weak entropy entropy-dissipation law, converge exponentially fast to the unique strictly positive complex balance equilibria.

**Theorem 1.1** (Exponential convergence to equilibrium for complex balanced systems without boundary equilibria).

Let \( \Omega \) be a bounded smooth (\( C^2 \)) domain of \( \mathbb{R}^n \). Consider a (positive definite) diffusion matrix \( D = \text{diag}(d_1, \ldots, d_N) \) with \( d_i > 0 \) for \( i = 1, 2, \ldots, N \). Let \( R \) be a reaction term describing a complex balanced chemical reaction network without boundary equilibria (coming out of the mass action law, cf. (1.2)), and fix a strictly positive initial mass vector \( M \in \mathbb{R}_+^N \). We recall that such a network possesses a unique strictly positive equilibrium \( c_\infty \).

Then, there exists a constant \( \lambda > 0 \) such that for all nonnegative measurable functions \( c = (c_1, \ldots, c_N) : \Omega \to \mathbb{R}^N_+ \) satisfying the mass conservation laws \( Q\tau = M \) (cf. (1.4) for the definition of \( Q \)) and the upper bound \( \mathcal{E}(\tau|c_\infty) \leq K \), the following entropy-entropy dissipation inequality holds:

\[
\mathcal{D}(c) \geq \lambda E(c|c_\infty),
\]

(1.13)

where \( E(c|c_\infty) \) and \( \mathcal{D}(c) \) are defined in (1.10) and (1.11), respectively, and \( \lambda \) solely depends on \( \Omega, d_i, M \), the coefficients appearing in \( R \), and an upper bound \( K \) on the relative entropy \( \mathcal{E}(\tau|c_\infty) \) of \( \tau \) w.r.t. \( c_\infty \).

Moreover, following [Fis15], for any nonnegative initial data with finite initial mass and relative entropy, i.e. \( c_0 \in (L^1(\Omega))^N \) and \( E(c_0|c_\infty) < +\infty \), the system (1.1)–(1.3) possesses a global renormalised solution \( c(x, t) = (c_1(x, t), \ldots, c_N(x, t)) \) (cf. Rmk. 1.1 for a precise definition of this solution concept).

Then, each such renormalised solution which moreover satisfies the weak entropy entropy-dissipation law

\[
E(c|c_\infty)(t) + \int_{t}^{\infty} \mathcal{D}(c)(r)dr \leq E(c|c_\infty)(s), \quad a.a. \quad t \geq s > 0,
\]

(1.14)

converges in \( L^1 \)-norm exponentially fast to the (unique, strictly positive) complex balance equilibrium. That is, for a constant \( C > 0 \) depending on the same parameters as \( \lambda \) and on \( c_0 \), there holds

\[
\sum_{i=1}^{N} ||c_i(t) - c_i,\infty||_{L^1(\Omega)}^2 \leq C e^{-\lambda t}, \quad \forall t > 0.
\]

(1.15)
Remark 1.1. The statement of Thm. 1.1 is formulated for renormalised solutions, which is the only available concept of global solutions for general systems (1.1)–(1.3). We recall here that a vector of measurable nonnegative components \( \mathbf{c}(x,t) = (c_1(x,t), \ldots, c_N(x,t)) \) is said to be a renormalised solution when \( c_i \in L^\infty_{\text{loc}}(\mathbb{R}_+; L^1(\Omega)) \), \( \sqrt{T} \in L^2_{\text{loc}}(\mathbb{R}_+; H^1(\Omega)) \), and when for any smooth function \( \xi : \mathbb{R}^N_+ \rightarrow \mathbb{R} \) with compactly supported derivative \( \partial^\alpha \xi \) and every \( \psi \in C^\infty(\overline{\Omega} \times \mathbb{R}_+) \), the equation

\[
\int_\Omega \xi(\mathbf{c}(\cdot,T))\psi(\cdot,T) \, dx - \int_\Omega \xi(\mathbf{c}_0)\psi(\cdot,0) \, dx - \int_0^T \int_\Omega \xi(\mathbf{c}) \frac{d}{dt} \psi \, dx dt \\
= - \sum_{i,j=1}^N \int_0^T \int_\Omega \psi \partial_j \xi(\mathbf{c})(d_i \nabla c_i) \cdot \nabla c_j \, dx dt \\
- \sum_{i=1}^N \int_0^T \int_\Omega \partial_i \xi(\mathbf{c})(d_i \nabla c_i) \cdot \nabla \psi \, dx dt + \sum_{i=1}^N \int_0^T \int_\Omega \partial_i \xi(\mathbf{c}) R_i(\mathbf{c}) \psi \, dx dt
\]  

(1.16)

holds for almost every \( T > 0 \), with \( \mathbf{R}(\mathbf{c}) = (R_1(\mathbf{c}), \ldots, R_N(\mathbf{c})) \).

We notice that the assumed weak entropy entropy-dissipation law (1.14), which holds at the formal level, is in fact not easy to prove in general for renormalised solutions, because of the lacking integrability of the reaction terms in the entropy dissipation functional (1.11). In Thm. 1.2 below, exponential equilibration of renormalised solutions (without any extra assumptions) in the case of a cyclic reaction network is obtained thanks to estimates which are uniform w.r.t. an approximation process.

For classical and sufficiently integrable weak solutions of (1.1)–(1.3), however, the entropy entropy-dissipation law (1.14) can be verified rigorously (even with an equality sign). This was done, for instance, in [DFP07] for weak \((L \log L)^2\)-solutions of a system with quadratic nonlinearities. More recently, sufficiently integrable weak solutions were shown to exist for some special systems of the form (1.1)–(1.3) when a (space dimension-dependent) “closeness” assumption on the diffusion coefficients of the system is made, see e.g. [CDF14, FL16]. Imposing an even stronger “closeness” assumption on the diffusion coefficients allows in fact to show the existence of classical solutions, see e.g. [FLS16].

Remark 1.2. In the above theorem, the constant \( C \) in (1.15) can be explicitly estimated. In fact, one can take \( C = C_{\text{CKP}}^{-1} \mathcal{E}(c_0|c_\infty) \), where \( C_{\text{CKP}} \) is the constant in a Csiszár-Kullback-Pinsker inequality (cf. Lemma 2.6 below).

It is however not possible with our method used in Thm. 1.1 to estimate \( \lambda \) in a completely explicit way. It is nevertheless possible to define the constant \( \lambda \) by a finite-dimensional minimisation problem (cf. the proof of Thm. 1.1). In particular, one needn’t use any abstract infinite-dimensional compactness argument.

The proof of Thm. 1.1 exploits an additivity property of the entropy functional, the Logarithmic Sobolev inequality (allowing to avoid the use of \( L^\infty \)-bounds, which are unavailable for general systems) and then uses a convexification argument presented in [MHH15], which allows to further reduce the functional inequality (1.8) into a finite dimensional inequality. The proof of this finite-dimensional inequality is based on Taylor expansions around the equilibrium (unfortunately not yielding an explicit bound for \( \lambda \), since the actual value of \( \lambda > 0 \) might be obtained at states far from equilibrium).

The second main result on complex balanced systems without boundary equilibria is stated in Theorem 1.2 below. Since the method of convexification used in Theorem 1.1 does not yield explicit estimates for the convergence rate \( \lambda \) (at least for general systems), we also propose a constructive method, inspired by [DF08, FL16, FT], to prove an entropy entropy-dissipation estimate with computable constants. Another advantage of this method is that it can be applied uniformly w.r.t. the approximating systems used for constructing renormalised solutions, and thus yields exponential convergence to equilibrium of renormalised solutions without having to assume the weak entropy entropy-dissipation law (1.14) as in Thm. 1.1.

The proposed method consists of four steps. The first three are proven for general complex balanced systems, but the last one, which crucially depends on the structure of the conservation laws, is rather a proof of concept which can be detailed only once a specific system is given. We demonstrate the details of this last step for the specific case of cycles of reactions connecting an arbitrary number of chemical substances:

\[
\alpha_1 A_1 \xrightarrow{k_1} \alpha_2 A_2 \xrightarrow{k_2} \cdots \xrightarrow{k_{N-1}} \alpha_N A_N \xrightarrow{k_N} \alpha_1 A_1,
\]
where \( \alpha_i \in \mathbb{N}_{>0} \), and \( k_i > 0 \) for \( i = 1, 2, \ldots, N \). The corresponding reaction-diffusion system writes as

\[
\begin{align*}
\partial_t a_1 &- d_1 \Delta a_1 = a_1(-k_1 a_1^\alpha_1 + k_N a_N^\alpha_N), & x \in \Omega, & t > 0, \\
\partial_t a_2 &- d_2 \Delta a_2 = a_2(-k_2 a_2^\alpha_2 + k_1 a_1^\alpha_1), & x \in \Omega, & t > 0, \\
\vdots & & & \\
\partial_t a_N &- d_N \Delta a_N = a_N(-k_N a_N^\alpha_N + k_{N-1} a_{N-1}^\alpha_{N-1}), & x \in \Omega, & t > 0,
\end{align*}
\]

(1.17)

 altogether with homogeneous Neumann boundary conditions and initial data (1.3) (with \( c \) replaced by \( a := (a_1, \ldots, a_N) \)). The system (1.17) has one mass conservation law corresponding to \( Q = \left( \frac{1}{\alpha_1}, \ldots, \frac{1}{\alpha_N} \right) \), i.e.

\[
\sum_{i=1}^N \frac{1}{\alpha_i} \int_\Omega a_i(x,t)dx = M := \sum_{i=1}^N \frac{1}{\alpha_i} \int_\Omega a_{i,0}(x)dx > 0 \quad \text{for all} \quad t > 0,
\]

(1.18)

where the initial total mass \( M \) is assumed to be strictly positive. As a consequence, the system (1.17) has a unique strictly positive complex balance equilibrium \( a_\infty = (a_1, \ldots, a_N) \) (cf. Lemma 2.7).

By denoting \( a = (a_1, a_2, \ldots, a_N) \), \( a_0 = (a_{1,0}, \ldots, a_{N,0}) \) and by using the periodic notation \( a_{N+1} := a_1 \), \( a_{N+1,\infty} := a_1, \infty \) and \( a_{N+1,0} := a_{1,0} \), the quantities (1.10) and (1.11) write in this specific case as

\[
\mathcal{E}(a|a_\infty) = \sum_{i=1}^N \int_\Omega \left( a_i \log \frac{a_i}{a_i,\infty} - a_i + a_i,\infty \right) dx,
\]

(1.19)

and

\[
\mathcal{D}(a) = \sum_{i=1}^N d_i \int_\Omega \frac{\nabla a_i^2}{a_i} + \sum_{i=1}^N k_i a_i^\alpha_i \sum_{\alpha_i} \int_\Omega \Psi \left( \frac{a_i^\alpha_i}{a_i,\infty^\alpha_i}, \frac{a_{i+1}^\alpha_i}{a_{i+1,\infty}^\alpha_i} \right) dx,
\]

(1.20)

where we recall that \( \Psi \) is defined by (1.12).

We now state the

**Theorem 1.2** (Explicit convergence to equilibrium for a cyclic reaction).

Let \( \Omega \) be a bounded smooth \((C^2)\) domain of \( \mathbb{R}^n \), \( d_i > 0 \), \( \alpha_i \in \mathbb{N}_{>0} \), and \( k_i > 0 \) for all \( i = 1, 2, \ldots, N \). We finally fix a strictly positive initial mass \( M > 0 \).

Then, there exists a constant \( \lambda > 0 \) which can be explicitly estimated in terms of \( \Omega, d_i, M, \alpha_i \) and \( k_i \), such that for any nonnegative measurable functions \( a_i : \Omega \rightarrow \mathbb{R}_+ \) \((i = 1, 2, \ldots, N)\) satisfying the mass conservation law \( \sum_{i=1}^N \frac{1}{\alpha_i} \int_\Omega a_i(x)dx = M \), we have

\[
\mathcal{D}(a) \geq \lambda \mathcal{E}(a|a_\infty),
\]

where \( a_\infty \) is the unique strictly positive complex balance equilibrium determined by \( M \), and \( \mathcal{E}(a|a_\infty) \) and \( \mathcal{D}(a) \) are defined in (1.19) and (1.20), respectively.

As a consequence, for any nonnegative initial data \( a_0 \) with positive mass and finite relative entropy, i.e. \( \sum_{i=1}^N \frac{1}{\alpha_i} \int_\Omega a_{i,0}(x)dx = M > 0 \) and \( \mathcal{E}(a_0|a_\infty) < +\infty \), the renormalised solutions of system (1.17) (as constructed in [Fis15] and defined in the same manner as in (1.16)), (1.3) (with \( c \) replaced by \( a \)) converge exponentially fast in \( L^1 \)-norm to the equilibrium \( a_\infty \):

\[
\sum_{i=1}^N \|a_i(t) - a_i,\infty\|_{L^1(\Omega)}^2 \leq C_{CKP}^{-1} e^{-\lambda t} \mathcal{E}(a_0|a_\infty) \quad \text{for all} \quad t > 0,
\]

where \( C_{CKP} \) is defined in Lemma 2.6 below.

**The second part of this paper**, presented in Section 3, is devoted to the study of complex balanced systems featuring boundary equilibria. We recall that even in the ODE setting (1.6), the convergence to the unique strictly positive complex balance equilibrium for general reaction networks with boundary equilibria involves the study of the Global Attractor Conjecture.

Moreover, we emphasise that due to the presence of boundary equilibria, we cannot expect an entropy dissipation rate similar to that in Thm. 1.1. Instead, it is easy to verify (see also Section 3) that the entropy dissipation \( \mathcal{D}(c) \) tends to zero for sequences of constant states, which converge to a boundary equilibrium and satisfy the mass conservation laws, yet the relative entropy to the positive complex balance equilibrium \( \mathcal{E}(c|c_\infty) \) remains, of course, far from 0 for such states.
In Section 3, we prove two quite general results concerning the convergence to the unique strictly positive complex balance equilibrium for the following two reaction-diffusion networks with boundary equilibria:

\[
\begin{align*}
2A & \xrightarrow{k_1/k_2} A + B \\
2B & \xrightarrow{k_3/k_2} B + C
\end{align*}
\]

The first system is described by the following 2 × 2 mass action law reaction-diffusion system

\[
\begin{aligned}
a_t - d_a \Delta a &= -a^2 + ab, & a \in \Omega, \quad t > 0, \\
b_t - d_b \Delta b &= a^2 - ab, & b \in \Omega, \quad t > 0, \\
\nab \cdot \nabla a &= \nabla b \cdot \nu = 0, & a \in \partial \Omega, \quad t > 0, \\

a(x, 0) &= a_0(x), & b(x, 0) &= b_0(x), & x \in \Omega,
\end{aligned}
\]

(1.21)

where we assume normalised reaction rate constants \( k_1 = k_2 = 1 \) (without loss of generality since different rates can be recovered thanks to a suitable scaling).

It is easily checked that the above 2 × 2 system is complex balanced, and even detailed balanced, but features a boundary equilibrium. We nevertheless are able to prove for this system the

**Proposition 1.3** (Convergence to detailed balance equilibrium for a 2 × 2 system with boundary equilibrium).

Let \( \Omega \) be a bounded smooth \((C^2)\) domain of \( \mathbb{R}^n \) and \( d_a, d_b > 0 \). Assume that the initial data \((a_0, b_0) \) is \( L^\infty(\Omega) \times L^\infty(\Omega) \) satisfying 0 \( < \varepsilon^2 \leq a_0(x), b_0(x) \leq \Lambda < +\infty \) for a.a. \( x \in \Omega \).

Then, the unique global classical \((C^2)\) solution \((a, b)\) to (1.21) conserves the total mass

\[
\int_\Omega (a(x, t) + b(x, t)) \, dx = M := \int_\Omega (a_0(x) + b_0(x)) \, dx > 0 \quad \text{for all} \quad t > 0,
\]

and converges exponentially fast to the unique strictly positive detailed balance equilibrium \((a_\infty, b_\infty)\):

\[
\|a(\cdot,t) - a_\infty\|_{L^2(\Omega)}^2 + \|b(\cdot,t) - b_\infty\|_{L^2(\Omega)}^2 \leq e^{-\lambda_\varepsilon t} \|a_0 - a_\infty\|_{L^2(\Omega)}^2 + \|b_0 - b_\infty\|_{L^2(\Omega)}^2 \quad \text{for all} \quad t > 0,
\]

where \( \lambda_\varepsilon \) is an explicit constant depending only on \( \Omega, M, d_a, d_b, \) and \( \varepsilon \). Moreover, one can choose \( \lambda_\varepsilon = O(\varepsilon^2) \) as \( \varepsilon \to 0 \).

This proposition can be used to get a result in the case when the initial data are not bounded below:

**Corollary 1.4**. Let \( \Omega, d_a, d_b \) be as in Prop. 1.3. Assume that the non-negative initial data \((a_0, b_0)\) belong to \( C(\Omega) \times C(\Omega) \), and that \( a_0 \) is not the trivial initial state 0.

Then, the unique global classical \((C^2)\) solution \((a, b)\) to (1.21) conserves the total mass (as in Prop. 1.3) and converges exponentially fast to the unique strictly positive detailed balance equilibrium, that is, for some \( C, \Lambda > 0 \),

\[
\|a(\cdot,t) - a_\infty\|_{L^2(\Omega)}^2 + \|b(\cdot,t) - b_\infty\|_{L^2(\Omega)}^2 \leq C e^{-\lambda_\varepsilon t} \quad \text{for all} \quad t > 0.
\]

Note that on the other hand, if \( a_0 \equiv 0 \) on \( \Omega \) (so that the initial total mass is concentrated in \( b_0 \)), then the classical solutions to (1.21) converge exponentially fast to the boundary equilibrium \((0, M)\).

In the proof of Prop. 1.3, the particular structure of system (1.21) allows to show via a comparison principle argument that solutions propagate the assumed positive lower and upper \( L^\infty \)-bounds of the initial data. As a consequence, the solutions remain bounded away from the boundary equilibrium. Also the proof of Cor. 1.4 applies the minimum principle. However, such comparison principle arguments (like the propagation of lower and upper \( L^\infty \)-bounds) are no longer true for more general systems such as the second example (described in the following).

Applying the mass action law to the second example leads to the below 3 × 3 reaction-diffusion system

\[
\begin{aligned}
a_t - d_a \Delta a &= -k_1 a + k_3 b^2, & a \in \Omega, \quad t > 0, \\
b_t - d_b \Delta b &= k_1 a - 2k_3 b^2, & b \in \Omega, \quad t > 0, \\
c_t - d_c \Delta c &= k_1 a - 2k_2 c, & c \in \Omega, \quad t > 0, \\
\nab \cdot \nabla a &= \nabla b \cdot \nu = \nabla c \cdot \nu = 0, & a \in \partial \Omega, \quad t > 0, \\

a(x, 0) &= a_0(x), & b(x, 0) &= b_0(x), & c(x, 0) &= c_0(x), & x \in \Omega.
\end{aligned}
\]

(1.22)
For system (1.22), we first address the existence of global classical solution. Due to the lack of comparison principle arguments (which held for the $2 \times 2$ system), we are able to show the existence of global classical solutions in dimensions $N \leq 5$, or, when $N \geq 6$, when the diffusion coefficients are assumed to be sufficiently close to each other (see Lemma 3.2). The $3 \times 3$ system (1.22) also features a boundary equilibrium. Since no uniform-in-time upper or lower positive a priori $L^\infty$-bounds can be obtained for system (1.22), the question of instability of the boundary equilibrium turns out be significantly more tricky.

We can nevertheless prove the

**Theorem 1.5** (Convergence to complex balance equilibrium for a $3 \times 3$ system with boundary equilibrium).

Let $\Omega$ be a bounded smooth ($C^2$) domain of $\mathbb{R}^n$, and $d_a, d_b, d_c > 0$, $k_1, k_2, k_3 > 0$. Assume that the initial data $(a_0, b_0, c_0) \in (L^\infty(\Omega))^3$ are such that $b_0$ is a.e. bounded below by a strictly positive constant, i.e. $\|a_0\|_{L^\infty(\Omega)} < \infty$.

Then, there exist explicit constants $C > 0$ and $\lambda > 0$ such that any classical solutions of (1.22) converge exponentially fast to the unique strictly positive complex balance equilibrium $(a_\infty, b_\infty, c_\infty)$ with the rate $\lambda$, that is

$$\|a(t) - a_\infty\|_{L^1(\Omega)} + \|b(t) - b_\infty\|_{L^1(\Omega)}^2 + \|c(t) - c_\infty\|_{L^1(\Omega)}^2 \leq C e^{-\lambda t}, \quad \text{for all } t > 0,$$

where the constant $C$ depends only on the initial relative entropy and $C_{CKP}$, and $\lambda$ depends only on $\Omega, M, d_a, d_b, d_c, k_1, k_2, k_3$ and $\|b_0\|_{L^\infty(\Omega)}$.

In order to prove this result, we first observe that solutions to system (1.22), which initially satisfy $\|A(0)\|_{L^\infty(\Omega)} < \infty$, obey a specific lower bound of the form $\|A(t)\|_{L^\infty(\Omega)} \leq C (1 + t)$. This lower bound shows that the dissipation of the relative entropy (with respect to the unique strictly positive equilibrium $(a_\infty, b_\infty, c_\infty)$) decays at most like $(1 + t)^{-1}$ times the relative entropy.

As a consequence, we can show algebraically fast convergence to $(a_\infty, b_\infty, c_\infty)$ via a Gronwall argument. The essence of this Gronwall argument is illustrated by the following contradiction argument: Suppose that the relative entropy to $(a_\infty, b_\infty, c_\infty)$ would remain bounded below by a strictly positive constant uniformly-in-time, then the entropy dissipation would also remain bounded below by a positive constant times $(1 + t)^{-1}$. The function $(1 + t)^{-1}$ is however non-integrable over the interval $[0, +\infty)$ and thus yields a contradiction to the fact that the time-integral of the entropy dissipation over $[0, +\infty)$ is bounded by the initial relative entropy.

Moreover in a second step, the obtained algebraic convergence allows to recover exponential convergence towards the strictly positive complex balance equilibrium $(a_\infty, b_\infty, c_\infty)$. In fact, the rate of the convergence can be estimated explicitly as in Thm. 1.2 (see Rmk. 3.5).

**Remark 1.3** (Convergence to boundary equilibria). The condition on $b_0$ in Thm 1.5 is clearly stronger than the assumption that the nonnegative initial datum $(a_0, b_0, c_0)$ is different from the boundary equilibrium. In fact, there is a class of initial data (strictly larger than the boundary equilibrium) for which solutions converge to the boundary equilibrium, see Rmk. 3.4. Such a class of initial data was also presented in Cor. 1.4 for the $2 \times 2$ system. Moreover, Cor. 1.4 allowed to precisely classify for continuous initial data, which solutions converge to the boundary equilibrium and which solutions converge to the positive complex balance equilibrium.

**Remark 1.4.** We would like to point out a significant difference between our method here and the linearisation techniques: when using linearisation, one needs to wait until the trajectory is trapped into a small enough neighbourhood of $c_\infty = (a_\infty, b_\infty, c_\infty)$. Our method, however, already allows to use the standard entropy method (and obtain exponential convergence to $c_\infty$) as soon as the solution is not close to the boundary equilibria.

Note also that the arguments used in the $3 \times 3$ system are different from the ones used in the so-called “slowly growing a priori bounds” variant of the entropy method (cf. [ToVi], [DF08]).

The main novelties of this work are the following:

(i) Up to our knowledge, this paper seems to be the first quantitative study of convergence to equilibrium for nonlinear complex balanced reaction-diffusion systems.

Previous related results on the large time behaviour of complex balanced systems in the ODE setting stated mostly qualitative results, that is convergence to equilibrium, but without
convergence rates, see e.g. [And11, CNN13, GMS14, Pan12] and references therein. The only quantitative results for ODE complex balanced systems known to us were proven by [SJ], in which the authors obtained convergence rates close to equilibrium via a linearisation technique.

The PDE setting (1.1) for complex balanced systems is even less studied. Concerning the convergence to equilibrium for detailed balanced systems in the context of semi-conductor models, we refer to [Grö92, GGH96, GH97], where convergence rates were proven, but with non-constructive constants coming out of an abstract compactness argument. In [DF06], the authors obtained the first results of convergence to equilibrium for nonlinear detailed balance reaction-diffusion systems with explicit rates and constants. Related results were later derived in e.g. [DF08, GZ10, DF14, MHM15, FT]. We point out that all these references consider reaction-diffusion systems which satisfy a detailed balance condition. Finally, a recent result on the quantitative convergence to equilibrium was proven in [FPT] in the special case of linear complex balanced systems for PDEs (coming out of first order chemical reaction networks).

(ii) Secondly, we study two specific complex balanced systems featuring boundary equilibria.

One of the main difficulties in proving the Global Attractor Conjecture for ODEs is the appearance of boundary equilibria, leading to the possibility of \( \omega \)-limit sets having nonempty intersections with the boundary \( \partial \mathbb{R}^N_+ \). Resolving this problem in the PDE setting (1.1) is, up to the best of our knowledge, completely open. Here, we investigate two particular systems featuring boundary equilibria and study in which modified sense the GAC could be investigated: First, we note that in the PDE setting, there exist nontrivial classes of solutions which converge to the boundary equilibrium. Secondly, we point out the observation that for some specific examples of complex balanced systems, the nonlinear reaction terms which gives rise to boundary equilibria, also lead to specifically decaying \textit{a priori} lower bounds which (in combination with an entropy dissipation argument) are sufficient to prove instability of those boundary equilibria.

\textbf{Outline:} The rest of the paper is organised as follows: In Section 2 contains the results concerning complex balanced systems without boundary equilibria. The two specific systems featuring boundary equilibria are then analysed in Section 3.

\textbf{Notations:} With a slight abuse of notation, we recall the following convention whenever a single letter \( y \) is used for a complex, then \( y_i \) (for \( i = 1, \ldots, N \)) denotes the stoichiometric coefficient of the \( i \)-th specie \( S_i \) in the complex \( y \). If the complex is written as \( y_r \) (or \( y_r' \)), then it denotes the source (or target) complex of the \( r \)-th reaction, and the \( i \)-th stoichiometric coefficient is denoted by \( y_{r,i} \) (or \( y'_{r,i} \)).

- For \( c, y \in \mathbb{R}^N \), we denote
  \[
  e^c = \prod_{i=1}^N c_i^{y_i} \quad \text{and} \quad \frac{c}{y} = \left( \frac{c_1}{y_1}, \frac{c_2}{y_2}, \ldots, \frac{c_N}{y_N} \right) \in \mathbb{R}^N.
  \]
  The scalar product is denoted by
  \[
  c \cdot y = \sum_{i=1}^N c_i y_i.
  \]
- For a function \( f : \mathbb{R} \to \mathbb{R} \) and \( c \in \mathbb{R}^N \), we denote
  \[
  f(c) = (f(c_1), f(c_2), \ldots, f(c_N)) \in \mathbb{R}^N.
  \]
- We denote by \( \| \cdot \| \) the usual norm in \( L^2(\Omega) \).
- We can assume (w.l.o.g. by rescaling the space variable \( x \)) that \( \Omega \) has normalised volume
  \[
  |\Omega| = 1.
  \]

Note that we shall systematically make this assumption in all the following proofs of the theorems concerning PDEs.

- In Section 2, we shall use the notations \( c, \bar{c}, c_\infty, \) etc. whenever a result is proven for general complex balanced systems without boundary equilibrium. On the other hand, for results specific to a cycle of reactions and systems (1.17), we shall use the notations \( a, \bar{a}, a_\infty, \) etc. instead. In this case, it is also convenient to use the periodic notation \( a_{N+1} := a_1 \) or \( a_{N+1,\infty} := a_{1,\infty} \) e.t.c. Moreover, we shall introduce capital letters as a short hand notation for square roots of quantities, e.g. \( c_i := \sqrt{c_i}, \bar{c}_i := \sqrt{\bar{c}_i}, c_{i,\infty} := \sqrt{c_{i,\infty}}, A_i := \sqrt{a_i}, \) etc.
2. Complex balanced systems without boundary equilibria

Before beginning the proof of Thm. 1.1, we first state and prove a proposition devoted to the study of ODEs coming out of a chemical reaction network, since the tools developed in this proof can then be used in the study of the PDEs.

2.1. Convergence to equilibrium for ODE systems.

Throughout this paper, we frequently use the two variables function \( \Psi \) defined by (1.12), and the elementary estimate

\[
\Psi(x, y) \geq (\sqrt{x} - \sqrt{y})^2.
\]  

(2.1)

It is well known for reaction terms \( R \) describing complex balanced chemical networks without boundary equilibrium that solutions \( c := c(t) \) of the ODE system (1.6) (with nonnegative initial data \( c_0 \geq 0 \) corresponding to a strictly positive mass vector \( M \in \mathbb{R}^m_{>0} \)) converge towards the unique strictly positive complex balance equilibrium \( c_\infty \). Moreover, the relative entropy functional

\[
E(c(t)|c_\infty) := \sum_{i=1}^N \Psi(c_i(t); c_{i, \infty}) = \sum_{i=1}^N \left( c_i(t) \log \frac{c_i(t)}{c_{i, \infty}} - c_i + c_{i, \infty} \right)
\]  

(2.2)

is decreasing along solution trajectories (see e.g. [HJ72, Fei79, And14, Gop13]). The corresponding entropy dissipation, however, seems not to be written down. We provide the following proposition to establish the explicit expression of the entropy dissipation:

**Proposition 2.1.** We consider a reaction rate defined by (1.2) and \( c_\infty \) a strictly positive complex balance equilibrium of a chemical reaction network (cf. Defs. 1.1 and 1.2). Then,

\[
D(c) := -R(c) \cdot \log \frac{c}{c_\infty} \geq \sum_{r=1}^{|R|} k_r \Psi \left( \frac{e^{y_r} c_{\infty}^{y_r}}{c_r^{y_r}}, \frac{e^{y_r} c_r^{y_r}}{c_{\infty}^{y_r}} \right) \geq 0.
\]  

(2.3)

Moreover \( D(c_\ast) = 0 \) for some \( c_\ast \in \mathbb{R}_{\geq 0}^N \) if and only if \( c_\ast \) is a complex balanced equilibrium (that is, (1.5) holds). Finally, if the chemical reaction network has no boundary equilibria, then \( c_\ast = c_\infty \).

**Proof.** We compute

\[
-R(c) \cdot \log \frac{c}{c_\infty} = \left( \sum_{r=1}^{|R|} k_r e^{y_r} (y_r - y'_r) \right) \cdot \log \frac{c}{c_\infty} = \sum_{r=1}^{|R|} k_r e^{y_r} \log \frac{e^{y_r} c_r^{y_r}}{c_\infty^{y_r}} - \sum_{r=1}^{|R|} k_r e^{y_r} - k_r e^{y_r} e^{y_r} c_r^{y_r} c_\infty^{y_r}
\]  

(2.4)

It remains to prove that

\[
\sum_{r=1}^{|R|} k_r e^{y_r} - k_r e^{y_r} e^{y_r} c_r^{y_r} c_\infty^{y_r} = 0.
\]  

(2.5)

Indeed, by using the properties of Def. 1.1,

\[
\sum_{r=1}^{|R|} k_r e^{y_r} - k_r e^{y_r} e^{y_r} c_r^{y_r} c_\infty^{y_r} = \sum_{y \in \mathcal{C}} \left( \sum_{\{r : y_r = y\}} k_r e^{y_r} - \sum_{\{s : y'_s = y\}} k_s e^{y'_s} e^{y'_s} c_s^{y'_s} c_\infty^{y'_s} \right)
\]  

(2.6)

where we have used the complex balance condition (1.5) in the last step.

Now assume that \( D(c_\ast) = 0 \) for some \( c_\ast \in \mathbb{R}_{\geq 0}^N \). Since \( \Psi(x, y) \geq 0 \) for all \( x \geq 0, y > 0 \) and \( \Psi(x, y) = 0 \) if and only if \( x = y \), it follows from \( D(c_\ast) = 0 \) that

\[
\frac{c_r^{y_r}}{c_\infty^{y_r}} = \frac{c_s^{y'_s}}{c_\infty^{y'_s}}, \text{ or equivalently } \frac{c_s^{y'_s}}{c_\infty^{y'_s}} = c_r^{y_r}.
\]  

(2.7)
for all \( r = 1, \ldots, |\mathcal{R}| \). Thus for any \( y \in \mathcal{C} \), we have
\[
\sum_{\{r: y_r = y\}} k_r c_r^{y_r} = \frac{e^{\theta r}}{e^\infty} \sum_{\{r: y_r = y\}} k_r c_r^{y_r} = \frac{e^{\theta r}}{e^\infty} \sum_{\{s: y_s' = y\}} k_s c_s^{y_s'} = \sum_{\{s: y_s' = y\}} k_s c_s^{y_s'}.
\]
Therefore, \( c_* \) is a complex balance equilibrium due to (1.5). Furthermore, if the considered chemical reactions network has no boundary equilibria, then obviously \( c_* \equiv c_\infty \).

**Remark 2.1 (Explicit Entropy-Dissipation).**
In the case of linear reaction networks, a recent result [FPT] shows explicitly that the relative entropy between any two solutions decays. Such a strong results does not seem true for nonlinear reaction networks, for which we suspect that only the relative entropy (2.2) is a Lyapunov functional in general. For linear systems, the existence of a dissipative relative entropy between any two solutions is due to a fundamental property of linear systems, see e.g. [FJ16]. It can also be derived from the General Entropy Principle for finite Markov chains, see [PMP06].

Thanks to the explicit expression of the entropy dissipation in Prop. 2.1, we are now in position to prove the following quantitative result about the decay towards equilibrium of solutions of the ODE systems (1.6) (for complex balanced chemical reactions networks without boundary equilibria).

**Proposition 2.2 (Exponential convergence to equilibrium for ODE systems without boundary equilibria).**
We consider a complex balanced chemical reaction network without boundary equilibria, the corresponding reaction term (1.2) and, for a given mass vector \( M \in \mathbb{R}_+^N \), the unique strictly positive complex balanced equilibrium \( c_\infty \) together with the ODE system (1.6).

Then, there exists a strictly positive constant \( \lambda > 0 \) depending only on the coefficients appearing in \( \mathbf{R} \), such that any solution \( c := c(t) \) to the ODE system (1.6) (with nonzero nonnegative initial data) satisfies the following exponentially fast decay of the relative entropy:
\[
E(c(t) | c_\infty) \leq E(c_0 | c_\infty) e^{-\lambda t} \quad \text{for all} \quad t > 0.
\]
As a consequence, these solutions \( c \) converge exponentially fast to the unique strictly positive complex balance equilibrium \( c_\infty \):
\[
\sum_{i=1}^N |c_i(t) - c_{i,\infty}|^2 \leq CE(c_0 | c_\infty) e^{-\lambda t} \quad \text{for all} \quad t > 0,
\]
where \( C = (\sqrt{K} + \max\{\sqrt{c_i, \infty}\})^2 \) and \( K \) is defined as in (2.22).

**Proof.** We first observe that
\[
- \frac{d}{dt} E(c|c_\infty) = - \sum_{i=1}^N \left( \frac{d}{dt} c_i \right) \log \frac{c_i}{c_i,\infty} = - \mathbf{R}(c) \cdot \log \frac{c}{c_\infty},
\]
so that thanks to (2.3),
\[
- \frac{d}{dt} E(c|c_\infty) = D(c) = \sum_{r=1}^{|\mathcal{R}|} k_r c_r^{y_r} \psi \left( \frac{e^{\psi_r}}{e^\infty} \frac{e^{y_r}}{e^\infty} \right).
\]
Thus, if we are able to show the following **entropy entropy-dissipation estimate**
\[
D(c(t)) \geq \lambda E(c(t)|c_\infty),
\]
for some \( \lambda > 0 \) and for all \( t > 0 \), we see that (2.8) will follow from a Gronwall argument. We consider
\[
\mathcal{C}_M = \{ \xi \in \mathbb{R}^N : \mathbb{Q} \xi = M \quad \text{and} \quad E(\xi|c_\infty) \leq E(c_0|c_\infty) \}.
\]
Thanks to (2.11), we see that \( c(t) \in \mathcal{C}_M \) for all \( t > 0 \). Therefore, we will prove the estimate (2.12) thanks to the

**Lemma 2.3.** Let \( E(\cdot|c_\infty) \), \( D(\cdot) \) and \( \mathcal{C}_M \) be defined in (2.2), (2.3) and (2.13) respectively. Then
\[
\lambda := \inf_{\xi \in \mathcal{C}_M} \frac{D(\xi)}{E(\xi|c_\infty)} > 0.
\]
Proof. Firstly, since \(D(\xi) \geq 0\) and \(E(\xi|c_\infty) \geq 0\), we see that \(\lambda \geq 0\). Secondly, it follows from definition (2.13) of \(E_M\) that the denominator \(E(\xi|c_\infty)\) is bounded above. On the other hand, thanks to Prop. 2.1, \(D(\xi) = 0\) if and only if \(\xi = c_\infty\). Thus, \(\lambda\) can approach zero only in cases when \(\xi \to c_\infty\) on \(E_M\) (where we recall that \(c_\infty\) is strictly positive). Therefore, in order to prove (2.14), it is sufficient to show that

\[
\liminf_{\xi \to c_\infty} \frac{D(\xi)}{E(\xi|c_\infty)} > 0.
\]

(2.15)

Thanks to \(D(c_\infty) = 0\), \(\nabla D(c_\infty) = 0\), and

\[
(\xi - c_\infty)^T \nabla^2 D(c_\infty)(\xi - c_\infty) = \sum_{r=1}^{\lceil R \rceil} k_r \left( \frac{y_{r,i} - y'_{r,i}}{c_{i,\infty}} (\xi - c_{i,\infty}) \right)^2,
\]

which follows from direct computations, we can Taylor expand around \(c_\infty\) to obtain

\[
\Lambda := \liminf_{\xi \to c_\infty} \frac{D(\xi)}{E(\xi|c_\infty)} = \inf_{\xi \in E_M} \left\{ \left( \sum_{r=1}^{\lceil R \rceil} k_r \left( \frac{y_{r,i} - y'_{r,i}}{c_{i,\infty}} (\xi - c_{i,\infty}) \right)^2 \right) \right\}.
\]

(2.16)

Denoting \(\mu = \xi - c_\infty\), we see that (2.16) becomes

\[
\Lambda = \inf_{\mu + c_\infty \in E_M} \left\{ \sum_{r=1}^{\lceil R \rceil} k_r \left( \frac{y_{r,i} - y'_{r,i}}{c_{i,\infty}} \mu_i \right)^2 \right\}.
\]

(2.17)

Since both numerator and denominator of the above fraction are of homogeneity two, it is sufficent to estimate (2.17) for \(\mu\) in the unit ball, i.e. \(|\mu| = 1\). By putting \(\eta_i = \frac{\mu_i}{c_{i,\infty}}\), we have \(\mu = \text{diag}(c_\infty) \eta\).

Hence, the numerator of (2.17) becomes

\[
\sum_{r=1}^{\lceil R \rceil} k_r \left( (y_{r,i} - y'_{r,i}) \cdot \eta \right)^2.
\]

(2.18)

This term can only be zero when \(\eta \in \ker(W)\), where we recall that \(W\) is the Wegscheider matrix

\[
W = [(y'_{r,i} - y_{r,i})_{r=1,...,\lceil R \rceil}]^T \in \mathbb{R}^{\lceil R \rceil \times N}.
\]

(2.19)

We now consider two cases, related to the number of conservation laws \(m = \dim \ker(W)\):

**Case 1.** \(m > 0\): Since the rows of \(Q\) form a basis of \(\ker(W)\), there exists \(a \in \mathbb{R}^m\) such that

\[
\eta = Q^T a.
\]

(2.20)

From \(Q \mu = Q \xi - Q c_\infty = M - M = 0\), we get

\[
Q \text{ diag}(c_\infty) Q^T a = Q \text{ diag}(c_\infty) \eta = Q \mu = 0.
\]

(2.21)

It follows from rank\((Q) = m\) and \(c_{i,\infty} > 0\) (for all \(i = 1, 2, \ldots, N\)) that \(Q \text{ diag}(c_\infty) Q^T\) is a positive definite \(m \times m\) matrix. Hence, (2.21) implies \(a = 0\), which shows \(\eta = 0\), and consequently \(|\mu| = 1\), which contradicts the fact that \(|\mu| = 1\).

**Case 2.** \(m = 0\): In this case it is obvious that \(\Lambda = 0\) if and only if \(\eta = 0\), since \(\ker(W) = 0\).

In conclusion, we have shown that \(\Lambda > 0\) and consequently (2.14).

A direct consequence of Lemma 2.3 is that (2.12) holds and thus (2.8) holds thanks to a classic Gronwall lemma. From \(E(c(t)|c_\infty) \leq E(c_0|c_\infty)\) and the inequality \(\Psi(x;y) \geq (\sqrt{x} - \sqrt{y})^2\) we get

\[
E(c_0|c_\infty) \geq \sum_{i=1}^{N} \Psi(c_i(t); c_{i,\infty}) \geq \sum_{i=1}^{N} \left( \sqrt{c_i(t)} - \sqrt{c_{i,\infty}} \right)^2 \geq \frac{1}{2} \sum_{i=1}^{N} c_i(t) - \sum_{i=1}^{N} c_{i,\infty}.
\]

Thus for all \(i = 1, \ldots, N\) and all \(t > 0\),

\[
c_i(t) \leq \tilde{K} := 2 \left( E(c_0|c_\infty) + \sum_{i=1}^{N} c_{i,\infty} \right).
\]

(2.22)
We can therefore show that
\[
E(c(t)|c_{\infty}) \geq \sum_{i=1}^{N} \left( \sqrt{c_i(t)} - \sqrt{c_i,_{\infty}} \right)^2 = \sum_{i=1}^{N} \frac{(c_i(t) - c_i,_{\infty})^2}{(\sqrt{c_i(t)} + \sqrt{c_i,_{\infty}})^2} \geq \frac{1}{(\sqrt{K} + \max \{\sqrt{c_i,_{\infty}}\})^2} \sum_{i=1}^{N} |c_i(t) - c_i,_{\infty}|^2
\]
and finally get (2.9).

\[\blacksquare\]

2.2. Convergence to equilibrium for PDE systems.

We now begin the proof of Theorem 1.1. We first rewrite its right hand side by using the additivity property of the relative entropy, that is (remembering that \(\overline{c_i} = \int_\Omega c_i\) and \(c_{\infty}\) do not depend upon \(x\)):
\[
\mathcal{E}(c|c_{\infty}) = \sum_{i=1}^{N} \int_\Omega c_i \log \frac{c_i}{\overline{c_i}} \, dx + \sum_{i=1}^{N} \left( \overline{c_i} \log \frac{\overline{c_i}}{c_i,_{\infty}} - \overline{c_i} + c_i,_{\infty} \right) = \mathcal{E}(\overline{c}|\overline{c}) + \mathcal{E}(\overline{c}|c_{\infty}).
\]
In order to treat the first term \(\mathcal{E}(\overline{c}|\overline{c})\) in the above identity, we apply the Logarithmic Sobolev Inequality
\[
\int_\Omega \frac{\|f\|_{L^2}^2}{\|f\|_{L^1}^2} \, dx \geq C_{LSI} \int_\Omega f \log \frac{f}{\mathcal{E}(\overline{c})} \, dx
\]
to estimate
\[
\frac{1}{2} \mathcal{D}(c) \geq \frac{1}{2} \sum_{i=1}^{N} \int_\Omega d_i \frac{\|\nabla c_i\|_{L^2}^2}{c_i} \, dx \geq \lambda_1 \mathcal{E}(\overline{c}|\overline{c}), \tag{2.23}
\]
with \(\lambda_1 = \frac{1}{2} C_{LSI} \min_i \{d_i\}\). By denoting
\[
G(c) := \frac{1}{2} \sum_{r=1}^{[R]} k_r \log \frac{c_{\infty}^{\Psi}}{c_{\infty}^{\Psi},_r}, \quad F(c) := \lambda_1 \sum_{i=1}^{N} \Psi(c_i; c_i,_{\infty}), \tag{2.24}
\]
we estimate (using again the Logarithmic Sobolev Inequality and the additivity of the relative entropy)
\[
\frac{1}{2} \mathcal{D}(c) \geq \lambda_1 \mathcal{E}(\overline{c}|\overline{c}) + \int_\Omega G(c) \, dx = \int_\Omega \underbrace{[F(c) + G(c)]}_{=\Phi(c)} \, dx - F(\overline{c})
\]
\[
= : \int_\Omega \Phi(c) \, dx - F(\overline{c}). \tag{2.25}
\]
Let \(\hat{\Phi}\) be the convexification of \(\Phi\), that is, \(\hat{\Phi}\) is the supremum of all affine functions below \(\Phi\). By noticing that \(\Phi \geq \hat{\Phi}\) and that \(\hat{\Phi}\) is convex, we can now use Jensen’s inequality (recall that \([R] = 1\), as stated in the paragraph devoted to notations at the end of the introduction) and get
\[
\frac{1}{2} \mathcal{D}(c) \geq \int_\Omega \hat{\Phi}(c) \, dx - F(\overline{c}) \geq \hat{\Phi}(\overline{c}) - F(\overline{c}). \tag{2.26}
\]
It remains to show the following finite-dimensional inequality:
\[
\hat{\Phi}(\overline{c}) - F(\overline{c}) \geq \lambda_2 \mathcal{E}(\overline{c}|c_{\infty}),
\]
for some positive constant \(\lambda_2 > 0\). By using Jensen’s inequality, we see that
\[
\mathcal{E}(\overline{c}|c_{\infty}) \leq \mathcal{E}(c|c_{\infty}) \leq \mathcal{E}(c_0|c_{\infty}).
\]
We define, for some given \(K > 0\),
\[
\mathcal{C}_M := \{ \xi \in \mathbb{R}_+^N : \xi = M \text{ and } \mathcal{E}(\xi|c_{\infty}) \leq K \}.
\]
Our goal is now to prove the
Lemma 2.4.

\[ \lambda_2 := \inf_{\xi \in \mathcal{C}_M} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_\infty)} > 0. \]  

Proof. We first observe that \( \lambda_2 \geq 0 \). Indeed, using the inequality \( f + g \geq \hat{f} + \hat{g} \) (see e.g. [WCHL96]), we have

\[ \hat{\Phi}(\xi) - F(\xi) \geq \hat{F}(\xi) + \hat{G}(\xi) - F(\xi) = \hat{G}(\xi) \geq 0, \]

thanks to the convexity of \( F \) (and the nonnegativity of \( G \)). On the other hand, since \( \mathcal{E}(\xi|c_\infty) \) is bounded above in \( \mathcal{C}_M \), then \( \lambda_2 \) can be equal to zero only when

\[ \hat{\Phi}(\xi) - F(\xi) \to 0 \quad \text{or equivalently} \quad \hat{G}(\xi) \to 0. \]

Now using the fact that (when \( \xi = M \)) \( G(\xi) = 0 \) if and only if \( \xi = c_\infty \), it follows that \( \hat{G}(\xi) = 0 \) if and only if \( \xi = c_\infty \), thanks to the definition of the convexification. Therefore \( \lambda_2 \) could only vanish for states \( \mathcal{C}_M \ni \xi \to c_\infty \). In other words, it is sufficient to prove that

\[ \liminf_{\mathcal{C}_M \ni \xi \to c_\infty} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_\infty)} > 0. \]  

We then use the

Lemma 2.5. [MHM15]

Let \( \Phi \) be defined by (2.25) and (2.24).

Then, there exists \( \delta > 0 \) such that for all \( \xi \in \mathbb{R}^N_+ \cap B(c_\infty, \delta) \) (ball centered at \( c_\infty \) with radius \( \delta \)) the identity \( \hat{\Phi}(\xi) = \Phi(\xi) \) holds.

Proof. The proof of this Lemma is given in [MHM15, Lemma 3.4]. For the sake of completeness and the convenience of the reader, we also provide a proof in the Appendix of this paper.

We point out that all estimates in the proof of Lemma 2.5 are explicit. This means that it is in principle possible to actually estimate the radius \( \delta > 0 \) of the ball \( B(c_\infty, \delta) \), on which \( \Phi \) and its convexification \( \hat{\Phi} \) are identical.

We continue the proof of Lemma 2.4. Thanks to Lemma 2.5 above, we see that \( \hat{\Phi} \equiv \Phi \) in a neighborhood of \( c_\infty \). Thus, (2.28) is equivalent to

\[ \liminf_{\mathcal{C}_M \ni \xi \to c_\infty} \frac{G(\xi)}{\mathcal{E}(\xi|c_\infty)} > 0, \]  

which was already proven in Proposition 2.2 (see (2.15) and note that \( D \) in (2.15) corresponds to \( G \) in (2.29)), up to a small modification (\( E(c_0|c_\infty) \) is replaced by some arbitrary \( K > 0 \)).

Therefore, from (2.23), (2.26) and (2.27), we obtain the main estimate (1.13) with \( \lambda = \min \{ \lambda_1, \lambda_2 \} \).

We first note that for the renormalised solutions considered in our theorem (that is, satisfying (1.14)), the entropy inequality \( \mathcal{E}(c(t)|c_\infty) \leq \mathcal{E}(c_0|c_\infty) \) holds, so that we can use estimate (1.13). We can then directly apply an integral form of Gronwall’s lemma, see e.g. [Wil] or more specifically [FL16], to finally ensure that

\[ \mathcal{E}(c(t)|c_\infty) \leq e^{-\lambda t} \mathcal{E}(c_0|c_\infty). \]

Finally, we conclude the proof of Theorem 1.1 by using the

Lemma 2.6 (Csiszár-Kullback-Pinsker inequality, [AMTU01]).

Assume that \( c : \Omega \to \mathbb{R}_+ \) is nonnegative and measurable and \( c_\infty \in \mathbb{R}^N_0 \) such that \( Q \sigma = Q c_\infty \).

Then, there exists a constant \( C_{CKP} > 0 \) depending only on \( \Omega \) and \( Q \) such that

\[ C_{CKP} \sum_{i=1}^N \| c_i - c_i,\infty \|_{L^1(\Omega)}^2 \leq \mathcal{E}(c|c_\infty). \]
2.3. A constructive entropy method and its application to a cycle of reactions.

Theorem 1.1 shows exponential convergence to equilibrium for the system (1.1), (1.3) when it is complex balanced and has no boundary equilibria. However, it does not give a quantitative estimate for the convergence rate $\lambda$. The reason is that the strictly positive limit in (2.29), which can be computed explicitly, is not necessary the positive infinite $\lambda_2$ in (2.27). Though this infinite comes out of an optimisation problem in finite dimension (and not from an abstract compactness argument in infinite dimension), it remains not explicitly computable. Therefore, in this subsection, we propose a constructive method to prove the entropy-dissipation estimate (1.13) and to explicitly estimate the convergence rate $\lambda$.

This method builds on some recent ideas developed in [FT] for nonlinear detailed balance systems. It consists of four steps. The first three ones can be proven for any general complex balanced systems without boundary equilibria. The last step, being based on the structure of the mass conservation laws, is proven only for a specific network consisting of a cycle of reactions with arbitrary number of species, described in Thm. 1.2.

We begin therefore the

Proof of Thm. 1.2: Since the three first steps of the proof hold for complex balance general systems without boundary equilibria, we use the notations $\rho, \overline{\rho}, \rho_{\infty}$. For step four we change the notation to $a, \bar{a}, a_{\infty}$ to underline that this step is specific to system (1.17).

Step 1 (Additivity of the relative entropy and Logarithmic Sobolev Inequality)

As in the proof of Thm. 1.1, we use the additivity property (for $\mathcal{E}$ given by formula (1.10)):

$$\mathcal{E}(\rho;\rho_{\infty}) = \mathcal{E}(\rho;\overline{\rho}) + \mathcal{E}(\overline{\rho};\rho_{\infty}),$$

and control the term $\mathcal{E}(\rho;\overline{\rho})$ via the Logarithmic Sobolev Inequality as in (2.23) (with $\mathcal{D}$ defined by formula (1.11)), i.e.

$$\frac{1}{2} \mathcal{D}(\rho) \geq \lambda_1 \mathcal{E}(\rho;\overline{\rho}),$$

for an explicit constant $\lambda_1 > 0$. It remains therefore to control $\mathcal{E}(\overline{\rho};\rho_{\infty})$.

Step 2 (Upper and lower bounds using quadratic terms of square root concentrations)

We estimate $\mathcal{D}(\rho)$ and $\mathcal{E}(\overline{\rho};\rho_{\infty})$ by quadratic terms involving the square roots of concentrations, which are significantly easier to handle than logarithmic terms. Recalling the short hand notation $C_i := \sqrt{c_i}$ and $C_{i,\infty} := \sqrt{c_{i,\infty}}$ for square roots of concentrations, we estimate first

$$\frac{1}{2} \mathcal{D}(\rho) \geq 2 \min_j \{d_j\} \sum_{i=1}^N \|\nabla C_i\|^2 + \frac{1}{2} \sum_{j=1}^R k_j c_i \rho_{\infty} \|c_i\|^2 \left( \frac{C_{i,\infty}^{y_j}}{C_i} - \frac{c_{i,\infty}^{y_j}}{C_{i,\infty}} \right)^2,$$

thanks to inequality (2.1). Now, by recalling the conservation of mass (1.18) and the nonnegativity of the solution, we see that $c_i \leq M$ for all $i = 1, \ldots, N$. Moreover, since $\Psi(\cdot, y)/(\sqrt{x} - \sqrt{y})^2$ is increasing for each fixed $y > 0$, the elementary estimate

$$\frac{\Psi(x; y)}{(\sqrt{x} - \sqrt{y})^2} \leq \frac{\Psi(M; y)}{(\sqrt{M} - \sqrt{y})^2}$$

holds for all $x \in (0, M]$ and $y > 0$, and we can estimate $\mathcal{E}(\overline{\rho};\rho_{\infty})$ as follows:

$$\mathcal{E}(\overline{\rho};\rho_{\infty}) \leq \sum_{i=1}^N \frac{\Psi(M; c_{i,\infty})}{(\sqrt{M} - \sqrt{c_{i,\infty}})^2} \left( \sqrt{C_i}^2 - C_{i,\infty} \right)^2 \leq K_2 \sum_{i=1}^N \left( \sqrt{C_i}^2 - C_{i,\infty} \right)^2,$$

where $K_2$ depends only on $M$ and $c_{\infty}$.

Step 3 (Control of the reaction dissipation term via a reaction dissipation term for spatial averages)

As another step towards exploiting the mass conservation laws, we estimate $\mathcal{D}(\rho)$ further by

$$\frac{1}{2} \mathcal{D}(\rho) \geq K_1 \left( \sum_{i=1}^N \|\nabla C_i\|^2 + \sum_{i=1}^N \left( \frac{C_{i,\infty}^{y_j}}{C_i} - \frac{c_{i,\infty}^{y_j}}{C_{i,\infty}} \right)^2 \right),$$

for an explicit constant $K_1 > 0$. This technical step follows from [FT, Lemma 2.6] with slight modifications. For the sake of completeness, we recall here the main arguments of the proof. The proof makes use of a domain decomposition corresponding to the deviation to the averages of $C_i$, i.e. by denoting $\delta_i(x) = C_i(x) - \overline{C}_i$, we consider the decomposition

$$\Omega = S \cup S^c$$
where $S = \{ x \in \Omega : |\delta_i(x)| \leq L \text{ for all } i = 1, \ldots, N \}$ with a fixed constant $L > 0$ (which can be chosen arbitrarily). On the set $S$, thanks to a Taylor expansion, we have
\[
\frac{C^{y_i}}{C^{y_\infty}} = \frac{1}{C^{y_\infty}} \prod_{i=1}^{N} (C_i + \delta_i)^{y_i} = \frac{C^{y_i}}{C^{y_\infty}} + \tilde{R}(C_1, \ldots, C_N, \delta_1, \ldots, \delta_N) \sum_{i=1}^{N} \delta_i
\]
with $|\tilde{R}(C_i, \delta_i)| \leq C(L)$ thanks to the upper bounds $C_i \leq \sqrt{C_i} \leq \sqrt{K}$ (obtained by a computation similar to (2.22)) and $|\delta_i(x)| \leq L$ in $S$. It thus follows that
\[
\frac{1}{2} \sum_{i=1}^{R} k_i e_i^{y_i} \left( \frac{C^{y_i}}{C^{y_\infty}} - \frac{C^{y_i}}{C^{y_\infty}} \right)^2 \geq \beta_1 \sum_{i=1}^{R} \left( \frac{C^{y_i}}{C^{y_\infty}} - \frac{C^{y_i}}{C^{y_\infty}} \right)^2 |S| - \beta_2 \sum_{i=1}^{N} \|\delta_i\|_{L^2(S)}^2
\]
for some constants $\beta_1, \beta_2 > 0$. On the other hand, by using the lower bounds $|\delta_i| \geq L$ on $S^c$ for some $1 \leq i \leq N$, the boundedness of $C_i$, and Poincaré’s inequality, it follows that
\[
\frac{1}{2} \min_{j} \{ d_j \} \sum_{i=1}^{N} \| \nabla C_i \|^2 \geq \beta_3 \sum_{i=1}^{N} \| \delta_i \|_{L^2(\Omega)}^2 \geq \beta_3 N L^2 |S^c| \geq \beta_4 \sum_{i=1}^{R} \left( \frac{C^{y_i}}{C^{y_\infty}} - \frac{C^{y_i}}{C^{y_\infty}} \right)^2 |S| \geq \beta_5 \sum_{i=1}^{R} \left( \frac{C^{y_i}}{C^{y_\infty}} - \frac{C^{y_i}}{C^{y_\infty}} \right)^2 \left( \frac{\| \nabla C_i \|^2}{\| \nabla C_i \|^2} \right)
\]
for constants $\beta_3, \beta_4 > 0$. Note that all constants $\beta_1, \beta_2, \beta_3, \beta_4$ are independent of $S$. Now, a combination of (2.30), (2.33) and (2.34) leads to, for any $\theta \in (0, 1)$,
\[
\frac{D(\varepsilon)}{2} \geq \min_{j} \{ d_j \} \sum_{i=1}^{N} \| \nabla C_i \|^2 + \frac{\beta_3}{2} N \sum_{i=1}^{N} \| \delta_i \|^2_{L^2(S)} + \min_{j} \{ \delta_{j1}, \delta_{j2} \} \sum_{i=1}^{N} \| \nabla C_i \|^2 + \frac{\theta}{2} \sum_{i=1}^{R} k_i e_i^{y_i} \left( \frac{C^{y_i}}{C^{y_\infty}} - \frac{C^{y_i}}{C^{y_\infty}} \right)^2
\]
where we recall that $|S| + |S^c| = |\Omega| = 1$. Hence, (2.32) follows by choosing $L, 0 < \theta < \min \{ \beta_3, \beta_2^{-1}, 1 \}$ and $K_1 = \min \{ \theta \beta_1, \beta_4, \min_{j} \{ d_j \} \}$.  

**Step 4** (Conservation laws and deviations of averaged concentration around the equilibrium)

Since this step is specific to the cyclic reaction (1.17), we change the notation to $a, \bar{a}, a_\infty, A, \varepsilon$, etc. in the remainder of the proof. Moreover, we recall the notations $A_i = \sqrt{a_i}, A_i = \sqrt{\bar{a}_i}$, and use the periodic notation $a_{N+1} := a_1$ or $a_{N+1} := a_\infty$, etc.

First, we recall that the system (1.17) is indeed a complex balanced system satisfying exactly one conservation law (corresponding to the conservation of the total mass): $\sum_{i=1}^{N} \frac{1}{\gamma_i} \int_{0}^{\infty} a_i = M$. Indeed, the following lemma holds:

**Lemma 2.7.** For any strictly positive initial mass $M > 0$, there exists a unique strictly positive complex balance equilibrium $a_\infty = (a_{1,\infty}, a_{2,\infty}, \ldots, a_{N,\infty})$ to the system (1.17), which solves
\[
\begin{cases}
  k_1 a_{1,\infty}^2 = k_2 a_{2,\infty}^2 = \ldots = k_N a_{N,\infty}^2, \\
  a_{1,\infty} + a_{2,\infty}^2 + \ldots + a_{N,\infty}^2 = M.
\end{cases}
\]

**Proof.** Using the first equation of (2.35), we see that $a_1 = \left( \frac{k_1}{k_2} a_{1,\infty}^2 \right)^{1/\alpha_1}$ for all $i = 2, 3, \ldots, N$. Hence, $a_{1,\infty}$ solves
\[
\frac{a_{1,\infty}}{\alpha_1} + \frac{1}{\alpha_2} \left( \frac{k_1}{k_2} a_{1,\infty}^2 \right)^{1/\alpha_2} + \ldots + \frac{1}{\alpha_N} \left( \frac{k_1}{k_N} a_{1,\infty}^2 \right)^{1/\alpha_N} = M.
\]

Denoting the left hand side of this equation by $f(a_{1,\infty})$, we see that $z \mapsto f(z)$ is a strictly increasing function on $[0, +\infty)$ with $f(0) = 0$ and $\lim_{z \to +\infty} f(z) = +\infty$. Thus, for any $M > 0$, there exists a unique solution $a_{1,\infty}$ to the equation $f(z) = M$, which completes the proof of the lemma.

In the case of a cyclic reaction (1.17), the relative entropy is specifically defined by (1.19), while the corresponding entropy dissipation is given by (1.20).

Moreover, following the above general method, it will be convenient for the readability of the proof to briefly restate the required key estimates of Steps 1-3: First, we see that Step 1 becomes
\[
\mathcal{E}(a|a_\infty) = \mathcal{E}(a|\bar{a}) + \mathcal{E}(\bar{a}|a_\infty),
\]
(2.36)
and, thanks to the logarithmic Sobolev inequality,

$$\frac{1}{2} \mathcal{D}(\mathbf{a}) \geq \lambda_1 \mathcal{E}(\mathbf{a} | \mathbf{a}^\infty).$$

(2.37)

Step 2 shows that $\mathcal{D}(\mathbf{a})$ and $\mathcal{E}(\mathbf{a}^\infty)$ are controlled by quadratic terms of square roots of concentrations

$$\mathcal{D}(\mathbf{a}) \geq 4 \min_j \left\{ d_j \right\} \sum_{i=1}^{N} \| \nabla A_i \|^2 + \sum_{i=1}^{N} k_i a_i^{\infty} \left\| \frac{A_i^{\infty}}{A_i^\infty} - \frac{A_i^{\infty+1}}{A_i^{\infty+1}} \right\|^2,$$

(2.38)

and, thanks to the logarithmic Sobolev inequality,

$$\mathcal{E}(\mathbf{a} | \mathbf{a}^\infty) \leq K_0 \sum_{i=1}^{N} \left( \sqrt{A_i^2 - A_i^{\infty+1}} \right)^2.$$ 

(2.39)

In Step 3, we obtain moreover,

$$\mathcal{D}(\mathbf{a}) \geq K_1 \left( \sum_{i=1}^{N} \| \nabla A_i \|^2 + \sum_{i=1}^{N} \left( \frac{A_i^{\infty}}{A_i^\infty} - \frac{A_i^{\infty+1}}{A_i^{\infty+1}} \right)^2 \right).$$

(2.40)

In order to complete the proof of Thm. 1.2 in the final Step 4, it therefore remains to show that

$$\sum_{i=1}^{N} \| \nabla A_i \|^2 + \sum_{i=1}^{N} \left( \frac{A_i^{\infty}}{A_i^\infty} - \frac{A_i^{\infty+1}}{A_i^{\infty+1}} \right)^2 \geq K_2 \sum_{i=1}^{N} \left( \sqrt{A_i^2 - A_i^{\infty}} \right)^2.$$ 

(2.41)

We recall the definition $\delta_i(x) = A_i(x) - \overline{A_i}$, which implies $\| \delta_i \|^2 = \overline{A_i^2} - \overline{A_i}^2$ for $i = 1, \ldots, N$. By using the ansatz

$$\overline{A_i^2} = A_i^{\infty} (1 + \mu_i)^2 \quad \text{with} \quad \mu_i \in [-1, +\infty) \quad \text{for} \quad i = 1, \ldots, N,$$

(2.42)

we compute

$$\overline{A_i} = A_i^{\infty} (1 + \mu_i) - \frac{\| \delta_i \|^2}{\sqrt{\overline{A_i^2} + \overline{A_i}}} =: A_i^{\infty} (1 + \mu_i) - \| \delta_i \|^2 R(A_i) \quad \text{for all} \quad i = 1, 2, \ldots, N,$$

where we have defined $R(A_i) := \left( \overline{A_i^2} + \overline{A_i} \right)^{-1}$. Moreover, thanks to the mass conservation law

$$\sum_{i=1}^{N} \frac{1}{\alpha_i} \int_{\Omega} a_i(x) \, dx = M = \sum_{i=1}^{N} \frac{1}{\alpha_i} A_i^{\infty},$$

(2.43)

we know that

$$\sum_{i=1}^{N} \frac{1}{\alpha_i} A_i^{\infty} \mu_i (\mu_i + 2) = 0.$$ 

Next, we can always fix an $\varepsilon > 0$ small enough such that

$$\gamma := \min_{i,j \in \{1, \ldots, N\}} \left\{ \frac{1}{2 \alpha_i^j+1} A_i^{\infty j} \left( \frac{M - \varepsilon^2/\alpha_i^j}{N-1} \right)^{\alpha_i^j} - \frac{\varepsilon^{2 \alpha_i}}{A_i^{\infty}} > 0, \right. \right.$$ 

(2.44)

and consider two cases.

**Case 4.1:** If $\overline{A_i^2} \geq \varepsilon^2$ for all $i = 1, \ldots, N$, then $R(A_i) \leq \varepsilon^{-1}$ for $i = 1, 2, \ldots, N$. Note that

$$\mu_i = -1 + \sqrt{\frac{\overline{A_i^2}}{A_i^{\infty}}} \leq -1 + \frac{\sqrt{\alpha_i M}}{A_i^{\infty}},$$

thanks to (2.42) and $\| \delta_i \| \leq \sqrt{\overline{A_i^2}} \leq \sqrt{\alpha_i M}$. Moreover, we use the ansatz (2.42) and the Taylor expansion

$$\frac{\overline{A_i^{\infty j}}}{A_i^{\infty j}} = (1 + \mu_i) \left( \frac{\| \delta_i \|^2 R(A_i)}{A_i^{\infty}} \right)^{\alpha_i} = (1 + \mu_i)^{\alpha_i} - R_i(\mu_i, R(A_i), \| \delta_i \|, \alpha_i, A_i^{\infty}) \| \delta_i \|$$
with a bounded remainder term $\tilde{R}_i (\mu_i, R(A_i), \|\delta_i\|, \alpha_i, A_i, \infty) \leq C(\varepsilon)$ thanks to the bounds on $\mu_i$, $R(A_i)$ and $\|\delta_i\|$. This yields for a $\theta \in (0,1)$ (to be chosen) the lower bound

$$\text{LHS of (2.41)} \geq \sum_{i=1}^{N} \|\nabla A_i\|^2 + \theta \sum_{i=1}^{N} \left[ (1 + \mu_i)^{\alpha_i} - \tilde{R}_i \|\delta_i\| \right] - \left( (1 + \mu_{i+1})^{\alpha_{i+1}} - \tilde{R}_{i+1} \|\delta_{i+1}\| \right)^2 \geq \sum_{i=1}^{N} \|\nabla A_i\|^2 + \theta \sum_{i=1}^{N} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_{i+1})^{\alpha_{i+1}} \right)^2 - \theta C(\varepsilon) \sum_{i=1}^{N} \|\delta_i\|^2 \geq \theta \sum_{i=1}^{N} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_{i+1})^{\alpha_{i+1}} \right)^2,$$

(2.45)

where we have applied the Poincaré’s inequality and chosen $\theta$ small enough in the last step. Moreover, the right hand side of (2.41) is bounded above by

$$K_3 \max_{i=1,\ldots,N} \left\{ A_{i,\infty}^2 \right\} \sum_{i=1}^{N} \mu_i^2.$$

(2.46)

Thanks to Lemma 2.8 below, we obtain (2.41) from (2.45) and (2.46) by choosing

$$K_3 \leq \frac{\theta}{2NN \max_{i=1,\ldots,N} \left\{ A_{i,\infty}^2 \right\}}.$$

(2.47)

**Lemma 2.8.** Assume that $\mu_i \in [-1, +\infty)$ and $\alpha_i \in \mathbb{N}_{>0}$, $i = 1, 2, \ldots, N$ satisfy (for some $A_{i, \infty} \geq 0$),

$$\sum_{i=1}^{N} \frac{1}{\alpha_i} A_{i,\infty}^2 (\mu_i^2 + 2\mu_i) = 0.$$

(2.48)

Then,

$$\sum_{i=1}^{N} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_{i+1})^{\alpha_{i+1}} \right)^2 \geq \frac{1}{NN} \sum_{i=1}^{N} \mu_i^2.$$

(2.49)

**Proof.** Note that the left hand side of (2.49) can be bounded below by

$$\sum_{i=1}^{N} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_{i+1})^{\alpha_{i+1}} \right)^2 \geq \frac{1}{NN} \sum_{i \neq j} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_j)^{\alpha_j} \right)^2,$$

(2.50)

thanks to the elementary inequality $a^2 + b^2 \geq \frac{1}{2} (a + b)^2$. Thanks to the mass conservation (2.48), there exists $\emptyset \neq I_0 \subset \{1, 2, \ldots, N\}$ such that $\mu_i \geq 0$ for $i \in I_0$ and $\mu_j < 0$ for $j \in I_0^C$, where also the complement $I_0^C$ is not empty. Thus, for $i \in I_0$ and $j \in I_0^C$, we have

$$(1 + \mu_i)^{\alpha_i} - (1 + \mu_j)^{\alpha_j} \geq (1 + \mu_i) - (1 + \mu_j) = \mu_i - \mu_j \geq 0.$$ \nonumber

Then

$$\left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_j)^{\alpha_j} \right)^2 \geq (\mu_i - \mu_j)^2 \geq \mu_i^2 + \mu_j^2$$

for all $i \in I_0$ and $j \in I_0^C$. We can thus continue to estimate (2.50)

$$\sum_{i=1}^{N} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_{i+1})^{\alpha_{i+1}} \right)^2 \geq \frac{1}{NN} \sum_{i \neq j} \left( (1 + \mu_i)^{\alpha_i} - (1 + \mu_j)^{\alpha_j} \right)^2 \geq \frac{1}{NN} \sum_{i \in I_0, j \in I_0^C} (\mu_i^2 + \mu_j^2) \geq \frac{1}{NN} \sum_{i=1}^{N} \mu_i^2.$$\nonumber

(2.51)

This ends the proof of Lemma 2.8 and Case 4.1 of Thm. 1.2.

**Remark 2.2.** The proof of the above Lemma 2.8 used the conservation law (2.48) in a rather weak way. Indeed, it only uses the fact that not all $\mu_i$ have the same sign. We believe that the ideas used in the proof of Lemma 2.8 can be generalised to many complex balanced systems, once the structure of the conservation laws is explicitly given.
**Case 4.2:** We now consider cases where there exists (at least) one small relative entropy towards the equilibrium) is bounded above by a constant due to the following distance from the equilibrium state. Hence, we expect that the entropy dissipation (and its lower bound on the left-hand-side of (2.41)) should be bounded below by a positive constant, which expresses this distance from the equilibrium state.

At first, however, we note that the right hand side of (2.41) (which is an upper bound of the relative entropy towards the equilibrium) is bounded above by a constant due to the following estimate:

\[ K_3 \sum_{i=1}^{N} \left( \sqrt{A_i^2} - A_i,\infty \right)^2 \leq 2K_3 \sum_{i=1}^{N} (\alpha_i + a_i,\infty) \leq 2K_3 \left( M \max\{\alpha_i\} + \sum_{i=1}^{N} a_i,\infty \right). \tag{2.52} \]

Hence, in order to prove (2.41), it remains to show that the left hand side of (2.41) is bounded below by a positive constant. To show that, we define

\[ \eta := \min_i \frac{M - \epsilon^2/\alpha_i}{2(N-1)}, \tag{2.53} \]

and distinguish two further subcases representing states with large and small spatial variations:

**Subcase 1:** There exists \( i_0 \in \{1, 2, \ldots, N\} \) such that \( \|\delta_{i_0}\|^2 \geq \eta > 0 \). Then, we can estimate directly thanks to (2.52) and Poincaré-Wirtinger’s inequality:

\[ \text{LHS of (2.41)} \geq \sum_{i=1}^{N} C_P \|\delta_i\|^2 \geq C_P \eta \geq K_3 \sum_{i=1}^{N} \left( \sqrt{A_i^2} - A_i,\infty \right)^2, \tag{2.54} \]

by choosing

\[ K_3 \leq \frac{C_P \eta}{2 \left( M \max\{\alpha_i\} + \sum_{i=1}^{N} c_i,\infty \right)}. \tag{2.55} \]

**Subcase 2:** For all \( i = 1, 2, \ldots, N \), we have \( \|\delta_i\|^2 \leq \eta \). In this case concerning states away from equilibrium yet with small spatial variation, we expect that the reaction terms represent in large parts the distance from the equilibrium. This can be quantified from firstly recalling the conservation law \( \sum_{i=1}^{N} \frac{1}{\alpha_i} A_i^2 = M \), and from observing the estimate

\[ \sum_{i \neq i_0} \frac{1}{\alpha_i} A_i^2 = M - \frac{1}{\alpha_{i_0}} A_{i_0}^2 \geq M - \frac{\epsilon^2}{\alpha_{i_0}}, \]

Thus, there exists \( j_0 \neq i_0 \) such that

\[ A_{j_0}^2 \alpha_{j_0} \geq \frac{M - \frac{\epsilon^2}{\alpha_{i_0}}}{N-1} \geq \frac{M - \frac{\epsilon^2}{\alpha_{i_0}}}{N-1} \]

since \( \alpha_{j_0} \geq 1 \). Hence, by using (2.53),

\[ \frac{\epsilon_{j_0}}{\alpha_{j_0}} A_{j_0}^2 - \|\delta_{j_0}\|^2 \geq \frac{M - \frac{\epsilon^2}{\alpha_{i_0}}}{N-1} - \eta \geq \frac{M - \frac{\epsilon^2}{\alpha_{i_0}}}{2(N-1)}. \]

Now, with \( \gamma \) defined in (2.44), by using the triangle inequality and Young’s inequality, we get

\[ \text{LHS of (2.41)} \geq \sum_{i=1}^{N} \left( \frac{A_{i_{0}}}{A_{i_{0},\infty}} \frac{A_{i+1}}{A_{i+1,\infty}} - \frac{A_{i+1}}{A_{i+1,\infty}} \frac{A_{i_{0}}}{A_{i_{0},\infty}} \right)^2 \geq \frac{1}{\left| j_0 - i_0 \right| + 1} \left( \frac{A_{i_{0}}^{\alpha_{i_0}}}{A_{i_{0},\infty}^{\alpha_{i_0}} - A_{j_0}^{\alpha_{i_0}} - A_{j_0}^{\alpha_{i_0}} A_{i_{0},\infty}^{\alpha_{i_0}} - \frac{A_{j_0}^{\alpha_{i_0}}}{A_{j_0}^{\alpha_{i_0}} A_{i_{0},\infty}^{\alpha_{i_0}}} \right)^2 \geq \frac{1}{N} \left( \frac{\epsilon_{j_0}}{\alpha_{j_0}} A_{j_0}^2 \right) - \frac{1}{N} \left( \frac{M - \frac{\epsilon^2}{\alpha_{i_0}}}{N-1} - \frac{\epsilon_{j_0}}{\alpha_{j_0}} A_{j_0}^2 \right) \]

\[ \geq \frac{\gamma}{N} \geq K_3 \sum_{i=1}^{N} \left( \sqrt{A_i^2} - A_i,\infty \right)^2, \tag{2.56} \]

whenever

\[ K_3 \leq \frac{\gamma}{2N \left( M \max\{\alpha_i\} + \sum_{i=1}^{N} c_i,\infty \right)}. \tag{2.57} \]
From (2.47), (2.55) and (2.57), we obtain (2.41) with an explicit $K_3$. Then by combining (2.36), (2.37) and (2.41), we can conclude the desired entropy entropy-dissipation inequality. The second part of the proof of Thm. 1.2 concerns the exponential convergence of the renormalised solution towards the strictly positive complex balance equilibrium.

Firstly, if the solution of the system (1.17) is regular enough in order to rigorously satisfy the weak entropy entropy-dissipation law (1.14), i.e., for a.a. $t \geq s > 0$,

\[ \mathcal{E}(a|a_\infty)(t) + \int_s^t \mathcal{D}(a)(r)dr \leq \mathcal{E}(a|a_\infty)(s), \]

then exponential convergence to equilibrium in relative entropy $\mathcal{E}(a|a_\infty)$ follows from a suitable Gronwall argument, see e.g. [Wil] or more specifically [FL16]. Moreover, the Csiszár-Kullback-Pinsker inequality yields exponential convergence of solutions to the complex balance equilibrium in $L^1$.

However, for renormalised solution as introduced in [Fis15], it is unclear if the weak entropy entropy-dissipation law (1.14) holds due to the lacking integrability of the nonlinear reaction terms in the entropy-dissipation. We will resolve this issue by proving exponential convergence to equilibrium with a uniform dissipation law (1.14) holds due to the lacking integrability of the nonlinear reaction terms in the entropy-dissipation law. From (2.47), (2.55) and (2.57), we obtain (2.41) with an explicit

\[ K_3 \]

where $\eta > 0$. It was proven in [Fis15] that (2.58) has a unique weak solution $c^\varepsilon$, with $c^\varepsilon \in L^\infty([0, +\infty]; \log L(\Omega))$ and, as $\varepsilon \to 0$,

\[ c^\varepsilon \to c \ \ \text{almost everywhere on } \Omega \times [0, +\infty), \]

(2.59)

where $c$ is the renormalised solution of (1.1). Note that (1.1) and (2.58) share the same complex balance equilibrium $c_\infty$. Moreover, we use the same relative entropy for (2.58) as for the limiting system:

\[ \mathcal{E}(c^\varepsilon|c_\infty) = \sum_{i=1}^N \int_{\Omega} \left( c_i^\varepsilon \log \frac{c_i^\varepsilon}{c_i^\varepsilon + c_{i,\infty}} - c_i^\varepsilon + c_{i,\infty} \right) dx. \]

By using Proposition 2.1 and the notation $G(c)$ in (2.24), we compute the entropy-dissipation for (2.58) as

\[ \mathcal{D}(c^\varepsilon) = -\frac{d}{dt} \mathcal{E}(c^\varepsilon|c_\infty) = \sum_{i=1}^N \int_{\Omega} \left( \frac{\nabla c_i^\varepsilon}{c_i^\varepsilon} \right)^2 dx + \int_{\Omega} \frac{G(c^\varepsilon)}{1 + \varepsilon |R(c^\varepsilon)|} dx. \]

We can now follow the first part of the proof of Thm. 1.2 to show that

\[ \mathcal{D}(c^\varepsilon) \geq \eta \mathcal{E}(c^\varepsilon|c_\infty) \]

(2.60)

where $\eta > 0$ is independent of $\varepsilon$. Indeed, in the first step, we write

\[ \mathcal{E}(c^\varepsilon|c_\infty) = \mathcal{E}(c^\varepsilon|\overline{c}) + \mathcal{E}(\overline{c}|c_\infty), \]

in which the first term is controlled by $\mathcal{D}(c^\varepsilon)$ thanks to the Logarithmic Sobolev inequality

\[ \frac{1}{2} \mathcal{D}(c^\varepsilon) \geq \lambda_1 \mathcal{E}(c^\varepsilon|\overline{c}), \]

where $\lambda_1$ is independent of $\varepsilon$. In the second step, we directly estimate $\mathcal{D}(c^\varepsilon)$ below by

\[ \mathcal{D}(c^\varepsilon) \geq 4 \min \{ d_j \} \sum_{i=1}^N \| \nabla C_i^\varepsilon \|^2 + \int_{\Omega} \frac{H(C^\varepsilon)}{1 + \varepsilon |R(c^\varepsilon)|} dx, \]

where

\[ H(C^\varepsilon) = \sum_{r=1}^{\lvert R \rvert} k_r \varepsilon^\alpha \left( \frac{(C^\varepsilon)^{\nu_r}}{C^\varepsilon_{\infty}^{\nu_r}} - \frac{(C^\varepsilon_{\infty})^{\nu_r}}{C^\varepsilon_{\infty}^{\nu_r}} \right)^2. \]

Following Step 3 and using the domain decomposition $\Omega = S \cup S^c$, we observe first that on $S = \{ x \in \Omega : \delta_i(x) \leq L \text{ for all } i = 1, \ldots, N \}$. Secondly, we note that for cyclic reactions satisfying the mass conservation law (1.18), we estimate $\overline{C_i^\varepsilon} \leq \varepsilon^\alpha M$ thanks to Jensen’s inequality. Therefore, we have
$|C_i^\varepsilon| \leq |\delta_i| + C^\varepsilon$ and thus $|R(e^\varepsilon)| \leq C(L, M)$ for some constant $C(L, M)$ depending on $L$ and $M$ (and the rates appearing in $R$). Note that for general systems, we can equally apply the upper bounds (2.22) instead of (1.18) and replace $M$ by $K$.

Hence, we obtain for $0 \leq \varepsilon \leq 1$ the uniform lower bound

$$\frac{1}{1 + \varepsilon |R(e^\varepsilon)|} \geq \frac{1}{1 + C(L, M)}.$$  

As a consequence, with this simple modification, we can show as in the proof of Thm. 1.2 that

$$D^\varepsilon(e^\varepsilon) \geq K_1 \left( \sum_{i=1}^N \|\nabla C_i^\varepsilon\|^2 + \sum_{i=1}^N \left( \frac{C_i^\varepsilon \rightarrow v_i^\varepsilon}{C_i^\varepsilon \rightarrow \infty} - \frac{C_i^\varepsilon \rightarrow v_i^\varepsilon}{C_i^\varepsilon \rightarrow \infty} \right)^2 \right)$$

where $K_1 > 0$ is independent of $\varepsilon$. Note that the statement involving $S^\varepsilon$ holds without change.

Moreover, we can now repeat all the arguments in Step 4 to finally prove estimate (2.60). Hence, by applying the Gronwall lemma, we obtain convergence to equilibrium in relative entropy, i.e.

$$E(e^\varepsilon(t)|e_\infty) \leq e^{-\eta t} E(e_0^\varepsilon|e_\infty),$$

(2.61)

for all $t > 0$ and with a rate $\eta$ which is independent of $\varepsilon$. Thanks to the almost everywhere convergence (2.59) and the convexity of $E$, we can pass to the limit $\varepsilon \to 0$ in (2.61), and end up with

$$E(e(t)|e_\infty) \leq \liminf_{\varepsilon \to 0} E(e^\varepsilon(t)|e_\infty) \leq e^{-\eta t} E(e_0|e_\infty),$$

which, combined with the Csiszár-Kullback-Pinsker inequality, allows us to finish the proof of Thm. 1.2.

3. Complex balanced systems with boundary equilibria

In the previous section, we have shown that for a complex balanced reaction-diffusion system without boundary equilibria, all solution trajectories converge exponentially fast to the unique strictly positive complex balance equilibrium thanks to a functional entropy-entropy-dissipation inequality.

It was also pointed out in the introduction that if a system has boundary equilibria, then such an entropy entropy-dissipation estimate cannot hold with the same generality. Indeed, if a trajectory approaches a boundary equilibrium, then the entropy dissipation tends to zero while the relative entropy functional towards the positive complex balance equilibrium remains strictly positive, see (3.9). The question of the large time behaviour for general complex balanced reaction-diffusion systems possessing boundary equilibria is therefore an open problem.

In this section, we give some partial answers. We first consider the two by two model (1.21), and then the three by three model (1.22). In the two by two model, the boundedness of solutions away from the boundary equilibrium $(0, M)$ follows from a comparison-principle argument. The instability of the boundary equilibrium $(0, 0, M)$ for the three by three model, however, turns out be tricky and makes system (1.22) a more interesting example.

In fact, we do not obtain the instability directly. To establish Thm. 1.5, we instead prove that if a trajectory should approach the boundary equilibrium, then the rate of this convergence process cannot be faster than $1/(1 + t)$. That (too) slow convergence is sufficient to apply an entropy entropy-dissipation-like estimate which yields instead convergence to the unique strictly positive complex balance equilibrium with an algebraic rate.

In a second step, thanks to this algebraically fast convergence, it follows after some positive time $T > 0$ that such trajectories remain outside of a neighbourhood of the boundary equilibrium. Hence, another entropy entropy-dissipation estimate can be proven, which implies exponential convergence to the unique strictly positive complex balance equilibrium.

**Remark 3.1.** It is easy to see that the result of Theorem 1.5 can be extended to the following class of complex balanced reaction networks:

$$A \xrightarrow{k_1} \alpha B + C$$

$$\xrightarrow{k_3} (\alpha + 1)B$$

(3.1)

with $\alpha \in \mathbb{N}_{>0}$. All these systems possess the boundary equilibrium $(0, 0, M)$. 
It is important to notice a substantial difference between the ODE and the PDE settings of (1.21) and (1.22). In the ODE setting it can be easily verified that any trajectory starting away from the boundary equilibria (even when starting at the boundary $\partial \mathbb{R}^d_+$ or $\partial \mathbb{R}^d_-$, respectively), will enter the interior of the positive quadrant and thus eventually converge to the unique strictly positive complex balance equilibrium. The PDE setting, however, allows the existence of a non-trivial class of initial data, for which solutions converge to the boundary equilibrium. We present such an example for system (1.22) in Rmk. 3.4.

We now start the

**Proof of Prop. 1.3.** We show first the existence of a unique positive equilibrium and a boundary equilibrium for (1.21). Indeed, the system (1.21) satisfies the mass conservation

$$\int_{\Omega} (a(x, t) + b(x, t)) dx = M := \int_{\Omega} (a_0(x) + b_0(x)) dx \quad \forall t > 0.$$  

Thus, an equilibrium of (1.21) solves

$$\begin{cases}
a_\infty^2 = a_\infty b_\infty, \\
a_\infty + b_\infty = M,
\end{cases}$$

so that there exists a unique strictly positive equilibrium $(a_\infty, b_\infty) = (M/2, M/2)$, and there also exists a boundary equilibrium $(a^*, b^*) = (0, M)$.

Next, we show the uniform propagation of lower and upper bounds for solutions to (1.21). While comparison principles cannot be expected to hold for general systems of parabolic equations, the specific structure of the $2 \times 2$ system (1.21) allows to prove the following *a priori* bounds and global existence of classical solutions:

**Lemma 3.1** (Uniform propagation of positive lower and upper bounds and existence of global classical solutions). Under the assumption of Prop. 1.3, there exists a unique global classical solution $(a, b)$ to (1.21) satisfying the same bounds

$$\varepsilon^2 \leq a(t, x), b(t, x) \leq \Lambda \quad \text{for all} \quad x \in \Omega, \quad t > 0. \quad (3.2)$$

**Proof.** The propagation of the claimed strictly positive lower and upper bounds to solutions of system (3.1) follows, for instance, from comparison principle arguments. Following e.g. [Kir90], one can argue that if one of the (possibly regularised) solutions $a$ or $b$ should violate these positive lower or upper bounds at some position $x_0 \in \Omega$ for a first time at $t_0 > 0$, then the right hand side of the corresponding equations (for the concentration minus the lower/upper bound) has the proper sign to yield a contradiction with the parabolic minimum/maximum principle for such extremal points.

Alternatively, the following argument (which works for the same structural reasons as the comparison principle argument) equally allows to prove the propagation of the positive lower and upper bounds as limits of a hierarchy of corresponding $L^p$-bounds. We formally compute (the formal computation can easily be made rigorous via approximating systems with strictly positive solutions e.g. by adding vanishing positive source terms) the quantity leading to the lower bound (the upper bound can be proven in a similar way):

$$\frac{d}{dt} \int_{\Omega} \left( \frac{1}{pa^p} + \frac{1}{pb^p} \right) dx$$

$$\quad = - \int_{\Omega} d_a(p + 1)a^{-p-2}|\nabla a|^2 dx - \int_{\Omega} d_b(p + 1)b^{-p-2}|\nabla b|^2 dx + \int_{\Omega} a(a^{-p-1} - b^{-p-1})(a - b) dx \leq 0.$$

Thus,

$$\sup_{t > 0} \left\{ \int_{\Omega} \left( \frac{1}{pa^p} + \frac{1}{pb^p} \right) dx \right\} \leq \int_{\Omega} \left( \frac{1}{a_0^p} + \frac{1}{b_0^p} \right) dx \xrightarrow{p \to \infty} \sup_{t \geq 0} \left\{ \frac{1}{a(t)} _{L^p} + \frac{1}{b(t)} _{L^p} \right\} \leq \frac{1}{a_0} _{L^p} + \frac{1}{b_0} _{L^p}.$$

Finally, due to the uniform-in-time $L^\infty$-bounds (3.2), the existence of a unique global classical solution to (1.21) (which satisfies the positive lower and upper bounds point-wise in $x$) follows from classical arguments.

Lemma 3.1 implies that initial data $(a_0, b_0)$, which are a.e. bounded away from the boundary equilibrium $(0, M)$, yield solutions $(a(\cdot, t), b(\cdot, t))$, which are uniformly-in-time bounded away from the same boundary equilibrium.
Then, direct computations show that
\[
\frac{1}{2} \frac{d}{dt} \left( \|a(\cdot,t) - a_\infty\|^2 + \|b(\cdot,t) - b_\infty\|^2 \right) = -d_a \int_\Omega |\nabla a|^2 dx - d_b \int_\Omega |\nabla b|^2 dx - \int_\Omega (a - b)^2 dx
\]
\[
\leq -d_a \int_\Omega |\nabla a|^2 dx - d_b \int_\Omega |\nabla b|^2 dx - \varepsilon^2 \int_\Omega (a - b)^2 dx,
\]
thanks to Lemma 3.1. Due to the mass conservation $\overline{a}(t) + \overline{b}(t) = M$ and previously established entropy-dissipation estimates for detailed balance reaction-diffusion systems without boundary equilibria (see e.g. [DF06, DF08, FT]), we obtain, for all $t > 0$,
\[
d_a \int_\Omega |\nabla a|^2 dx + d_b \int_\Omega |\nabla b|^2 dx + \varepsilon^2 \int_\Omega (a - b)^2 dx \geq \frac{\lambda_c}{2} (\|a(\cdot,t) - a_\infty\|^2 + \|b(\cdot,t) - b_\infty\|^2),
\]
where $\lambda_c$ depends only on $\Omega$, $M$, $d_a$, $d_b$ (and $\varepsilon$) and moreover $\lambda_c = O(\varepsilon^2)$. By inserting the above estimate into (3.3), exponential convergence to equilibrium follows from a standard Gronwall inequality. ■

**Proof of Cor. 1.4.** First, we remark that classical existence results imply for each continuous nonnegative initial data a unique global classical nonnegative (in the sense that both components are nonnegative) solution $(a, b)$ to the system (1.21). Moreover, these solutions satisfy an $L^\infty$ upper bound as in the proof of Prop. 1.3. Thus, we get the estimate
\[
\begin{cases}
a_t - d_a \Delta a \geq -\|a\|_{L^\infty} a, \\
a_0(x) \geq 0 \quad \text{with} \quad \int_\Omega a_0(x) \, dx > 0.
\end{cases}
\]
Hence, the strong minimum principle for the heat-equation with Neumann boundary conditions (see e.g. [Pao92]) implies for classical solutions that for any $\tau > 0$ there exists an $\varepsilon_a(\tau) > 0$ such that $a(x, \tau) \geq \varepsilon_a(\tau)$. Otherwise, any point where $a(x, \tau) = 0$ (with $x \in \Omega$) would be a minimum and yield a contradiction with the minimum principle except if $a_0(x) \equiv 0$ on $\Omega$, which we have excluded of the set of admissible initial data (note that for homogeneous Dirichlet data, we would indeed only have $a(x, \tau) > 0$ on $\Omega$). The same argument holds for $b(x, t)$ provided that $\int_\Omega b_0(x) \, dx > 0$. In the special case when $\int_\Omega b_0(x) \, dx = 0$, we observe that
\[
\frac{d}{dt} \int_\Omega b(x, t) \, dx \bigg|_{t=0} \geq \int_\Omega a_0(x) \, dx > 0.
\]
Hence, for any (sufficiently small) $\tau/2 > 0$, we see that $\int_\Omega b(x, \tau/2) \, dx > 0$ and, via an analog minimum principle argument, that $b(x, \tau) \geq \varepsilon_b(\tau)$ for some $\varepsilon_b(\tau) > 0$. As a consequence, the statement of Corollary 1.4 follows from applying Proposition 1.3 for $t \geq \tau > 0$.

For the second statement of Cor. 1.4, we observe that $a_0 \equiv 0$ and (1.21) imply that the solution $a(x, t) \equiv 0$ for all times, and thus
\[
\begin{cases}
b_t - d_b \Delta b = 0, & x \in \Omega, \quad t > 0, \\
b_0(x) \geq 0, & x \in \Omega, \quad \int_\Omega b_0(x) \, dx = M.
\end{cases}
\]
Therefore, by standard results for the heat equation and for all $t > 0$, we have
\[
\|b(t) - M\|^2 \leq e^{-d_b t \lambda_P} \|b_0 - M\|^2,
\]
where $\lambda_P$ is the best constant in Poincaré-Wirtinger’s inequality $\|\nabla f\|^2 \geq \lambda_P \|f - \overline{f}\|^2$ on the domain $\Omega$. ■

**Remark 3.2** (Higher regularity). By using the $L^p$ bounds on $a$ and $b$, we can actually obtain exponential convergence to equilibrium in any $L^p$-norm with $1 \leq p < +\infty$ via interpolation. Moreover, following e.g. [DF08], one can verify that any Sobolev norm of classical solutions to systems (1.21) (which matches the assumed regularity of the boundary $\partial \Omega$) will grow at most polynomially-in-time, so that exponential convergence to equilibrium in a (slightly lower) Sobolev norm follows again via interpolation with (1.3).

**Remark 3.3** (Generalised $2 \times 2$ systems). The above arguments can be generalised to obtain exponential convergence to equilibrium for the following class of systems
\[
\begin{cases}
a_t - d_a \Delta a = -\varphi(a)(r(a) - r(b)), & x \in \Omega, \quad t > 0, \\
b_t - d_b \Delta b = \varphi(a)(r(a) - r(b)), & x \in \Omega, \quad t > 0,
\end{cases}
\]
supplemented by homogeneous Neumann boundary condition and initial data satisfying the bounds (3.2), where $\varphi : [0, +\infty) \to [0, +\infty)$ is continuously increasing and $r : [0, +\infty) \to [0, +\infty)$ and its derivative $r'$ are continuously increasing.
As example, we can consider reactions of the form
\[(n + m)A \rightleftharpoons nA + mB\]
with \(n \geq 0\) and \(m \geq 1\), which are described by the mass action law systems
\[
\begin{aligned}
  a_x - d_x \Delta a &= -m a^n(a^m - b^m), & x &\in \Omega, & t &> 0, \\
  b_x - d_b \Delta b &= m a^n(a^m - b^m), & x &\in \Omega, & t &> 0,
\end{aligned}
\]
corresponding to the cases \(\varphi(z) = mz^n\) and \(r(z) = z^m\).

We now consider the complex-balanced chemical reaction network defined by (1.22). In strict contrast to the 2 \times 2 system (1.21), general reaction-diffusion systems like (1.22) do not allow to show uniform-in-time propagation of upper and positive lower a priori \(L^{\infty}\)-estimates.

The existence of global classical solutions to (1.22) in space dimensions \(N \leq 5\) follows from bootstrap arguments (in the spirit of e.g. [Rot84]) thanks to the specific structure of (1.22), and since the nonlinear reaction terms are at most quadratic. In higher space dimensions \(N \geq 6\), existence of global classical solutions can be shown under a “closeness” condition on the diffusion coefficients by following e.g. [CDF14]. The (rather standard) bootstrap argument along with the complete proof of the following Lemma stating existence of global, classical solutions is given in the Appendix.

**Lemma 3.2** (Global classical solutions).
Let \(\Omega\) be a bounded smooth (\(C^2\)) domain of \(\mathbb{R}^N\), and \(d_a, d_b, d_c > 0\), \(k_1, k_2, k_3 > 0\). Assume nonnegative initial data \(a_0, b_0, c_0 \in L^{\infty}(\Omega)\), and denote
\[
\delta := \max\{d_a, d_b, d_c\} - \min\{d_a, d_b, d_c\}.
\]
Consider either \(1 \leq N \leq 5\) or \(N \geq 6\) and \(\delta > 0\) sufficiently small (depending on \(N\) as specified in the Appendix).

Then, there exists a unique, nonnegative, global classical solution \((a, b, c)\) to system (1.22), which satisfies the following \(L^{\infty}\)-bound for all \(T > 0\):
\[
\|a\|_{L^{\infty}(\Omega_T)} + \|b\|_{L^{\infty}(\Omega_T)} + \|c\|_{L^{\infty}(\Omega_T)} \leq C(T),
\]
where \(C(T)\) grows at most polynomially with respect to \(T\).

Moreover, \(a, b, c\) satisfy the conservation of mass
\[
\int_{\Omega} (2a(x,t) + b(x,t) + c(x,t)) \, dx = M := \int_{\Omega} (2a_0(x) + b_0(x) + c_0(x)) \, dx \quad \text{for all} \quad t > 0. \quad (3.4)
\]

Considering system (1.22), we observe that any equilibrium \((a_\infty, b_\infty, c_\infty)\) solves the stationary state conditions
\[
\begin{aligned}
  -k_1 a_\infty + k_3 b_\infty^2 &= 0, \\
  k_1 a_\infty + k_2 b_\infty c_\infty - 2k_3 b_\infty^2 &= 0, \\
  k_1 a_\infty - k_2 b_\infty c_\infty &= 0,
\end{aligned}
\]
and satisfies the mass conservation law
\[2a_\infty + b_\infty + c_\infty = M.\]

Thus, it follows that system (1.22) features two equilibria, first the unique positive complex-balance equilibrium
\[(a_\infty, b_\infty, c_\infty) \quad \text{where} \quad a_\infty = \frac{k_3}{k_1} b_\infty^2, \quad b_\infty = \frac{-k_1 + \sqrt{k_1(k_1 + 2k_2 M)}}{2k_2} \quad \text{and} \quad c_\infty = \frac{k_3}{k_2} b_\infty, \quad (3.5)\]
and secondly the boundary equilibrium
\[(a^*, b^*, c^*) = (0, 0, M).\]

For notational convenience, we shall write \(c = (a, b, c)\) and \(c_\infty = (a_\infty, b_\infty, c_\infty)\). At some places we shall also use (coherently with the first sections of this paper) \(c_1 := a\), \(c_2 := b\) and \(c_3 := c\) (and similarly for \(c_{i, \infty}\) with \(i = 1, 2, 3\)).

The relative entropy writes as (cf. (1.10))
\[
\mathcal{E}(c|c_\infty) = \sum_{i=1}^{3} \int_{\Omega} \left( c_i \log \frac{c_i}{c_{i, \infty}} - c_i + c_{i, \infty} \right) \, dx, \quad (3.6)
\]
and its entropy dissipation is given by (cf. Prop. 2.1)
\[
D(c) = d_a \int_\Omega \frac{\nabla a}{a}^2 dx + d_b \int_\Omega \frac{\nabla b}{b}^2 dx + d_c \int_\Omega \frac{\nabla c}{c}^2 dx
+ \int_\Omega \left[ k_1 a_{\infty} \Psi \left( \frac{a}{a_{\infty}} \right) + k_2 b_{\infty} c_{\infty} \Psi \left( \frac{bc}{b_{\infty} c_{\infty}} \right) + k_3 b_{\infty} \Psi \left( \frac{b_3}{b_{\infty}} \right) \right] dx. \tag{3.7}
\]

The uniform-in-time \textit{a priori} estimates for the average \( \bar{\rho}(t) = \int_\Omega \rho(x,t) \, dx \) are a natural consequence of the mass conservation law (3.4) and the nonnegativity of the solutions:
\[
2\bar{\rho}(t) + \bar{\eta}(t) + \rho(t) = M \quad \text{for all } \quad t \geq 0.
\tag{3.8}
\]

It is easy to see that if \( \{c_k \} = (a_k, b_k, c_k) \) is a space homogeneous sequence satisfying (3.8) and converging to \((0,0,M)\), then
\[
\lim_{k \to \infty} D(c_k) = 0 \quad \text{and} \quad \lim_{k \to \infty} \mathcal{E}(c_k|c_{\infty}) = a_{\infty} + b_{\infty} + \Psi(M; c_{\infty}) > 0,
\tag{3.9}
\]
that means there does not exist any \( \lambda > 0 \) such that \( D(c) \geq \lambda \mathcal{E}(c|c_{\infty}) \) for all \( c \) satisfying the mass conservation (3.8). Hence, the convergence of solutions to (1.22) towards the strictly positive complex balance equilibrium is not clear. In fact, similarly to the Cor.1.4, the following example shows that there is a non-trivial class of initial data, for which solutions to (1.22) will converge to the boundary equilibrium.

**Remark 3.4 (Convergence towards boundary equilibrium).**
Consider initial data \( a_0 = b_0 = c_0 \equiv M \), yet with \( c_0 \neq M \) (so that the initial data \( (a_0, b_0, c_0) \) is not the boundary equilibrium \((0,0,M)\)).

By arguing similarly to the second part of Cor. 1.4, we obtain that \( a(t) = b(t) \equiv 0 \) (for all \( t \geq 0 \)), while \( c := c(t) \) is given by the solution of the homogeneous heat equation (with homogeneous Neumann conditions), and thus converges exponentially fast to \( M \) as
\[
||c(t) - M||^2 \leq e^{-\lambda_P t} ||c_0 - M||^2,
\]
where \( \lambda_P \) is the constant in Poincaré-Wirtinger’s inequality \( ||\nabla f||^2 \geq \lambda_P \|f - \bar{f}\|^2 \) on the domain \( \Omega \).

We now start the

**Proof of Theorem 1.5.** We consider in this proof a classical solution \( c := c(x,t) \) of system (1.22).

We first use the assumption \( \|b_0 \|_{L^\infty(\Omega)} < \infty \) to show that
\[
\inf_{x \in \Omega} b(x,t) \geq h(t) := \frac{1}{\|b_0 \|_{L^\infty(\Omega)} + 2k_3 t}, \quad \text{for all} \quad t \geq 0. \tag{3.10}
\]
Indeed thanks to the continuity of classical solutions, there exists a \( t^* > 0 \) such that \( \|b(t,\cdot)^{-1}\|_{L^\infty} \leq 2 \|b_0^{-1}\|_{L^\infty} \) for all \( t \in [0,t^*] \). Thus, for all \( t \in [0,t^*] \), we can compute
\[
\partial_t \left( \frac{1}{b} \right) - d_b \Delta \left( \frac{1}{b} \right) = - \frac{k_1 a}{2b} \frac{2k_3 c}{b^2} + 2k_3 - 2d_b \frac{\nabla b}{b} \leq 2k_3.
\]
Thus, using the maximum principle, we obtain
\[
\left\| \frac{1}{b(t)} \right\|_{L^\infty(\Omega)} \leq \left\| \frac{1}{b_0} \right\|_{L^\infty(\Omega)} + 2k_3 t, \quad \text{for all} \quad t \in [0,t^*],
\tag{3.11}
\]
which implies the desired estimate for the time-interval \([0,t^*]\). Furthermore, it is easy to see that this argument can be iterated in time, and eventually yields the bound (3.11) for all \( t > 0 \).

We see therefore that if \( b \) decays to zero, then it decays at most with a rate of the form \((1 + t)^{-1}\). We will see that this lower bound, in combination with the entropy entropy-dissipation structure, is sufficient to show that it is impossible for the solution to converge to the boundary equilibrium.

As previously, \( \mathcal{E} \) defined by (3.6) satisfies
\[
\mathcal{E}(c|c_{\infty}) = \mathcal{E}(c|\bar{c}) + \mathcal{E}(\bar{c}|c_{\infty}),
\tag{3.12}
\]
and via the Logarithmic Sobolev Inequality, we get
\[
\frac{1}{2} D(c) \geq \lambda_1 \mathcal{E}(c|\bar{c}). \tag{3.13}
\]
We denote \( F(c) = \lambda_1 \sum_{i=1}^3 \Psi(c_i; c_{i,\infty}) \). By using estimate (3.11) (for all \( t \geq 0 \), the Logarithmic Sobolev Inequality and the additivity property (3.12), we see that (with \( h(t) \) being defined in (3.10))
\[
\frac{1}{2} D(c) \geq \lambda_1 \left( \mathcal{E}(c; c_{\infty}) - \mathcal{E}(\Phi(c) - F(\Phi)) \right) + \frac{1}{2} \int_\Omega \left[ k_1 a, \Psi \left( \frac{a}{a_{\infty}} - \frac{bc}{b_{\infty} c_{\infty}} \right) + k_2 c, \h(t) \Psi \left( \frac{c}{c_{\infty}} - \frac{b}{b_{\infty}} \right) + k_3 b^2, \Psi \left( \frac{b^2}{b_{\infty}^2} - \frac{a}{a_{\infty}} \right) \right] dx \\
\geq \alpha h(t) \left\{ \int_\Omega \left[ F(c) + \Psi \left( \frac{a}{a_{\infty}} - \frac{bc}{b_{\infty} c_{\infty}} \right) + \Psi \left( \frac{c}{c_{\infty}} - \frac{b}{b_{\infty}} \right) + \Psi \left( \frac{b^2}{b_{\infty}^2} - \frac{a}{a_{\infty}} \right) \right] dx - F(\Phi) \right\} := \phi(c) 
\]
where \( \alpha = \| a_{\infty} \|_{L^\infty} \) and \( : \Phi := \Psi \left( \frac{a}{a_{\infty}} - \frac{bc}{b_{\infty} c_{\infty}} \right) + \Psi \left( \frac{c}{c_{\infty}} - \frac{b}{b_{\infty}} \right) + \Psi \left( \frac{b^2}{b_{\infty}^2} - \frac{a}{a_{\infty}} \right) \).

We also recall that \( \hat{\Phi} \) denotes the convexification of \( \Phi \). Define
\[
\mathcal{E}_M = \{ \xi = (a, b, c) \in \mathbb{R}^3_+: 2a + b + c = M \text{ and } \mathcal{E}(\xi|c_{\infty}) \leq \mathcal{E}(c_0|c_{\infty}) \}.
\]
the set of all concentrations satisfying the mass conservation (3.4) and suitably bounded relative entropy. Note that
\[
\xi \in \mathcal{E}_M \text{ and } Q(\xi) = 0 \iff \xi \equiv c_{\infty},
\]
and
\[
\hat{\Phi}(\xi) = F(\xi) \geq \hat{F}(\xi) + \hat{Q}(\xi) = \hat{Q}(\xi) \geq 0.
\]

Hence,
\[
\inf_{\xi \in \mathcal{E}_M} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_{\infty})} > 0 \text{ if } \liminf_{\mathcal{E}_M} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_{\infty})} > 0.
\]

On the other hand, we can estimate by using Lemma 2.5 and Taylor expansion around \( c_{\infty} \) (see Proposition 2.2 or [MHM15, Proposition 3.3])
\[
\inf_{\xi \in \mathcal{E}_M} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_{\infty})} \geq \liminf_{\mathcal{E}_M} \frac{\hat{Q}(\xi)}{\mathcal{E}(\xi|c_{\infty})} = \liminf_{\mathcal{E}_M} \frac{Q(\xi)}{\mathcal{E}(\xi|c_{\infty})} > 0.
\]

Finally, we see that
\[
\inf_{\xi \in \mathcal{E}_M} \frac{\hat{\Phi}(\xi) - F(\xi)}{\mathcal{E}(\xi|c_{\infty})} \geq \beta > 0,
\]
and consequently
\[
\frac{1}{2} D(c) \geq \alpha \beta h(t) \mathcal{E}(\Phi(c) - F(\Phi)) \quad \text{for all } t > 0,
\]
(3.14)
which, combining with (3.13), implies
\[
D(c) \geq \gamma h(t) \mathcal{E}(\Phi(c) - F(\Phi)) \quad \text{for all } t > 0,
\]
(3.15)
where \( \gamma = \min \left\{ \alpha ; \lambda_1 \left\| \frac{1}{b} \right\|_{L^\infty(\Omega)} \right\} \). Then, thanks to Gronwall’s lemma, we see that for all \( t \geq 0 \),
\[
\mathcal{E}(c(t)|c_{\infty}) \leq \mathcal{E}(c_0|c_{\infty}) \left( \left\| \frac{1}{b_0} \right\|_{L^\infty(\Omega)} + 2k_3 t \right)^{-\gamma/2k_3}
\]
and the relative entropy with respect to the complex balance equilibrium decays therefore to zero with the algebraic rate \( \gamma/2k_3 \). Moreover, by using a Csizsár-Kullback-Pinsker inequality (see e.g. Lemma 2.6), we get the estimate
\[
\| a(t) - a_{\infty} \|^2_{L^1(\Omega)} + \| b(t) - b_{\infty} \|^2_{L^1(\Omega)} + \| c(t) - c_{\infty} \|^2_{L^1(\Omega)} \leq \mathcal{E}(c_0|c_{\infty}) \left( \left\| \frac{1}{b_0} \right\|_{L^\infty(\Omega)} + 2k_3 t \right)^{-\gamma/2k_3}.
\]

As a consequence, there exists for any (sufficiently small) \( \varepsilon > 0 \) a time \( T_\varepsilon \) such that solutions are bounded away from the boundary equilibrium in the sense of \( L^1 \), i.e.
\[
\| a(t) \|^2_{L^1(\Omega)} \geq \varepsilon^2, \quad \| b(t) \|^2_{L^1(\Omega)} \geq \varepsilon^2
\]
and \( ||c(t)||_{L^1(\Omega)} \geq \varepsilon^2 \) for all \( t \geq T_\varepsilon \). These \( L^1 \)-bounds away from the boundary equilibrium allow to apply a specialised entropy entropy-dissipation estimate proven in Lemma 3.3 below. Finally, via another Gronwall arguments, we obtain
\[
\mathcal{E}(c(t)\varepsilon) \leq e^{-\lambda_t \varepsilon} \mathcal{E}(c_0\varepsilon) \quad \text{for all} \quad t \geq T_\varepsilon
\]
which consequently, together with another use of a Csiszár-Kullback-Pinsker inequality, implies Theorem 1.5.

**Remark 3.5.** We remark that since we prove estimate (3.14) via the convexification technique, we get a non-explicit constant \( \gamma \) in (3.15). In fact, \( \gamma \) can be explicitly estimated by using the constructive method in Subsection 2.3, at the price of a much longer proof.

**Lemma 3.3** (Entropy-entropy-dissipation estimate). Consider a positive initial mass \( M > 0 \).

Then, for any nonnegative, measurable functions \( a, b, c : \Omega \to \mathbb{R}_+ \) satisfying the mass conservation law
\[
\int_{\Omega} (2a(x) + b(x) + c(x)) \, dx = M,
\]
and the lower bounds \( ||a||_{L^1(\Omega)} \geq \varepsilon^2 \), \( ||b||_{L^1(\Omega)} \geq \varepsilon^2 \) and \( ||c||_{L^1(\Omega)} \geq \varepsilon^2 \), the functional inequality
\[
\mathcal{D}(\varepsilon) \geq \lambda_c \mathcal{E}(\varepsilon c_\infty)
\]
holds, for some explicit constant \( \lambda_c > 0 \), which depends only on the initial mass \( M \), the domain \( \Omega \), the diffusion coefficients \( d_a, d_b, d_c \), the reaction rates \( k_1, k_2, k_3 \), and on \( \varepsilon \). Here \( \mathcal{E}(\varepsilon c_\infty) \) and \( \mathcal{D}(\varepsilon) \) are defined by (3.6) and (3.7), while \( c_\infty \) is defined by (3.5). Moreover, we have
\[
\lambda_c = O(\varepsilon^6) \quad \text{as} \quad \varepsilon \to 0.
\]

**Proof.** We will apply the method described in Subsection 2.3. We already know that (3.13) (that is, Step 1), holds.

In Step 2, we bound the entropy dissipation \( \mathcal{D}(\varepsilon) \) below and the relative entropy \( \mathcal{E}(\varepsilon c_\infty) \) above by quadratic terms for the square root concentrations \( C := (A, B, C) \):
\[
\mathcal{D}(\varepsilon) \geq K_0 \left( \|A\|^2 + \|B\|^2 + \|C\|^2 + \left\| \frac{A}{A_\infty} - \frac{B^2}{B_{\infty}^2} \right\|^2 + \left\| \frac{B^2}{B_{\infty}^2} - \frac{BC}{B_{\infty}C_{\infty}} \right\|^2 + \left\| \frac{BC}{B_{\infty}C_{\infty}} - \frac{A}{A_\infty} \right\|^2 \right),
\]
and
\[
\mathcal{E}(\varepsilon c_\infty) \leq K_2 \left( \left( \sqrt{A^2} - A_\infty \right)^2 + \left( \sqrt{B^2} - B_\infty \right)^2 + \left( \sqrt{C^2} - C_\infty \right)^2 \right).
\]

Step 3 enables to estimate the entropy dissipation \( \mathcal{D}(\varepsilon) \) in terms of reaction terms for averaged quantities:
\[
\mathcal{D}(\varepsilon) \geq K_1 \left[ \|\nabla A\|^2 + \|\nabla B\|^2 + \|\nabla C\|^2 + \left( \frac{A}{A_\infty} - \frac{B^2}{B_{\infty}^2} \right)^2 + \left( \frac{B^2}{B_{\infty}^2} - \frac{BC}{B_{\infty}C_{\infty}} \right)^2 + \left( \frac{BC}{B_{\infty}C_{\infty}} - \frac{A}{A_\infty} \right)^2 \right].
\]

It remains to find \( K_3 > 0 \) such that
\[
\|\nabla A\|^2 + \|\nabla B\|^2 + \|\nabla C\|^2 + \left( \frac{A}{A_\infty} - \frac{B^2}{B_{\infty}^2} \right)^2 + \left( \frac{B^2}{B_{\infty}^2} - \frac{BC}{B_{\infty}C_{\infty}} \right)^2 + \left( \frac{BC}{B_{\infty}C_{\infty}} - \frac{A}{A_\infty} \right)^2 \geq K_3 \left( \left( \sqrt{A^2} - A_\infty \right)^2 + \left( \sqrt{B^2} - B_\infty \right)^2 + \left( \sqrt{C^2} - C_\infty \right)^2 \right).
\]

We exploit the ansatz
\[
\overline{A} = A_\infty(1 + \mu_A)^2, \quad \overline{B} = B_{\infty}(1 + \mu_B)^2, \quad \overline{C} = C_\infty(1 + \mu_C)^2, \quad \text{with} \quad \mu_A, \mu_B, \mu_C \in [-1, +\infty).
\]

Thanks to the natural upper bounds for \( \overline{A}^2, \overline{B}^2 \) and \( \overline{C}^2 \), we have more precisely
\[
-1 \leq \mu_A, \mu_B, \mu_C \leq \mu_{\max} < +\infty,
\]
for some \( \mu_{\max} > 0 \). We again denote \( \delta_A(x) = A(x) - \overline{A} \), \( \delta_B(x) = B(x) - \overline{B} \) and \( \delta_C(x) = C(x) - \overline{C} \), and recall that
\[
\overline{A} = A_\infty(1 + \mu_A) - \frac{||\delta_A||^2}{\sqrt{A^2 + \overline{A}}} =: A_\infty(1 + \mu_A) - R(A)||\delta_A||^2.
\]
Similarly,
\[
\overline{B} = B_{\infty}(1 + \mu_B) - R(B)||\delta_B||^2 \quad \text{and} \quad \overline{C} = C_{\infty}(1 + \mu_C) - R(C)||\delta_C||^2.
\]
Thanks to the assumption $\overline{A^2} \geq \varepsilon^2$, $\overline{B^2} \geq \varepsilon^2$ and $\overline{C^2} \geq \varepsilon^2$, we see that 
$$\|R(A), R(B), R(C)\| \leq \frac{1}{\varepsilon}.$$ 
By using computations similar to (2.45), we get 
$$\left(\frac{\overline{A}}{A_\infty} - \frac{\overline{B}}{B_\infty}\right)^2 \geq \frac{1}{2} \left[ (1 + \mu_A) - (1 + \mu_B) \right]^2 - 3\left(\frac{\|\delta A\|}{A_\infty}\right)^2 (1 + \mu_B)^2 R(B)^2 - 3\|\delta B\|^2 R(B)^2 - 3\|\delta B\|^2 \|\delta B\|^2 - 3\|\delta B\|^8 R(B)^4.$$ 
Then, thanks to the boundedness of $\|\delta A\|$ and $\|\delta B\|$ and $1/\varepsilon^2 \leq 1/\varepsilon^4$ for $\varepsilon \in [0, 1]$, we can estimate 
$$\left(\frac{\overline{A}}{A_\infty} - \frac{\overline{B}}{B_\infty}\right)^2 \geq \frac{1}{2} \left[ (1 + \mu_A) - (1 + \mu_B) \right]^2 - \frac{K}{\varepsilon^4}\left(\|\delta A\|^2 + \|\delta B\|^2\right),$$ 
where $K$ is independent of $\varepsilon$. Similarly, 
$$\left(\frac{\overline{B}}{B_\infty} - \frac{\overline{BC}}{BC_\infty}\right)^2 \geq \frac{1}{2} \left[ (1 + \mu_B) \right]^2 (\mu_B - \mu_C)^2 - \frac{K}{\varepsilon^4}\left(\|\delta B\|^2 + \|\delta C\|^2\right)$$ 
and 
$$\left(\frac{\overline{BC}}{BC_\infty} - \frac{\overline{A}}{A_\infty}\right)^2 \geq \frac{1}{2} \left[ (1 + \mu_B)(1 + \mu_C) - (1 + \mu_A) \right]^2 - \frac{K}{\varepsilon^4}\left(\|\delta A\|^2 + \|\delta B\|^2 + \|\delta C\|^2\right).$$ 
The constant $K$ above may vary but is always independent of $\varepsilon$. Therefore, the left hand side of (3.17) is bounded below by 
$$K_4(\varepsilon) \left( B\right) \left[ (1 + \mu_A) - (1 + \mu_B)^2 \right] + (1 + \mu_B)^2 (\mu_B - \mu_C)^2 + (1 + \mu_B)(1 + \mu_C) - (1 + \mu_A)^2,$$ 
with $K_4(\varepsilon) = O(\varepsilon^4)$, while the right hand side of (3.17) is bounded above by 
$$K_3 \max\{A_\infty^2, B_\infty^2, C_\infty^2\} \left(\mu_A^2 + \mu_B^2 + \mu_C^2\right).$$ 
Therefore, we only have to show for some $K_5 > 0$ (and $K_5 = O(\varepsilon^2)$ as $\varepsilon \rightarrow 0$), that 
$$\left[ (1 + \mu_A) - (1 + \mu_B)^2 \right] + (1 + \mu_B)^2 (\mu_B - \mu_C)^2 + (1 + \mu_B)(1 + \mu_C) - (1 + \mu_A)^2 \geq K_5 \left(\mu_A^2 + \mu_B^2 + \mu_C^2\right),$$ 
under the constraint imposed by the mass conservation law 
$$2A_\infty^2 \mu_A(\mu_A + 2) + B_\infty^2 \mu_B(\mu_B + 2) + C_\infty^2 \mu_C(\mu_C + 2) = 0.$$ 
(3.19) 
To prove (3.18), we first use $\overline{B^2} \geq \varepsilon^2$, which leads to $(1 + \mu_B)^2 \geq \frac{\varepsilon^2}{B_\infty^2}$ and 
$$2A_\infty^2 \mu_A(\mu_A + 2) + B_\infty^2 \mu_B(\mu_B + 2) + C_\infty^2 \mu_C(\mu_C + 2) = 0.$$ 
(3.19) 
The mass conservation law (3.19) implies then that only the following three cases concerning the signs of $\mu_A, \mu_B$ and $\mu_C$ can appear: 
(I) $\mu_A$ and $\mu_B$ have different signs, 
(II) $\mu_A \geq 0$, $\mu_B \geq 0$ and $\mu_C \leq 0$, 
(III) $\mu_A \leq 0$, $\mu_B \leq 0$ and $\mu_C \geq 0$. 
We will treat each case separately. 
Case (I): $\mu_A$ and $\mu_B$ have different signs. In this case, we see that 
$$\left[ (1 + \mu_A) - (1 + \mu_B)^2 \right] \geq \mu_A^2 + \mu_B^2.$$ 
Indeed, if $\mu_A \geq 0$ and $\mu_B \leq 0$, 
$$\left[ (1 + \mu_A) - (1 + \mu_B)^2 \right] = (\mu_A - \mu_B - \mu_B(1 + \mu_B))^2 \geq (\mu_A - \mu_B)^2 \geq \mu_A^2 + \mu_B^2$$ 
since $\mu_A - \mu_B \geq 0$ and $-\mu_B(1 + \mu_B) \geq 0$. If $\mu_A \leq 0$ and $\mu_B \geq 0$, then 
$$\left[ (1 + \mu_A) - (1 + \mu_B)^2 \right] = (\mu_B - \mu_A + \mu_B(1 + \mu_B))^2 \geq (\mu_B - \mu_A)^2 \geq \mu_B^2 + \mu_A^2.$$
From (3.20) and (3.21), we get
\[
\text{LHS of (3.18)} \geq \mu_A^2 + \mu_B^2 + \frac{\varepsilon^2}{B_\infty^2} (\mu_B - \mu_C)^2 \geq \frac{1}{4} \min \left\{ \frac{\varepsilon^2}{B_\infty^2}, 1 \right\} (\mu_A^2 + \mu_B^2 + \mu_C^2),
\]
and thus obtain (3.18) with
\[
K_5 = \frac{1}{4} \min \left\{ \frac{\varepsilon^2}{B_\infty^2}, 1 \right\}.
\]

Case (II): \( \mu_A \geq 0, \mu_B \geq 0 \) and \( \mu_C \leq 0 \). In this case, we first estimate further (3.20) as
\[
(1 + \mu_B)^2 (\mu_B - \mu_C)^2 \geq \frac{\varepsilon^2}{B_\infty^2} (\mu_B - \mu_C)^2 \geq \frac{\varepsilon^2}{B_\infty^2} (\mu_A^2 + \mu_C^2).
\] (3.22)
There are two possibilities: \( \mu_A \geq \mu_B \) or \( \mu_B \geq \mu_A \). If \( \mu_A \geq \mu_B \), then
\[
[(1 + \mu_B)(1 + \mu_C) - (1 + \mu_A)]^2 = (\mu_A - \mu_B - \mu_C(1 + \mu_B))^2 \geq (\mu_A - \mu_B)^2.
\] (3.23)
If \( \mu_B \geq \mu_A \), then
\[
[(1 + \mu_A) - (1 + \mu_B)]^2 = (\mu_B - \mu_A + \mu_B(1 + \mu_B))^2 \geq (\mu_B - \mu_A)^2.
\] (3.24)
Therefore, thanks to (3.22), (3.23) and (3.24), we can obtain (3.18) in case (II) with
\[
K_5 = \frac{1}{4} \min \left\{ \frac{\varepsilon^2}{B_\infty^2}, 1 \right\}.
\]

Case (III): \( \mu_A \leq 0, \mu_B \leq 0 \) and \( \mu_C \geq 0 \). This case can be treated in the same way as case (II) by using
\[
(1 + \mu_B)^2 (\mu_B - \mu_C)^2 \geq \frac{\varepsilon^2}{B_\infty^2} (\mu_C - \mu_B)^2 \geq \frac{\varepsilon^2}{B_\infty^2} (\mu_A^2 + \mu_B^2),
\]
and considering \( \mu_A \geq \mu_B \) or \( \mu_A \leq \mu_B \). We get from this case (3.18), with again
\[
K_5 = \frac{1}{4} \min \left\{ \frac{\varepsilon^2}{B_\infty^2}, 1 \right\}.
\]

In conclusion, we have proven (3.18), with
\[
K_5 = \frac{1}{4} \min \left\{ \frac{\varepsilon^2}{B_\infty^2}, 1 \right\},
\]
and consequently finished the proof of Lemma 3.3. \( \blacksquare \)

**Remark 3.6.** Obtaining a general stability statement of the strictly positive complex balance equilibrium for general (complex balanced) reaction-diffusion systems featuring boundary equilibria looks quite involved, and remains open for future investigation. We would just like to point out that the difficulty appears only when a trajectory \( t \mapsto \mathbf{c}_\infty(t) \) gets too close to a boundary equilibrium.

More precisely, let us consider a complex balanced system, and denote
\[
\mathcal{B}_E = \{ \mathbf{c}^* \in \partial \mathbb{R}^N_{\geq 0} : \mathbf{c}^* \text{ is a complex balance equilibrium of (1.1)} \}
\]
and
\[
\mathcal{N}_\varepsilon(\mathcal{B}_E) = \{ \mathbf{c} \in \mathbb{R}^N_{> 0} : d(\mathbf{c}, \mathcal{B}_E) \geq \varepsilon \}
\]
an \( \varepsilon \)-neighbourhood of \( \mathcal{B}_E \), where \( d \) is the Euclidean distance in \( \mathbb{R}^n \). Let us assume that there exists \( \varepsilon > 0 \) such that each trajectory of (1.1) starting outside of \( \mathcal{N}_\varepsilon(\mathcal{B}_E) \) remains outside \( \mathcal{N}_\varepsilon(\mathcal{B}_E) \) for all \( t \geq 0 \), that is
\[
\mathbf{c}(t) \in \mathbb{R}^N_{> 0} \setminus \mathcal{N}_\varepsilon(\mathcal{B}_E) \quad \text{for all} \quad t \geq 0.
\]
Then one can show that there exists a constant \( \lambda_\varepsilon > 0 \) depending on the domain \( \Omega \), the diffusion coefficients \( D \), the stoichiometric coefficients, and additionally the parameter \( \varepsilon \), such that
\[
D(\mathbf{c}(t)) \geq \lambda_\varepsilon \mathcal{E}(\mathbf{c}(t)|\mathbf{c}_\infty).
\]
4. Appendix

In this Appendix, we prove Lemma 3.2 about the global existence of classical solution to the system (1.22) and Lemma 2.5. We first need the following classical lemma on the regularity of solutions to the heat equation.

**Lemma 4.1.** [see e.g. [CDF14]] Assume that \( u_0 \in L^\infty(\Omega) \), \( f \in L^p(\Omega_T) \), \( d > 0 \), and that \( u \) is the solution to

\[
\begin{cases}
    u_t - d\Delta u = f, & x \in \Omega, \quad t > 0, \\
    \nabla u \cdot \nu = 0, & x \in \partial\Omega, \quad t > 0, \\
    u(x,0) = u_0(x), & x \in \Omega.
\end{cases}
\]

If \( 1 < p < \frac{N+2}{2} \), then \( u \in L^s(\Omega_T) \) for all \( 1 \leq s < \frac{p(N+2)}{N+2-2p} \). If \( p \geq \frac{N+2}{2} \), then \( u \in L^\infty(\Omega_T) \). Moreover, the corresponding norms grow at most polynomially w.r.t \( T \).

**Proof of Lemma 3.2.** When \( 1 \leq N \leq 2 \) or \( N \geq 6 \) and \( \delta \) is small enough, the results follow from [CDF14]. Here we prove that when \( 3 \leq N \leq 5 \), the system (1.22) possesses global classical solution without assumption on the diffusion coefficients. First we observe that

\[ (2a + b + c)t - \Delta(2d_a a + d_bb + d_cc) = 0, \]

or equivalently

\[ z_t - \Delta(Mz) = 0, \]

with \( z = 2a + b + c \), and

\[ 0 < \min\{d_a, d_b, d_c\} \leq M(x,t) := \frac{2d_a a + d_bb + d_cc}{2a + b + c} \leq \max\{d_a, d_b, d_c\} < +\infty. \]

It follows from e.g. [Pie10] that \( z \in L^2(\Omega_T) \) for all \( T > 0 \) (with norms at most polynomially growing w.r.t \( T \)), which in combination with the nonnegativity of the concentrations leads to \( a, b, c \in L^2(\Omega_T) \). Secondly, we obtain from (1.22) that

\[
\begin{align*}
    a_t - d_a \Delta a &\leq k_3b^2, \quad (4.1a) \\
    b_t - d_b \Delta b &\leq k_4a + k_2bc, \quad (4.1b) \\
    c_t - d_c \Delta c &\leq k_1a. \quad (4.1c)
\end{align*}
\]

The strategy of the proof is the following: By using Lemma 4.1 and the regularity of \( a \), we get from (4.1c) an improved regularity of \( c \). This yields an improved regularity of \( a + bc \) and thus from (4.1b) an improvement of the regularity of \( b \). This information is used in (4.1a) to deduce the improved regularity of \( a \). Graphically, we have the following bootstrap iteration:

[Diagram of the bootstrap iteration]

This iteration terminates whenever we achieve \( a, b, c \in L^\infty(\Omega_T) \). In the rest of the proof, we will show how this iteration works when \( N = 3, 4, 5 \) and also explain why it fails when \( N \geq 6 \). To avoid unnecessary long explanations, we will present the iterations for each case in a table. We shall use the notation \( f \in L^{\alpha-0} \) if \( f \in L^\alpha(\Omega_T) \) for all \( p < \alpha \) and \( f \in L^\infty -0 \) if \( f \in L^\infty(\Omega_T) \) for all \( p < +\infty \) (with norms at most polynomially growing w.r.t \( T \)).

**Case \( N = 3 \):**

|   |   |   |   |   |
|---|---|---|---|---|
|   | \( L^2 \) | \( L^2 \) | \( - \) | \( L^2 \) |
| Step 0 | \( L^2 \) | \( L^{1-0} \) | \( L^{1/3-0} \) | \( L^{5-0} \) |
| Step 1 | \( L^\infty -0 \) | \( L^\infty \) | \( L^{5-0} \) | \( L^\infty \) |
| Step 2 | \( L^\infty -0 \) | \( - \) | \( - \) | \( - \) |
| Step 3 | \( L^\infty \) | \( - \) | \( - \) | \( - \) |

**Case \( N = 4 \):**


In order to do so, recall from (2.25) that $\Phi(a)$ is an approximation of $\Phi$ at point $a$. In the following, we will prove that if there exists such an $a \in B_{\delta}(c_{\infty})$ for $\delta$ small enough, then the linear approximation of $\Phi$ at $a$, which we denote by $L_{a}\Phi$, is such an affine function.

It is obvious that $\Phi(a) = L_{a}\Phi(a)$. It remains to show that

$$\Phi(c) \geq L_{a}\Phi(c) \quad \text{for all} \quad c \in \mathbb{R}^{N}_{\geq 0}. \quad (4.2)$$

In order to do so, recall from (2.25) that $\Phi(c) = F(c) + G(c)$. We will treat $F$ and $G$ separately.

First, by using the notations $\rho = |c - a|$, $\delta = |a - c_{\infty}|$, and direct computation of the Taylor expansion, we have thanks to the convexity of $F$,

$$F(c) - L_{a}F(c) \geq \frac{\nu \rho^{2}}{1 + \rho},$$

for some constant $\nu > 0$.

Next, thanks to the smoothness of $G$, and $G(c_{\infty}) = DG(c_{\infty}) = 0$ and $D^{2}G(c_{\infty}) \geq 0$, we get for $a \in B_{\varepsilon_{*}}(c_{\infty})$ (with $\varepsilon_{*}$ small enough)

$$|G(a)| \leq M\delta^{2}, \quad |DG(a)| \leq M\delta, \quad D^{2}G(a) \geq -M\delta, \quad |D^{3}G(a)| \leq M. \quad (4.3)$$

By using Taylor expansion again, we see that locally:

$$G(c) - L_{a}G(c) \geq -M\delta \rho^{2} - M\rho^{3}. \quad (4.3)$$

When $c$ is far away from $a$ and $\rho$ is not small, then by using the fact that $a$ is close to $c_{\infty}$ and $G(c_{\infty})$ is zero only at point $c_{\infty}$, we obtain that $G(c) - L_{a}G(c)$ is bounded below. Thus, (4.3) holds also in this case for some suitable $M > 0$. On the other hand, we can also estimate

$$G(c) - L_{a}G(c) \geq -M\delta^{2} - M\rho. \quad (4.4)$$

Hence, we obtain from (4.3) and (4.4) the lower bound

$$G(c) - L_{a}G(c) \geq -M(\delta + \rho) \min\{\delta, \rho^{2}\} \geq -\frac{4M\sqrt{\delta}\rho^{2}}{1 + \rho}. \quad (4.5)$$

Now, by choosing $\delta = \min\{\varepsilon_{*}, \nu^{2}/(4M)^{2}\}$, we have for all $a \in B_{\delta}(c_{\infty})$,

$$\Phi(c) - L_{a}\Phi(c) \geq \frac{(\nu - 4M\sqrt{\delta})\rho^{2}}{1 + \rho} \geq 0. \quad (4.6)$$

This proves (4.2) and thus completes the proof of the Lemma.
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References
[And11] D.F. Ardenson, A proof of the Global Attractor Conjecture in the single linkage class case, SIAM J. Appl. Math., 71 (2011) pp. 1487–1508.
[And14] D.F. Ardenson, A short note on the Lyapunov function for complex-balanced chemical reaction networks, online notes https://www.math.wisc.edu/~anderson/CRNT_Lyapunov.pdf.
[AMTU01] A. Arnold, P. Markowich, G. Toscani, A. Unterreiter, On convex Sobolev inequalities and the rate of convergence to equilibrium for Fokker-Planck type equations, Comm. Partial Differential Equations 26 (2001) 43–100.
[Bol1896] L. Boltzmann, Gastheorie, (1896) Leipzig, J. A. Barth.
[Bol1887] L. Boltzmann, Neuer Beweis zweier Sätze über das Wärmegleichgewicht unter mehratomigen Gasmolekülen, Sitzungsberichte der Kaiserlichen Akademie der Wissenschaften in Wien. 95 (1887), 153–164.
[CDF14] J.A. Cañizo, L. Desvillettes and K. Fellner, Improved duality estimates and applications to reaction-diffusion equations, Comm. Partial Differential Equations, 39 (2014) pp. 1185–1204.
[CNN13] G. Craciun, F. Nazarov, C. Pantea, Persistence and permanence of mass-action and power-law dynamical systems, SIAM J. Appl. Math., 73 (2013) pp. 305–329.
[Cra] G. Craciun, Toric Differential Inclusions and a Proof of the Global Attractor Conjecture, arXiv:1501.02860.
[CAAB09] G. Craciun, A. Dickenstein, A. Shiu, B. Sturmfels, Toric dynamical systems, J. Symbolic Comput. 44, no. 11 (2009) 1551–1565.
[DF06] L. Desvillettes, K. Fellner, Exponential decay toward equilibrium via entropy methods for reaction–diffusion equations, J. Math. Anal. Appl., 319 (2006), pp. 157–176.
[DF08] L. Desvillettes, K. Fellner, Entropy methods for reaction–diffusion equations: slowly growing a priori bounds, Rev. Mat. Iberoamericana, 24 (2008), pp. 407–431.
[DF14] L. Desvillettes, K. Fellner, Exponential Convergence to Equilibrium for a Nonlinear Reaction-Diffusion System with Detailed Balance Condition, arXiv:1601.05992.
[DFP07] L. Desvillettes, K. Fellner, M. Pierre, J. Vovelle, About Global existence of quadratic systems of reaction-diffusion equations. J. Adv. Nonlinear Stud. 7 (2007), 491–511.
[Fei79] M. Feinberg, Lectures on Chemical Reaction Networks, University of Wisconsin-Madison, 1979, https://crnt.osu.edu/LecturesOnReactionNetworks.
[Fei87] 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), pp. 2229–2268.
[FFH74] M. Feinberg and F. J. M. Horn, Dynamics of open chemical systems and the algebraic structure of the underlying reaction network, Chem. Eng. Sci., 29 (1974), pp. 775–777.
[FL16] K. Fellner, E–H. Laamri, Exponential decay towards equilibrium and global classical solutions for nonlinear reaction–diffusion systems, J. Evol. Equ., 16 (2016), pp. 681–704.
[FLS16] K. Fellner, E. Latos and T. Suzuki, Global classical solutions for mass-conserving, (super-)quadratic reaction–diffusion systems in three and higher space dimensions, to appear in Discrete Contin. Dyn. Syst. Ser. B.
[FLT] K. Fellner, E. Latos and B. Tang, Well-posedness and exponential equilibration of a volume-surface reaction-diffusion system with nonlinear boundary coupling, to appear in Ann. Inst. H. Poincaré Anal. Non Linéaire.
[FPT] K. Fellner, W. Prager and B.Q. Tang, The entropy method for reaction-diffusion systems without detailed balance: first order chemical reaction networks, Preprint.
[FT] K. Fellner and B.Q. Tang, Explicit exponential convergence to equilibrium for mass action reaction-diffusion systems with detailed balance condition, arXiv:1601.05992.
[Fis15] J. Fischer, Global Existence of Renormalized Solutions to Entropy-Dissipating Reaction-Diffusion Systems, Arch. Rational Mech. Anal. 218 (2015) pp. 553–587.
[FJ16] J. Fontbona, B. Jourdain, A trajecotorial interpretation of the dissipations of entropy and Fisher information for stochastic differential equations, Ann. Probab., 44 (2016) pp. 131–170.
[GHH96] A. Glitzky, K. Gröger, R. Hünlich, Free energy and dissipation rate for reaction–diffusion processes of electrically charged species, Appl. Anal. 60 (1996), 201–217.
[GH97] A. Glitzky, R. Hünlich, Energetic estimates and asymptotics for electro-reaction-diffusion systems, Z. Angew. Math. Mech. 77 (1997), 823–832.
[Grö92] K. Gröger, Free energy estimates and asymptotic behaviour of reaction-diffusion processes, Preprint 20, Institut für Angewandte Analysis und Stochastik, Berlin, 1992.
[Gun03] J. Gunawardena, Chemical Reaction Network Theory in In-Silico Biologists, 2003.
[Gop13] M. Gopalkrishnan, On the Lyapunov function for complex-balanced mass-action systems, arXiv:1312.3043.
[GMS14] M. Gopalkrishnan, E. Miller, A. Shiu, A geometric approach to the global attractor conjecture, SIAM J. Appl. Dyn. Syst., 13 (2014), pp. 758–797.
[GZ10] I. Gentil, B. Zegarlinski, Asymptotic behaviour of a general reversible chemical reaction-diffusion equation, Kinet. Relat. Models. 3 (2010) pp. 427–444.
[Hor72] F.J.M. Horn, Necessary and sufficient conditions for complex balancing in chemical kinetics, Arch. Rational Mech. Anal., 49 (1972), pp. 172–186.
[Ho74] F.J.M. Horn, The dynamics of open reaction systems, in SIAM-AMS Proceedings, Vol. VIII, SIAM, Philadelphia, 1974, pp. 125–137.

[HHJ72] F.J.M. Horn and R. Jackson, General mass action kinetics, Arch. Rational Mech. Anal., 47 (1972), pp. 81–116.

[Kir90] M. Kirane, On stabilization of solutions of the system of parabolic differential equations describing the kinetics of an autocatalytic reversible chemical reaction, Bull. Inst. Mat. Acad. Sin. 18, no. 4 (1990), pp. 369-377.

[MHM15] A. Mielke, J. Haskovec, P. A. Markowich, On uniform decay of the entropy for reaction-diffusion systems, J. Dynam. Differential Equations 27 (2015) 897–928.

[Pa92] C. V. Pao, Nonlinear Parabolic and Elliptic Equations, Springer (1992).

[Pan12] C. Pantea, On the persistence and global stability of mass-action systems, SIAM J. Math. Anal., 44 (2012) pp. 1636–1673.

[Pie10] M. Pierre, Global existence in reaction-diffusion systems with control of mass: a survey, Milan J. Math. 78.2 (2010), pp. 417–455.

[PMP06] Philippe Michel, Stéphane Mischler, and Benoît Perthame, General relative entropy inequality: an illustration on growth models, J. Math. Pures Appl. 84 (2005) pp. 1235–1260.

[Rot84] F. Rothe, Global Solutions of Reaction-Diffusion Systems, Lecture Notes in Mathematics, Springer, Berlin, (1984).

[SJ80] D. Siegel, M.D. Johnston, Linearization of complex balanced reaction systems, Preprint 2008.

[SM00] D. Siegel, D. MacLean, Global stability of complex balanced mechanisms, J. Math. Chem. 27 (2000) pp. 89–110.

[TeVi90] G. Toscani, C. Villani, On the trend to equilibrium for some dissipative systems with slowly increasing a priori bounds, J. Statist. Phys. 98, no. 5-6 (2000), pp. 1279-1309.

[Vol72] A.I. Volpert, Differential equations on graphs, Mat. Sb. 88 no.130 (1972) 578–588 (in Russian) Math. USSR-Sb. 17 (1972) 571–582 (in English).

[Weg1901] R. Wegscheider, Über simultane Gleichgewichte und die Beziehungen zwischen Thermodynamik und Reactionskinetik homogener Systeme, Monatshfte für Chemie 32 (1901), 849–906.

[Wil] D. Willett, A linear generalization of Gronwall’s inequality Proc. Amer. Math. Soc. 16 (1965) pp. 774–778.

[WCHL96] C. Wu, L. Cheng, M. Ha, E.S. Lee, Convexification of nonconvex functions and application to minimum and maximum principles for nonconvex sets, Computers Math. Appl. 31 (1996) 27–36.

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