# **The SYMBIONT Project: Symbolic Methods for Biological Networks**

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**SYMBIONT** PROJECT

#### **Overview**

SYMBIONT ranges from mathematics via computer science to systems biology, with a balanced team of researchers from those fields. At the present stage the project has a clear focus on fundamental research on mathematical methods and prototypes in software. Results are systematically benchmarked against models from computational biology databases. We summarize the motivation and aims for the project, and report on some existing results.

# Main Objectives

- 1. Solving high dimensional parametric models for fixed points
- 2. Order reduction of ordinary differential equations
- 3. Improved techniques for parameter fitting with respect to observed data
- 4. Applications in systems biology and medicine

# **Methods and Results**

**Computer algebra methods for solving** systems of equations and inequalities

For polynomial or rational vector fields, classical topics in the qualitative theory of ordinary differential equations such as the existence and stability of periodic solutions and equilibria can be reduced to real quantifier elimination problems. These questions are important for chemical reaction networks (CRN). However, they turn out difficult for medium size and large networks when using generic methods. Based on ideas from Stoichiometric Network Analysis (SNA) specifically virtual substitution and regular chain techniques allow the symbolic computation of Hopf bifurcation fixed

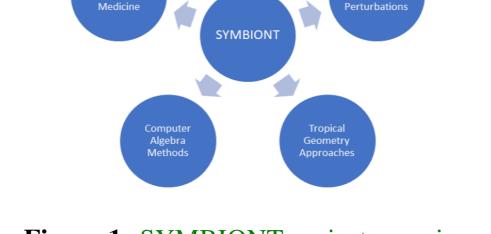


Figure 1: SYMBIONT project overview

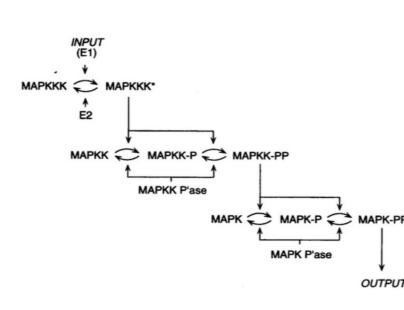
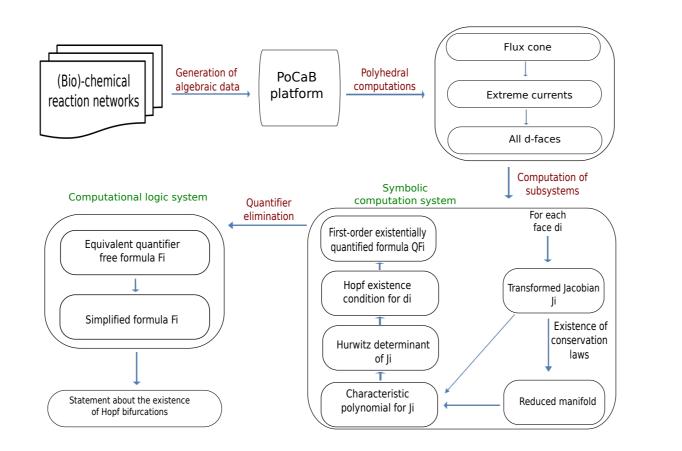


Figure 2: MAPK is an important signalling network. Reduced models serve as test cases.



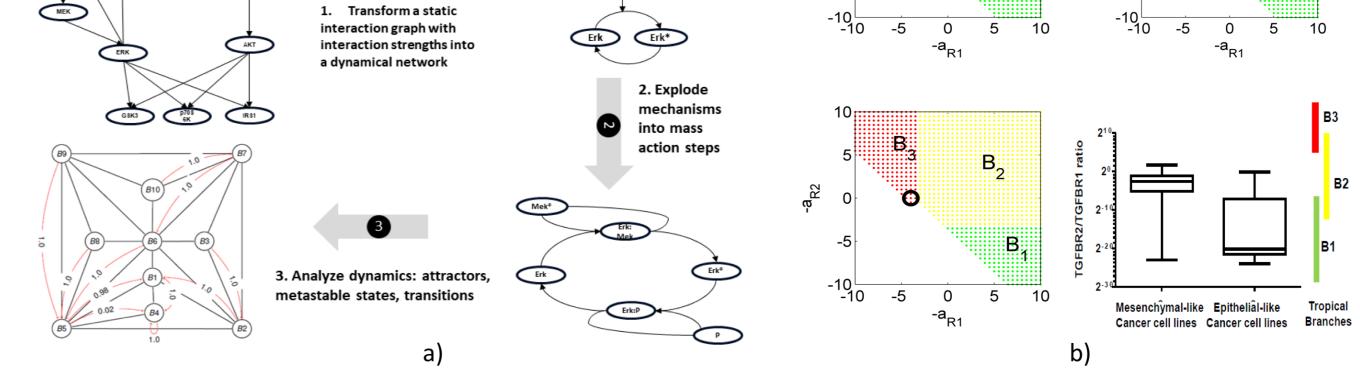


Figure 6: a) Flowchart indicating reconstruction and analysis of dynamical networks used as models for cell processes. Static networks are directed graphs; Dynamic networks are systems of polynomial ODEs; Branches B1, ..., B10 of tropical equilibrations indicate metastable regimes; Transitions between metastable regimes form a non-deterministic automaton. b) Tropical equilibration branches provide proxys for metastable regimes of large biochemical networks.

#### **Parameter fitting**

We aim to develop symbolic and numeric methods dedicated to the parameter estimation problem for models formulated by means of nonlinear integrodifferential equations (IDE). In particular, we address the computation of the model input-output equation and the numerical integration of IDE systems (cf. [1] and Fig. 7).

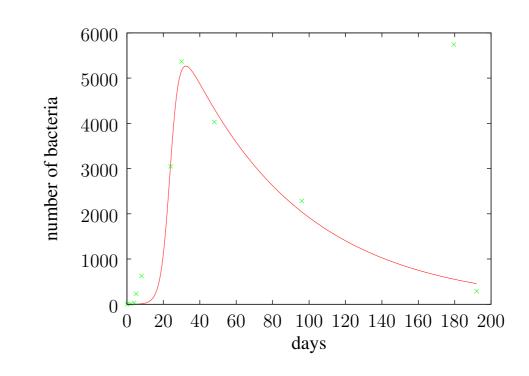


Figure 7: Fitting of parameters (coefficients) of the Volterra–Kostitzin integro-differential model, described as  $\dot{p}(t) = \varepsilon p(t) - hp^2(t) - bp^2(t)$  $cp(t) \int_0^t K(t-\tau)p(\tau) d\tau$  where p(t) denotes the number of bacteria.

Figure 3: Schematic view of HoCoQ—an algorithm for computing Hopf bifurcations using convex coordinates and quantifier elimination.

points. Some crucial structural features of CRN are used to derive corresponding heuristics for quantifier elimination [2]. "Complex Balancing," which is frequently encountered in CRNs, leads to parameterized binomial systems. For solving such systems, we are using polynomial time solution methods [4].

#### Model order reduction and qualitative dynamics

Model order reduction often relies on quasi-steady state phenomena that lead to slow-fast variable decomposition. This is related to the classical singular perturbation scheme due to Tikhonov and Fenichel [3]. The aim is to determine critical manifolds, i.e., parameter values resulting in stationary points with positive dimension. Figure 4 illustrates typical results. In addition, we use scaling and more generally Lie symmetries for variable and parameter nondimensionalisation and identification of small parameters.

Tropical geometry offers an alternative framework. For polynomial ODEs, we use tropical methods to rank monomial 7 complex invariant(s) terms according to their orders of mag-No (more) complete equilibration nitude and to identify dominant terms. in [25]: tropicalize(tropical\_max\_degree: 3 E1\_KKK+E2\_P\_KKK+KKK+P\_KKK+P\_KKK\_KK+P\_KKK\_P\_KK E1+E1 KKK When there is only one dominant term E2+E2 P KKH KK+KKPase\_PP\_KK+KKPase\_P\_KK+PP\_KK\_PP\_KK\_P\_KK\_P\_KK+P\_KKK\_KK+P\_KKK\_P\_KK KKPase+KKPase\_PP\_KK+KKPase\_P\_KK K+KPase\_PP\_K+KPase\_P\_K+PP\_KK\_K+PP\_KK\_P\_K+P\_K or when the dominant terms all have the KPase+KPase\_PP\_K+KPase\_P\_K 7 complex invariant(s) found a complete equilibration leading to the rescaling: E1' =  $\epsilon^{-}$  5 \* E1 E1\_KKK' =  $\epsilon^{-}$  7 \* E1\_KKK E2' =  $\epsilon^{-}$  4 \* E2 same sign, the dynamics is fast and the  $E2\_P\_KKK' = \varepsilon^{-} 7 * E2\_P\_KKK$  K' = Ksystem tends rapidly towards a region in KK' = KK  $KKK' = \epsilon^{-} 3 * KKK$ KKPase' =  $\epsilon^{-}$  5 \* KKPase phase space, where at least two domi-KKPase\_PP\_KK' =  $\epsilon^{-}$  6 \* KKPase\_PP\_KK nant terms of opposite signs are equili-Figure 5: Computation of tropical equilibrations using Conbrated. We call a solution to the latter straint Logic Programming in BIOCHAM-4. problem a tropical equilibration, which we can compute automatically (cf. [5] and Fig. 5). We show that tropical equilibrations can be algorithmically grouped into branches and put into correspondence with metastable dynamic regimes of chemical reaction networks (cf. Fig. 6a). Each tropical equilibration branch leads to a different reduced model.

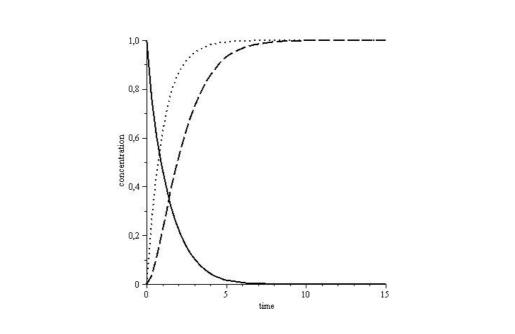


Figure 4: An instance where quasi-steady state reduction provides incorrect results (Michaelis-Menten system with small parameter  $k_{-1}$ ). The dashed line represents the concentration of product correctly, the dotted line represents the quasi-steady state approximation.

In [24]:	<pre>tropicalize(tropical_max_degree: 2).</pre>
Out[24]:	E1_KKK+E2_P_KKK+KKK+P_KKK+P_KKK_KK+P_KKK_P_KK
	E1+E1_KKK
	E2+E2_P_KKK
	KK+KKPase_PP_KK+KKPase_P_KK+PP_KK+PP_KK_K+PP_KK_P_KK+P_KKK+P_KKK+P_KKK_P_KK
	KKPase+KKPase_PP_KK+KKPase_P_KK
	K+KPase_PP_K+KPase_P_K+PP_K+PP_KK_K+PP_KK_P_K+P_K
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#### **Applications in systems biology and medicine**

**SYMBIONT** eventually aims at investigating challenging models for systems biology and medicine applying the various methods described here. It will compute attractors, dominant subsystems, and metastable regimes as well as tipping points and transitions between such regimes (cf. Fig. 6b, Fig. 8). Our analysis could provide hints about the onset of a disease and support the mechanistic understanding of the disease

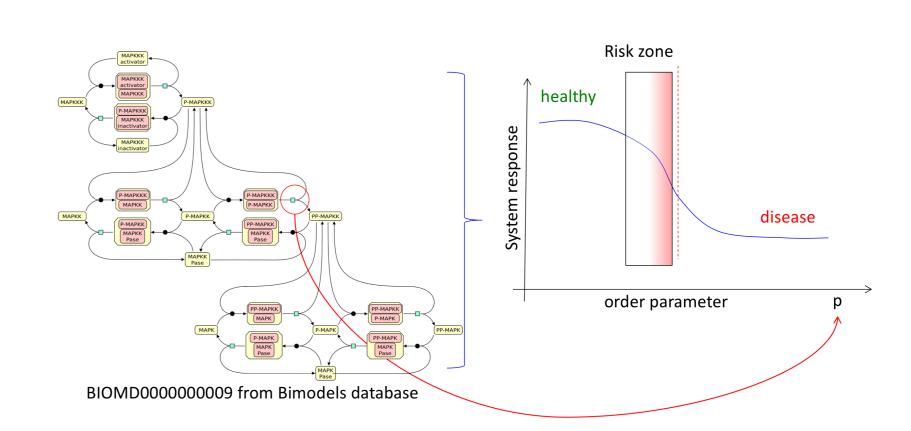


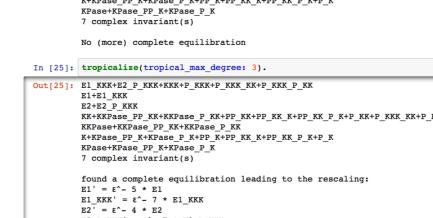
Figure 8: A representation of critical transitions in biological systems. The deviation of few "critical" parameters in the network (left) can shift the system's response resulting in the transition from "healthy" to "disease" state (right).

process. For instance, network dynamic regimes are related to tumor cell phenotypes in the case of TGF- $\beta$  signalling [6].

# Challenges

- A rigorous proof of the appearance of Hopf bifurcations, e.g., in the MAPK cascade
- Using hidden toric structures of steady states in combination with linear conservation constraints
- Model reduction for ODEs with sums of fractions
- Additional constraints (e.g. stability constraints) to the tropical equilibration branches
- Tipping points and dynamical regimes of large biological networks symbolically

### References



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We are seeking PhD students and postdoctoral researchers for our project. We also organize regular workshops open for external researchers fostering collaborations. Find information on our website.



