Time-dependent propagators for stochastic models of gene expression: an analytical method

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Abstract

The inherent stochasticity of gene expression in the context of regulatory networks profoundly influences the dynamics of the involved species. Mathematically speaking, the propagators which describe the evolution of such networks in time are typically defined as solutions of the corresponding Chemical Master Equation (CME). However, it is not possible in general to obtain exact solutions to the CME in closed form, which is due largely to its high dimensionality. In the present article, we propose an analytical method for the efficient approximation of these propagators. We illustrate our method on the basis of two categories of stochastic models for gene expression that have been discussed in the literature. The requisite procedure consists of three steps: a probability-generating function is introduced which transforms the CME into (a system of) partial differential equations (PDEs); application of the method of characteristics then yields (a system of) ordinary differential equations (ODEs) which can be solved using dynamical systems techniques, giving closed-form expressions for the generating function; finally, propagator probabilities can be reconstructed numerically from these expressions via the Cauchy integral formula. The resulting 'library' of propagators lends itself naturally to implementation in a Bayesian parameter inference scheme, and can be generalised systematically to related categories of stochastic models beyond the ones considered here.

1 Introduction

1.1 Motivation

Understanding the process of gene expression in the context of gene regulatory networks is indispensable for gaining insight into the fundamentals of numerous biological processes. However, gene expression can be highly stochastic in nature, both in prokaryotic and in eukaryotic organisms; see e.g. the work by Elowitz et al. 2002; Raj and Oudenaarden 2008; Shahrezaei and Swain 2008b, and the references therein. This inherent stochasticity has a profound influence on the dynamics of the involved species, in particular when their abundance is low. Therefore, gene expression is often appropriately described by stochastic models (Bressloff 2014; Karlebach and Shamir 2008; Thattai and Oudenaarden 2001; Wilkinson 2009). A schematic of the canonical model for gene expression is depicted in Figure 1. Here, the processes of transcription, translation, and degradation are approximated by single rates.

To test the validity of such stochastic models, a comparison with experimental data needs to be performed. The development of experimental techniques, such as time-lapse fluorescence microscopy (Coutu and Schroeder 2013; Elowitz et al. 2002; Larson et al. 2011; Muzzey and Oudenaarden 2009; Raj et al. 2006; Young et al. 2011), allows for real-time tracking of gene expression dynamics in single cells, providing mRNA or protein abundance time series, such

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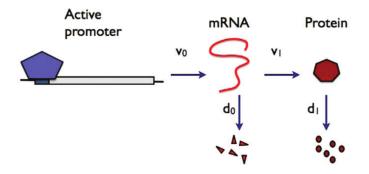


Figure 1: The canonical model of gene expression. Transcription of mRNA occurs with rate v_0 ; mRNA is translated to protein with rate v_1 . Both mRNA and protein decay, with rates d_0 and d_1 , respectively. (Figure courtesy of Shahrezaei and Swain 2008a.¹)

as those depicted in Figure 2. To select between competing hypotheses on the underlying regulatory networks given measurement data, as well as to infer the values of the corresponding model parameters, we can apply Bayesian inference theory to calculate the likelihood of a given model, which is constructed as follows.

The abundance of a protein, denoted by n, is sampled at times t_i , see Figure 2, yielding a list of measurement pairs (t_i, n_i) from which *transitions* $(\Delta t, n_i \rightarrow n_{i+1})$ between states can be extracted; here, $\Delta t = t_{i+1} - t_i$ is the regular, fixed, sampling interval. Next, the underlying stochastic model with parameter set Θ is used to calculate the probabilities of these transitions, which are denoted by $P_{n_{i+1}|n_i}(\Delta t)$. These so-called *propagators* give the probability of n_{i+1} protein being present after time Δt , given an initial protein abundance of n_i . The log-likelihood $L(\Theta|D)$ of the parameter set Θ , given the observed data D, is now defined in terms of the propagators as

$$L(\Theta|D) = \sum_{i} \log P_{n_{i+1}|n_i}(\Delta t).$$
(1.1)

To infer the values of parameters in the model, propagators are calculated for a wide range of parameter combinations, resulting in a 'log-likelihood landscape'; the maximal value of the log-likelihood as a function of the model parameters yields the most likely parameter values, given the experimental data. An example realisation of the above procedure can e.g. be found in the work by Feigelman, Popović and Marr 2015.

To calculate accurately the log-likelihood in (1.1), it is imperative that the values of the propagators can be extracted from the underlying stochastic model for any desired combination of parameters in Θ . In particular, we need to be able to calculate the propagator $P_{n|n_0}(t)$ as a function of time *t* for any initial protein number n_0 . Unfortunately, the highly complex nature of the stochastic models involved makes it very difficult to obtain explicit expressions for these probabilities. Some analytical progress can be made when a steady-state approximation is performed, i.e. when it is assumed that the system is allowed to evolve for a sufficiently long time, such that it converges to a time-independent state. However, the sampling interval Δt used for obtaining experimental data, as seen in Figure 2, is often *short* with respect to the protein life time. As that life time represents a natural time scale for the system dynamics, it follows that the evolution of the probabilities $P_{n|n_0}(t)$ should be studied over short times, in contradiction with the steady-state or long-evolution-time approximations which have previously been employed to derive analytical results (Bokes et al. 2012a; Hornos et al. 2005; Iyer-Biswas and Jayaprakash 2014; Shahrezaei and Swain 2008a).

The complex nature of stochastic models for gene expression has led to the widespread use of stochastic simulation techniques, such as Gillespie's algorithm (Gillespie 1977), with the aim

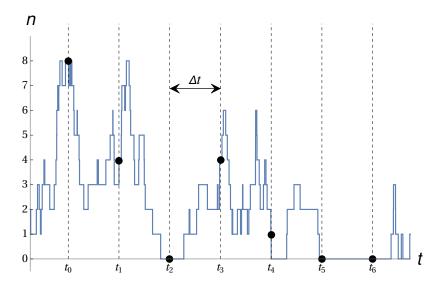


Figure 2: Sketch of a potential time series of protein abundance *n*, sampled at times t_i , with regular, fixed, sampling interval Δt .

of predicting values for the associated propagators from these models; see Feigelman et al. 2016 for recent work combining stochastic simulation with a particle filtering approach. However, these approaches can still be very time-consuming, due to the (relatively) high dimensionality of the model parameter space, combined with the fact that, for each combination of parameter values, the stochastic model has to be simulated sufficiently many times to yield a probability distribution that can be used to infer the corresponding propagator. For that reason, it is desirable to be able to obtain explicit expressions for the propagator $P_{n|n_0}(t)$ directly in terms of the model parameters, if necessary in an appropriate approximation.

1.2 Analytical method

In the present article, we develop an analytical method for the efficient evaluation of timedependent propagators in stochastic gene expression models, for arbitrary values of the model parameters. The results of our analysis can be implemented in a straightforward fashion in a Bayesian parameter inference framework, as outlined above.

To demonstrate our approach, we analyse two different stochastic models for gene expression. The first model, henceforth referred to as 'model A', is a model that incorporates autoregulation, where transcription and translation are approximated by single rates and protein can either stimulate or inhibit its own production by influencing the activity of DNA; see Figure 3. That model was first studied by Iyer-Biswas and Jayaprakash 2014 via a steady-state approximation. The second model, henceforth referred to as 'model B', models both mRNA and protein explicitly and again incorporates DNA switching between an active and an inactive state; see Figure 4. That model was first studied by Shahrezaei and Swain 2008a in a long-evolution-time approximation.

Both model A and model B are formulated in terms of the Chemical Master Equation (CME), which is the accepted mathematical representation of stochastic gene expression in the context of the model categories considered here; cf. Iyer-Biswas and Jayaprakash 2014 and Shahrezaei and Swain 2008a, respectively. Mathematically speaking, the CME is an infinitedimensional system of linear ordinary differential equations (ODEs) that describes the evolution in time of the probabilities of observing a specific state in the system, given some initial state. Numerous approaches have been suggested for the (approximate) solution of the CME;

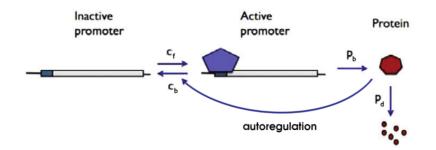


Figure 3: Schematic of model A, a gene expression model with autoregulation. (Base figure courtesy of Shahrezaei and Swain $2008a.^1$)

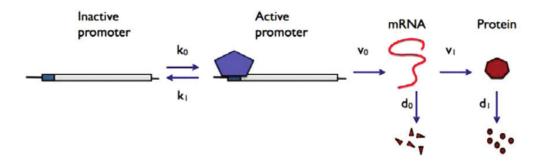


Figure 4: Schematic of model B, a gene expression model that explicitly incorporates transcription (Shahrezaei and Swain 2008a). (Figure courtesy of Shahrezaei and Swain 2008a.¹)

see e.g. Popović, Marr and Swain 2016 and the references therein for details. Our method relies on a combination of various techniques from the theory of differential equations and dynamical systems; specifically, we perform three consecutive steps, as follows.

- 1. **CME system** → **PDE system:** We introduce a probability-generating function to convert the CME into a (system of) partial differential equations (PDEs).
- 2. **PDE system** \rightarrow **ODE system:** Applying the method of characteristics combined, if necessary, with perturbation techniques we transform the system of PDEs obtained in step 1 into a dynamical system, that is, a system of ODEs.
- 3. **ODE system** \rightarrow **Explicit solution:** Making use of either special functions (model A) or multiple-time-scale analysis (model B), we obtain explicit solutions to the dynamical system found in step 2.

We emphasise that the 'characteristic system' of ODEs which is obtained in step 2 is lowdimensional, in contrast to the underlying CME system, as well as that it exhibits additional structure, allowing for the derivation of a closed-form analytical approximation for the associated generating function.

To convert the results of the above procedure into solutions to the original stochastic model, the three steps involved in our analysis have to be reverted. To that end, we require the following three ingredients:

- 1. Initial conditions are originally stated in terms of the CME, and first have to be reformulated in terms of the corresponding system of PDEs to ensure well-posedness; then, initial conditions can be extracted for the dynamical system that was obtained via the method of characteristics, reverting step 3.
- 2. To transform solutions to the characteristic system into solutions of the underlying PDE system, the associated 'characteristic transformation' has to be inverted, reverting step 2.
- 3. Lastly, solutions of the CME have to be extracted from solutions to the resulting PDE system, reverting step 1. Although the correspondence between the two sets of solutions is exact, theoretically speaking, the complexity of the expressions involved precludes the efficient analytical reconstruction of propagators from their generating functions. Therefore, we propose a novel hybrid analytical-numerical approach which relies on the Cauchy integral formula.

The various steps in our analytical method, as indicated above, are represented in Figure 5. It is important to mention that the implementation of Bayesian parameter inference, as outlined in Section 1.1, is not a topic for the present article; rather, the aim here is to describe our method, and to present analytical results which can readily be implemented in the context of parameter inference. The article hence realises the first stage of our research programme; the natural next stage, which is precisely that implementation, will be the subject of a follow-up article by the same authors.

1.3 Outline

The present article is organised as follows. In Section 2, we apply the analytical method outlined in Section 1.2 to model A, the gene expression model with autoregulation. Here, we use a perturbative approach to incorporate the autoregulatory aspects of the model; the resulting dynamical system can be solved in terms of confluent hypergeometric functions, see §13 in *NIST Digital Library of Mathematical Functions*. In Section 3, the same method is applied to model B, the model that explicitly incorporates transcription. We also indicate how autoregulation can be added to that model, and how the resulting extended model can be analysed on the basis

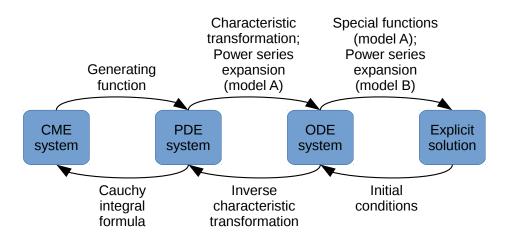


Figure 5: Schematic overview of the analytical method.

of our treatment of model A. The analysis carried out in Sections 2 and 3 yields a 'library' of explicit asymptotic expressions for the probability-generating functions associated to the underlying stochastic models. To obtain quantifiable expressions for their propagators, we introduce a novel hybrid analytical-numerical approach in Section 4, which can be readily implemented in the Bayesian parameter inference framework that provided the motivation for our analysis; see Section 1.1. We conclude with a discussion of our results, and an outlook to future work, in Section 5.

2 Model A: Gene Expression with Autoregulation

We first demonstrate our analytical method in the context of an autoregulatory stochastic gene expression model, as presented by Iyer-Biswas and Jayaprakash 2014; see also Figure 3. In the original article (Iyer-Biswas and Jayaprakash 2014), a Poisson representation was used to obtain analytical descriptions for time-independent solutions to the model. For a visual guide to the upcoming analysis, the reader is referred to Figure 5.

2.1 Stochastic model and CME

The basic stochastic model for gene expression is represented by the reaction scheme

$$D \stackrel{c_f}{\underset{c_b}{\leftarrow}} D^* \qquad (DNA \text{ switching}),$$

$$D^* \stackrel{p_b}{\rightarrow} D^* + P \qquad (\text{production of protein}),$$

$$P \stackrel{p_d}{\rightarrow} \emptyset \qquad (\text{decay of protein}).$$

$$(2.1)$$

The gene can hence switch between the inactive state D and the active state D^* , with switching rates c_f and c_b , respectively. The active gene produces protein (P) with rate p_b , while protein decays with rate p_d .

The autoregulatory part of the model is implemented as either positive or negative feedback:

$$D + P \xrightarrow{a} D^* + P$$
 (autoactivation), (2.2a)

$$D^* + P \xrightarrow{\prime} D + P$$
 (autorepression). (2.2b)

In the case of autoactivation, viz. (2.2a), protein induces activation of the gene with activation rate *a*, thereby accelerating its own production; in the case of autorepression, viz. (2.2b), protein deactivates the active gene with repression rate *r*, impeding its own production.

The CME system that is associated to the reaction scheme in (2.1), with autoactivation as in (2.2a), is given by

$$\frac{\mathrm{d}P_n^{(0)}}{\mathrm{d}t} = -\left(\kappa_f + \frac{a}{p_d}n\right)P_n^{(0)} + \kappa_b P_n^{(1)} + \left[(n+1)P_{n+1}^{(0)} - nP_n^{(0)}\right],\tag{2.3a}$$

$$\frac{\mathrm{d}P_n^{(1)}}{\mathrm{d}t} = \left(\kappa_f + \frac{a}{p_d}n\right)P_n^{(0)} - \kappa_b P_n^{(1)} + \left[(n+1)P_{n+1}^{(1)} - nP_n^{(1)}\right] + \lambda \left[P_{n-1}^{(1)} - P_n^{(1)}\right].$$
 (2.3b)

Here, $P_n^{(j)}(t)$ (j = 0, 1) represents the probability of *n* protein being present at time *t* while the gene is either inactive (0) or active (1). The time variable is nondimensionalised by the protein decay rate p_d ; other model parameters are scaled as

$$\kappa_f = \frac{c_f}{p_d}, \quad \kappa_b = \frac{c_b}{p_d}, \quad \text{and} \quad \lambda = \frac{p_b}{p_d}.$$
(2.4)

Analogously, the CME system for the case of autorepression, as defined in (2.2b), is given by

$$\frac{\mathrm{d}P_n^{(0)}}{\mathrm{d}t} = -\kappa_f P_n^{(0)} + \left(\kappa_b + \frac{r}{p_d}n\right) P_n^{(1)} + \left[(n+1)P_{n+1}^{(0)} - nP_n^{(0)}\right],\tag{2.5a}$$

$$\frac{\mathrm{d}P_n^{(1)}}{\mathrm{d}t} = \kappa_f P_n^{(0)} - \left(\kappa_b + \frac{r}{p_d}n\right) P_n^{(1)} + \left[(n+1)P_{n+1}^{(1)} - nP_n^{(1)}\right] + \lambda \left(P_{n-1}^{(1)} - P_n^{(1)}\right).$$
(2.5b)

Remark 2.1. A priori, it is possible to incorporate both autoactivation and autorepression in a single model, by merging systems (2.3) and (2.5). However, since autoactivation and autorepression precisely counteract each other, a partial cancellation would ensue, resulting in *effective* activation or repression. It can hence be argued that the simultaneous inclusion of both effects would introduce superfluous terms and parameters, which could be considered as poor model-ling practice. Therefore, we choose to model the two autoregulation mechanisms separately.

2.2 Generating function PDE

Rather than investigating the dynamics of (2.3) and (2.5) numerically, using stochastic simulation, we aim to employ an analytical approach. To that end, we define the probability-generating functions $F^{(j)}(z,t)$ (j = 0, 1) as follows; see e.g. Gardiner 2009:

$$F^{(j)}(z,t) = \sum_{n=0}^{\infty} z^n P_n^{(j)}(t).$$
(2.6)

In the case of autoactivation, the generating functions $F^{(j)}(z, t)$ can be seen to satisfy

$$\partial_t F^{(0)} + (z-1)\partial_z F^{(0)} + \frac{a}{p_d} z \partial_z F^{(0)} = -\kappa_f F^{(0)} + \kappa_b F^{(1)}, \qquad (2.7a)$$

$$\partial_t F^{(1)} + (z-1)\partial_z F^{(1)} - \frac{a}{p_d} z \partial_z F^{(0)} = \kappa_f F^{(0)} - \kappa_b F^{(1)} + \lambda (z-1) F^{(1)}$$
(2.7b)

if the coefficients $P_n^{(j)}(t)$ in (2.6) obey the CME system (2.3); likewise, in the case of autorepression, (2.3) gives rise to

$$\partial_t F^{(0)} + (z-1)\partial_z F^{(0)} - \frac{r}{p_d} z \partial_z F^{(1)} = -\kappa_f F^{(0)} + \kappa_b F^{(1)}, \qquad (2.8a)$$

$$\partial_t F^{(1)} + (z-1)\partial_z F^{(1)} + \frac{r}{p_d} z \partial_z F^{(1)} = \kappa_f F^{(0)} - \kappa_b F^{(1)} + \lambda(z-1)F^{(1)}.$$
 (2.8b)

Both (2.7) and (2.8) are systems of coupled, linear, first-order, hyperbolic partial differential equations (PDEs). Systems of this type are typically difficult to analyse; existing techniques only provide general results (Courant and Hilbert 1962; Taylor 2011).

To allow for an explicit analysis of systems (2.7) and (2.8), we make the following assumption:

Assumption 2.2. We assume that the autoactivation rate a in (2.2) is small in comparison with the other model parameters; specifically, we write

$$a = \alpha p_d \,\delta,\tag{2.9}$$

where $0 < \delta < 1$ is sufficiently small. Likewise, we assume that the autorepression rate r is small in comparison with the other model parameters, writing

$$r = \rho p_d \,\delta. \tag{2.10}$$

Previous work on the inclusion of autoregulatory effects in model selection by Feigelman et al. 2016 suggests that, in the context of Nanog expression in mouse embryonic stem cells, autoregulation rates are indeed small compared to other model parameters.

Based on Assumption 2.2, we can expand the generating functions $F^{(j)}$ (j = 0, 1) as power series in δ :

$$F^{(j)}(z,t) = \sum_{m=0}^{\infty} \delta^m F_m^{(j)}(z,t).$$
(2.11)

Substitution of (2.11) into (2.7) yields

$$\partial_t F_m^{(0)} + (z-1)\partial_z F_m^{(0)} = -\kappa_f F_m^{(0)} + \kappa_b F_m^{(1)} - \alpha \, z \partial_z F_{m-1}^{(0)}, \tag{2.12a}$$

$$\partial_t F_m^{(1)} + (z-1)\partial_z F_m^{(1)} = \kappa_f F_m^{(0)} - \kappa_b F_m^{(1)} + \lambda(z-1)F_m^{(1)} + \alpha z \partial_z F_{m-1}^{(0)}; \qquad (2.12b)$$

analogously, we substitute (2.11) into (2.8) to find

$$\partial_t F_m^{(0)} + (z-1)\partial_z F_m^{(0)} = -\kappa_f F_m^{(0)} + \kappa_b F_m^{(1)} + \rho \, z \partial_z F_{m-1}^{(1)}, \tag{2.13a}$$

$$\partial_t F_m^{(1)} + (z-1)\partial_z F_m^{(1)} = \kappa_f F_m^{(0)} - \kappa_b F_m^{(1)} + \lambda(z-1)F_m^{(1)} - \rho \, z \partial_z F_{m-1}^{(1)}.$$
(2.13b)

We observe that, in (2.12) and (2.13), the same leading-order differential operator acts on both $F_m^{(0)}$ and $F_m^{(1)}$, which allows us to apply the method of characteristics to solve the equations for $F_m^{(j)}$ (j = 0, 1) simultaneously. In particular, we emphasise that, mathematically speaking, the resulting perturbation is *regular* in the perturbation parameter δ .

2.3 Dynamical systems analysis

In this section, we apply the method of characteristics to derive the 'characteristic equations' that are associated to the PDE systems (2.12) and (2.13), respectively; the former are systems of ODEs, which are naturally analysed in the language of dynamical systems.

2.3.1 Autoactivation

We first consider the case of autoactivation; to that end, we rewrite system (2.12) as

$$(\partial_t + v\partial_v)F_m^{(0)} + \kappa_f F_m^{(0)} - \kappa_b F_m^{(1)} = -\alpha(v+1)\partial_v F_{m-1}^{(0)}, \qquad (2.14a)$$

$$(\partial_t + v\partial_v)F_m^{(1)} - \kappa_f F_m^{(0)} + \kappa_b F_m^{(1)} - \lambda v F_m^{(1)} = \alpha (v+1)\partial_v F_{m-1}^{(0)}, \qquad (2.14b)$$

where we have introduced the new variable

$$v = z - 1.$$
 (2.15)

The differential operator $\partial_t + v \partial_v$ in Equation (2.14) gives rise to characteristics $\xi(s; v_0)$ that obey the characteristic equation

$$\frac{\partial v}{\partial s} = v; \tag{2.16}$$

these characteristics can thus be expressed as

$$\xi(s; v_0) = (s, v_0 e^s). \tag{2.17}$$

Since the partial differential operators in (2.14) transform into

$$\partial_t + v\partial_v = \partial_s$$
 and (2.18a)

$$\alpha(\nu+1)\partial_{\nu} = \alpha \left(\nu_0 + e^{-s}\right)\partial_{\nu_0},\tag{2.18b}$$

we arrive at the following system:

$$\partial_s F_m^{(0)} + \kappa_f F_m^{(0)} - \kappa_b F_m^{(1)} = -\alpha \left(v_0 + e^{-s} \right) \partial_{v_0} F_{m-1}^{(0)}, \qquad (2.19a)$$

$$\partial_{s} F_{m}^{(1)} - \kappa_{f} F_{m}^{(0)} + \kappa_{b} F_{m}^{(1)} - \lambda v_{0} e^{s} F_{m}^{(1)} = \alpha \left(v_{0} + e^{-s} \right) \partial_{v_{0}} F_{m-1}^{(0)}.$$
(2.19b)

Note that Equation (2.19) is a recursive (nonhomogeneous) system of ordinary differential equations for $F_m^{(j)}$ (j = 0, 1). Henceforth, we will therefore refer to (2.19) as such, while retaining the use of partial derivatives ∂_s due to the presence of ∂_{v_0} in the corresponding right-hand sides. To solve system (2.19), we rewrite it as a second-order ODE for $F_m^{(0)}$: we hence obtain

$$\left[\partial_s^2 + (\kappa_f + \kappa_b - \lambda v_0 \mathbf{e}^s)\partial_s - \kappa_f \lambda v_0 \mathbf{e}^s\right] F_m^{(0)} = \left[\lambda v_0 \mathbf{e}^s - \partial_s\right] \alpha (v_0 + \mathbf{e}^{-s})\partial_{v_0} F_{m-1}^{(0)}, \tag{2.20}$$

which can be solved recursively to determine $F_m^{(0)}$ for any $m \ge 0$. To simplify (2.20), we introduce the variable

$$w = \lambda v_0 e^s, \tag{2.21}$$

which transforms the partial derivatives ∂_s and ∂_{v_0} into

$$\partial_s \to w \partial_w$$
 and $\partial_{v_0} \to \frac{w}{v_0} \partial_w + \partial_{v_0};$ (2.22)

Equation (2.20) hence reads

$$\left[(w\partial_w)^2 + (\kappa_f + \kappa_b - w)(w\partial_w) - \kappa_f w \right] F_m^{(0)} = (w - w\partial_w) \alpha (w + \lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_{m-1}^{(0)}.$$
(2.23)

Using (2.19a), we can express the second component $F_m^{(1)}$ in terms of $F_m^{(0)}$ as

$$F_{m}^{(1)} = \frac{1}{\kappa_{b}} \left[\partial_{s} F_{m}^{(0)} + \kappa_{f} F_{m}^{(0)} + \alpha \left(v_{0} + e^{-s} \right) \partial_{v_{0}} F_{m-1}^{(0)} \right]$$

$$= \frac{1}{\kappa_{b}} \left[w \partial_{w} F_{m}^{(0)} + \kappa_{f} F_{m}^{(0)} + \alpha \left(w + \lambda \right) \left(\partial_{w} + \frac{v_{0}}{w} \partial_{v_{0}} \right) F_{m-1}^{(0)} \right].$$
(2.24)

At leading order, i.e. for m = 0, (2.23) reduces to

$$[(w\partial_w)^2 + (\kappa_f + \kappa_b - w)(w\partial_w) - \kappa_f w]F_0^{(0)} = 0, \qquad (2.25)$$

the solutions of which can be expressed in terms of the confluent hypergeometric function ${}_{1}F_{1}$, see §13 of NIST Digital Library of Mathematical Functions, to yield

$$F_0^{(0)}(w) = c_{11}F_1(\kappa_f, 1 + \kappa_f + \kappa_b, w) + c_2 w^{-\kappa_f - \kappa_b} F_1(-\kappa_b, 1 - \kappa_f - \kappa_b, w).$$
(2.26)

Using (2.24), we can determine

$$F_0^{(1)}(w) = \frac{1}{\kappa_b} \left(w \partial_w F_0^{(0)} + \kappa_f F_0^{(0)} \right), \tag{2.27}$$

with $F_0^{(0)}$ as given in (2.26); for an explicit expression, see Equation (A.1) in Appendix A. The expression for $F_0^{(0)}(w)$ in (2.26) allows us to determine the first-order correction $F_1^{(0)}(w)$: substituting m = 1 in (2.23), we obtain

$$\left[(w\partial_w)^2 + (\kappa_f + \kappa_b - w)(w\partial_w) - \kappa_f w \right] F_1^{(0)} = (w - w\partial_w) \alpha (w + \lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_0^{(0)}.$$
(2.28)

Next, we apply the method of variation of constants to express the solution to (2.28) as

$$F_{1}^{(0)}(w) = \frac{1}{\kappa_{f} + \kappa_{b}} \bigg[{}_{1}F_{1}(\kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, w) \int_{c_{3}}^{w} F_{1}(-\kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, \hat{w}) g(\hat{w}) d\hat{w}$$
(2.29)
$$- w^{-\kappa_{f} - \kappa_{b}} {}_{1}F_{1}(-\kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, w) \int_{c_{4}}^{w} F_{1}(\kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, \hat{w}) \frac{g(\hat{w})}{\hat{w}^{-\kappa_{f} - \kappa_{b}}} d\hat{w} \bigg],$$

where

$$g(w) = \frac{e^{-w}}{w}(w - w\partial_w) \alpha (w + \lambda) \left(\partial_w + \frac{v_0}{w}\partial_{v_0}\right) F_0^{(0)},$$
(2.30)

with $F_0^{(0)}$ as given in (2.26). Finally, we may again use (2.24) to determine

$$F_1^{(1)}(w) = \frac{1}{\kappa_b} \left[w \partial_w F_1^{(0)} + \kappa_f F_1^{(0)} + \alpha (w + \lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_0^{(0)} \right],$$
(2.31)

with $F_0^{(0)}$ and $F_1^{(0)}$ as given in (2.26) and (2.29), respectively. At this point in the analysis, the constants c_1 and c_2 in (2.26) and the integration limits c_3 and c_4 in (2.29) remain undetermined. To fix these constants, and thereby determine a unique solution to (2.19), we have to prescribe appropriate initial conditions.

2.3.2 Autorepression

Given the analysis of autoactivation in the previous subsection, the case of autorepression can be analysed in an analogous manner. Employing the same characteristics as before, recall (2.17), we obtain

$$\partial_s F_m^{(0)} + \kappa_f F_m^{(0)} - \kappa_b F_m^{(1)} = \rho \left(v_0 + e^{-s} \right) \partial_{v_0} F_{m-1}^{(1)}, \qquad (2.32a)$$

$$\partial_s F_m^{(1)} - \kappa_f F_m^{(0)} + \kappa_b F_m^{(1)} - \lambda v_0 e^s F_m^{(1)} = -\rho \left(v_0 + e^{-s} \right) \partial_{v_0} F_{m-1}^{(1)}$$
(2.32b)

from system (2.13); cf. Equation (2.19). Next, we rewrite (2.32) as a second-order ODE for $F_m^{(1)}$, using again the variable transformation in (2.21):

$$\left[(w\partial_w)^2 + \left(\kappa_f + \kappa_b - w \right) (w\partial_w) - (1 + \kappa_f) w \right] F_m^{(1)} = - (w\partial_w) \rho \left(w + \lambda \right) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_{m-1}^{(1)},$$
(2.33)

(2.33) which can be solved recursively to obtain $F_m^{(1)}$ for any $m \ge 1$. The first component $F_m^{(0)}$ can be expressed in terms of $F_m^{(1)}$ as

$$F_m^{(0)} = \frac{1}{\kappa_f} \left[w \partial_w F_m^{(1)} + (\kappa_b - w) F_m^{(1)} + \rho \left(w + \lambda \right) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_{m-1}^{(1)} \right]; \tag{2.34}$$

to leading order, we thus obtain

$$F_0^{(1)}(w) = \hat{c}_{1\,1}F_1(1+\kappa_f, 1+\kappa_f+\kappa_b, w) + \hat{c}_2 w^{-\kappa_f-\kappa_b} {}_1F_1(1-\kappa_b, 1-\kappa_f-\kappa_b, w)$$
(2.35)

and

$$F_0^{(0)} = \frac{1}{\kappa_f} \left[w \partial_w F_0^{(1)} + (\kappa_b - w) F_0^{(1)} \right].$$
(2.36)

The corresponding equation for the first-order correction $F_1^{(1)}$ reads

$$\left[(w\partial_w)^2 + (\kappa_f + \kappa_b - w)(w\partial_w) - (1 + \kappa_f)w \right] F_1^{(1)} = -(w\partial_w)\rho(w + \lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_0^{(1)}, \quad (2.37)$$

which can be solved via the method of variation of constants to give

$$F_{1}^{(1)}(w) = \frac{1}{\kappa_{f} + \kappa_{b}} \bigg[{}_{1}F_{1}(1 + \kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, w) \int_{\hat{c}_{3}}^{w} F_{1}(1 - \kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, \hat{w}) h(\hat{w}) d\hat{w} - w^{-\kappa_{f} - \kappa_{b}} {}_{1}F_{1}(1 - \kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, w) \int_{\hat{c}_{4}}^{w} F_{1}(1 + \kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, \hat{w}) \frac{h(\hat{w})}{\hat{w}^{-\kappa_{f} - \kappa_{b}}} d\hat{w} \bigg];$$

$$(2.38)$$

here,

$$h(w) = -\frac{e^{-w}}{w} (w\partial_w) \rho (w + \lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0}\right) F_0^{(1)}, \qquad (2.39)$$

with $F_0^{(1)}$ as in (2.35). The first-order correction to the first component $F_0^{(0)}$ is hence given by

$$F_1^{(0)}(w) = \frac{1}{\kappa_f} \left[w \partial_w F_1^{(1)} + (\kappa_b - w) F_1^{(1)} + \rho \left(w + \lambda \right) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_0^{(1)} \right], \tag{2.40}$$

with $F_0^{(1)}$ and $F_1^{(1)}$ as in (2.35) and (2.38), respectively. As in the case of autoactivation, the constants \hat{c}_1 and \hat{c}_2 in (2.35) and the integration limits \hat{c}_3 and \hat{c}_4 in (2.38) remain undetermined, and have to be fixed through suitable initial conditions.

2.4 Initial conditions

To determine appropriate initial conditions for the dynamical systems (2.19) and (2.32), we consider the original CME systems (2.3) and (2.5), respectively.

At time t = 0, we impose an initial protein number $n = n_0$, which implies

$$P_n^{(0)}(0) + P_n^{(1)}(0) = \delta_{n,n_0}$$
(2.41)

for the probabilities $P_n^{(j)}(t)$ (j = 0, 1); here, δ_{n,n_0} denotes the standard Kronecker symbol, with $\delta_{n,n_0} = 1$ for $n = n_0$ and $\delta_{n,n_0} = 0$ otherwise. Using the definition of the generating functions $F^{(j)}(z, t)$ in (2.6), we find that

$$F^{(0)}(z,0) + F^{(1)}(z,0) = z^{n_0} = (v+1)^{n_0},$$
(2.42)

taking into account the change of variables in (2.15). Thus, (2.42) provides an initial condition for the PDE systems (2.7) and (2.8). Given the power series expansion in (2.11), we infer that the coefficients $F_m^{(j)}(v,t)$ (j = 0, 1) satisfy

$$F_0^{(0)}(v,0) + F_0^{(1)}(v,0) = (v+1)^{n_0}$$
 and (2.43a)

$$F_m^{(0)}(v,0) + F_m^{(1)}(v,0) = 0 \quad \text{for all } m \ge 1,$$
(2.43b)

which holds for both (2.12) and (2.13).

To be able to interpret the initial conditions in (2.43) in the context of the dynamical systems (2.19) and (2.32), we revisit the method of characteristics, which was used to map the PDE systems (2.12) and (2.13) to the former, respectively.

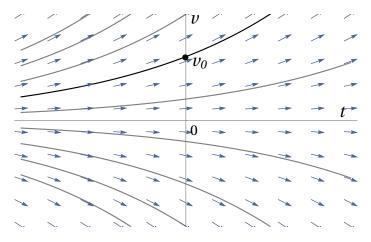


Figure 6: The (t, v)-coordinate plane, on which the PDE systems (2.12) and (2.13) are solved. The characteristics are integral curves of the vector field (1, v), indicated in blue. The black characteristic curve intersects the *v*-axis at $v = v_0$.

The characteristics of the differential operator $\partial_t + v \partial_v (= \partial_s)$ in (2.19) and (2.32) are the integral curves of the vector field (1, v). Geometrically speaking, these characteristic curves foliate the (t, v)-plane over the *v*-axis. Therefore, each characteristic can be identified by its base point, which is the point where the characteristic curve intersects the *v*-axis, at $v = v_0$; see Equation (2.17) and Figure 6.

Equivalently, each characteristic can be represented as a graph over the *t*-axis. Indeed, the differential equation for the *v*-component of a characteristic curve (2.16) can be solved to obtain the natural parametric description of a characteristic 'fibre' with base point $v = v_0$, which is given by (s, v_0e^s) . Here, the parameter *s* along the characteristic is chosen such that its point of intersection with the *v*-axis lies at s = 0. Given that choice, it is natural to identify the parameter along the characteristic (s) with the time variable (t). Hence, the initial conditions in (2.43), which determine a relation between $F_m^{(0)}$ and $F_m^{(1)}$ on the *v*-axis, can be interpreted on every characteristic as

$$\left[F_0^{(0)} + F_0^{(1)}\right]_{s=0} = (v_0 + 1)^{n_0} \quad \text{and} \tag{2.44a}$$

$$\left[F_m^{(0)} + F_m^{(1)}\right]_{s=0} = 0 \quad \text{for all } m \ge 1,$$
(2.44b)

which again holds for both (2.19) and (2.32).

Remark 2.3. For CME systems such as (2.3) and (2.5), it is customary to impose a 'normalisation' condition of the form

$$\sum_{m=0}^{\infty} P_m^{(0)}(t) + P_m^{(1)}(t) = 1,$$
(2.45)

as $P_m^{(0)}(t)$ and $P_m^{(1)}(t)$ represent probabilities. Recast in the framework of generating functions, recall (2.6), the above normalisation condition yields the boundary condition

$$\left[F^{(0)}(z,t) + F^{(1)}(z,t)\right]_{z=1} = \left[F^{(0)}(v,t) + F^{(1)}(v,t)\right]_{v=0} = 1.$$
 (2.46)

It is worthwhile to note that (2.46) is automatically satisfied whenever (2.43) is imposed: by adding the two equations in system (2.12) – or, equivalently, in (2.13) – one sees that $F_m^{(0)} + F_m^{(1)}$ satisfies

$$\partial_s F_m^{(0)} + \partial_s F_m^{(1)} = \lambda v F_m^{(1)}.$$
 (2.47)

The line {v = 0} is represented by the 'trivial' characteristic, which is identified by $v_0 = 0$; therefore, on that characteristic, we impose (2.43) to find

$$\partial_s F_m^{(0)} + \partial_s F_m^{(1)} = 0, (2.48a)$$

$$\left[F_0^{(0)} + F_0^{(1)}\right]_{s=0} = 1, \quad \text{and}$$
(2.48b)

$$\left[F_m^{(0)} + F_m^{(1)}\right]_{s=0} = 0 \quad \text{for all } m \ge 1,$$
(2.48c)

which implies that $F_0^{(0)} + F_0^{(1)} = 1$ for all *s*, as well as that $F_m^{(0)} + F_m^{(1)}$ vanishes identically for all $m \ge 1$. Substituting these results into the power series representation for $F^{(j)}(z, t)$ (j = 0, 1) in (2.11), we obtain (2.46).

At this point, it is important to observe that Equation (2.43) determines a *line* of initial conditions in the phase spaces of the dynamical systems (2.19) and (2.32). Therefore, at every order in δ , we can only fix *one* of the two free parameters that arise in the solution of the corresponding differential equations. In particular, (2.43) only fixes either c_1 or c_2 in (2.26), and either c_3 or c_4 in (2.29), and so forth. That indeterminacy motivates us to introduce a new parameter χ , which is defined as follows:

Definition 2.4. Consider the CME systems (2.3) and (2.5). We define $\chi(n_0)$ to be the probability that the gene modelled by the reaction scheme in (2.1) is switched off at time t = 0, given an initial protein number n_0 .

Definition 2.4 immediately specifies initial conditions for systems (2.3) and (2.5) via

$$P_n^{(0)}(0) = \chi(n_0) \,\delta_{n,n_0}$$
 and $P_n^{(1)}(0) = (1 - \chi(n_0)) \,\delta_{n,n_0};$ (2.49)

the above expression, in turn, provides us with a complete set of initial conditions for the PDE systems (2.7) and (2.8), to wit

$$F^{(0)}(v,0) = \chi(n_0) (v+1)^{n_0}$$
 and (2.50a)

$$F^{(1)}(v,0) = (1 - \chi(n_0))(v+1)^{n_0}.$$
(2.50b)

Here, we allow for the fact that $\chi(n_0)$ may depend on other model parameters, and in particular on the autoregulation rates *a* and *r*; for a discussion of alternative options, see Section 4.1. We therefore expand $\chi(n_0)$ as a power series in δ ,

$$\chi(n_0) = \sum_{m=0}^{\infty} \delta^m \chi_m(n_0).$$
 (2.51)

The above expansion can be used to infer a complete set of initial conditions for the PDE systems (2.12) and (2.13), yielding

$$F_m^{(0)}(v,0) = \chi_m (v+1)^{n_0} \quad \text{for all } m \ge 0,$$
(2.52a)

$$F_0^{(1)}(v,0) = (1-\chi_0)(v+1)^{n_0}$$
, and (2.52b)

$$F_m^{(1)}(v,0) = -\chi_m (v+1)^{n_0} \quad \text{for all } m \ge 1.$$
(2.52c)

By the same reasoning that inferred (2.44) from (2.43), we can conclude that the complete set of initial conditions for the dynamical systems (2.19) and (2.32) is given by

$$\left[F_m^{(0)}\right]_{s=0} = \chi_m (v_0 + 1)^{n_0} \quad \text{for all } m \ge 0,$$
(2.53a)

$$\left[F_{0}^{(1)}\right]_{s=0} = (1 - \chi_{0})(v_{0} + 1)^{n_{0}}, \text{ and}$$
(2.53b)

$$\left[F_m^{(1)}\right]_{s=0} = -\chi_m (v_0 + 1)^{n_0} \quad \text{for all } m \ge 1.$$
(2.53c)

We can now use the conditions in (2.53) to determine the free constants c_1 and c_2 in (2.26), which yields

$$c_{1} = (1 + v_{0})^{n_{0}} e^{-\lambda v_{0}} \frac{\kappa_{b}}{\kappa_{f} + \kappa_{b}} \left({}_{1}F_{1}(-\kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, \lambda v_{0}) + \frac{\chi_{0}\lambda v_{0}}{1 - \kappa_{f} - \kappa_{b}} {}_{1}F_{1}(1 - \kappa_{b}, 2 - \kappa_{f} - \kappa_{b}, \lambda v_{0}) \right) \text{ and } (2.54a)$$

$$c_{2} = (1 + v_{0})^{n_{0}} e^{-\lambda v_{0}} \frac{(\lambda v_{0})^{\kappa_{f} + \kappa_{b}}}{\kappa_{f} + \kappa_{b}} \Biggl\{ [\chi_{0}(\kappa_{f} + \kappa_{b}) - \kappa_{b}]_{1} F_{1}(\kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, \lambda v_{0}) + \frac{\kappa_{f} \chi_{0} \lambda v_{0}}{1 + \kappa_{f} + \kappa_{b}} {}_{1} F_{1}(1 + \kappa_{f}, 2 + \kappa_{f} + \kappa_{b}, \lambda v_{0}) \Biggr\}.$$
(2.54b)

Analogously, for \hat{c}_1 and \hat{c}_2 in (2.35), we obtain

$$\hat{c}_1 = c_1 \frac{\kappa_f}{\kappa_b}$$
 and $\hat{c}_2 = c_2$, (2.55)

with c_1 and c_2 as in (2.54). Using conversion formulas found in §13.3(i) of *NIST Digital Library* of *Mathematical Functions*, one can show that the expressions resulting from (2.35) and (2.36) match those in (2.27) and (2.26), respectively, as expected; see also Equation (A.1). Explicit expressions for the integration limits c_3 and c_4 in (2.29), as well as for the corresponding limits \hat{c}_3 and \hat{c}_4 in (2.38), can be found in Appendix A.

2.5 Inverse transformation

The final step towards providing explicit solutions to the PDE systems (2.7) and (2.8) consists in interpreting the solutions to the dynamical systems (2.19) and (2.32), with initial conditions as in (2.53), as solutions to the PDE systems (2.12) and (2.13), respectively. To that end, we again consider the corresponding characteristics from a geometric viewpoint.

As mentioned in Section 2.4, the (t, v)-coordinate plane is foliated by characteristics, which are the integral curves of the vector field (1, v); recall Figure 6. Hence, any point (t, v) lies on a unique characteristic; flowing backward along that characteristic to its intersection with the *v*-axis, we can determine the corresponding base point v_0 by the inverse transformation

$$(t, v) \mapsto v_0(t, v) = v e^{-t},$$
 (2.56)

since we have identified the parameter along the characteristic (s) with the time variable (t).

To determine the value of $F_m^{(j)}(v,t)$ (j = 0, 1), interpreted as a solution to the PDE system (2.12) or (2.13), we proceed as follows. We first apply the inverse transformation in (2.56) to establish on which characteristic the coordinate pair (t, v) lies. For that characteristic, identified by its base point v_0 , we then find the solution to the dynamical system (2.19) or (2.32), which is a function of s and v_0 . Next, we substitute s = t and $v_0 = ve^{-t}$ into that solution to obtain an explicit expression for the solution to the PDE system (2.12) or (2.13):

$$F_m^{(j)}(v,t)$$
 [as solution to (2.12) or (2.13)] = $\left[F_m^{(j)}\right]_{(s,v_0)=(t,ve^{-t})}$, (2.57)

where $F_m^{(j)}$ (j = 0, 1) on the right-hand side denotes the solution to (2.19) or (2.32), with initial conditions as in (2.53). Lastly, we substitute $F_m^{(j)}(v, t)$ into the power series in (2.11) to obtain an explicit solution to (2.7) or (2.8), to satisfactory order in δ .

Remark 2.5. The geometric interpretation of characteristics that was used to motivate the inverse transformation in (2.56) also shows that the introduction of the new system parameter χ in Definition 2.4 is *necessary* for the generating functions $F^{(j)}$ to be determined uniquely as

solutions to (2.7) or (2.8), even if we are only interested in their sum $F^{(0)}(v, t) + F^{(1)}(v, t)$. The crucial point is that any free constants obtained in the process of solving the dynamical systems (2.19) and (2.32) – or, equivalently, their second-order one-dimensional counterparts (2.20) and (2.33), respectively – are constant in *s*. In other words, they are constant along the particular characteristic on which the dynamical system is solved. These constants – such as e.g. c_1 and c_2 in (2.26) – can, and generally will, depend on the base point v_0 of the characteristic; see for example (2.54). The inverse transformation in (2.56) that is used to reconstruct the solution to the original PDE from that of the corresponding dynamical system would then yield undetermined functions $c(v e^{-t})$ in the resulting solutions to (2.7) and (2.8), respectively.

2.6 Summary of main result

To summarise Section 2, we combine the analysis of the previous subsections to state our main result.

Main result: The PDE system (2.7) can be solved for sufficiently small autoactivation rates *a*; see Assumption 2.2. Its solutions $F^{(j)}(z,t)$ (j = 0,1) are expressed as power series in the small parameter δ ; recall (2.11). The coefficients $F_m^{(j)}(z,t)$ in these series, written in terms of the shifted variable *v* defined in (2.15), can be found by

- (1) solving recursively the second-order ODE (2.20) and using the identity in (2.24), incorporating the initial conditions in (2.53);
- (2) and, subsequently, applying the inverse transformation in (2.56) to the resulting solutions.

Likewise, we can solve the PDE system (2.8) for sufficiently small autorepression rates r; see Assumption 2.2. Its solutions $F^{(j)}(z, t)$ are again expressed as power series in the small parameter δ ; cf. (2.11). The coefficients $F_m^{(j)}(z, t)$ in these series, written in terms of the shifted variable v defined in (2.15), can be found by

- (1) solving recursively the second-order ODE (2.33) and using the identity in (2.34), incorporating the initial conditions in (2.53);
- (2) and, subsequently, applying the inverse transformation in (2.56) to the resulting solutions.

To illustrate the procedure described above, we state the resulting explicit expressions for the leading-order solution to (2.7) – or, equivalently, to (2.8) – in terms of the original variables

z and t:

$$F_{0}^{(0)}(z,t) = \left[1 + (z-1)e^{-t}\right]^{n_{0}}e^{-\lambda(z-1)e^{-t}}\frac{\kappa_{b}}{\kappa_{f}+\kappa_{b}}\left[{}_{1}F_{1}(-\kappa_{b},1-\kappa_{f}-\kappa_{b},\lambda(z-1)e^{-t}) \quad (2.58a) + \frac{\chi_{0}\lambda(z-1)e^{-t}}{1-\kappa_{f}-\kappa_{b}}{}_{1}F_{1}(1-\kappa_{b},2-\kappa_{f}-\kappa_{b},\lambda(z-1)e^{-t})\right]{}_{1}F_{1}(\kappa_{f},1+\kappa_{f}+\kappa_{b},\lambda(z-1)) - \left[1 + (z-1)e^{-t}\right]^{n_{0}}e^{-\lambda(z-1)e^{-t}}\frac{e^{-(\kappa_{f}+\kappa_{b})t}}{\kappa_{f}+\kappa_{b}}\left\{\left[\chi_{0}(\kappa_{f}+\kappa_{b})-\kappa_{b}\right]{}_{1}F_{1}(\kappa_{f},1+\kappa_{f}+\kappa_{b},\lambda(z-1)e^{-t}) + \frac{\kappa_{f}\chi_{0}\lambda(z-1)e^{-t}}{1+\kappa_{f}+\kappa_{b}}{}_{1}F_{1}(1+\kappa_{f},2+\kappa_{f}+\kappa_{b},\lambda(z-1)e^{-t})\right\}{}_{1}F_{1}(-\kappa_{b},1-\kappa_{f}-\kappa_{b},\lambda(z-1)) \quad \text{and}$$

$$F_{0}^{(1)}(z,t) = \left[1 + (z-1)e^{-t}\right]^{n_{0}}e^{-\lambda(z-1)e^{-t}}\frac{\kappa_{f}}{\kappa_{f}+\kappa_{b}}\left[{}_{1}F_{1}(-\kappa_{b},1-\kappa_{f}-\kappa_{b},\lambda(z-1)e^{-t}) - \frac{\chi_{0}\lambda(z-1)e^{-t}}{1+\kappa_{f}+\kappa_{b}}\lambda(z-1)e^{-t}\right]{}_{1}F_{1}(1-\kappa_{b},2-\kappa_{f}-\kappa_{b},\lambda(z-1)e^{-t})\right]$$

$$+ \left[1 + (z - 1)e^{-t}\right]^{n_0}e^{-\lambda(z-1)e^{-t}}\frac{e^{-(\kappa_f + \kappa_b)t}}{\kappa_f + \kappa_b} \left\{ \left[\chi_0(\kappa_f + \kappa_b) - \kappa_b\right]_1 F_1(\kappa_f, 1 + \kappa_f + \kappa_b, \lambda(z - 1)e^{-t}) + \frac{\kappa_f \chi_0 \lambda(z - 1)e^{-t}}{1 + \kappa_f + \kappa_b} {}_1 F_1(1 + \kappa_f, 2 + \kappa_f + \kappa_b, \lambda(z - 1)e^{-t}) \right\}_1 F_1(1 - \kappa_b, 1 - \kappa_f - \kappa_b, \lambda(z - 1))$$

Note that similar expressions were derived by Iyer-Biswas, Hayot and Jayaprakash 2009, where an analogous generating function approach was applied under the assumption that the gene is initially inactive – i.e. that $\chi(n_0) = 1$, see Definition 2.4 – as well as that the initial protein number n_0 is zero. With these choices, the expression for $F_0^{(0)} + F_0^{(1)}$ from(2.58) can be seen to coincide with that found in Equations (5) through (7) of Iyer-Biswas, Hayot and Jayaprakash 2009, using §13.3.4 and §13.2.39 of *NIST Digital Library of Mathematical Functions*.

3 Model B: Gene Expression with Explicit Transcription

In this section, we apply our analytical method to model B, a stochastic gene expression model presented by Shahrezaei and Swain 2008a which explicitly incorporates the transcription stage in the expression process, as well as DNA switching; see also Figure 4. In the original article by Shahrezaei and Swain 2008a, a generating function approach was used to obtain analytical expressions for the time-independent ('stationary') solution to the model. For a visual guide to the upcoming analysis, the reader is again referred to Figure 5.

3.1 Stochastic model and CME

The model for stochastic gene expression considered here is given by the reaction scheme

D	$\stackrel{k_0}{\underset{k_1}{\rightleftharpoons}}$	D^*	(DNA switching),	
		$D^* + M$	(transcription of DNA to mRNA),	
М	$\xrightarrow{\nu_1}$	M + P	(translation of mRNA to protein),	(3.1)
	$\xrightarrow{d_0}$		(decay of mRNA),	
Р	$\xrightarrow{d_1}$	Ø	(decay of protein).	

The modelled gene can hence switch between inactive and active states which are denoted by D and D^* , respectively, with corresponding switching rates k_0 and k_1 . The active gene is transcribed to mRNA (M) with rate v_0 ; mRNA is translated to protein (P) with rate v_1 . Finally, mRNA decays with rate d_0 , while protein decays with rate d_1 .

As in model A, autoregulatory terms can be added to the core reaction scheme in (3.1). Since both mRNA and protein are modelled explicitly, we can identify four distinct autoregulatory mechanisms, in analogy to those in (2.2a) and (2.2b):

$$D + M \xrightarrow{a_M} D^* + M$$
 (autoactivation through mRNA), (3.2a)

$$D^* + M \xrightarrow{r_M} D + M$$
 (autorepression through mRNA), (3.2b)

$$D + P \xrightarrow{a_P} D^* + P$$
 (autoactivation through protein), (3.2c)

$$D^* + P \xrightarrow{r_P} D + P$$
 (autorepression through protein). (3.2d)

Autoactivation can be achieved either by mRNA or by protein, with rates a_M or a_P , respectively; similarly, autorepression can occur either through mRNA or through protein, with respective rates r_M or r_P .

The CME system associated to the reaction scheme in (3.1) is given by

$$\frac{\mathrm{d}P_{m,n}^{(0)}}{\mathrm{d}t} = -\kappa_0 P_{m,n}^{(0)} + \kappa_1 P_{m,n}^{(1)} + \gamma \left[(m+1) P_{m+1,n}^{(0)} - m P_{m,n}^{(0)} \right] + \left[(n+1) P_{m,n+1}^{(0)} - n P_{m,n}^{(0)} \right] + \gamma \mu \left(m P_{m,n-1}^{(0)} - m P_{m,n}^{(0)} \right),$$
(3.3a)

$$\frac{\mathrm{d}P_{m,n}^{(0)}}{\mathrm{d}t} = \kappa_0 P_{m,n}^{(0)} - \kappa_1 P_{m,n}^{(1)} + \gamma \left[(m+1)P_{m+1,n}^{(1)} - mP_{m,n}^{(1)} \right] + \left[(n+1)P_{m,n+1}^{(1)} - nP_{m,n}^{(1)} \right] + \gamma \mu \left(mP_{m,n-1}^{(1)} - mP_{m,n}^{(1)} \right) + \lambda \left(P_{m-1,n}^{(1)} - P_{m,n}^{(1)} \right).$$
(3.3b)

Here, $P_{m,n}^{(j)}(t)$ (j = 0, 1) represents the probability of *m* mRNA and *n* protein being present at time *t* while the gene is either inactive (0) or active (1). As in (2.3) and (2.5), the time variable is nondimensionalised by the protein decay rate d_1 ; other model parameters are scaled as

$$\kappa_0 = \frac{k_0}{d_1}, \quad \kappa_1 = \frac{k_1}{d_1}, \quad \gamma = \frac{d_0}{d_1}, \quad \lambda = \frac{\nu_0}{d_1}, \quad \text{and} \quad \mu = \frac{\nu_1}{d_0}.$$
(3.4)

We note that the above scaling was also used by Shahrezaei and Swain 2008a. The effects of incorporating the autoregulatory mechanisms in (3.2) into the CME system (3.3) are specified in Table 1.

Autoregulation type	Contribution to (3.3a)	Contribution to (3.3b)
mRNA autoactivation	$-\frac{a_M}{d_1}mP_{m,n}^{(0)}$	$+\frac{a_{M}}{d_{1}}mP_{m,n}^{(0)}$
mRNA autorepression	$+\frac{r_{M}}{d_{1}}mP_{m,n}^{(1)}$	$-\frac{r_M}{d_1}mP_{m,n}^{(1)}$
protein autoactivation	$-\frac{a_P}{d_1}nP_{m,n}^{(0)}$	$+\frac{a_P}{d_1}nP_{m,n}^{(0)}$
protein autorepression	$+\frac{r_P}{d_1}nP_{m,n}^{(1)}$	$-\frac{r_P}{d_1}nP_{m,n}^{(1)}$

Table 1: Contribution to the right-hand sides of (3.3) that is due to incorporation of the autoregulatory mechanisms in (3.2).

3.2 Generating function PDE

Since the probabilities $P_{m,n}^{(j)}(t)$ (j = 0, 1) in (3.3) depend on both the mRNA number *m* and the protein number *n*, we introduce probability-generating functions that are defined by double asymptotic series:

$$F^{(j)}(w,z,t) = \sum_{m=0}^{\infty} \sum_{n=0}^{\infty} w^m z^n P^{(j)}_{m,n}(t) \quad \text{for } j = 0, 1.$$
(3.5)

For coefficients $P_{m,n}^{(j)}(t)$ that obey the CME system (3.3), the associated generating functions $F^{(j)}(w, z, t)$ satisfy

$$\partial_t F^{(0)} + (z-1)\partial_z F^{(0)} + \gamma(w-1)\partial_w F^{(0)} - \gamma\mu(z-1)w\partial_w F^{(0)} = -\kappa_0 F^{(0)} + \kappa_1 F^{(1)}, \qquad (3.6a)$$

$$\partial_t F^{(1)} + (z-1)\partial_z F^{(1)} + \gamma(w-1)\partial_w F^{(1)} - \gamma\mu(z-1)w\partial_w F^{(1)} = \kappa_0 F^{(0)} - \kappa_1 F^{(1)}$$

$$+\lambda(w-1)F^{(1)}$$
. (3.6b)

The effects of incorporating the autoregulatory mechanisms in (3.2) into system (3.6) are specified in Table 2.

Autoregulation type	Contribution to (3.6a)	Contribution to (3.6b)
mRNA autoactivation	$-\frac{a_M}{d_1}w\partial_w F^{(0)}$	$+\frac{a_M}{d_1}w\partial_w F^{(0)}$
mRNA autorepression	$+\frac{r_M}{d_1}w\partial_w F^{(1)}$	$-\frac{\dot{r}_{M}}{d_{1}}w\partial_{w}F^{(1)}$
protein autoactivation	$-\frac{a_P}{d_1}z\partial_z F^{(0)}$	$+\frac{a_P}{d_1}z\partial_z F^{(0)}$
protein autorepression	$+\frac{\dot{r_P}}{d_1}z\partial_z F^{(1)}$	$-\frac{\widetilde{r}_{P}}{d_{1}}z\partial_{z}F^{(1)}$

Table 2: Contribution to the right-hand sides of (3.6) that is due to incorporation of the autoregulatory mechanisms in (3.2).

3.3 Dynamical systems analysis

As before, the PDE system (3.6) for the generating function can be reformulated as a system of ODEs via the method of characteristics. The differential operator $\partial_t + (z - 1)\partial_z + \gamma(w - 1)\partial_w - \gamma\mu(z - 1)w\partial_w$ now gives rise to the characteristic system

$$\dot{u} = \gamma [u - \mu v(u+1)], \qquad (3.7a)$$

$$\dot{v} = v, \tag{3.7b}$$

where $\dot{u} = \frac{du}{ds}$ is the derivative along the characteristic, which is parametrised by *s*. For simplicity, we have introduced the new variables *u* and *v*, which are defined as

$$u = w - 1$$
 and $v = z - 1$, (3.8)

respectively. On the resulting characteristics, Equation (3.6) yields the system of ODEs

$$\dot{F}^{(0)} = -\kappa_0 F^{(0)} + \kappa_1 F^{(1)}, \tag{3.9a}$$

$$\dot{F}^{(1)} = \kappa_0 F^{(0)} - \kappa_1 F^{(1)} + \lambda u F^{(1)}.$$
(3.9b)

First, we note that the characteristic system (3.7) can be solved explicitly in terms of the incomplete Gamma function $\Gamma(a, z)$, see §8.2(i) of *NIST Digital Library of Mathematical Functions*), yielding

$$u(s) = e^{\gamma(s-\mu\nu_0 e^s)} \Big\{ u_0 e^{\gamma\mu\nu_0} + (-\gamma\mu\nu_0)^{\gamma} \left[\Gamma(1-\gamma, -\gamma\mu\nu_0) - \Gamma(1-\gamma, -\gamma\mu\nu_0 e^s) \right] \Big\},$$
(3.10a)

$$v(s) = v_0 e^s.$$
 (3.10b)

However, due to its complex nature, the expression for u(s) given in (3.10a) cannot be used to obtain explicit expressions for the generating functions $F^{(j)}$ that solve system (3.9). Therefore, inspired by the analysis by Shahrezaei and Swain 2008a and Popović, Marr and Swain 2016, we make the following assumption:

Assumption 3.1. We assume that the decay rate of protein (d_1) is smaller than the decay rate of mRNA (d_0) ; specifically, we write

$$\frac{d_0}{d_1} = \gamma = \frac{1}{\varepsilon},\tag{3.11}$$

where $0 < \varepsilon < 1$ is sufficiently small.

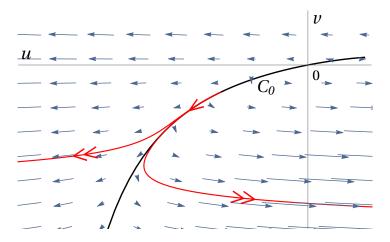


Figure 7: Phase space dynamics of systems (3.12) and (3.13). The slow flow along C_0 is indicated by single arrows; the fast dynamics transverse to C_0 are denoted by double arrows.

The resulting scale separation between mRNA and protein decay rates is well-documented in many microbial organisms, including in bacteria and yeast (Shahrezaei and Swain 2008a; Yu et al. 2006).

For clarity of presentation, we make an additional assumption here.

Assumption 3.2. We assume that all other model parameters κ_0 , κ_1 , λ , and μ , as defined in (3.4), are O(1) in ε .

Remark 3.3. Although Assumption 3.2 is not strictly necessary for the upcoming analysis, it is beneficial. It is worthwhile to note that the analytical scheme presented in this section can be applied in a straightforward fashion in cases where Assumption 3.2 fails, which is particularly relevant in relation to previous work (Shahrezaei and Swain 2008a; Feigelman, Popović and Marr 2015), where the CME system (3.3) is studied for parameter values far beyond the range implied by Assumption 3.2.

Using Assumption 3.1, we can write the characteristic system (3.7) as

$$\varepsilon \dot{u} = u - \mu v(u+1), \tag{3.12a}$$

$$\dot{v} = v. \tag{3.12b}$$

Since ε is assumed to be small, we can classify (3.12) as a *singularly* perturbed *slow-fast* system in standard form; see Kuehn 2015. A comprehensive slow-fast analysis of Equation (3.12) was carried out by Popović, Marr and Swain 2016; we highlight some relevant aspects of that analysis here.

System (3.12) gives rise to a critical manifold $C_0 = \{(u, v) \mid u = \frac{\mu v}{1 - \mu v}\}$. If ε is asymptotically small, orbits of (3.12) can be separated into slow and fast segments, using Fenichel's geometric singular perturbation theory (Kuehn 2015). The critical manifold C_0 is normally repelling; in other words, orbits converge to C_0 in backward time at an exponential rate. For initial conditions asymptotically close to C_0 , orbits initially follow C_0 closely for some time, after which they move away from C_0 under the fast dynamics; see also Figure 7.

To analyse the fast dynamics of system (3.12), we introduce the fast variable $\sigma = \frac{s}{\varepsilon}$; in terms of σ , (3.12) is hence expressed as

$$u' = u - \mu v(u+1),$$
 (3.13a)

$$v' = \varepsilon v, \tag{3.13b}$$

where $u' = \frac{du}{d\sigma}$. We can solve (3.13b) explicitly and write the result as a power series in ε , which yields

$$v(\sigma) = v_0 e^{\varepsilon \sigma} = v_0 \sum_{n=0}^{\infty} \varepsilon^n \frac{\sigma^n}{n!}.$$
(3.14)

Expressing the solution to (3.13a) as a power series in ε , as well, i.e. writing

$$u(\sigma) = \sum_{n=0}^{\infty} \varepsilon^n \hat{u}_n(\sigma), \qquad (3.15)$$

substituting (3.15) into (3.13a), and making use of (3.14), we find

$$\frac{\mathrm{d}\hat{u}_n}{\mathrm{d}\sigma} = \hat{u}_n - \mu v_0 \frac{\sigma^n}{n!} \left[1 + \sum_{l=0}^n \frac{\sigma^{-l} n!}{(n-l)!} \hat{u}_l \right],\tag{3.16}$$

with initial conditions

$$\hat{u}_0(0) = u_0$$
 and $\hat{u}_n(0) = 0$ for $n \ge 1$. (3.17)

The first two terms in (3.15) are thus given by

$$\hat{u}_0(\sigma) = \frac{\mu v_0}{1 - \mu v_0} + \left(u_0 - \frac{\mu v_0}{1 - \mu v_0}\right) e^{(1 - \mu v_0)\sigma} \quad \text{and}$$
(3.18a)

$$\hat{u}_1(\sigma) = \frac{\mu v_0}{(1-\mu v_0)^2} \left[\frac{1-e^{(1-\mu v_0)\sigma}}{1-\mu v_0} + \sigma \right] - \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \mu v_0 \frac{\sigma^2}{2} e^{(1-\mu v_0)\sigma}.$$
 (3.18b)

We now employ the expansion in (3.15) to obtain explicit expressions for the generating functions $F^{(j)}$ (j = 0, 1). In the fast variable $\sigma = \frac{s}{\varepsilon}$, system (3.9) takes the form

$$F^{(0)\prime} = -\varepsilon \kappa_0 F^{(0)} + \varepsilon \kappa_1 F^{(1)}, \qquad (3.19a)$$

$$F^{(1)\prime} = \varepsilon \kappa_0 F^{(0)} - \varepsilon \kappa_1 F^{(1)} + \varepsilon \lambda u F^{(1)}.$$
(3.19b)

As in the analysis of model A, recall Section 2.3, we rewrite system (3.19) as a second-order ODE for $F^{(0)}$ to find

$$\frac{\mathrm{d}^2}{\mathrm{d}\sigma^2}F^{(0)} + \varepsilon(\kappa_0 + \kappa_1 - \lambda u)\frac{\mathrm{d}}{\mathrm{d}\sigma}F^{(0)} - \varepsilon^2\lambda\kappa_0 uF^{(0)} = 0.$$
(3.20)

Next, we use (3.19a) to express $F^{(1)}$ in terms of $F^{(0)}$ as

$$F^{(1)} = \frac{1}{\varepsilon \kappa_1} \left(F^{(0)\prime} + \varepsilon \kappa_0 F^{(0)} \right). \tag{3.21}$$

To incorporate the expansion for u in (3.15), we also expand $F^{(j)}$ (j = 0, 1) in powers of ε , writing

$$F^{(j)}(\sigma) = \sum_{n=0}^{\infty} \varepsilon^n F_n^{(j)}(\sigma).$$
(3.22)

Substitution of (3.22) into (3.20) then yields

$$\frac{d^2}{d\sigma^2}F_0^{(0)} = 0, \qquad (3.23a)$$

$$\frac{\mathrm{d}^2}{\mathrm{d}\sigma^2}F_1^{(0)} + (\kappa_0 + \kappa_1 - \lambda\hat{u}_0)\frac{\mathrm{d}}{\mathrm{d}\sigma}F_0^{(0)} = 0, \qquad (3.23\mathrm{b})$$

and

$$\frac{\mathrm{d}^2}{\mathrm{d}\sigma^2} F_n^{(0)} + (\kappa_0 + \kappa_1 - \lambda \hat{u}_0) \frac{\mathrm{d}}{\mathrm{d}\sigma} F_{n+1}^{(0)} - \lambda \sum_{k=0}^n \left(\hat{u}_{n+1-k} \frac{\mathrm{d}}{\mathrm{d}\sigma} + \kappa_0 \, \hat{u}_{n-k} \right) F_k^{(0)} = 0$$
(3.24)

for $n \ge 2$. By combining (3.21) with (3.22), we obtain

$$\frac{d}{d\sigma}F_0^{(0)} = 0 \tag{3.25}$$

and

$$F_n^{(1)} = \frac{\kappa_0}{\kappa_1} F_n^{(0)} + \frac{1}{\kappa_1} \frac{\mathrm{d}}{\mathrm{d}\sigma} F_{n+1}^{(0)}$$
(3.26)

for $n \ge 0$. We can solve Equations (3.23) and (3.24) iteratively – taking into account the additional condition on $F_0^{(0)}$ in (3.25) – to find

$$F_0^{(0)} = f_0, (3.27a)$$

$$F_1^{(0)} = f_1 + g_1 \sigma$$
, and (3.27b)

$$F_{2}^{(0)} = f_{2} + g_{2}\sigma + \frac{\lambda(f_{0}\kappa_{0} + g_{1})}{1 - \mu\nu_{0}} \left[\frac{e^{(1 - \mu\nu_{0})\sigma}}{1 - \mu\nu_{0}} \left(u_{0} - \frac{\mu\nu_{0}}{1 - \mu\nu_{0}} \right) + \mu\nu_{0}\frac{\sigma^{2}}{2} \right] - g_{1}(\kappa_{0} + \kappa_{1})\frac{\sigma^{2}}{2} \quad (3.27c)$$

for the first three terms which, upon substitution into (3.26), yields

$$F_0^{(1)} = \frac{f_0 \kappa_0 + g_1}{\kappa_1}$$
 and (3.28a)

$$F_{1}^{(1)} = \frac{f_{1}\kappa_{0} + g_{2}}{\kappa_{1}} + \frac{\lambda(f_{0}\kappa_{0} + g_{1})}{\kappa_{1}(1 - \mu v_{0})} \left[e^{(1 - \mu v_{0})\sigma} \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) + \sigma \mu v_{0} \right] - \sigma g_{1};$$
(3.28b)

here, f_i and g_i are free constants, to be determined by initial conditions; see Section 3.4.

Contrary to common practice in the study of slow-fast systems such as (3.19), we forego a detailed analysis of the slow system (3.9), continuing our discussion with the determination of appropriate initial conditions; cf. Section 2.4. For details on why the slow dynamics is disregarded, the reader is referred to Remark 3.5.

3.4 Initial conditions and inverse transformation

To complete our analytical method, we discuss the determination of initial conditions and the reconstruction of the solution to the original PDE system (3.6), as was done for model A in Sections 2.4 and 2.5, respectively.

3.4.1 Initial conditions

We follow the reasoning of Section 2.4, and determine appropriate initial conditions by considering the original CME system (3.3).

At time t = 0, we prescribe initial mRNA and protein numbers m_0 and n_0 , respectively. As in Section 2.4, we again introduce the parameter χ , which is defined as follows; compare Definition 2.4:

Definition 3.4. Consider the CME system (3.3). We define $\chi(m_0, n_0)$ to be the probability that the gene modelled by the reaction scheme in (3.1) is switched off at time t = 0, given initial mRNA and protein numbers m_0 and n_0 , respectively.

For the probabilities $P_{m,n}^{(j)}(t)$ (j = 0, 1), Definition 3.4 implies that

$$P_{m,n}^{(0)}(0) = \chi \,\delta_{m,m_0} \,\delta_{n,n_0}$$
 and (3.29a)

$$P_{m,n}^{(1)}(0) = (1 - \chi)\delta_{m,m_0}\delta_{n,n_0}.$$
(3.29b)

Using the definition of the generating functions $F^{(j)}(w, z, t)$ in (3.5), we find

$$F^{(0)}(w, z, 0) = \chi w^{m_0} z^{n_0} = \chi (1+u)^{m_0} (1+v)^{n_0}$$
 and (3.30a)

$$F^{(1)}(w,z,0) = (1-\chi)w^{m_0}z^{n_0} = (1-\chi)(1+u)^{m_0}(1+v)^{n_0},$$
(3.30b)

taking into account the change of variables in (3.8).

Next, we can infer e.g. from (3.10) and well-known properties of the incomplete Gamma function – see §8.2 of *NIST Digital Library of Mathematical Functions* – that solutions to (3.7) exist globally, i.e. for all *s*. In particular, given an arbitrary triple (s_*, u_*, v_*) , we can apply the inverse flow of (3.7) to flow backward (u_*, v_*) for time s_* . We conclude that the characteristics of the operator $\partial_t + (z-1)\partial_z + \gamma(w-1)\partial_w - \gamma\mu(z-1)w\partial_w = \partial_t + v\partial_v + \gamma u\partial_u - \gamma\mu(u+1)v\partial_u$, which are given by the orbits of system (3.7), foliate the (t, u, v)-coordinate space over the $\{t = 0\}$ -plane when the parameter along the characteristic – interpreted as a fibre over the $\{t = 0\}$ -plane – by its base point. Because orbits of (3.7) provide a parametrisation of the underlying characteristics, the coordinates of that base point are given by $(0, u_0, v_0)$, where (u_0, v_0) are the initial values of the corresponding orbit of (3.7). Hence, the conditions in (3.30) yield, on each characteristic,

$$\left[F^{(0)}\right]_{s=0} = \chi (1+u_0)^{m_0} (1+v_0)^{n_0} \quad \text{and} \tag{3.31a}$$

$$\left[F^{(1)}\right]_{s=0} = (1-\chi)(1+u_0)^{m_0}(1+v_0)^{n_0}.$$
(3.31b)

As s = 0 implies $\sigma = 0$, we can apply the initial conditions in (3.31) to the solutions of the ODE system (3.19) in order to determine the free constants f_i and g_i in (3.27) and (3.28). Combining the power series expansion in (3.22) with (3.31), we obtain

$$\left[F_0^{(0)}\right]_{\sigma=0} = \chi (1+u_0)^{m_0} (1+v_0)^{n_0}, \qquad (3.32a)$$

$$\left[F_{1}^{(1)}\right]_{\sigma=0} = (1-\chi)(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}, \text{ and}$$
(3.32b)

$$\left[F_{n}^{(j)}\right]_{\sigma=0} = 0 \quad \text{for all } n \ge 1, \text{ with } j = 0, 1,$$
 (3.32c)

which implies

$$f_0 = \chi (1 + u_0)^{m_0} (1 + v_0)^{n_0}, \tag{3.33a}$$

$$f_1 = 0,$$
 (3.33b)

$$f_2 = -(1-\chi)(1+u_0)^{m_0}(1+v_0)^{n_0}\frac{\lambda\kappa_1}{(1-\mu v_0)^2}\left(u_0 - \frac{\mu v_0}{1-\mu v_0}\right),$$
(3.33c)

$$g_1 = (\kappa_1(1-\chi) - \kappa_0\chi)(1+u_0)^{m_0}(1+v_0)^{n_0}, \text{ and}$$
 (3.33d)

$$g_2 = -(1-\chi)(1+u_0)^{m_0}(1+v_0)^{n_0}\frac{\lambda\kappa_1}{1-\mu v_0}\left(u_0 - \frac{\mu v_0}{1-\mu v_0}\right)$$
(3.33e)

in (3.27) and (3.28).

3.4.2 Inverse transformation

Since the (t, u, v)-coordinate space is foliated by the characteristics of the operator $\partial_t + v\partial_v + \gamma u\partial_u - \gamma \mu(u+1)v\partial_u$, any point (t, u, v) lies on a unique characteristic. Flowing backward along that characteristic to its intersection with the $\{t = 0\}$ -plane, we can determine the corresponding base point $(0, u_0, v_0)$ by inverting the relations in (3.10). Since the dynamics of the *v*-coordinate do not depend on *u*, we may use (2.56) to express v_0 in terms of *t* and *v* only. Taking the resulting expression as input for inverting (3.10a), we obtain the inverse characteristic transformation

$$(t, u, v) \mapsto (u_0(t, u, v), v_0(t, v)),$$
 (3.34)

$$u_{0}(t, u, v) = e^{-\gamma(t+\mu ve^{-t})} \left\{ u e^{\gamma \mu v} - (-\gamma \mu v)^{\gamma} \left[\Gamma(1-\gamma, -\gamma \mu ve^{-t}) - \Gamma(1-\gamma, -\gamma \mu v) \right] \right\} \text{ and } (3.35a)$$

$$v_{0}(t, v) = ve^{-t}. \tag{3.35b}$$

Under Assumption 3.1, we can employ the power series expansion in (3.15), in combination with the recursive set of ODEs in (3.16) and initial conditions as in (3.17), as an alternative to (3.35a) to obtain u_0 as a function of u, v (or v_0), and σ . Rewriting the result as a power series in ε , we find

$$u_{0}(\sigma, u, v) = \frac{\mu v}{1 - \mu v} + e^{(\mu v - 1)\sigma} \left(u - \frac{\mu v}{1 - \mu v} \right) - \varepsilon \mu v \left[\frac{e^{(\mu v - 1)\sigma} - 1}{(1 - \mu v)^{3}} + \frac{\sigma}{(1 - \mu v)^{2}} + \frac{\sigma^{2}}{2} e^{(\mu v - 1)\sigma} \left(u - \frac{\mu v}{1 - \mu v} \right) \right] + O(\varepsilon^{2}); \quad (3.36)$$

naturally, (3.14) gives rise to

$$v_0(\sigma, v) = v_0 \sum_{n=0}^{\infty} \varepsilon^n \frac{(-\sigma)^n}{n!}.$$
 (3.37)

Since the independent variable in (3.16) is σ , and since *s* is naturally identified with *t*, we replace σ with $\frac{t}{s}$ in (3.36) to obtain a perturbative expansion for $u_0(t, u, v)$.

The solution to the PDE system (3.6) can now be found to satisfactory order in ε by applying the inverse transformation in (3.34) to the solutions given in (3.27) and (3.28), taking into account the values of f_i and g_i in (3.33). In other words,

$$F^{(j)}(u, v, t) \text{ [as solution to (3.6)]} = \left[\sum_{n=0}^{\infty} \varepsilon^n \left[F_n^{(j)}\right]_{(u_0, v_0) = (u_0(\sigma, u, v), ve^{-\varepsilon\sigma})}\right]_{\sigma = \frac{1}{\varepsilon}},$$
(3.38)

where $F_n^{(j)}$ (j = 0, 1) on the right-hand side of the above expression denotes the solution to Equations (3.23a) through (3.26), with initial conditions as in (3.32).

Remark 3.5. The absence of any detailed analysis of the characteristic system in its slow formulation, Equation (3.12), can be argued as follows, by considering the corresponding phase space, as depicted in Figure 7.

- (a) For arbitrary initial conditions (u_0, v_0) , the dominant dynamics are fast, since the critical manifold C_0 is normally repelling. In other words, solutions are generally repelled away from C_0 under the fast dynamics.
- (b) All orbits that have their initial conditions on the same Fenichel fibre are *exponentially* close (in ε) to each other near the slow manifold C_{ε} that is associated to the critical manifold C_0 . Therefore, flowing backward from (u, v) to (u_0, v_0) as expressed through the inverse transformation in (3.34) that yields the corresponding PDE solution may introduce *exponentially large* terms in the transformation, precluding any sensible series expansion.

Thus, although the construction of a composite '(initially) slow – (ultimately) fast' expression of $F^{(j)}$ as a solution to systems (3.12) and (3.13) certainly makes sense from a dynamical systems perspective, the extreme lack of sensitivity of orbits on their initial conditions (u_0 , v_0) may prevent such a composite expansion from being useful for obtaining solutions to the original PDE system (3.6).

Remark 3.6. In Section 3.3, explicit expressions are given for the expansion of $F^{(0)}$ up to $O(\varepsilon^2)$ only, cf. (3.27); similarly, $F^{(1)}$ is approximated to $O(\varepsilon)$ in (3.28), for the sake of brevity. It is worthwhile to note that a lower bound on the order of the expansion is stipulated by the

with

application; recall Section 1.1: the sampling time Δt can be considered as a *minimum* time interval over which the results of our analysis should be (reasonably) accurate. To that end, we have to compare Δt with $\varepsilon = \frac{1}{\gamma}$, the parameter defining the fast time scale on which the above analytical results have been derived. We can then apply the classical theory of Poincaré expansions (Verhulst 2000) to infer that, if for example $\Delta t = O(1)$ – which implies $\Delta t = O(\varepsilon^{-1})$ in the fast time variable σ – the generating functions $F^{(j)}$ should at least be expanded up to $O(\varepsilon^2)$ for the resulting approximation to be accurate to $O(\varepsilon)$.

3.5 Autoregulation

The inclusion of any type of autoregulation into system (3.6) – which is equivalent to the addition of model terms from Table 2 to the right-hand sides of the corresponding equations – precludes the direct application of the method of characteristics, as the resulting partial differential operators in Equations (3.6a) and (3.6b) do not coincide anymore. To resolve that complication, we follow the approach of Section 2, making the following assumption:

Assumption 3.7. We assume that the autoregulation rates a_M , r_M , a_P , and r_P , as defined in (3.2), are small in comparison with the protein decay rate d_1 ; specifically, we write

$$a_M = \alpha_M d_1 \delta, \quad r_M = \rho_M d_1 \delta, \quad a_P = \alpha_P d_1 \delta, \quad and \quad r_P = \rho_P d_1 \delta, \quad (3.39)$$

where $0 < \delta < 1$ is sufficiently small.

Next, we expand the generating functions $F^{(j)}$ (j = 0, 1) as power series in δ ; recall (2.11):

$$F^{(j)}(z,w,t) = \sum_{m=0}^{\infty} \delta^m F_m^{(j)}(z,w,t).$$
(3.40)

To demonstrate the procedure, we include mRNA autoactivation in (3.6), see again Table 2; the analysis of the remaining autoregulatory mechanisms can be performed in a similar fashion. Substitution of (3.40) now yields

$$\left[\partial_{t} + (z-1)\partial_{z} + \gamma(w-1)\partial_{w} - \gamma\mu(z-1)w\partial_{w} \right] F_{m}^{(0)} = -\kappa_{0}F_{m}^{(0)} + \kappa_{1}F_{m}^{(1)} - \alpha_{M}w\partial_{w}F_{m-1}^{(0)}, \quad (3.41a)$$

$$\left[\partial_{t} + (z-1)\partial_{z} + \gamma(w-1)\partial_{w} - \gamma\mu(z-1)w\partial_{w} \right] F_{m}^{(1)} = \kappa_{0}F_{m}^{(0)} - \kappa_{1}F_{m}^{(1)} + \lambda(w-1)F_{m}^{(1)} + \alpha_{M}w\partial_{w}F_{m-1}^{(0)}; \quad (3.41b)$$

cf. (2.12). System (3.41) is then amenable to the method of characteristics; in fact, employing the same characteristics as in the unperturbed setting, recall (3.7), we find

$$\partial_t + (z-1)\partial_z + \gamma(w-1)\partial_w - \gamma\mu(z-1)w\partial_w = \partial_s, \qquad (3.42)$$

while the partial differential operators in Table 2 transform into

$$w\partial_w = (u+1) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0}$$
 and (3.43a)

$$z\partial_z = (v+1)\left(\frac{\partial v}{\partial v_0}\right)^{-1} \left[-\left(\frac{\partial u}{\partial v_0}\right) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} + \partial_{v_0} \right];$$
(3.43b)

here, $u(s; u_0, v_0)$ and $v(s; v_0)$ are as given in (3.10). Thus, the mRNA autoactivation system (3.41) transforms to

$$\partial_s F_m^{(0)} = -\kappa_0 F_m^{(0)} + \kappa_1 F_m^{(1)} - \alpha_M (u+1) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} F_{m-1}^{(0)}, \qquad (3.44a)$$

$$\partial_s F_m^{(1)} = \kappa_0 F_m^{(0)} - \kappa_1 F_m^{(1)} + \lambda \, u F_m^{(1)} + \alpha_M (u+1) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} F_{m-1}^{(0)}. \tag{3.44b}$$

To obtain explicit solutions to system (3.44), we adopt Assumption 3.1 and revert to the fast time scale $\sigma = \frac{s}{\varepsilon}$ to write the dynamical system (3.44) as a second-order ODE for $F_m^{(0)}$, which yields

$$\left[\partial_{\sigma}^{2} + \varepsilon(\kappa_{0} + \kappa_{1} - \lambda u)\partial_{\sigma} - \varepsilon^{2}\kappa_{0}\lambda u\right]F_{m}^{(0)} = (\varepsilon^{2}\lambda u - \varepsilon\partial_{\sigma})\alpha_{M}(u+1)\left(\frac{\partial u}{\partial u_{0}}\right)^{-1}\partial_{u_{0}}F_{m-1}^{(0)}.$$
 (3.45)

Using (3.44), we can express $F_m^{(1)}$ in terms of $F_m^{(0)}$ as

$$F_m^{(1)} = \frac{1}{\varepsilon \kappa_1} \left[\partial_\sigma F_m^{(0)} + \varepsilon \kappa_0 F_m^{(0)} + \varepsilon \alpha_M (u+1) \left(\frac{\partial u}{\partial u_0} \right)^{-1} \partial_{u_0} F_{m-1}^{(0)} \right].$$
(3.46)

To solve (3.45) (recursively), we expand $F_m^{(j)}$ (j = 0, 1) in powers of ε :

$$F_m^{(j)}(\sigma) = \sum_{n=0}^{\infty} \varepsilon^n F_{m,n}^{(j)}(\sigma); \qquad (3.47)$$

recall Equation (3.22). Together with the series expansion for u in (3.15), we thus obtain

$$\partial_{\sigma}^2 F_{m,0}^{(0)} = 0, \tag{3.48}$$

$$\partial_{\sigma}^{2} F_{m,1}^{(0)} = -(\kappa_{0} + \kappa_{1} - \lambda \hat{u}_{0}) \partial_{\sigma} F_{m,0}^{(0)} - \partial_{\sigma} G_{m-1,0}, \qquad (3.49)$$

and

$$\partial_{\sigma}^{2} F_{m,n}^{(0)} + (\kappa_{0} + \kappa_{1} - \lambda \hat{u}_{0}) \partial_{\sigma} F_{m,n+1}^{(0)} - \lambda \sum_{k=0}^{n} \left(\hat{u}_{n+1-k} \partial_{\sigma} + \kappa_{0} \, \hat{u}_{n-k} \right) F_{m,k}^{(0)} = -\partial_{\sigma} G_{m-1,n+1}^{(0)} + \lambda \sum_{k=0}^{n} \hat{u}_{n-k} G_{m-1,k}^{(0)}$$
(3.50)

for $n \ge 2$; compare with Equations (3.23a), (3.23b), and (3.24). The coefficients $G_{m,n}^{(0)}$ in the above expression are defined from an expansion of the autoregulation term as

$$\alpha_M(u+1) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} F_{m-1}^{(0)} = \sum_{n=0}^{\infty} \varepsilon^n G_{m-1,n}^{(0)}(u_0,\sigma).$$
(3.51)

From (3.46), we obtain

$$\partial_{\sigma} F_{m,0}^{(0)} = 0$$
 (3.52)

and

$$F_{m,n}^{(1)} = \frac{1}{\kappa_1} \partial_\sigma F_{m,n+1}^{(0)} + \frac{\kappa_0}{\kappa_1} F_{m,n}^{(0)} + \frac{1}{\kappa_1} G_{m-1,n}^{(0)}$$
(3.53)

for $n \ge 0$; recall (3.25) and (3.26). To solve Equations (3.48) through (3.53) iteratively, we fix m – the order of the expansion in δ – and determine the solution to satisfactory order in n, the order of the expansion in ε . Then, we increase m to m + 1 and take the result as input for the dynamics at order m+1. The resulting repeated iteration procedure yields an explicit expression for the generating functions $F^{(j)}$ (j = 0, 1) as double asymptotic series in both δ and ε .

The determination of appropriate initial conditions is largely analogous to the non-autoregulated case; see Section 3.4.1. However, with the inclusion of autoregulation into model B, we need to incorporate the possibility that $\chi(m_0, n_0)$ depends on the corresponding autoregulation rates. As in the case of model A, we expand $\chi(m_0, n_0)$ as a power series in δ :

$$\chi(m_0, n_0) = \sum_{m=0}^{\infty} \delta^m \chi_m(m_0, n_0);$$
(3.54)

recall (2.51). In that case, the initial conditions for $F_{m,n}^{(j)}$ can be inferred from (3.32) to give

$$\left[F_{m,0}^{(0)}\right]_{\sigma=0} = \chi_m (1+u_0)^{m_0} (1+v_0)^{n_0} \quad \text{for all } m \ge 0,$$
(3.55a)

$$\left[F_{0,1}^{(1)}\right]_{\sigma=0} = (1-\chi_0)(1+u_0)^{m_0}(1+v_0)^{n_0}, \qquad (3.55b)$$

$$\left[F_{m,1}^{(1)}\right]_{\sigma=0} = -\chi_m (1+u_0)^{m_0} (1+v_0)^{n_0} \quad \text{for all } m \ge 1, \text{ and}$$
(3.55c)

$$\left[F_{m,n}^{(j)}\right]_{\sigma=0} = 0$$
 for all $m \ge 0$ and $n \ge 1$, with $j = 0, 1$. (3.55d)

Solutions to Equations (3.48) through (3.53) that incorporate the conditions in (3.55) for all types of autoregulation introduced in (3.2) can be found in Appendix B. Finally, our previous results on the inverse characteristic transformation in the non-autoregulated case from Section 3.4.2 can now be applied in a straightforward fashion to give solutions to the PDE system (3.6) with added autoregulation.

3.6 Summary of main result

To summarise Section 3, we combine the analysis of the previous subsections to state our main result.

Main result: The PDE system (3.6) can be solved for sufficiently large values of γ ; see Assumption 3.1. Its solutions $F^{(j)}(w, z, t)$ (j = 0, 1) are expressed as power series in the small parameter $\varepsilon = \frac{1}{\gamma}$; cf. (3.22). The coefficients $F_n^{(j)}(w, z, t)$ in these series, written in terms of the shifted variables *u* and *v* defined in (3.8), can be found by

- (1) solving recursively the second-order ODEs (3.23a) through (3.24) and using the identities in (3.25) and (3.26), incorporating the initial conditions in (3.32);
- (2) subsequently applying the inverse transformations in (3.36) and (3.37) to the resulting solutions;
- (3) and, finally, substituting $\sigma = \frac{t}{s}$.

To illustrate the procedure described above, we state the resulting explicit expressions for the leading-order solution to (3.6) in terms of the original variables *w*, *z*, and *t* here:

$$F_0^{(0)}(w,z,t) = \chi \left[1 + (z-1)e^{-t}\right]^{n_0} \left\{ \frac{1}{1+\mu(1-z)} + e^{-\left[1+\mu(1-z)\right]\frac{t}{\varepsilon}} \left[w - \frac{1}{1+\mu(1-z)}\right] \right\}^{m_0} \quad \text{and}$$
(3.56a)

$$F_0^{(1)}(w,z,t) = \frac{1-\chi}{\chi} F_0^{(0)}(w,z,t).$$
(3.56b)

Note that the sum $F_0^{(0)} + F_0^{(1)}$ corresponds precisely to the leading-order fast expansion found in Equation (21) of Bokes et al. 2012b.

If autoregulation as in (3.2) is incorporated into model B, the main result can be formulated as follows.

Main result (autoregulatory extension): The PDE system (3.6) incorporating any one type of autoregulation from Table 2 can be solved as long as γ is sufficiently large and δ is sufficiently small; see Assumptions 3.1 and 3.7, respectively. Its solutions $F^{(j)}(w, z, t)$ (j = 0, 1) are expressed as double power series in the small parameters δ and ε , viz.

$$F^{(j)} = \sum_{m=0}^{\infty} \sum_{n=0}^{\infty} \delta^m \varepsilon^n F^{(j)}_{m,n}.$$
 (3.57)

The coefficients $F_{m,n}^{(j)}$ in these series, written in terms of the shifted variables *u* and *v* defined in (3.8), can be found by

- (1a) solving recursively the second-order ODEs (3.48) through (3.50) for fixed *m* and using the identities in (3.52) and (3.53), incorporating the initial conditions in (3.55);
- (1b) increasing m to m + 1, and repeating step (1a) until a sufficient accuracy in δ (and ε) is attained;
- (2) subsequently applying the inverse transformations in (3.36) and (3.37) to the resulting solutions;
- (3) and, finally, substituting $\sigma = \frac{t}{\epsilon}$.

4 From Generating Function to Propagator

The final step in our analytical method consists in reconstructing the probabilities $P_n^{(j)}$ (model A) and $P_{m,n}^{(j)}$ (model B), respectively, from the explicit expressions for the associated generating functions $F^{(j)}$ (j = 0, 1), which were the main analytical outcome of Sections 2 and 3.

In principle, the relation between probabilities and probability-generating functions is clear from the definition of the latter, and is given in (2.6) and (3.5), respectively. Specifically, probabilities can be expressed in terms of derivatives of their generating functions as follows:

$$P_{n}^{(j)}(t) = \frac{1}{n!} \left[\frac{\partial^{n}}{\partial z^{n}} F^{(j)}(z, t) \right]_{z=0}$$
(model A), (4.1a)

$$P_{m,n}^{(j)}(t) = \frac{1}{m!n!} \left[\frac{\partial^{m+n}}{\partial w^m \partial z^n} F^{(j)}(w, z, t) \right]_{(w,z)=(0,0)}$$
(model B). (4.1b)

However, explicit expressions for the *n*-th and (m + n)-th order derivatives, respectively, of these generating functions become progressively unwieldy with increasing *m* and *n*. Indeed, from the expressions for the generating functions $F^{(j)}$ obtained previously, which combine (2.26) for specific initial conditions as in (2.54) with the inverse characteristic transformation in (2.56), it is clear that finding explicit expressions for derivatives of arbitrary order is very difficult indeed, if it is possible at all.²

To complete successfully the final step towards approximating propagators for parameter inference in the present setting, we abandon the requirement of deriving explicit expressions for the probabilities $P_n^{(j)}$ and $P_{m,n}^{(j)}$. Instead, we use the standard Cauchy integral formula for derivatives of holomorphic functions to write

$$P_n^{(j)}(t) = \frac{1}{2\pi i} \oint_{\gamma_A} \frac{F^{(j)}(z,t)}{z^{n+1}} \,\mathrm{d}z \qquad (\text{model A}), \tag{4.2a}$$

$$P_{m,n}^{(j)}(t) = \frac{1}{(2\pi i)^2} \bigoplus_{\gamma_B} \frac{F^{(j)}(w, z, t)}{w^{m+1} z^{n+1}} \, \mathrm{d}w\mathrm{d}z \qquad (\text{model B}); \tag{4.2b}$$

here, γ_A is a suitably chosen contour around z = 0, while γ_B is a (double) contour around (w, z) = (0, 0).

The above expression of probabilities as integrals is well suited for an efficient numerical implementation, which is naturally incorporated into the realisation of the parameter inference scheme discussed in Section 1. From a numerical perspective, the integral formula in (4.2) has the additional advantage that the values of $F^{(j)}$ on (a discretisation of) the integral contours γ_A and γ_B , respectively, only have to be determined once to yield propagators for any values of *m*

²In some specific cases, however, such explicit expressions can be found; see e.g. Popović, Marr and Swain 2016.

and n and fixed initial states m_0 and n_0 . Moreover, we are free to choose the integration contours γ_A and γ_B , which allows us to accelerate the calculation of these integrals; see Bornemann 2011. Here, we note that the choice of circular integration contours with unit radius, and subsequent discretisation of those contours as *M*-sided and *N*-sided polygons, respectively, coincides with the 'Fourier mode' approach, as presented by Bokes et al. 2012a.

Remark 4.1. By introducing the Cauchy integral formula for derivatives of holomorphic functions in (4.2), we implicitly assume that the integration contours γ_A and γ_B are chosen such that they lie completely within the open neighbourhoods of the origin in \mathbb{C} and \mathbb{C}^2 , respectively, where the canonical complex extensions of the generating functions $F^{(j)}$ – which exist by the Cauchy-Kowalevski theorem – are holomorphic. In other words, γ_A and γ_B must be chosen such that any poles of $F^{(j)}$ lie outside of these integration contours. The expansion for u_0 in (3.36) shows that this is not a moot point: the generating functions $F^{(j)}$ resulting for model B, as established in Section 3, will generically have a pole at $v = \frac{1}{\mu}$, i.e. at $z = 1 + \frac{1}{\mu}$. As μ is positive, by (3.4), choosing the z-contour of γ_B to be a circle with at most unit radius allows us to avoid that pole.

4.1 **Incorporation** of χ

In the course of the analysis presented in Sections 2 and 3, the introduction of the parameter χ was necessary to obtain definite, explicit expressions for the generating functions as solutions to the PDE systems (2.7), (2.8), and (3.6); see Definitions 2.4 and 3.4. The successful implementation of these expressions in a parameter inference scheme requires us to decide how to incorporate that new parameter. We identify three options here.

- 1. Before implementing parameter inference, we can marginalise over the new parameter χ to eliminate it altogether, using a predetermined measure $d\mu(\chi)$, which adds an additional integration step to the requisite numerical scheme.
- 2. We can make a choice for χ that is based on the specifics of the model under consideration. Thus, exploiting the Markov property of the stochastic models underlying (2.3), (2.5), and (3.3), we may use the switching rates and any autoregulation rates to express χ in model A as

$$\chi(n_0) = \frac{c_b}{c_b + c_f + an_0}$$
 (autoactivation), (4.3a)
$$\chi(n_0) = \frac{c_b + rn_0}{c_b + rn_0 + c_f}$$
 (autorepression); (4.3b)

the corresponding expressions for model B read

 k_1

$$\chi(m_0, n_0) = \frac{k_1}{k_0 + k_1}$$
 (no autoregulation), (4.4a)

$$\chi(m_0, n_0) = \frac{\kappa_1}{k_0 + k_1 + a_M m_0}$$
(mRNA autoactivation), (4.4b)

$$\chi(m_0, n_0) = \frac{k_1 + r_M m_0}{k_0 + k_1 + r_M m_0}$$
(mRNA autorepression), (4.4c)
$$\chi(m_0, n_0) = \frac{k_1}{k_0 + k_1 + a_P n_0}$$
(protein autoactivation), (4.4d)
$$k_1 + r_P n_0$$

$$\chi(m_0, n_0) = \frac{\kappa_1 + r_P n_0}{k_0 + k_1 + r_P n_0}$$
 (protein autorepression). (4.4e)

3. We can determine χ 'experimentally' by including the latter in the parameter set that is to be inferred in the (numerical) process of parameter inference.

Note that step 2 has been anticipated in the analysis of model A, by introducing the series expansion in (2.51). Indeed, by Assumption 2.2, we can expand $\chi(n_0)$ as

$$\chi = \frac{\kappa_b}{\kappa_b + \kappa_f + \delta\alpha n_0} = \frac{\kappa_b}{\kappa_f + \kappa_b} \sum_{m=0}^{\infty} \delta^m \left(\frac{-\alpha n_0}{\kappa_f + \kappa_b}\right)^m$$
(autoactivation), (4.5a)

$$\chi = \frac{\kappa_b + \delta\rho n_0}{\kappa_b + \kappa_f + \delta\rho n_0} = \frac{\kappa_b}{\kappa_f + \kappa_b} - \frac{\kappa_f}{\kappa_f + \kappa_b} \sum_{m=1}^{\infty} \delta^m \left(\frac{-\rho n_0}{\kappa_f + \kappa_b}\right)^m \qquad \text{(autorepression).} \tag{4.5b}$$

Likewise, when autoregulation is added to model B, Assumption 3.7 implies an expansion for $\chi(m_0, n_0)$ of the form

$$\chi = \frac{\kappa_1}{\kappa_0 + \kappa_1 + \delta \alpha_M m_0} = \frac{\kappa_1}{\kappa_0 + \kappa_1} \sum_{m=0}^{\infty} \delta^m \left(\frac{-\alpha_M m_0}{\kappa_0 + \kappa_1}\right)^m$$
(mRNA autoactivation),
(4.6a)

$$\chi = \frac{\kappa_1 + \delta \rho_M m_0}{\kappa_0 + \kappa_1 + \delta \rho_M m_0} = \frac{\kappa_1}{\kappa_0 + \kappa_1} - \frac{\kappa_0}{\kappa_0 + \kappa_1} \sum_{m=1}^{\infty} \delta^m \left(\frac{-\rho_M m_0}{\kappa_0 + \kappa_1}\right)^m \quad \text{(mRNA autorepression)},$$
(4.6b)

$$\chi = \frac{\kappa_1}{\kappa_0 + \kappa_1 + \delta \alpha_P n_0} = \frac{\kappa_1}{\kappa_0 + \kappa_1} \sum_{m=0}^{\infty} \delta^m \left(\frac{-\alpha_P n_0}{\kappa_0 + \kappa_1}\right)^m$$
(protein autoactivation),
(4.6c)

$$\chi = \frac{\kappa_1 + \delta\rho_P n_0}{\kappa_0 + \kappa_1 + \delta\rho_P n_0} = \frac{\kappa_1}{\kappa_0 + \kappa_1} - \frac{\kappa_0}{\kappa_0 + \kappa_1} \sum_{m=1}^{\infty} \delta^m \left(\frac{-\rho_P n_0}{\kappa_0 + \kappa_1}\right)^m \qquad \text{(protein autorepression).}$$
(4.6d)

5 Discussion and Outlook

In the present article, we have developed an analytical method for obtaining explicit, fully timedependent expressions for the probability-generating functions that are associated to models for stochastic gene expression. Moreover, we have presented a computationally efficient approach which allows us to derive model predictions (in the form of propagators) from these generating functions, using the Cauchy integral formula. It is important to note that our method does not make any steady-state or long-evolution-time approximations. On the contrary, the perturbative nature of our approach naturally optimises its applicability over relatively short (or intermediate) time scales; see also Remark 3.6. As is argued in Section 1.1, such relatively short evolution times naturally occur in the calculation of quantities such as the log-likelihood, as defined in Equation (1.1). Therefore, our analytical approach is naturally suited to an implementation in a Bayesian parameter inference scheme, such as is outlined in Section 1.1.

As mentioned in Sections 2.2 and 3.3, the introduction of Assumptions 2.2 and 3.7 in our analysis of the systems of PDEs and ODEs that are obtained via the generating function approach is *necessary* for determining explicit expressions for the generating functions themselves. Therefore, we can only be certain of the validity of our approach if we assume that the autoregulation rates are small in comparison with other model parameters, as is done there. Moreover, in the analysis of model B, we have to assume that the protein decay rate is smaller than the decay rate of mRNA; recall Assumption 3.1. That assumption is valid for a large class of (microbial) organisms (Shahrezaei and Swain 2008a; Yu et al. 2006); however, it is by no means generic, as the two decay rates are often comparable in mammalian cells (Schwanhäusser et al. 2011; Vogel and Marcotte 2012). Since the accuracy of approximation of the explicit expressions for the generating functions derived here is quantified in terms of orders of the perturbation parameter(s),

see e.g. Remark 3.6, violation of Assumptions 2.2, 3.1, or 3.7 will decrease the predictive power of the results obtained by the application of the analytical method developed in the article.

The method which is described in Section 1.2, and outlined visually in Figure 5, hence provides a generic framework for the analysis of stochastic gene expression models such as model A (Figure 3 and Section 2.1) and model B (Figure 4 and Section 3.1). Note that, for example, the steady-state and long-evolution-time approximations derived by Shahrezaei and Swain 2008a could be extended to autoregulatory systems via the same approach. However, as is apparent from the (differences between the) analysis presented in Sections 2 and 3, the 'path to an explicit solution' is highly model-dependent. The decision on which analytical techniques to apply, such as the perturbative expansion postulated in (3.22), has to be made on a case-bycase basis. The success of the method presented in the article fully depends on whether the resulting dynamical systems can be solved *explicitly*. To that end, it is highly beneficial that the systems (2.19) and (3.19) obtained here are linear, which is a direct consequence of the fact that all reactions described in the reaction schemes in (2.1) and (2.2), as well as in (3.1) and (3.2), are of first order. Inclusion of second-order reactions would introduce both nonlinear terms and second-order differential operators in the PDE systems for the corresponding generating functions, which would severely increase the complexity of these systems, thus preventing us from obtaining explicit solutions.

The method presented in this article, and the results thus obtained, can be seen, first and foremost, as the natural extension of previous work by Popović, Marr and Swain 2016. Analytical results for the classes of models studied here can be found in several earlier articles. We mention the article by Shahrezaei and Swain 2008a, where a leading-order approximation was obtained in a long-evolution-time and steady-state limit. Bokes et al. 2012a derived analytical expressions for stationary distributions in a model that is equivalent to that considered by Popović, Marr and Swain 2016. Also for that model, a time-scale separation was exploited by Bokes et al. 2012b, in a manner that is similar to the present article, to obtain leading-order analytical expressions on both time scales. The model that is referred to as Model A in Section 2 was analysed in a steady-state setting by Iyer-Biswas and Jayaprakash 2014 via the Poisson integral transform. A similar model was studied by Hornos et al. 2005, were a generating function approach was used; making a steady-state Ansatz, the authors were able to obtain an explicit solution for the generating function in terms of special (Kummer) functions; see also NIST Digital Library of Mathematical Functions. The same model was later solved in a fully time-dependent context by Ramos, Innocentini and Hornos 2011, after a cleverly chosen variable substitution, in terms of another class of special (Heun) functions; cf. again NIST Digital Library of Mathematical Functions.

Other authors have attempted to solve several classes of CMEs directly, i.e. without resorting to generating function techniques or integral transforms. A noteworthy example is the work of Jahnke and Huisinga 2007 on monomolecular systems. Another, more recent example can be found in the work by Iserles and MacNamara 2017, where exact solutions are determined for explicitly time-dependent isomerisation models.

It is important to emphasise that the 'time dependence' referred to in the title of the present article is solely due to the dynamic nature of the underlying stochastic process, and that it hence manifests exclusively through time derivatives in the associated CMEs, such as e.g. in (2.3). In particular, none of the model parameters are time-dependent, as opposed to, for example, the system studied by Iserles and MacNamara 2017. The inclusion of such explicitly time-dependent parameters would be a starting point for incorporating the influence of (extrinsic) noise in the context of the model categories considered in the article.

The availability of analytical expressions for generating functions does, in principle, allow one to try to obtain insight in the underlying processes by studying the explicit form of said expressions, as has been done e.g. by Bokes et al. (Bokes et al. 2012a; Bokes et al. 2012b). However, the complex nature of the processes we analyse here seems to preclude such insights. For example, the integrals over confluent hypergeometric functions, which appear in (2.29), cannot themselves be efficiently expressed in terms of (other) special functions. Still, that complication does not necessarily pose an obstacle to the application we ultimately have in mind, i.e. to Bayesian parameter inference. As the last step in our method – the extraction of propagators from generating functions, see Section 4 – is numerical, the precise functional form of the generating function is not of importance. The mere fact that an explicit expression can be obtained is sufficient for the application of the Cauchy integral formula, where these generating functions enter into the calculation of the appropriate integrals; see again Section 4.

The analytical approach explored in the article does not, of course, represent the only feasible way of obtaining numerical values for propagator probabilities, which can, in turn, serve as input for a Bayesian parameter inference scheme. For an example of a direct numerical method in which the Cauchy integral plays a central role, the reader is referred to the work by Mac-Namara 2015. Our main motivation for pursuing an analytical alternative is reducing the need for potentially lengthy numerical simulations. An efficient implementation of the resulting expressions can result in (significantly) reduced computation times; see, for example, the work by Bornemann 2011. The optimisation of the underlying numerical procedures is, however, beyond the scope of the present article in particular, and of our research programme in general.

The analytical results obtained thus far, as presented in the article, are ready for implementation in a Bayesian parameter inference framework. An analysis of the performance of the resulting approximations to the associated generating functions in the spirit of the article by Feigelman, Popović and Marr 2015, where parameter inference is tested on simulated data based on specific stochastic models, is ongoing work. Moreover, the successful application of our analytical method to specific model categories, such as are represented by model A and model B, suggests several feasible expansions of the 'model library' for which explicit expressions for the corresponding generating functions can be constructed. Thus, stochastic models comprised of multiple proteins represent a natural next stage, bringing the analysis of toggle switch-type models within reach. In addition, one could begin exploring the vast field of gene regulatory networks by considering a simple two-protein system with, for example, activator-inhibitor interaction. Under the assumption of small interaction rates, the resulting PDE system for the associated generating function would be directly amenable to the analytical method described in the article. The analysis of these and similar systems could be a topic for future research.

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A Model A: Explicit Expressions

An explicit expression for (2.27) can be stated as follows:

$$F_{0}^{(1)}(w) = \frac{1}{\kappa_{b}} \left(w \frac{\mathrm{d}}{\mathrm{d}w} F_{0}^{(0)} + \kappa_{f} F_{0}^{(0)} \right)$$

$$= c_{1} \frac{\kappa_{f}}{\kappa_{b}} \left({}_{1}F_{1}(\kappa_{f}, 1 + \kappa_{f} + \kappa_{b}, w) + \frac{w}{1 + \kappa_{f} + \kappa_{b}} {}_{1}F_{1}(1 + \kappa_{f}, 2 + \kappa_{f} + \kappa_{b}, w) \right)$$
(A.1)
$$- c_{2} w^{-\kappa_{f} - \kappa_{b}} \left({}_{1}F_{1}(-\kappa_{b}, 1 - \kappa_{f} - \kappa_{b}, w) + \frac{w}{1 - \kappa_{f} - \kappa_{b}} {}_{1}F_{1}(1 - \kappa_{b}, 2 - \kappa_{f} - \kappa_{b}, w) \right),$$

which can be shown to be equal to (2.35) using §13.3(i) of *NIST Digital Library of Mathematical Functions* when $\hat{c}_1 = c_1 \frac{\kappa_f}{\kappa_b}$ and $\hat{c}_2 = c_2$. Explicit expressions involving the integration limits c_3 and c_4 in (2.29), respectively, are obtained as

$$\int_{c_3}^{\lambda v_0} F_1(-\kappa_b, 1-\kappa_f-\kappa_b, \hat{w}) g(\hat{w}) d\hat{w} = e^{-\lambda v_0} \left\{ \chi_1 (1+v_0)^{n_0} \frac{\lambda v_0 \kappa_b}{1-\kappa_f-\kappa_b} {}_1F_1 (1-\kappa_b, 2-\kappa_f-\kappa_b, \lambda v_0) - {}_1F_1 (-\kappa_b, 1-\kappa_f-\kappa_b, \lambda v_0) \left[\alpha (w+\lambda) \left(\partial_w + \frac{v_0}{w} \partial_{v_0} \right) F_0^{(0)} \right]_{w=\lambda v_0} \right\}$$
(A.2)

and

$$\int_{c_4}^{\lambda v_0} F_1(\kappa_f, 1+\kappa_f+\kappa_b, \hat{w}) \frac{g(\hat{w})}{\hat{w}^{-\kappa_f-\kappa_b}} d\hat{w} = -e^{-\lambda v_0} (\lambda v_0)^{\kappa_f+\kappa_b} \bigg\{ \chi_1 (1+v_0)^{n_0} (\kappa_f+\kappa_b)_1 F_1(\kappa_f, 1+\kappa_f+\kappa_b, \lambda v_0) + \chi_1 (1+v_0)^{n_0} \frac{\lambda v_0 \kappa_f}{1+\kappa_f+\kappa_b} {}_1 F_1 (1+\kappa_f, 2+\kappa_f+\kappa_b, \lambda v_0) + {}_1 F_1(\kappa_f, 1+\kappa_f+\kappa_b, \lambda v_0) \bigg[\alpha (w+\lambda) \bigg(\partial_w + \frac{v_0}{w} \partial_{v_0} \bigg) F_0^{(0)} \bigg]_{w=\lambda v_0} \bigg\}, \quad (A.3)$$

respectively, with $F_0^{(0)}$ as in (2.26). Similarly, explicit expression that involve the integration limits \hat{c}_3 and \hat{c}_4 in (2.38), respectively, read

$$\begin{split} \int_{\hat{c}_{3}}^{\lambda v_{0}} F_{1}(1-\kappa_{b},1-\kappa_{f}-\kappa_{b},\hat{w}) h(\hat{w}) d\hat{w} &= -e^{-\lambda v_{0}} \bigg\{ \chi_{1}(1+v_{0})^{n_{0}} \lambda v_{0} {}_{1}F_{1}(1-\kappa_{b},1-\kappa_{f}-\kappa_{b},\lambda v_{0}) \\ &-\chi_{1}(1+v_{0})^{n_{0}} \frac{\lambda v_{0}(1-\kappa_{b})}{1-\kappa_{f}-\kappa_{b}} {}_{1}F_{1}(2-\kappa_{b},2-\kappa_{f}-\kappa_{b},\lambda v_{0}) \\ &+{}_{1}F_{1}(1-\kappa_{b},1-\kappa_{f}-\kappa_{b},\lambda v_{0}) \bigg[\rho(w+\lambda) \bigg(\partial_{w} + \frac{v_{0}}{w} \partial_{v_{0}} \bigg) F_{0}^{(1)} \bigg]_{w=\lambda v_{0}} \bigg\} \quad (A.4) \end{split}$$

and

$$\begin{split} \int_{\hat{c}_{4}}^{\lambda v_{0}} F_{1}(\kappa_{f}, 1+\kappa_{f}+\kappa_{b}, \hat{w}) \frac{h(\hat{w})}{\hat{w}^{-\kappa_{f}-\kappa_{b}}} \, d\hat{w} &= e^{-\lambda v_{0}} (\lambda v_{0})^{\kappa_{f}+\kappa_{b}} \left\{ \chi_{1}(1+v_{0})^{n_{0}} (\kappa_{f}+\kappa_{b}-\lambda v_{0})_{1} F_{1}(1+\kappa_{f}, 1+\kappa_{f}+\kappa_{b}, \lambda v_{0}) \right. \\ &+ \chi_{1}(1+v_{0})^{n_{0}} \frac{\lambda v_{0}(1+\kappa_{f})}{1+\kappa_{f}+\kappa_{b}} {}_{1}F_{1}(2+\kappa_{f}, 2+\kappa_{f}+\kappa_{b}, \lambda v_{0}) \\ &+ {}_{1}F_{1}(1+\kappa_{f}, 1+\kappa_{f}+\kappa_{b}, \lambda v_{0}) \left[\rho(w+\lambda) \left(\partial_{w} + \frac{v_{0}}{w} \partial_{v_{0}} \right) F_{0}^{(1)} \right]_{w=\lambda v_{0}} \right\}, \quad (A.5) \end{split}$$

respectively, with $F_0^{(1)}$ as in (2.35).

B Model B with Autoregulation: Explicit Expressions

To leading order in δ , i.e. for m = 0, the right-hand side of Equation (3.45) vanishes; hence, the expressions in Equations (3.20) and (3.21) apply. In other words, we can identify $F_{0,n}^{(j)}$ with $F_n^{(j)}$ as given in (3.27) and (3.28), for j = 0, 1 and n = 0, 1, 2. The free constants f_i and g_i are as stated in (3.33), with χ replaced with χ_0 . In the following sections, we present explicit expressions for the first-order correction in δ – i.e. for m = 1 – for all autoregulatory scenarios listed in (3.2).

B.1 mRNA autoactivation

We can solve (3.45) to second order in ε , i.e. for n = 2, applying the initial conditions in (3.55) to obtain

$$F_{1,0}^{(0)} = h_{0,M},\tag{B.1}$$

$$F_{1,1}^{(0)} = h_{1,M} + j_{1,M}\sigma + \alpha_M \chi_0 m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} \frac{e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^2}, \quad \text{and}$$
(B.2)

$$F_{1,2}^{(0)} = h_{2,M} + j_{2,M}\sigma \tag{B.3}$$

$$+ \alpha_{M}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\kappa_{1}\left[\frac{e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}}\left(\frac{1}{1-\mu v_{0}}+\sigma\right)-\frac{\sigma^{2}}{2}(1-\mu v_{0})\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)\right] \\ + \alpha_{M}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\chi_{0}\left\{\frac{e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}}\left[\frac{3\mu v_{0}}{(1-\mu v_{0})^{2}}-\sigma\left(\kappa_{0}+\kappa_{1}-\frac{2\mu v_{0}}{1-\mu v_{0}}\right)+\frac{\sigma^{2}}{2}\mu v_{0}\right] \\ + \sigma^{2}(\kappa_{0}+\kappa_{1})(1-\mu v_{0})\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)\right\} \\ - (1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\chi_{1}\left[\frac{\lambda \kappa_{1}e^{(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{2}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)+\frac{\sigma^{2}}{2}\left(\frac{\lambda \kappa_{1}\mu v_{0}}{1-\mu v_{0}}-(\kappa_{0}+\kappa_{1})^{2}\right)\right], \tag{B.4}$$

which yields

$$F_{1,0}^{(1)} = \frac{1}{\kappa_1} \left[j_{1,M} + \kappa_0 h_{0,M} + \alpha_M m_0 \chi_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \right] \quad \text{and} \tag{B.5}$$

$$F_{1,1}^{(1)} = \frac{j_{2,M} + \kappa_0(h_{1,M} + j_{1,M}\sigma)}{\kappa_1}$$
(B.6)

$$-\frac{\chi_{1}}{\kappa_{1}}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\left[\frac{\lambda\kappa_{1}e^{(1-\mu\nu_{0})\sigma}}{1-\mu\nu_{0}}\left(u_{0}-\frac{\mu\nu_{0}}{1-\mu\nu_{0}}\right)+\sigma\left(\frac{\lambda\kappa_{1}\mu\nu_{0}}{1-\mu\nu_{0}}-(\kappa_{0}+\kappa_{1})^{2}\right)\right]\\-\alpha_{M}\frac{\chi_{0}}{\kappa_{1}}m_{0}(1+u_{0})^{m_{0}-1}(1+\nu_{0})^{n_{0}}\left[\frac{\mu\nu_{0}}{(1-\mu\nu_{0})^{3}}+\frac{\kappa_{1}e^{-(1-\mu\nu_{0})\sigma}}{(1-\mu\nu_{0})^{2}}-\sigma(\kappa_{0}+\kappa_{1})\left(u_{0}-\frac{\mu\nu_{0}}{1-\mu\nu_{0}}\right)\right],$$

where

$$h_{0,M} = \chi_1 (1 + u_0)^{m_0} (1 + v_0)^{n_0},$$
(B.7)

$$h_{1,M} = -\alpha_M \chi_0 m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} \frac{1}{(1-\mu v_0)^2},$$
(B.8)

$$j_{1,M} = -\chi_1(\kappa_0 + \kappa_1)(1 + u_0)^{m_0}(1 + v_0)^{n_0}$$

$$+ \alpha_{M} \chi_0 m_0 (1 + u_0)^{m_0 - 1} (1 + v_0)^{n_0} \left(u_0 - \frac{\mu v_0}{2} \right)$$
(B.9)

$$+ \alpha_{M} \chi_{0} m_{0} (1 + u_{0})^{m_{0}} (1 + v_{0})^{n_{0}} \left(u_{0} - \frac{1}{1 - \mu v_{0}} \right),$$

$$h_{2,M} = \chi_{1} (1 + u_{0})^{m_{0}} (1 + v_{0})^{n_{0}} \frac{\lambda \kappa_{1}}{(1 - \mu v_{0})^{2}} \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right)$$
(B.10)
$$\alpha_{1} m_{0} (1 + u_{0})^{m_{0}-1} (1 + v_{0})^{n_{0}} \frac{1}{(1 - \mu v_{0})^{2}} \left(u_{0} - \frac{\lambda \kappa_{1}}{1 - \mu v_{0}} \right)$$
(B.10)

$$-\alpha_{M}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\frac{1}{(1-\mu v_{0})^{3}}\left[\kappa_{1}+\frac{1}{1-\mu v_{0}}\right], \text{ and}$$

$$j_{2,M} = \chi_{1}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\frac{\lambda \kappa_{1}}{1-\mu v_{0}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)$$

$$+\alpha_{M}\chi_{0}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\frac{1}{(1-\mu v_{0})^{2}}\left(\kappa_{0}+\kappa_{1}+\frac{\mu v_{0}}{1-\mu v_{0}}\right). \tag{B.11}$$

B.2 mRNA autorepression

For the case of mRNA autorepression, we replace $\alpha_M(u+1)\left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} F_{m-1}^{(0)}$ in Equations (3.45) and (3.46) with

$$-\rho_M(u+1)\left(\frac{\partial u}{\partial u_0}\right)^{-1}\partial_{u_0}F_{m-1}^{(1)},\tag{B.12}$$

which yields

$$F_{1,0}^{(0)} = k_{0,M},$$
(B.13)
$$e^{-(1-\mu\nu_0)\sigma}$$

$$F_{1,1}^{(0)} = k_{1,M} + l_{1,M}\sigma - \rho_M(1-\chi_0)m_0(1+u_0)^{m_0-1}(1+v_0)^{n_0}\frac{e^{-(1-\mu_0)\sigma}}{(1-\mu_0)^2}, \quad \text{and}$$
(B.14)
$$F_{1,2}^{(0)} = k_{2,M} + l_{2,M}\sigma$$
(B.15)

$$\begin{aligned} F_{1,2}^{(0)} &= k_{2,M} + l_{2,M}\sigma \end{aligned} \tag{B.15} \\ &- \rho_M m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} \kappa_0 \left[\frac{e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^2} \left(\frac{1}{1-\mu v_0} + \sigma \right) - \frac{\sigma^2}{2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \right] \\ &+ \rho_M m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} (1-\chi_0) \left\{ \frac{\lambda e^{(1-\mu v_0)\sigma}}{(1-\mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)^2 \right. \\ &+ \frac{\lambda e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^3} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) - \frac{(\lambda+3)\mu v_0 e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^4} \\ &- \sigma \frac{e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^2} \left[\frac{(\lambda+2)\mu v_0}{1-\mu v_0} - (\kappa_0+\kappa_1) \right] \\ &- \frac{\sigma^2}{2} \left[\frac{\mu v_0 e^{-(1-\mu v_0)\sigma}}{(1-\mu v_0)^2} + \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \left(2(\kappa_0+\kappa_1) - \frac{\lambda \mu v_0}{1-\mu v_0} \right) \right] \right\} \\ &- (1+u_0)^{m_0} (1+v_0)^{n_0} \chi_1 \left[\frac{\lambda \kappa_1 e^{(1-\mu v_0)\sigma}}{(1-\mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) + \frac{\sigma^2}{2} \left(\frac{\lambda \kappa_1 \mu v_0}{1-\mu v_0} - (\kappa_0+\kappa_1)^2 \right) \right] \\ &+ \rho_M (1+u_0)^{m_0} (1+v_0)^{n_0} (1-\chi_0) \frac{\lambda}{(1-\mu v_0)^2} \left[e^{(1-\mu v_0)\sigma} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) + \frac{e^{-(1-\mu v_0)\sigma}}{1-\mu v_0} \right]. \end{aligned}$$

Hence, it follows that

$$\begin{aligned} F_{1,0}^{(1)} &= \frac{1}{\kappa_1} \left[l_{1_M} + \kappa_0 k_{0,M} - \rho_M m_0 (1 - \chi_0) (1 + u_0)^{m_0 - 1} (1 + v_0)^{n_0} \left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right) \right] & \text{and} \quad (B.16) \\ F_{1,1}^{(1)} &= \frac{l_{2,M} + \kappa_0 (k_{1,M} + l_{1,M} \sigma)}{\kappa_1} \\ &\quad - \frac{\chi_1}{\kappa_1} (1 + u_0)^{m_0} (1 + v_0)^{n_0} \left[\frac{\lambda \kappa_1 e^{(1 - \mu v_0)\sigma}}{1 - \mu v_0} \left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right) + \sigma \left(\frac{\lambda \kappa_1 \mu v_0}{1 - \mu v_0} - (\kappa_0 + \kappa_1)^2 \right) \right] \\ &\quad + \rho_M (1 + u_0)^{m_0} (1 + v_0)^{n_0} (1 - \chi_0) \frac{\lambda}{\kappa_1 (1 - \mu v_0)} \left[\left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right) - \frac{1}{1 - \mu v_0} \right] \\ &\quad + \rho_M (1 - \chi_0) m_0 (1 + u_0)^{m_0 - 1} (1 + v_0)^{n_0} \left[\frac{e^{-(1 - \mu v_0)\sigma}}{(1 - \mu v_0)^2} - \frac{\lambda}{\kappa_1 (1 - \mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right) \right] \\ &\quad + \frac{\mu v_0}{\kappa_1 (1 - \mu v_0)^3} + \frac{\lambda}{\kappa_1 (1 - \mu v_0)} \left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right)^2 - \frac{\sigma}{\kappa_1} (\kappa_0 + \kappa_1) \left(u_0 - \frac{\mu v_0}{1 - \mu v_0} \right) \right], \end{aligned}$$

where

$$k_{0,M} = \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0}, \tag{B.18}$$

$$k_{1,M} = \rho_M (1 - \chi_0) m_0 (1 + u_0)^{m_0 - 1} (1 + v_0)^{n_0} \frac{1}{(1 - \mu v_0)^2},$$
(B.19)

$$l_{1,M} = -\chi_1(\kappa_0 + \kappa_1)(1 + u_0)^{m_0}(1 + v_0)^{n_0}$$
(B.20)

$$-\rho_{M}(1-\chi_{0})m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right),$$

$$k_{2,M} = \chi_{1}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\frac{\lambda \kappa_{1}}{(1-\mu v_{0})^{2}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)$$

$$-\rho_{M}(1+u_{0})^{m_{0}+1}(1+v_{0})^{n_{0}}(1-\chi_{0})\frac{\lambda}{(1-\mu v_{0})^{2}}$$

$$+\rho_{M}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}m_{0}\frac{\kappa_{0}}{(1-\mu v_{0})^{3}}$$

$$+\rho_{M}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}(1-\chi_{0})\frac{1}{(1-\mu v_{0})^{2}}\left[\frac{(\lambda+3)\mu v_{0}}{(1-\mu v_{0})^{2}}+\frac{\lambda}{1-\mu v_{0}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)\right]$$

$$-\lambda\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)^{2}\right], \text{ and}$$

$$l_{2,M} = \chi_{1}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\frac{\lambda \kappa_{1}}{1-\mu v_{0}}\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)$$

$$(B.22)$$

$$\begin{split} l_{2,M} &= \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda \kappa_1}{1-\mu v_0} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \\ &- \rho_M (1+u_0)^{m_0} (1+v_0)^{n_0} (1-\chi_0) \frac{\lambda}{1-\mu v_0} \left[\left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) - \frac{1}{1-\mu v_0} \right] \\ &- \rho_M (1+u_0)^{m_0-1} (1+v_0)^{n_0} m_0 (1-\chi_0) \frac{1}{1-\mu v_0} \left[\frac{\kappa_0 + \kappa_1 - 1}{1-\mu v_0} + \frac{1}{(1-\mu v_0)^2} \right] \\ &- \frac{\lambda}{1-\mu v_0} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) + \lambda \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)^2 \right]. \end{split}$$
(B.22)

B.3 Protein autoactivation

For the case of protein autoactivation, we replace $\alpha_M(u+1)\left(\frac{\partial u}{\partial u_0}\right)^{-1}\partial_{u_0}F_{m-1}^{(0)}$ in Equations (3.45) and (3.46) with

$$\alpha_P(\nu+1) \left(\frac{\partial \nu}{\partial \nu_0}\right)^{-1} \left[-\left(\frac{\partial u}{\partial \nu_0}\right) \left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} + \partial_{\nu_0} \right] F_{m-1}^{(0)}, \tag{B.23}$$

which yields

$$\begin{split} F_{1,0}^{(0)} &= h_{0,P}, & (B.24) \\ F_{1,1}^{(0)} &= h_{1,P} + j_{1,P}\sigma - \alpha_{P\chi_{0}}m_{0}\mu(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1} \bigg[\frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{3}} + \frac{\sigma^{2}}{2} \bigg(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \bigg) \bigg], & \text{and} \\ & (B.25) \\ F_{1,2}^{(0)} &= h_{2,P} + j_{2,P}\sigma & (B.26) \\ &+ \alpha_{P}n_{0}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\mu \frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{4}} \bigg[\kappa_{1}(1+v_{0}) + \chi_{0} \bigg] \\ &- \alpha_{P}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\mu \frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{4}} \bigg[\kappa_{1}(1+v_{0}) + \chi_{0} \frac{1+2v_{0}+\mu v_{0}(5+4v_{0})}{1-\mu v_{0}} \bigg] \\ &+ \alpha_{P}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\sigma\mu \frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{3}} \bigg\{ -\chi_{0} \frac{(1+(3+2v_{0})\mu]v_{0}}{1-\mu v_{0}} + [(1-\chi_{0})\kappa_{1}-\chi_{0}\kappa_{0}](1+v_{0})\bigg\} \\ &+ \alpha_{P}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\mu \frac{\sigma^{2}}{2} \frac{1}{(1-\mu v_{0})^{2}} \bigg\{ -\chi_{0}(1+v_{0})\frac{\mu v_{0}e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}} + (1+v_{0})[2(\kappa_{0}+\kappa_{1})\chi_{0}-\kappa_{1}] + \chi_{0}\frac{1+\mu v_{0}^{2}}{1-\mu v_{0}}\bigg\} \\ &+ \alpha_{P}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}}\mu \frac{\sigma^{3}}{6} \bigg\{ u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \bigg\} ((1+v_{0})[3\chi_{0}(\kappa_{0}+\kappa_{1}) - 2\kappa_{1}] + \chi_{0}(1-v_{0})\} \\ &- (1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\chi_{1} \bigg\{ \frac{\lambda \kappa_{1}e^{(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{2}} \bigg\{ u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \bigg\} + \frac{\sigma^{2}}{2} \bigg[\frac{\lambda \kappa_{1}\mu v_{0}}{1-\mu v_{0}} - (\kappa_{0}+\kappa_{1})^{2} \bigg] \bigg\}. \end{split}$$

Hence, it follows that

where

$$h_{0,P} = \chi_1 (1 + u_0)^{m_0} (1 + v_0)^{n_0}, \tag{B.30}$$

$$h_{1,P} = \alpha_P \chi_0 m_0 \mu (1+u_0)^{m_0-1} (1+v_0)^{n_0+1} \frac{1}{(1-\mu v_0)^3},$$
(B.31)

$$j_{1,P} = -\chi_1(\kappa_0 + \kappa_1)(1 + u_0)^{m_0}(1 + v_0)^{n_0}$$
(B.32)

$$-\alpha_{P}\chi_{0}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\left[n_{0}+\mu m_{0}\frac{1+v_{0}}{1+u_{0}}\frac{1}{(1-\mu v_{0})^{2}}\right],$$

$$P=\chi_{1}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}}\frac{\lambda\kappa_{1}}{(1-\mu v_{0})^{2}}\left(u_{0}-\frac{\mu v_{0}}{1+\mu v_{0}}\right)$$
(B.33)

$$h_{2,P} = \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda \kappa_1}{(1-\mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)$$
(B.33)
+ $\alpha_P m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} - \frac{\mu}{(1+v_0)\kappa_1 + v_0} \left[(1+v_0)\kappa_1 + v_0 \frac{1+2v_0 + \mu v_0(5+4v_0)}{1-\mu v_0} \right],$ and

$$j_{2,P} = \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda \kappa_1}{1-\mu v_0} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)$$

$$+ \alpha_P \chi_0 \mu m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0+1} \frac{1}{(1-\mu v_0)^3} \left[2 - (\kappa_0 + \kappa_1) - \frac{3}{1-\mu v_0} \right].$$
(B.34)

B.4 Protein autorepression

For the case of protein autorepression, we replace $\alpha_M(u+1)\left(\frac{\partial u}{\partial u_0}\right)^{-1} \partial_{u_0} F_{m-1}^{(0)}$ in Equations (3.45) and (3.46) with

$$-\rho_P(\nu+1)\left(\frac{\partial\nu}{\partial\nu_0}\right)^{-1} \left[-\left(\frac{\partial u}{\partial\nu_0}\right)\left(\frac{\partial u}{\partial u_0}\right)^{-1}\partial_{u_0} + \partial_{\nu_0}\right] F_{m-1}^{(1)}, \qquad (B.35)$$

which yields

$$\begin{split} F_{1,0}^{(0)} &= k_{0,P}, & (B.36) \\ F_{1,1}^{(0)} &= k_{1,P} + l_{1,P}\sigma + \rho_{P}(1-\chi_{0})m_{0}\mu(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1} \left[\frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{3}} + \frac{\sigma^{2}}{2} \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) \right] \right], & \text{and} \\ & (B.37) \\ F_{1,2}^{(0)} &= k_{2,P} + l_{2,P}\sigma & (B.38) \\ &+ \rho_{P}(1-\chi_{0})\lambda\mu(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}+1} \frac{1}{(1-\mu v_{0})^{3}} \left[e^{(1-\mu v_{0})\sigma} \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) - \frac{e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}} \right] \\ &+ \rho_{P}(1-\chi_{0})\lambda_{0}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}+1} \frac{1}{(1-\mu v_{0})^{3}} \left[e^{(1-\mu v_{0})\sigma} \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) - \frac{e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}} \right] \\ &+ \rho_{P}(1-\chi_{0})\lambda_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1} \frac{\lambda \mu e^{-(1-\mu v_{0})\sigma}}{1-\mu v_{0}} \left[u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right] \left[\frac{1}{1-\mu v_{0}} - \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) \right] \\ &+ \rho_{P}m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1} \frac{\lambda e^{(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{3}} \left[u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right] \left[\frac{1}{1-\mu v_{0}} - \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) \right] \\ &+ \rho_{P}\sigma \lambda \mu m_{0}(1-\chi_{0})(1+u_{0})^{m_{0}-1}(1+v_{0})m_{0} + 1 \frac{e^{(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{2}} \left[u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right]^{2} \\ &+ \rho_{P}\sigma \mu m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})m_{0} \frac{e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{2}} \left[x_{0}(1+v_{0})(\kappa_{0}+\kappa_{1}) - (1+v_{0})\kappa_{1} \right] \\ &+ (1-\chi_{0})\frac{v_{0}}{+\mu v_{0}(3+2v_{0}) + \mu v_{0}\lambda(1+v_{0})}{1-\mu v_{0}} \right] \\ &+ \rho_{P}\frac{\sigma^{2}}{2}(1+u_{0})^{m_{0}}(1+v_{0})^{n_{0}+1}(1-\chi_{0})\frac{\lambda \mu}{1-\mu v_{0}}} \left[\frac{1}{1-\mu v_{0}} - \left(u_{0} - \frac{\mu v_{0}}{1-\mu v_{0}} \right) \right] \\ &+ \rho_{P}\frac{\sigma^{2}}{2}\mu m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1}(1-\chi_{0})\frac{\mu u_{0}e^{-(1-\mu v_{0})\sigma}}{(1-\mu v_{0})^{3}} \\ &+ \rho_{P}\frac{\sigma^{2}}{2}\mu m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}} \left\{ \left[-(1-2\chi_{0})(\kappa_{0}+\kappa_{1}) - \kappa_{1} \right] \frac{1+v_{0}}{(1-\mu v_{0})^{3}} \\ &+ \rho_{P}\frac{\sigma^{2}}{2}\mu m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}} \left\{ \left[-(1-2\chi_{0})(\kappa_{0}+\kappa_{1}) - \kappa_{1} \right] \frac{1+v_{0}}{(1-\mu v_{0})^{2}} \\ &- (1-\chi_{0})\frac{1+\mu v_{0}^{2}}{(1-\mu v_{0})^{3}} + (1-\chi_{0})(1+v_{0}) \left[\frac{\mu v_{0}}{(1-\mu v_{0})^{3}} \\ &+ \rho_{P}\frac$$

It follows that

$$\begin{aligned} F_{1,0}^{(1)} &= \frac{1}{\kappa_{1}} \left(l_{1,P} + \kappa_{0} k_{0,P} \right) & (B.39) \\ &- \rho_{P} m_{0} (1 - \chi_{0}) (1 + u_{0})^{m_{0}-1} (1 + v_{0})^{n_{0}} \frac{1}{\kappa_{1}} \left\{ \frac{m_{0} \mu (1 + v_{0})}{(1 - \mu v_{0})^{2}} + n_{0} \left[\frac{1}{1 - \mu v_{0}} + \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) \right] \right\} & \text{and} \\ F_{1,1}^{(1)} &= \frac{l_{2,M} + \kappa_{0} (k_{1,M} + l_{1,M} \sigma)}{\kappa_{1}} & (B.40) \\ &+ \rho_{P} (1 - \chi_{0}) (1 + u_{0})^{m_{0}+1} (1 + v_{0})^{n_{0}+1} \frac{\lambda \mu}{\kappa_{1}} \frac{1}{(1 - \mu v_{0})^{2}} \\ &+ \rho_{P} n_{0} (1 - \chi_{0}) (1 + u_{0})^{m_{0}} (1 + v_{0})^{n_{0}} \frac{1}{\kappa_{1}} \left[\frac{\lambda}{1 - \mu v_{0}} \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) - (\kappa_{0} + \kappa_{1}) \sigma \right] \\ &+ \rho_{P} \mu m_{0} (1 - \chi_{0}) (1 + u_{0})^{m_{0}-1} (1 + v_{0})^{n_{0}+1} \frac{1}{\kappa_{1}} \left\{ - \frac{3}{(1 - \mu v_{0})^{4}} + \frac{1}{(1 - \mu v_{0})^{3}} \left[-\kappa_{1} e^{-(1 - \mu v_{0})\sigma} + 2 + 2\lambda \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) \right] \\ &- \sigma \frac{\kappa_{0} + \kappa_{1}}{(1 - \mu v_{0})^{2}} - \frac{\sigma^{2}}{2} \kappa_{1} \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) \right\} \\ &- \frac{\chi_{1}}{\kappa_{1}} (1 + u_{0})^{m_{0}} (1 + v_{0})^{n_{0}} \left[\frac{\lambda \kappa_{1} e^{(1 - \mu v_{0})\sigma}}{1 - \mu v_{0}} \left(u_{0} - \frac{\mu v_{0}}{1 - \mu v_{0}} \right) + \sigma \left(\frac{\lambda \kappa_{1} \mu v_{0}}{1 - \mu v_{0}} - (\kappa_{0} + \kappa_{1})^{2} \right) \right], \end{aligned}$$

where

$$k_{0,P} = \chi_1 (1 + u_0)^{m_0} (1 + v_0)^{n_0}, \tag{B.41}$$

$$k_{1,P} = -\rho_P (1 - \chi_0) m_0 (1 + u_0)^{m_0 - 1} (1 + v_0)^{n_0 + 1} \frac{\mu}{(1 - \mu v_0)^3},$$
(B.42)

$$l_{1,P} = -\chi_1(\kappa_0 + \kappa_1)(1 + u_0)^{m_0}(1 + v_0)^{n_0}$$

$$-\rho_P(1 - \chi_0)(1 + u_0)^{m_0}(1 + v_0)^{n_0} \left[n_0 + \mu m_0 \frac{1 + v_0}{1 + u_0} \frac{1}{(1 - \mu v_0)^2} \right],$$
(B.43)

$$k_{2,P} = \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda \kappa_1}{(1-\mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)$$
(B.44)
+ $\rho_P (1-\chi_0) (1+u_0)^{m_0} (1+v_0)^{n_0+1} \frac{\lambda \mu}{(1-\mu v_0)^3} \left[\frac{1}{1-\mu v_0} - \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \right]$
- $\rho_P (1-\chi_0) n_0 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda}{(1-\mu v_0)^2} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right) \right]$
+ $\rho_P (1-\chi_0) m_0 \mu (1+u_0)^{m_0-1} (1+v_0)^{n_0+1} \frac{\lambda}{(1-\mu v_0)^3} \left[\left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)^2 - \frac{\mu v_0}{(1-\mu v_0)^2} \right]$
- $\rho_P m_0 (1+u_0)^{m_0-1} (1+v_0)^{n_0} \frac{\mu}{(1-\mu v_0)^4} \left\{ \frac{1-\chi_0}{1-\mu v_0} [1+2v_0+\mu v_0(5+4v_0)] + (1+v_0)\kappa_0 \right\}, \text{ and}$
 $l_{2,P} = \chi_1 (1+u_0)^{m_0} (1+v_0)^{n_0} \frac{\lambda \kappa_1}{1-\mu v_0} \left(u_0 - \frac{\mu v_0}{1-\mu v_0} \right)$ (B.45)
 $- \rho_P (1+u_0)^{m_0+1} (1+v_0)^{n_0+1} (1-\chi_0) \frac{\lambda \mu}{(1-\mu v_0)^2}$

$$+\rho_{P}\mu m_{0}(1+u_{0})^{m_{0}-1}(1+v_{0})^{n_{0}+1}(1-\chi_{0})\frac{1}{(1-\mu v_{0})^{3}}\left[-2-2\lambda\left(u_{0}-\frac{\mu v_{0}}{1-\mu v_{0}}\right)+\frac{3}{1-\mu v_{0}}+\kappa_{0}+\kappa_{1}\right].$$