## **SBML Model Report**

# Model name: "Tyson1999\_CircClock"



May 6, 2016

#### 1 General Overview

This is a document in SBML Level 2 Version 1 format. This model was created by Nicolas Le Novre<sup>1</sup> at July seventh 2005 at 3:19 p.m. and last time modified at February 25<sup>th</sup> 2015 at 1:24 p.m. Table 1 gives an overview of the quantities of all components of this model.

Table 1: Number of components in this model, which are described in the following sections.

Element	Quantity	Element	Quantity
compartment types	0	compartments	2
species types	0	species	3
events	0	constraints	0
reactions	5	function definitions	0
global parameters	1	unit definitions	1
rules	0	initial assignments	0

#### **Model Notes**

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To cite BioModels Database, please use: Li C, Donizelli M, Rodriguez N, Dharuri H, Endler L, Chelliah V, Li L, He E, Henry A, Stefan MI, Snoep JL, Hucka M, Le Novre N, Laibe C (2010) BioModels Database: An enhanced, curated and annotated resource for published quantitative kinetic models. BMC Syst Biol., 4:92.

#### 2 Unit Definitions

This is an overview of five unit definitions of which four are predefined by SBML and not mentioned in the model.

#### 2.1 Unit time

Name hour

**Definition** 3600 s

#### 2.2 Unit substance

**Notes** Mole is the predefined SBML unit for substance.

**Definition** mol

#### 2.3 Unit volume

**Notes** Litre is the predefined SBML unit for volume.

**Definition** 1

#### 2.4 Unit area

**Notes** Square metre is the predefined SBML unit for area since SBML Level 2 Version 1.

**Definition** m<sup>2</sup>

### 2.5 Unit length

**Notes** Metre is the predefined SBML unit for length since SBML Level 2 Version 1.

**Definition** m

## 3 Compartments

This model contains two compartments.

Table 2: Properties of all compartments.

			1	I			
Id	Name	SBO	Spatial Dimensions	Size	Unit	Constant	Outside
			Difficilisions				
default			3	1	litre		
CYTOPLASM			3	1	litre		

## 3.1 Compartment default

This is a three dimensional compartment with a constant size of one litre.

## 3.2 Compartment CYTOPLASM

This is a three dimensional compartment with a constant size of one litre.

# 4 Species

This model contains three species. The boundary condition of one of these species is set to true so that this species' amount cannot be changed by any reaction. Section 7 provides further details and the derived rates of change of each species.

Table 3: Properties of each species.

		P			
Id	Name	Compartment	Derived Unit	Constant	Boundary Condi- tion
EmptySet M P	Pt	default CYTOPLASM CYTOPLASM	$\begin{array}{c} \operatorname{mol} \cdot l^{-1} \\ \operatorname{mol} \cdot l^{-1} \\ \operatorname{mol} \cdot l^{-1} \end{array}$	<b>⊿</b> ⊟ ⊟	<b>☑</b> ⊟

## **5 Parameter**

This model contains one global parameter.

Table 4: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
N_A		6	$0.0221367 \cdot 10^{2}$	23	$ \mathbf{Z} $

## 6 Reactions

This model contains five reactions. All reactions are listed in the following table and are subsequently described in detail. If a reaction is affected by a modifier, the identifier of this species is written above the reaction arrow.

Table 5: Overview of all reactions

Nº	Id	Name	Reaction Equation	SBO
1	TC	transcription	$\text{EmptySet} \xrightarrow{P} M$	
2	TL	translation	EmptySet $\xrightarrow{\mathbf{M}}$ P	
3	mRNAD	mRNA degradation	$M \longrightarrow EmptySet$	
4	ProteinD	protein degradation	$P \longrightarrow EmptySet$	
5	DBT	Per phosphorylation and degradation	$P \longrightarrow EmptySet$	

#### 6.1 Reaction TC

This is an irreversible reaction of one reactant forming one product influenced by one modifier.

Name transcription

## **Reaction equation**

$$EmptySet \xrightarrow{P} M \tag{1}$$

#### Reactant

Table 6: Properties of each reactant.

Id	Name	SBO
EmptySet		

#### **Modifier**

Table 7: Properties of each modifier.

Id	Name	SBO
P	Pt	

#### **Product**

Table 8: Properties of each product.

Id	Name	SBO
М		

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{1} = \text{vol}\left(\text{CYTOPLASM}\right) \cdot \frac{\text{Vm}}{1 + \left(\frac{\left[P\right] \cdot \left(1 - \frac{2}{1 + \left(1 + 8 \cdot \text{Keq} \cdot \left[P\right]\right)^{0.5}}\right)}{2 \cdot \text{Pcrit}}\right)^{2}}\right)}$$
(2)

Table 9: Properties of each parameter.

		1 1	
Id	Name	SBO Value Unit	Constant
Vm		1.0	$\square$
Pcrit		0.1	
Keq		200.0	

## 6.2 Reaction TL

This is an irreversible reaction of one reactant forming one product influenced by one modifier.

Name translation

## **Reaction equation**

$$EmptySet \xrightarrow{M} P$$
 (3)

#### Reactant

Table 10: Properties of each reactant.

Id	Name	SBO
EmptySet		

#### **Modifier**

Table 11: Properties of each modifier.

Id	Name	SBO
М		

## **Product**

Table 12: Properties of each product.

Id	Name	SBO
Р	Pt	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_2 = V \cdot [M] \cdot vol(CYTOPLASM) \tag{4}$$

Table 13: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
V	Vp	0.5	$\overline{Z}$

## 6.3 Reaction mRNAD

This is an irreversible reaction of one reactant forming one product.

Name mRNA degradation

## **Reaction equation**

$$M \longrightarrow EmptySet$$
 (5)

#### Reactant

Table 14: Properties of each reactant.

Id	Name	SBO
М		

#### **Product**

Table 15: Properties of each product.

Id	Name	SBO
EmptySet		

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_3 = D \cdot [M] \cdot vol(CYTOPLASM)$$
 (6)

Table 16: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
D	km	0.1	

#### **6.4 Reaction ProteinD**

This is an irreversible reaction of one reactant forming one product.

Name protein degradation

## **Reaction equation**

$$P \longrightarrow EmptySet$$
 (7)

#### Reactant

Table 17: Properties of each reactant.

Id	Name	SBO
Р	Pt	

#### **Product**

Table 18: Properties of each product.

Id	Name	SBO
EmptySet		

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_4 = D \cdot [P] \cdot vol(CYTOPLASM) \tag{8}$$

Table 19: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
D	kp3	0.1	

#### 6.5 Reaction DBT

This is an irreversible reaction of one reactant forming one product.

Name Per phosphorylation and degradation

#### **Reaction equation**

$$P \longrightarrow EmptySet$$
 (9)

#### Reactant

Table 20: Properties of each reactant.

Id	Name	SBO
Р	Pt	

#### **Product**

Table 21: Properties of each product.

Id	Name	SBO
EmptySet		

#### **Kinetic Law**

Derived unit contains undeclared units

$$v_5 = \text{vol}\left(\text{CYTOPLASM}\right) \cdot \frac{\text{k1} \cdot [\text{P}] \cdot \frac{2}{1 + (1 + 8 \cdot \text{Keq} \cdot [\text{P}])^{0.5}} + \text{k2} \cdot [\text{P}]}{\text{J} + [\text{P}]}$$
 (10)

Table 22: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
k1	kp1	10.00	$\overline{Z}$
k2	kp2	0.03	$\square$
J	Jp	0.05	
Keq		200.00	$\square$

# 7 Derived Rate Equations

When interpreted as an ordinary differential equation framework, this model implies the following set of equations for the rates of change of each species.

Identifiers for kinetic laws highlighted in gray cannot be verified to evaluate to units of SBML substance per time. As a result, some SBML interpreters may not be able to verify the consistency of the units on quantities in the model. Please check if

- parameters without an unit definition are involved or
- volume correction is necessary because the hasOnlySubstanceUnits flag may be set to false and spacialDimensions > 0 for certain species.

### 7.1 Species EmptySet

#### Initial amount 0 mol

This species takes part in five reactions (as a reactant in TC, TL and as a product in mRNAD, ProteinD, DBT), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{EmptySet} = 0\tag{11}$$

### 7.2 Species M

#### Initial amount 1 mol

This species takes part in three reactions (as a reactant in mRNAD and as a product in TC and as a modifier in TL).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{M} = |v_1| - |v_3| \tag{12}$$

#### 7.3 Species P

#### Name Pt

## Initial amount 1 mol

This species takes part in four reactions (as a reactant in ProteinD, DBT and as a product in TL and as a modifier in TC).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{P} = |v_2| - |v_4| - |v_5| \tag{13}$$

SBML2LATEX was developed by Andreas Dräger<sup>a</sup>, Hannes Planatscher<sup>a</sup>, Dieudonné M Wouamba<sup>a</sup>, Adrian Schröder<sup>a</sup>, Michael Hucka<sup>b</sup>, Lukas Endler<sup>c</sup>, Martin Golebiewski<sup>d</sup> and Andreas Zell<sup>a</sup>. Please see http://www.ra.cs.uni-tuebingen.de/software/SBML2LaTeX for more information.

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