# SBML Model Report

# Model identifier: "MPhase"



February 28, 2009

## 1 General Overview

This is a document in SBML Level 2 Version 1 format. Table 1 gives an overview of the quantities of all components of this model.

Table 1: The SBML components in this model. All components are described in more detail in the following sections.

8			
Element	Quantity	Element	Quantity
compartment types	0	compartments	2
species types	0	species	20
events	0	constraints	0
reactions	11	function definitions	0
global parameters	0	unit definitions	0
rules	0	initial assignments	0

## 2 Unit Definitions

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

#### 2.1 Unit substance

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

**Definition** mol

## 2.2 Unit volume

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

**Definition** 1

#### 2.3 Unit area

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

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

## 2.4 Unit length

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

**Definition** m

#### 2.5 Unit time

**Notes** Second is the predefined SBML unit for time.

**Definition** s

## 3 Compartments

This model contains two compartments.

Table 2: Properties of all compartments.

			*				
Id	Name	SBO	Spatial	Size	Unit	Constant	Outside
			Dimensions				
default			3	1	litre	<b></b>	
c1	cell		3	1	litre	$\overline{\mathbb{Z}}$	default

## 3.1 Compartment default

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

## 3.2 Compartment c1

This is a three-dimensional compartment with a constant size of one litre that is surrounded by default.

Name cell

# 4 Species

This model contains 20 species. Section 6 provides further details and the derived rates of change of each species.

Table 3: Properties of each species.

Id	Name	Compartment	Derived Unit	Constant	Boundary Condi-
- 2	ContinD		$\operatorname{mol} \cdot \operatorname{l}^{-1}$		tion
s3	CyclinB	c1	$mol \cdot l^{-1}$		
s5	PP2A	c1			
s6	Kinase X	c1	$\text{mol} \cdot 1^{-1}$		
s7	CAK	c1	$\text{mol} \cdot 1^{-1}$	$\Box$	
s10	Nim1	c1	$\operatorname{mol} \cdot 1^{-1}$		$\Box$
s11	Lamin	c1	$\text{mol} \cdot l^{-1}$		
s12	M-Phase	c1	$\text{mol} \cdot l^{-1}$	$\Box$	$\Box$
s21	Lamin	c1	$\mathrm{mol}\cdot\mathrm{l}^{-1}$		
s4	Cdc25	c1	$\mathrm{mol}\cdot\mathrm{l}^{-1}$		$\Box$
s8	Mik1	c1	$\operatorname{mol} \cdot 1^{-1}$		$\Box$
s9	Wee1	c1	$\operatorname{mol} \cdot \operatorname{l}^{-1}$		$\Box$
s22	Cdc25	c1	$\operatorname{mol} \cdot 1^{-1}$		$\Box$
s24	Wee1	c1	$\operatorname{mol} \cdot 1^{-1}$		
s25	Mik1	c1	$\mathrm{mol}\cdot\mathrm{l}^{-1}$		
s26	a33_degraded	c1	$\mathrm{mol}\cdot\mathrm{l}^{-1}$		$\Box$
s2	Cdc2	c1	$\mathrm{mol}\cdot\mathrm{l}^{-1}$		$\Box$
s27	Complex(CyclinB,Cdc2)	c1	$\text{mol} \cdot l^{-1}$		$\Box$
s28	Complex(CyclinB,Cdc2)	c1	$\operatorname{mol} \cdot 1^{-1}$		$\Box$
s29	Complex(CyclinB,Cdc2)	c1	$\operatorname{mol} \cdot 1^{-1}$		$\Box$
<b>s</b> 30	Complex(CyclinB,Cdc2)	c1	$\text{mol} \cdot 1^{-1}$		

# **5 Reactions**

This model contains eleven reactions. All reactions are listed in the following table and are subsequently described in detail. If a reaction is affected by one or more modifiers, the identifiers of the modifier species are written above the reaction arrow.

Table 4: Overview of all reactions

No	Id	Name	Reaction Equation	SBO
1	r1		$s21 \xrightarrow{s29} s11$	
2	r2		$\mathtt{s11} \longrightarrow \mathtt{s12}$	
3	<b>r</b> 7		$s4 \xrightarrow{s5} s22$	
4	r8		$\mathtt{s22} \xrightarrow{\mathtt{s6}} \mathtt{s4}$	
5	r11		$s25 \xrightarrow{s10} s8$	
6	r12		$s24 \xrightarrow{s10} s9$	
7	r13		$s27 \xrightarrow{s7} s28$	
8	r14		$s28 \xrightarrow{s22} s29$	
9	r15		$s30 \xrightarrow{s24, s25} s27$	
10	r17		$s29 \longrightarrow s2 + s26$	
11	r18		$s2 + s3 \longrightarrow s30$	

#### **5.1 Reaction** r1

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

## **Reaction equation**

$$s21 \xrightarrow{s29} s11 \tag{1}$$

#### Reactant

Table 5: Properties of each reactant.

Id	Name	SBO
s21	Lamin	

#### **Modifier**

Table 6: Properties of each modifier.

Id	Name	SBO
s29	Complex(CyclinB,Cdc2)	

#### **Product**

Table 7: Properties of each product.

Id	Name	SBO
s11	Lamin	

## **Kinetic Law**

$$v_1 = \text{not specified}$$
 (2)

## 5.2 Reaction r2

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

## **Reaction equation**

$$s11 \longrightarrow s12$$
 (3)

## Reactant

Table 8: Properties of each reactant.

Id	Name	SBO
s11	Lamin	

## **Product**

Table 9: Properties of each product.

Id	Name	SBO
s12	M-Phase	

## **Kinetic Law**

$$v_2 = \text{not specified}$$
 (4)

## **5.3 Reaction** r7

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

## **Reaction equation**

$$s4 \xrightarrow{s5} s22$$
 (5)

#### Reactant

Table 10: Properties of each reactant.

Id	Name	SBO
s4	Cdc25	

#### **Modifier**

Table 11: Properties of each modifier.

Id	Name	SBO
s5	PP2A	

#### **Product**

Table 12: Properties of each product.

Id	Name	SBO
s22	Cdc25	

#### **Kinetic Law**

$$v_3 = \text{not specified}$$
 (6)

## 5.4 Reaction r8

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

## **Reaction equation**

$$s22 \xrightarrow{s6} s4$$
 (7)

#### Reactant

Table 13: Properties of each reactant.

Id	Name	SBO
s22	Cdc25	

## **Modifier**

Table 14: Properties of each modifier.

Id	Name	SBO
s6	Kinase X	

#### **Product**

Table 15: Properties of each product.

Id	Name	SBO
s4	Cdc25	

#### **Kinetic Law**

$$v_4 = \text{not specified}$$
 (8)

#### 5.5 Reaction r11

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

## **Reaction equation**

$$s25 \xrightarrow{s10} s8 \tag{9}$$

#### Reactant

Table 16: Properties of each reactant.

Id	Name	SBO
s25	Mik1	

#### **Modifier**

Table 17: Properties of each modifier.

Id	Name	SBO
s10	Nim1	

#### **Product**

Table 18: Properties of each product.

Id	Name	SBO
s8	Mik1	

#### **Kinetic Law**

$$v_5 = \text{not specified}$$
 (10)

#### 5.6 Reaction r12

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

## **Reaction equation**

$$s24 \xrightarrow{s10} s9 \tag{11}$$

#### Reactant

Table 19: Properties of each reactant.

Id	Name	SBO
s24	Wee 1	

#### **Modifier**

Table 20: Properties of each modifier.

Id	Name	SBO
s10	Nim1	

#### **Product**

Table 21: Properties of each product.

Id	Name	SBO
s9	Wee1	

## **Kinetic Law**

$$v_6 = \text{not specified}$$
 (12)

## 5.7 Reaction r13

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

## **Reaction equation**

$$s27 \xrightarrow{s7} s28 \tag{13}$$

#### Reactant

Table 22: Properties of each reactant.

Id	Name	SBO
s27	Complex(CyclinB,Cdc2)	

#### **Modifier**

Table 23: Properties of each modifier.

Id	Name	SBO
s7	CAK	

#### **Product**

Table 24: Properties of each product.

Id	Name	SBO
s28	Complex(CyclinB,Cdc2)	

#### **Kinetic Law**

$$v_7 = \text{not specified}$$
 (14)

#### 5.8 Reaction r14

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

## **Reaction equation**

$$s28 \xrightarrow{s22} s29 \tag{15}$$

#### Reactant

Table 25: Properties of each reactant.

Tueste 25: 1 repetites et euen reuetant.		
Id	Name	SBO
s28	Complex(CyclinB,Cdc2)	

#### **Modifier**

Table 26: Properties of each modifier.

Id	Name	SBO
s22	Cdc25	

## **Product**

Table 27: Properties of each product.

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Id	Name	SBO
s29	Complex(CyclinB,Cdc2)	

## **Kinetic Law**

$$v_8 = \text{not specified}$$
 (16)

## 5.9 Reaction r15

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

## **Reaction equation**

$$s30 \xrightarrow{s24, s25} s27 \tag{17}$$

#### Reactant

Table 28: Properties of each reactant.

	Name	SBO
s30	Complex(CyclinB,Cdc2)	

#### **Modifiers**

Table 29: Properties of each modifier.

Id	Name	SBO
s24	Wee1	
s25	Mik1	

#### **Product**

Table 30: Properties of each product.

Id	Name	SBO
s27	Complex(CyclinB,Cdc2)	

## **Kinetic Law**

$$v_9 = \text{not specified}$$
 (18)

#### **5.10 Reaction** r17

This is an irreversible reaction of one reactant forming two products.

## **Reaction equation**

$$s29 \longrightarrow s2 + s26$$
 (19)

#### Reactant

Table 31: Properties of each reactant.

	Name	SBO
s29	Complex(CyclinB,Cdc2)	

#### **Products**

Table 32: Properties of each product.

Id	Name	SBO
s2	Cdc2	
s26	a33_degraded	

#### **Kinetic Law**

$$v_{10} = \text{not specified}$$
 (20)

## **5.11 Reaction** r18

This is an irreversible reaction of two reactants forming one product.

## **Reaction equation**

$$s2 + s3 \longrightarrow s30$$
 (21)

#### **Reactants**

Table 33: Properties of each reactant.

Id	Name	SBO
s2	Cdc2	
s3	CyclinB	

#### **Product**

Table 34: Properties of each product.

Id	Name	SBO
s30	Complex(CyclinB,Cdc2)	

#### **Kinetic Law**

$$v_{11} = \text{not specified}$$
 (22)

# **6 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.

The identifiers for reactions, which are not defined properly or which are lacking a kinetic equation, are highlighted in red.

## **6.1 Species** s3

## Name CyclinB

#### Initial amount 0 mol

This species takes part in one reaction (as a reactant in r18).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}3 = -\nu_{11} \tag{23}$$

## 6.2 Species s5

#### Name PP2A

#### Initial amount 0 mol

This species takes part in one reaction (as a modifier in r7).

$$\frac{\mathrm{d}}{\mathrm{d}t}s5 = 0\tag{24}$$

## **6.3 Species** s6

Name Kinase X

Initial amount 0 mol

This species takes part in one reaction (as a modifier in r8).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}\mathbf{6} = 0\tag{25}$$

## **6.4 Species** s7

Name CAK

Initial amount 0 mol

This species takes part in one reaction (as a modifier in r13).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}7 = 0\tag{26}$$

## **6.5 Species** s10

Name Nim1

Initial amount 0 mol

This species takes part in two reactions (as a modifier in r11, r12).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}\mathbf{10} = 0\tag{27}$$

## 6.6 Species s11

Name Lamin

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r2 and as a product in r1).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}\mathbf{1}\mathbf{1} = \mathbf{v}_1 - \mathbf{v}_2 \tag{28}$$

## **6.7 Species** s12

Name M-Phase

**Initial amount** 0 mol

This species takes part in one reaction (as a product in r2).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}12 = \mathbf{v}_2\tag{29}$$

#### 6.8 Species s21

Name Lamin

Initial amount 0 mol

This species takes part in one reaction (as a reactant in r1).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{s}21 = -\nu_1 \tag{30}$$

## 6.9 Species s4

Name Cdc25

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r7 and as a product in r8).

$$\frac{\mathrm{d}}{\mathrm{d}t} \mathbf{s} 4 = \mathbf{v_4} - \mathbf{v_3} \tag{31}$$

## 6.10 Species s8

Name Mik1

Initial amount 0 mol

This species takes part in one reaction (as a product in r11).

$$\frac{\mathrm{d}}{\mathrm{d}t} s8 = v_5 \tag{32}$$

## **6.11 Species** s9

Name Wee1

Initial amount 0 mol

This species takes part in one reaction (as a product in r12).

$$\frac{\mathrm{d}}{\mathrm{d}t} s9 = v_6 \tag{33}$$

## **6.12 Species** s22

Name Cdc25

Initial amount 0 mol

This species takes part in three reactions (as a reactant in r8 and as a product in r7 and as a modifier in r14).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{s}22 = v_3 - v_4 \tag{34}$$

## **6.13 Species** s24

Name Wee1

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r12 and as a modifier in r15).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{s}24 = -v_6 \tag{35}$$

## **6.14 Species** s25

Name Mik1

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r11 and as a modifier in r15).

$$\frac{\mathrm{d}}{\mathrm{d}t} s25 = -v_5 \tag{36}$$

## **6.15 Species** s26

Name a33\_degraded

Initial amount 0 mol

This species takes part in one reaction (as a product in r17).

$$\frac{\mathrm{d}}{\mathrm{d}t} s26 = v_{10} \tag{37}$$

### 6.16 Species s2

Name Cdc2

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r18 and as a product in r17).

$$\frac{d}{dt}s2 = v_{10} - v_{11} \tag{38}$$

## **6.17 Species** s27

Name Complex(CyclinB,Cdc2)

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r13 and as a product in r15).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}27 = \mathbf{v}_9 - \mathbf{v}_7\tag{39}$$

### **6.18 Species** s28

Name Complex(CyclinB,Cdc2)

**Initial amount** 0 mol

This species takes part in two reactions (as a reactant in r14 and as a product in r13).

$$\frac{\mathrm{d}}{\mathrm{d}t} s28 = v_7 - v_8 \tag{40}$$

#### **6.19 Species** s29

Name Complex(CyclinB,Cdc2)

Initial amount 0 mol

This species takes part in three reactions (as a reactant in r17 and as a product in r14 and as a modifier in r1).

$$\frac{d}{dt}s29 = v_8 - v_{10} \tag{41}$$

#### **6.20 Species** s30

Name Complex(CyclinB,Cdc2)

Initial amount 0 mol

This species takes part in two reactions (as a reactant in r15 and as a product in r18).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathbf{s}30 = \mathbf{v}_{11} - \mathbf{v}_{9} \tag{42}$$

 $\mathfrak{BML2}^{d}$ X 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>, Wolfgang Müller<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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