

## SBML Model Report

**Model name: “Qu2003\_CellCycle”**



May 6, 2016

### 1 General Overview

This is a document in SBML Level 2 Version 4 format. This model was created by the following two authors: Enuo He<sup>1</sup> and Lukas Endler<sup>2</sup> at September 29<sup>th</sup> 2008 at one o’ clock in the afternoon. and last time modified at February 25<sup>th</sup> 2015 at 11:11 a. 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	1
species types	0	species	15
events	0	constraints	0
reactions	22	function definitions	0
global parameters	30	unit definitions	2
rules	2	initial assignments	0

### Model Notes

This model is from the article:

**Dynamics of the cell cycle: checkpoints, sizers, and timers.**

Qu Z, MacLellan WR, Weiss JN Biophys. J.2003 Dec; 85(6): 3600-11 [14645053](#),

**Abstract:**

We have developed a generic mathematical model of a cell cycle signaling network in higher

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eukaryotes that can be used to simulate both the G1/S and G2/M transitions. In our model, the positive feedback facilitated by CDC25 and wee1 causes bistability in cyclin-dependent kinase activity, whereas the negative feedback facilitated by SKP2 or anaphase-promoting-complex turns this bistable behavior into limit cycle behavior. The cell cycle checkpoint is a Hopf bifurcation point. These behaviors are coordinated by growth and division to maintain normal cell cycle and size homeostasis. This model successfully reproduces sizer, timer, and the restriction point features of the eukaryotic cell cycle, in addition to other experimental findings.

Figure6B has been reproduced by both SBMLodeSolver online and MathSBML. We do not include the synthesis of cyclins is proportional to cell size (Equation 2 in Page3604 of the paper) in this model. The author of the paper keep all the variables and parameters dimensionless. But in the model, we choose to use default units of SBML.

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## 2 Unit Definitions

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

### 2.1 Unit substance

**Name** arbitrary subst.

**Definition** dimensionless

### 2.2 Unit volume

**Name** arbitrary vol

**Definition** dimensionless

### 2.3 Unit area

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

**Definition**  $\text{m}^2$

## 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 Compartment

This model contains one compartment.

Table 2: Properties of all compartments.

Id	Name	SBO	Spatial Dimensions	Size	Unit	Constant	Outside
<code>cell</code>	<code>cell</code>		3	1	dimensionless	<input checked="" type="checkbox"/>	

## 3.1 Compartment `cell`

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

**Name** `cell`

## 4 Species

This model contains 15 species. Section 8 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 Condition
x1	Inactive Cyclin:CDK complex	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
x	Active Cyclin:CDK complex	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
c0	Total CDK	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
c	Free CDK	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
z0	Unphosphorylated CDC25	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
z1	One-site phosphorylated CDC25	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
z2	Two-site phosphorylated CDC25	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
w0	Unphosphorylated wee1	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
w1	phosphorylated wee1	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
u	Active SKP2 or APC	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖
i	Free CKI	cell	dimensionless dimensionless <sup>-1</sup>	· ⊖	⊖

Id	Name	Compartment	Derived Unit	Constant	Boundary Condition
ix	Cyclin:CDK:CKI complex with CKI un-phosphorylated	cell	dimensionless · dimensionless <sup>-1</sup>	<input type="checkbox"/>	<input type="checkbox"/>
ixp	Cyclin:CDK:CKI complex with CKI phosphorylated	cell	dimensionless · dimensionless <sup>-1</sup>	<input type="checkbox"/>	<input type="checkbox"/>
y	Free cyclin	cell	dimensionless · dimensionless <sup>-1</sup>	<input type="checkbox"/>	<input type="checkbox"/>
totalCyclin		cell	dimensionless · dimensionless <sup>-1</sup>	<input type="checkbox"/>	<input type="checkbox"/>

## 5 Parameters

This model contains 30 global parameters.

Table 4: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k1			300.0		<input checked="" type="checkbox"/>
k2			5.0		<input checked="" type="checkbox"/>
k3			30.0		<input checked="" type="checkbox"/>
k4			30.0		<input checked="" type="checkbox"/>
k5			0.1		<input checked="" type="checkbox"/>
k6			1.0		<input checked="" type="checkbox"/>
k7			10.0		<input checked="" type="checkbox"/>
k8			100.0		<input checked="" type="checkbox"/>
k9			1.0		<input checked="" type="checkbox"/>
k10			10.0		<input checked="" type="checkbox"/>
k11			1.0		<input checked="" type="checkbox"/>
k12			0.0		<input checked="" type="checkbox"/>
k13			1.0		<input checked="" type="checkbox"/>
k14			1.0		<input checked="" type="checkbox"/>
k15			1.0		<input checked="" type="checkbox"/>
k16			2.0		<input checked="" type="checkbox"/>
k2u			50.0		<input checked="" type="checkbox"/>
k7u			0.0		<input checked="" type="checkbox"/>
k16u			25.0		<input checked="" type="checkbox"/>
a			4.0		<input checked="" type="checkbox"/>
az			10.0		<input checked="" type="checkbox"/>
aw			10.0		<input checked="" type="checkbox"/>
ai			10.0		<input checked="" type="checkbox"/>
bz			0.1		<input checked="" type="checkbox"/>
bw			0.1		<input checked="" type="checkbox"/>
bi			0.1		<input checked="" type="checkbox"/>
cz			1.0		<input checked="" type="checkbox"/>
cw			1.0		<input checked="" type="checkbox"/>
ci			1.0		<input checked="" type="checkbox"/>
Tau			25.0		<input checked="" type="checkbox"/>

## 6 Rules

This is an overview of two rules.

### 6.1 Rule `c`

Rule `c` is an assignment rule for species `c`:

$$c = \frac{[c0] - ([x] + [x1] + [ix] + [ixp])}{[c0] \cdot \text{vol}(\text{cell})} \quad (1)$$

**Derived unit** dimensionless

**Notes** Free CDK ( normalized with `c0` ) $c=(c0-x-x1-ix-ixp)/c0$ .

### 6.2 Rule `totalCyclin`

Rule `totalCyclin` is an assignment rule for species `totalCyclin`:

$$\text{totalCyclin} = [x] + [x1] + [y] \quad (2)$$

**Derived unit** dimensionless<sup>-1</sup>

**Notes** Total cyclin ( $x+x1+y$ )

## 7 Reactions

This model contains 22 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	reaction- _0000001	Free cyclin synthesis	$\emptyset \longrightarrow y$	
2	reaction- _0000002	Cyclin binding with CDK	$y \xrightleftharpoons{c} x1$	
3	reaction- _0000003	Cyclin degradation	$y \xrightarrow{u} \emptyset$	
4	reaction- _0000004	z0 phosphorylation	$z0 \xrightleftharpoons{x} z1$	
5	reaction- _0000005	x phosphorylation	$x \xrightleftharpoons{z2, w0} x1$	
6	reaction- _0000006	z1 phosphorylation	$z1 \xrightleftharpoons{x} z2$	
7	reaction- _0000007	unphosphorylated wee1 synthesis	$\emptyset \longrightarrow w0$	
8	reaction_000008	wee1 phosphorylation	$w0 \xrightleftharpoons{x} w1$	
9	reaction_000009	SKP2 synthesis	$\emptyset \xrightarrow{x} u$	
10	reaction- _0000010	CKI synthesis	$\emptyset \longrightarrow i$	
11	reaction- _0000011	CKI binding to Active Cyclin:CDK complex	$i + x \rightleftharpoons ix$	



Nº	Id	Name	Reaction Equation	SBO
12	reaction- _0000014	ix phosphorylation	$ix \xrightarrow{x} ixp$	
13	reaction- _0000013	SKP2 degradation	$u \longrightarrow \emptyset$	
14	reaction- _0000015	CKI degradation	$i \longrightarrow \emptyset$	
15	reaction- _0000016	Cyclin:CDK:CKIp degradation	$ixp \longrightarrow x$	
16	reaction- _0000017	unphosphorylated wee1 degradation	$w0 \longrightarrow \emptyset$	
17	reaction- _0000018	phosphorylated wee1 degradation	$w1 \longrightarrow \emptyset$	
18	reaction- _0000019	Unphosphorylated CDC25 degradation	$z0 \longrightarrow \emptyset$	
19	reaction- _0000020	One site phosphorylated CDC25 degradaton	$z1 \longrightarrow \emptyset$	
20	reaction- _0000021	Two site phosphorylated CDC25 degradation	$z2 \rightleftharpoons \emptyset$	
21	reaction- _0000022	Unphosphorylation CDC25 synthesis	$\emptyset \longrightarrow z0$	
22	reaction- _0000023	Active cyclin:CDK complex degradation	$x \xrightarrow{u} \emptyset$	

### 7.1 Reaction [reaction\\_0000001](#)

This is an irreversible reaction of no reactant forming one product.

**Name** Free cyclin synthesis

#### Reaction equation



#### Product

Table 6: Properties of each product.

Id	Name	SBO
y	Free cyclin	

#### Kinetic Law

**Derived unit** contains undeclared units

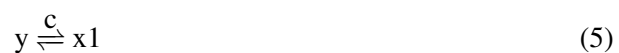
$$v_1 = k_1 \cdot \text{vol}(\text{cell}) \quad (4)$$

### 7.2 Reaction [reaction\\_0000002](#)

This is a reversible reaction of one reactant forming one product influenced by one modifier.

**Name** Cyclin binding with CDK

#### Reaction equation



#### Reactant

Table 7: Properties of each reactant.

Id	Name	SBO
y	Free cyclin	

#### Modifier

Table 8: Properties of each modifier.

Id	Name	SBO
c	Free CDK	

## Product

Table 9: Properties of each product.

Id	Name	SBO
x1	Inactive Cyclin:CDK complex	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_2 = \text{vol}(\text{cell}) \cdot (k_3 \cdot [c] \cdot [y] - [x1] \cdot k_4) \quad (6)$$

## 7.3 Reaction `reaction_0000003`

This is an irreversible reaction of one reactant forming no product influenced by one modifier.

**Name** Cyclin degradation

## Reaction equation



## Reactant

Table 10: Properties of each reactant.

Id	Name	SBO
y	Free cyclin	

## Modifier

Table 11: Properties of each modifier.

Id	Name	SBO
u	Active SKP2 or APC	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_3 = \text{vol}(\text{cell}) \cdot (k_2 + k_{2u} \cdot [u]) \cdot [y] \quad (8)$$

## 7.4 Reaction `reaction_0000004`

This is a reversible reaction of one reactant forming one product influenced by one modifier.

**Name** z0 phosphorylation

### Reaction equation



### Reactant

Table 12: Properties of each reactant.

Id	Name	SBO
z0	Unphosphorylated CDC25	

### Modifier

Table 13: Properties of each modifier.

Id	Name	SBO
x	Active Cyclin:CDK complex	

### Product

Table 14: Properties of each product.

Id	Name	SBO
z1	One-site phosphorylated CDC25	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_4 = \text{vol}(\text{cell}) \cdot ((bz + cz \cdot [x]) \cdot [z0] - [z1] \cdot az) \quad (10)$$

## 7.5 Reaction [reaction\\_0000005](#)

This is a reversible reaction of one reactant forming one product influenced by two modifiers.

**Name** x phosphorylation

### Reaction equation



### Reactant

Table 15: Properties of each reactant.

Id	Name	SBO
x	Active Cyclin:CDK complex	

### Modifiers

Table 16: Properties of each modifier.

Id	Name	SBO
z2	Two-site phosphorylated CDC25	
w0	Unphosphorylated wee1	

### Product

Table 17: Properties of each product.

Id	Name	SBO
x1	Inactive Cyclin:CDK complex	

### Kinetic Law

**Derived unit** contains undeclared units

$$v_5 = \text{vol}(\text{cell}) \cdot ((k6 + [w0]) \cdot [x] - (k5 + [z2]) \cdot [x1]) \quad (12)$$

## 7.6 Reaction [reaction\\_0000006](#)

This is a reversible reaction of one reactant forming one product influenced by one modifier.

**Name** z1 phosphorylation

### Reaction equation



### Reactant

Table 18: Properties of each reactant.

Id	Name	SBO
z1	One-site phosphorylated CDC25	

### Modifier

Table 19: Properties of each modifier.

Id	Name	SBO
x	Active Cyclin:CDK complex	

### Product

Table 20: Properties of each product.

Id	Name	SBO
z2	Two-site phosphorylated CDC25	

### Kinetic Law

**Derived unit** contains undeclared units

$$v_6 = \text{vol}(\text{cell}) \cdot ((bz + cz \cdot [x]) \cdot [z1] - az \cdot [z2]) \quad (14)$$

### 7.7 Reaction [reaction\\_0000007](#)

This is an irreversible reaction of no reactant forming one product.

**Name** unphosphorylated wee1 synthesis

### Reaction equation



### Product

Table 21: Properties of each product.

Id	Name	SBO
w0	Unphosphorylated wee1	

**Kinetic Law**

**Derived unit** contains undeclared units

$$v_7 = k_{10} \cdot \text{vol}(\text{cell}) \quad (16)$$

**7.8 Reaction** `reaction_000008`

This is a reversible reaction of one reactant forming one product influenced by one modifier.

**Name** wee1 phosphorylation

**Reaction equation****Reactant**

Table 22: Properties of each reactant.

Id	Name	SBO
w0	Unphosphorylated wee1	

**Modifier**

Table 23: Properties of each modifier.

Id	Name	SBO
x	Active Cyclin:CDK complex	

**Product**

Table 24: Properties of each product.

Id	Name	SBO
w1	phosphorylated wee1	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_8 = \text{vol}(\text{cell}) \cdot ((bw + cw \cdot [x]) \cdot [w0] - aw \cdot [w1]) \quad (18)$$

## 7.9 Reaction `reaction_000009`

This is an irreversible reaction of no reactant forming one product influenced by one modifier.

**Name** SKP2 synthesis

### Reaction equation



### Modifier

Table 25: Properties of each modifier.

Id	Name	SBO
x	Active Cyclin:CDK complex	

### Product

Table 26: Properties of each product.

Id	Name	SBO
u	Active SKP2 or APC	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_9 = \frac{\frac{\text{vol}(\text{cell}) \cdot [x]^2}{a^2 + [x]^2}}{\text{Tau}} \quad (20)$$

## 7.10 Reaction `reaction_0000010`

This is an irreversible reaction of no reactant forming one product.

**Name** CKI synthesis



### Reaction equation



### Product

Table 27: Properties of each product.

Id	Name	SBO
i	Free CKI	

### Kinetic Law

**Derived unit** contains undeclared units

$$v_{10} = k_{12} \cdot \text{vol}(\text{cell}) \quad (22)$$

### 7.11 Reaction [reaction\\_0000011](#)

This is a reversible reaction of two reactants forming one product.

**Name** CKI binding to Active Cyclin:CDK complex

### Reaction equation



### Reactants

Table 28: Properties of each reactant.

Id	Name	SBO
i	Free CKI	
x	Active Cyclin:CDK complex	

### Product

Table 29: Properties of each product.

Id	Name	SBO
ix	Cyclin:CDK:CKI complex with CKI unphosphorylated	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_{11} = (k_{14} \cdot [x] \cdot [i] - k_{15} \cdot [ix]) \cdot \text{vol}(\text{cell}) \quad (24)$$

## 7.12 Reaction [reaction\\_0000014](#)

This is a reversible reaction of one reactant forming one product influenced by one modifier.

**Name** ix phosphorylation

### Reaction equation



### Reactant

Table 30: Properties of each reactant.

Id	Name	SBO
ix	Cyclin:CDK:CKI complex with CKI unphosphorylated	

### Modifier

Table 31: Properties of each modifier.

Id	Name	SBO
x	Active Cyclin:CDK complex	

### Product

Table 32: Properties of each product.

Id	Name	SBO
ixp	Cyclin:CDK:CKI complex with CKI phosphorylated	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_{12} = \text{vol}(\text{cell}) \cdot ((b_i + c_i \cdot [x]) \cdot [ix] - a_i \cdot [ixp]) \quad (26)$$

7.13 Reaction [reaction\\_0000013](#)

This is an irreversible reaction of one reactant forming no product.

**Name** SKP2 degradation

Reaction equation



Reactant

Table 33: Properties of each reactant.

Id	Name	SBO
u	Active SKP2 or APC	

Kinetic Law

**Derived unit** contains undeclared units

$$v_{13} = \frac{\text{vol}(\text{cell}) \cdot [u]}{\text{Tau}}$$

(28)

7.14 Reaction [reaction\\_0000015](#)

This is an irreversible reaction of one reactant forming no product.

**Name** CKI degradation

Reaction equation



Reactant

Table 34: Properties of each reactant.

Id	Name	SBO
i	Free CKI	

Kinetic Law

**Derived unit** contains undeclared units

$$v_{14} = \text{vol}(\text{cell}) \cdot k_{13} \cdot [i] \quad (30)$$

### 7.15 Reaction `reaction_0000016`

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

**Name** Cyclin:CDK:CKIp degradation

#### Reaction equation



#### Reactant

Table 35: Properties of each reactant.

Id	Name	SBO
ixp	Cyclin:CDK:CKI complex with CKI phosphorylated	

#### Product

Table 36: Properties of each product.

Id	Name	SBO
x	Active Cyclin:CDK complex	

#### Kinetic Law

**Derived unit** contains undeclared units

$$v_{15} = \text{vol}(\text{cell}) \cdot k_{16} \cdot k_{16u} \cdot [\text{ixp}] \quad (32)$$

### 7.16 Reaction `reaction_0000017`

This is an irreversible reaction of one reactant forming no product.

**Name** unphosphorylated wee1 degradation

#### Reaction equation



#### Reactant

Table 37: Properties of each reactant.

Id	Name	SBO
w0	Unphosphorylated wee1	

#### Kinetic Law

**Derived unit** contains undeclared units

$$v_{16} = \text{vol}(\text{cell}) \cdot [\text{w0}] \cdot k_{11} \quad (34)$$

#### 7.17 Reaction [reaction\\_0000018](#)

This is an irreversible reaction of one reactant forming no product.

**Name** phosphorylated wee1 degradation

#### Reaction equation



#### Reactant

Table 38: Properties of each reactant.

Id	Name	SBO
w1	phosphorylated wee1	

#### Kinetic Law

**Derived unit** contains undeclared units

$$v_{17} = \text{vol}(\text{cell}) \cdot k_{11} \cdot [\text{w1}] \quad (36)$$

#### 7.18 Reaction [reaction\\_0000019](#)

This is an irreversible reaction of one reactant forming no product.

**Name** Unphosphorylated CDC25 degradation

#### Reaction equation



**Reactant**

Table 39: Properties of each reactant.

Id	Name	SBO
z0	Unphosphorylated CDC25	

#### Kinetic Law

**Derived unit** contains undeclared units

$$v_{18} = \text{vol}(\text{cell}) \cdot k_9 \cdot [z0] \quad (38)$$

#### 7.19 Reaction `reaction_0000020`

This is an irreversible reaction of one reactant forming no product.

**Name** One site phosphorylated CDC25 degradaton

#### Reaction equation



#### Reactant

Table 40: Properties of each reactant.

Id	Name	SBO
z1	One-site phosphorylated CDC25	

#### Kinetic Law

**Derived unit** contains undeclared units

$$v_{19} = \text{vol}(\text{cell}) \cdot k_9 \cdot [z1] \quad (40)$$

#### 7.20 Reaction `reaction_0000021`

This is a reversible reaction of one reactant forming no product.

**Name** Two site phosphorylated CDC25 degradation

#### Reaction equation



**Reactant**



Table 41: Properties of each reactant.

Id	Name	SBO
z2	Two-site phosphorylated CDC25	

### Kinetic Law

**Derived unit** contains undeclared units

$$v_{20} = \text{vol}(\text{cell}) \cdot k_9 \cdot [z2] \quad (42)$$

### 7.21 Reaction [reaction\\_0000022](#)

This is an irreversible reaction of no reactant forming one product.

**Name** Unphosphorylation CDC25 synthesis

### Reaction equation



### Product

Table 42: Properties of each product.

Id	Name	SBO
z0	Unphosphorylated CDC25	

### Kinetic Law

**Derived unit** contains undeclared units

$$v_{21} = \text{vol}(\text{cell}) \cdot k_8 \quad (44)$$

### 7.22 Reaction [reaction\\_0000023](#)

This is an irreversible reaction of one reactant forming no product influenced by one modifier.

**Name** Active cyclin:CDK complex degradation

### Reaction equation



## Reactant

Table 43: Properties of each reactant.

Id	Name	SBO
x	Active Cyclin:CDK complex	

## Modifier

Table 44: Properties of each modifier.

Id	Name	SBO
u	Active SKP2 or APC	

## Kinetic Law

**Derived unit** contains undeclared units

$$v_{22} = \text{vol}(\text{cell}) \cdot (k_7 + k_7 u \cdot [u]) \cdot [x] \quad (46)$$

## 8 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.

### 8.1 Species x1

**Name** Inactive Cyclin:CDK complex

**Initial concentration**  $0.1 \text{ dimensionless} \cdot \text{dimensionless}^{-1}$

This species takes part in two reactions (as a product in [reaction\\_0000002](#), [reaction\\_0000005](#)).

$$\frac{d}{dt}x1 = v_2 + v_5 \quad (47)$$

## 8.2 Species $x$

**Name** Active Cyclin:CDK complex

**Initial concentration**  $0.1 \text{ dimensionless} \cdot \text{dimensionless}^{-1}$

This species takes part in nine reactions (as a reactant in [reaction\\_0000005](#), [reaction\\_0000011](#), [reaction\\_0000023](#) and as a product in [reaction\\_0000016](#) and as a modifier in [reaction\\_0000004](#), [reaction\\_0000006](#), [reaction\\_000008](#), [reaction\\_000009](#), [reaction\\_0000014](#)).

$$\frac{d}{dt}x = v_{15} - v_5 - v_{11} - v_{22} \quad (48)$$

## 8.3 Species $c0$

**Name** Total CDK

**Initial concentration**  $200 \text{ dimensionless} \cdot \text{dimensionless}^{-1}$

This species does not take part in any reactions. Its quantity does hence not change over time:

$$\frac{d}{dt}c0 = 0 \quad (49)$$

## 8.4 Species $c$

**Name** Free CDK

**Initial concentration**  $0 \text{ dimensionless} \cdot \text{dimensionless}^{-1}$

**Involved in rule** [c](#)

This species takes part in one reaction (as a modifier in [reaction\\_0000002](#)) and is also involved in one rule which determines this species' quantity.

## 8.5 Species $z0$

**Name** Unphosphorylated CDC25

**Initial concentration**  $0 \text{ dimensionless} \cdot \text{dimensionless}^{-1}$

This species takes part in three reactions (as a reactant in [reaction\\_0000004](#), [reaction\\_0000019](#) and as a product in [reaction\\_0000022](#)).

$$\frac{d}{dt}z0 = v_{21} - v_4 - v_{18} \quad (50)$$

## 8.6 Species $z1$

**Name** One-site phosphorylated CDC25

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in three reactions (as a reactant in [reaction\\_0000006](#), [reaction\\_0000020](#) and as a product in [reaction\\_0000004](#)).

$$\frac{d}{dt}z1 = v_4 - v_6 - v_{19} \quad (51)$$

## 8.7 Species $z2$

**Name** Two-site phosphorylated CDC25

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in three reactions (as a reactant in [reaction\\_0000021](#) and as a product in [reaction\\_0000006](#) and as a modifier in [reaction\\_0000005](#)).

$$\frac{d}{dt}z2 = v_6 - v_{20} \quad (52)$$

## 8.8 Species $w0$

**Name** Unphosphorylated wee1

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in four reactions (as a reactant in [reaction\\_000008](#), [reaction\\_0000017](#) and as a product in [reaction\\_0000007](#) and as a modifier in [reaction\\_0000005](#)).

$$\frac{d}{dt}w0 = v_7 - v_8 - v_{16} \quad (53)$$

## 8.9 Species $w1$

**Name** phosphorylated wee1

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in two reactions (as a reactant in [reaction\\_0000018](#) and as a product in [reaction\\_000008](#)).

$$\frac{d}{dt}w1 = v_8 - v_{17} \quad (54)$$

### 8.10 Species $u$

**Name** Active SKP2 or APC

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in four reactions (as a reactant in [reaction\\_0000013](#) and as a product in [reaction\\_000009](#) and as a modifier in [reaction\\_0000003](#), [reaction\\_0000023](#)).

$$\frac{d}{dt}u = v_9 - v_{13} \quad (55)$$

### 8.11 Species $i$

**Name** Free CKI

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in three reactions (as a reactant in [reaction\\_0000011](#), [reaction\\_0000015](#) and as a product in [reaction\\_0000010](#)).

$$\frac{d}{dt}i = v_{10} - v_{11} - v_{14} \quad (56)$$

### 8.12 Species $ix$

**Name** Cyclin:CDK:CKI complex with CKI unphosphorylated

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in two reactions (as a reactant in [reaction\\_0000014](#) and as a product in [reaction\\_0000011](#)).

$$\frac{d}{dt}ix = v_{11} - v_{12} \quad (57)$$

### 8.13 Species $ixp$

**Name** Cyclin:CDK:CKI complex with CKI phosphorylated

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in two reactions (as a reactant in [reaction\\_0000016](#) and as a product in [reaction\\_0000014](#)).

$$\frac{d}{dt}ixp = v_{12} - v_{15} \quad (58)$$

## 8.14 Species `y`

**Name** Free cyclin

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

This species takes part in three reactions (as a reactant in [reaction\\_0000002](#), [reaction\\_0000003](#) and as a product in [reaction\\_0000001](#)).

$$\frac{d}{dt}y = v_1 - v_2 - v_3 \quad (59)$$

## 8.15 Species `totalCyclin`

**Initial concentration** 0 dimensionless · dimensionless<sup>-1</sup>

**Involved in rule** `totalCyclin`

One rule which determines this species' quantity.

SBML2<sup>A</sup>TeX 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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