

## SBML Model Report

**Model name:**  
**“Kofahl2004\_PheromonePathway”**



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: Harish Dharuri<sup>1</sup> and Jacky L Snoep<sup>2</sup> at June eighth 2005 at 2:45 p. m. and last time modified at April eighth 2016 at 3:19 p. m. Table 1 shows 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	37
events	0	constraints	0
reactions	47	function definitions	0
global parameters	0	unit definitions	6
rules	0	initial assignments	0

### Model Notes

This a model from the article:

**Modelling the dynamics of the yeast pheromone pathway.**

Kofahl B, Klipp E Yeast[2004 Jul; Volume: 21 (Issue: 10 )] Page info: 831-50 [15300679](#),

**Abstract:**

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We present a mathematical model of the dynamics of the pheromone pathways in haploid yeast cells of mating type MATa after stimulation with pheromone alpha-factor. The model consists of a set of differential equations and describes the dynamics of signal transduction from the receptor via several steps, including a G protein and a scaffold MAP kinase cascade, up to changes in the gene expression after pheromone stimulation in terms of biochemical changes (complex formations, phosphorylations, etc.). The parameters entering the models have been taken from the literature or adapted to observed time courses or behaviour. Using this model we can follow the time course of the various complex formation processes and of the phosphorylation states of the proteins involved. Furthermore, we can explain the phenotype of more than a dozen well-characterized mutants and also the graded response of yeast cells to varying concentrations of the stimulating pheromone.

The model was updated on 21<sup>st</sup> October 2010, by Vijayalakshmi Chelliah. The following changes were made: 1) The model has been converted to SBML l2v4.2) The model has been recurated and the curation figure was updated (units are in nanoMolar; but the publication has units in microMolar). Simulations were done using Copasi v4.6 (Build 32).3) Notes have been added.4) Annotation for one of the species has been corrected (Complex M).

**SBML level 2 code generated for the JWS Online project by Jacky Snoep using PySCeS**  
**Run this model online at <http://jjj.biochem.sun.ac.za>**  
**To cite JWS Online please refer to: Olivier, B.G. and Snoep, J.L. (2004) [Web-based modelling using JWS Online](#), *Bioinformatics*, 20:2143-2144**

The following are the four major differences between the original publication by Kofahl et al and the model that actually is able to replicate the results as depicted in the publication (those corrections have been made in agreement with the authors):

1. Bar1 is the inactive protease present inside the cell but the publication wrongly mentions that Bar1 is also the protease that is present on the extracellular surface. The model correctly names the protease in it's different forms by calling inactive Bar1 within the cell as Bar1, active Bar1 within the cell as Bar1a and extracellular Bar1 as Bar1aex
2. The initial amount of Alpha-factor is given as 1000nM but the model uses a value of 100nM.
3. The value of the parameter k8 is given as 0.33 but the model uses a value of 0.033.
4. The value of the parameter k41 is given as 0.002 but the model uses a value of 0.02.

This model originates from BioModels Database: A Database of Annotated Published Models (<http://www.ebi.ac.uk/biomodels/>). It is copyright (c) 2005-2010 The BioModels.net Team.

For more information see the [terms of use](#).

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 nine unit definitions of which three are predefined by SBML and not mentioned in the model.

### 2.1 Unit `substance`

**Name** nanomole (default)

**Definition** nmol

### 2.2 Unit `time`

**Name** minute (default)

**Definition** 60 s

### 2.3 Unit `nanoMolar`

**Name** nanoMolar

**Definition**  $\text{nmol} \cdot \text{l}^{-1}$

### 2.4 Unit `min_inv`

**Name** min\_inv

**Definition**  $(60 \text{ s})^{-1}$

### 2.5 Unit `min_inv_nM_inv`

**Name** min\_inv\_nM\_inv

**Definition**  $\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$

### 2.6 Unit `nM_min_inv`

**Name** nM\_min\_inv

**Definition**  $\text{nmol} \cdot \text{l}^{-1} \cdot (60 \text{ s})^{-1}$

### 2.7 Unit `volume`

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

**Definition** 1

## 2.8 Unit area

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

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

## 2.9 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.

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

### 3.1 Compartment Extracellular

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

### 3.2 Compartment compartment

This is a three dimensional compartment with a constant size of one litre, which is surrounded by Extracellular.

**Name** cell

## 4 Species

This model contains 37 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 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- tion
alpha	-factor	Extracellular	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste2		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste2a	Ste2active	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Gabc	G	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
GaGTP	GGTP	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Gbc	G	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
GaGDP	GGDP	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexC		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexD		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste5		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste11		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexA		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste7		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Fus3		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexB		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste20		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexE		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexF		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexG		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexH		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexI		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>

Id	Name	Compartment	Derived Unit	Constant	Boundary Condi- tion
complexL		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Fus3PP		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexK		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste12		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Ste12a	Ste12active	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Bar1		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Bar1a	Bar1active	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Bar1aex	Bar1activeEx	Extracellular	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Far1		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Far1PP		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Far1U	Far1ubiquitin	compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexM		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
complexN		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Cdc28		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
Sst2		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input type="checkbox"/>
p		compartment	$\text{nmol} \cdot \text{l}^{-1}$	<input type="checkbox"/>	<input checked="" type="checkbox"/>

## 5 Reactions

This model contains 47 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 4: Overview of all reactions

Nº	Id	Name	Reaction Equation	SBO
1	v1	alpha degradation	$\text{alpha} \xrightarrow{\text{Bar1aex}} \emptyset$	
2	v2		$\text{Ste2} \xrightleftharpoons{\text{alpha}} \text{Ste2a}$	
3	v3		$\text{Ste2a} \rightleftharpoons \text{Ste2}$	
4	v4		$\text{Ste2a} \rightleftharpoons \text{p}$	
5	v5		$\text{Ste2} \rightleftharpoons \emptyset$	
6	v6		$\text{Gabc} \xrightleftharpoons{\text{Ste2a}} \text{GaGTP} + \text{Gbc}$	
7	v7		$\text{GaGTP} \rightleftharpoons \text{GaGDP}$	
8	v8		$\text{GaGTP} \xrightleftharpoons{\text{Sst2}} \text{GaGDP}$	
9	v9		$\text{GaGDP} + \text{Gbc} \rightleftharpoons \text{Gabc}$	
10	v10		$\text{Gbc} + \text{complexC} \rightleftharpoons \text{complexD}$	
11	v11		$\text{complexD} \rightleftharpoons \text{Gbc} + \text{complexC}$	
12	v12		$\text{Ste11} + \text{Ste5} \rightleftharpoons \text{complexA}$	
13	v13		$\text{complexA} \rightleftharpoons \text{Ste11} + \text{Ste5}$	
14	v14		$\text{Fus3} + \text{Ste7} \rightleftharpoons \text{complexB}$	
15	v15		$\text{complexB} \rightleftharpoons \text{Fus3} + \text{Ste7}$	
16	v16		$\text{complexA} + \text{complexB} \rightleftharpoons \text{complexC}$	
17	v17		$\text{complexC} \rightleftharpoons \text{Fus3} + \text{Ste11} + \text{Ste7} + \text{Ste5}$	
18	v18		$\text{complexD} + \text{Ste20} \rightleftharpoons \text{complexE}$	
19	v19		$\text{complexE} \rightleftharpoons \text{complexD} + \text{Ste20}$	
20	v20		$\text{complexE} \rightleftharpoons \text{complexF}$	

Nº	Id	Name	Reaction Equation	SBO
21	v21		$\text{complexE} \rightleftharpoons \text{Gbc} + \text{Ste7} + \text{Ste5} + \text{Fus3} + \text{Ste20} + \text{Ste11}$	
22	v22		$\text{complexF} \rightleftharpoons \text{complexG}$	
23	v23		$\text{complexF} \rightleftharpoons \text{Gbc} + \text{Ste7} + \text{Ste5} + \text{Fus3} + \text{Ste20} + \text{Ste11}$	
24	v24		$\text{complexG} \rightleftharpoons \text{complexH}$	
25	v25		$\text{complexG} \rightleftharpoons \text{Gbc} + \text{Ste7} + \text{Ste5} + \text{Fus3} + \text{Ste20} + \text{Ste11}$	
26	v26		$\text{complexH} \rightleftharpoons \text{complexI}$	
27	v27		$\text{complexH} \rightleftharpoons \text{Gbc} + \text{Ste7} + \text{Ste5} + \text{Fus3} + \text{Ste20} + \text{Ste11}$	
28	v28		$\text{complexI} \rightleftharpoons \text{complexL} + \text{Fus3PP}$	
29	v29		$\text{complexL} + \text{Fus3} \rightleftharpoons \text{complexK}$	
30	v30		$\text{complexK} \rightleftharpoons \text{complexL} + \text{Fus3}$	
31	v31		$\text{complexK} \rightleftharpoons \text{complexI}$	
32	v32		$\text{complexL} \rightleftharpoons \text{Gbc} + \text{Ste7} + \text{Ste5} + \text{Ste20} + \text{Ste11}$	
33	v33		$\text{Fus3PP} \rightleftharpoons \text{Fus3}$	
34	v34		$\text{Ste12} + \text{Fus3PP} \rightleftharpoons \text{Ste12a}$	
35	v35		$\text{Ste12a} \rightleftharpoons \text{Ste12} + \text{Fus3PP}$	
36	v36		$\text{Bar1} \xrightleftharpoons{\text{Ste12a}} \text{Bar1a}$	
37	v37		$\text{Bar1a} \rightleftharpoons \text{Bar1}$	
38	v38		$\text{Bar1a} \rightleftharpoons \text{Bar1aex}$	
39	v39		$\text{Far1} \xrightleftharpoons{\text{Fus3PP}} \text{Far1PP}$	
40	v40		$\text{Far1PP} \rightleftharpoons \text{Far1}$	
41	v41		$\text{Far1} \xrightleftharpoons{\text{Cdc28}} \text{Far1U}$	
42	v42		$\text{Gbc} + \text{Far1PP} \rightleftharpoons \text{complexM}$	
43	v43		$\text{complexM} \rightleftharpoons \text{Gbc} + \text{Far1PP}$	
44	v44		$\text{complexN} \rightleftharpoons \text{Cdc28} + \text{Far1PP}$	



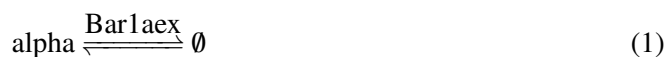
Nº	Id	Name	Reaction Equation	SBO
45	v45		$\text{Cdc28} + \text{Far1PP} \rightleftharpoons \text{complexN}$	
46	v46		$\text{p} \xrightleftharpoons{\text{Fus3PP}} \text{Sst2}$	
47	v47		$\text{Sst2} \rightleftharpoons \text{p}$	

## 5.1 Reaction v1

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

**Name** alpha degradation

### Reaction equation



### Reactant

Table 5: Properties of each reactant.

Id	Name	SBO
alpha	-factor	

### Modifier

Table 6: Properties of each modifier.

Id	Name	SBO
Bar1aex	Bar1activeEx	

### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_1 = \text{vol}(\text{Extracellular}) \cdot [\text{alpha}] \cdot [\text{Bar1aex}] \cdot k_1 \quad (2)$$

Table 7: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k1			0.03	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.2 Reaction v2

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

## Reaction equation



## Reactant

Table 8: Properties of each reactant.

Id	Name	SBO
Ste2		

## Modifier

Table 9: Properties of each modifier.

Id	Name	SBO
alpha	-factor	

## Product

Table 10: Properties of each product.

Id	Name	SBO
Ste2a	Ste2active	

## Kinetic Law

**Derived unit**  $9.99999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_2 = \text{vol}(\text{compartment}) \cdot [\text{Ste2}] \cdot [\alpha] \cdot k_2 \quad (4)$$

Table 11: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k2			0.001	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.3 Reaction v3

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 12: Properties of each reactant.

Id	Name	SBO
Ste2a	Ste2active	

### Product

Table 13: Properties of each product.

Id	Name	SBO
Ste2		

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_3 = \text{vol}(\text{compartment}) \cdot [\text{Ste2a}] \cdot k_3 \quad (6)$$

Table 14: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k3			0.6	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.4 Reaction v4

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 15: Properties of each reactant.

Id	Name	SBO
Ste2a	Ste2active	

## Product

Table 16: Properties of each product.

Id	Name	SBO
p		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_4 = \text{vol}(\text{compartment}) \cdot [\text{Ste2a}] \cdot k_4 \quad (8)$$

Table 17: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k4			0.24	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.5 Reaction v5

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

### Reaction equation



## Reactant

Table 18: Properties of each reactant.

Id	Name	SBO
Ste2		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_5 = \text{vol}(\text{compartment}) \cdot [\text{Ste2}] \cdot k_5 \quad (10)$$

Table 19: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k5			0.024	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.6 Reaction v6

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

### Reaction equation



## Reactant

Table 20: Properties of each reactant.

Id	Name	SBO
Gabc	G	

## Modifier

Table 21: Properties of each modifier.

Id	Name	SBO
Ste2a	Ste2active	

## Products

Table 22: Properties of each product.

Id	Name	SBO
GaGTP	GGTP	
Gbc	G	

Id	Name	SBO
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### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_6 = \text{vol}(\text{compartment}) \cdot [\text{Ste2a}] \cdot [\text{Gabc}] \cdot k_6 \quad (12)$$

Table 23: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k6			0.004	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.7 Reaction $v_7$

This is a reversible reaction of one reactant forming one product.

#### Reaction equation



#### Reactant

Table 24: Properties of each reactant.

Id	Name	SBO
GaGTP	GGTP	

#### Product

Table 25: Properties of each product.

Id	Name	SBO
GaGDP	GGDP	

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_7 = \text{vol}(\text{compartment}) \cdot [\text{GaGTP}] \cdot k_7 \quad (14)$$

Table 26: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k7			0.24	$(60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.8 Reaction v8

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

### Reaction equation



### Reactant

Table 27: Properties of each reactant.

Id	Name	SBO
GaGTP	GGTP	

### Modifier

Table 28: Properties of each modifier.

Id	Name	SBO
Sst2		

### Product

Table 29: Properties of each product.

Id	Name	SBO
GaGDP	GGDP	

### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60\text{ s})^{-1}$

$$v_8 = \text{vol}(\text{compartment}) \cdot [\text{GaGTP}] \cdot [\text{Sst2}] \cdot k_8 \quad (16)$$



Table 30: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k8			0.033	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.9 Reaction v9

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

### Reaction equation



### Reactants

Table 31: Properties of each reactant.

Id	Name	SBO
GaGDP	GGDP	
Gbc	G	

### Product

Table 32: Properties of each product.

Id	Name	SBO
Gabc	G	

### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_9 = \text{vol}(\text{compartment}) \cdot [\text{GaGDP}] \cdot [\text{Gbc}] \cdot k_9 \quad (18)$$

Table 33: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k9			2000.0	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.10 Reaction v10

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

### Reaction equation



### Reactants

Table 34: Properties of each reactant.

Id	Name	SBO
Gbc	G	
complexC		

### Product

Table 35: Properties of each product.

Id	Name	SBO
complexD		

### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{10} = \text{vol}(\text{compartment}) \cdot [\text{Gbc}] \cdot [\text{complexC}] \cdot k_{10} \quad (20)$$

Table 36: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k10			0.1	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.11 Reaction v11

This is a reversible reaction of one reactant forming two products.

### Reaction equation



## Reactant

Table 37: Properties of each reactant.

Id	Name	SBO
complexD		

## Products

Table 38: Properties of each product.

Id	Name	SBO
Gbc	G	
complexC		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{11} = \text{vol}(\text{compartment}) \cdot [\text{complexD}] \cdot k_{11} \quad (22)$$

Table 39: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k11			5.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.12 Reaction v12

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

## Reaction equation



## Reactants

Table 40: Properties of each reactant.

Id	Name	SBO
Ste11		

Id	Name	SBO
Ste5		

## Product

Table 41: Properties of each product.

Id	Name	SBO
complexA		

## Kinetic Law

**Derived unit**  $9.99999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{12} = \text{vol}(\text{compartment}) \cdot [\text{Ste5}] \cdot [\text{Ste11}] \cdot k_{12} \quad (24)$$

Table 42: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k12			1.0	$\text{nmol}^{-1} \cdot 1 \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.13 Reaction v13

This is a reversible reaction of one reactant forming two products.

### Reaction equation



## Reactant

Table 43: Properties of each reactant.

Id	Name	SBO
complexA		

## Products

Table 44: Properties of each product.

Id	Name	SBO
Ste11		
Ste5		

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{13} = \text{vol}(\text{compartment}) \cdot [\text{complexA}] \cdot k_{13} \quad (26)$$

Table 45: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k13			3.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.14 Reaction v14

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

### Reaction equation



### Reactants

Table 46: Properties of each reactant.

Id	Name	SBO
Fus3		
Ste7		

### Product

Table 47: Properties of each product.

Id	Name	SBO
complexB		

## Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{14} = \text{vol}(\text{compartment}) \cdot [\text{Ste7}] \cdot [\text{Fus3}] \cdot k_{14} \quad (28)$$

Table 48: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k14			1.0	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.15 Reaction v15

This is a reversible reaction of one reactant forming two products.

### Reaction equation



### Reactant

Table 49: Properties of each reactant.

Id	Name	SBO
complexB		

### Products

Table 50: Properties of each product.

Id	Name	SBO
Fus3		
Ste7		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{15} = \text{vol}(\text{compartment}) \cdot [\text{complexB}] \cdot k_{15} \quad (30)$$

Table 51: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k15			3.0	$(60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.16 Reaction v16

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

#### Reaction equation



#### Reactants

Table 52: Properties of each reactant.

Id	Name	SBO
	complexA	
	complexB	

#### Product

Table 53: Properties of each product.

Id	Name	SBO
	complexC	

#### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60\text{ s})^{-1}$

$$v_{16} = \text{vol}(\text{compartment}) \cdot [\text{complexA}] \cdot [\text{complexB}] \cdot k16 \quad (32)$$

Table 54: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k16			3.0	$\text{nmol}^{-1} \cdot \text{l} \cdot (60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.17 Reaction v17

This is a reversible reaction of one reactant forming four products.

#### Reaction equation



#### Reactant

Table 55: Properties of each reactant.

Id	Name	SBO
complexC		

#### Products

Table 56: Properties of each product.

Id	Name	SBO
Fus3		
Ste11		
Ste7		
Ste5		

#### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{17} = \text{vol}(\text{compartment}) \cdot [\text{complexC}] \cdot k_{17} \quad (34)$$

Table 57: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k17			100.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.18 Reaction v18

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



### Reaction equation



### Reactants

Table 58: Properties of each reactant.

Id	Name	SBO
complexD		
Ste20		

### Product

Table 59: Properties of each product.

Id	Name	SBO
complexE		

### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{18} = \text{vol}(\text{compartment}) \cdot [\text{complexD}] \cdot [\text{Ste20}] \cdot k18 \quad (36)$$

Table 60: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k18			5.0	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.19 Reaction v19

This is a reversible reaction of one reactant forming two products.

### Reaction equation



### Reactant

Table 61: Properties of each reactant.

Id	Name	SBO
complexE		

## Products

Table 62: Properties of each product.

Id	Name	SBO
complexD		
Ste20		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{19} = \text{vol}(\text{compartment}) \cdot [\text{complexE}] \cdot k_{19} \quad (38)$$

Table 63: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k19			1.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.20 Reaction v20

This is a reversible reaction of one reactant forming one product.

### Reaction equation



## Reactant

Table 64: Properties of each reactant.

Id	Name	SBO
complexE		

## Product

Table 65: Properties of each product.

Id	Name	SBO
complexF		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{20} = \text{vol}(\text{compartment}) \cdot [\text{complexE}] \cdot k_{20} \quad (40)$$

Table 66: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k20			10.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.21 Reaction v21

This is a reversible reaction of one reactant forming six products.

## Reaction equation



## Reactant

Table 67: Properties of each reactant.

Id	Name	SBO
complexE		

## Products

Table 68: Properties of each product.

Id	Name	SBO
Gbc	G	
Ste7		

Id	Name	SBO
Ste5		
Fus3		
Ste20		
Ste11		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{21} = \text{vol}(\text{compartment}) \cdot [\text{complexE}] \cdot k_{21} \quad (42)$$

Table 69: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k21			5.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.22 Reaction v22

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 70: Properties of each reactant.

Id	Name	SBO
complexF		

### Product

Table 71: Properties of each product.

Id	Name	SBO
complexG		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{22} = \text{vol}(\text{compartment}) \cdot [\text{complexF}] \cdot k_{22} \quad (44)$$

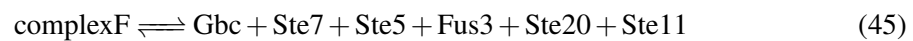
Table 72: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k22			47.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.23 Reaction v23

This is a reversible reaction of one reactant forming six products.

### Reaction equation



## Reactant

Table 73: Properties of each reactant.

Id	Name	SBO
complexF		

## Products

Table 74: Properties of each product.

Id	Name	SBO
Gbc	G	
Ste7		
Ste5		
Fus3		
Ste20		
Ste11		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{23} = \text{vol}(\text{compartment}) \cdot [\text{complexF}] \cdot k_{23} \quad (46)$$

Table 75: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k23			5.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.24 Reaction v24

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 76: Properties of each reactant.

Id	Name	SBO
	complexG	

### Product

Table 77: Properties of each product.

Id	Name	SBO
	complexH	

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{24} = \text{vol}(\text{compartment}) \cdot [\text{complexG}] \cdot k_{24} \quad (48)$$

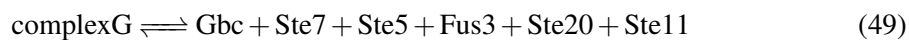
Table 78: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k24			345.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.25 Reaction v25

This is a reversible reaction of one reactant forming six products.

#### Reaction equation



#### Reactant

Table 79: Properties of each reactant.

Id	Name	SBO
complexG		

#### Products

Table 80: Properties of each product.

Id	Name	SBO
Gbc	G	
Ste7		
Ste5		
Fus3		
Ste20		
Ste11		

#### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{25} = \text{vol}(\text{compartment}) \cdot [\text{complexG}] \cdot k_{25} \quad (50)$$

Table 81: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k25			5.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.26 Reaction v26

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 82: Properties of each reactant.

Id	Name	SBO
complexH		

### Product

Table 83: Properties of each product.

Id	Name	SBO
complexI		

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{26} = \text{vol}(\text{compartment}) \cdot [\text{complexH}] \cdot k_{26} \quad (52)$$

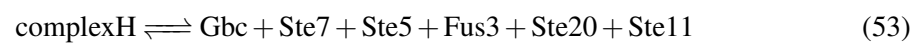
Table 84: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k26			50.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.27 Reaction v27

This is a reversible reaction of one reactant forming six products.

### Reaction equation



### Reactant



Table 85: Properties of each reactant.

Id	Name	SBO
complexH		

## Products

Table 86: Properties of each product.

Id	Name	SBO
Gbc	G	
Ste7		
Ste5		
Fus3		
Ste20		
Ste11		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{27} = \text{vol}(\text{compartment}) \cdot [\text{complexH}] \cdot k_{27} \quad (54)$$

Table 87: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k27			5.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.28 Reaction v28

This is a reversible reaction of one reactant forming two products.

## Reaction equation



## Reactant

Table 88: Properties of each reactant.

Id	Name	SBO
complexI		

## Products

Table 89: Properties of each product.

Id	Name	SBO
complexL		
Fus3PP		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{28} = \text{vol}(\text{compartment}) \cdot [\text{complexI}] \cdot k_{28} \quad (56)$$

Table 90: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k28			140.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.29 Reaction v29

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

### Reaction equation



## Reactants

Table 91: Properties of each reactant.

Id	Name	SBO
complexL		
Fus3		

## Product

Table 92: Properties of each product.

Id	Name	SBO
complexK		

## Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{29} = \text{vol}(\text{compartment}) \cdot [\text{complexL}] \cdot [\text{Fus3}] \cdot k_{29} \quad (58)$$

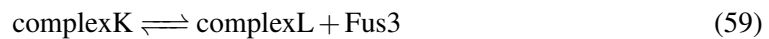
Table 93: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k29			10.0	$\text{nmol}^{-1} \cdot 1 \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.30 Reaction v30

This is a reversible reaction of one reactant forming two products.

## Reaction equation



## Reactant

Table 94: Properties of each reactant.

Id	Name	SBO
complexK		

## Products

Table 95: Properties of each product.

Id	Name	SBO
complexL		
Fus3		

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{30} = \text{vol}(\text{compartment}) \cdot [\text{complexK}] \cdot k_{30} \quad (60)$$

Table 96: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k30			1.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.31 Reaction v31

This is a reversible reaction of one reactant forming one product.

#### Reaction equation



#### Reactant

Table 97: Properties of each reactant.

Id	Name	SBO
complexK		

#### Product

Table 98: Properties of each product.

Id	Name	SBO
complexI		

### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{31} = \text{vol}(\text{compartment}) \cdot [\text{complexK}] \cdot k_{31} \quad (62)$$

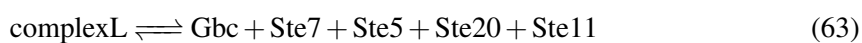
Table 99: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k31			250.0	$(60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.32 Reaction v32

This is a reversible reaction of one reactant forming five products.

#### Reaction equation



#### Reactant

Table 100: Properties of each reactant.

Id	Name	SBO
complexL		

#### Products

Table 101: Properties of each product.

Id	Name	SBO
Gbc	G	
Ste7		
Ste5		
Ste20		
Ste11		

#### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60\text{ s})^{-1}$

$$v_{32} = \text{vol}(\text{compartment}) \cdot [\text{complexL}] \cdot k_{32} \quad (64)$$

Table 102: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k32			5.0	$(60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.33 Reaction v33

This is a reversible reaction of one reactant forming one product.

#### Reaction equation



#### Reactant

Table 103: Properties of each reactant.

Id	Name	SBO
Fus3PP		

#### Product

Table 104: Properties of each product.

Id	Name	SBO
Fus3		

#### Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{33} = \text{vol}(\text{compartment}) \cdot [\text{Fus3PP}] \cdot k_{33} \quad (66)$$

Table 105: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k33			50.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.34 Reaction v34

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

#### Reaction equation



## Reactants

Table 106: Properties of each reactant.

Id	Name	SBO
Ste12		
Fus3PP		

## Product

Table 107: Properties of each product.

Id	Name	SBO
Ste12a	Ste12active	

## Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{34} = \text{vol}(\text{compartment}) \cdot [\text{Ste12}] \cdot [\text{Fus3PP}] \cdot k_{34} \quad (68)$$

Table 108: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k34			18.0	$\text{nmol}^{-1} \cdot 1 \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.35 Reaction v35

This is a reversible reaction of one reactant forming two products.

## Reaction equation



## Reactant

Table 109: Properties of each reactant.

Id	Name	SBO
Ste12a	Ste12active	

## Products

Table 110: Properties of each product.

Id	Name	SBO
Ste12		
Fus3PP		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{35} = \text{vol}(\text{compartment}) \cdot [\text{Ste12a}] \cdot k_{35} \quad (70)$$

Table 111: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k35			10.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.36 Reaction v36

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

## Reaction equation



## Reactant

Table 112: Properties of each reactant.

Id	Name	SBO
Bar1		

## Modifier

Table 113: Properties of each modifier.

Id	Name	SBO
Ste12a	Ste12active	



Id	Name	SBO
----	------	-----

## Product

Table 114: Properties of each product.

Id	Name	SBO
Bar1a	Bar1active	

## Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{36} = \text{vol}(\text{compartment}) \cdot [\text{Ste12a}] \cdot [\text{Bar1}] \cdot k_{36} \quad (72)$$

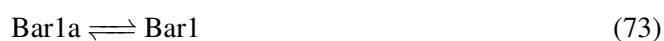
Table 115: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k36			0.1	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.37 Reaction $v_{37}$

This is a reversible reaction of one reactant forming one product.

## Reaction equation



## Reactant

Table 116: Properties of each reactant.

Id	Name	SBO
Bar1a	Bar1active	

## Product

Table 117: Properties of each product.

Id	Name	SBO
Bar1		

**Kinetic Law****Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$ 

$$v_{37} = \text{vol}(\text{compartment}) \cdot [\text{Bar1a}] \cdot k_{37} \quad (74)$$

Table 118: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k37			0.1	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

**5.38 Reaction v38**

This is a reversible reaction of one reactant forming one product.

**Reaction equation****Reactant**

Table 119: Properties of each reactant.

Id	Name	SBO
Bar1a	Bar1active	

**Product**

Table 120: Properties of each product.

Id	Name	SBO
Bar1aex	Bar1activeEx	

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{38} = \text{vol}(\text{compartment}) \cdot [\text{Bar1a}] \cdot k_{38} \quad (76)$$

Table 121: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k38			0.01	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.39 Reaction v39

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

### Reaction equation



## Reactant

Table 122: Properties of each reactant.

Id	Name	SBO
Far1		

## Modifier

Table 123: Properties of each modifier.

Id	Name	SBO
Fus3PP		

## Product

Table 124: Properties of each product.

Id	Name	SBO
Far1PP		

## Kinetic Law

**Derived unit** contains undeclared units

$$v_{39} = \text{vol}(\text{compartment}) \cdot \frac{[\text{Far1}] \cdot [\text{Fus3PP}] \cdot [\text{Fus3PP}]}{100 \cdot 100 + [\text{Fus3PP}] \cdot [\text{Fus3PP}]} \cdot k_{39} \quad (78)$$

Table 125: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k39			18.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.40 Reaction v40

This is a reversible reaction of one reactant forming one product.

### Reaction equation



### Reactant

Table 126: Properties of each reactant.

Id	Name	SBO
Far1PP		

### Product

Table 127: Properties of each product.

Id	Name	SBO
Far1		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{40} = \text{vol}(\text{compartment}) \cdot [\text{Far1PP}] \cdot k_{40} \quad (80)$$

Table 128: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k40			1.0	$(60\text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.41 Reaction $v_{41}$

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

#### Reaction equation



#### Reactant

Table 129: Properties of each reactant.

Id	Name	SBO
Far1		

#### Modifier

Table 130: Properties of each modifier.

Id	Name	SBO
Cdc28		

#### Product

Table 131: Properties of each product.

Id	Name	SBO
Far1U	Far1ubiquitin	

#### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60\text{ s})^{-1}$

$$v_{41} = \text{vol}(\text{compartment}) \cdot [\text{Far1}] \cdot [\text{Cdc28}] \cdot k_{41} \quad (82)$$

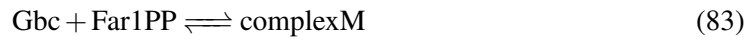
Table 132: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k41			0.02	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

### 5.42 Reaction v42

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

#### Reaction equation



#### Reactants

Table 133: Properties of each reactant.

Id	Name	SBO
Gbc	G	
Far1PP		

#### Product

Table 134: Properties of each product.

Id	Name	SBO
complexM		

#### Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{42} = \text{vol}(\text{compartment}) \cdot [\text{Gbc}] \cdot [\text{Far1PP}] \cdot k_{42} \quad (84)$$

Table 135: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k42			0.1	$\text{nmol}^{-1} \cdot \text{l} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

5.43 Reaction v43

This is a reversible reaction of one reactant forming two products.

Reaction equation



Reactant

Table 136: Properties of each reactant.

Id	Name	SBO
complexM		

Products

Table 137: Properties of each product.

Id	Name	SBO
Gbc	G	
Far1PP		

Kinetic Law

Derived unit  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$v_{43} = \text{vol}(\text{compartment}) \cdot [\text{complexM}] \cdot k_{43}$

(86)

Table 138: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k43			0.01	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

5.44 Reaction v44

This is a reversible reaction of one reactant forming two products.

Reaction equation



## Reactant

Table 139: Properties of each reactant.

Id	Name	SBO
complexN		

## Products

Table 140: Properties of each product.

Id	Name	SBO
Cdc28		
Far1PP		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{44} = \text{vol}(\text{compartment}) \cdot [\text{complexN}] \cdot k_{44} \quad (88)$$

Table 141: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k44			0.01	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.45 Reaction $v_{45}$

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

## Reaction equation



## Reactants

Table 142: Properties of each reactant.

Id	Name	SBO
Cdc28		



Id	Name	SBO
Far1PP		

## Product

Table 143: Properties of each product.

Id	Name	SBO
complexN		

## Kinetic Law

**Derived unit**  $9.999999999999998 \cdot 10^{-10} \text{ mol} \cdot (60 \text{ s})^{-1}$

$$v_{45} = \text{vol}(\text{compartment}) \cdot [\text{Far1PP}] \cdot [\text{Cdc28}] \cdot k_{45} \quad (90)$$

Table 144: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k45			0.1	$\text{nmol}^{-1} \cdot 1 \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.46 Reaction v46

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

### Reaction equation



## Reactant

Table 145: Properties of each reactant.

Id	Name	SBO
p		

## Modifier

Table 146: Properties of each modifier.

Id	Name	SBO
Fus3PP		

## Product

Table 147: Properties of each product.

Id	Name	SBO
Sst2		

## Kinetic Law

**Derived unit** contains undeclared units

$$v_{46} = \text{vol}(\text{compartment}) \cdot \frac{[\text{Fus3PP}]^2}{4^2 + [\text{Fus3PP}]^2} \cdot k_{46} \quad (92)$$

Table 148: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k46			200.0	$\text{nmol} \cdot \text{l}^{-1} \cdot (60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 5.47 Reaction v47

This is a reversible reaction of one reactant forming one product.

## Reaction equation



## Reactant

Table 149: Properties of each reactant.

Id	Name	SBO
Sst2		

## Product

Table 150: Properties of each product.

Id	Name	SBO
p		

## Kinetic Law

**Derived unit**  $\text{nmol} \cdot (60 \text{ s})^{-1}$

$$v_{47} = \text{vol}(\text{compartment}) \cdot [\text{Sst2}] \cdot k_{47} \quad (94)$$

Table 151: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
k47			1.0	$(60 \text{ s})^{-1}$	<input checked="" type="checkbox"/>

## 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.

### 6.1 Species `alpha`

**Name** -factor

**Initial concentration**  $100 \text{ nmol} \cdot \text{l}^{-1}$

This species takes part in two reactions (as a reactant in [v1](#) and as a modifier in [v2](#)).

$$\frac{d}{dt} \text{alpha} = -v_1 \quad (95)$$

### 6.2 Species `Ste2`

**Initial concentration**  $1666.6666667 \text{ nmol} \cdot \text{l}^{-1}$

This species takes part in three reactions (as a reactant in [v2](#), [v5](#) and as a product in [v3](#)).

$$\frac{d}{dt} \text{Ste2} = v_3 - v_2 - v_5 \quad (96)$$

### 6.3 Species Ste2a

**Name** Ste2active

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in four reactions (as a reactant in v3, v4 and as a product in v2 and as a modifier in v6).

$$\frac{d}{dt}\text{Ste2a} = v_2 - v_3 - v_4 \quad (97)$$

### 6.4 Species Gabc

**Name** G

**Initial concentration** 1666.6666667 nmol · l<sup>-1</sup>

This species takes part in two reactions (as a reactant in v6 and as a product in v9).

$$\frac{d}{dt}\text{Gabc} = v_9 - v_6 \quad (98)$$

### 6.5 Species GaGTP

**Name** GGTP

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in v7, v8 and as a product in v6).

$$\frac{d}{dt}\text{GaGTP} = v_6 - v_7 - v_8 \quad (99)$$

### 6.6 Species Gbc

**Name** G

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in eleven reactions (as a reactant in v9, v10, v42 and as a product in v6, v11, v21, v23, v25, v27, v32, v43).

$$\frac{d}{dt}\text{Gbc} = v_6 + v_{11} + v_{21} + v_{23} + v_{25} + v_{27} + v_{32} + v_{43} - v_9 - v_{10} - v_{42} \quad (100)$$

## 6.7 Species GaGDP

**Name** GGDP

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in v9 and as a product in v7, v8).

$$\frac{d}{dt}\text{GaGDP} = v_7 + v_8 - v_9 \quad (101)$$

## 6.8 Species complexC

**Initial concentration** 235.724935791903 nmol · l<sup>-1</sup>

This species takes part in four reactions (as a reactant in v10, v17 and as a product in v11, v16).

$$\frac{d}{dt}\text{complexC} = v_{11} + v_{16} - v_{10} - v_{17} \quad (102)$$

## 6.9 Species complexD

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in four reactions (as a reactant in v11, v18 and as a product in v10, v19).

$$\frac{d}{dt}\text{complexD} = v_{10} + v_{19} - v_{11} - v_{18} \quad (103)$$

## 6.10 Species Ste5

**Initial concentration** 158.33176608789 nmol · l<sup>-1</sup>

This species takes part in eight reactions (as a reactant in v12 and as a product in v13, v17, v21, v23, v25, v27, v32).

$$\frac{d}{dt}\text{Ste5} = v_{13} + v_{17} + v_{21} + v_{23} + v_{25} + v_{27} + v_{32} - v_{12} \quad (104)$$

## 6.11 Species Ste11

**Initial concentration** 158.33176608789 nmol · l<sup>-1</sup>

This species takes part in eight reactions (as a reactant in v12 and as a product in v13, v17, v21, v23, v25, v27, v32).

$$\frac{d}{dt}\text{Ste11} = v_{13} + v_{17} + v_{21} + v_{23} + v_{25} + v_{27} + v_{32} - v_{12} \quad (105)$$

### 6.12 Species `complexA`

**Initial concentration** 105.943298120207 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in `v13`, `v16` and as a product in `v12`).

$$\frac{d}{dt}\text{complexA} = v_{12} - v_{13} - v_{16} \quad (106)$$

### 6.13 Species `Ste7`

**Initial concentration** 36.3997016405141 nmol · l<sup>-1</sup>

This species takes part in eight reactions (as a reactant in `v14` and as a product in `v15`, `v17`, `v21`, `v23`, `v25`, `v27`, `v32`).

$$\frac{d}{dt}\text{Ste7} = v_{15} + v_{17} + v_{21} + v_{23} + v_{25} + v_{27} + v_{32} - v_{14} \quad (107)$$

### 6.14 Species `Fus3`

**Initial concentration** 686.399701640513 nmol · l<sup>-1</sup>

This species takes part in ten reactions (as a reactant in `v14`, `v29` and as a product in `v15`, `v17`, `v21`, `v23`, `v25`, `v27`, `v30`, `v33`).

$$\frac{d}{dt}\text{Fus3} = v_{15} + v_{17} + v_{21} + v_{23} + v_{25} + v_{27} + v_{30} + v_{33} - v_{14} - v_{29} \quad (108)$$

### 6.15 Species `complexB`

**Initial concentration** 77.8753625675829 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in `v15`, `v16` and as a product in `v14`).

$$\frac{d}{dt}\text{complexB} = v_{14} - v_{15} - v_{16} \quad (109)$$

### 6.16 Species `Ste20`

**Initial concentration** 1000 nmol · l<sup>-1</sup>

This species takes part in seven reactions (as a reactant in `v18` and as a product in `v19`, `v21`, `v23`, `v25`, `v27`, `v32`).

$$\frac{d}{dt}\text{Ste20} = v_{19} + v_{21} + v_{23} + v_{25} + v_{27} + v_{32} - v_{18} \quad (110)$$

### 6.17 Species `complexE`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in four reactions (as a reactant in [v19](#), [v20](#), [v21](#) and as a product in [v18](#)).

$$\frac{d}{dt}\text{complexE} = v_{18} - v_{19} - v_{20} - v_{21} \quad (111)$$

### 6.18 Species `complexF`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in [v22](#), [v23](#) and as a product in [v20](#)).

$$\frac{d}{dt}\text{complexF} = v_{20} - v_{22} - v_{23} \quad (112)$$

### 6.19 Species `complexG`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in [v24](#), [v25](#) and as a product in [v22](#)).

$$\frac{d}{dt}\text{complexG} = v_{22} - v_{24} - v_{25} \quad (113)$$

### 6.20 Species `complexH`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in [v26](#), [v27](#) and as a product in [v24](#)).

$$\frac{d}{dt}\text{complexH} = v_{24} - v_{26} - v_{27} \quad (114)$$

### 6.21 Species `complexI`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in [v28](#) and as a product in [v26](#), [v31](#)).

$$\frac{d}{dt}\text{complexI} = v_{26} + v_{31} - v_{28} \quad (115)$$

### 6.22 Species `complexL`

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in four reactions (as a reactant in [v29](#), [v32](#) and as a product in [v28](#), [v30](#)).

$$\frac{d}{dt}\text{complexL} = v_{28} + v_{30} - v_{29} - v_{32} \quad (116)$$

### 6.23 Species Fus3PP

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

This species takes part in six reactions (as a reactant in v33, v34 and as a product in v28, v35 and as a modifier in v39, v46).

$$\frac{d}{dt}\text{Fus3PP} = v_{28} + v_{35} - v_{33} - v_{34} \quad (117)$$

### 6.24 Species complexK

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

This species takes part in three reactions (as a reactant in v30, v31 and as a product in v29).

$$\frac{d}{dt}\text{complexK} = v_{29} - v_{30} - v_{31} \quad (118)$$

### 6.25 Species Ste12

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

This species takes part in two reactions (as a reactant in v34 and as a product in v35).

$$\frac{d}{dt}\text{Ste12} = v_{35} - v_{34} \quad (119)$$

### 6.26 Species Ste12a

**Name** Ste12active

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

This species takes part in three reactions (as a reactant in v35 and as a product in v34 and as a modifier in v36).

$$\frac{d}{dt}\text{Ste12a} = v_{34} - v_{35} \quad (120)$$

### 6.27 Species Bar1

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

This species takes part in two reactions (as a reactant in v36 and as a product in v37).

$$\frac{d}{dt}\text{Bar1} = v_{37} - v_{36} \quad (121)$$



### 6.28 Species Bar1a

**Name** Bar1active

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in v37, v38 and as a product in v36).

$$\frac{d}{dt}\text{Bar1a} = v_{36} - v_{37} - v_{38} \quad (122)$$

### 6.29 Species Bar1aex

**Name** Bar1activeEx

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in two reactions (as a product in v38 and as a modifier in v1).

$$\frac{d}{dt}\text{Bar1aex} = v_{38} \quad (123)$$

### 6.30 Species Far1

**Initial concentration** 500 nmol · l<sup>-1</sup>

This species takes part in three reactions (as a reactant in v39, v41 and as a product in v40).

$$\frac{d}{dt}\text{Far1} = v_{40} - v_{39} - v_{41} \quad (124)$$

### 6.31 Species Far1PP

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in six reactions (as a reactant in v40, v42, v45 and as a product in v39, v43, v44).

$$\frac{d}{dt}\text{Far1PP} = v_{39} + v_{43} + v_{44} - v_{40} - v_{42} - v_{45} \quad (125)$$

### 6.32 Species Far1U

**Name** Far1ubiquitin

**Initial concentration** 0 nmol · l<sup>-1</sup>

This species takes part in one reaction (as a product in v41).

$$\frac{d}{dt}\text{Far1U} = v_{41} \quad (126)$$

### 6.33 Species `complexM`

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

This species takes part in two reactions (as a reactant in [v43](#) and as a product in [v42](#)).

$$\frac{d}{dt}\text{complexM} = v_{42} - v_{43} \quad (127)$$

### 6.34 Species `complexN`

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

This species takes part in two reactions (as a reactant in [v44](#) and as a product in [v45](#)).

$$\frac{d}{dt}\text{complexN} = v_{45} - v_{44} \quad (128)$$

### 6.35 Species `Cdc28`

**Initial concentration**  $300 \text{ nmol} \cdot \text{l}^{-1}$

This species takes part in three reactions (as a reactant in [v45](#) and as a product in [v44](#) and as a modifier in [v41](#)).

$$\frac{d}{dt}\text{Cdc28} = v_{44} - v_{45} \quad (129)$$

### 6.36 Species `Sst2`

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

This species takes part in three reactions (as a reactant in [v47](#) and as a product in [v46](#) and as a modifier in [v8](#)).

$$\frac{d}{dt}\text{Sst2} = v_{46} - v_{47} \quad (130)$$

### 6.37 Species `p`

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

This species takes part in three reactions (as a reactant in [v46](#) and as a product in [v4](#), [v47](#)), which do not influence its rate of change because this species is on the boundary of the reaction system:

$$\frac{d}{dt}p = 0 \quad (131)$$

SBML<sup>2</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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