# **SBML Model Report**

# Model name: "Reyes-Palomares2012 - a combined model hepatic polyamine and sulfur aminoacid metabolism - version2"



May 5, 2016

#### 1 General Overview

This is a document in SBML Level 2 Version 4 format. This model was created by the following two authors: Armando Reyes-Palomares<sup>1</sup> and Vijayalakshmi Chelliah<sup>2</sup> at May nineth 2013 at 11:39 a.m. and last time modified at April fourth 2014 at 2:32 a.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	4
species types	0	species	53
events	0	constraints	0
reactions	74	function definitions	70
global parameters	178	unit definitions	2
rules	17	initial assignments	2

## **Model Notes**

Reyes-Palomares2012 - a combined model hepatic polyamine and sulfur aminoacid metabolism - version2

<sup>&</sup>lt;sup>1</sup>University of M?laga. CIBERER (Enfermedades Raras), armando@uma.es

<sup>&</sup>lt;sup>2</sup>EMBL-EBI, viji@ebi.ac.uk

Mammalian polyamine metabolism consists of a bi-cycle with two required entrances, omithine and S-adenosyl methionine (SAM), and several alternative exists. The relevant regulatory roles of the short half-life enzymes ornithine decarboxylase (ODC), S-adenosyl methione decarboxylase (SAMDC) and spermindine/spermine acetyl transferase (SSAT) in polyamine metabolism are well studied, and has been modelled here.

This model is described in the article: A combined model of hepatic polyamine and sulfur amino acid metabolism to analyze S-adenosyl methionine availability. Reyes-Palomares A, Montaez R, Snchez-Jimnez F, Medina MAAmino Acids, February 2012, Volume 42, Issue 2-3, pp 597-610

Abstract:

Many molecular details remain to be uncovered concerning the regulation of polyamine metabolism. A previous model of mammalian polyamine metabolism showed that S-adenosyl methionine availability could play a key role in polyamine homeostasis. To get a deeper insight in this prediction, we have built a combined model by integration of the previously published polyamine model and one-carbon and glutathione metabolism model, published by different research groups. The combined model is robust and it is able to achieve physiological steady-state values, as well as to reproduce the predictions of the individual models. Furthermore, a transition between two versions of our model with new regulatory factors added properly simulates the switch in methionine adenosyl transferase isozymes occurring when the liver enters in proliferative conditions. The combined model is useful to support the previous prediction on the role of S-adenosyl methionine availability in polyamine homeostasis. Furthermore, it could be easily adapted to get deeper insights on the connections of polyamines with energy metabolism.

#### Notes by the author:

This model combines BIOMD000000190 and BIOMD0000000268 from BioModels Database, both models include corrections respect to their originals publications.

To simulate a MATI/MATIII switch to MATII in proliferating liver:

We set to 0 the Vmax parameters of MATI and MATIII

We included MATII reaction equation.

We add a regulation factor dependent of SAM levels in ODC and SAMDCe rates of synthesis (66.5/[SAM]).

H2O2 was increased in a 50 % according to an initial state of proliferating and regenerating liver.

This model is hosted on BioModels Database and identified by: MODEL1305060001.

To cite BioModels Database, please use: BioModels Database: An enhanced, curated and annotated resourcefor published quantitative kinetic models.

To the extent possible under law, all copyright and related orneighbouring rights to this encoded model have been dedicated to the publicdomain worldwide. Please refer to CCO Public DomainDedication for more information.

#### 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 time

**Definition** 3600 s

#### 2.2 Unit substance

**Definition** µmol

#### 2.3 Unit volume

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

**Definition** 1

#### 2.4 Unit area

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

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

## 2.5 Unit length

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

**Definition** m

# 3 Compartments

This model contains four compartments.

Table 2: Properties of all compartments.

Id	Name	SBO	Spatial Dimensions	Size	Unit	Constant	Outside
blood	blood		3	1	litre		
cytosol	cytosol		3	1	litre	$\overline{\mathbf{Z}}$	
mito	mitochondrion		3	1	litre	$\overline{\mathbf{Z}}$	
cell	cell		3	1	litre	$\overline{\mathbb{Z}}$	

## 3.1 Compartment blood

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

Name blood

# **3.2 Compartment** cytosol

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

Name cytosol

# 3.3 Compartment mito

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

Name mitochondrion

## 3.4 Compartment cell

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

Name cell

# 4 Species

This model contains 53 species. The boundary condition of eleven of these species is set to true so that these species' amount cannot be changed by any reaction. Section 10 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-
	1.36.41		1 1-1		tion
b_met	b_Methionine	blood	$\mu$ mol·l <sup>-1</sup>		Z
b_ser	b_Serine	blood	$\mu$ mol·l <sup>-1</sup>		
b_gly	b_Glycine	blood	$\mu$ mol·l <sup>-1</sup>		
b_glu	b_Glutamate	blood	$\mu$ mol·l <sup>-1</sup>	$\Box$	
b_cys	b_Cysteine	blood	$\mu mol \cdot l^{-1}$	$\Box$	
b_gsg	b_GSSG	blood	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
b_gsh	b_GSH	blood	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
GAR	GAR	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
NADPH	NADPH	cytosol	$\mu mol \cdot l^{-1}$		
BET	Betaine	cytosol	$\mu mol \cdot l^{-1}$		
DUMP	dUMP	cytosol	$\mu mol \cdot l^{-1}$		
H202	H2O2	cytosol	$\mu mol \cdot l^{-1}$		
$c_{-}$ thf	c_THF	cytosol	$\mu mol \cdot l^{-1}$		
$c_5mf$	c_5-methyl-THF	cytosol	$\mu mol \cdot l^{-1}$		
$c_2cf$	c_5-10-methylene-THF	cytosol	$\mu mol \cdot l^{-1}$	$\Box$	
$c_{-}1cf$	c_5-10-methenyl-THF	cytosol	$\mu mol \cdot l^{-1}$		
$c_{-}10f$	$c_10$ -formyl-THF	cytosol	$\mu mol \cdot l^{-1}$		
$c_dhf$	c_DHF	cytosol	$\mu$ mol·l <sup>-1</sup>		
aic	AICAR	cytosol	$\mu$ mol· $1^{-1}$		
c_glu	c_Glutamate	cytosol	$\mu$ mol· $1^{-1}$		
c_cys	c_Cysteine	cytosol	$\mu mol \cdot l^{-1}$		

Id	Name	Compartment	Derived Unit	Constant	Boundary Condi- tion
glc	Glutamyl-Cysteine	cytosol	$\mu$ mol·l <sup>-1</sup>	$\Box$	
c_gly	c_Glycine	cytosol	$\mu \text{mol} \cdot l^{-1}$		
c_gsg	c_GSSG	cytosol	$\mu mol \cdot l^{-1}$	$\Box$	
c_gsh	c_GSH	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
cyt	Cystathionine	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
hcy	Homocysteine	cytosol	$\mu$ mol·l <sup>-1</sup>		
c_ser	c_Serine	cytosol	$\mu$ mol·l <sup>-1</sup>		
sah	SAH	cytosol	$\mu$ mol·l <sup>-1</sup>		
sam	SAM	cytosol	$\mu$ mol·l <sup>-1</sup>		
met	c_Methionine	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
C_C00	c_Formate	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
species_1	dcSAM	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
species_2	Putrescine	cytosol	$\mu mol \cdot l^{-1}$	$\Box$	
species_3	Spermine	cytosol	$\mu mol \cdot l^{-1}$	$\Box$	
species_4	Spermidine	cytosol	$\mu$ mol·l <sup>-1</sup>		
species_5	Acetylspermine	cytosol	$\mu mol \cdot l^{-1}$		
species_6	Acetylspermidine	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
species_7	Ornithine	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
species_8	Acetyl-CoA	cytosol	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\Box$	
species_9	CoA	cytosol	$\mu mol \cdot l^{-1}$	$\Box$	
CO	CO2	mito	$\mu mol \cdot l^{-1}$		
$\mathtt{m\_thf}$	$m_{-}THF$	mito	$\mu mol \cdot l^{-1}$		
$m_2cf$	m_5-10-methylene-THF	mito	$\mu mol \cdot l^{-1}$	$\Box$	
$m_1cf$	m_5-10-methenyl-THF	mito	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
$m_10f$	m_10-formyl-THF	mito	$\mu mol \cdot l^{-1}$	$\Box$	
${\tt m\_ser}$	m_Serine	mito	$\mu$ mol·l <sup>-1</sup>		
$m_{-}gly$	m_Glycine	mito	$\mu \text{mol} \cdot l^{-1}$		

Id	Name	Compartment	Derived Unit	Constant	Boundary Condi-
					tion
m_coo	m_Formate	mito	$\mu$ mol·l <sup>-1</sup>		$\Box$
Fol	Folate	cell	$\mu mol \cdot l^{-1}$	$\square$	$\square$
HCHO	Formaldehyde	cell	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$	$\square$	$\square$
src	Sarcosine	cell	$\mu \mathrm{mol} \cdot \mathrm{l}^{-1}$		
dmg	Dimethylglycine	cell	$\mu mol \cdot l^{-1}$		$\Box$

# **5 Parameters**

This model contains 178 global parameters.

Table 4: Properties of each parameter.

Id	Name	SBO	Value	Unit	Constant
tot_cfol	tot_cfol		13.400		
tot_mfol	tot_mfol		40.200		
V_oCys_b	V_oCys_b		70.000		
$V_{\tt oGly\_b}$	$V_oGly_b$		630.000		
$V_{\tt oGlu\_b}$	V_oGlu_b		273.000		
$V_{\mathtt{gshHb}}$	$V_gshHb$		150.000		
${ t K\_gshHb}$	$K_gshHb$		150.000		$   \overline{\checkmark} $
$V_{\mathtt{gshLb}}$	$V_g sh Lb$		1100.000		$\overline{\mathbf{Z}}$
$h\_{\sf gshLb}$	$h\_gshLb$		3.000		$\overline{\mathbf{Z}}$
$K_{\mathtt{gshLb}}$	$K_gshLb$		3000.000		
$V_{\mathtt{gsgHb}}$	$V_{-}gsgHb$		40.000		
${ t K\_gsgHb}$	$K_gsgHb$		1250.000		$\overline{\mathbf{Z}}$
$V_{\mathtt{gsgLb}}$	$V_{gsgLb}$		4025.000		
$K_{-}$ gsg $L$ b	$K_{-}gsgLb$		7100.000		$\overline{\mathbf{Z}}$
$V_{ extsf{b}}$ cysc	V_bcysc		14950.000		$\overline{\mathbf{Z}}$
${\tt K\_bcysc}$	K_bcysc		2100.000		$\overline{\mathbf{Z}}$
$V_{ extbf{bglutc}}$	V_bglutc		28000.000		$\overline{\mathbf{Z}}$
${ t K\_bglutc}$	K_bglutc		300.000		
$k\_out\_glu$	k_out_glu		1.000		
${\tt V\_bglyc}$	$V_bglyc$		4600.000		
${\tt K\_bglyc}$	K_bglyc		150.000		
$k\_{\tt out\_gly}$	k_out_gly		1.000		
$V_{\mathtt{bserc}}$	V_bserc		2700.000		
${\tt K\_bserc}$	K_bserc		150.000		
$k\_out\_ser$	k_out_ser		1.000		
${\tt V\_bmetc}$	V_bmetc		913.400		
${\tt K\_bmetc}$	K_bmetc		150.000		
$k\_{\tt out\_met}$	$k_out_met$		1.000		
$Vm\_mFTD$	Vm_mFTD		1050.000		
$K_10f_FTD$	K_10f_FTD		20.000		
${\tt Vf\_mSHMT}$	$Vf_mSHMT$		11440.000		
$K_{thf\_SHMT}$	K_thf_SHMT		50.000		$\mathbf{Z}$
$K\_ser\_SHMT$	K_ser_SHMT		600.000		$   \overline{\mathbf{Z}} $
${\tt Vr\_mSHMT}$	$Vr\_mSHMT$		$3 \cdot 10^7$		$   \overline{\mathbf{Z}} $
$K_gly_SHMT$	$K_gly_SHMT$		10000.000		$\overline{\mathbf{Z}}$
K_2cf_SHMT	K_2cf_SHMT		3200.000		$\overline{\mathbf{Z}}$
Vf_mFTS	Vf_mFTS		2000.000		$\overline{\mathbb{Z}}$

Id	Name	SBO	Value	Unit	Constant
K_thf_mFTS	K_thf_mFTS		3.000		Ø
K_coo_mFTS	K_coo_mFTS		43.000		$\overline{Z}$
$Vr\_mFTS$	$Vr\_mFTS$		6300.000		$\overline{\mathbf{Z}}$
$K_10f_mFTS$	$K_10f_mFTS$		22.000		$\overline{\mathbf{Z}}$
k1_mNE	k1_mNE		0.030		<u> </u>
k2_mNE	k2_mNE		20.000		$\overline{\mathbf{Z}}$
$Vm\_GDC$	$Vm\_GDC$		15000.000		$\overline{\mathbf{Z}}$
$K_{thf_GDC}$	K_thf_GDC		50.000		$\overline{\mathbf{Z}}$
$K_gly_GDC$	K_gly_GDC		3400.000		$ \overline{\mathbf{Z}} $
$Vm\_SDH$	VmSDH		15000.000		$\overline{\mathbf{Z}}$
$K_{thf\_SDH}$	K_thf_SDH		50.000		$\overline{\mathbf{Z}}$
$K\_src\_SDH$	K_src_SDH		320.000		$\overline{\mathbf{Z}}$
$Vm\_DMGD$	Vm_DMGD		15000.000		$\overline{\mathbf{Z}}$
$K_{thf_DMGD}$	K_thf_DMGD		50.000		$\overline{\mathbf{Z}}$
$K_dmg_DMGD$	K_dmg_DMGD		50.000		$\overline{\mathbf{Z}}$
Vf_mMTD	Vf_mMTD		180000.000		$\overline{\mathbf{Z}}$
K_2cf_MTD	K_2cf_MTD		2.000		$\overline{\mathbb{Z}}$
$Vr\_MTD$	$Vr\_MTD$		594000.000		$\overline{\mathbf{Z}}$
$K_1cf_MTD$	K_1cf_MTD		10.000		$\overline{\mathbf{Z}}$
Vf_mMTCH	Vf_mMTCH		790000.000		$\overline{\mathbf{Z}}$
$K_1cf_MTCH$	K_1cf_MTCH		250.000		$ \overline{\checkmark} $
$Vr\_MTCH$	Vr_MTCH		20000.000		$ \overline{\checkmark} $
$K_10f_MTCH$	K_10f_MTCH		100.000		
V_mser	V_mser		10000.000		$ \overline{\checkmark} $
K_mser	K_mser		5700.000		$ \overline{\mathbf{Z}} $
$V_{\tt cser}$	V_cser		10000.000		$ \overline{\mathbf{Z}} $
$K_{\mathtt{cser}}$	K_cser		5700.000		
k_in_coo	k_in_coo		100.000		
k_out_coo	k_out_coo		100.000		
$V_{mgly}$	$V_{mgly}$		10000.000		
$K_mgly$	$K_mgly$		5700.000		$ \overline{\checkmark} $
$V_{\tt cgly}$	V_cgly		10000.000		$ \overline{\mathbf{Z}} $
$K_{\mathtt{cgly}}$	K_cgly		5700.000		
$Vm\_MS$	$Vm\_MS$		500.000		
$K_5mf_MS$	K_5mf_MS		25.000		
$K_hcy_MS$	K_hcy_MS		1.000		$\overline{\mathbf{Z}}$
ssH2O2	ssH2O2		0.010		$   \overline{\checkmark} $
Ki_MS	Ki_MS		0.010		$\overline{\mathbf{Z}}$
Vm_DHFR	Vm_DHFR		2000.000		$\overline{\mathbb{Z}}$
$K_dhf_DHFR$	K_dhf_DHFR		0.500		$\overline{\mathbb{Z}}$
K_NADPH_DHFR	K_NADPH_DHFR		4.000		$\overline{\mathbf{Z}}$
${\tt Vm\_cFTD}$	Vm_cFTD		500.000		$\overline{\mathbf{Z}}$

Id	Name	SBO Value	Unit	Constant
Vm_PGT	Vm_PGT	24300.0	000	<b>✓</b>
$K_10f_PGT$	K_10f_PGT	4.9	900	
$K\_GAR\_PGT$	K_GAR_PGT	520.0	000	
${\tt Vm\_cFTS}$	Vm_cFTS	3900.0	000	
$K_{thf_cFTS}$	K_thf_cFTS	3.0	000	$   \overline{\mathscr{A}} $
$K_{coo_cFTS}$	K_coo_cFTS	43.0	000	
$Vf_cSHMT$	Vf_cSHMT	5200.0	000	
${\tt Vr\_cSHMT}$	Vr_cSHMT	1.5 · 1	$10^{7}$	
$k1_cNE$	k1_cNE	0.0	030	$ \mathbf{Z} $
$k2$ _cNE	k2₋cNE	22.0	000	$\overline{\mathbf{Z}}$
$Vm_{-}TS$	$Vm_{-}TS$	5000.0	000	$\overline{\mathbf{Z}}$
K_DUMP_TS	K_DUMP_TS	6.3	300	$\overline{\mathbf{Z}}$
$K_2cf_TS$	K_2cf_TS	14.0	000	$\overline{\mathbf{Z}}$
Vm_MTHFR	Vm_MTHFR	5300.0	000	$\overline{\mathbf{Z}}$
K_2cf_MTHFR	K_2cf_MTHFR	50.0	000	$\overline{\mathbf{Z}}$
K_NADPH-	K_NADPH-	16.0	000	$\overline{\mathbf{Z}}$
_MTHFR	_MTHFR			
$Vf\_cMTD$	Vf_cMTD	80000.0	000	
$Vr\_cMTD$	Vr_cMTD	600000.0	000	$\overline{\mathscr{L}}$
Vf_cMTCH	Vf_cMTCH	500000.0	000	$\overline{\mathscr{L}}$
$Vm\_ART$	$Vm\_ART$	55000.0	000	$\overline{\mathbf{Z}}$
K_10f_ART	$K_10f_ART$	5.9	900	$\overline{\mathbf{Z}}$
$K_aic_ART$	K_aic_ART	100.0	000	$\overline{\mathscr{A}}$
$Vm\_BHMT$	$Vm_BHMT$	2160.0	000	$\overline{\mathbf{Z}}$
K_hcy_BHMT	K_hcy_BHMT	12.0	000	$\mathbf{Z}$
K_bet_BHMT	K_bet_BHMT	100.0	000	$\overline{\mathscr{L}}$
Ki_BHMT	Ki_BHMT	0.0	010	$\overline{\mathbf{Z}}$
Vm_MAT1	Vm_MAT1	0.0	000	$\overline{\mathbf{Z}}$
Km_MAT1	Km_MAT1	41.0	000	$\overline{\mathbf{Z}}$
Ki_MAT1	Ki_MAT1	2140.0	000	$\overline{\mathscr{A}}$
Vm_MAT3	Vm_MAT3	0.0	000	$\overline{\mathscr{L}}$
Km_MAT3	Km_MAT3	300.0	000	$\mathbf{Z}$
Ka_MAT3	Ka_MAT3	360.0	000	$\overline{\mathbf{Z}}$
Ki_MAT3	Ki_MAT3	4030.0	000	$\overline{\mathbf{Z}}$
$Vm\_GNMT$	$Vm_{-}GNMT$	260.0	000	$\overline{\mathbf{Z}}$
$K_sam_GNMT$	K_sam_GNMT	63.0	000	<b>Z</b>
$K_gly_GNMT$	K_gly_GNMT	130.0	000	$\mathbf{Z}$
$Ki_GNMT$	Ki_GNMT	18.0	000	<b>Z</b>
$Vm_DNMT$	Vm_DNMT	180.0	000	<b>Z</b>
Km_DNMT	Km_DNMT	1.4	100	<b>Z</b>
Ki_DNMT	Ki_DNMT	1.4	100	<b>Z</b>
Vf_SAHH	Vf_SAHH	320.0	000	$\overline{\mathbf{Z}}$

Id	Name	SBO	Value	Unit	Constant
$K_sah_SAHH$	K_sah_SAHH		6.500		
${\tt Vr\_SAHH}$	Vr_SAHH		4530.000		$\overline{\mathbf{Z}}$
$K_hcy_SAHH$	K_hcy_SAHH		150.000		
$Vm_CBS$	Vm_CBS	42	20000.000		$\overline{\mathscr{L}}$
$K_hcy_CBS$	K_hcy_CBS		1000.000		$\overline{\mathscr{L}}$
$K\_ser\_CBS$	K_ser_CBS		2000.000		$   \overline{\mathscr{L}} $
$Ka\_CBS$	Ka_CBS		0.035		$   \overline{\mathscr{L}} $
${\tt Vm\_CTGL}$	Vm_CTGL		1500.000		$ \overline{\checkmark} $
$K_{\mathtt{cyt}}$	K_cyt_CTGL		500.000		
$Vm\_GCS$	$Vm\_GCS$		3600.000		$   \overline{\mathscr{L}} $
$Ke\_GCS$	Ke_GCS		5597.000		$   \overline{\mathscr{L}} $
$K_{cys}_{GCS}$	K_cys_GCS		100.000		$   \overline{\mathscr{L}} $
$K_glu_GCS$	K_glu_GCS		1900.000		
${\tt Ki\_GCS}$	Ki_GCS		8200.000		$   \overline{\mathscr{L}} $
${\tt Kp\_GCS}$	Kp_GCS		300.000		
$Ka\_GCS$	Ka_GCS		0.010		$   \overline{\mathscr{L}} $
${\tt Vm\_GS}$	$Vm_{-}GS$		5400.000		$   \overline{\mathscr{L}} $
$Ke\_GS$	Ke_GS		5600.000		$   \overline{\mathscr{L}} $
$K_gly_GS$	K_gly_GS		300.000		$   \overline{\mathscr{L}} $
$K_glc_GS$	K_glc_GS		22.000		$   \overline{\mathscr{L}} $
${\tt Kp\_GS}$	Kp_GS		30.000		
$Vm\_GPX$	$Vm_{-}GPX$		4500.000		
${ t K\_gsh\_GPX}$	K_gsh_GPX		1330.000		
$K_H202_GPX$	K_H2O2_GPX		0.090		$   \overline{\mathscr{L}} $
$Vm_{-}GR$	$Vm_{-}GR$		892.500		$   \overline{\mathscr{L}} $
$K_{gsg}GR$	K_gsg_GR		107.000		$ \overline{\checkmark} $
$K_NADPH_GR$	K_NADPH_GR		10.400		$\mathbf{Z}$
dinner	dinner		1.000		$\mathbf{Z}$
lunch	lunch		1.000		$\mathbf{Z}$
breakfast	breakfast		1.000		$\mathbf{Z}$
fasting	fasting		1.000		$\mathbf{Z}$
daytime	daytime		0.000		
$\mathtt{aa\_input}$	Aminoacid_input		1.000		
b_met_basal	b_met_basal		30.000		
$b_ser_basal$	b_ser_basal		150.000		
$V_oGly_b-$	V_oGly_b_basal		630.000		$\mathbf{Z}$
_basal					
V_oGlu_b- _basal	V_oGlu_b_basal		273.000		$\square$
V_oCys_b- _basal	V_oCys_b_basal		70.000		
k_out_cys	k_out_cys		1.000		

Id	Name	SBO	Value	Unit	Constant
parameter_1	Vm_ODC		72.256		
$parameter_2$	$Vm_SSAT$		42.285		
$parameter_3$	$Vm\_SAMDC$		21.134		
$parameter_4$	ANTZ		0.585		
$parameter_5$	Ke_POL		1.000		
$parameter_6$	Kd_ODC		3.000		
$parameter_7$	Ks_ODC		300.000		$\overline{\mathbf{Z}}$
$parameter_8$	Kd_SSAT		12.000		
$parameter_9$	Ks_SSAT		0.060		
$parameter_10$	Kd_SAMDC		1.200		$\overline{\checkmark}$
$parameter_11$	Ks_SAMDC		60.000		$\overline{\mathbf{Z}}$
$parameter_12$	Kd_ANTZ		0.020		$\overline{\mathbf{Z}}$
$parameter_13$	Ks_ANTZ		0.020		$\overline{\mathbf{Z}}$
$parameter_14$	K_accoa		0.240		
$parameter_15$	K_coa		0.720		
parameter_16	R		0.240		
parameter_17	S+D		140.951		
parameter_18	R_percent		100.000		
parameter_19	Vm_MATII		220.000		

# 6 Initialassignments

This is an overview of two initialassignments.

## **6.1 Initialassignment** c\_thf

Derived unit contains undeclared units

$$\label{eq:math} \mbox{Math} \ \ \frac{[Fol]}{\frac{2\cdot 3}{4}} - \left( [c\_5mf] + [c\_2cf] + [c\_1cf] + [c\_10f] + [c\_dhf] \right)$$

## 6.2 Initialassignment m\_thf

Derived unit contains undeclared units

$$\label{eq:math} \mbox{Math} \ \ \frac{[Fol]}{\frac{2\cdot 1}{4}} - \left( [m\_2cf] + [m\_1cf] + [m\_10f] \right)$$

## 7 Function definitions

This is an overview of 70 function definitions.

## 7.1 Function definition function\_4\_b\_glu\_loss\_1

 $\textbf{Name} \ \ function\_4\_b\_glu\_loss\_1$ 

Argument [b\_glu]

## **Mathematical Expression**

$$0.1 \cdot [b\_glu] \tag{1}$$

## 7.2 Function definition function\_1

Name Constant flux (reversible)

Argument v

## **Mathematical Expression**

#### 7.3 Function definition function\_4\_VmHCOOHc\_1

Name function\_4\_VmHCOOHc\_1

**Arguments** [c\_coo], vol (cytosol), k\_in\_coo, k\_out\_coo, [m\_coo], vol (mito)

## **Mathematical Expression**

$$\frac{k\_in\_coo \cdot [m\_coo] \cdot vol \, (mito)}{3} - k\_out\_coo \cdot [c\_coo] \cdot vol \, (cytosol) \tag{3}$$

## **7.4 Function definition** function\_4\_b\_cys\_cystine\_conv\_1

Name function\_4\_b\_cys\_cystine\_conv\_1

**Argument** [b\_cys]

#### **Mathematical Expression**

$$0.25 \cdot [b\_cys] \tag{4}$$

## **7.5 Function definition** function\_4\_b\_gsg\_decomp\_1

Name function\_4\_b\_gsg\_decomp\_1

Argument [b\_gsg]

$$67.5 \cdot [b\_gsg] \tag{5}$$

## **7.6 Function definition** function\_4\_b\_cys\_loss\_1

Name function\_4\_b\_cys\_loss\_1

**Argument** [b\_cys]

## **Mathematical Expression**

$$0.1 \cdot [b\_cys] \tag{6}$$

## 7.7 Function definition function\_4\_b\_gsh\_decomp\_1

Name function\_4\_b\_gsh\_decomp\_1

Argument [b\_gsh]

## **Mathematical Expression**

$$90 \cdot [b\_gsh] \tag{7}$$

#### 7.8 Function definition function\_4\_V\_GDC\_1

Name function\_4\_V\_GDC\_1

Arguments K\_gly\_GDC, K\_thf\_GDC, Vm\_GDC, [m\_gly], [m\_thf]

#### **Mathematical Expression**

$$\frac{Vm\_GDC \cdot [m\_thf] \cdot [m\_gly]}{(K\_thf\_GDC + [m\_thf]) \cdot (K\_gly\_GDC + [m\_gly])} \tag{8}$$

## **7.9 Function definition** function\_4\_b\_gsg\_loss\_1

Name function\_4\_b\_gsg\_loss\_1

Argument [b\_gsg]

#### **Mathematical Expression**

$$7.5 \cdot [b_{-}gsg] \tag{9}$$

#### 7.10 Function definition function\_4\_VmNE\_1

Name function\_4\_VmNE\_1

**Arguments** [HCHO], k1\_mNE, k2\_mNE, [m\_2cf], [m\_thf], vol (mito)

$$vol(mito) \cdot (k1\_mNE \cdot [m\_thf] \cdot [HCHO] - k2\_mNE \cdot [m\_2cf])$$
 (10)

## **7.11 Function definition** function\_4\_b\_gsh\_loss\_1

Name function\_4\_b\_gsh\_loss\_1

Argument [b\_gsh]

**Mathematical Expression** 

$$0.7 \cdot [b\_gsh] \tag{11}$$

## **7.12 Function definition** function\_4\_b\_gly\_loss\_1

Name function\_4\_b\_gly\_loss\_1

Argument [b\_gly]

**Mathematical Expression** 

$$0.1 \cdot [b\_gly] \tag{12}$$

#### 7.13 Function definition function\_4\_VmFTD\_1

Name function\_4\_VmFTD\_1

Arguments K\_10f\_FTD, Vm\_mFTD, [m\_10f]

**Mathematical Expression** 

$$\frac{Vm\_mFTD \cdot [m\_10f]}{K\_10f\_FTD + [m\_10f]}$$
 (13)

#### 7.14 Function definition MM\_twosubst

Name MM\_twosubst

Arguments Vmax, Km1, Km2, S1, S2

**Mathematical Expression** 

$$\frac{Vmax \cdot S1 \cdot S2}{(Km1 + S1) \cdot (Km2 + S2)} \tag{14}$$

#### 7.15 Function definition MM

Name MM

 $\textbf{Arguments}\ \ Vmax,\,Km,\,S$ 

$$\frac{V \max \cdot S}{K m + S} \tag{15}$$

## **7.16 Function definition** function\_4\_V\_c\_gshLb\_1

Name function\_4\_V\_c\_gshLb\_1

Arguments K\_gshLb, V\_gshLb, [c\_gsh], vol(cytosol), h\_gshLb

#### **Mathematical Expression**

$$\frac{\text{vol}\left(\text{cytosol}\right) \cdot \text{V\_gshLb} \cdot [\text{c\_gsh}]^{\text{h\_gshLb}}}{\text{K\_gshLb}^{\text{h\_gshLb}} + [\text{c\_gsh}]^{\text{h\_gshLb}}} \tag{16}$$

#### 7.17 Function definition function\_4\_V\_MS\_1

Name function\_4\_V\_MS\_1

**Arguments** [H2O2], K\_5mf\_MS, K\_hcy\_MS, Ki\_MS, Vm\_MS, [c\_5mf], [hcy], ssH2O2

#### **Mathematical Expression**

$$\frac{MM\_twosubst(Vm\_MS,K\_5mf\_MS,K\_hcy\_MS,[c\_5mf],[hcy])\cdot(ssH2O2+Ki\_MS)}{[H2O2]+Ki\_MS}$$
 (17)

#### 7.18 Function definition function\_4\_V\_DHFR\_1

Name function\_4\_V\_DHFR\_1

Arguments K\_NADPH\_DHFR, K\_dhf\_DHFR, [NADPH], Vm\_DHFR, [c\_dhf]

## **Mathematical Expression**

$$\frac{Vm\_DHFR \cdot [c\_dhf] \cdot [NADPH]}{(K\_dhf\_DHFR + [c\_dhf]) \cdot (K\_NADPH\_DHFR + [NADPH])}$$
(18)

#### **7.19 Function definition** function\_4\_VcFTD\_1

Name function\_4\_VcFTD\_1

**Arguments** K\_10f\_FTD, Vm\_cFTD, [c\_10f]

$$\frac{Vm\_cFTD \cdot [c\_10f]}{K\_10f\_FTD + [c\_10f]}$$
(19)

## **7.20 Function definition** function\_4\_V\_PGT\_1

Name function\_4\_V\_PGT\_1

Arguments [GAR], K\_10f\_PGT, K\_GAR\_PGT, Vm\_PGT, [c\_10f]

#### **Mathematical Expression**

$$\frac{Vm\_PGT \cdot [c\_10f] \cdot [GAR]}{(K\_10f\_PGT + [c\_10f]) \cdot (K\_GAR\_PGT + [GAR])}$$
(20)

#### 7.21 Function definition function\_4\_VcFTS\_1

Name function\_4\_VcFTS\_1

Arguments K\_coo\_cFTS, K\_thf\_cFTS, Vm\_cFTS, [c\_coo], [c\_thf]

#### **Mathematical Expression**

$$\frac{Vm\_cFTS \cdot [c\_thf] \cdot [c\_coo]}{(K\_thf\_cFTS + [c\_thf]) \cdot (K\_coo\_cFTS + [c\_coo])}$$
(21)

#### **7.22 Function definition** function\_4\_VcSHMT\_1

Name function\_4\_VcSHMT\_1

**Arguments** K\_2cf\_SHMT, K\_gly\_SHMT, K\_ser\_SHMT, K\_thf\_SHMT, Vf\_cSHMT, Vr\_cSHMT, [c\_2cf], [c\_gly], [c\_ser], [c\_thf]

## **Mathematical Expression**

#### **7.23 Function definition** function\_4\_VcNE\_1

Name function\_4\_VcNE\_1

Arguments [HCHO], [c\_2cf], [c\_thf], vol (cytosol), k1\_cNE, k2\_cNE

$$vol(cytosol) \cdot (k1\_cNE \cdot [c\_thf] \cdot [HCHO] - k2\_cNE \cdot [c\_2cf])$$
 (23)

## **7.24 Function definition** function\_4\_V\_TS\_1

Name function\_4\_V\_TS\_1

Arguments [DUMP], K\_2cf\_TS, K\_DUMP\_TS, Vm\_TS, [c\_2cf]

#### **Mathematical Expression**

$$\frac{Vm\_TS \cdot [DUMP] \cdot [c\_2cf]}{(K\_DUMP\_TS + [DUMP]) \cdot (K\_2cf\_TS + [c\_2cf])}$$
(24)

#### **7.25 Function definition** function 4 V MTHFR 1

Name function\_4\_V\_MTHFR\_1

Arguments K\_2cf\_MTHFR, K\_NADPH\_MTHFR, [NADPH], Vm\_MTHFR, [c\_2cf], [sah], [sam]

## **Mathematical Expression**

$$\frac{\text{MM\_twosubst}(\text{Vm\_MTHFR}, \text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, [\text{c\_2cf}], [\text{NADPH}]) \cdot 72}{10 + [\text{sam}] - [\text{sah}]} \tag{25}$$

#### **7.26 Function definition** function\_4\_VcMTD\_1

Name function\_4\_VcMTD\_1

Arguments K\_1cf\_MTD, K\_2cf\_MTD, Vf\_cMTD, Vr\_cMTD, [c\_1cf], [c\_2cf]

#### **Mathematical Expression**

$$MM(Vf\_cMTD, K\_2cf\_MTD, [c\_2cf]) - MM(Vr\_cMTD, K\_1cf\_MTD, (26)$$

$$[c\_1cf])$$

## 7.27 Function definition function\_4\_VcMTCH\_1

Name function\_4\_VcMTCH\_1

Arguments K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_cMTCH, Vr\_MTCH, [c\_10f], [c\_1cf]

$$MM\left(Vf\_cMTCH,K\_1cf\_MTCH,[c\_1cf]\right) - MM\left(Vr\_MTCH,K\_10f\_MTCH,[c\_10f]\right) \tag{27}$$

#### 7.28 Function definition function\_4\_V\_ART\_1

Name function\_4\_V\_ART\_1

Arguments K\_10f\_ART, K\_aic\_ART, Vm\_ART, [aic], [c\_10f]

#### **Mathematical Expression**

$$\frac{Vm\_ART \cdot [c\_10f] \cdot [aic]}{(K\_10f\_ART + [c\_10f]) \cdot (K\_aic\_ART + [aic])}$$
(28)

#### 7.29 Function definition function\_4\_V\_BHMT\_1

Name function\_4\_V\_BHMT\_1

Arguments [BET], [H2O2], K\_bet\_BHMT, K\_hcy\_BHMT, Ki\_BHMT, Vm\_BHMT, vol (cytosol), [hcy], [sah], [sam], ssH2O2

#### **Mathematical Expression**

$$vol (cytosol) \cdot exp (0.0021 \cdot ([sam] + [sah])) \cdot exp (0.0021 \cdot 102.6)$$

$$\cdot MM\_twosubst (Vm\_BHMT, K\_hcy\_BHMT,$$

$$K\_bet\_BHMT, [hcy], [BET]) \cdot \frac{ssH2O2 + Ki\_BHMT}{[H2O2] + Ki\_BHMT}$$

$$(29)$$

## 7.30 Function definition function\_4\_V\_MATI\_1

Name function\_4\_V\_MATI\_1

Arguments Ki\_MAT1, Km\_MAT1, Vm\_MAT1, [c\_gsg], [met], [sam]

## **Mathematical Expression**

$$Vm\_MAT1 \cdot \frac{[met]}{Km\_MAT1 + [met]} \cdot (0.23 + 0.8 \cdot exp (0.0026 \cdot [sam])) \cdot \frac{Ki\_MAT1 + 66.71}{Ki\_MAT1 + [c\_gsg]} \tag{30}$$

#### 7.31 Function definition function\_4\_V\_MATIII\_1

Name function\_4\_V\_MATIII\_1

**Arguments** Ka\_MAT3, Ki\_MAT3, Km\_MAT3, Vm\_MAT3, [c\_gsg], [met], [sam]

$$Vm\_MAT3 \cdot \frac{[met]^{1.21}}{Km\_MAT3 + [met]^{1.21}} \cdot \left(1 + \frac{7.2 \cdot [sam]^2}{Ka\_MAT3^2 + [sam]^2}\right) \cdot \frac{Ki\_MAT3 + 66.71}{Ki\_MAT3 + [c\_gsg]} \tag{31}$$

#### **7.32 Function definition** function\_4\_V\_GNMT\_1

Name function\_4\_V\_GNMT\_1

 $\label{eq:conditional} \mbox{\bf Arguments} \ \ \mbox{\bf K\_gly\_GNMT}, \mbox{\bf K\_sam\_GNMT}, \mbox{\bf Ki\_GNMT}, \mbox{\bf Vm\_GNMT}, \mbox{\bf [c\_5mf]}, \mbox{\bf [c\_gly]}, \mbox{\bf vol} \mbox{\bf (cytosol)}, \mbox{\bf [sah]}, \mbox{\bf [sam]}$ 

#### **Mathematical Expression**

$$\begin{array}{l} vol\left(cytosol\right) \\ \cdot MM\_twosubst\left(Vm\_GNMT, K\_sam\_GNMT, K\_gly\_GNMT, [sam], [c\_gly]\right) \\ \cdot \frac{1}{1 + \frac{[sah]}{Ki\_GNMT}} \cdot \frac{4.8}{0.35 + [c\_5mf]} \\ \end{array}$$

#### 7.33 Function definition function\_4\_V\_DNMT\_1

Name function\_4\_V\_DNMT\_1

Arguments Ki\_DNMT, Km\_DNMT, Vm\_DNMT, [sah], [sam]

## **Mathematical Expression**

$$Vm\_DNMT \cdot \frac{[sam]}{Km\_DNMT \cdot \left(1 + \frac{[sah]}{Ki\_DNMT}\right) + [sam]} \tag{33}$$

## 7.34 Function definition function\_4\_V\_SAHH\_1

Name function\_4\_V\_SAHH\_1

**Arguments** K\_hcy\_SAHH, K\_sah\_SAHH, Vf\_SAHH, Vr\_SAHH, [hcy], [sah]

#### **Mathematical Expression**

$$MM(Vf\_SAHH, K\_sah\_SAHH, [sah]) - MM(Vr\_SAHH, K\_hcy\_SAHH, (34) [hcy])$$

#### **7.35 Function definition** function\_4\_gluconeogenesis\_ser\_1

Name function\_4\_gluconeogenesis\_ser\_1

**Argument** [c\_ser]

$$1.2 \cdot [c\_ser] \tag{35}$$

#### 7.36 Function definition function\_4\_V\_CBS\_1

Name function\_4\_V\_CBS\_1

Arguments [H2O2], K\_hcy\_CBS, K\_ser\_CBS, Ka\_CBS, Vm\_CBS, [c\_ser], [hcy], [sah], [sam], ssH2O2

#### **Mathematical Expression**

$$\begin{aligned} & \text{MM\_twosubst} \left( \text{Vm\_CBS}, \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, [\text{hcy}], [\text{c\_ser}] \right) \\ & \cdot \frac{\left( \frac{30}{102.59} \right)^2 + 1}{\left( \frac{30}{[\text{sam}] + [\text{sah}]} \right)^2 + 1} \cdot \frac{[\text{H2O2}] + \text{Ka\_CBS}}{\text{ssH2O2} + \text{Ka\_CBS}} \end{aligned}$$

#### 7.37 Function definition function\_4\_V\_CTGL\_1

Name function\_4\_V\_CTGL\_1

**Arguments** K\_cyt\_CTGL, Vm\_CTGL, [cyt]

## **Mathematical Expression**

$$\frac{Vm\_CTGL \cdot [cyt]}{K\_cyt\_CTGL + [cyt]}$$
(37)

#### 7.38 Function definition function\_4\_V\_GCS\_1

Name function\_4\_V\_GCS\_1

**Arguments** [H2O2], K\_cys\_GCS, K\_glu\_GCS, Ka\_GCS, Ke\_GCS, Ki\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gsh], [glc], ssH2O2

#### **Mathematical Expression**

$$\frac{Vm\_GCS \cdot \left( [c\_cys] \cdot [c\_glu] - \frac{[glc]}{Kc\_GCS} \right)}{K\_cys\_GCS \cdot K\_glu\_GCS + [c\_glu] \cdot K\_cys\_GCS + [c\_cys] \cdot \left( K\_glu\_GCS \cdot \left( 1 + \frac{[c\_gsh]}{Ki\_GCS} \right) + [c\_glu] \right) + \frac{[glc]}{Kp\_GCS} + \frac{[c\_glu]}{Kp\_GCS} + \frac{[c\_gl$$

#### 7.39 Function definition function\_4\_V\_GS\_1

Name function\_4\_V\_GS\_1

Arguments K\_glc\_GS, K\_gly\_GS, Ke\_GS, Kp\_GS, Vm\_GS, [c\_gly], [c\_gsh], [glc]

#### **Mathematical Expression**

$$\frac{Vm\_GS \cdot \left( [c\_gly] \cdot [glc] - \frac{[c\_gsh]}{Ke\_GS} \right)}{K\_gly\_GS \cdot K\_glc\_GS + [glc] \cdot K\_gly\_GS + [c\_gly] \cdot \left( K\_glc\_GS + [glc] \right) + \frac{[c\_gsh]}{Kp\_GS}}$$

# **7.40 Function definition** function\_4\_c\_glu\_usage\_1

Name function\_4\_c\_glu\_usage\_1

Argument [c\_glu]

#### **Mathematical Expression**

$$0.07 \cdot [c_{glu}] \tag{40}$$

## **7.41 Function definition** function\_4\_cys\_usage\_1

Name function\_4\_cys\_usage\_1

Argument [c\_cys]

#### **Mathematical Expression**

$$\frac{0.35 \cdot [c\_cys]^2}{200} \tag{41}$$

#### 7.42 Function definition function\_4\_V\_GPX\_1

Name function\_4\_V\_GPX\_1

Arguments [H2O2], K\_H2O2\_GPX, K\_gsh\_GPX, Vm\_GPX, [c\_gsh]

#### **Mathematical Expression**

$$Vm\_GPX \cdot \left(\frac{[c\_gsh]}{K\_gsh\_GPX + [c\_gsh]}\right)^2 \cdot \frac{[H2O2]}{K\_H2O2\_GPX + [H2O2]} \tag{42}$$

#### 7.43 Function definition function\_4\_V\_GR\_1

Name function\_4\_V\_GR\_1

Arguments K\_NADPH\_GR, K\_gsg\_GR, [NADPH], Vm\_GR, [c\_gsg]

#### **Mathematical Expression**

22

$$\frac{Vm\_GR \cdot [c\_gsg] \cdot [NADPH]}{(K\_gsg\_GR + [c\_gsg]) \cdot (K\_NADPH\_GR + [NADPH])}$$
(43)

## **7.44 Function definition** function\_4\_c\_gsh\_degr\_1

Name function\_4\_c\_gsh\_degr\_1

**Argument** [c\_gsh]

#### **Mathematical Expression**

$$0.0020 \cdot [c\_gsh] \tag{44}$$

## **7.45 Function definition** function\_4\_c\_gsg\_degr\_1

Name function\_4\_c\_gsg\_degr\_1

Argument [c\_gsg]

## **Mathematical Expression**

$$0.1 \cdot [c\_gsg] \tag{45}$$

## **7.46 Function definition** function\_4\_V\_ODC\_1

Name function\_4\_V\_ODC\_1

**Arguments** Kipodc, Kmodc, parameter\_1, [species\_2], [species\_7]

#### **Mathematical Expression**

$$\frac{\text{parameter\_1} \cdot [\text{species\_7}]}{\text{Kmodc} \cdot \left(1 + \frac{[\text{species\_2}]}{\text{Kipodc}}\right) + [\text{species\_7}]}$$
(46)

#### 7.47 Function definition function\_4\_V\_SAMDC\_1

Name function\_4\_V\_SAMDC\_1

Arguments Kapsamdc, Kiasamdc, Kissamdc, Kmsamdc, parameter\_3, [sam], [species\_1], [species\_2], [species\_3]

$$\frac{\frac{\text{parameter.}^{3}}{1 + \frac{[\text{species.}3]}{\text{Kissamdc}}} \cdot [\text{sam}]}{\text{Kmsamdc} \cdot \left(1 + \frac{\text{Kapsamdc}}{[\text{species.}2]} + \frac{[\text{species.}1]}{\text{Kiasamdc}}\right) + [\text{sam}]}$$
(47)

## 7.48 Function definition function\_4\_V\_SSAT\_S\_1

Name function\_4\_V\_SSAT\_S\_1

**Arguments** C, Kmaccoassat, Kmcoassat, Kmdssat, Kmsssat, parameter\_2, [species\_3], [species\_4], [species\_8], [species\_9]

## **Mathematical Expression**

$$\frac{\frac{1}{C} \cdot parameter\_2 \cdot [species\_3] \cdot [species\_8]}{Kmsssat \cdot \left(1 + \frac{[species\_4]}{Kmdssat}\right) \cdot Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) + Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot [species\_3] + Kmsssat}$$

#### 7.49 Function definition function\_4\_V\_SSAT\_D\_1

Name function\_4\_V\_SSAT\_D\_1

**Arguments** Kmaccoassat, Kmcoassat, Kmdssat, Kmsssat, parameter\_2, [species\_3], [species\_4], [species\_8], [species\_9]

## **Mathematical Expression**

$$\frac{parameter\_2 \cdot [species\_4] \cdot [species\_8]}{Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmsssat}\right) \cdot Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) + Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot [species\_4] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmdsat \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right$$

#### 7.50 Function definition function\_4\_V\_PAO\_AS\_1

Name function\_4\_V\_PAO\_AS\_1

**Arguments** Kmadpao, Kmaspao, Kmdpao, Kmspao, Vmpao, [species\_3], [species\_4], [species\_5], [species\_6]

#### **Mathematical Expression**

$$\frac{\text{Vmpao} \cdot [\text{species\_5}]}{\text{Kmaspao} \cdot \left(1 + \frac{[\text{species\_6}]}{\text{Kmadpao}} + \frac{[\text{species\_5}]}{\text{Kmaspao}} + \frac{[\text{species\_4}]}{\text{Kmdpao}} + \frac{[\text{species\_3}]}{\text{Kmspao}}\right)}$$
(50)

#### 7.51 Function definition function\_4\_V\_PAO\_AD\_1

Name function\_4\_V\_PAO\_AD\_1

**Arguments** Kmadpao, Kmaspao, Kmdpao, Kmspao, Vmpao, [species\_3], [species\_4], [species\_5], [species\_6]

$$\frac{Vmpao \cdot [species\_6]}{Kmadpao \cdot \left(1 + \frac{[species\_6]}{Kmadpao} + \frac{[species\_5]}{Kmaspao} + \frac{[species\_4]}{Kmdpao} + \frac{[species\_3]}{Kmspao}\right)}$$
(51)

#### 7.52 Function definition function\_4\_V\_SPDS\_1

Name function\_4\_V\_SPDS\_1

Arguments KaSpds, Kiaspds, Kidspds, Kpspds, Vmspds, [species\_1], [species\_2], [species\_4]

#### **Mathematical Expression**

$$\frac{V m s p d s \cdot [s pecies\_1] \cdot [s pecies\_2]}{K i a s p d s \cdot K p s p d s \cdot \left(1 + \frac{[s pecies\_4]}{K i d s p d s}\right) + K p s p d s \cdot [s pecies\_1] + K a S p d s \cdot \left(1 + \frac{[s pecies\_4]}{K i d s p d s}\right) \cdot [s pecies\_2] + [s pecies\_1] \cdot [s pecies\_2]}$$

#### 7.53 Function definition function\_4\_V\_SPMS\_1

Name function\_4\_V\_SPMS\_1

Arguments Kaspms, Kdspms, Kiaspms, Kisspms, Vmspms, [species\_1], [species\_3], [species\_4]

## **Mathematical Expression**

$$\frac{V m spms \cdot [species\_1] \cdot [species\_4]}{Kiaspms \cdot K d spms \cdot \left(1 + \frac{[species\_3]}{Kisspms}\right) + K d spms \cdot [species\_1] + K a spms \cdot \left(1 + \frac{[species\_3]}{Kisspms}\right) \cdot [species\_4] + [species\_1]}$$

#### 7.54 Function definition function\_4\_V\_MATII

Name function\_4\_V\_MATII

**Arguments** Ki, Km, [met], parameter\_19, [sam]

#### **Mathematical Expression**

$$\frac{\text{parameter}_{-19}}{1 + \left(\frac{\text{Km} \cdot \left(1 + \frac{[\text{sam}]}{\text{Ki}}\right)}{[\text{met}]}\right)^{0.76}}$$
(54)

#### 7.55 Function definition function\_4\_V\_SDH\_1

Name function\_4\_V\_SDH\_1

Arguments K\_src\_SDH, K\_thf\_SDH, Vm\_SDH, [m\_thf], vol (mito), [src]

$$vol\left(mito\right) \cdot MM\_twosubst\left(Vm\_SDH, K\_thf\_SDH, K\_src\_SDH, [m\_thf], \\ [src]\right)$$

#### 7.56 Function definition function\_4\_V\_b\_GLU\_c\_1

Name function\_4\_V\_b\_GLU\_c\_1

Arguments K\_bglutc, V\_bglutc, [b\_glu], [c\_glu], vol (cytosol), k\_out\_glu

#### **Mathematical Expression**

$$vol\left(cytosol\right)\cdot\left(MM\left(V\_bglutc,K\_bglutc,[b\_glu]\right)-k\_out\_glu\cdot[c\_glu]\right) \quad (56)$$

## 7.57 Function definition function\_4\_VmGLYc\_1

Name function\_4\_VmGLYc\_1

**Arguments** K\_cgly, K\_mgly, V\_cgly, V\_mgly, [c\_gly], vol (cytosol), [m\_gly], vol (mito)

#### **Mathematical Expression**

$$MM (V_{mgly}, K_{mgly}, [m_{gly}]) \cdot vol(mito) \cdot \frac{1}{3}$$

$$- MM (V_{cgly}, K_{cgly}, [c_{gly}]) \cdot vol(cytosol)$$
(57)

#### 7.58 Function definition function\_4\_V\_b\_SER\_c\_1

Name function\_4\_V\_b\_SER\_c\_1

Arguments K\_bserc, V\_bserc, [b\_ser], [c\_ser], vol (cytosol), k\_out\_ser

#### **Mathematical Expression**

$$vol(cytosol) \cdot (MM(V\_bserc, K\_bserc, [b\_ser]) - k\_out\_ser \cdot [c\_ser])$$
 (58)

#### 7.59 Function definition function\_4\_VmSHMT\_1

Name function\_4\_VmSHMT\_1

## **7.60 Function definition** function\_4\_V\_c\_gshHb\_1

Name function\_4\_V\_c\_gshHb\_1

**Arguments** K\_gshHb, V\_gshHb, [c\_gsh], vol (cytosol)

#### **Mathematical Expression**

$$vol(cytosol) \cdot MM(V_gshHb, K_gshHb, [c_gsh])$$
 (60)

#### **7.61 Function definition** function\_4\_V\_b\_MET\_c\_1

Name function\_ $4_V_b_MET_c_1$ 

**Arguments** K\_bmetc, V\_bmetc, [b\_met], vol (cytosol), k\_out\_met, [met]

#### **Mathematical Expression**

$$vol(cytosol) \cdot (MM(V\_bmetc, K\_bmetc, [b\_met]) - k\_out\_met \cdot [met])$$
 (61)

#### **7.62 Function definition** function\_4\_V\_b\_GLY\_c\_1

Name function\_4\_V\_b\_GLY\_c\_1

**Arguments** K\_bglyc, V\_bglyc, [b\_gly], [c\_gly], vol(cytosol), k\_out\_gly

#### **Mathematical Expression**

$$vol(cytosol) \cdot (MM(V_bglyc, K_bglyc, [b_gly]) - k_out_gly \cdot [c_gly])$$
 (62)

#### 7.63 Function definition function\_4\_VmFTS\_1

Name function\_4\_VmFTS\_1

**Arguments** K\_10f\_mFTS, K\_coo\_mFTS, K\_thf\_mFTS, Vf\_mFTS, Vr\_mFTS, [m\_10f], [m\_coo], [m\_thf]

#### **Mathematical Expression**

$$MM_{twosubst}(Vf_{mFTS}, K_{thf_{mFTS}}, K_{coo_{mFTS}}, K_{coo_{mFTS}}, K_{loo_{mFTS}}, K_{$$

#### 7.64 Function definition function\_4\_VmMTD\_1

Name function\_4\_VmMTD\_1

Arguments K\_1cf\_MTD, K\_2cf\_MTD, Vf\_mMTD, Vr\_MTD, [m\_1cf], [m\_2cf]

$$MM (Vf\_mMTD, K\_2cf\_MTD, [m\_2cf]) - MM (Vr\_MTD, K\_1cf\_MTD, (64)$$

$$[m\_1cf])$$

#### **7.65 Function definition** function\_4\_V\_DMGD\_1

Name function\_4\_V\_DMGD\_1

**Arguments** K\_dmg\_DMGD, K\_thf\_DMGD, Vm\_DMGD, [dmg], [m\_thf], vol (mito)

## **Mathematical Expression**

 $vol\left(mito\right) \cdot MM\_twosubst\left(Vm\_DMGD, K\_thf\_DMGD, K\_dmg\_DMGD, [m\_thf], \\ [dmg]\right)$ 

(65)

#### 7.66 Function definition function\_4\_VmMTCH\_1

Name function\_4\_VmMTCH\_1

Arguments K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_mMTCH, Vr\_MTCH, [m\_10f], [m\_1cf]

## **Mathematical Expression**

 $\label{eq:mmtch} \begin{aligned} &MM\left(Vf\_mMTCH,K\_1cf\_MTCH,[m\_1cf]\right) - MM\left(Vr\_MTCH,K\_10f\_MTCH,\\ &[m\_10f]\right) \end{aligned}$ 

(66)

## **7.67 Function definition** function\_4\_V\_b\_CYS\_c\_1

Name function\_4\_V\_b\_CYS\_c\_1

**Arguments** K\_bcysc, V\_bcysc, [b\_cys], vol (cytosol)

#### **Mathematical Expression**

$$vol(cytosol) \cdot MM(V\_bcysc, K\_bcysc, [b\_cys])$$
 (67)

## **7.68 Function definition** function\_4\_V\_c\_gsgLb\_1

Name function\_4\_V\_c\_gsgLb\_1

**Arguments** K\_gsgLb, V\_gsgLb, [c\_gsg], vol (cytosol)

$$vol(cytosol) \cdot MM(V_gsgLb, K_gsgLb, [c_gsg])$$
 (68)

#### **7.69 Function definition** function\_4\_VmSERc\_1

Name function\_4\_VmSERc\_1

**Arguments** K\_cser, K\_mser, V\_cser, V\_mser, [c\_ser], vol (cytosol), [m\_ser], vol (mito)

## **Mathematical Expression**

$$\left( \frac{\text{MM} \left( \text{V\_mser}, \text{K\_mser}, [\text{m\_ser}] \right) \cdot \text{vol} \left( \text{mito} \right)}{3} \right.$$

$$\left. - \text{MM} \left( \text{V\_cser}, \text{K\_cser}, [\text{c\_ser}] \right) \right) \cdot \text{vol} \left( \text{cytosol} \right)$$

$$(69)$$

## **7.70 Function definition** function\_4\_V\_c\_gsgHb\_1

Name function\_4\_V\_c\_gsgHb\_1

**Arguments** K\_gsgHb, V\_gsgHb, [c\_gsg], vol(cytosol)

#### **Mathematical Expression**

$$vol(cytosol) \cdot MM(V_gsgHb, K_gsgHb, [c_gsg])$$
 (70)

#### 8 Rules

This is an overview of 17 rules.

#### 8.1 Rule tot\_cfol

Rule tot\_cfol is an assignment rule for parameter tot\_cfol:

$$tot\_cfol = [c\_5mf] + [c\_2cf] + [c\_1cf] + [c\_10f] + [c\_dhf] + [c\_thf]$$
(71)

**Derived unit**  $\mu mol \cdot l^{-1}$ 

#### 8.2 Rule tot\_mfol

Rule tot\_mfol is an assignment rule for parameter tot\_mfol:

$$tot\_mfol = [m\_thf] + [m\_2cf] + [m\_1cf] + [m\_10f]$$
(72)

**Derived unit**  $\mu mol \cdot l^{-1}$ 

#### 8.3 Rule daytime

Rule daytime is an assignment rule for parameter daytime:

$$daytime = time - 24 \cdot \left\lfloor \frac{time}{24} \right\rfloor \tag{73}$$

## 8.4 Rule aa\_input

Rule aa\_input is an assignment rule for parameter aa\_input:

#### 8.5 Rule b\_met

Rule b\_met is an assignment rule for species b\_met:

$$b\_met = aa\_input \cdot b\_met\_basal$$
 (75)

#### 8.6 Rule b\_ser

Rule b\_ser is an assignment rule for species b\_ser:

$$b\_ser = aa\_input \cdot b\_ser\_basal$$
 (76)

## 8.7 Rule V\_oGly\_b

Rule V\_oGly\_b is an assignment rule for parameter V\_oGly\_b:

$$V_oGly_b = aa\_input \cdot V_oGly_b\_basal$$
 (77)

## 8.8 Rule V\_oCys\_b

Rule V\_oCys\_b is an assignment rule for parameter V\_oCys\_b:

$$V_oCys_b = aa\_input \cdot V_oCys_b\_basal$$
 (78)

## 8.9 Rule V\_oGlu\_b

Rule V\_oGlu\_b is an assignment rule for parameter V\_oGlu\_b:

$$V_oGlu_b = aa_input \cdot V_oGlu_b_basal$$
 (79)

#### 8.10 Rule parameter\_14

Rule parameter\_14 is an assignment rule for parameter parameter\_14:

$$parameter_14 = parameter_16$$
 (80)

## **8.11 Rule** parameter\_15

Rule parameter\_15 is an assignment rule for parameter parameter\_15:

$$parameter_{15} = 3 \cdot parameter_{16}$$
 (81)

## **8.12 Rule** parameter\_17

Rule parameter\_17 is an assignment rule for parameter parameter\_17:

$$parameter_17 = [species_3] + [species_4]$$
 (82)

Derived unit  $\mu mol \cdot l^{-1}$ 

## 8.13 Rule parameter\_18

Rule parameter\_18 is an assignment rule for parameter parameter\_18:

$$parameter_{18} = 100 \cdot \frac{parameter_{16}}{0.24}$$
 (83)

## 8.14 Rule parameter\_1

Rule parameter\_1 is a rate rule for parameter parameter\_1:

$$\frac{d}{dt}parameter\_1 = 60 \cdot \frac{65.06}{[sam]} \cdot parameter\_7 \cdot \frac{1}{1 + parameter\_5 \cdot ([species\_4] + [species\_3])}$$

$$- parameter\_6 \cdot parameter\_4 \cdot parameter\_1$$
(84)

#### 8.15 Rule parameter\_2

Rule parameter\_2 is a rate rule for parameter parameter\_2:

$$\frac{d}{dt} parameter\_2 = 60 \cdot parameter\_9 \cdot \left(1 - \frac{1}{1 + parameter\_5 \cdot ([species\_4] + [species\_3])}\right) - parameter\_8 \cdot \frac{1}{1 + parameter\_5 \cdot ([species\_4] + [species\_3])} \cdot parameter\_2$$
(85)

## 8.16 Rule parameter\_3

Rule parameter\_3 is a rate rule for parameter parameter\_3:

$$\frac{d}{dt} parameter_{3} = 60 \cdot \frac{65.06}{[sam]} \cdot parameter_{1} \cdot \frac{1}{1 + parameter_{5} \cdot ([species_{4}] + [species_{3}])}$$
(86)
$$- parameter_{1} \cdot 0 \cdot parameter_{3}$$

# 8.17 Rule parameter\_4

Rule parameter\_4 is a rate rule for parameter parameter\_4:

$$\frac{d}{dt} parameter\_4 = 1 \cdot parameter\_13 \cdot \left(1 - \frac{1}{1 + parameter\_5 \cdot 0.01 \cdot ([species\_4] + [species\_3])}\right) - parameter\_12 \cdot parameter\_4$$
(87)

# 9 Reactions

This model contains 74 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
	IU	ivanic		300
1	$b_{gsh_decomp}$	b_gsh_decomp	$b_g sh \xrightarrow{b_g sh, b_g sh} b_c cys + b_g ly + b_g lu$	
2	$b\_gsg\_decomp$	b_gsg_decomp	$b_{gsg} = \frac{b_{gsg}, b_{gsg}}{b_{gsg}} 2b_{gsg} + 2b_{gly} + 2b_{glu}$	
3	b_cys_cystine- _conv	b_cys_cystine_conv	b_cys \( \frac{b_cys, b_cys}{\limits}  \	
4	b_cys_loss	b_cys_loss	$b_{\text{cys}} \stackrel{b_{\text{cys}}, b_{\text{cys}}}{\longleftarrow} \emptyset$	
5	b_glu_loss	b_glu_loss	b_glu \(\frac{b_glu, b_glu}{\tau}\) \(\phi\)	
6	b_gly_loss	b_gly_loss	$b_{-gly} \stackrel{b_{-gly}, b_{-gly}}{\longleftarrow} \emptyset$	
7	b_gsh_loss	b_gsh_loss	$b_g sh \xrightarrow{b_g sh} \emptyset$	
8	b_gsg_loss	b_gsg_loss	$b_{-gsg} \xrightarrow{b_{-gsg}, b_{-gsg}} \emptyset$	
9	$b\_cys\_import$	b_cys_import	Ø <del>====</del> b_cys	
10	$b_gly_import$	b_gly_import	$\emptyset \rightleftharpoons b_g ly$	
11	$b_glu_import$	b_glu_import	Ø <del>====</del> b_glu	
12	$V_{-}c_{-}gshHb$	$V_{c_gshHb}$	$c_gsh \xrightarrow{c_gsh} b_gsh$	
13	$V_c_gshLb$	V_c_gshLb	$c_gsh \xrightarrow{c_gsh, c_gsh} b_gsh$	
14	$V_c_gsgHb$	$V_c_gsgHb$	$c_{-gsg} = \frac{c_{-gsg}, c_{-gsg}}{c_{-gsg}} b_{-gsg}$	
15	$V_c_gsgLb$	$V_{-}c_{-}gsgLb$	$c\_gsg \xrightarrow{c\_gsg, c\_gsg} b\_gsg$	

34	No	Id	Name	Reaction Equation	SBO
	16	V_b_CYS_c	V_b_CYS_c	$b_{\text{-cys}} \xrightarrow{b_{\text{-cys}}} c_{\text{-cys}}$	
	17	$V_b_GLU_c$	$V_b_GLU_c$	b_glu \(\frac{b_glu, c_glu, b_glu, c_glu}{\tau}\) c_glu	
	18	$V_b_GLY_c$	V_b_GLY_c	$b_{-gly} \xrightarrow{b_{-gly}, c_{-gly}, b_{-gly}, c_{-gly}} c_{-gly}$	
	19	V_b_SER_c	V_b_SER_c	b_ser \(\frac{b_ser, c_ser, b_ser, c_ser}{c_ser}\)	
	20	V_b_MET_c	V_b_MET_c	b_met b_met, met, b_met, met met	
	21	VmFTD	VmFTD	$m_10f \xrightarrow{m_10f, m_10f} m_thf$	
Produ	22	VmSHMT	VmSHMT	m_thf+m_ser = m_2cf, m_gly, m_ser, m_thf, m_2cf, m_grader m_2cf	=====================================
iced l	23	VmFTS	VmFTS	$m_{thf} + m_{coo} = \frac{m_{1}0f, m_{coo}, m_{thf}, m_{1}0f, m_{coo}, m_{thf}}{m_{thf}}$	$\frac{\text{n\_thf}}{\text{m}}$ m_10f
9 <i>y</i> <b>9</b> 8	24	VmNE	VmNE	m_thf+HCHO HCHO, m_2cf, m_thf, HCHO, m_2cf,	$\frac{\text{m\_thf}}{\text{m\_2cf}}$
Produced by SBML2l <sup>ET</sup> EX	25	$V_{-}GDC$	V_GDC	$\begin{array}{l} \text{m\_thf} + \text{m\_gly} \xleftarrow{\text{m\_gly, m\_thf, m\_gly, m\_thf}} \text{m\_2cf} + \\ \text{CO} \end{array}$	
' <del>'</del> ×	26	$V_{-}SDH$	V_SDH	$m_{thf} + src \xrightarrow{m_{thf}, src, m_{thf}, src} m_{2}cf + m_{gly}$	
	27	V_DMGD	V_DMGD	$m_{thf} + dmg \xrightarrow{dmg, m_{thf}, dmg, m_{thf}} m_{2}cf + src$	
	28	VmMTD	VmMTD	$m_2cf \xrightarrow{m_1cf, m_2cf, m_1cf, m_2cf} m_1cf$	
	29	VmMTCH	VmMTCH	$m_{-}1cf \xrightarrow{m_{-}10f, m_{-}1cf, m_{-}10f, m_{-}1cf} m_{-}10f$	
	30	VmSERc	VmSERc	$3 \text{ m\_ser} \xrightarrow{\text{c\_ser, m\_ser, c\_ser, m\_ser}} \text{c\_ser}$	
	31	VmHCOOHc	VmHCOOHc	3 m_coo	
	32	VmGLYc	VmGLYc	$3 \text{ m\_gly} \stackrel{\text{c\_gly, m\_gly, c\_gly, m\_gly}}{\longleftarrow} \text{c\_gly}$	

Nº Id	Name	Reaction Equation SBO	
33 V_MS	V_MS	c_5mf+hcy H2O2, c_5mf, hcy, H2O2, c_5mf, hcy c_thf+	
	V _IVIS	met C_JIII + IICY \ C_LIII +	
34 V_DHFR	V_DHFR	$c_dhf + NADPH \xrightarrow{NADPH, c_dhf, NADPH, c_dhf} c_thf$	
04 V_Dnrk	V-DHFK		
35 VcFTD	VcFTD	$c_{-10f} \stackrel{c_{-10f}, c_{-10f}}{\longleftarrow} c_{-thf}$	
36 V_PGT	V_PGT	$c_{-}10f + GAR \xrightarrow{GAR, c_{-}10f, GAR, c_{-}10f} aic + c_{-}thf$	
37 VcFTS	VcFTS	$c_{-thf} + c_{-coo} \xleftarrow{c_{-coo}, c_{-thf}, c_{-coo}, c_{-thf}} c_{-10f}$	
38 VcSHMT	VcSHMT	$c_{ser} + c_{thf} = \frac{c_2cf, c_gly, c_{ser}, c_thf, c_2cf, c_gly, c_{ser}, c_thf}{c_gly}$	
Jo Vebinii	VESTIMI	c_2cf	
39 VcNE	VcNE	$c_{-}$ thf+HCHO $\leftarrow$ C_2cf, c_thf, HCHO, c_2cf, c_thf $\leftarrow$ c_2cf	
39 VCNE			
$40 \text{ V}_{-}\text{TS}$	$V_{-}TS$	$DUMP + c_2cf \xrightarrow{DUMP, c_2cf, DUMP, c_2cf} c_dhf$	
41 V_MTHFR	V_MTHFR	c_2cf+NADPH sah, sam, NADPH, c_2cf, sah, sam, NADPH, c_2cf, sah, sa	$\stackrel{\text{n}}{=}$ c
42 VcMTD	VcMTD	$c_{-}2cf \xrightarrow{c_{-}1cf, c_{-}2cf, c_{-}1cf, c_{-}2cf} c_{-}1cf + NADPH$	
43 VcMTCH	VcMTCH	$c_{-1}cf = \frac{c_{-1}0f, c_{-1}cf, c_{-1}0f, c_{-1}cf}{c_{-1}0f}$	
44 V_ART	$V\_ART$	$c_{-}10f + aic \stackrel{aic, c_{-}10f, aic, c_{-}10f}{\rightleftharpoons} c_{-}thf$	
		H2O2, sah, sam, BET, H2O2, hcv, sah, sam, BET, H2O2, hcv, s	ah,
45 V_BHMT	$V\_BHMT$	hcy+BET	
		dmg	
46 V_MATI	V_MATI	$met \xrightarrow{c\_gsg, c\_gsg, met, sam, c\_gsg, met, sam} sam$	
47 V_MATIII	V_MATIII	$met \xleftarrow{c\_gsg, c\_gsg, met, sam, c\_gsg, met, sam} sam$	
48 V_GNMT	$V_{-}GNMT$	$sam+c\_gly \stackrel{c\_5mf, c\_5mf, c\_gly, sah, sam, c\_5mf, c\_gly, sah, sam}{=} sah+$	
10 V_MIIII	V_01\1111	stri + c_gry \	

36	No	Id	Name	Reaction Equation	SBO
Produced by SBMI2 <sup>E</sup>	49	V_DNMT	V_DNMT	sam sah, sam, sah, sam sah	
	50	V_SAHH	$V_SAHH$	$sah \xrightarrow{hcy, sah, hcy, sah} hcy$	
	51	gluconeogenesis- _ser	- gluconeogenesis_ser	$c_{ser} \stackrel{c_{ser}, c_{ser}}{\longleftarrow} \emptyset$	
	52	V_CBS	V_CBS	hcy+c_ser H2O2, sah, sam, H2O2, c_ser, hcy, sah, s	sam, H2O2, c_ser, hcy, sah, sam
	53	$V_{-}CTGL$	V_CTGL	$cyt \xrightarrow{cyt, cyt} c\_cys$	
	54	V_GCS	V_GCS	$c\_cys + c\_glu = H2O2, c\_gsh, H2O2, c\_cys, c\_glu, c\_gsh$	h, glc, H2O2, c_cys, c_glu, c_gsh,
	55	cys_usage	cys_usage	$c\_{cys} \stackrel{c\_{cys}, c\_{cys}}{\longleftarrow} \emptyset$	
	56	c_glu_usage	c_glu_usage	$c_{glu} \stackrel{c_{glu}, c_{glu}}{\longleftarrow} \emptyset$	
	57	$V_{-}GS$	V_GS	$glc + c_gly \xrightarrow{c_gly, c_gsh, glc, c_gly, c_gsh, glc} c_gsh$	
	58	$V_{-}GPX$	$V_{-}GPX$	$2 c_{gsh} + H2O2 \xrightarrow{H2O2, c_{gsh}, H2O2, c_{gsh}} c_{gsg}$	
	59	$V_{-}GR$	$V_{-}GR$	c_gsg+NADPH NADPH, c_gsg, NADPH, c_gsg 2 c	z_gsh
	60	c_gsh_degr	c_gsh_degr	$c_g sh \xrightarrow{c_g sh} \theta$	1
	61	c_gsg_degr	c_gsg_degr	$c_{-gsg} \stackrel{c_{-gsg}, c_{-gsg}}{\longleftarrow} \emptyset$	
	62	${\tt reaction\_1}$	V_ODC	species_7, species_7, species_2, species_7 s	
	63	reaction_2	V_SAMDC	sam species_3, species_2, sam, species_1, species_2, s	
	64	reaction_3	V_SSAT_S	species_3 + species_8 species_4, species_3, species_4, species_9	
	65	reaction_4	V_SSAT_D	species_4+species_8 species_3, species_3, species_4, species_9	, species_8, species_9, species_3, s

N₀	Id	Name	Reaction Equation	SBO
66	reaction_5	V_PAO_AD	species_6 species_5, species_4, species_3, species_6	s_3, species_4, species_5, species_6, sp
			species_5, species_3, species_3, species_5	s_4, species_5, species_6, species_3, sp
67	reaction_6	V_PAO_AS		
68	$reaction_{-}7$	V_SPDS	species_1 + species_2 species_1, species_2, species	
69	reaction_8	V_SPMS	species_1+species_4 species_3, species_3	es_4, species_1, species_3, species_4 species_4
70	reaction_9	V_COA	species_8 species_8, species_8 species_9	
71	reaction_10	V_ACCOA	species_9 $\xrightarrow{\text{species}\_9}$ species_8	
72	reaction_11	V_PUT_efflux	species_2 $\xrightarrow{\text{species}\_2}$ , $\xrightarrow{\text{species}\_2}$ $\emptyset$	
73	reaction_12	V_AD_efflux	species_6 $\xrightarrow{\text{species}\_6}$ , $\xrightarrow{\text{species}\_6}$ $\emptyset$	
74	reaction_13	V_MATII	$met \stackrel{sam, met, sam, met, sam}{\longleftarrow} sam$	

# 9.1 Reaction b\_gsh\_decomp

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

Name b\_gsh\_decomp

# **Reaction equation**

$$b\_gsh \xrightarrow{b\_gsh, b\_gsh} b\_cys + b\_gly + b\_glu \tag{88}$$

#### Reactant

Table 6: Properties of each reactant.

Id	Name	SBO
b_gsh	b_GSH	

### **Modifiers**

Table 7: Properties of each modifier.

Id	Name	SBO
	b_GSH	
b_gsh	b_GSH	

### **Products**

Table 8: Properties of each product.

Id	Name	SBO
b_cys b_gly b_glu	b_Cysteine b_Glycine b_Glutamate	

### **Kinetic Law**

$$v_1 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_gsh\_decomp\_1}([\text{b\_gsh}])$$
 (89)

$$function_4_b_g sh_d e comp_1([b_g sh]) = 90 \cdot [b_g sh]$$
(90)

function\_4\_b\_gsh\_decomp\_1 ([b\_gsh]) = 
$$90 \cdot [b_gsh]$$
 (91)

# 9.2 Reaction b\_gsg\_decomp

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

Name b\_gsg\_decomp

# **Reaction equation**

$$b\_gsg \xrightarrow{b\_gsg, b\_gsg} 2b\_cys + 2b\_gly + 2b\_glu$$
 (92)

#### Reactant

Table 9: Properties of each reactant.

Id	Name	SBO
b_gsg	b_GSSG	

## **Modifiers**

Table 10: Properties of each modifier.

Id	Name	SBO
b_gsg	b_GSSG	
b_gsg	b_GSSG	

## **Products**

Table 11: Properties of each product.

Id	Name	SBO
b_cys b_gly b_glu	b_Cysteine b_Glycine b_Glutamate	

### **Kinetic Law**

$$v_2 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_gsg\_decomp\_1}([\text{b\_gsg}])$$
 (93)

function\_4\_b\_gsg\_decomp\_1 ([b\_gsg]) = 
$$67.5 \cdot [b_gsg]$$
 (94)

function\_4\_b\_gsg\_decomp\_1 ([b\_gsg]) = 
$$67.5 \cdot [b_gsg]$$
 (95)

# **9.3 Reaction** b\_cys\_cystine\_conv

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

Name b\_cys\_cystine\_conv

## **Reaction equation**

$$b_{-cys} \xrightarrow{b_{-cys}, b_{-cys}} \emptyset$$
 (96)

#### Reactant

Table 12: Properties of each reactant.

Id	Name	SBO
b_cys	b_Cysteine	

### **Modifiers**

Table 13: Properties of each modifier.

Id	Name	SBO
b_cys	b_Cysteine	
b_cys	b_Cysteine	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_3 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_cys\_cystine\_conv\_1}([\text{b\_cys}])$$
 (97)

function\_4\_b\_cys\_cystine\_conv\_1 ([b\_cys]) = 
$$0.25 \cdot [b_cys]$$
 (98)

function\_4\_b\_cys\_cystine\_conv\_1 ([b\_cys]) = 
$$0.25 \cdot [b_cys]$$
 (99)

# **9.4 Reaction** b\_cys\_loss

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

Name b\_cys\_loss

## **Reaction equation**

$$b\_cys \xrightarrow{b\_cys, b\_cys} \emptyset$$
 (100)

### Reactant

Table 14: Properties of each reactant.

Id	Name	SBO
b_cys	b_Cysteine	

## **Modifiers**

Table 15: Properties of each modifier.

Id	Name	SBO
b_cys b_cys		

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_4 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_cys\_loss\_1}([\text{b\_cys}])$$
 (101)

$$function\_4\_b\_cys\_loss\_1([b\_cys]) = 0.1 \cdot [b\_cys]$$
 (102)

$$function\_4\_b\_cys\_loss\_1([b\_cys]) = 0.1 \cdot [b\_cys]$$

$$(103)$$

# 9.5 Reaction b\_glu\_loss

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

Name b\_glu\_loss

# **Reaction equation**

$$b\_glu \xrightarrow{b\_glu, b\_glu} \emptyset$$
 (104)

### Reactant

Table 16: Properties of each reactant.

Id	Name	SBO
b_glu	b_Glutamate	

### **Modifiers**

Table 17: Properties of each modifier.

Id	Name	SBO
_	b_Glutamate b_Glutamate	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_5 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_glu\_loss\_1}([\text{b\_glu}])$$
 (105)

$$function\_4\_b\_glu\_loss\_1([b\_glu]) = 0.1 \cdot [b\_glu]$$

$$(106)$$

$$function\_4\_b\_glu\_loss\_1([b\_glu]) = 0.1 \cdot [b\_glu]$$

$$(107)$$

# **9.6 Reaction** b\_gly\_loss

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

Name b\_gly\_loss

## **Reaction equation**

$$b_{gly} = 0$$

$$b_{gly} = 0$$

$$(108)$$

## Reactant

Table 18: Properties of each reactant.

Id	Name	SBO
b_gly	b_Glycine	

### **Modifiers**

Table 19: Properties of each modifier.

Id	Name	SBO
b_gly	b_Glycine	

Id	Name	SBO
b_gly	b_Glycine	

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_6 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_gly\_loss\_1}([\text{b\_gly}])$$
 (109)

$$function\_4\_b\_gly\_loss\_1([b\_gly]) = 0.1 \cdot [b\_gly]$$

$$(110)$$

$$function\_4\_b\_gly\_loss\_1([b\_gly]) = 0.1 \cdot [b\_gly]$$

$$(111)$$

# 9.7 Reaction b\_gsh\_loss

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

Name b\_gsh\_loss

## **Reaction equation**

$$b\_gsh \xrightarrow{b\_gsh, b\_gsh} \emptyset$$
 (112)

### Reactant

Table 20: Properties of each reactant.

Id	Name	SBO
b_gsh	b_GSH	

#### **Modifiers**

Table 21: Properties of each modifier.

Id	Name	SBO
•	b_GSH	
b_gsh	b_GSH	

### **Kinetic Law**

$$v_7 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_gsh\_loss\_1}([\text{b\_gsh}])$$
 (113)

function\_4\_b\_gsh\_loss\_1 ([b\_gsh]) = 
$$0.7 \cdot [b_gsh]$$
 (114)

function\_4\_b\_gsh\_loss\_1([b\_gsh]) = 
$$0.7 \cdot [b_gsh]$$
 (115)

# 9.8 Reaction b\_gsg\_loss

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

Name b\_gsg\_loss

## **Reaction equation**

$$b_{-gsg} \stackrel{b_{-gsg}, b_{-gsg}}{\longleftarrow} \emptyset$$
 (116)

#### Reactant

Table 22: Properties of each reactant.

Id	Name	SBO
b_gsg	b_GSSG	

### **Modifiers**

Table 23: Properties of each modifier.

Id	Name	SBO
b_gsg	b_GSSG	
b_gsg	b_GSSG	

### **Kinetic Law**

$$v_8 = \text{vol}(\text{blood}) \cdot \text{function\_4\_b\_gsg\_loss\_1}([\text{b\_gsg}])$$
 (117)

function\_4\_b\_gsg\_loss\_1([b\_gsg]) = 
$$7.5 \cdot [b_gsg]$$
 (118)

function\_4\_b\_gsg\_loss\_1([b\_gsg]) = 
$$7.5 \cdot [b_gsg]$$
 (119)

# 9.9 Reaction b\_cys\_import

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

Name b\_cys\_import

# **Reaction equation**

$$\emptyset \rightleftharpoons b_{-}cys$$
 (120)

## **Product**

Table 24: Properties of each product.

Id	Name	SBO
b_cys	b_Cysteine	

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_9 = \text{vol}(\text{blood}) \cdot \text{function}_1(V_oCys_b)$$
 (121)

$$function_{-}1(v) = v (122)$$

$$function_{-1}(v) = v (123)$$

# 9.10 Reaction b\_gly\_import

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

Name b\_gly\_import

# **Reaction equation**

$$\emptyset \rightleftharpoons b_g ly$$
 (124)

### **Product**

Table 25: Properties of each product.

Id	Name	SBO
b_gly	b_Glycine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{10} = \text{vol}(\text{blood}) \cdot \text{function}_1(V_oGly_b)$$
 (125)

$$function_{-}1(v) = v (126)$$

$$function_{-}1(v) = v (127)$$

# 9.11 Reaction b\_glu\_import

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

Name b\_glu\_import

## **Reaction equation**

$$\emptyset \rightleftharpoons b_{-g}lu$$
 (128)

### **Product**

Table 26: Properties of each product.

Id	Name	SBO
b_glu	b_Glutamate	

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{11} = \text{vol}(\text{blood}) \cdot \text{function}_1(\text{V}_{-0}\text{Glu}_{-b})$$
 (129)

$$function_{-1}(v) = v \tag{130}$$

$$function_{-}1(v) = v (131)$$

# 9.12 Reaction V\_c\_gshHb

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

Name V\_c\_gshHb

$$c_{-gsh} \xleftarrow{c_{-gsh}, c_{-gsh}} b_{-gsh}$$
 (132)

#### Reactant

Table 27: Properties of each reactant.

Id	Name	SBO
c_gsh	c_GSH	

#### **Modifiers**

Table 28: Properties of each modifier.

Id	Name	SBO
•	c_GSH c_GSH	

### **Product**

Table 29: Properties of each product.

Id	Name	SBO
b_gsh	b_GSH	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{12} = \text{function\_4\_V\_c\_gshHb\_1} (K\_gshHb, V\_gshHb, [c\_gsh], vol (cytosol))$$
 (133)

# 9.13 Reaction V\_c\_gshLb

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

Name V\_c\_gshLb

$$c_{-gsh} \xleftarrow{c_{-gsh}, c_{-gsh}} b_{-gsh}$$
 (135)

## Reactant

Table 30: Properties of each reactant.

Id	Name	SBO
c_gsh	c_GSH	

#### **Modifiers**

Table 31: Properties of each modifier.

Id	Name	SBO
c_gsh	c_GSH	
$c\_gsh$	c_GSH	

### **Product**

Table 32: Properties of each product.

Id	Name	SBO
b_gsh	b_GSH	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_{13} = \text{function\_4\_V\_c\_gshLb\_1} (K\_\text{gshLb}, V\_\text{gshLb}, [c\_\text{gsh}], \text{vol}(\text{cytosol}), h\_\text{gshLb})$$
 (136)

$$\begin{split} & \text{function\_4\_V\_c\_gshLb\_1}\left(K\_gshLb, V\_gshLb, [c\_gsh], \text{vol}\left(\text{cytosol}\right), h\_gshLb\right) \\ &= \frac{\text{vol}\left(\text{cytosol}\right) \cdot V\_gshLb \cdot [c\_gsh]^{h\_gshLb}}{K\_gshLb^{h\_gshLb} + [c\_gsh]^{h\_gshLb}} \end{split} \tag{137}$$

# 9.14 Reaction V\_c\_gsgHb

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

Name V\_c\_gsgHb

$$c_{-gsg} = \underbrace{c_{-gsg}, c_{-gsg}}_{b_{-gsg}} b_{-gsg}$$
 (138)

#### Reactant

Table 33: Properties of each reactant.

Id	Name	SBO
c_gsg	c_GSSG	

#### **Modifiers**

Table 34: Properties of each modifier.

Id	Name	SBO
	c_GSSG c_GSSG	

#### **Product**

Table 35: Properties of each product.

Id	Name	SBO
b_gsg	b_GSSG	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{14} = \text{function\_4\_V\_c\_gsgHb\_1} (K\_gsgHb, V\_gsgHb, [c\_gsg], vol (cytosol))$$
 (139)

$$\begin{aligned} & \text{function\_4\_V\_c\_gsgHb\_1} \left( \text{K\_gsgHb}, \text{V\_gsgHb}, [\text{c\_gsg}], \text{vol} \left( \text{cytosol} \right) \right) \\ &= & \text{vol} \left( \text{cytosol} \right) \cdot \text{MM} \left( \text{V\_gsgHb}, \text{K\_gsgHb}, [\text{c\_gsg}] \right) \end{aligned} \tag{140}$$

# 9.15 Reaction V\_c\_gsgLb

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

Name  $V_c_gsgLb$ 

$$c\_gsg \xrightarrow{c\_gsg, c\_gsg} b\_gsg$$
 (141)

#### Reactant

Table 36: Properties of each reactant.

Id	Name	SBO
c_gsg	c_GSSG	

#### **Modifiers**

Table 37: Properties of each modifier.

Id	Name	SBO
c_gsg c_gsg	c_GSSG c_GSSG	

#### **Product**

Table 38: Properties of each product.

Id	Name	SBO
b_gsg	b_GSSG	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_{15} = \text{function\_4\_V\_c\_gsgLb\_1} (K\_gsgLb, V\_gsgLb, [c\_gsg], vol (cytosol))$$
 (142)

## 9.16 Reaction V\_b\_CYS\_c

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

Name V\_b\_CYS\_c

$$b_{\text{cys}} \stackrel{b_{\text{cys}}, b_{\text{cys}}}{=\!=\!=\!=\!=} c_{\text{cys}}$$
 (144)

#### Reactant

Table 39: Properties of each reactant.

Id	Name	SBO
b_cys	b_Cysteine	

#### **Modifiers**

Table 40: Properties of each modifier.

Id	Name	SBO
b_cys	b_Cysteine	
b_cys	b_Cysteine	

### **Product**

Table 41: Properties of each product.

	1	1
Id	Name	SBO
c_cys	c_Cysteine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{16} = \text{function\_4\_V\_b\_CYS\_c\_1} (K\_bcysc, V\_bcysc, [b\_cys], vol(cytosol))$$
 (145)

$$function\_4\_V\_b\_CYS\_c\_1(K\_bcysc, V\_bcysc, [b\_cys], vol(cytosol))$$

$$= vol(cytosol) \cdot MM(V\_bcysc, K\_bcysc, [b\_cys])$$

$$(146)$$

## 9.17 Reaction V\_b\_GLU\_c

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

Name V\_b\_GLU\_c

$$b\_glu \xrightarrow{b\_glu, c\_glu, b\_glu, c\_glu} c\_glu \tag{147}$$

#### Reactant

Table 42: Properties of each reactant.

Id	Name	SBO
b_glu	b_Glutamate	

### **Modifiers**

Table 43: Properties of each modifier.

Id	Name	SBO
b_glu	b_Glutamate	
c_glu	$c_Glutamate$	
b_glu	b_Glutamate	
c_glu	c_Glutamate	

### **Product**

Table 44: Properties of each product.

Id	Name	SBO
c_glu	c_Glutamate	-

#### **Kinetic Law**

Derived unit contains undeclared units

$$v_{17} = function\_4\_V\_b\_GLU\_c\_1 (K\_bglutc, V\_bglutc, [b\_glu], [c\_glu], vol (cytosol), k\_out\_glu)$$

$$(148)$$

# 9.18 Reaction V\_b\_GLY\_c

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

Name V\_b\_GLY\_c

$$b\_gly \xleftarrow{b\_gly, c\_gly, b\_gly, c\_gly} c\_gly \tag{150}$$

#### Reactant

Table 45: Properties of each reactant.

Id	Name	SBO
b_gly	b_Glycine	

#### **Modifiers**

Table 46: Properties of each modifier.

Id	Name	SBO
b_gly	b_Glycine	
$c\_{\tt gly}$	c_Glycine	
$b_gly$	b_Glycine	
$c\_{gly}$	c_Glycine	

### **Product**

Table 47: Properties of each product.

Id	Name	SBO
c_gly	c_Glycine	

#### **Kinetic Law**

Derived unit contains undeclared units

$$v_{18} = function\_4\_V\_b\_GLY\_c\_1 (K\_bglyc, V\_bglyc, [b\_gly], [c\_gly], vol (cytosol), k\_out\_gly) \tag{151}$$

## 9.19 Reaction V\_b\_SER\_c

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

Name V\_b\_SER\_c

$$b\_ser \xrightarrow{b\_ser, c\_ser, b\_ser, c\_ser} c\_ser$$
 (153)

### Reactant

Table 48: Properties of each reactant.

Id	Name	SBO
b_ser	b_Serine	

### **Modifiers**

Table 49: Properties of each modifier.

Id	Name	SBO
$b_{-}ser$	$b_{-}$ Serine	
c_ser	c_Serine	
$b_ser$	b_Serine	
c_ser	c_Serine	

## **Product**

Table 50: Properties of each product.

Id	Name	SBO
c_ser	c_Serine	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_{19} = \text{function\_4\_V\_b\_SER\_c\_1} (K\_\text{bserc}, V\_\text{bserc}, [b\_\text{ser}], [c\_\text{ser}], \text{vol} (\text{cytosol}), \text{k\_out\_ser})$$
(154)

$$\begin{array}{l} function\_4\_V\_b\_SER\_c\_1 \left( K\_bserc, V\_bserc, [b\_ser], [c\_ser], vol \left( cytosol \right), k\_out\_ser \right) \\ = vol \left( cytosol \right) \cdot \left( MM \left( V\_bserc, K\_bserc, [b\_ser] \right) - k\_out\_ser \cdot [c\_ser] \right) \end{array} \tag{155}$$

## 9.20 Reaction V\_b\_MET\_c

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

Name V\_b\_MET\_c

$$b_{met} \xrightarrow{b_{met}, met, b_{met}, met} met$$
 (156)

### Reactant

Table 51: Properties of each reactant.

Id	Name	SBO
b_met	b_Methionine	

### **Modifiers**

Table 52: Properties of each modifier.

Id	Name	SBO
b_met	b_Methionine	
met	c_Methionine	
b_met	b_Methionine	
met	c_Methionine	

## **Product**

Table 53: Properties of each product.

Id	Name	SBO
met	c_Methionine	

### **Kinetic Law**

### Derived unit contains undeclared units

$$v_{20} = \text{function\_4\_V\_b\_MET\_c\_1} (K\_bmetc, V\_bmetc, [b\_met], vol (cytosol), k\_out\_met, [met])$$
(157)

$$\begin{array}{l} function\_4\_V\_b\_MET\_c\_1\left(K\_bmetc,V\_bmetc,[b\_met],vol\left(cytosol\right),k\_out\_met,[met]\right) \\ = vol\left(cytosol\right)\cdot\left(MM\left(V\_bmetc,K\_bmetc,[b\_met]\right) - k\_out\_met\cdot[met]\right) \end{array}$$

# 9.21 Reaction VmFTD

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

### Name VmFTD

$$m_{-}10f \xrightarrow{m_{-}10f, m_{-}10f} m_{-}thf$$
 (159)

## Reactant

Table 54: Properties of each reactant.

Id	Name	SBO
m_10f	m_10-formyl-THF	

## **Modifiers**

Table 55: Properties of each modifier.

Id	Name	SBO
	m_10-formyl-THF m_10-formyl-THF	

### **Product**

Table 56: Properties of each product.

Id	Name	SBO
m_thf	m_THF	

## **Kinetic Law**

Derived unit contains undeclared units

$$v_{21} = \text{vol}(\text{mito}) \cdot \text{function\_4\_VmFTD\_1}(\text{K\_10f\_FTD}, \text{Vm\_mFTD}, [\text{m\_10f}])$$
 (160)

$$function\_4\_VmFTD\_1\left(K\_10f\_FTD,Vm\_mFTD,[m\_10f]\right) = \frac{Vm\_mFTD\cdot[m\_10f]}{K\_10f\_FTD+[m\_10f]} \quad (161)$$

$$function\_4\_VmFTD\_1\left(K\_10f\_FTD,Vm\_mFTD,[m\_10f]\right) = \frac{Vm\_mFTD\cdot[m\_10f]}{K\_10f\_FTD+[m\_10f]} \quad (162)$$

### 9.22 Reaction VmSHMT

This is a reversible reaction of two reactants forming two products influenced by eight modifiers.

### Name VmSHMT

## **Reactants**

Table 57: Properties of each reactant.

Id	Name	SBO
$\mathtt{m}_{-}\mathtt{thf}$	$m_{-}THF$	
${\tt m\_ser}$	m_Serine	

### **Modifiers**

Table 58: Properties of each modifier.

Table 30. I roperties of each mounter.		
Id	Name	SBO
m_2cf	m_5-10-methylene-THF	
$m\_gly$	m_Glycine	
${\tt m\_ser}$	m_Serine	
$\mathtt{m\_thf}$	$m_{-}THF$	
$m_{-}2cf$	m_5-10-methylene-THF	
$m\_{\tt gly}$	m_Glycine	
${\tt m\_ser}$	m_Serine	
${\tt m\_thf}$	m_THF	

## **Products**

Table 59: Properties of each product.

Id	Name	SBO
0 0	m_Glycine m_5-10-methylene-THF	

## **Kinetic Law**

$$v_{22} = vol(mito) \cdot function\_4\_VmSHMT\_1 (K\_2cf\_SHMT, K\_gly\_SHMT, K\_ser\_SHMT, K\_thf\_SHMT, Vf\_mSHMT, Vr\_mSHMT, [m\_2cf], [m\_gly], [m\_ser], [m\_thf])$$

$$(164)$$

$$\begin{split} & function\_4\_VmSHMT\_1\left(K\_2cf\_SHMT,K\_gly\_SHMT,K\_ser\_SHMT,K\_thf\_SHMT, \\ & Vf\_mSHMT,Vr\_mSHMT,[m\_2cf],[m\_gly],[m\_ser],[m\_thf]\right) = MM\_twosubst\left(Vf\_mSHMT,K\_ser\_SHMT,[m\_thf],[m\_ser]\right) \\ & - MM\_twosubst\left(Vr\_mSHMT,K\_gly\_SHMT,K\_2cf\_SHMT,[m\_gly],[m\_2cf]\right) \end{aligned} \tag{165}$$

$$\begin{split} & function\_4\_VmSHMT\_1\left(K\_2cf\_SHMT,K\_gly\_SHMT,K\_ser\_SHMT,K\_thf\_SHMT, \\ & Vf\_mSHMT,Vr\_mSHMT,[m\_2cf],[m\_gly],[m\_ser],[m\_thf]\right) = MM\_twosubst\left(Vf\_mSHMT,K\_ser\_SHMT,[m\_thf],[m\_ser]\right) \\ & - MM\_twosubst\left(Vr\_mSHMT,K\_gly\_SHMT,K\_2cf\_SHMT,[m\_gly],[m\_2cf]\right) \end{aligned} \tag{166}$$

### 9.23 Reaction VmFTS

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

#### Name VmFTS

## **Reaction equation**

$$m_{thf} + m_{coo} \xrightarrow{m_{l}10f, m_{coo}, m_{thf}, m_{l}10f, m_{coo}, m_{thf}} m_{l}10f$$

$$(167)$$

## **Reactants**

Table 60: Properties of each reactant.

Id	Name	SBO
m_thf	m_THF	
$m\_coo$	m_Formate	

#### **Modifiers**

Table 61: Properties of each modifier.

Id	Name	SBO
m_10f	m_10-formyl-THF	
${\tt m\_coo}$	m_Formate	
$\mathtt{m}_{-}\mathtt{thf}$	$m_{-}THF$	
$\mathtt{m}_{-}\mathtt{10f}$	m_10-formyl-THF	
${\tt m\_coo}$	m_Formate	
$\mathtt{m\_thf}$	m_THF	

#### **Product**

Table 62: Properties of each product.

14010 02	. Troperties of each	producti
Id	Name	SBO
$m_{-}10f$	m_10-formyl-THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{23} = \text{vol}(\text{mito}) \cdot \text{function\_4\_VmFTS\_1} (\text{K\_10f\_mFTS}, \text{K\_coo\_mFTS}, \text{K\_thf\_mFTS}, \text{Vf\_mFTS}, \text{Vr\_mFTS}, [\text{m\_10f}], [\text{m\_coo}], [\text{m\_thf}])$$

$$(168)$$

$$\begin{array}{l} \text{function\_4\_VmFTS\_1} \left( \text{K\_10f\_mFTS}, \text{K\_coo\_mFTS}, \text{K\_thf\_mFTS}, \text{Vf\_mFTS}, \text{Vr\_mFTS}, \\ [\text{m\_10f}], [\text{m\_coo}], [\text{m\_thf}] \right) = \text{MM\_twosubst} \left( \text{Vf\_mFTS}, \text{K\_thf\_mFTS}, \text{K\_coo\_mFTS}, \\ [\text{m\_thf}], [\text{m\_coo}] \right) - \text{MM} \left( \text{Vr\_mFTS}, \text{K\_10f\_mFTS}, [\text{m\_10f}] \right) \end{array}$$

$$\begin{array}{l} \text{function\_4\_VmFTS\_1} \left( \text{K\_10f\_mFTS}, \text{K\_coo\_mFTS}, \text{K\_thf\_mFTS}, \text{Vf\_mFTS}, \text{Vr\_mFTS}, \\ [\text{m\_10f}], [\text{m\_coo}], [\text{m\_thf}] \right) = \text{MM\_twosubst} \left( \text{Vf\_mFTS}, \text{K\_thf\_mFTS}, \text{K\_coo\_mFTS}, \\ [\text{m\_thf}], [\text{m\_coo}] \right) - \text{MM} \left( \text{Vr\_mFTS}, \text{K\_10f\_mFTS}, [\text{m\_10f}] \right) \end{array}$$

## 9.24 Reaction VmNE

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

### Name VmNE

### **Reaction equation**

$$m_{thf} + HCHO \xrightarrow{HCHO, m_{thf}, HCHO, m_{thf}, HCHO, m_{thf}, m_{thf}} m_{thf}$$

$$(171)$$

#### **Reactants**

Table 63: Properties of each reactant.

Id	Name	SBO
m_thf HCHO	m_THF Formaldehyde	

### **Modifiers**

Table 64: Properties of each modifier.

Id	Name	SBO
НСНО	Formaldehyde	
$m_2cf$	m_5-10-methylene-THF	
$\mathtt{m\_thf}$	$m_{-}THF$	
HCHO	Formaldehyde	
$m_2cf$	m_5-10-methylene-THF	
${\tt m\_thf}$	m_THF	

## **Product**

Table 65: Properties of each product.

Id	Name	SBO
m_2cf	m_5-10-methylene-THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{24} = \text{function\_4\_VmNE\_1} ([\text{HCHO}], \text{k1\_mNE}, \text{k2\_mNE}, [\text{m\_2cf}], [\text{m\_thf}], \text{vol} (\text{mito}))$$
 (172)

$$\begin{array}{l} \text{function\_4\_VmNE\_1}\left([\text{HCHO}], \text{k1\_mNE}, \text{k2\_mNE}, [\text{m\_2cf}], [\text{m\_thf}], \text{vol}\left(\text{mito}\right)\right) \\ = \text{vol}\left(\text{mito}\right) \cdot \left(\text{k1\_mNE} \cdot [\text{m\_thf}] \cdot [\text{HCHO}] - \text{k2\_mNE} \cdot [\text{m\_2cf}]\right) \end{array}$$

# 9.25 Reaction V\_GDC

This is a reversible reaction of two reactants forming two products influenced by four modifiers.

### Name V\_GDC

## **Reaction equation**

$$m_{thf} + m_{gly} \xrightarrow{m_{gly}, m_{thf}, m_{gly}, m_{thf}} m_{2cf} + CO$$

$$(174)$$

### **Reactants**

Table 66: Properties of each reactant.

Id	Name	SBO
m_thf	m_THF	
$m_{-} \texttt{gly}$	m_Glycine	

## **Modifiers**

Table 67: Properties of each modifier.

Id	Name	SBO
m_gly m_thf	m_Glycine m_THF	
m_gly m_thf	m_Glycine m_THF	

#### **Products**

Table 68: Properties of each product.

Id	Name	SBO
m_2cf	m_5-10-methylene-THF	

### **Kinetic Law**

$$v_{25} = \text{vol}(\text{mito}) \cdot \text{function\_4\_V\_GDC\_1} (\text{K\_gly\_GDC}, \text{K\_thf\_GDC}, \text{Vm\_GDC}, [\text{m\_gly}], [\text{m\_thf}])$$
(175)

$$\begin{split} & \text{function\_4\_V\_GDC\_1}\left(\text{K\_gly\_GDC}, \text{K\_thf\_GDC}, \text{Vm\_GDC}, [\text{m\_gly}], [\text{m\_thf}]\right) \\ & = \frac{\text{Vm\_GDC} \cdot [\text{m\_thf}] \cdot [\text{m\_gly}]}{\left(\text{K\_thf\_GDC} + [\text{m\_thf}]\right) \cdot \left(\text{K\_gly\_GDC} + [\text{m\_gly}]\right)} \end{split} \tag{176}$$

$$\begin{split} & \text{function\_4\_V\_GDC\_1}\left(\text{K\_gly\_GDC}, \text{K\_thf\_GDC}, \text{Vm\_GDC}, [\text{m\_gly}], [\text{m\_thf}]\right) \\ & = \frac{\text{Vm\_GDC} \cdot [\text{m\_thf}] \cdot [\text{m\_gly}]}{\left(\text{K\_thf\_GDC} + [\text{m\_thf}]\right) \cdot \left(\text{K\_gly\_GDC} + [\text{m\_gly}]\right)} \end{split} \tag{177}$$

# 9.26 Reaction V\_SDH

This is a reversible reaction of two reactants forming two products influenced by four modifiers.

## Name V\_SDH

# **Reaction equation**

$$m_{thf} + src \xrightarrow{m_{thf}, src, m_{thf}, src} m_{2}cf + m_{gly}$$
 (178)

## **Reactants**

Table 69: Properties of each reactant.

Id	Name	SBO
m_thf	m_THF	
src	Sarcosine	

## **Modifiers**

Table 70: Properties of each modifier.

Id	Name	SBO
$m_{-}$ thf	$m_{-}THF$	
src	Sarcosine	
$\mathtt{m\_thf}$	$m_{-}THF$	
src	Sarcosine	

### **Products**

Table 71: Properties of each product.

Id	Name	SBO
	m_5-10-methylene-THF	
$m_gly$	m_Glycine	

## **Kinetic Law**

$$v_{26} = \text{function\_4\_V\_SDH\_1} \left( \text{K\_src\_SDH}, \text{K\_thf\_SDH}, \text{Vm\_SDH}, [\text{m\_thf}], \text{vol} \left( \text{mito} \right), [\text{src}] \right)$$

$$(179)$$

$$\begin{array}{l} function\_4\_V\_SDH\_1\left(K\_src\_SDH,K\_thf\_SDH,Vm\_SDH,[m\_thf],vol\left(mito\right),[src]\right) \\ = vol\left(mito\right)\cdot MM\_twosubst\left(Vm\_SDH,K\_thf\_SDH,K\_src\_SDH,[m\_thf],[src]\right) \end{array}$$

# 9.27 Reaction V\_DMGD

This is a reversible reaction of two reactants forming two products influenced by four modifiers.

# Name V\_DMGD

# **Reaction equation**

$$m\_thf + dmg \xrightarrow{dmg, m\_thf, dmg, m\_thf} m\_2cf + src$$
 (181)

## **Reactants**

Table 72: Properties of each reactant.

Id	Name	SBO
m_thf dmg	m_THF Dimethylglycine	

## **Modifiers**

Table 73: Properties of each modifier.

Id	Name	SBO
dmg m_thf dmg m_thf	Dimethylglycine m_THF Dimethylglycine m_THF	

### **Products**

Table 74: Properties of each product.

Id	Name	SBO
m_2cf	m_5-10-methylene-THF	
src	Sarcosine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{27} = \text{function\_4\_V\_DMGD\_1} (\text{K\_dmg\_DMGD}, \text{K\_thf\_DMGD}, \text{Vm\_DMGD}, [\text{dmg}], \\ [\text{m\_thf}], \text{vol} (\text{mito}))$$
 (182)

$$\begin{aligned} & \text{function\_4\_V\_DMGD\_1}\left(K\_dmg\_DMGD, K\_thf\_DMGD, \\ & \text{Vm\_DMGD}, [dmg], [m\_thf], vol\left(mito\right)\right) = vol\left(mito\right) \\ & \cdot \text{MM\_twosubst}\left(\text{Vm\_DMGD}, K\_thf\_DMGD, K\_dmg\_DMGD, [m\_thf], [dmg]\right) \end{aligned} \tag{183}$$

### 9.28 Reaction VmMTD

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

## Name VmMTD

# **Reaction equation**

$$m_2cf \xrightarrow{m_1lcf, m_2cf, m_1lcf, m_2cf} m_1lcf$$
 (184)

#### Reactant

Table 75: Properties of each reactant.

Id	Name	SBO
m_2cf	m_5-10-methylene-THF	

#### **Modifiers**

Table 76: Properties of each modifier.

Id	Name	SBO
${\tt m\_1cf}$	m_5-10-methenyl-THF	
$m_2cf$	m_5-10-methylene-THF	
${\tt m\_1cf}$	m_5-10-methenyl-THF	
$m_2cf$	m_5-10-methylene-THF	

### **Product**

Table 77: Properties of each product

Id	Name	SBO
m_1cf	m_5-10-methenyl-THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{28} = \text{vol}(\text{mito}) \cdot \text{function\_4\_VmMTD\_1} (\text{K\_1cf\_MTD}, \text{K\_2cf\_MTD}, \text{Vf\_mMTD}, \\ \text{Vr\_MTD}, [\text{m\_1cf}], [\text{m\_2cf}])$$
 (185)

$$\begin{aligned} & \text{function\_4\_VmMTD\_1} \left( \text{K\_1cf\_MTD}, \text{K\_2cf\_MTD}, \text{Vf\_mMTD}, \text{Vr\_MTD}, [\text{m\_1cf}], [\text{m\_2cf}] \right) \\ &= \text{MM} \left( \text{Vf\_mMTD}, \text{K\_2cf\_MTD}, [\text{m\_2cf}] \right) - \text{MM} \left( \text{Vr\_MTD}, \text{K\_1cf\_MTD}, [\text{m\_1cf}] \right) \end{aligned} \tag{186}$$

$$\begin{aligned} & \text{function\_4\_VmMTD\_1}\left(K\_1\text{ef\_MTD}, K\_2\text{ef\_MTD}, Vf\_mMTD, Vr\_MTD, [m\_1\text{ef}], [m\_2\text{ef}]\right) \\ &= & \text{MM}\left(Vf\_mMTD, K\_2\text{ef\_MTD}, [m\_2\text{ef}]\right) - & \text{MM}\left(Vr\_MTD, K\_1\text{ef\_MTD}, [m\_1\text{ef}]\right) \end{aligned} \tag{187}$$

## 9.29 Reaction VmMTCH

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

#### Name VmMTCH

## **Reaction equation**

$$m_{-}1cf \xrightarrow{m_{-}10f, m_{-}1cf, m_{-}10f, m_{-}1cf} m_{-}10f$$
 (188)

# Reactant

Table 78: Properties of each reactant.

Id	Name	SBO
m_1cf	m_5-10-methenyl-THF	

### **Modifiers**

Table 79: Properties of each modifier.

	I	
Id	Name	SBO
$m_{-}10f$	m_10-formyl-THF	
${\tt m\_1cf}$	m_5-10-methenyl-THF	7
$\mathtt{m}_{-}\mathtt{10f}$	m_10-formyl-THF	
$m_{-}1cf$	m_5-10-methenyl-THF	<del>-</del>

#### **Product**

Table 80: Properties of each product.

Id	Name	SBO
m_10f	m_10-formyl-THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{29} = \text{vol} (\text{mito}) \cdot \text{function\_4\_VmMTCH\_1} (K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_mMTCH, Vr\_MTCH, [m\_10f], [m\_1ef])$$
 (189)

$$\begin{split} & \text{function\_4\_VmMTCH\_1}\left(K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_mMTCH, \\ & \text{Vr\_MTCH}, [m\_10f], [m\_1cf]\right) = \text{MM}\left(\text{Vf\_mMTCH}, K\_1cf\_MTCH, [m\_1cf]\right) \\ & - \text{MM}\left(\text{Vr\_MTCH}, K\_10f\_MTCH, [m\_10f]\right) \end{split}$$

$$\begin{split} & \text{function\_4\_VmMTCH\_1}\left(K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_mMTCH, \\ & \text{Vr\_MTCH}, [m\_10f], [m\_1cf]\right) = \text{MM}\left(\text{Vf\_mMTCH}, K\_1cf\_MTCH, [m\_1cf]\right) \\ & - \text{MM}\left(\text{Vr\_MTCH}, K\_10f\_MTCH, [m\_10f]\right) \end{split}$$

## 9.30 Reaction VmSERc

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

Name VmSERc

## **Reaction equation**

$$3 \text{ m\_ser} \xleftarrow{\text{c\_ser, m\_ser, c\_ser, m\_ser}} \text{c\_ser}$$
 (192)

## Reactant

Table 81: Properties of each reactant.

Id	Name	SBO
m_ser	m_Serine	

## **Modifiers**

Table 82: Properties of each modifier.

Id	Name	SBO
c_ser	c_Serine	
${\tt m\_ser}$	m_Serine	
c_ser	c_Serine	
m_ser	m_Serine	

#### **Product**

Table 83: Properties of each product.

Id	Name	SBO
c_ser	c_Serine	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_{30} = \text{function\_4\_VmSERc\_1} (K\_\text{cser}, K\_\text{mser}, V\_\text{cser}, V\_\text{mser}, [c\_\text{ser}], \text{vol} (\text{cytosol}),$$

$$[m\_\text{ser}], \text{vol} (\text{mito}))$$
(193)

$$\begin{aligned} \text{function\_4\_VmSERc\_1} \left( \text{K\_cser}, \text{K\_mser}, \text{V\_cser}, \text{V\_mser}, [\text{c\_ser}], \text{vol}\left(\text{cytosol}\right), [\text{m\_ser}], \\ \text{vol}\left(\text{mito}\right) \right) &= \left( \frac{\text{MM}\left(\text{V\_mser}, \text{K\_mser}, [\text{m\_ser}]\right) \cdot \text{vol}\left(\text{mito}\right)}{3} - \text{MM}\left(\text{V\_cser}, \text{K\_cser}, \frac{194}{3}\right) \right) \\ &= \left[ \text{c\_ser} \right] \right) \right) \cdot \text{vol}\left(\text{cytosol}\right) \end{aligned}$$

## 9.31 Reaction VmHCOOHc

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

## Name VmHCOOHc

$$3 \text{ m\_coo} \xleftarrow{\text{c\_coo, m\_coo, c\_coo, m\_coo}} \text{c\_coo}$$
 (195)

#### Reactant

Table 84: Properties of each reactant.

Id	Name	SBO
m_coo	m_Formate	

### **Modifiers**

Table 85: Properties of each modifier.

Id	Name	SBO
C_C00	c_Formate	
$m\_coo$	m_Formate	
C_C00	c_Formate	
m_coo	m_Formate	

### **Product**

Table 86: Properties of each product.

Id	Name	SBO
C_C00	c_Formate	

# **Kinetic Law**

$$v_{31} = \text{function\_4\_VmHCOOHc\_1}([\text{c\_coo}], \text{vol}(\text{cytosol}), \text{k\_in\_coo}, \text{k\_out\_coo}, [\text{m\_coo}], \text{vol}(\text{mito}))$$

$$\begin{aligned} & \text{function\_4\_VmHCOOHc\_1}\left([\text{c\_coo}], \text{vol}\left(\text{cytosol}\right), \text{k\_in\_coo}, \text{k\_out\_coo}, [\text{m\_coo}], \\ & \text{vol}\left(\text{mito}\right)\right) = \frac{\text{k\_in\_coo} \cdot [\text{m\_coo}] \cdot \text{vol}\left(\text{mito}\right)}{3} - \text{k\_out\_coo} \cdot [\text{c\_coo}] \cdot \text{vol}\left(\text{cytosol}\right) \end{aligned}$$

# 9.32 Reaction VmGLYc

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

Name VmGLYc

# **Reaction equation**

$$3 \text{ m\_gly} \xrightarrow{\text{c\_gly, m\_gly, c\_gly, m\_gly}} \text{c\_gly}$$
 (198)

### Reactant

Table 87: Properties of each reactant.

Id	Name	SBO
m_gly	m_Glycine	

### **Modifiers**

Table 88: Properties of each modifier.

Id	Name	SBO
$c_gly$	c_Glycine	
$m\_{\tt gly}$	m_Glycine	
$c\_{\tt gly}$	c_Glycine	
m_gly	m_Glycine	

### **Product**

Table 89: Properties of each product.

Id	Name	SBO
c_gly	c_Glycine	

### **Kinetic Law**

$$v_{32} = function\_4\_VmGLYc\_1 (K\_cgly, K\_mgly, V\_cgly, V\_mgly, [c\_gly], vol (cytosol), (199)$$

$$[m\_gly], vol (mito))$$

$$\begin{split} & \text{function\_4\_VmGLYc\_1}\left(K\_\text{cgly}, K\_\text{mgly}, V\_\text{cgly}, V\_\text{mgly}, [c\_\text{gly}], \\ & \text{vol}\left(\text{cytosol}\right), [m\_\text{gly}], \text{vol}\left(\text{mito}\right)\right) = MM\left(V\_\text{mgly}, K\_\text{mgly}, [m\_\text{gly}]\right) \\ & \cdot \text{vol}\left(\text{mito}\right) \cdot \frac{1}{3} - MM\left(V\_\text{cgly}, K\_\text{cgly}, [c\_\text{gly}]\right) \cdot \text{vol}\left(\text{cytosol}\right) \end{split} \tag{200}$$

## 9.33 Reaction V\_MS

This is a reversible reaction of two reactants forming two products influenced by seven modifiers.

## Name V\_MS

# **Reaction equation**

$$c\_5mf + hcy \xrightarrow{H2O2, H2O2, c\_5mf, hcy, H2O2, c\_5mf, hcy} c\_thf + met \tag{201}$$

### **Reactants**

Table 90: Properties of each reactant.

Id	Name	SBO
c_5mf hcy	c_5-methyl-THF Homocysteine	

## **Modifiers**

Table 91: Properties of each modifier.

Id	Name	SBO
H202	H2O2	
H202	H2O2	
$\texttt{c\_5mf}$	c_5-methyl-THF	
hcy	Homocysteine	
H202	H2O2	
$\texttt{c}_{-}\texttt{5mf}$	c_5-methyl-THF	
hcy	Homocysteine	

### **Products**

Table 92: Properties of each product.

Id	Name	SBO
$c_{-}$ thf	c_THF	
met	c_Methionine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{33} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_MS\_1}([\text{H2O2}], \text{K\_5mf\_MS}, \text{K\_hcy\_MS}, \text{Ki\_MS}, \\ \text{Vm\_MS}, [\text{c\_5mf}], [\text{hcy}], \text{ssH2O2})$$

$$\begin{aligned} & \text{function\_4\_V\_MS\_1}\left([\text{H2O2}], \text{K\_5mf\_MS}, \text{K\_hcy\_MS}, \text{Ki\_MS}, \text{Vm\_MS}, [\text{c\_5mf}], [\text{hcy}], \text{ssH2O2}\right) \\ &= \frac{\text{MM\_twosubst}\left(\text{Vm\_MS}, \text{K\_5mf\_MS}, \text{K\_hcy\_MS}, [\text{c\_5mf}], [\text{hcy}]\right) \cdot \left(\text{ssH2O2} + \text{Ki\_MS}\right)}{[\text{H2O2}] + \text{Ki\_MS}} \end{aligned} \tag{203}$$

$$\begin{split} & \text{function\_4\_V\_MS\_1} \left( [\text{H2O2}], \text{K\_5mf\_MS}, \text{K\_hcy\_MS}, \text{Ki\_MS}, \text{Vm\_MS}, [\text{c\_5mf}], [\text{hcy}], \text{ssH2O2} \right) \\ &= \frac{\text{MM\_twosubst} \left( \text{Vm\_MS}, \text{K\_5mf\_MS}, \text{K\_hcy\_MS}, [\text{c\_5mf}], [\text{hcy}] \right) \cdot \left( \text{ssH2O2} + \text{Ki\_MS} \right) }{[\text{H2O2}] + \text{Ki\_MS}} \end{aligned}$$

## 9.34 Reaction V\_DHFR

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

### Name V\_DHFR

### **Reaction equation**

$$c_-dhf + NADPH \xrightarrow{NADPH, c_-dhf, NADPH, c_-dhf} c_-thf$$
 (205)

#### **Reactants**

Table 93: Properties of each reactant.

Id	Name	SBO
0_4111	c_DHF NADPH	

### **Modifiers**

Table 94: Properties of each modifier.

Id	Name	SBO
NADPH	NADPH	
$c_dhf$	c_DHF	
NADPH	NADPH	
$c_{-}dhf$	c_DHF	

#### **Product**

Table 95: Properties of each product.

Id	Name	SBO
$c_{-}$ thf	c_THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{34} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_DHFR\_1} (K\_NADPH\_DHFR, K\_dhf\_DHFR, [c\_dhf])$$
 (206)

$$\begin{aligned} & \text{function\_4\_V\_DHFR\_1} \left( \text{K\_NADPH\_DHFR}, \text{K\_dhf\_DHFR}, [\text{NADPH}], \text{Vm\_DHFR}, \\ & \left[ \text{c\_dhf} \right] \right) = \frac{\text{Vm\_DHFR} \cdot [\text{c\_dhf}] \cdot [\text{NADPH}]}{\left( \text{K\_dhf\_DHFR} + [\text{c\_dhf}] \right) \cdot \left( \text{K\_NADPH\_DHFR} + [\text{NADPH}] \right)} \end{aligned} \tag{207}$$

$$\begin{aligned} & \text{function\_4\_V\_DHFR\_1} \ (K\_NADPH\_DHFR, K\_dhf\_DHFR, [NADPH], Vm\_DHFR,} \\ & [c\_dhf]) = \frac{Vm\_DHFR \cdot [c\_dhf] \cdot [NADPH]}{(K\_dhf\_DHFR + [c\_dhf]) \cdot (K\_NADPH\_DHFR + [NADPH])} \end{aligned}$$

# 9.35 Reaction VcFTD

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

Name VcFTD

## **Reaction equation**

$$c_{-}10f \underbrace{c_{-}10f, c_{-}10f}_{} c_{-}thf$$
 (209)

# Reactant

Table 96: Properties of each reactant.

Id	Name	SBO
c_10f	c_10-formyl-THF	

Table 97: Properties of each modifier.

Id	Name	SBO
	c_10-formyl-THF c_10-formyl-THF	

## **Product**

Table 98: Properties of each product.

Id	Name	SBO
$c_{-}thf$	c_THF	

## **Kinetic Law**

Derived unit contains undeclared units

$$v_{35} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_VcFTD\_1}(\text{K\_10f\_FTD}, \text{Vm\_cFTD}, [\text{c\_10f}])$$
 (210)

$$function\_4\_VcFTD\_1\left(K\_10f\_FTD,Vm\_cFTD,[c\_10f]\right) = \frac{Vm\_cFTD\cdot[c\_10f]}{K\_10f\_FTD+[c\_10f]} \quad (211)$$

$$function\_4\_VcFTD\_1\left(K\_10f\_FTD,Vm\_cFTD,[c\_10f]\right) = \frac{Vm\_cFTD\cdot[c\_10f]}{K\_10f\_FTD+[c\_10f]} \quad (212)$$

## 9.36 Reaction V\_PGT

This is a reversible reaction of two reactants forming two products influenced by four modifiers.

Name V\_PGT

## **Reaction equation**

$$c_{-}10f + GAR \xrightarrow{GAR, c_{-}10f, GAR, c_{-}10f} aic + c_{-}thf$$
 (213)

## Reactants

Table 99: Properties of each reactant.

Id	Name	SBO
c_10f GAR	c_10-formyl-THF GAR	

Table 100: Properties of each modifier.

	*	
Id	Name	SBO
GAR	GAR	
$c_{-}10f$	$c_10$ -formyl-THF	
GAR	GAR	
$c_{-}10f$	$c_{-}10$ -formyl-THF	

#### **Products**

Table 101: Properties of each product.

Id	Name	SBO
aic	AICAR	
c_thf	c_THF	

#### **Kinetic Law**

$$v_{36} = vol\left(cytosol\right) \cdot function\_4\_V\_PGT\_1\left([GAR], K\_10f\_PGT, K\_GAR\_PGT, Vm\_PGT, [c\_10f]\right) \tag{214}$$

$$\begin{split} & \text{function\_4\_V\_PGT\_1}\left([GAR], K\_10f\_PGT, K\_GAR\_PGT, Vm\_PGT, [c\_10f]\right) \\ & = \frac{Vm\_PGT \cdot [c\_10f] \cdot [GAR]}{\left(K\_10f\_PGT + [c\_10f]\right) \cdot \left(K\_GAR\_PGT + [GAR]\right)} \end{split} \tag{215}$$

$$\begin{aligned} & \text{function\_4\_V\_PGT\_1}\left([GAR], K\_10f\_PGT, K\_GAR\_PGT, Vm\_PGT, [c\_10f]\right) \\ & = \frac{Vm\_PGT \cdot [c\_10f] \cdot [GAR]}{(K\_10f\_PGT + [c\_10f]) \cdot (K\_GAR\_PGT + [GAR])} \end{aligned}$$

## 9.37 Reaction VcFTS

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

### Name VcFTS

# **Reaction equation**

$$c_{-}thf + c_{-}coo \xrightarrow{c_{-}coo, c_{-}thf, c_{-}coo, c_{-}thf} c_{-}10f$$
(217)

## **Reactants**

Table 102: Properties of each reactant.

Id	Name	SBO
$c_{-}$ thf	c_THF	
C_C00	c_Formate	

### **Modifiers**

Table 103: Properties of each modifier.

Id	Name	SBO
C_C00		
$c_{-} thf$	c_THF	
C_C00	$c_Formate$	
c_thf	c_THF	

### **Product**

Table 104: Properties of each product.

Id	Name	SBO
c_10f	c_10-formyl-THF	

## **Kinetic Law**

$$v_{37} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_VcFTS\_1} (\text{K\_coo\_cFTS}, \text{K\_thf\_cFTS}, \text{Vm\_cFTS}, [\text{c\_coo}], [\text{c\_thf}])$$
(218)

$$\begin{split} & \text{function\_4\_VcFTS\_1} \left( \text{K\_coo\_cFTS}, \text{K\_thf\_cFTS}, \text{Vm\_cFTS}, [\text{c\_coo}], [\text{c\_thf}] \right) \\ &= \frac{\text{Vm\_cFTS} \cdot [\text{c\_thf}] \cdot [\text{c\_coo}]}{\left( \text{K\_thf\_cFTS} + [\text{c\_thf}] \right) \cdot \left( \text{K\_coo\_cFTS} + [\text{c\_coo}] \right)} \end{aligned} \tag{219}$$

$$\begin{aligned} & \text{function\_4\_VcFTS\_1} \left( \text{K\_coo\_cFTS}, \text{K\_thf\_cFTS}, \text{Vm\_cFTS}, [\text{c\_coo}], [\text{c\_thf}] \right) \\ &= \frac{\text{Vm\_cFTS} \cdot [\text{c\_thf}] \cdot [\text{c\_coo}]}{\left( \text{K\_thf\_cFTS} + [\text{c\_thf}] \right) \cdot \left( \text{K\_coo\_cFTS} + [\text{c\_coo}] \right)} \end{aligned} \tag{220}$$

### 9.38 Reaction VcSHMT

This is a reversible reaction of two reactants forming two products influenced by eight modifiers.

## Name VcSHMT

### **Reaction equation**

$$c\_ser + c\_thf \xrightarrow{c\_2cf, c\_gly, c\_ser, c\_thf, c\_2cf, c\_gly, c\_ser, c\_thf} c\_gly + c\_2cf$$
 (221)

## **Reactants**

Table 105: Properties of each reactant.

Id	Name	SBO
0_00_	c_Serine c_THF	

Table 106: Properties of each modifier.

Id	Name	SBO
c_2cf	c_5-10-methylene-THF	
$c\_{\tt gly}$	c_Glycine	
c_ser	c_Serine	
$c_{-}thf$	c_THF	
$c_2cf$	c_5-10-methylene-THF	
$c\_{gly}$	c_Glycine	
c_ser	c_Serine	
c_thf	c_THF	

## **Products**

Table 107: Properties of each product.

Tueste 1071 11 repetition of out-in production		
Id	Name	SBO
0 0	c_Glycine c_5-10-methylene-THF	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$\begin{aligned} v_{38} &= \text{vol}\left(\text{cytosol}\right) \cdot \text{function\_4\_VcSHMT\_1}\left(\text{K\_2cf\_SHMT}, \text{K\_gly\_SHMT}, \text{K\_ser\_SHMT}, \\ &\quad \text{K\_thf\_SHMT}, \text{Vf\_cSHMT}, \text{Vr\_cSHMT}, [\text{c\_2cf}], [\text{c\_gly}], [\text{c\_ser}], [\text{c\_thf}]\right) \end{aligned} \tag{222}$$

$$\begin{split} & function\_4\_VcSHMT\_1\left(K\_2cf\_SHMT,K\_gly\_SHMT,K\_ser\_SHMT,K\_thf\_SHMT,Vf\_cSHMT,Vr\_cSHMT,[c\_2cf],[c\_gly],[c\_ser],[c\_thf]\right) = & MM\_twosubst\left(Vf\_cSHMT,K\_thf\_SHMT,K\_thf\_SHMT,K\_ser\_SHMT,[c\_thf],[c\_ser]\right) \\ & - & MM\_twosubst\left(Vr\_cSHMT,K\_gly\_SHMT,K\_2cf\_SHMT,[c\_gly],[c\_2cf]\right) \end{split} \label{eq:control_state} \tag{223}$$

$$\begin{split} & function\_4\_VcSHMT\_1\left(K\_2cf\_SHMT,K\_gly\_SHMT,K\_ser\_SHMT,K\_thf\_SHMT,Vf\_cSHMT,Vr\_cSHMT,[c\_2cf],[c\_gly],[c\_ser],[c\_thf]\right) = & MM\_twosubst\left(Vf\_cSHMT,K\_thf\_SHMT,K\_thf\_SHMT,K\_thf\_SHMT,F_c\_thf],[c\_ser]\right) \\ & - & MM\_twosubst\left(Vr\_cSHMT,K\_gly\_SHMT,K\_2cf\_SHMT,[c\_gly],[c\_2cf]\right) \end{split} \label{eq:control_shape} \tag{224}$$

## 9.39 Reaction VcNE

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

Name VcNE

### **Reaction equation**

$$c_{-}thf + HCHO \xrightarrow{HCHO, c_{-}2cf, c_{-}thf, HCHO, c_{-}2cf, c_{-}thf} c_{-}2cf$$
 (225)

#### Reactants

Table 108: Properties of each reactant.

Id	Name	SBO
c_thf	c_THF	
HCH0	Formaldehyde	

Table 109: Properties of each modifier.

Id	Name	SBO
нсно	Formaldehyde	
$c_2cf$	c_5-10-methylene-THF	
$c_{-}thf$	c_THF	
HCHO	Formaldehyde	
$c_2cf$	c_5-10-methylene-THF	
c_thf	c_THF	

### **Product**

Table 110: Properties of each product.

Id	Name	SBO
	c_5-10-methylene-THF	

## **Kinetic Law**

### **Derived unit** contains undeclared units

$$v_{39} = \text{function\_4\_VcNE\_1} ([\text{HCHO}], [\text{c\_2cf}], [\text{c\_thf}], \text{vol}(\text{cytosol}), \text{k1\_cNE}, \text{k2\_cNE})$$
 (226)

$$function\_4\_VcNE\_1 ([HCHO], [c\_2cf], [c\_thf], vol (cytosol), k1\_cNE, k2\_cNE)$$

$$= vol (cytosol) \cdot (k1\_cNE \cdot [c\_thf] \cdot [HCHO] - k2\_cNE \cdot [c\_2cf])$$

$$(227)$$

### 9.40 Reaction V\_TS

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

Name V\_TS

# **Reaction equation**

$$DUMP + c_2cf \xrightarrow{DUMP, c_2cf, DUMP, c_2cf} c_dhf$$
 (228)

### **Reactants**

Table 111: Properties of each reactant.

SBO
hylene-THF

#### **Modifiers**

Table 112: Properties of each modifier.

Id	Name	SBO
DUMP	dUMP	
$c_2cf$	c_5-10-methylene-THF	
DUMP	dUMP	
$c_2cf$	c_5-10-methylene-THF	

### **Product**

Table 113: Properties of each product.

Id	Name	SBO
c_dhf	c_DHF	

### **Kinetic Law**

$$v_{40} = vol\left(cytosol\right) \cdot function\_4\_V\_TS\_1\left([DUMP], K\_2cf\_TS, K\_DUMP\_TS, Vm\_TS, [c\_2cf]\right) \tag{229}$$

$$\begin{aligned} & \text{function\_4\_V\_TS\_1}\left([\text{DUMP}], \text{K\_2cf\_TS}, \text{K\_DUMP\_TS}, \text{Vm\_TS}, [\text{c\_2cf}]\right) \\ & = \frac{\text{Vm\_TS} \cdot [\text{DUMP}] \cdot [\text{c\_2cf}]}{(\text{K\_DUMP\_TS} + [\text{DUMP}]) \cdot (\text{K\_2cf\_TS} + [\text{c\_2cf}])} \end{aligned}$$

$$\begin{split} & \text{function\_4\_V\_TS\_1}\left([\text{DUMP}], \text{K\_2cf\_TS}, \text{K\_DUMP\_TS}, \text{Vm\_TS}, [\text{c\_2cf}]\right) \\ & = \frac{\text{Vm\_TS} \cdot [\text{DUMP}] \cdot [\text{c\_2cf}]}{\left(\text{K\_DUMP\_TS} + [\text{DUMP}]\right) \cdot \left(\text{K\_2cf\_TS} + [\text{c\_2cf}]\right)} \end{split} \tag{231}$$

# 9.41 Reaction V\_MTHFR

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

### Name V\_MTHFR

# **Reaction equation**

$$c\_2cf + NADPH \xrightarrow{sah, sam, NADPH, c\_2cf, sah, sam, NADPH, c\_2cf, sah, sam} c\_5mf \quad (232)$$

## **Reactants**

Table 114: Properties of each reactant.

Id	Name	SBO
	c_5-10-methylene-THF NADPH	

## **Modifiers**

Table 115: Properties of each modifier.

Id	Name	SBO
sah	SAH	
sam	SAM	
NADPH	NADPH	
$c_2cf$	c_5-10-methylene-THF	
sah	SAH	
sam	SAM	
NADPH	NADPH	
$c_2cf$	c_5-10-methylene-THF	
sah	SAH	
sam	SAM	

# **Product**

Table 116: Properties of each product.

Id	Name	SBO
$c_5mf$	c_5-methyl-THF	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{41} = \text{vol}\left(\text{cytosol}\right) \cdot \text{function\_4\_V\_MTHFR\_1}\left(\text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, \\ [\text{NADPH}], \text{Vm\_MTHFR}, [\text{c\_2cf}], [\text{sah}], [\text{sam}]\right)$$

$$\text{function\_4\_V\_MTHFR\_1}\left(\text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, \\ [\text{NADPH}], \text{Vm\_MTHFR}, [\text{c\_2cf}], [\text{sah}], [\text{sam}]\right)$$

$$= \frac{\text{MM\_twosubst}\left(\text{Vm\_MTHFR}, \text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, [\text{c\_2cf}], [\text{NADPH}]\right) \cdot 72}{10 + [\text{sam}] - [\text{sah}]}$$

$$\text{function\_4\_V\_MTHFR\_1}\left(\text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, \\ [\text{NADPH}], \text{Vm\_MTHFR}, [\text{c\_2cf}], [\text{sah}], [\text{sam}]\right)$$

$$= \frac{\text{MM\_twosubst}\left(\text{Vm\_MTHFR}, \text{K\_2cf\_MTHFR}, \text{K\_NADPH\_MTHFR}, [\text{c\_2cf}], [\text{NADPH}]\right) \cdot 72}{10 + [\text{sam}] - [\text{sah}]}$$

### 9.42 Reaction VcMTD

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

### Name VcMTD

## **Reaction equation**

$$c_2cf \xrightarrow{c_2lef, c_2lef, c_2lef} c_1lef + NADPH$$
 (236)

(235)

### Reactant

Table 117: Properties of each reactant.

Id	Name	SBO
c_2cf	c_5-10-methylene-THF	

Table 118: Properties of each modifier.

	I	
Id	Name	SBO
c_1cf	c_5-10-methenyl-THF	
$c_2cf$	c_5-10-methylene-THF	
$c_1cf$	c_5-10-methenyl-THF	
$c_2cf$	c_5-10-methylene-THF	

#### **Products**

Table 119: Properties of each product

Table 117. I Toperties of each product.		
Id	Name	SBO
	c_5-10-methenyl-THF NADPH	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{42} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_VcMTD\_1} (\text{K\_1cf\_MTD}, \text{K\_2cf\_MTD}, \text{Vf\_cMTD}, \\ \text{Vr\_cMTD}, [\text{c\_1cf}], [\text{c\_2cf}])$$
 (237)

$$\begin{aligned} & \text{function\_4\_VcMTD\_1} \left( \text{K\_1cf\_MTD}, \text{K\_2cf\_MTD}, \text{Vf\_cMTD}, \text{Vr\_cMTD}, [\text{c\_1cf}], [\text{c\_2cf}] \right) \\ &= \text{MM} \left( \text{Vf\_cMTD}, \text{K\_2cf\_MTD}, [\text{c\_2cf}] \right) - \text{MM} \left( \text{Vr\_cMTD}, \text{K\_1cf\_MTD}, [\text{c\_1cf}] \right) \end{aligned} \tag{238}$$

### 9.43 Reaction VcMTCH

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

#### Name VcMTCH

## **Reaction equation**

$$c_{-1}cf \xrightarrow{c_{-1}0f, c_{-1}cf, c_{-1}0f, c_{-1}cf} c_{-1}0f$$
 (240)

## Reactant

Table 120: Properties of each reactant.

Id	Name	SBO
c_1cf	c_5-10-methenyl-THF	

Table 121: Properties of each modifier.

ruore r	ruote 121. I roperties et euen moumen.		
Id	Name	SBO	
$c_{-}1cf$	c_10-formyl-THF c_5-10-methenyl-THF c_10-formyl-THF		
	c_5-10-methenyl-THF		

### **Product**

Table 122: Properties of each product.

Id	Name	SBO
c_10f	c_10-formyl-THF	

#### **Kinetic Law**

## **Derived unit** contains undeclared units

$$v_{43} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_VcMTCH\_1} (\text{K\_10f\_MTCH}, \text{K\_1cf\_MTCH}, \text{Vf\_cMTCH}, \text{Vr\_MTCH}, \text{[c\_10f]}, \text{[c\_1cf]})$$
 (241)

$$\begin{array}{l} \mbox{function\_4\_VcMTCH\_1}\left(K\_10f\_MTCH, K\_1cf\_MTCH, Vf\_cMTCH, \\ \mbox{Vr\_MTCH}, [c\_10f], [c\_1cf]\right) = \mbox{MM}\left(Vf\_cMTCH, K\_1cf\_MTCH, [c\_1cf]\right) \\ - \mbox{MM}\left(Vr\_MTCH, K\_10f\_MTCH, [c\_10f]\right) \end{array}$$

#### 9.44 Reaction V\_ART

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

## Name V\_ART

## **Reaction equation**

$$c_{-}10f + aic \xrightarrow{aic, c_{-}10f, aic, c_{-}10f} c_{-}thf$$
 (244)

## **Reactants**

Table 123: Properties of each reactant.

Id	Name	SBO
c_10f aic	c_10-formyl-THF AICAR	

## **Modifiers**

Table 124: Properties of each modifier.

Id	Name	SBO
aic	AICAR	-
$c_10f$	c_10-formyl-THF	
aic	AICAR	
$c_{-}10f$	$c_10$ -formyl-THF	

## **Product**

Table 125: Properties of each product.

Id	Name	SBO
$c_{-}$ thf	c_THF	

#### **Kinetic Law**

$$v_{44} = vol\left(cytosol\right) \cdot function\_4\_V\_ART\_1\left(K\_10f\_ART, K\_aic\_ART, Vm\_ART, [aic], [c\_10f]\right) \tag{245}$$

$$\begin{split} & \text{function\_4\_V\_ART\_1}\left(K\_10f\_ART, K\_aic\_ART, Vm\_ART, [aic], [c\_10f]\right) \\ & = \frac{Vm\_ART \cdot [c\_10f] \cdot [aic]}{\left(K\_10f\_ART + [c\_10f]\right) \cdot \left(K\_aic\_ART + [aic]\right)} \end{split} \tag{246}$$

$$\begin{split} & \text{function\_4\_V\_ART\_1} \left( \text{K\_10f\_ART}, \text{K\_aic\_ART}, \text{Vm\_ART}, [\text{aic}], [\text{c\_10f}] \right) \\ & = \frac{\text{Vm\_ART} \cdot [\text{c\_10f}] \cdot [\text{aic}]}{(\text{K\_10f\_ART} + [\text{c\_10f}]) \cdot (\text{K\_aic\_ART} + [\text{aic}])} \end{split} \tag{247}$$

## 9.45 Reaction V\_BHMT

This is a reversible reaction of two reactants forming two products influenced by 13 modifiers.

### Name V\_BHMT

## **Reaction equation**

$$hcy + BET \xrightarrow{\mbox{H2O2, sah, sam, BET, H2O2, hcy, sah, sam, BET, H2O2, hcy, sah, sam} \mbox{met} + dmg \end{math}$$

### **Reactants**

Table 126: Properties of each reactant.

Id	Name	SBO
hcy BET	Homocysteine Betaine	

Table 127: Properties of each modifier.

Id	Name	SBO
H202	H2O2	
sah	SAH	
sam	SAM	
BET	Betaine	
H202	H2O2	
hcy	Homocysteine	
sah	SAH	
sam	SAM	
BET	Betaine	
H202	H2O2	
hcy	Homocysteine	
sah	SAH	
sam	SAM	

## **Products**

Table 128: Properties of each product.

Id	Name	SBO
	c_Methionine Dimethylglycine	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{45} = \text{function\_4\_V\_BHMT\_1} ([BET], [H2O2], K\_bet\_BHMT, K\_hcy\_BHMT, \\ \text{Ki\_BHMT}, \text{Vm\_BHMT}, \text{vol} (\text{cytosol}), [\text{hcy}], [\text{sah}], [\text{sam}], \text{ssH2O2}) \\ \\ \text{function\_4\_V\_BHMT\_1} ([BET], [H2O2], K\_bet\_BHMT, K\_hcy\_BHMT, \\ \text{Ki\_BHMT}, \text{Vm\_BHMT}, \text{vol} (\text{cytosol}), [\text{hcy}], [\text{sah}], [\text{sam}], \text{ssH2O2}) = \text{vol} (\text{cytosol}) \\ \cdot \text{exp} (0.0021 \cdot ([\text{sam}] + [\text{sah}])) \cdot \text{exp} (0.0021 \cdot 102.6) \\ \cdot \text{MM\_twosubst} (\text{Vm\_BHMT}, \text{K\_hcy\_BHMT}, \text{K\_bet\_BHMT}, [\text{hcy}], [BET]) \\ \cdot \frac{\text{ssH2O2} + \text{Ki\_BHMT}}{[\text{H2O2}] + \text{Ki\_BHMT}} \\ \end{cases}$$

### 9.46 Reaction V\_MATI

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

### Name V\_MATI

## **Reaction equation**

$$met \xleftarrow{c\_gsg, \ c\_gsg, \ met, \ sam, \ c\_gsg, \ met, \ sam} sam \tag{251}$$

#### Reactant

Table 129: Properties of each reactant.

Id	Name	SBO
met	c_Methionine	

Table 130: Properties of each modifier.

Id	Name	SBO
c_gsg	c_GSSG	
c_gsg	c_GSSG	
met	c_Methionine	
sam	SAM	
c_gsg	$c\_GSSG$	
met	c_Methionine	
sam	SAM	

#### **Product**

Table 131: Properties of each product.

Id	Name	SBO
sam	SAM	

### **Kinetic Law**

Derived unit contains undeclared units

$$v_{46} = vol\left(cytosol\right) \\ \cdot function\_4\_V\_MATI\_1\left(Ki\_MAT1, Km\_MAT1, Vm\_MAT1, [c\_gsg], [met], [sam]\right)$$
 (252)

$$function\_4\_V\_MATI\_1\left(Ki\_MAT1,Km\_MAT1,Vm\_MAT1,[c\_gsg],[met],[sam]\right)$$

$$= Vm\_MAT1 \cdot \frac{[met]}{Km\_MAT1 + [met]} \cdot (0.23 + 0.8 \cdot exp(0.0026 \cdot [sam])) \cdot \frac{Ki\_MAT1 + 66.71}{Ki\_MAT1 + [c\_gsg]}$$
(253)

$$function\_4\_V\_MATI\_1\left(Ki\_MAT1,Km\_MAT1,Vm\_MAT1,[c\_gsg],[met],[sam]\right)$$

$$= Vm\_MAT1 \cdot \frac{[met]}{Km\_MAT1 + [met]} \cdot (0.23 + 0.8 \cdot exp(0.0026 \cdot [sam])) \cdot \frac{Ki\_MAT1 + 66.71}{Ki\_MAT1 + [c\_gsg]} \tag{254}$$

### 9.47 Reaction V\_MATIII

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

## Name V\_MATIII

# **Reaction equation**

$$met \xrightarrow{c\_gsg, \ c\_gsg, \ met, \ sam, \ c\_gsg, \ met, \ sam} sam \tag{255}$$

## Reactant

Table 132: Properties of each reactant.

Id	Name	SBO
met	c_Methionine	

## **Modifiers**

Table 133: Properties of each modifier.

Id	Name	SBO
c_gsg	c_GSSG	
c_gsg	$c\_GSSG$	
met	c_Methionine	
sam	SAM	
c_gsg	$c\_GSSG$	
met	$c\_Methionine$	
sam	SAM	

## **Product**

Table 134: Properties of each product.

Id	Name	SBO
sam	SAM	

### **Kinetic Law**

$$\begin{split} & \text{function\_4\_V\_MATIII\_1} \left( \text{Ka\_MAT3}, \text{Ki\_MAT3}, \text{Km\_MAT3}, \text{Vm\_MAT3}, [\text{c\_gsg}], [\text{met}], [\text{sam}] \right) \\ & = \text{Vm\_MAT3} \cdot \frac{[\text{met}]^{1.21}}{\text{Km\_MAT3} + [\text{met}]^{1.21}} \cdot \left( 1 + \frac{7.2 \cdot [\text{sam}]^2}{\text{Ka\_MAT3}^2 + [\text{sam}]^2} \right) \cdot \frac{\text{Ki\_MAT3} + 66.71}{\text{Ki\_MAT3} + [\text{c\_gsg}]} \end{aligned} \tag{257}$$

$$\begin{split} & \text{function\_4\_V\_MATIII\_1} \left( \text{Ka\_MAT3}, \text{Ki\_MAT3}, \text{Km\_MAT3}, \text{Vm\_MAT3}, [\text{c\_gsg}], [\text{met}], [\text{sam}] \right) \\ & = \text{Vm\_MAT3} \cdot \frac{[\text{met}]^{1.21}}{\text{Km\_MAT3} + [\text{met}]^{1.21}} \cdot \left( 1 + \frac{7.2 \cdot [\text{sam}]^2}{\text{Ka\_MAT3}^2 + [\text{sam}]^2} \right) \cdot \frac{\text{Ki\_MAT3} + 66.71}{\text{Ki\_MAT3} + [\text{c\_gsg}]} \end{aligned}$$

## 9.48 Reaction V\_GNMT

This is a reversible reaction of two reactants forming two products influenced by nine modifiers.

### Name V\_GNMT

## **Reaction equation**

$$sam + c\_gly \xleftarrow{c\_5mf, c\_5mf, c\_gly, sah, sam, c\_5mf, c\_gly, sah, sam} sah + src \tag{259}$$

#### Reactants

Table 135: Properties of each reactant.

Id	Name	SBO
sam	SAM	
$c_gly$	c_Glycine	

Table 136: Properties of each modifier.

Id	Name	SBO
c_5mf	c_5-methyl-THF	
$c_5mf$	c_5-methyl-THF	
$c_gly$	c_Glycine	
sah	SAH	
sam	SAM	
$\texttt{c\_5mf}$	c_5-methyl-THF	
$c\_{\tt gly}$	c_Glycine	
sah	SAH	
sam	SAM	

|--|

### **Products**

Table 137: Properties of each product.

Id	Name	SBO
sah	SAH	
src	Sarcosine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$\begin{aligned} v_{48} &= \text{function\_4\_V\_GNMT\_1} \left( \text{K\_gly\_GNMT}, \text{K\_sam\_GNMT}, \text{Ki\_GNMT}, \text{Vm\_GNMT}, \\ & & & & & & & & & & & & & & \\ & & & & & & & & & & & & \\ & & & & & & & & & & \\ & & & & & & & & & & \\ & & & & & & & & & \\ & & & & & & & & \\ & & & & & & & \\ & & & & & & & \\ & & & & & & & \\ & & & & & & \\ & & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & &$$

# 9.49 Reaction V\_DNMT

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

### Name V\_DNMT

## **Reaction equation**

$$\operatorname{sam} \frac{\operatorname{sah, sam, sah, sam}}{\operatorname{sah}} \operatorname{sah}$$
 (262)

### Reactant

Table 138: Properties of each reactant.

Id	Name	SBO
sam	SAM	

Table 139: Properties of each modifier.

Id	Name	SBO
sah	SAH	
sam	SAM	
sah	SAH	
sam	SAM	

#### **Product**

Table 140: Properties of each product.

Id	Name	SBO
sah	SAH	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{49} = \text{vol}\left(\text{cytosol}\right) \cdot \text{function\_4\_V\_DNMT\_1}\left(\text{Ki\_DNMT}, \text{Km\_DNMT}, \text{Vm\_DNMT}, [\text{sah}], [\text{sam}]\right)$$
(263)

$$\begin{split} & \text{function\_4\_V\_DNMT\_1}\left(\text{Ki\_DNMT}, \text{Km\_DNMT}, \text{Vm\_DNMT}, [\text{sah}], [\text{sam}]\right) \\ & = \text{Vm\_DNMT} \cdot \frac{[\text{sam}]}{\text{Km\_DNMT} \cdot \left(1 + \frac{[\text{sah}]}{\text{Ki\_DNMT}}\right) + [\text{sam}]} \end{split}$$

$$\begin{split} & \text{function\_4\_V\_DNMT\_1} \left( \text{Ki\_DNMT}, \text{Km\_DNMT}, \text{Vm\_DNMT}, [\text{sah}], [\text{sam}] \right) \\ & = \text{Vm\_DNMT} \cdot \frac{[\text{sam}]}{\text{Km\_DNMT} \cdot \left( 1 + \frac{[\text{sah}]}{\text{Ki\_DNMT}} \right) + [\text{sam}]} \end{aligned}$$

## 9.50 Reaction V\_SAHH

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

# Name V\_SAHH

# Reaction equation

$$sah \xrightarrow{hcy, sah, hcy, sah} hcy$$
 (266)

## Reactant

Table 141: Properties of each reactant.

Id	Name	SBO
sah	SAH	

## **Modifiers**

Table 142: Properties of each modifier.

Id	Name	SBO
hcy	Homocysteine	
$\mathtt{sah}$	SAH	
hcy	Homocysteine	
sah	SAH	

#### **Product**

Table 143: Properties of each product.

Id	Name	SBO
hcy	Homocysteine	

#### **Kinetic Law**

$$v_{50} = vol\left(cytosol\right) \cdot function\_4\_V\_SAHH\_1\left(K\_hcy\_SAHH, K\_sah\_SAHH, Vf\_SAHH, [hcy], [sah]\right)$$
 
$$Vr\_SAHH\_[hcy], [sah]\right)$$
 
$$function\_4\_V\_SAHH\_1\left(K\_hcy\_SAHH, K\_sah\_SAHH, Vf\_SAHH, Vr\_SAHH, [hcy], [sah]\right)$$
 
$$= MM\left(Vf\_SAHH, K\_sah\_SAHH, [sah]\right) - MM\left(Vr\_SAHH, K\_hcy\_SAHH, [hcy]\right)$$
 
$$(268)$$
 
$$function\_4\_V\_SAHH\_1\left(K\_hcy\_SAHH, K\_sah\_SAHH, Vf\_SAHH, Vr\_SAHH, [hcy], [sah]\right)$$
 
$$= MM\left(Vf\_SAHH, K\_sah\_SAHH, [sah]\right) - MM\left(Vr\_SAHH, K\_hcy\_SAHH, [hcy]\right)$$
 
$$(269)$$

# 9.51 Reaction gluconeogenesis\_ser

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

Name gluconeogenesis\_ser

# **Reaction equation**

$$c\_ser \xrightarrow{c\_ser, c\_ser} \emptyset$$
 (270)

### Reactant

Table 144: Properties of each reactant.

Id	Name	SBO
c_ser	c_Serine	

## **Modifiers**

Table 145: Properties of each modifier.

Id	Name	SBO
	c_Serine c_Serine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{51} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_gluconeogenesis\_ser\_1}([\text{c\_ser}])$$
 (271)

function\_4\_gluconeogenesis\_ser\_1([c\_ser]) = 
$$1.2 \cdot [c_ser]$$
 (272)

function\_4\_gluconeogenesis\_ser\_1 ([c\_ser]) = 
$$1.2 \cdot [c_ser]$$
 (273)

## 9.52 Reaction V\_CBS

This is a reversible reaction of two reactants forming one product influenced by 13 modifiers.

Name V\_CBS

# **Reaction equation**

$$hcy + c\_ser \xleftarrow{H2O2, sah, sam, H2O2, c\_ser, hcy, sah, sam, H2O2, c\_ser, hcy, sah, sam} cyt$$

# **Reactants**

Table 146: Properties of each reactant.

Id	Name	SBO
hcy c_ser	Homocysteine c_Serine	

# **Modifiers**

Table 147: Properties of each modifier.

Id	Name	SBO
	Traine	500
H202	H2O2	
$\mathtt{sah}$	SAH	
sam	SAM	
H202	H2O2	
c_ser	c_Serine	
hcy	Homocysteine	
$\mathtt{sah}$	SAH	
sam	SAM	
H202	H2O2	
c_ser	c_Serine	
hcy	Homocysteine	
sah	SAH	
sam	SAM	

# **Product**

Table 148: Properties of each product.

Id	Name	SBO
cyt	Cystathionine	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{52} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_CBS\_1}([\text{H2O2}], \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, \text{Ka\_CBS}, \text{Vm\_CBS}, [\text{c\_ser}], [\text{hcy}], [\text{sah}], [\text{sam}], \text{ssH2O2})$$
 (275)

$$\begin{split} \text{function\_4\_V\_CBS\_1} \, ([\text{H2O2}], \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, \text{Ka\_CBS}, \text{Vm\_CBS}, [\text{c\_ser}], \\ [\text{hcy}], [\text{sah}], [\text{sam}], \text{ssH2O2}) &= \text{MM\_twosubst} \, (\text{Vm\_CBS}, \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, \\ [\text{hcy}], [\text{c\_ser}]) \cdot \frac{\left(\frac{30}{102.59}\right)^2 + 1}{\left(\frac{30}{[\text{sam}] + [\text{sah}]}\right)^2 + 1} \cdot \frac{[\text{H2O2}] + \text{Ka\_CBS}}{\text{ssH2O2} + \text{Ka\_CBS}} \end{split} \eqno(276)$$

$$\begin{split} \text{function\_4\_V\_CBS\_1} & ([\text{H2O2}], \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, \text{Ka\_CBS}, \text{Vm\_CBS}, [\text{c\_ser}], \\ & [\text{hcy}], [\text{sah}], [\text{sam}], \text{ssH2O2}) = \text{MM\_twosubst} (\text{Vm\_CBS}, \text{K\_hcy\_CBS}, \text{K\_ser\_CBS}, \\ & [\text{hcy}], [\text{c\_ser}]) \cdot \frac{\left(\frac{30}{102.59}\right)^2 + 1}{\left(\frac{30}{[\text{sam}] + [\text{sah}]}\right)^2 + 1} \cdot \frac{[\text{H2O2}] + \text{Ka\_CBS}}{\text{ssH2O2} + \text{Ka\_CBS}} \end{split} \eqno(277)$$

### 9.53 Reaction V\_CTGL

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

Name V\_CTGL

### **Reaction equation**

$$cyt \xleftarrow{cyt, cyt} c\_cys$$
 (278)

#### Reactant

Table 149: Properties of each reactant.

Id	Name	SBO
cyt	Cystathionine	

Table 150: Properties of each modifier.

Id	Name	SBO
cyt cyt	Cystathionine Cystathionine	

### **Product**

Table 151: Properties of each product.

Id	Name	SBO
c_cys	c_Cysteine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{53} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_CTGL\_1}(\text{K\_cyt\_CTGL}, \text{Vm\_CTGL}, [\text{cyt}])$$
 (279)

$$function\_4\_V\_CTGL\_1\left(K\_cyt\_CTGL,Vm\_CTGL,[cyt]\right) = \frac{Vm\_CTGL\cdot[cyt]}{K\_cyt\_CTGL+[cyt]} \quad (280)$$

$$function\_4\_V\_CTGL\_1\left(K\_cyt\_CTGL,Vm\_CTGL,[cyt]\right) = \frac{Vm\_CTGL\cdot[cyt]}{K\_cyt\_CTGL+[cyt]} \quad (281)$$

## 9.54 Reaction V\_GCS

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

## Name V\_GCS

### **Reaction equation**

$$c\_cys + c\_glu \xleftarrow{H2O2, c\_gsh, H2O2, c\_cys, c\_glu, c\_gsh, glc, H2O2, c\_cys, c\_glu, c\_gsh, glc} glc \tag{282}$$

### **Reactants**

Table 152: Properties of each reactant.

Cysteine Glutamate	
	Cysteine Glutamate

Table 153: Properties of each modifier.

Id	Name	SBO
H202	H2O2	
$c_gsh$	c_GSH	
H202	H2O2	
c_cys	c_Cysteine	
$c\_{\tt glu}$	c_Glutamate	
$c\_gsh$	c_GSH	
glc	Glutamyl-Cysteine	
H202	H2O2	
c_cys	c_Cysteine	
$c\_{\tt glu}$	$c_Glutamate$	
$c\_gsh$	c_GSH	
glc	Glutamyl-Cysteine	

# **Product**

Table 154: Properties of each product.

	1	1
Id	Name	SBO
glc	Glutamyl-Cysteine	

## **Kinetic Law**

$$\begin{split} \nu_{54} &= vol\left(cytosol\right) \cdot function\_4\_V\_GCS\_1\left([H2O2], K\_cys\_GCS, K\_glu\_GCS, Ka\_GCS, \right. \\ &\quad \left. (283) \right. \\ &\quad \left. \left. \left. \left. \left. \left( E\_GCS, Ki\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gsh], [glc], ssH2O2 \right) \right. \\ &\quad \left. \left. \left( E\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gsh], [glc], ssH2O2 \right) \right. \\ &\quad \left. \left. \left( E\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gsh], [glc], ssH2O2 \right) \right. \\ &\quad \left. \left( E\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gsh], [glc], ssH2O2 \right) \right. \\ &\quad \left. \left( E\_GCS, Kp\_GCS, Vm\_GCS, [c\_cys], [c\_glu], [c\_gl$$

$$\begin{split} & \text{function\_4\_V\_GCS\_1}\left([\text{H2O2}], \text{K\_cys\_GCS}, \text{K\_glu\_GCS}, \text{Ka\_GCS}, \text{Ke\_GCS}, \\ & \text{Ki\_GCS}, \text{Kp\_GCS}, \text{Vm\_GCS}, [\text{c\_cys}], [\text{c\_glu}], [\text{c\_gsh}], [\text{glc}], \text{ssH2O2}\right) \\ & = \frac{\text{Vm\_GCS} \cdot \left([\text{c\_cys}] \cdot [\text{c\_glu}] - \frac{[\text{glc}]}{\text{Ke\_GCS}}\right)}{\text{K\_cys\_GCS} \cdot \text{K\_glu\_GCS} + [\text{c\_glu}] \cdot \text{K\_cys\_GCS} + [\text{c\_cys}] \cdot \left(\text{K\_glu\_GCS} \cdot \left(1 + \frac{[\text{c\_gsh}]}{\text{Ki\_GCS}}\right) + [\text{c\_glu}]\right) + \frac{[\text{glc}]}{\text{Kp\_GCS}} + \frac{[\text{H2O2}] + \text{Ka\_GCS}}{\text{ssH2O2} + \text{Ka\_GCS}}} \\ & \frac{[\text{H2O2}] + \text{Ka\_GCS}}{\text{ssH2O2} + \text{Ka\_GCS}} \\ & \text{function\_4\_V\_GCS\_1}\left([\text{H2O2}], \text{K\_cys\_GCS}, \text{K\_glu\_GCS}, \text{Ka\_GCS}, \text{Ke\_GCS}, \\ & \text{Ki\_GCS}, \text{Kp\_GCS}, \text{Vm\_GCS}, [\text{c\_cys}], [\text{c\_glu}], [\text{c\_gsh}], [\text{glc}], \text{ssH2O2}}\right) \\ & = \frac{\text{Vm\_GCS} \cdot \left([\text{c\_cys}] \cdot [\text{c\_glu}] - \frac{[\text{glc}]}{\text{Ke\_GCS}}\right)}{\text{K\_cys\_GCS} \cdot \text{K\_glu\_GCS} + [\text{c\_glu}] \cdot \text{K\_cys\_GCS} + [\text{c\_cys}] \cdot \left(\text{K\_glu\_GCS} \cdot \left(1 + \frac{[\text{c\_gsh}]}{\text{Ki\_GCS}}\right) + [\text{c\_glu}]\right) + \frac{[\text{glc}]}{\text{Kp\_GCS}} + \\ & [\text{H2O2}] + \text{Ka\_GCS} \end{aligned}$$

## 9.55 Reaction cys\_usage

 $ssH2O2 + Ka\_GCS$ 

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

Name cys\_usage

# **Reaction equation**

$$c_{\text{cys}} \stackrel{c_{\text{cys}}, c_{\text{cys}}}{\longleftarrow} \emptyset$$
 (286)

#### Reactant

Table 155: Properties of each reactant.

Id	Name	SBO
c_cys	c_Cysteine	

Table 156: Properties of each modifier.

Id	Name	SBO
c_cys	c_Cysteine	
c_cys	c_Cysteine	

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{55} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_cys\_usage\_1}([\text{c\_cys}])$$
 (287)

function\_4\_cys\_usage\_1 ([c\_cys]) = 
$$\frac{0.35 \cdot [c_cys]^2}{200}$$
 (288)

function\_4\_cys\_usage\_1 ([c\_cys]) = 
$$\frac{0.35 \cdot [c_cys]^2}{200}$$
 (289)

# 9.56 Reaction c\_glu\_usage

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

Name c\_glu\_usage

### **Reaction equation**

$$c_{glu} \stackrel{c_{glu}, c_{glu}}{====} \emptyset$$
 (290)

## Reactant

Table 157: Properties of each reactant.

Id	Name	SBO
c_glu	$c_{-}Glutamate$	

## **Modifiers**

Table 158: Properties of each modifier.

Id	Name	SBO
_	c_Glutamate c_Glutamate	

### **Kinetic Law**

$$v_{56} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_c\_glu\_usage\_1}([\text{c\_glu}])$$
 (291)

function\_4\_c\_glu\_usage\_1 ([c\_glu]) = 
$$0.07 \cdot [c_glu]$$
 (292)

function\_4\_c\_glu\_usage\_1 ([c\_glu]) = 
$$0.07 \cdot [c_glu]$$
 (293)

# 9.57 Reaction V\_GS

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

## Name V\_GS

## **Reaction equation**

$$glc + c\_gly \xleftarrow{c\_gly, c\_gsh, glc, c\_gly, c\_gsh, glc} c\_gsh$$
 (294)

## **Reactants**

Table 159: Properties of each reactant.

Id	Name	SBO
glc c_gly	Glutamyl-Cysteine c_Glycine	

# **Modifiers**

Table 160: Properties of each modifier.

	2. 1 Toperties of each in	
Id	Name	SBO
c_gly	c_Glycine	
$c\_gsh$	c_GSH	
glc	Glutamyl-Cysteine	
$c\_{gly}$	c_Glycine	
$c\_gsh$	c_GSH	
glc	Glutamyl-Cysteine	

## **Product**

Table 161: Properties of each product.

Id	Name	SBO
c_gsh	c_GSH	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{57} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_V\_GS\_1} (\text{K\_glc\_GS}, \text{K\_gly\_GS}, \text{Ke\_GS}, \text{Kp\_GS}, \text{Vm\_GS}, \\ [\text{c\_gly}], [\text{c\_gsh}], [\text{glc}])$$
(295)

$$\begin{split} & \text{function\_4\_V\_GS\_1} \left( \text{K\_glc\_GS}, \text{K\_gly\_GS}, \text{Ke\_GS}, \text{Kp\_GS}, \text{Vm\_GS}, [\text{c\_gly}], [\text{c\_gsh}], [\text{glc}] \right) \\ & = \frac{\text{Vm\_GS} \cdot \left( [\text{c\_gly}] \cdot [\text{glc}] - \frac{[\text{c\_gsh}]}{\text{Ke\_GS}} \right)}{\text{K\_gly\_GS} \cdot \text{K\_glc\_GS} + [\text{glc}] \cdot \text{K\_gly\_GS} + [\text{c\_gly}] \cdot \left( \text{K\_glc\_GS} + [\text{glc}] \right) + \frac{[\text{c\_gsh}]}{\text{Kp\_GS}}} \end{split}$$

$$\begin{split} & \text{function\_4\_V\_GS\_1} \left( \text{K\_glc\_GS}, \text{K\_gly\_GS}, \text{Ke\_GS}, \text{Kp\_GS}, \text{Vm\_GS}, [\text{c\_gly}], [\text{c\_gsh}], [\text{glc}] \right) \\ & = \frac{\text{Vm\_GS} \cdot \left( [\text{c\_gly}] \cdot [\text{glc}] - \frac{[\text{c\_gsh}]}{\text{Ke\_GS}} \right)}{\text{K\_gly\_GS} \cdot \text{K\_glc\_GS} + [\text{glc}] \cdot \text{K\_gly\_GS} + [\text{c\_gly}] \cdot \left( \text{K\_glc\_GS} + [\text{glc}] \right) + \frac{[\text{c\_gsh}]}{\text{Kp\_GS}}} \end{split}$$

### 9.58 Reaction V\_GPX

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

### Name V\_GPX

## **Reaction equation**

$$2c\_gsh + H2O2 \xrightarrow{H2O2, c\_gsh, H2O2, c\_gsh} c\_gsg$$
 (298)

#### **Reactants**

Table 162: Properties of each reactant.

Id	Name	SBO
c_gsh H2O2	c_GSH H2O2	

Table 163: Properties of each modifier.

Id	Name	SBO
H202	H2O2	
$c\_gsh$	$c\_GSH$	
H202	H2O2	
$c_gsh$	$c_GSH$	

#### **Product**

Table 164: Properties of each product.

Id	Name	SBO
c_gsg	c_GSSG	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{58} = \text{vol}(\text{cytosol})$$

$$\cdot \text{function\_4\_V\_GPX\_1}([\text{H2O2}], \text{K\_H2O2\_GPX}, \text{K\_gsh\_GPX}, \text{Vm\_GPX}, [\text{c\_gsh}])$$
(299)

$$\begin{split} & \text{function\_4\_V\_GPX\_1}\left([\text{H2O2}], \text{K\_H2O2\_GPX}, \text{K\_gsh\_GPX}, \text{Vm\_GPX}, [\text{c\_gsh}]\right) \\ & = \text{Vm\_GPX} \cdot \left(\frac{[\text{c\_gsh}]}{\text{K\_gsh\_GPX} + [\text{c\_gsh}]}\right)^2 \cdot \frac{[\text{H2O2}]}{\text{K\_H2O2\_GPX} + [\text{H2O2}]} \end{split} \tag{300}$$

$$\begin{split} & \text{function\_4\_V\_GPX\_1}\left([\text{H2O2}], \text{K\_H2O2\_GPX}, \text{K\_gsh\_GPX}, \text{Vm\_GPX}, [\text{c\_gsh}]\right) \\ & = \text{Vm\_GPX} \cdot \left(\frac{[\text{c\_gsh}]}{\text{K\_gsh\_GPX} + [\text{c\_gsh}]}\right)^2 \cdot \frac{[\text{H2O2}]}{\text{K\_H2O2\_GPX} + [\text{H2O2}]} \end{aligned} \tag{301}$$

### 9.59 Reaction V\_GR

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

### Name V\_GR

## **Reaction equation**

$$c\_gsg + NADPH \xrightarrow{NADPH, c\_gsg, NADPH, c\_gsg} 2c\_gsh$$
 (302)

### **Reactants**

Table 165: Properties of each reactant.

Id	Name	SBO
c_gsg NADPH	c_GSSG NADPH	

Table 166: Properties of each modifier.

Id	Name	SBO
NADPH	NADPH	
c_gsg	$c\_GSSG$	
NADPH	NADPH	
c_gsg	c_GSSG	

#### **Product**

Table 167: Properties of each product.

Id	Name	SBO
Iu	Name	SDO
c_gsh	c_GSH	

### **Kinetic Law**

$$v_{59} = \text{vol}(\text{cytosol})$$

$$\cdot \text{function\_4\_V\_GR\_1}(\text{K\_NADPH\_GR}, \text{K\_gsg\_GR}, [\text{NADPH}], \text{Vm\_GR}, [\text{c\_gsg}])$$

$$(303)$$

$$\begin{split} & \text{function\_4\_V\_GR\_1}\left(\text{K\_NADPH\_GR}, \text{K\_gsg\_GR}, [\text{NADPH}], \text{Vm\_GR}, [\text{c\_gsg}]\right) \\ & = \frac{\text{Vm\_GR} \cdot [\text{c\_gsg}] \cdot [\text{NADPH}]}{\left(\text{K\_gsg\_GR} + [\text{c\_gsg}]\right) \cdot \left(\text{K\_NADPH\_GR} + [\text{NADPH}]\right)} \end{aligned} \tag{304}$$

$$\begin{aligned} & \text{function\_4\_V\_GR\_1} \left( \text{K\_NADPH\_GR}, \text{K\_gsg\_GR}, [\text{NADPH}], \text{Vm\_GR}, [\text{c\_gsg}] \right) \\ & = \frac{\text{Vm\_GR} \cdot [\text{c\_gsg}] \cdot [\text{NADPH}]}{(\text{K\_gsg\_GR} + [\text{c\_gsg}]) \cdot (\text{K\_NADPH\_GR} + [\text{NADPH}])} \end{aligned} \tag{305}$$

# 9.60 Reaction c\_gsh\_degr

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

Name c\_gsh\_degr

## **Reaction equation**

$$c\_gsh \xrightarrow{c\_gsh, c\_gsh} \emptyset$$
 (306)

#### Reactant

Table 168: Properties of each reactant.

Id	Name	SBO
c_gsh	c_GSH	

### **Modifiers**

Table 169: Properties of each modifier.

Id	Name	SBO
•	c_GSH	
c_gsh	c_GSH	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{60} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_c\_gsh\_degr\_1}([\text{c\_gsh}])$$
 (307)

function\_4\_c\_gsh\_degr\_1 ([c\_gsh]) = 
$$0.0020 \cdot [c_gsh]$$
 (308)

function\_4\_c\_gsh\_degr\_1([c\_gsh]) = 
$$0.0020 \cdot [c_gsh]$$
 (309)

# 9.61 Reaction c\_gsg\_degr

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

Name c\_gsg\_degr

### **Reaction equation**

$$c_{-gsg} = \underbrace{c_{-gsg}, c_{-gsg}}_{(310)} \emptyset$$

### Reactant

Table 170: Properties of each reactant.

Id	Name	SBO
c_gsg	c_GSSG	

## **Modifiers**

Table 171: Properties of each modifier.

Id	Name	SBO
c_gsg c_gsg	c_GSSG c_GSSG	

## **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{61} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_c\_gsg\_degr\_1}([\text{c\_gsg}])$$
 (311)

function\_4\_c\_gsg\_degr\_1 ([c\_gsg]) = 
$$0.1 \cdot [c_gsg]$$
 (312)

function\_4\_c\_gsg\_degr\_1 ([c\_gsg]) = 
$$0.1 \cdot [c_gsg]$$
 (313)

## 9.62 Reaction reaction\_1

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

## Name $V\_ODC$

# **Reaction equation**

### Reactant

Table 172: Properties of each reactant.

Id	Name	SBO
species_7	Ornithine	

Table 173: Properties of each modifier.

Id	Name	SBO
species_2	Putrescine	
species_7	Ornithine	
species_2	Putrescine	
${\tt species\_7}$	Ornithine	

# **Product**

Table 174: Properties of each product.

Id	Name	SBO
species_2	Putrescine	

### **Kinetic Law**

$$v_{62} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_V\_ODC\_1} (\text{Kipodc}, \text{Kmodc}, \text{parameter\_1}, [\text{species\_2}], [\text{species\_7}])$$
(315)

$$\begin{aligned} & \text{function\_4\_V\_ODC\_1} \left( \text{Kipodc}, \text{Kmodc}, \text{parameter\_1}, [\text{species\_2}], [\text{species\_7}] \right) \\ &= \frac{\text{parameter\_1} \cdot [\text{species\_7}]}{\text{Kmodc} \cdot \left( 1 + \frac{[\text{species\_2}]}{\text{Kipodc}} \right) + [\text{species\_7}]} \end{aligned}$$

Table 175: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
Kipodc	Kipodc	1300.0	
Kmodc	Kmodc	60.0	$   \overline{\mathcal{L}} $

# 9.63 Reaction reaction\_2

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

## Name V\_SAMDC

# **Reaction equation**

## Reactant

Table 176: Properties of each reactant.

Id	Name	SBO
sam	SAM	

### **Modifiers**

Table 177: Properties of each modifier.

Id	Name	SBO
species_3	Spermine	
species_2	Putrescine	
sam	SAM	
species_1	dcSAM	
species_2	Putrescine	
species_3	Spermine	
sam	SAM	
species_1	dcSAM	
species_2	Putrescine	
species_3	Spermine	

### **Product**

Table 178: Properties of each product.

Id	Name	SBO
species_1	dcSAM	

#### **Kinetic Law**

#### **Derived unit** contains undeclared units

$$v_{63} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_V\_SAMDC\_1} (\text{Kapsamdc}, \text{Kiasamdc}, \text{Kissamdc}, \text{Kissamdc}, \text{parameter\_3}, [\text{sam}], [\text{species\_1}], [\text{species\_2}], [\text{species\_3}])$$
(319)

function\_4\_V\_SAMDC\_1 (Kapsamdc, Kiasamdc, Kissamdc,

$$\begin{split} & \text{Kmsamdc}, \text{parameter\_3}, [\text{sam}], [\text{species\_1}], [\text{species\_2}], \\ & \underbrace{\frac{\text{parameter\_3}}{1 + \frac{[\text{species\_3}]}{\text{Kissamdc}}} \cdot [\text{sam}]}_{1 + \frac{[\text{species\_1}]}{\text{Kissamdc}}} + \underbrace{[\text{species\_1}]}_{\text{Kiasamdc}}) + [\text{sam}] \end{split} \tag{320}$$

function\_4\_V\_SAMDC\_1 (Kapsamdc, Kiasamdc, Kissamdc, Kmsamdc, parameter\_3, [sam], [species\_1], [species\_2],

parameter\_3 . [sam]

$$[species\_3]) = \frac{\frac{parameter\_3}{1 + \frac{[species\_3]}{Kissamdc}} \cdot [sam]}{Kmsamdc \cdot \left(1 + \frac{Kapsamdc}{[species\_2]} + \frac{[species\_1]}{Kiasamdc}\right) + [sam]}$$
(321)

Table 179: Properties of each parameter.

Name	SBO	Value	Unit	Constant
Kapsamdc		0.5		
Kiasamdc		2.5		
Kissamdc		500.0		
Kmsamdc		50.0		
	Kapsamdc Kiasamdc Kissamdc	Kapsamdc Kiasamdc Kissamdc	Kapsamdc0.5Kiasamdc2.5Kissamdc500.0	Kapsamde0.5Kiasamde2.5Kissamde500.0

### 9.64 Reaction reaction\_3

This is a reversible reaction of two reactants forming two products influenced by nine modifiers.

Name V\_SSAT\_S

### **Reaction equation**

#### **Reactants**

Table 180: Properties of each reactant.

Id	Name	SBO
species_3 species_8	Spermine Acetyl-CoA	

### **Modifiers**

Table 181: Properties of each modifier.

Id	Name	SBO
species_4	Spermidine	
species_3	Spermine	
${ t species\_4}$	Spermidine	
species_8	Acetyl-CoA	
species_9	CoA	
species_3	Spermine	
${\tt species\_4}$	Spermidine	
species_8	Acetyl-CoA	
${\tt species\_9}$	CoA	

#### **Products**

Table 182: Properties of each product.

Id	Name	SBO
species_5 species_9	Acetylspermine CoA	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{64} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_V\_SSAT\_S\_1} (C, \text{Kmaccoassat}, \text{Kmcoassat}, \text{Kmdssat}, \\ \text{Kmsssat}, \text{parameter\_2}, [\text{species\_3}], [\text{species\_4}], [\text{species\_8}], [\text{species\_9}])$$

$$\text{function\_4\_V\_SSAT\_S\_1} (C, \text{Kmaccoassat}, \text{Kmcoassat}, \text{Kmdssat}, \\ \text{Kmsssat}, \text{parameter\_2}, [\text{species\_3}], [\text{species\_4}], [\text{species\_9}])$$

$$= \frac{\frac{1}{C} \cdot \text{parameter\_2} \cdot [\text{species\_3}] \cdot [\text{species\_8}]}{\text{Kmsssat} \cdot \left(1 + \frac{[\text{species\_9}]}{\text{Kmdssat}}\right) \cdot \text{Kmaccoassat} \cdot \left(1 + \frac{[\text{species\_9}]}{\text{Kmcoassat}}\right) \cdot [\text{species\_3}] + \text{Kmssat}}$$

$$\frac{1}{C} \cdot parameter_2 \cdot [species_3] \cdot [species_8]$$

$$= \frac{\frac{1}{C} \cdot parameter\_2 \cdot [species\_3] \cdot [species\_8]}{Kmsssat \cdot \left(1 + \frac{[species\_4]}{Kmdssat}\right) \cdot Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) + Kmaccoassat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot [species\_3] + Kmssat \cdot \left(1 + \frac{[species\_4]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left(1 + \frac{[species\_9]}{Kmcoassat}\right) \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmsat \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right] + Kmssat \cdot \left[1 + \frac{[species\_9]}{Kmcoassat}\right$$

Table 183: Properties of each parameter.

Id	Name	SBO Value	Unit	Constant
С	C	4.44		
Kmaccoassat	Kmaccoassat	1.50		
Kmcoassat	Kmcoassat	40.00		
Kmdssat	Kmdssat	130.00		
Kmsssat	Kmsssat	35.00		Ø

### 9.65 Reaction reaction\_4

This is a reversible reaction of two reactants forming two products influenced by nine modifiers.

Name V\_SSAT\_D

### **Reaction equation**

### **Reactants**

Table 184: Properties of each reactant.

Id	Name	SBO
species_4 species_8	Spermidine Acetyl-CoA	

Table 185: Properties of each modifier.

Id	Name	SBO
species_3	Spermine	
species_3	Spermine	

Id	Name	SBO
species_4 species_8 species_9	Spermidine Acetyl-CoA CoA	
species_3 species_4 species_8 species_9	Spermine Spermidine Acetyl-CoA CoA	

### **Products**

Table 186: Properties of each product.

	1 1	
Id	Name	SBO
species_6 species_9	Acetylspermidine CoA	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{65} = \text{vol} (\text{cytosol}) \cdot \text{function\_4\_V\_SSAT\_D\_1} (\text{Kmaccoassat}, \text{Kmcoassat}, \text{Kmdssat}, \\ \text{Kmsssat}, \text{parameter\_2}, [\text{species\_3}], [\text{species\_4}], [\text{species\_8}], [\text{species\_9}])$$

$$\text{function\_4\_V\_SSAT\_D\_1} (\text{Kmaccoassat}, \text{Kmcoassat}, \text{Kmdssat}, \\ \text{Kmsssat}, \text{parameter\_2}, [\text{species\_3}], [\text{species\_4}], [\text{species\_8}], [\text{species\_9}])$$

$$= \frac{\text{parameter\_2} \cdot [\text{species\_4}] \cdot [\text{species\_8}]}{\text{Kmdssat} \cdot \left(1 + \frac{[\text{species\_3}]}{\text{Kmsssat}}\right) \cdot \text{Kmaccoassat} \cdot \left(1 + \frac{[\text{species\_9}]}{\text{Kmcoassat}}\right) \cdot [\text{species\_4}] + \text{Kmd}}$$

$$\text{function\_4\_V\_SSAT\_D\_1} (\text{Kmaccoassat}, \text{Kmcoassat}, \text{Kmdssat}, \\ \text{Kmsssat}, \text{parameter\_2}, [\text{species\_3}], [\text{species\_4}], [\text{species\_9}])$$

$$= \frac{\text{parameter\_2} \cdot [\text{species\_4}] \cdot [\text{species\_8}]}{\text{Kmdssat} \cdot \left(1 + \frac{[\text{species\_9}]}{\text{Kmcoassat}}\right) \cdot \text{Kmaccoassat} \cdot \left(1 + \frac{[\text{species\_9}]}{\text{Kmcoassat}}\right) \cdot [\text{species\_4}] + \text{Kmd}}$$

Table 187: Properties of each parameter.

Id	Name	SBO Value	Unit	Constant
Kmaccoassat	Kmaccoassat	1.5		
Kmcoassat	Kmcoassat	40.0		
Kmdssat	Kmdssat	130.0		$\overline{\mathbf{Z}}$
Kmsssat	Kmsssat	35.0	,	$\overline{\checkmark}$
112	r	COURCE BY SOIVIES FIE	1	

# 9.66 Reaction reaction\_5

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

Name V\_PAO\_AD

# **Reaction equation**

species\_6, species\_4, species\_3, species\_4, species\_5, species\_6, species\_6, species\_4, species\_5

(330)

### Reactant

Table 188: Properties of each reactant.

Id	Name	SBO
species_6	Acetylspermidine	

### **Modifiers**

Table 189: Properties of each modifier.

	I	
Id	Name	SBO
species_5	Acetylspermine	
species_4	Spermidine	
species_3	Spermine	
species_3	Spermine	
species_4	Spermidine	
species_5	Acetylspermine	
species_6	Acetylspermidine	
species_3	Spermine	
species_4	Spermidine	
species_5	Acetylspermine	
species_6	Acetylspermidine	

### **Product**

Table 190: Properties of each product.

Id	Name	SBO
species_2	Putrescine	

### **Kinetic Law**

#### **Derived unit** contains undeclared units

$$v_{66} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_PAO\_AD\_1} \text{ (Kmadpao, Kmaspao, Kmdpao, Kmspao, Vmpao, [species\_3], [species\_4], [species\_5], [species\_6])}$$

$$[\text{species\_6}]) = \frac{\text{Vmpao} \cdot [\text{species\_6}]}{\text{Kmadpao} \cdot \left(1 + \frac{[\text{species\_6}]}{\text{Kmadpao}} + \frac{[\text{species\_5}]}{\text{Kmaspao}} + \frac{[\text{species\_4}]}{\text{Kmdpao}} + \frac{[\text{species\_3}]}{\text{Kmspao}}\right)}$$
(332)

function\_4\_V\_PAO\_AD\_1 (Kmadpao, Kmaspao, Kmdpao,

$$[\text{species\_6}]) = \frac{\text{Vmpao} \cdot [\text{species\_6}]}{\text{Kmadpao} \cdot \left(1 + \frac{[\text{species\_6}]}{\text{Kmadpao}} + \frac{[\text{species\_5}]}{\text{Kmaspao}} + \frac{[\text{species\_4}]}{\text{Kmdpao}} + \frac{[\text{species\_3}]}{\text{Kmspao}}\right)}$$
(333)

Table 191: Properties of each parameter.

		1	
Id	Name	SBO Value Unit	Constant
Kmadpao	Kmadpao	14.0	$ \mathcal{L} $
Kmaspao	Kmaspao	0.6	
Kmdpao	Kmdpao	50.0	
Kmspao	Kmspao	15.0	$\square$
Vmpao	Vmpao	621.0	$\square$

### 9.67 Reaction reaction\_6

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

Name V\_PAO\_AS

#### Reaction equation

# Reactant

Table 192:	Properties o	of each	reactant.
------------	--------------	---------	-----------

Tuble 192: Troperties of cuent reactant.			
Id	Name	SBO	
species_5	Acetylspermine		

# **Modifiers**

Table 193: Properties of each modifier.

	1	
Id	Name	SBO
species_6	Acetylspermidine	
species_3	Spermine	
species_3	Spermine	
${\tt species\_4}$	Spermidine	
species_5	Acetylspermine	
species_6	Acetylspermidine	
species_3	Spermine	
species_4	Spermidine	
species_5	Acetylspermine	
${\tt species\_6}$	Acetylspermidine	

### **Product**

Table 194: Properties of each product.

Id	Name	SBO
species_4	Spermidine	

# **Kinetic Law**

# **Derived unit** contains undeclared units

$$v_{67} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_PAO\_AS\_1}(\text{Kmadpao}, \text{Kmaspao}, \text{Kmdpao}, \text{Kmspao}, \text{Vmpao}, [\text{species\_4}], [\text{species\_5}], [\text{species\_6}])$$

$$\begin{split} & \text{function\_4\_V\_PAO\_AS\_1} \text{ (Kmadpao, Kmaspao, Kmdpao,} \\ & \text{Kmspao, Vmpao, [species\_3], [species\_4], [species\_5],} \\ & \text{[species\_6])} = \frac{\text{Vmpao} \cdot [\text{species\_5}]}{\text{Kmaspao} \cdot \left(1 + \frac{[\text{species\_6}]}{\text{Kmadpao}} + \frac{[\text{species\_5}]}{\text{Kmaspao}} + \frac{[\text{species\_3}]}{\text{Kmdpao}} + \frac{[\text{species\_3}]}{\text{Kmspao}} \right)} \end{split}$$

$$\begin{aligned} & \text{function\_4\_V\_PAO\_AS\_1} \text{ (Kmadpao, Kmaspao, Kmdpao,} \\ & \text{Kmspao, Vmpao, [species\_3], [species\_4], [species\_5],} \\ & \text{[species\_6])} = \frac{\text{Vmpao} \cdot [\text{species\_5}]}{\text{Kmaspao} \cdot \left(1 + \frac{[\text{species\_6}]}{\text{Kmadpao}} + \frac{[\text{species\_5}]}{\text{Kmaspao}} + \frac{[\text{species\_4}]}{\text{Kmdpao}} + \frac{[\text{species\_3}]}{\text{Kmspao}} \right)} \end{aligned} \tag{337}$$

Table 195: Properties of each parameter.

		1	
Id	Name	SBO Value Unit	Constant
Kmadpao	Kmadpao	14.0	
Kmaspao	Kmaspao	0.6	
Kmdpao	Kmdpao	50.0	
Kmspao	Kmspao	15.0	
Vmpao	Vmpao	621.0	$\square$

# 9.68 Reaction reaction\_7

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

# Name V\_SPDS

# **Reaction equation**

#### **Reactants**

Table 196: Properties of each reactant.

Id	Name	SBO
species_1 species_2		

Table 197: Properties of each modifier.

Id	Name	SBO
species_1	dcSAM	

Id	Name	SBO
species_2	Putrescine	
${ t species\_4}$	Spermidine	
${ t species\_1}$	dcSAM	
species_2	Putrescine	
species_4	Spermidine	

### **Product**

Table 198: Properties of each product.

Id	Name	SBO
species_4	Spermidine	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{68} = \text{vol}\left(\text{cytosol}\right) \cdot \text{function\_4\_V\_SPDS\_1}\left(\text{KaSpds}, \text{Kiaspds}, \text{Kidspds}, \text{Kpspds}, \text{Vmspds}, [\text{species\_2}], [\text{species\_2}], [\text{species\_4}]\right)}$$

$$\text{function\_4\_V\_SPDS\_1}\left(\text{KaSpds}, \text{Kiaspds}, \text{Kidspds}, \text{Kidspds}, \text{Kpspds}, [\text{species\_2}], [\text{species\_4}]\right)$$

$$= \frac{\text{Vmspds} \cdot [\text{species\_4}] \cdot [\text{species\_4}]}{\text{Kiaspds} \cdot \text{Kpspds} \cdot (1 + \frac{[\text{species\_4}]}{\text{Kidspds}})} + \text{Kpspds} \cdot [\text{species\_1}] + \text{KaSpds} \cdot \left(1 + \frac{[\text{species\_4}]}{\text{Kidspds}}\right) \cdot [\text{species\_2}] + [\text{species\_1}]}$$

$$\text{function\_4\_V\_SPDS\_1}\left(\text{KaSpds}, \text{Kiaspds}, \text{Kidspds}, \text{Kidspds}, \text{Kidspds}, \text{Kidspds}, \text{Kpspds}, [\text{species\_2}], [\text{species\_4}]\right)}$$

$$= \frac{\text{Vmspds} \cdot [\text{species\_4}]}{\text{Kiaspds} \cdot \text{Kpspds} \cdot [\text{species\_4}]} \cdot [\text{species\_4}] \cdot [\text{species\_4}]$$

$$= \frac{\text{Vmspds} \cdot [\text{species\_4}]}{\text{Kiaspds} \cdot \text{Kpspds} \cdot [\text{species\_4}]} \cdot [\text{species\_4}] \cdot [\text{species\_4}]} \cdot [\text{species\_2}] + [\text{species\_1}]}$$

Table 199: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
KaSpds	KaSpds	0.3	
Kiaspds	Kiaspds	0.8	$\square$
Kidspds	Kidspds	100.0	$\square$
Kpspds	Kpspds	40.0	$\square$
Vmspds	Vmspds	657.0	$\square$

# 9.69 Reaction reaction\_8

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

### Name V\_SPMS

# **Reaction equation**

### **Reactants**

Table 200: Properties of each reactant.

Id	Name	SBO
species_1 species_4	dcSAM Spermidine	

# **Modifiers**

Table 201: Properties of each modifier.

Id	Name	SBO
species_1	dcSAM	
species_3	Spermine	
species_4	Spermidine	
species_1	dcSAM	
species_3	Spermine	
${\tt species\_4}$	Spermidine	

# **Product**

Table 202: Properties of each product.

Id	Name	SBO
species_3	Spermine	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{69} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_SPMS\_1}(\text{Kaspms}, \text{Kdspms}, \text{Kiaspms}, \text{Kisspms}, \text{Vmspms}, [\text{species\_4}], [\text{species\_4}])$$
 (343)

(344)

Kisspms, Vmspms, [species\_1], [species\_3], [species\_4])

 $Vmspms \cdot [species\_1] \cdot [species\_4]$ 

$$\frac{\text{Vmspms} \cdot [\text{species\_1}] \cdot [\text{species\_4}]}{\text{Kiaspms} \cdot \text{Kdspms} \cdot \left(1 + \frac{[\text{species\_3}]}{\text{Kisspms}}\right) + \text{Kdspms} \cdot [\text{species\_1}] + \text{Kaspms} \cdot \left(1 + \frac{[\text{species\_3}]}{\text{Kisspms}}\right) \cdot [\text{species\_4}] + [\text{species\_4}]}$$

function\_4\_V\_SPMS\_1 (Kaspms, Kdspms, Kiaspms,

(345)

Kisspms, Vmspms, [species\_1], [species\_3], [species\_4])

$$\frac{V m spms \cdot [species\_1] \cdot [species\_4]}{Kiaspms \cdot K d spms \cdot \left(1 + \frac{[species\_3]}{Kisspms}\right) + K d spms \cdot [species\_1] + K a spms \cdot \left(1 + \frac{[species\_3]}{Kisspms}\right) \cdot [species\_4] + [species\_4]}$$

Table 203: Properties of each parameter.

Id	Name	SBO Value	Unit Constant
Kaspms	Kaspms	0.10	<b>Z</b>
Kdspms	Kdspms	60.00	
Kiaspms	Kiaspms	0.06	
Kisspms	Kisspms	25.00	
Vmspms	Vmspms	193.80	

### 9.70 Reaction reaction\_9

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

# Name V\_COA

### **Reaction equation**

species\_8 
$$\xrightarrow{\text{species}\_8, \text{ species}\_8}$$
 species\_9 (346)

#### Reactant

Table 204: Properties of each reactant.

Id	Name	SBO
species_8	Acetyl-CoA	

### **Modifiers**

Table 205: Properties of each modifier.

Id	Name	SBO
-	Acetyl-CoA Acetyl-CoA	

### **Product**

Table 206: Properties of each product.

Id	Name	SBO
species_9	CoA	

# **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{70} = \text{vol}(\text{cytosol}) \cdot \text{parameter} \cdot 15 \cdot [\text{species} \cdot 8]$$
 (347)

# 9.71 Reaction reaction\_10

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

# Name V\_ACCOA

# **Reaction equation**

species\_9 
$$\xrightarrow{\text{species}\_9, \text{ species}\_9}$$
 species\_8 (348)

# Reactant

Table 207: Properties of each reactant.

Id	Name	SBO
species_9	CoA	

Table 208: Properties of each modifier.

Name	SBO
CoA	
	Name CoA CoA

# **Product**

Table 209: Properties of each product.

Id	Name	SBO
species_8	Acetyl-CoA	

### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{71} = \text{vol}(\text{cytosol}) \cdot \text{parameter}\_14 \cdot [\text{species}\_9]$$
 (349)

# 9.72 Reaction reaction\_11

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

Name V\_PUT\_efflux

# **Reaction equation**

species\_2 
$$\xrightarrow{\text{species}\_2, \text{ species}\_2} \emptyset$$
 (350)

# Reactant

Table 210: Properties of each reactant.

Id	Name	SBO
species_2	Putrescine	

Table 211: Properties of each modifier.

Id	Name	SBO
species_2 species_2		

# **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{72} = \text{vol}(\text{cytosol}) \cdot \text{k1} \cdot [\text{species}\_2]$$
 (351)

Table 212: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
k1	k1	0.6	

# 9.73 Reaction reaction\_12

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

# Name $V\_AD\_efflux$

# **Reaction equation**

species\_6 
$$\xrightarrow{\text{species}\_6, \text{ species}\_6} \emptyset$$
 (352)

# Reactant

Table 213: Properties of each reactant.

Id	Name	SBO
species_6	Acetylspermidine	

Table 214: Properties of each modifier.

Id	Name	SBO
species_6	Acetylspermidine	
species_6	Acetylspermidine	

Id	Name	SBO

# **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{73} = \text{vol}(\text{cytosol}) \cdot \text{k1} \cdot [\text{species\_6}]$$
 (353)

Table 215: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
k1	k1	0.6	

# 9.74 Reaction reaction\_13

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

# Name V\_MATII

# **Reaction equation**

$$met \xrightarrow{sam, met, sam, met, sam} sam$$
 (354)

# Reactant

Table 216: Properties of each reactant.

Id	Name	SBO
met	c_Methionine	

Table 217: Properties of each modifier.

Id	Name	SBO
sam	SAM	
met	c_Methionine	
sam	SAM	
met	c_Methionine	
sam	SAM	

### **Product**

Table 218: Properties of each product.

Id	Name	SBO
sam	SAM	

#### **Kinetic Law**

**Derived unit** contains undeclared units

$$v_{74} = \text{vol}(\text{cytosol}) \cdot \text{function\_4\_V\_MATII}(\text{Ki}, \text{Km}, [\text{met}], \text{parameter\_19}, [\text{sam}])$$
 (355)

$$function\_4\_V\_MATII(Ki,Km,[met],parameter\_19,[sam]) = \frac{parameter\_19}{1 + \left(\frac{Km \cdot \left(1 + \frac{[sam]}{Ki}\right)}{[met]}\right)^{0.76}}$$
(356)

$$function\_4\_V\_MATII(Ki, Km, [met], parameter\_19, [sam]) = \frac{parameter\_19}{1 + \left(\frac{Km \cdot \left(1 + \frac{[sam]}{Ki}\right)}{[met]}\right)^{0.76}} \quad (357)$$

Table 219: Properties of each parameter.

Id	Name	SBO Value Unit	Constant
Ki	Ki	50.0	$ \overline{\checkmark} $
Km	Km	4.0	$\overline{\checkmark}$

# 10 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.

# 10.1 Species b\_met

Name b\_Methionine

Initial concentration  $30 \mu mol \cdot l^{-1}$ 

Involved in rule b\_met

This species takes part in three reactions (as a reactant in V\_b\_MET\_c and as a modifier in V\_b\_MET\_c, V\_b\_MET\_c). Not these but one rule determines the species' quantity because this species is on the boundary of the reaction system.

### 10.2 Species b\_ser

Name b\_Serine

Initial concentration  $150 \mu mol \cdot l^{-1}$ 

Involved in rule b ser

This species takes part in three reactions (as a reactant in V\_b\_SER\_c and as a modifier in V\_b\_SER\_c, V\_b\_SER\_c). Not these but one rule determines the species' quantity because this species is on the boundary of the reaction system.

# 10.3 Species b\_gly

Name b\_Glycine

Initial concentration  $218.733171504338 \, \mu \text{mol} \cdot 1^{-1}$ 

This species takes part in nine reactions (as a reactant in b\_gly\_loss, V\_b\_GLY\_c and as a product in b\_gsh\_decomp, b\_gsg\_decomp, b\_gly\_import and as a modifier in b\_gly\_loss, b\_gly\_loss, V\_b\_GLY\_c, V\_b\_GLY\_c).

$$\frac{d}{dt}b_{-g}ly = v_1 + 2v_2 + v_{10} - v_6 - v_{18}$$
 (358)

# 10.4 Species b\_glu

Name b\_Glutamate

Initial concentration  $60.4651616225031 \ \mu mol \cdot l^{-1}$ 

This species takes part in nine reactions (as a reactant in b\_glu\_loss, V\_b\_GLU\_c and as a product in b\_gsh\_decomp, b\_gsg\_decomp, b\_glu\_import and as a modifier in b\_glu\_loss, b\_glu\_loss, V\_b\_GLU\_c, V\_b\_GLU\_c).

$$\frac{\mathrm{d}}{\mathrm{d}t} b_{-} \mathrm{glu} = |v_1| + 2|v_2| + |v_{11}| - |v_5| - |v_{17}| \tag{359}$$

# 10.5 Species b\_cys

Name b\_Cysteine

Initial concentration  $183.099466381356 \ \mu mol \cdot l^{-1}$ 

This species takes part in twelve reactions (as a reactant in b\_cys\_cystine\_conv, b\_cys\_loss, V\_b\_CYS\_c and as a product in b\_gsh\_decomp, b\_gsg\_decomp, b\_cys\_import and as a modifier in b\_cys\_cystine\_conv, b\_cys\_cystine\_conv, b\_cys\_loss, b\_cys\_loss, V\_b\_CYS\_c, V\_b-\_CYS\_c).

$$\frac{d}{dt}b_{-}cys = v_1 + 2v_2 + v_9 - v_3 - v_4 - v_{16}$$
(360)

# 10.6 Species b\_gsg

Name b\_GSSG

Initial concentration  $0.472632922783833 \ \mu mol \cdot l^{-1}$ 

This species takes part in eight reactions (as a reactant in b\_gsg\_decomp, b\_gsg\_loss and as a product in V\_c\_gsgHb, V\_c\_gsgLb and as a modifier in b\_gsg\_decomp, b\_gsg\_decomp, b\_gsg\_loss, b\_gsg\_loss).

$$\frac{d}{dt}b_{-}gsg = v_{14} + v_{15} - v_{2} - v_{8}$$
 (361)

# 10.7 Species b\_gsh

Name b\_GSH

Initial concentration  $12.5470655822207 \, \mu \text{mol} \cdot 1^{-1}$ 

This species takes part in eight reactions (as a reactant in b\_gsh\_decomp, b\_gsh\_loss and as a product in V\_c\_gshHb, V\_c\_gshLb and as a modifier in b\_gsh\_decomp, b\_gsh\_decomp, b\_gsh\_loss, b\_gsh\_loss).

$$\frac{d}{dt}b_{-}gsh = |v_{12}| + |v_{13}| - |v_{1}| - |v_{7}|$$
(362)

# 10.8 Species GAR

Name GAR

Initial concentration  $10 \ \mu mol \cdot l^{-1}$ 

This species takes part in three reactions (as a reactant in V\_PGT and as a modifier in V\_PGT, V\_PGT), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{GAR} = 0\tag{363}$$

# 10.9 Species NADPH

Name NADPH

Initial concentration  $50 \ \mu mol \cdot l^{-1}$ 

This species takes part in ten reactions (as a reactant in V\_DHFR, V\_MTHFR, V\_GR and as a product in VcMTD and as a modifier in V\_DHFR, V\_DHFR, V\_MTHFR, V\_MTHFR, V\_GR, V\_GR), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{NADPH} = 0\tag{364}$$

# 10.10 Species BET

Name Betaine

Initial concentration  $50 \ \mu mol \cdot l^{-1}$ 

This species takes part in three reactions (as a reactant in V\_BHMT and as a modifier in V\_BHMT, V\_BHMT), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{BET} = 0\tag{365}$$

### 10.11 Species DUMP

Name dUMP

Initial concentration  $20 \mu mol \cdot l^{-1}$ 

This species takes part in three reactions (as a reactant in V\_TS and as a modifier in V\_TS, V\_TS), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{DUMP} = 0\tag{366}$$

# **10.12 Species** H202

Name H2O2

Initial concentration  $0.015 \ \mu mol \cdot l^{-1}$ 

This species takes part in 15 reactions (as a reactant in V\_GPX and as a modifier in V\_MS, V\_MS, V\_MS, V\_BHMT, V\_BHMT, V\_BHMT, V\_CBS, V\_CBS, V\_CBS, V\_GCS, V\_GCS, V\_GCS, V\_GPX, V\_GPX), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{H2O2} = 0\tag{367}$$

# 10.13 Species c\_thf

Name c\_THF

Initial concentration  $4.0250933022918 \ \mu mol \cdot l^{-1}$ 

### Initial assignment c\_thf

This species takes part in 14 reactions (as a reactant in VcFTS, VcSHMT, VcNE and as a product in V\_MS, V\_DHFR, VcFTD, V\_PGT, V\_ART and as a modifier in VcFTS, VcFTS, VcSHMT, VcNE, VcNE, VcNE).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}thf = |v_{33}| + |v_{34}| + |v_{35}| + |v_{36}| + |v_{44}| - |v_{37}| - |v_{38}| - |v_{39}|$$
(368)

# 10.14 Species c\_5mf

**Name** c\_5-methyl-THF

Initial concentration  $5.36079859230553 \ \mu mol \cdot l^{-1}$ 

This species takes part in seven reactions (as a reactant in V\_MS and as a product in V\_MTHFR and as a modifier in V\_MS, V\_MS, V\_GNMT, V\_GNMT, V\_GNMT).

$$\frac{d}{dt}c_{-}5mf = v_{41} - v_{33} \tag{369}$$

# 10.15 Species c\_2cf

Name c\_5-10-methylene-THF

Initial concentration  $0.463962655701761 \ \mu mol \cdot l^{-1}$ 

This species takes part in 15 reactions (as a reactant in V\_TS, V\_MTHFR, VcMTD and as a product in VcSHMT, VcNE and as a modifier in VcSHMT, VcNE, VcNE, V\_TS, V\_MTHFR, V\_MTHFR, VcMTD, VcMTD).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{2}cf = v_{38} + v_{39} - v_{40} - v_{41} - v_{42} \tag{370}$$

### 10.16 Species c\_1cf

Name c\_5-10-methenyl-THF

Initial concentration  $0.259497676806752 \mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in VcMTCH and as a product in VcMTD and as a modifier in VcMTD, VcMTD, VcMTCH, VcMTCH).

$$\frac{d}{dt}c_{-}1cf = |v_{42}| - |v_{43}| \tag{371}$$

# **10.17 Species** c\_10f

Name c\_10-formyl-THF

Initial concentration  $3.25539505319571 \ \mu mol \cdot l^{-1}$ 

This species takes part in 13 reactions (as a reactant in VcFTD, V\_PGT, V\_ART and as a product in VcFTS, VcMTCH and as a modifier in VcFTD, VcFTD, V\_PGT, VcMTCH, VcMTCH, VcMTCH, V\_ART, V\_ART).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}10f = v_{37} + v_{43} - v_{35} - v_{36} - v_{44} \tag{372}$$

# 10.18 Species c\_dhf

Name c\_DHF

Initial concentration  $0.0352527196984464 \mu mol \cdot l^{-1}$ 

This species takes part in four reactions (as a reactant in V\_DHFR and as a product in V\_TS and as a modifier in V\_DHFR, V\_DHFR).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{c}_{-}\mathrm{dhf} = v_{40} - v_{34} \tag{373}$$

### 10.19 Species aic

Name AICAR

Initial concentration  $0.944676738309717 \mu mol \cdot l^{-1}$ 

This species takes part in four reactions (as a reactant in V\_ART and as a product in V\_PGT and as a modifier in V\_ART, V\_ART).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{aic} = v_{36} - v_{44} \tag{374}$$

# 10.20 Species c\_glu

Name c\_Glutamate

Initial concentration  $3236.78229408139 \, \mu \text{mol} \cdot l^{-1}$ 

This species takes part in nine reactions (as a reactant in V\_GCS, c\_glu\_usage and as a product in V\_b\_GLU\_c and as a modifier in V\_b\_GLU\_c, V\_b\_GLU\_c, V\_GCS, V\_GCS, c\_glu\_usage, c\_glu\_usage).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}\mathrm{glu} = |v_{17}| - |v_{54}| - |v_{56}| \tag{375}$$

# 10.21 Species c\_cys

Name c\_Cysteine

Initial concentration  $179.792196767939 \ \mu mol \cdot l^{-1}$ 

This species takes part in eight reactions (as a reactant in V\_GCS, cys\_usage and as a product in V\_b\_CYS\_c, V\_CTGL and as a modifier in V\_GCS, V\_GCS, cys\_usage, cys\_usage).

$$\frac{d}{dt}c\_cys = |v_{16}| + |v_{53}| - |v_{54}| - |v_{55}| \tag{376}$$

# 10.22 Species glc

Name Glutamyl-Cysteine

Initial concentration  $9.60706615144005~\mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in V\_GS and as a product in V\_GCS and as a modifier in V\_GCS, V\_GCS, V\_GS, V\_GS).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{glc} = v_{54} - v_{57} \tag{377}$$

# 10.23 Species c\_gly

Name c\_Glycine

Initial concentration  $927.56013101536 \ \mu mol \cdot l^{-1}$ 

This species takes part in 15 reactions (as a reactant in V\_GNMT, V\_GS and as a product in V\_b\_GLY\_c, VmGLYc, VcSHMT and as a modifier in V\_b\_GLY\_c, V\_b\_GLY\_c, VmGLYc, VmGLYc, VcSHMT, VcSHMT, V\_GNMT, V\_GNMT, V\_GS, V\_GS).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}\mathrm{gly} = v_{18} + v_{32} + v_{38} - v_{48} - v_{57} \tag{378}$$

# 10.24 Species c\_gsg

Name c\_GSSG

Initial concentration  $59.8062682413464 \, \mu mol \cdot l^{-1}$ 

This species takes part in 19 reactions (as a reactant in V\_c\_gsgHb, V\_c\_gsgLb, V\_GR, c\_gsg\_degr and as a product in V\_GPX and as a modifier in V\_c\_gsgHb, V\_c\_gsgHb, V\_c\_gsgLb, V\_c\_gsgLb, V\_MATI, V\_MATI, V\_MATIII, V\_MATIII, V\_MATIII, V\_MATIII, V\_GR, V\_GR, c\_gsg\_degr, c\_gsg\_degr).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}gsg = |v_{58}| - |v_{14}| - |v_{15}| - |v_{59}| - |v_{61}|$$
(379)

# 10.25 Species c\_gsh

Name c\_GSH

Initial concentration  $6272.51492720171 \ \mu mol \cdot l^{-1}$ 

This species takes part in 19 reactions (as a reactant in V\_c\_gshHb, V\_c\_gshLb, V\_GPX, c\_gsh\_degr and as a product in V\_GS, V\_GR and as a modifier in V\_c\_gshHb, V\_c\_gshHb, V\_c\_gshLb, V\_c\_gshLb, V\_GCS, V\_GCS, V\_GS, V\_GS, V\_GPX, V\_GPX, c\_gsh\_degr, c\_gsh\_degr).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}gsh = |v_{57}| + 2|v_{59}| - |v_{12}| - |v_{13}| - 2|v_{58}| - |v_{60}|$$
(380)

### 10.26 Species cyt

Name Cystathionine

Initial concentration  $32.3047581589977 \ \mu mol \cdot l^{-1}$ 

This species takes part in four reactions (as a reactant in V\_CTGL and as a product in V\_CBS and as a modifier in V\_CTGL, V\_CTGL).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{cyt} = v_{52} - v_{53} \tag{381}$$

# 10.27 Species hcy

Name Homocysteine

Initial concentration  $1.02425194863179 \ \mu mol \cdot l^{-1}$ 

This species takes part in twelve reactions (as a reactant in V\_MS, V\_BHMT, V\_CBS and as a product in V\_SAHH and as a modifier in V\_MS, V\_BHMT, V\_BHMT, V\_SAHH, V\_SAHH, V\_CBS, V\_CBS).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{hcy} = |v_{50}| - |v_{33}| - |v_{45}| - |v_{52}| \tag{382}$$

# 10.28 Species c\_ser

Name c\_Serine

Initial concentration  $571.225285800751 \, \mu \text{mol} \cdot l^{-1}$ 

This species takes part in 15 reactions (as a reactant in VcSHMT, gluconeogenesis\_ser, V\_CBS and as a product in V\_b\_SER\_c, VmSERc and as a modifier in V\_b\_SER\_c, V\_b\_SER\_c, VmSERc, VmSERc, VcSHMT, VcSHMT, gluconeogenesis\_ser, gluconeogenesis\_ser, V\_CBS, V\_CBS).

$$\frac{\mathrm{d}}{\mathrm{d}t}c_{-}\mathrm{ser} = |v_{19}| + |v_{30}| - |v_{38}| - |v_{51}| - |v_{52}| \tag{383}$$

# 10.29 Species sah

Name SAH

Initial concentration  $15.5626693211634 \ \mu mol \cdot l^{-1}$ 

This species takes part in 18 reactions (as a reactant in V\_SAHH and as a product in V\_GNMT, V\_DNMT and as a modifier in V\_MTHFR, V\_MTHFR, V\_MTHFR, V\_BHMT, V\_BHMT, V\_BHMT, V\_GNMT, V\_DNMT, V\_DNMT, V\_DNMT, V\_SAHH, V\_CBS, V\_CBS, V\_CBS,

$$\frac{d}{dt} \operatorname{sah} = |v_{48}| + |v_{49}| - |v_{50}| \tag{384}$$

### 10.30 Species sam

Name SAM

Initial concentration  $65.0613824844853 \, \mu mol \cdot l^{-1}$ 

This species takes part in 28 reactions (as a reactant in V\_GNMT, V\_DNMT, reaction\_2 and as a product in V\_MATI, V\_MATIII, reaction\_13 and as a modifier in V\_MTHFR, V\_MTHFR, V\_MTHFR, V\_BHMT, V\_BHMT, V\_BHMT, V\_MATI, V\_MATIII, V\_MATIII, V\_MATIII, V\_GNMT, V\_DNMT, V\_CBS, V\_CBS, V\_CBS, reaction\_2, reaction\_13, reaction\_13, reaction\_13).

$$\frac{d}{dt}sam = v_{46} + v_{47} + v_{74} - v_{48} - v_{49} - v_{63}$$
 (385)

### 10.31 Species met

Name c\_Methionine

Initial concentration  $50.6006924474774 \ \mu mol \cdot l^{-1}$ 

This species takes part in 14 reactions (as a reactant in V\_MATI, V\_MATIII, reaction\_13 and as a product in V\_b\_MET\_c, V\_MS, V\_BHMT and as a modifier in V\_b\_MET\_c, V\_b\_MET\_c, V\_MATI, V\_MATIII, V\_MATIII, reaction\_13, reaction\_13).

$$\frac{d}{dt} met = |v_{20}| + |v_{33}| + |v_{45}| - |v_{46}| - |v_{47}| - |v_{74}|$$
(386)

### 10.32 Species c\_coo

Name c\_Formate

Initial concentration  $13.9758110645557 \, \mu \text{mol} \cdot 1^{-1}$ 

This species takes part in six reactions (as a reactant in VcFTS and as a product in VmHCOOHc and as a modifier in VmHCOOHc, VmHCOOHc, VcFTS, VcFTS).

$$\frac{d}{dt}c\_coo = v_{31} - v_{37} \tag{387}$$

# 10.33 Species species\_1

Name dcSAM

Initial concentration  $0.011178638638793 \ \mu mol \cdot l^{-1}$ 

This species takes part in nine reactions (as a reactant in reaction\_7, reaction\_8 and as a product in reaction\_2 and as a modifier in reaction\_2, reaction\_2, reaction\_7, reaction\_7, reaction\_8, reaction\_8).

$$\frac{d}{dt} \text{species}_{1} = |v_{63}| - |v_{68}| - |v_{69}| \tag{388}$$

### **10.34 Species** species\_2

Name Putrescine

Initial concentration  $98.2036296721139 \, \mu mol \cdot l^{-1}$ 

This species takes part in 13 reactions (as a reactant in reaction\_7, reaction\_11 and as a product in reaction\_1, reaction\_5 and as a modifier in reaction\_1, reaction\_1, reaction\_2, reaction\_2, reaction\_7, reaction\_7, reaction\_11, reaction\_11).

$$\frac{d}{dt} \text{species}_2 = |v_{62}| + |v_{66}| - |v_{68}| - |v_{72}| \tag{389}$$

### 10.35 Species species\_3

Name Spermine

Initial concentration  $61.3595114874529 \, \mu \text{mol} \cdot l^{-1}$ 

This species takes part in 18 reactions (as a reactant in reaction\_3 and as a product in reaction\_8 and as a modifier in reaction\_2, reaction\_2, reaction\_2, reaction\_3, reaction\_4, reaction\_4, reaction\_4, reaction\_5, reaction\_5, reaction\_5, reaction\_6, reaction\_6, reaction\_8, reaction\_8).

$$\frac{d}{dt}$$
 species\_3 =  $|v_{69}| - |v_{64}|$  (390)

# 10.36 Species species\_4

Name Spermidine

Initial concentration  $79.5917525310194 \, \mu \text{mol} \cdot l^{-1}$ 

This species takes part in 18 reactions (as a reactant in reaction\_4, reaction\_8 and as a product in reaction\_6, reaction\_7 and as a modifier in reaction\_3, reaction\_3, reaction\_4, reaction\_4, reaction\_5, reaction\_5, reaction\_5, reaction\_6, reaction\_6, reaction\_7, reaction\_7, reaction\_8, reaction\_8).

$$\frac{d}{dt} \text{species} = v_{67} + v_{68} - v_{65} - v_{69}$$
(391)

# 10.37 Species species\_5

Name Acetylspermine

Initial concentration  $0.0272742537826481 \ \mu mol \cdot l^{-1}$ 

This species takes part in seven reactions (as a reactant in reaction\_6 and as a product in reaction\_3 and as a modifier in reaction\_5, reaction\_5, reaction\_5, reaction\_6, reaction\_6.

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{species}_{.5} = |v_{64}| - |v_{67}| \tag{392}$$

# 10.38 Species species\_6

Name Acetylspermidine

Initial concentration  $0.903751442137693 \ \mu mol \cdot l^{-1}$ 

This species takes part in ten reactions (as a reactant in reaction\_5, reaction\_12 and as a product in reaction\_4 and as a modifier in reaction\_5, reaction\_5, reaction\_6, reaction\_6, reaction\_12, reaction\_12).

$$\frac{d}{dt} \text{species}_{.6} = |v_{65}| - |v_{66}| - |v_{73}| \tag{393}$$

### **10.39 Species** species\_7

Name Ornithine

Initial concentration  $300 \, \mu mol \cdot l^{-1}$ 

This species takes part in three reactions (as a reactant in reaction\_1 and as a modifier in reaction\_1, reaction\_1), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{species}_{.7} = 0 \tag{394}$$

### 10.40 Species species\_8

Name Acetyl-CoA

Initial concentration  $38.833390762794 \, \mu \text{mol} \cdot l^{-1}$ 

This species takes part in ten reactions (as a reactant in reaction\_3, reaction\_4, reaction\_9 and as a product in reaction\_10 and as a modifier in reaction\_3, reaction\_3, reaction\_4, reaction\_9, reaction\_9).

$$\frac{d}{dt} \text{species}_{8} = |v_{71}| - |v_{64}| - |v_{65}| - |v_{70}| \tag{395}$$

# 10.41 Species species\_9

Name CoA

Initial concentration  $160.666609237206 \, \mu mol \cdot l^{-1}$ 

This species takes part in ten reactions (as a reactant in reaction\_10 and as a product in reaction\_3, reaction\_4, reaction\_9 and as a modifier in reaction\_3, reaction\_3, reaction\_4, reaction\_10, reaction\_10).

$$\frac{d}{dt} \text{species}_{9} = |v_{64}| + |v_{65}| + |v_{70}| - |v_{71}|$$
(396)

### 10.42 Species CO

Name CO2

Initial concentration  $0 \ \mu mol \cdot l^{-1}$ 

This species takes part in one reaction (as a product in V\_GDC), which does not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{CO} = 0\tag{397}$$

### 10.43 Species m\_thf

Name m\_THF

Initial concentration  $20.989363966111 \, \mu mol \cdot l^{-1}$ 

Initial assignment m\_thf

This species takes part in 19 reactions (as a reactant in VmSHMT, VmFTS, VmNE, V\_GDC, V\_SDH, V\_DMGD and as a product in VmFTD and as a modifier in VmSHMT, VmSHMT, VmFTS, VmFTS, VmNE, VmNE, V\_GDC, V\_GDC, V\_SDH, V\_DMGD, V\_DMGD).

$$\frac{\mathrm{d}}{\mathrm{d}t} \mathbf{m}_{t} \mathbf{h} \mathbf{f} = |v_{21}| - |v_{22}| - |v_{23}| - |v_{24}| - |v_{25}| - |v_{26}| - |v_{27}|$$
(398)

### 10.44 Species m\_2cf

Name m\_5-10-methylene-THF

Initial concentration  $1.66729681864136 \ \mu mol \cdot l^{-1}$ 

This species takes part in twelve reactions (as a reactant in VmMTD and as a product in VmSHMT, VmNE, V\_GDC, V\_SDH, V\_DMGD and as a modifier in VmSHMT, VmSHMT, VmNE, VmNE, VmMTD, VmMTD).

$$\frac{\mathrm{d}}{\mathrm{d}t} \text{m.2cf} = |v_{22}| + |v_{24}| + |v_{25}| + |v_{26}| + |v_{27}| - |v_{28}|$$
(399)

# 10.45 Species m\_1cf

Name m\_5-10-methenyl-THF

Initial concentration  $1.55022131198345 \ \mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in VmMTCH and as a product in VmMTD and as a modifier in VmMTD, VmMTD, VmMTCH, VmMTCH).

$$\frac{d}{dt}m_{-}1cf = v_{28} - v_{29} \tag{400}$$

# **10.46** Species m\_10f

Name m\_10-formyl-THF

Initial concentration  $15.9931179032642 \ \mu mol \cdot l^{-1}$ 

This species takes part in nine reactions (as a reactant in VmFTD and as a product in VmFTS, VmMTCH and as a modifier in VmFTD, VmFTD, VmFTS, VmFTS, VmMTCH, VmMTCH).

$$\frac{\mathrm{d}}{\mathrm{d}t} \mathbf{m}_{-} 10\mathbf{f} = |v_{23}| + |v_{29}| - |v_{21}| \tag{401}$$

# 10.47 Species m\_ser

Name m\_Serine

Initial concentration  $2150.05782513217 \, \mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in VmSHMT, VmSERc and as a modifier in VmSHMT, VmSHMT, VmSERc, VmSERc).

$$\frac{d}{dt} \text{m\_ser} = -v_{22} - 3v_{30} \tag{402}$$

# 10.48 Species m\_gly

Name m\_Glycine

Initial concentration  $2043.470284255 \ \mu mol \cdot l^{-1}$ 

This species takes part in ten reactions (as a reactant in V\_GDC, VmGLYc and as a product in VmSHMT, V\_SDH and as a modifier in VmSHMT, VmSHMT, V\_GDC, VmGLYc, VmGLYc).

$$\frac{d}{dt}m_{-}gly = |v_{22}| + |v_{26}| - |v_{25}| - 3|v_{32}|$$
(403)

# 10.49 Species m\_coo

Name m\_Formate

Initial concentration  $58.3710073680041 \ \mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in VmFTS, VmHCOOHc and as a modifier in VmFTS, VmHCOOHc, VmHCOOHc).

$$\frac{d}{dt}m_{-}coo = -v_{23} - 3v_{31} \tag{404}$$

# 10.50 Species Fol

Name Folate

Initial concentration  $20.1 \ \mu mol \cdot l^{-1}$ 

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{Fol} = 0\tag{405}$$

# 10.51 Species HCHO

Name Formaldehyde

Initial concentration  $500 \ \mu mol \cdot l^{-1}$ 

This species takes part in six reactions (as a reactant in VmNE, VcNE and as a modifier in VmNE, VmNE, VcNE, VcNE), which do not influence its rate of change because this constant species is on the boundary of the reaction system:

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{HCHO} = 0\tag{406}$$

# 10.52 Species src

Name Sarcosine

Initial concentration  $8.25411023033989 \, \mu \text{mol} \cdot 1^{-1}$ 

This species takes part in five reactions (as a reactant in V\_SDH and as a product in V\_DMGD, V\_GNMT and as a modifier in V\_SDH, V\_SDH).

$$\frac{d}{dt}src = |v_{27}| + |v_{48}| - |v_{26}| \tag{407}$$

# 10.53 Species dmg

Name Dimethylglycine

Initial concentration  $0.677559463168903 \ \mu mol \cdot l^{-1}$ 

This species takes part in four reactions (as a reactant in V\_DMGD and as a product in V\_BHMT and as a modifier in V\_DMGD, V\_DMGD).

$$\frac{\mathrm{d}}{\mathrm{d}t}\mathrm{dmg} = |v_{45}| - |v_{27}| \tag{408}$$

 $\mathfrak{BML2}^{lA}$  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.

<sup>&</sup>lt;sup>a</sup>Center for Bioinformatics Tübingen (ZBIT), Germany

<sup>&</sup>lt;sup>b</sup>California Institute of Technology, Beckman Institute BNMC, Pasadena, United States

<sup>&</sup>lt;sup>c</sup>European Bioinformatics Institute, Wellcome Trust Genome Campus, Hinxton, United Kingdom

<sup>&</sup>lt;sup>d</sup>EML Research gGmbH, Heidelberg, Germany