The Systems Biology Workbench (SBW) Version 1.0: Framework and Modules

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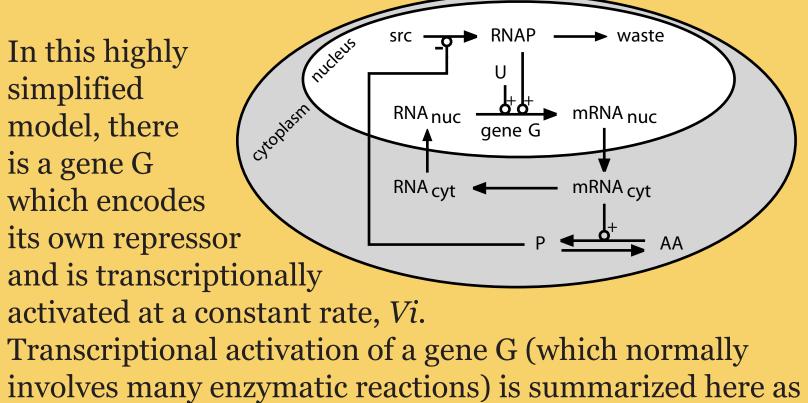
Introduction

Progress in molecular biotechnology has fueled an explosion in the development of software tools. Regrettably, developers often end up recreating similar facilities in separate software packages.

In an effort to make it more attractive for developers to share rather than reimplement resources, we have implemented the Systems Biology Workbench (SBW), a free, open-source, application integration environment. Our goal has been to create a framework simple enough that software authors find it easier to provide an SBW interface than to recreate functionality available in other tools. By doing so, we hope **developers** can concentrate on creating best-of-breed solutions in their areas of expertise.

SBW in Action: A Sample Session

Here is an example of using several SBW-enabled tools to create and simulate a two-compartment model of a hypothetical single-gene oscillatory circuit.



the production of active *RNAP* from source material, *src*, and degradation to ____

Reaction

 $src \rightarrow RNAP$

 $RNAP \rightarrow waste$

 $RNA_{nuc} \rightarrow mRNA_{nuc}$

 $mRNA_{nuc} \rightarrow mRNA_{cyt}$

 $mRNA_{cyt} \rightarrow RNA_{cyt}$

 $RNA_{cyt} \rightarrow RNA_{nuc}$

 $AA \rightarrow P$

 $P \longrightarrow AA$

Rate

 V_{i}

 $1 + P/K_i$

 $V_{kd} \cdot RNAP$

 $k_1 \cdot mRNA_{nuc}$

 $k_2 \cdot RNA_{cyt}$

 $V_{m4} \cdot P$

 $\overline{P+K_{m4}}$

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 $V_{m2} \cdot mRNA_{cyt}$

 $mRNA_{cyt} + K_{m2}$

 $V_{m3} \cdot \textit{mRNA}_{cyt} \cdot \textit{AA}$

 $AA + K_{m3}$

\$2,14

 $V_{m1} \cdot RNAP \cdot RNA_{nuc}$

 $K_{m1} + RNA_{nuc}$

Third-Party Modules Available For SBW

- Jarnac, a biochemical simulation package for Windows
- JDesigner, a visual biochemical network layout tool
- **Pasadena Twain**, a simple interactive ODE solver
- A stochastic simulator based on Gillespie's algorithm
- A **bifurcation analysis** module
- An **optimization** module
- An **SBML validator** for checking SBML model files
- An **inspector** that lists running modules & their services

Coming Attractions

More open-source developers are joining the SBW project, and together we are enhancing and extending SBW in many ways. Here is a preview of some coming attractions:

What Does SBW Provide?

SBW provides libraries for enabling applications to learn about and communicate with each other. The applications may be running on separate computers.

SBW lets heterogevneous packages connect to each other using a **remote procedure call** mechanism based on a message-passing network protocol. The interfaces to SBW are encapsulated in client libraries for different languages.

The SBW Broker

and coordinates

communications

A Broker is started

automatically for

the user if one is

the first SBW

not running when

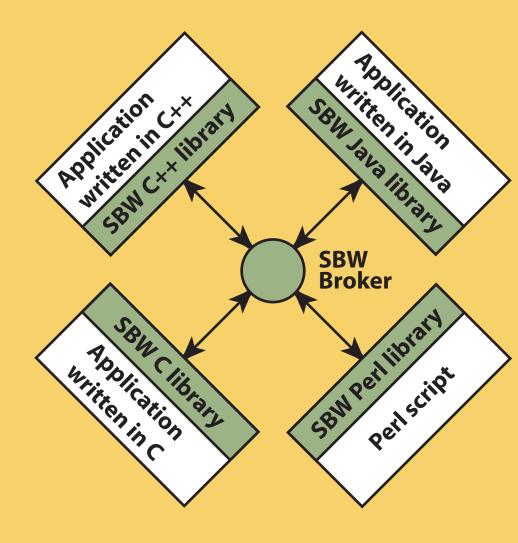
application starts.

on demand,

on a given

computer.

starts applications



Features of SBW Version 1.0

- Languages supported: C, C++, Delphi, Java, Perl, and **Python**.
- Windows (98, 2000, XP) and Linux supported, with **MacOS X** planned in the near future.
- Secure, distributed operation via SSH, featuring remote startup of brokers and applications.

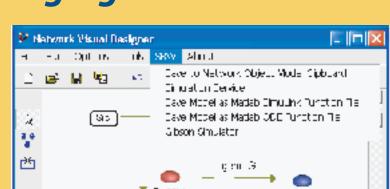
waste. Transcribed *mRNA* is then transported from the nucleus into the cytoplasm, where it is translated into the product *P* from constituent amino acids AA and where it is also subject to degradation.

Model Capture: **Using a Visual Editor**

Using the SBW-enabled JDesigner biochemical network editor, a user can create the model using a graphical interface.



An application such as JDesigner can dynamically create a menu of tools with which it can interact, by querying SBW to find all installed SBW-enabled packages that provide services for processing biochemical models.



- Support for JDK 1.4
- Support for MacOS X
- MATLAB scripting interface
- New modules, including:
 - Graphical browser for the SBW environment
 - Rewritten SBML parser/writer for SBML Level 2
 - Improved generic GUI for simulators
 - Improved, full-featured plot module
 - New simulation engines

How to Get Started with SBW

The SBW version 1.0 package and extensive documentation are available from the project web site,

http://www.sbw-sbml.org/.

SBW is distributed under the terms of the GNU LGPL.

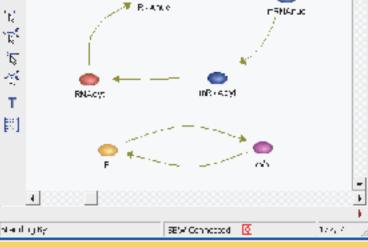
Acknowledgments

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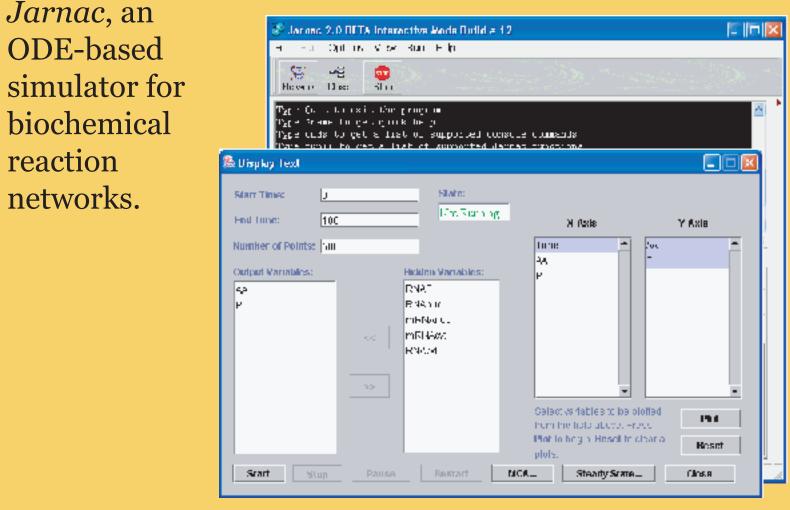
We thank the following groups engaged in developing software tools for systems biology, for their feedback and guidance in developing SBW:

- BioSpice (Arkin et al.)
- Cellerator (Mjolsness et al.)
- DBsolve (Goryanin et al.)

- CORBA gateway for bidirectional communication between SBW-based apps and CORBA-based apps.
- Collection of basic applications provided with the SBW distribution, including:
 - A simple **stochastic simulator** based on the Gibson-Bruck variant of the Gillespie algorithm
 - An SBML-to-MATLAB ODE & Simulink translator
 - An **SBML reader tool** that allows a program to extract (via an API) components of an SBML model
 - A "clipboard" module that stores an SBML model description, and allows the easy transfer of models between separate modules
 - A "browser" module that allows querying SBW for registered modules and producing descriptions of each module's interface in Java or CORBA IDL
 - A simple **plotting module** for time-series data
 - A generic simulation control GUI interface.
 - A collection of tutorial **example modules** in C, C++, Delphi and Java
- Extensive documentation—in addition to overview documents and published papers, every language library has its own programmer's manual and API

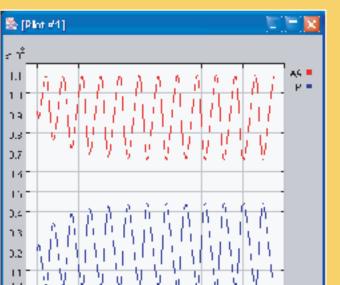


Here, picking the menu item "Simulation Service" invokes a generic simulation control GUI, which in turn invokes



Model Visualization and Analysis

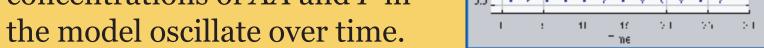
Setting the run parameters for the simulation and selecting the output variables in the simulation control GUI allows the user to plot the values of quantities over time. The plot at the right shows how the concentrations of AA and P in



- E-CELL (Tomita et al.)
- Gepasi (Mendes et al.)
- ProMoT/DIVA (Gilles et al.)
- StochSim (Bray et al.)
- Virtual Cell (Loew et al.)

We also thank the following individuals for their comments, suggestions and assistance: Tau-Mu Yi, Mineo Morohashi, Pieter van der Zee, Fred Livingston, Paul Shannon, Deanne Taylor, Andrew de Laix, Venkat Jagadish, Lukasz Salwinski, Mark Johnson, Richard Giuli, Greg Riddick, Michael Vanier.





A user can also perform other analyses on the model via







open source