

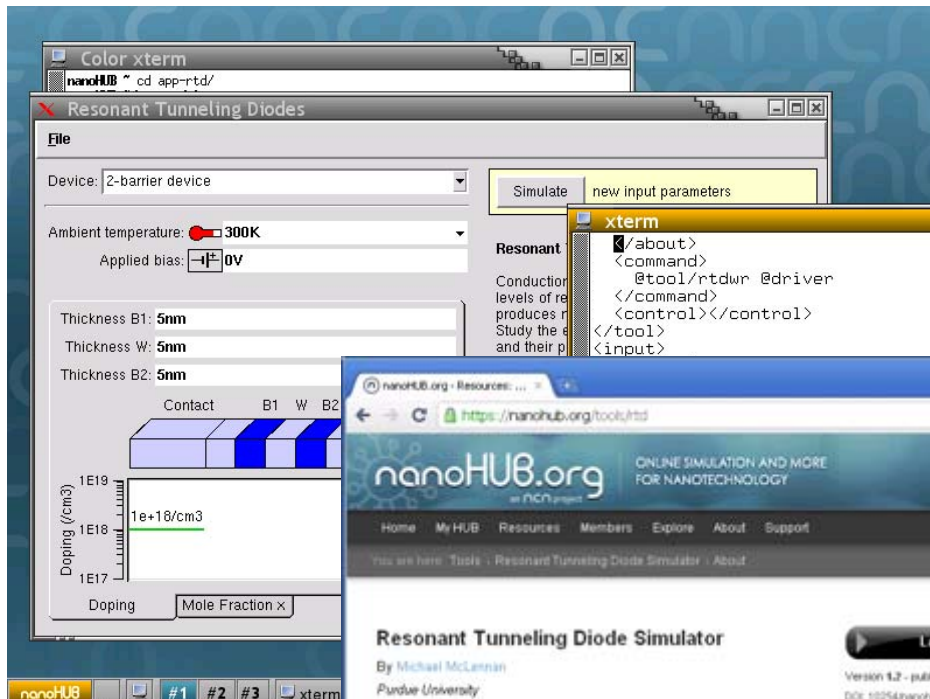
Regression Testing

Michael McLennan

HUBzero® Platform for Scientific Collaboration

Purdue University

Build the tool...



Install, approve...

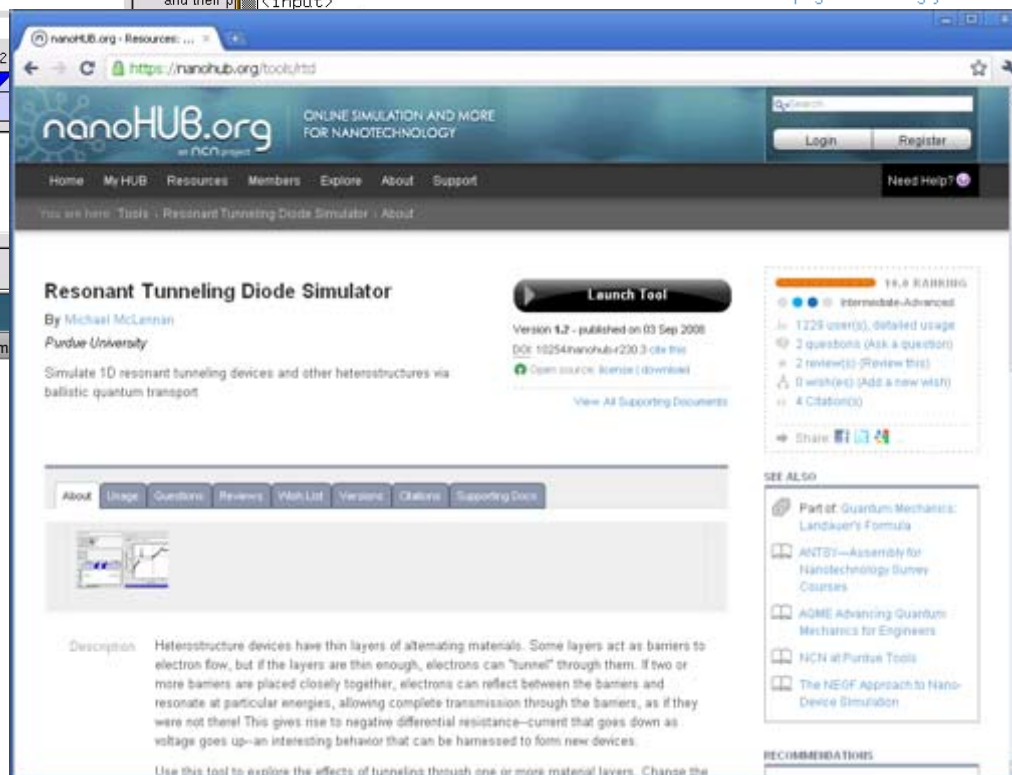
What's next?

Your latest code is installed and ready on nanoHUB.org. Please test your tool by clicking the button below to make sure that everything is working properly, as well as verify that the page describing your tool is created and displays correct information:

→ Test your application:



→ Review the page describing your tool



is working properly, click here to

d in your latest fixes, click here to

the latest version for testing and

Published!

Usually doesn't end there...

[Allow download of 3d simulation data](#)

Proposed by Golam Rabbani on 31 May 2009, 2 Comments

+7 0

+5

257439 [When I launch this tool, it crashes right away and collapses down to a...](#)

by Michael McLennan ([mmc](#))

open

Wishes

Bugs

New physics

$$\frac{m^* v_{emit}^2}{2} \approx \hbar \omega_{phonon}(opt.) \quad \mu = m_{sat}^2 \frac{2L}{W} \frac{1}{C_i}$$

```

Color xterm
real hbar, hbar2, Kb, m0, eVtoJ, pi, q
real ks, mtocm
parameter ( hbar=6.58216e-16,
+ hbar2=hbar**2,
+ Kb=8.61738e-5,
+ m0=9.1095e-31,
+ eVtoJ=1.60218e-19,
+ pi=3.141592654,
+ q=eVtoJ,
+ ks=12.846,
+ mtocm=100.0 )
c
c DEFINITIONS --
c ltor ... electrons traveling left to right
c rtor ... electrons traveling right to left
    
```

What's next?

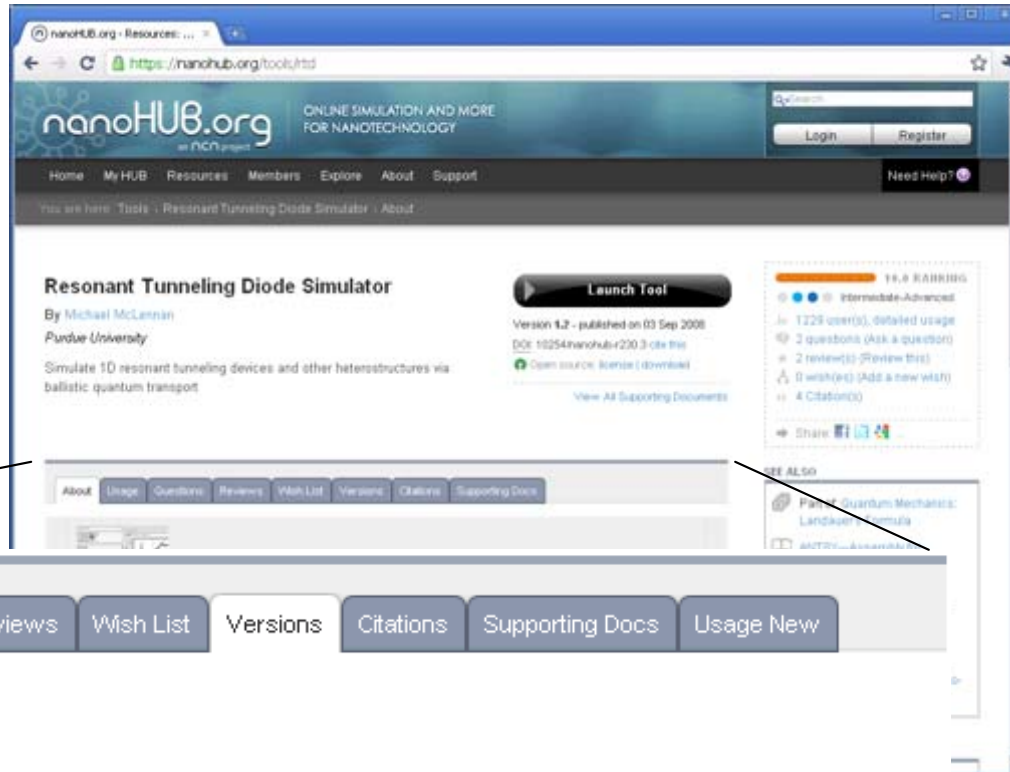
Your tool has been published and is now open to the public. Users will find out about your tool by reading the following page. Be sure to keep this information up to date:

<http://nanoHUB.org/tools/rtd>

Your options:

- > I've made changes [Please install the latest code for testing and approval.](#)





Versions

Version	Released	DOI Handle	Published
3.01	03 Dec 2010	10254/nanohub-r1307.7	✓
3.0	03 Nov 2010	10254/nanohub-r1307.6	✗
2.1	17 Jul 2009	10254/nanohub-r1307.5	✓
2.02	25 Sep 2008	10254/nanohub-r1307.4	✗
2.0.1	08 May 2008	10254/nanohub-r1307.3	✗

You can publish multiple versions and keep more than one active

- Build up a suite of test cases
- Run them as you go along to make sure they still work
- Investigate failures and track down the problem

The screenshot shows the Rappture Regression Tester window. The left pane displays a tree view of test cases under the heading 'Result' and 'Test Case'. The tree structure is as follows:

- helium
 - 2eV (failed)
 - 1eV (failed)
 - 0eV (failed)
 - fail (failed)
- multiple
 - levels
 - spaces are allowed... (passed)
- roomtemp (selected)
 - 1V (failed)
 - 0eV (failed)
- nitrogen
 - 1eV (passed)
 - 0eV (passed)

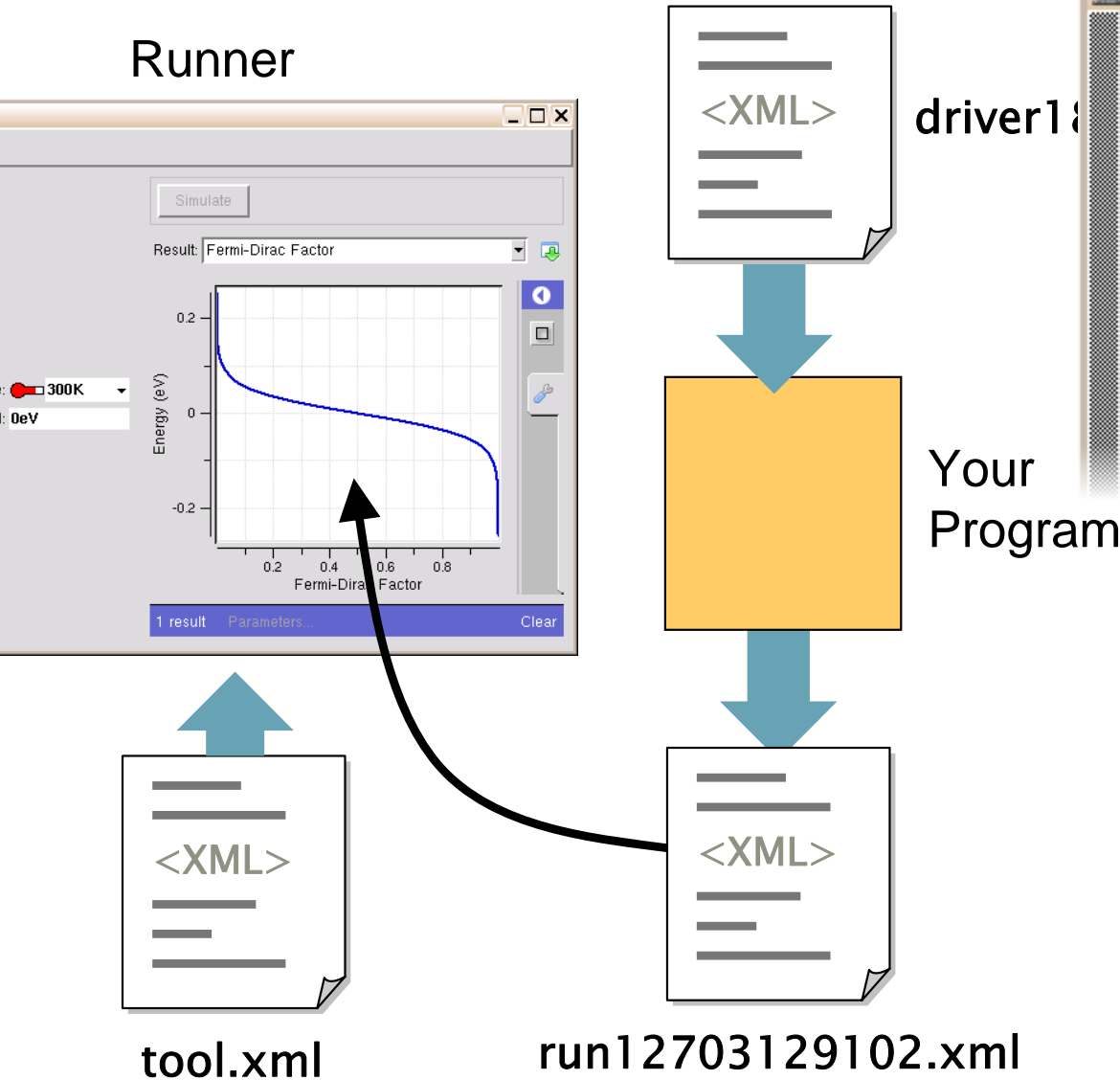
The right pane shows details for the selected test case, 'Test: 1V', which has failed. A red 'X' icon and the text 'Test failed' are displayed. Below this, a message states: 'This test should generate a warning due to differences in the input parameters.' A 'Run' button is visible in the top right corner of the right pane.

Differences:

- Input: Control value**
input.boolean(extra)
Test case has this extra input value
- Output: Extra number in output**
output.number(extra)
Result is missing from current output
- Output: Femi-Dirac Factor**
output.curve(f12)
Result differs from expected value

At the bottom of the window, there are 'Select: All None' buttons and a '<< New golden standard' button.

1 Run the desired test case...



2 Move into tests directory

```

Color xterm
$ ls
docs  spiro.m  tool.xml
$ mkdir tests
$ unset SESSIONDIR
$ rapture
$ ls
docs  run12703129102.xml  spiro.m  tes
$ mv run12703129102.xml tests
$ █
    
```

Results are moved out of the current working directory to the "results" directory **unless you unset SESSIONDIR**



~/data/results/\$SESSION

3 Add label/description to each test

```
Color xterm  
$ vi tests/run12703129102.xml
```

```
<?xml version="1.0"?>  
<run>  
  <test>  
    <label>roomtemp|0eV</label>  
    <description>Should work at room temperature and 0eV.</description>  
  </test>  
  <tool>  
    <about>Press Simulate to view results.</about>  
    <command>tclsh @tool/fermi.tcl @driver</command>  
  </tool>  
  <input>  
    <number id="temperature">  
      <about>  
        <label>Ambient temperature</label>  
        <description>Temperature of the environment.</description>  
      </about>
```

separator for test folders

X Rapture Regression Tester

Result	Test Case
	helium
✗	2eV
✗	1eV
✗	0eV
✗	fail
	multiple
	levels
✓	spaces are allowed...
	roomtemp
✗	1V
✗	0eV
	nitrogen
✓	1eV
✓	0eV

Test: 1V

Test failed

This test should generate a warning due to differences in the input parameters.

Run

Differences:

- ⚠ **Input: Control value**
 input.boolean(extra)
 Test case has this extra input value
- ✗ **Output: Extra number in output**
 output.number(extra)
 Result is missing from current output
- ✗ **Output: Fermi-Dirac Factor**
 output.curve(f12)
 Result differs from expected value

Select: All None

<< New golden standard

Output results:

- ✘ Output value has changed
- ✘ Output value is missing
- ✘ Output value is extra (not supposed to be there)

Input values:

- ✘ Input value has changed--label, units, etc.
- ✘ Input value is missing from test case
- ✘ Input value is extra (not supposed to be there)

Fix the tool

```
Color xterm
real hbar, hbar2, Kb, n0, eVtoJ, pi, q
real ks, ntoen
parameter ( hbar=6.582116e-16,
+          hbar2=hbar**2,
+          Kb=8.61730e-5,
+          n0=9.1095e-31,
+          eVtoJ=1.60218e-19,
+          pi=3.141592654,
+          q=eVtoJ,
+          ks=12.946,
+          ntoen=100,0 )
c
c DEFINITIONS --
c ltor ... electrons traveling left to right
c rtor ... electrons traveling right to left
```

Fix the test

<< New golden standard

Create a test suite with these cases

Fancy cross

$$n_1 = 13$$

$$n_2 = -7$$

$$n_3 = -3$$

Flower

$$n_1 = 19$$

$$n_2 = -13$$

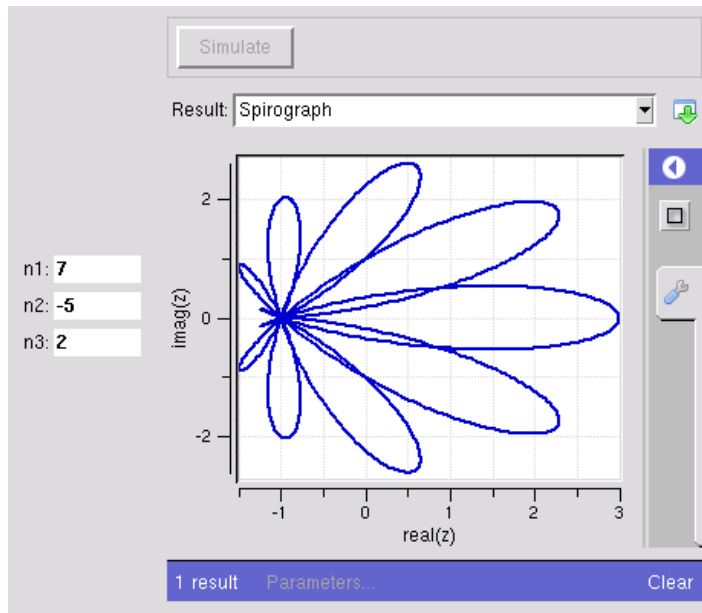
$$n_3 = 3$$

Palm Branch

$$n_1 = 7$$

$$n_2 = -5$$

$$n_3 = 2$$



- Run the regression tests (should run cleanly)
- Edit one of the tests and delete some numbers
- Run the tests, discover the error, then regoldenize