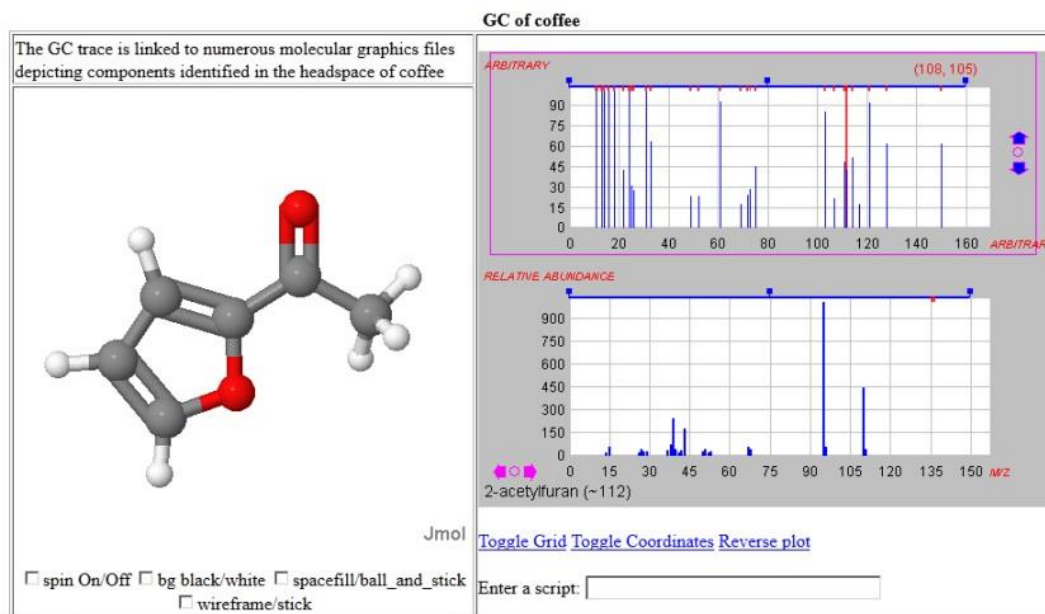


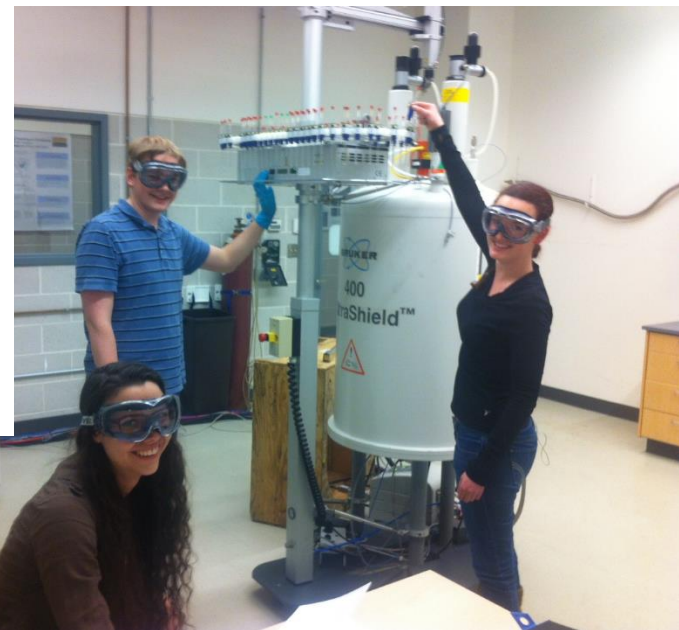
SwingJS -- Resurrecting the interactive functionality of Java on the web as JavaScript

Bob Hanson
Department of Chemistry
St. Olaf College



The top display is the GC trace and the lower display is for the MS spectra.
Clicking on a peak in the GC will load the appropriate Mol file and MS.

Reference: [Google Docs](#)



MSCS Colloquium
St. Olaf College
April 15, 2019

Acknowledgments

Summer, 2016

Nadia El Mouldi

Andreas Raduege



Acknowledgments

Summer, 2017

Andrew Lee

Tahir Ahsan

Nikesh Yadav



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Zhou Renjian

Shanghai, China



early SwingJS
collaborators

Paul Falstad

Minneapolis, MN



Udo Borkowski

Aachen, Germany



Outline of presentation

Historical context

The problem and its solution

Where we are now

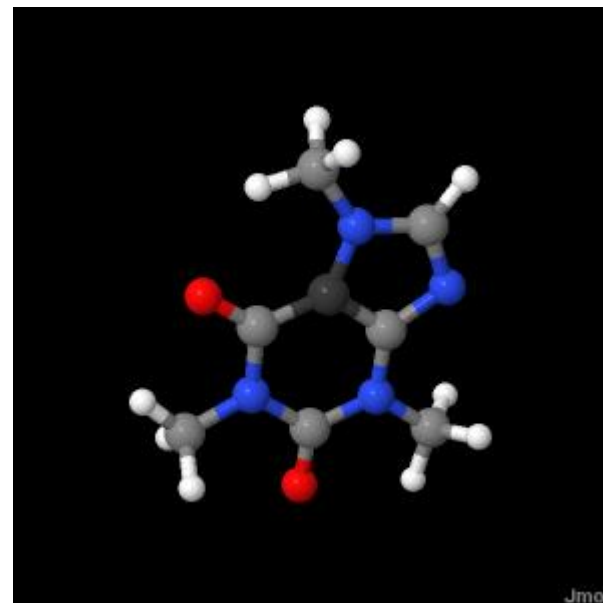
Where we are headed

Jmol

Mission: The high-quality, real-time visualization of molecular structure, dynamics, and energetics.

open source Java
cross-disciplinary
development driven by user input

Principal developer since 2006



Jmol

amylopectin

amylose

cellulose

periodic table

glycogen

glucose

protein

sucrose

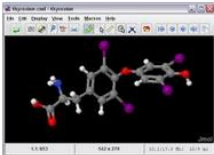
color

peptidyl transferase center


software

structure


icon



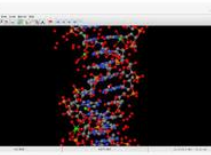
Jmol Wiki - Jmol Application
wiki.jmol.org




Jmol - Wikipedia
en.wikipedia.org



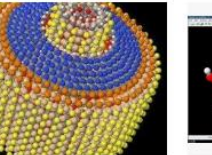
JSmol and Jmol
rcoo.org



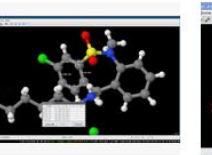
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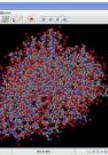
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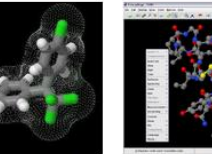
crystal structure - How to get Jmol to ...
chemistry.stackexchange.com



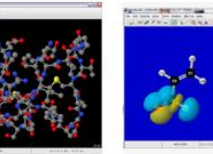
Jmol Tutorial
chemistry.uneche.maine.edu




Jmol Tutorial
chemistry.uneche.maine.edu



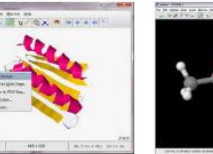
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sourceforge.net



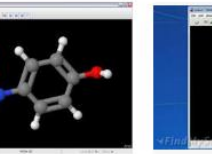
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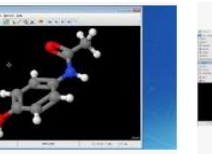
NBOP ro@Jmol
roo7.chem.wisc.edu



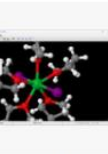
Jmol Training Guide - MSOE Center...
cbm.msoc.edu



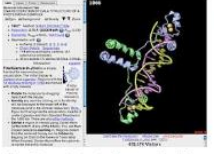
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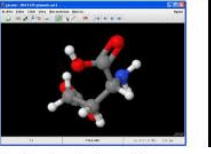
First Glance in Jmol - Proteopedia, lif...
proteopedia.org



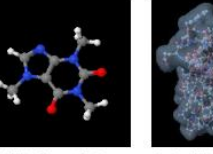
Jmol - Workbench
wiki.c2i2.columbia.edu




Jmol/JSmol Interactive Script Doc...
chemapps.stola.edu



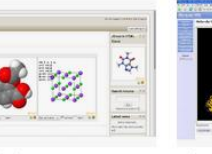
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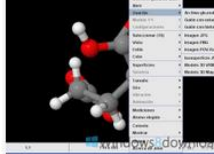
Jmol - Wikipedia
en.wikipedia.org



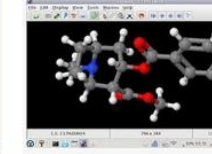
Moodle plugins directory: Jmol
moodle.org



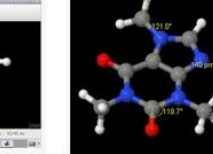
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Jmol - Windows & Downloads
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repo.openpandora.org Repository ...
repo.openpandora.org



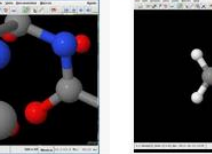
Screenshots
jmol.sourceforge.net



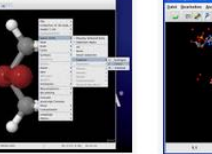
Getting 3D model
biostar.org



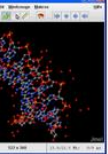
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wiki.jmol.org



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Jmol - Download
jmol.de.sofpatic.com

Demo Jmol and Jspecview

<https://chemapps.stolaf.edu/jmol/jsmol/simple2.htm>

https://chemapps.stolaf.edu/jmol/jsmol/nmr_predict_1H.htm

Jmol

Symmetry @ Otterbein

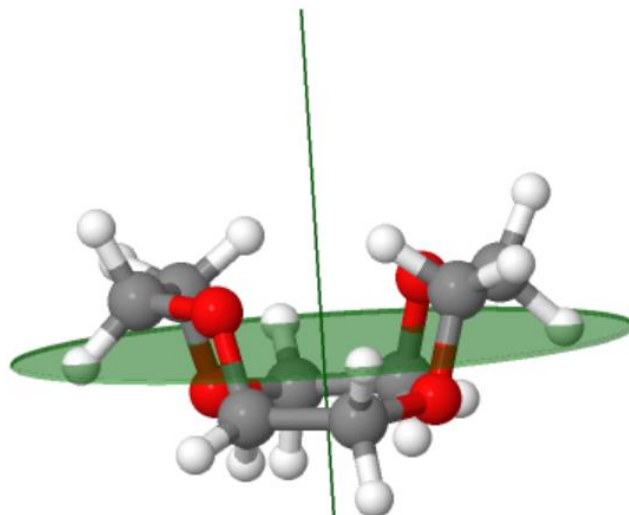
[Home](#)[Tutorial](#)[Gallery](#)[Challenge](#)[Info](#)[Feedback](#)

Point Group Type:

All

Select Molecule:

- $\Lambda(\lambda,\lambda)$ -*cis*-[Co(en)₂Cl₂]⁺
- (δ,δ)-*trans*-[Co(en)₂Cl₂]⁺
- (δ,λ)-*trans*-[Co(en)₂Cl₂]⁺
- 1,1-dichloroethylene
- 1,2-dibromobenzene
- 1,2-dichloroethylene (cis)
- 1,2-dichloroethylene (trans)
- 1,4-dibromobenzene
- 12-crown-4
- 18-crown-6
- trans*-cyclooctene
- closo*-[Ge₉]²⁻
- nido*-[Ge₉]⁴⁻

[Preferences](#)[HTML5](#) • [Java](#)Point Group = S₄S₄

JSmol

Element Operation

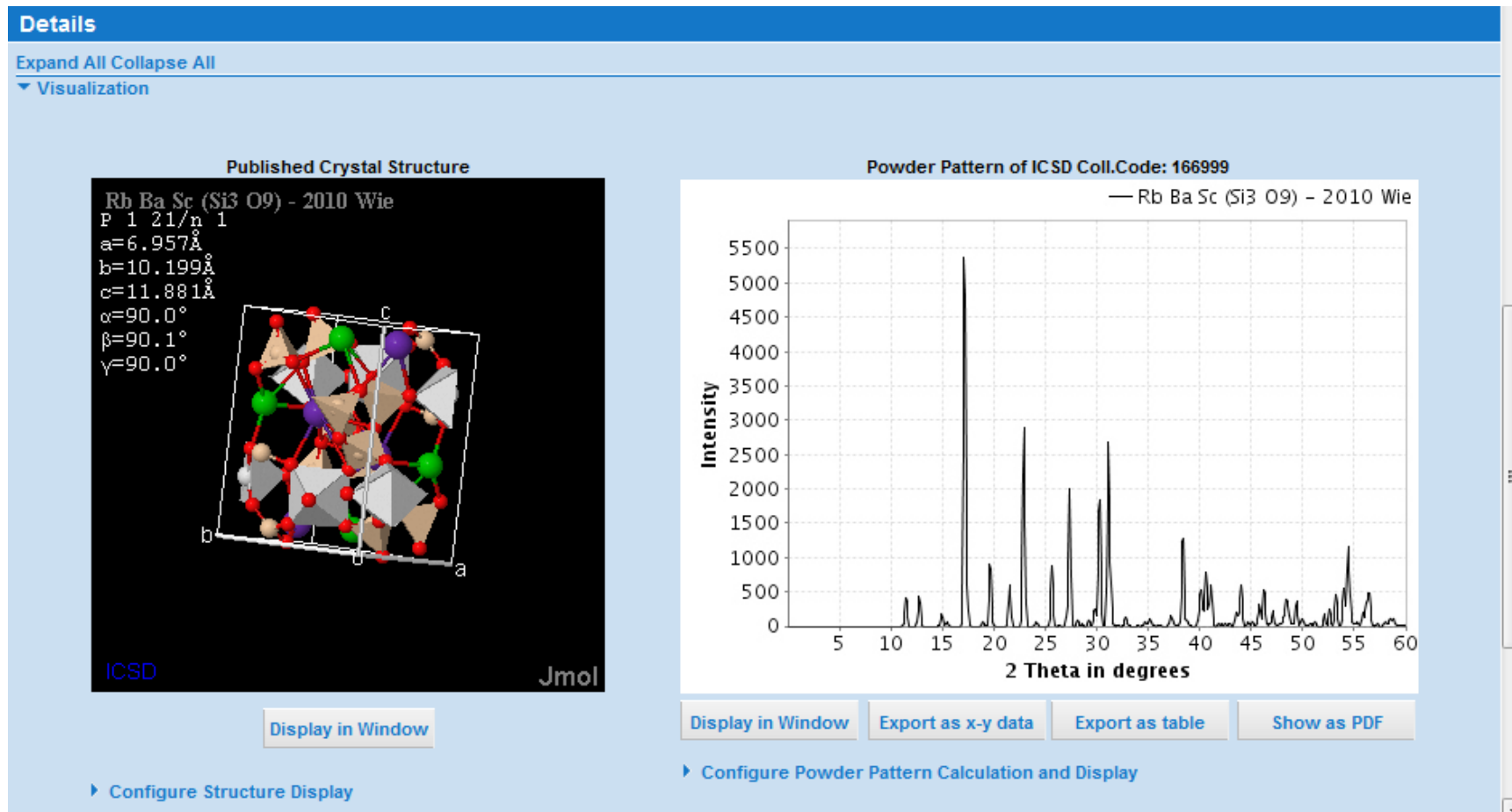
☐ C₂ axis [Rotate](#)☒ S₄ axis [Rotate](#)

Reference: Al-Jallal, N. A.; Al-Kahtani, A. A.; El-Azhary, A. A. J. *Phys. Chem. A* **2005**, *109*, 3694. DOI: [10.1021/jp050133c](https://doi.org/10.1021/jp050133c)

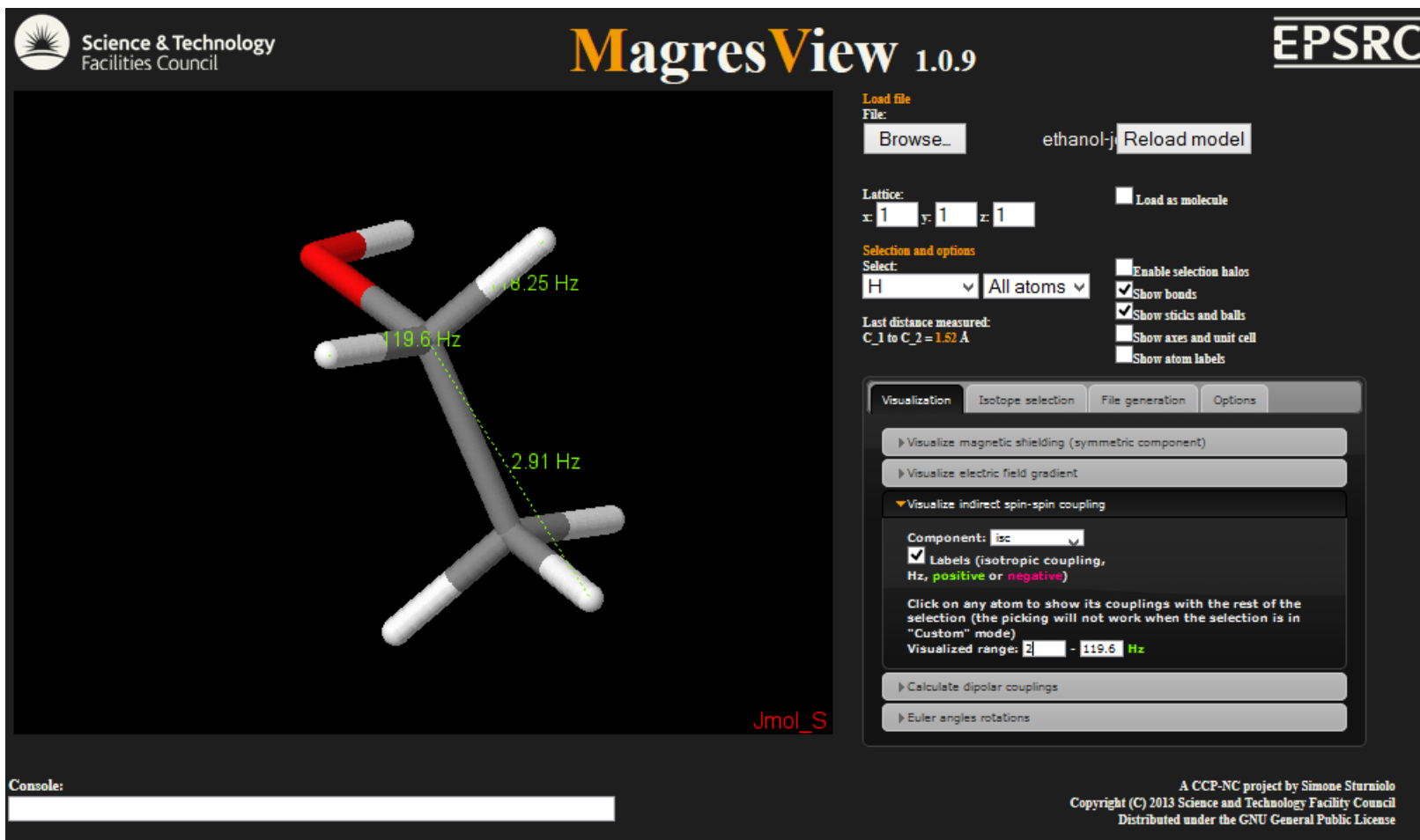
Demo Otterbein Symmetry Gallery

<http://symmetry.otterbein.edu/gallery/>

Jmol

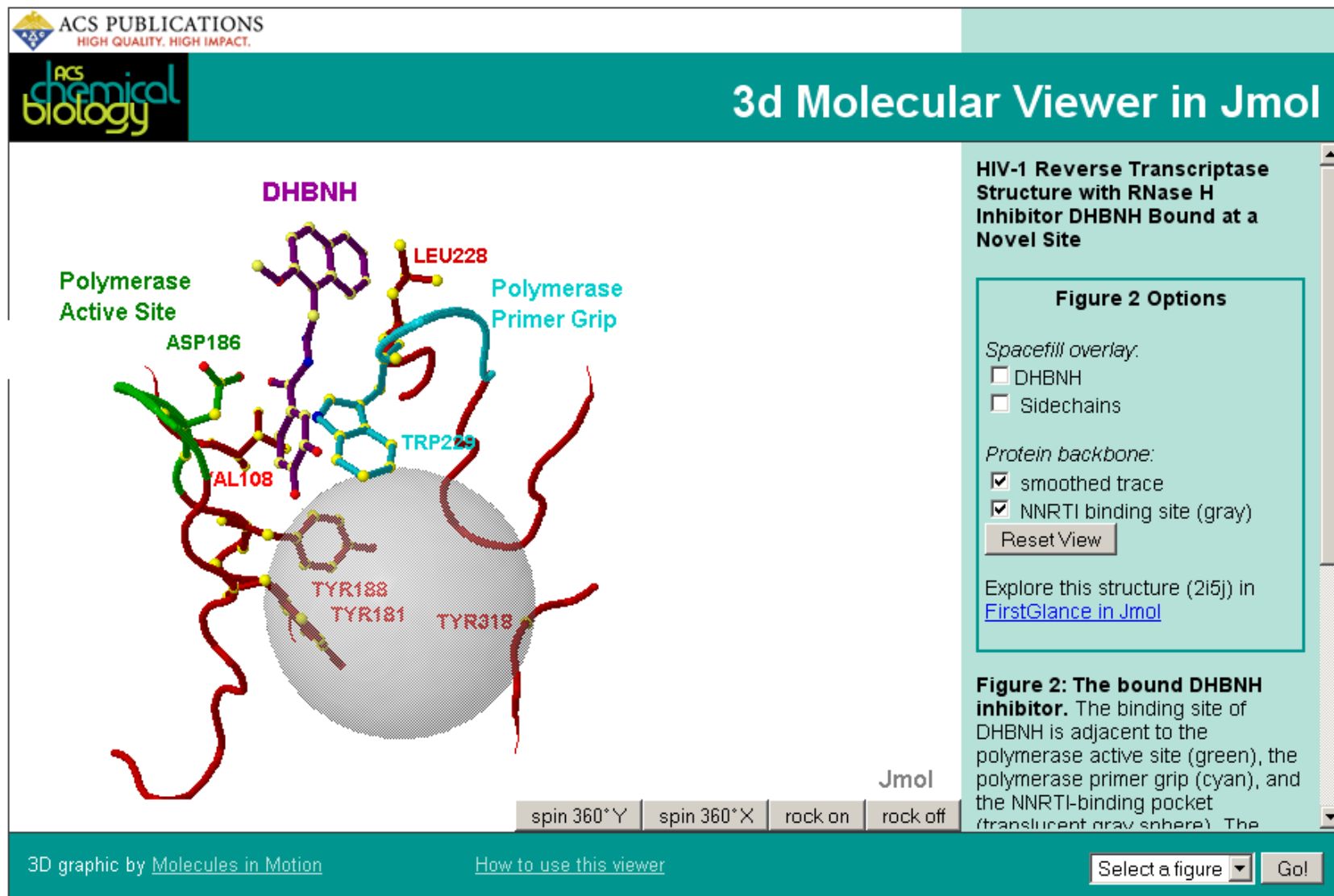


Jmol



Simone Sturniolo and Jonathan Yates, Department of Materials, University of Oxford
https://www.ccpnc.ac.uk/magresview/magresview/magres_view.html

Jmol



Historical context

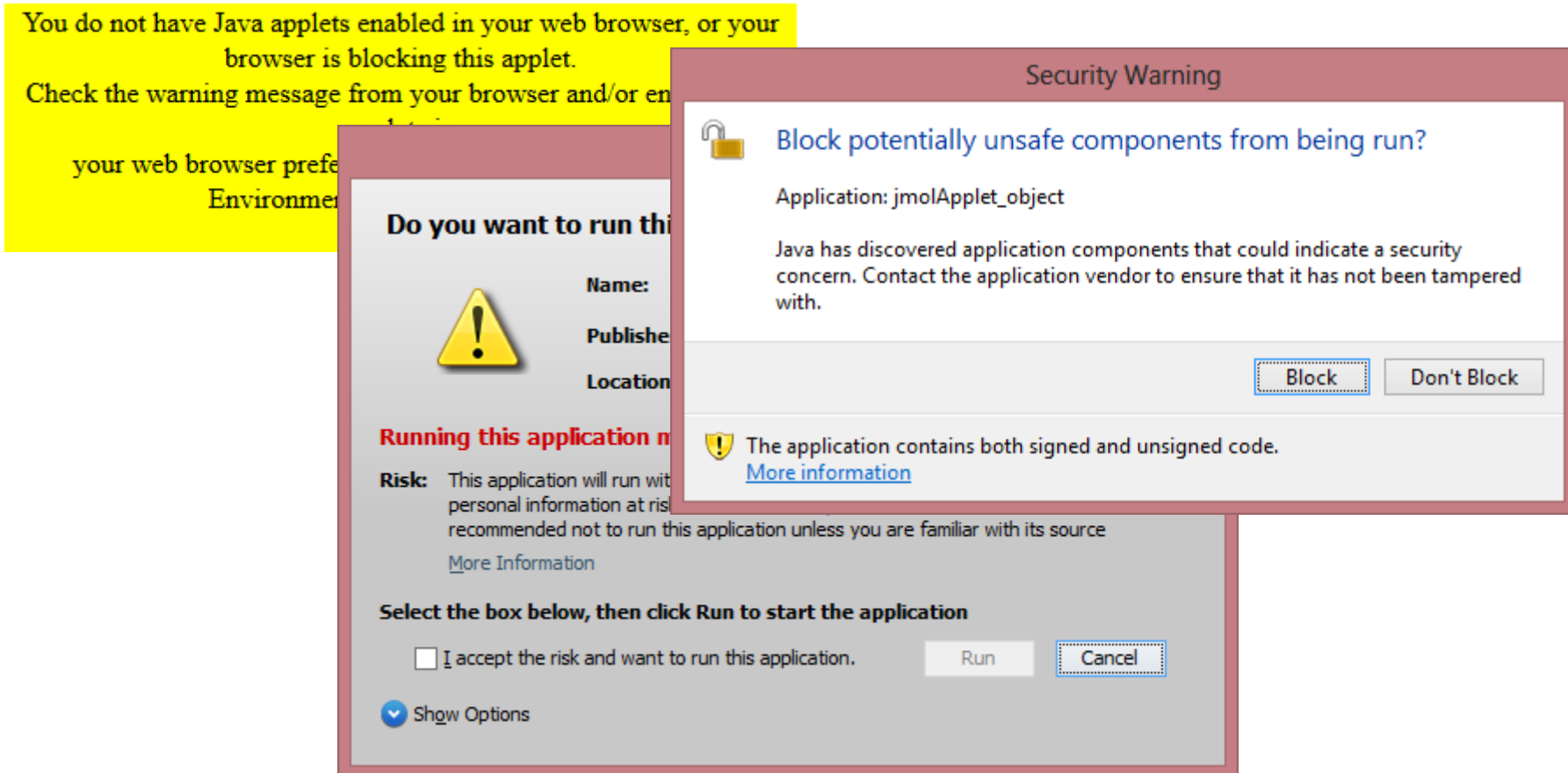
The problem and its solution

Where we are now

Where we are headed

Death of the Java applet

During the summer of 2012 it became apparent that Java applets were no longer suitable for general web use.



The Problem – Java Availability

Most notably, a series of security concerns surrounding Java applets during the fall and early winter of 2012 resulted in an announcement on Jan. 10, 2013, by the U.S. Department of Homeland Security with this suggestion:

- ***Disable Java in web browsers***
- *This and previous Java vulnerabilities have been widely targeted by attackers, and new Java vulnerabilities are likely to be discovered. To defend against this and future Java vulnerabilities, **consider disabling Java in web browsers until adequate updates are available**. As with any software, unnecessary features should be disabled or removed as appropriate for your environment.*

<https://www.us-cert.gov/ncas/alerts/TA13-010A>

java2script to the rescue!

Open-Source Plug-In for Eclipse

Works as a *transpiler* (effectively)

Allows *simultaneous co-creation* of
Java “byte code” class files and
equivalent JavaScript files



Zhou Renjian

java2script to the rescue!

Open-Source Plug-In for Eclipse

Includes a purely JavaScript run-time library that emulates the Java Virtual Machine

Prototype stage of development by Zhou Renjian – just a few demo implementations – never distributed



Zhou Renjian

java2script to the rescue!

Sept – Dec. 2012 “JSmol” work begins at St. Olaf

Intense Jmol implementation effort (150,000 lines of code)

Sept/Oct Collaborative development (fixing transpiler bugs!)

Nov/Dec Extensive Java/JavaScript performance optimization

Jan 2013 – released!

JSmol Success! 16,000+ web pages

For the month of
February, 2014

Visits

112,721



Unique Visitors

79,827



Pageviews

255,522

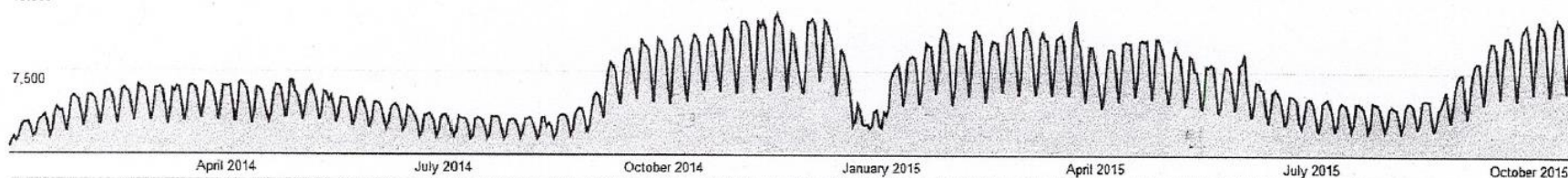


Jan 2014 - Oct, 2015

Sessions

15,000

7,500



Sessions

3,938,294



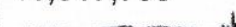
Users

2,407,160



Pageviews

10,017,793



Pages / Session

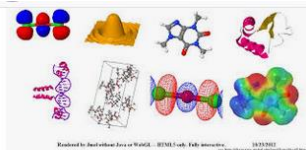
2.54



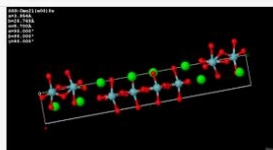
New Visitor Returning Visitor



JSmol Success! 16,000+ web pages



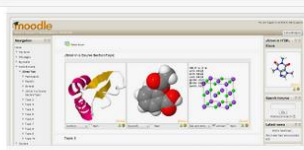
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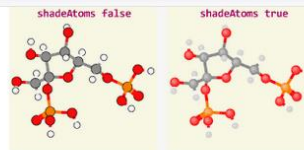
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JSmol and Jmol
rcsb.org



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moodle.org



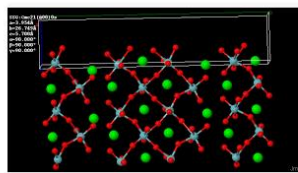
Jmol Wiki - Lightweight JSmol
wiki.jmol.org



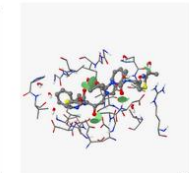
Jmol Wiki - Jmol JavaScript Object
wiki.jmol.org



Plugins/Blocktype/Jmol - Mahara Wiki
wiki.mahara.org



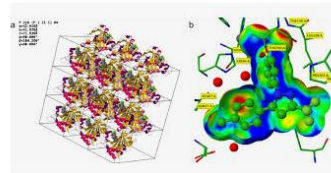
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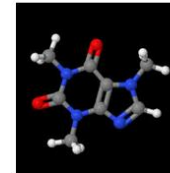
jmolApplet0 - Jmol 14.29.18
chemapps.stolaf.edu

System	Browser	JSmol (No Java)	Jmol Java applet
Internet Explorer	Compatibility mode	Good	
Firefox	OK	Good	
Chrome	Best!	Good	
Opera	Rotation not good	Good	
Safari	OK	Good	
Firefox	OK	Good	
Mac OS X	Best!	Incompatible with Java	
Opera	Rotation not good	Good	

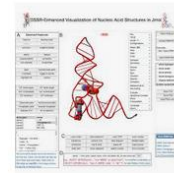
Jmol Wiki - JSmol pros and cons
wiki.jmol.org



Two new custom rendering options on the ...
researchgate.net



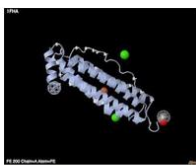
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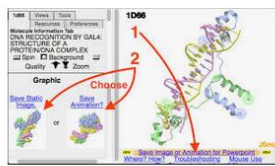
DSSR-JSmol of web interfac...
researchgate.net



Jmol - Wikipedia
en.wikipedia.org



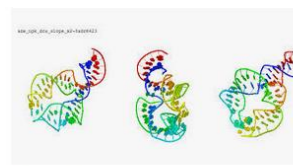
Using FirstGlance in JSmol - C...
sites.google.com



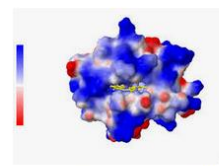
Versions of FirstGlance in Jmol
bioinformatics.org



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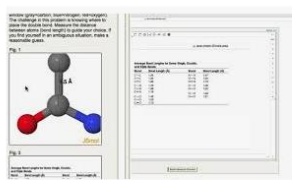
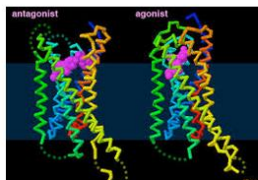
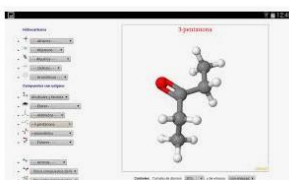
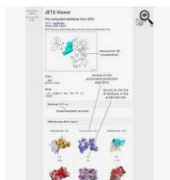
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Jmol hashtag on Twitter
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JSmol

AFLOW
Automatic - FLOW for Materials Discovery

The Brillouin Zone

TRI MCL MCLC ORC **ORCC** ORCI ORCF TET BCT RHL HEX CUB BCC FCC

Showing 2996 of 2996 structures.
The colored buttons below group together sets of elements to search for.

CLEAR **NOT** [Purple Box] + [Dark Blue Box] + [Light Blue Box] **ONLY** ?

Search Options

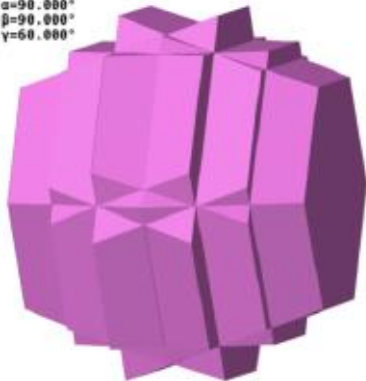
1A	2A	3B	4B	5B	6B	7B	8	1B	2B	3A	4A	5A	6A	7A	NOBLE		
1 H															He		
2 Li	Be									B	C	N	O	F	Ne		
3 Na	Mg									Al	Si	P	S	Cl	Ar		
4 K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr
5 Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe
6 Cs	Ba	La - Lu	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn
7 Fr	Ra	Ac - Lr	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Cn	Nh	Fl	Mc	Lv	Ts	Og

Collaboration with Duke University and US Naval Research Laboratory

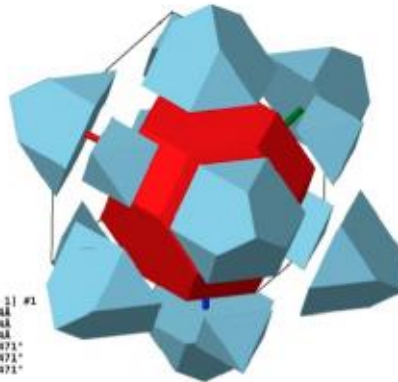
JSmol



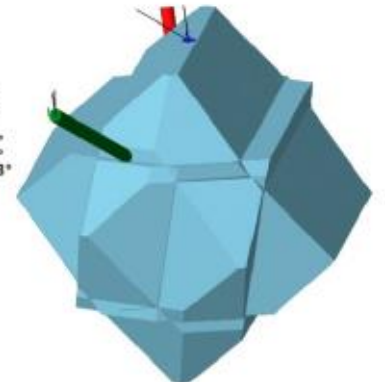
a=2.483Å
b=2.483Å
c=0.064Å
α=90.000°
β=90.000°
γ=60.000°



P 1 [P 1] #1
a=2.614Å
b=2.614Å
c=2.614Å
α=189.471°
β=189.471°
γ=189.471°



a=2.627Å
b=2.627Å
c=1.246Å
α=58.889°
β=58.889°
γ=110.298°



Collaboration with Duke University and US Naval Research Laboratory

Historical context

The problem and its solution

Where we are now

Where we are headed

JSmol

Release 2.2.0 by P. Canepa and R. Hanson - Updated Tue 11 Sep 2018 07:33:23 +08- powered by [JSmol](#)

[Help](#) | [Acknowledgments](#)

FILE

CELL

SHOW

EDIT

SYM BUILD

MEASURE

ORIENT

POLY

SURFACE

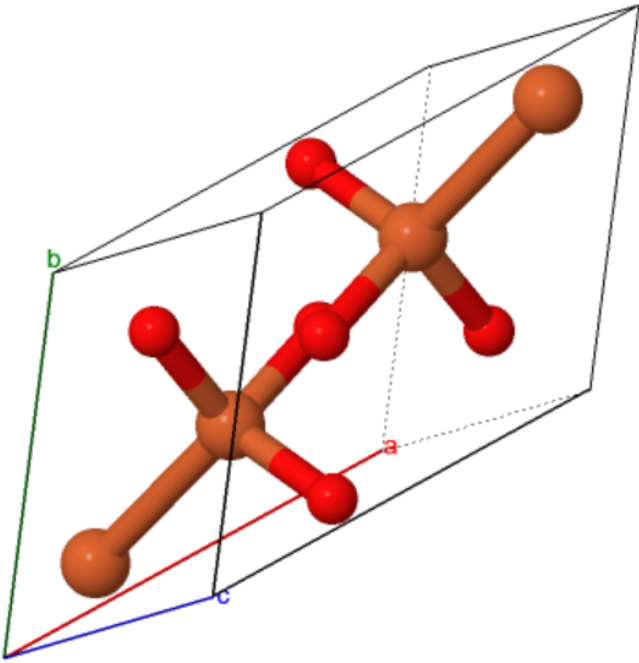
OPTIMIZE

SPECTRA

E&M

OTHER

P 1 [P 1] #1
a=5.461Å
b=5.461Å
c=5.461Å
α=55.773°
β=55.773°
γ=55.773°



STRUCTURE APPEARANCE

Select atom/s by:

by element

select

☐ within a plane

select All

unselect All

Halo/s off

Label On

Label Off

Atom/s & bond/s style

Atom/s colour:

Default colour

Bond colour:

Default colour

Atom/s & bond/s finish ☒ opaque ☐ translucent

select

Dot surface

select

☐ off

Atom/s & bond/s Size

Stick & Ball

Stick

Wire

Ball

CPK

Ionic

wireframe 0.2 Å

vdW radii 20 %

H-bonds: ☐ on ☒ off / solid H-bond ☐ on ☒ off

H-bond colour:

Default colour

View / Hide Hydrogen/s ☒

☐ Show Commands

Clear

History

Scripting Help

Collaboration with the University of Turin and Singapore National University

Beyond JSmol: SwingJS

Summer, 2016

Nadia El Mouldi

Andreas Raduege



Beyond JSmol: SwingJS

Summer, 2016

Q: Could this idea be generalized?

Q: Or does it end with JSmol?

Beyond JSmol: SwingJS

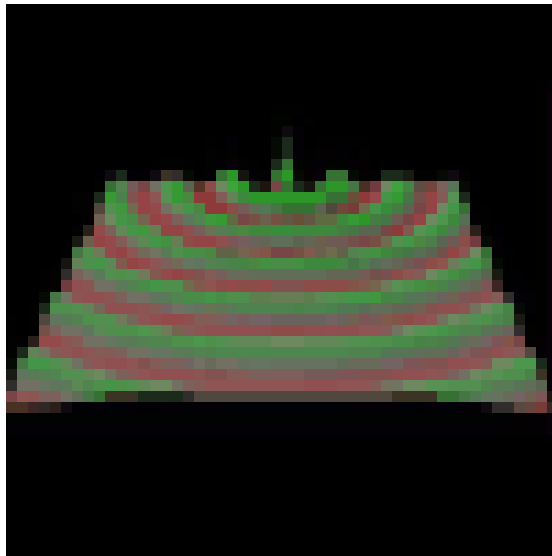
Summer, 2016

Task: Explore scope and limitations of java2script

Task: Contact possible collaborators

Task: Implement a few Java applets in JavaScript

Beyond JSmol: SwingJS



Paul Falstad

Minneapolis, MN



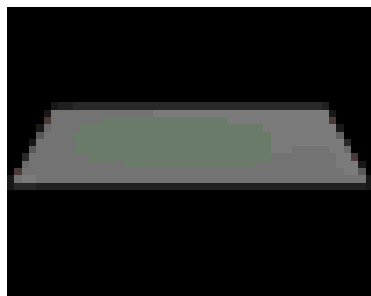
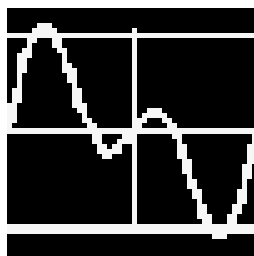
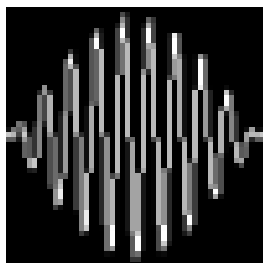
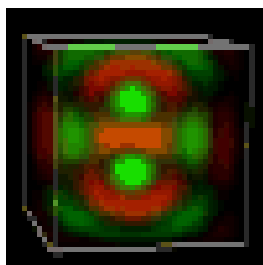
Ripple Tank (2-D Waves) Applet

Ripple tank simulation that demonstrates wave motion, interference, diffraction, refraction, Doppler effect, etc.

Beyond JSmol: SwingJS

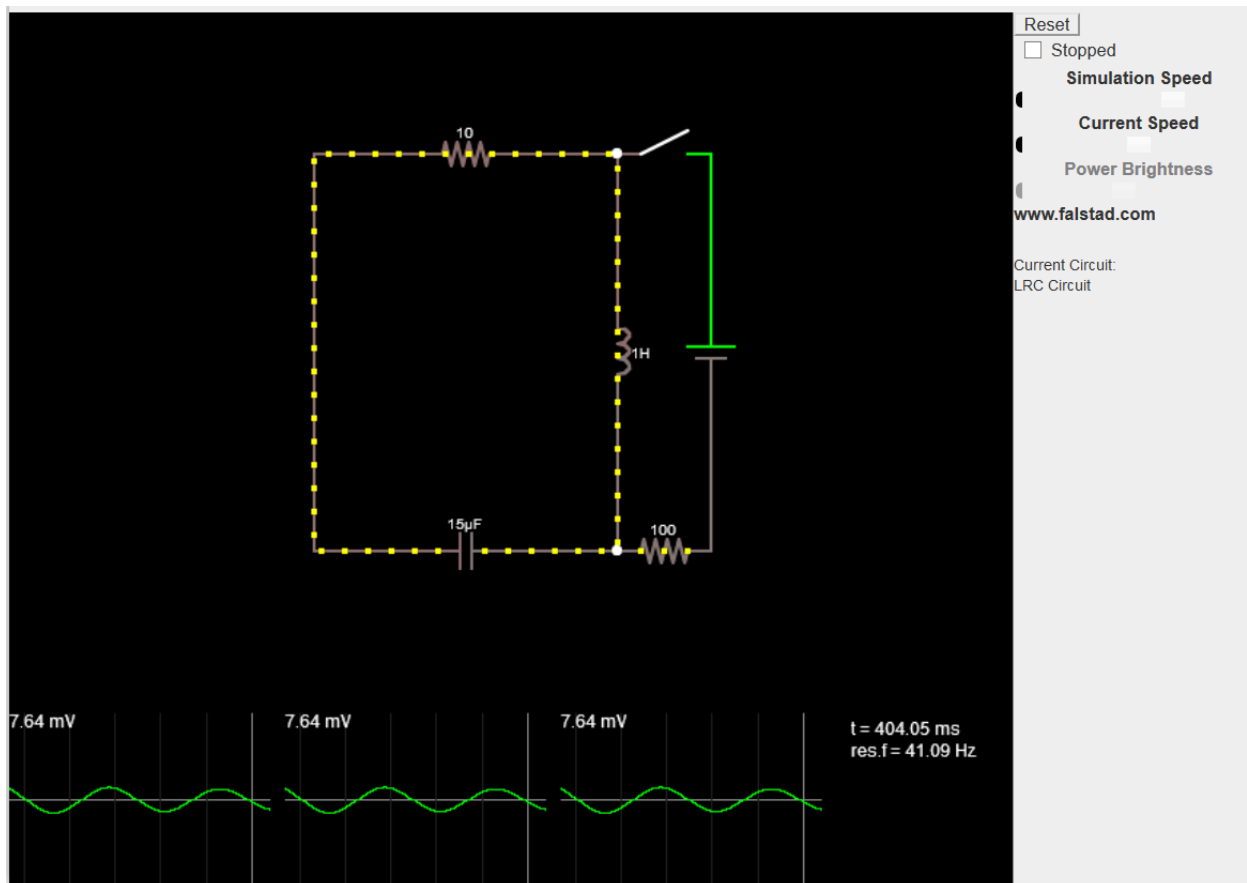
Paul Falstad

Minneapolis, MN



Beyond JSmol: SwingJS

circuit



Paul Falstad

Minneapolis, MN



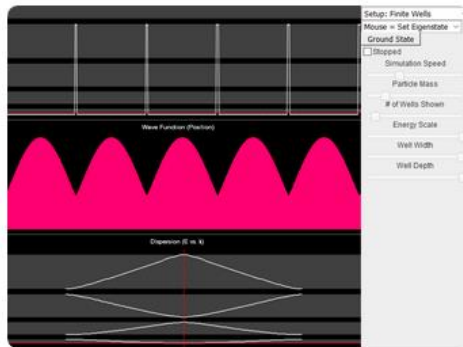
Demo Falstad Circuit

https://chemapps.stolaf.edu/swingjs/phet/site/swingjs/examples/applets/_Circuit_.html

Beyond JSmol: SwingJS

Physics - Solid State

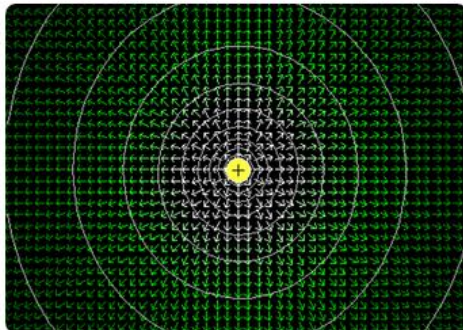
1-D Quantum Crystal



Periodic potentials in one dimension

Physics - Electromagnetics

2-D Electrostatics



Demonstrates static electric fields and steady-state current

Physics - Quantum

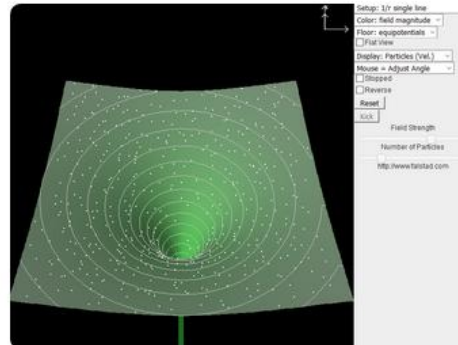
1-D Quantum States



Single-particle quantum mechanics states in one dimension

Mathematics

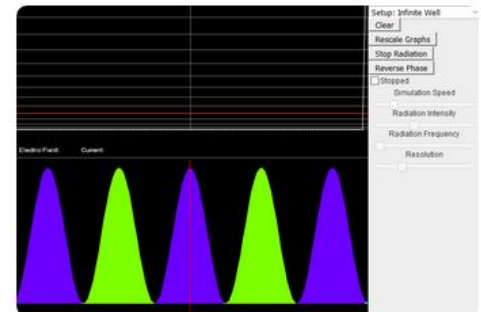
2-D Vector Fields



Demonstrates various properties of vector fields, including divergence and curl, etc.

Physics - Quantum

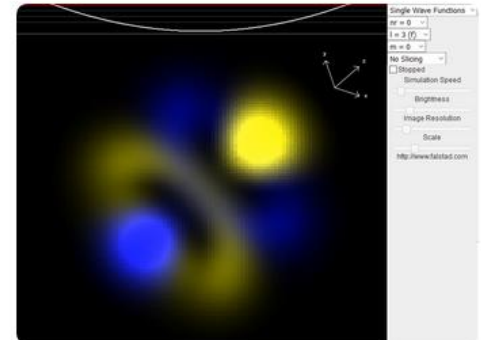
1-D Quantum Transitions



Radiative transitions (absorption and stimulated emission) in one dimension

Physics

3-D Quantum Harmonic Oscillator

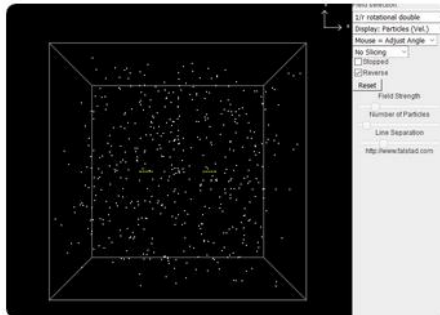


Harmonic oscillator in three dimensions

Beyond JSmol: SwingJS

Mathematics

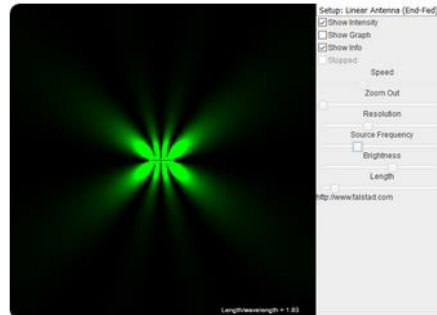
3-D Vector Fields



Demonstrates vector fields in 3 dimensions.
Includes the Lorenz Attractor and Rossler Attractor

Physics - Electronics

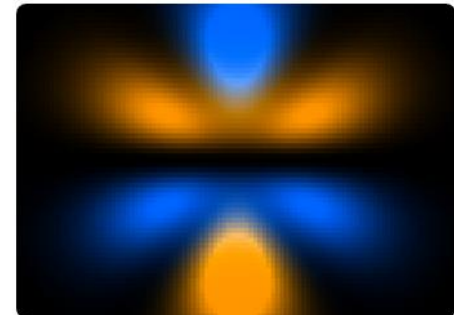
Antenna



Generates antenna radiation patterns

Physics - Atomic

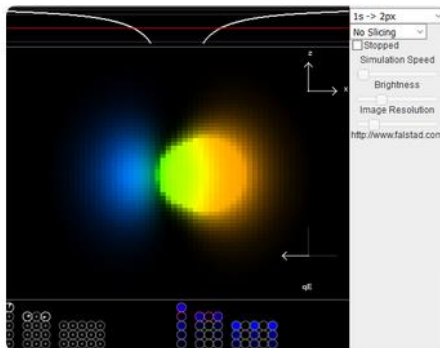
AtomViewer



Displays the wave functions (orbitals) of the hydrogen atom in 3-D

Physics - Atomic

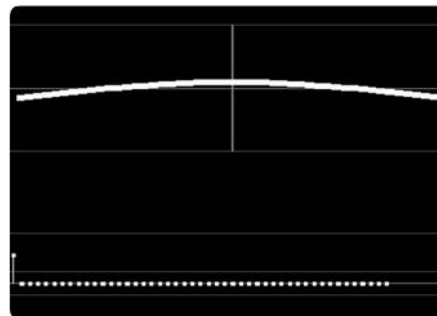
Atomic Dipole Transitions



Radiative transitions (absorption and stimulated emission) in atoms

Physics - Waves

BarWaves



A simulation that demonstrates standing flexural waves in a bar

Mathematics

Boids

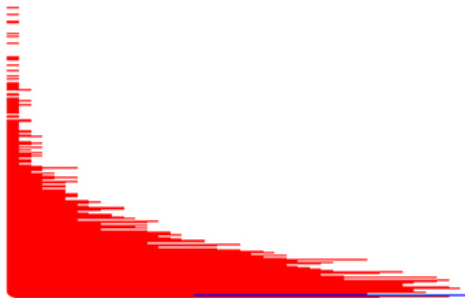


Illustrates a flocking algorithm

Beyond JSmol: SwingJS

Chemistry

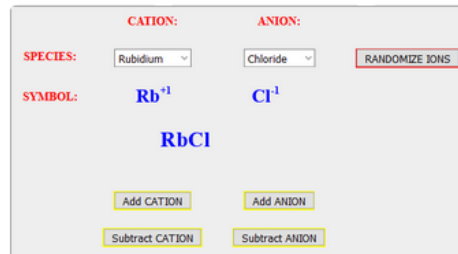
Boltzmann



A simulation of the Boltzmann distribution.

Chemistry

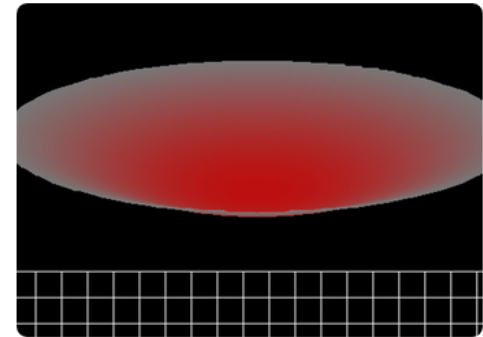
Chemical Charge

A web interface for creating chemical formulas. It has two main sections: 'CATION:' and 'ANION:'. Under 'CATION:', there is a 'SPECIES:' dropdown menu set to 'Rubidium', a 'SYMBOL:' label, and the text 'Rb⁺'. Under 'ANION:', there is an 'ANION:' dropdown menu set to 'Chloride', a 'SYMBOL:' label, and the text 'Cl⁻'. Below these, the text 'RbCl' is displayed. There are four buttons: 'Add CATION', 'Add ANION', 'Subtract CATION', and 'Subtract ANION'. A 'RANDOMIZE IONS' button is also present.

Combine charges to create a proper chemical formula

Physics - Mechanics

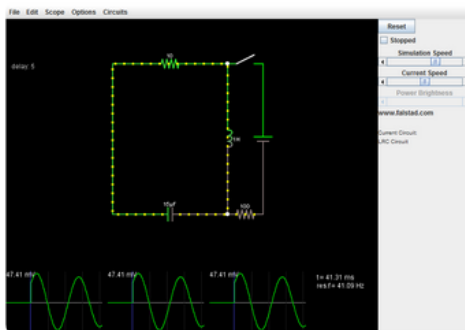
CircOsc



A simulation of waves in a circular membrane, showing its various vibrational modes.

Physics - Electronics

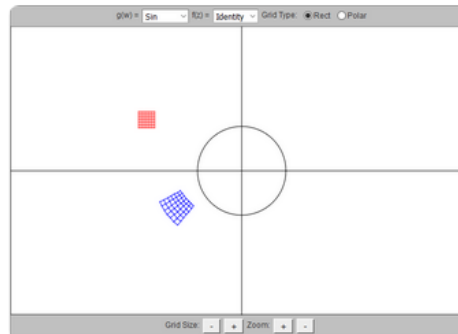
Circuit



An electronic circuit simulator. When the

Mathematics

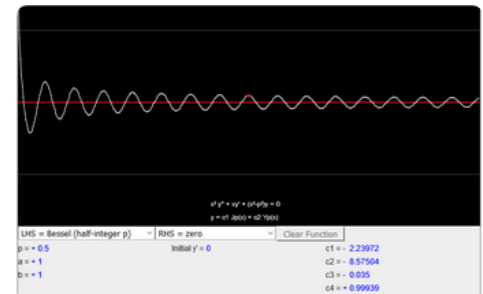
ComplexViewer



A viewer of mappings in the complex plane.

Mathematics

DiffEq

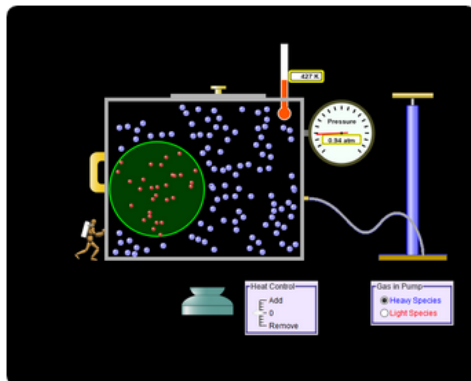


Visual differential equation solver.

Beyond JSmol: SwingJS

Chemistry - Ideal Gases

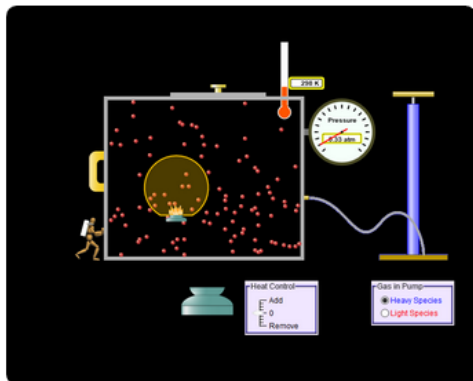
PhET Helium Balloon



Interactive ideal gas simulation involving a helium-filled balloon

Chemistry - Ideal Gases

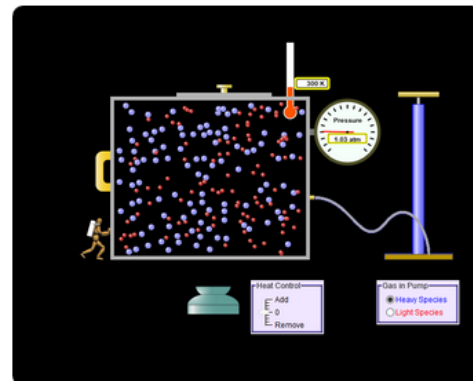
PhET Hot Air Balloon



Interactive ideal gas simulation involving a hot air balloon

Chemistry - Ideal Gases

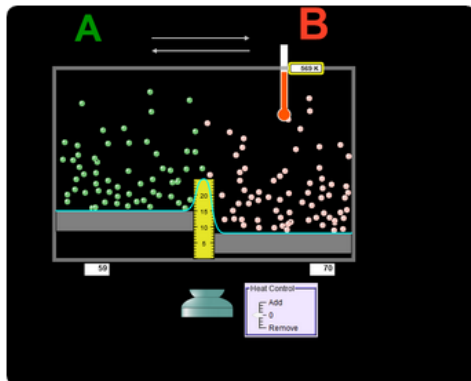
PhET Ideal Gas Properties



Ideal Gas Simulation

Chemistry - Reactions

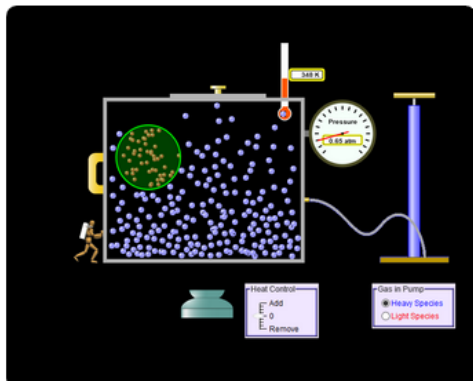
PhET Reversible Reactions



Interactive simulation of a reversible reaction

Chemistry - Ideal Gases

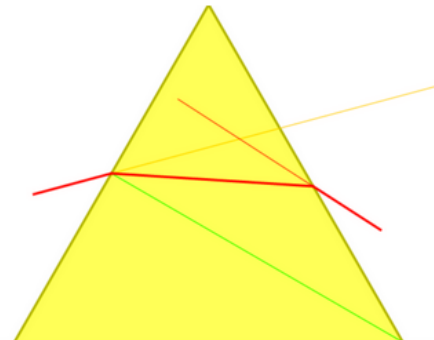
PhET Rigid Sphere



Interactive ideal gas simulation involving a

Physics - Optics

Prism

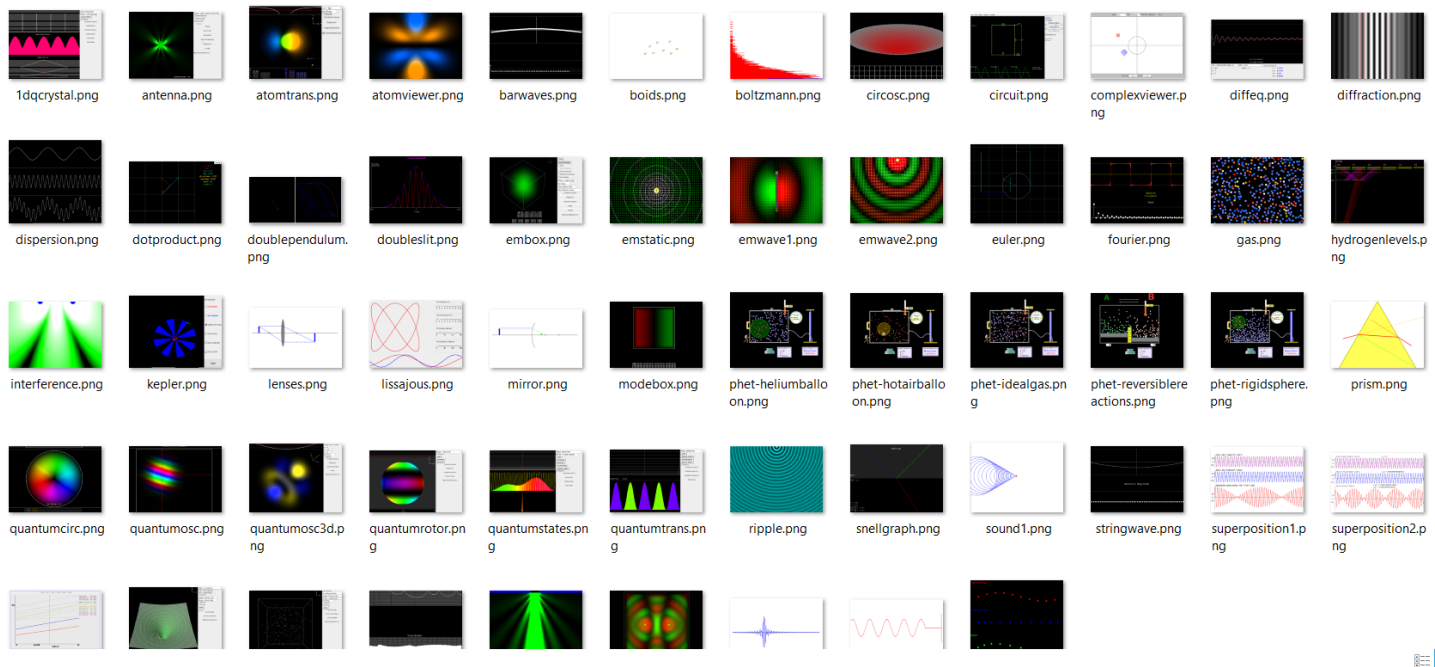


Illustrates the refraction of light by a prism

Beyond JSmol: SwingJS

Summer, 2016

– generated more than 50 applets



– discovered numerous transpiler and run-time issues

Further Generalization

Summer, 2017

Andrew Lee

Tahir Ahsan



Nikesh Yadav



Further Generalization

Zhou Renjian's first-generation transpiler was error-prone and still experimental, requiring many manually-generated work-arounds, primarily due to:

- minimal Java class implementation
- inaccurate class loading
- lack of numerical typing in JavaScript
byte, short, int, long, float, double
- no array typing
String[] int[] float[][] ...
- no method signature overloading
read(int) read(int, int) read(double) ...

Further Generalization

enter... Udo Borkowski

independent consultant

contributor to java2script

diplom in computer science – U. of Bonn

special interest in compiler construction



“Bob, it’s all about method binding. That has to be done by the transpiler, not at run time. I’m certain it can be done.”

Further Generalization

summer of 2017 into spring of 2017

- Completely rewritten transpiler now *100% consistent with Java class loading*
- Full implementation of the Java event queue
- Full numerical typing implemented for *char, byte, short, integer*, and (to the extent possible) *long*
- All method overloading issues resolved
- JavaScript single-thread issue solved

Beyond JSmol: SwingJS

Java:

```
package test;

import java.awt.Font;
import java.awt.Label;

import javax.swing.JApplet;

public class HelloWorld extends JApplet {

    @Override
    public void init() {

        Label label = new Label("Hello, World!");
        label.setFont(new Font(Font.SANS_SERIF, Font.PLAIN, 32));
        label.setAlignment(Label.CENTER);
        add(label);

    }

}
```

Hello, World!

Beyond JSmol: SwingJS

JavaScript:



Hello, World!



Hello, World!

```
var P$=Clazz.newPackage("test"),I$=[[0,'java.awt.Label','java.awt.Font']];
var C$=Clazz.newClass(P$, "HelloWorld", null, 'javax.swing.JApplet');

Clazz.newMeth(C$, ['init'], function () {
  var label=Clazz.new_($I$(1).c$$S,["Hello, World!"]);
  label.setFont$java_awt_Font(CClazz.new_($I$(2).c$$S$I$I,["SansSerif", 0, 32]));
  label.setAlignment$I(1);
  this.add$java_awt_Component(label);
});
```

Demo HelloWorld and Eclipse

Current Collaborations

spring/summer of 2018

- present

Physlet Physics

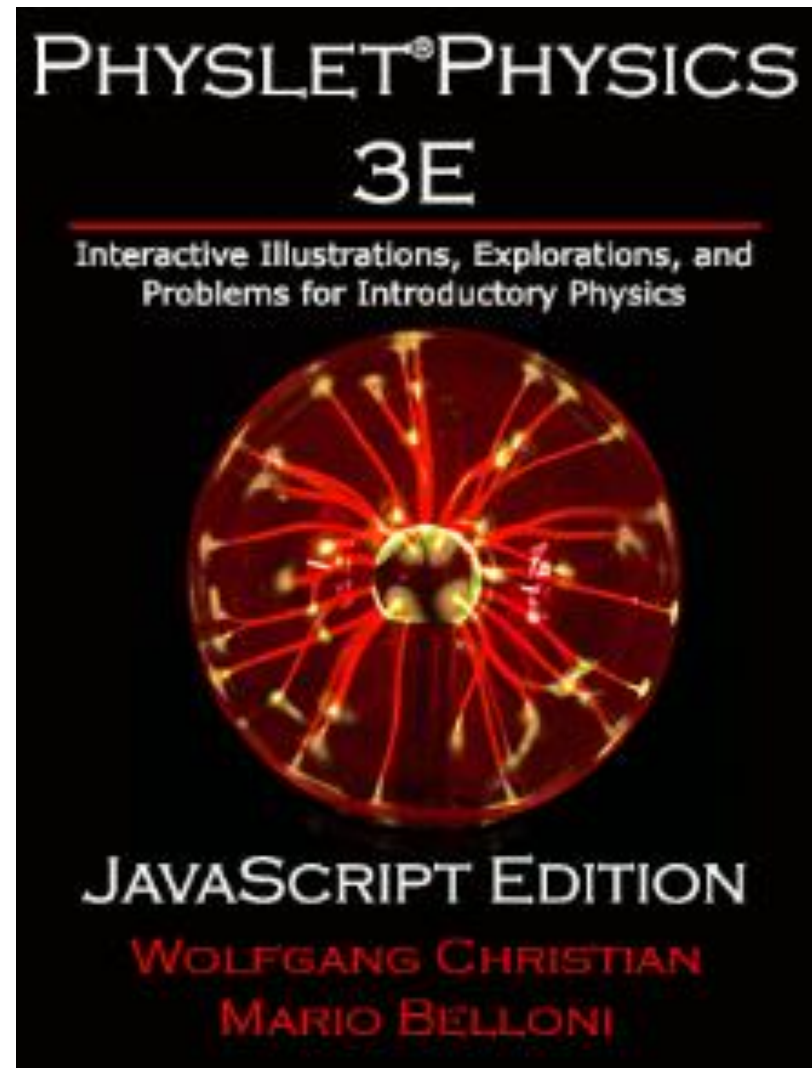
Wolfgang Christian

Davidson College

39 chapters of interactive
on-line introductory physics

30 applets; 800+ simulations

1500 Java classes



Current Collaborations

summer of 2018 –
present

Jalview

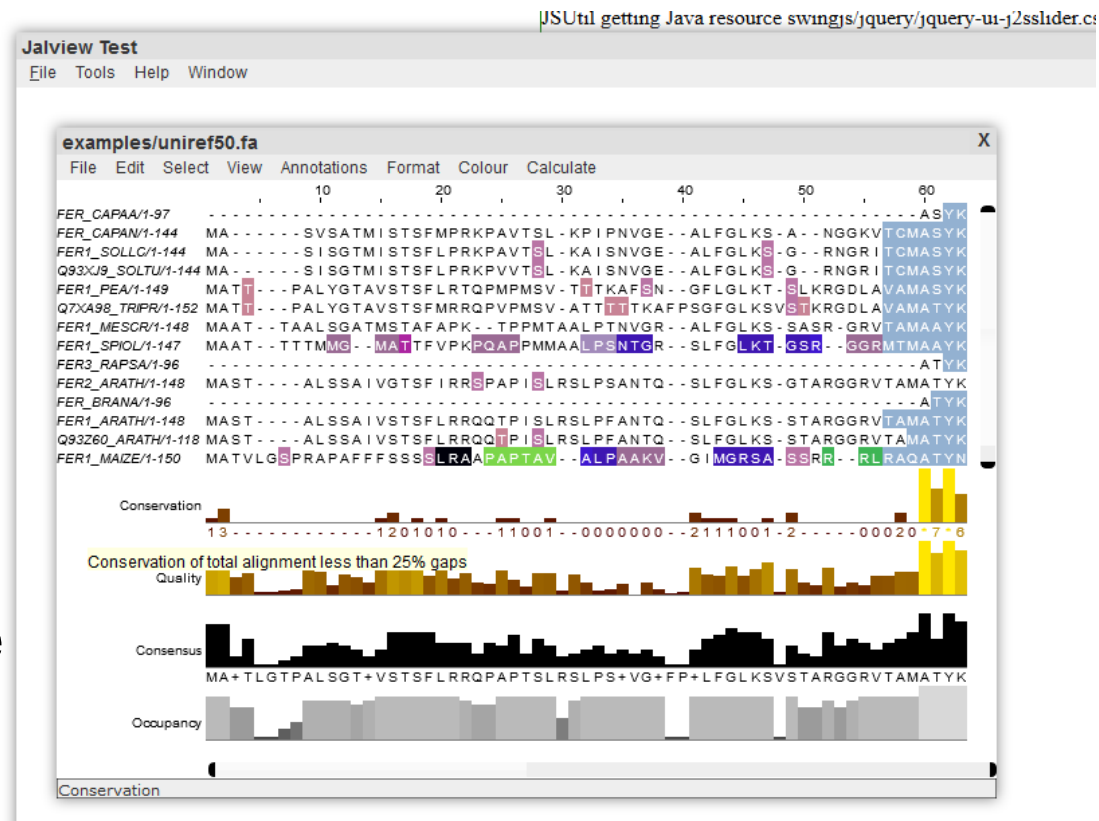
Geoff Barton

University of Dundee

desktop application

200,000 lines of code

3000 Java classes



Current Collaborations

fall of 2018 - present

Math^e(Prism)^a

Mathe-Prisma

Karsten Blankenagel
Wuppertal University

53 modules

500+ applets

1000+ web pages

4400 Java classes

Algebra
Arithmetik
Analysis
Geometrie
Graphentheorie
Stochastik
Informatik

Historical Context

The Problem and its Solution

Where we are now

Where we are headed

The Cost of Progress

Symmetry @ Otterbein

Home

Tutorial

Gallery

Challenge

Info

Feedback

Point Group Type:
All

Select Molecule:

- $\Lambda(\lambda,\lambda)$ -*cis*-[Co(en)₂Cl₂]⁺
- (δ,δ)-*trans*-[Co(en)₂Cl₂]⁺
- (δ,λ)-*trans*-[Co(en)₂Cl₂]⁺
- 1,1-dichlorethylene
- 1,2-dibromobenzene
- 1,2-dichlorethylene (cis)
- 1,2-dichloroethylene (tran
- 1,4-dibromobenzene
- 12-crown-4
- 18-crown-6
- trans*-cyclooctene
- closo*-[Ge₉]²⁻
- nido*-[Ge₉]⁴⁻

Preferences

[HTML5](#) • [Java](#)

Point Group = S₄

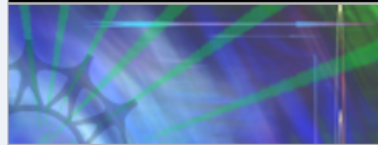
JSmol

Reference: Al-Jallal, N. A.; Al-Kahtani, A. A.; El-Azhary, A. A. J. *Phys. Chem. A* **2005**, *109*, 3694. DOI: [10.1021/jp050133c](https://doi.org/10.1021/jp050133c)

Element	Operation
<input type="checkbox"/> C ₂ axis	<input type="button" value="Rotate"/>
<input checked="" type="checkbox"/> S ₄ axis	<input type="button" value="Rotate"/>

The Cost of Progress

Awards



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Award Abstract #0536710

Visual-Spatial Learning: Development of an Interactive Web-Based Symmetry Tutorial

NSF Org:	DUE Division of Undergraduate Education
Initial Amendment Date:	February 24, 2006
Latest Amendment Date:	February 24, 2006
Award Number:	0536710
Award Instrument:	Standard Grant
Program Manager:	Eileen L. Lewis DUE Division of Undergraduate Education EHR Directorate for Education & Human Resources
Start Date:	April 1, 2006
End Date:	August 31, 2008 (Estimated)
Awarded Amount to Date:	\$57,641.00
Investigator(s):	Dean Johnston DJohnston@otterbein.edu (Principal Investigator)
Sponsor:	Otterbein College 1 S. Grove St. Westerville, OH 43081-2006 (614)823-1846

...and the reward for success

Awards

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Award Abstract #0536710
Visual-Spatial Learning: Development of an Interactive Web-Based Symmetry Tutorial

NSF Org:	DUE Division of Undergraduate Education
Initial Amendment Date:	February 24, 2006
Latest Amendment Date:	February 24, 2006

Johnston, Dean

2:35 PM (15 minutes ago)


to me ▼

Yes - infinite thanks to you and all your work to make the things I have done still useful after all these years...

[Policy Office Website](#)


End Date:	August 31, 2008 (Estimated)
Awarded Amount to Date:	\$57,641.00
Investigator(s):	Dean Johnston DJohnston@otterbein.edu (Principal Investigator)
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The Cost of Progress



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Physlet® Physics 3E

Interactive Illustrations, Explorations, and Problems for Introductory Physics

Wolfgang Christian
Mario Belloni

With contributing authors:
Anne Cox, Melissa H. Dancy, and Aaron Titus

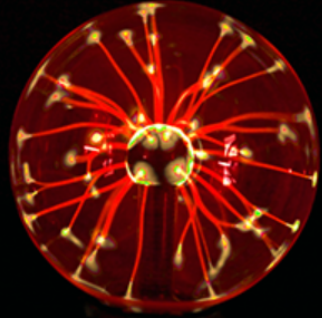
Exploration Worksheets by:
Thomas M. Colbert

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PHYSLET® PHYSICS

3E

Interactive Illustrations, Explorations, and Problems for Introductory Physics



JAVASCRIPT EDITION

WOLFGANG CHRISTIAN
MARIO BELLONI

NSF Awards – “physlets”

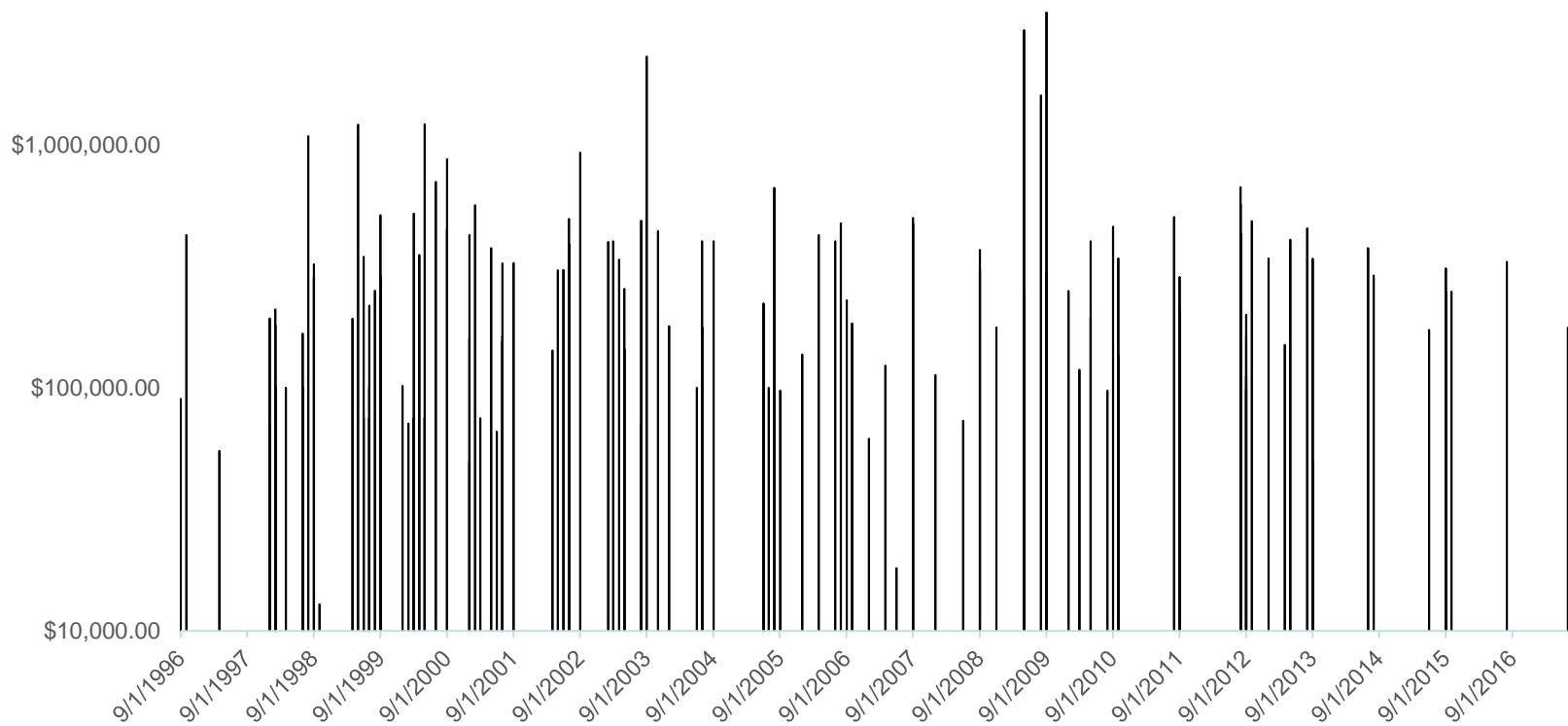
\$2,272,543

Title	Award
Ubiquitous Contextual Access to STEM Educational Resources (UCASTER)	\$521,702
Cross-Linked Models (XLM)	\$149,977
OPTIC: Open Physics Technology for Interactive Curricula	\$450,000
Revising intermediate optics for greater conceptual understanding and the development of complex problem solving skills	\$60,566
Open-Source Physics EducationN: OPEN	\$494,715
Creation of a Computation and Visualization Laboratory	\$53,466
Multimedia-focused Kinematics Questions: An Innovative Approach to Teaching Kinematics	\$74,410
Computerizing Introductory Physics Laboratories to Improve Student Learning	\$31,545
Implementing the Workshop Model and other Research-based Instructional Strategies in Physics & Mathematics Courses	\$81,207
Studio Based Modern Physics	\$50,205
WebPhysics	\$304,750

NSF Awards – “applets”

\$47,191,650

\$10,000,000.00 Total NSF "applet" awards 1996-2017: 134 (\$47,191,650)



NSF Awards – “applets”

\$47,191,650



Summary

- SwingJS “proof of concept” is complete – around 10,000 Java classes implemented and working smoothly.
- Developer tools are now in place so that others can proceed with minimal assistance (GitHub, Eclipse).
- Still a few Java Swing classes to fully implement (e.g. JTree). Testing/debugging continues.
- Not totally automatic. A certain amount of Java redesign may be necessary, and performance optimization is critical.

Summary

- The success of Jmol/JSmol has been generalized to a wide range of Java applets and applications across the fields of mathematics, physics, chemistry, and biology.
- Java may be gone from the web, but, oddly enough, that is not the case for *Java applets!* *They can live on as JavaScript.*
- java2script/SwingJS allows for re-enabling \$Ms of superb resources that have been lost in the past few years.
- Our technology allows for a novel “Java+JavaScript” option for web app and stand-alone application developers.

Thanks!

St. Olaf Students

Nadia El Mouldi
Andreas Raduege

Andrew Lee
Tahir Ahsan
Nikesh Yadav

Thanks!

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Thanks!

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Karsten Blankenagel

Friends at St. Olaf

Matt Richey
Dick Brown
Olaf Hall-Holt

MSCS for use of anansi server
(<https://chemapps.stolaf.edu>)

Thanks!

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(<https://chemapps.stolaf.edu>)

and you, for your attention