**Chemical Structure Simulator (CSS)** Documentation

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# Introduction

## Purpose of the documentation

The purpose of this document is to allow the user to fully understand all the features of the application and how to effectively make use of them.

## How to use this document

The documentation is split into different sections. Each section will contain a single feature. Each section will be further divided into multiple subsections: general summary of the feature, how to use it, what to expect from it, and a more technical description of the feature. At the end of this document, there will be a list of shortcuts that can be used in the application to speed up some tasks. In the appendix, there will be the features of the algorithms.

# Prerequisites

## Getting this application

The entirety of this application, along with the optional source code, is available on GitHub. To install the executable and/or the source code, please visit and clone:

<https://github.com/volovikariel/ChemicalSimulator>

## Dependencies

### Java

To run this application natively on Java (using the jar file instead of the executable), the computer running this application must also have JRE (Java Runtime Environment) installed. Otherwise, the user must use the executable (\*.exe, \*.out).

## Specifications

Since the algorithm this application is based on is a very broad and thus very computationally expensive, it is highly recommended to have at least 2GB of memory available on the device.

# Features

The features that follow are all features that affect the user directly, sparing the overly technical descriptions of the algorithms, which are accessible in the appendix.

## Periodic Table

### Summary

The Periodic Table is available as a form of reference, and as a form of input for the user.

### How to use

The user can drag and drop individual atoms into the top portion of the screen, this will add it to the list of atoms that will be considered for the algorithm. The user can also right click to remove them, or alternatively drag and drop them back into the periodic table. The user can also delete individual atoms by pressing backspace after having added it.

### What to expect

The user can expect intuitive controls for the addition and removal of atoms. The user can also expect to have a periodic table which looks familiar to the one they’re used to, that is to say: it is ordered in the standard form so that the user knows what column the atom is situated in, it is colour coded to facilitate finding metals and non-metals and other subgroups, it shows the atom’s atomic number. [POSSIBLY ADD MORE INFO LIKE ATOMIC MASS WHEN IT MATTERS IN ALGORITHM]

### Technical Description

## Text Input

### Summary

The text input is a way for the user to input the set of atoms for which the solutions will be computed. If the user intends to write an equation with many of the same atom, it bypasses the monotony of manually placing the same atom over and over by dragging and dropping from the periodic table.

### How to use

The user can simply start typing the chemical compound’s equation and a text field will pop up, allowing the user to keep track of what has been written. The user can at any time press backspace, and an atom will be removed from the list. The user has **two main ways to input the equation**. The user can type an atom at a time, **HOH**, or they can type **H2O**, both will be properly interpreted. The input is **case sensitive**, this is to avoid bizarre cases such as inputting, CO instead of Co (Carbon Monoxide vs Cobalt).

### What to expect

The user can expect the input to be rather intuitive and responsive. The user can expect to have a seamless transition between using the periodic table to add elements and the text input, for instance, if the user wants to add several carbons, the user can drag and drop a single carbon and then using their keyboard, write how many carbons they want.

### Technical Description

## Loading Screen

### Summary

The loading screen is visible whilst the algorithm is calculating the solutions. This is after the user has pressed enter and is ready to see their results, but before the results have been finalized.

### How to use

The user can leave the application whilst the algorithm is doing its job, the user will receive a notification in the bottom right of their screen once the algorithm has finished.

### What to expect

If the user wishes to look at the loading screen, they will see molecules swirling around.

### Technical Description

## Tabs

### Summary

The tabs are what the Lewis Structure and the Three-Dimensional Display are contained in. The user can choose which tab they want selected by clicking the numbered tabs at the top of the screen.

### How to use

The user can play around with the Lewis Structure and Three-Dimensional Display in one tab, and then go to the next. The user can, easily navigate between tabs by tapping the left and right arrow keys.

### What to expect

The tabs are generally **ordered in terms of likeliness** to appear, which is calculated by the algorithm, so the solution in tab 1 is generally more likely to appear than in tab 10. Do note that this isn’t entirely accurate because there are many things to consider. Once the user has rotated or moved the *group*, it remains in that position, so the user can compare their angles and such by jumping between tabs.

### Technical Description

## Lewis Structure

### Summary

### How to use

### What to expect

### Technical Description

## Three-Dimensional Display

### Summary

The Three-Dimensional Display is a 3D representation of the chemical compound (the *group*).

### How to use

The *group* can be **rotated** by pressing down the primary mouse button and moving the cursor.

The *group* can be **translated**/**moved** by pressing down the secondary mouse button (usually the right) and moving the cursor to where you want the group to be.

*group* can be **zoomed in and out** by using the scroll wheel.

A *group’s* atom can be hovered over and its **symbol** will be **displayed** at the bottom of the screen.

### What to expect

The user can expect the angles between atoms to be close to what they are theoretically.

The user can expect to see the proper bond counts for each solution.

The user can expect to see the formal charge on each atom if it is not neutral.

The atoms are distinguishable one from the other, for instance hydrogen atoms as seen in the group are of a lighter colour and of a smaller size. The user can also display its symbol by hovering over it.

### Technical Description

# Shortcuts

# Appendix

## Structures

## Sorting