

Every public MolDraw page

266 public pages with their web addresses, grouped by section.

MolDraw (<https://www.moldraw.com>) · Help centre: <https://www.moldraw.com/support/> · Updated 2026-10-17

Start (1)

- <https://www.moldraw.com/> — <https://www.moldraw.com/>
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Documentation (36)

- [MolDraw Documentation | 2D Editor Guide](https://www.moldraw.com/documentation/) — <https://www.moldraw.com/documentation/>
- [MolDraw Docs: how to use the MolDraw chemistry editor](https://www.moldraw.com/docs/) — <https://www.moldraw.com/docs/>
- [Introduction to MolDraw](https://www.moldraw.com/docs/get-started/what-is-moldraw-free-online-chemical-structure-editor/) — <https://www.moldraw.com/docs/get-started/what-is-moldraw-free-online-chemical-structure-editor/>
- [Draw your first structure](https://www.moldraw.com/docs/get-started/how-to-draw-your-first-chemical-structure-in-moldraw/) — <https://www.moldraw.com/docs/get-started/how-to-draw-your-first-chemical-structure-in-moldraw/>
- [The MolDraw screen](https://www.moldraw.com/docs/get-started/moldraw-editor-interface-toolbar-and-panels-explained/) — <https://www.moldraw.com/docs/get-started/moldraw-editor-interface-toolbar-and-panels-explained/>
- [Save and open your work](https://www.moldraw.com/docs/get-started/how-to-save-and-open-chemical-drawings-in-moldraw/) — <https://www.moldraw.com/docs/get-started/how-to-save-and-open-chemical-drawings-in-moldraw/>
- [Keyboard shortcuts](https://www.moldraw.com/docs/get-started/moldraw-keyboard-shortcuts-list/) — <https://www.moldraw.com/docs/get-started/moldraw-keyboard-shortcuts-list/>
- [Draw atoms](https://www.moldraw.com/docs/drawing/how-to-draw-atoms-and-change-elements/) — <https://www.moldraw.com/docs/drawing/how-to-draw-atoms-and-change-elements/>
- [Draw bonds](https://www.moldraw.com/docs/drawing/how-to-draw-single-double-triple-and-wedge-bonds/) — <https://www.moldraw.com/docs/drawing/how-to-draw-single-double-triple-and-wedge-bonds/>
- [Rings and structures](https://www.moldraw.com/docs/drawing/how-to-draw-benzene-cyclohexane-and-fused-rings/) — <https://www.moldraw.com/docs/drawing/how-to-draw-benzene-cyclohexane-and-fused-rings/>
- [Atom labels and R-groups](https://www.moldraw.com/docs/drawing/how-to-add-atom-labels-and-r-groups/) — <https://www.moldraw.com/docs/drawing/how-to-add-atom-labels-and-r-groups/>
- [Stereochemistry](https://www.moldraw.com/docs/drawing/how-to-draw-stereochemistry-wedge-hash-and-r-s-labels/) — <https://www.moldraw.com/docs/drawing/how-to-draw-stereochemistry-wedge-hash-and-r-s-labels/>
- [Lone pairs, radicals and charges](https://www.moldraw.com/docs/drawing/how-to-add-lone-pairs-radicals-and-formal-charges/) — <https://www.moldraw.com/docs/drawing/how-to-add-lone-pairs-radicals-and-formal-charges/>
- [Polymer brackets](https://www.moldraw.com/docs/drawing/how-to-draw-polymer-repeat-unit-brackets/) — <https://www.moldraw.com/docs/drawing/how-to-draw-polymer-repeat-unit-brackets/>
- [Text, shapes and images](https://www.moldraw.com/docs/drawing/how-to-add-text-shapes-images-and-lab-glassware/) — <https://www.moldraw.com/docs/drawing/how-to-add-text-shapes-images-and-lab-glassware/>
- [Style and display](https://www.moldraw.com/docs/drawing/how-to-change-structure-style-colors-and-display/) — <https://www.moldraw.com/docs/drawing/how-to-change-structure-style-colors-and-display/>
- [Reaction arrows](https://www.moldraw.com/docs/reactions/how-to-draw-reaction-arrows-with-reagents/) — <https://www.moldraw.com/docs/reactions/how-to-draw-reaction-arrows-with-reagents/>
- [Build a reaction scheme](https://www.moldraw.com/docs/reactions/how-to-make-a-chemical-reaction-scheme/) — <https://www.moldraw.com/docs/reactions/how-to-make-a-chemical-reaction-scheme/>
- [Library and templates](https://www.moldraw.com/docs/library/how-to-use-the-structure-library-and-templates/) — <https://www.moldraw.com/docs/library/how-to-use-the-structure-library-and-templates/>

- [Copy and paste SMILES](https://www.moldraw.com/docs/import-export/how-to-copy-and-paste-smiles-inchi-and-mol/) — <https://www.moldraw.com/docs/import-export/how-to-copy-and-paste-smiles-inchi-and-mol/>
- [Download and file formats](https://www.moldraw.com/docs/import-export/how-to-export-png-svg-pdf-mol-and-cdxml-files/) — <https://www.moldraw.com/docs/import-export/how-to-export-png-svg-pdf-mol-and-cdxml-files/>
- [Search PubChem](https://www.moldraw.com/docs/import-export/how-to-search-pubchem-by-name-or-cas-number/) — <https://www.moldraw.com/docs/import-export/how-to-search-pubchem-by-name-or-cas-number/>
- [Names, formula and properties](https://www.moldraw.com/docs/analyze/how-to-find-iupac-name-molecular-formula-and-weight/) — <https://www.moldraw.com/docs/analyze/how-to-find-iupac-name-molecular-formula-and-weight/>
- [3D viewer](https://www.moldraw.com/docs/analyze/how-to-view-molecules-in-3d/) — <https://www.moldraw.com/docs/analyze/how-to-view-molecules-in-3d/>
- [AI assistant](https://www.moldraw.com/docs/ai/how-to-use-the-ai-chemistry-assistant/) — <https://www.moldraw.com/docs/ai/how-to-use-the-ai-chemistry-assistant/>
- [Plugins overview](https://www.moldraw.com/docs/plugins/moldraw-plugins-overview-and-how-to-install/) — <https://www.moldraw.com/docs/plugins/moldraw-plugins-overview-and-how-to-install/>
- [NMR Analyser plugin](https://www.moldraw.com/docs/plugins/nmr-analyser-plugin-predict-1h-13c-and-2d-nmr-spectra/) — <https://www.moldraw.com/docs/plugins/nmr-analyser-plugin-predict-1h-13c-and-2d-nmr-spectra/>
- [ChEMBL plugin](https://www.moldraw.com/docs/plugins/chembl-plugin-bioactivity-and-drug-target-lookup/) — <https://www.moldraw.com/docs/plugins/chembl-plugin-bioactivity-and-drug-target-lookup/>
- [Electrostatic Potential \(MEP\) plugin](https://www.moldraw.com/docs/plugins/mep-plugin-molecular-electrostatic-potential-surface/) — <https://www.moldraw.com/docs/plugins/mep-plugin-molecular-electrostatic-potential-surface/>
- [Compare Molecules plugin](https://www.moldraw.com/docs/plugins/compare-molecules-plugin-tanimoto-similarity/) — <https://www.moldraw.com/docs/plugins/compare-molecules-plugin-tanimoto-similarity/>
- [Protein Viewer plugin](https://www.moldraw.com/docs/plugins/protein-viewer-plugin-view-pdb-structures-in-3d/) — <https://www.moldraw.com/docs/plugins/protein-viewer-plugin-view-pdb-structures-in-3d/>
- [Smart Draw plugin](https://www.moldraw.com/docs/plugins/smart-draw-plugin-hand-drawn-sketch-to-structure/) — <https://www.moldraw.com/docs/plugins/smart-draw-plugin-hand-drawn-sketch-to-structure/>
- [AI Spectroscopy plugin](https://www.moldraw.com/docs/plugins/ai-spectroscopy-plugin-predict-nmr-ms-and-uv-vis/) — <https://www.moldraw.com/docs/plugins/ai-spectroscopy-plugin-predict-nmr-ms-and-uv-vis/>
- [Themes \(Ball-and-Stick\) plugin](https://www.moldraw.com/docs/plugins/themes-plugin-ball-and-stick-2d-style/) — <https://www.moldraw.com/docs/plugins/themes-plugin-ball-and-stick-2d-style/>
- [Symmetry and Point Groups plugin](https://www.moldraw.com/docs/plugins/symmetry-plugin-point-group-finder-3d-symmetry-operations/) — <https://www.moldraw.com/docs/plugins/symmetry-plugin-point-group-finder-3d-symmetry-operations/>
- [pKa and Microspecies plugin](https://www.moldraw.com/docs/plugins/pka-plugin-microspecies-protonation-states-at-any-ph/) — <https://www.moldraw.com/docs/plugins/pka-plugin-microspecies-protonation-states-at-any-ph/>

Learn (course) (39)

- [Learn MolDraw: free hands-on chemistry drawing lessons](https://www.moldraw.com/learn/) — <https://www.moldraw.com/learn/>
- [How to add atoms, elements and formal charges | Learn MolDraw](https://www.moldraw.com/learn/how-to-add-atoms-elements-and-formal-charges/) — <https://www.moldraw.com/learn/how-to-add-atoms-elements-and-formal-charges/>
- [How to add ligands and functional groups | Learn MolDraw](https://www.moldraw.com/learn/how-to-add-ligands-and-functional-groups-from-the-library/) — <https://www.moldraw.com/learn/how-to-add-ligands-and-functional-groups-from-the-library/>
- [How to add text and shapes to a drawing | Learn MolDraw](https://www.moldraw.com/learn/how-to-add-text-and-shapes-to-a-chemical-drawing/) — <https://www.moldraw.com/learn/how-to-add-text-and-shapes-to-a-chemical-drawing/>
- [How to align, arrange and space out structures | Learn MolDraw](https://www.moldraw.com/learn/how-to-align-arrange-and-distribute-structures-on-the-canvas/) — <https://www.moldraw.com/learn/how-to-align-arrange-and-distribute-structures-on-the-canvas/>
- [How to change atom colours and colour a structure | Learn MolDraw](https://www.moldraw.com/learn/how-to-change-atom-colours-and-colour-structures-in-moldraw/) — <https://www.moldraw.com/learn/how-to-change-atom-colours-and-colour-structures-in-moldraw/>
- [How to change font size and atom label style | Learn MolDraw](https://www.moldraw.com/learn/how-to-change-font-size-and-atom-label-style-in-chemical-drawings/) — <https://www.moldraw.com/learn/how-to-change-font-size-and-atom-label-style-in-chemical-drawings/>

- [How to clean up a chemical structure | Learn MolDraw](https://www.moldraw.com/learn/how-to-clean-up-a-chemical-structure-drawing/) — <https://www.moldraw.com/learn/how-to-clean-up-a-chemical-structure-drawing/>
- [How to compare two molecules for similarity | Learn MolDraw](https://www.moldraw.com/learn/how-to-compare-two-molecules-for-similarity/) — <https://www.moldraw.com/learn/how-to-compare-two-molecules-for-similarity/>
- [How to turn a name or SMILES into a structure | Learn MolDraw](https://www.moldraw.com/learn/how-to-convert-a-name-or-smiles-to-a-structure/) — <https://www.moldraw.com/learn/how-to-convert-a-name-or-smiles-to-a-structure/>
- [How to draw a reaction scheme with arrows | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-a-chemical-reaction-scheme-with-arrows/) — <https://www.moldraw.com/learn/how-to-draw-a-chemical-reaction-scheme-with-arrows/>
- [How to draw a benzene ring and phenol | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-benzene-ring-and-phenol-structure/) — <https://www.moldraw.com/learn/how-to-draw-benzene-ring-and-phenol-structure/>
- [How to draw double, triple and wedge bonds | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-chemical-bonds-double-triple-and-wedge-bonds/) — <https://www.moldraw.com/learn/how-to-draw-chemical-bonds-double-triple-and-wedge-bonds/>
- [How to draw curved arrow mechanisms | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-curved-arrow-reaction-mechanisms/) — <https://www.moldraw.com/learn/how-to-draw-curved-arrow-reaction-mechanisms/>
- [How to draw lab apparatus and glassware | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-lab-apparatus-and-glassware/) — <https://www.moldraw.com/learn/how-to-draw-lab-apparatus-and-glassware/>
- [How to draw lone pairs and radicals | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-lone-pairs-and-radicals-on-atoms/) — <https://www.moldraw.com/learn/how-to-draw-lone-pairs-and-radicals-on-atoms/>
- [How to draw polymer repeat unit brackets | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-polymer-repeat-unit-brackets/) — <https://www.moldraw.com/learn/how-to-draw-polymer-repeat-unit-brackets/>
- [How to draw rings, chairs and fused rings | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-rings-cyclohexane-chair-and-fused-rings/) — <https://www.moldraw.com/learn/how-to-draw-rings-cyclohexane-chair-and-fused-rings/>
- [How to draw stereochemistry with wedge and hash bonds | Learn MolDraw](https://www.moldraw.com/learn/how-to-draw-stereochemistry-with-wedge-and-hash-bonds/) — <https://www.moldraw.com/learn/how-to-draw-stereochemistry-with-wedge-and-hash-bonds/>
- [How to export a structure for a paper or slides | Learn MolDraw](https://www.moldraw.com/learn/how-to-export-a-chemical-structure-for-a-paper-or-slides/) — <https://www.moldraw.com/learn/how-to-export-a-chemical-structure-for-a-paper-or-slides/>
- [How to export MOL, SVG and other file formats | Learn MolDraw](https://www.moldraw.com/learn/how-to-export-mol-svg-and-other-chemical-file-formats/) — <https://www.moldraw.com/learn/how-to-export-mol-svg-and-other-chemical-file-formats/>
- [How to find the formula, weight and IUPAC name | Learn MolDraw](https://www.moldraw.com/learn/how-to-find-molecular-formula-weight-and-iupac-name/) — <https://www.moldraw.com/learn/how-to-find-molecular-formula-weight-and-iupac-name/>
- [How to look up a molecule in ChEMBL | Learn MolDraw](https://www.moldraw.com/learn/how-to-look-up-a-molecule-in-chembl/) — <https://www.moldraw.com/learn/how-to-look-up-a-molecule-in-chembl/>
- [How to make publication quality structure figures | Learn MolDraw](https://www.moldraw.com/learn/how-to-make-publication-quality-chemical-structure-figures/) — <https://www.moldraw.com/learn/how-to-make-publication-quality-chemical-structure-figures/>
- [How to map atoms in a reaction | Learn MolDraw](https://www.moldraw.com/learn/how-to-map-atoms-in-a-reaction/) — <https://www.moldraw.com/learn/how-to-map-atoms-in-a-reaction/>
- [How to open the AI chemistry assistant | Learn MolDraw](https://www.moldraw.com/learn/how-to-open-the-ai-chemistry-assistant/) — <https://www.moldraw.com/learn/how-to-open-the-ai-chemistry-assistant/>
- [How to predict a mass spectrum | Learn MolDraw](https://www.moldraw.com/learn/how-to-predict-a-mass-spectrum/) — <https://www.moldraw.com/learn/how-to-predict-a-mass-spectrum/>
- [How to predict an NMR spectrum from a structure | Learn MolDraw](https://www.moldraw.com/learn/how-to-predict-an-nmr-spectrum-from-a-structure/) — <https://www.moldraw.com/learn/how-to-predict-an-nmr-spectrum-from-a-structure/>
- [How to search PubChem by name or CAS number | Learn MolDraw](https://www.moldraw.com/learn/how-to-search-pubchem-by-name-or-cas-number/) — <https://www.moldraw.com/learn/how-to-search-pubchem-by-name-or-cas-number/>
- [How to set bond length, line width and bond style | Learn MolDraw](https://www.moldraw.com/learn/how-to-set-bond-length-line-width-and-bond-style/) — <https://www.moldraw.com/learn/how-to-set-bond-length-line-width-and-bond-style/>
- [How to show an electrostatic potential map | Learn MolDraw](https://www.moldraw.com/learn/how-to-show-an-electrostatic-potential-map/) — <https://www.moldraw.com/learn/how-to-show-an-electrostatic-potential-map/>
- [How to use Community, Dashboard and your profile | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-community-dashboard-and-profile/) — <https://www.moldraw.com/learn/how-to-use-community-dashboard-and-profile/>

- [How to use keyboard shortcuts to draw faster | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-keyboard-shortcuts-to-draw-faster/) — <https://www.moldraw.com/learn/how-to-use-keyboard-shortcuts-to-draw-faster/>
 - [How to use R groups and labels like Ph and COOH | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-r-groups-and-atom-labels-like-ph-and-cooh/) — <https://www.moldraw.com/learn/how-to-use-r-groups-and-atom-labels-like-ph-and-cooh/>
 - [How to use ring templates and the structure library | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-ring-templates-and-the-structure-library/) — <https://www.moldraw.com/learn/how-to-use-ring-templates-and-the-structure-library/>
 - [How to sketch structures with Smart Draw | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-smart-draw-to-sketch-structures/) — <https://www.moldraw.com/learn/how-to-use-smart-draw-to-sketch-structures/>
 - [How to use themes and ball and stick style | Learn MolDraw](https://www.moldraw.com/learn/how-to-use-themes-and-ball-and-stick-style-for-molecules/) — <https://www.moldraw.com/learn/how-to-use-themes-and-ball-and-stick-style-for-molecules/>
 - [How to view a molecule in 3D | Learn MolDraw](https://www.moldraw.com/learn/how-to-view-a-molecule-in-3d/) — <https://www.moldraw.com/learn/how-to-view-a-molecule-in-3d/>
 - [How to view a protein structure from the PDB | Learn MolDraw](https://www.moldraw.com/learn/how-to-view-a-protein-structure-from-the-pdb/) — <https://www.moldraw.com/learn/how-to-view-a-protein-structure-from-the-pdb/>
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Spectroscopy guides (4)

- [Spectroscopy Guides: How to Read Mass, IR and NMR Spectra](https://www.moldraw.com/guides/) — <https://www.moldraw.com/guides/>
 - [How to Read an IR Spectrum: Step-by-Step Guide & Examples](https://www.moldraw.com/guides/how-to-read-ir-spectrum/) — <https://www.moldraw.com/guides/how-to-read-ir-spectrum/>
 - [How to Read a Mass Spectrum: Step-by-Step Guide & Examples](https://www.moldraw.com/guides/how-to-read-mass-spectrum/) — <https://www.moldraw.com/guides/how-to-read-mass-spectrum/>
 - [How to Read NMR Spectra \(\$^1\text{H}\$ & \$^{13}\text{C}\$ \): Shifts, Splitting, DEPT](https://www.moldraw.com/guides/how-to-read-nmr-spectrum/) — <https://www.moldraw.com/guides/how-to-read-nmr-spectrum/>
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Video tutorials (1)

- [MolDraw Tutorials | Video Guides for the Free Chemistry Drawing App](https://www.moldraw.com/tutorials/) — <https://www.moldraw.com/tutorials/>
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Interactive demos (1)

- [Interactive MolDraw demos | Step-by-step, no video](https://www.moldraw.com/demos/) — <https://www.moldraw.com/demos/>
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Blog and how-to articles (41)

- [MolDraw Blog | Chemical Draw, Reaction Drawing and ChemDraw Alternative Guides](https://www.moldraw.com/blog/) — <https://www.moldraw.com/blog/>
- [Predict NMR, Mass and UV-Vis Spectra with AI: Spectroscopy Plugin](https://www.moldraw.com/blog/ai-spectroscopy-plugin-nmr-ms-uv-prediction.html) — <https://www.moldraw.com/blog/ai-spectroscopy-plugin-nmr-ms-uv-prediction.html>
- [Look Up Drug Targets and Bioactivity from Your Drawing: ChEMBL Plugin](https://www.moldraw.com/blog/chembl-plugin-bioactivity-lookup.html) — <https://www.moldraw.com/blog/chembl-plugin-bioactivity-lookup.html>
- [ChemDraw vs MolDraw: Free ChemDraw Alternative \(2026\)](https://www.moldraw.com/blog/chemdraw-vs-moldraw-free-alternative-guide.html) — <https://www.moldraw.com/blog/chemdraw-vs-moldraw-free-alternative-guide.html>
- [Draw Chemical Structures Online | Chemical Structure Drawing App](https://www.moldraw.com/blog/chemical-structure-drawing-app-seo-optimized.html) — <https://www.moldraw.com/blog/chemical-structure-drawing-app-seo-optimized.html>

- [Chemistry Draw Online Guide | Chemical Drawing Workflows](https://www.moldraw.com/blog/chemistry-drawing-keyword-guide.html) — <https://www.moldraw.com/blog/chemistry-drawing-keyword-guide.html>
- [Compare Analogues with Tanimoto Similarity: Compare Molecules Plugin](https://www.moldraw.com/blog/compare-molecules-plugin-tanimoto-similarity.html) — <https://www.moldraw.com/blog/compare-molecules-plugin-tanimoto-similarity.html>
- [Read Cyclic Voltammograms Online: E_{1/2}, ΔE_p, HOMO/LUMO and Scan Rate](https://www.moldraw.com/blog/cyclic-voltammetry-viewer-guide.html) — <https://www.moldraw.com/blog/cyclic-voltammetry-viewer-guide.html>
- [Download MolDraw for Windows \(Free\) | Official Install Guide](https://www.moldraw.com/blog/download-moldraw-windows.html) — <https://www.moldraw.com/blog/download-moldraw-windows.html>
- [Embed MolDraw on Your Site | Free Molecule Editor Iframe](https://www.moldraw.com/blog/embed-moldraw.html) — <https://www.moldraw.com/blog/embed-moldraw.html>
- [Free Chemical Databases: PubChem, ChEMBL & More](https://www.moldraw.com/blog/free-chemical-databases-guide.html) — <https://www.moldraw.com/blog/free-chemical-databases-guide.html>
- [Free Chemical Drawing Software and Chemistry Drawing Program](https://www.moldraw.com/blog/free-chemical-drawing-software-iupac-guide.html) — <https://www.moldraw.com/blog/free-chemical-drawing-software-iupac-guide.html>
- [Free Chemistry Tools and Calculators](https://www.moldraw.com/blog/free-chemistry-tools-and-calculators.html) — <https://www.moldraw.com/blog/free-chemistry-tools-and-calculators.html>
- [How to Add Lone Pairs in MolDraw](https://www.moldraw.com/blog/how-to-add-lone-pair-in-moldraw.html) — <https://www.moldraw.com/blog/how-to-add-lone-pair-in-moldraw.html>
- [How to Turn On the Grid in MolDraw \(Interactive Demo\)](https://www.moldraw.com/blog/how-to-turn-on-the-grid-in-moldraw.html) — <https://www.moldraw.com/blog/how-to-turn-on-the-grid-in-moldraw.html>
- [Read and Compare IR Spectra in Your Browser](https://www.moldraw.com/blog/ir-spectrum-viewer-guide.html) — <https://www.moldraw.com/blog/ir-spectrum-viewer-guide.html>
- [Ketcher Online Molecule Editor Guide | Chemical Structure Drawing](https://www.moldraw.com/blog/ketcher-online-molecule-editor-guide.html) — <https://www.moldraw.com/blog/ketcher-online-molecule-editor-guide.html>
- [chemfig Mechanism Samples: Multi-Step, Curly Arrows and Radical Fish-Hooks](https://www.moldraw.com/blog/chemfig-mechanism-samples.html) — <https://www.moldraw.com/blog/chemfig-mechanism-samples.html>
- [chemfig and MolDraw: Paste chemfig, Edit Visually, Export LaTeX](https://www.moldraw.com/blog/chemfig-latex-moldraw-guide.html) — <https://www.moldraw.com/blog/chemfig-latex-moldraw-guide.html>
- [Draw Lab Glassware and Apparatus Setups: The MolDraw Apparatus Library](https://www.moldraw.com/blog/lab-glassware-apparatus-library-guide.html) — <https://www.moldraw.com/blog/lab-glassware-apparatus-library-guide.html>
- [Check Isotope Patterns, Adducts and ppm Error in Your Browser](https://www.moldraw.com/blog/mass-spectrum-viewer-guide.html) — <https://www.moldraw.com/blog/mass-spectrum-viewer-guide.html>
- [See Where a Molecule Is Electron-Rich: Electrostatic Potential \(MEP\) Plugin](https://www.moldraw.com/blog/mep-plugin-electrostatic-potential-map.html) — <https://www.moldraw.com/blog/mep-plugin-electrostatic-potential-map.html>
- [Complete Guide to All MolDraw Free Chemistry Tools](https://www.moldraw.com/blog/moldraw-all-free-chemistry-tools-guide.html) — <https://www.moldraw.com/blog/moldraw-all-free-chemistry-tools-guide.html>
- [Molecule Structures and Properties](https://www.moldraw.com/blog/molecules/) — <https://www.moldraw.com/blog/molecules/>
- [Aspirin Structure, Formula, Molecular Weight, SMILES & 3D Viewer](https://www.moldraw.com/blog/molecules/aspirin.html) — <https://www.moldraw.com/blog/molecules/aspirin.html>
- [Caffeine Structure, Formula, Molecular Weight, SMILES & 3D Viewer](https://www.moldraw.com/blog/molecules/caffeine.html) — <https://www.moldraw.com/blog/molecules/caffeine.html>
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- [Paracetamol Structure, Formula, Molecular Weight, SMILES & 3D Viewer](https://www.moldraw.com/blog/molecules/paracetamol.html) — <https://www.moldraw.com/blog/molecules/paracetamol.html>
- [Predict ¹H, ¹³C and 2D NMR from Your Drawing: NMR Analyser Plugin](https://www.moldraw.com/blog/nmr-analyser-plugin-predict-nmr-spectra.html) — <https://www.moldraw.com/blog/nmr-analyser-plugin-predict-nmr-spectra.html>
- [View, Integrate and Clean Up NMR Spectra in Your Browser](https://www.moldraw.com/blog/nmr-spectrum-viewer-guide.html) — <https://www.moldraw.com/blog/nmr-spectrum-viewer-guide.html>

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- [View PDB Protein Structures in 3D in Your Browser: Protein Viewer Plugin](https://www.moldraw.com/blog/protein-viewer-plugin-pdb-3d.html) — <https://www.moldraw.com/blog/protein-viewer-plugin-pdb-3d.html>
- [Clean Up and Read Raman Spectra Online: Shift, Fluorescence and Spikes](https://www.moldraw.com/blog/raman-spectrum-viewer-guide.html) — <https://www.moldraw.com/blog/raman-spectrum-viewer-guide.html>
- [Sketch a Molecule by Hand and Get a Clean Structure: Smart Draw Plugin](https://www.moldraw.com/blog/smart-draw-plugin-sketch-to-structure.html) — <https://www.moldraw.com/blog/smart-draw-plugin-sketch-to-structure.html>
- [Find a Molecule's Point Group and See Its Symmetry in 3D: Symmetry Plugin](https://www.moldraw.com/blog/symmetry-plugin-point-groups-3d.html) — <https://www.moldraw.com/blog/symmetry-plugin-point-groups-3d.html>
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- [Make Ball-and-Stick Figures for Slides and Posters: Themes Plugin](https://www.moldraw.com/blog/themes-plugin-ball-and-stick-style.html) — <https://www.moldraw.com/blog/themes-plugin-ball-and-stick-style.html>
- [Read a UV-Vis Spectrum Online: \$\lambda_{\text{max}}\$, Beer-Lambert, Colour and Tauc Band Gap](https://www.moldraw.com/blog/uv-vis-spectrum-viewer-guide.html) — <https://www.moldraw.com/blog/uv-vis-spectrum-viewer-guide.html>
- [XPS Peak Fitting in Your Browser: Charge Correction, Shirley and Atomic %](https://www.moldraw.com/blog/xps-viewer-peak-fitting-guide.html) — <https://www.moldraw.com/blog/xps-viewer-peak-fitting-guide.html>

Free tools and spectrum viewers (87)

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- [Free Cyclic Voltammetry Viewer Online: E1/2, \$\Delta E_p\$, HOMO/LUMO](https://www.moldraw.com/tools/cv-viewer/index.html) — <https://www.moldraw.com/tools/cv-viewer/index.html>
- [MolDraw Tools Hub | Free Chemistry Tools](https://www.moldraw.com/tools/free-chem-tools/) — <https://www.moldraw.com/tools/free-chem-tools/>
- [3D Molecular Model Viewer – Ball and Stick & Space-Filling Models](https://www.moldraw.com/tools/free-chem-tools/3d-molecular-model-viewer.html) — <https://www.moldraw.com/tools/free-chem-tools/3d-molecular-model-viewer.html>
- [AP Chemistry Score Calculator \(2026\) – Estimate Your 1-5](https://www.moldraw.com/tools/free-chem-tools/ap-chemistry-score-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/ap-chemistry-score-calculator.html>
- [Build a Molecule Online – Molecule Builder with 3D & Challenges](https://www.moldraw.com/tools/free-chem-tools/build-a-molecule.html) — <https://www.moldraw.com/tools/free-chem-tools/build-a-molecule.html>
- [CAS Number to SMILES Converter](https://www.moldraw.com/tools/free-chem-tools/cas-number-to-smiles-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/cas-number-to-smiles-converter.html>
- [CAS Number to Structure | Chemical Structure Lookup](https://www.moldraw.com/tools/free-chem-tools/cas-number-to-structure.html) — <https://www.moldraw.com/tools/free-chem-tools/cas-number-to-structure.html>
- [Chemical Equation Balancer - Balance Online Free](https://www.moldraw.com/tools/free-chem-tools/chemical-equation-balancer.html) — <https://www.moldraw.com/tools/free-chem-tools/chemical-equation-balancer.html>
- [Draw a Chemical Reaction Free | Reaction Drawer](https://www.moldraw.com/tools/free-chem-tools/chemical-reaction-drawer.html) — <https://www.moldraw.com/tools/free-chem-tools/chemical-reaction-drawer.html>
- [CID to SMILES Converter](https://www.moldraw.com/tools/free-chem-tools/cid-to-smiles-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/cid-to-smiles-converter.html>
- [Degree of Unsaturation Calculator \(IHD / DBE\) – Rings + \$\pi\$ Bonds](https://www.moldraw.com/tools/free-chem-tools/degree-of-unsaturation-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/degree-of-unsaturation-calculator.html>
- [Dilution Calculator \(C1V1 = C2V2\) - Free](https://www.moldraw.com/tools/free-chem-tools/dilution-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/dilution-calculator.html>
- [Druglikeness Calculator | Lipinski and Veber Checks](https://www.moldraw.com/tools/free-chem-tools/druglikeness-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/druglikeness-calculator.html>

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- [Empirical Formula Calculator - % to Formula](https://www.moldraw.com/tools/free-chem-tools/empirical-formula-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/empirical-formula-calculator.html>
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- [Grams to Moles Calculator | Moles to Grams](https://www.moldraw.com/tools/free-chem-tools/grams-to-moles-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/grams-to-moles-calculator.html>
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- [LogP Calculator | XLogP from SMILES or Name](https://www.moldraw.com/tools/free-chem-tools/logp-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/logp-calculator.html>
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- [Molecular Formula to Structure Search | Formula to Molecule](https://www.moldraw.com/tools/free-chem-tools/molecular-formula-to-structure.html) — <https://www.moldraw.com/tools/free-chem-tools/molecular-formula-to-structure.html>
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- [Protein Molecular Weight Calculator - kDa & ε280](https://www.moldraw.com/tools/free-chem-tools/protein-molecular-weight-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/protein-molecular-weight-calculator.html>
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- [SMILES to 3D Structure Converter - View & Download](https://www.moldraw.com/tools/free-chem-tools/smiles-to-3d-structure-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-3d-structure-converter.html>
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- [SMILES to IUPAC Name Converter - Free & Instant](https://www.moldraw.com/tools/free-chem-tools/smiles-to-iupac-name-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-iupac-name-converter.html>
- [SMILES to MOL Converter](https://www.moldraw.com/tools/free-chem-tools/smiles-to-mol-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-mol-converter.html>
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- [SMILES to PDB Converter](https://www.moldraw.com/tools/free-chem-tools/smiles-to-pdb-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-pdb-converter.html>

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- [SMILES to SDF Converter](https://www.moldraw.com/tools/free-chem-tools/smiles-to-sdf-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-sdf-converter.html>
- [SMILES to Structure Converter Online | Draw 2D Molecules](https://www.moldraw.com/tools/free-chem-tools/smiles-to-structure.html) — <https://www.moldraw.com/tools/free-chem-tools/smiles-to-structure.html>
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- [Structure to IUPAC Name - Draw & Get Name Free](https://www.moldraw.com/tools/free-chem-tools/structure-to-iupac-name.html) — <https://www.moldraw.com/tools/free-chem-tools/structure-to-iupac-name.html>
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- [VSEPR Chart & Molecular Geometry Calculator \(3D\)](https://www.moldraw.com/tools/free-chem-tools/vsepr-molecular-geometry-calculator.html) — <https://www.moldraw.com/tools/free-chem-tools/vsepr-molecular-geometry-calculator.html>
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- [XYZ to SMILES Converter | Coordinate File to SMILES](https://www.moldraw.com/tools/free-chem-tools/xyz-to-smiles-converter.html) — <https://www.moldraw.com/tools/free-chem-tools/xyz-to-smiles-converter.html>
- [Free IR Spectrum Viewer Online: Open JCAMP-DX, CSV and SPC Files](https://www.moldraw.com/tools/ir-viewer/index.html) — <https://www.moldraw.com/tools/ir-viewer/index.html>
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- [Free Raman Spectrum Viewer Online: Baseline, Peaks, Bands](https://www.moldraw.com/tools/raman-viewer/index.html) — <https://www.moldraw.com/tools/raman-viewer/index.html>
- [Free TGA and DSC Viewer Online: DTG, Onset, T_g, ΔH, Crystallinity](https://www.moldraw.com/tools/tga-dsc-viewer/index.html) — <https://www.moldraw.com/tools/tga-dsc-viewer/index.html>
- [Free UV-Vis Spectrum Viewer: λ_{max}, Beer-Lambert, Colour and Tauc Band Gap](https://www.moldraw.com/tools/uv-vis-viewer/index.html) — <https://www.moldraw.com/tools/uv-vis-viewer/index.html>
- [Free XPS Viewer and Peak Fitting Online: Shirley, Doublets, Atomic %](https://www.moldraw.com/tools/xps-viewer/index.html) — <https://www.moldraw.com/tools/xps-viewer/index.html>

Named reactions (8)

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 - [Grignard Reaction Guide](https://www.moldraw.com/reaction/grignard-reaction) — <https://www.moldraw.com/reaction/grignard-reaction>
 - [Heck Reaction Guide](https://www.moldraw.com/reaction/heck-reaction) — <https://www.moldraw.com/reaction/heck-reaction>
 - [Suzuki Coupling Reaction Guide](https://www.moldraw.com/reaction/suzuki-coupling) — <https://www.moldraw.com/reaction/suzuki-coupling>
 - [Wittig Reaction Guide](https://www.moldraw.com/reaction/wittig-reaction) — <https://www.moldraw.com/reaction/wittig-reaction>
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Molecule pages (30)

- [3D Molecule Models – 29 Interactive Ball-and-Stick Molecular Models](https://www.moldraw.com/molecules/) — <https://www.moldraw.com/molecules/>
- [Acetone Molecule Model – Interactive 3D Structure \(C₃H₆O\)](https://www.moldraw.com/molecules/acetone.html) — <https://www.moldraw.com/molecules/acetone.html>
- [Adrenaline Molecule Model – Interactive 3D Structure \(C₉H₁₃NO₃\)](https://www.moldraw.com/molecules/adrenaline.html) — <https://www.moldraw.com/molecules/adrenaline.html>
- [Ammonia Molecule Model – Interactive 3D Structure \(NH₃\)](https://www.moldraw.com/molecules/ammonia.html) — <https://www.moldraw.com/molecules/ammonia.html>
- [Aspirin Molecule Model – Interactive 3D Structure \(C₉H₈O₄\)](https://www.moldraw.com/molecules/aspirin.html) — <https://www.moldraw.com/molecules/aspirin.html>
- [ATP Molecule Model – Interactive 3D Structure \(C₁₀H₁₆N₅O₁₃P₃\)](https://www.moldraw.com/molecules/atp.html) — <https://www.moldraw.com/molecules/atp.html>
- [Benzene Molecule Model – Interactive 3D Structure \(C₆H₆\)](https://www.moldraw.com/molecules/benzene.html) — <https://www.moldraw.com/molecules/benzene.html>
- [Caffeine Molecule Model – Interactive 3D Structure \(C₈H₁₀N₄O₂\)](https://www.moldraw.com/molecules/caffeine.html) — <https://www.moldraw.com/molecules/caffeine.html>
- [Carbon dioxide Molecule Model – Interactive 3D Structure \(CO₂\)](https://www.moldraw.com/molecules/carbon-dioxide.html) — <https://www.moldraw.com/molecules/carbon-dioxide.html>
- [Cholesterol Molecule Model – Interactive 3D Structure \(C₂₇H₄₆O\)](https://www.moldraw.com/molecules/cholesterol.html) — <https://www.moldraw.com/molecules/cholesterol.html>
- [Citric acid Molecule Model – Interactive 3D Structure \(C₆H₈O₇\)](https://www.moldraw.com/molecules/citric-acid.html) — <https://www.moldraw.com/molecules/citric-acid.html>
- [Estradiol Molecule Model – Interactive 3D Structure \(C₁₈H₂₄O₂\)](https://www.moldraw.com/molecules/estradiol.html) — <https://www.moldraw.com/molecules/estradiol.html>
- [Ethanol Molecule Model – Interactive 3D Structure \(C₂H₆O\)](https://www.moldraw.com/molecules/ethanol.html) — <https://www.moldraw.com/molecules/ethanol.html>
- [Fructose Molecule Model – Interactive 3D Structure \(C₆H₁₂O₆\)](https://www.moldraw.com/molecules/fructose.html) — <https://www.moldraw.com/molecules/fructose.html>
- [Glucose Molecule Model – Interactive 3D Structure \(C₆H₁₂O₆\)](https://www.moldraw.com/molecules/glucose.html) — <https://www.moldraw.com/molecules/glucose.html>
- [Glycerol Molecule Model – Interactive 3D Structure \(C₃H₈O₃\)](https://www.moldraw.com/molecules/glycerol.html) — <https://www.moldraw.com/molecules/glycerol.html>
- [Glycine Molecule Model – Interactive 3D Structure \(C₂H₅NO₂\)](https://www.moldraw.com/molecules/glycine.html) — <https://www.moldraw.com/molecules/glycine.html>
- [Hydrogen chloride Molecule Model – Interactive 3D Structure \(HCl\)](https://www.moldraw.com/molecules/hydrogen-chloride.html) — <https://www.moldraw.com/molecules/hydrogen-chloride.html>
- [Melatonin Molecule Model – Interactive 3D Structure \(C₁₃H₁₆N₂O₂\)](https://www.moldraw.com/molecules/melatonin.html) — <https://www.moldraw.com/molecules/melatonin.html>
- [Methane Molecule Model – Interactive 3D Structure \(CH₄\)](https://www.moldraw.com/molecules/methane.html) — <https://www.moldraw.com/molecules/methane.html>
- [Methanol Molecule Model – Interactive 3D Structure \(CH₄O\)](https://www.moldraw.com/molecules/methanol.html) — <https://www.moldraw.com/molecules/methanol.html>

- [Oxygen Molecule Model - Interactive 3D Structure \(O2\)](https://www.moldraw.com/molecules/oxygen.html) — <https://www.moldraw.com/molecules/oxygen.html>
 - [Ozone Molecule Model - Interactive 3D Structure \(O3\)](https://www.moldraw.com/molecules/ozone.html) — <https://www.moldraw.com/molecules/ozone.html>
 - [Propane Molecule Model - Interactive 3D Structure \(C3H8\)](https://www.moldraw.com/molecules/propane.html) — <https://www.moldraw.com/molecules/propane.html>
 - [Serotonin Molecule Model - Interactive 3D Structure \(C10H12N2O\)](https://www.moldraw.com/molecules/serotonin.html) — <https://www.moldraw.com/molecules/serotonin.html>
 - [Sucrose Molecule Model - Interactive 3D Structure \(C12H22O11\)](https://www.moldraw.com/molecules/sucrose.html) — <https://www.moldraw.com/molecules/sucrose.html>
 - [Testosterone Molecule Model - Interactive 3D Structure \(C19H28O2\)](https://www.moldraw.com/molecules/testosterone.html) — <https://www.moldraw.com/molecules/testosterone.html>
 - [Vanillin Molecule Model - Interactive 3D Structure \(C8H8O3\)](https://www.moldraw.com/molecules/vanillin.html) — <https://www.moldraw.com/molecules/vanillin.html>
 - [Vitamin C Molecule Model - Interactive 3D Structure \(C6H8O6\)](https://www.moldraw.com/molecules/vitamin-c.html) — <https://www.moldraw.com/molecules/vitamin-c.html>
 - [Water Molecule Model - Interactive 3D Structure \(H2O\)](https://www.moldraw.com/molecules/water.html) — <https://www.moldraw.com/molecules/water.html>
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About, help and policies (12)

- [About MolDraw - Free Chemical Structure Drawing App](https://www.moldraw.com/pages/about.html) — <https://www.moldraw.com/pages/about.html>
 - [How to Use AI Assistant](https://www.moldraw.com/pages/ai-help.html) — <https://www.moldraw.com/pages/ai-help.html>
 - [Authors and Team](https://www.moldraw.com/pages/authors.html) — <https://www.moldraw.com/pages/authors.html>
 - [Community Guidelines](https://www.moldraw.com/pages/community-guidelines.html) — <https://www.moldraw.com/pages/community-guidelines.html>
 - [Contact](https://www.moldraw.com/pages/contact.html) — <https://www.moldraw.com/pages/contact.html>
 - [Editorial Guidelines](https://www.moldraw.com/pages/editorial-guidelines.html) — <https://www.moldraw.com/pages/editorial-guidelines.html>
 - [MolDraw FAQ - Frequently Asked Questions | Free ChemDraw Alternative](https://www.moldraw.com/pages/faq.html) — <https://www.moldraw.com/pages/faq.html>
 - [MolDraw MCP privacy and security](https://www.moldraw.com/pages/mcp-privacy.html) — <https://www.moldraw.com/pages/mcp-privacy.html>
 - [Privacy Policy](https://www.moldraw.com/pages/privacy-policy.html) — <https://www.moldraw.com/pages/privacy-policy.html>
 - [HTML Sitemap](https://www.moldraw.com/pages/sitemap.html) — <https://www.moldraw.com/pages/sitemap.html>
 - [Terms of Service](https://www.moldraw.com/pages/terms-of-service.html) — <https://www.moldraw.com/pages/terms-of-service.html>
 - [Updates](https://www.moldraw.com/pages/updates.html) — <https://www.moldraw.com/pages/updates.html>
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Other pages (6)

- [MolDraw Addons | PowerPoint and Word \(Coming Soon\)](https://www.moldraw.com/addons) — <https://www.moldraw.com/addons>
- [MolDraw Help Centre: Support, Guides and Knowledge Base](https://www.moldraw.com/support/) — <https://www.moldraw.com/support/>
- [ChemSketch Alternative: Free Online Chemical Structure Editor](https://www.moldraw.com/chemsketch) — <https://www.moldraw.com/chemsketch>
- [ChemSketch Download: Try MolDraw in Your Browser](https://www.moldraw.com/chemsketch-download) — <https://www.moldraw.com/chemsketch-download>
- [ChemSketch Free & ACD ChemSketch Freeware Alternative](https://www.moldraw.com/chemsketch-free) — <https://www.moldraw.com/chemsketch-free>
- [MolDraw Chrome Addon | Snip Any Molecule Image to an Editable Structure \(Coming Soon\)](https://www.moldraw.com/chrome-addon) — <https://www.moldraw.com/chrome-addon>