

Resource Summary Report

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VMD

RRID:SCR_024368

Type: Tool

Proper Citation

VMD (RRID:SCR_024368)

Resource Information

URL: <http://www.ks.uiuc.edu/Research/vmd/>

Proper Citation: VMD (RRID:SCR_024368)

Description: Software tool as molecular visualization program for displaying, animating, and analyzing large biomolecular systems using 3-D graphics and built-in scripting. VMD supports computers running MacOS X, Unix, or Windows, is distributed free of charge, and includes source code.

Synonyms: vmd

Resource Type: software resource, software application

Keywords: molecular visualization, displaying, animating, analyzing, large biomolecular systems, .

Availability: Free, Available for download, Freely available,

Resource Name: VMD

Resource ID: SCR_024368

Old URLs: <https://sources.debian.org/src/vmd/>

Record Creation Time: 20230830T050217+0000

Record Last Update: 20260801T191121+0000

Ratings and Alerts

No rating or validation information has been found for VMD.

No alerts have been found for VMD.

Data and Source Information

Source: [SciCrunch Registry](#)

Usage and Citation Metrics

We found 33 mentions in open access literature.

Listed below are recent publications. The full list is available at [ASWG](#).

Camargo PG, et al. (2025) In silico evaluation of N-aryl-1,10-phenanthroline-2-amines as potential inhibitors of *T. cruzi* GP63 zinc-metalloprotease by docking and molecular dynamics simulations. *Scientific reports*, 15(1), 6036.

Akulov V, et al. (2025) Phosphorylation-Regulated Conformational Diversity and Topological Dynamics of an Intrinsically Disordered Nuclear Receptor. *The journal of physical chemistry. B*, 129(30), 7719.

Garcia Jimenez D, et al. (2025) Linker Methylation as a Strategy to Enhance PROTAC Oral Bioavailability: Insights from Molecular Properties and Conformational Analysis. *Journal of medicinal chemistry*, 68(15), 16666.

Camargo PG, et al. (2025) In vitro assays identified thiohydantoins with anti-trypanosomatid activity and molecular modelling studies indicated possible selective CYP51 inhibition. *Scientific reports*, 15(1), 465.

Singh AK, et al. (2025) Conformational Landscape of the Di- and Tripeptide Permease A Transport Cycle. *Journal of chemical information and modeling*, 65(12), 6198.

Ghavi Z, et al. (2025) Molecular dynamics insights into non-covalent functionalization of boron nitride nanotubes for doxorubicin delivery. *Scientific reports*, 15(1), 30213.

Motuziuk O, et al. (2025) The effect of C60 fullerene on the musculus gastrocnemius contraction in chronically alcohol-exposed rats and the potential mechanism of its action. *Scientific reports*, 15(1), 32651.

Zan B, et al. (2025) The difference between Melp5 and melittin membrane poration. *Scientific reports*, 15(1), 7442.

Mohebbi A, et al. (2025) A new method based on binary mixture concept for prediction of ionic liquids critical properties using molecular dynamics simulation. *Scientific reports*, 15(1), 7545.

Yu S, et al. (2025) c-di-GMP inhibits rRNA methylation and impairs ribosome assembly in the presence of kanamycin. *EMBO reports*, 26(5), 1367.

Konečná K, et al. (2025) Functional characterization of 16 variants found in the LDL receptor gene. *Journal of lipid research*, 66(9), 100873.

Doran MH, et al. (2025) The hypertrophic cardiomyopathy-associated A331P actin variant enhances basal contractile activity and elicits resting muscle dysfunction. *iScience*, 28(2),

111816.

Dalal K, et al. (2025) Interpreting regulatory mechanisms of Hippo signaling through a deep learning sequence model. *Cell genomics*, 5(4), 100821.

Tang J, et al. (2025) Targeting N-Myc in neuroblastoma with selective Aurora kinase A degraders. *Cell chemical biology*, 32(2), 352.

Schumann J, et al. (2024) Ten-electron count rule for the binding of adsorbates on single-atom alloy catalysts. *Nature chemistry*, 16(5), 749.

Camargo PG, et al. (2024) Py-CoMFA, docking, and molecular dynamics simulations of *Leishmania (L.) amazonensis* arginase inhibitors. *Scientific reports*, 14(1), 11575.

Tourkova IL, et al. (2024) Chloride/proton antiporters ClC3 and ClC5 support bone formation in mice. *Bone reports*, 21, 101763.

Hossain A, et al. (2024) Role of cold shock proteins B and D in *Aeromonas salmonicida* subsp. *salmonicida* physiology and virulence in lumpfish (*Cyclopterus lumpus*). *Infection and immunity*, 92(8), e0001124.

Bakhtiari MA, et al. (2024) Investigation the behavior of different fullerenes on graphene surface. *Scientific reports*, 14(1), 18220.

Nelic D, et al. (2024) Agonist-selective activation of individual G-proteins by muscarinic receptors. *Scientific reports*, 14(1), 9652.