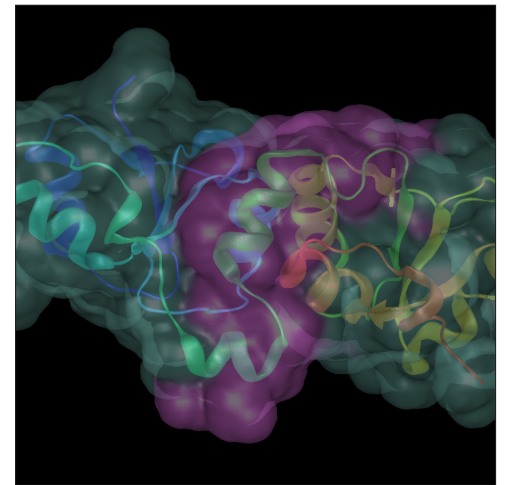
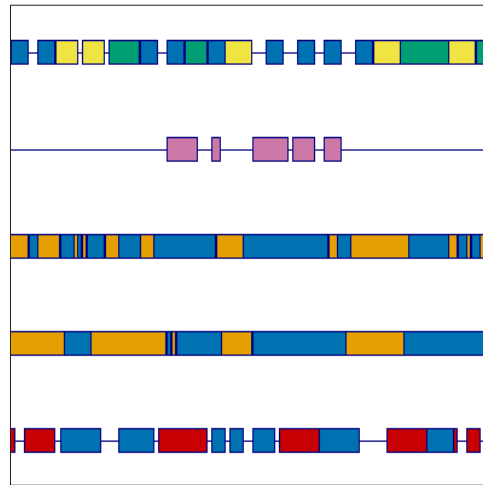
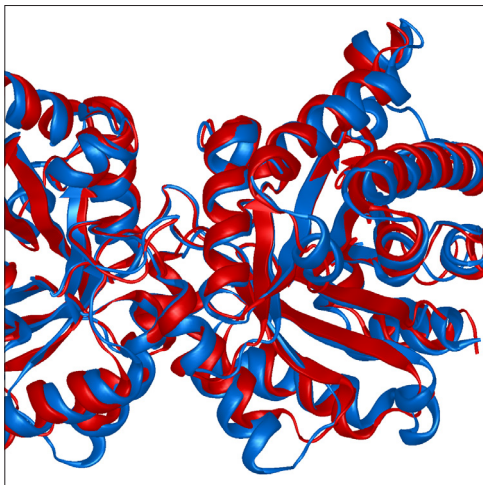


LASERGENE PROTEIN AND NOVA APPLICATIONS

Software for protein modeling and analysis



PROTEIN SEQUENCE ANALYSIS

- Utilize integrated views and analysis methods for sequence, secondary structure, and tertiary structure
- Predict secondary structure characteristics

PROTEIN STRUCTURE ANALYSIS

- Predict B-cell epitopes
- Create molecular and solvent accessible surfaces to visualize predicted epitopes
- Align entire structures or selected regions
- Create publication quality graphics
- Visualize conformational changes of nearly 400 animated macromolecular structures

PROTEIN MODELING

- Predict 3D structure for any protein sequence using the renowned **AlphaFold 2**, **AlphaFold Multimer**, or **Boltz-2** algorithms
- Model antibody structures and identify antibody/antigen binding sites
- Predict protein function, ligand binding sites, and enzyme activity
- Model docking for any receptor and ligand pair
- Predict binding interactions and energy
- Create and model variants on protein structures
- Perform hot-spot scans and improve fold stability with protein design tools

Comprehensive tools for protein modeling

Structure Prediction with NovaFold AI Boltz

- NovaFold AI Boltz utilizes the Boltz-2 protein structure prediction algorithm, an open-source analogue of AlphaFold 3. NovaFold AI Boltz delivers fast, accurate, and cost-effective results.
- NovaFold AI uses the AlphaFold 2 algorithm from DeepMind to predict distance and create dihedral maps using deep multiple sequence alignments as input.
- NovaFold AI-Multimer uses the AlphaFold-Multimer algorithm to predict the structure of a multimeric protein assembly.

Protein-Protein Docking with NovaDock

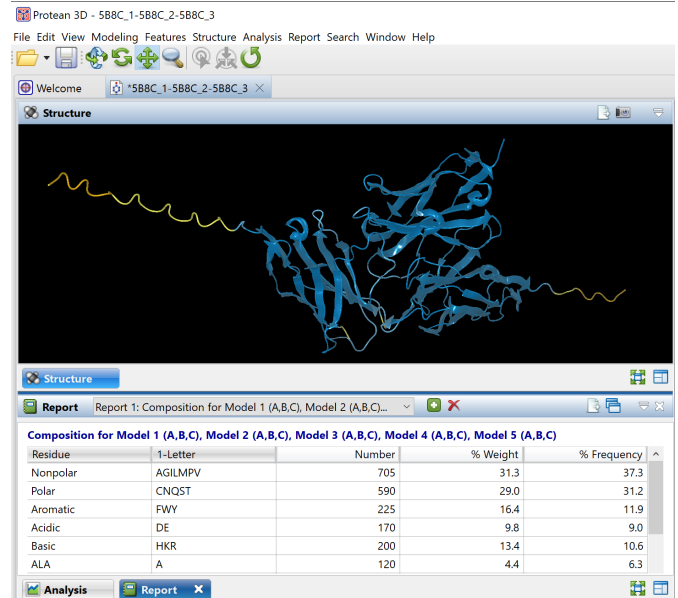
- Based on SwarmDock, a high-resolution docking algorithm
- Model protein docking and binding interactions
- Explore protein flexibility during docking

Antibody Modeling with NovaFold Antibody

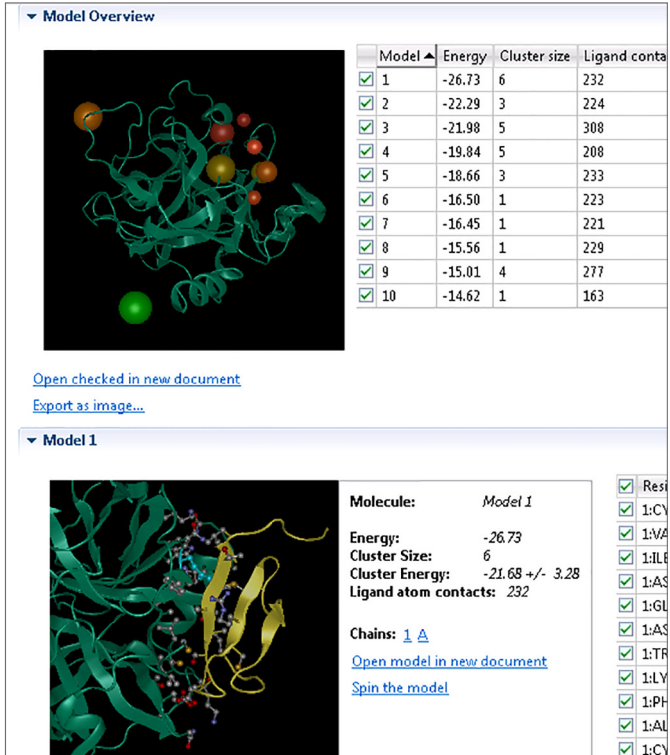
- Model Fv, Fab, VH, sdAb in minutes
- Search a library of antibody frameworks, or provide custom templates
- *Ab initio* loop modeling for H3 - up to 15 residues
- Automated annotation of CDR loops

Advanced Protein Design

- Create, model, and analyze variants on structure
- Calculate energy changes caused by mutations
- Perform serine and alanine variant scans
- Improve protein fold stability



The structure view and the composition report for a NovaFold AI-Multimer prediction. Many other customizable views and reports are available.



The screenshot shows the NovaDock report. The 'Model Overview' section displays a table of docking models. The 'Model 1' section shows a detailed view of the top docking model, including a 3D visualization of the protein and ligand, and a table of properties.

Model	Energy	Cluster size	Ligand conta
1	-26.73	6	232
2	-22.29	3	224
3	-21.98	5	308
4	-19.84	5	208
5	-18.66	3	233
6	-16.50	1	223
7	-16.45	1	221
8	-15.56	1	229
9	-15.01	4	277
10	-14.62	1	163

Model 1

Molecule: Model 1

Energy: -26.73

Cluster Size: 6

Cluster Energy: -21.68 +/- 3.28

Ligand atom contacts: 232

Chains: 1 A

Open model in new document

Spin the model

The NovaDock report showing the top ligand-receptor docking models for a completed prediction.



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