

# CS612 - Algorithms in Bioinformatics

Visualization and Representation

February 19, 2019

# The PDB File Format

		Amino Acid		Chain name		Sequence Number		-----Coordinates-----		
		Element						X	Y	Z (etc.)
ATOM	1	N	ASP	L	1			4.060	7.307	5.186 ...
ATOM	2	CA	ASP	L	1			4.042	7.776	6.553 ...
ATOM	3	C	ASP	L	1			2.668	8.426	6.644 ...
ATOM	4	O	ASP	L	1			1.987	8.438	5.606 ...
ATOM	5	CB	ASP	L	1			5.090	8.827	6.797 ...
ATOM	6	CG	ASP	L	1			6.338	8.761	5.929 ...
ATOM	7	OD1	ASP	L	1			6.576	9.758	5.241 ...
ATOM	8	OD2	ASP	L	1			7.065	7.759	5.948 ...

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Element position within amino acid

# Modeling and Visualization of Protein Structures

- Computer graphics, scientific visualization and geometry to create a 3-D visual model of molecular structures.
- Facilitates structure, dynamic and function analysis.
- The PDB represents molecular structures as a set of Cartesian coordinates in 3-D.
- Internal coordinates – bond lengths, angles and dihedrals (more about them later...).
- Possible to switch back and forth.

# Cartesian Coordinates

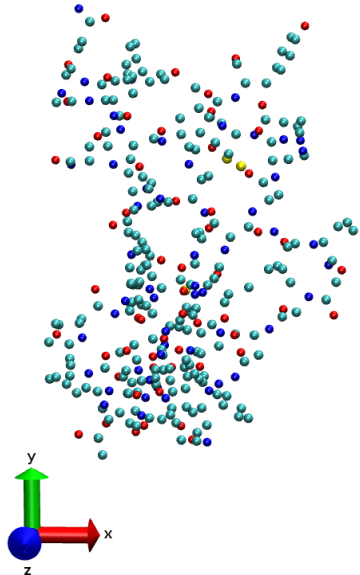
- Representing each atom as a set of  $(x, y, z)$  coordinates
- Easy and convenient to render on screen.
- Every line in a PDB file represents the 3-D location of the atom center with respect to some (arbitrary) axis system.
- Using a graphics package, simply draw a dot for each atom.

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# Visualization of Protein Structures

- Numerous tools are available for visualizing the structures stored in the PDB and other repositories.
- Most such tools allow a detailed examination of the molecule in a variety of rendering modes.
- For example, sometimes it may be useful to have a detailed image of the surface of the molecule as experienced by a molecule of water.
- For other purposes, a simple, cartoonish representation of the major structural features may be sufficient.

- “Full featured” academic packages
  - UCSF Chimera (<http://chimera.ucsf.edu/>)
  - PyMOL (<http://pymol.sourceforge.net/>)
  - VMD (<http://www.ks.uiuc.edu/Research/vmd/>)
- Viewers
  - RasMol/Chime (<http://www.openrasmol.org/>)
  - Jmol (<http://jmol.sourceforge.net/>)
  - SwissProt PDB-Viewer (DeepView) (<http://www.expasy.org/spdbv/>)
  - RCSB Protein Workshop (<http://www.rcsb.org/>)

# Modeling Tools

- Amber (<http://amber.scripps.edu/>)
- Charmm (<http://www.charmm.org/>)
- NAMD (<http://www.ks.uiuc.edu/Research/namd/>)
- Gaussian (<http://www.gaussian.com/>)
- ModBase (<http://modbase.compbio.ucsf.edu/>)
- Modeller (<http://www.salilab.org/modeller/>)
- DOCK (<http://dock.compbio.ucsf.edu/>)
- Many, many more (see <http://cmm.info.nih.gov/>)



# Comparison of Visualization Packages

- Jmol
  - Best-in-class viewer
  - Web enabled
  - Scriptable
  - Input only
  - Compatible with RasMol scripts
  - Limited analytical capabilities
  - Mostly through Javascript wrappers

# Comparison of Visualization Packages

- PyMol
  - Best-in-class for peptidometrics, speed
  - Single-screen interface (+command line)
  - Extensible
  - Some modeling capabilities
  - Good publication tools (built-in ray tracer)
  - Scriptable

# Comparison of Visualization Packages

- VMD
  - Best-in-class for MD (integrated with NAMD) and other analysis tools
  - Scriptable
  - Excellent stereo capabilities
  - Embedded ray tracer (Tachyon)
  - Extensible
  - Now supports Python, previously was only TCL/TK

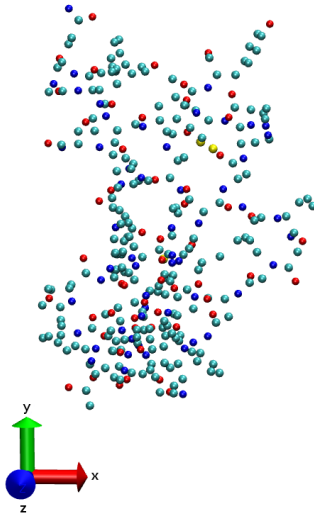
# Comparison of Visualization Packages

- Chimera
  - Best-in-class for visualizing very large structures
  - Multiscale extension, volume viewer
  - Focus on extensibility, broad functionality
  - Primarily analytical interface
  - Familiar GUI interface (+command line)
  - Scriptable
  - Reasonable tools for publication & presentation
  - Embedded ray tracer (POV-Ray)
  - Excellent sequence/structure capabilities
  - Reasonable interface to modeling programs

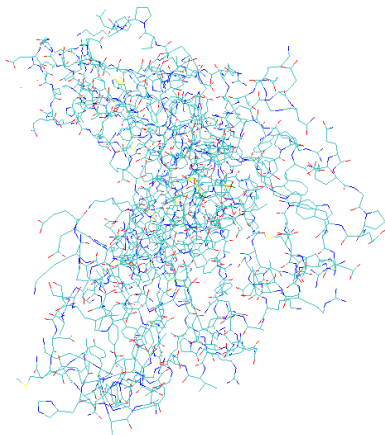
# Comparison of Visualization Packages

- There is no “best” package for everything (IMHO)
- YMMV (Your Mileage May Vary)
  - What we think is easy, you may think is hard
  - What we think is hard, you may think is easy
- Choosing the best package for you
  - Does what you need
  - Good documentation
  - Good support (either local or from the authors)

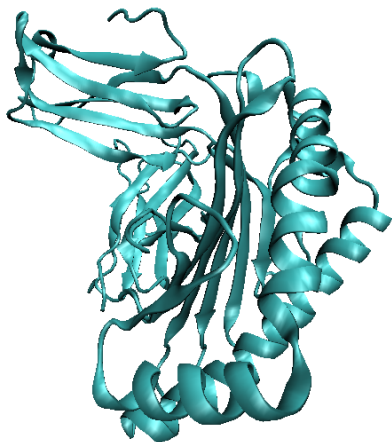
# Representation of Protein Structures – Set of Dots



# Representation of Protein Structures – Wireframe

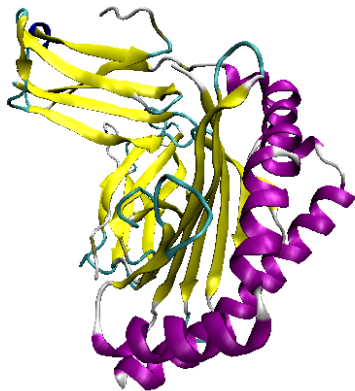


# Representation of Protein Structures – Cartoon

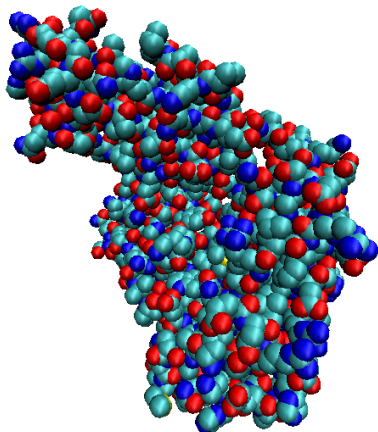




# Representation of Protein Structures – Color by Secondary Structure



# Representation of Protein Structures – Spheres



# Representation of Protein Structures – Solvent Accessible Surface, Colored by Chain

