

# Protein Structure Viewers

## Hot Topics – March 2009

### Swiss-PdbViewer demo

1. Open (or import) cholera toxin (1xtc) PDB file.
2. Browse main interface
  - Rotate, translate zoom
  - Display → Render in Solid 3D
  - Wind (windows) → Control Panel, Alignment
3. Select by chain or atoms, and/or using Select menu
4. Browse Control Panel
  - Chain, secondary structure, group, show (ball and stick), side (chain), label, surface (::v), ribbon, color, backbone/side chain/etc.
5. Preferences (Prefs) menu
6. Change 5 chains of ring (DEFGH) into ribbons
  - Prefs → colors
7. Change chain C to orange space-filled
8. Make molecular surface of chains AC
  - Prefs → Surfaces
    - Check "Ignore Selected Residues"
    - Transparency (1 - 100) → 90
  - Tools → Compute Molecular Surface
9. Select active site residues and show as red space-filled atoms
10. Save layer (File →)
11. Save Pov-Ray Scene (File →)
12. Open (or import) E. coli enterotoxin (1tii) PDB file.
13. Browse new display options:
  - Control Panel → pull-down menu for structure selection
  - Visible? Can move?
14. Fit → Iterative Magic Fit (and then separate)
15. Select active residues and change to space-filling
  - Look at Alignment
10. Save Project (File →)
11. Save Pov-Ray Scene (File →)
12. Go to POV-Ray (or open \*.pov file) and Run, generating image of desired size.