Subject: Re: Molecular Modelling really possible with pv-wave?
Posted by hewat on Fri, 09 Feb 1996 08:00:00 GMT
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Bernd Ihmels <bihmels1@gwdg.de> writes:

> So I wonder, if anyone of you ever wrote a program do draw large molecules,
> with a fast algorithm, so that you won't have to wait several minutes for just
> one image? Or even better, a program, able to turn a molecule ones created, so
> that you don't have to create a new image for every new view-angle?
>
> Or is something like that impossible to do with pv-wave?

Why would you want to do that with IDL :-) Take a look at VRML based applications-
http://ws05.pc.chemie.th-darmstadt.de/vrml/pdb2vrml.html and then at RasMol
http://www.umass.edu/microbio/rasmol to see the competition you are up against :-)

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