

<b>Course name:</b> Introduction to structural bioinformatics	<b>Credits:</b> 5 ECTS
<b>Class type:</b> On-line lectures + individual practice	<b>Hours per week:</b>
<b>Type of the exam:</b> Project work	
<b>Prerequisites</b> (if exist): basic biochemistry and molecular biology. Ability of downloading and installing the free UCSF Chimera software	
<p><b>Course description:</b> Basics of search, manipulation and analyses of structures of large biological molecules, especially proteins.</p> <ul style="list-style-type: none"> <li>• Basics of protein structures and structure determination. Simple validation of models by Ramachandran plots. Basic use of molecular graphics software</li> <li>• Molecular graphics: illustrating and highlighting molecular details on screen and print; generating molecular surfaces</li> <li>• Comparison of structures: overlaying molecules and measuring their structural similarity</li> <li>• Molecular animations</li> <li>• Theory of protein modeling and protein dynamics</li> <li>• Validation and analysis of models</li> </ul>	
<b>Required reading:</b>	
<b>Recommended reading:</b> Gu, J & Bourne, P.E. (eds, 2009): Structural bioinformatics, 2nd ed. Wiley-Blackwell	
<b>Lecturer</b> ( <i>name, position, degree</i> ): Zoltán Gáspári, associate professor,, PhD.	
<b>Additional lecturers</b> , if exist:	