UnityMoI: Visualization with a game engine

Z. Lv^{1,2}, A. Tek², M. Chavent³ and M. Baaden²

¹Ocean University of China, QingDao, China.²Institut de Biologie Physico-Chimique, Laboratoire de Biochimie Théorique, CNRS UPR 9080, 13, rue Pierre et Marie Curie, F-75005 Paris, France. ²CEA, DAM, DIF, 91297 Arpaion, France.



Introduction

In every biology fields technological progress continously increases the amount of experimental and theoretical data. Hence, building visualization application to retrieve relevant information is more and more complex and requires higher and higher programming skills. This is not achievable by most of non-computer science teams, which then struggle to analyze new data.

In the other hand, the video game market is full of high-quality and visually appealing games playable by everyone and created by small independant companies with low resources. The needs in terms of graphics and interactivity are often comparable to science visualization application.

Can the scientific community use game development methods and tools to ease the creation and diffusion of visual applications?



We chose the Unity3D game engine (<u>http://unity3d.com/</u>) as it proposes a free license and many useful features :

- ✓Multi-platform (Windows, MacOS, web, mobile devices) applications with minimal programming effort
- ✓ User-friendly interface for creation of 3D application using object oriented concepts
- ✓ Well featured and documented API available in C#, JavaScript and Boo/Python including physics and 3D interaction routines
- ✓ Shader Lab to create and debug shader programs
- ✓ Active and helpful web community

Unity



We created UnityMol, a molecular viewer prototype, using the basics features of Unity and see how it is reliable to build science applications



• References: [1] M. Chavent, A. Vanel, A. Tek, B. Levy, B. Raffin, S. Robert, M. Baaden, GPU-accelerated atom and dynamic bond visualization using HyperBalls, a unified algorithm for balls, sticks and hyperboloids, accepted in Journal of Computational Chemistry.