

Molecular Modeling in Process Engineering

2023/2024

Project 8

Preliminary steps

Download and install the software AIMAll.

The program needs the wfn file from QM calculations (see manual).

Electron Density Analysis

- 1) Calculate at the B3LYP/6-31G** level the electron density of the following systems: ethane, ethene, ethyne, and benzene (optimize the geometry of the systems)

By AIMAll search for:

- Critical Points
- Molecular graphs
- BCPs sono quelli verdi
- Density at the BCPs
- Order of the C-C bonds
- Ellipticity of the bonds

- 2) Calculate at the B3LYP/6-31G** level the electron density of the systems: water and water dimer (optimize the geometry of the systems)

By AIMAll search for:

- Critical Points
- Molecular graphs
- BCPs
- Density at the BCPs

- 3) Calculate at the B3LYP/6-31G** level the electron density of the systems: H₂, HF and LiF (optimize the geometry of the systems)

By AIMAll search for:

- Critical Points
- Molecular graphs
- BCPs
- Density at the BCPs
- Laplacian of the density at the BCPs
- Plot the L=-Laplacian of the density