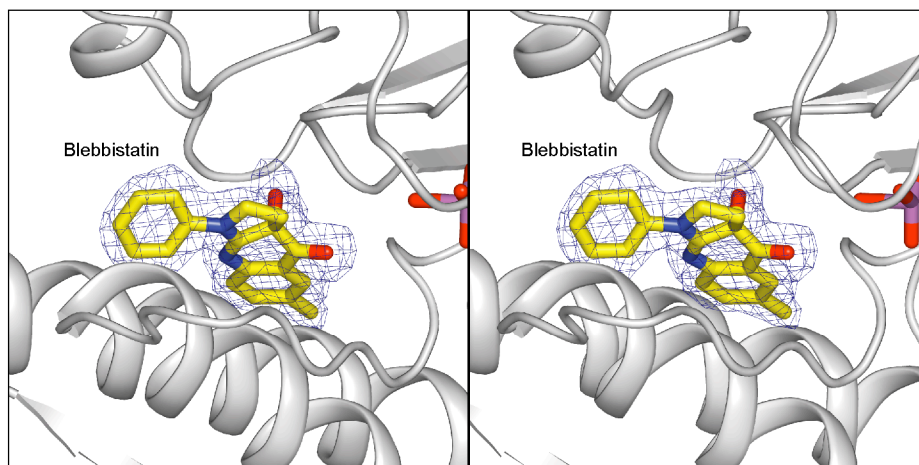


### Supplementary Figure 1. (Rayment)



**Supplementary Figure 1.** Stereo view of the electron density for blebbistatin. The map was calculated with coefficients of the form  $F_o - F_c$  where the inhibitor was omitted from the refinement and phase calculation. Electron density was contoured at  $3 \sigma$ .