

Supporting Information

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Synthesis and molecular docking study of ethyl piperidine-1-carboxylate derivative Schiff bases

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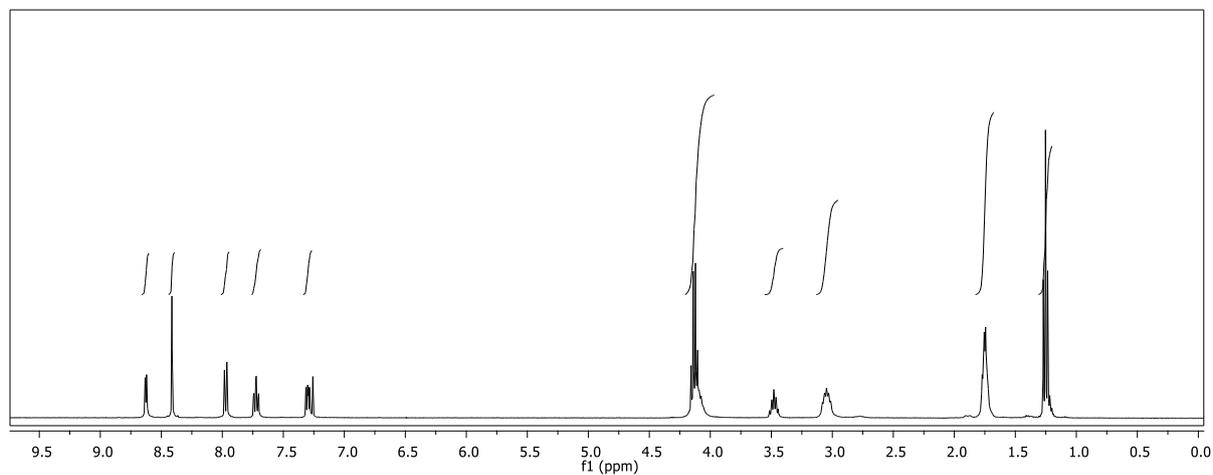
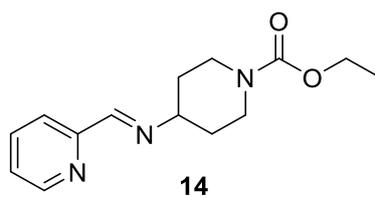


Figure S1: 400 MHz ^1H NMR spectrum of Ethyl (*E*)-4-((pyridin-2-ylimino)methyl)piperidine-1-carboxylate (**14**) in CDCl_3

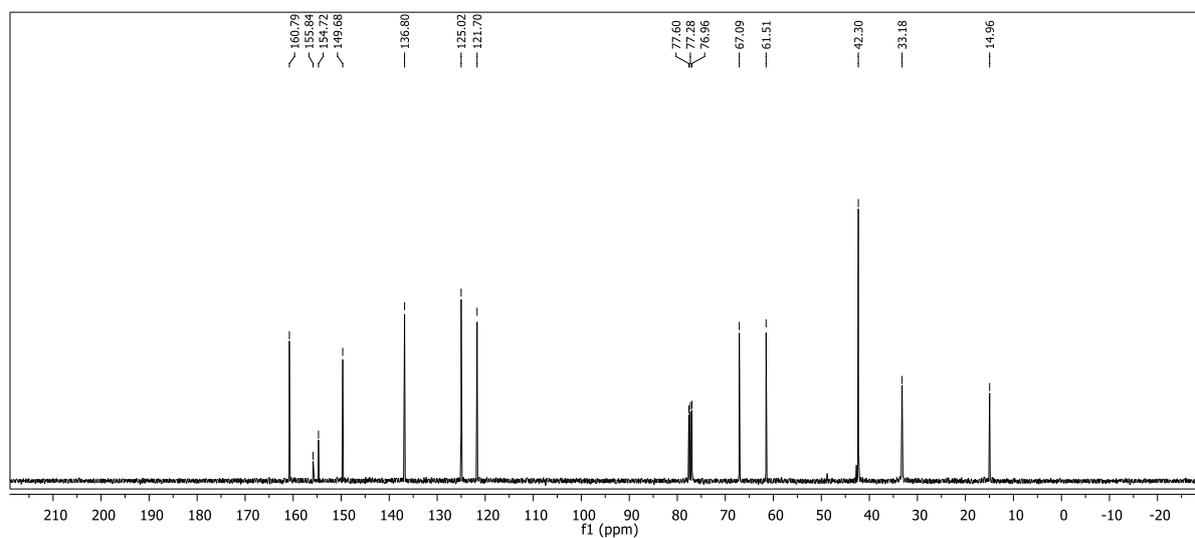


Figure S2: 100 MHz ^{13}C -NMR spectrum of Ethyl (*E*)-4-((pyridin-2-ylimino)methyl)piperidine-1-carboxylate (**14**) in CDCl_3

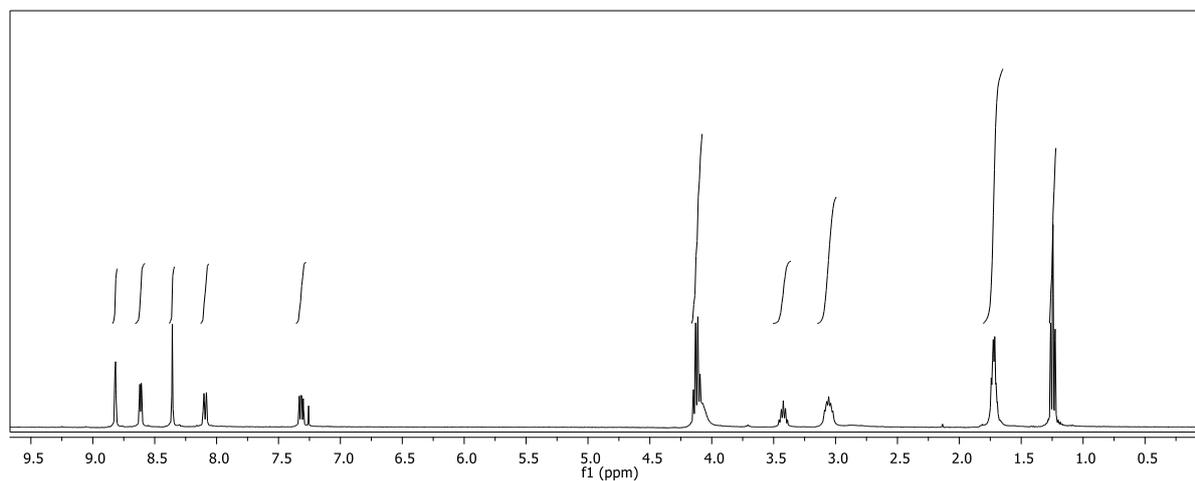
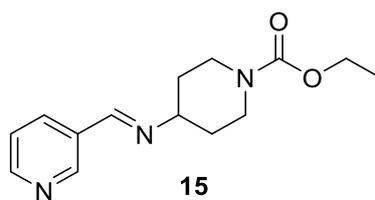


Figure S3: 400 MHz ^1H NMR spectrum of Ethyl (*E*)-4-((pyridin-3-ylimino)methyl)piperidine-1-carboxylate (**15**) in CDCl_3

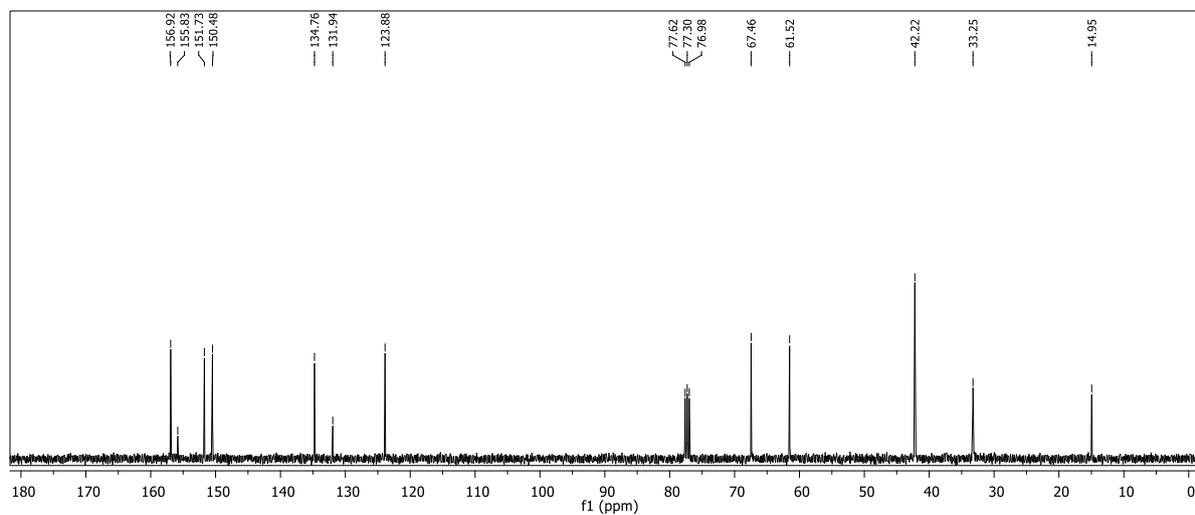


Figure S4: 100 MHz ^{13}C -NMR spectrum of Ethyl (*E*)-4-((pyridin-3-ylimino)methyl)piperidine-1-carboxylate (**15**) in CDCl_3

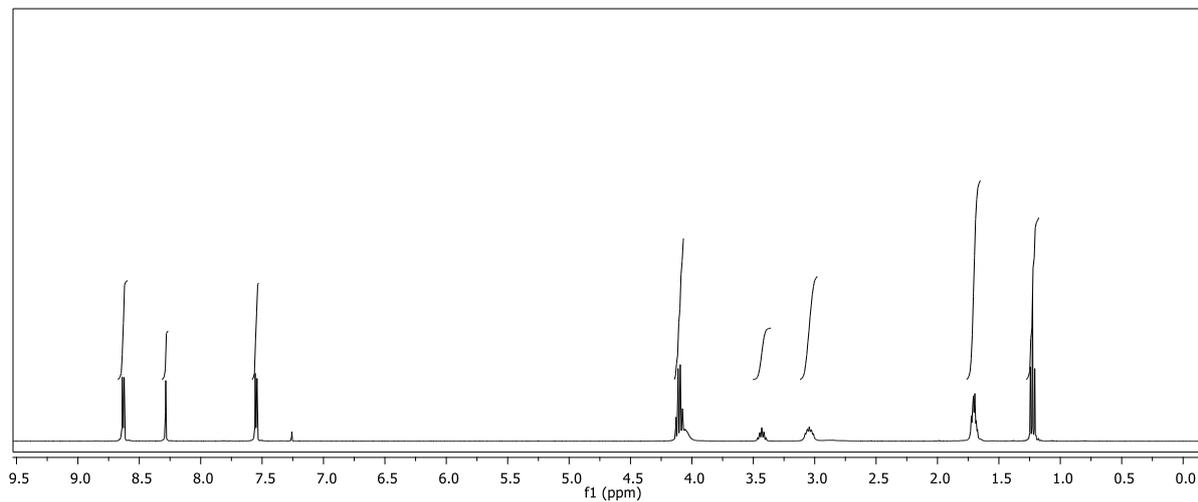
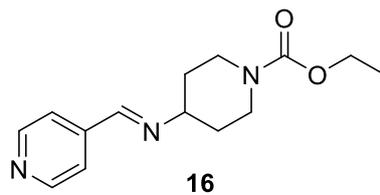


Figure S5: 400 MHz ^1H NMR spectrum of Ethyl (*E*)-4-((pyridin-4-ylimino)methyl)piperidine-1-carboxylate (**16**) in CDCl_3

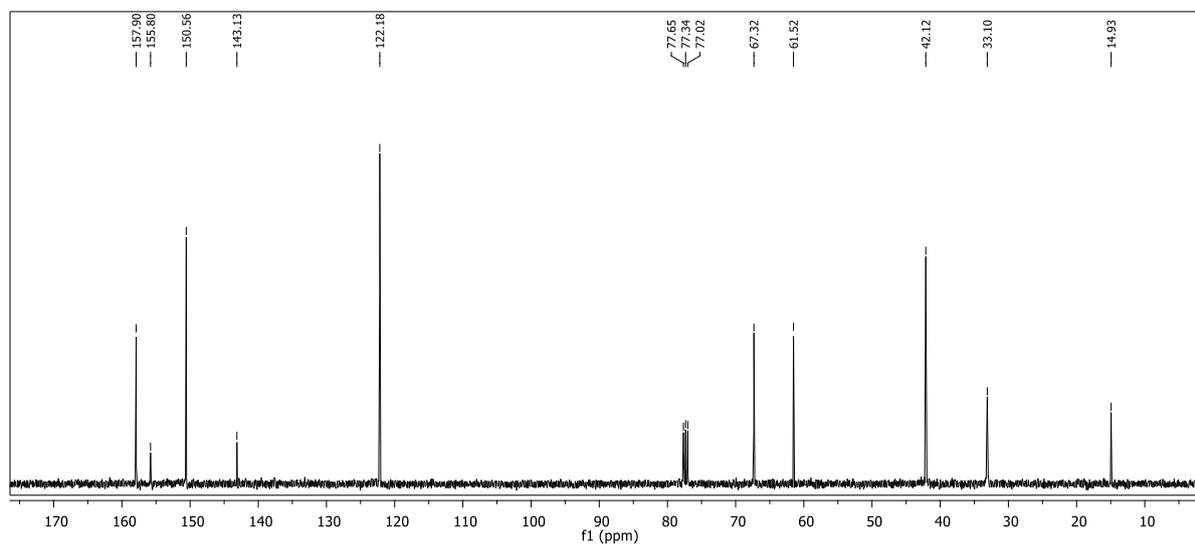


Figure S6: 100 MHz ^{13}C -NMR spectrum of Ethyl (*E*)-4-((pyridin-4-ylimino)methyl)piperidine-1-carboxylate (**16**) in CDCl_3

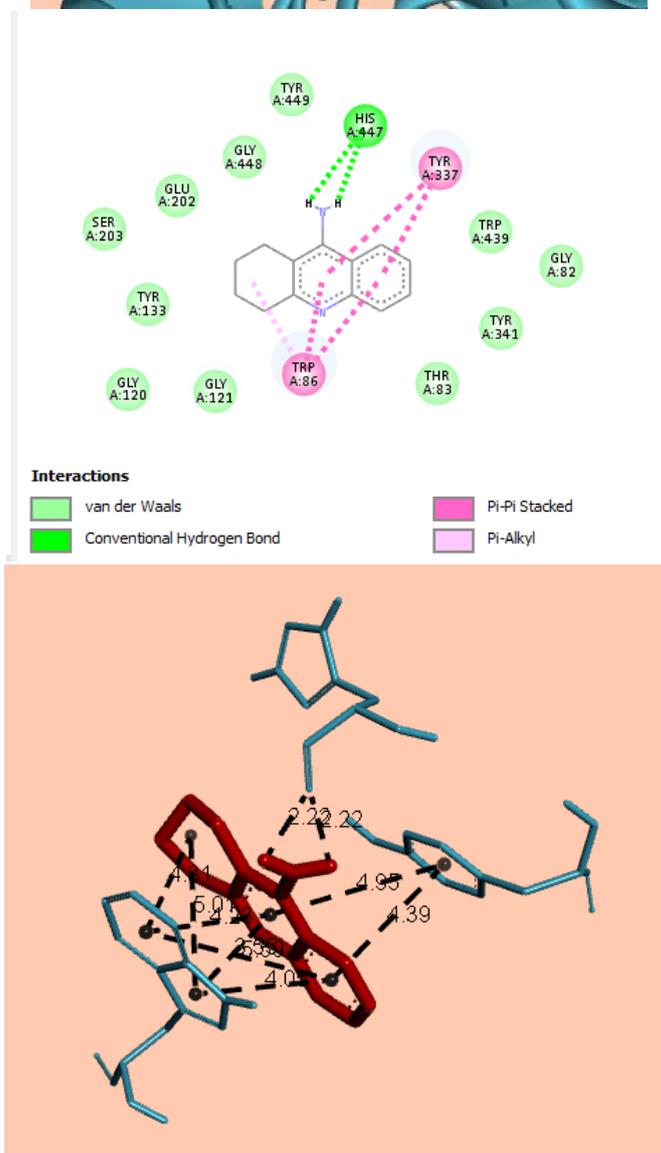
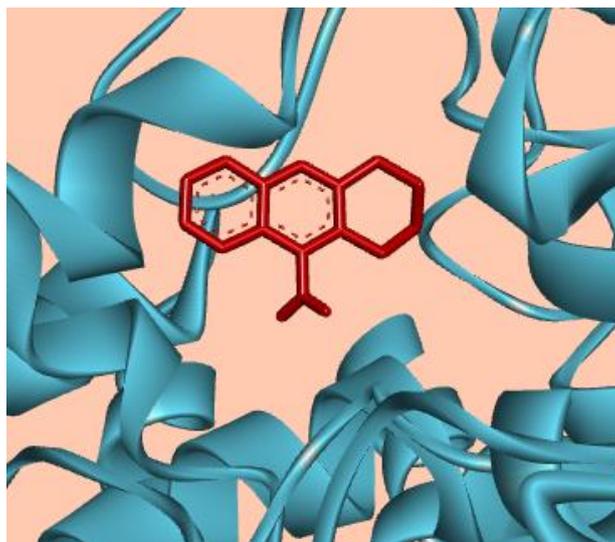


Figure S7: Docking interactions between the protein 7XN1 and the Tacrine

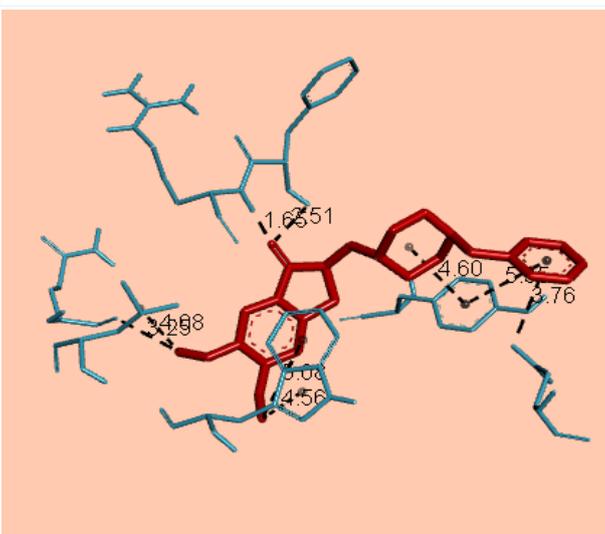
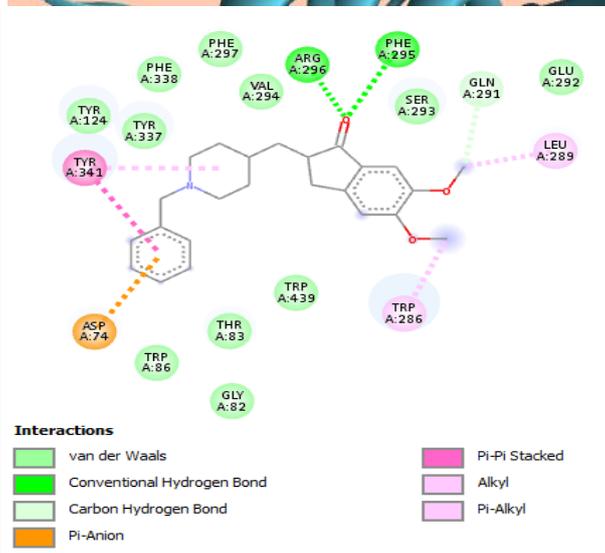
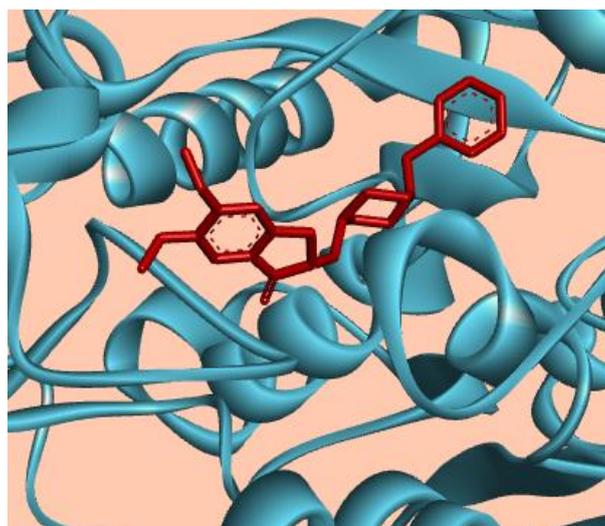
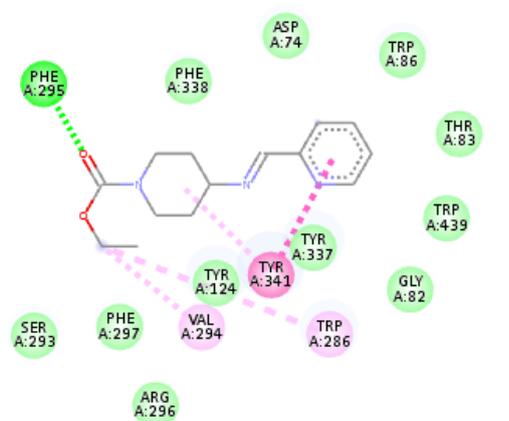
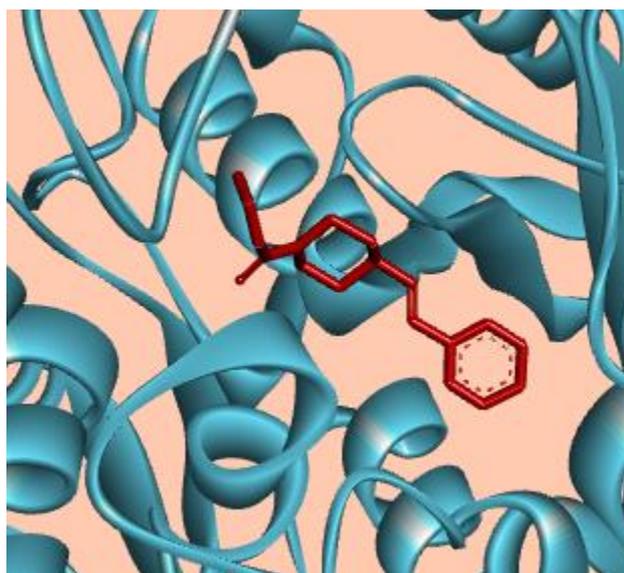


Figure S8: Docking interactions between the protein 7XN1 and the Donepezil



Interactions

- | | |
|--|--|
|  van der Waals |  Alkyl |
|  Conventional Hydrogen Bond |  Pi-Alkyl |
|  Pi-Pi Stacked | |

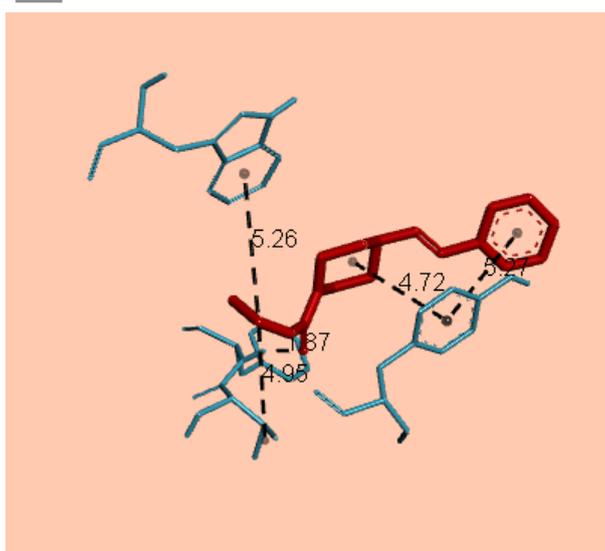


Figure S9: Docking interactions between the protein 7XN1 and compound 14

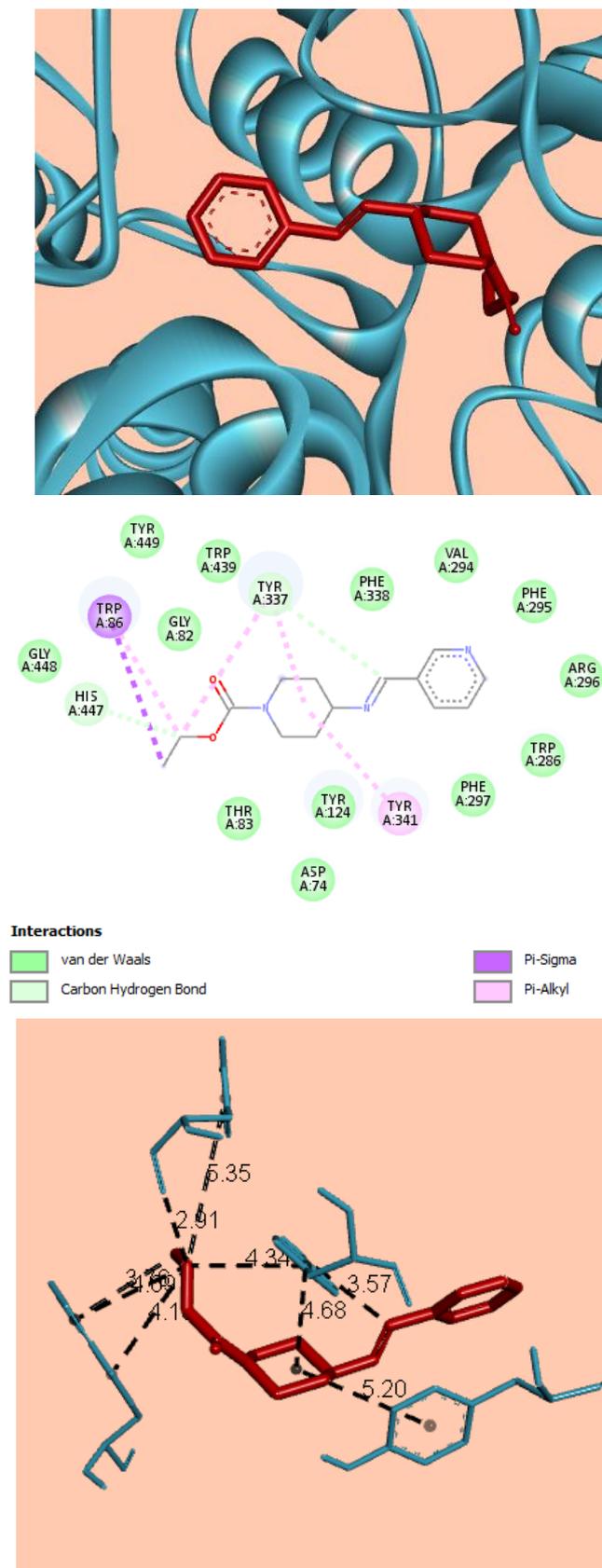


Figure S10: Docking interactions between the protein 7XN1 and Compound 15

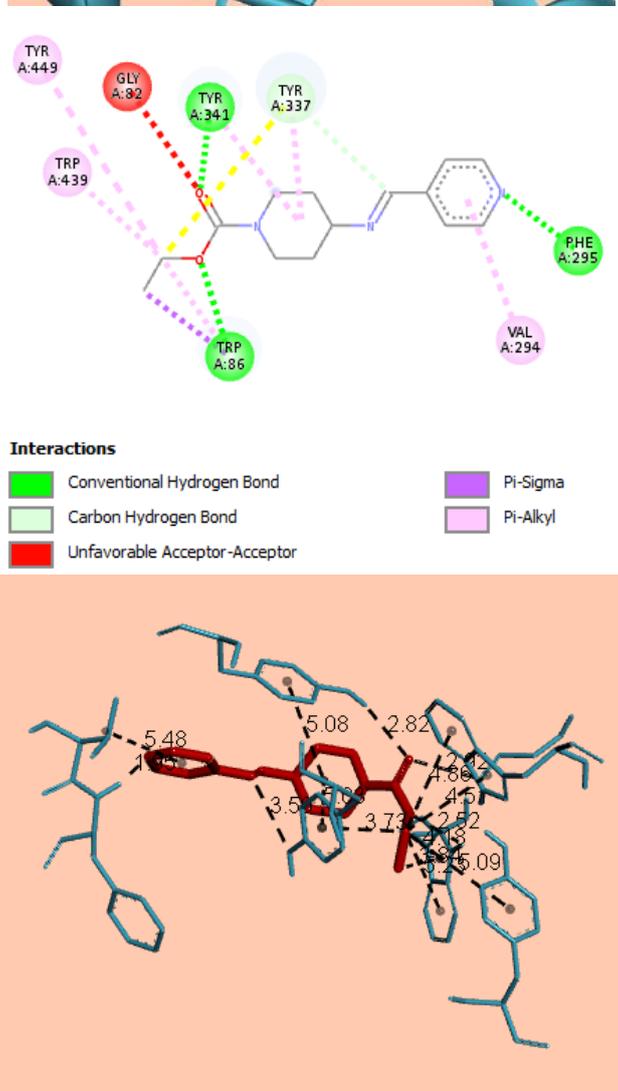
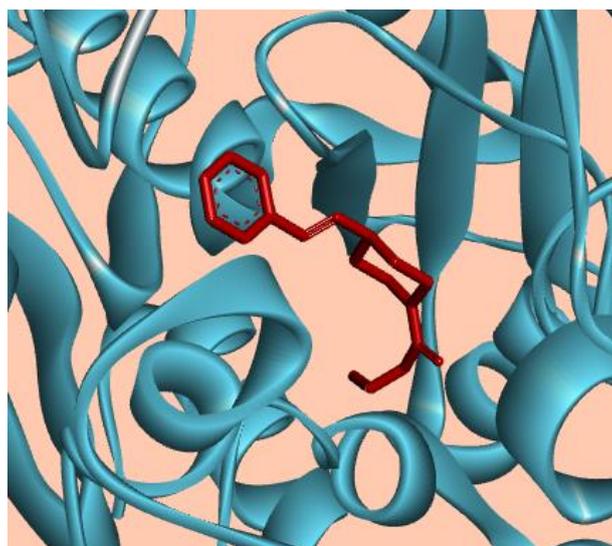


Figure S11: Docking interactions between the protein 7XN1 and the compound **16**.