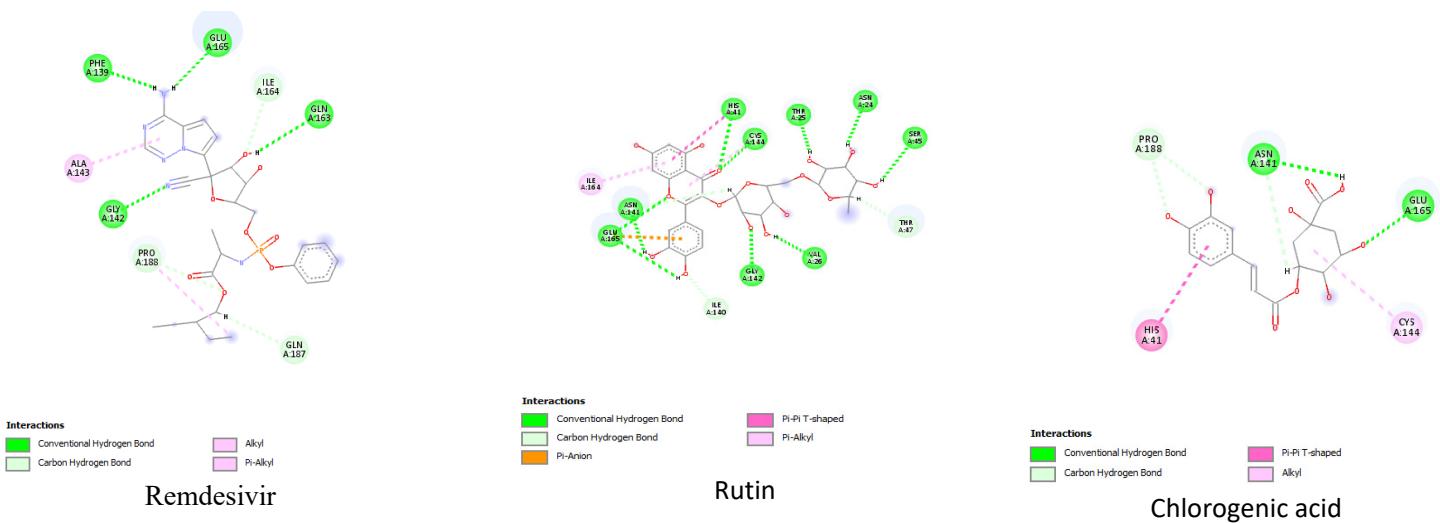
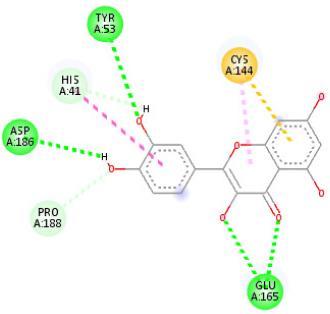


Figure S1. Frontier molecular orbitals (HOMOs, LUMOs) and Molecular electrostatic potential (MEP) surface of all detected phenolic compounds by using DFT/B3LYP/6-311++G(2d,2p) method





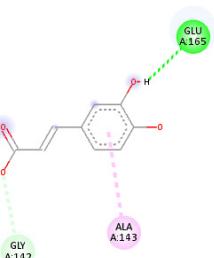
Interactions

- Conventional Hydrogen Bond
- Pi-Sulfur
- Carbon Hydrogen Bond
- Pi-Pi T-shaped
- Pi-Donor Hydrogen Bond
- Pi-Alkyl



Interactions

- Conventional Hydrogen Bond
- Pi-Sulfur
- Carbon Hydrogen Bond
- Pi-Pi T-shaped
- Pi-Sulfur



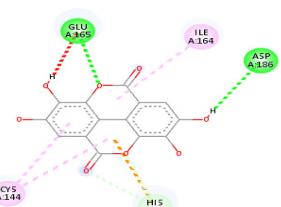
Interactions

- Conventional Hydrogen Bond
- Carbon Hydrogen Bond
- Pi-Alkyl

Quercetin

Kaempferol

Caffeic acid



Interactions

- Conventional Hydrogen Bond
- Pi-Cation
- Carbon Hydrogen Bond
- Unfavorable Donor-Donor
- Pi-Alkyl

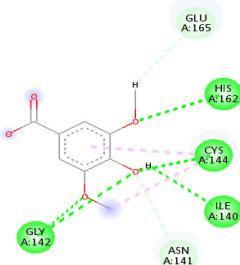
Ellagic acid



Interactions

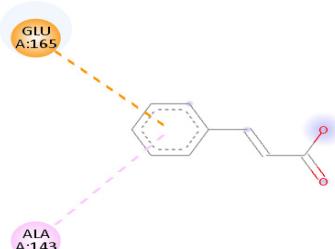
- Conventional Hydrogen Bond
- Carbon Hydrogen Bond
- Pi-Alkyl

Syringic acid



Interactions

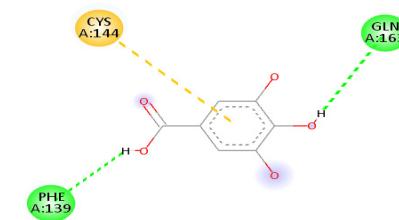
- Conventional Hydrogen Bond
- Carbon Hydrogen Bond
- Alkyl
- Pi-Alkyl



Interactions

- Pi-Anion
- Pi-Alkyl

Cinnamic acid



Interactions

- Conventional Hydrogen Bond
- Pi-Sulfur

Methyl gallate

Interactions

- Conventional Hydrogen Bond
- Alkyl
- Pi-Alkyl

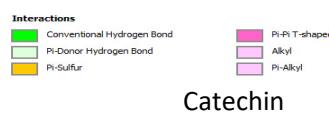
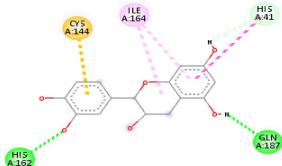
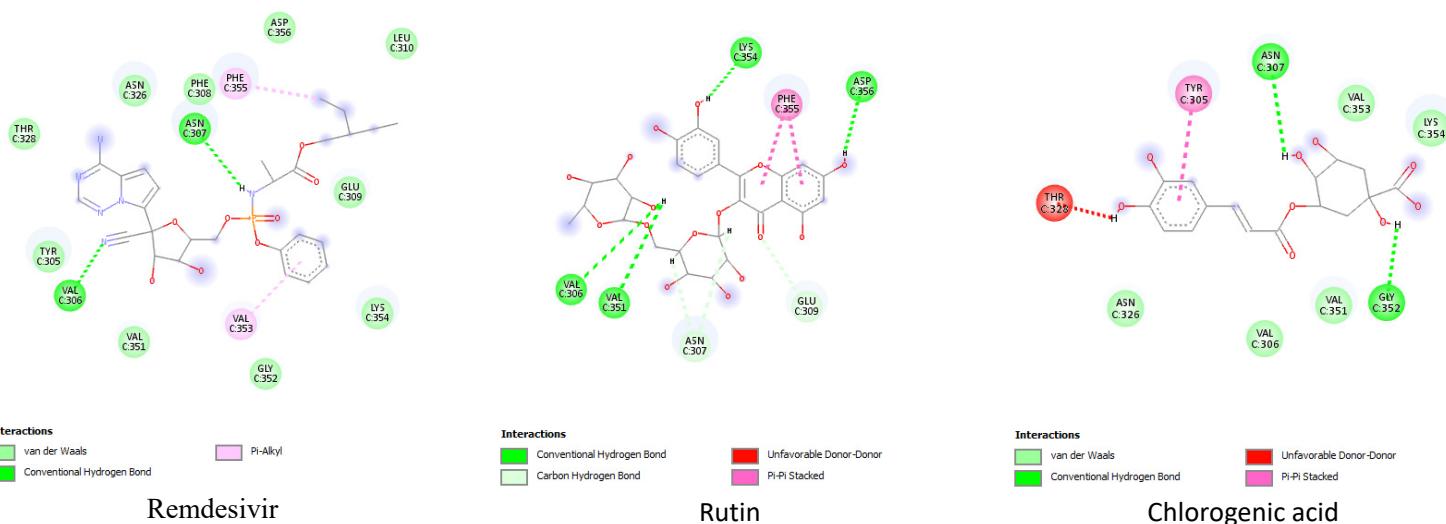
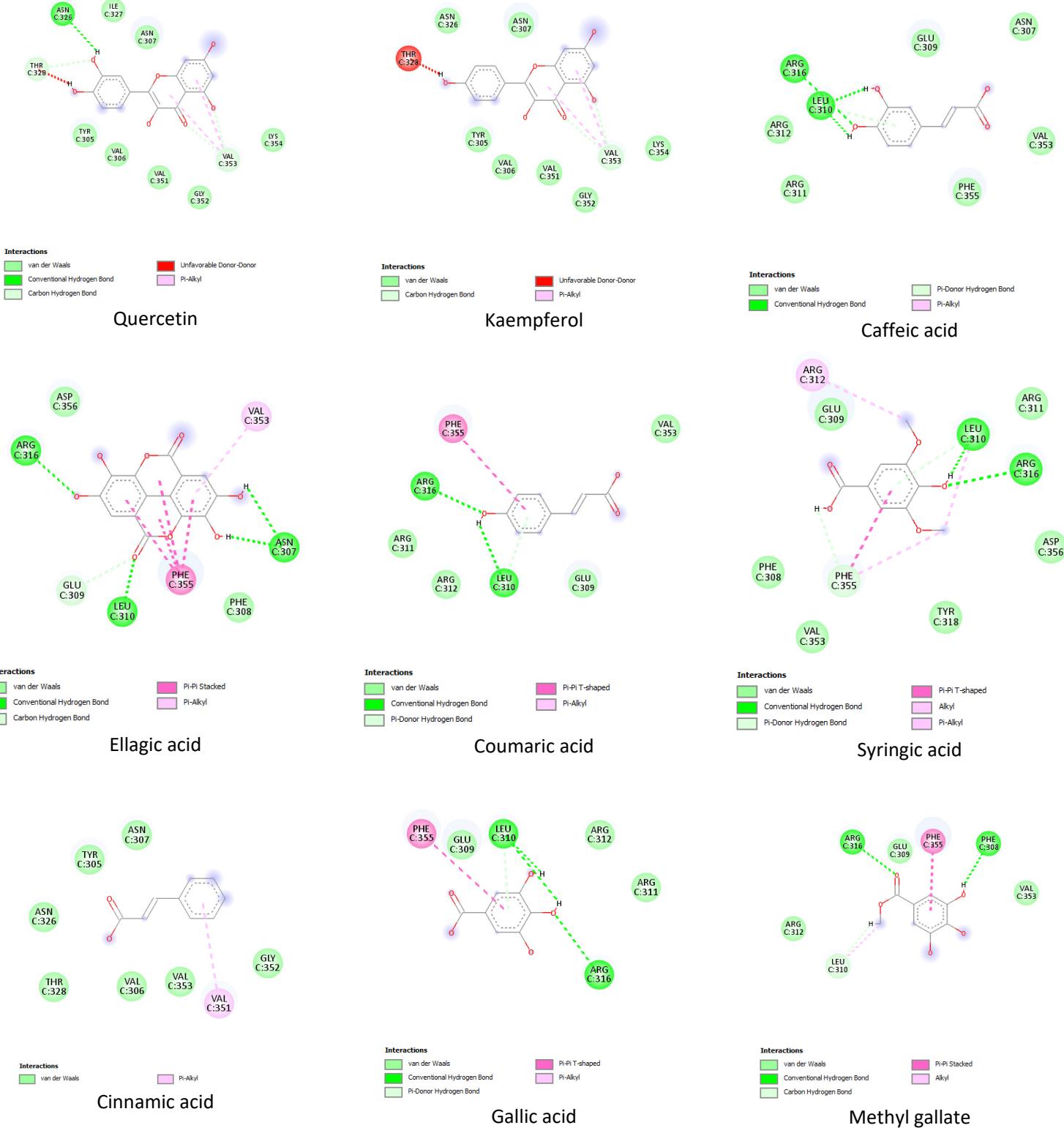


Figure S2. 2D interactions 12 phenolic compounds of GCAE and Remdesivir docked on HCoV-229E M^{pro} (PDB ID: 2ZU2)





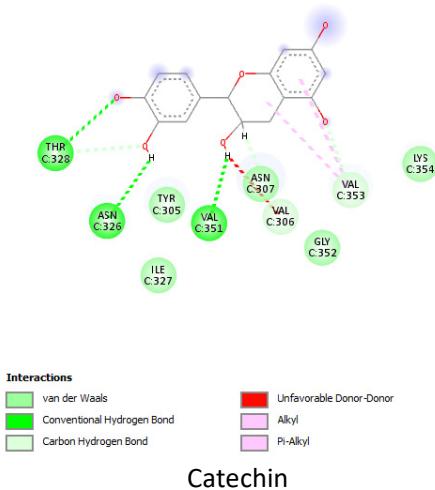
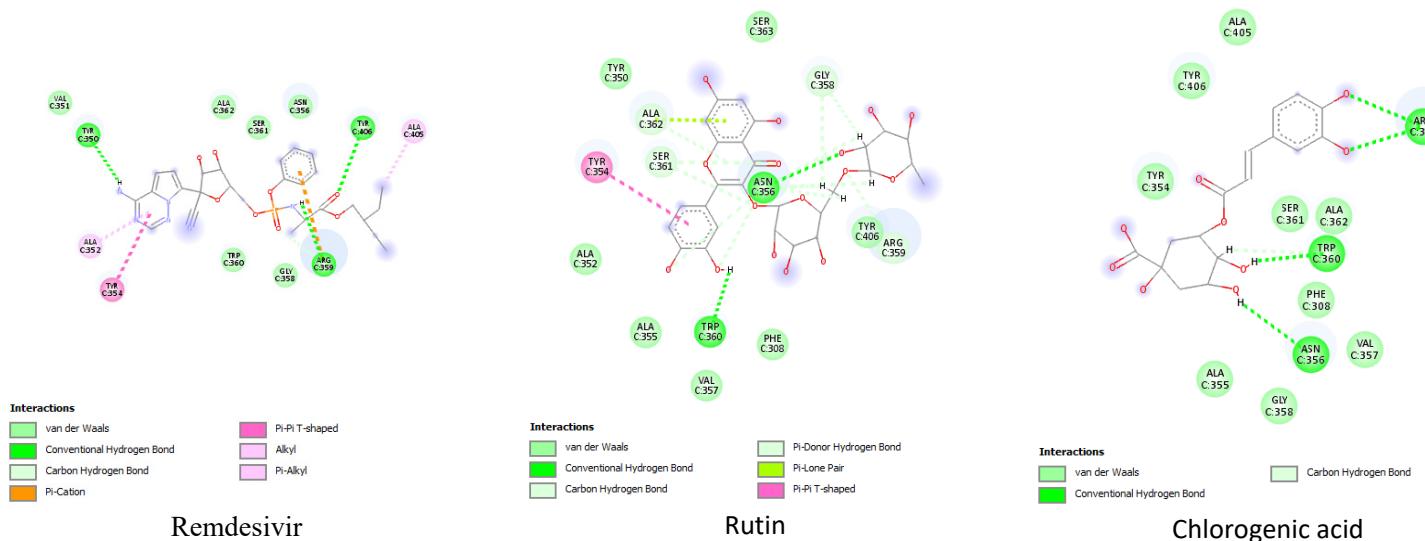
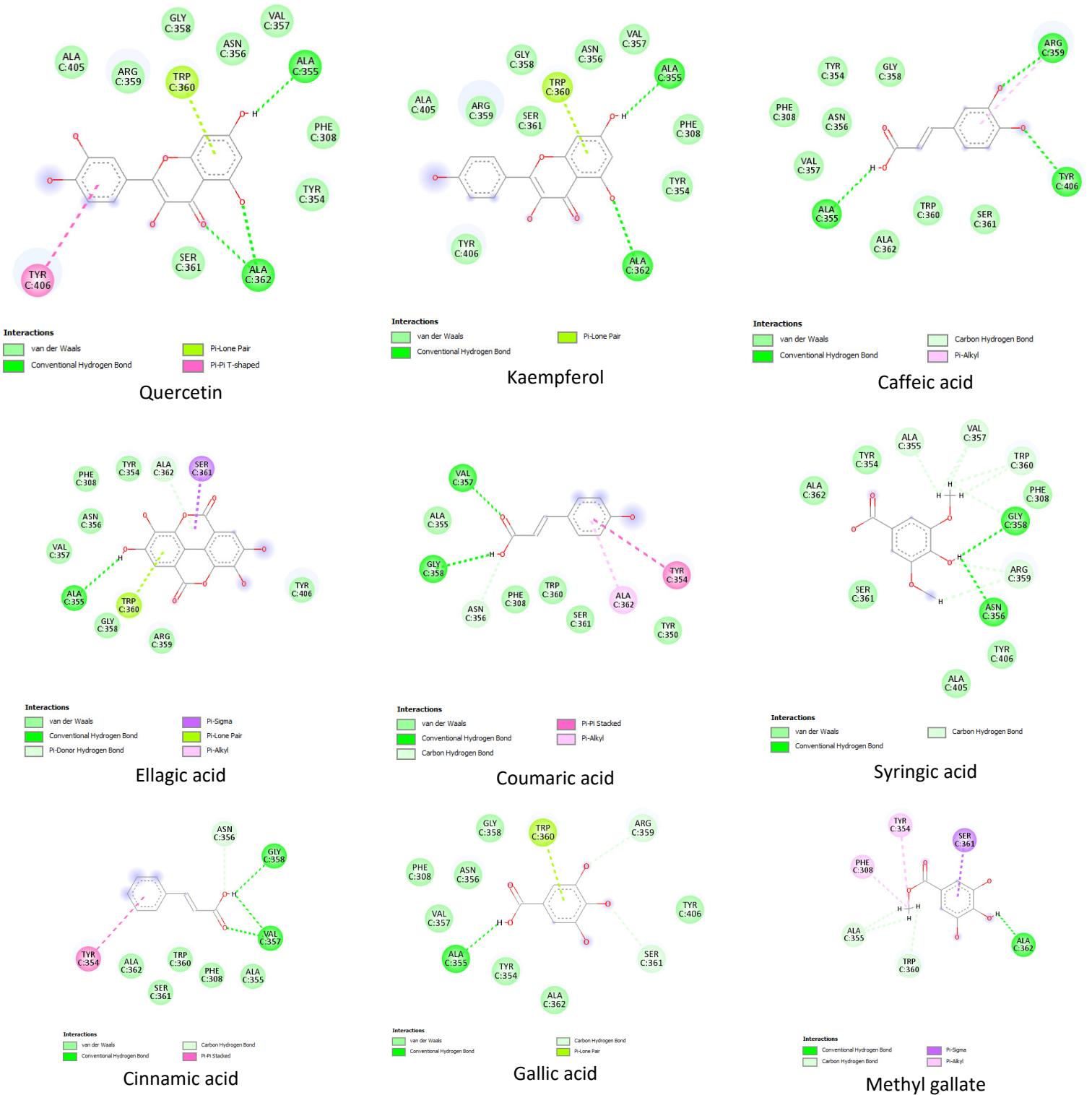


Figure S3. 2D interactions 12 phenolic compounds of GCAE and Remdesivir docked on HCoV-229E RBD Class V (PDB ID: 6U7G)





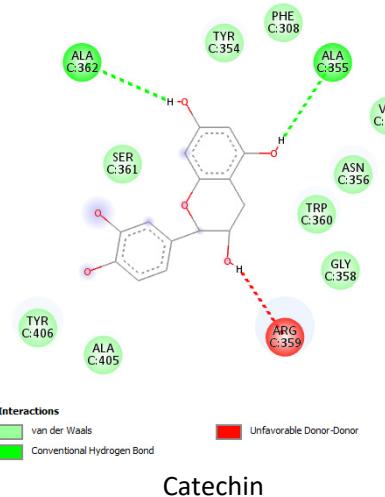
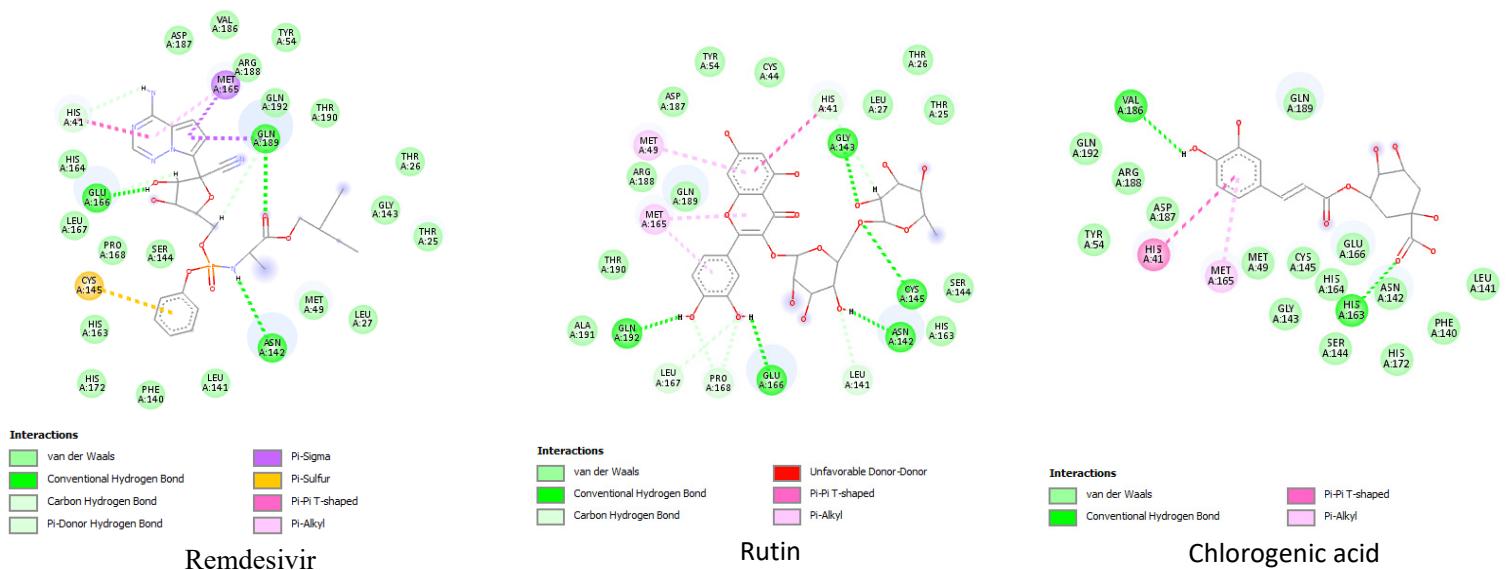
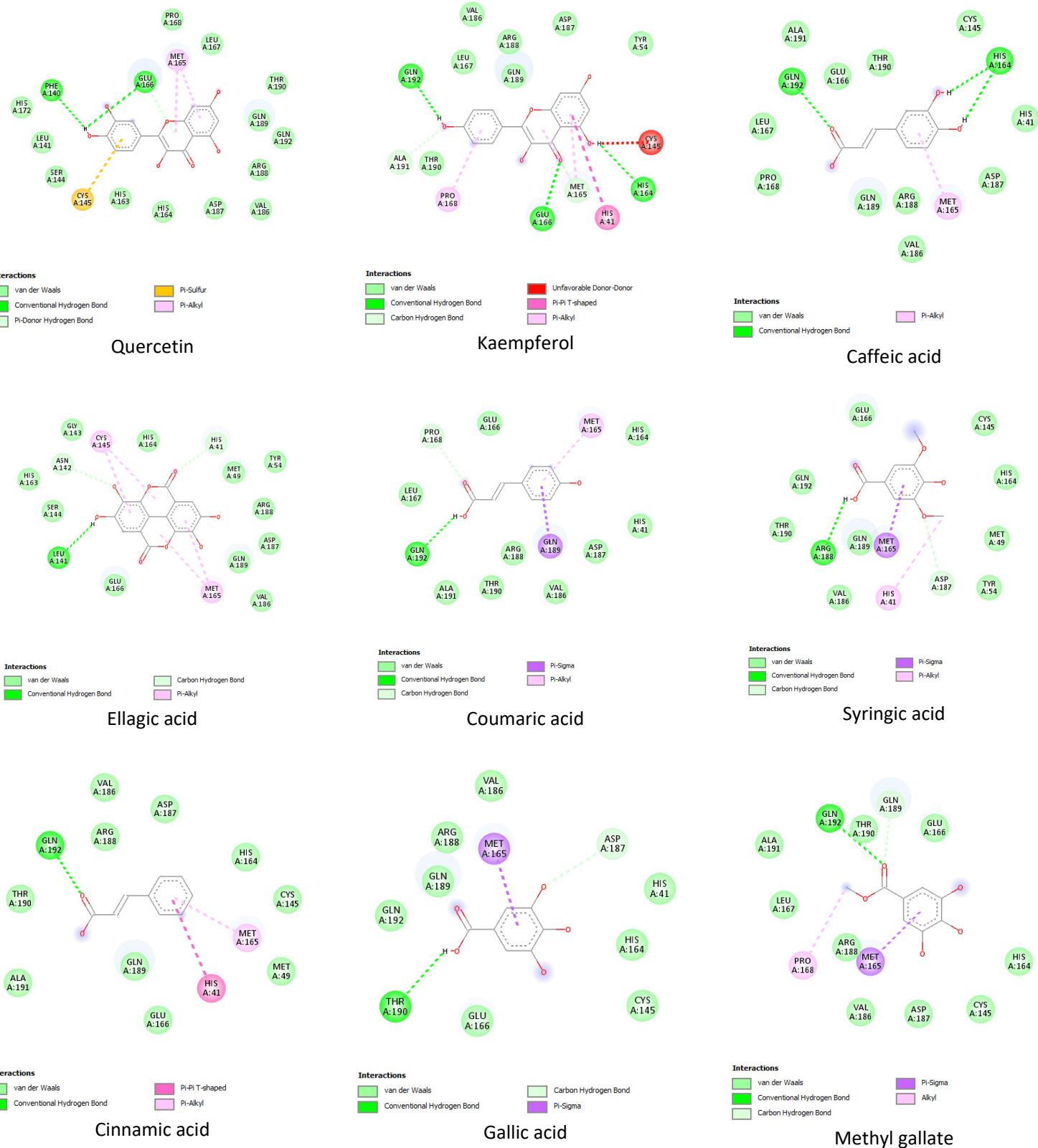


Figure S4. 2D interactions 12 phenolic compounds of GCAE and Remdesivir docked on HCoV-229E spike protein (PDB ID: 7VN9)





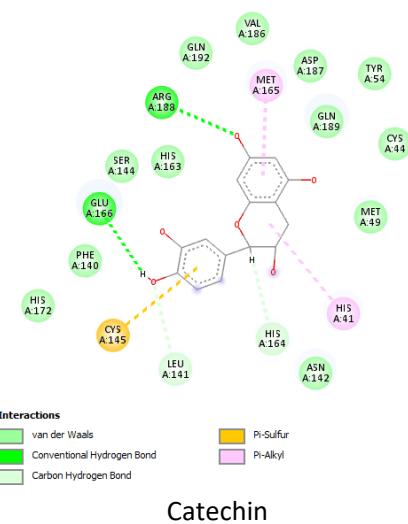


Figure S5. 2D interactions 12 phenolic compounds of GCAE and Remdesivir docked on SARS-CoV-2 M^{pro} (PDB ID: 6WTT)