# Discovery of GABA aminotransferase Inhibitors via Molecular Docking, Molecular Dynamic Simulation, and Biological Evaluation

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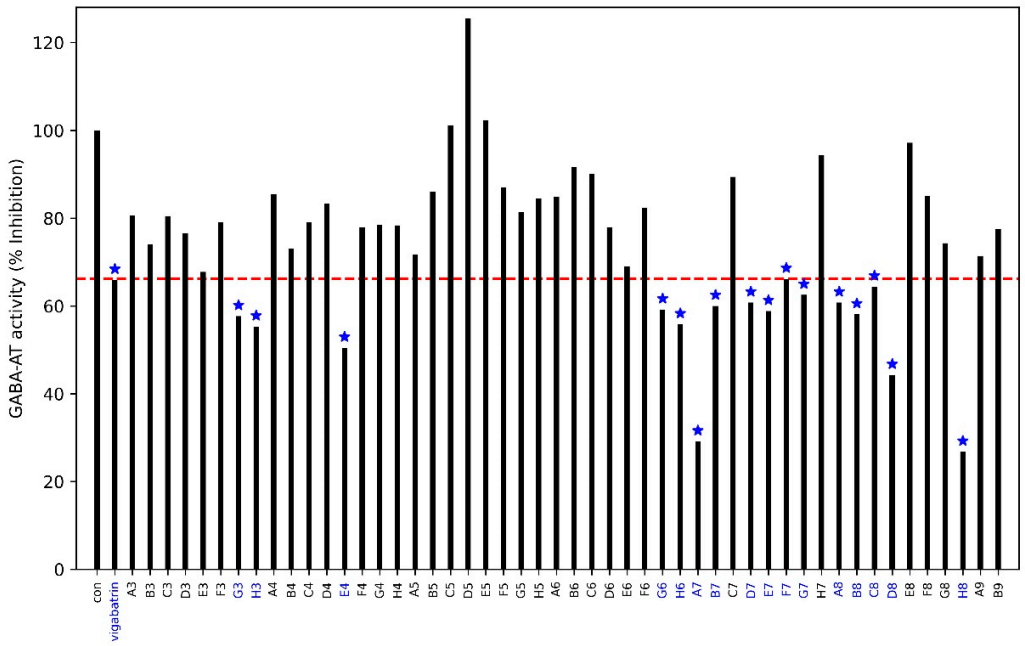



**Figure S1.** Preliminary GABA-AT inhibitory activity screening of pharmacophore-filtered KCB 50 compounds

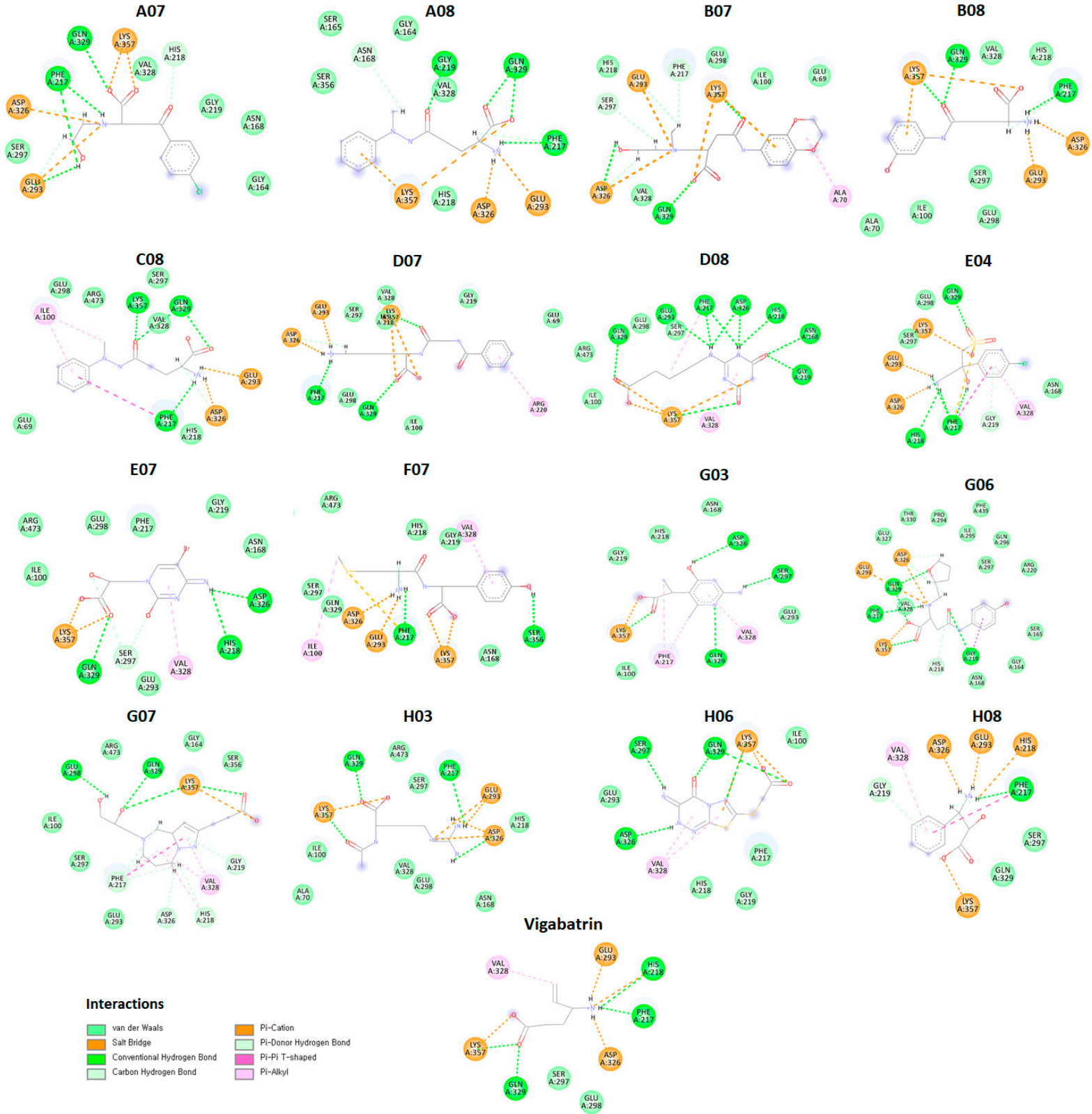



**Figure S2.** The interaction analysis of sixteen docked complexes.