Supplementary data

Synthesis and Molecular Docking Studies of Novel Bi-heterocyclic Propanamides as Anti-diabetic Agents having Mild Cytotoxicity

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Figure S1: Aromatic region of ¹H-NMR spectrum of 8r.



Figure S2: Aliphatic region of ¹H-NMRspectrum of **8r**.



Figure S3: Downfield region of ¹³C-NMR spectrum of 8r.



Figure S4: Upfield region of ¹³C-NMR spectrum of 8r.



Figure S5: EI-MS spectrum of 8r.



Figure S6: Fragmentation pattern of 8r.



Figure S7: Ramachandran graph.



Figure S8: Docking of 8a with α -glucosidase.



Figure S9: Docking of 8b with α -glucosidase.



Figure S10: Docking of 8c with α -glucosidase.



Figure S11: Docking of 8d with α -glucosidase.



Figure S12: Docking of 8e with α -glucosidase.



Figure S13: Docking of 8f with α -glucosidase.



Figure S15: Docking of 8h with α -glucosidase.



Figure S17: Docking of 8j with α -glucosidase.



Figure S18: Docking of 8k with α -glucosidase.



Figure S19: Docking of 8m with $\alpha\mbox{-glucosidase}.$



Figure S21: Docking of 80 with α -glucosidase.



Figure S22: Docking of 8p with α -glucosidase.



Figure S23: Docking of 8q with α -glucosidase.



Figure S25: Docking of 8s with α -glucosidase.

Asp202

Lys203