Supporting Information

Design, Synthesis and Molecular Docking Analysis of Flavonoid Derivatives as Potential Telomerase Inhibitors

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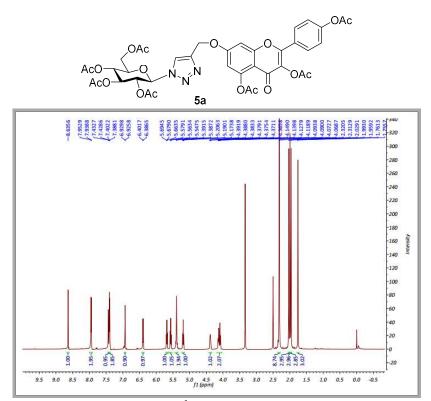


Figure S1. The ¹H-NMR of compound 5a

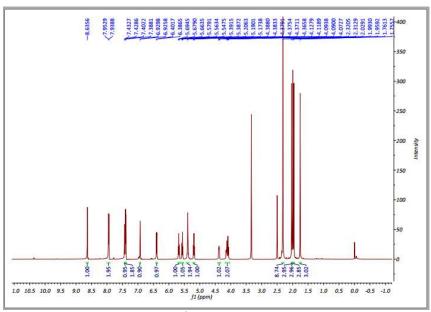
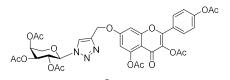


Figure S2. The ¹H-NMR of compound 5b



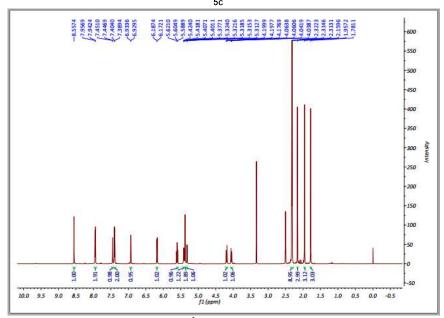
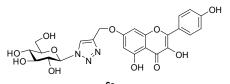


Figure S3. The ¹H-NMR of compound 5c



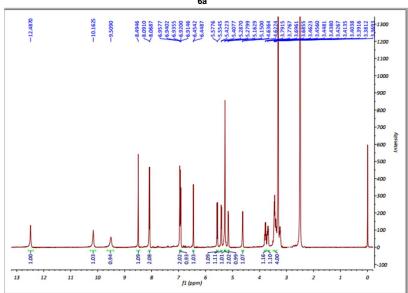


Figure S4. The ¹H-NMR of compound 6a

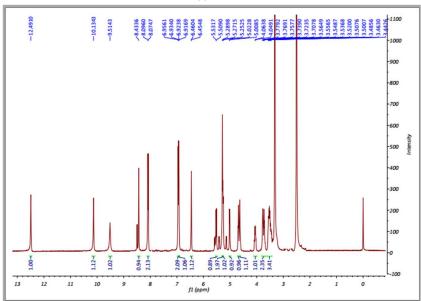


Figure S5. The ¹H-NMR of compound 6b

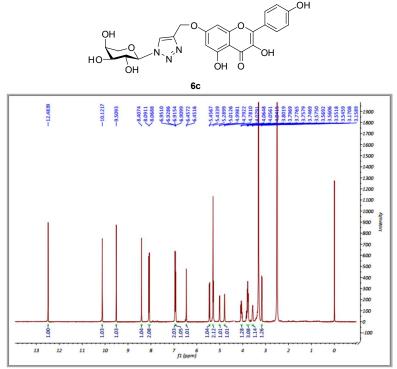


Figure S6. The ¹H-NMR of compound 6c

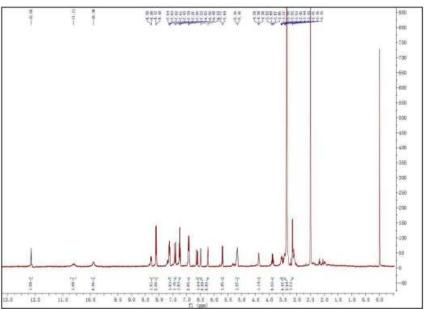


Figure S7. The ¹H-NMR of compound 14a

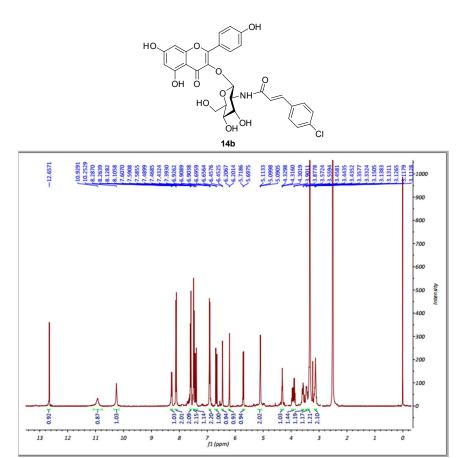
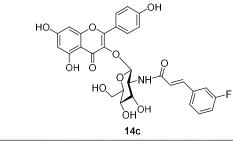


Figure S8. The ¹H-NMR of compound 14b



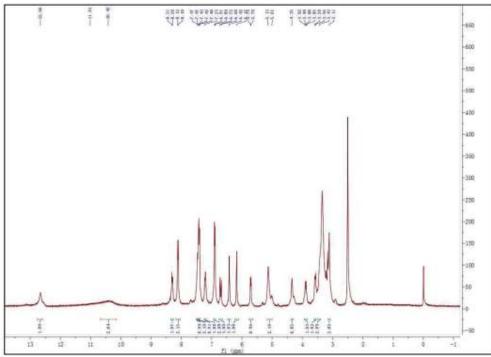


Figure S9. The ¹H-NMR of compound 14c

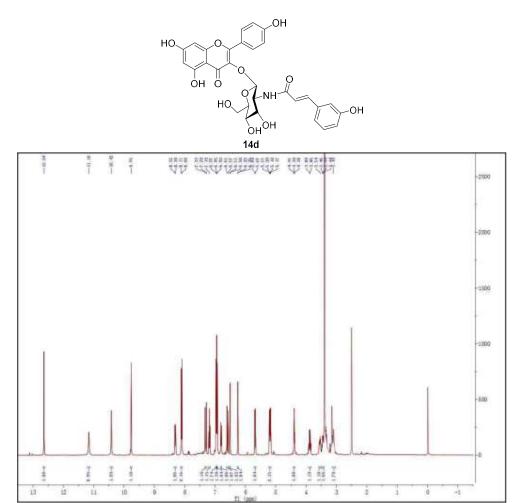


Figure S10. The ¹H-NMR of compound 14d

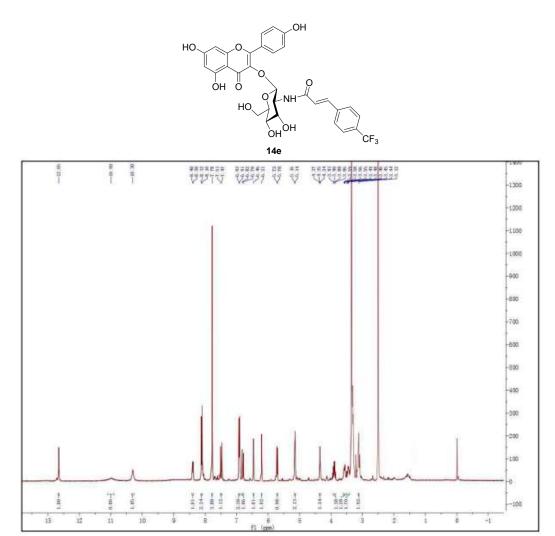


Figure S11. The ¹H-NMR of compound 14e

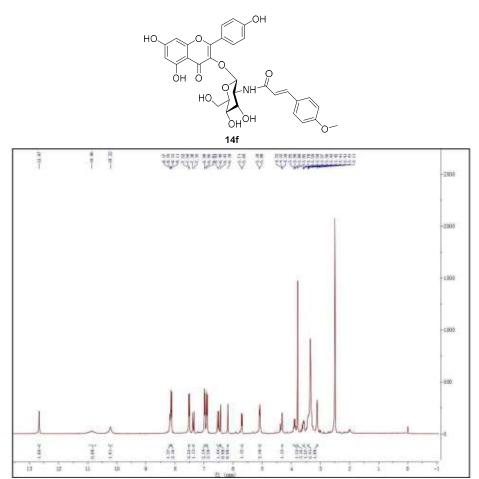


Figure S12. The ¹H-NMR of compound 14f