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Figure S1. Mass spectrum of compound 3 (4-((7-chloroquinolin-4-yl)oxy)-3-

ethoxybenzaldehyde) in MeOH showing prominent parent ion peak at 328.0742 (m/z) which corresponds to $[M+H]^+$.



Figure S2. Mass spectrum of compound 5a (N-(4-((7-chloroquinolin-4-yl)oxy)-3-

ethoxybenzyl)aniline) in MeOH showing prominent parent ion peak at 405.1387 (m/z) which corresponds to $[M+H]^+$.



Figure S3. Mass spectrum of compound 5b (N-(4-((7-chloroquinolin-4-yl)oxy)-3-

ethoxybenzyl)-4-methylaniline) in MeOH showing prominent parent ion peak at 419.1523 (m/z) which corresponds to $[M+H]^+$.



Figure S4. Mass spectrum of compound **5c** (N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-6-methylpyridin-2-amine) in MeOH showing prominent parent ion peak at 420.1480 (m/z) which corresponds to $[M+H]^+$.



Figure S5. Mass spectrum of compound 5d (N-(4-((7-chloroquinolin-4-yl)oxy)-3-

ethoxybenzyl)-4-methylaniline) in MeOH showing prominent parent ion peak at 406.1273 (m/z) which corresponds to $[M+H]^+$.



Figure S6. Mass spectrum of compound **5e** (N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-3-(trifluoromethyl)aniline) in MeOH showing prominent parent ion peak at 473.1248 (m/z) which corresponds to $[M+H]^+$.



Figure S7. Mass spectrum of compound **5f** (N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-2,4-difluoroaniline) in MeOH showing prominent parent ion peak at 441.0981 (m/z) which corresponds to $[M+H]^+$.



Figure S8. Mass spectrum of compound **5g** (N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-4-methoxyaniline) in MeOH showing prominent parent ion peak at 435.1469 (m/z) which corresponds to $[M+H]^+$.



Figure S9. ¹H NMR spectrum of compound 3 in CDCl₃.



Figure S10. ¹H NMR spectrum of compound **5a** in CDCl₃.



Figure S11. ¹H NMR spectrum of compound 5b in CDCl₃.



Figure S12. ¹H NMR spectrum of compound 5c in CDCl₃.



Figure S13. ¹H NMR spectrum of compound 5d in CDCl₃.



Figure S14. ¹H NMR spectrum of compound 5e in CDCl₃.



Figure S15. ¹H NMR spectrum of compound 5f in CDCl₃.



Figure S16. ¹H NMR spectrum of compound 5g in CDCl₃.



Figure S17. ¹³C NMR spectrum of compound 3 in CDCl₃.



Figure S18. ¹³C NMR spectrum of compound 5a in CDCl₃.



Figure S19. ¹³C NMR spectrum of compound 5b in CDCl₃.



Figure S20. ¹³C NMR spectrum of compound **5c** in d₆-DMSO.



Figure S21. ¹³C NMR spectrum of compound 5d in CDCl₃.



Figure S22. ¹³C NMR spectrum of compound 5e in d₆-DMSO.



Figure S23. ¹³C NMR spectrum of compound 5f in CDCl₃.



Figure S24. ¹³C NMR spectrum of compound 5g in CDCl₃.



Figure S25. Unit cell packing diagram of compound **5b**, viz. N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-4-methylaniline. Color code: Cl, Purple; N, green; C, black; O, blue; H, white.



Figure S26. Unit cell packing diagram of synthesized molecule **5b**, N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-4-methylaniline, showing short contact bonding (color code: Cl, Purple; N, green; C, black; O, blue; H, white).



Figure S27. Unit cell packing diagram of synthesized molecule **5f**, viz. N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-2,4-difluoroaniline. Color code: Cl, Purple; F, yellow; N, green; C, black; O, blue; H, white.



Figure S28. Unit cell packing diagram of synthesized molecule **5f**, viz. N-(4-((7-chloroquinolin-4-yl)oxy)-3-ethoxybenzyl)-2,4-difluoroaniline, showing short contact bonding (color code: Cl, Purple; F, yellow; N, green; C, black; O, blue; H, white).