

Supplementary Information

Synthesis and Evaluation of Fenofibric Acid Ester Derivatives: Studies of Different Formulation with Their Bioavailability and Absorption Conditions

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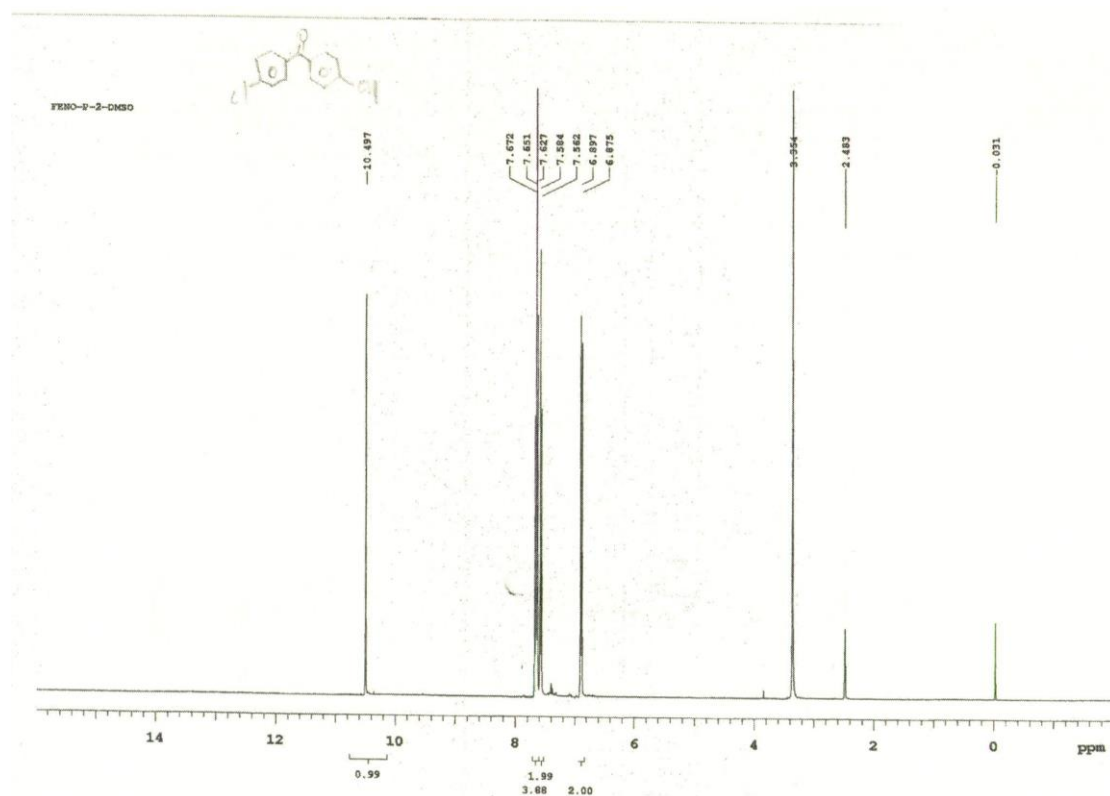


Figure S1. ¹H NMR spectrum (300 MHz, DMSO-*d*₆) of (4-chlorophenyl)(4-hydroxyphenyl)methanone.

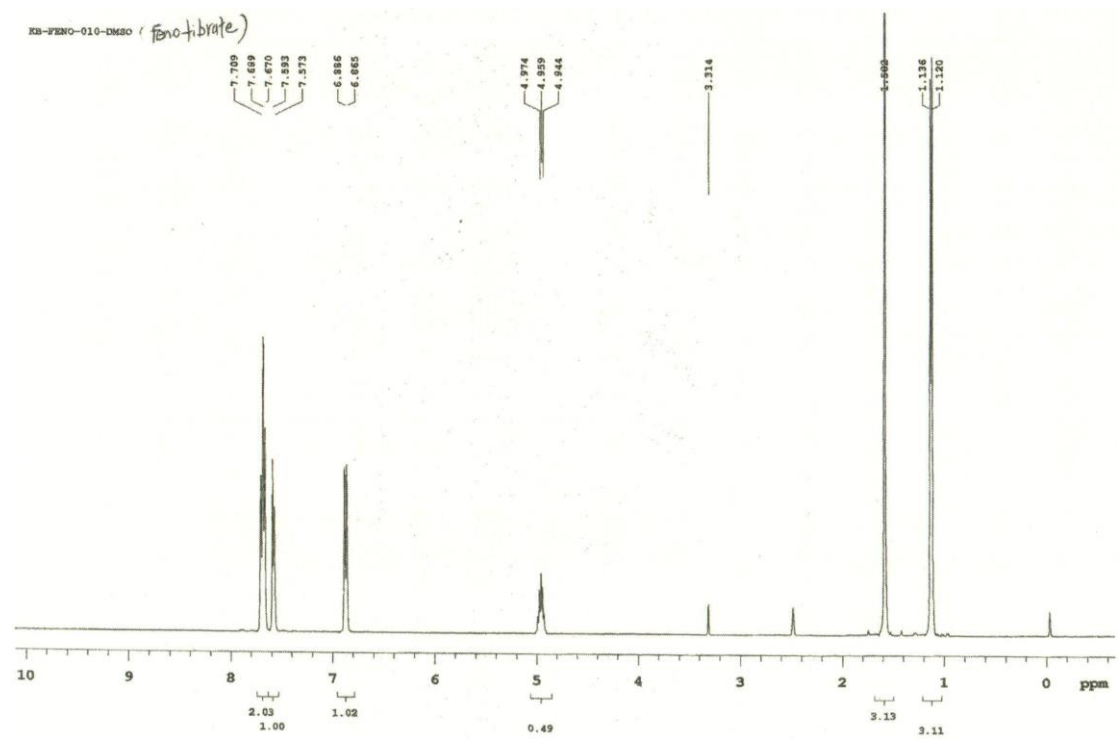


Figure S2. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of fenofibrate.

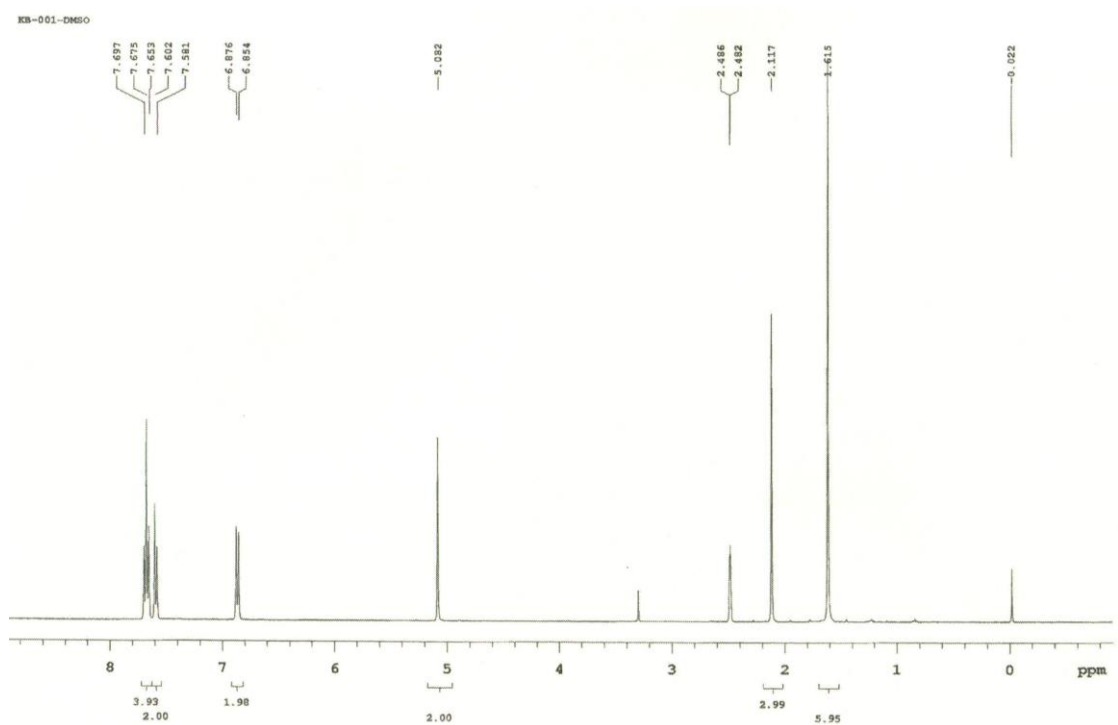


Figure S3. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of (5-methyl-2-oxo-1,3-dioxol-4-yl)-methyl-2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoate (**JF-1**).

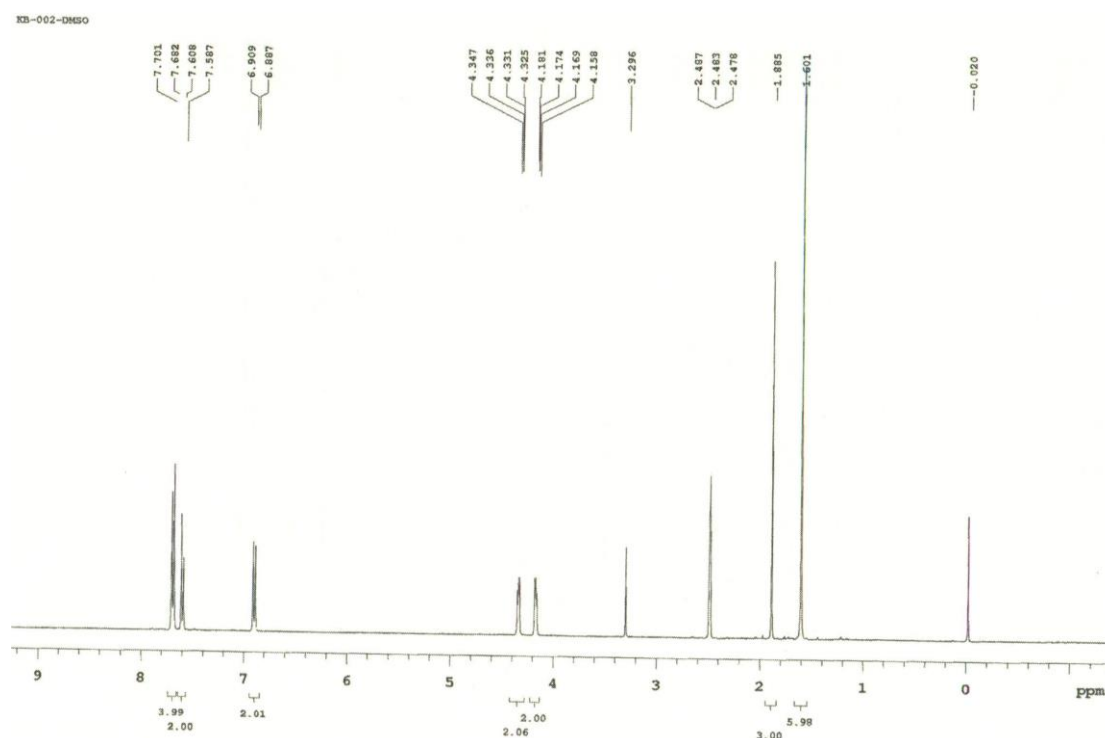


Figure S4. ¹H NMR spectrum (300 MHz, DMSO-*d*₆) of 2-acetoxyethyl-2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoate (**JF-2**).

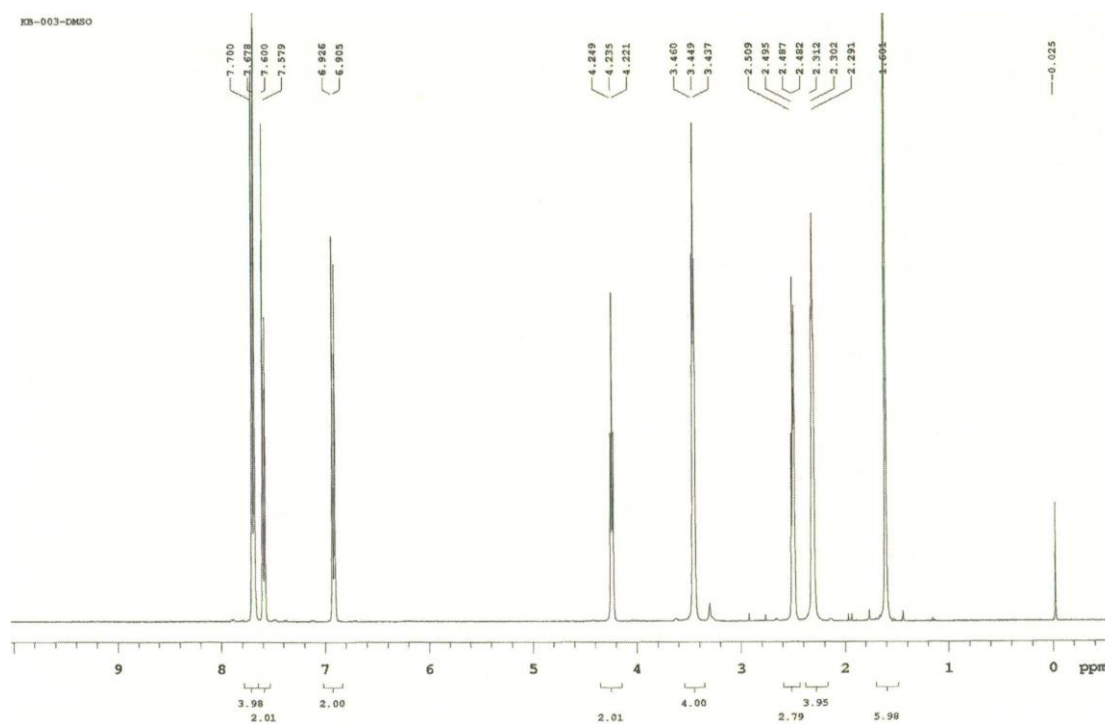


Figure S5. ¹H NMR spectrum (300 MHz, DMSO-*d*₆) of 2-morpholinoethyl-2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoate (**JF-3**).

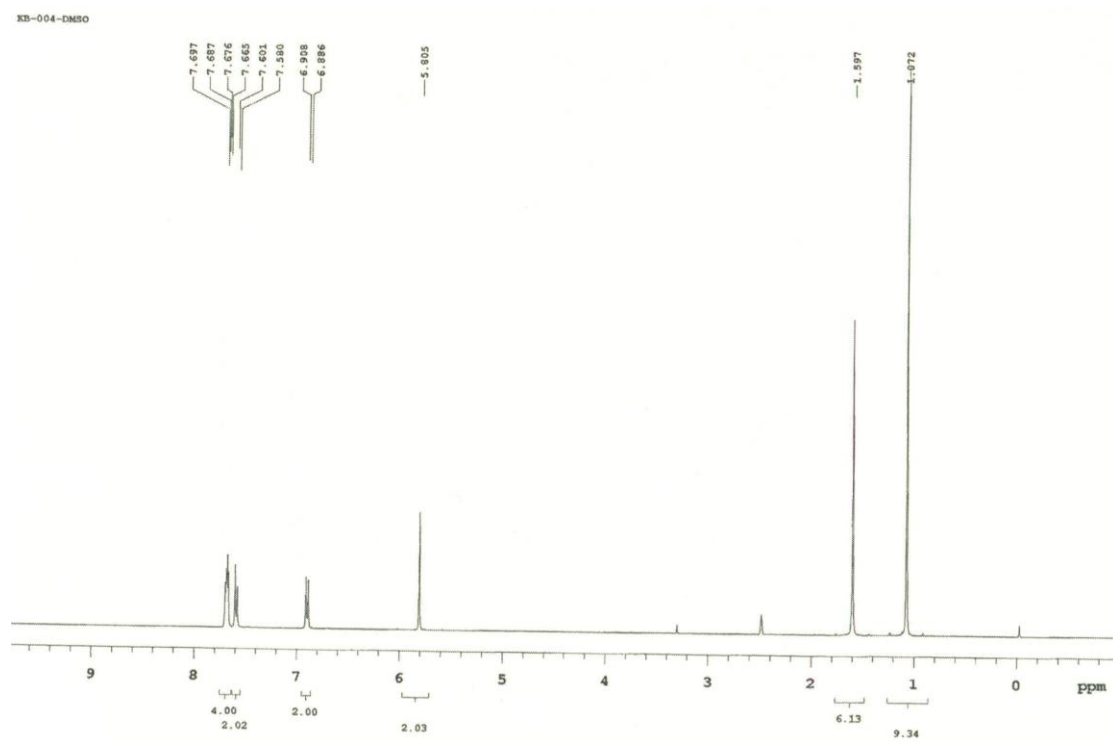


Figure S6. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) methylpivalate (**JF-4**).

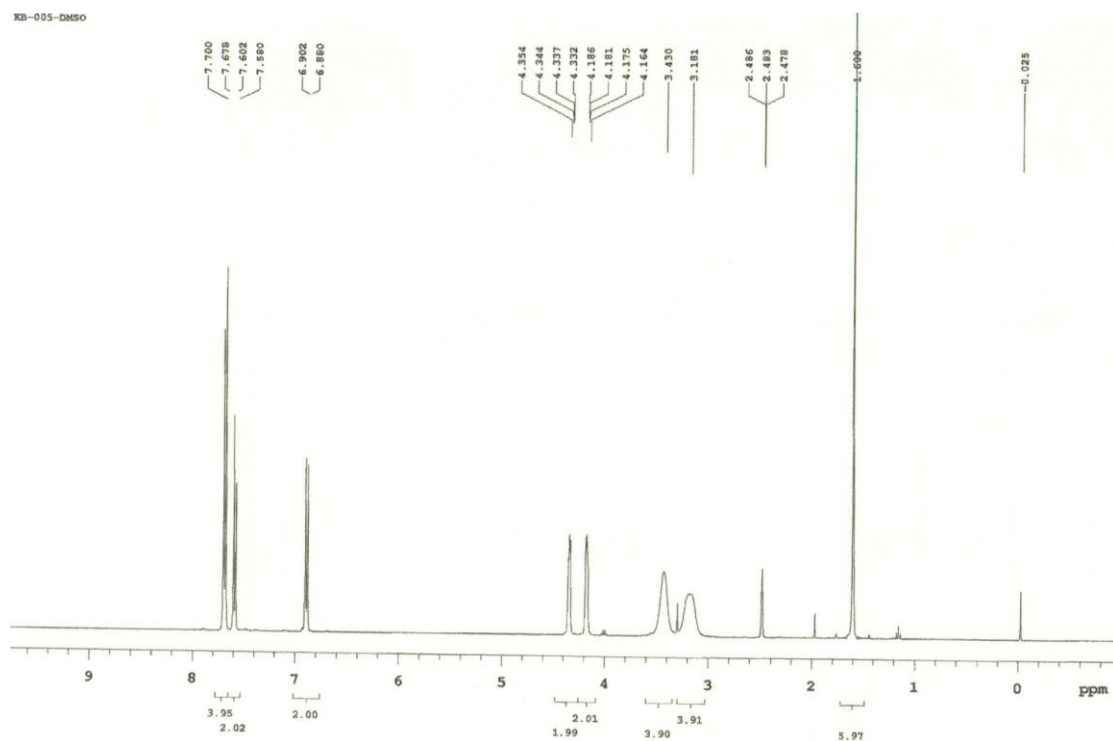


Figure S7. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethylmorpholine-4-carboxylate (**JF-5**).

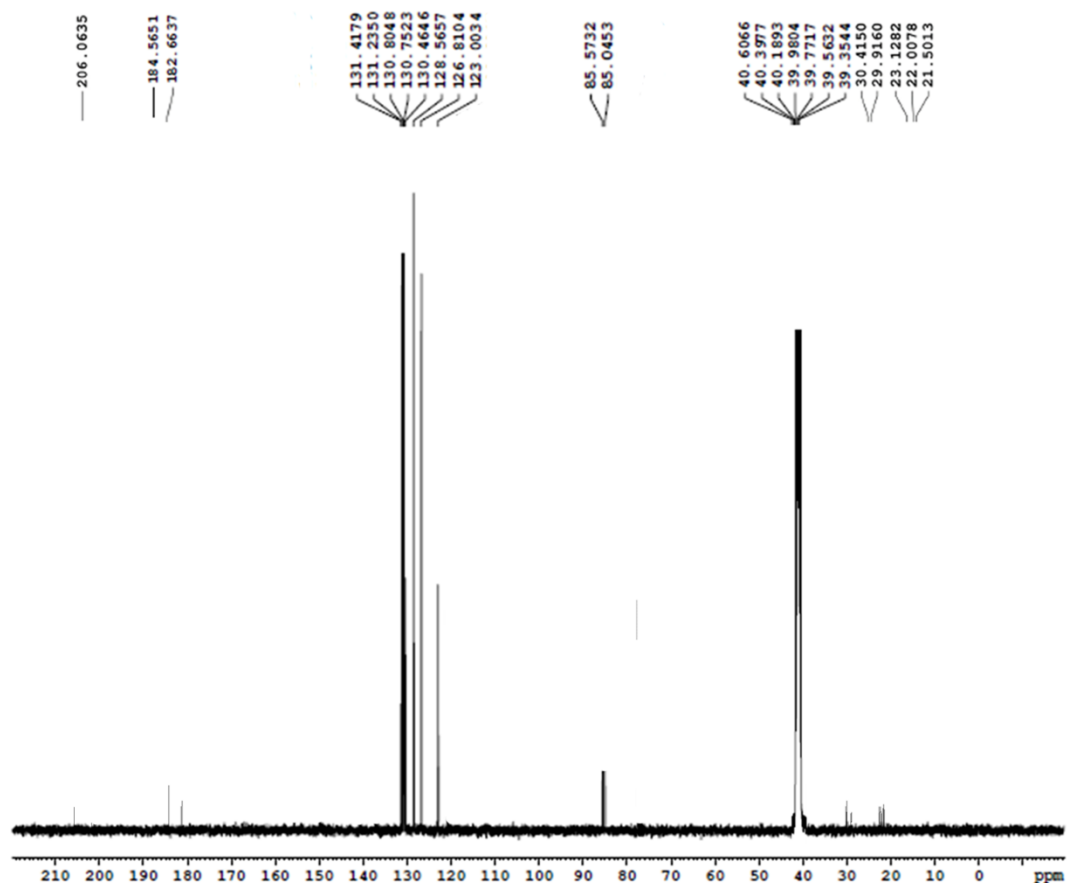


Figure S8. ^{13}C NMR spectrum (75 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethylmorpholine-4-carboxylate (**JF-5**).

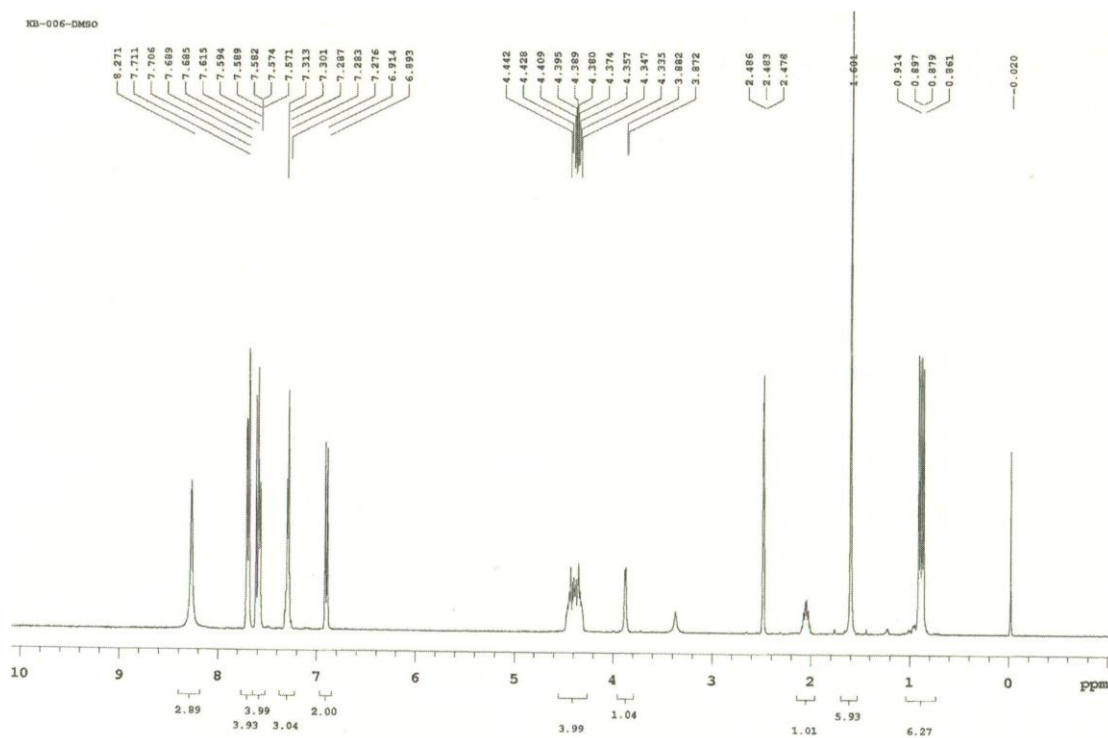


Figure S9. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethyl-2-amino-3-methylbutanionate mono-benzenesulfonate (**JF-6**).

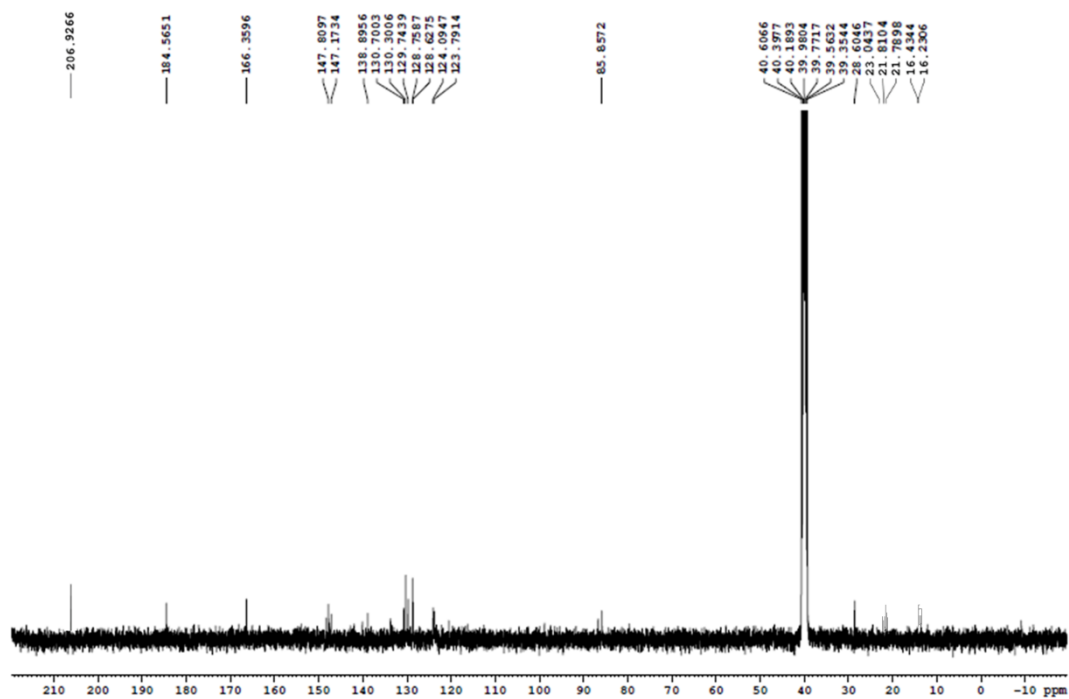


Figure S10. ^{13}C NMR spectrum (75 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethyl-2-amino-3-methylbutanonate mono-benzenesulfonate (**JF-6**).

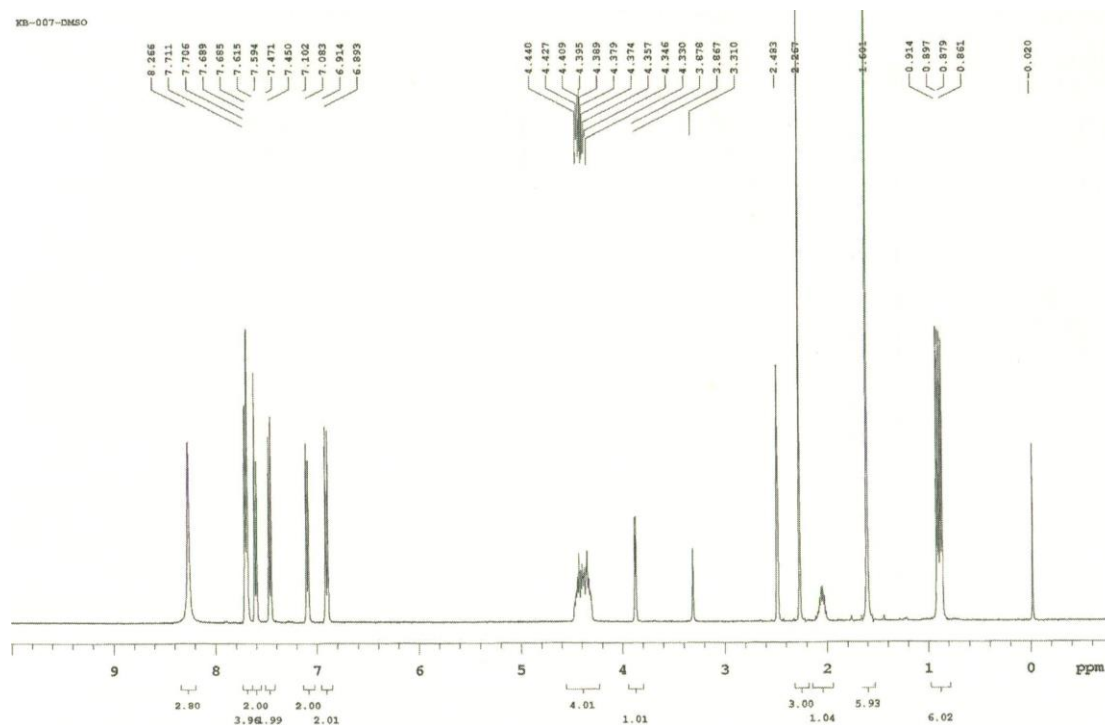


Figure S11. ^1H NMR spectrum (300 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethyl-2-amino-3-methylbutanonate mono-tosylate (**JF-7**).

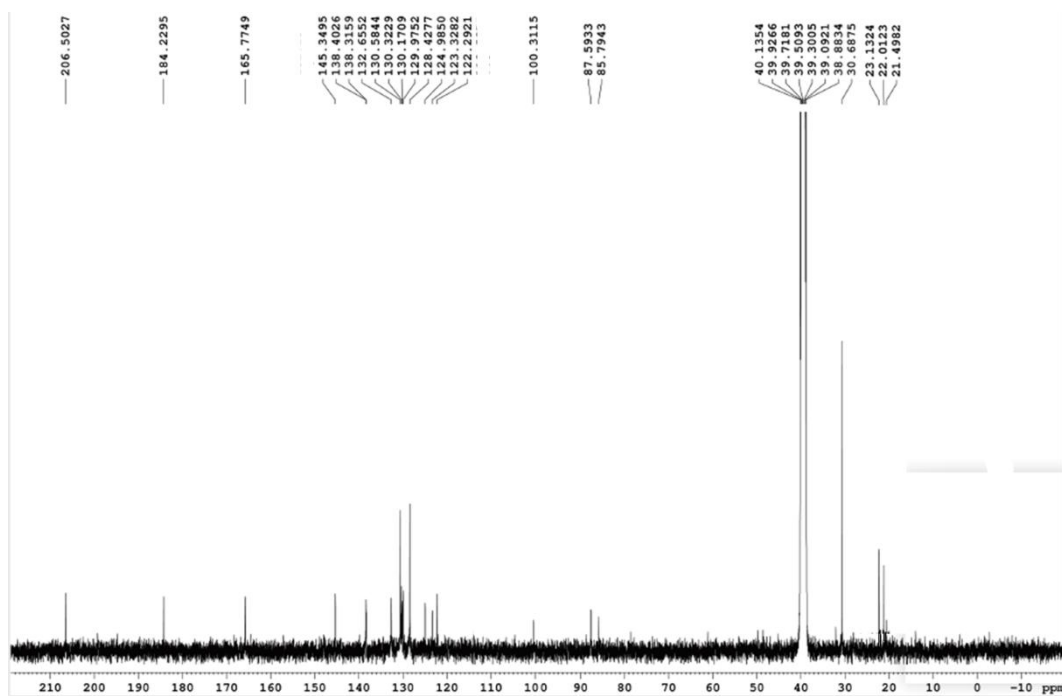


Figure S12. ^{13}C NMR spectrum (75 MHz, $\text{DMSO-}d_6$) of (2-(4-(4-chlorobenzoyl)phenoxy)-2-methylpropanoyloxy) ethyl-2-amino-3-methylbutanonate mono-tosylate (**JF-7**).