4-[(4-Chlorophenyl)carbamoyl]butanoic acid

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*** SUPPLEMENTARY MATERIALS ***

Figure S1: The FTIR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**).

Figure S2: The ^{1}H NMR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**).

Figure S3: The ^{13}C{^{1}H} NMR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**).

Figure S4: The UV spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**).

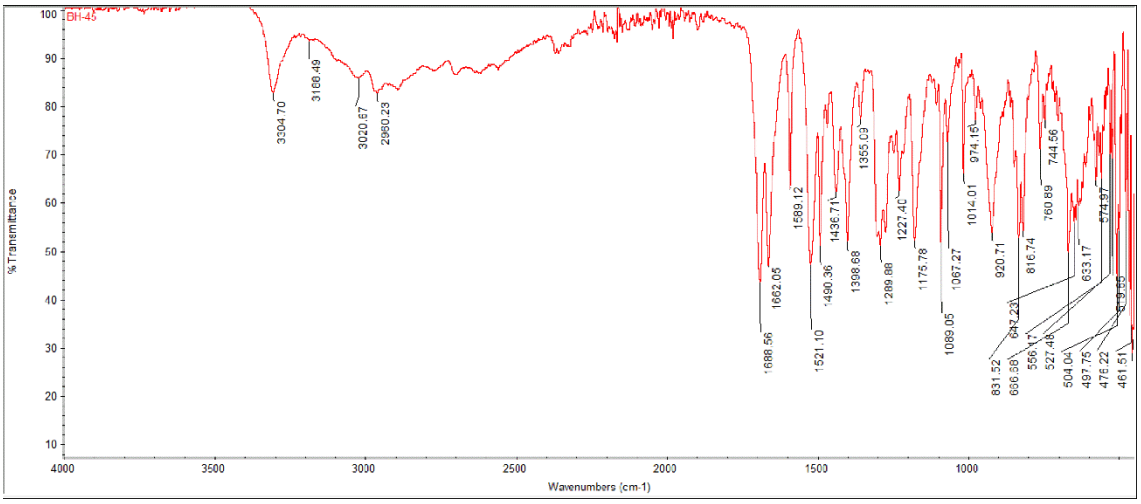

Figure S1: The FTIR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**). The FTIR spectrum was measured on a Thermo Nicolet-6700 spectrophotometer (Vienna, Austria) from 4000 to 450 cm^{-1}.

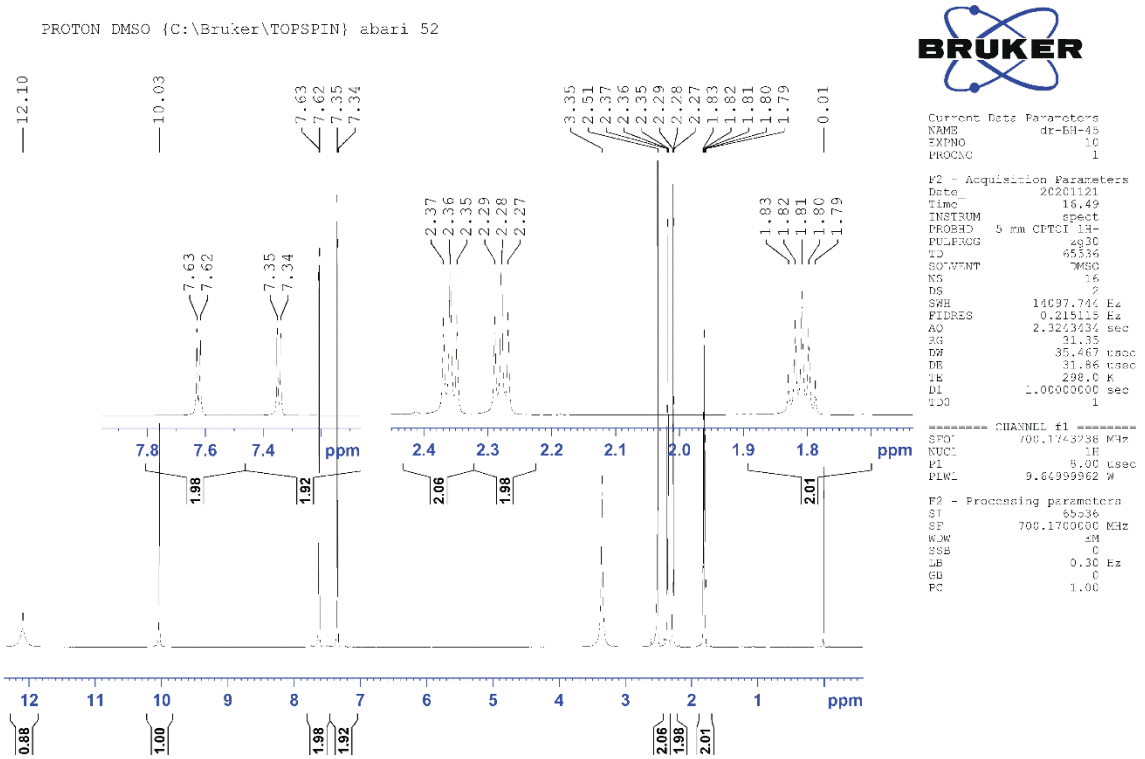

Figure S2: The ^{1}H NMR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**). The ^{1}H NMR spectrum were recorded in DMSO-d_6 solution on a Bruker Avance 500 MHz NMR (Billerica, MA, USA) spectrometer with chemical shifts relative to tetramethylsilane.

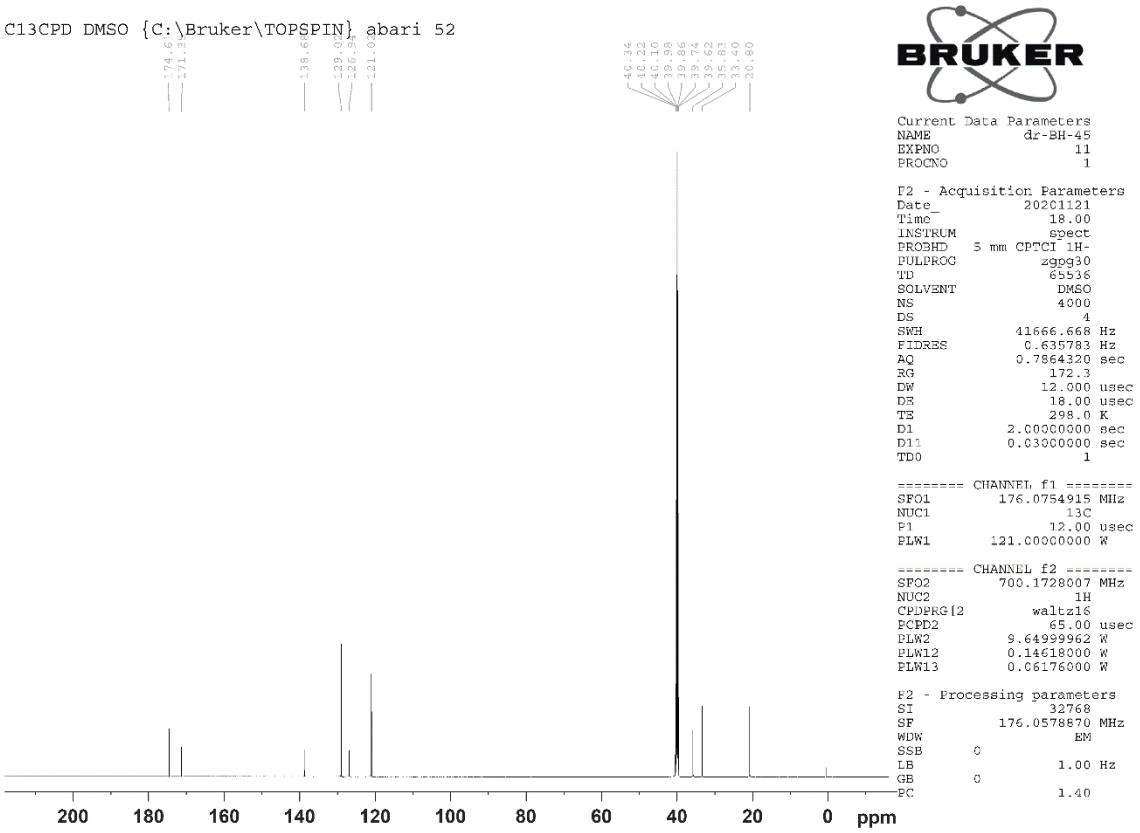

Figure S3: The $^{13}C\{^{1}H\}$ NMR spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**). The $^{13}C\{^{1}H\}$ NMR spectrum was recorded in DMSO-d_6 solution on a Bruker Avance 500 MHz NMR (Billerica, MA, USA) spectrometer with chemical shifts relative to tetramethylsilane.

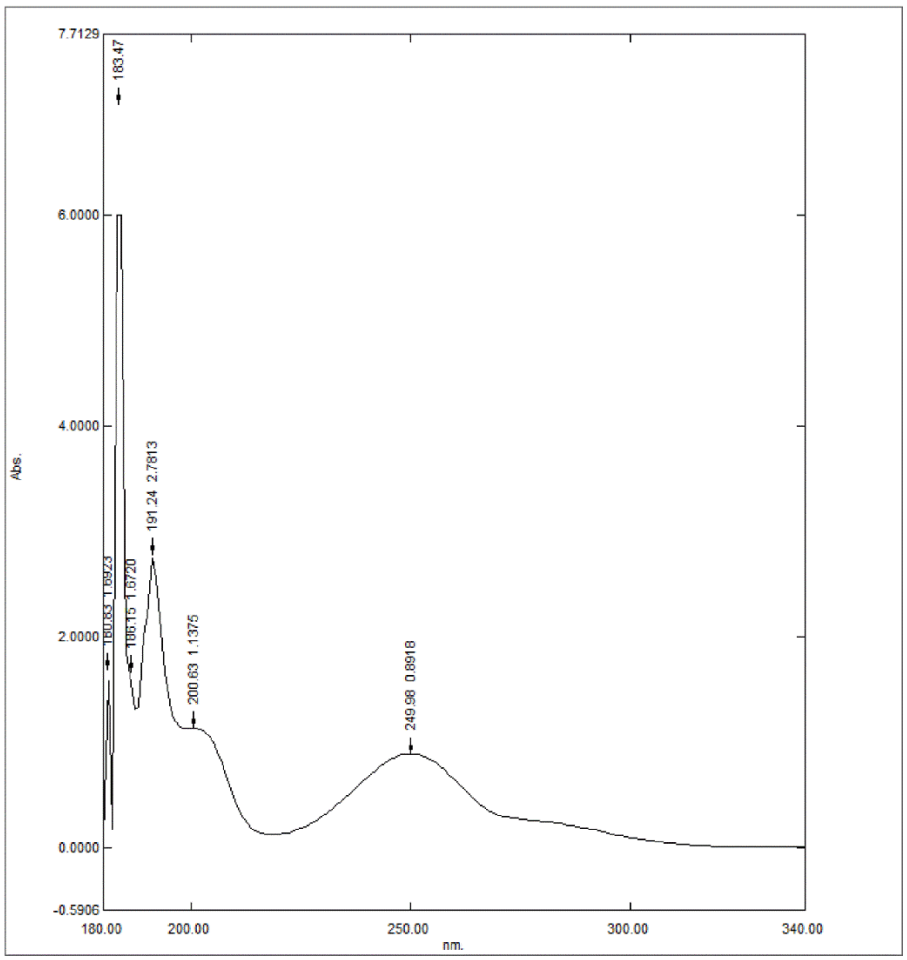

Figure S4: The UV spectrum of 4-[(4-chlorophenyl)carbamoyl]butanoic acid (**1**). The optical absorption spectrum was obtained from a 4.14 × 10^{-5} M acetonitrile solution of (**1**) in the range 185–340 nm on a Shimadzu UV-3600 plus UV/VIS/NIR (Shimadzu Corporation, Kyoto Prefecture, Japan) spectrophotometer.