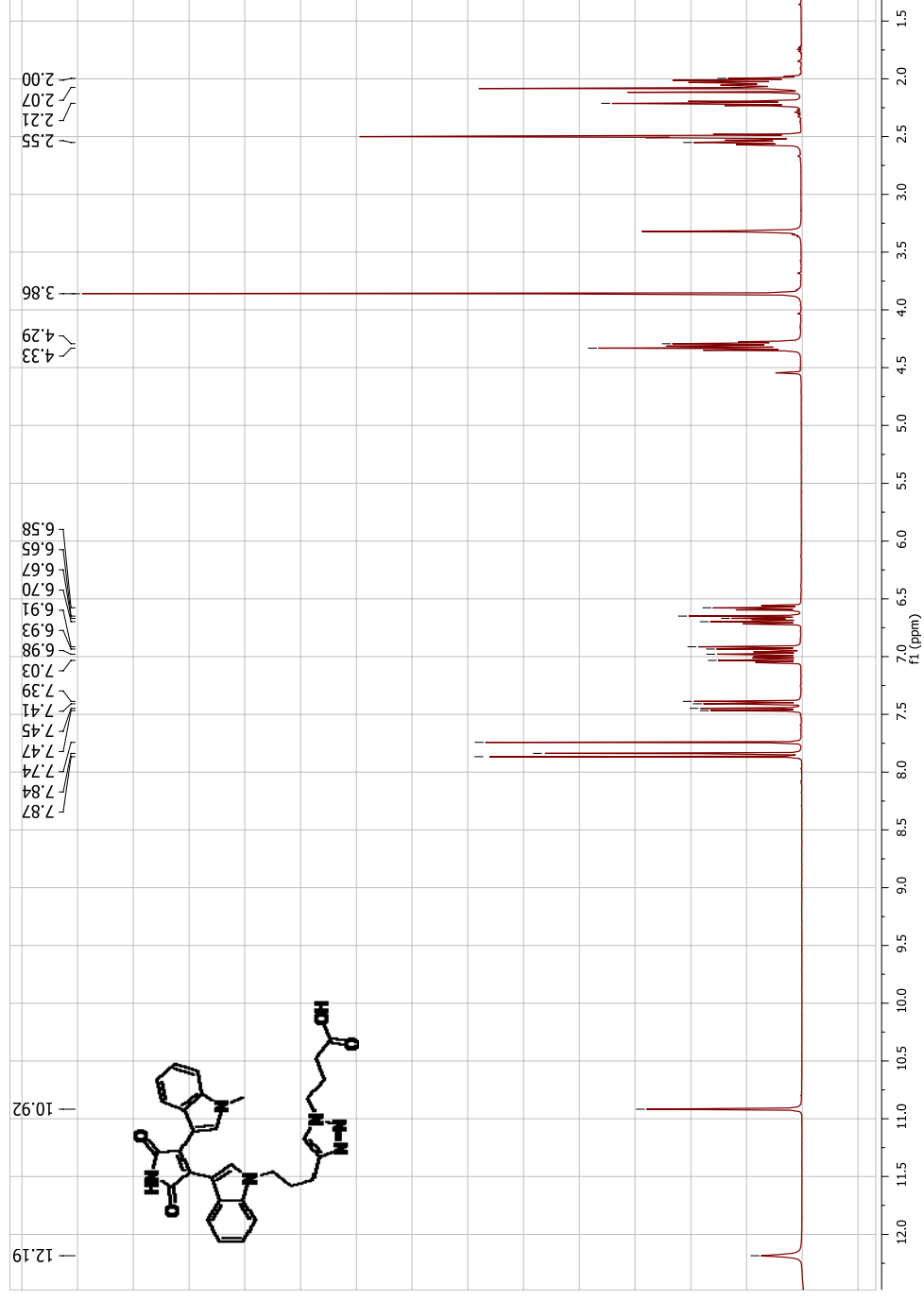
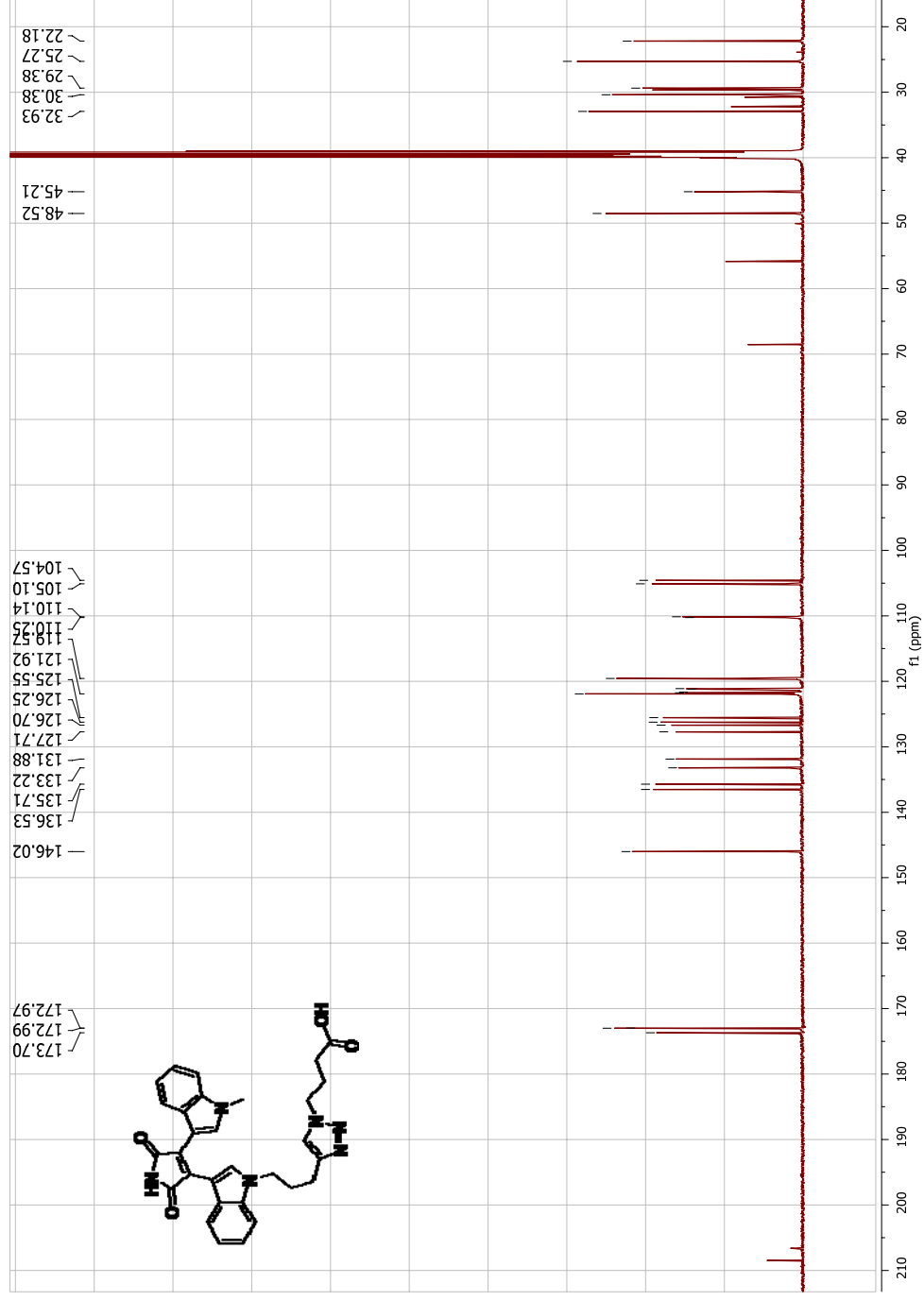


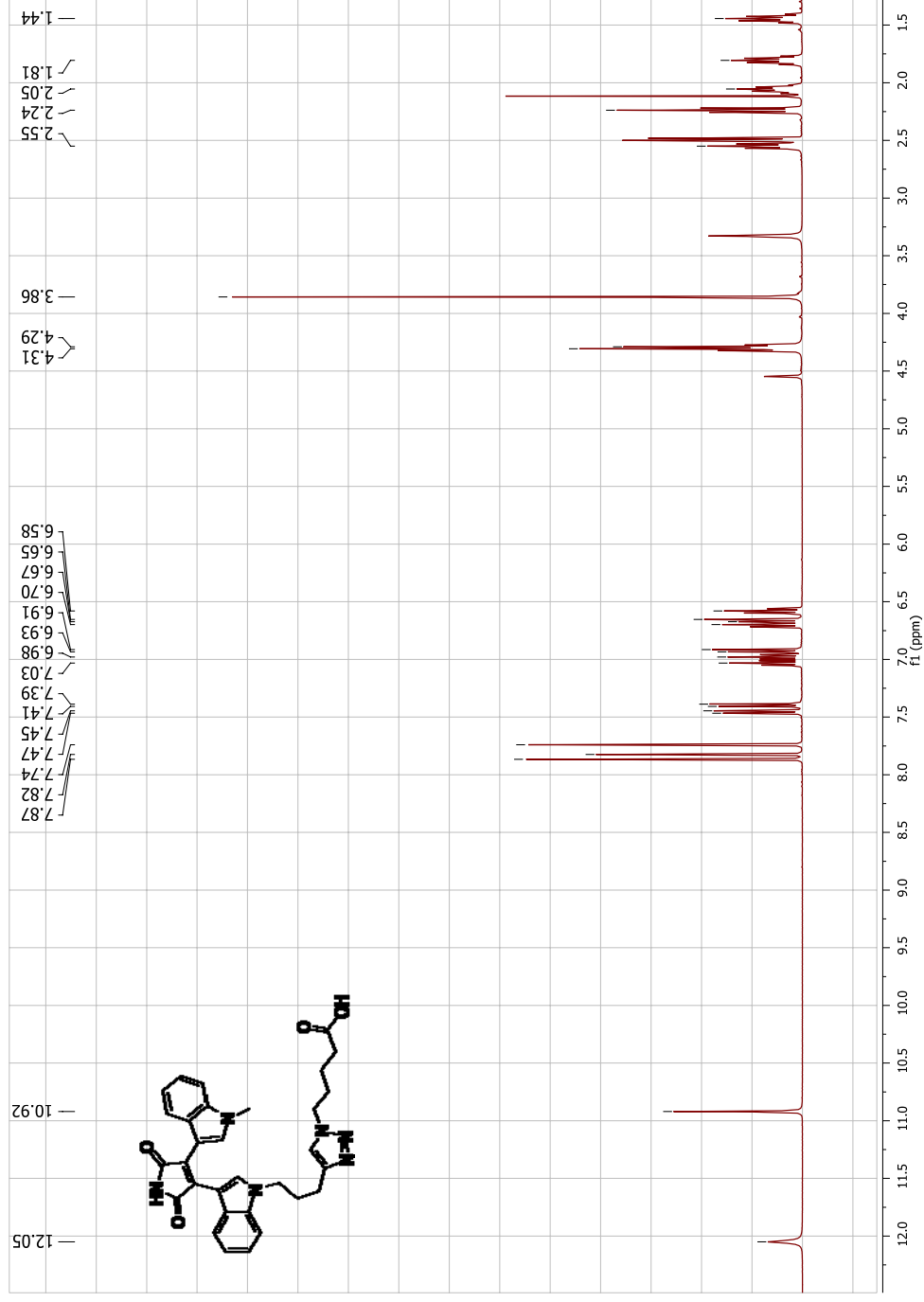
¹H NMR of 4-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-(1,2,3-triazol-1-yl)-butyric acid (36)



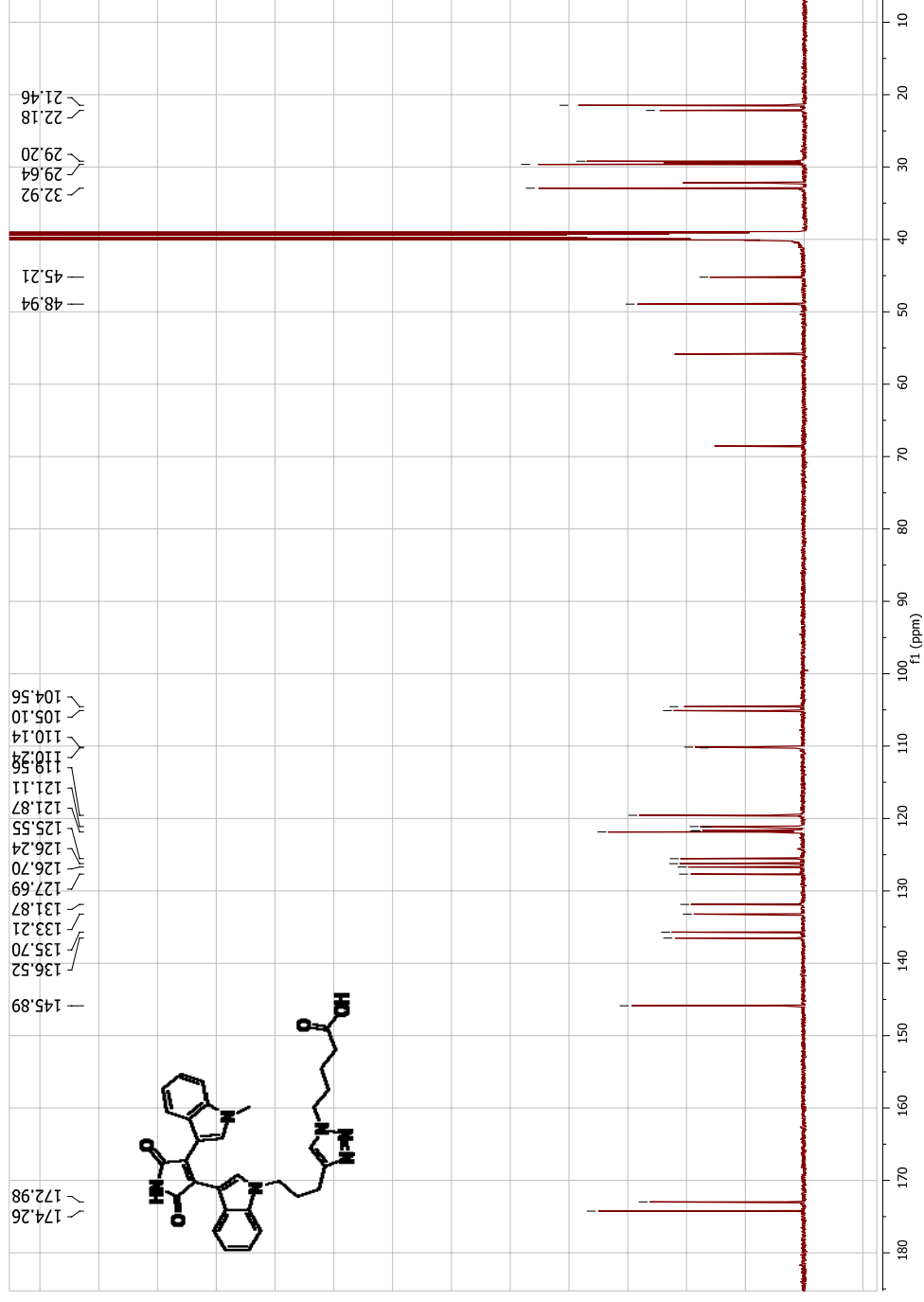
¹³C NMR of 4-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-(1,2,3-triazol-1-yl)-butyric acid (36)



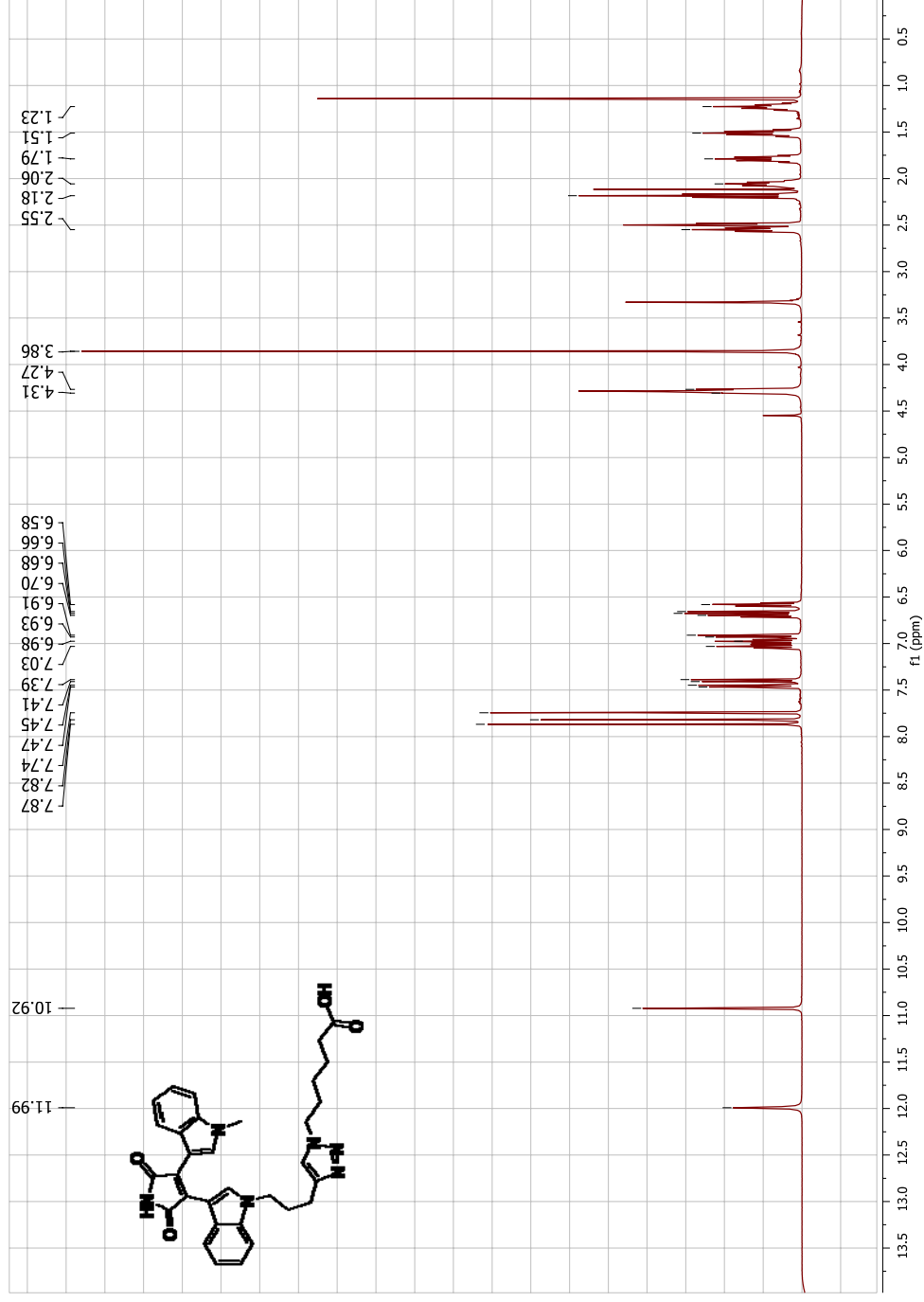
**¹H NMR of 5-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-
-(1,2,3)triazol-1-yl)-pentanoic acid (37)**



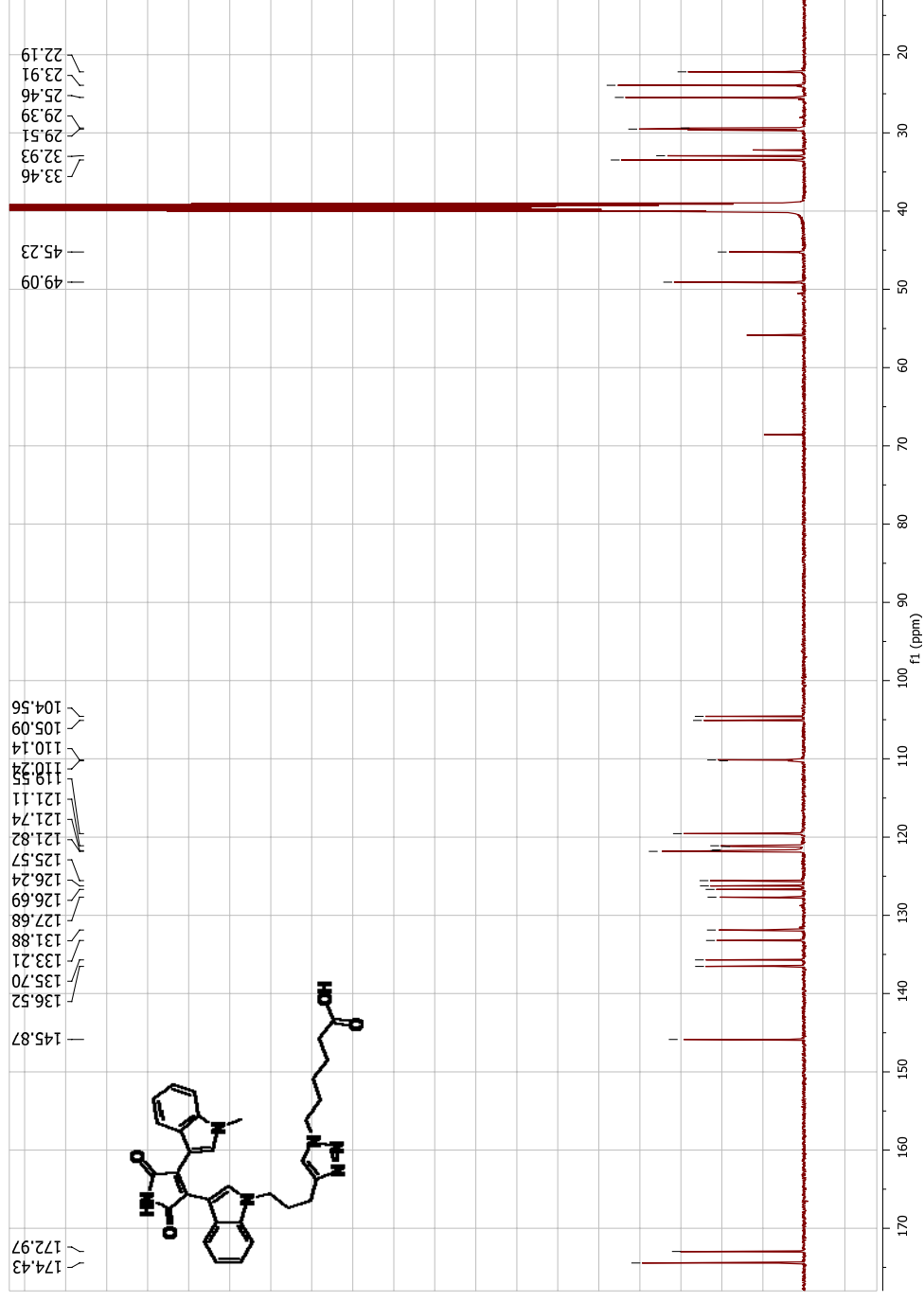
**¹³C NMR of 5-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-
-(1,2,3)triazol-1-yl)-pentanoic acid (37)**



**¹H NMR of 6-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-
-(1,2,3)triazol-1-yl)-hexanoic acid (38)**



**¹³C NMR of 6-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-
-(1,2,3)triazol-1-yl)-hexanoic acid (38)**



¹H NMR spectrum (DMSO-d₆) of compound 10.

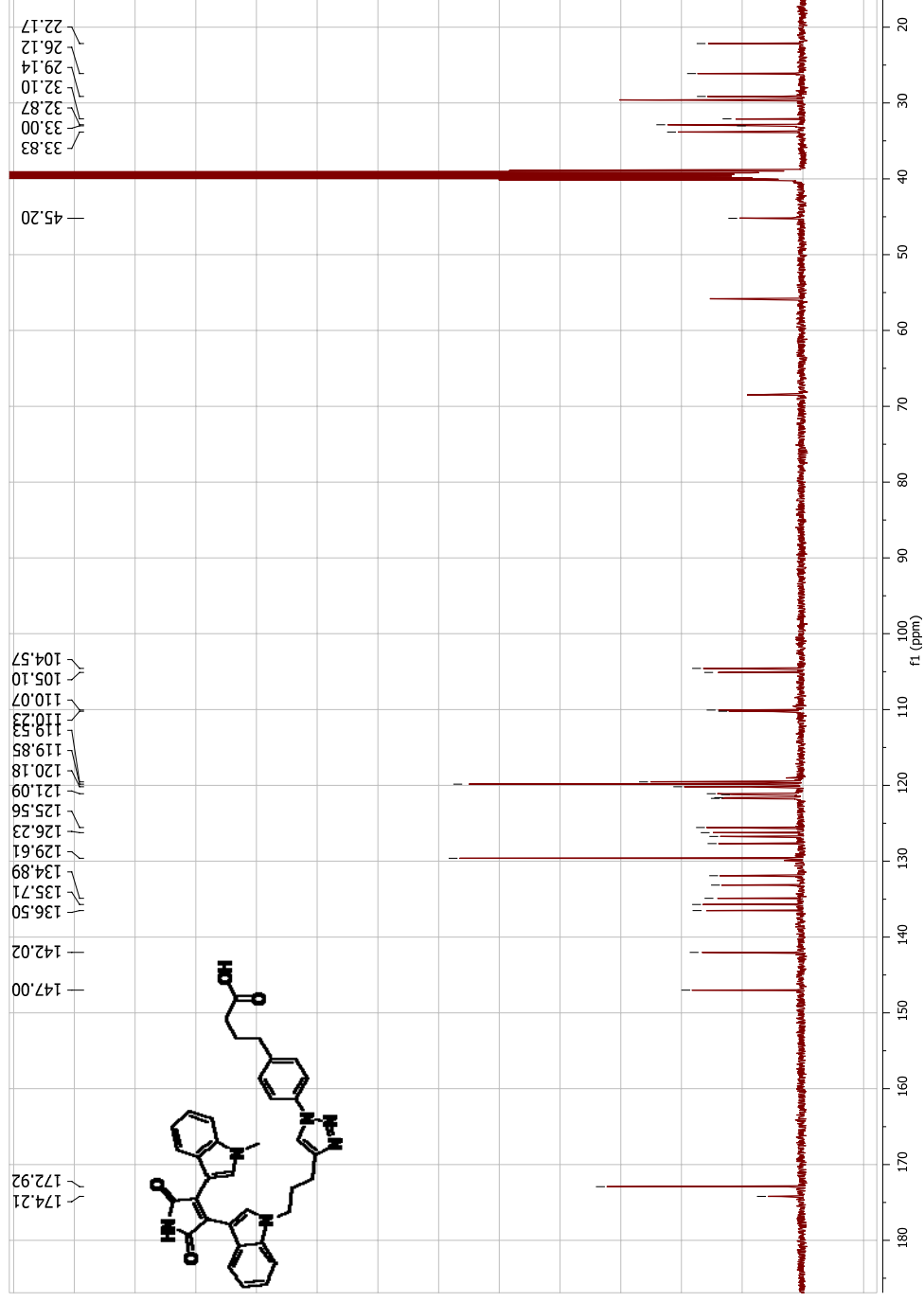
Chemical structure of compound 10:

O=C(O)CCc1ccc(N2C=NC(CCC3C4C=CC=CC=C4C(=O)NC3=O)CC2)cc1

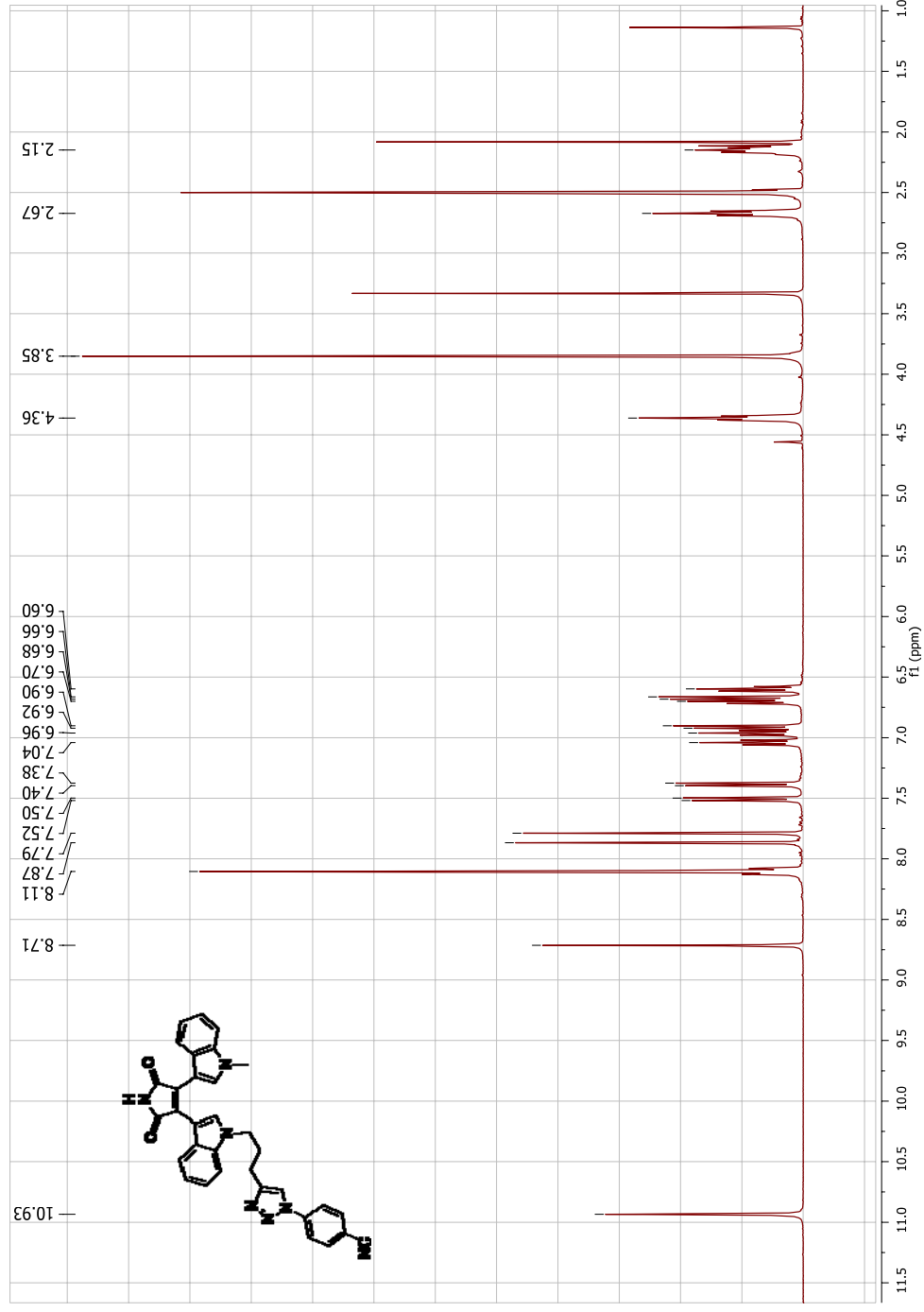
Peak list (ppm):

Chemical Shift (ppm)
12.11
10.94
8.49
7.87
7.79
7.76
7.52
7.50
7.41
7.40
7.39
7.38
7.04
6.97
6.93
6.91
6.70
6.69
6.67
6.60
4.35
3.85
3.50
2.67
2.65
2.24
2.14
1.83

¹³C NMR of 4-(4-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-(1,2,3)triazol-1-yl)-phenyl)-butyric acid (39)



¹H NMR of 4-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-(1,2,3)triazol-1-yl)-benzonitrile (40)



**¹³C NMR of 4-(4-(3-(4-(1-methyl-1*H*-indol-3-yl)-2,5-dioxo-2,5-dihydro-1*H*-pyrrol-3-yl)-indol-1-yl)-propyl)-
-(1,2,3)triazol-1-yl)-benzonitrile (40)**

