SUPPORTING INFORMATION (*New J. Chem*)

Understanding multivalent effects in glycosidase inhibition using *C*-glycoside click clusters as molecular probes

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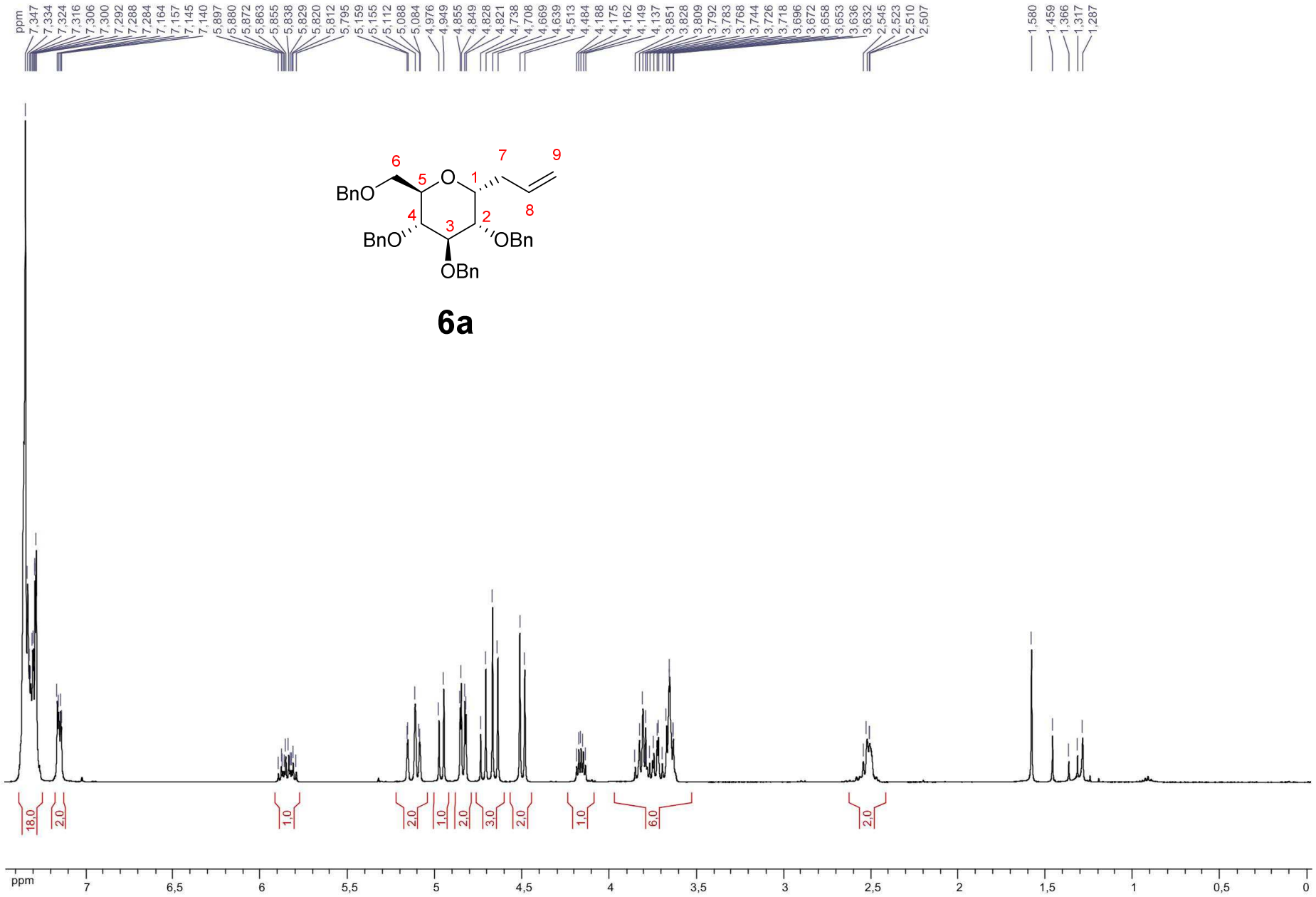
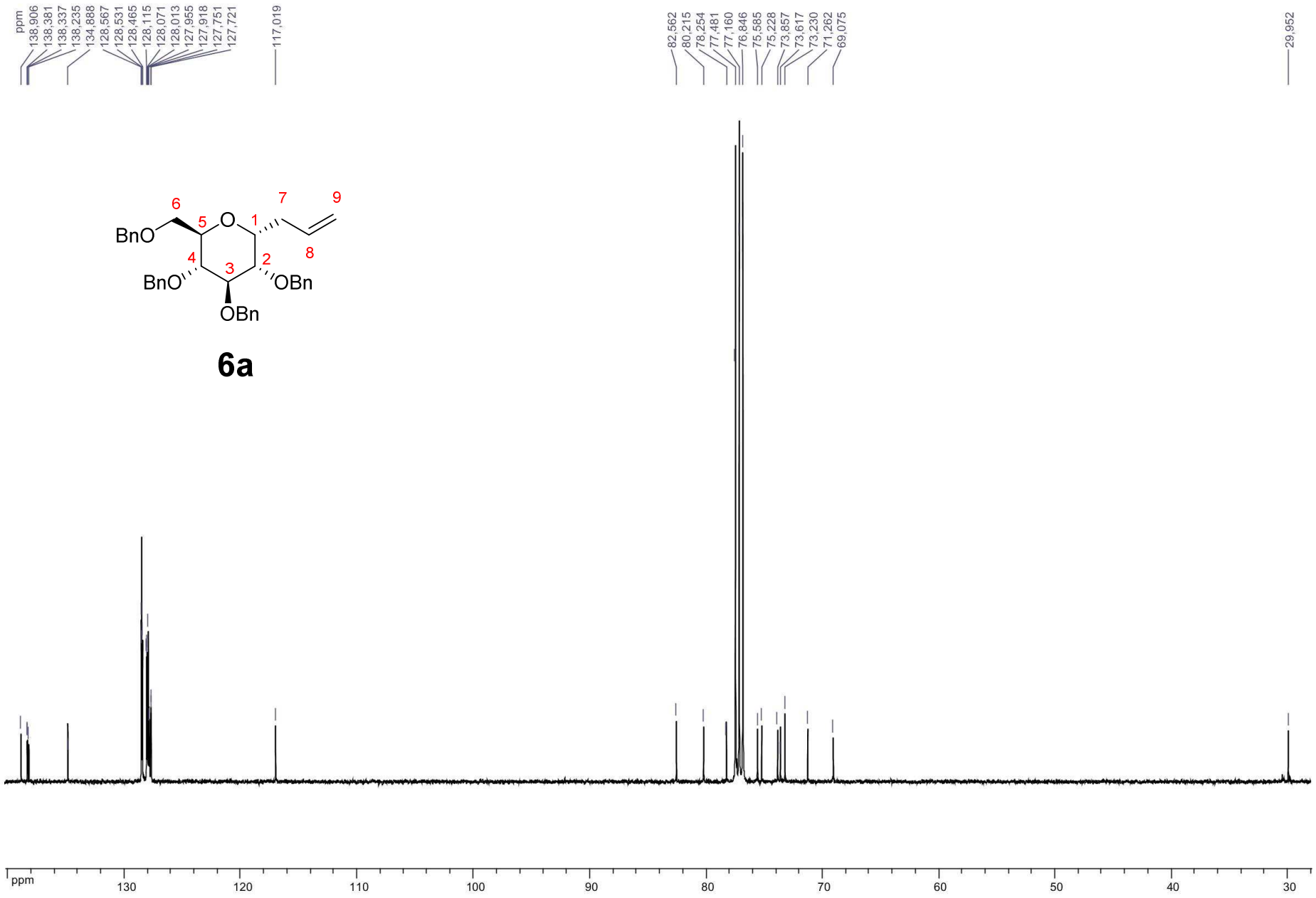
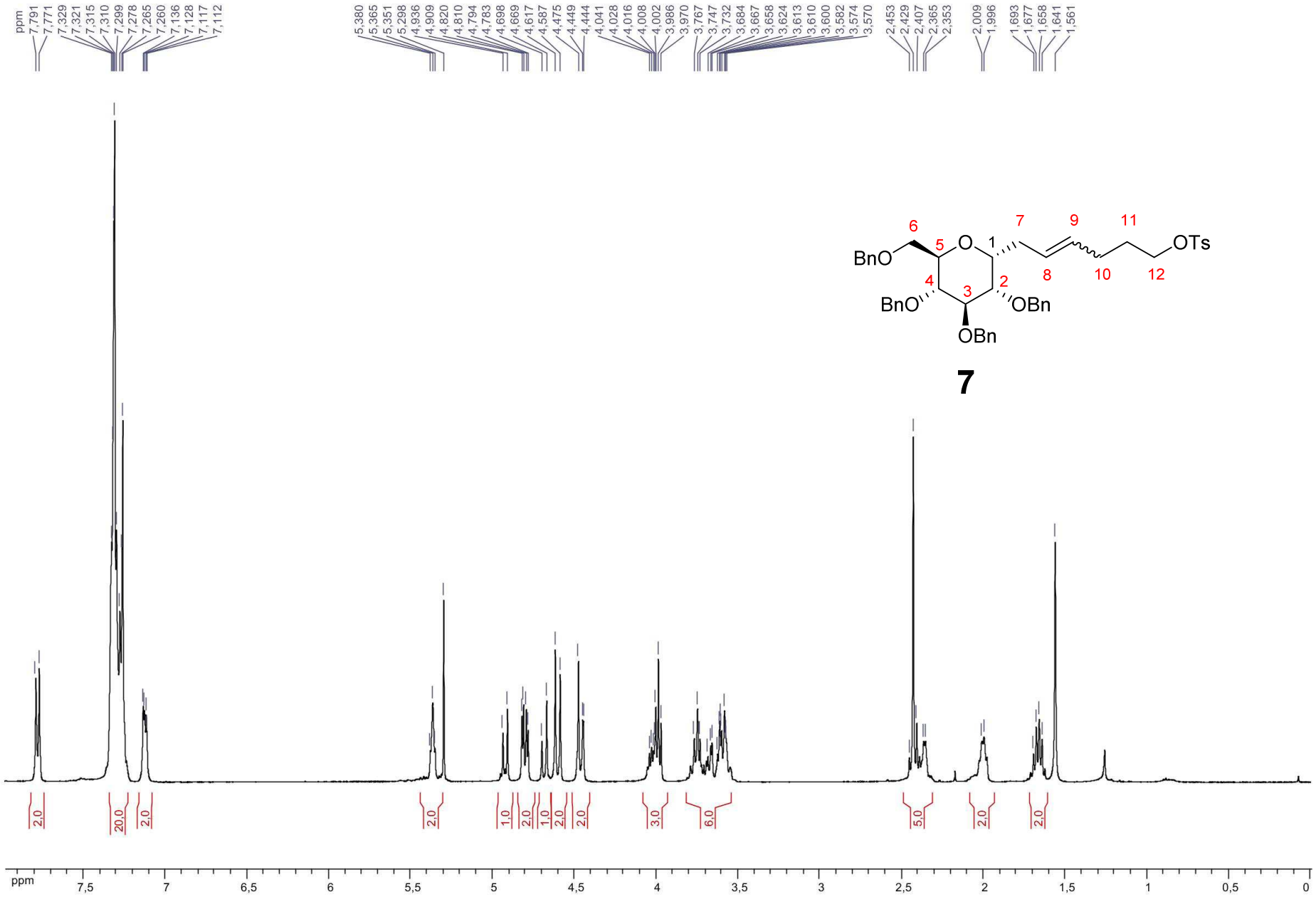
^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 6a:



^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 6a:

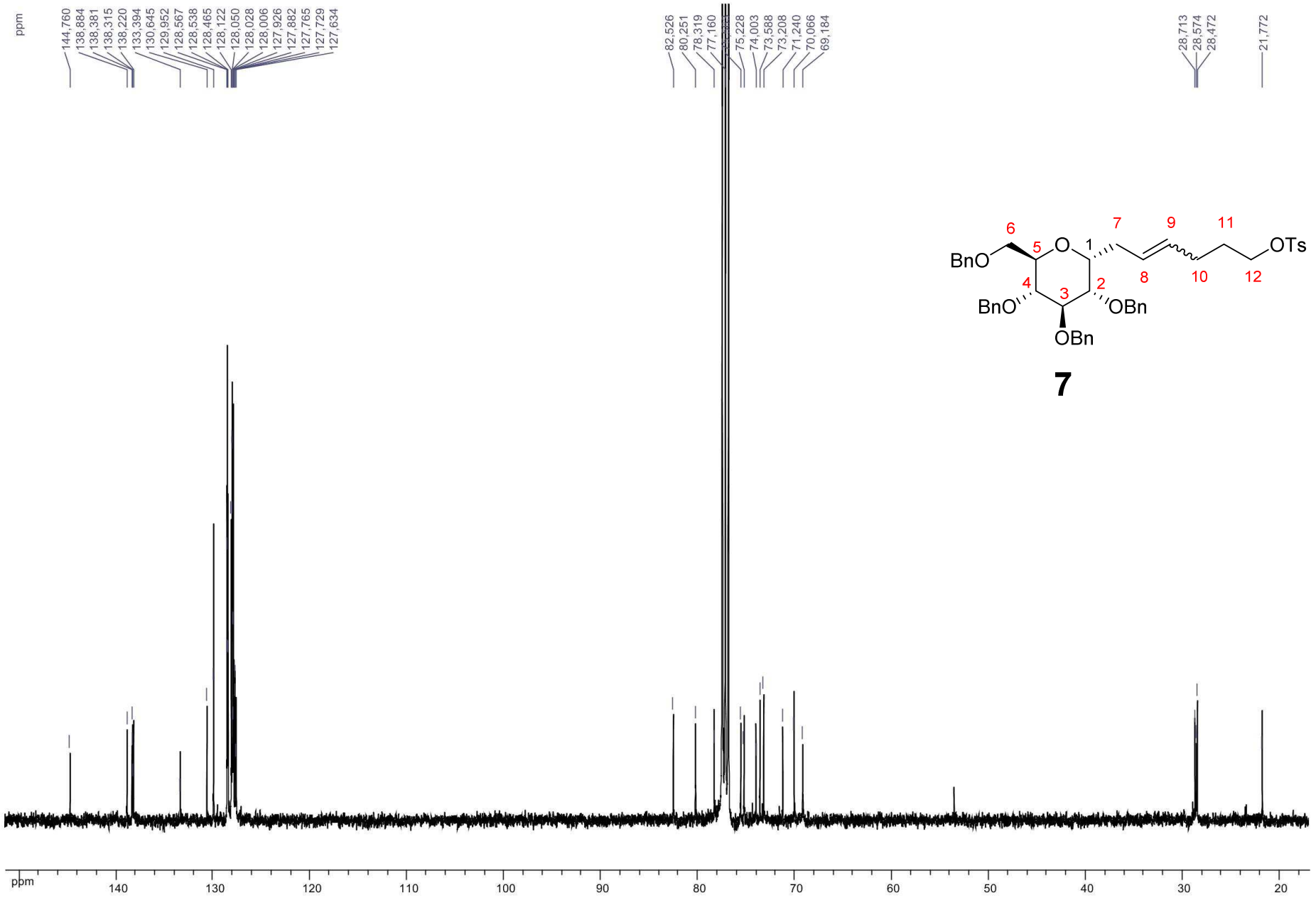


^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 7:

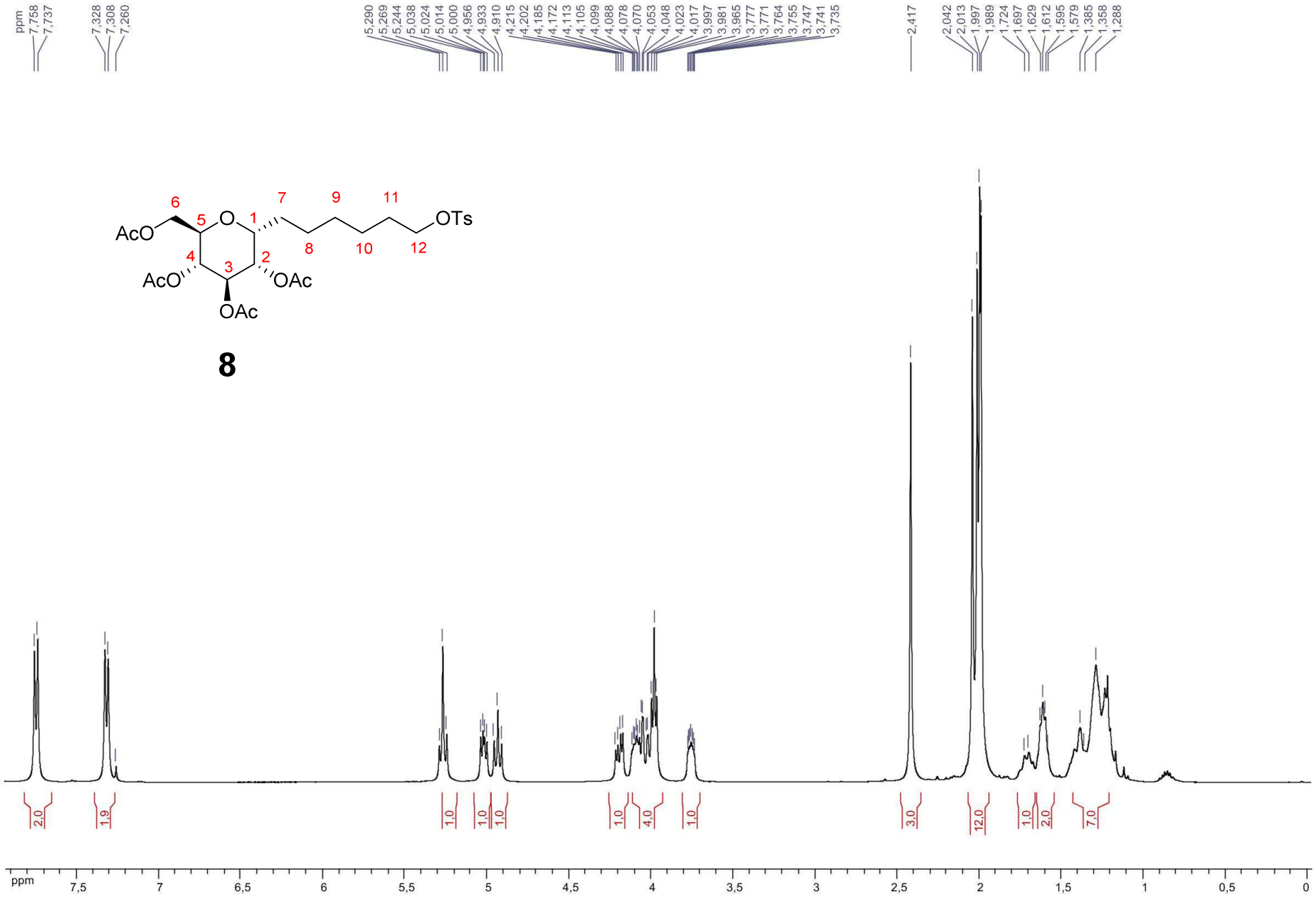


7

^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 7:

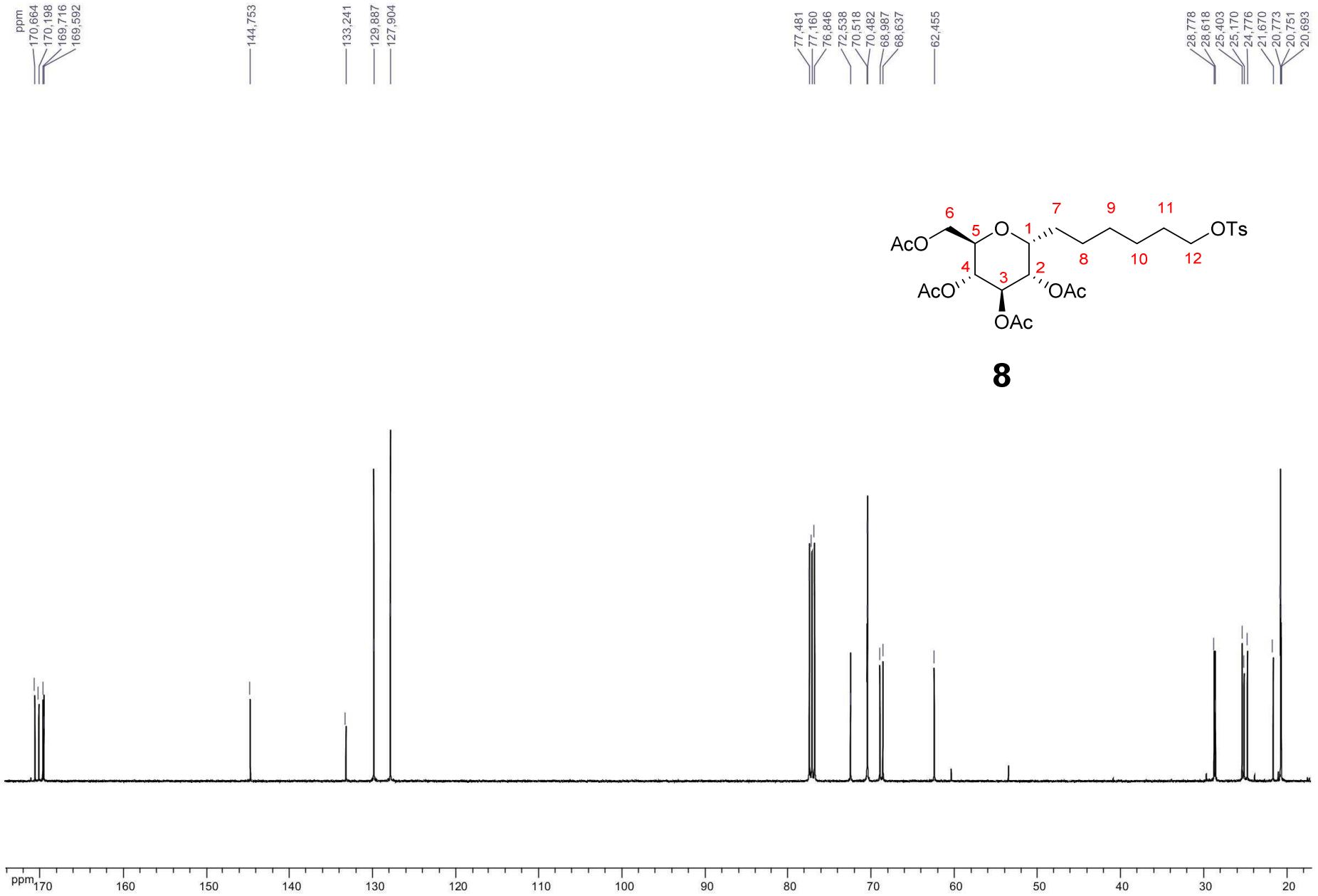


^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 8:



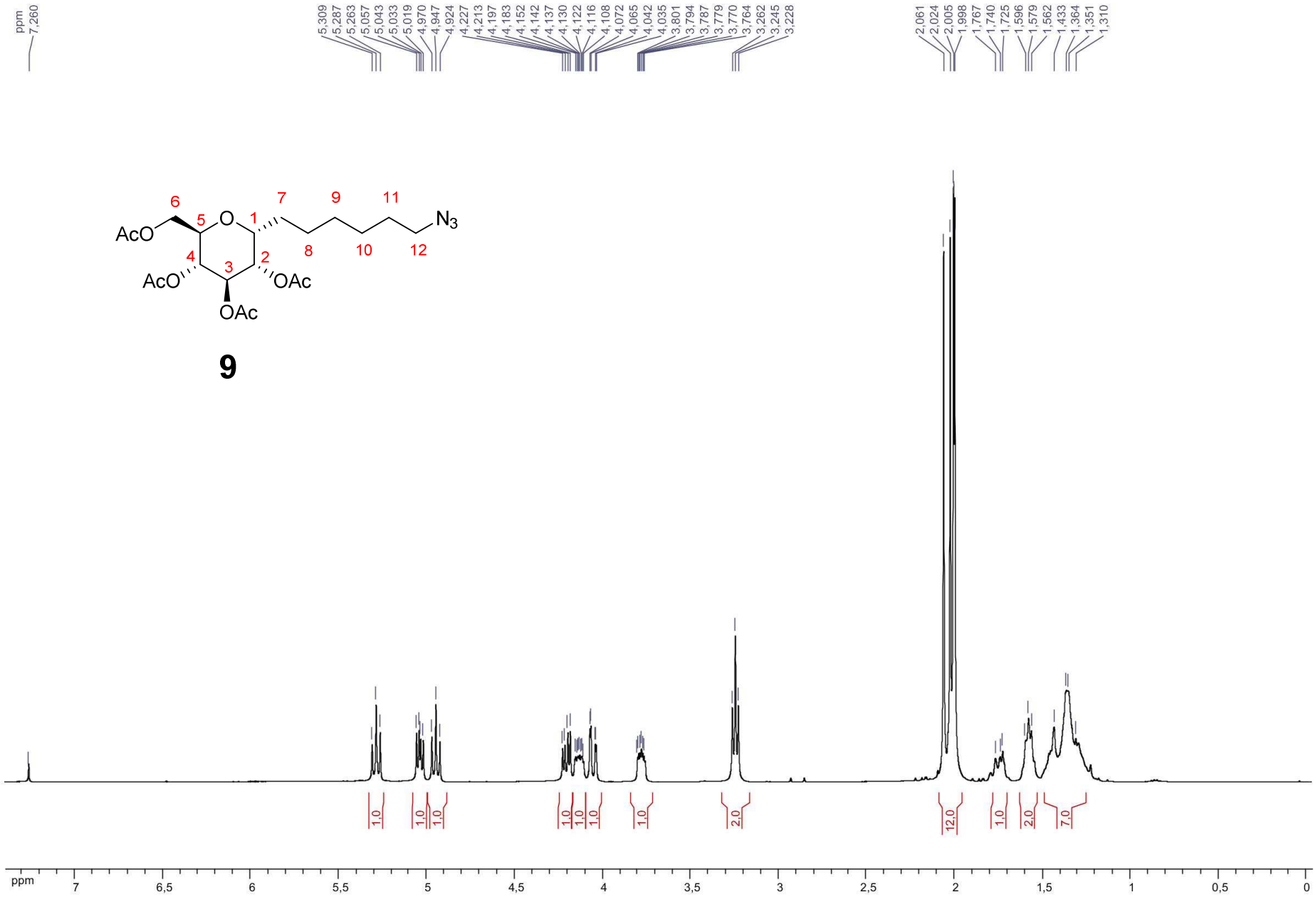
8

^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 8:



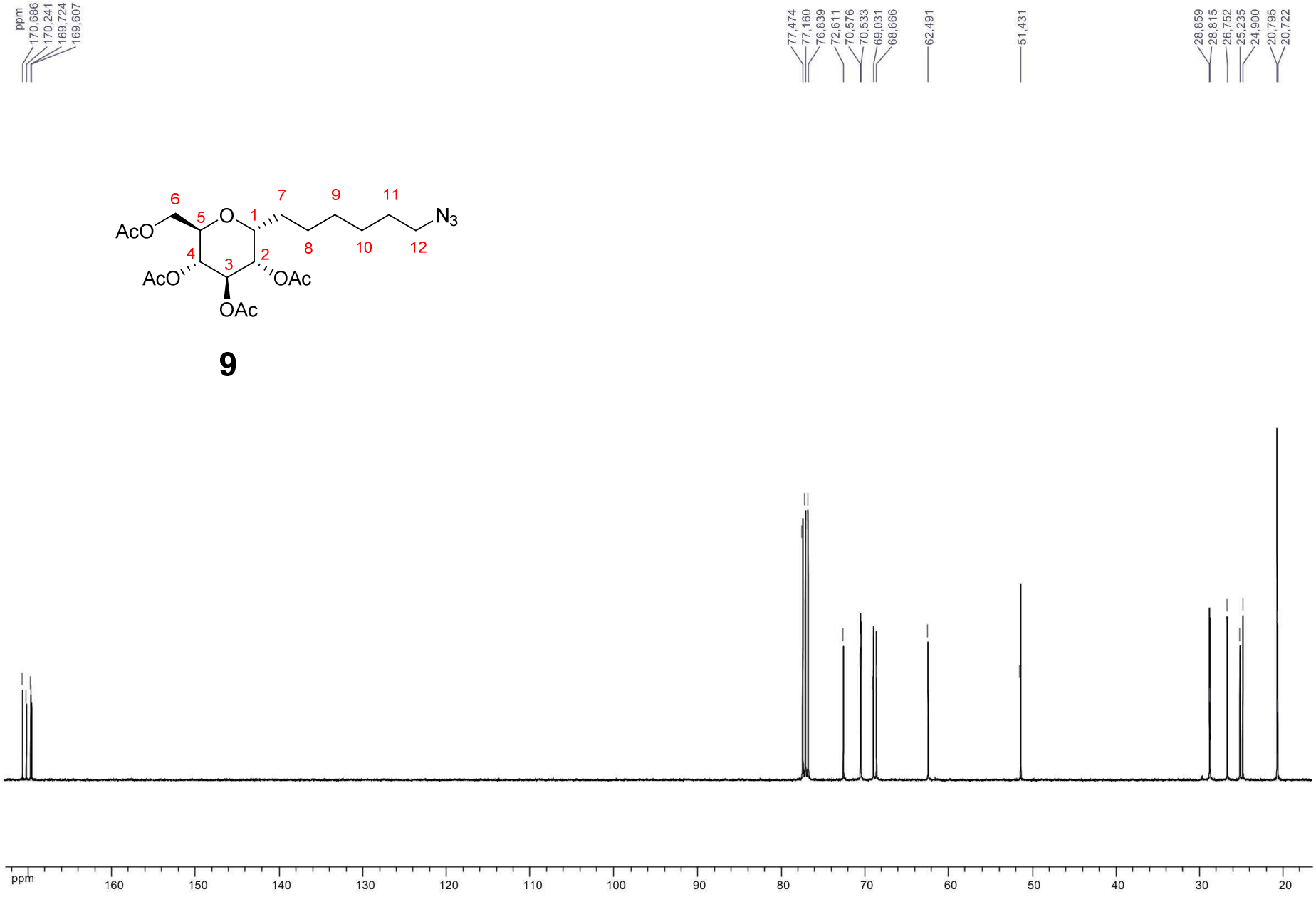
8

^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 9:



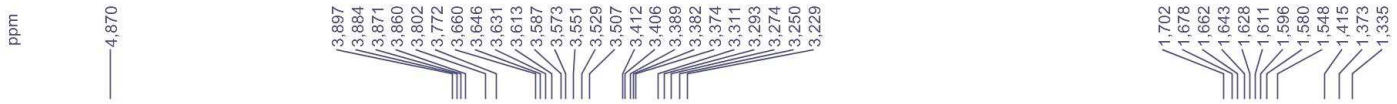
9

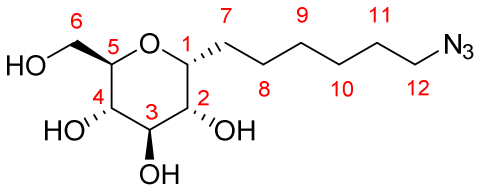
^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 9:



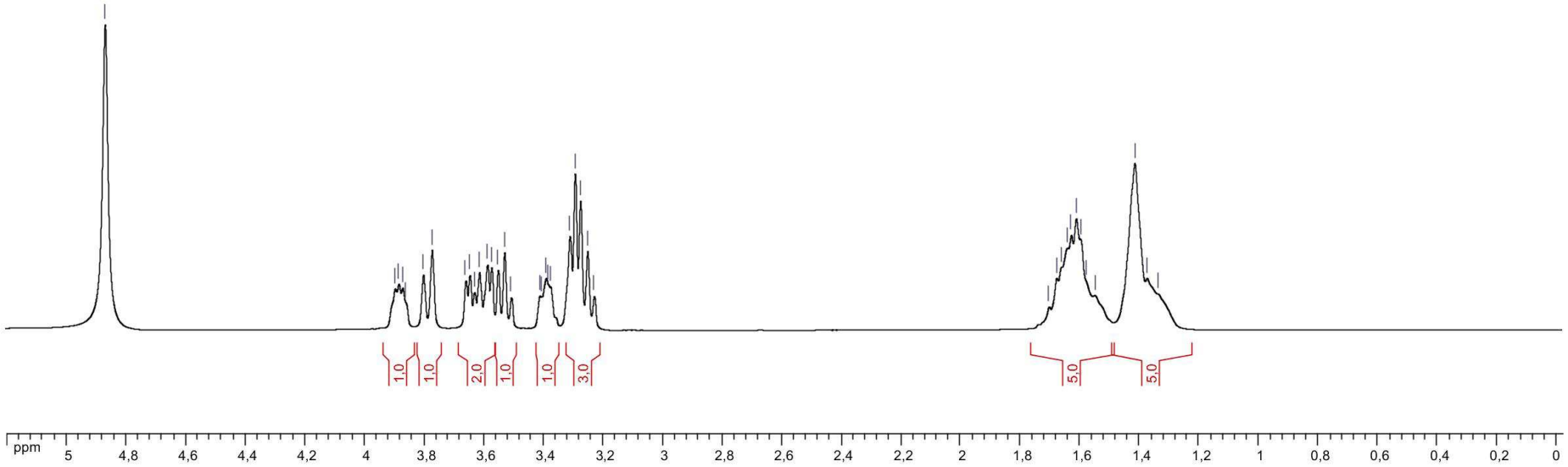
9

^{1}H NMR spectrum (400 MHz, CD_3OD) of compound 10:

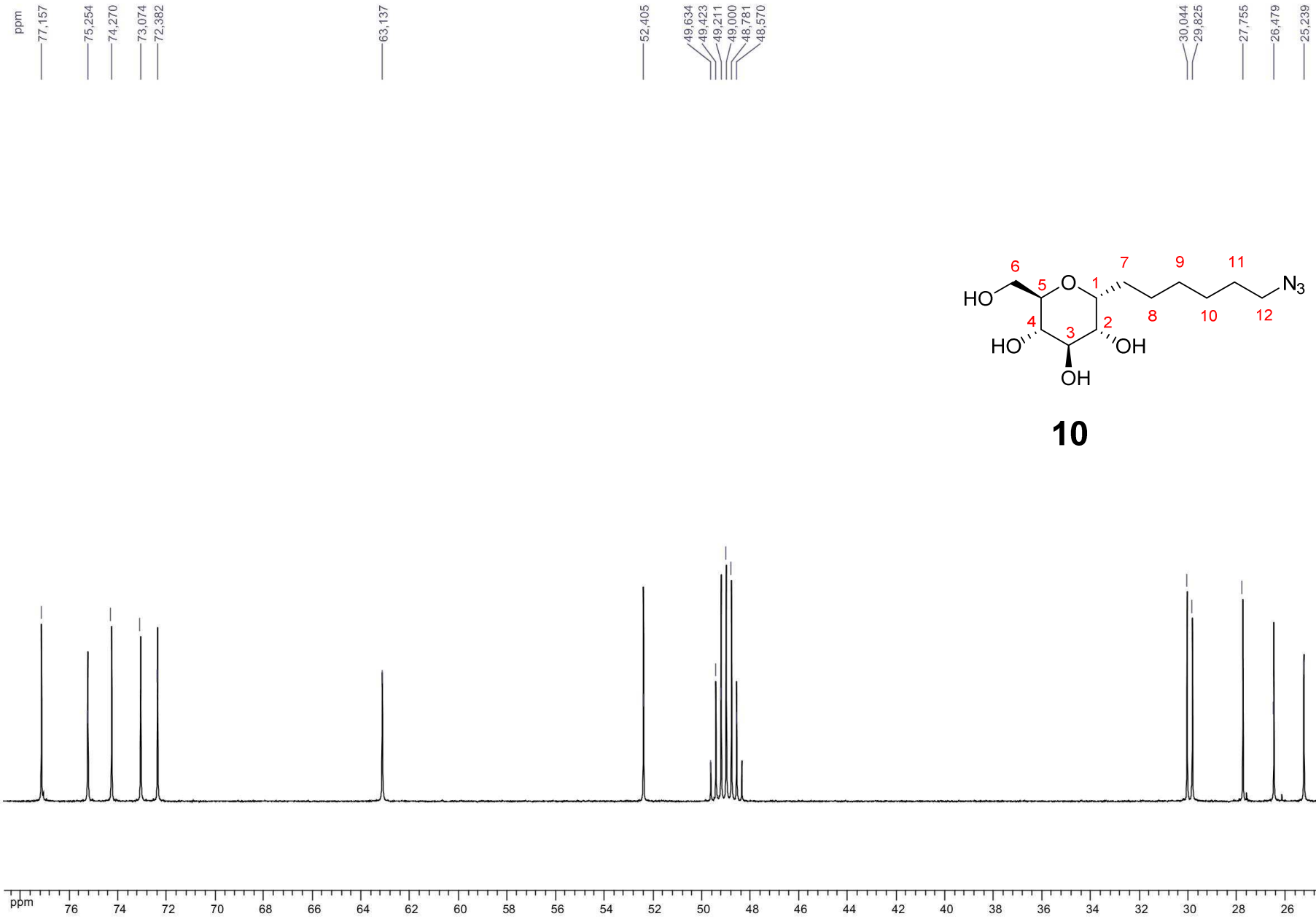




10

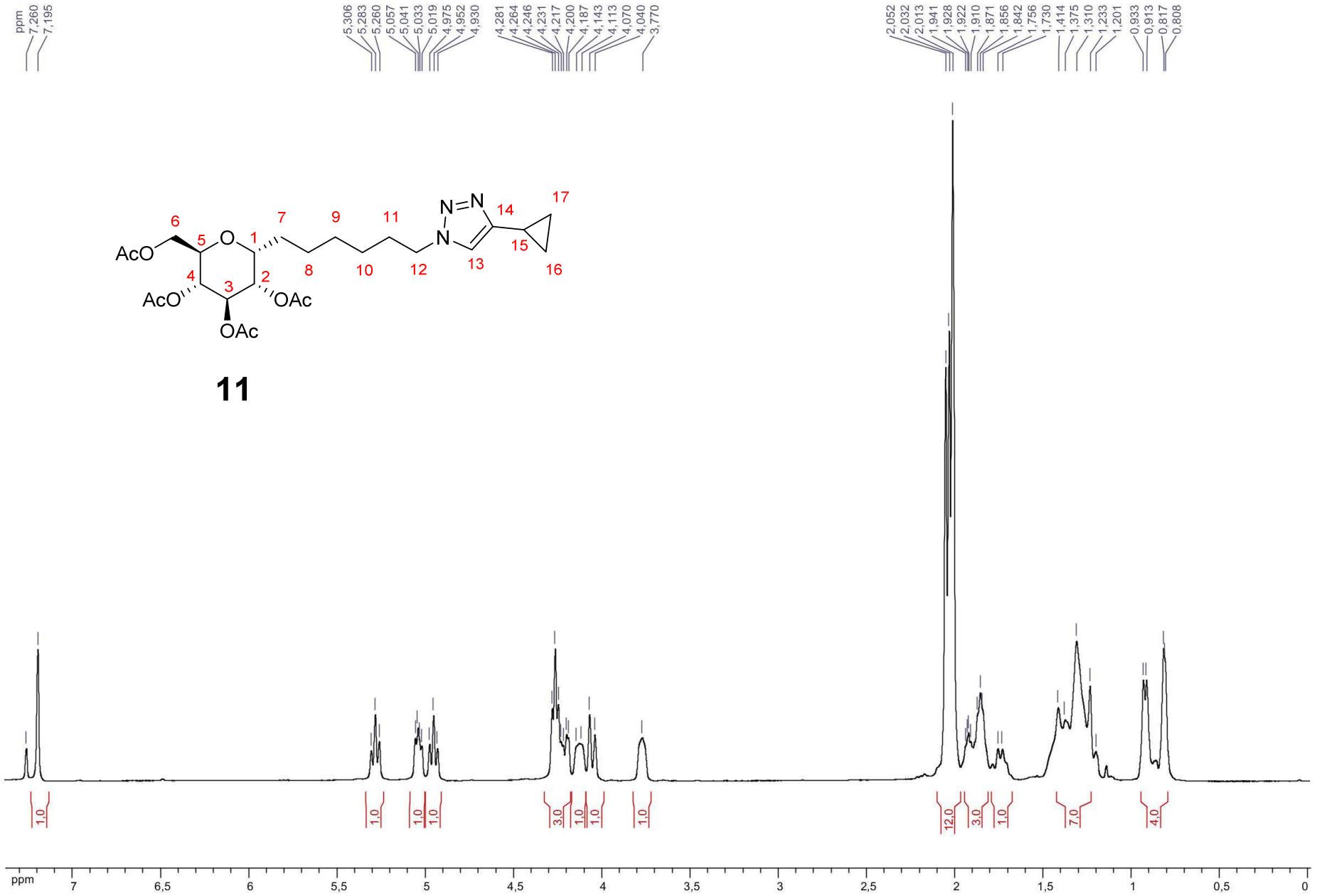


^{13}C NMR spectrum (100 MHz, CD_3OD) of compound 10:



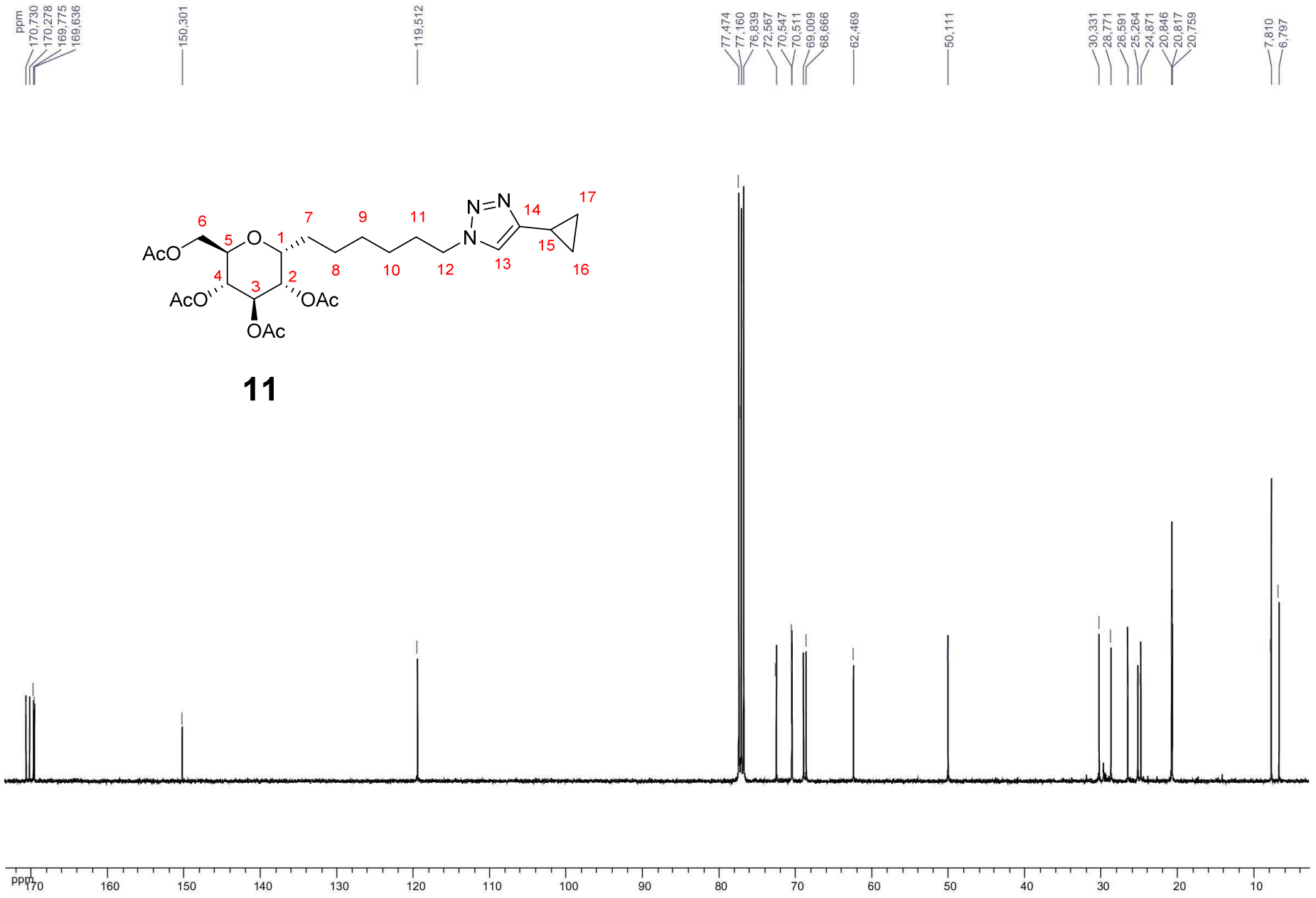
n

^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 11:



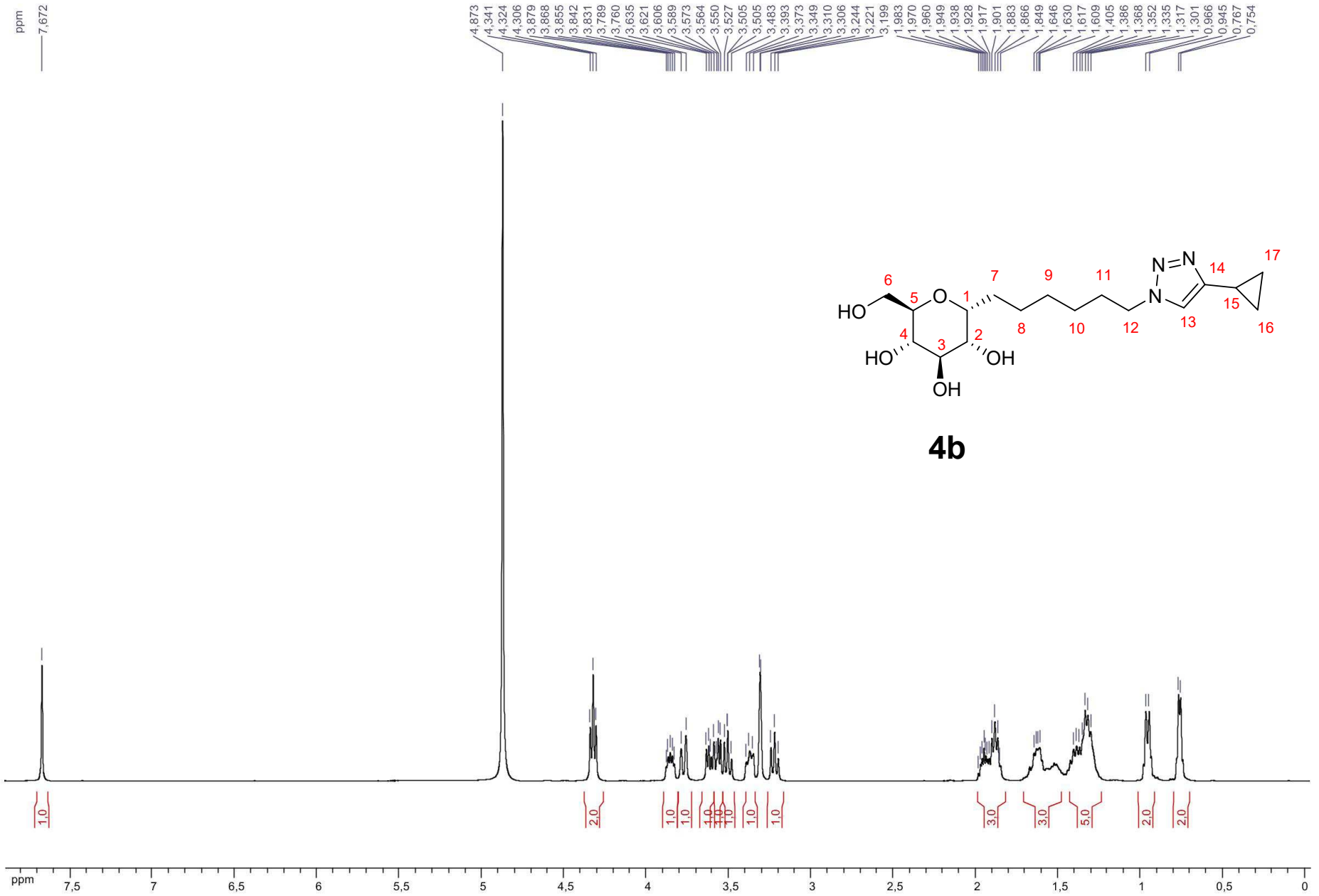
11

^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 11:



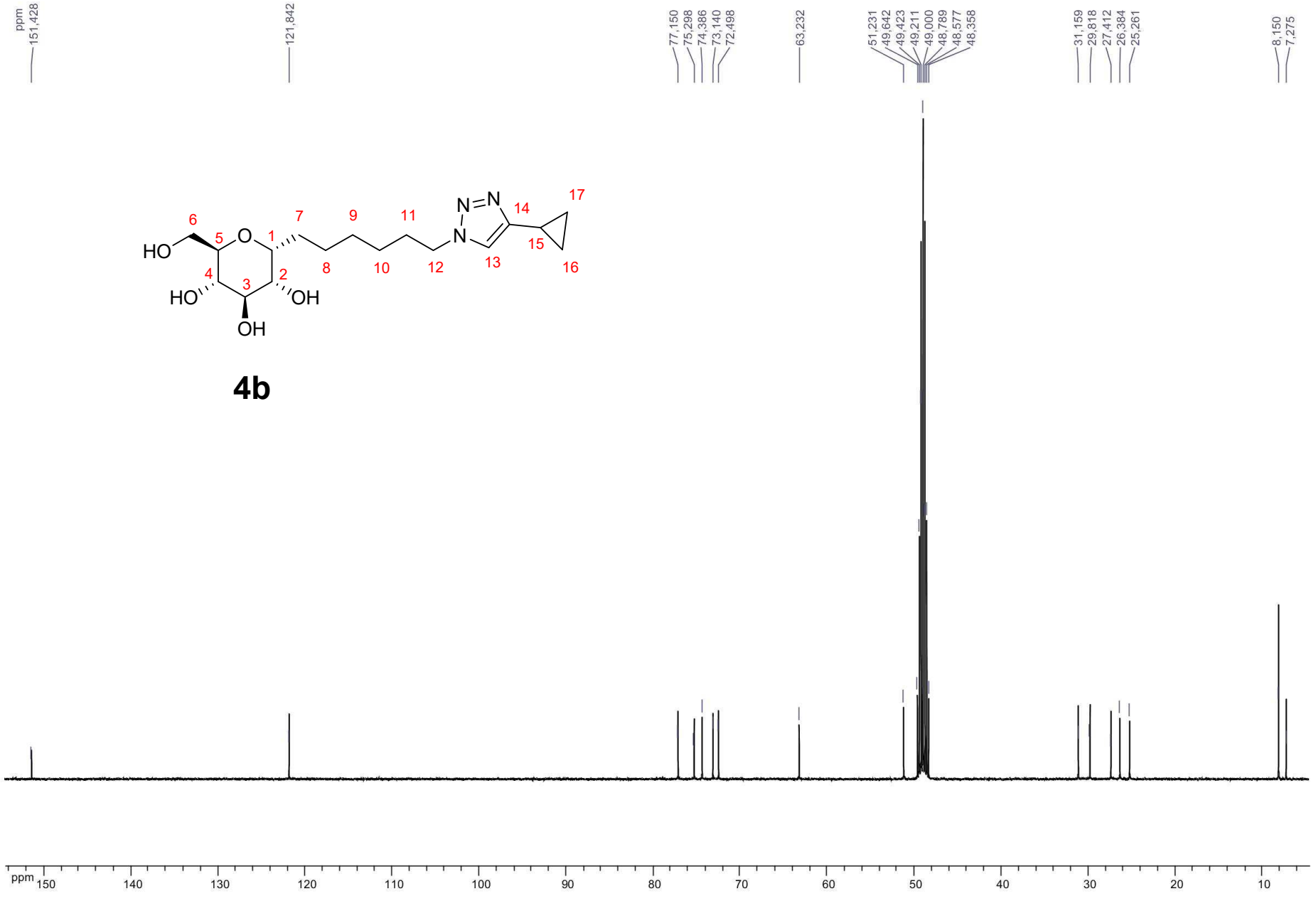
11

^{1}H NMR spectrum (400 MHz, CD_3OD) of compound 4b:



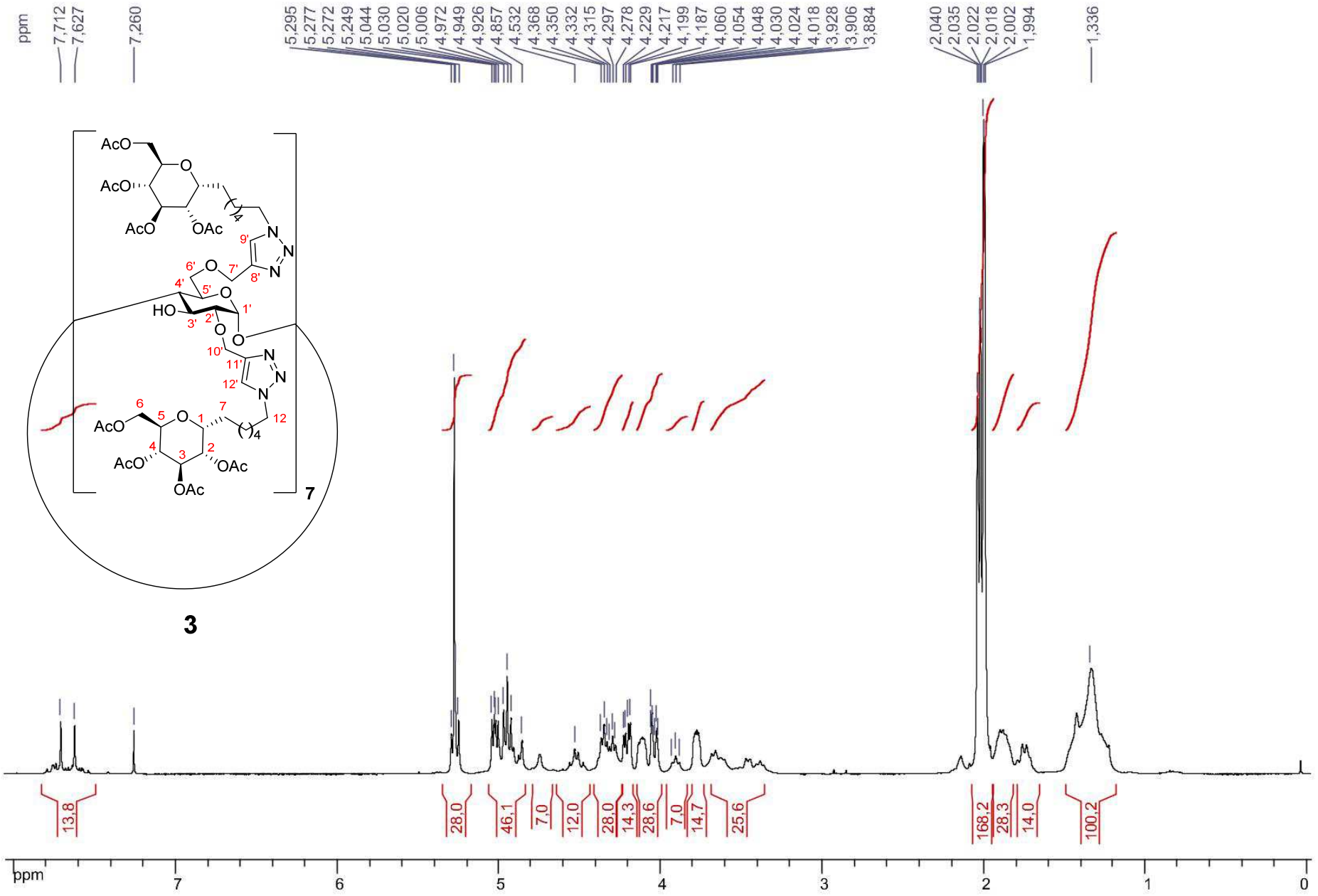
4b

^{13}C NMR spectrum (100 MHz, CD_3OD) of compound 4b:

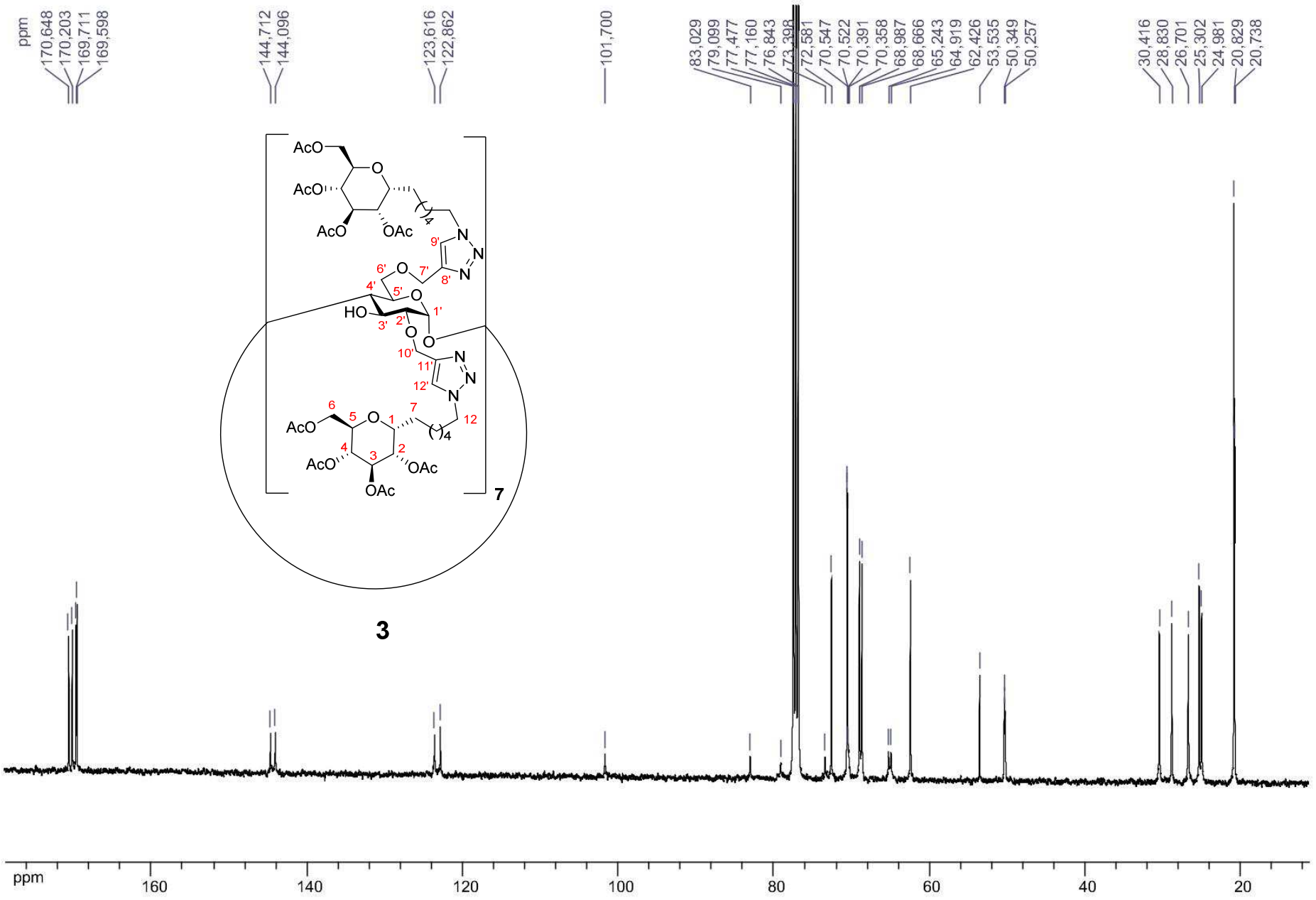


4b

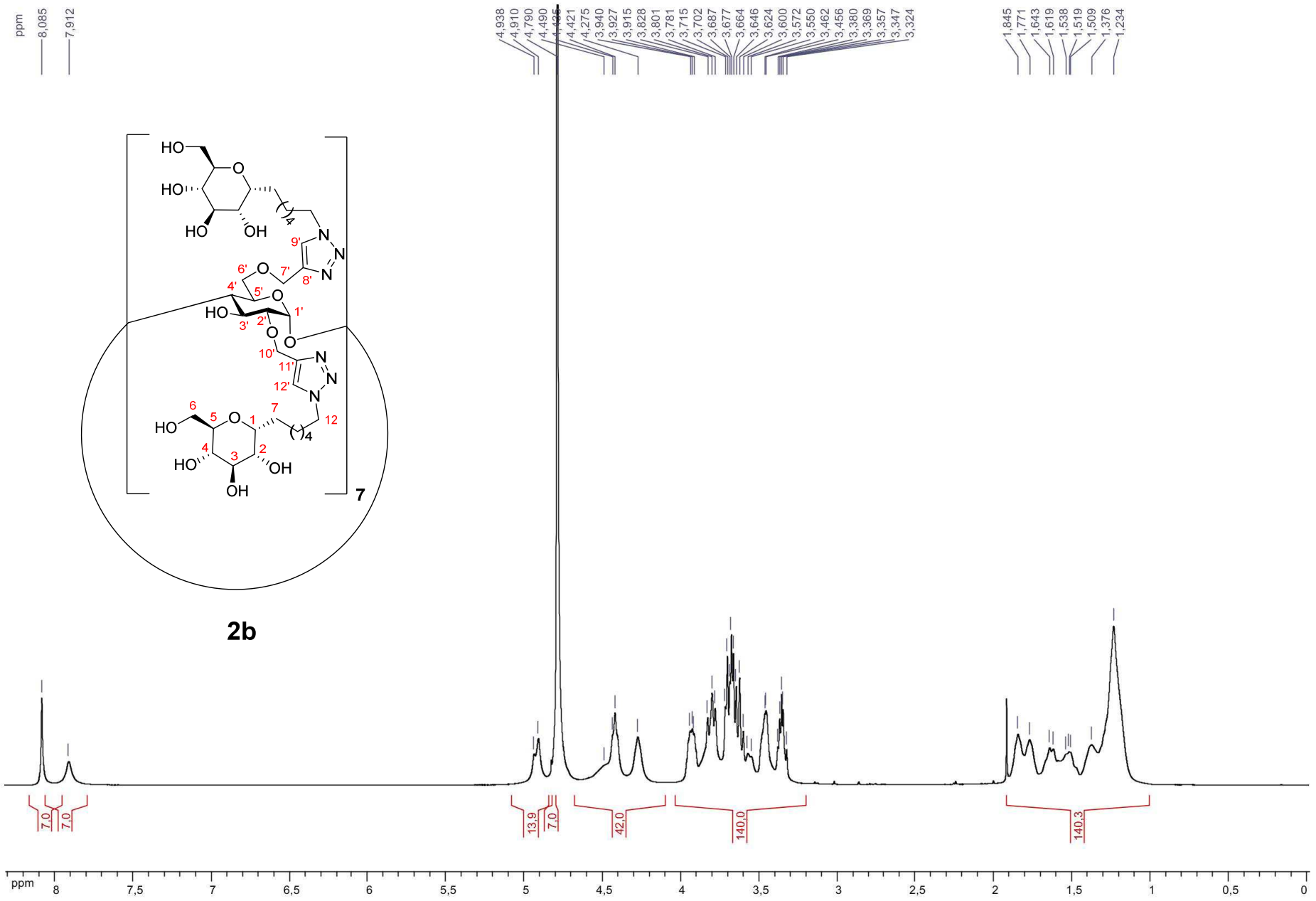
^{1}H NMR spectrum (400 MHz, $CDCl_3$) of compound 3:



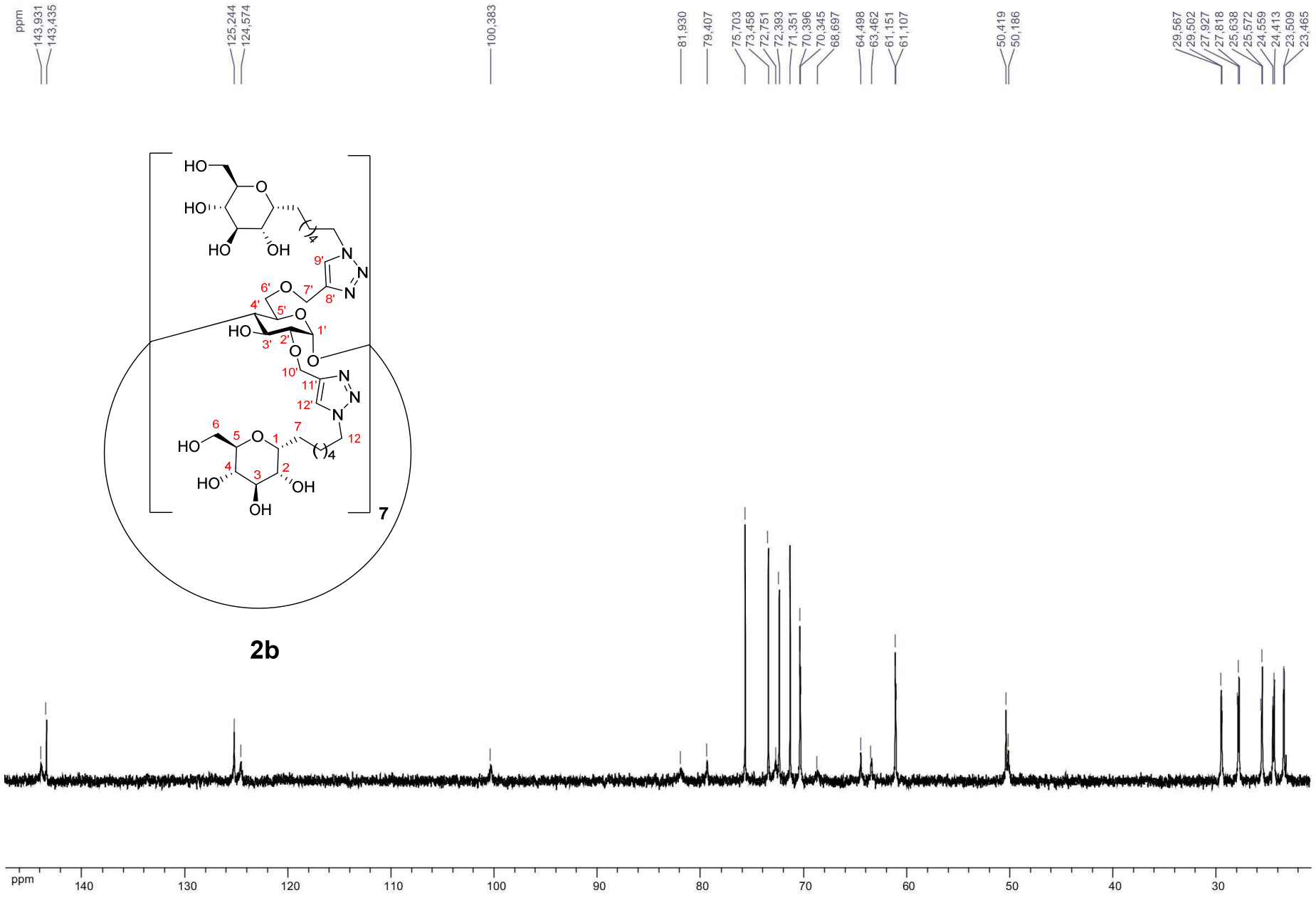
^{13}C NMR spectrum (100 MHz, $CDCl_3$) of compound 3:



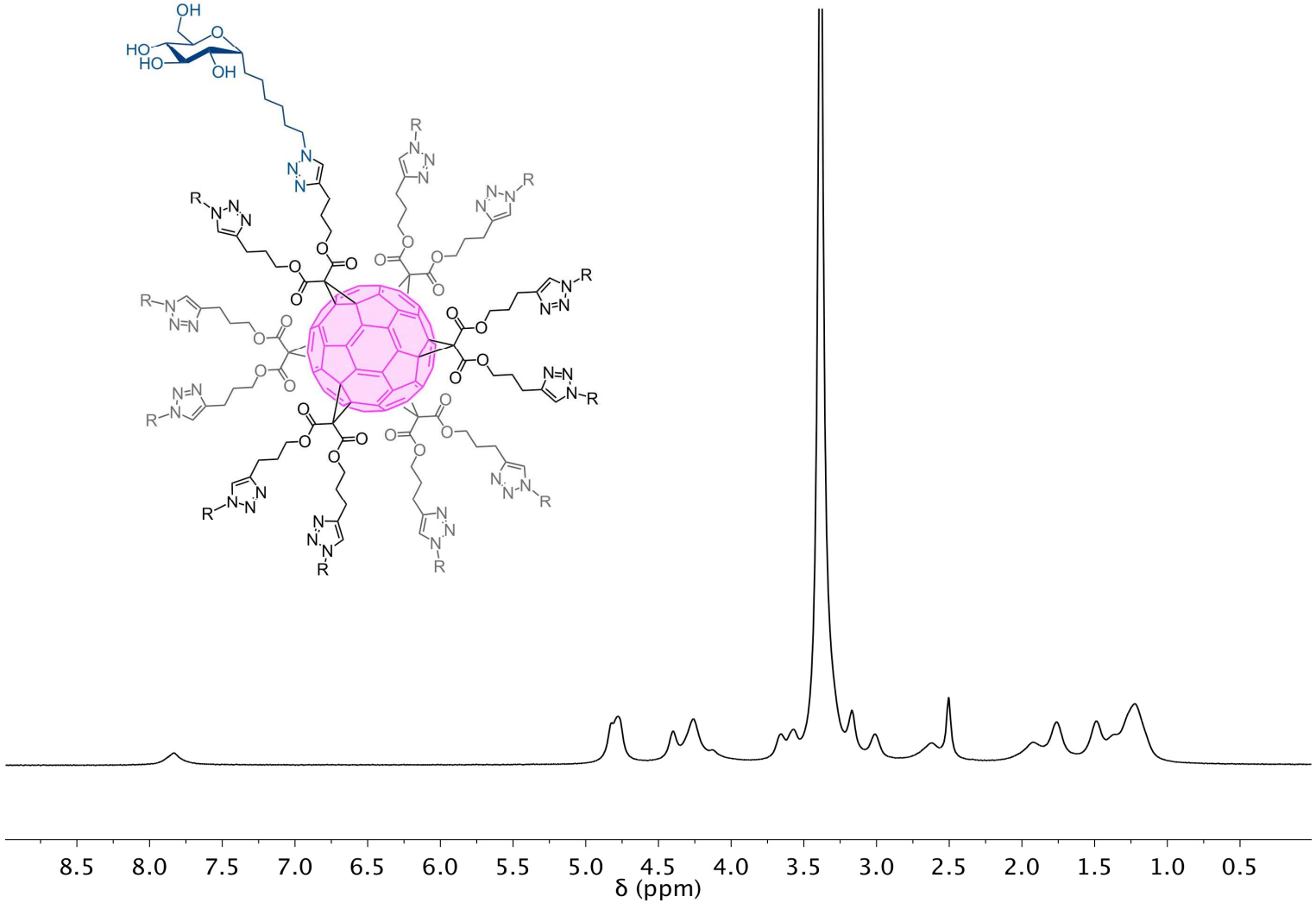
^{1}H NMR spectrum (300 MHz, D_2O) of compound 2b:



^{13}C NMR spectrum (75.5 MHz, D_2O) of compound 2b:



^{1}H NMR spectrum (400 MHz, D_2O/DMSO-d_6, 4:1) of compound 2a:



^{13}C NMR spectrum (100 MHz, D_2O/DMSO-d_6, 4:1) of compound 2a:

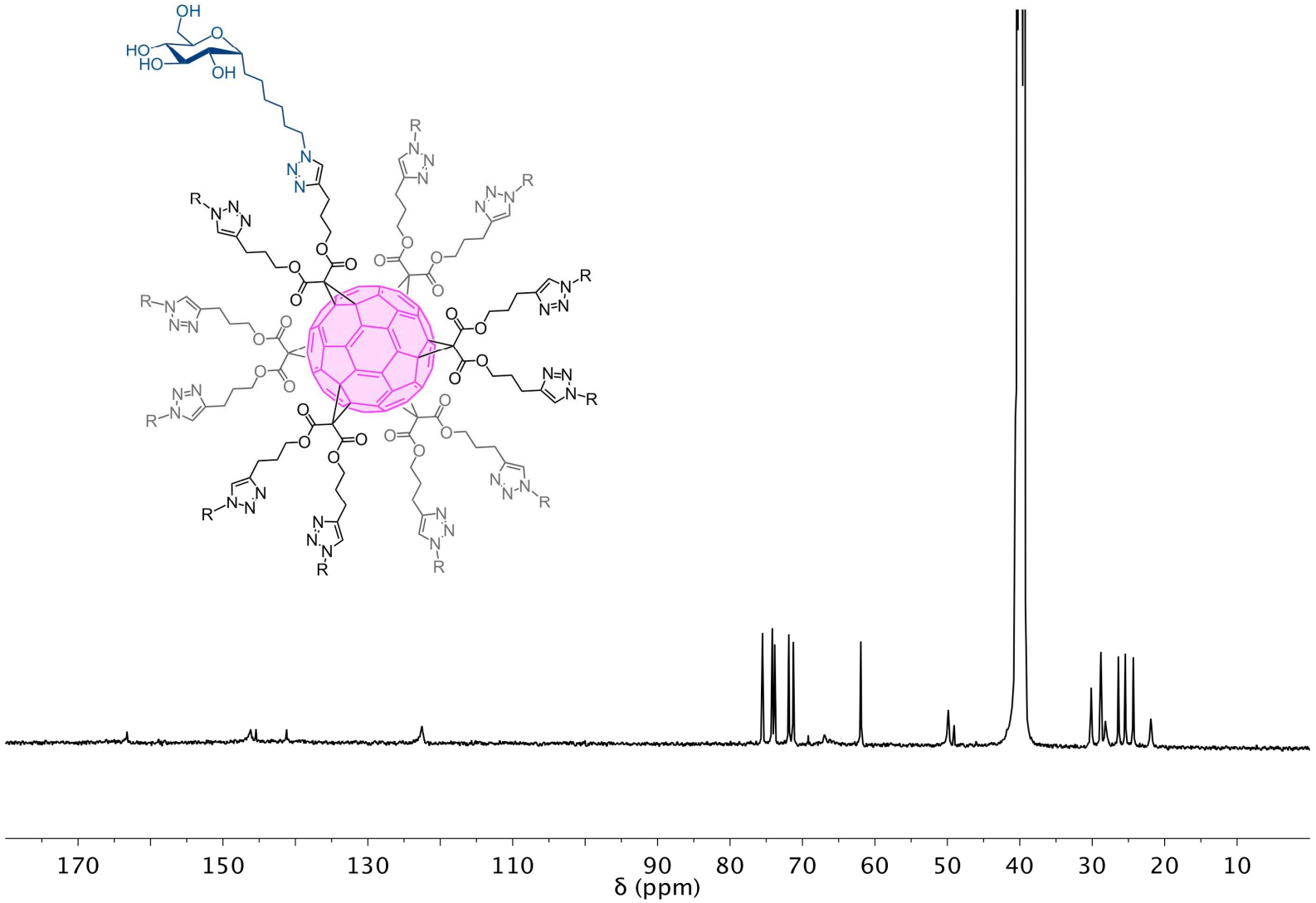


Table S1. Optimization of the olefin metathesis reaction of compound **6a**.[a]

Entry	Catalyst	*T* (°C)	Time	Additive	Product **7**
1	Hoveyda-Grubbs II (20 mol%)	40	18 h	-	~ 42%[b]
2	Grubbs I (10 mol%)	40	26 h	1,4-BQ (20 mol%)	35%
3	Grubbs II (15 mol%)	rt	23 h	1,4-BQ (20 mol%)	55%
4	Grubbs II (15 mol%)	rt	70 h	1,4-BQ (20 mol%)	68%

[a] Reactions were performed in CH_2Cl_2 under argon. [b] Contaminated with analogues of **7** bearing C_5 or C_7 alkyl spacer as judged by MS analysis.