

Figure S1. NMR spectra of Sop_3 .

(A) ^1H -NMR, (B) ^{13}C -NMR, (C) DQF-COSY, (D) TOCSY, (E) HSQC and (F) HMBC. I, II, and III denote first, second, and third glucose residues from reducing end, respectively. Letters in parenthesis represent position of hydroxyl group on the anomeric carbon. Arabic numbers shown with roman numbers represent positions of carbons and protons in sugar rings.

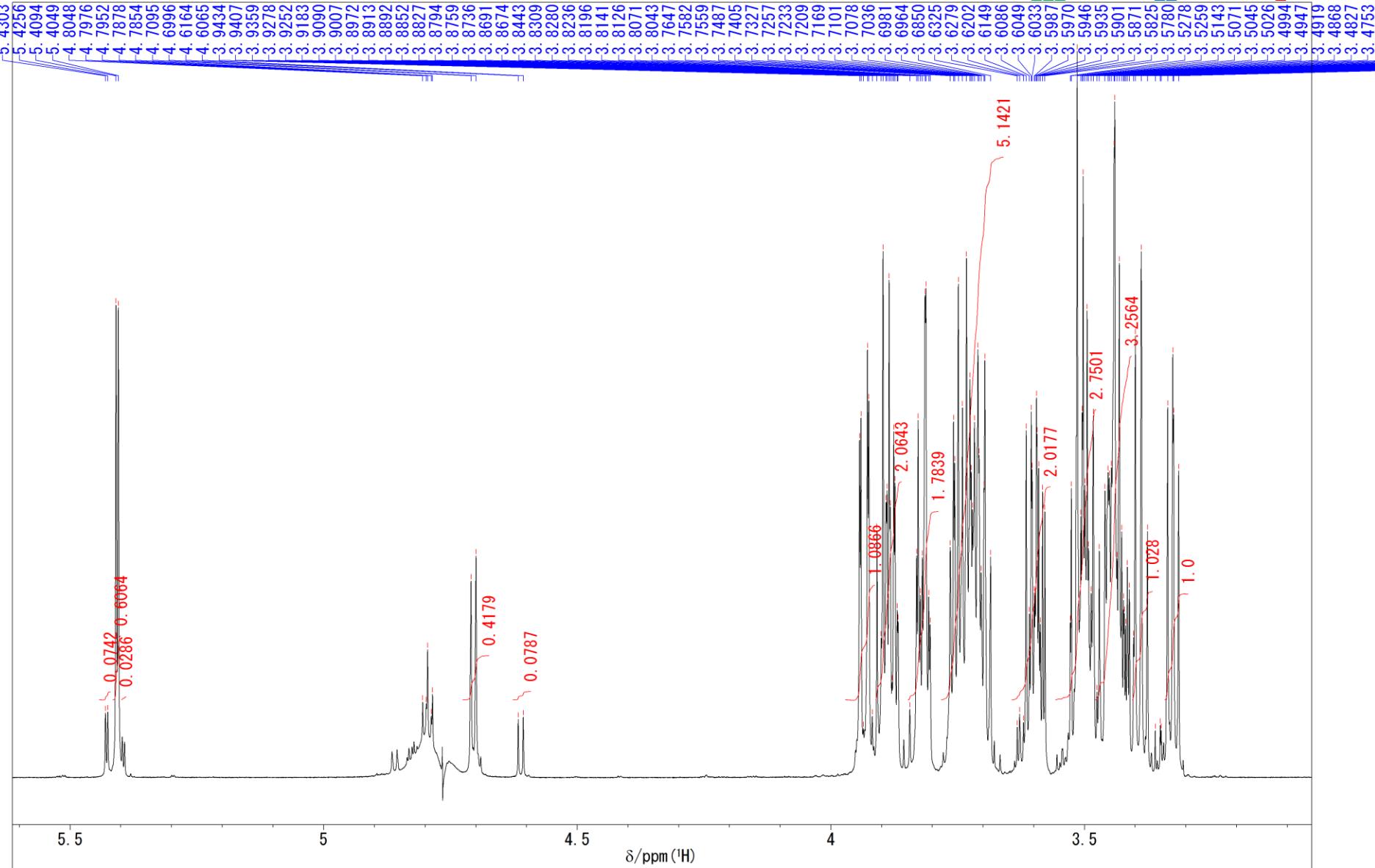
(A) ^1H 1D

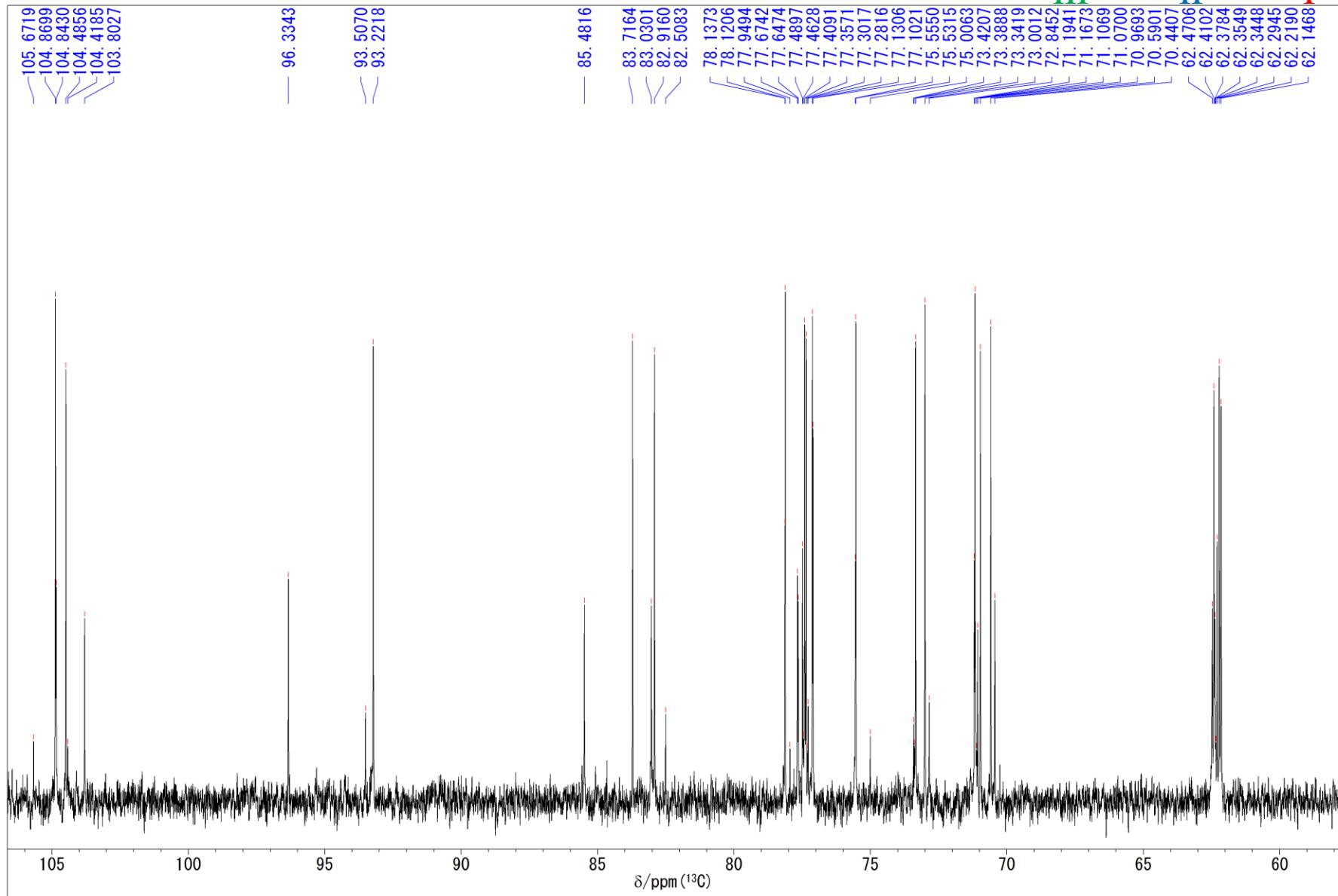
Glc β 1,2Glc β 1,2Glc

III

II

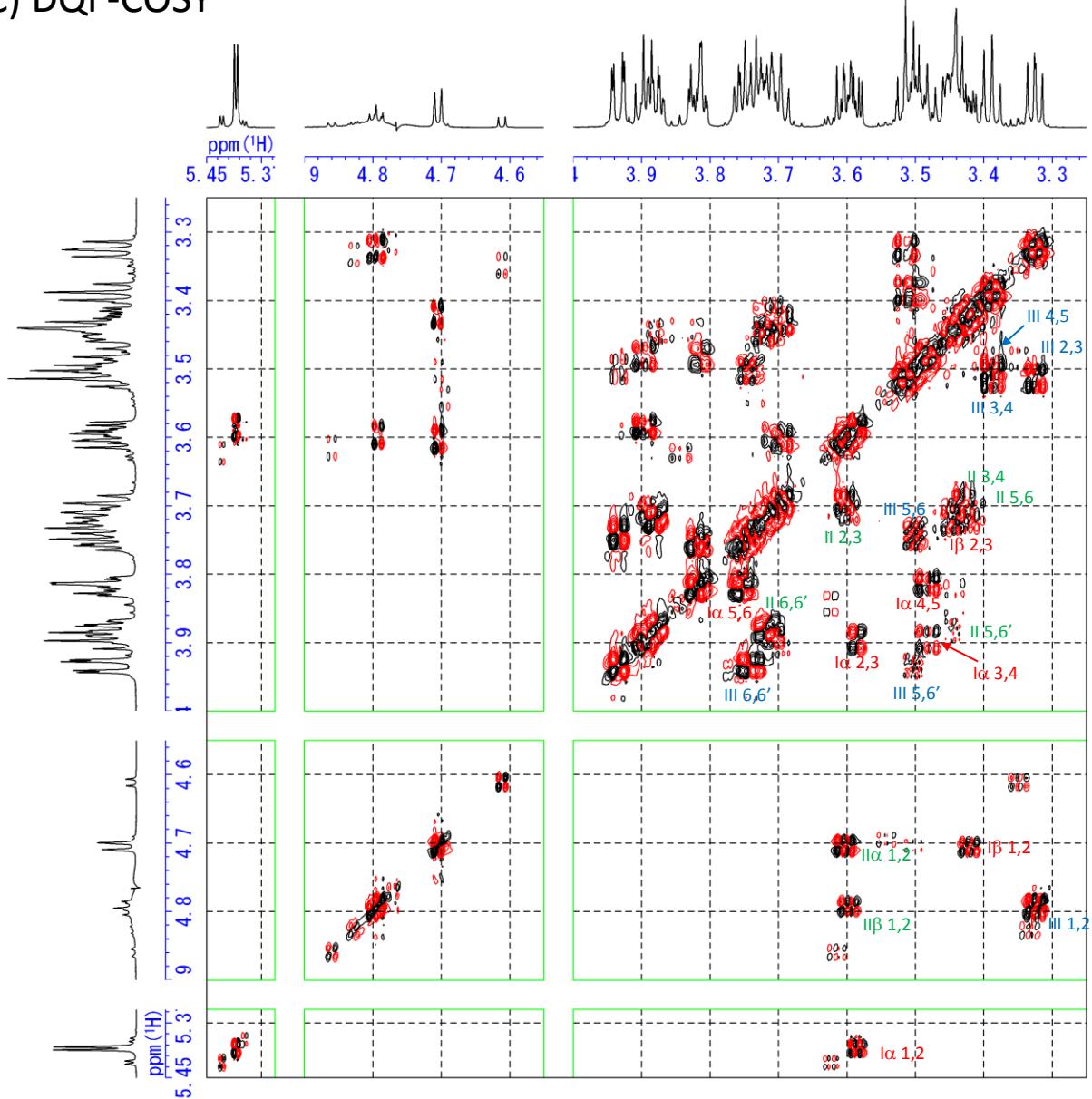
I



(B) ^{13}C 1D

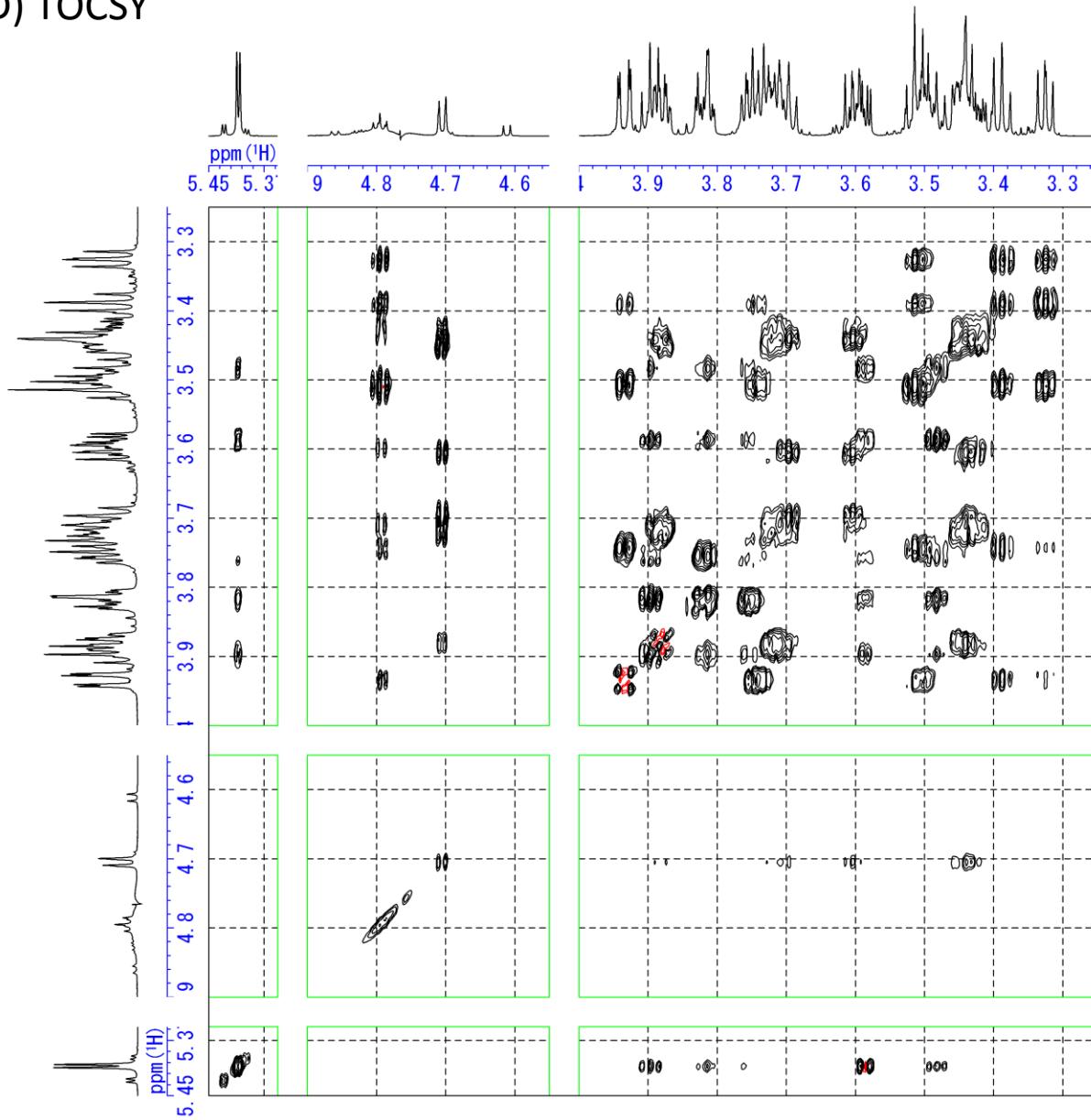
(C) DQF-COSY

Glc β 1,2Glc β 1,2Glc
III II I

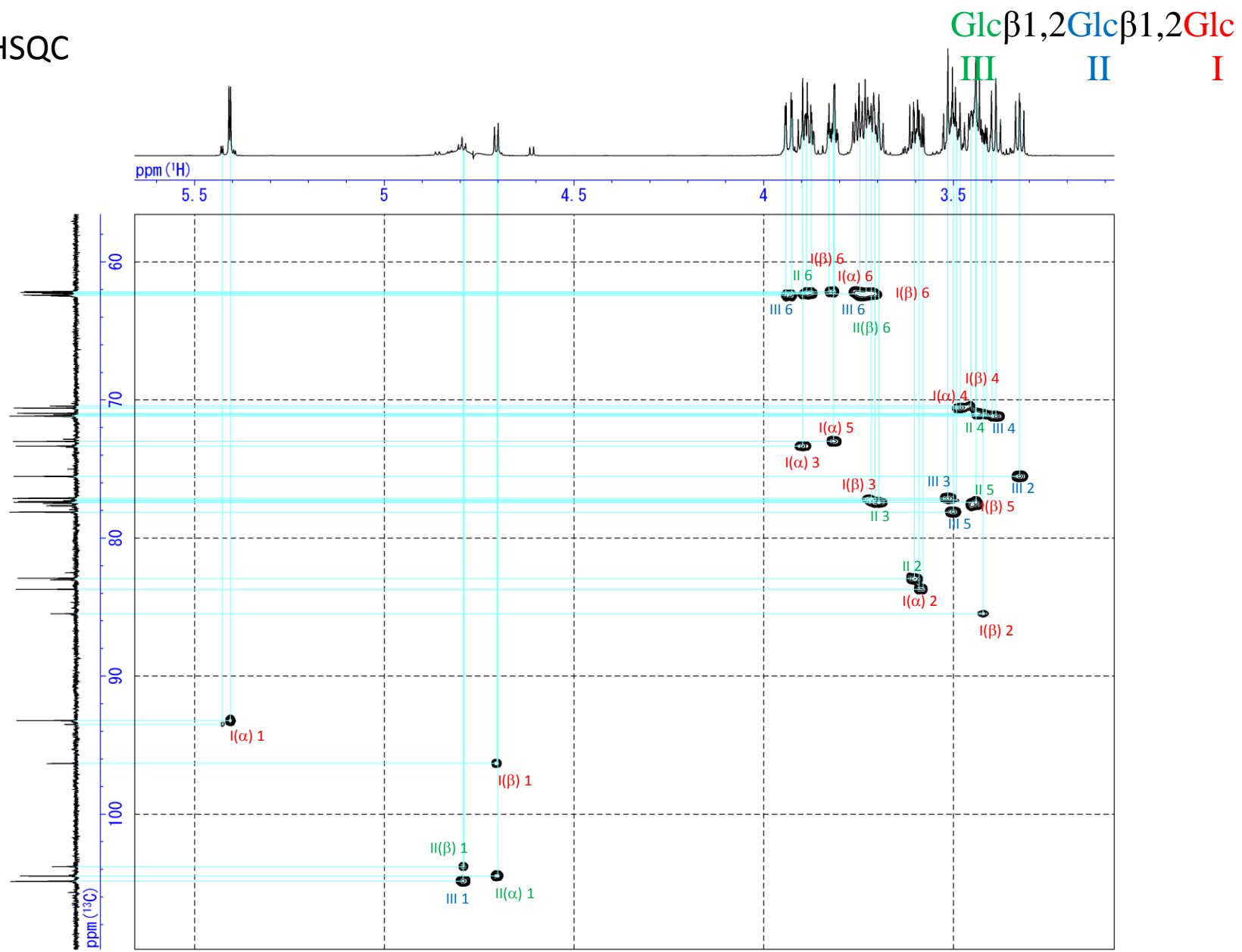


(D) TOCSY

Glc β 1,2Glc β 1,2Glc
III II I



(E) HSQC



(F) HMBC

