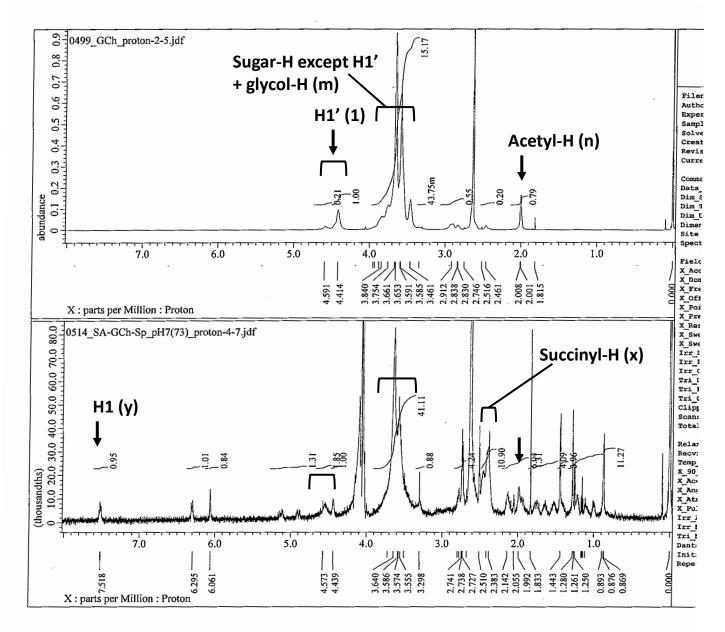
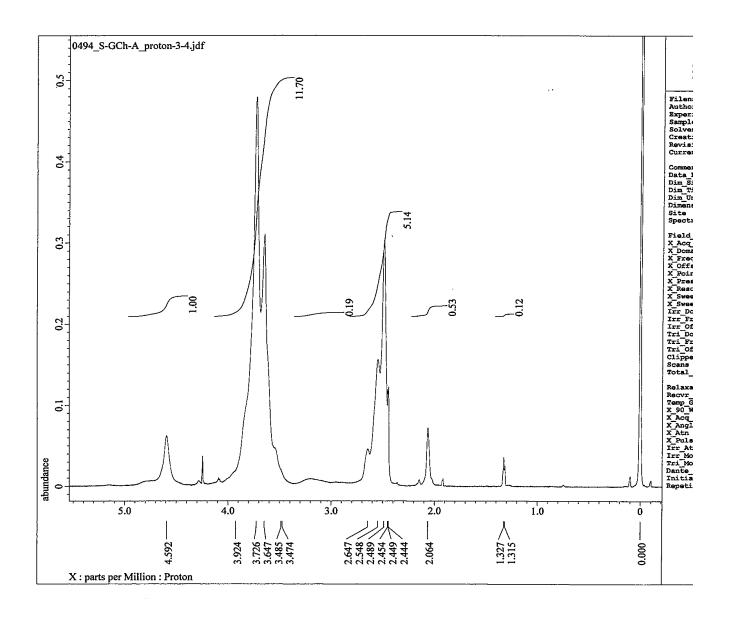


Sapplementary file 1. Proposed structure of S-S-GCh-SP for composition analysis from <sup>1</sup>H-NMR.



Sapplementary file 2. Assignment of each proton of S-GCh-SP. The character or number in the charts shows a molar number in the structure in suuplementary file 1.



Supplementary file 3. The <sup>1</sup>H-NMR spectrum of S-GCh synthesized in the similar reaction.

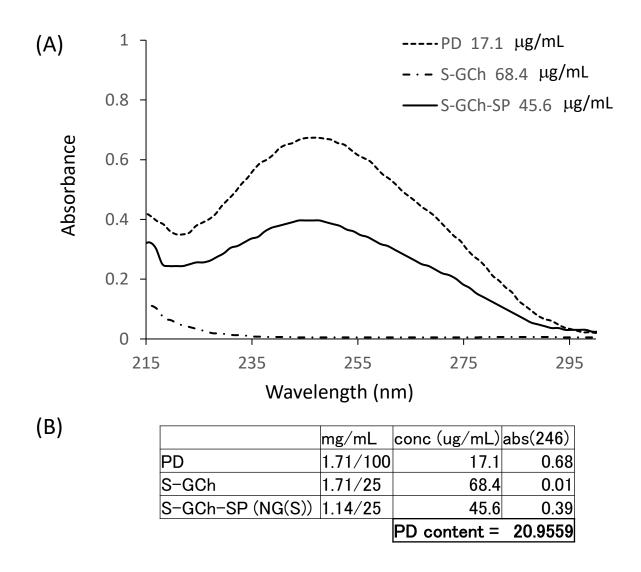
		cal	obs	normalize(mol)		mol ratio
GCh	H1'	1	1.21	1		
	other H'	6 + 4m	15.17	12.53719008	m=	1.63429752
	СОСН3	3n	0.79	0.652892562	n=	0.21763085

		cal	obs	normalize			mol ratio	
S-GCh-SP	H1'	1	1.85	0.564315252				
	other H'	6 + 4m	41.11	12.54	•	m=	1.635	
	СОСН3	3n	1.5	0.457552907		n=	0.15251764	
	Suc	4x	10.9	3.324884456		x=	0.83	
	PD	у	0.95	0.289783508		y=	0.28	
						x-y =	0.55	Succinylation d

0.28 mol PD part amount PD MW=360

PD content 23.42116269 (%, w/w)

Supplementary file 4. Determination of molar ratio of each part (m, n, x, y) and PD content calculated from the 1H-NMR spectra. The m value over 1.0 means glycol group is attached to the parts other C6' position. The PD content = 23.4 % (w/w).



Supplementary file 5. UV absorption profile of PD, S-GCh and S-GCh-SP (treated at 45 oC for 10 min) in 0.1 M NaOH (A), and PD content calculated from the UV absorption at 246 nm. The PD content = 21.0 % (w/w). The PD content was estimated to be 20.8 % (w/w) in triplicated measurements.