

## Supplementary data

### Chondroitin sulfate/dermatan sulfate hybrid chains from swim bladder:

### Isolation, structural analysis, and anticoagulant activity

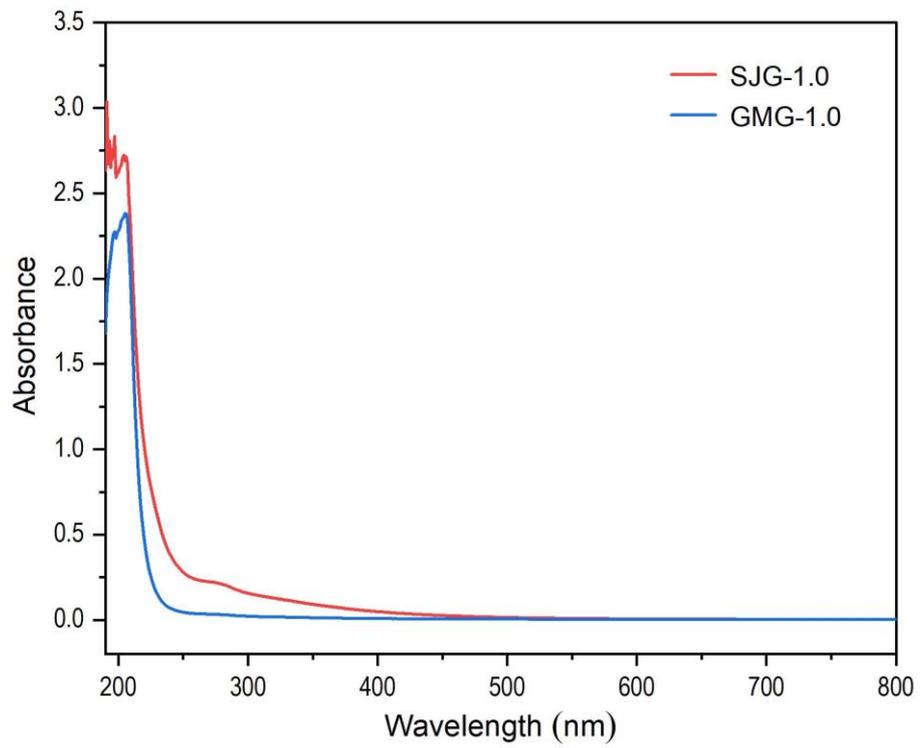
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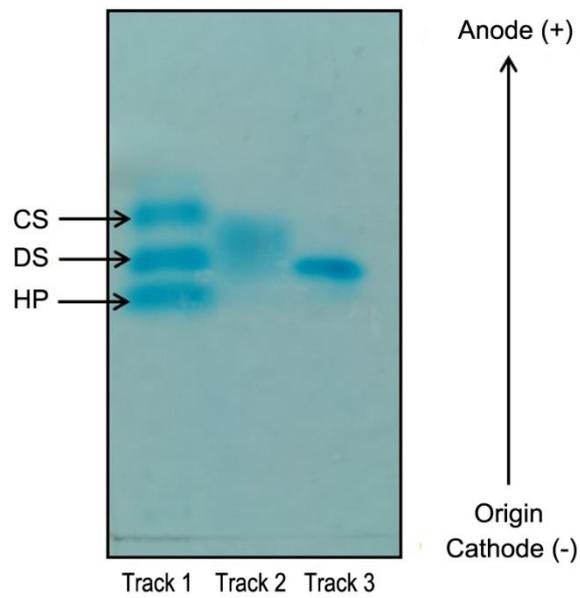
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**Figure S1.** UV spectrum of SJG-1.0 and GMG-1.0



**Figure S2.** Electrophoretogram of SJG-1.0 (Track 2) and GMG-1.0 (Track 3)



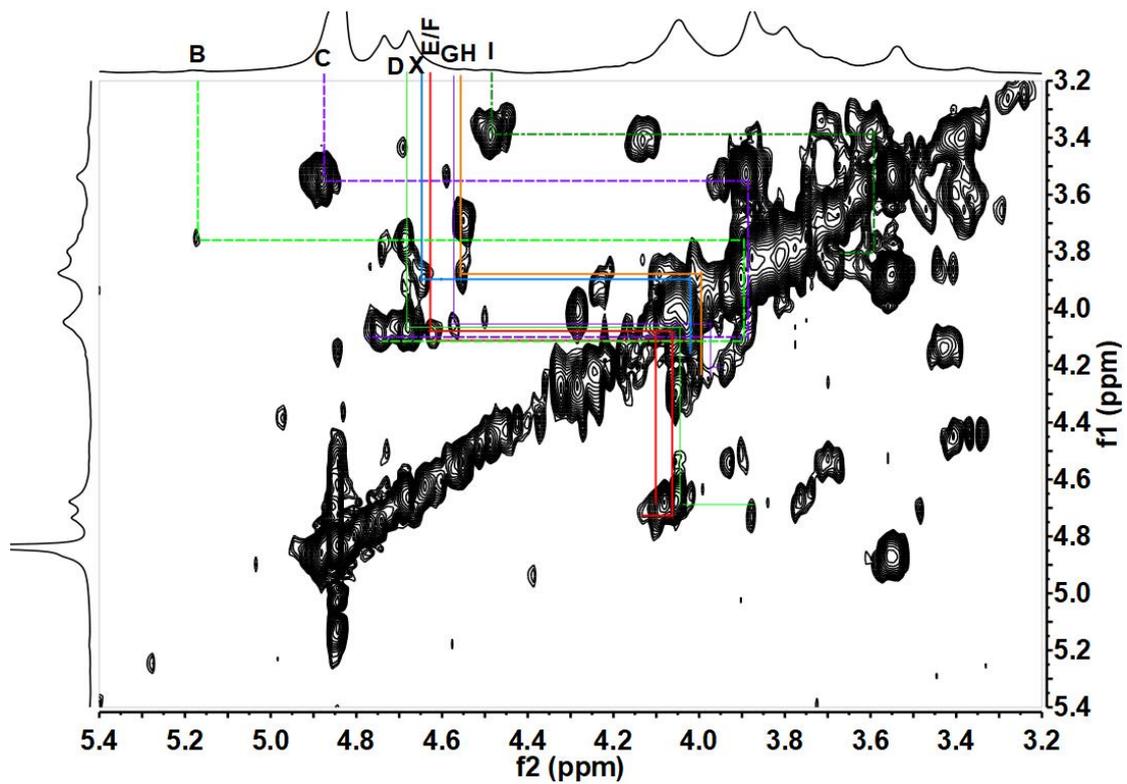


Figure S5.  $^1\text{H}$ - $^1\text{H}$  COSY spectrum of GMG-1.0

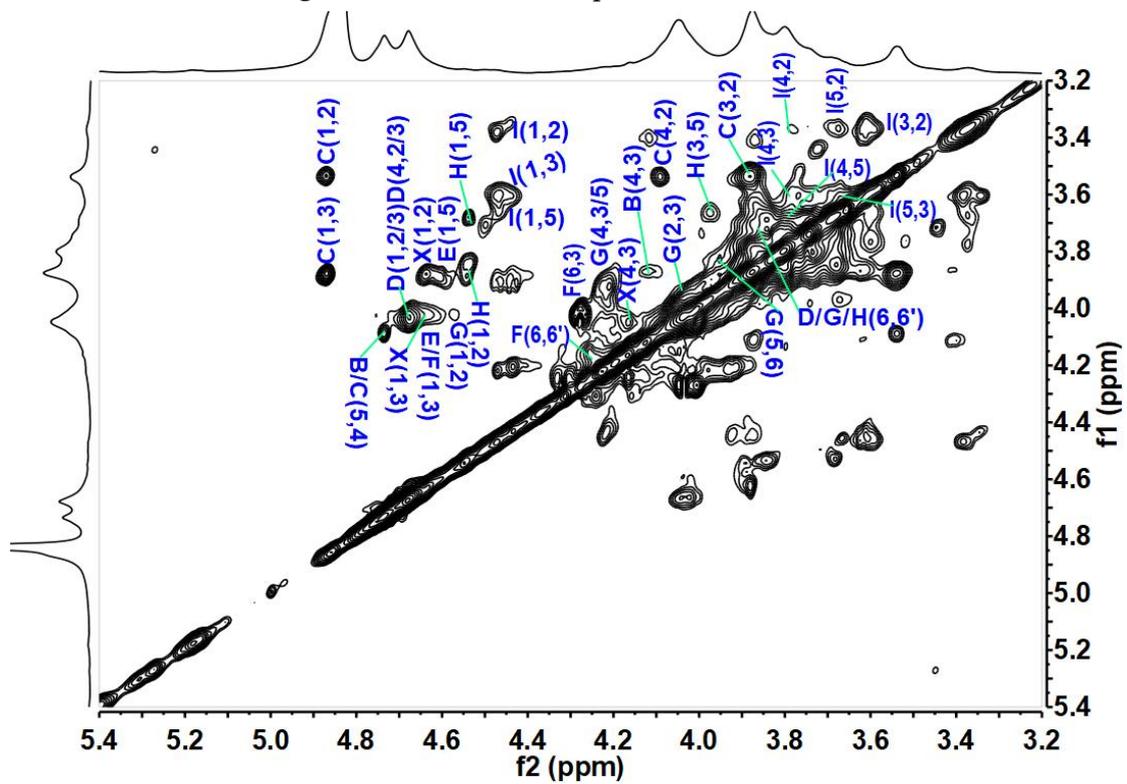


Figure S6.  $^1\text{H}$ - $^1\text{H}$  TOCSY spectrum of GMG-1.0

Table S1. Assignment of <sup>1</sup>H and <sup>13</sup>C NMR signals of GMG-1.0.

Residues	H/C	Chemical Shifts ( $\delta$ , ppm) <sup>a</sup>							
		1	2	3	4	5	6	7	8
A	H	5.27	<u>3.85</u> <sup>b</sup>	n.d.	<b>4.10</b> <sup>c</sup>	4.76			
$\alpha$ -D-IdoA <sub>2S</sub>	C	103.6	<u>79.0</u>	n.d.	<b>83.0</b>	72.4	176.8		
B	H	5.18	<u>3.76</u>	3.87	<b>4.12</b>	4.74			
$\alpha$ -D-IdoA <sub>2S</sub>	C	103.2	<u>79.1</u>	73.5	<b>83.7</b>	72.7	176.8		
C	H	4.87	3.54	3.89	<b>4.08</b>	4.74			
$\alpha$ -D-IdoA	C	106.2	72.4	74.5	<b>83.7</b>	72.7	176.8		
D	H	4.67	4.08	<b>4.04</b>	<u>4.68</u>	3.87	3.74/3.84		2.08
$\beta$ -D-GalNAc <sub>4S</sub>	C	105.0	54.3	<b>78.1</b>	<u>78.9</u>	77.4	63.8	178.1	25.0
X	H	4.64	<u>3.89</u>	<b>4.02</b>	4.16	3.85	<u>4.17/4.84</u>		2.08
$\beta$ -D-GalN <sub>2S6S</sub>	C	105.3	<u>55.0</u>	<b>78.9</b>	70.8	77.4	<u>70.0</u>	178.1	25.0
E	H	4.62	4.10	<b>4.08</b>	<u>4.65</u>	3.88	3.73/3.84		2.05
$\beta$ -D-GalNAc <sub>4S</sub>	C	105.0	54.7	<b>78.6</b>	<u>79.1</u>	77.7	63.9	177.9	25.5
F	H	4.61	4.08	<b>4.04</b>	<u>4.73</u>	4.15	<u>4.17/4.25</u>		2.04
$\beta$ -D-GalNAc <sub>4S6S</sub>	C	105.0	54.3	<b>78.9</b>	<u>79.4</u>	75.7	<u>70.0</u>	177.9	25.5
G	H	4.56	4.02	<b>3.96</b>	4.21	3.93	3.74/3.84		2.04
$\beta$ -D-GalNAc	C	105.0	54.5	<b>78.6</b>	70.3	77.4	63.9	177.4	25.5
H	H	4.54	<u>3.88</u>	<b>4.00</b>	4.19	3.69	3.70/3.80		
$\beta$ -D-GalN <sub>2S</sub>	C	105.4	<u>55.0</u>	<b>78.4</b>	70.8	77.4	63.8		
I	H	4.48	3.36	3.59	<b>3.78</b>	3.67			
$\beta$ -D-GlcA	C	106.5	75.1	76.4	<b>83.3</b>	79.1	177.4		

n.d. -not detected.

<sup>a</sup> The 600 MHz NMR spectra were recorded at 298 K. All chemical shifts are relative to TSP at 0 ppm.

<sup>b, c</sup> Values with underline and in boldface indicate sulfated and glycosylated positions, respectively.