SUPPORTING INFORMATION

Synthesis and Characterization of Thiourea and Urea Linked Glycolipids as Low Molecular-Weight Hydrogelators

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	 glucopyranosyl]-N^{**}-(n-dodecyl)-thioureidoacetamide (55)



Figure S1 Mode of molecular packing a) monosaccharide b) Lactose (β (1-4) linkage) b) Maltose (α (1-4) linkage).



Figure S2 SEM images of the air dried thiourea-linked glycolipids (0.5 wt % water) a) 44 (GlcTUC8) b) 45 (GlcTUC12) c) 52 (LacTUC8) d) 53 (LacTUC12)



Figure S3 Differential scanning calorimetry cycles of **57** (GalUC8) in nitrogen atmosphere (2 cycles, 5 $^{\circ}C m^{-1}$) a) heating, b) cooling



Differential scanning calorimetry of **58** (ManUC8), **62** (D-AraUC8) and **63** (L-AraUC8) in nitrogen atmosphere (10 $^{\circ}$ C m⁻¹) Figure S4



Data collection and refinement statistics for 56 (GlcUC8) Table S1

Parameter	GlcUC8 (56)
Empirical Formula	$C_{34}H_{73}N_6O_{18.5}$
Formula weight	861.98
Wavelength	0.71073 Å
Crystal system	Monoclinic
Space group	<i>C</i> 2
Cell Dimensions	a = 30.174(10) Å, b = 4.6629(19) Å, c = 32.708(12) Å, $\alpha = 90^{\circ}, \beta = 90^{\circ}, \gamma = 90^{\circ}$
Volume (Å ³)	4543(3)
Z, calculated density (Mg/m ³)	4, 1.260
Absorption coefficient (mm ⁻¹)	0.102
F(000)	1868
Crystal size (mm)	0.3 x 0.2 x 0.2
Theta range (°)	1.71 to 19.90
Index ranges	-28<=h<=28, -3<=k<=4, -26<=l<=31
Reflections collected /unique	6429 / 3183 [R(int) = 0.0556]
Data / restraints / parameters	3183 / 14 / 547
Goodness-of-fit on F ²	1.048
Final R indices [I > 2sigma (I)]	R1 = 0.0689, wR2 = 0.1749
R indices (all data)	R1 = 0.1059, wR2 = 0.2031



Figure S6 ¹³C-NMR (100 MHz, CDCl₃) of *N*-(2,3,4-tri-*O*-acetyl- α -D-arabinopyranosyl)azidoacetamide (17)



Figure S8 ¹³C-NMR (100 MHz, CDCl₃) of N-(2,3,4-tri-O-acetyl- α -L-arabinopyranosyl) azidoacetamide (**18**)









octyl)-thioureidoacetamide (23)

Figure S17 ¹H-NMR (400 MHz, CDCl₃) of 1-*N*-(2,3,4,6-tetra-*O*-acetyl-β-D-mannopyranosyl)-*N*''-(*n*-octyl)-thioureidoacetamide (**24**)

octyl)-thioureidoacetamide (24)

octyl)-thioureidoacetamide (**26**)

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re S34 ¹³C-NMR (100 MHz, CDCl₃) of $1-N-[4-O-(2',3',4',6'-tetra-O-acetyl-\alpha-D-glucopyranosyl)-2,3,6-tri-O-acetyl-\beta-D-glucopyranosyl]-N''-(n-octyl)-thioureidoacetamide ($ **32**)

octyl)-ureidoacetamide (35)

glucopyranosyl)-*N*''-(*n*-octyl)-ureidoacetamide (**37**)

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Figure S88 ¹³C-NMR (100 MHz, DMSO-D⁶) of 1-*N*-(2-deoxy-2-acetamido-β-D-glucopyranosyl)-*N*^{''}-(*n*-octyl)-ureidoacetamide (**59**)

