

**Supplementary Table S1. Data collection and refinement statistics**

PDB code	7XXU	7XXV	7XXW	7XXX	7XXY	7XXZ	<u>human Gal-10</u> <u>5XRH</u>
Resolution (Å)	19.99-1.80 (1.84-1.80)	19.97-1.60 (1.63-1.60)	19.96-1.63 (1.66-1.63)	19.99-1.94 (1.99-1.94)	19.47-1.92 (1.97-1.92Å)	19.54-1.83 (1.87-1.83)	<u>19.50-1.55 (1.58-1.55)</u>
Space group	P6 <sub>5</sub> 22	P6 <sub>5</sub> 22	P6 <sub>5</sub> 22	P6 <sub>5</sub> 22	P6 <sub>5</sub> 22	P6 <sub>5</sub> 22	<u>P6<sub>5</sub>22</u>
Unit cell parameters (a, b, c) (Å), ( $\alpha$ , $\beta$ , $\gamma$ ) (°)	(48.53, 48.53, 258.91), (90.00, 90.00, 120.00)	(48.53, 48.53, 258.57), (90.00, 90.00, 120.00)	(48.86, 48.86, 259.47), (90.00, 90.00, 120.00)	(48.59, 48.59, 258.91), (90.00, 90.00, 120.00)	(48.60, 48.62, 256.27), (90.00, 90.00, 119.99)	(48.72, 48.72, 259.20), (90.00, 90.00, 120.00)	<u>(48.56, 48.56, 260.33), (90.00, 90.00, 120.00)</u>
No. of measured reflections	289574 (17117)	4266594 (7893)	429428 (18387)	261243 (18307)	264291 (17271)	157494 (9846)	<u>489727 (17756)</u>
No. of unique reflections	17996 (1029)	24980 (1152)	24277 (1145)	14536 (962)	14830 (941)	17232 (1034)	<u>27731 (1288)</u>
Completeness (%)	99.9 (100.0)	99.7 (95.7)	99.9 (100.0)	99.9 (100.0)	99.9 (100.0)	99.6 (99.8)	<u>99.7 (96.9)</u>
Multiplicity	16.1 (16.6)	17.1 (6.9)	17.7 (16.1)	18.0 (19.0)	17.8 (18.4)	9.1 (9.5)	<u>17.7 (13.8)</u>
$R_{\text{merge}}$ (%)	9.3 (37.9)	6.8 (76.8)	6.5 (46.9)	9.5 (57.8)	10.3 (74.1)	10.0 (65.0)	<u>8.0 (45.3)</u>
$\langle I/\delta(I) \rangle$	20.1 (6.1)	26.5 (2.6)	26.5 (5.0)	22.6 (5.5)	21.1 (4.6)	14.0 (3.0)	<u>24.3 (6.7)</u>
$R_{\text{model}}$ (%)	18.43	17.57	17.71	17.60	17.39	17.70	<u>17.15</u>

$R_{\text{free}}$ (%)	21.72	19.29	20.54	21.21	20.90	20.77	<u>19.57</u>
Rmsd bond lengths (Å)	0.006	0.006	0.005	0.006	0.006	0.006	<u>0.006</u>
Rmsd bond angles (°)	0.853	0.92	0.853	0.886	0.934	0.908	<u>0.905</u>
Number of atoms							
Protein	1129	1129	1129	1132	1132	1132	<u>1123</u>
Ligand	-	45	6	-	23	6	=
Water	189	171	166	190	31	204	<u>188</u>
Average B factors (Å <sup>2</sup> ) (protein/ligand/solvent)	18.97/-/30.45	16.85/24.16/29.42	19.57/32.55/30.80	21.632/-/33.06	24.47/33.92/34.18	19.08/30.38/31.03	<u>16.71/-/29/64</u>
Ramachandran plot <sup>f</sup> residues in favored regions (%)	97.78	97.78	97.78	97.06	97.79	97.79	<u>97.90</u>
Ramachandran plot <sup>f</sup> residues in allowed regions (%)	1.48	1.48	1.48	2.21	1.47	1.47	<u>1.4</u>
Ramachandran plot <sup>f</sup> residues in outliers regions (%)	0.74	0.74	0.74	0.74	0.74	0.74	<u>0.7</u>
Substrate/Ligand	-	Lactose	Glycerol	-	Lactose	Glycerol	=

