



**Figure S1. Glycolipid receptor of Cry5B.**

Top, One of the glycolipid receptors of Cry5B, previously identified as band B (Griffitts et al., 2001). Middle, a more minimal glycan receptor of Cry5B, previously identified as band E (Griffitts et al., 2001). Bottom, glycan previously identified as band D (Griffitts et al., 2001), which is not bound by Cry5B. The arthroseries tetrasaccharide core is boxed in dotted lines.

**Table S1.** Data collection, phasing and refinement statistics

	Native	SeMet
<b>Data collection</b>		
Space group	P6 <sub>3</sub>	P6 <sub>3</sub>
Cell dimensions		
<i>a</i> , <i>b</i> , <i>c</i> (Å)	a=b=115.5 c=110.3	a=b=114.5 c=111.0
$\alpha$ , $\beta$ , $\gamma$ (°)	$\alpha=\beta=90$ , $\gamma=120$	$\alpha=\beta=90$ , $\gamma=120$
		<i>Peak</i>
Wavelength (Å)	0.97930	0.97865
Resolution (Å)	50.00-2.30 (2.34- 2.30)	50.00-3.15 (3.20- 3.15)
<i>R</i> <sub>merge</sub>	15.3 (70.5)	20.6 (75.5)
<i>I</i> / $\sigma$ <sub><i>I</i></sub>	17.5 (3.2)	41.9 (7.3)
Completeness (%)	99.9 (100)	100 (100)
Redundancy	6.3 (6.3)	42.7 (35.9)
<b>Refinement</b>		
Resolution (Å)	39.9-2.3	
No. reflections	37,191	
<i>R</i> <sub>work</sub> / <i>R</i> <sub>free</sub>	0.173/0.218	
No. atoms		
Protein	4,624	
Ligand/ion		
Water	481	
B-factors (Å <sup>2</sup> )		
Protein	32.3	
Ligand/ion		
Water	33.0	
R.m.s deviations		
Bond lengths (Å)	0.008	
Bond angles (°)	1.132	

\*Highest resolution shell is shown in parenthesis.