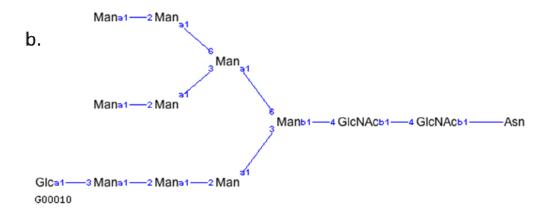
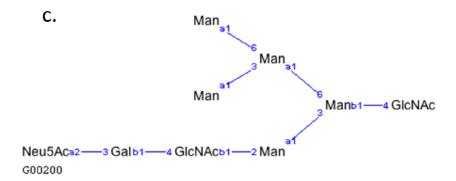
Figures

a. **ENTRY** G00010 Glycan NODE 13 Asn 27 -2 GlcNAc 18 -2 1 Asn GICNAC 9 -2 3 Man 5 Man -6 5 Man Man -13 10 Man 9 Man -13 -9 10 Man -20 10 11 Man -20 0 12 Man -20 -9 -27 -9 13 Glc **EDGE** 12 2:b1 1 3:b1 2:4 4:b1 3:4 5:a1 4:6 4:3 6:a1 7:a1 5:6 8:a1 5:3 9:a1 6:2 10:a1 7:2 10 11:a1 8:2 11 12:a1 9:2 12 13:a1 12:3 ///

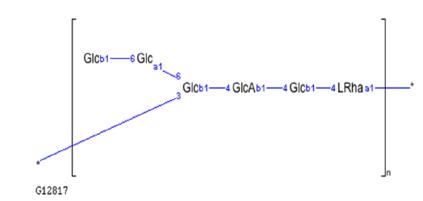




d.

*
$$4$$
 Glc $_{b1}$ 4 LRha $_{a1}$ 3 Glc $_{b1}$ 4 2dGlcA $_{b1}$ * G11448

e.



f.

GlCa1
$$\longrightarrow$$
 2 GlCa1 \longrightarrow 2 GlC + H $_2$ O \longrightarrow GlC + GlCa1 \longrightarrow 2 GlC G00833 \bigcirc G00404

Supplementary Figure 1: KEGG Glycans. a-e: Examples of different types of glycans found in the KEGG Glycans database. (a) A textual representation of the KCF graph for glycan G00010. (b) A visual representation of the KCF graph for glycan G00010. (c) A regular glycan. (d) A Linear repeating

glycan. (e) A non linear repeating glycan. (f) Reconstruction of GlyDeR reactions. The GlyDeR algorithm receives a glycan graph structure and an EC number as input and generates the appropriate glycan products, while considering the following rules: Glycosidic linkages hydrolyzed, Endo vs. exo acting enzyme, Degree of polymerization preference, Reducing vs. non-reducing end preference, Contained sub-glycan, Glycan Released. In the example depicted in the figure, these fields had the following values respectively: Glc a1-2 Glc, exo, 10+, non-reducing, unknown, TAU00015 (the glycan ID of glucose).