## Ruthenium complexes carrying a disialo complex-type oligosaccharide: enzymatic synthesis and its application to a luminescent probe to detect influenza viruses

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Molecular dynamics (MD) calculations of YDS-adducts. MD calculations of YDS-adducts were carried out using InsightII/Discover program and esff force field (Temp. = 300 K and  $\varepsilon$  = 80.0). After 500 ps equilibration, 500 ps dynamics was carried out. Figure S1 shows the most stable conformations of YDS<sub>1</sub>- $\Delta$ Ru during dynamics step. The complex core is highly shielded from outer space by the densely packed saccharide appendages including YDS.

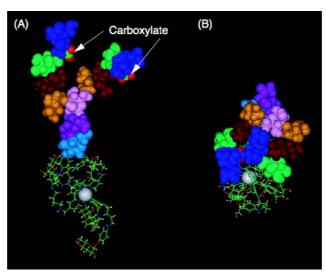


Fig. S1 The (A) initial and (B) most stable conformation of YDS-ΔRu. The each saccharide residues in YDS-moieties are highlighted in CPK