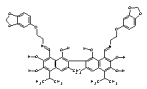

Formula: C48H48N2O12

MW: 844.92

LogP: 2.45

LogS: -3.62

Acceptors: 12

Donors: 6

Rotation Bonds: 19

Chiral Centers: 0

N+O: 14

LIPINSKY: 1

IUPAC: 8-((1E)-4-(2H-benzo[d]1,3-dioxolan-5-yloxy)-2-azabut-1-enyl)-2-[8-((1E)-4-(2H-benzo[d]1,3-dioxolan-5-yloxy)-2-azabut-1-enyl)-1,6,7-trihydroxy-3-methyl-5-(methylethyl)(2-naphthyl)]-3-methyl-5-(methylethyl)naphthalene-1,6,7-triol

Smiles:
c1(c2c(c3c(/C=NCCOc4cc5c(cc4)OCO5)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=NCCOc3cc4c(cc3)OCO4)c(O)c(c(c2cc1C)C(C)C)O)O