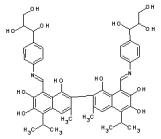

Formula: C48H52N2O12

MW: 848.95

LogP: 5.12

LogS: -5.12

Acceptors: 12

Donors: 12

Rotation Bonds: 21

Chiral Centers: 4

N+O: 14

LIPINSKY: 1

IUPAC: 8-{(1E)-2-[4-(1,2,3-trihydroxypropyl)phenyl]-2-azavinyl}-2-(8-{(1E)-2-[4-(1,2, 3-trihydroxypropyl)phenyl]-2-azavinyl}-1,6,7-trihydroxy-3-methyl-5-(methylethy l)(2-naphthyl))-3-methyl-5-(methylethyl)naphthalene-1,6,7-triol

Smiles:
c1(c(O)c(c2cc(C)c(c3c(c4c(cc3C)c(c(O)c(O)c4/C=Nc3ccc(cc3)C(C(O)CO)O)C(C)C)O)c(c2c1/C =Nc1ccc(cc1)C(O)C(CO)O)O)C(C)C)O