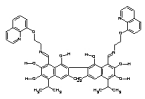

Formula: C52H50N4O8

MW: 858.99

LogP: 11.41

LogS: -7.5

Acceptors: 8

Donors: 6

Rotation Bonds: 19

Chiral Centers: 0

N+O: 12

LIPINSKY: 1

IUPAC: 8-((1E)-4-(8-quinolyloxy)-2-azabut-1-enyl)-2-[8-((1E)-4-(8-quinolyloxy)-2-azab ut-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=NCCOc4c5c(ccc4)cccn5)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=NCCOc3c4c(ccc 3)cccn4)c(O)c(c(c2cc1C)C(C)C)O)O