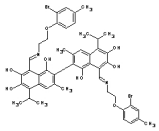

Formula: C48H50Br2N2O8

MW: 942.74

LogP: -3.28

LogS: -1.84

Acceptors: 8

Donors: 6

Rotation Bonds: 19

Chiral Centers: 0

N+O: 10

LIPINSKY: 2

IUPAC: 8-[(1E)-4-(2-bromo-4-methylphenoxy)-2-azabut-1-enyl]-2-{8-[(1E)-4-(2-bromo-4-methylphenoxy)-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(O)c(c(c3cc2C)C(C)C)O)/C=NCCOc2c(cc(cc2)C)Br)O)c(c2c(c(O)c(c(c2cc1C)C(C)C)O)/C=NCCOc1c(cc(cc1)C)Br)O