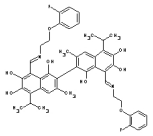

Formula: C46H46F2N2O8

MW: 792.88

LogP: -5.14

LogS: -2.27

Acceptors: 8

Donors: 6

Rotation Bonds: 19

Chiral Centers: 0

N+O: 10

LIPINSKY: 2

IUPAC: 8-[(1E)-4-(2-fluorophenoxy)-2-azabut-1-enyl]-2-{8-[(1E)-4-(2-fluorophenoxy)-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(F)cccc2)O)c(c2c(c(O)c(c(c2cc1C)C(C)C)O)/C=NCCOc1c(F)cccc1)O