



Formula: C₅₀H₅₆N₂O₈

MW: 813



LogP: 2.29

LogS: -3.42

Acceptors: 8

Donors: 6

Rotation Bonds: 17

Chiral Centers: 0

N+O: 10

LIPINSKY: 2

IUPAC: 8-[(1E)-4-(2,6-dimethylphenoxy)-2-azabut-1-enyl]-2-{8-[(1E)-4-(2,6-dimethylphe noxy)-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=NCCOc4c(cccc4C)C)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=NCCOc3c(cccc3C)C)c(O)c(c(c2cc1C)C(C)C)O)O