



Formula: C₅₀H₅₂Cl₄N₂O₈

MW: 950.78



LogP: 3.29

LogS: -4.1

Acceptors: 8

Donors: 6

Rotation Bonds: 19

Chiral Centers: 0

N+O: 10

LIPINSKY: 2

IUPAC: 8-[(1E)-4-(2,4-dichloro-3,5-dimethylphenoxy)-2-azabut-1-enyl]-2-{8-[(1E)-4-(2, 4-dichloro-3,5-dimethylphenoxy)-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(c(C)c(c(c2)C)Cl)Cl)O)c(c2c(c(O)c(c(c2cc1C)C(C)C)O)/C=NCCOc1c(c(C)c(c(c1)C)Cl)Cl)O