



Formula: C₅₄H₆₂Cl₂N₂O₈

MW: 938



LogP: 0.86

LogS: -4.26

Acceptors: 8

Donors: 6

Rotation Bonds: 21

Chiral Centers: 0

N+O: 10

LIPINSKY: 2

IUPAC: 8-{(1E)-4-[4-chloro-5-methyl-2-(methylethyl)phenoxy]-2-azabut-1-enyl}-2-(8-{(1 E)-4-[4-chloro-5-methyl-2-(methylethyl)phenoxy]-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(c(c2)C)Cl)C(C)C)O)c(c2c(c(O)c(c(c2cc1)C(C)C)O)/C=NCCOc1c(cc(c(c1)C)Cl)C(C)C)O