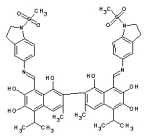

Formula: C48H50N4O10S2

MW: 907.08

LogP: 13.8

LogS: -9.24

Acceptors: 10

Donors: 6

Rotation Bonds: 11

Chiral Centers: 0

N+O: 14

LIPINSKY: 1

IUPAC: 5-{(1E)-2-[7-(8-{(1E)-2-[1-(methylsulfonyl)indolin-5-yl]-2-azavinyl}-1,6,7-tri hydroxy-3-methyl-5-(methylethyl)(2-naphthyl))-2,3,8-trihydroxy-6-methyl-4-(met hylethyl)naphthyl]-1-azavinyl}-1-(methylsulfonyl)indoline

Smiles:
c1(c(c(C(C)C)c2c(c1/C=Nc1ccc3c(c1)CCN3S(C)(=O)=O)c(O)c(c1c(cc3c(c1O)c(/C=Nc1cc4c(cc 1)N(CC4)S(=O)(=O)C)c(O)c(O)c3C(C)C)C)c(C)c2)O)O