



Formula: C₄₂H₄₂N₄O₁₀

MW: 762.82



LogP: -1.05

LogS: -2.94

Acceptors: 10

Donors: 8

Rotation Bonds: 15

Chiral Centers: 0

N+O: 14

LIPINSKY: 2

IUPAC: N-{(1E)-2-[7-(8-{(1E)-2-[(2-methyl(3-furyl))carbonylamino]-2-azaviny}-1,6,7-t rihydroxy-3-methyl-5-(methylethyl)(2-naphthyl))-2,3,8-trihydroxy-6-methyl-4-(m ethylethyl)naphthyl]-1-azaviny}-2-methyl(3-furyl))carboxamide

Smiles:

c1c(c2c(c3c(/C=NNC(=O)c4ccoc4C)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=NNC(=O)c3ccoc3C)c(O)c(c(c2cc1C)C(C)C)O)O