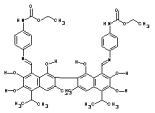

Formula: C48H50N4O10

MW: 842.95

MDL: MFCD12965944

TNP:

LogP: -4.33

LogS: -1.47

Acceptors: 10

Donors: 8

Rotation Bonds: 15

Chiral Centers: 0

N+O: 14

LIPINSKY: 2

IUPAC: N-(4-{(1E)-2-[7-(8-{(1E)-2-[4-(ethoxycarbonylamino)phenyl]-2-azavinyl}-1,6,7-t rihydroxy-3-methyl-5-(methylethyl)(2-naphthyl))-2,3,8-trihydroxy-6-methyl-4-(m ethylethyl)naphthyl]-1-azavinyl}phenyl)ethoxycarboxamide

Smiles:
c1(c2c(c3c(/C=Nc4ccc(cc4)NC(=O)OCC)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=Nc3ccc(cc3)NC (=O)OCC)c(O)c(c(c2cc1C)C(C)C)O)O