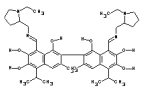

Formula: C44H58N4O6

MW: 738.97

LogP: 4.39

LogS: -4.55

Acceptors: 6

Donors: 6

Rotation Bonds: 17

Chiral Centers: 2

N+O: 10

LIPINSKY: 2

IUPAC: 8-[(1E)-3-(1-ethylpyrrolidin-2-yl)-2-azaprop-1-enyl]-2-{8-[(1E)-3-(1-ethylpyrr olidin-2-yl)-2-azaprop-1-enyl]-1,6,7-trihydroxy-3-methyl-5-(methylethyl)(2-nap hthyl)}-3-methyl-5-(methylethyl)naphthalene-1,6,7-triol

Smiles:
c1(c2c(c3c(/C=NCC4N(CCC4)CC)c(O)c(c(c3cc2C)C(C)C)O)O)c(c2c(/C=NCC3CCCN3CC)c(O) c(c(c2cc1C)C(C)C)O)O