



CAS Registry Number: 38990-09-9

Molecular Formula: $C_{20}H_{24}N_2O_2$

Molecular Weight: 324.422

Accurate Mass: 324.183778

Percentage Composition: C 74.05%; H 7.46%; N 8.63%; O 9.86%

Biological Source: Alkaloid from the stem bark of *Pleiocarpa talbotii* (Apocynaceae)

Physical Description: Cryst. (Et₂O/pentane)

Melting Point: Mp 152-154°

Optical Rotation: $[\alpha]_D^{23} -190$ (c, 0.95 in CHCl₃)

pKa Value: pK_a 6.9 (MCS)

>>Derivative: 16-Epimer, picrate

Physical Description: Cryst. (MeOH)

Melting Point: Mp 187-189°

>>Derivative: 16-Epimer, N-Me

Synonym(s): N-Methyl-16-epiaffinine

CAS Registry Number: 58262-66-1

Molecular Formula: $C_{21}H_{26}N_2O_2$

Molecular Weight: 338.449

Accurate Mass: 338.199428

Percentage Composition: C 74.53%; H 7.74%; N 8.28%; O 9.45%

Biological Source: Alkaloid from *Tabernaemontana accedens* (Apocynaceae)

Physical Description: Plates (MeOH)

Melting Point: Mp 208-210 dec.°

Optical Rotation: $[\alpha]_D^{20} -243$ (c, 0.05 in CHCl₃)

>>Derivative: 16-Epimer, 10-hydroxy

Synonym(s): 10-Hydroxy-16-epiaffinine

CAS Registry Number: 82513-70-0

Molecular Formula: $C_{20}H_{24}N_2O_3$

Molecular Weight: 340.421

Accurate Mass: 340.178693

Percentage Composition: C 70.57%; H 7.11%; N 8.23%; O 14.10%

Biological Source: Alkaloid from *Hunteria zeylanica* (Apocynaceae)

Physical Description: Amorph.

Optical Rotation: $[\alpha]_D -122$ (c, 0.46 in CHCl₃)