



Accurate Mass: 385.188924
Percentage Composition: C 68.55%; H 7.06%; N 3.63%; O 20.75%
Biological Source: Alkaloid from *Colchicum speciosum*
Melting Point: Mp 158-160°
Optical Rotation: $[\alpha]_D +321$

>>Derivative: *O*²-De-Me

Synonym(s): 2-*O*-Demethyl-β-lumicolchicine

CAS Registry Number: 18172-24-2

Molecular Formula: C₂₁H₂₃NO₆
Molecular Weight: 385.416
Accurate Mass: 385.152539
Percentage Composition: C 65.44%; H 6.01%; N 3.63%; O 24.91%
Biological Source: Alkaloid from *Colchicum autumnale*, *Colchicum keselringii*, *Gloriosa superba* and *Wurmbea* spp.
Physical Description: Cryst. (EtOAc)
Melting Point: Mp 198-200°
Optical Rotation: $[\alpha]_D^{20} +337$ (c, 0.45 in CHCl₃)

>>Derivative: *O*³-De-Me

Synonym(s): 3-*O*-Demethyl-β-lumicolchicine

CAS Registry Number: 30632-52-1

Molecular Formula: C₂₁H₂₃NO₆
Molecular Weight: 385.416
Accurate Mass: 385.152539
Percentage Composition: C 65.44%; H 6.01%; N 3.63%; O 24.91%
Biological Source: Alkaloid from *Colchicum* spp., *Gloriosa superba* and *Merendera raddeana*
Melting Point: Mp 235°
Optical Rotation: $[\alpha]_D^{22} +316$ (CHCl₃)

>>Derivative: 3-*O*-De-Me, *N*-de-Ac, *N*-formyl

Synonym(s): *N*-Deacetyl-*N*-formyl-3-*O*-demethyl-β-lumicolchicine

CAS Registry Number: 18172-25-3

Molecular Formula: C₂₀H₂₁NO₆
Molecular Weight: 371.389
Accurate Mass: 371.136889
Percentage Composition: C 64.68%; H 5.70%; N 3.77%; O 25.85%
Biological Source: Alkaloid from *Colchicum decaisnei* and *Gloriosa superba* (Liliaceae)
Physical Description: Cryst. (MeOH/EtOAc)
Melting Point: Mp 229-230°
Optical Rotation: $[\alpha]_D +240$ (c, 0.12 in MeOH)

>>Derivative: 8,12-Diepimer

Synonym(s): γ-Lumicolchicine. Lumicolchicine. Lumicolchicine II. Substance J

CAS Registry Number: 6901-14-0

Molecular Formula: C₂₂H₂₅NO₆
Molecular Weight: 399.443
Accurate Mass: 399.168189
Percentage Composition: C 66.15%; H 6.31%; N 3.51%; O 24.03%
Biological Source: Alkaloid from *Colchicum* spp., *Merendera* spp. and others, as artifact of illumination of Colchicine [CFR10-M](#) and poss. as genuine alkaloid (Liliaceae)
Physical Description: Cryst. (EtOAc/Et₂O, dioxan or anisole)
Melting Point: Mp 268°
Optical Rotation: $[\alpha]_D^{18} -445$ (c, 1 in CHCl₃)
UV: [neutral] λ_{max} 265 () (EtOH) (Berdy)
Sigma: C0148

>>Derivative: 5,12-Diepimer, *N*-de-Ac, *N*-formyl

Synonym(s): *N*-Deacetyl-*N*-formyl-γ-lumicolchicine

CAS Registry Number: 18172-22-0