



Other Data: Abs. config. not detd., may belong to the (3*S*,6*S*)-form

>>Derivative: 3-*O*-Phenylacetyl

Synonym(s): Tropane-3,6-diol phenylacetate

CAS Registry Number: 104086-64-8

Molecular Formula: C₁₆H₂₁NO₃

Molecular Weight: 275.347

Accurate Mass: 275.152144

Percentage Composition: C 69.79%; H 7.69%; N 5.09%; O 17.43%

Biological Source: Alkaloid from root bark of *Erythroxylum hypericifolium* (Erythroxylaceae)

Physical Description: Cryst. (EtOH)

Melting Point: Mp 85°

Optical Rotation: $[\alpha]_{\text{D}}^{20} +4.3$ (c, 1.5 in EtOH)

>>Derivative: 3-*O*-Phenylacetyl, 6-Ac

Synonym(s): Tropane 3,6-diol 3-phenylacetate 6-acetate

CAS Registry Number: 97345-82-9

Molecular Formula: C₁₈H₂₃NO₄

Molecular Weight: 317.384

Accurate Mass: 317.162709

Percentage Composition: C 68.12%; H 7.30%; N 4.41%; O 20.16%

Biological Source: Alkaloid from root bark of *Erythroxylum hypericifolium* (Erythroxylaceae)

Melting Point: Mp 133° (as picrate)

>>Variant: (3*S*,6*S*)-form

CAS Registry Number: 4839-11-6

Molecular Formula: C₈H₁₅NO₂

Molecular Weight: 157.212

Accurate Mass: 157.110279

Percentage Composition: C 61.12%; H 9.62%; N 8.91%; O 20.35%

Biological Source: Alkaloid in *Anthocercis* spp. (Solanaceae)

Melting Point: Mp 212°

Optical Rotation: $[\alpha]_{\text{D}}^{27} -23.3$ (c, 2 in EtOH)

>>Derivative: 3,6-Di-Ac, hydrobromide

Melting Point: Mp 219-220°

>>Derivative: 3-*O*-(2-Methylpropanoyl)

Synonym(s): 3-Isobutyryloxytropan-6-ol

CAS Registry Number: 78886-99-4

Molecular Formula: C₁₂H₂₁NO₃

Molecular Weight: 227.303

Accurate Mass: 227.152144

Percentage Composition: C 63.41%; H 9.31%; N 6.16%; O 21.12%

Biological Source: Alkaloid from aerial parts of *Anthocercis albicans* (Solanaceae)

>>Derivative: 3-*O*-(3-Methylbutanoyl)

Synonym(s): Valeroidine

CAS Registry Number: 490-96-0

Molecular Formula: C₁₃H₂₃NO₃

Molecular Weight: 241.33

Accurate Mass: 241.167794

Percentage Composition: C 64.70%; H 9.61%; N 5.80%; O 19.89%

Biological Source: Alkaloid from *Datura sanguinea* and *Duboisia myoporoides* (Solanaceae)

Melting Point: Mp 85°

Optical Rotation: $[\alpha]_{\text{D}}^{20} -9$ (c, 5 in EtOH)