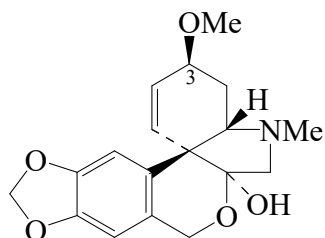




>>Entry Name: Tazettine

Synonym(s): Ungernine. Sekisanine. Sekisanoline



Absolute
Configuration

Molecular Formula: C₁₈H₂₁NO₅

Molecular Weight: 331.368

Accurate Mass: 331.141974

Percentage Composition: C 65.24%; H 6.39%; N 4.23%; O 24.14%

>>Variant: (+)-form

CAS Registry Number: 507-79-9

Molecular Formula: C₁₈H₂₁NO₅

Molecular Weight: 331.368

Accurate Mass: 331.141974

Percentage Composition: C 65.24%; H 6.39%; N 4.23%; O 24.14%

Biological Source: Isol. from *Narcissus tazetta* and a very large number of spp. in the Amaryllidaceae

Biological Use/Importance: Weak hypotensive agent

Melting Point: Mp 210-211°

Optical Rotation: [α]_D¹⁶+150 (CHCl₃)

Other Data: Has been shown to be an artifact produced by base-catalysed rearr. of Pretazettine [CGQ73-J](#) in two cases studied, and is probably an artifact in all cases

Hazard & Toxicity: LD₅₀ (dog, ivn) 71 mg/kg

>>Derivative: Picrate

Melting Point: Mp 205-208°

>>Derivative: O-Ac

Molecular Formula: C₂₀H₂₃NO₆

Molecular Weight: 373.405

Accurate Mass: 373.152539

Percentage Composition: C 64.33%; H 6.21%; N 3.75%; O 25.71%

Melting Point: Mp 125-126.5°

>>Derivative: O-De-Me

Synonym(s): O-Demethyltazettine. **Tazettinol**

CAS Registry Number: 560-31-6

Molecular Formula: C₁₇H₁₉NO₅

Molecular Weight: 317.341

Accurate Mass: 317.126324

Percentage Composition: C 64.34%; H 6.03%; N 4.41%; O 25.21%

Biological Source: Alkaloid from bulbs of *Hippeastrum equestre*

Melting Point: Mp 179-181°

Optical Rotation: [α]_D²²+85 (c, 1 in MeOH)

>>Derivative: O-De-Me, 3-O-(3-hydroxybutanoyl)

Synonym(s): 3-O-(3-Hydroxybutanoyl)tazettinol

Molecular Formula: C₂₁H₂₅NO₇

Molecular Weight: 403.431

Accurate Mass: 403.163104

Percentage Composition: C 62.52%; H 6.25%; N 3.47%; O 27.76%

Biological Source: Alkaloid from *Galanthus plicatus* ssp. *byzantinus*

Physical Description: Amorph. solid

Optical Rotation: [α]_D +63.5 (c, 0.22 in MeOH)

UV: [neutral] $\lambda_{\max}$ 215 (log ϵ 4.12); 234 (log ϵ 3.87); 292 (log ϵ 3.6) (MeOH)