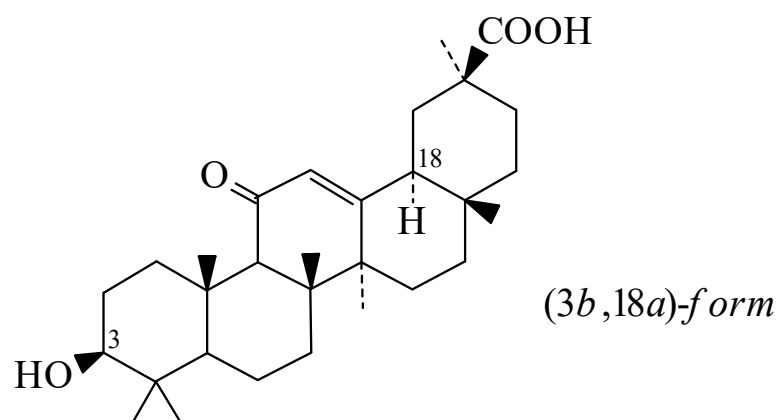




>>Entry Name: 3-Hydroxy-11-oxo-12-oleanen-30-oic acid



Molecular Formula: C₃₀H₄₆O₄

Molecular Weight: 470.691

Accurate Mass: 470.33961

Percentage Composition: C 76.55%; H 9.85%; O 13.60%

Calculated Log P Data: Log P 6.25 (calc)

>>Variant: (3β,18α)-form

Synonym(s): 18α-Glycyrrhetic acid. Isoglycyrrhetic acid. β-Glycyrrhetic acid. Uralenic acid

CAS Registry Number: 1449-05-4

Molecular Formula: C₃₀H₄₆O₄

Molecular Weight: 470.691

Accurate Mass: 470.33961

Percentage Composition: C 76.55%; H 9.85%; O 13.60%

Physical Description: Cryst. (CHCl₃/MeOH)

Melting Point: Mp 330-335°

Optical Rotation: [α]_D +98

Solubility: BERDY SOL: Sol. MeOH, CHCl₃; fairly sol. EtOH; poorly sol. H₂O

UV: [neutral] λ_{max} 242 (ε 12200) (MeOH) (Berdy)

Hazard & Toxicity: LD₅₀ (mus, orl) 560 mg/kg

Sigma: G8503

>>Variant: (3β,18β)-form

Synonym(s): Glycyrrhetic acid. α-Glycyrrhetic acid. Enoxolone, BAN, INN. Biogastrone acid. Biosone. Glycyrrhetin. Glycyrrhetic acid. Rhetinic acid. Uralenic acid

CAS Registry Number: 471-53-4

Molecular Formula: C₃₀H₄₆O₄

Molecular Weight: 470.691

Accurate Mass: 470.33961

Percentage Composition: C 76.55%; H 9.85%; O 13.60%

Biological Source: Aglycone from *Glycyrrhiza glabra* and some other plants

Biological Use/Importance: Antibacterial, antitussive agent. Used in treatment of noninfective inflammatory disorders (skin, mouth, etc.)

Physical Description: Cryst. (MeOH)

Melting Point: Mp 300-304°

Optical Rotation: [α]_D +161 (CHCl₃)

Calculated Log P Data: Log P 6.25 (calc)

Other Data: Sweet taste

Hazard & Toxicity: LD₅₀ (mus, ivn) 56 mg/kg, inhibits 11β-hydroxysteroid dehydrogenase

Aldrich: G1010-5

Fluka: 50510

Sigma: G5628