>>**Variant:** (7'*S*,8*R*,8'*R*,9*S*)-form

Molecular Formula: $C_{18}H_{18}O_6$
Molecular Weight: 330.337
Accurate Mass: 330.11034
Percentage Composition: C 65.45%; H 5.49%; O 29.06%

>>**Derivative:** 3',5,9-Tri-Me ether
Synonym(s): Formosanol. Tsugacetal

CAS Registry Number: 101312-79-2

Biological Source: Isol. from *Dacrydium intermedium* and *Juniperus formosana*
Melting Point: Mp 188-189°
Optical Rotation: $[\alpha]_D^{23}$ –81.3 (c, 1 in $CHCl_3$)

>>**Derivative:** 3',5,9-Tri-Me ether, di-Ac

Melting Point: Mp 201°
Optical Rotation: $[\alpha]_D^{20}$ –76.3 (c, 0.5 in $CHCl_3$)

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