# PRODUCT INFORMATION



## Probetaenone I

Item No. 33252

CAS Registry No.: 115473-44-4

Formal Name: 3-hydroxy-1-[(1R,2S,4aR,6S,8R,8aS)-

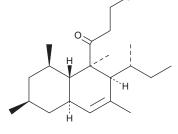
> 1,2,4a,5,6,7,8,8a-octahydro-1,3,6,8tetramethyl-2-[(1R)-1-methylpropyl]-1-

naphthalenyl]-1-propanone

MF:  $C_{21}H_{36}O_{2}$ FW: 320.5 **Purity:** ≥95% Supplied as: A solid Storage: -20°C Stability: ≥4 years

Item Origin: Fungus/Pleospora betae

Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.



## **Laboratory Procedures**

Probetaenone I is supplied as a solid. A stock solution may be made by dissolving the probetaenone I in the solvent of choice, which should be purged with an inert gas. Probetaenone I is soluble in methanol and DMSO.

## Description

Probetaenone I is an intermediate in the biosynthesis of the phytotoxin betaenone B from P. betae.1

### Reference

1. Oikawa, H., Ichihara, A., and Sakamura, S. Biosynthetic study of betaenone B: Origin of the oxygen atoms and accumulation of a deoxygenated intermediate using P-450 inhibitor. J. Chem. Soc. Chem. Commun. 9, 600-602 (1988).

WARNING
THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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