# PRODUCT INFORMATION



# (±)-epi CP 47,497 (exempt preparation)

Item No. 13801

CAS Registry No.: 70434-83-2

Formal Name: rel-5-(1,1-dimethylheptyl)-2-[(1R,3R)-

3-hydroxycyclohexyl]-phenol

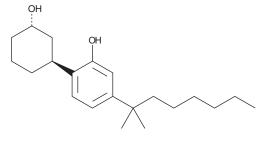
Synonym: 3-trans CP 47,497

MF:  $C_{21}H_{34}O_{2}$ FW: 318.5 **Purity:** ≥98%

UV/Vis.:  $\lambda_{max}$ : 219, 276 nm Supplied as: A solution in methanol

Storage: -20°C Stability: ≥2 years

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



# Description

(±)-epi CP 47,497 (exempt preparation) (Item No. 13801) is an analytical reference standard that is structurally categorized as a cannabinoid (CB). It is an epimer of (±)-CP 47,497 (Item Nos. 10913 | 16851). ( $\pm$ )-CP 47,497 is a synthetic cannabinoid that avidly binds the CB<sub>1</sub> receptor ( $K_i = 2.2$  nM) and is equivalent in analgesic potency to  $\Delta^9$ -THC.<sup>1,2</sup> (±)-epi CP 47,497 is regulated as a Schedule 1 drug in the United States. (±)-epi CP 47,497 (exempt preparation) (Item No. 13801) is provided as a DEA exempt preparation. This product is intended for research and forensic applications.

### References

- 1. Huffman, J.W., Thompson, A.L.S., Wiley, J.L., et al. Synthesis and pharmacology of 1-deoxy analogs of CP-47,497 and CP-55,940. Bioorg. Med. Chem. 16, 322-335 (2008).
- 2. Melvin, L.S., Milne, G.M., Johnson, M.R., et al. Structure-activity relationships for cannabinoid receptorbinding and analgesic activity: Studies of bicyclic cannabinoid analogs. Mol. Pharmacol. 44, 1008-1015 (1993).

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.

# WARRANTY AND LIMITATION OF REMEDY

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