Hexahydrocannabinols (HHCs)

Hydrogenation of tetrahydrocannabinols (Δ^8- or Δ^9-THC) leads to formation of a mixture of hexahydrocannabinols (HHCs) comprised of the 9(S)- and 9(R)-HHC diastereomers. These compounds retain some psychoactivity but avoid classification as THCs as well as any THC-related regulations. Separation of the HHC diastereomers is possible by chromatography (HPLC, GC), but identity can only be confirmed with verified reference standards. Cayman offers fully characterized reference standards to aid in the correct identification and differentiation of hydrogenated phytocannabinoids.

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ISO/IEC 17025 #AT-1773
ISO 17034 #AR-1774

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Characterization of Hexahydrocannabinol Diastereomers by NMR, HPLC, GC-MS, and TLC

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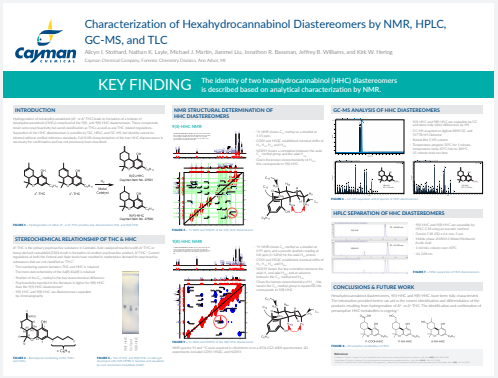
Discover a reliable method to differentiate two HHC diastereomers resulting from Δ^8- or Δ^9-THC hydrogenation.

View the full poster INSIDE


Characterization of Hexahydrocannabinol Diastereomers by NMR, HPLC, GC-MS, and TLC

KEY FINDING The identity of two hexahydrocannabinol (HHC) diastereomers is described based on analytical characterization by NMR.

INTRODUCTION

NMR STRUCTURAL DETERMINATION OF HHC DIASTEREOMERS

GC-MS ANALYSIS OF HHC DIASTEREOMERS

HPLC SEPARATION OF HHC DIASTEREOMERS

STEREOCHEMICAL RELATIONSHIP OF THC & HHC

CONCLUSIONS & FUTURE WORK

KEY FINDING

The identity of
is described ba

INTRODUCTION

Hydrogenation of tetrahydrocannabinols (Δ^8- or Δ^9-THC) leads to formation of a mixture of hexahydrocannabinols (HHCs) comprised of the 9(*S*)- and 9(*R*)-HHC diastereomers. These compounds retain some psychoactivity but avoid classification as THCs as well as any THC-related regulations. Separation of the HHC diastereomers is possible by TLC, HPLC, and GC-MS, but identity cannot be inferred without verified reference standards. Full NMR characterization of the two HHC diastereomers is necessary for confirmation and has not previously been described.

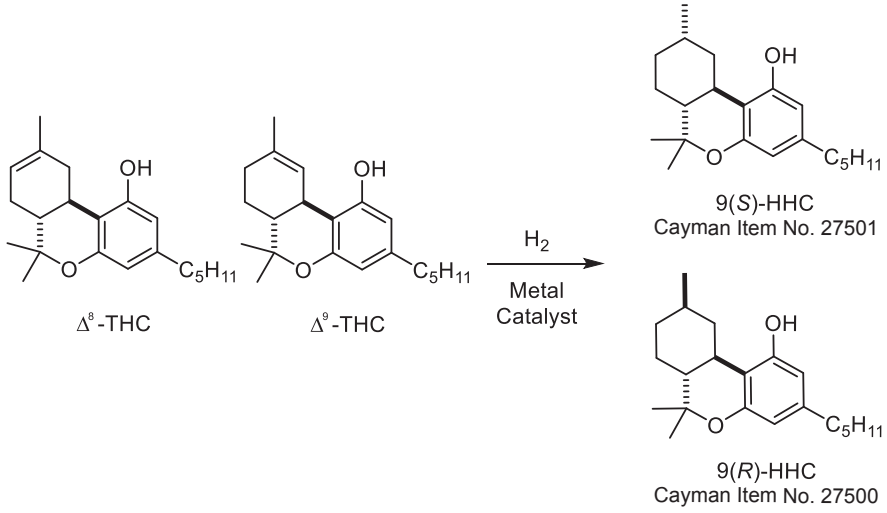


FIGURE 1 – Hydrogenation of either Δ^8- or Δ^9-THC provides two diastereomers 9(*S*)- and 9(*R*)-HHC

STEREOCHEMICAL RELATIONSHIP OF THC & HHC

Δ^9-THC is the primary psychoactive substance in *Cannabis*. Acid-catalyzed reactions with Δ^9-THC or hemp-derived cannabidiol (CBD) result in formation of another psychoactive product, Δ^8-THC.[1] Current regulations at both the Federal and State levels have resulted in marketplace demand for psychoactive substances that are not classified as "THC."

- The numbering system between THC and HHC is retained
- The *trans* stereochemistry of the 6a(*R*),10a(*R*) is retained
- Position of the C_{11}-methyl is the key stereochemical difference
- Psychoactivity reported in the literature is higher for 9(*R*)-HHC than the 9(*S*)-HHC diastereomer[2]
- 9(*S*)-HHC and 9(*R*)-HHC are diastereomers separable by chromatography

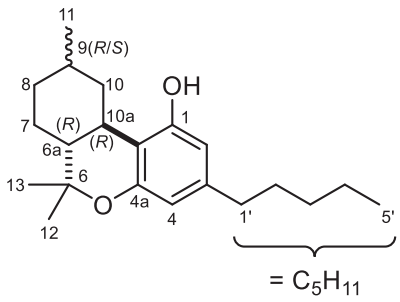


FIGURE 2 – Benzopyran numbering of the THCs and HHCs

FIGURE 3 – TLC of 9(*S*)- and 9(*R*)-HHC on silica gel developed with 10% MTBE in heptane and visualized by ceric ammonium molybdate (CAM)

NMR STRUCTURAL DETERMIN
HHC DIASTEREOMERS

9(*S*)-HHC NMR

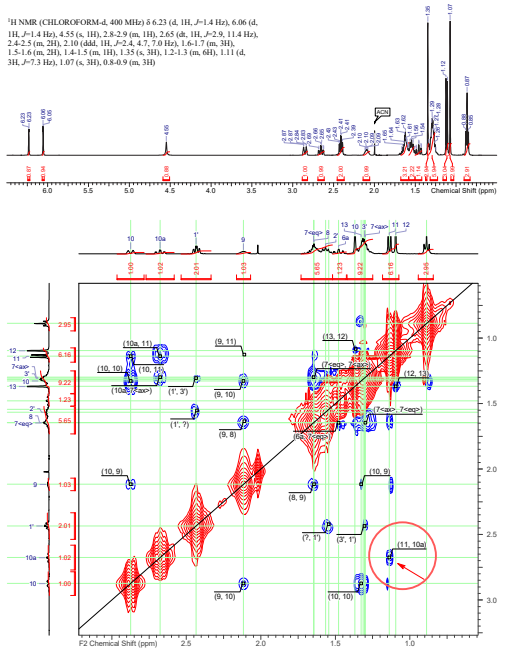


FIGURE 4 – ^{1}H-NMR and NOESY of the 9(*S*)-HHC diastereo

9(*R*)-HHC NMR

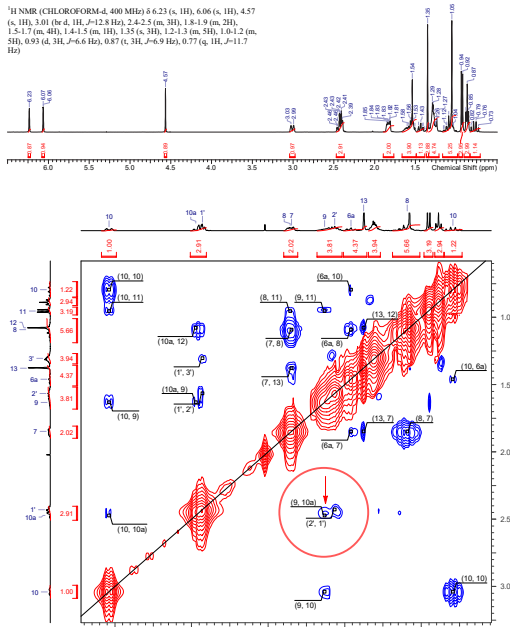


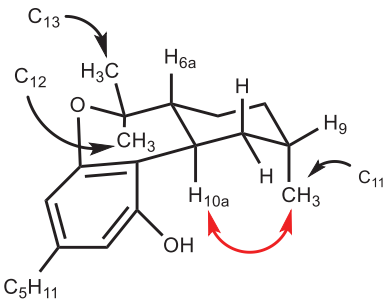
FIGURE 5 – ^{1}H-NMR and NOESY of the 9(*R*)-HHC diastereo

NMR spectra ^{1}H and ^{13}C were acquired in chloroform experiments included COSY, HSQC, and NOESY.

two hexahydrocannabinol (HHC) diastereomers sed on analytical characterization by NMR.

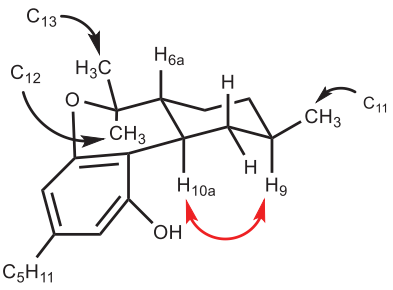
IATION OF

- ^{1}H-NMR shows C_{11}-methyl as a doublet at 1.15 ppm
- COSY and HSQC established chemical shifts of H_9, H_{10}, H_{6a}, and H_{10a}
- NOESY shows a correlation between the axial C_{11}-methyl group and the axial H_{10a}
- Given the known stereochemistry of H_{10a}, this corresponds to 9(*S*)-HHC



omer

- ^{1}H-NMR shows C_{11}-methyl as a doublet at 0.95 ppm, and a pseudo quartet coupling at 0.8 ppm (*J*=12Hz) for the axial H_{10} proton
- COSY and HSQC established chemical shifts of H_9, H_{10}, H_{6a}, and H_{10a}
- NOESY shows the key correlation between the axial H_9 and axial H_{10a}, and an absence between the C_{11}-methyl and H_{10a}
- Given the known stereochemistry of H_{10a}, this means the C_{11}-methyl group is equatorial, this corresponds to 9(*R*)-HHC



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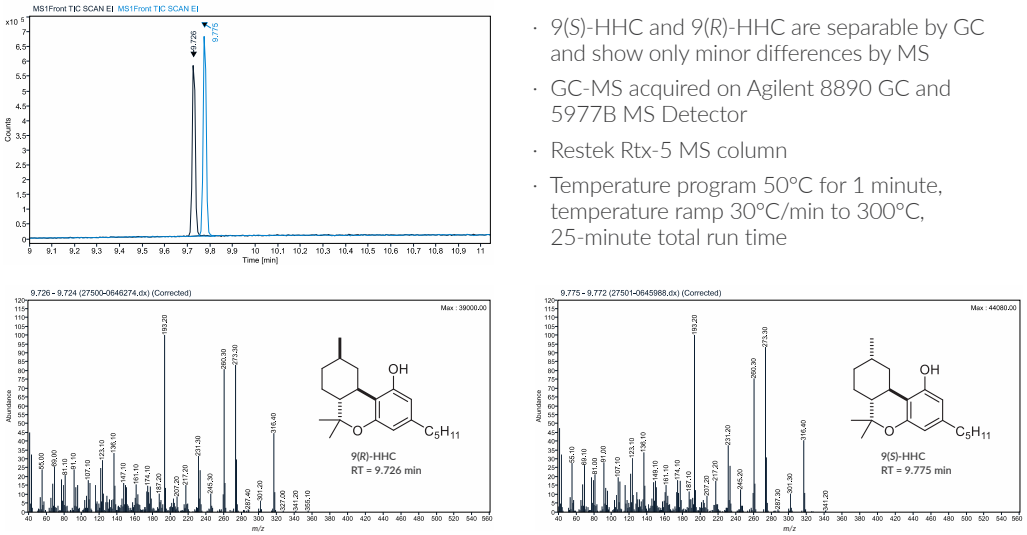
-d on a JEOL ECZ-400S spectrometer, 2D

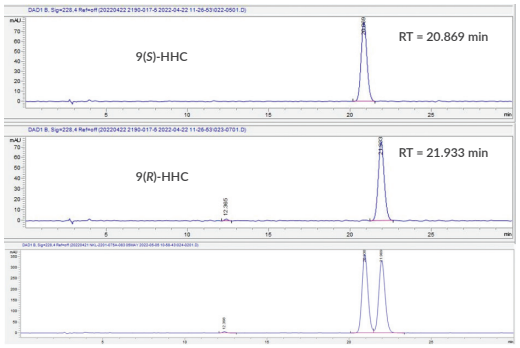
GC-MS ANALYSIS OF HHC DIASTEREOMERS



- 9(*S*)-HHC and 9(*R*)-HHC are separable by GC and show only minor differences by MS
- GC-MS acquired on Agilent 8890 GC and 5977B MS Detector
- Restek Rtx-5 MS column
- Temperature program 50°C for 1 minute, temperature ramp 30°C/min to 300°C, 25-minute total run time

FIGURE 6 – GC-MS separation and EI spectra of HHC diastereomers

HPLC SEPARATION OF HHC DIASTEREOMERS



- 9(*S*)-HHC and 9(*R*)-HHC are separable by HPLC-C18 using an isocratic method
- Gemini-C18 250 x 4.6 mm, 5 µm
- Mobile phase 20:80:0.1 Water/Methanol/ Acetic Acid
- 1 ml/min, column oven 40°C
- UV 228 nm

FIGURE 7 – HPLC separation of HHC diastereomers

CONCLUSIONS & FUTURE WORK

Hexahydrocannabinol diastereomers, 9(*S*)-HHC and 9(*R*)-HHC, have been fully characterized. The information provided herein can aid in the correct identification and differentiation of the products resulting from hydrogenation of Δ^8- or Δ^9-THC. The identification and confirmation of presumptive HHC metabolites is ongoing.[3]

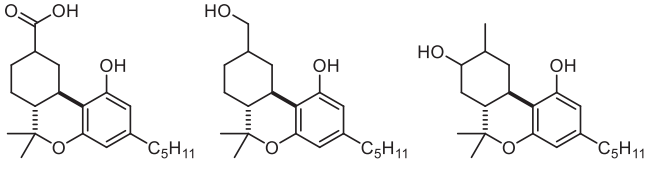


FIGURE 8 – Presumptive metabolites of HHC

References

1. Marzullo, P., Foschi, F., Coppini, D.A., *et al.* Cannabidiol as the substrate in acid-catalyzed intramolecular cyclization. *J. Nat. Prod.* **83(10)**, 2894-2901 (2020).
2. Mechoulam, R., Lander, N., Varkony, T.H., *et al.* Stereochemical requirements for cannabinoid activity. *J. Med. Chem.* **23(10)**, 1068–1072 (1980).
3. Harvey, D.J. and Brown, N.K. *In vitro* metabolism of the equatorial C_{11}-methyl isomer of hexahydrocannabinol in several mammalian species. *Drug Metab. Dispos.* **19(3)**, 714-716 (1991).

Hydrogenated Phytocannabinoids Available from Cayman

More than 20 reference standards for:

- Hexahydrocannabinols (HHCs)
- Hexahydrocannabiphorols (HHCPs)
- Tetrahydrocannabidiols (H4-CBDs)

Hexahydrocannabinols (HHCs)

Item No.	Product Name
27500	9(R)-Hexahydrocannabinol
37913	9(R)-Hexahydrocannabinol (CRM)
35368	9(R)-Hexahydrocannabinol Acetate
27501	9(S)-Hexahydrocannabinol
37942	9(S)-Hexahydrocannabinol (CRM)
35369	9(S)-Hexahydrocannabinol Acetate

Other HHCs of Interest

Item No.	Product Name
36256	(±)-9-nor-9α-hydroxy Hexahydrocannabinol
36257	(±)-9-nor-9β-hydroxy Hexahydrocannabinol
36129	(±)-9α-hydroxy Hexahydrocannabinol
35266	(±)-9β-hydroxy Hexahydrocannabinol
36353	11-hydroxy-9(R)-Hexahydrocannabinol
36354	11-hydroxy-9(S)-Hexahydrocannabinol
36355	11-nor-9(R)-carboxy-Hexahydrocannabinol
36356	11-nor-9(S)-carboxy-Hexahydrocannabinol
36250	8(R)-hydroxy-9(R)-Hexahydrocannabinol
36249	8(S)-hydroxy-9(S)-Hexahydrocannabinol

View all HHC cannabinoids at www.caymanchem.com

Hexahydrocannabiphorols (HHCPs)

Item No.	Product Name
36346	9(R)-Hexahydrocannabiphorol
37847	9(R)-Hexahydrocannabiphorol Acetate
36347	9(S)-Hexahydrocannabiphorol
37848	9(S)-Hexahydrocannabiphorol Acetate

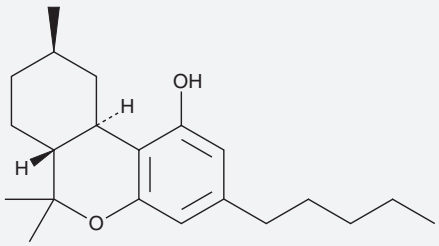
View all HHCP cannabinoids at www.caymanchem.com

Tetrahydrocannabidiols (H4-CBDs)

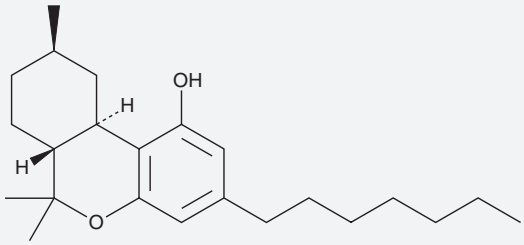
Item No.	Product Name
36350	1(R)-Tetrahydrocannabidiol (1(R)-H4-CBD)
36351	1(S)-Tetrahydrocannabidiol (1(S)-H4-CBD)

View all H4-CBD cannabinoids at www.caymanchem.com

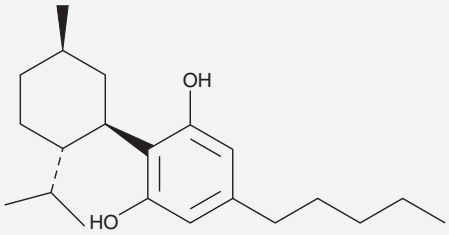
Major Hydrogenated Phytocannabinoids



9(R)-HHC
Item No. 27500



9(R)-HHCP
Item No. 36346



1(R)-H4-CBD
Item No. 36350

CAN'T FIND YOUR UNKNOWN?

Submit your GC-MS data (preferred) and any relevant information (suspected chemical class, matrix, solvent, GC parameters, etc.) to **techserv@caymanchem.com** and our scientists can help you solve your unknown.