

## 1. GENERAL INFORMATION

**IUPAC Name:** *rel*-2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol

**CFR:** Schedule I

**CAS#:** 70434-82-1

**Synonyms:** (±)-CP 47,497, (C7)-CP 47,497,  
5-(1,1-dimethylheptyl)-2-[(1R,3S)-3-hydroxycyclohexyl]-phenol

**Source:** DEA Reference Material Collection

**Appearance:** White powder

**Kovat's Index:** Pending

**$UV_{max}$  (nm):** 275.1

## 2. CHEMICAL AND PHYSICAL DATA

### 2.1 CHEMICAL DATA

Form	Chemical Formula	Molecular Weight	Melting Point (°C)
CP 47,497	C <sub>21</sub> H <sub>34</sub> O <sub>2</sub>	318	66.5-88.8

## 3. ADDITIONAL RESOURCES

[Forendex](#)

[Wikipedia](#)

Uchiyama, N.; Kikura-Hanajiri, R.; Ogata, J.; Goda, Y. Chemical analysis of synthetic cannabinoids as designer drugs in herbal products. *For. Sci. Int.* **2010**, *198*, 31-38.

## 4. QUALITATIVE DATA

### 4.1 NUCLEAR MAGNETIC RESONANCE

#### Method NMR CD<sub>3</sub>OD

*Sample Preparation:* Dilute analyte to ~5 mg/mL in CD<sub>3</sub>OD containing TMS for 0 ppm reference and maleic acid as quantitative internal standard.

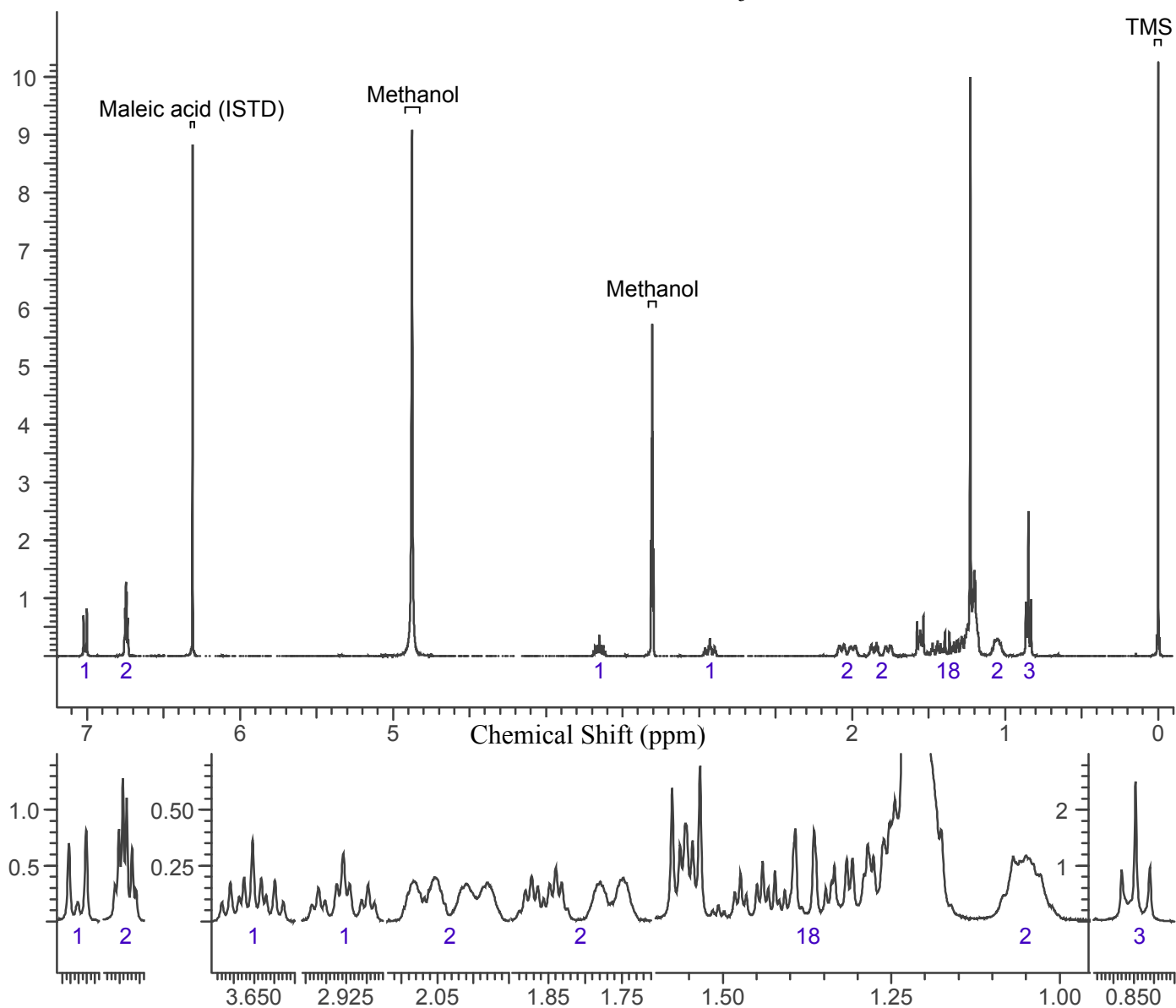
*Instrument:* 400 MHz NMR spectrometer

*Parameters:* Spectral width: at least containing -3 ppm through 13 ppm

Pulse angle: 90°

Delay between pulses: 45 seconds

<sup>1</sup>H NMR: CP 47,497; Lot 0419860-11; CD<sub>3</sub>OD; 400 MHz



## 4.2 Gas Chromatography/Mass Spectrometry

*Sample Preparation:* Dilute analyte ~ 1 mg/mL in methanol

***Instrument:*** Agilent gas chromatograph operated in split mode with MS detector

***Column:*** DB-1 MS (or equivalent); 30m x 0.25 mm x 0.25  $\mu$ m

***Carrier Gas:*** Helium at 1 mL/min

***Temperatures:*** Injector: 280°C

MSD transfer line: 280°C

MS Source: 230°C

MS Quad: 150°C

Oven program:

1) 100°C initial temperature for 1.0 min

2) Ramp to 300°C at 12 °C/min

3) Hold final temperature for 9.0 min

***Injection Parameters:*** Split Ratio = 20:1, 1  $\mu$ L injected

***MS Parameters:*** Mass scan range: 30-550 amu

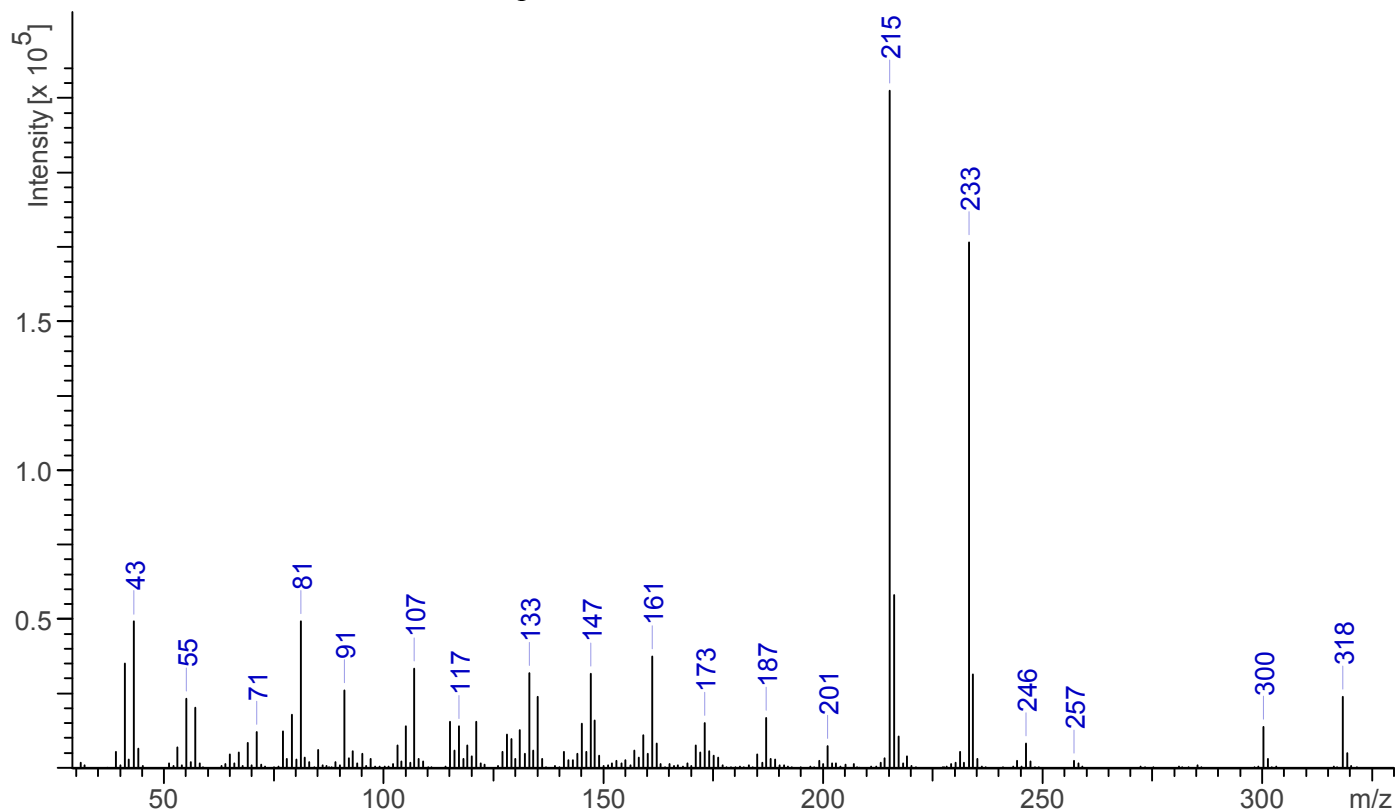
Threshold: 100

Tune file: stune.u

Acquisition mode: scan

***Retention Time:*** 16.294 min

EI Mass Spectrum: CP 47,497; Lot 0419860-11



### 4.3 INFRARED SPECTROSCOPY (FTIR)

**Instrument:** FTIR with diamond ATR attachment (3 bounce)

**Scan Parameters:**  
Number of scans: 32  
Number of background scans: 32  
Resolution: 4 cm<sup>-1</sup>  
Sample gain: 8  
Aperture: 150

FTIR ATR (Diamond, 3 Bounce): CP 47,497; Lot 0419860-11

