

# ISOLATION AND CHARACTERIZATION OF COMPOUNDS FROM CINNAMON OIL (CINNAMOMUM BURMANII) DISTILLATION RESIDU

*Maria Dewi Astuti, Lathifah Fauzi, Kamilia Mustikasari*

Program Studi Kimia FMIPA Universitas Lambung Mangkurat  
Jalan A Yani Km 36 Banjarbaru Kalimantan Selatan  
e-mail: [mdastuti@ulm.ac.id](mailto:mdastuti@ulm.ac.id)

## ABSTRACT

*This study aimed to isolate and characterize compounds from the distillation residue of cinnamon oil from Loksado, South Kalimantan. Cinnamon (*Cinnamomum burmanii*) distillation residue was extracted with methanol as solvent. The methanol extract was fractionated by liquid vacuum chromatography to obtain fractions A, B, C, and D. The crystals contained in fraction C were washed with cold n-hexane to obtain 5.4 mg of yellow isolate (FC1). FC1 isolates were characterized by UV-Vis, IR,  $^1\text{H-NMR}$ , and  $^{13}\text{C-NMR}$  spectrophotometers. UV spectra showed a maximum wavelength at 307, 316, and 321 nm indicating the presence of a conjugated or aromatic system. The infrared spectra showed  $\text{-C=O}$ ,  $\text{-OH}$ ,  $\text{C-O}$ ,  $\text{C-H}$ ,  $\text{C-N}$ , and  $\text{C=N}$  groups. The  $^1\text{H-NMR}$  spectra showed the presence of aromatic protons at 6.38 ppm (1H, d,  $J=9.5$  Hz), 7.67 ppm (1H, d,  $J=9.5$  Hz), 7.29 ppm (1H, d,  $J=8$  Hz), 7.44 ppm (1H, d,  $J=8$  Hz), and 7.49 ppm (1H, t) and there was a methyl proton (acetyl group) at H 2.13 ppm (3H,s). The  $^{13}\text{C-NMR}$  spectra showed the presence of a  $\text{C=O}$  ketone group at 207.26 ppm and there were 9  $\text{C-sp}^2$  at 116.9; 119.0; 124.6; 128.1; 132.0; 143.7; 154.3; 161.0 ppm, which  $\delta_{\text{C}}$  161.0 ppm was C-oxyaryl. Based on UV, IR,  $^1\text{H}$  and  $^{13}\text{C-NMR}$  spectra data, FC1 isolate was suggested as an isoquinoline alkaloid substituted by OH and acetyl groups.*

**Keywords:** *distillation residu, Cinnamomum burmanii, alkaloid, isoquinoline.*