

Article

Diarylheptanoid Glycosides of *Morella salicifolia* Bark

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Supplementary Material

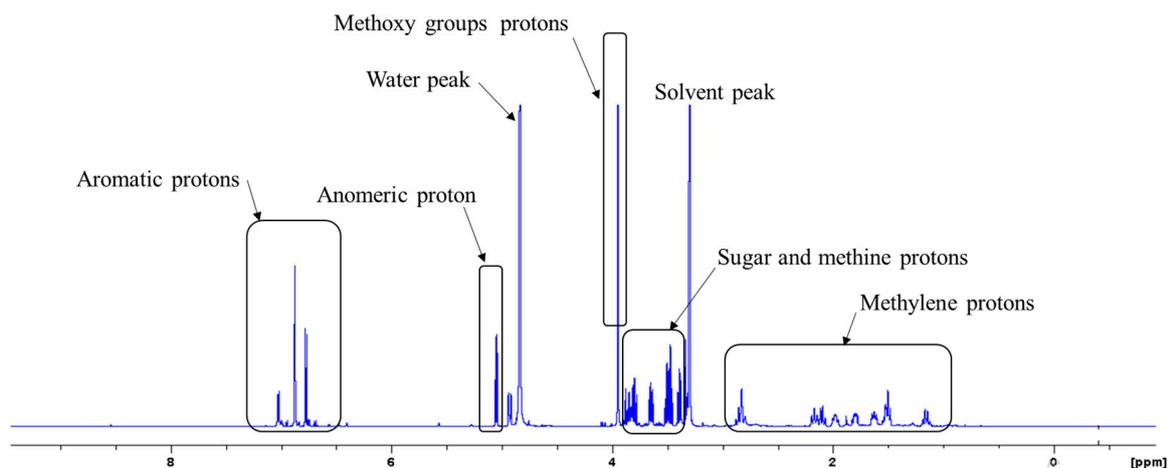


Figure S1. ¹H NMR spectrum (600 MHz, methanol-d₄, 298 K) of compound 1: salicimeckol (7-hydroxymyricanol 5-O-β-D-glucopyranoside).

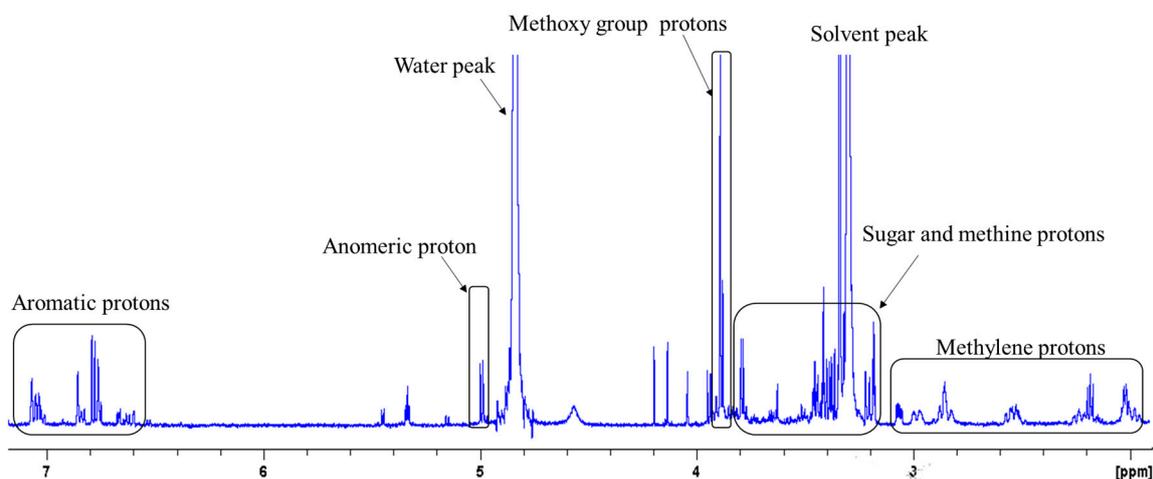


Figure S2. ¹H-NMR spectrum (600 MHz, methanol-d₄, 298 K) of compound 2: salicireneol A (juglanin B 3-O-β-D-glucopyranoside).

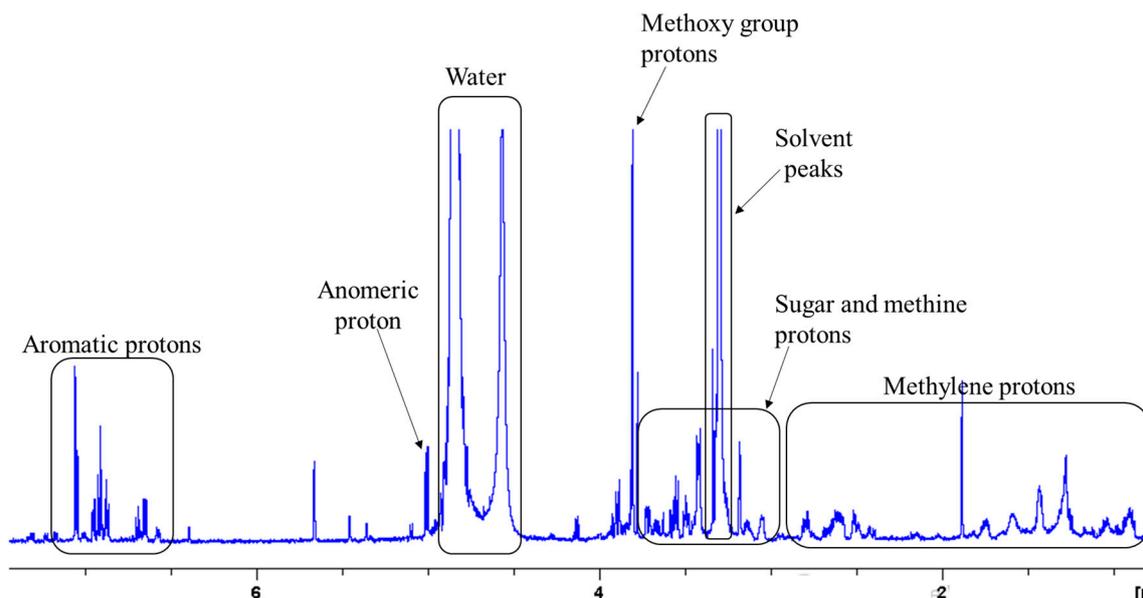


Figure S3. ¹H-NMR spectrum (600 MHz, methanol-d₄, 298 K) of compound 3: salicireneol B (16-hydroxyjuglanin B 17-O-β-D-glucopyranoside).

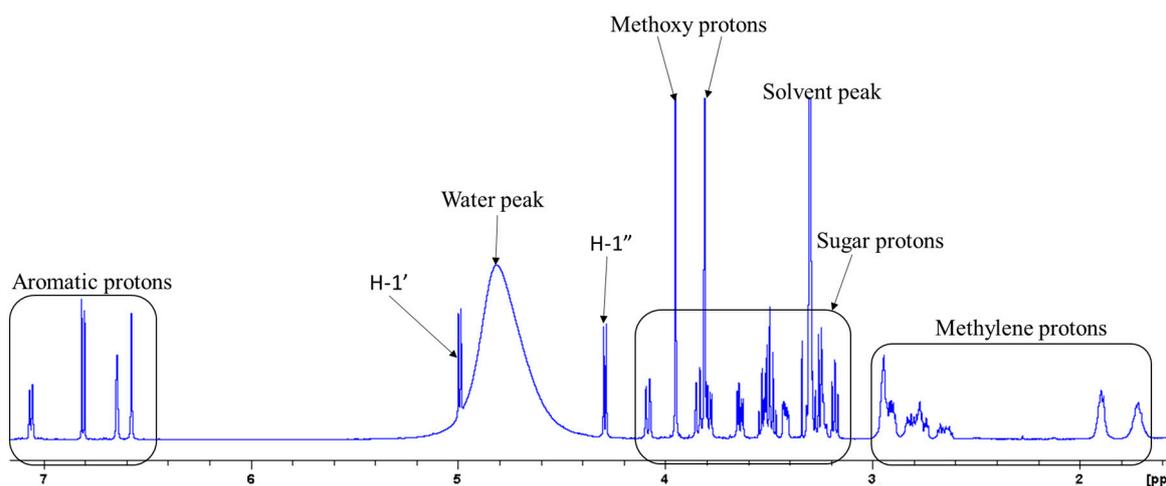


Figure S4. ¹H-NMR spectrum (600 MHz, methanol-d₄, 298 K) of compound 4: saliclaireone A (myricanone 5-O-β-D-glucopranosyl-(1-6)-β-D-glucopyranoside).

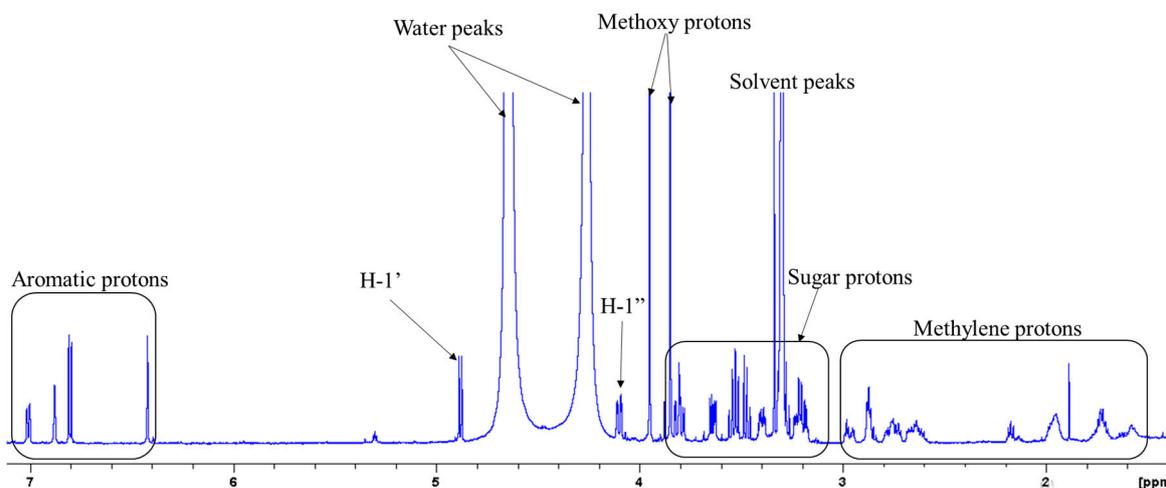


Figure S5. ¹H-NMR spectrum (600 MHz, methanol-d₄, 298 K) of compound 5: saliclaireone B (neomyricanone 5-O-β-D-glucopranosyl-(1-6)-β-D-glucopyranoside).

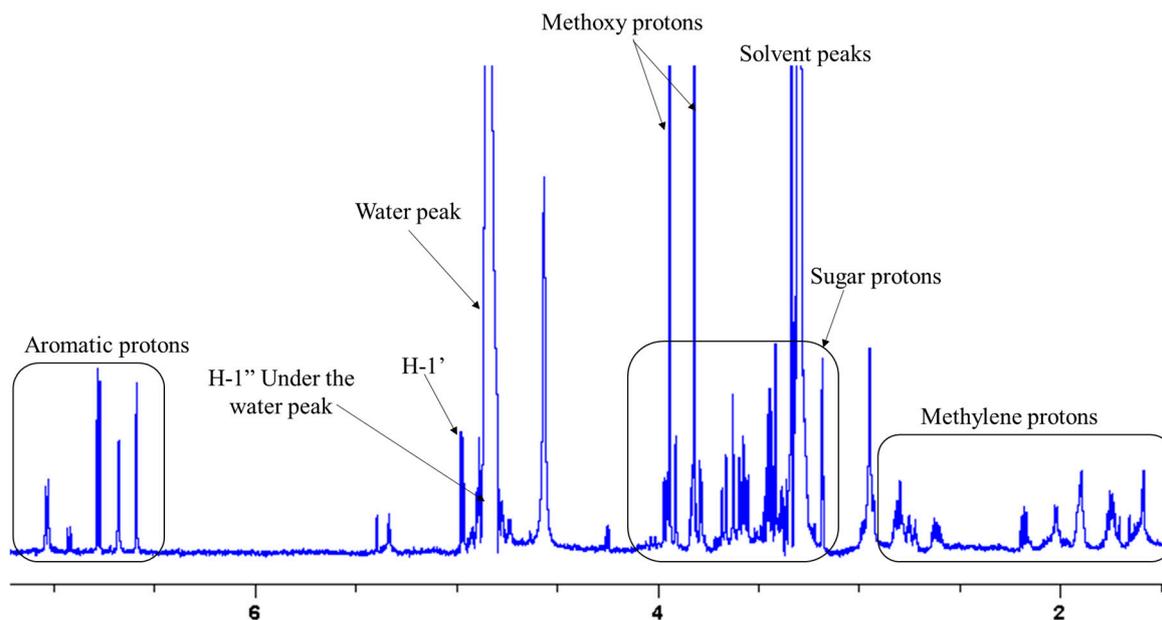


Figure S6. $^1\text{H-NMR}$ spectrum (600 MHz, methanol- d_4 , 298 K) of compound **6**: saliclaireone C (myricanone 17- O - α -L-arabinofuranosyl-(1-6)- β -D-glucopyranoside).

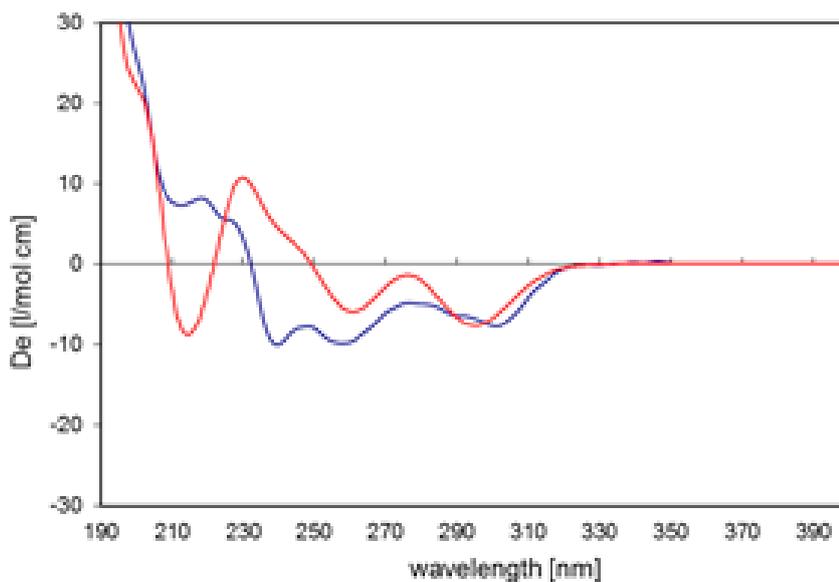


Figure S7. **Blue:** Experimental CD spectrum of myricanol. **Red:** Averaged CD spectrum for the *S,Sa* (87%) and *S,Ra* form (13%). TDDFT: RB3LYP/6-31G(d,p), nstates = 30. Calculated spectrum was red-shifted by -0.15 eV and scaled by a factor of 0.67.