Supplementary Materials

Semi-synthesis of C-ring Cyclopropyl Analogues of Fraxinellone and Their Insecticidal Activity against *Mythimna separata* Walker

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| Identification code | 3a |
|---|--|
| Empirical formula | C23H24O5 |
| Formula weight | 380.42 |
| Temperature/K | 293(2) |
| Space group | P212121 |
| a/Å | 8.2081(4) |
| b/Å | 14.3113(8) |
| c/Å | 16.7802(9) |
| α/° | 90.00 |
| β/° | 90.00 |
| γ/° | 90.00 |
| Volume/Å ³ | 1971.15(18) |
| Z | 4 |
| Qcalcg/cm ³ | 1.282 |
| µ/mm ⁻¹ | 0.731 |
| F(000) | 808.0 |
| Crystal size/mm ³ | $0.35 \times 0.13 \times 0.10$ |
| Radiation | $CuK\alpha~(\lambda = 1.54178)$ |
| 2⊖ range for data collection/° | 8.12 to 132.04 |
| Index ranges | $-9 \le h \le 4$, $-15 \le k \le 16$, $-19 \le l \le 13$ |
| Reflections collected | 4353 |
| Independent reflections | 2992 [$R_{int} = 0.0224$, $R_{sigma} = N/A$] |
| Data/restraints/parameters | 2992/0/257 |
| Goodness-of-fit on F ² | 1.081 |
| Final R indexes [I>=2σ (I)] | $R_1 = 0.0434$, $wR_2 = 0.1001$ |
| Final R indexes [all data] | $R_1 = 0.0568$, $wR_2 = 0.1097$ |
| Largest diff. peak/hole / e Å ⁻³ | 0.10/-0.12 |
| Flack parameter | 0.0(3) |

Table 1. Crystal data and structure refinement for **3a**.



Figure S1. ¹H and ¹³C NMR (MeOD) spectra of compound **2b**.





Figure S2. ¹H and ¹³C NMR (CDCl₃) spectra of compound 2c.

Figure S3. ¹H, DEPT135 and ¹³C NMR (CDCl₃) spectra of compound 3a.



Figure S4. ¹H, DEPT135 and ¹³C NMR (CDCl₃) spectra of compound **3b**.

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¹H NMR (500 MHz, Chloroform-d)



Figure S5. ¹H, DEPT135 and ¹³C NMR (CDCl₃) spectra of compound 3c.



Figure S6. HR-ESI-MS spectra of compound 3a.



+TOF MS *m*/*z*: Found 403.1496 [M+Na]⁺ (calcd for C₂₃H₂₄NaO₅, 403.1521).

Figure S7. HR-ESI-MS spectra of compound 3b.

+TOF MS *m*/*z*: Found 417.1647 [M+Na]⁺ (calcd for C₂₄H₂₆NaO₅, 417.1678).



Figure S8. HR-ESI-MS spectra of compound 3c.



