## **Supporting Information**

## Rec. Nat. Prod. 13:5 (2019) 424-428

## Polyacetylenes from the Roots of Aralia dumetorum

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| Table of Contents  | Page |
|--|------|
| Figure S1: 1H NMR spectrum of 1 in CDCl <sub>3</sub> (400 MHz).                | 2    |
| Figure S2: DEPT spectrum of 1 in CDCl <sub>3</sub> (100 MHz).                  | 3    |
| Figure S3: HSQC spectrum of 1 in CDCl <sub>3</sub> .                           | 4    |
| Figure S4: 1H-1H COSY spectrum of 1 in CDCl <sub>3</sub> .                     | 5    |
| Figure S5: HMBC spectrum of 1 in CDCl <sub>3</sub> .                           | 6    |
| Figure S6: ROESY spectrum of 1 in CDCl <sub>3</sub> .                          | 7    |
| Figure S7: IR spectrum of 1  | 8    |
| Figure S8: HRESIMS of 1  | 9    |
| Table S1. <sup>1</sup> H NMR and <sup>13</sup> C NMR data of compounds 1 and 2 | 10   |

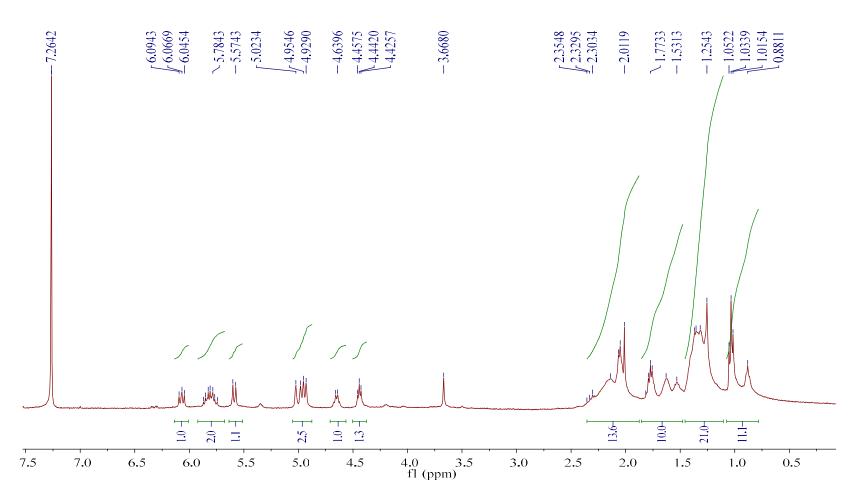


Figure S1: 1H NMR spectrum of 1 in CDCl<sub>3</sub> (400 MHz).

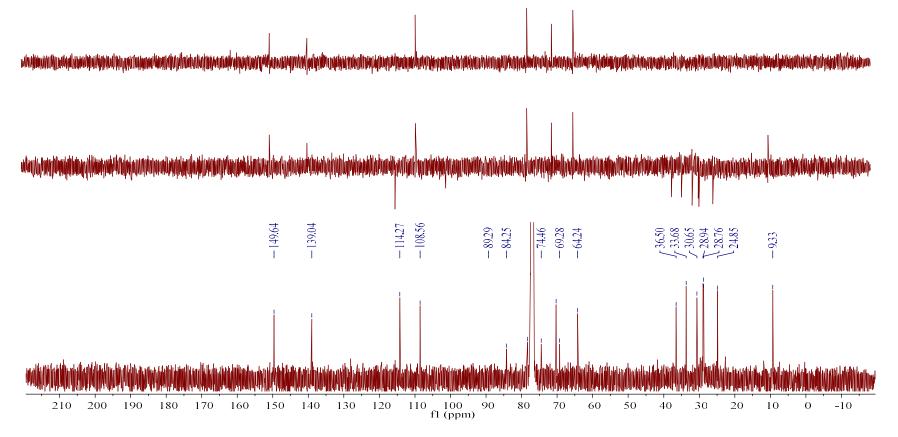


Figure S2: DEPT spectrum of 1 in CDCl<sub>3</sub> (100 MHz).

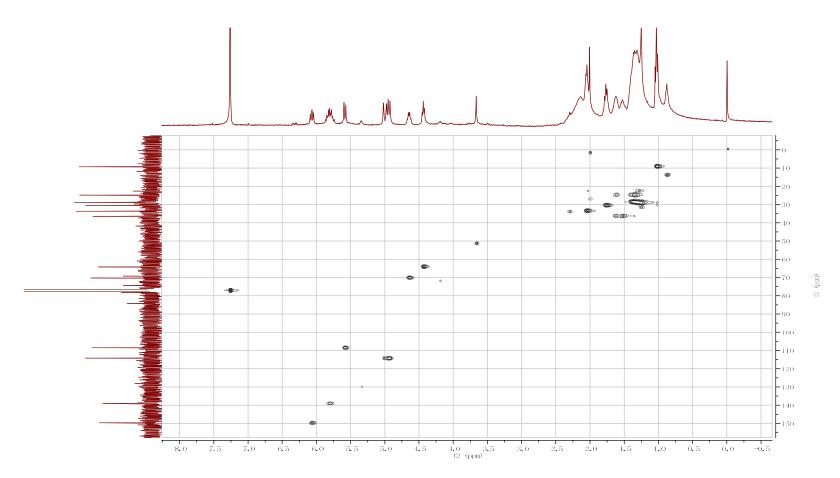


Figure S3: HSQC spectrum of 1 in CDCl<sub>3</sub>.

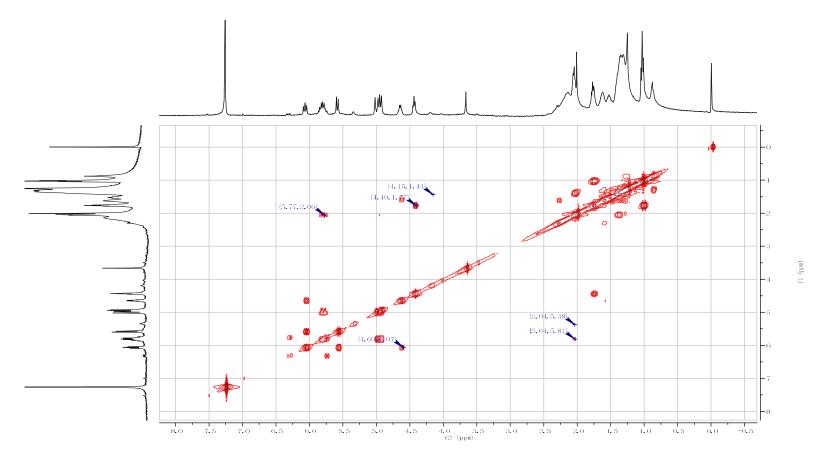
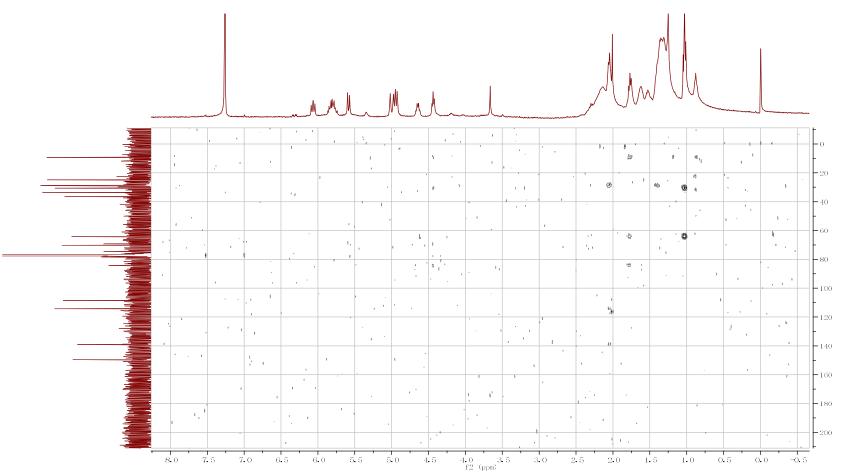


Figure S4: 1H-1H COSY spectrum of 1 in CDCl<sub>3</sub>.



fil (ppm)



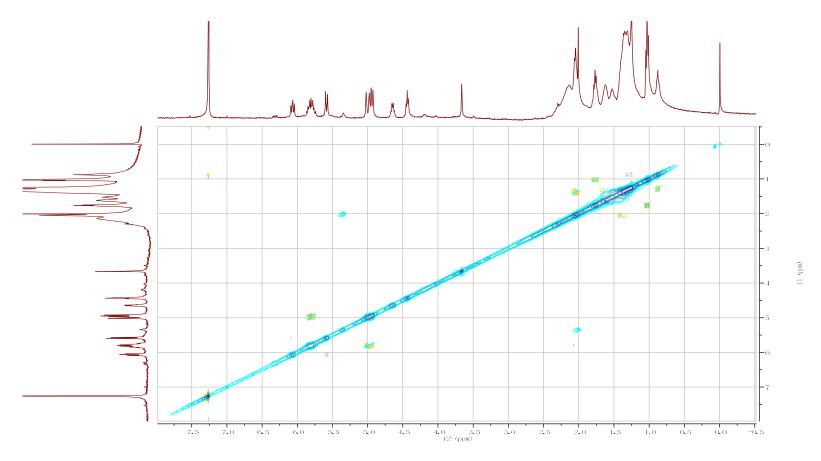


Figure S6: ROESY spectrum of 1 in CDCl<sub>3</sub>.

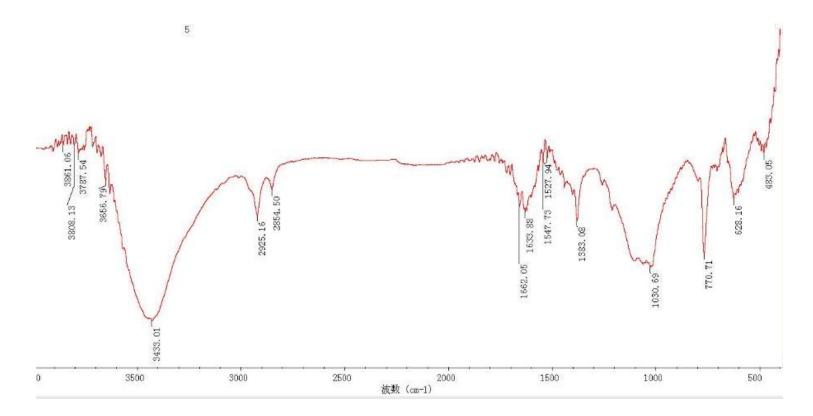


Figure S7: IR spectrum of 1

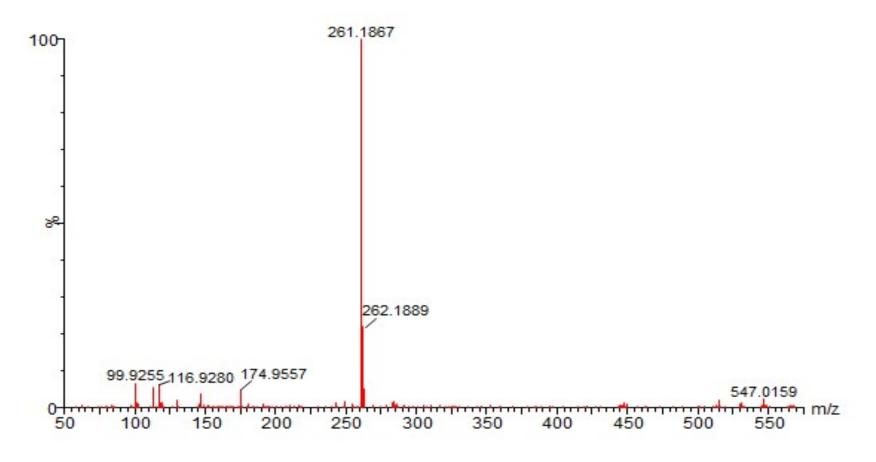


Figure S8: HRESIMS of 1

| Position | 1  |                          | 2  |                          |
|----------|--|--------------------------|--|--------------------------|
|          | <sup>1</sup> H NMR   | <sup>13</sup> C NMR      | <sup>1</sup> H NMR   | <sup>13</sup> C NMR      |
| 1        | 1.03 (3H, <i>t</i> , <i>J</i> = 7.2)   | 9.3 (CH <sub>3</sub> )   | 1.02 (3H, t, J = 6.9)  | 9.3 (CH <sub>3</sub> )   |
| 2        | 1.77 (2H, <i>m</i> )   | 30.6 (CH <sub>2</sub> )  | 1.74 (2H, <i>m</i> )   | 30.6 (CH <sub>2</sub> )  |
| 3        | 4.44 (1H, t, J = 6.0)  | 64.2 (CH)                | 4.41 (1H, <i>t</i> , <i>J</i> = 6.6)   | 64.3 (CH)                |
| 4        |  | 84.2 (C)                 |  | 82.9 (C)                 |
| 5        |  | 69.1 (C)                 |  | 70.3 (C)                 |
| 6        |  | 74.5 (C)                 |  | 73.6 (C)                 |
| 7        |  | 78.3 (C)                 |  | 77.9 (C)                 |
| 8        | 5.59 (1H, <i>d</i> , <i>J</i> = 10.8 Hz)   | 108.6 (CH)               | 5.75 (1H, <i>brd</i> , <i>J</i> = 15.9 Hz)   | 108.2 (CH)               |
| 9        | 6.06 (1H, <i>dd</i> , <i>J</i> = 10.8, 8.3)  | 149.6 (CH)               | 6.31 (1H, <i>dd</i> , <i>J</i> = 15.9, 6.0)  | 149.6 (CH)               |
| 10       | 4.65 (1H, <i>m</i> )   | 69.3 (CH)                | 4.19 (1H, <i>ddt</i> , <i>J</i> = 6.8, 6.0, 1.2)                                   | 72.0 (CH)                |
| 11       | 1.50 (2H, <i>m</i> )   | 36.5 (CH <sub>2</sub> )  | 1.50 (2H, q, J = 6.8)  | 36.8 (CH <sub>2</sub> )  |
| 12       | 1.30 (2H, <i>m</i> )   | 24.8 (CH <sub>2</sub> )  | 1.31 (2H, <i>m</i> )   | 25.0 (CH <sub>2</sub> )  |
| 13       | 1.31 (2H, m)   | 28.9 (CH <sub>2</sub> )  | 1.31 (2H, <i>m</i> )   | 29.0 (CH <sub>2</sub> )  |
| 14       | 1.35 (2H, <i>m</i> )   | 28.8 (CH <sub>2</sub> )  | 1.37 (2H, <i>m</i> )   | 28.8 (CH <sub>2</sub> )  |
| 15       | 2.05 (2H, <i>m</i> )   | 33.7 (CH <sub>2</sub> )  | 2.02 (2H, $q$ , $J = 7.0$ )  | 33.6 (CH <sub>2</sub> )  |
| 16       | 5.81 (1H, <i>m</i> )   | 139.0 (CH)               | 5.78 (1H, <i>ddt</i> , , <i>J</i> = 16.8, 10.4, 7.0)                               | 138.9 (CH)               |
| 17       | 5.00 (1H, <i>brd</i> , <i>J</i> = 17.2)<br>4.94 (1H, <i>brd</i> , <i>J</i> = 10.2) | 114.3 (CH <sub>2</sub> ) | 4.98 (1H, <i>brd</i> , <i>J</i> = 16.8)<br>4.94 (1H, <i>brd</i> , <i>J</i> = 10.4) | 114.4 (CH <sub>2</sub> ) |

Table S1. <sup>1</sup>H NMR and <sup>13</sup>C NMR data of compounds 1 and 2