Diphenyl Ethers from a Marine-Derived Aspergillus sydowii

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Figure S1. IR spectrum of compound 1



Figure S2. UV spectrum of compound 1



Figure S4. ¹³C NMR spectrum of compound 1 in DMSO-*d*₆ (150 MHz)



Figure S6. HSQC spectrum of compound 1 in DMSO-d6



Figure S7. COSY spectrum of compound 1 in DMSO-d6



Figure S8. HRESIMS of compound 1



Figure S9. IR spectrum of compound 2



Figure S10. UV spectrum of compound 2



Figure S12. ¹³C NMR spectrum of compound 2 in CD₃OD (150 MHz)



Figure S14. HSQC spectrum of compound **2** in CD3OD



Figure S8. HRESIMS of compound 2



Figure S17. IR spectrum of compound 3



Figure S18. UV spectrum of compound 3



Figure S20. ¹³C NMR spectrum of compound 3 in DMSO-*d*₆ (150 MHz)



Figure S22. HSQC spectrum of compound 3 in DMSO-d6



Figure S23. J Resolved HSQC spectrum of compound 3 in CD3OD



| m/z | | lon | Formula | Abundance |] | | | | | | | | | |
|-----|----------|-------------|----------------|----------------------------|----------------------------------|----------------------------------|--|---|--|---|--|--|---|---|
| 4 | 473.1403 | (M+Na)+ | C22 H26 Na O10 | 207149.1 |] | | | | | | | | | |
| Be | st | Formula (M) | Ion Formula | Score | Cross Sco | Mass | Calc Mass | Calc m/z | Diff (ppm) | Abs Diff (ppm) | Mass Match | Abund Match | Spacing Match | DBE |
| | G . | C22 H26 O10 | C22 H26 Na O10 | 99.79 | | 450.151 | 450.1526 | 473.1418 | 3.46 | 3.46 | 99.63 | 99.88 | 100 | 10 |
| | 4 | | C22 H26 010 | CZ2 H26 O10 CZ2 H26 Na O10 | C22 H26 O10 C22 H26 Na O10 99.79 | C22 H26 OT0 C22 H26 Na OT0 99.79 | C22 H26 010 C22 H26 Na 010 99.79 430.131 | C22 H26 O10 C22 H26 Na O10 99.79 430.131 430.1320 | CZZ H26 010 CZZ H26 Na 010 99.79 450.151 450.1520 475.1410 | C22 H26 O10 C22 H26 N8 O10 99.79 430.131 430.1320 473.1410 5.40 | CZZ H26 010 CZZ H26 Na 010 99.79 450.151 450.1520 475.1410 5.40 5.40 | C22 H26 OT0 C22 H26 Na OT0 99.79 450.151 450.1520 473.1418 5.40 5.40 55.00 | C22 H26 OT0 C22 H26 Na OT0 99.79 450.151 450.1520 475.1418 5.40 5.40 5.40 55.00 | CZ2 H26 010 CZ2 H26 N8 010 99.79 430.131 430.1320 475.1410 5.40 5.40 5.40 5.40 5.40 |

Figure S25. HRESIMS of compound 3



Figure S26. IR spectrum of compound 4



Figure S27. UV spectrum of compound 4



Figure S29. ¹³C NMR spectrum of compound **4** in CD₃OD (150 MHz)



Figure S31. HSQC spectrum of compound 4 in CD3OD



Figure S32. COSY spectrum of compound 4 in CD₃OD



Figure S33. HRESIMS of compound 4



Figure S34. IR spectrum of compound 5



Figure S35. UV spectrum of compound 5



Figure S37. ¹³C NMR spectrum of compound 5 in CD₃OD (150 MHz)

Figure S41. HRESIMS of compound 5

Figure S42. IR spectrum of compound 6

Figure S43. UV spectrum of compound 6

Figure S45. ¹³C NMR spectrum of compound 6 in CD₃OD (150 MHz)

Figure S47. HSQC spectrum of compound 6 in CD3OD

Figure S48. COSY spectrum of compound 6 in CD3OD

Figure S49. HRESIMS of compound 6

8

Figure S51. ¹³C NMR spectrum of compound 6a in CD₃OD (150 MHz)

Figure S53. HSQC spectrum of compound 6a in CD3OD

Figure S54. LC-ESI-MS analysis of derivative of sugar moiety in **1**. The sugar moiety in **1** was determined as D-ribose by ESI mass spectrum (extraction ion chromatogram at m/z 417, t_{R} 20.1 for L-ribose, t_{R} 22.3 for D-ribose).

Figure S55. LC-ESI-MS analysis of derivatives of sugar moiety in **2**. The sugar moiety in **2** was determined as D-ribose by ESI mass spectrum (extraction ion chromatogram at m/z 417, $t_{\rm R}$ 20.1 for L-ribose, $t_{\rm R}$ 22.3 for D-ribose).

Figure S56. LC-ESI-MS analysis of derivatives of sugar moiety in **3**. The sugar moiety in **3** was determined as D-glucose by ESI mass spectrum (extraction ion chromatogram at m/z 447, $t_{\rm R}$ 20.6 for L-glucose, $t_{\rm R}$ 21.0 for D-glucose).

Figure S57. The internal transcribed spacers (ITS) sequence of strain FNA026