Diphenyl Ethers from a Marine-Derived Aspergillus sydowii

Ya-Nan Wang ^{2,†}, Yan-Hua Mou ^{1,†}, Yu Dong ^{1,2,†}, Yan Wu ², Bing-Yu Liu ², Jian Bai ², Dao-Jiang Yan ², Le Zhang ², Dan-Qing Feng ³, Yue-Hu Pei ^{1,*}and You-Cai Hu ^{2,*}

- ¹ Shenyang Pharmaceutical University, Shenyang 110016, China; mu_hua_jj@sina.com (Y.-H.M.); allenn7@foxmail.com (Y.D.);
- ² State Key Laboratory of Bioactive Substance and Function of Natural Medicines, Institute of Materia Medica, Chinese Academy of Medical Sciences and Peking Union Medical College, Beijing 100050, China; wangyanan@imm.ac.cn (Y.-N.W.); wuyan@imm.ac.cn (Y.W.); liubingyu@imm.ac.cn (B.-Y.L.); baijian@imm.ac.cn (B.J.); yandj@imm.ac.cn (Y.-D.J.); zhangle@imm.ac.cn (L. Z.)
- ³ State-Province Joint Engineering Laboratory of Marine Bioproducts and Technology, College of Ocean & Earth Sciences, Xiamen University, Xiamen, 361102, China; dqfeng@xmu.edu.cn (D.-Q.F.)
- * Correspondence: peiyueh@vip.163.com (Y.-H.P.); huyoucai@imm.ac.cn (Y.-C.H.); Tel.: +86-024-23986485 (Y.-H.P.); +86-010- 61271883 (Y.-C.H.)
- * These authors contributed equally to this paper.

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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