

Supplementary Material

Stress-Driven Discovery of New Angucycline-Type Antibiotics from a Marine *Streptomyces pratensis* NA-ZhouS1

Najeeb Akhter ¹, Yaqin Liu ², Bibi Nazia Auckloo ¹, Yutong Shi ¹, Kuiwu Wang ³, Juanjuan Chen ⁴, Xiaodan Wu ⁵, and Bin Wu ^{1,*}

¹ Ocean College, Zhejiang University, Hangzhou 310058, China. na_memon@yahoo.com; naz22ia@hotmail.com; 11434028@zju.edu.cn; wubin@zju.edu.cn

² Department of Chemistry, Zhejiang University, Hangzhou 301000, China. yaqin86@zju.edu.cn

³ Department of Chemistry, Zhejiang Gongshang University, Hangzhou 310012, China. wkwnpc@zjgsu.edu.cn

⁴ Key Laboratory of Applied Marine Biotechnology, Ningbo University, Chinese Ministry of Education, Ningbo 315211, China. chenjuanjuan@nbu.edu.cn.

⁵ Centre of Analysis and Measurement, Zhejiang University, Hangzhou 310058, China. wxd_zju@163.com

* Correspondence: wubin@zju.edu.cn; Tel. /Fax: +86-05802092258

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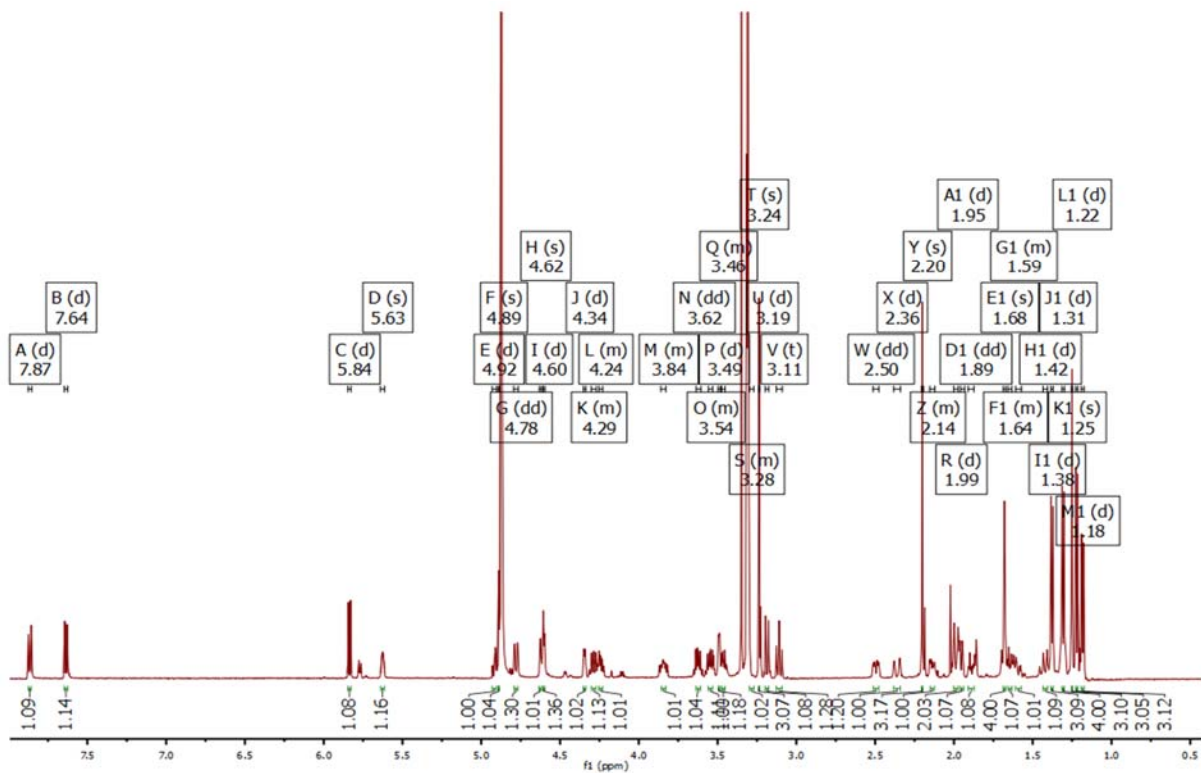


Figure S 1. ^1H NMR spectrum in MeOD for compound 1.

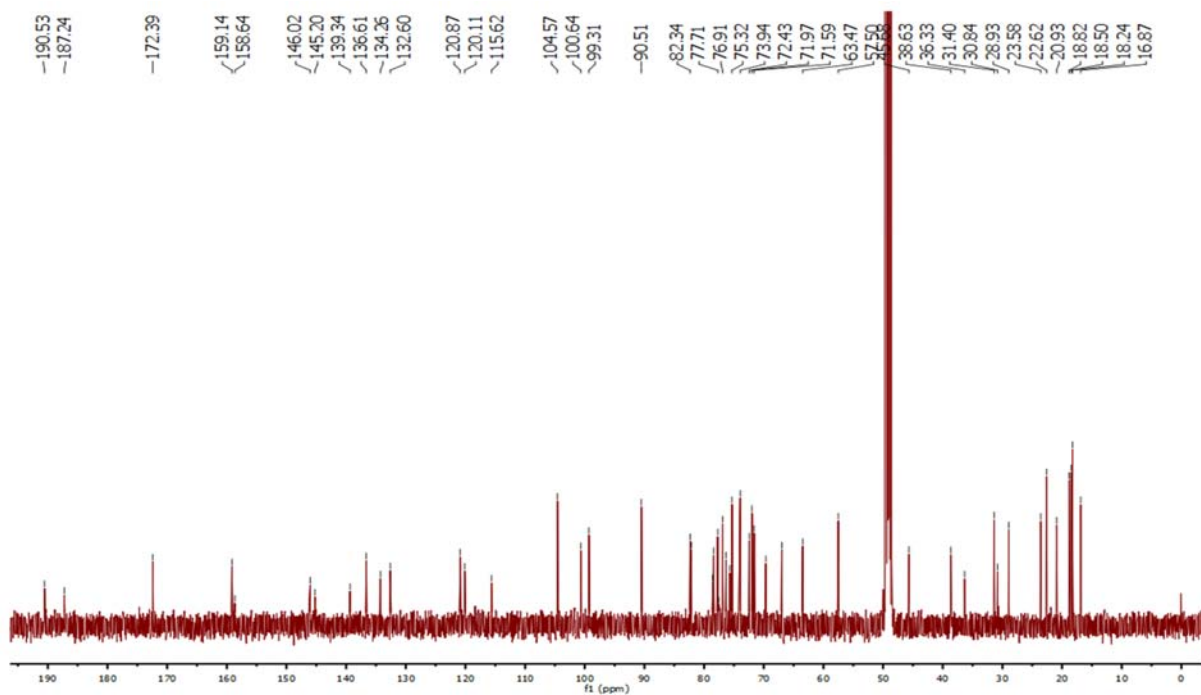


Figure S 2. ^{13}C NMR spectrum in MeOD for compound 1.

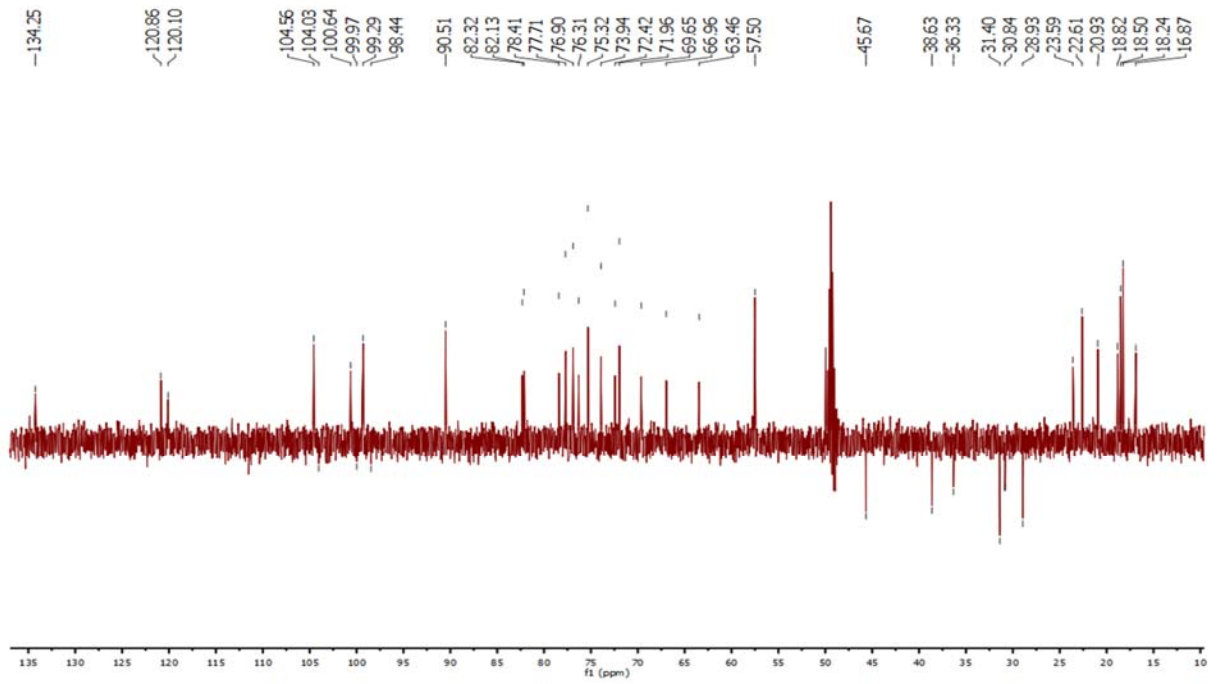


Figure S 3. DEPT-135 spectrum in MeOD for compound 1.

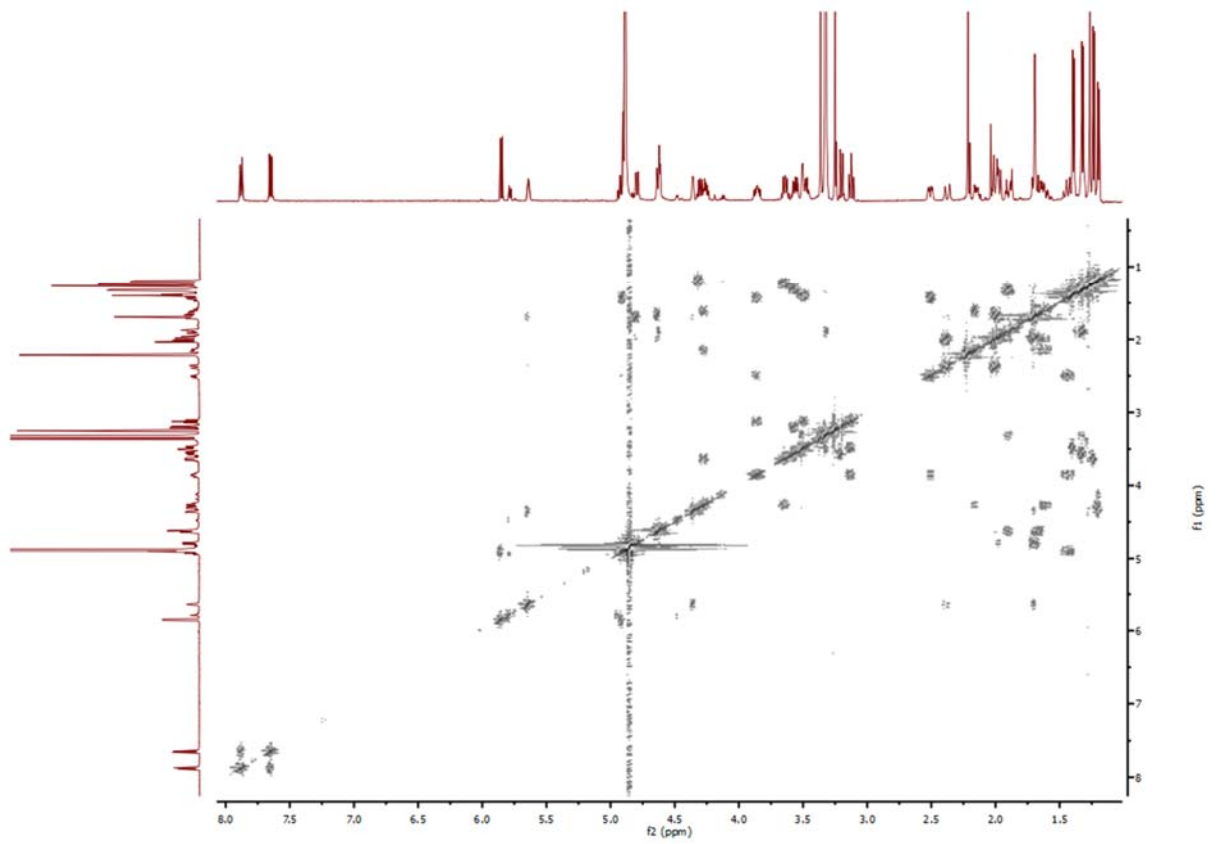


Figure S 4. COSY spectrum in MeOD for compound 1.

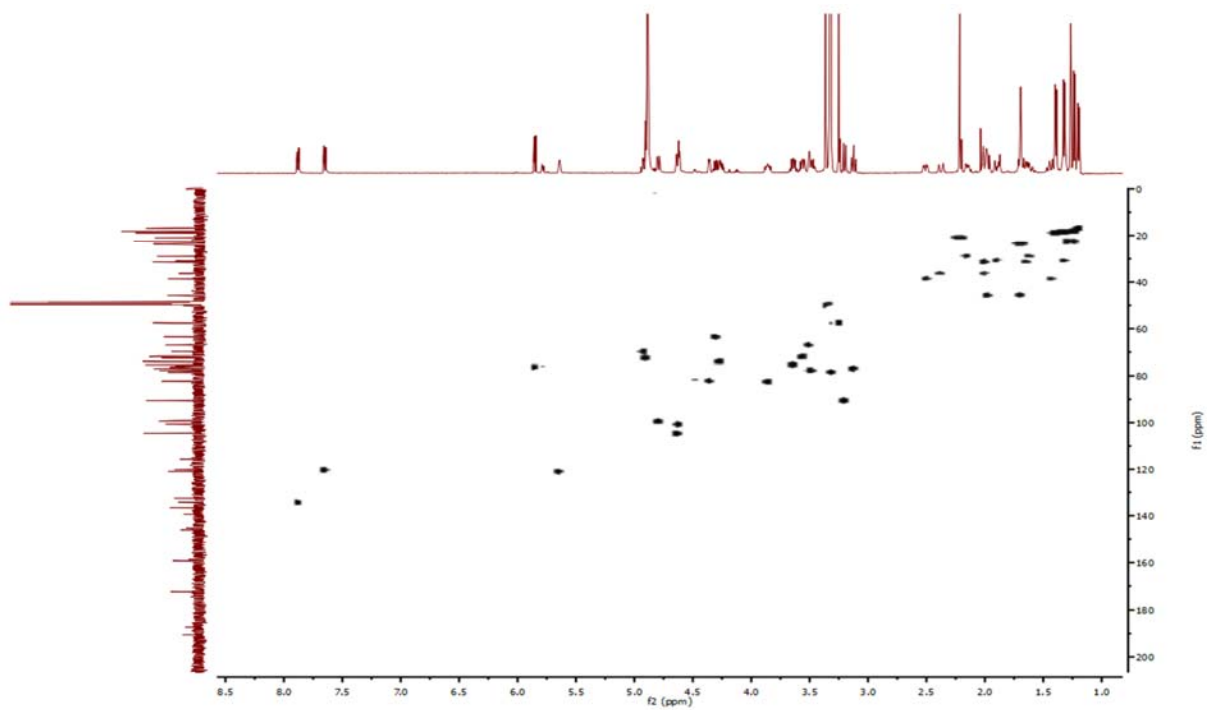


Figure S 5. HSQC spectrum in MeOD for compound 1.

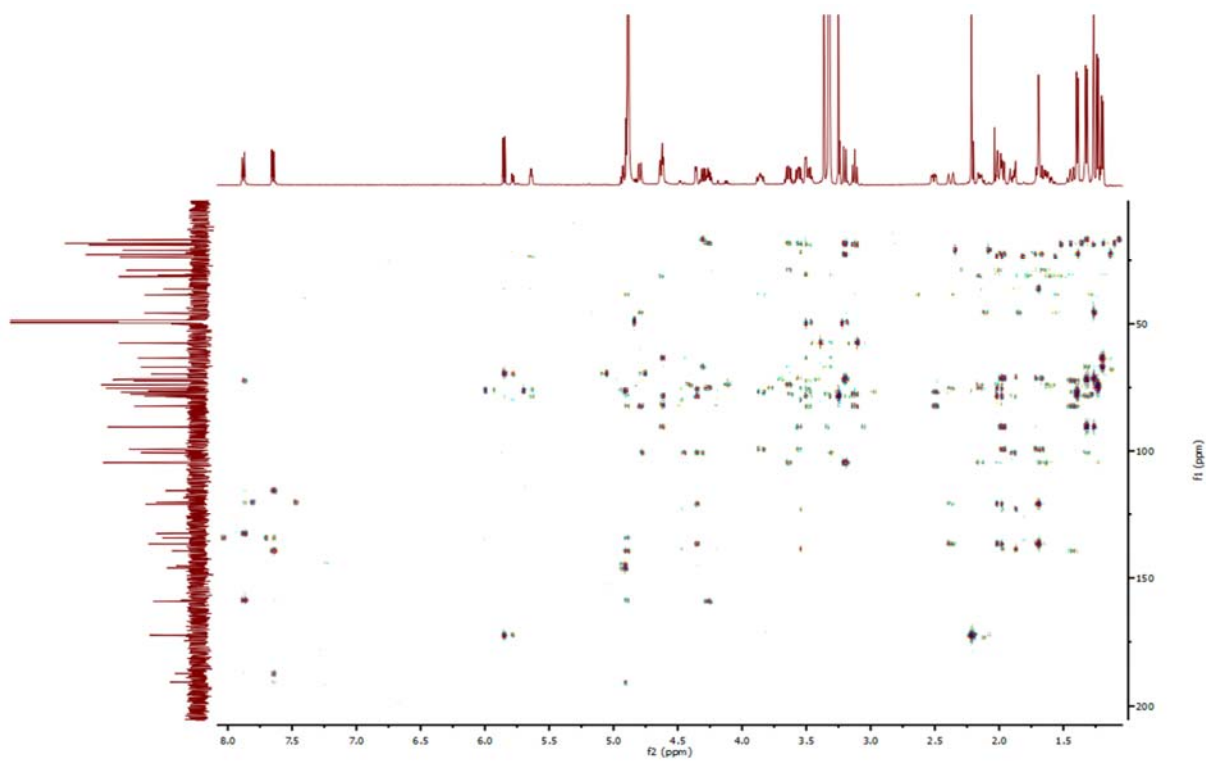


Figure S 6. HMBC spectrum in MeOD for compound 1.

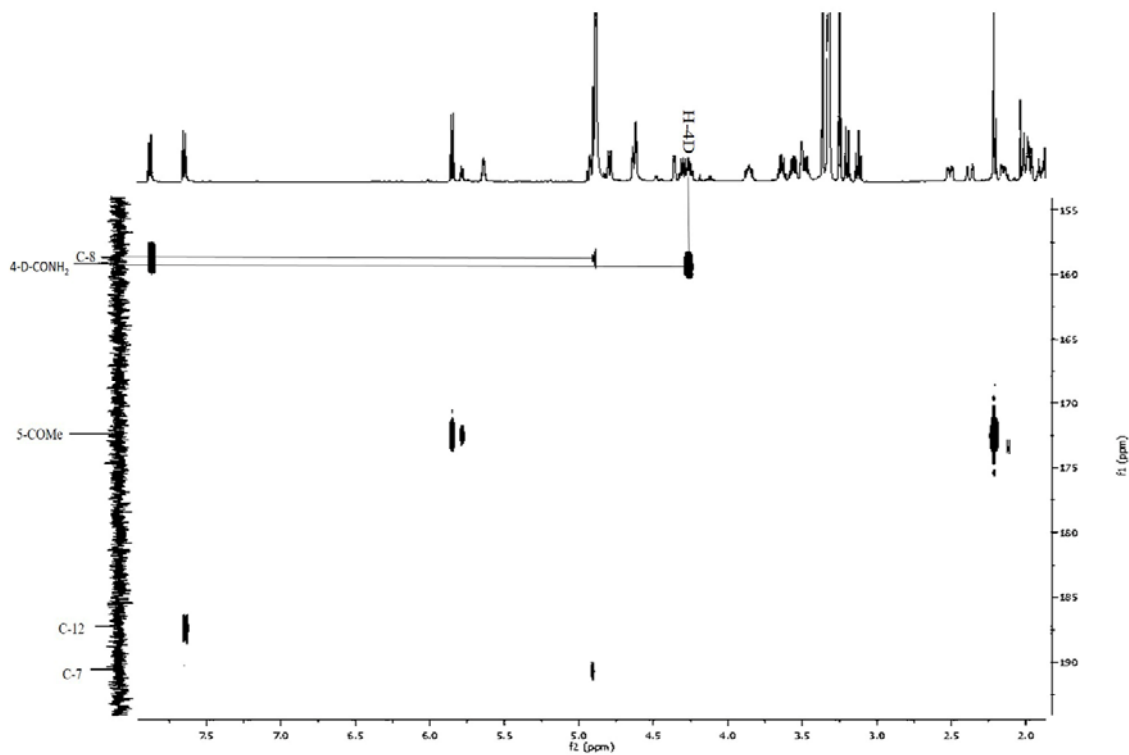


Figure S 7. HMBC-2 spectrum in MeOD for compound **1**. Enlarged view of the spectrum to clarify the cross peaks of H-4D to carbamoyl carbon resonated at δ c 159.6.

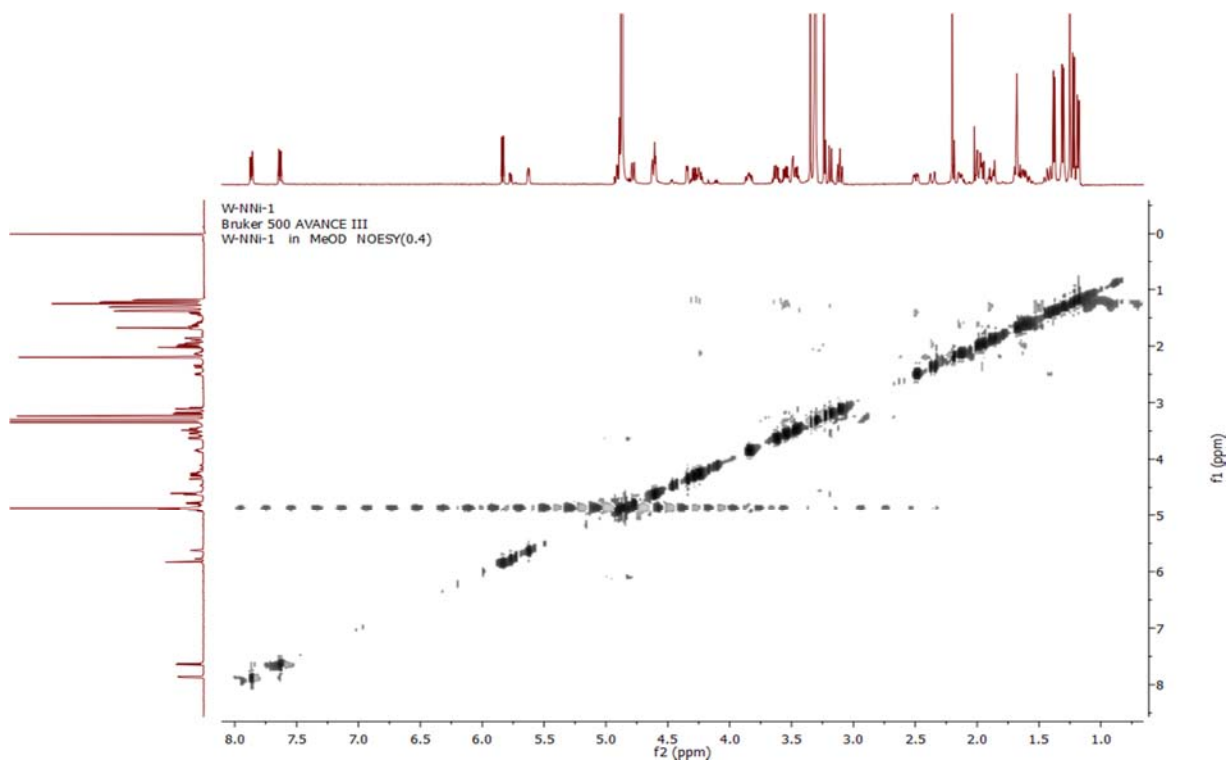


Figure S 8. NOESY spectrum in MeOD for compound **1**.

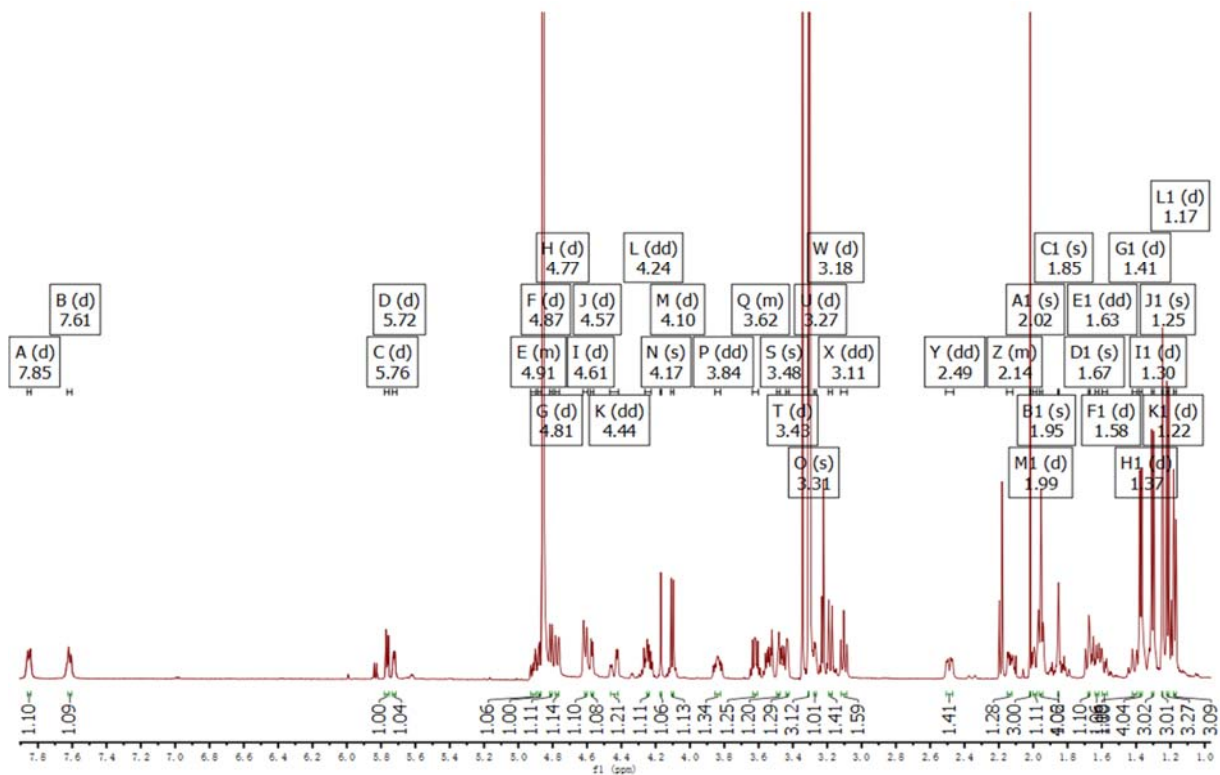


Figure S 9. ^1H NMR spectrum in MeOD for compound 2.

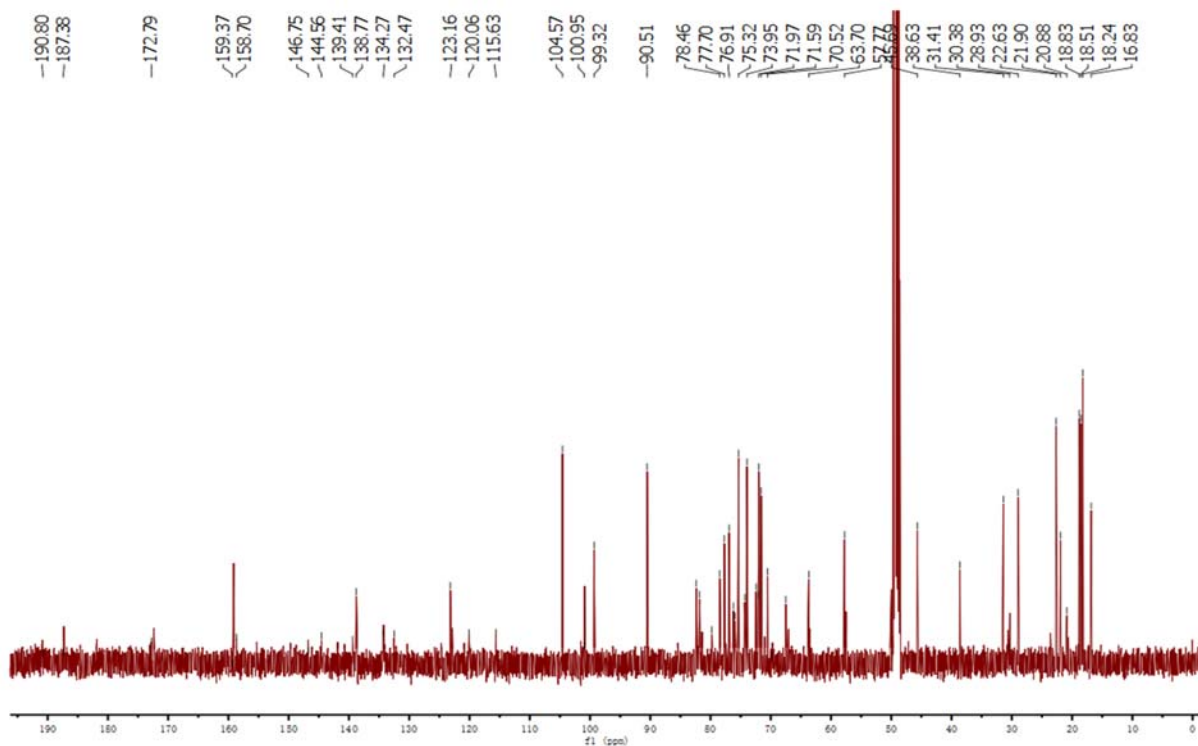


Figure S 10. ^{13}C NMR spectrum in MeOD for compound 2.

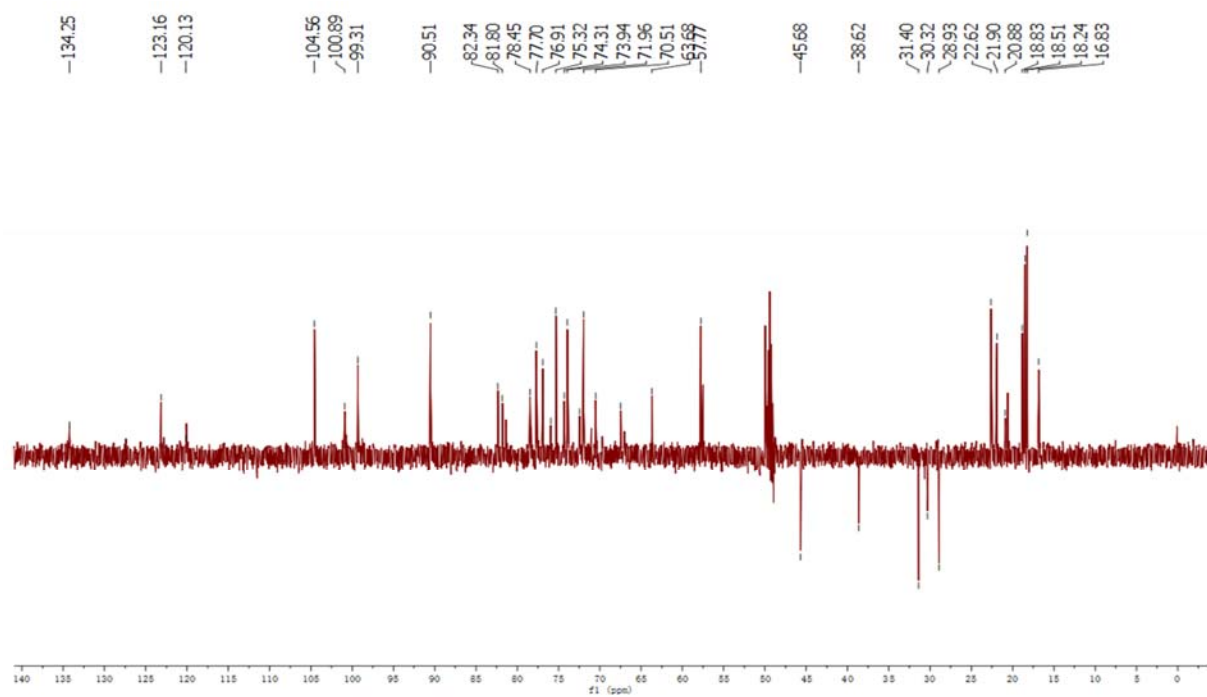


Figure S 11. DEPT-135 spectrum in MeOD for compound 2.

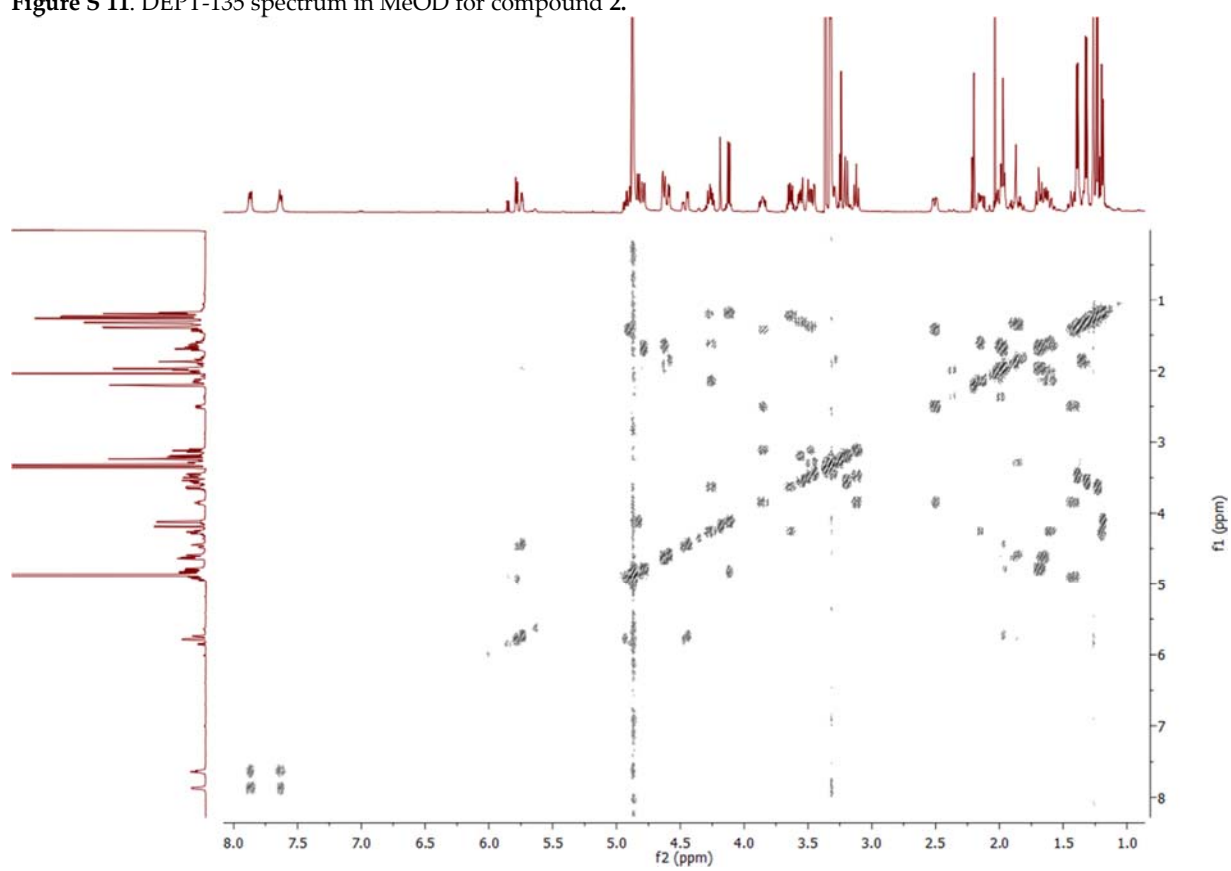


Figure S 12. COSY spectrum in MeOD for compound 2.

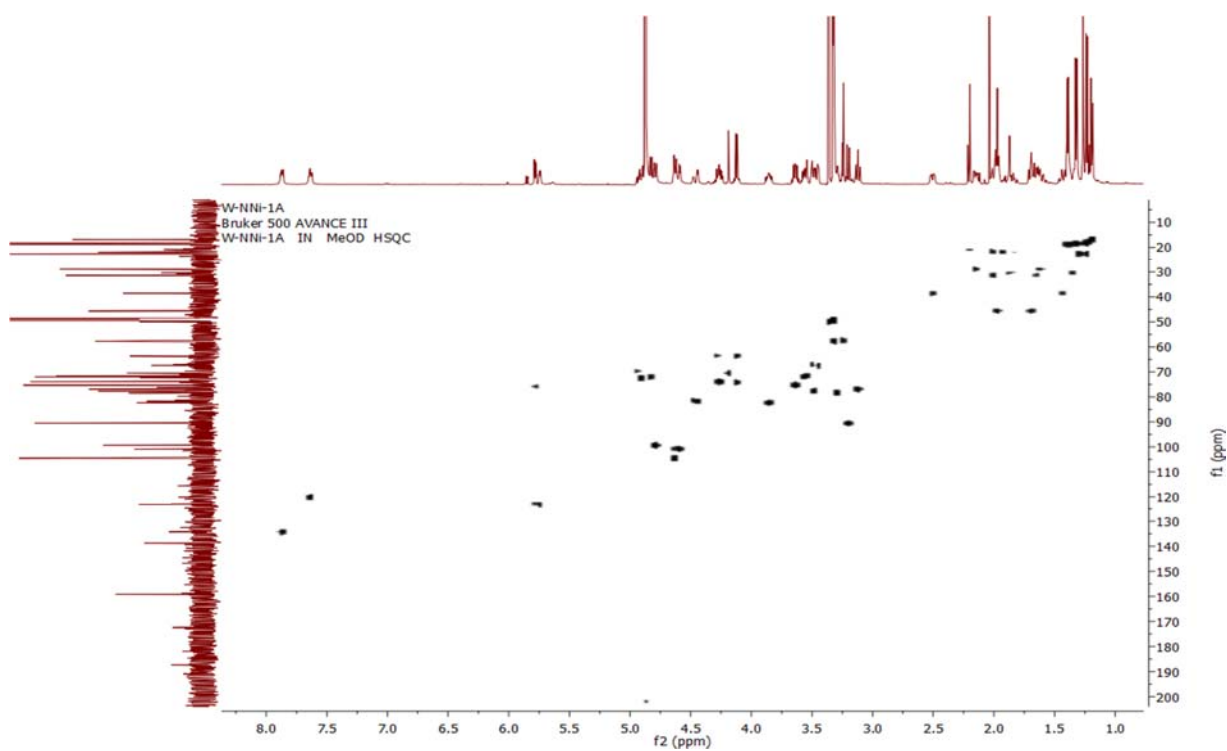


Figure S 13. HSQC spectrum in MeOD for compound 2.

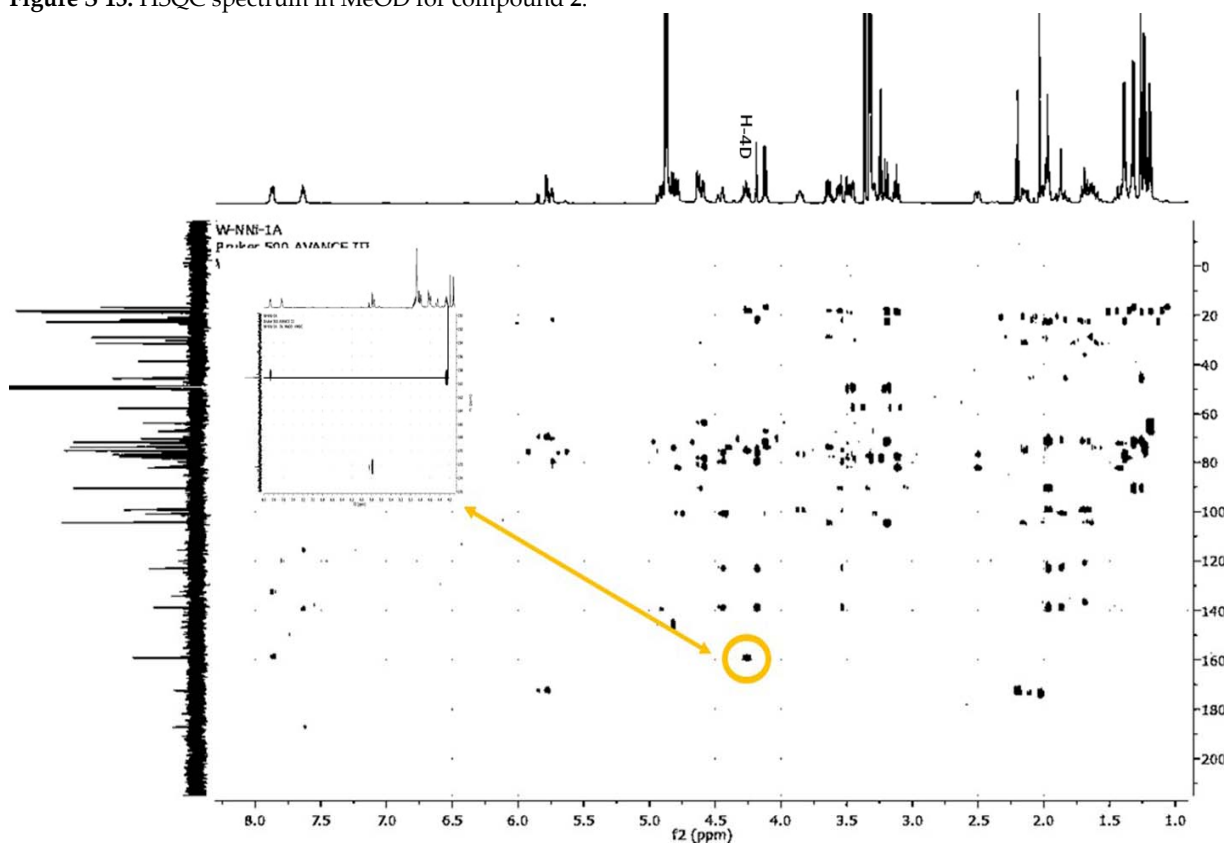


Figure S 14. HMBC spectrum in MeOD for compound 2.

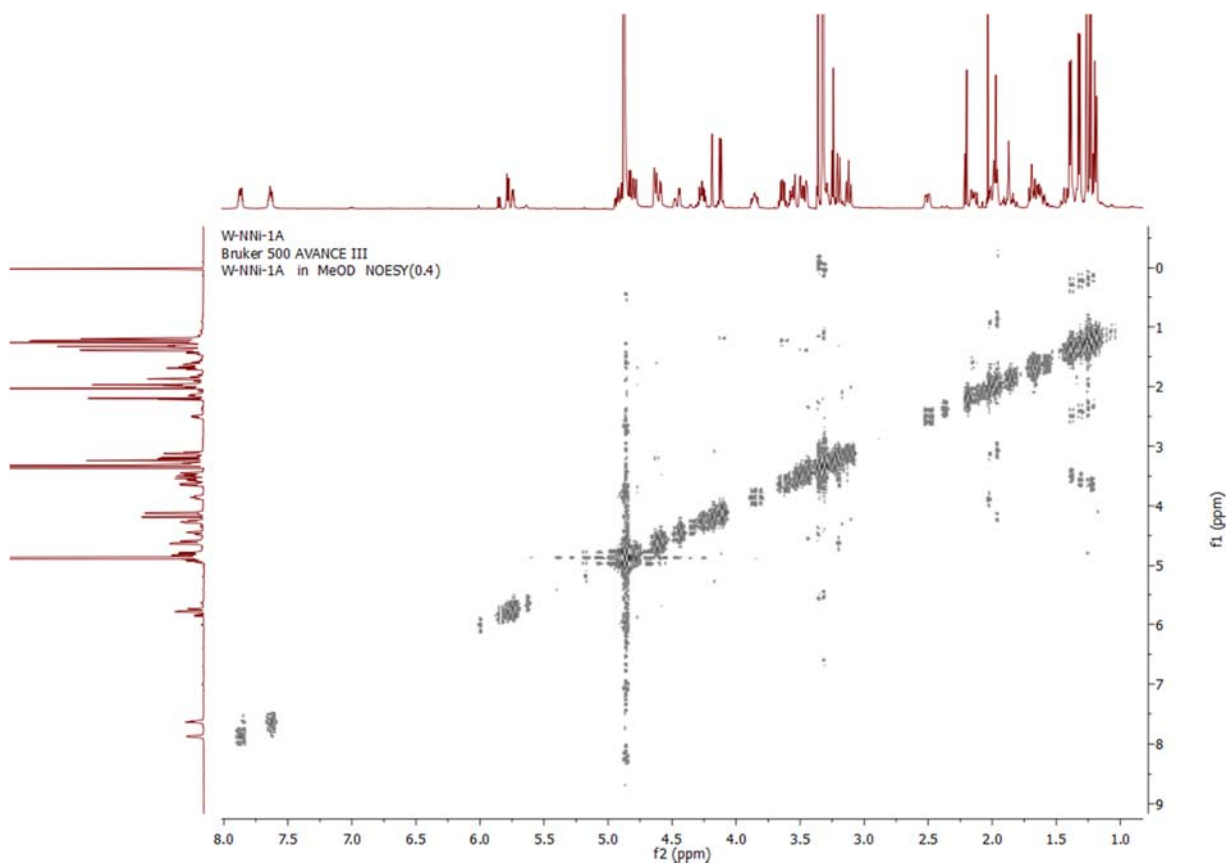


Figure S 15. NOESY spectrum in MeOD for compound 2.

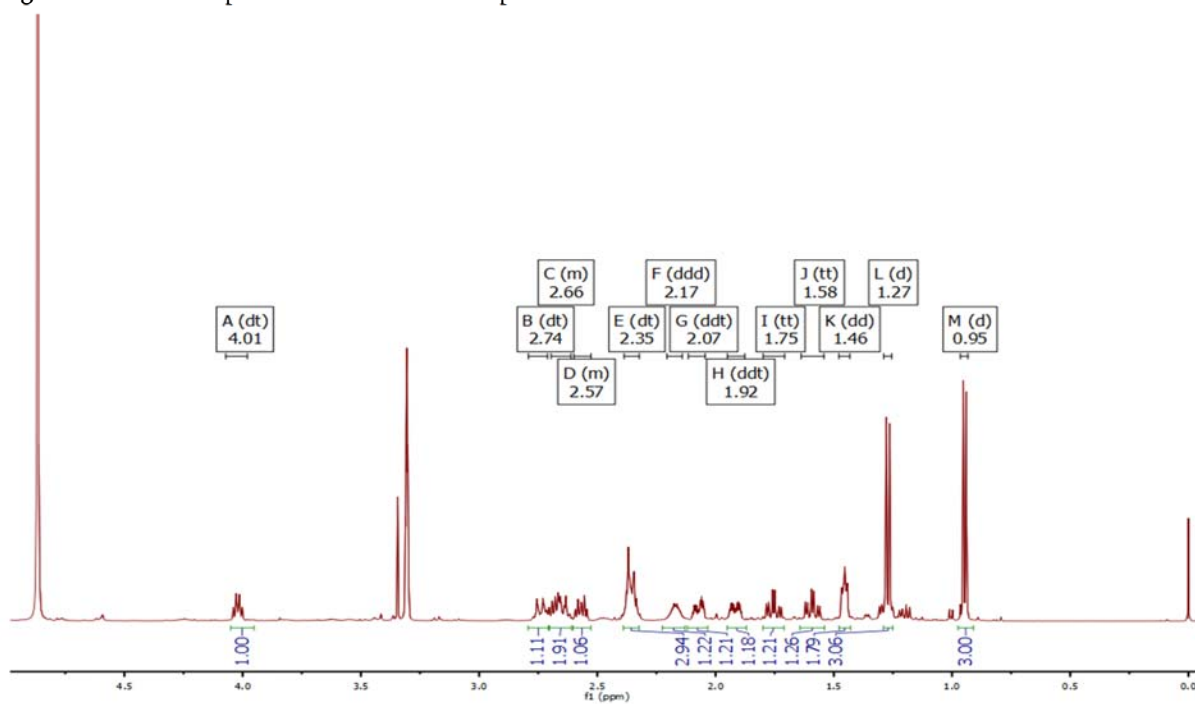


Figure S 16. ^1H NMR spectrum in MeOD for compound 3.

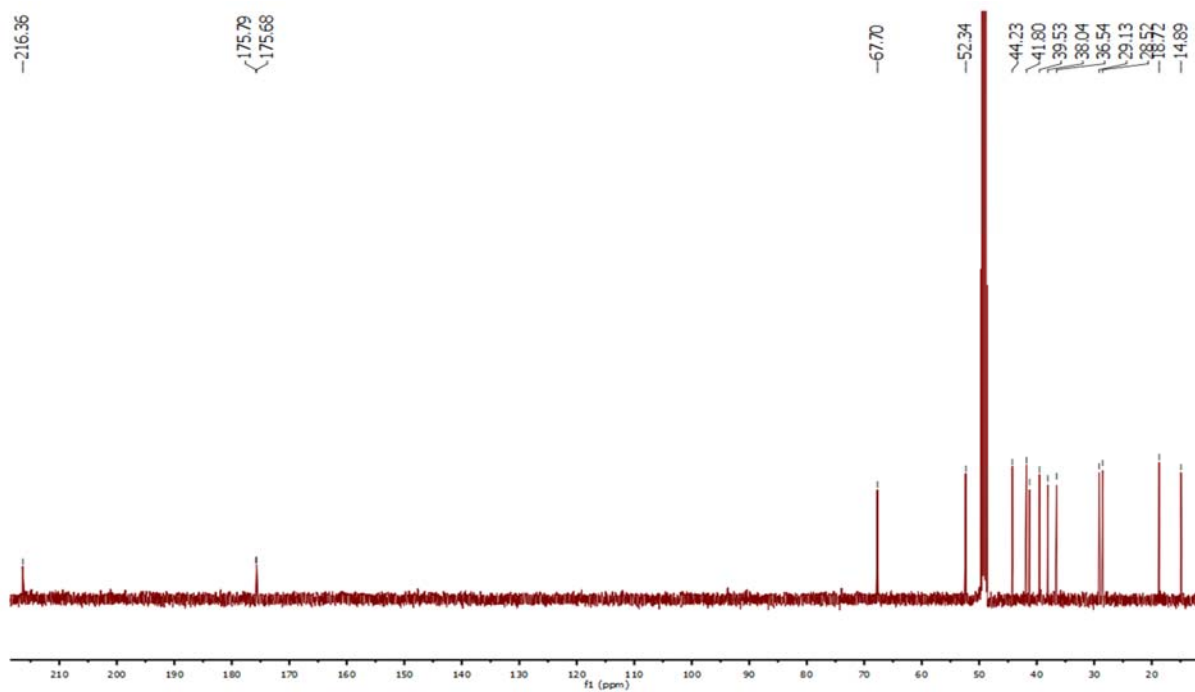


Figure S 17. ^{13}C NMR spectrum in MeOD for compound 3.

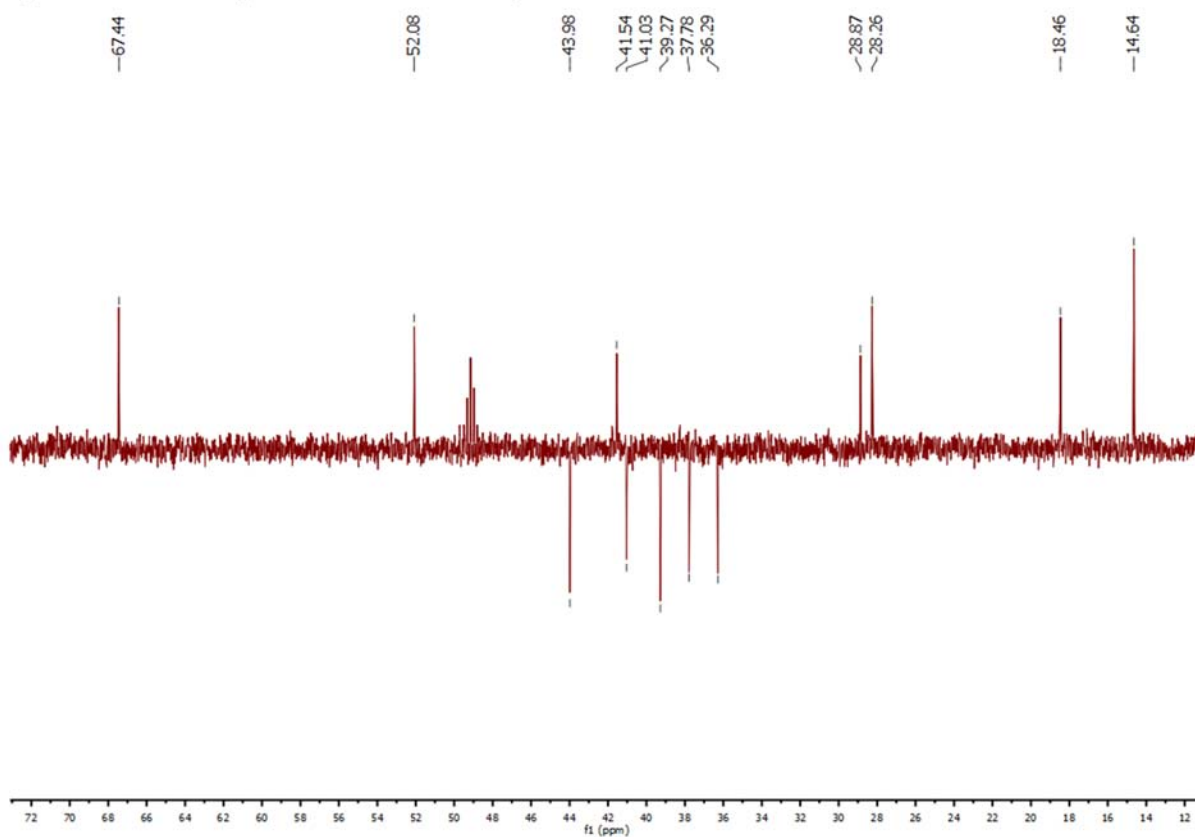


Figure S 18. DEPT-135 spectrum in MeOD for compound 3.

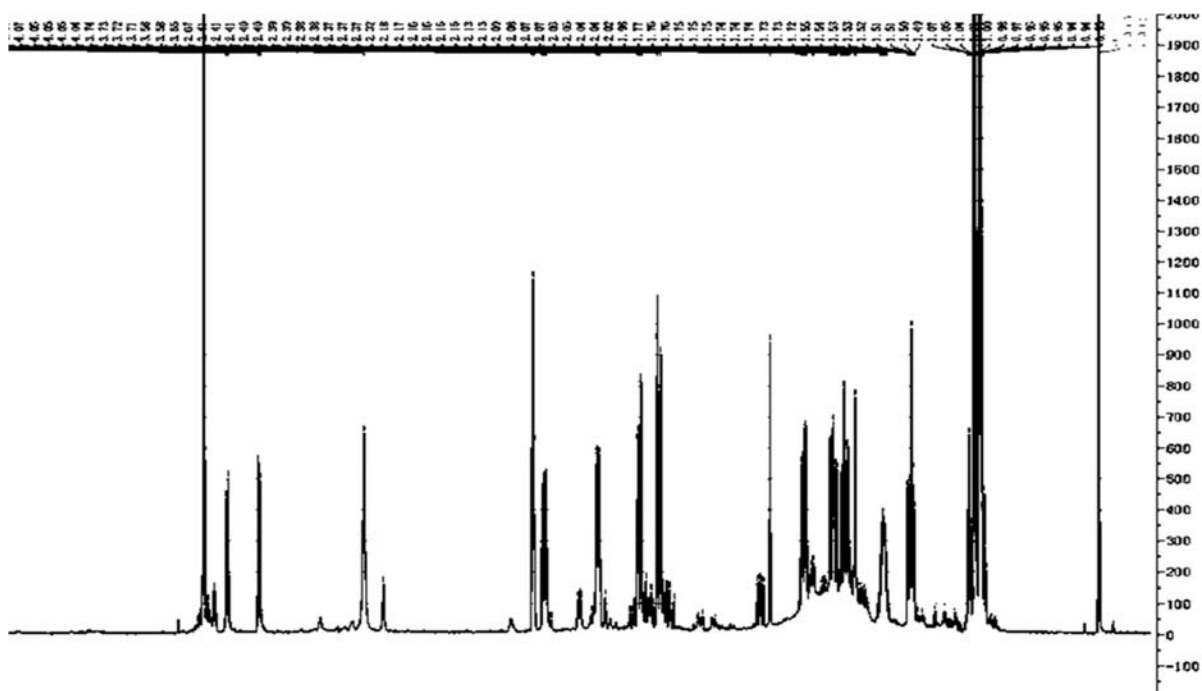


Figure S 19. ^1H NMR spectrum in CDCl_3 for compound 4.

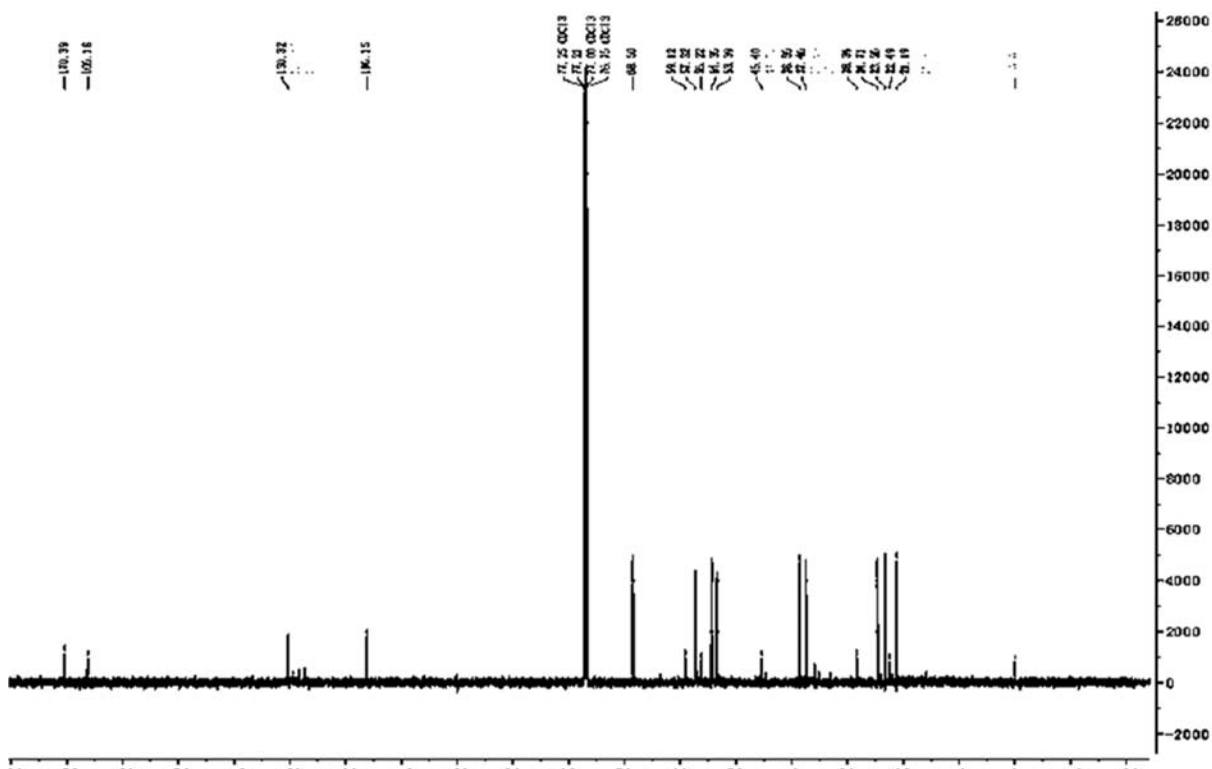


Figure S 20. ^{13}C NMR spectrum in CDCl_3 for compound 4.

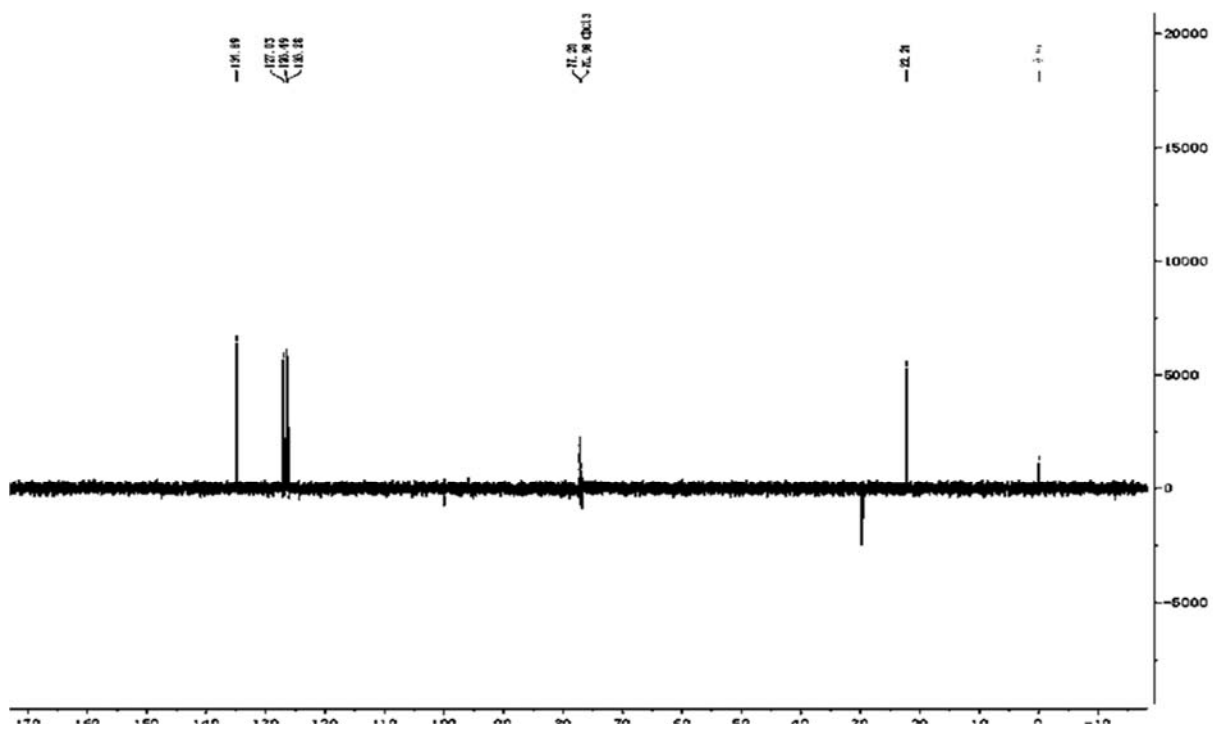


Figure S 23. DEPT spectrum in CDCl₃ for compound 5.

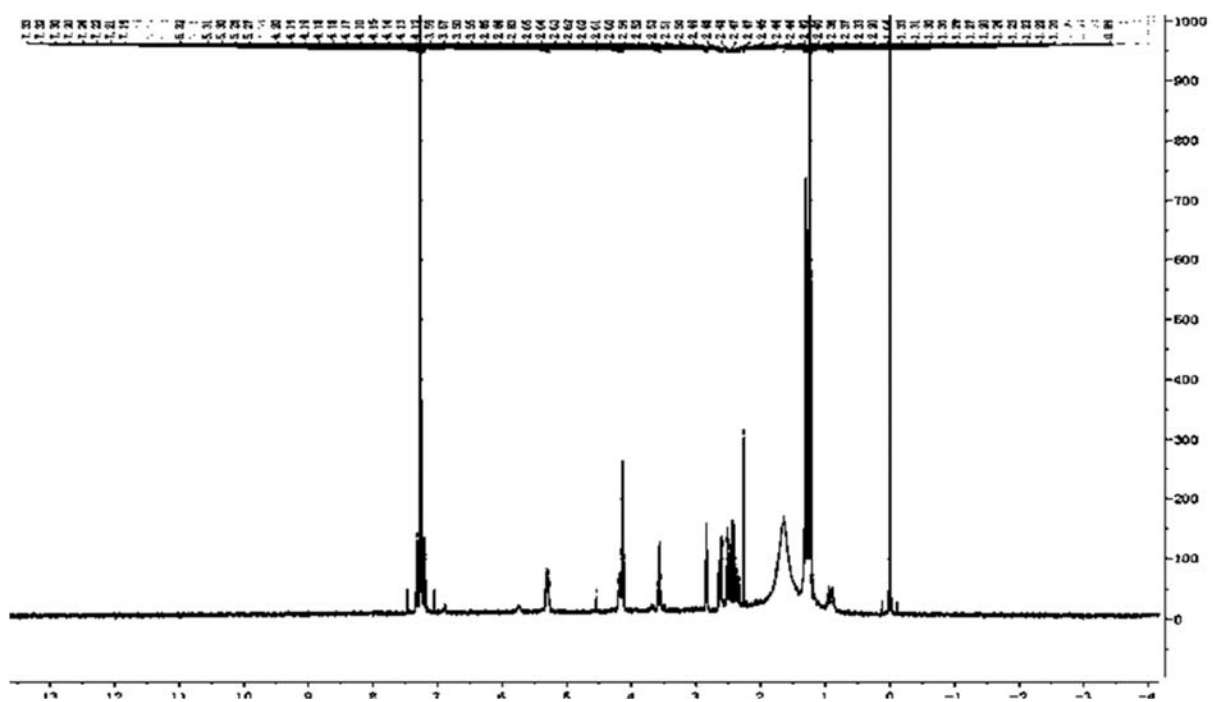


Figure S 24. ¹H NMR spectrum in CDCl₃ for compound 6.

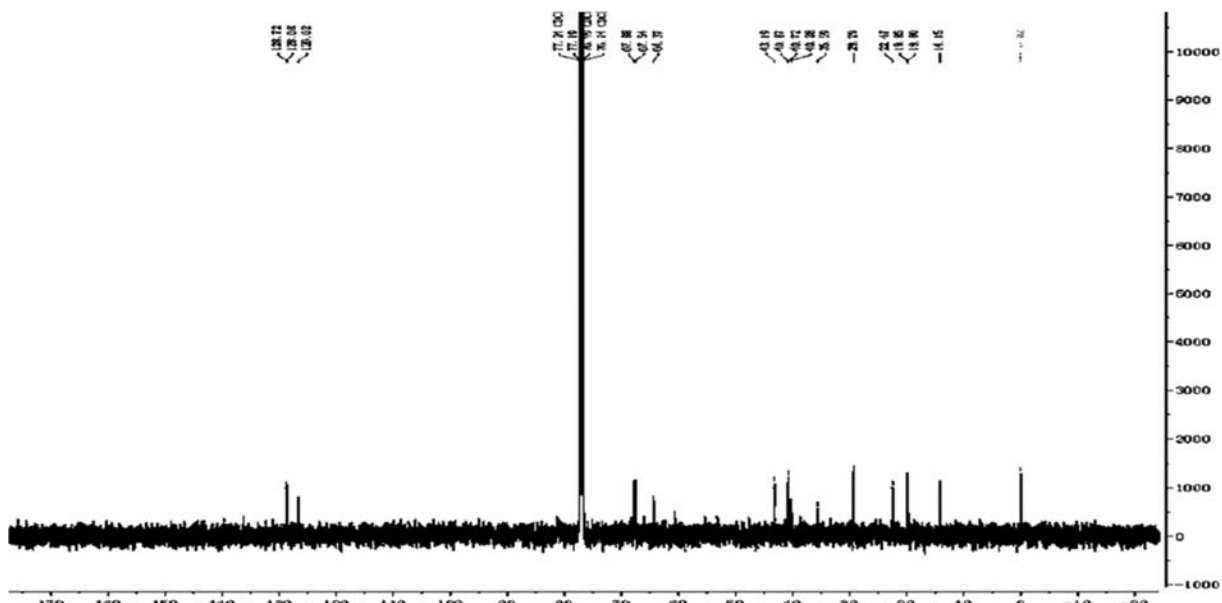


Figure S 25. ¹H NMR spectrum in CDCl₃ for compound 6.

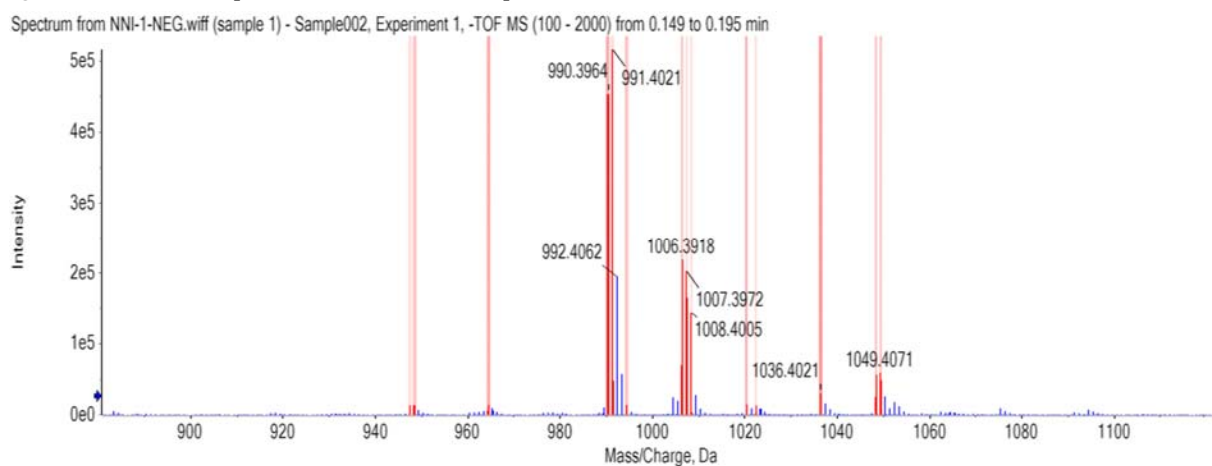


Figure S 26. HR-TOF-MS spectrum in negative mode for compound 1.

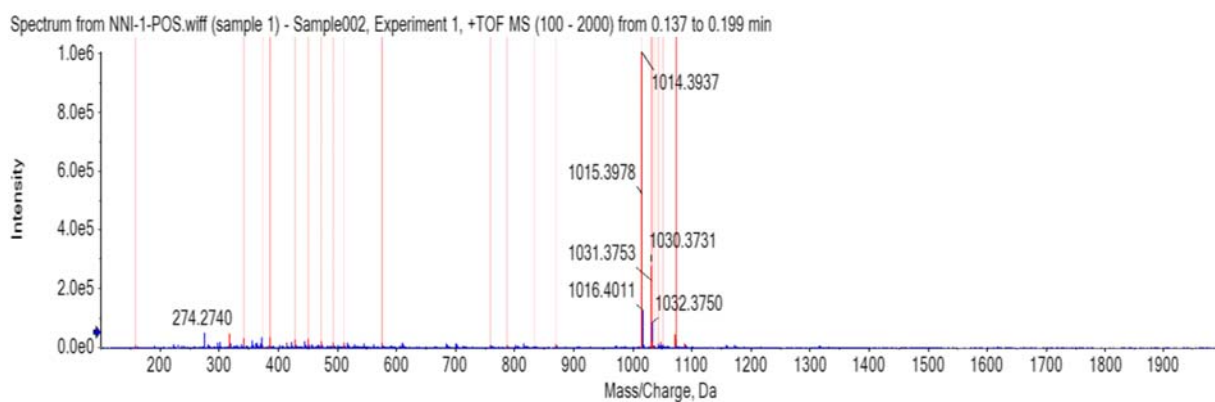


Figure S 27. HR-TOF-MS spectrum in positive mode for compound 1.

NNI-1_171017141246#126-135 RT: 1.80-1.98 AV: 8 NL: 2.02E3
T: ITMS + c ESI Full ms2 1014.50@cid25.00 [275.00-2000.00]

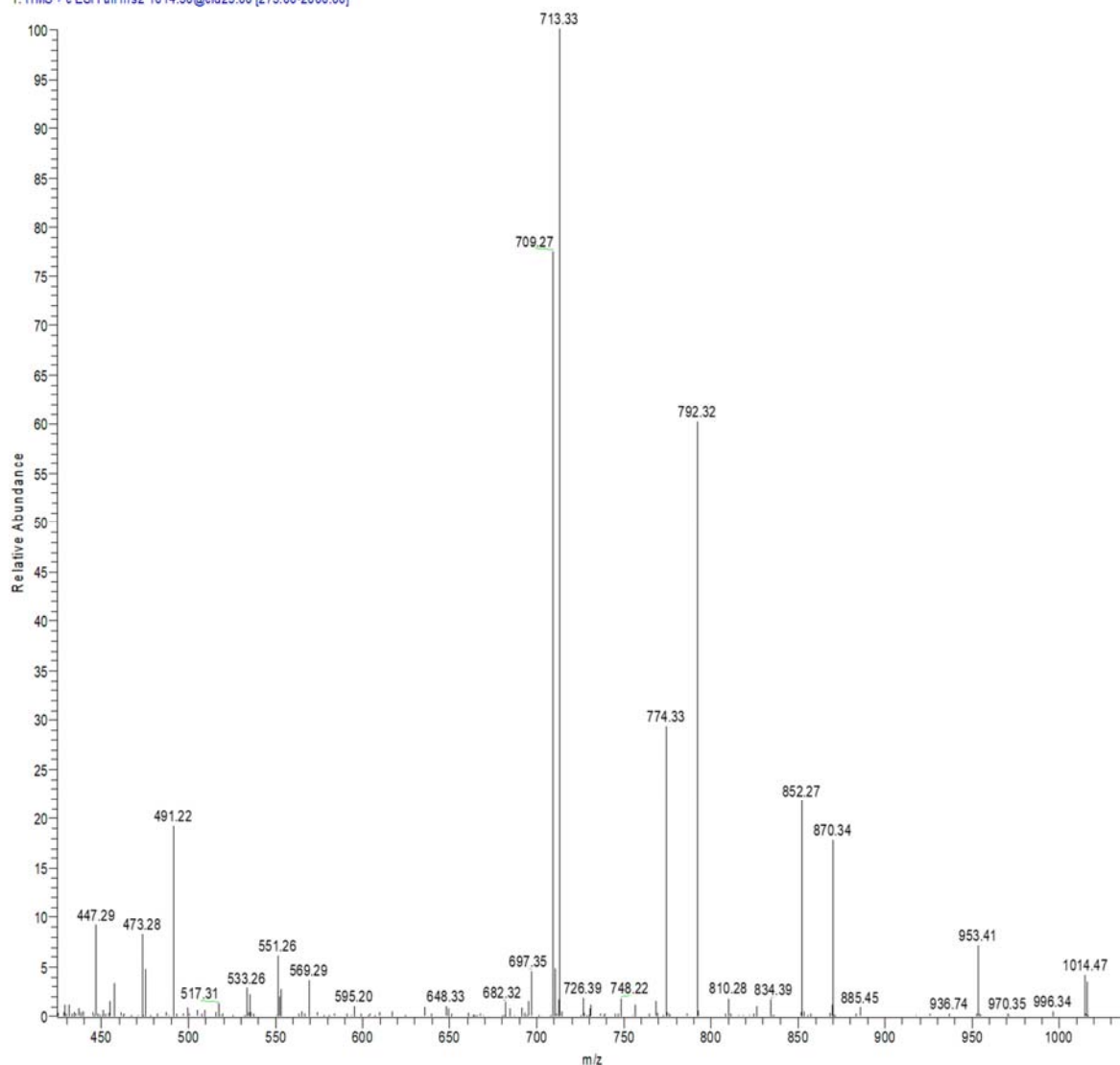


Figure S 28. MSⁿ Spectrum for compound 1.

Spectrum from NNI-1A-NEG.wiff (sample 1) - Sample004, Experiment 1, -TOF MS (100 - 2000) from 0.168 min

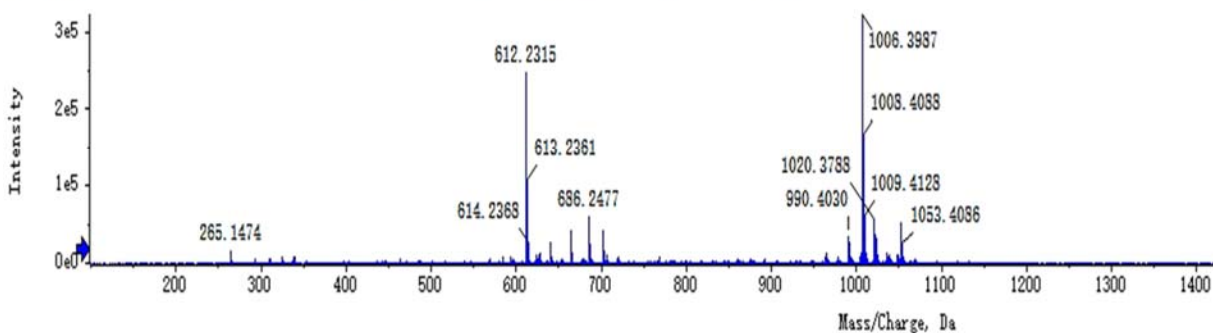


Figure S 29. HR-TOF-MS spectrum in negative mode for compound 2.

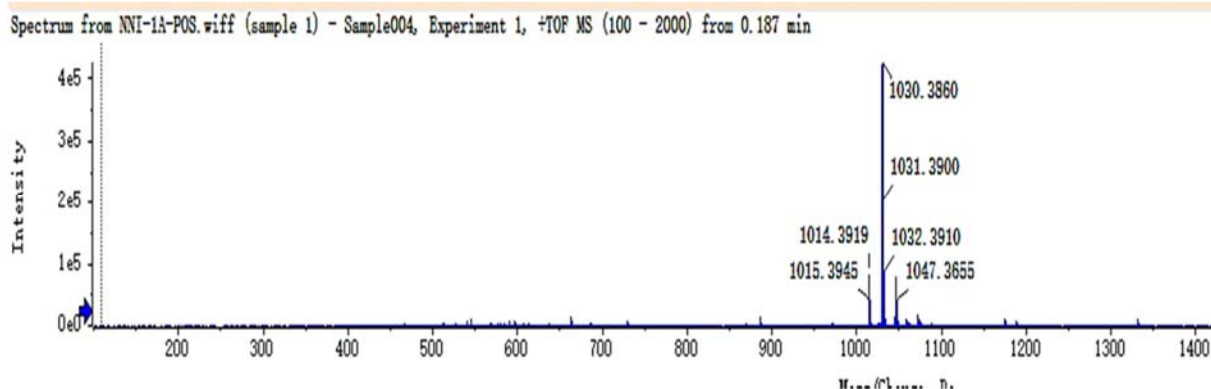


Figure S 30. HR-TOF-MS spectrum in negative mode for compound 2.

Table S 1. Biological activities of compounds (1-5) (MIC values are given in $\mu\text{g/mL}$)

Compounds	MRSA	<i>Pseudomonas aeruginosa</i>	<i>Klebsiella pneumoniae</i>	<i>Escherichia coli</i>	<i>Bacillus subtilis</i>
	MIC	MIC	MIC	MIC	MIC
1	16	16	16	16	8-16
2	16	16	16	16	8-16
3	16-32	16-32	16-32	-	-
4	32	16	16	16	16
5	32	16	16	16	16
Tetracycline	16	32	32	16	8

Strain	16S rDNA gene sequence
<i>Streptomyces pratensis</i> NA-ZhouS1	AGGGCGGGCGTGCTTACCATGTCAGTCGAACGATGAAGCCTTTCCG GGTGGATTAGTGGCGAACGGGTGAGTAACACGTGGGCAATCTGCC CTTCACTCTGGGACAAGCCCTGGAAACGGGGTCTAATACCGGATA ACACTCTGTCCCGCATGGGACGGGGTTAAAGCTCCGGCGGTGAA GGATGAGCCCGCGGCTATCAGCTTGTGGTGGGGTAATGGCCTAC CAAGGCGACGACGGGTAGCCGGCCTGAGAGGGCGACCCGCCACA CTGGGACTGAGACACGGCCAGACTCCTACGGGAGGCAGCATGG GGAATATGCAACAATGGGCGAAAGCCTGATGCAGCGACGCCGCT GAGGGATGACGGCCTTCGGGTGTAAACCTCTTTCAGCAGGGAAG AAGCGAAAGTGACGGTACCTGCAGAGAAAGCGCCGGCTAACTAC GTGCCAGCAGCCGGTAATACGTAGGGCGCAAGCGTTGTCCGGA ATTATTGGGCGTAAAGAGCTCGTAGGGCGCTTGTACGTCGGATGT GAAAGCCC GGCGCTTAACCCCGGGTCTGCATTCGATACGGGCTAG CTAGAGTGTGGTAGGGGAGATCGAATTCTGGTGTACGGTGAA ATGCGCAGATATCAGGAGGAACACCGGTGGCGAAGGCGGATCTCT GGGCCATTACTGACGCTGACGAGCGAAAGCGTGGGAGCGAACA GGATTAGATACCCTGGTAGTCCACGCCGTAACGTTGGGAAGTAC GTGTGGCGACATTCCACGTCGTCGGTGCCCGACGTAACGCATTAA GTTCCCCGCTGGGGAGTACGGCCGCAAGGCTAAAACTCAAAGGA ATTGACGGGGGCCCGCACAAGCAGCGGAGCATGTGGCTTAATTTCG ACGCAACCGGAAGAACCCTTACCAAGGCTTGACATATACCGGAAAG CATCAGAGATGGTGCCTTGTGGTTCGGTATACAGGTGGTGCAT GGCTGTCGTCAGCTCGTGTGAGATGTTGGGTTAAGTCCCGCAA CGAGCGCAACCCTTGTCTGTGTTGCCAGCATGCCCTTCGGGGTGA TGGGGACTCACAGGAGACTGCCGGGTCAACTCGGAGGAAGGTGG GGACGACGTCAAGTCAATCATGCCCTTATGTCTTGGGCTGCACAG TGCTACAATGGCCGGTACAATGAGCTGCGATGCCCGGAGCGGAG CGAATCTCAAAAAGCCGGTCTCAGTTCGGATTGGGGTCTGCAACTC GACCCCATGAAGTCGGAGTTGTAGTAATCGCAGATCAGCAATTGCT GCGGTGAATACGTTCCCGGGCTTGTACACACCCCGCTCACGTCA CGAAAGTCGGTAACACCCGAAAGCCGGTGGCCCAACCCTTGTGGG AGGAGCTTCGAAAGGTGTACGGCAAGTTCCCTT

Figure S 31. 16S ribosomal DNA gene, full sequence of *Streptomyces* sp. NA-ZhouS1

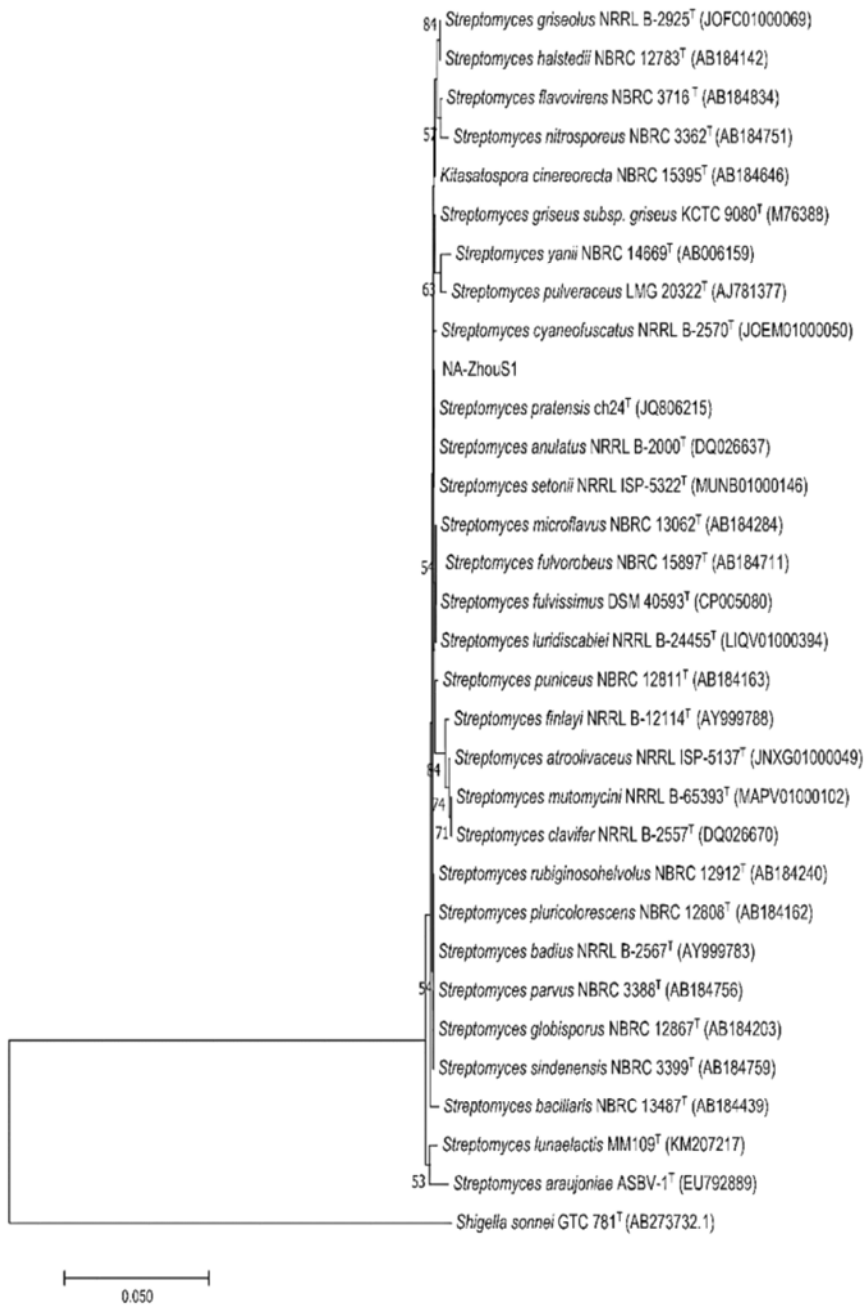


Figure S 32. Neighbor-joining phylogenetic tree of strain NA-ZhouS1 based on 16S rDNA sequences generated by Mega version 7. Numbers at branch points indicate levels of bootstrap (expressed as percentages of 1000 replicates), only values exceeding 50% are given. *Shigella sonnei* GTC 781^T (AB273732.1) was used as an outgroup. Bar, 0.050 substitutions per nucleotide position.