Supporting Information

Metabolites from the Fungal Endophyte *Aspergillus* austroafricanus in Axenic Culture and in Fungal-Bacterial Mixed Cultures

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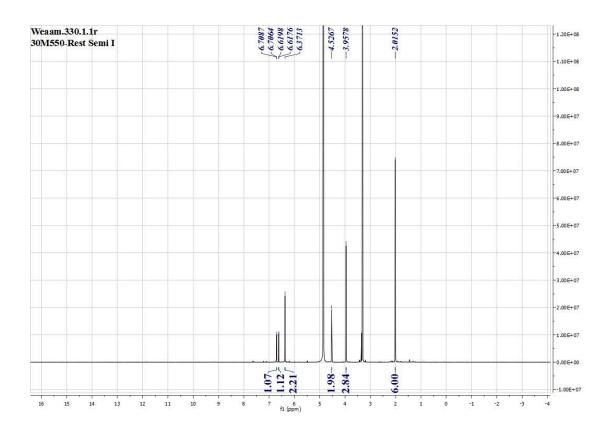
E-mail addresses: <u>weaam.ebrahim@uni-duesseldorf.de</u> (W.E.), <u>georgios.daletos@uni-duesseldorf.de</u> (G.D.), <u>proksch@uni-duesseldorf.de</u> (P.P.).

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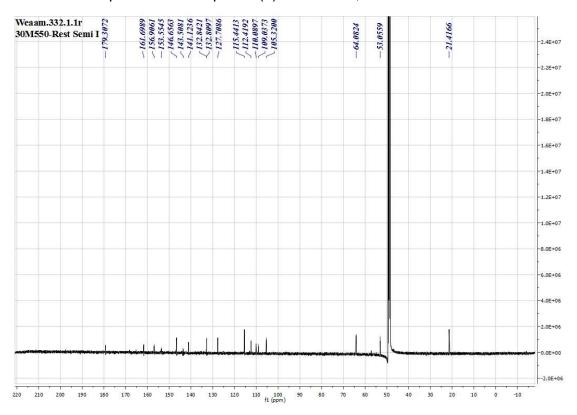
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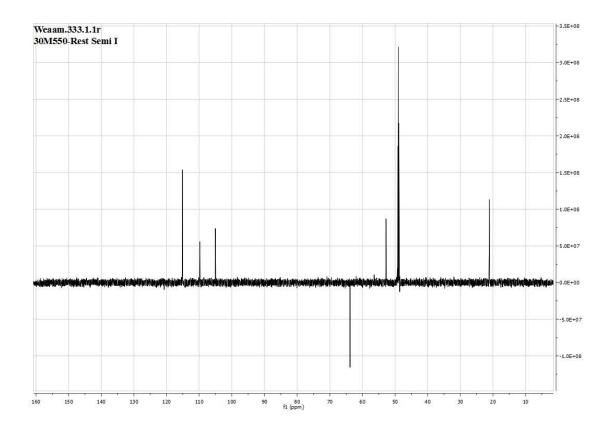
S1. ¹H NMR spectrum of compound (**1**) in MeOH-*d*₄

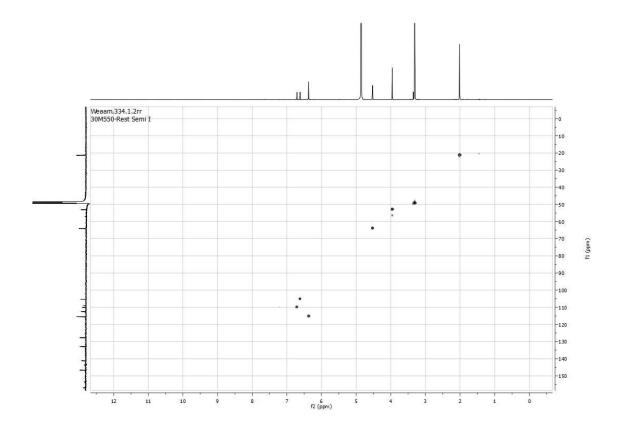


S2. 13 C NMR spectrum of compound (1) in MeOH- d_4

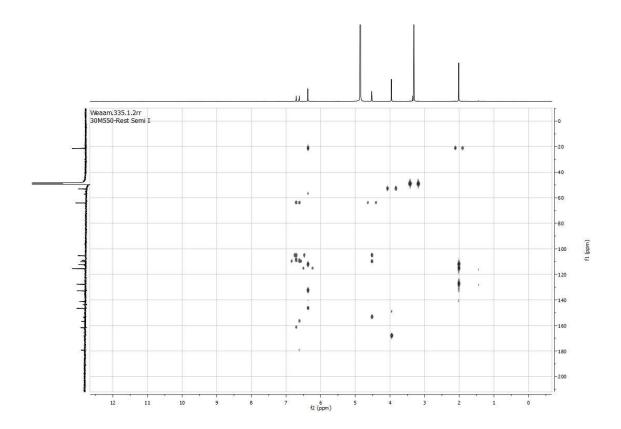


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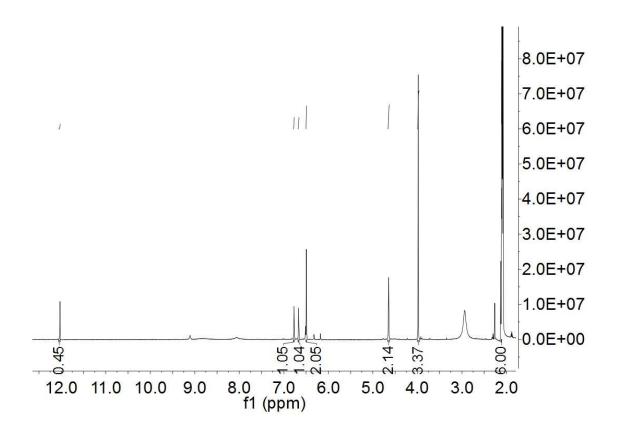




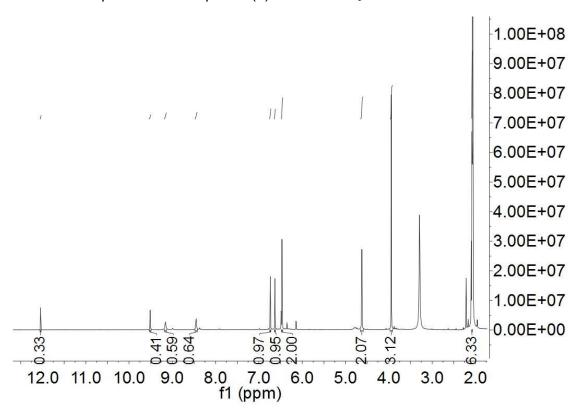
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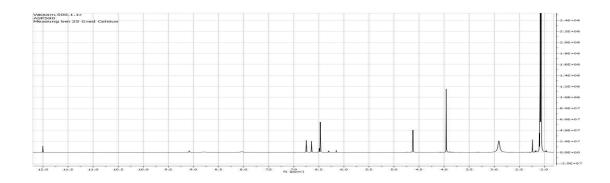
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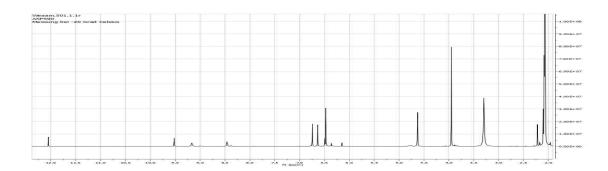


S7. ¹H NMR spectrum of compound (1) in acetone-*d*₆ at -20°C

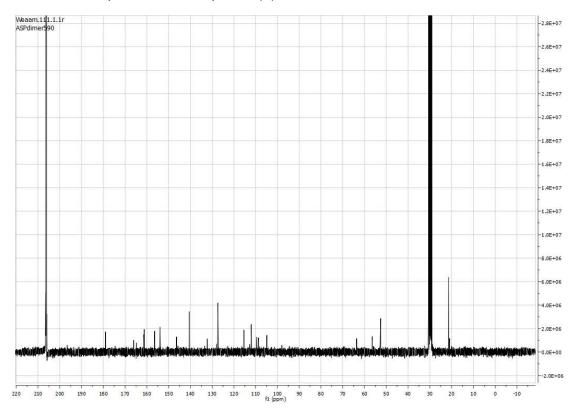


S8. Comparison of 1H NMR spectrum of compound (1) in acetone- d_6 at room temperature and at -20 $^\circ$ C

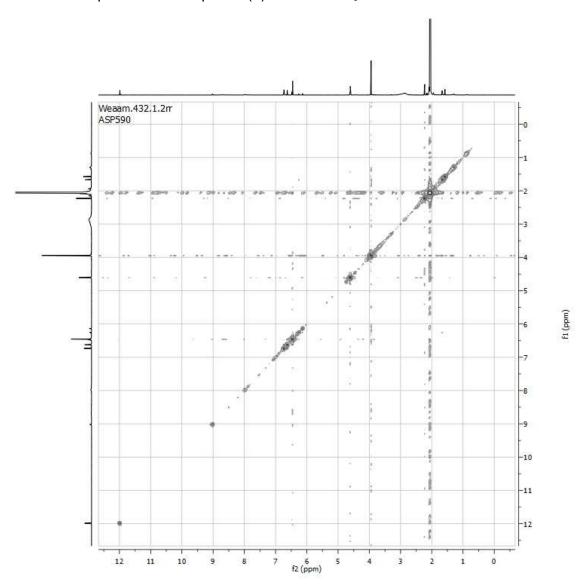




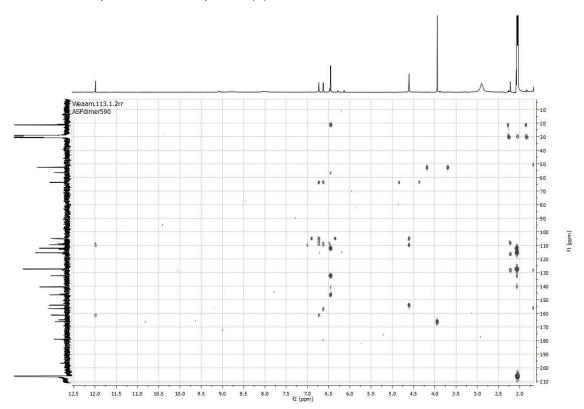
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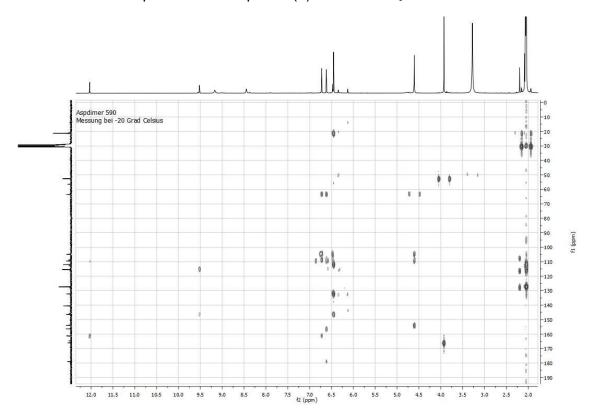
S10. COSY spectrum of compound (1) in acetone- d_6



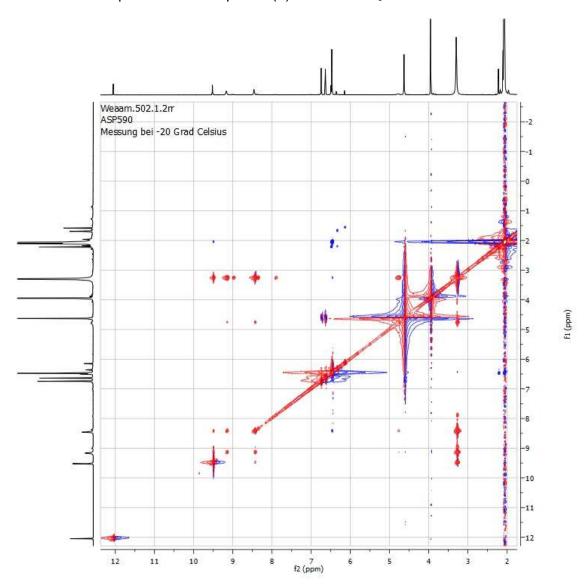
S11. HMBC spectrum of compound (1) in acetone- d_6



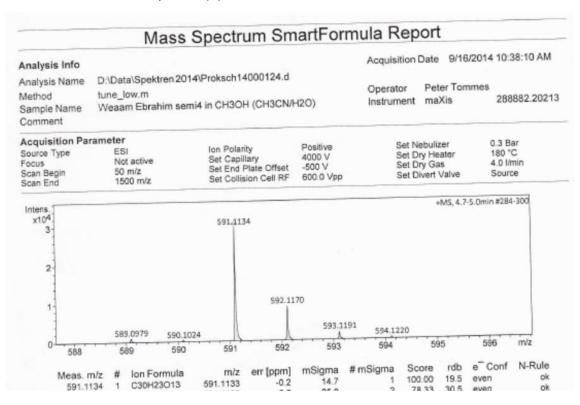
S12. HMBC NMR spectrum of compound (1) in acetone-d₆ at -20°C



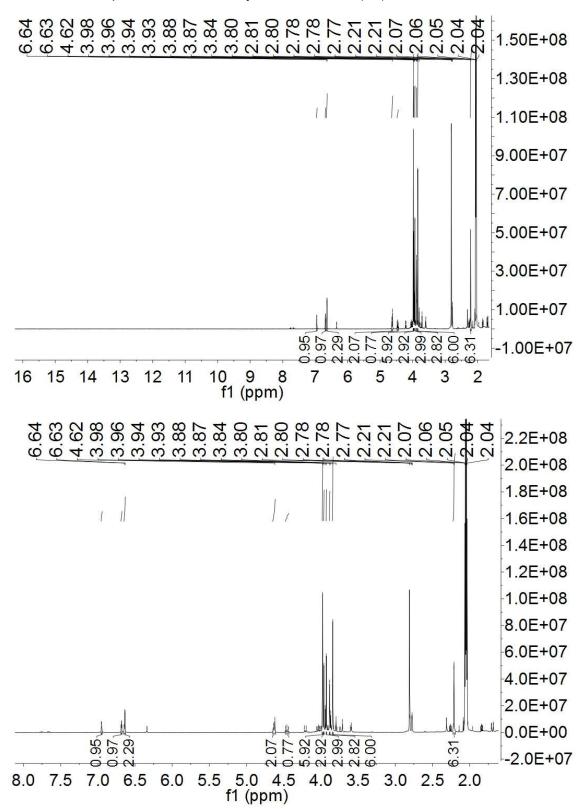
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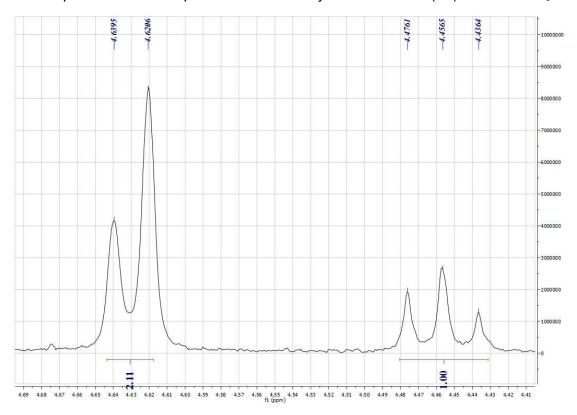
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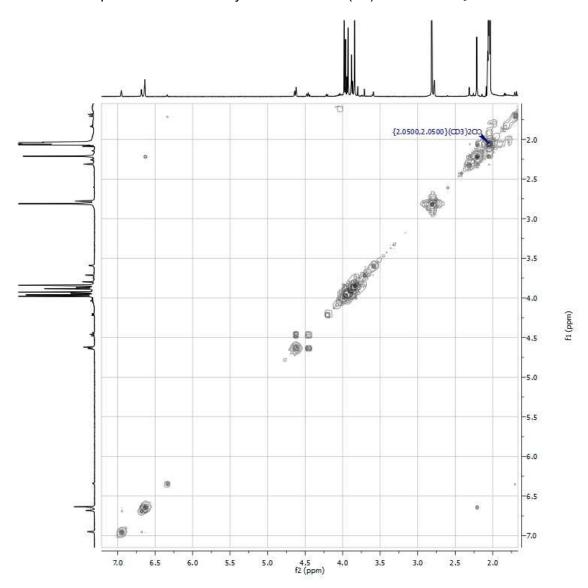
S15. ¹H NMR spectrum of the methylated derivative (**1b**) in acetone-d₆



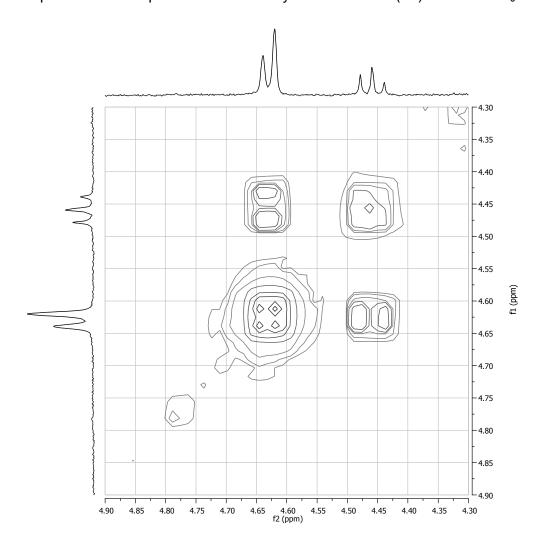
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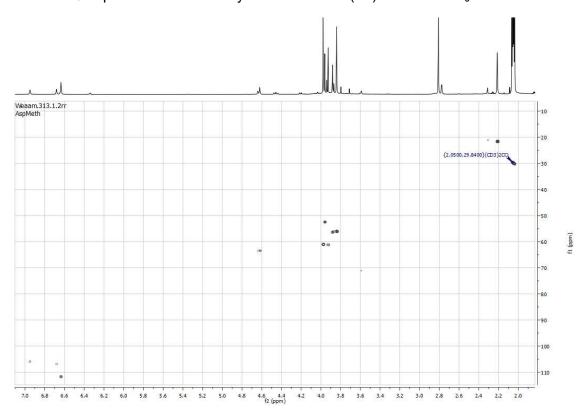
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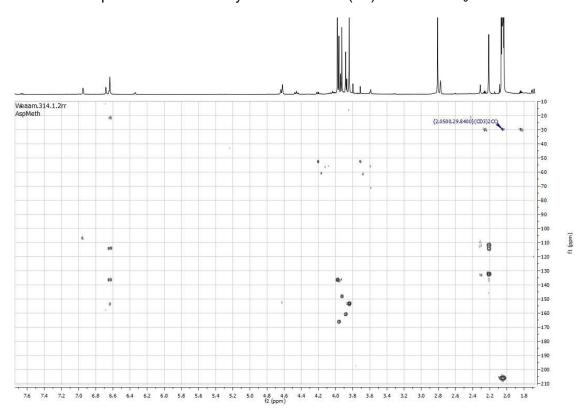
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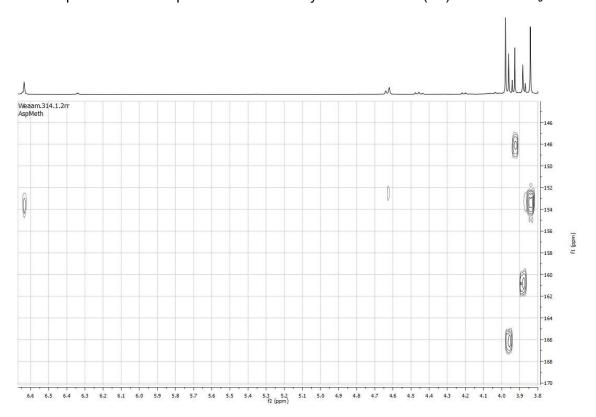
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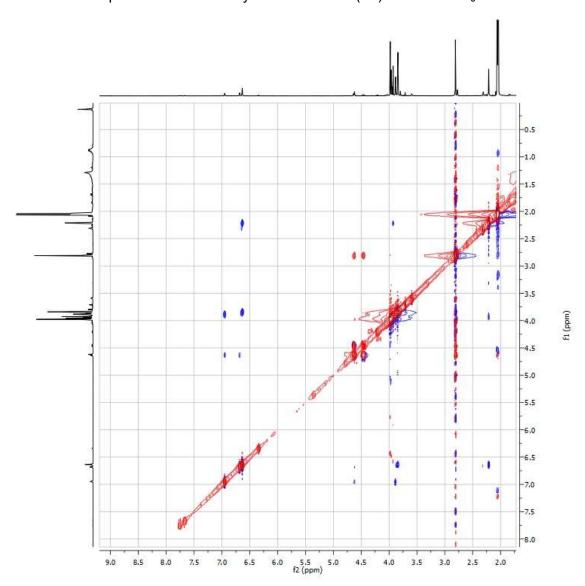
S20. HMBC spectrum of the methylated derivative (1b) in acetone- d_6



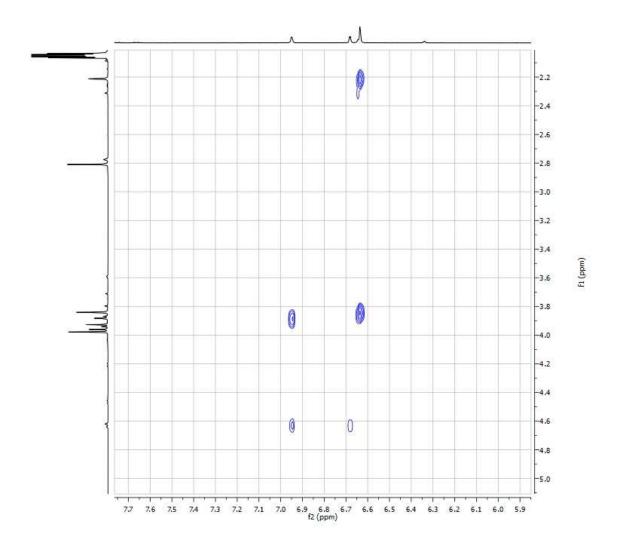
S21. Expanded HMBC spectrum of the methylated derivative (1b) in acetone- d_6



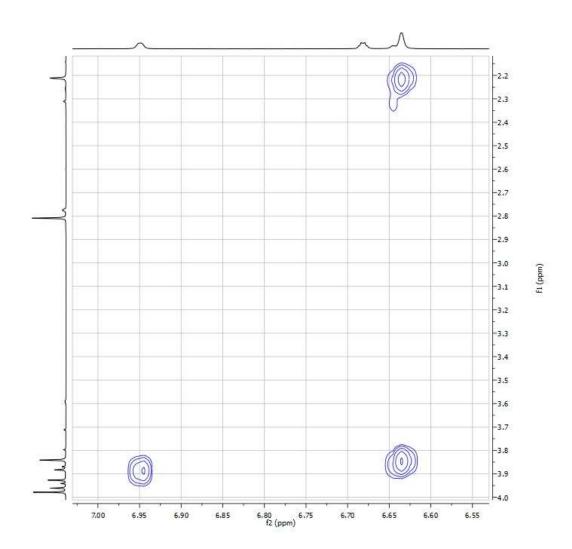
S22. ROESY spectrum of the methylated derivative (1b) in acetone- $d_{\rm 6}$



S23. Expanded ROESY spectrum of the methylated derivative (${f 1b}$) in acetone- ${\it d}_{\it 6}$



S24. Expanded ROESY spectrum of the methylated derivative (1b) in acetone- d_6

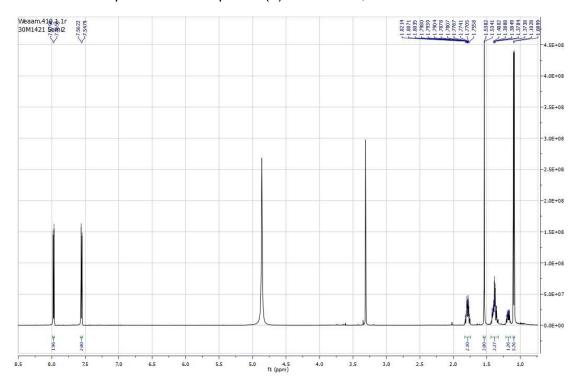


S25. HRESIMS of the methylated derivative (1b)

Mass Spectrum SmartFormula Report Analysis Info Acquisition Date 3/11/2015 9:40:32 AM Analysis Name D:\Data\Spektren 2015\Proksch15HR000088.d Method tune_low.m Peter Tommes Operator Sample Name L. Lewald Meth (CH3OH) Instrument maXis 288882.20213 13 ul auf 1000 ul Comment **Acquisition Parameter** ESI Not active 50 m/z Positive 4000 V -500 V Source Type Ion Polarity Set Nebulizer 0.3 Bar Focus Scan Begin Set Capillary Set End Plate Offset Set Dry Heater Set Dry Gas 180 °C 4.0 I/min Scan End Set Collision Cell RF 600.0 Vpp Set Divert Valve Source Intens. +MS, 3.6-3.7min #216-222 x104 675,2072 2-676.2105 677.2135 678.2159 678 675 676 677 679 m/z 675.2072 C36H35O13 675.2072 -0.0 43.8 100.00 19.5 even

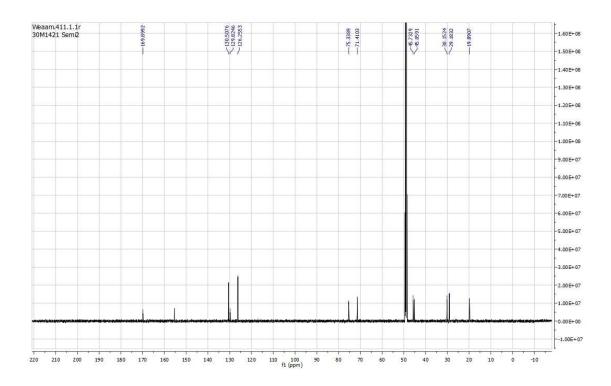
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S26. 1 H NMR spectrum of compound (2) in MeOH- d_{4}

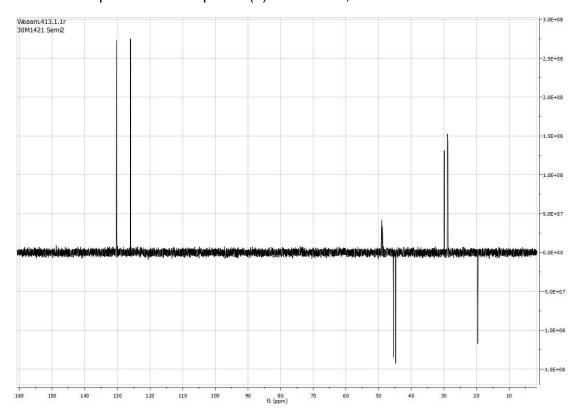


S27. 13 C NMR spectrum of compound (2) in MeOH- d_4

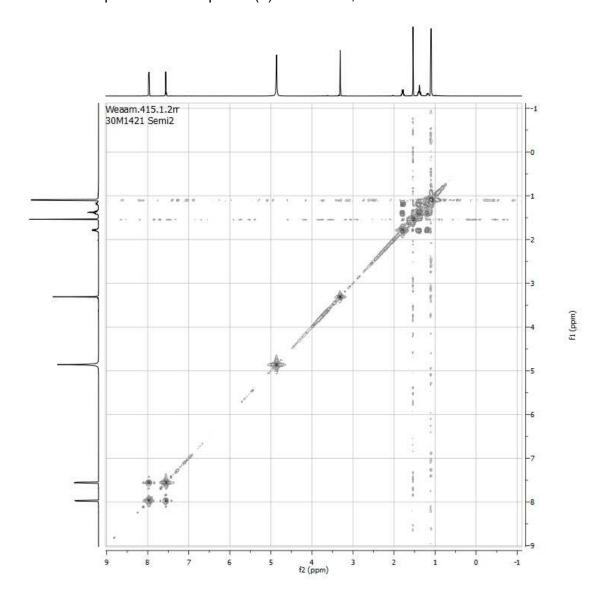
 $^{13}\text{C NMR (151 MHz, MeOD)} \ \delta \ 169.90, 130.51, 129.82, 126.26, 75.34, 71.41, 45.73, 45.06, 30.15, 29.18, 19.89.$



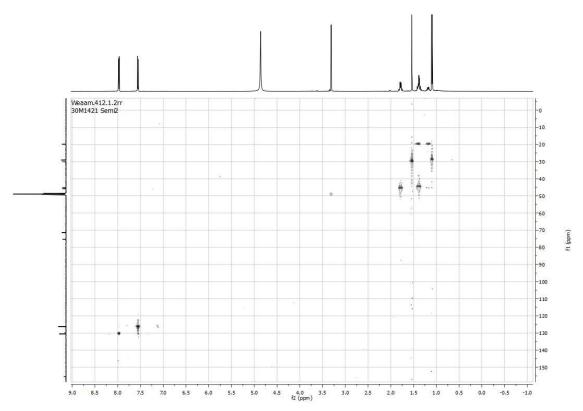
S28. DEPT spectrum of compound (2) in MeOH- d_4



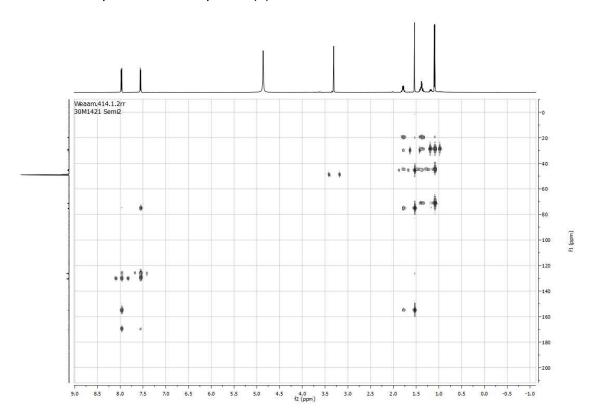
S29. COSY spectrum of compound (2) in MeOH- d_4



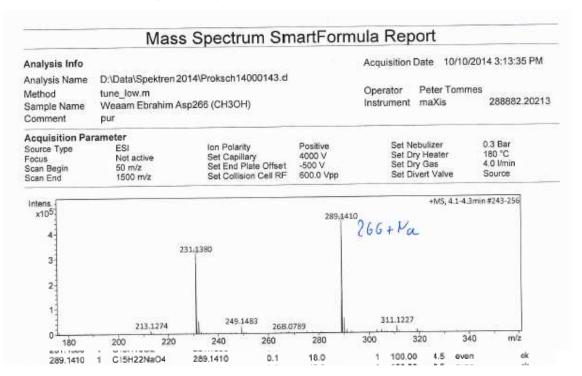
S30. HMQC spectrum of compound (2) in MeOH- d_4



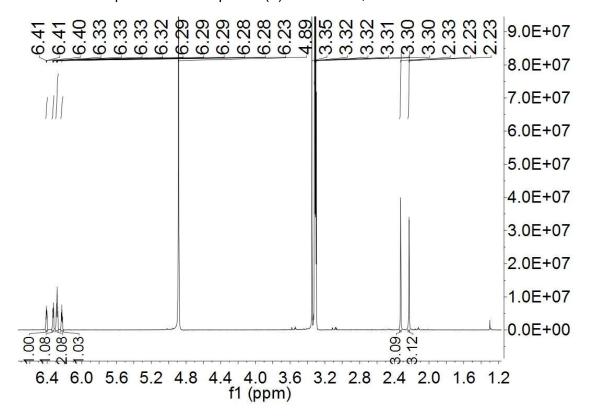
S31. HMBC spectrum of compound (2) in MeOH- d_4



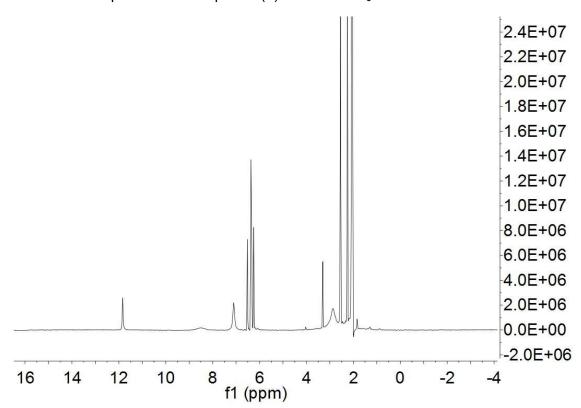
S32. HRESIMS of compound (2)



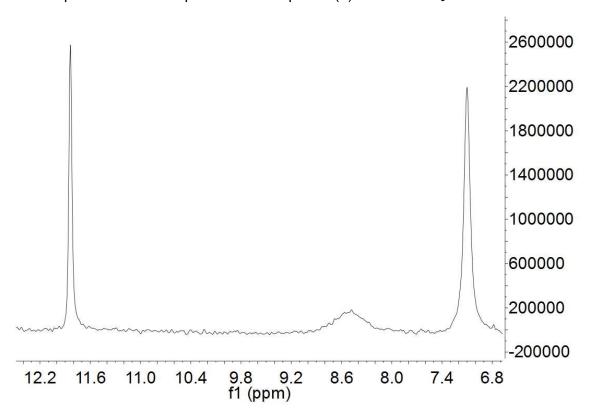
S33. ¹H NMR spectrum of compound (**8**) in MeOH-d₄



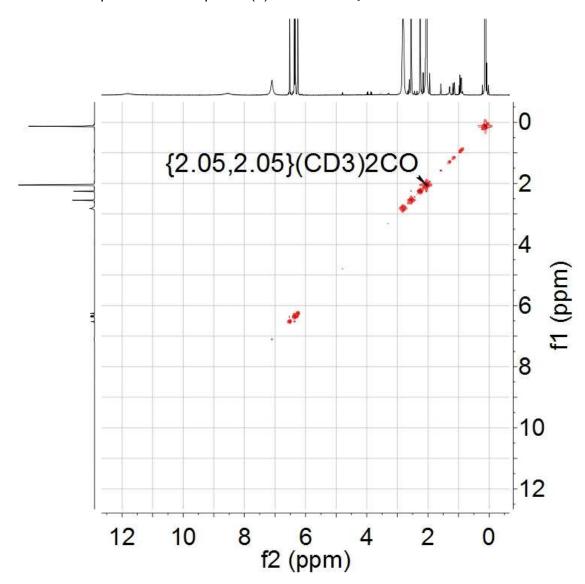
S34. ¹H NMR spectrum of compound (**8**) in acetone-*d*₆



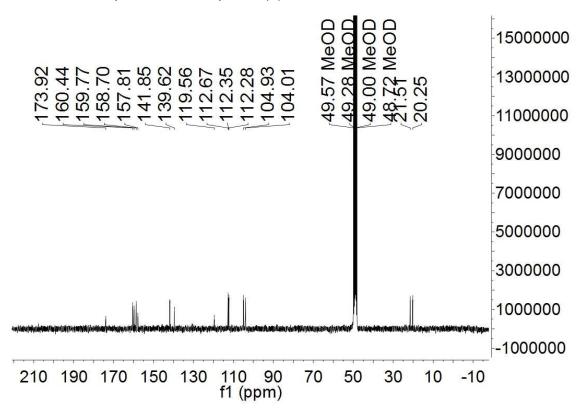
S35. Expanded ¹H NMR spectrum of compound (**8**) in acetone-*d*₆



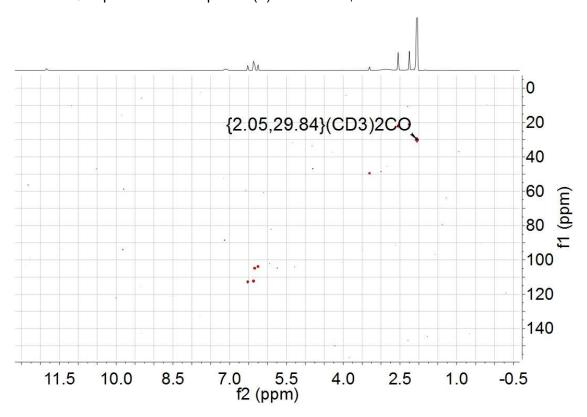
S36. COSY spectrum of compound (8) in acetone-d₆



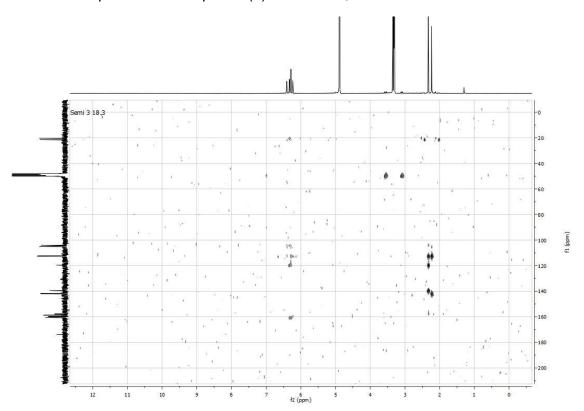
S37. ¹³C NMR spectrum of compound (8) in MeOH-d₄



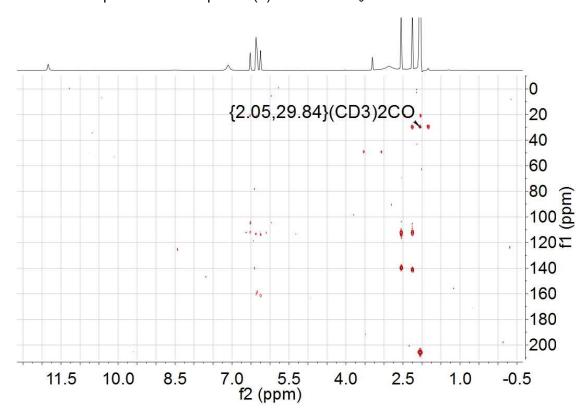
S38. HMQC spectrum of compound (8) in MeOH-d4



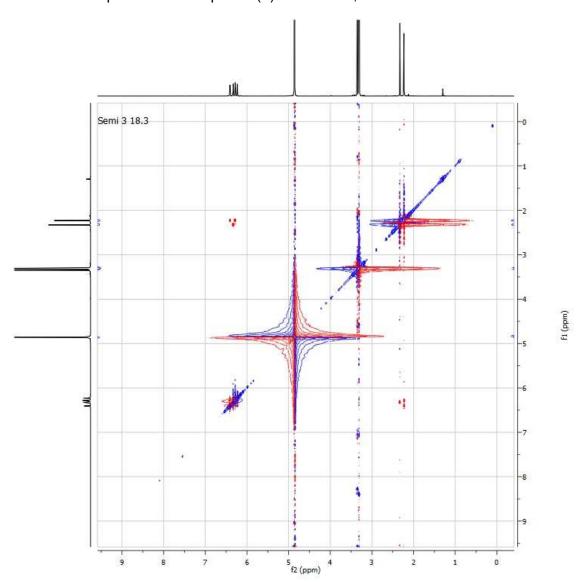
S39. HMBC spectrum of compound (8) in MeOH- d_4



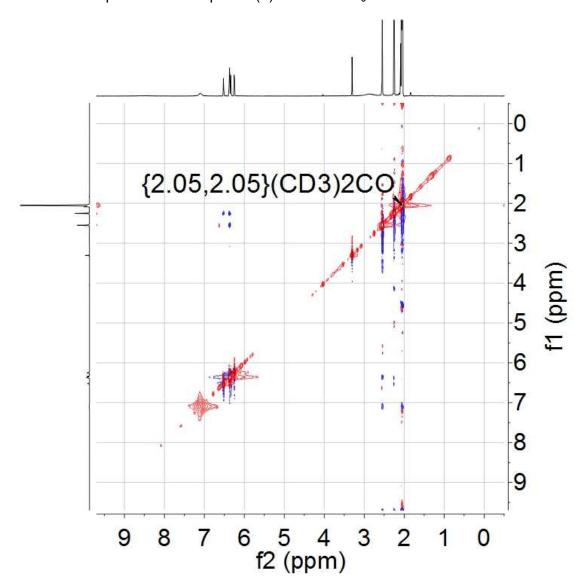
S40. HMBC spectrum of compound (8) in acetone- d_6



S41. ROESY spectrum of compound (8) in MeOH- d_4



S42. ROESY spectrum of compound (8) in acetone- d_6



S43. HRESIMS of compound (8)

