

Supplementary information

¹³C NMR spectroscopic data of compounds:

Artobiloxanthone (2) ¹³C NMR (125 MHz, acetone-*d*₆) δ 21.99 (C-12), 22.27 (C-9), 28.20 (C-18), 28.54 (C-17), 38.14 (C-10), 78.71 (C-16), 99.78 (C-6), 102.03 (C-8), 103.7 (C-3'), 105.51 (C-4a), 107.03 (C-1'), 111.43 (C-3), 111.91 (C-13), 116.14 (C-14), 127.88 (C-6'), 129.88 (C-15), 136.68 (C-5'), 145.35 (C-4'), 150.92 (C-11), 151.57 (C-2'), 152.31 (C-8a), 159.48 (C-7), 161.80 (C-5), 162.57 (C-2), 181.27 (C=O, C-4).

Artoindonesianin P (3) ¹³C NMR (125 MHz, acetone-*d*₆) δ 21.18 (C-9), 23.07 (C-12), 28.54 (C-13), 50.05 (C-10), 94.44 (C-11), 95.23 (C-8), 100.11 (C-6), 105.22 (C-1'), 105.27 (C-4a), 103.73 (C-3'), 112.96 (C-3), 133.92 (C-6'), 138.35 (C-5'), 147.88 (C-4'), 152.16 (C-2'), 158.57 (C-8a), 162.63 (C-5), 163.16 (C-2), 165.39 (C-7), 181.98 (C=O, C-4).

Cycloartobiloxanthone (4) ¹³C NMR (125 MHz, acetone-*d*₆) δ 20.61 (C-9), 23.10 (C-13), 28.51 (C-18), 28.53 (C-17), 28.71 (C-12), 47.78 (C-10), 78.99 (C-16), 93.97 (C-11), 100.19 (C-6), 102.14 (C-8), 105.13 (C-4a), 105.64 (C-1'), 105.64 (C-3'), 112.95 (C-3), 116.21 (C-14), 128.16 (C-15), 133.96 (C-6'), 138.17 (C-5'), 147.30 (C-4'), 151.84 (C-2'), 152.31 (C-8a), 159.77 (C-5), 161.78 (C-2), 162.88 (C-7), 181.69 (C=O, C-4).

Artonin E (6) ¹³C NMR (125 MHz, acetone-*d*₆) δ 18.02 (C-13), 25.03 (C-9), 26.19 (C-12), 28.62 (C-17 and C-18), 79.12 (C-16), 100.06 (C-6), 101.94 (C-8), 105.07 (C-3'), 105.94 (C-4a), 111.76 (C-1'), 115.79 (C-14), 117.40 (C-6'), 121.99 (C-3), 122.84 (C-10), 128.36 (C-15), 132.67 (C-11), 139.45 (C-5'), 149.83 (C-4'), 150.13 (C-2'), 153.61 (C-8a), 160.29 (C-7), 162.59 (C-5), 163.09 (C-2), 183.63 (C=O, C-4).

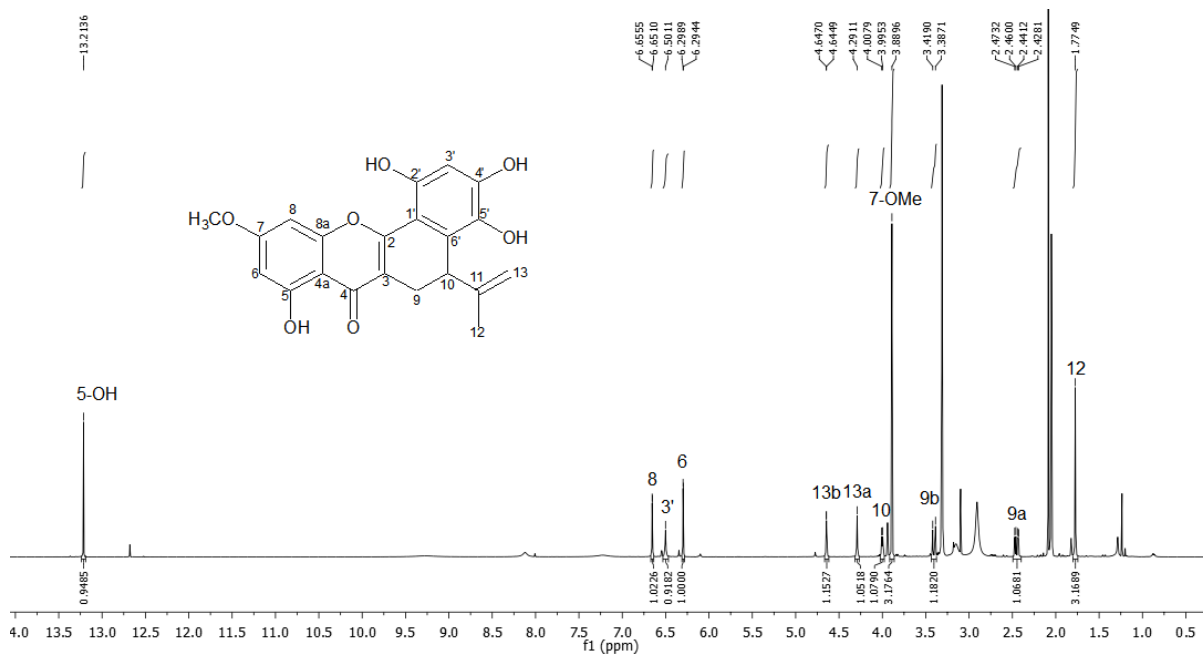


Figure S1. ¹H-NMR spectrum of compound 1 (500MHz, acetone-*d*₆)

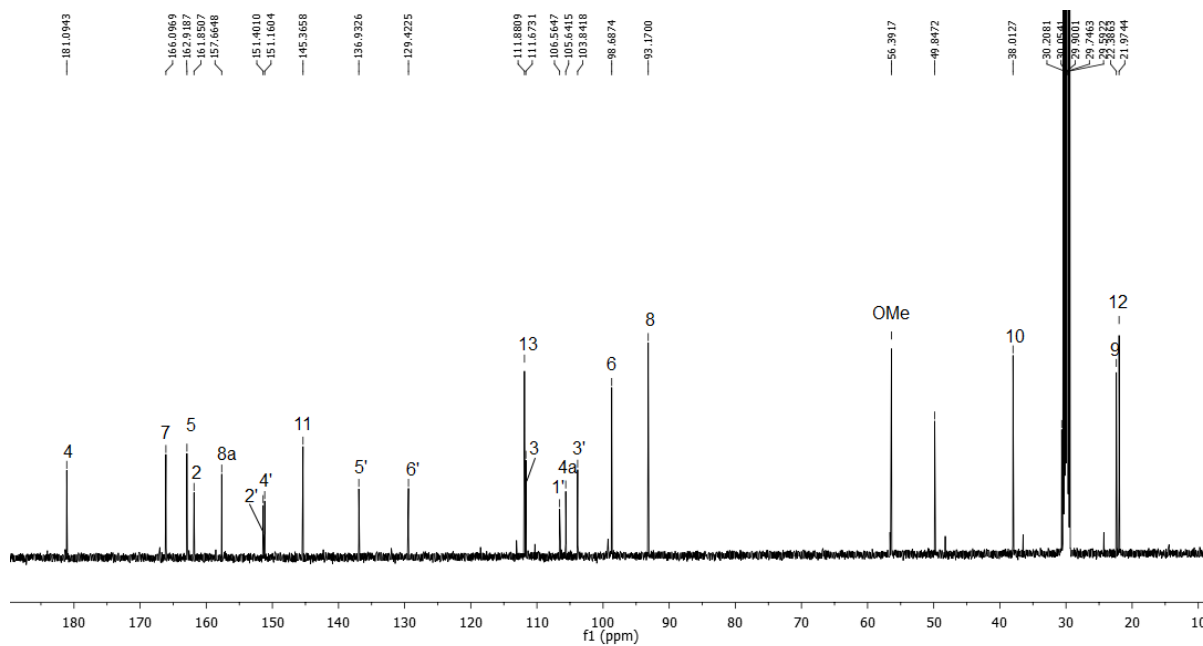


Figure S2. ¹³C-NMR spectrum of compound 1 (125MHz, acetone-*d*₆)

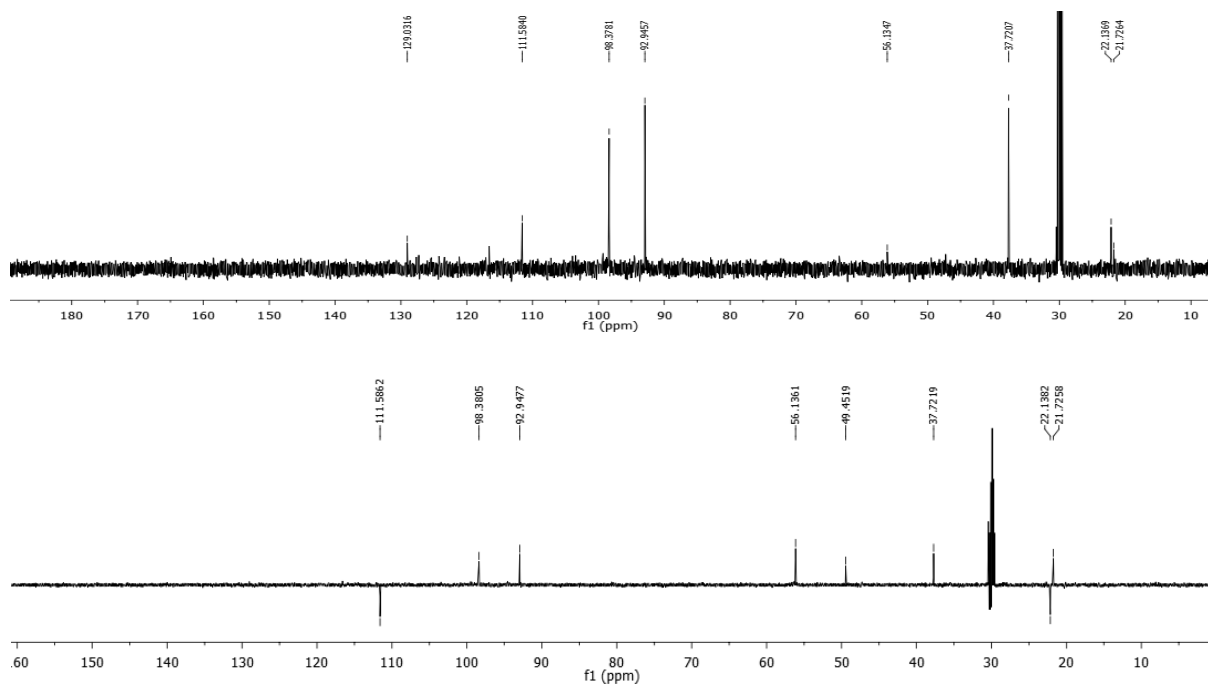


Figure S3. DEPT-90 and -135 spectrum of compound **1**.

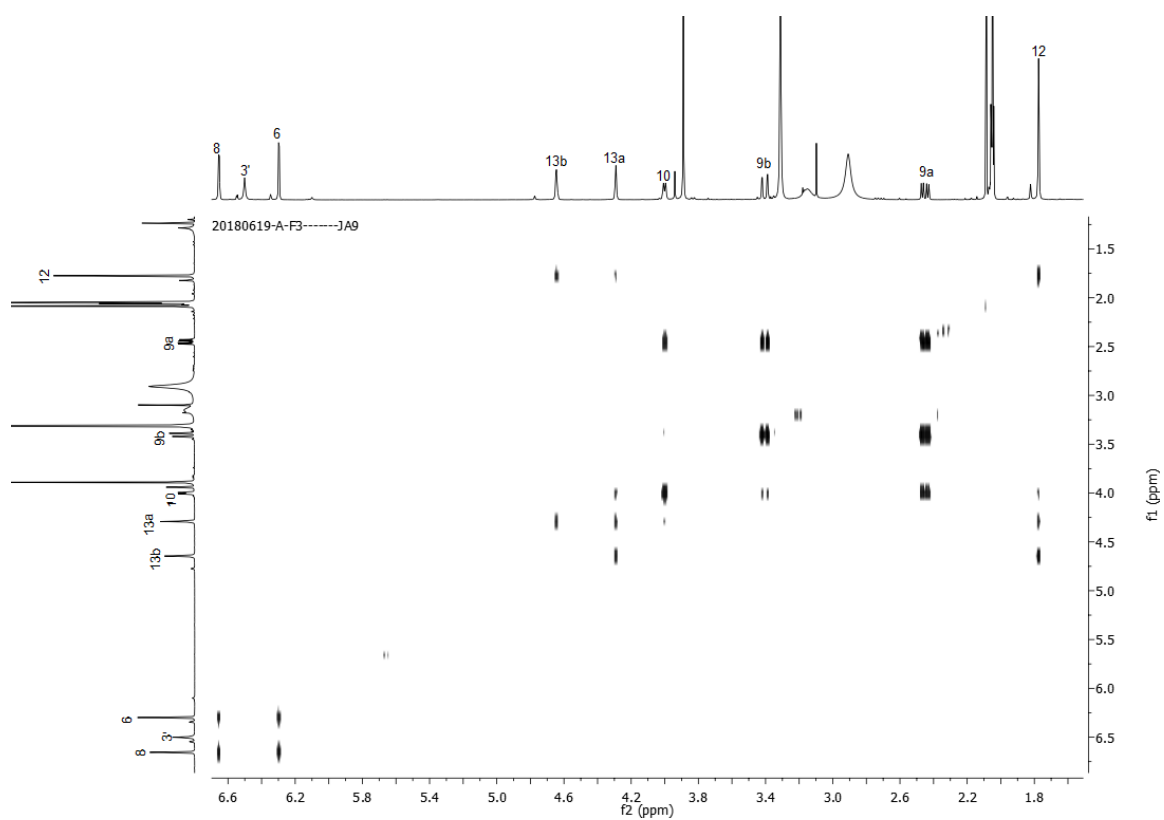


Figure S4. COSY spectrum of compound **1**.

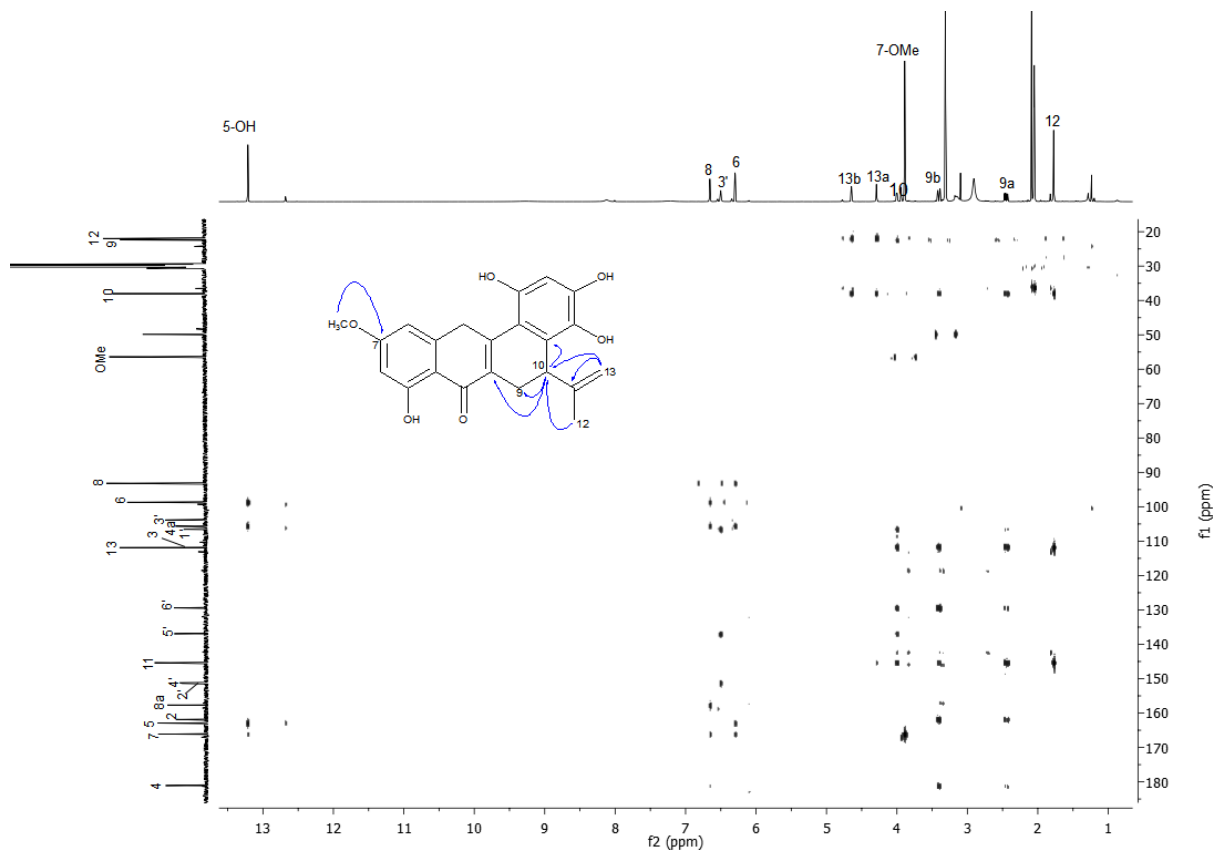


Figure S5. HMBC spectrum of compound 1.

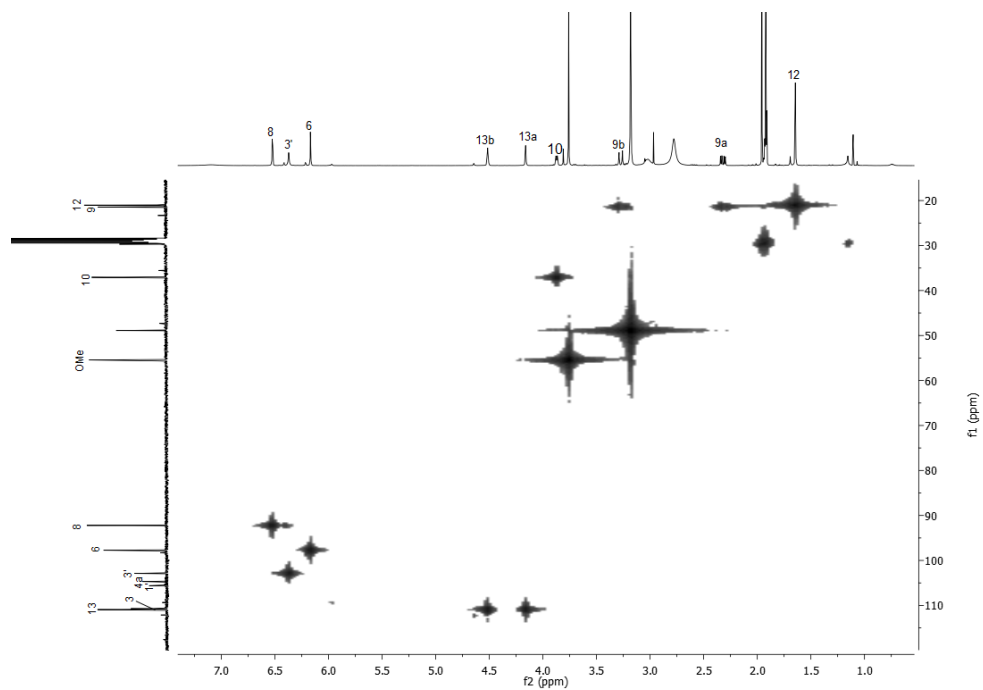
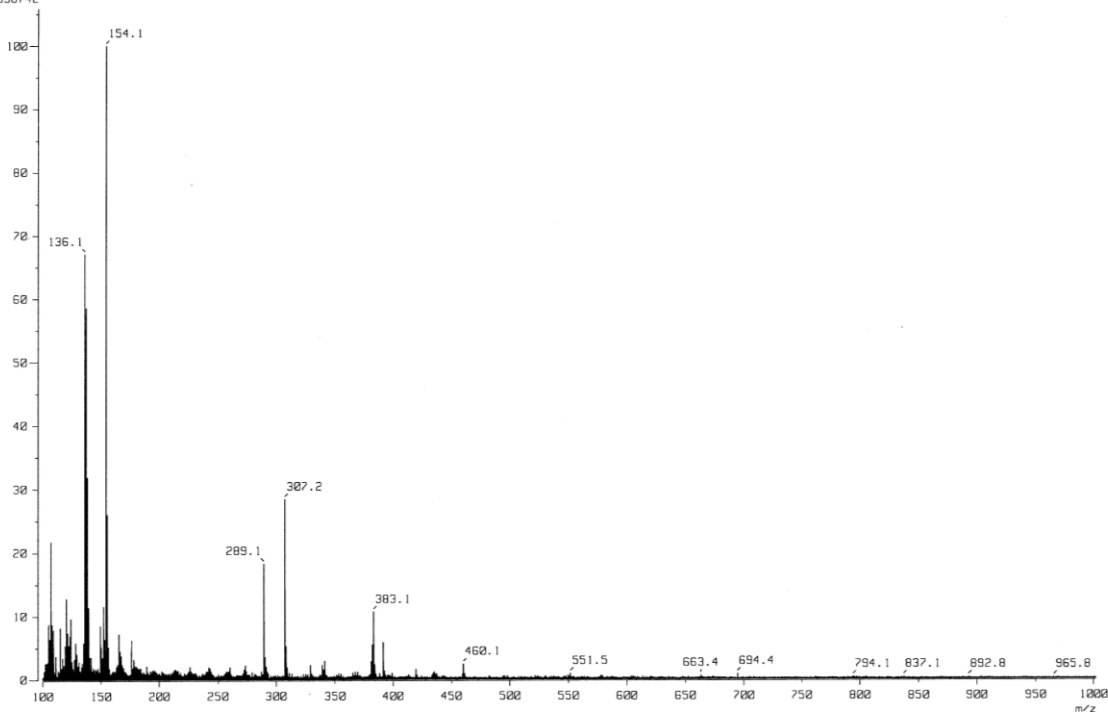


Figure S6. HMQC spectrum of compound 1.

[Mass Spectrum]
 Date : F3-1 Date : 21-Dec-2018 16:54
 Sample : -
 Note : -
 Inlet : Direct Ion Mode : FAB+
 Spectrum Type : Normal Ion [MF-Linear]
 RT : 0.45 min Scan# : 4
 BP : m/z 154.0629 Int. : 237.67
 Output m/z range : 100.0000 to 1002.0772 Cut Level : 0.00 %
 2638742



[Elemental Composition]

Pag

Data : A-F3_HR Date : 10-Jan-2019 16:29
 Sample : -
 Note : -
 Inlet : Direct Ion Mode : FAB+
 RT : 4.67 min Scan# : 57
 Elements : C 110/1, H 130/1, O 8/1
 Mass Tolerance : 100ppm, 3mmu if m/z > 30
 Unsaturation (U.S.) : 0.0 - 30.0

Observed m/z	Int%	Err [ppm / mmu]	U.S. Composition
383.1149	100.0	+4.8 / +1.9	12.5 C 21 H 19 O 7

Figure S7. FABMS and HRFABMS data of compound 1.

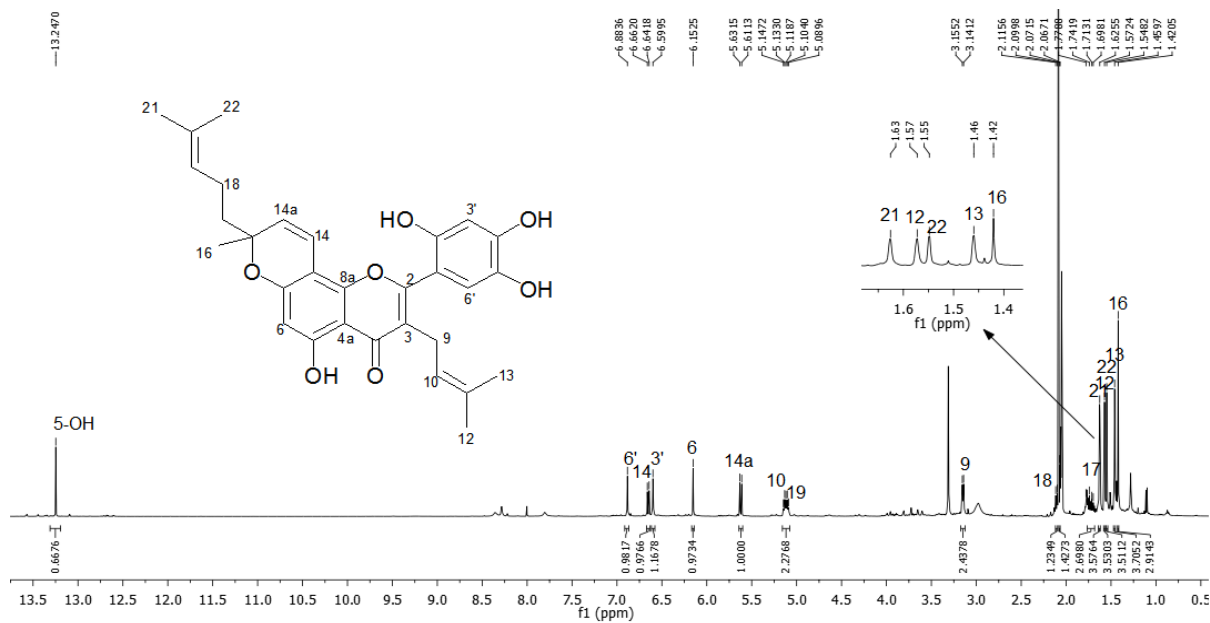


Figure S8. $^1\text{H-NMR}$ spectrum of compound **5** (500MHz, acetone- d_6)

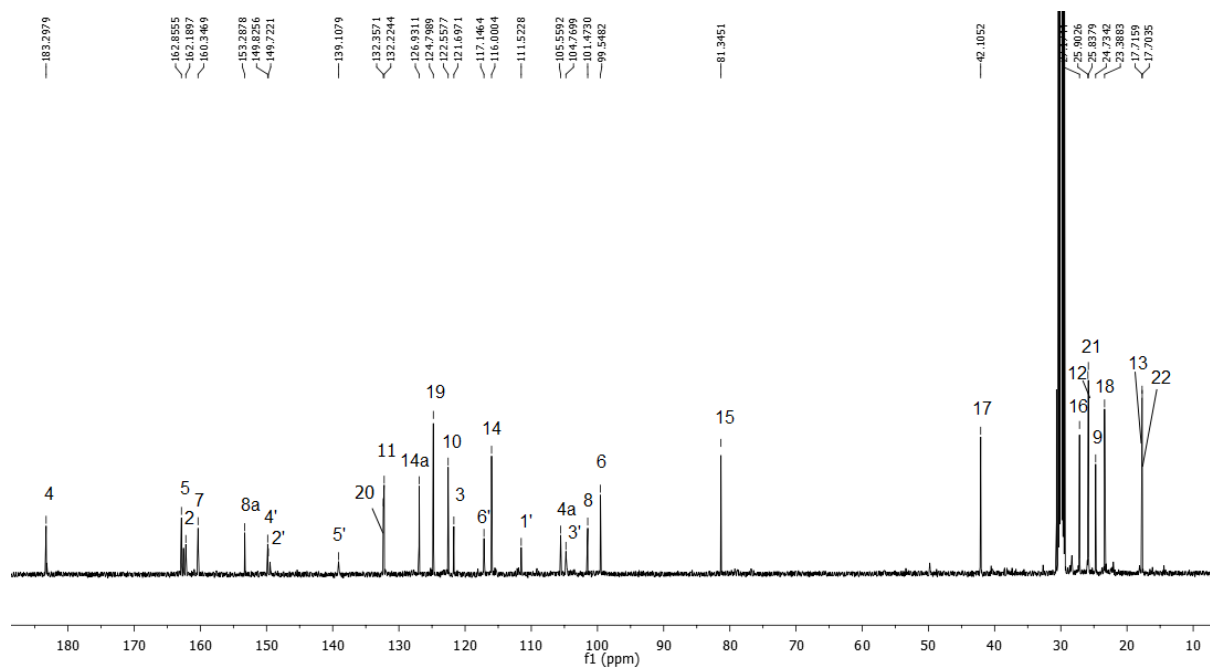


Figure S9. $^{13}\text{C-NMR}$ spectrum of compound **5** (125MHz, acetone- d_6)

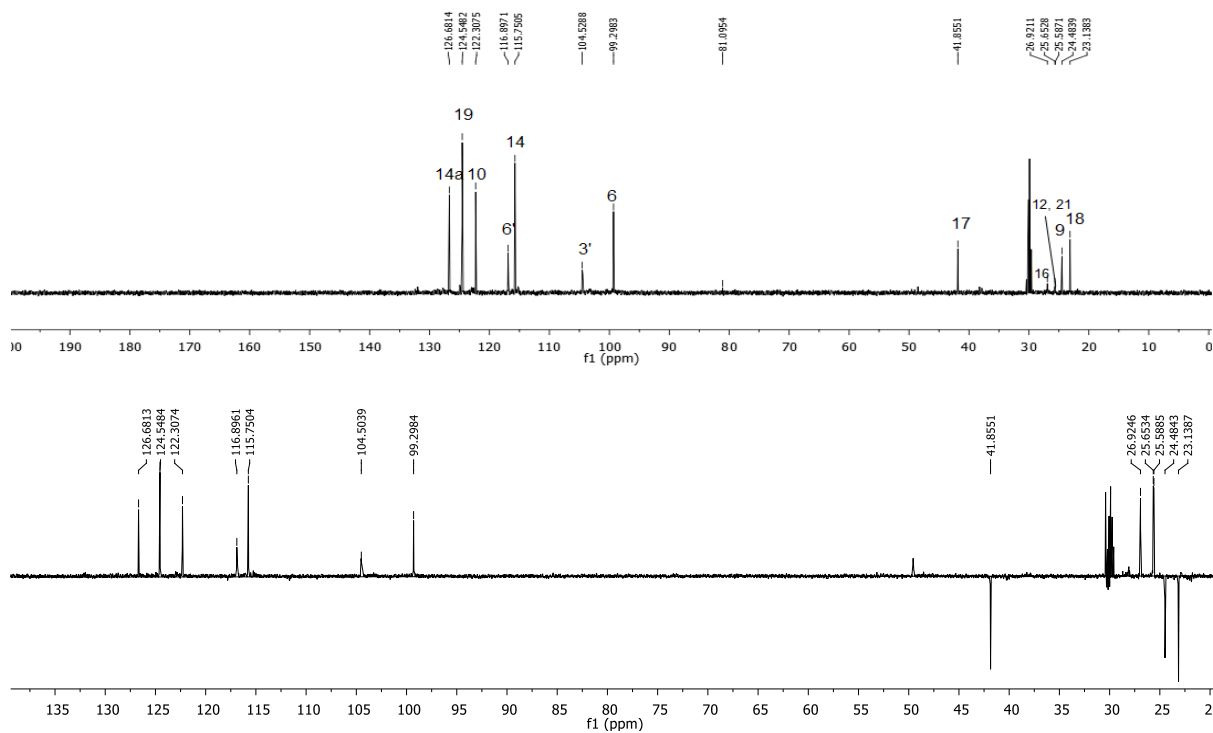


Figure S10. DEPT135 spectrum of compound 5.

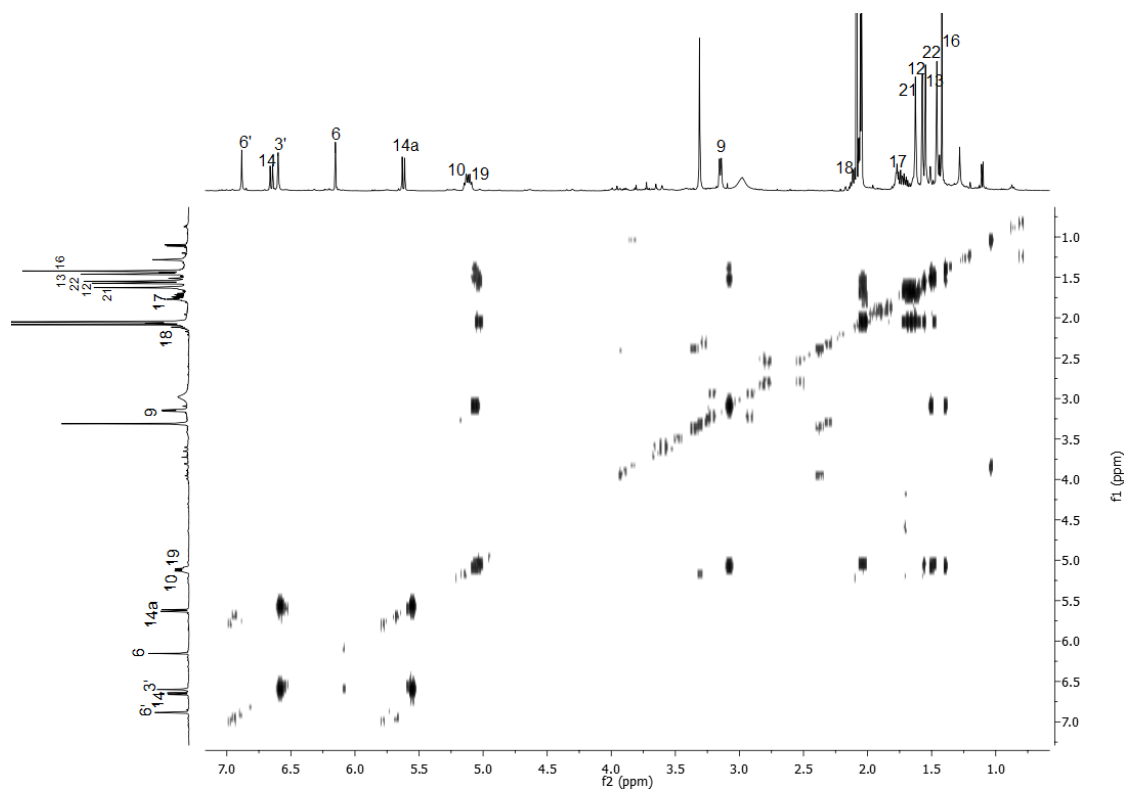


Figure S11. COSY spectrum of compound 5.

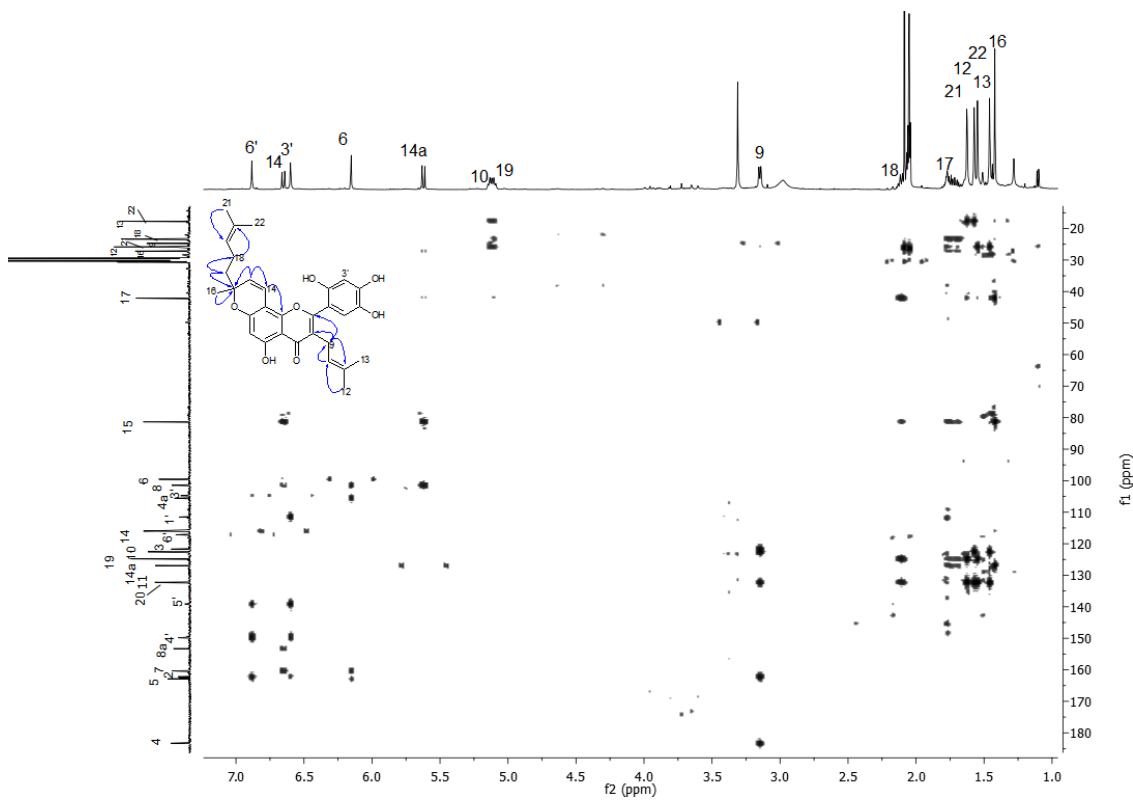


Figure S12. HMBC spectrum of compound 5.

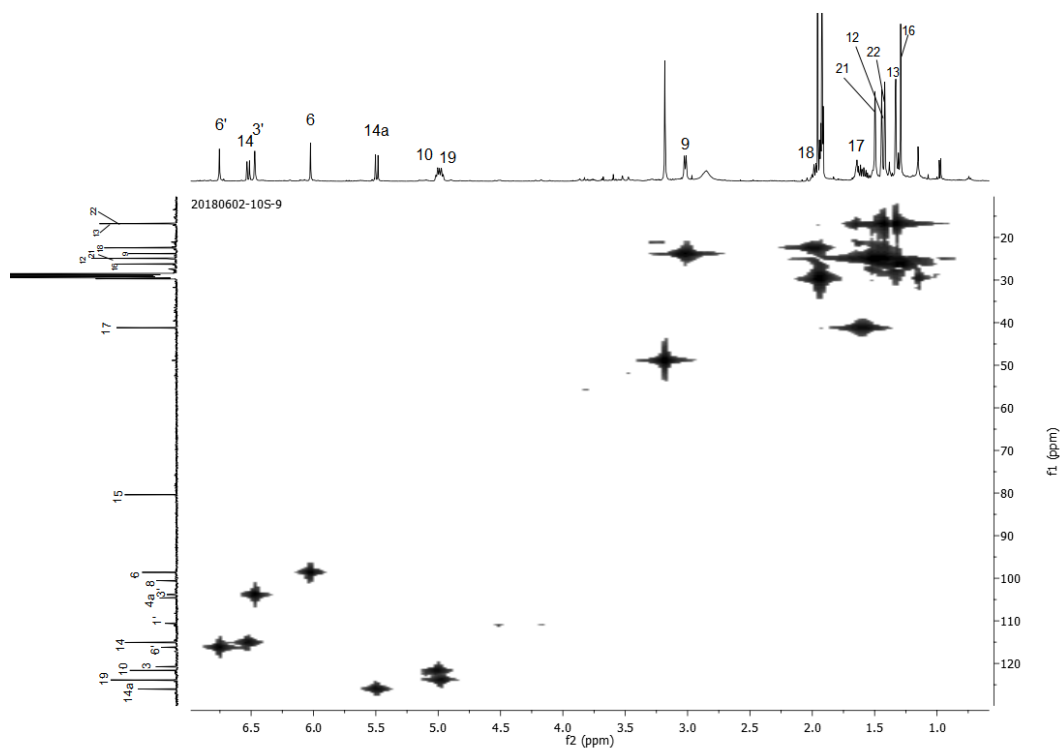
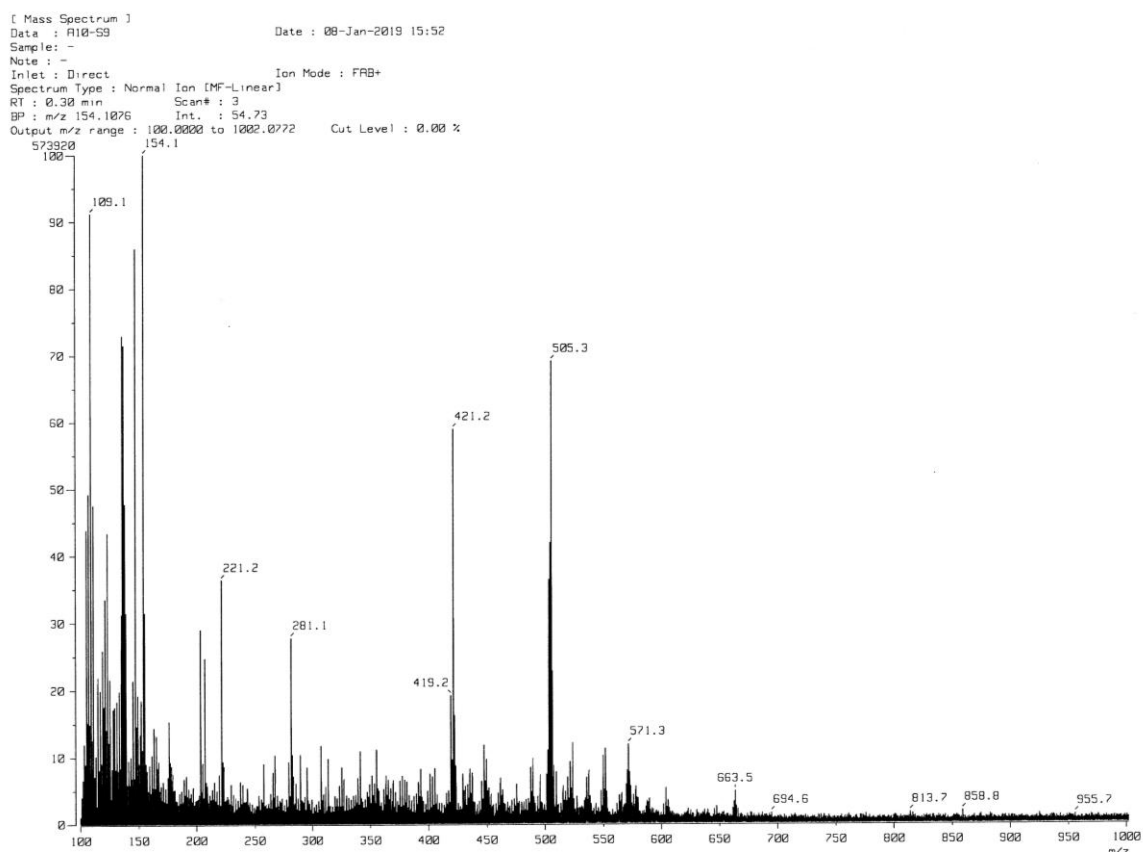


Figure S13. HMQC spectrum of compound 5.



[Elemental Composition]
 Data : A10-S9_HR Date : 08-Jan-2019 17:31
 Sample : -
 Note : -
 Inlet : Direct Ion Mode : FAB+
 RT : 4.84 min Scan# : 59
 Elements : C 110/1, H 130/1, O 8/1
 Mass Tolerance : 100ppm, 3mmu if m/z > 30
 Unsaturation (U.S.) : 0.0 - 30.0

Observed m/z	Int%	Err [ppm / mmu]	U.S.	Composition
505.2216	34.6	-1.9 / -1.0	14.5	C 30 H 33 O 7

Figure S14. FABMS and HRFABMS data of compound 5.

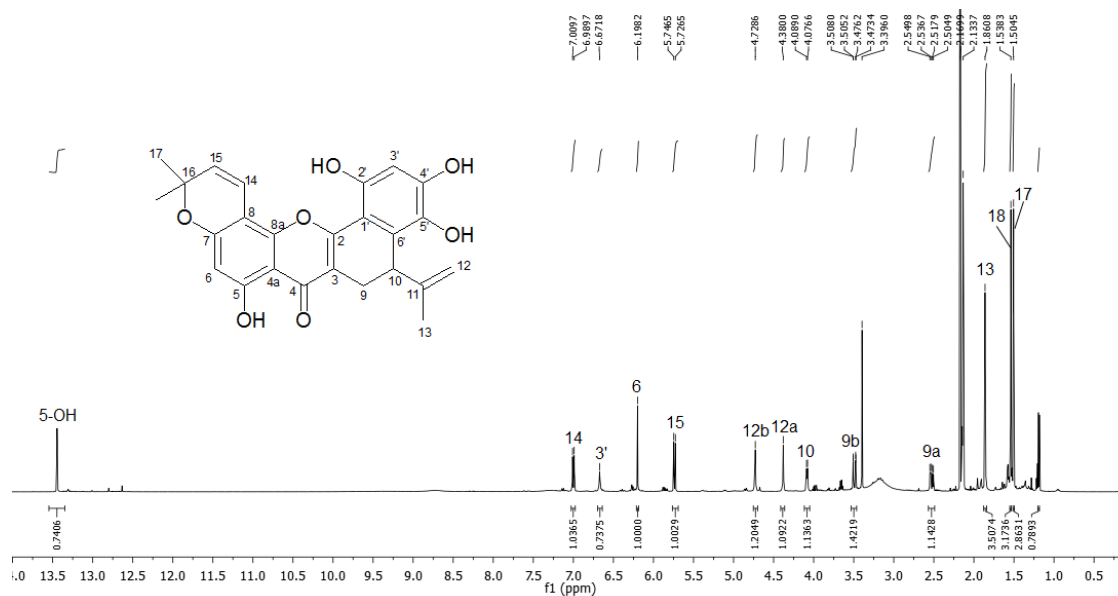


Figure S15. $^1\text{H-NMR}$ spectrum of compound **2** (500MHz, acetone- d_6)

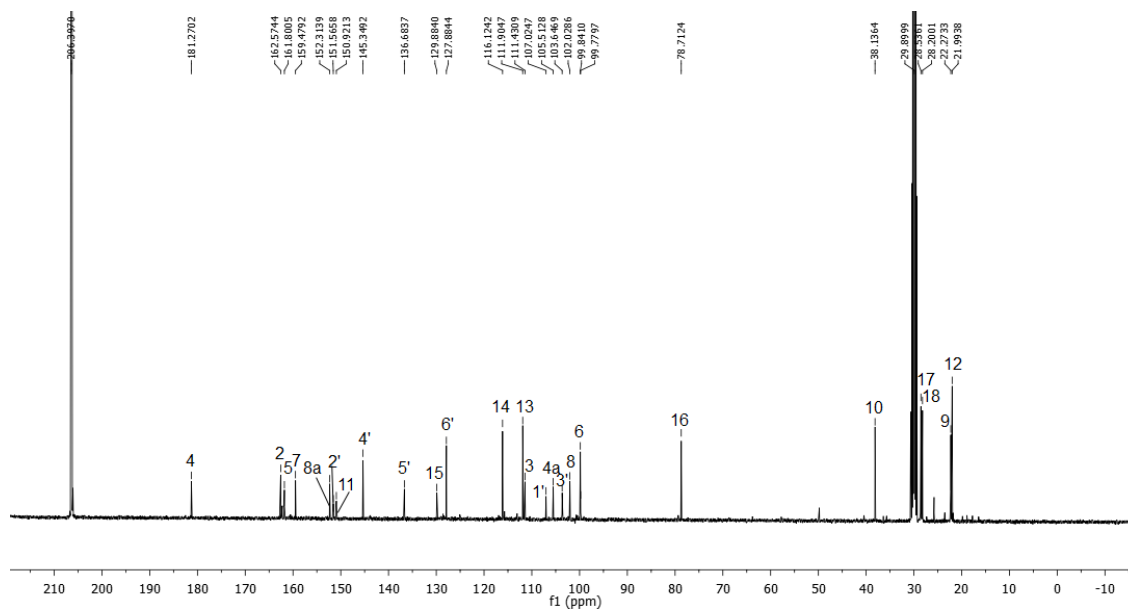


Figure S16. $^{13}\text{C-NMR}$ spectrum of compound **2** (125MHz, acetone- d_6)

Data: A-F2_HR Date: 19-Jun-2018 14:03
 Instrument: MStation
 Sample: -
 Note: -
 Inlet: Direct Ion Mode: EI+
 RT: 1.80 min Scan#: 28
 Elements: C 100/1, H 100/1, O 10/1
 Mass Tolerance : 1000ppm, 3mmu if m/z > 3
 Unsaturation (U.S.): -0.5 - 20.0

Observed m/z	Int%	Err [ppm / mmu]	U.S. Composition
1 434.1363	21.31	-0.6 / -0.3	15.0 C25 H22 O7

Figure S17. HR-ESI-MS data of compound **2**.

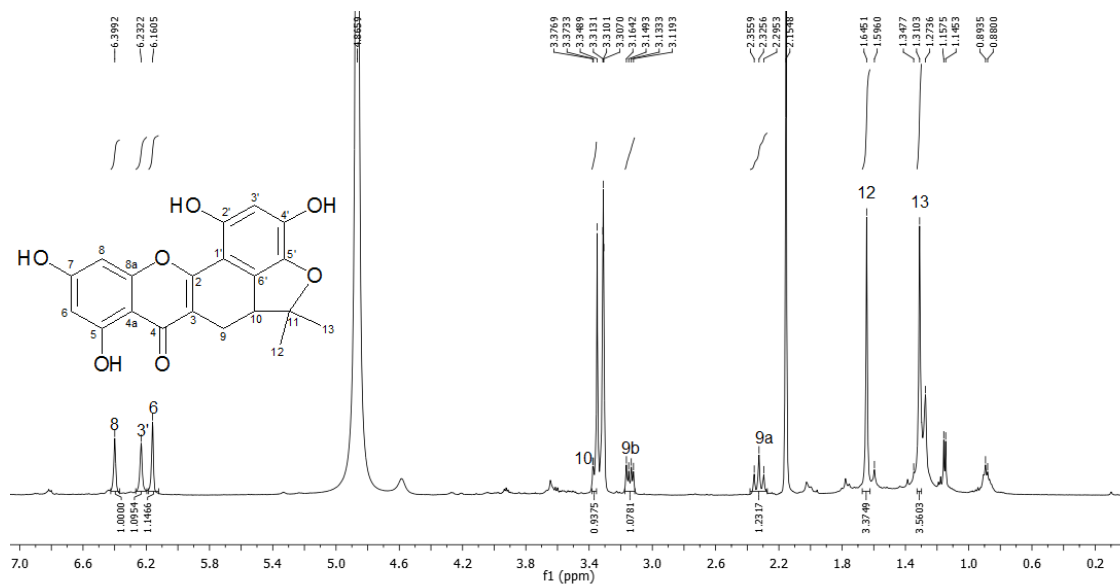


Figure S18. $^1\text{H-NMR}$ spectrum of compound **3** (500MHz, acetone- d_6)

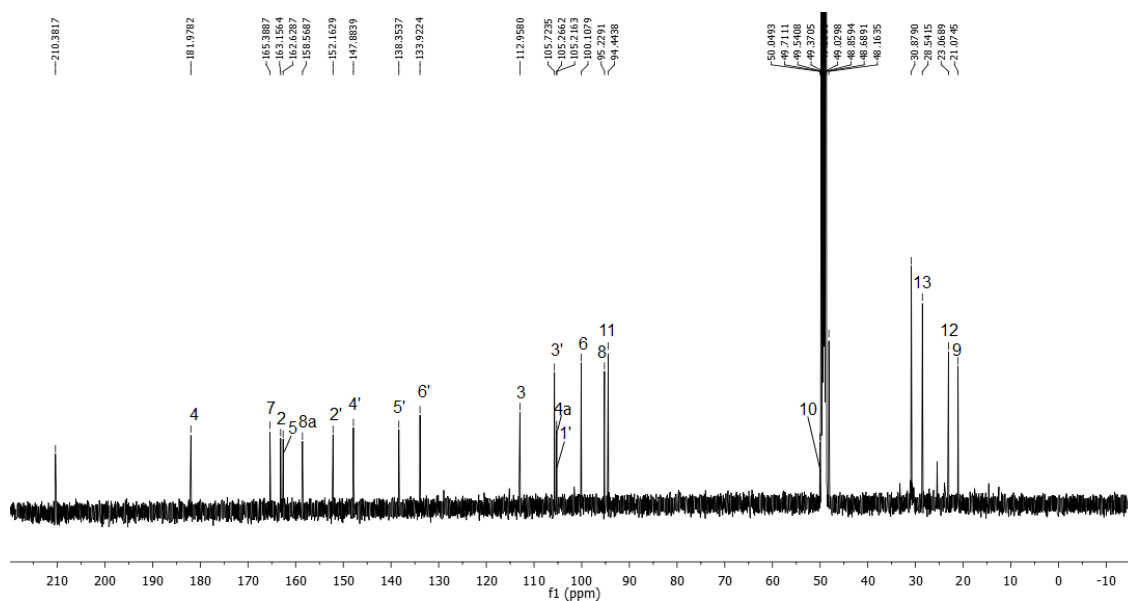


Figure S19. $^{13}\text{C-NMR}$ spectrum of compound **3** (125MHz, acetone- d_6)

Data : F3-4_HR Date : 19-Jun-2018 15:31
 Instrument : MStation
 Sample : -
 Note : -
 Inlet : Direct Ion Mode : EI+
 RT : 1.94 min Scan# : 30
 Elements : C 100/1, H 100/1, O 10/1
 Mass Tolerance : 1000ppm, 3mmu if m/z > 3
 Unsaturation (U.S.) : -0.5 - 20.0

	Observed m/z	Int%	Err [ppm / mmu]	U.S. Composition
1	368.0899	21.53	+0.8 / +0.3	13.0 C20 H16 O7

Figure S20. HR-ESI-MS data of compound **3**.

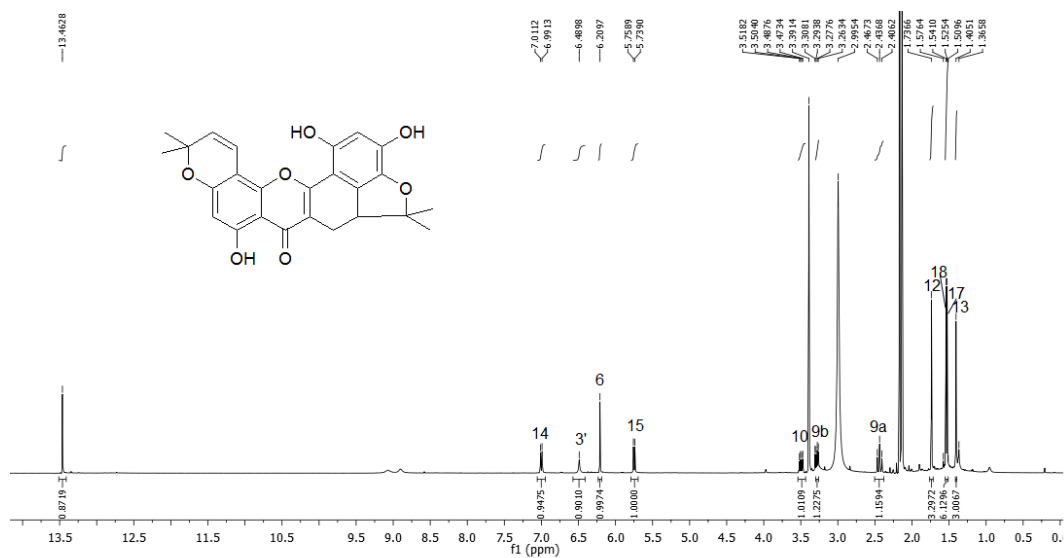


Figure S21. ¹H-NMR spectrum of compound 4 (500MHz, acetone-*d*₆)

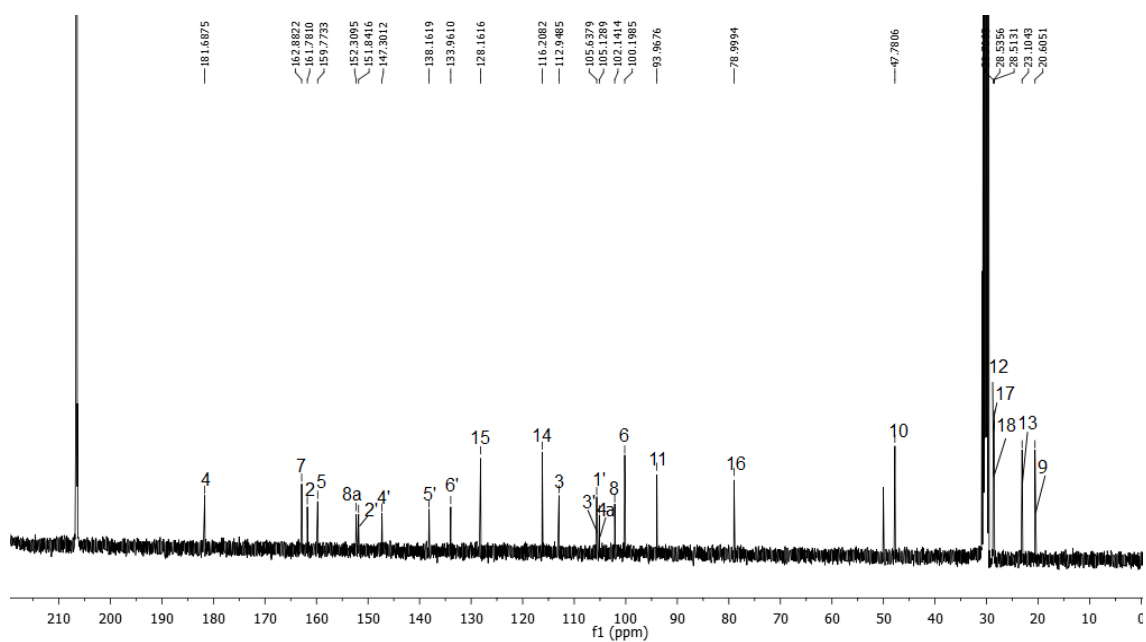


Figure S22. ¹³C-NMR spectrum of compound 4 (125MHz, acetone-*d*₆)

Data: F11_HR Date: 19-Jun-2018 14:15
 Instrument: MStation
 Sample: -
 Note: -
 Inlet: Direct Ion Mode: EI+
 RT: 3.14 min Scan#: 48
 Elements: C 100/1, H 100/1, O 10/1
 Mass Tolerance: 1000ppm, 3mmu if m/z > 3
 Unsaturation (U.S.): -0.5 - 20.0

Observed m/z	Int%	Err[ppm / mmu]	U.S. Composition
1 434.1364	22.03	-0.4 / -0.2	15.0 C25 H22 O7

Figure S23. HR-ESI-MS data of compound 4.

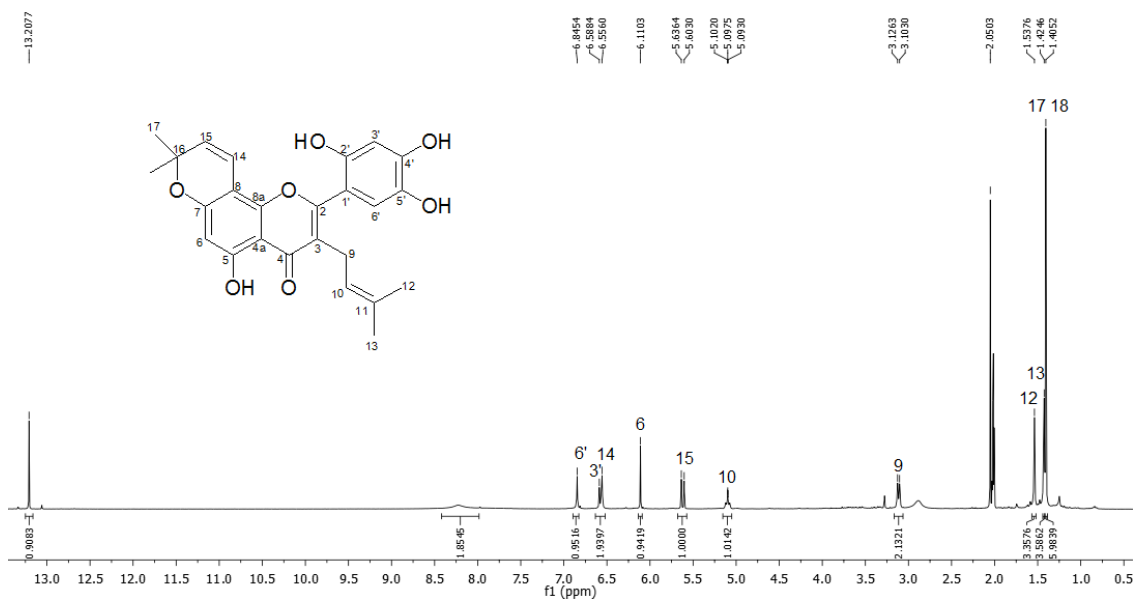


Figure S24. ^1H -NMR spectrum of compound **6** (500MHz, acetone- d_6)

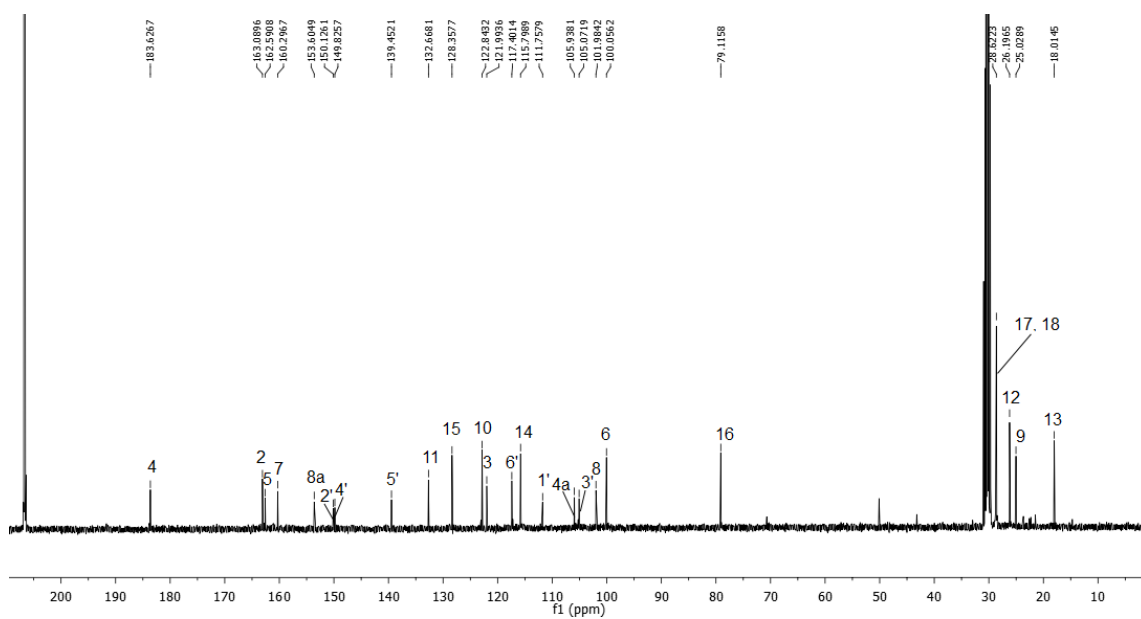


Figure S25. ^{13}C -NMR spectrum of compound **6** (125MHz, acetone- d_6)

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[ Elemental Composition ]
Data : A-F8_HR                               Date : 08-Jan-2019 17:22
Sample: -                                     Ion Mode : FAB+
Note : -                                       Scan#: 8
Inlet : Direct
RT : 0.59 min
Elements : C 110/1, H 130/1, O 8/1
Mass Tolerance : 100ppm, 5mmu if m/z > 50
Unsaturation (U.S.) : 0.0 - 30.0

Observed m/z Int% Err [ppm / mmu] U.S. Composition
437.1635 20.4 +7.8 / +3.4 13.5 C 25 H 25 O 7
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Figure S26. HR-ESI-MS data of compound **6**.

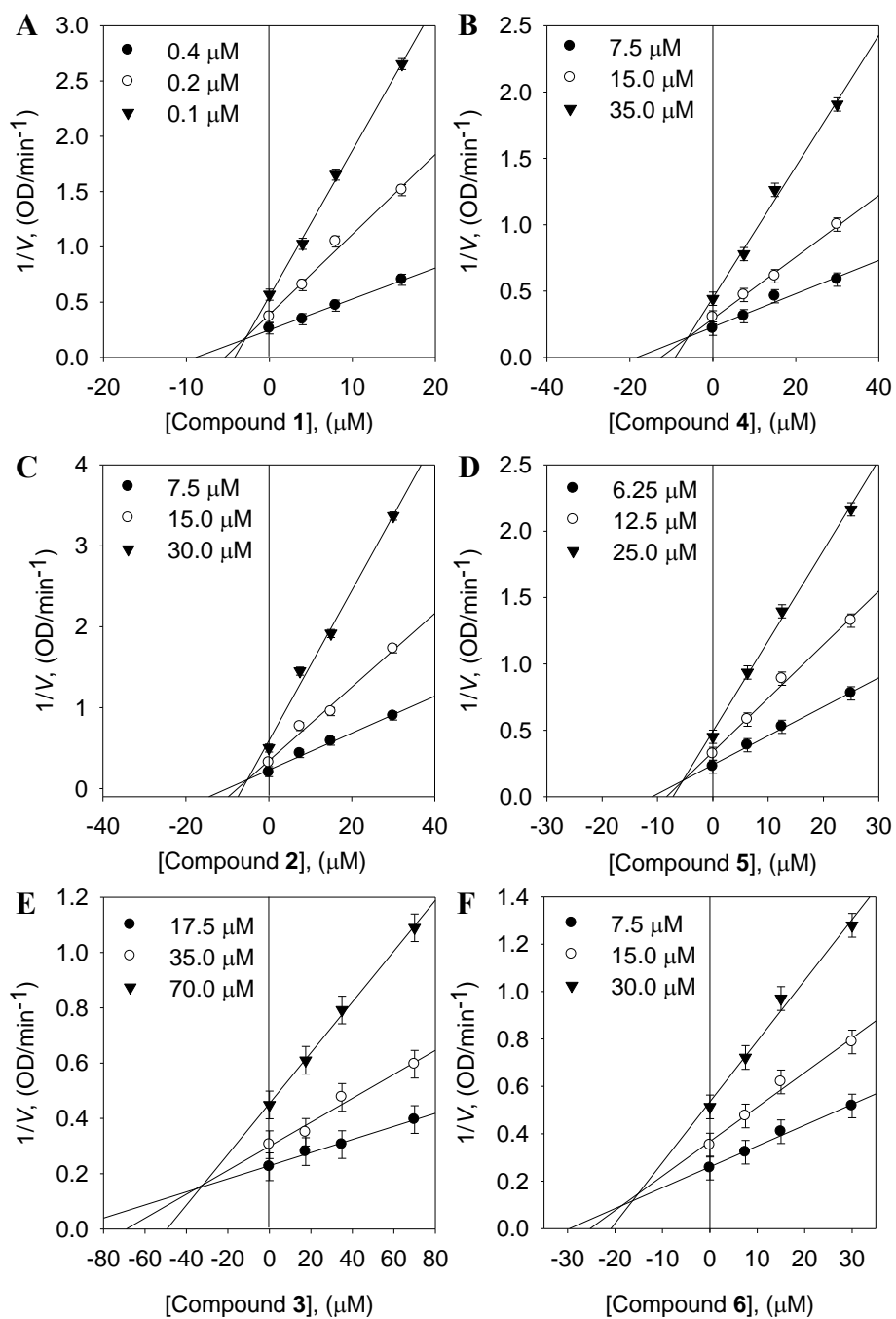


Figure S27. Dixon plots for α -glucosidase of compounds 1-6.

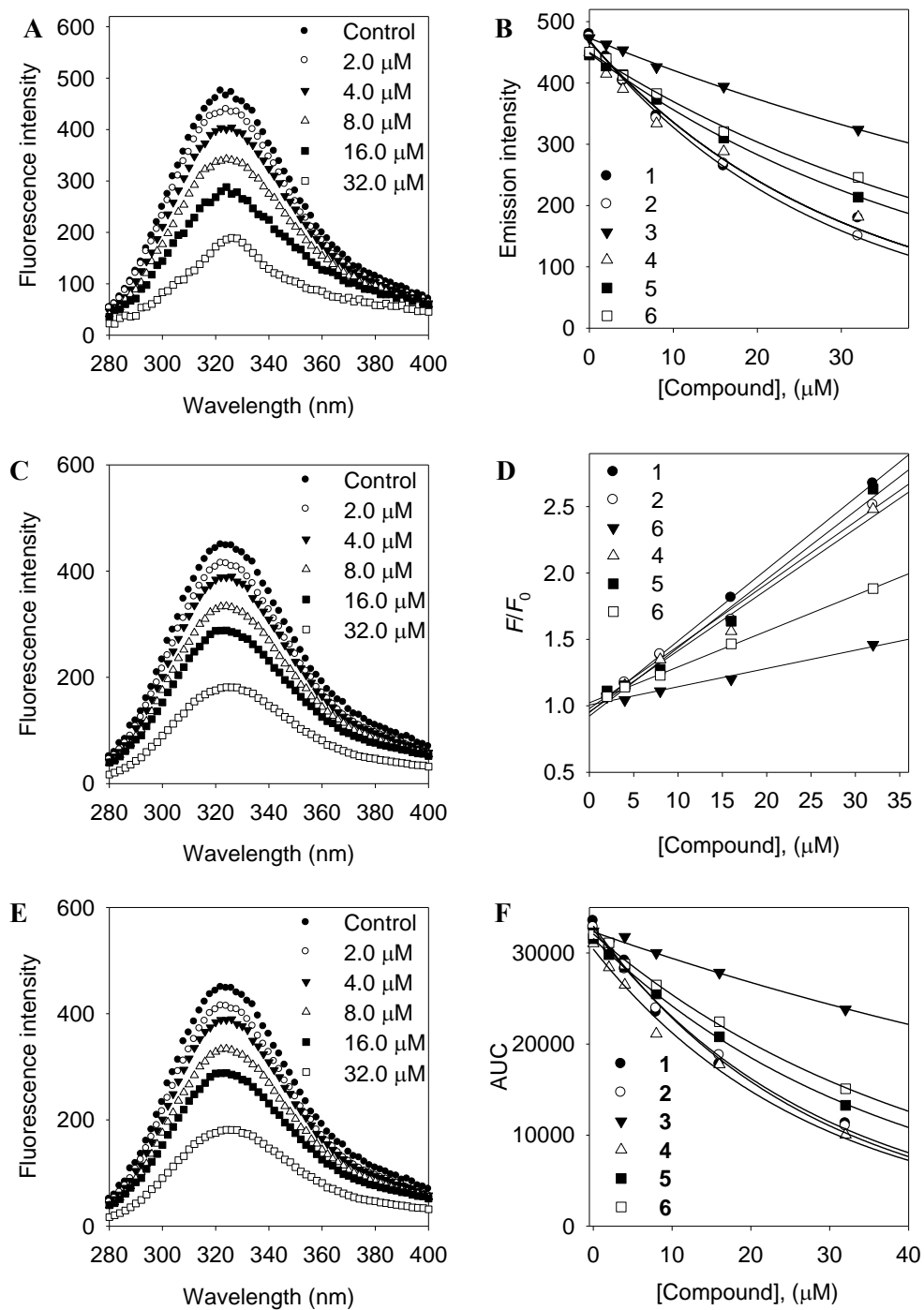


Figure S28. (A) The fluorescence effect of compound 2, (C) compound 4 and (E) compound 5 on fluorescence emission spectra of α -glucosidase, (B) Decrease in intensity of the emission plots of compounds, (D) Stern-Volmer plots of compounds at 37 $^{\circ}$ C, (F) Area under curve (AUC) of compounds.

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1  MTISDHPETE  PKWWKEATIY  QIYPASFKDS  NNDGWGDLKG  ITSKLQYIKD
51  LGVDAIWVCP  FYDSPQQDMG  YDISNYEKVW  PTYGTNEDCF  ELIDKTHKLG
101 MKFITDLVIN  HCSTEHEWFK  ESRSSKTNPK  RDWFFWRPPK  GYDAEGKPIP
151 PNNWKSFFGG  SAWTFDETTN  EFYLRLFASR  QVDLNWENED  CRRAIFESAV
201 GFWLDHGVDG  FRIDTAGLYS  KRPGLPDSPI  FDKTSKLQHP  NWGSHNGPRI
251 HEYHQELHRF  MKNRVKDGRE  IMTVGEVAHG  SDNALYTSAA  RYEVSEVFSF
301 THVEVGTSPF  FRYNIVPFTL  KQWKEAIASN  FLFINGTDSW  ATTYIENHDQ
351 ARSITRFADD  SPKYRKISGK  LLTLLECSLT  GTLYVYQQQE  IGQINFKEWP
401 IEKYEDVDVK  NNYEIIKSF  GKNSKEMKDF  FKGIALLSRD  HSRTPMPWTK
451 DKPNAGFTGP  DVKPWFLLNE  SFEQGINVEQ  ESRDDDSVLN  FWKRALQARK
501 KYKELMIYGY  DFQFIDLSD  QIFSFTKEYE  DKTLFAALNF  SGEEIEFSLP
551 REGASLSFIL  GNYDDTDVSS  RVLKPWEGRI  YLVK

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Figure S29. Amino acid sequence of *Saccharomyces cerevisiae* α -glucosidase. The Trp(W), Phe(F), and Tyr(Y) residues are shown as yellow, cyan, and pink respectively.

Trp(W): 20, Phe(F): 41, Tyr(Y): 26

Table S1. ^1H and ^{13}C spectroscopic data of (δ in ppm, J in Hz) **1** and **5** (acetone- d_6)

No.	1		5	
	$\delta_{\text{H}} J$ (Hz)	δ_{C} m	$\delta_{\text{H}} J$ (Hz)	δ_{C} m
2		161.9		162.2
3		111.7		121.7
4		181.1		183.3
4a		105.6		105.6
5		162.9		162.8
6	6.29 d (2.25)	98.7	6.15 s	99.6
7		166.1		160.4
8	6.65 d (2.25)	93.2		101.5
8a		157.7		153.2
9a	2.43 dd (15.98)	22.4	3.14 d (7)	24.7
9b	3.39 d (15.95)			
10	3.99 d (6.3)	38.0	5.09 m	122.6
11		145.4		132.2
12	1.78 s	22.0	1.57 s	25.9
13a	4.29 s	111.9	1.46 s	17.7
13b	4.65 s			
14			5.61 d (10.1)	116.0
14a			5.61 d (10.1)	126.9
15				81.4
16			1.42 s	27.2
17			1.69 m	42.1
18			2.07 m	23.4
19			5.09 m	124.8
20				132.4
21			1.63 s	25.8
22			1.54 s	17.7
1'		106.6		111.5
2'		151.4		149.7
3'	6.50 s	103.8	6.59 s	104.8
4'		151.2		149.8
5'		136.9		139.1
6'		129.4	6.88 s	117.2
OMe	3.89 s	56.4		

Table S2. Determination of K_{ik}/K_{iv} ratios by α -glucosidase enzyme inhibitory behaviours.

Compounds	[I] (μ M)	V_{max}	K_m	K_{ik}/K_{iv}
2	0	9.337	186.0878	-
	7.5	9.7056	740.5780	76.6587
	15	9.6712	868.3559	102.4604
	30	9.8012	2117.5032	209.3396
3	0	6.5274	193.6945	-
	17.5	6.6800	301.7435	23.8593
	35	6.7889	439.0428	31.6240
	70	6.7204	627.7285	75.7804
4	0	6.8027	202.3061	-
	7.5	6.3816	397.2304	16.6987
	15	7.3692	812.0855	11.4837
	30	7.4627	1318.5821	20.2662
6	0	6.8013	194.8085	-
	7.5	5.0352	265.5740	3.7049
	15	4.2992	319.4798	2.5098
	30	3.6219	365.4835	1.8101