

Supporting Information

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Isolation and Characterization of New Constituents from *Tricholepis eburnea*

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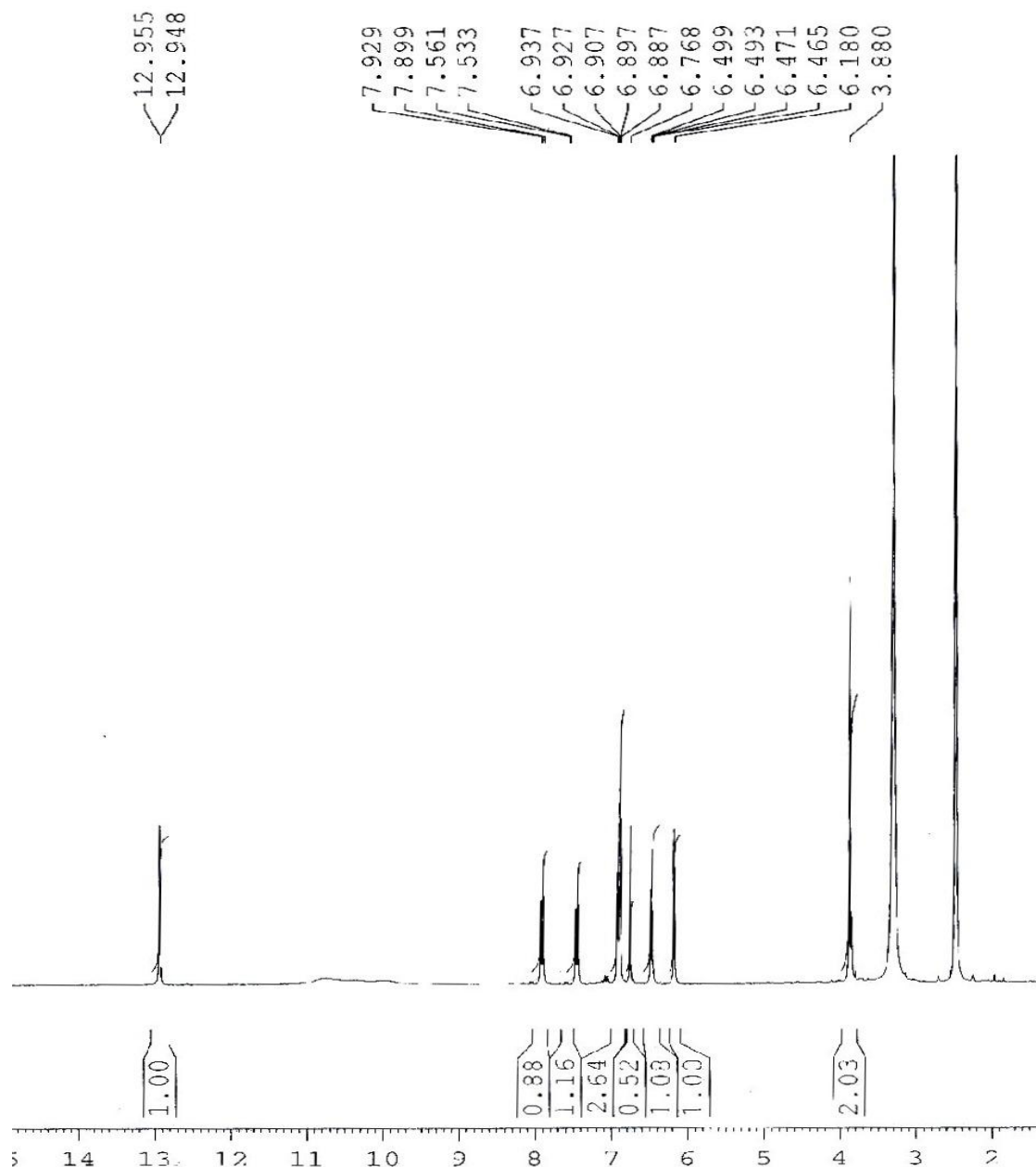
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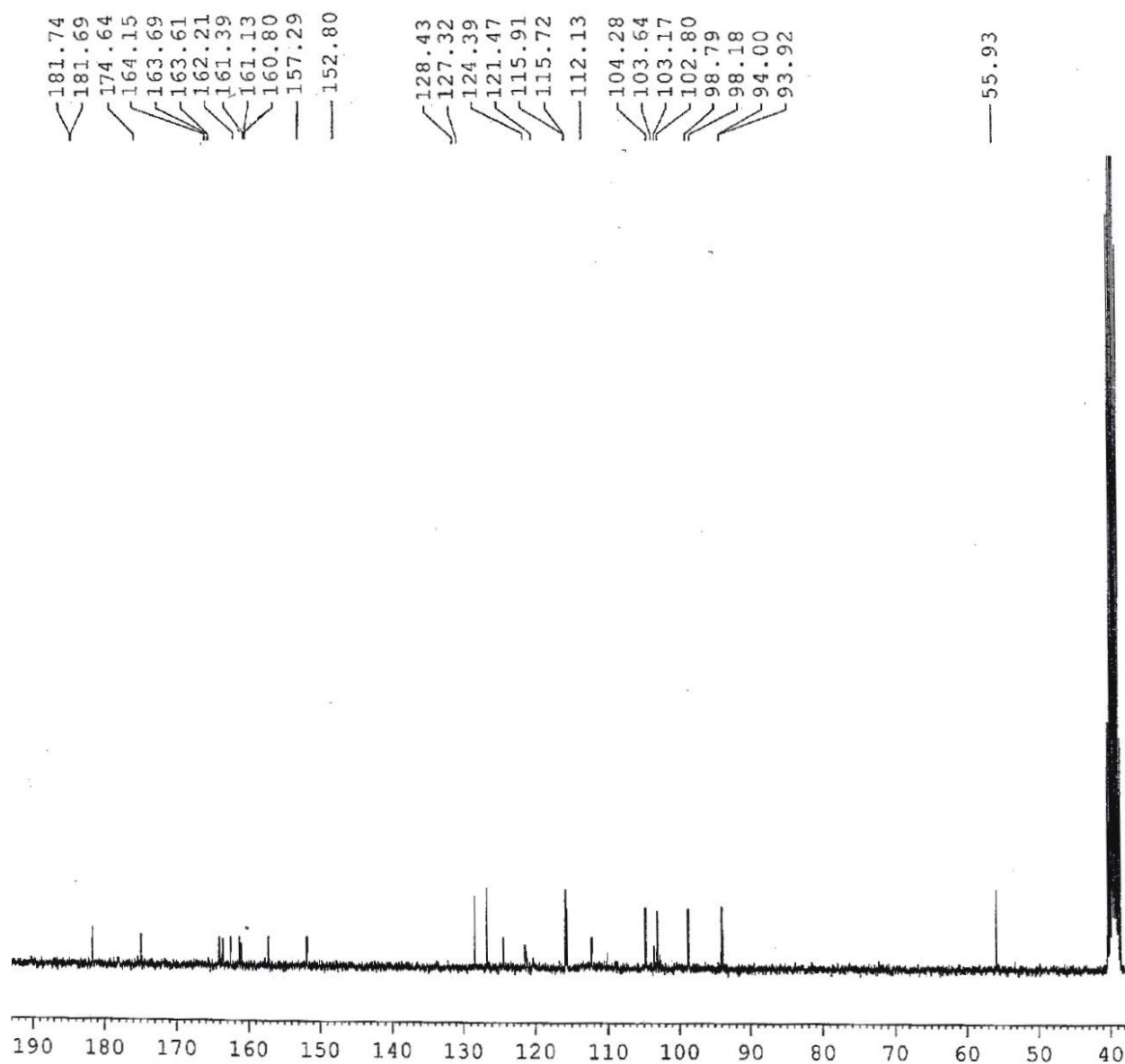
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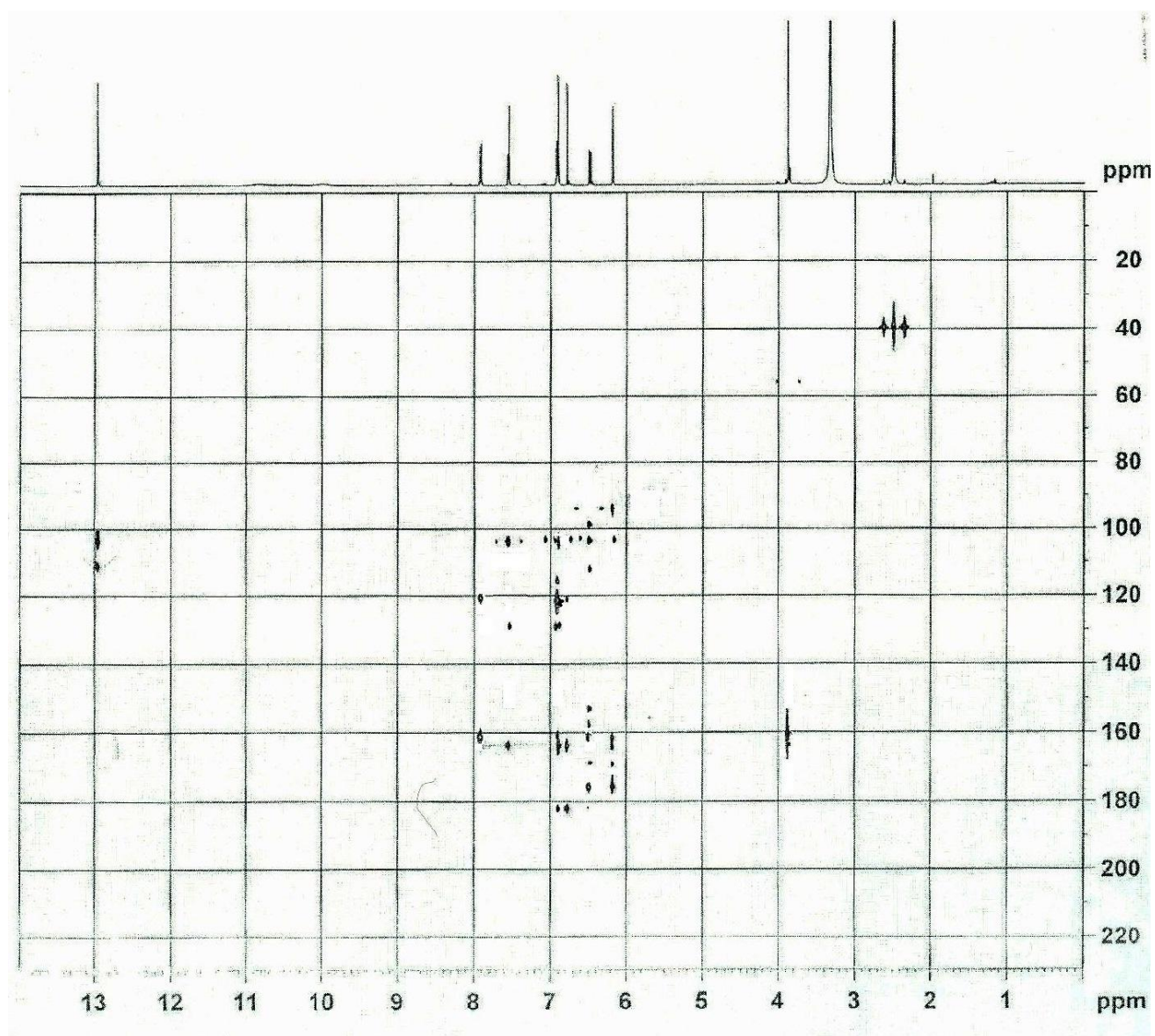
S1: ^1H -NMR (300 MHz, DMSO) Spectrum of Compound **1** (trichonoide A)

Trichonoide A (**1**): Yellow solid. IR ($\nu_{\max}$ cm^{-1}): 3389 (OH), 1676 (conjugated carbonyl), 1648 and 1481 cm^{-1} (aromatic moiety). UV ($\lambda_{\max}$ nm) 268 and 338 nm. ^1H -NMR (DMSO, 300 MHz), δ : 6.76 (1H, s, H-3), 6.18 (1H, d, overlapped, H-6), 6.46 (1H, d, $J = 1.8$, H-8), 7.92 (2H, d, $J = 9$, H-2', -6'), 6.93 (2H, d, overlapped, H-3', -5'), 6.88 (1H, s, H-3''), 6.18 (1H, d, overlapped, H-6''), 6.49 (1H, d, $J = 1.8$, H-8''), 7.56 (2H, d, $J = 8.4$, H-2''', -6'''), 6.90 (2H, d, overlapped, H-3''', -5'''), 3.95 (s, OCH_3 (4'')). ^{13}C -NMR (CDCl_3 , 75 MHz), δ : 163.7 (C-2), 103.1 (C-3), 181.7 (C-4), 103.6 (C-4a), 162.2 (C-5), 98.1 (C-6), 164.1 (C-7), 94.0 (C-8), 157.2 (C-8a), 121.1 (C-1'), 128.4 (C-2', -6'), 115.7 (C-3', -5'), 161.1 (C-4'), 163.6 (C-2''), 104.2 (C-3''), 181.6 (C-4''), 112.1 (C-4a''), 160.8 (C-5''), 98.7 (C-6''), 174.6 (C-7''),

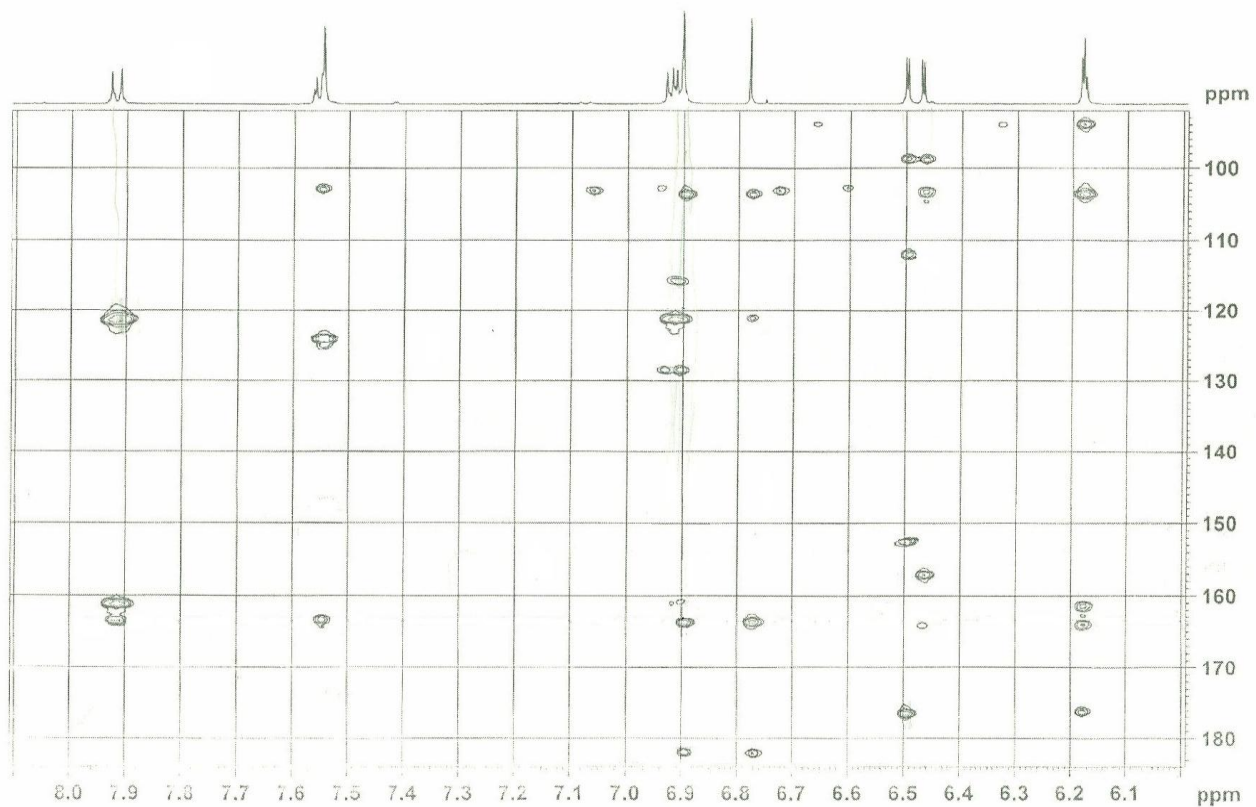
93.9 (C-8''), 152.8 (C-8a''), 124.4 (C-1'''), 127.3 (C-2''', -6'''), 115.9 (C-3''', -5'''), 161.3 (C-4'''), 55.9 (OCH₃(C-4''')). HR-ESI-MS: $m/z = 537.1180$ [M+H]⁺ for formula C₃₁H₂₁O₉.



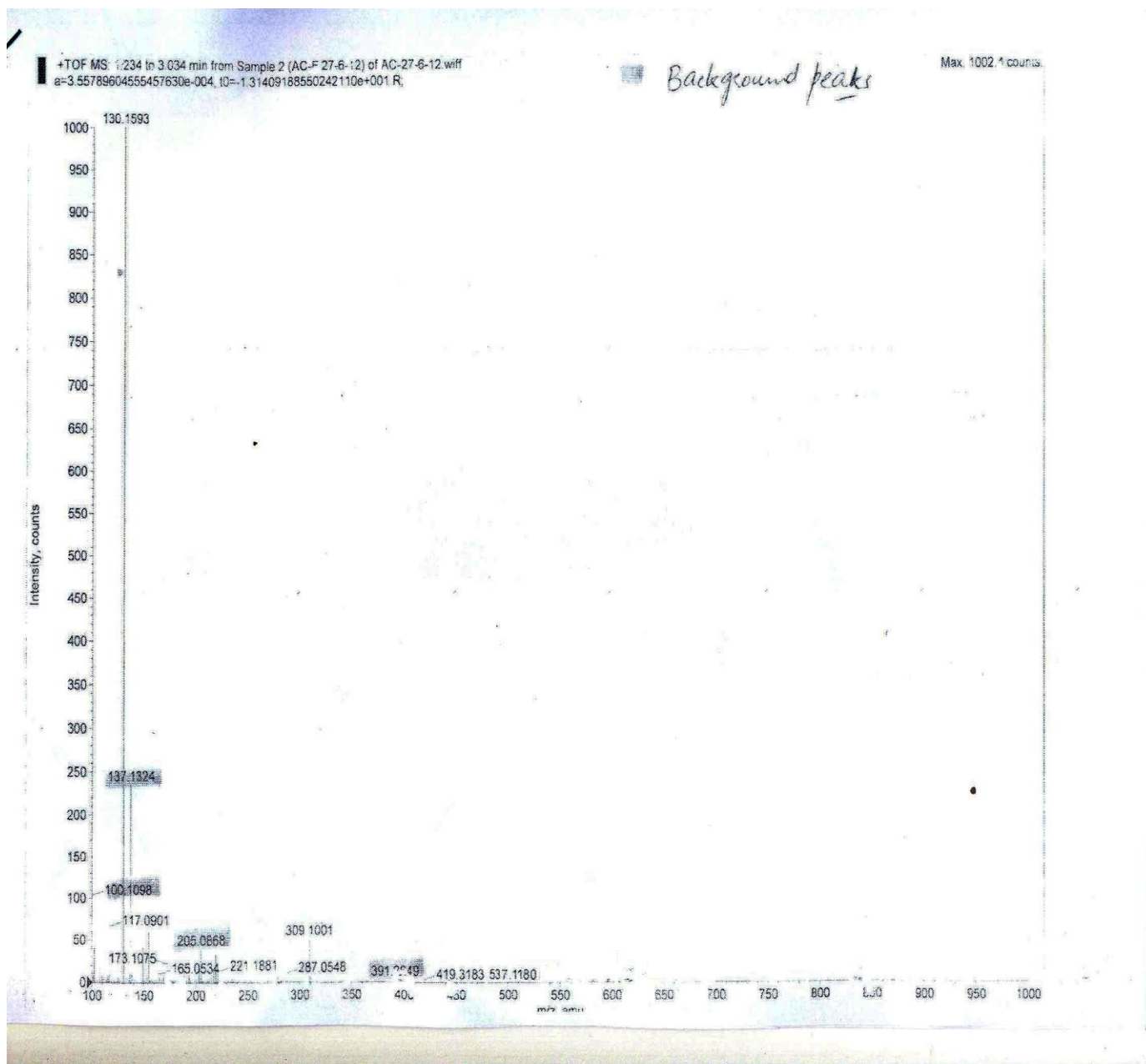
S2: ¹³C-NMR (75 MHz, CDCl₃) Spectrum of Compound **1** (trichonoide A)



S3: HMBC Spectrum of Compound **1** (trichonoide A)



S4: Extended-HMBC Spectrum of Compound **1** (trichonoide A)



S5: HRESIMS Spectrum of Compound 1 (trichonoide A)

Elemental composition calculator

Target m/z: +537.1180 amu

Tolerance: +10.0000 ppm

Result type: Elemental

Max num of results: 100

Min DBE: -0.5000 Max DBE: +50.0000

Electron state: OddAndEven

Num of charges: 0

Add water: N/A

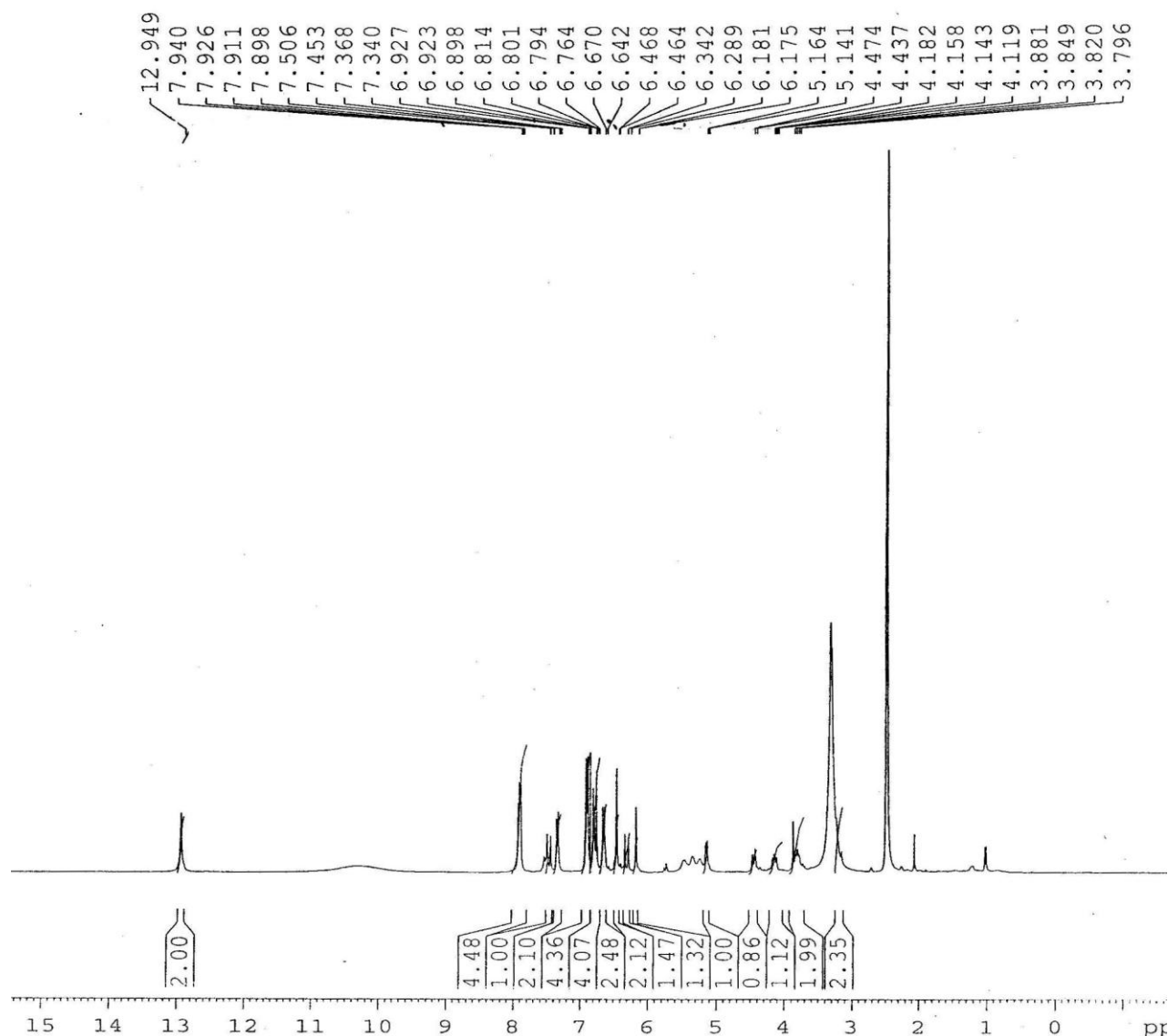
Add proton: N/A

File Name: XXXXXXXXXXXXXX

	Elements	Min Number	Max Number
1	H	0	30
2	O	0	10
3	C	0	35

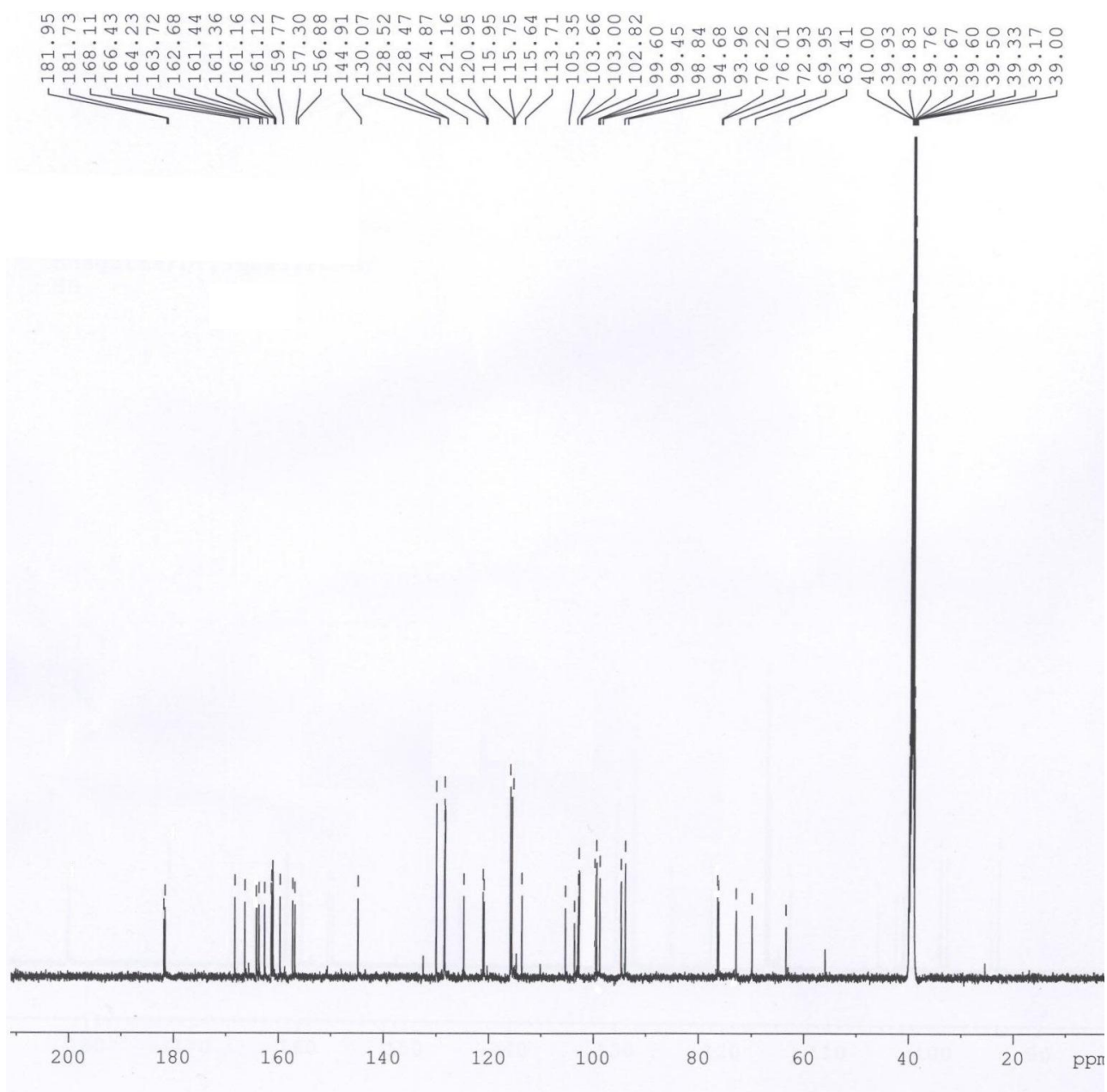
	Formula	Calculated m/z (amu)	mDa Error	PPM Error	DBE
1	C31 H21 O9	537.1185	-0.5575	-1.0381	21.5

S5: HRESIMS Spectrum of Compound 1 (trichonoide A)

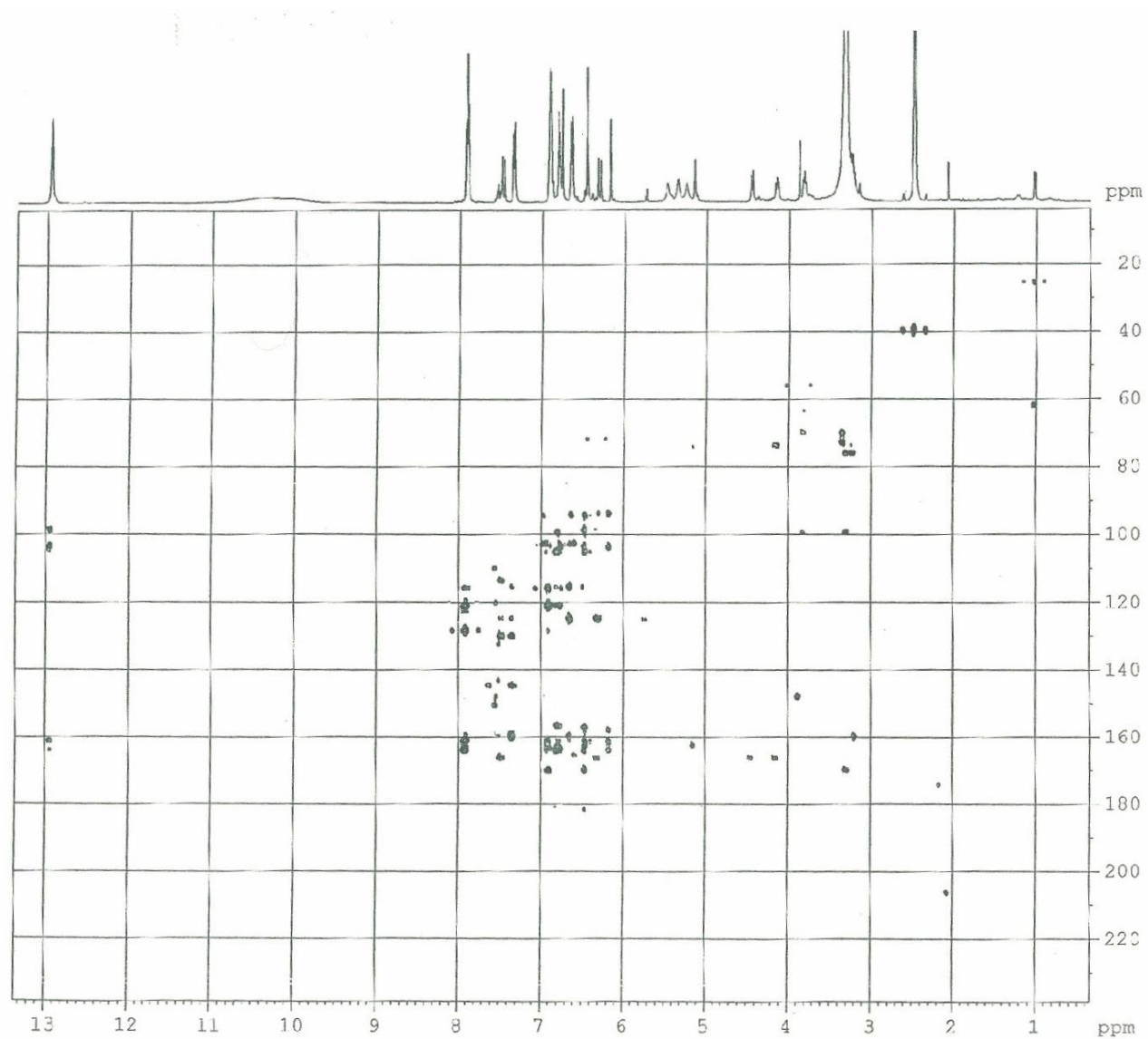


S6: ^1H -NMR (300 MHz, DMSO) Spectrum of Compound **2** (trichonoide B)

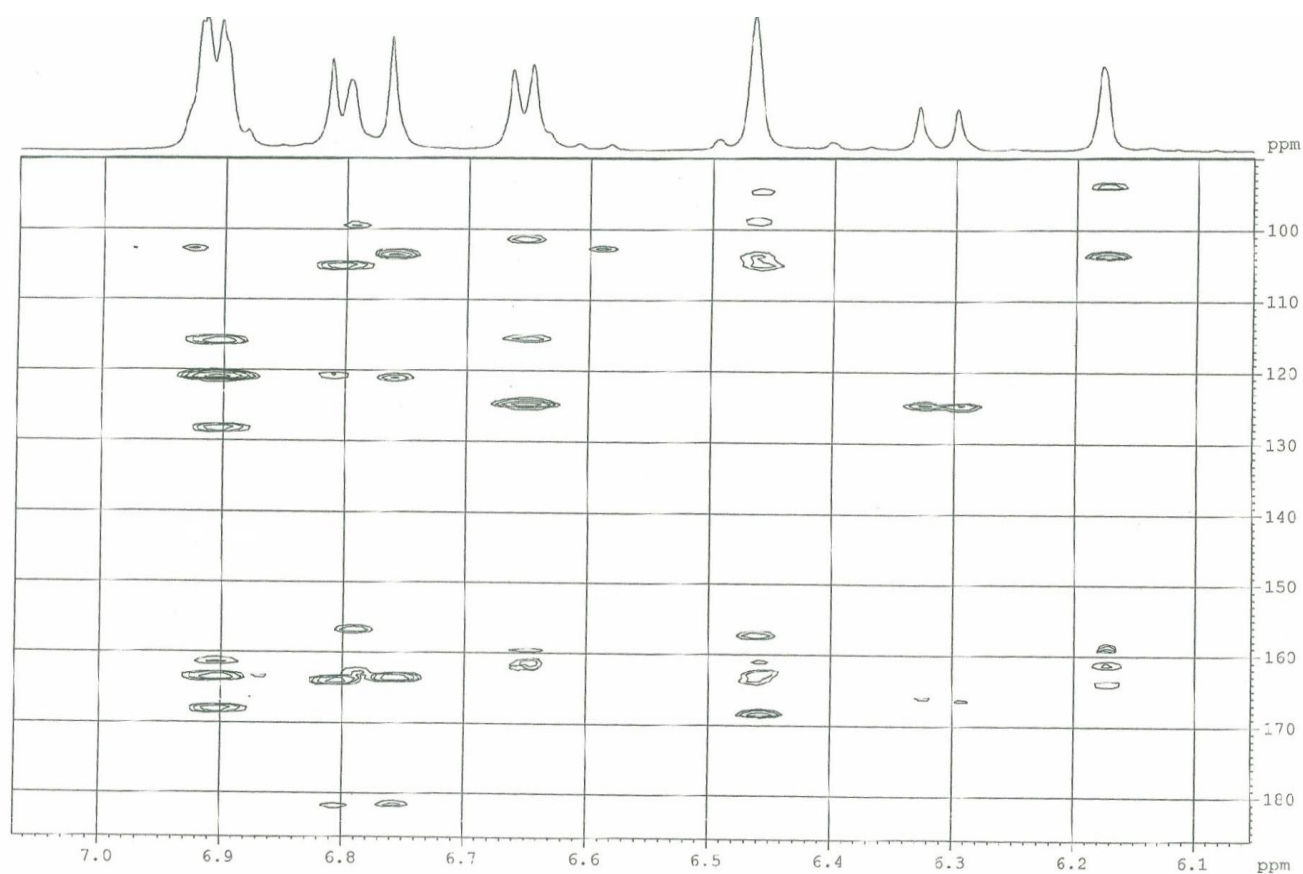
Trichonoide (2): Yellow solid. IR (ν_{max} cm^{-1}): 3385 (OH), 1755 (ester moiety), 1678 (conjugated carbonyl), 1642 and 1485 cm^{-1} (aromatic moiety). UV (λ_{max} nm) 269 and 345 nm. ^1H -NMR (DMSO, 300 MHz), δ : 6.81 (1H, s, H-3), 6.18 (1H, d, $J = 1.8$, H-6), 6.46 (1H, d, overlapped, H-8), 7.94, (2H, d, overlapped, H-2', -6'), 6.92 (2H, d, overlapped, H-3', -5'), 5.16 (1H, d, $J = 7.3$, H-1''), 3.20 (1H, m, H-2''), 3.53 (1H, m, H-3''), 3.15 (1H, m, 4''), 3.87 (1H, m, H-5''), 4.18 (1H_a, m, H-6''), 4.463 (1H_b, m, 6''), 6.67 (2H, d, $J = 8.4$, H-2''', -6'''), 7.36 (2H, d, $J = 8.4$, H-3''', -5'''), 7.50 (1H, d, $J = 15.9$ H-7'''), 6.34 (1H, d, $J = 15.9$, H-8'''), 6.76 (1H, s, H-3'''), 6.47 (1H, d, overlapped, H-6'''), 6.90 (1H, d, $J = 2.1$, H-8'''), 7.92, (2H, d, overlapped, H-2''', -6'''), 6.89 (2H, d, overlapped, H-3''', -5'''). ^{13}C -NMR (DMSO, 150 MHz), δ : 161.4 (C-2), 103.0 (C-3), 181.9 (C-4), 105.3 (C-4a), 159.7 (C-5), 98.8 (C-6), 164.2 (C-7), 93.9 (C-8), 157.3 (C-8a), 121.2 (C-1'), 128.5 (C-2', -6'), 115.8 (C-3', -5'), 162.6 (C-4'), 99.6 (C-1''), 76.0 (C-2''), 72.9 (C-3''), 69.9 (C-4''), 76.2 (C-5''), 63.4 (C-6''), 161.1 (C-1'''), 130.0 (C-2''', -6'''), 115.8 (C-3''', -5'''), 124.8 (C-4'''), 144.9 (C-7'''), 113.7 (C-8'''), 166.4 (C-9'''), 163.7 (C-2'''), 102.8 (C-3'''), 181.7 (C-4'''), 103.6 (C-4a'''), 161.2 (C-5'''), 99.4 (C-6'''), 168.1 (C-7'''), 94.6 (C-8'''), 156.9 (C-8a'''), 120.9 (C-1'''), 128.5 (C-2''', -6'''), 115.7 (C-3''', 5'''), 161.3 (C-4'''). HR-ESI-MS: $m/z = 831.1920$ $[\text{M}+\text{H}]^+$ for formula $\text{C}_{45}\text{H}_{35}\text{O}_{16}$.



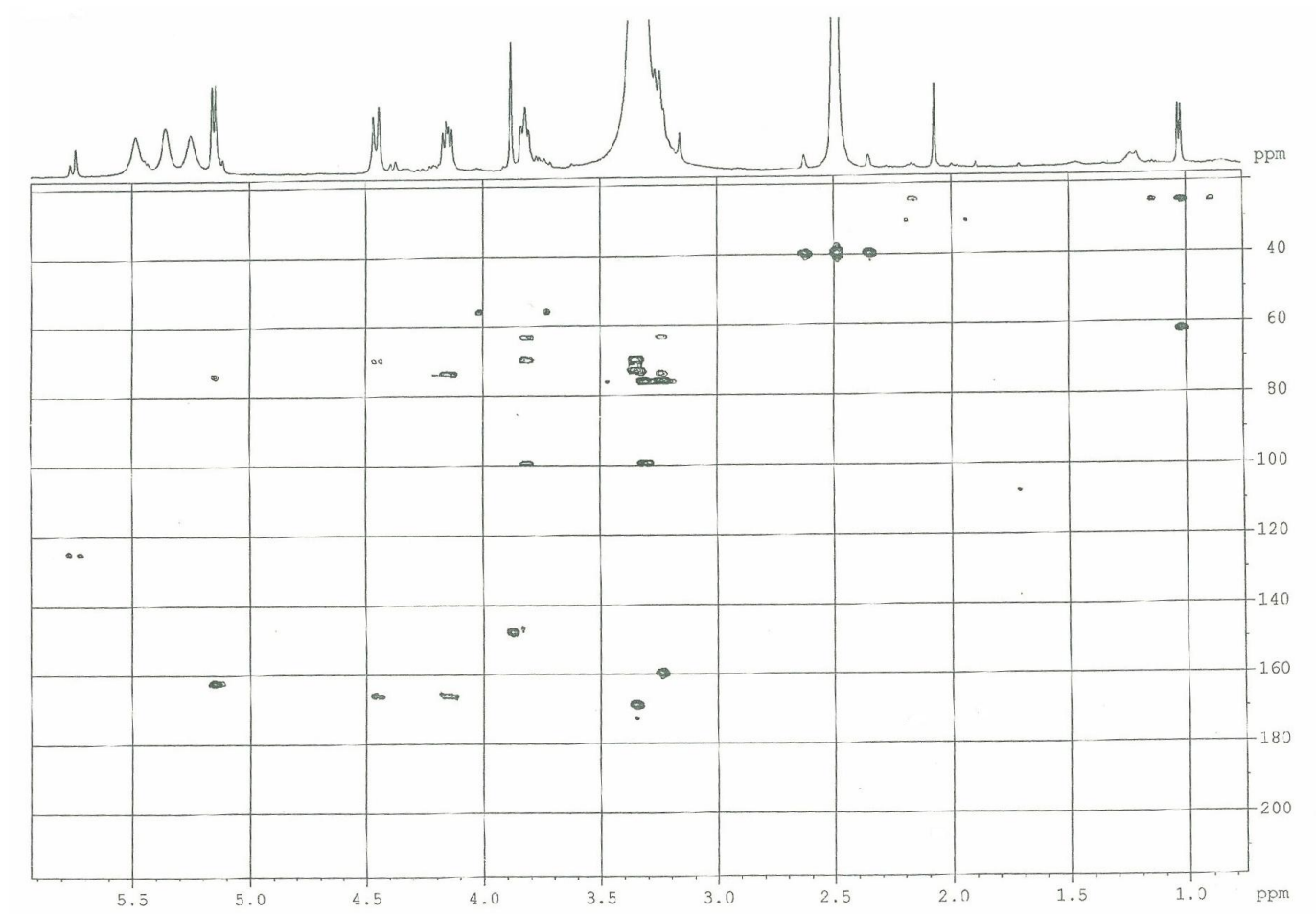
S7: ^{13}C NMR (125 MHz, DMSO) Spectrum of Compound **2** (trichonoide B)



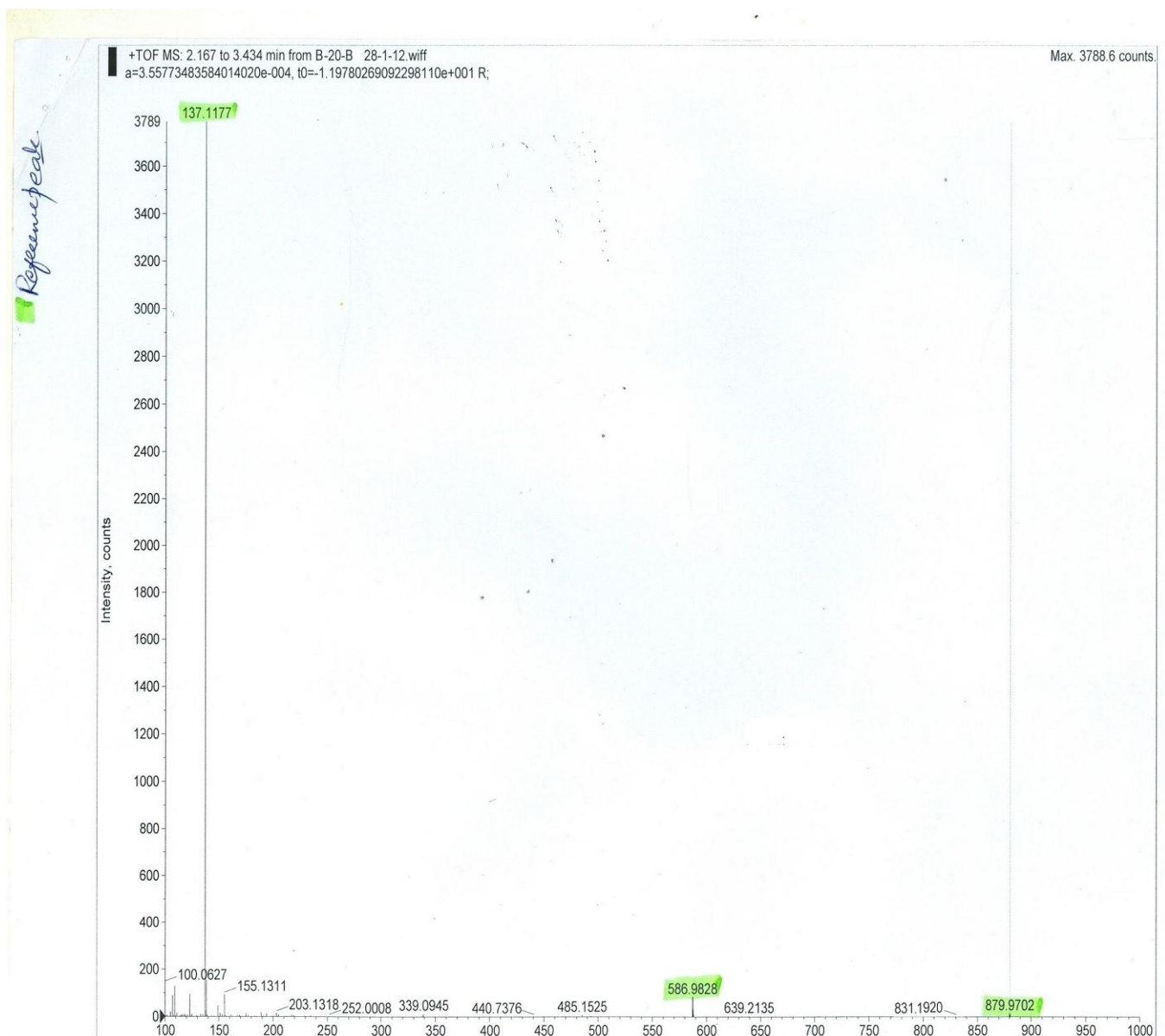
S8: HMBC Spectrum of Compound 2 (trichonoide B)



S9: Extended-HMBC Spectrum of Compound 2 (trichonoide B)



S9: Extended-HMBC Spectrum of Compound **2** (trichonoide B)



S10: HRESIMS Spectrum of Compound **2** (trichonoide B)

Elemental composition calculator

Target m/z: +831.1920 amu

Tolerance: +10.0000 ppm

Result type: Elemental

Max num of results: 100

Min DBE: -0.5000 Max DBE: +50.0000

Electron state: OddAndEven

Num of charges: 0

Add water: N/A

Add proton: N/A

File Name:

	Elements	Min Number	Max Number
1	H	0	40
2	O	0	20
3	C	0	50

	Formula	Calculated m/z (amu)	mDa Error	PPM Error	DBE
1	C45 H35 O16	831.1925	-0.5106	-0.6143	28.5

S10: HRESIMS Spectrum of Compound 2 (trichonoide B)