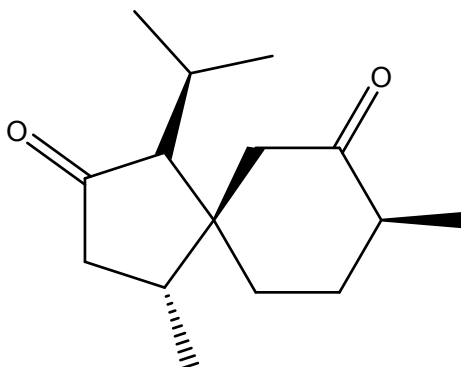


**Table S1: Voucher specimen numbers of collected plant material submitted at KASH Herbarium, Centre for Biodiversity and Taxonomy, University of Kashmir, India**

<b>S. No.</b>	<b>Sample collection site</b>	<b>Vocher No.</b>
01	Kawoosa	2438
02	Najan	2719
03	Palhalan Pattan	2720
04	Singpore Ganastan	2721
05	Yarikha Tangmarg	2722
06	Kakpora Pulwama	2723
07	Aarath Budgam	3071
08	Soibugh Budgam	3072
09	Wadwan ( cultivated)	3073
10	Tuli Rajouri	3074
11	Gotapora	3075
12	IIIM Srinagar	3076
13	Mattan Anantnag	3077
14	Zabigul Beerwah	3078

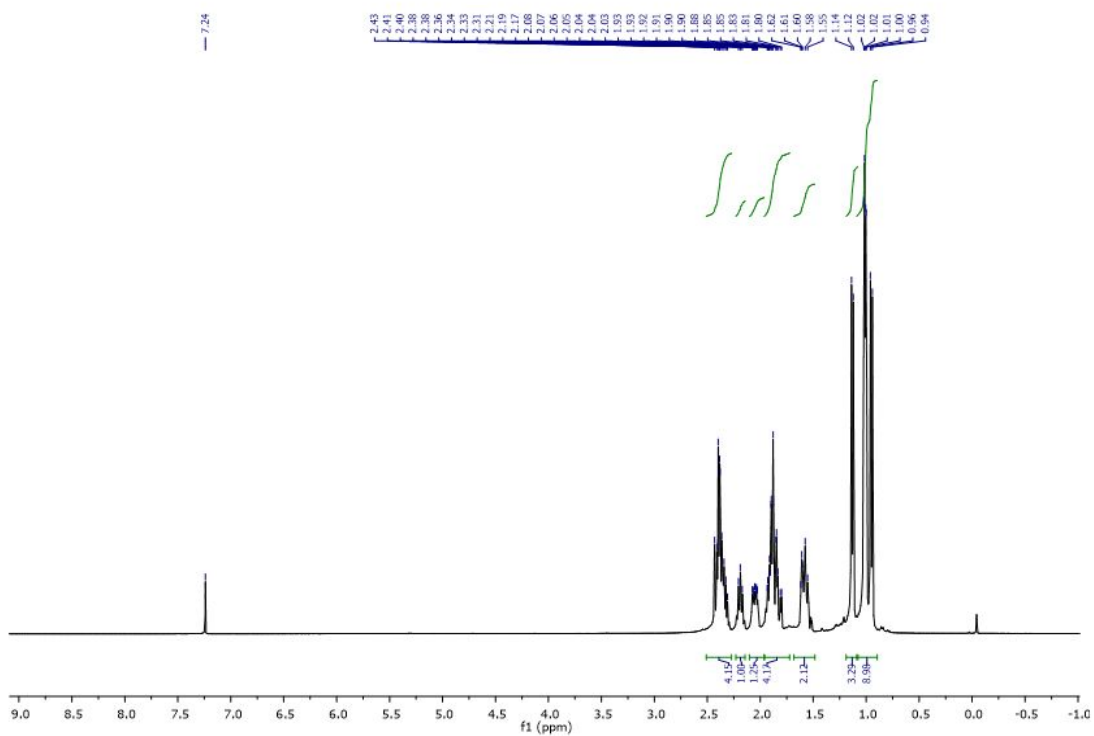
**Compound 1 (ACS01): Isoacorone**

**IUPAC Name:** (1S, 8S)-1-Isopropyl-4,8-dimethylspiro [4.5] decan-2,7-dione

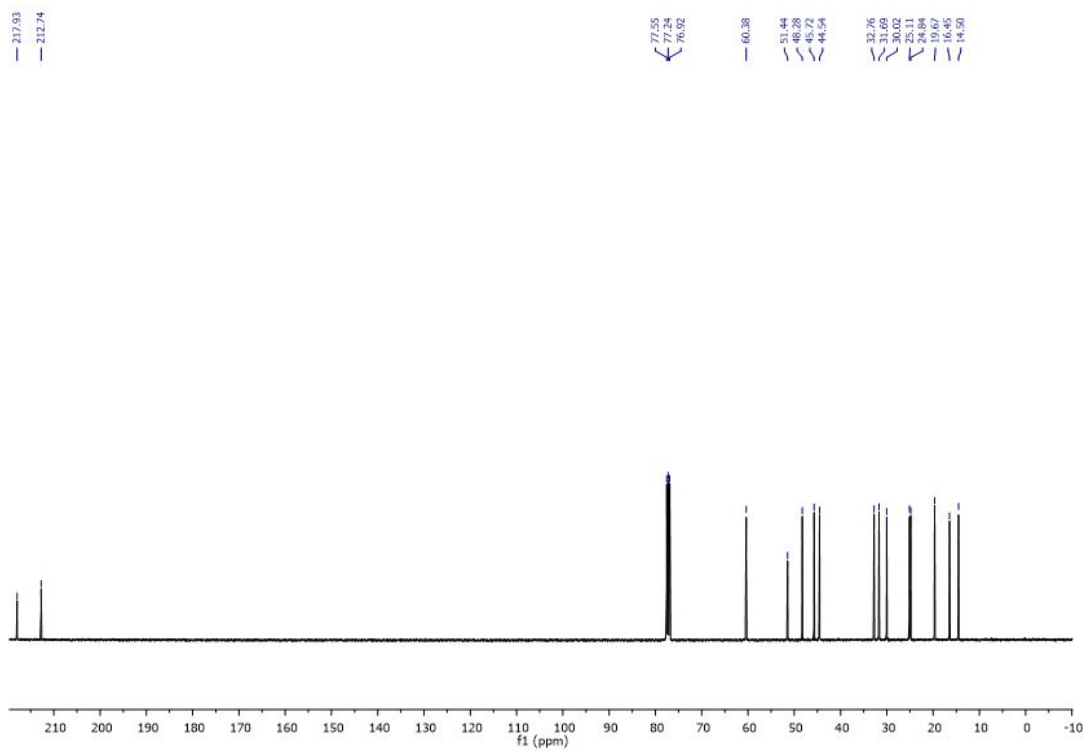


Isoacorone (ACS01)

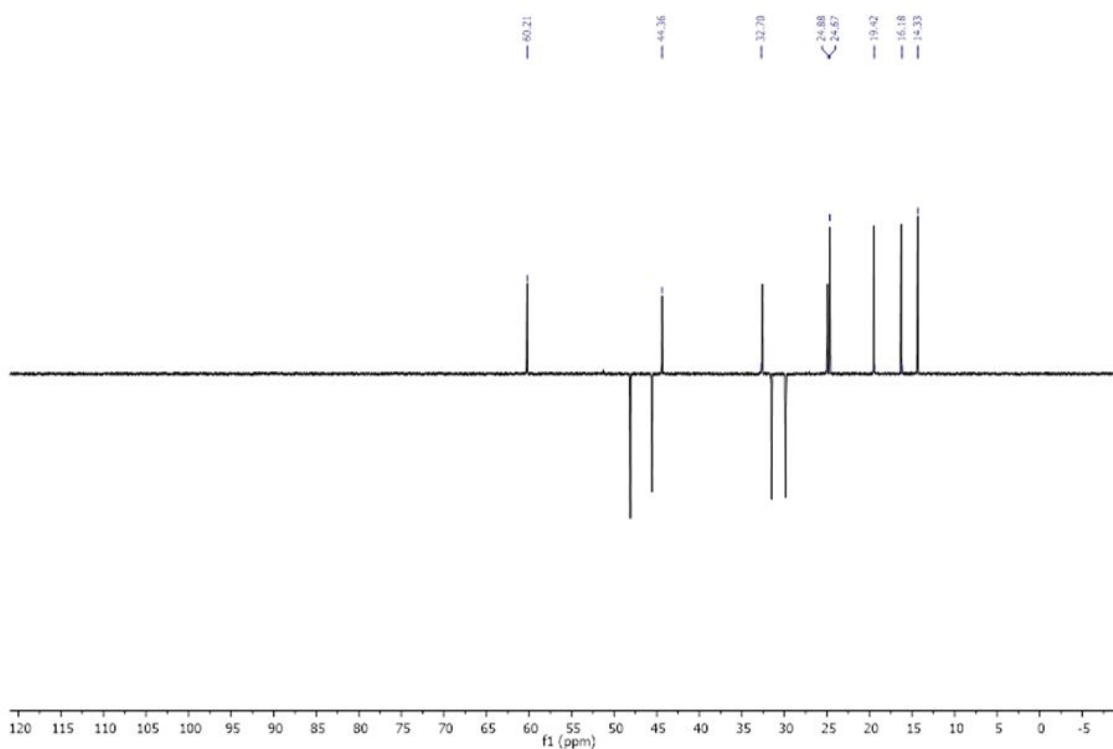
<b>Chemical formula</b>	:	$C_{15}H_{24}O_2$
<b>Exact mass</b>	:	236.18
<b>Molecular weight</b>	:	236.35 g/m
<b>Melting point</b>	:	96-97 °C
<b>Physical state</b>	:	Crystals
<b>Specific rotation</b>	:	-00.003 to -00.006
<b>M/z</b>	:	236.18 (100%), 237.18 (16.6%), 238.18 (1.6%)
<b>Elemental analysis</b>	:	C: 76.23, H: 10.24, O: 13.54



**<sup>1</sup>H NMR spectrum of Isoacorone**



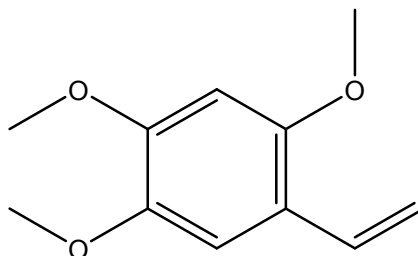
**<sup>13</sup>C NMR spectrum of Isoacorone**



### DEPT spectrum of Isoacorone

**Compound 2 (ACS02):** Cis-asarone ( $\beta$ -asarone)

**IUPAC Name** : 1, 2, 4 trimethoxy -5-(prop -1-en-1-yl) benzene



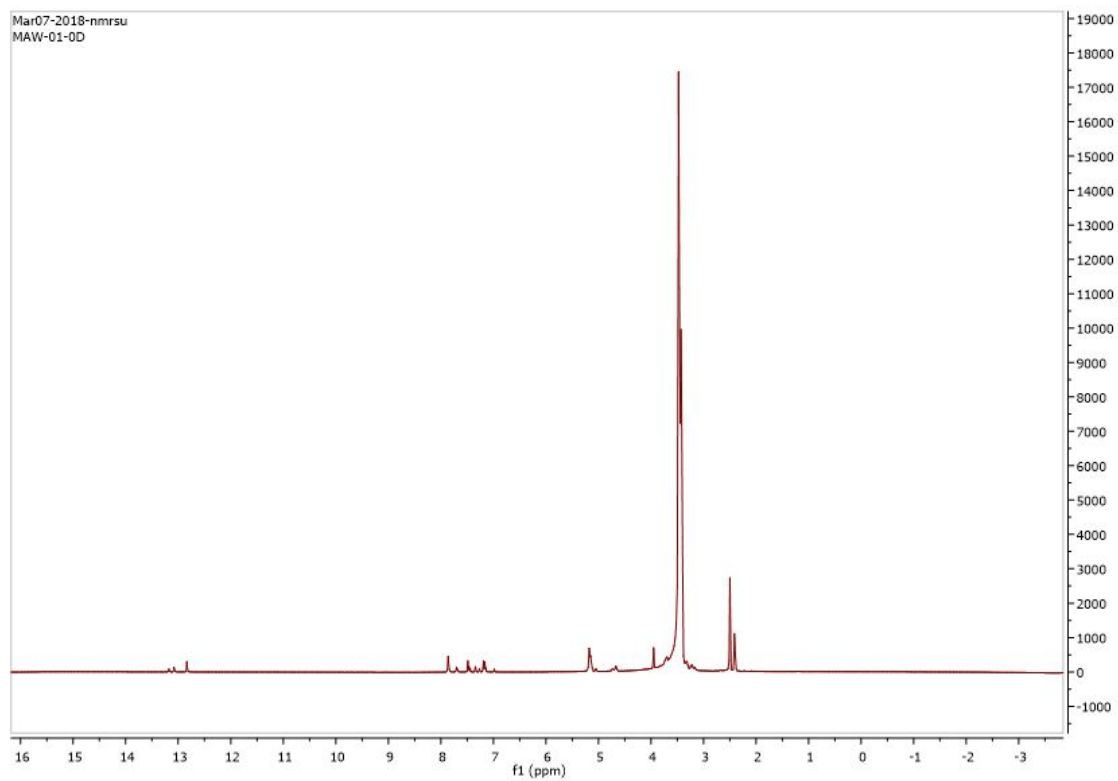
Beta-asarone (ACS02)

**Molecular Formula** :  $C_{12}H_{16}O_3$

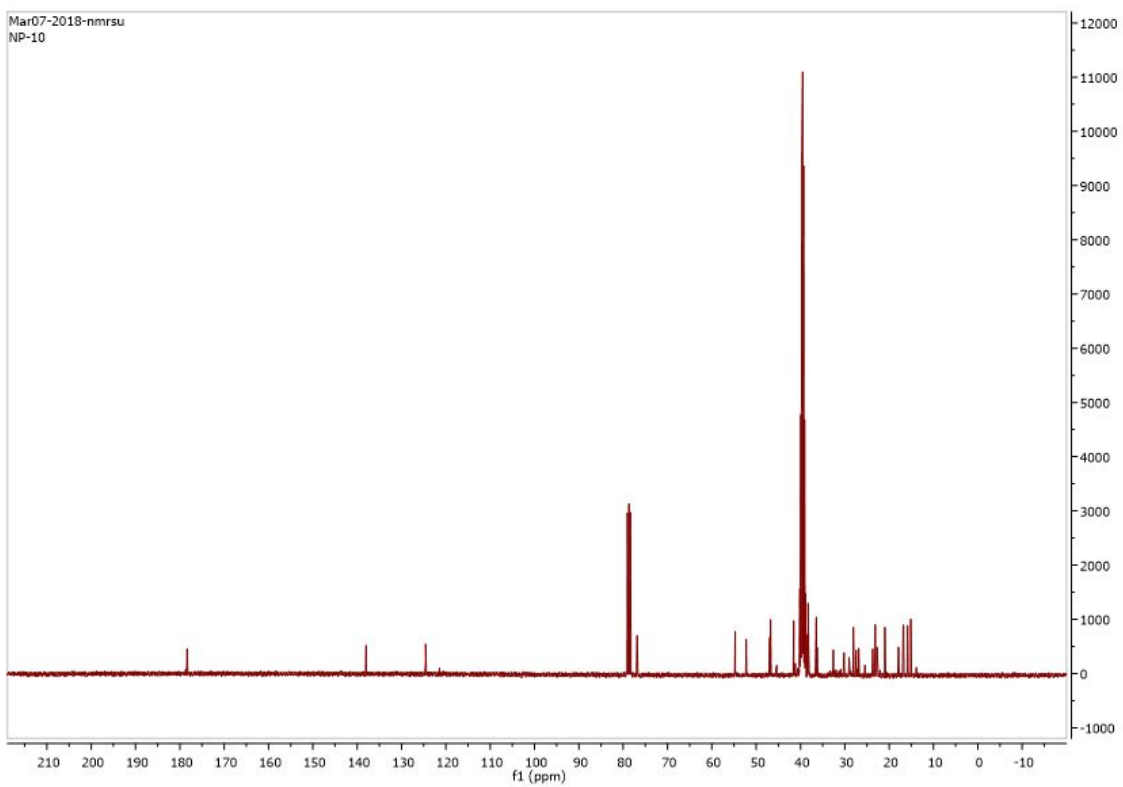
**Molecular weight** : 208.257 g/m

**Colour** : Yellowish oily liquid

**Melting point** : 62.00 to 63.00 °C.



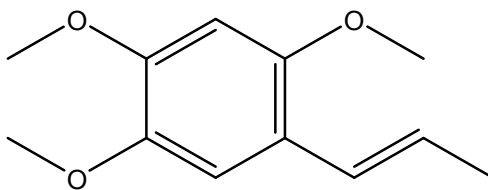
**<sup>1</sup>H NMR spectrum of  $\beta$ -Asarone**



**$^{13}\text{C}$  NMR spectrum of Beta asarone**

**Compound 3 (ACS03): Trans-asarone ( $\alpha$  - asarone)**

**IUPAC Name** : Trans-1, 2, 4-trimethoxy -5(2-propenyl) benzene



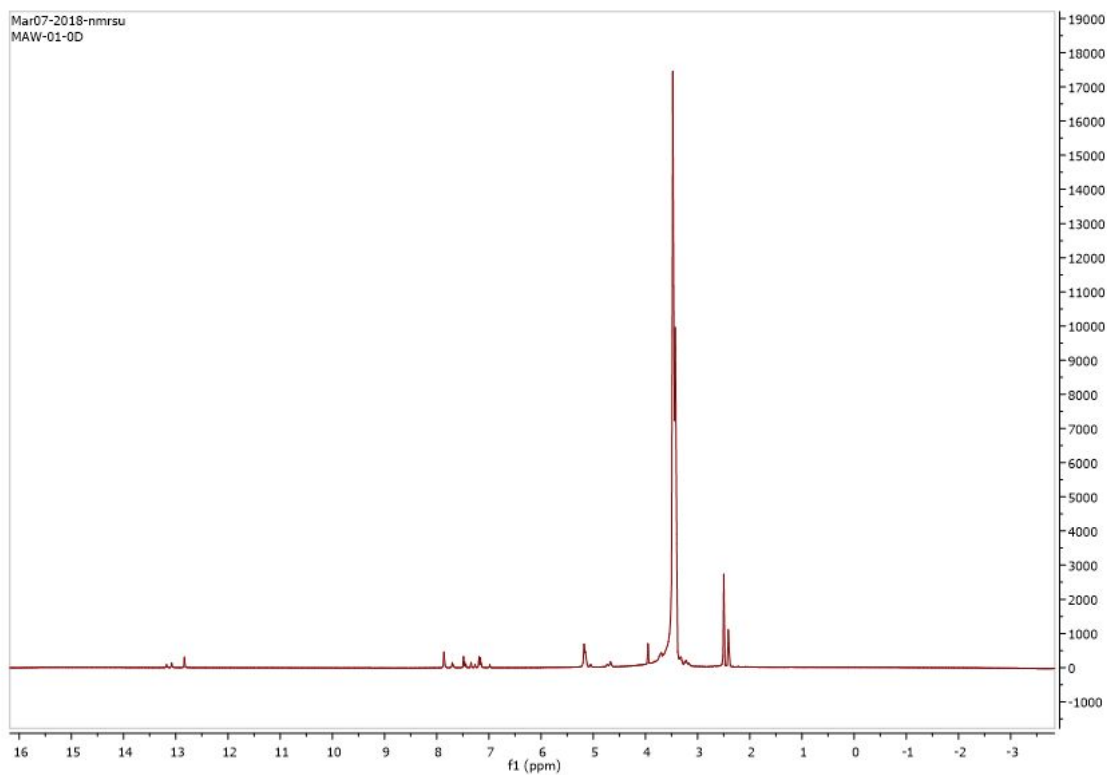
Trans-asarone (ACS03)

**Molecular Formula** : C<sub>12</sub>H<sub>16</sub>O<sub>3</sub>

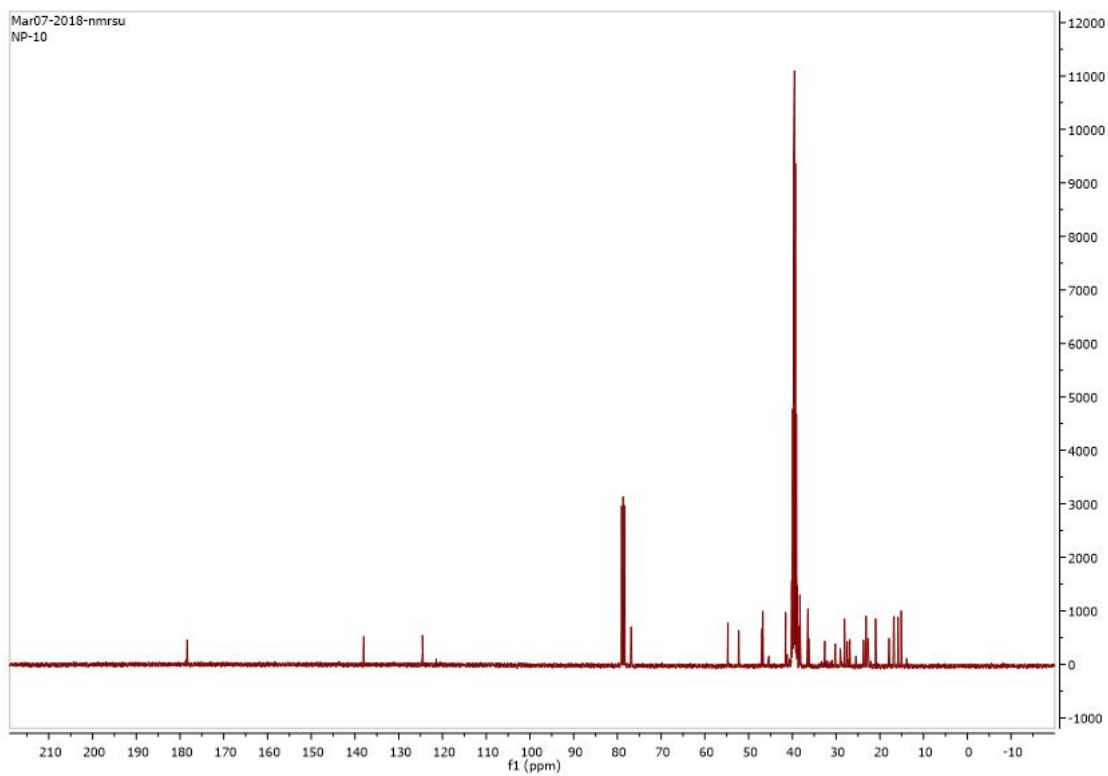
**Molecular Weight** : 208.25

**Colour** : Colourless crystalline solid

**Melting point** : 62 °C



**<sup>1</sup>H NMR Spectrum of  $\alpha$ - asarone**

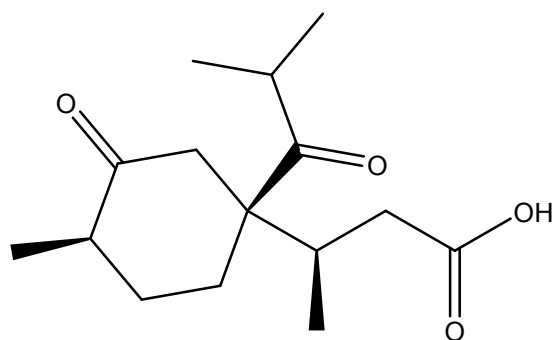


$^{13}\text{C}$  NMR Spectrum of  $\alpha$ -asarone



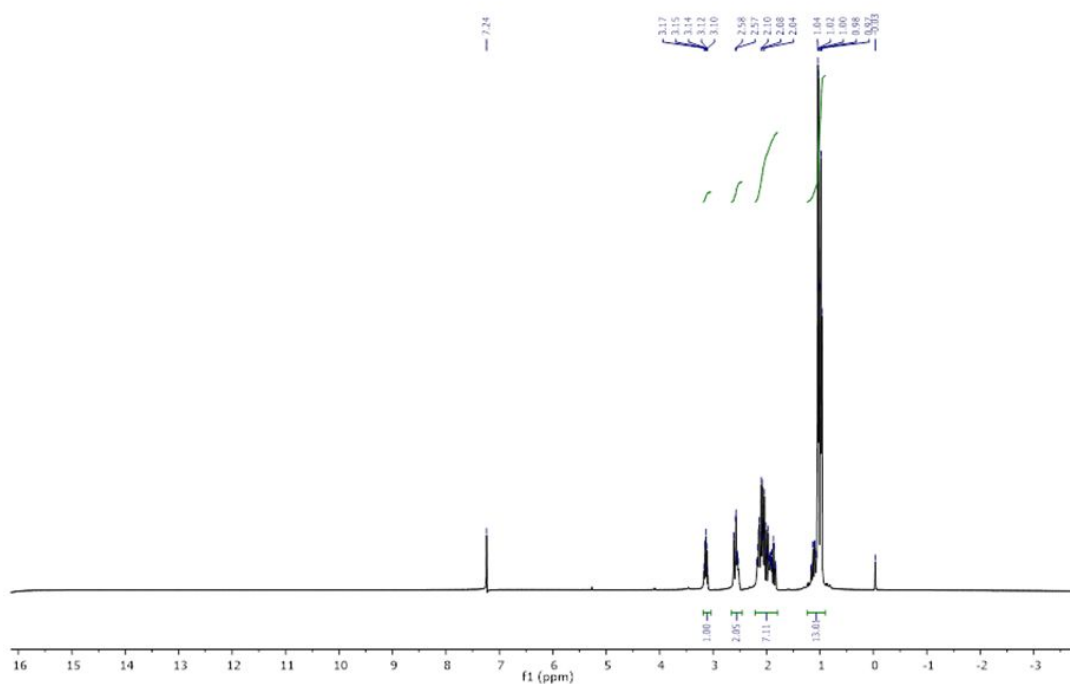
**Compound 4 (ACS04): Acoric acid**

**IUPAC Name** : (S)-3[(1S, 4R)-1-isobutyryl-4-methyl-3-oxocyclohexyl] butanoic Acid

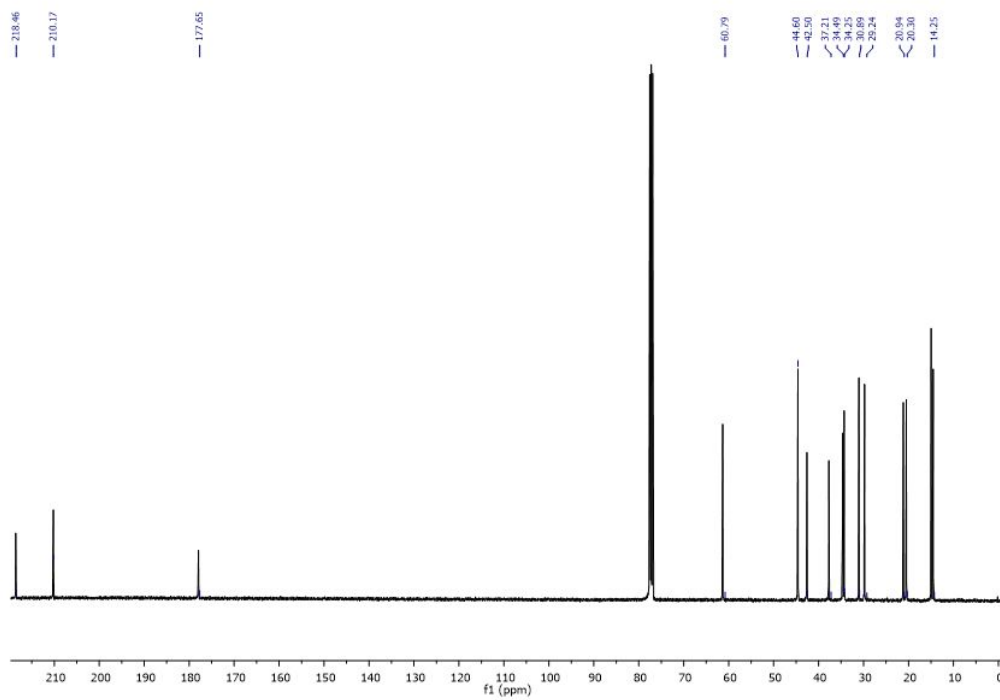


Acoric acid (ACS04)

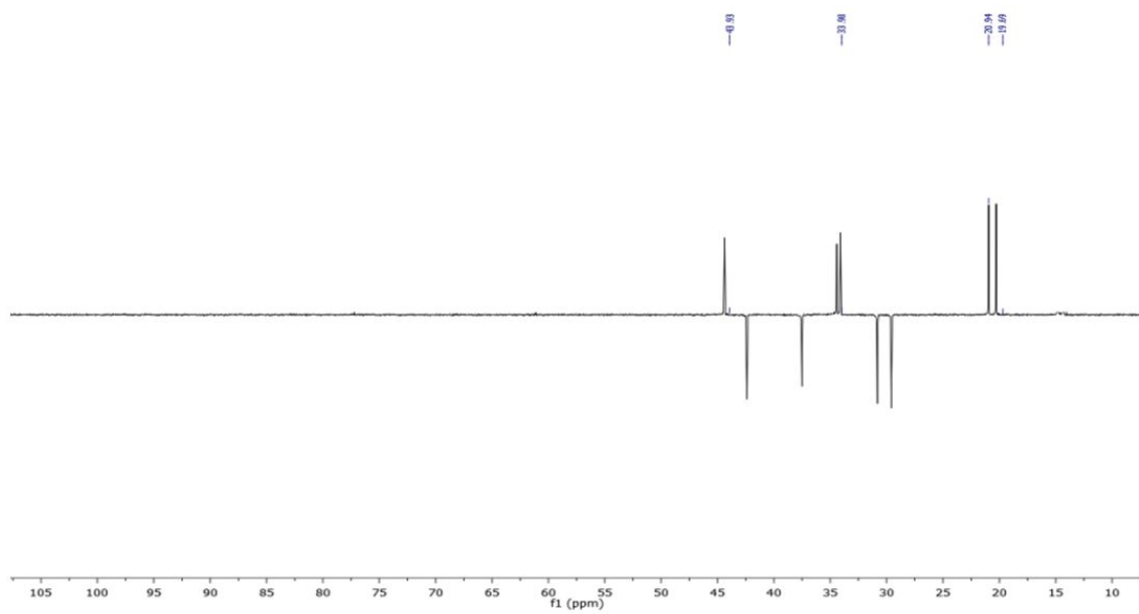
**Molecular formula** :  $C_{15}H_{24}O_4$   
**Molecular weight** : 268.353g/m  
**Exact Mass** : 268.17g/m  
**M/z** : 268.17(100%), 269.17(16.7%), 270.17(2.1%)  
**Elemental analysis** : C: 67.14, H: 9.01, O: 23.85  
**Melting point** : 166 °C - 168 °C  
**Physical state** : crystalline



**<sup>1</sup>H NMR Spectrum of Acoric acid**



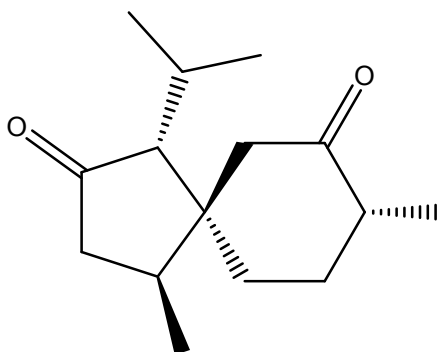
**<sup>13</sup>C NMR Spectrum of Acoric acid**



**DEPT Spectrum of Acoric acid**

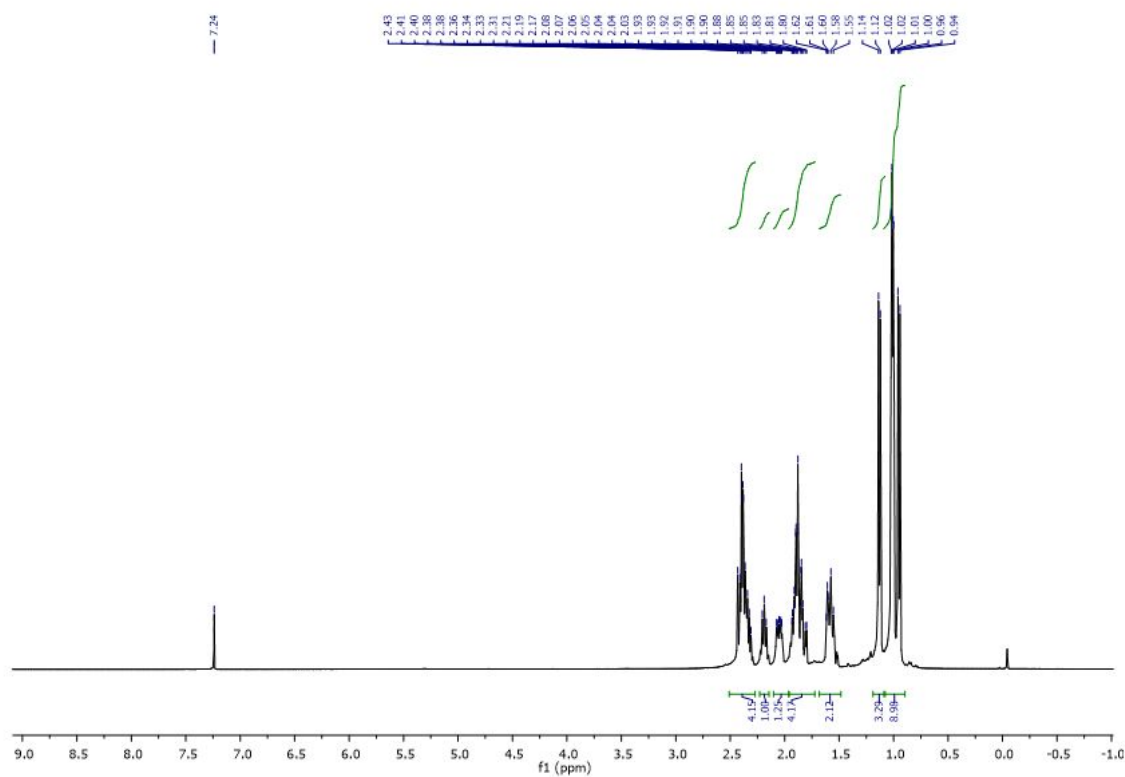
### Compound 5 (ACS05): Acorone

IUPAC Name: 1-Isopropyl-4, 8- dimethylspiro [4.5] decan-2, 7-dione

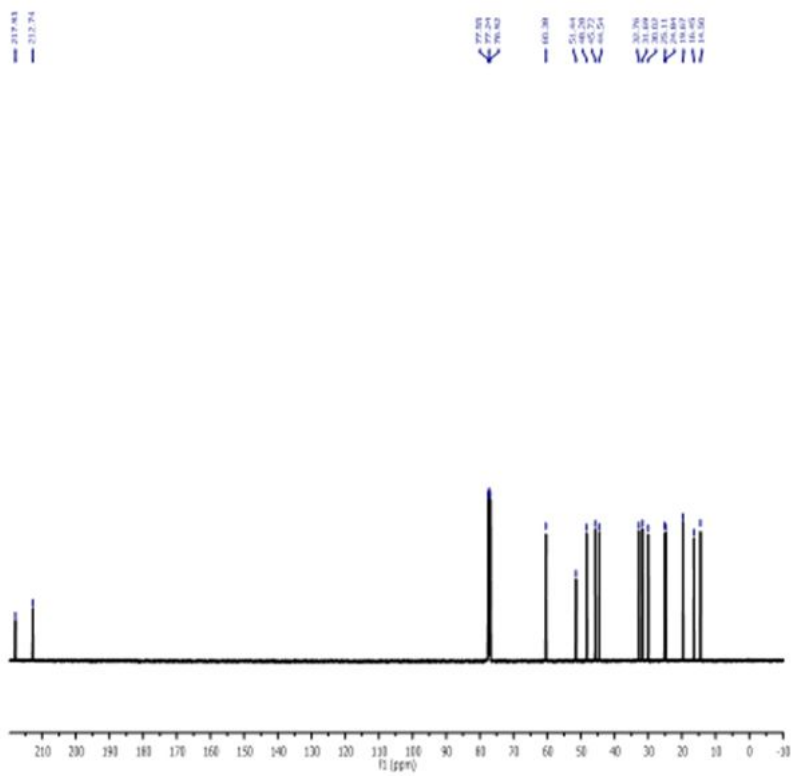


Acorone (ACS05)

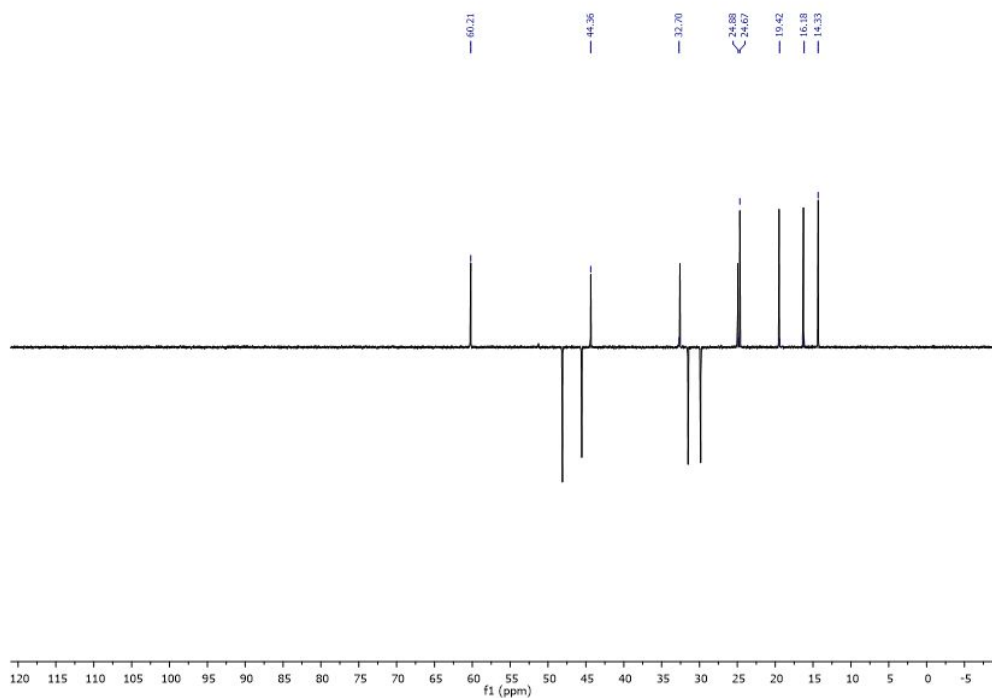
<b>Chemical formula</b>	:	$C_{15}H_{24}O_2$
<b>Molecular weight</b>	:	236 g/m
<b>Melting point</b>	:	100-101 °C
<b>Physical state</b>	:	Crystalline
<b>Specific rotation</b>	:	+00.006 to 0.008



<sup>1</sup>H NMR Spectrum of Acorone



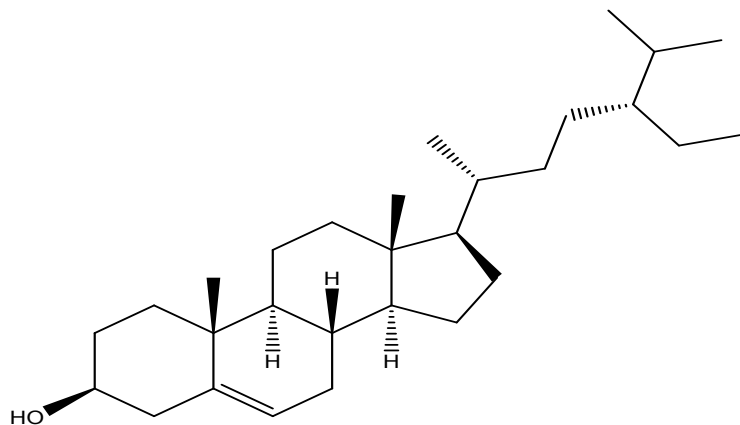
**<sup>13</sup>C NMR spectral data of compound Acorone**



**DEPT Spectral data of Acorone**

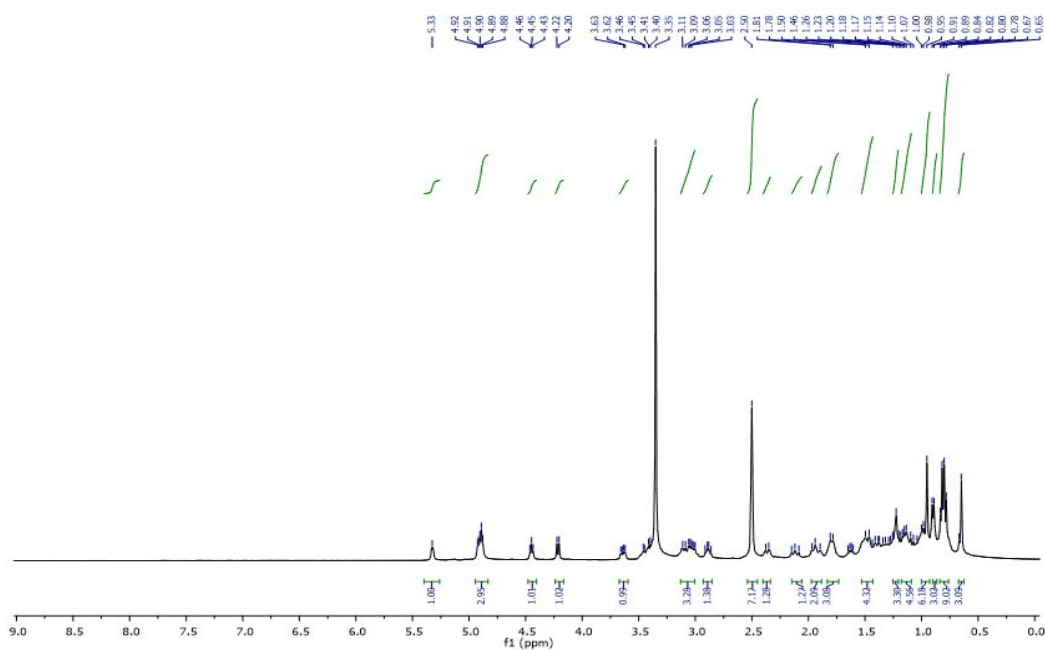
## Compound 6 (ACS6): Beta sitosterol glycoside

**IUPAC Name** : 2-((17-(5, 6-dimethylheptan-2-yl)-10, 13-dimethyl-2,3,4,7,8,9,10,11,12,13,14,15,16,17-tetradecahydro-1H-cyclopenta[a] phenanthren-3- yl)oxy) -6- (hydroxymethyl) tetrahydro-2H-pyran-3, 4, 5-triol

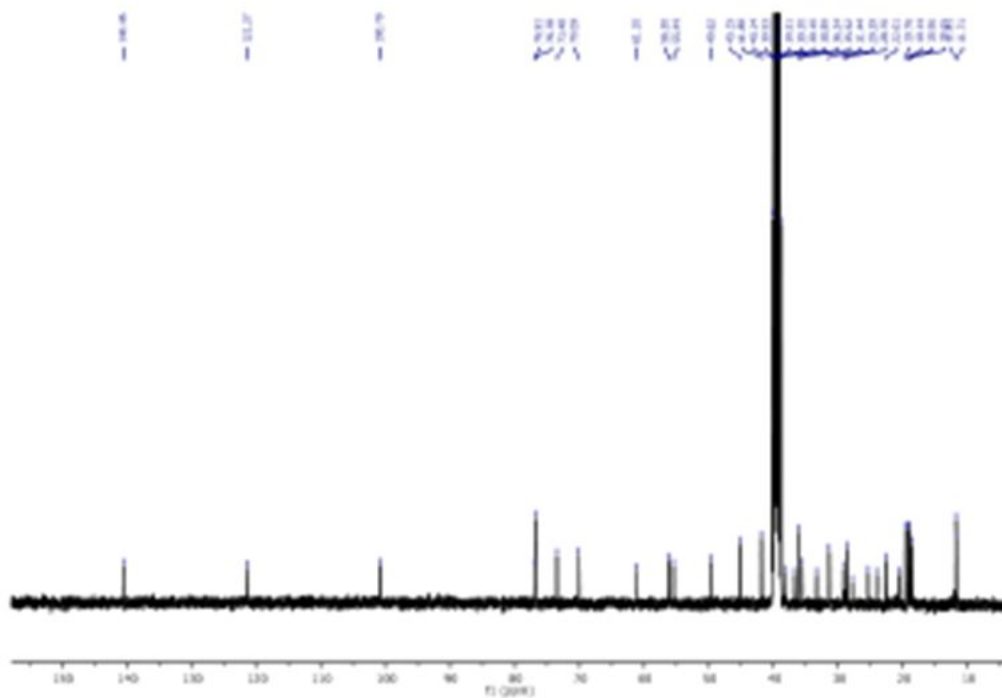


Beta sitosterol (ACS06)

**Chemical formula** :  $C_{34}H_{58}O_6$   
**Molecular weight** : 562.82g/m  
**Exact Mass** : 562.42  
**Physical state** : Solid white powder  
**M/z** : 562.42(100%), 563.43(37.7%), 564.43(8.1%), 565.43(1.2%)  
**Elemental analysis** : C: 72.56, H: 10.39, O: 17.06

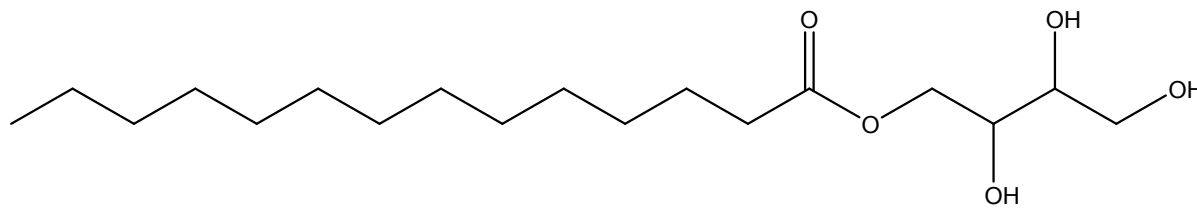


<sup>1</sup>H NMR Spectrum of compound β-Sitosterol glycoside



**Compound 7 (ACS07): 2, 3, 4-Trihydroxybutyl tetradecanoate**

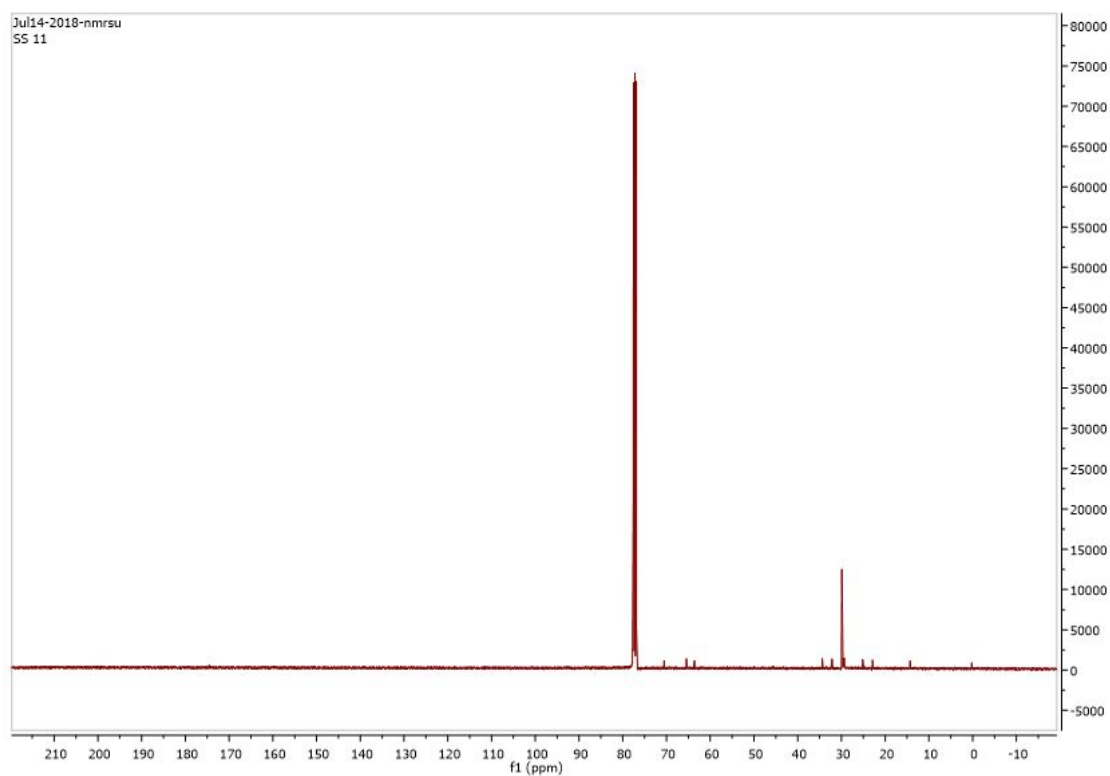
**IUPAC Name** : 2, 3, 4-Trihydroxybutyl tetradecanoate



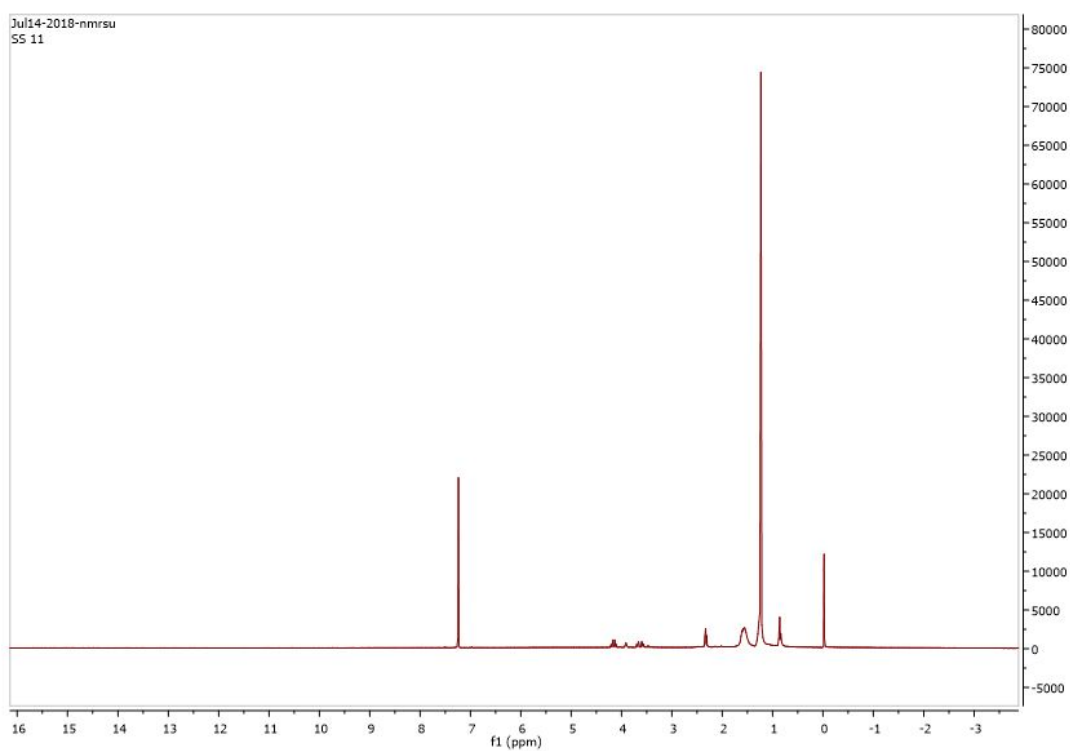
2, 3, 4-Trihydroxybutyl tetradecanoate (ACS07)

**Chemical formula** :  $C_{18}H_{36}O_5$

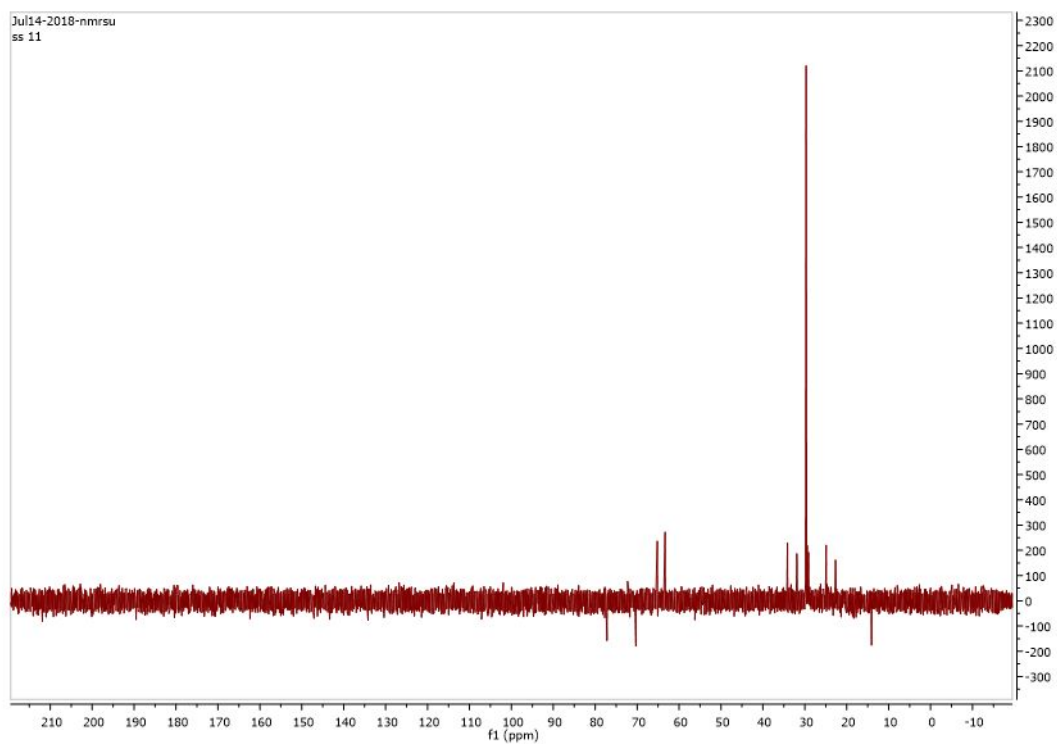
**Molecular weight** : 332.47 g/m



**$^1H$  NMR Spectrum of compound 2, 3, 4-Trihydroxybutyl tetradecanoate**



**$^{13}$  C NMR Spectrum of compound 2, 3, 4-Trihydroxybutyl tetradecanoate**

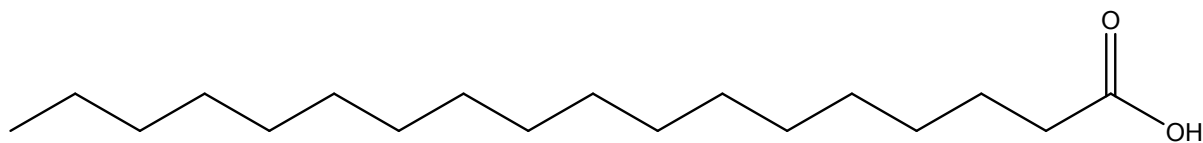


**DEPT Spectrum of compound 2, 3, 4-Trihydroxybutyl tetradecanoate**



**Compound 8 (ACS08): Stearic acid**

**IUPAC Name** : Octadenoic acid



Stearic acid (ACS08)

**Chemical formula** :  $C_{18}H_{36}O_2$  ,  $CH_3(CH_2)_{16}COOH$

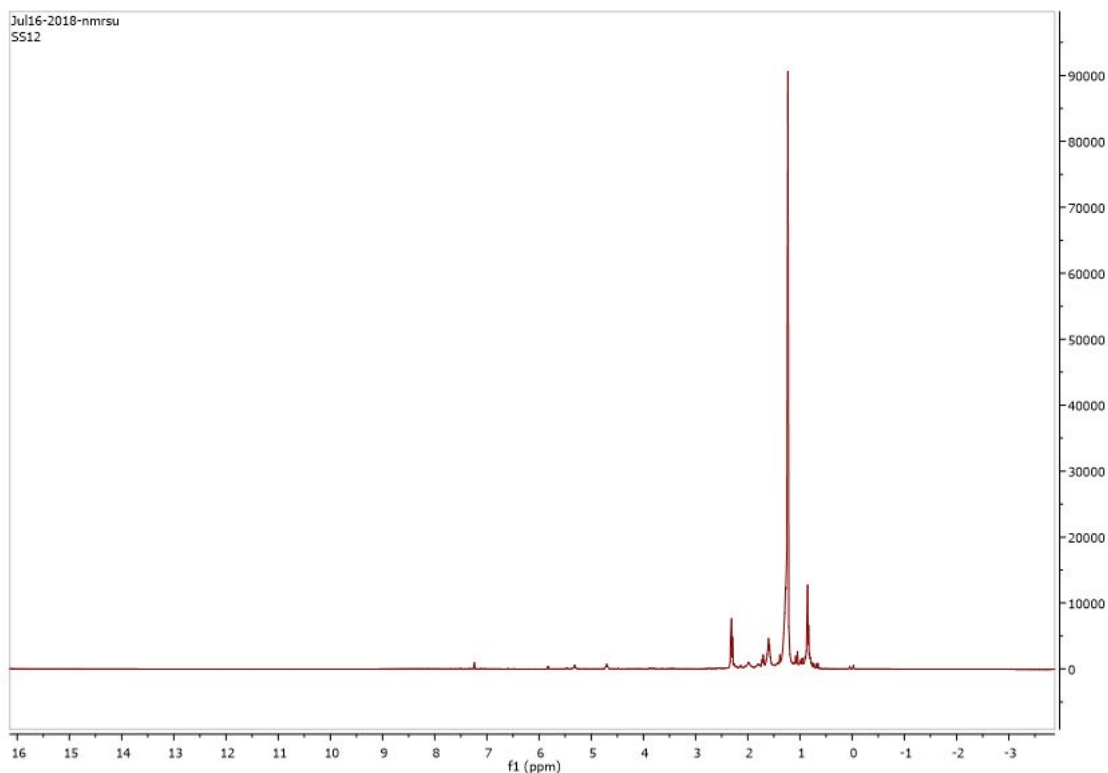
**Molecular weight** : 284.48 g/mol

**Melting point** : 69.3%

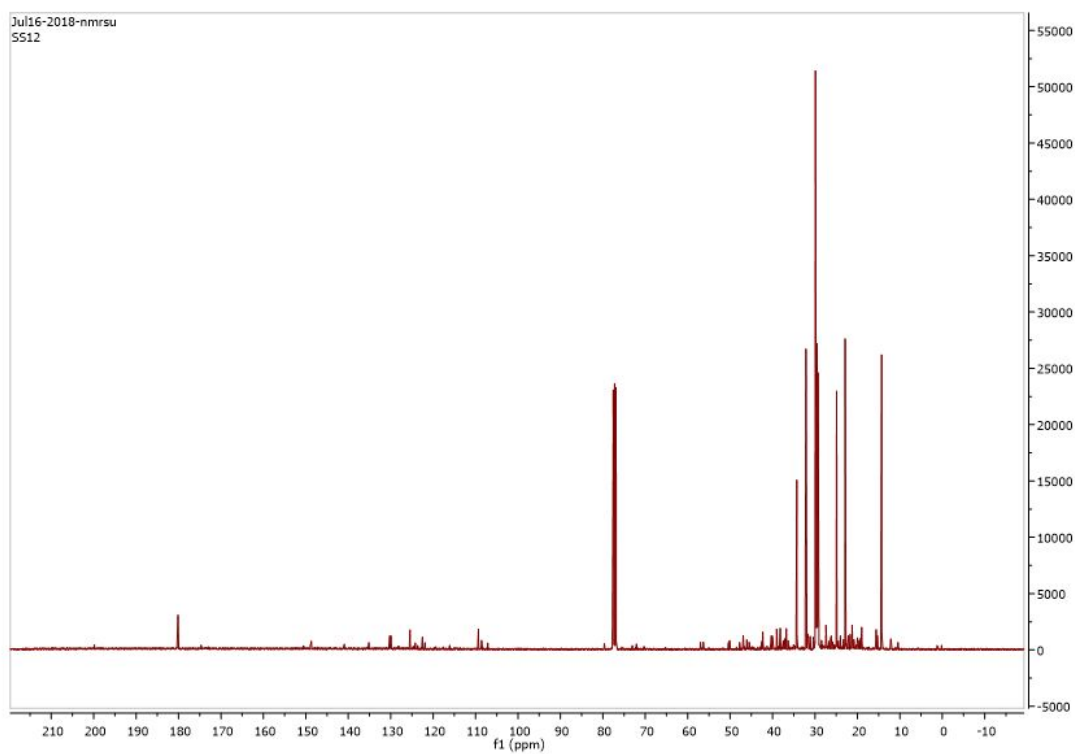
**Boiling point** : 361<sup>o</sup>C

**Physical state** : Saturated fatty acid

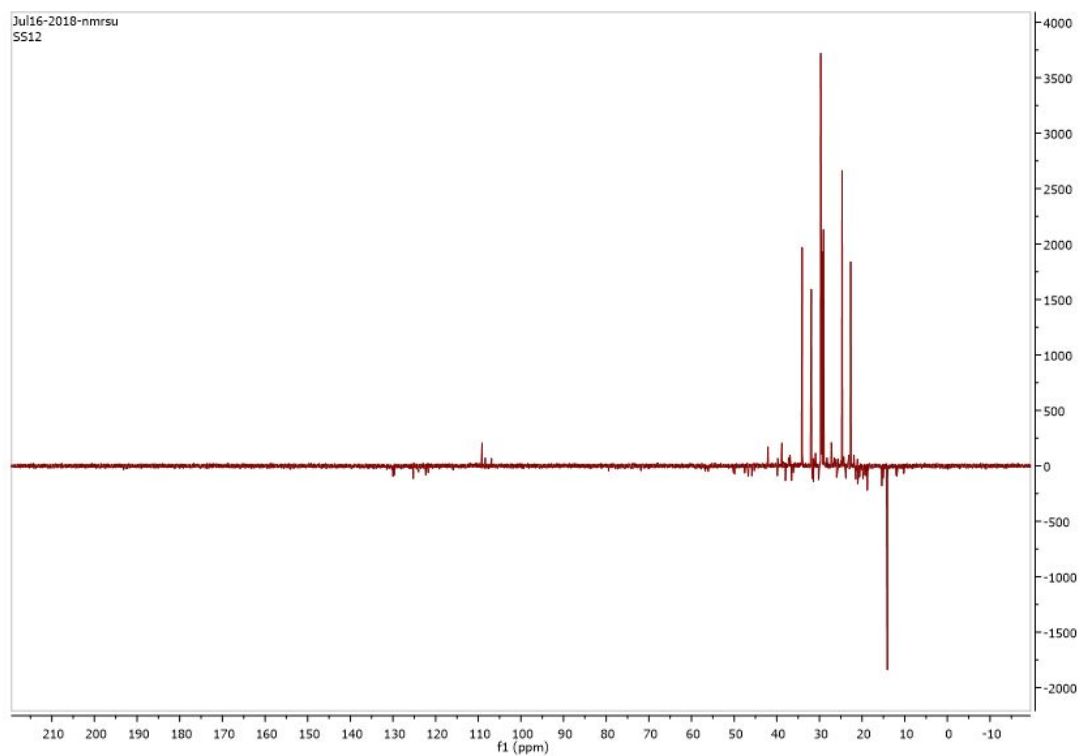
**Density** : 941kg/m



**<sup>1</sup>H NMR Spectrum of compound Stearic acid**



**$^{13}\text{C}$  NMR Spectrum of compound Stearic acid**



**DEPT Spectrum of compound stearic acid**