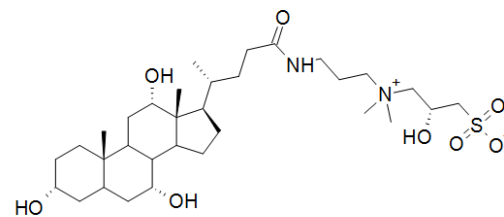


CHAPS

(3-[(3-cholamidopropyl)dimethylammonio]-1-propanesulfonate).



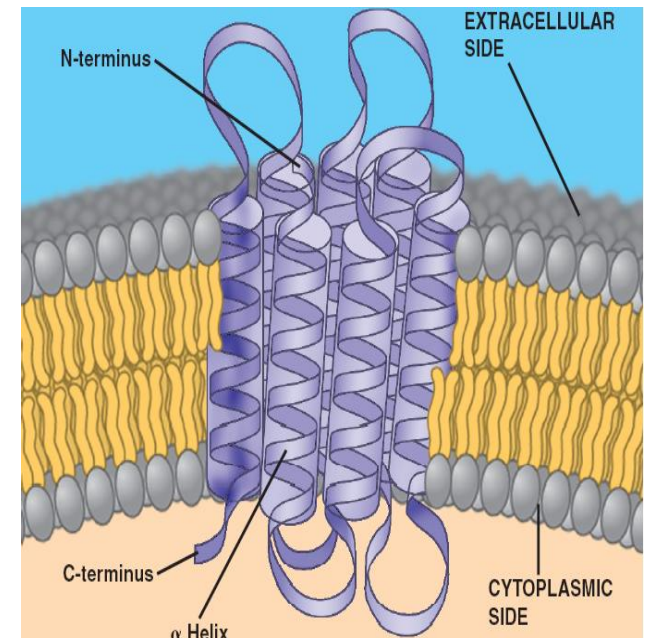
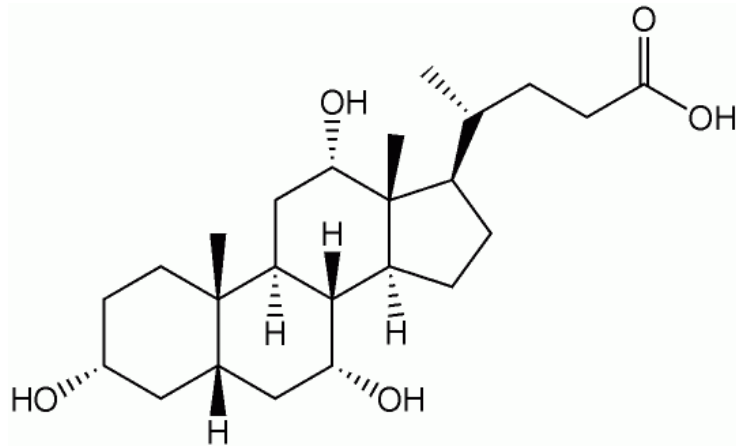
CHAPSO

3-[(3-cholamidopropyl)dimethylammonio]-2 hydroxyl-1-propanesulfonate

CHAPS/CHAPSO

By Martial Fotso and Amanda Vo

History and Description



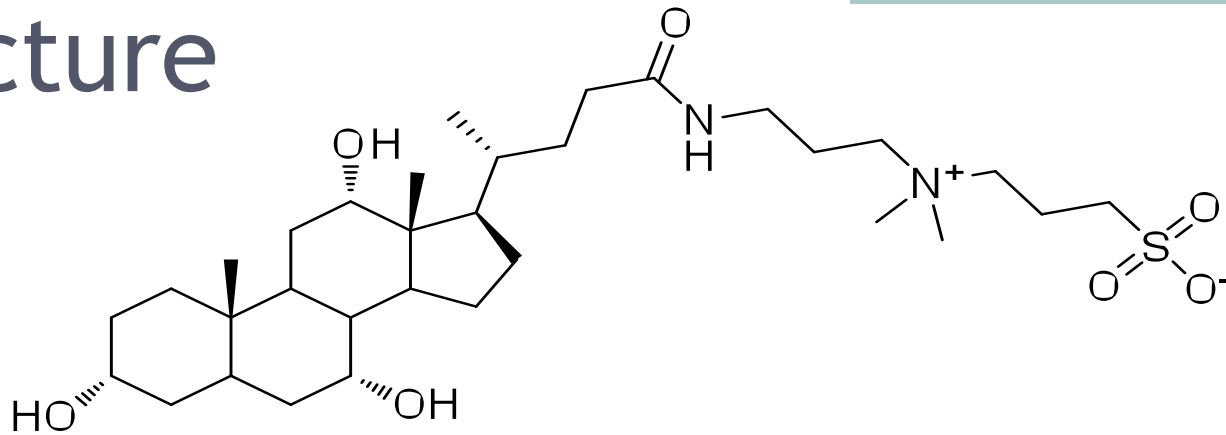
- Cholic acid derivative
- Small pH range in which net charge is present
- Mild surfactant
- Does not affect protein charge
- 1980s used in protein purification

Qin X., Liu M., Daiwen Yang D., and Zhang X. Concentration-Dependent Aggregation of CHAPS Investigated by NMR Spectroscopy. *J. Phys. Chem* **2010**, *114*, 3863–3868.

Biology 1152, <http://www.bio1152.nicerweb.net> (Accessed March 6 2012)

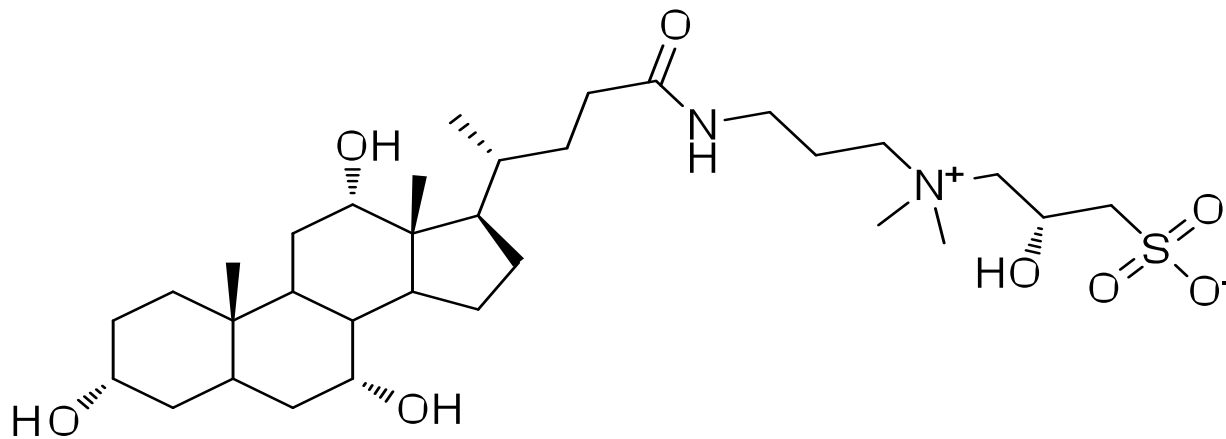
Wikipedia, <http://www.en.wikipedia.org> (Accessed March 6 2012)

Structure



CHAPS

(3-[(3-cholamidopropyl)dimethylammonio] 1-propanesulfonate).



CHAPSO

3-[(3-cholamidopropyl)dimethylammonio] 2 hydroxy-1-propane sulfonate

Uses

- Protein solubilization
- Disaggregation
- Eluting agent during hydrophobic interaction chromatography



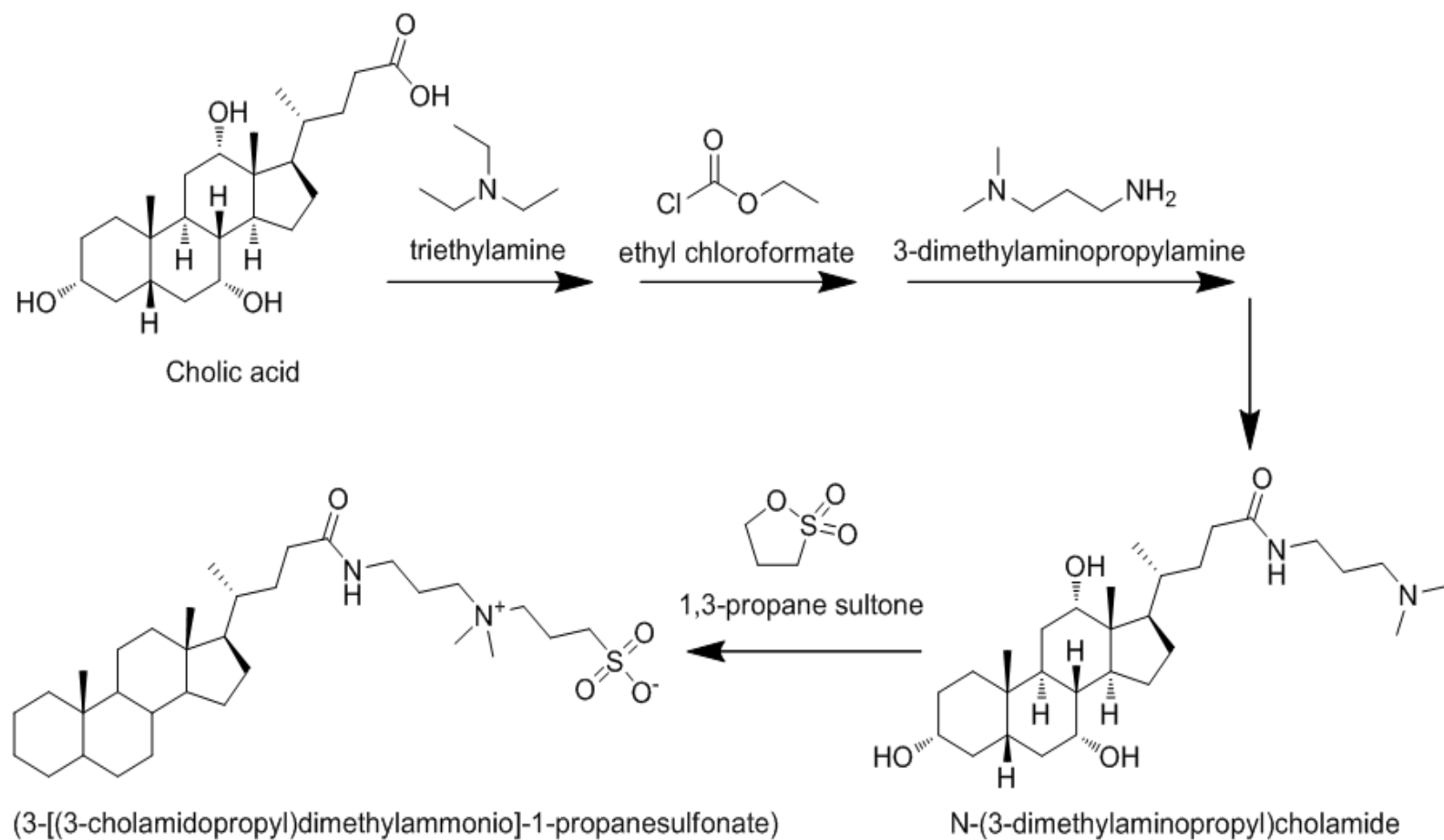
Giacomelli C., Arnouldus W. P., Vermeer, and Norde W. Micellization and Adsorption Characteristics of CHAPS, *Langmuir* **2000**, *16*, 4853-4858.
EMP Genetech, <http://www.empgenetech.com> (Accessed March 6 2012)

Hazards of CHAPS



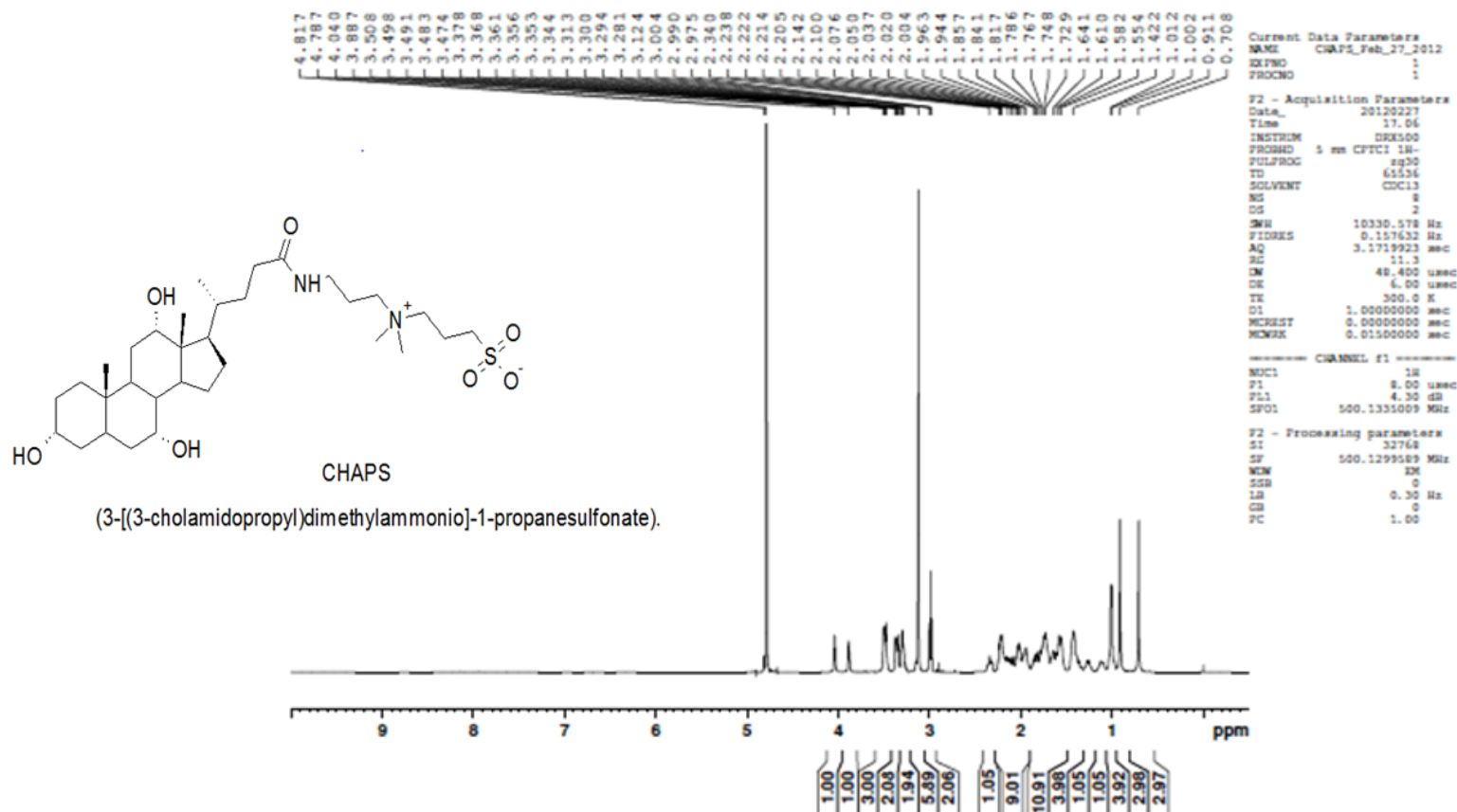
- Reactive with oxidizing reagents and acids
- Slightly combustible and should be kept away from flame
- Minimal skin contact
- Do not ingest (obviously)

Synthesis

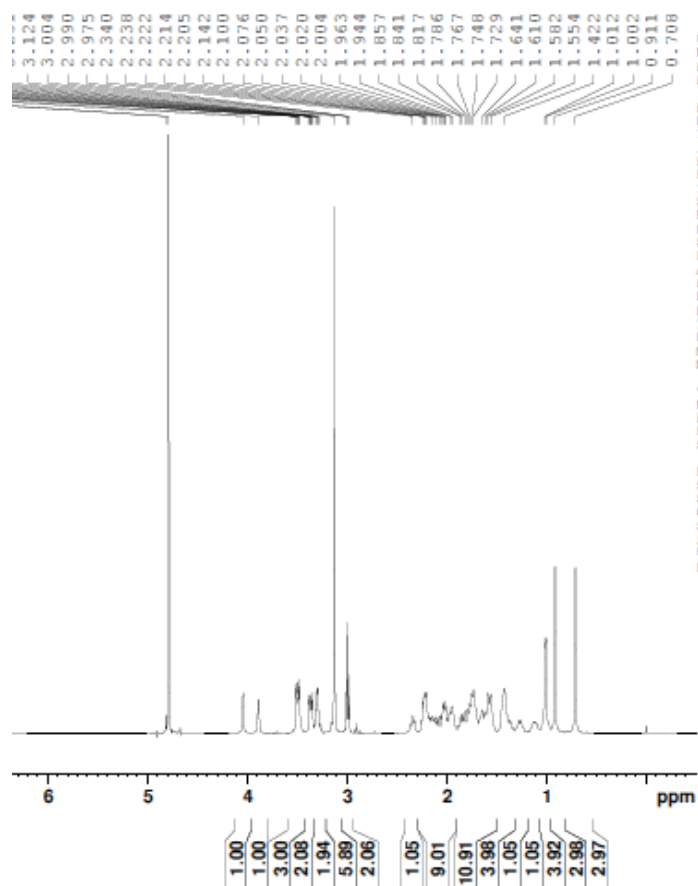


Hjelmeland, L. M. A Nondenaturing Zwitterionic detergent for Membrane Biochemistry Design and Synthesis. *Proc. Natl. Acad. Sci. USA* **1980**, 77, 6368-6370.

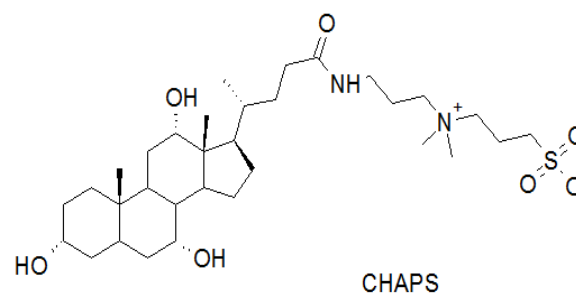
^1H 1-Dimensional Proton NMR



^1H 1-Dimensional Proton NMR

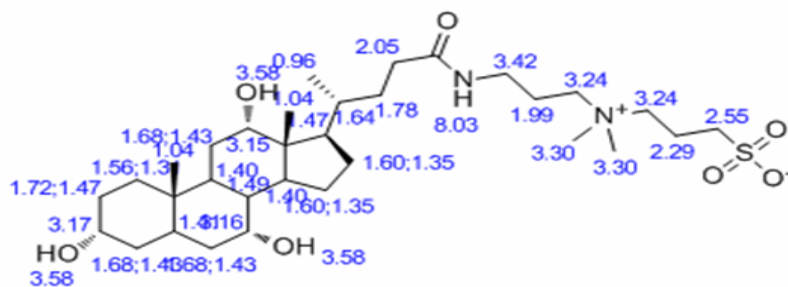


Current Data Parameters
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 PROCNO 1
 F2 - Acquisition Parameters
 Date_ 20120227
 Time 17.06
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 PROBHD 5 mm CPYCL 1H-
 PULPROG zg30
 TD 65536
 SOLVENT DMSO
 NS 8
 DS 2
 SWH 10330.578 Hz
 FIDRES 0.157632 Hz
 AQ 3.1719923 sec
 RG 11.3
 DM 48.400 use
 DE 6.00 use
 YS 300.0 K
 D1 1.0000000 sec
 MCRBST 0.0000000 sec
 MCRBK 0.0150000 sec
 ===== CHANNEL f1 =====
 NUC1 1H
 P1 8.00 use
 PL1 4.30 dB
 SFO1 500.1335009 MHz
 F2 - Processing parameters
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 SF 500.1299089 MHz
 WDM 82
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

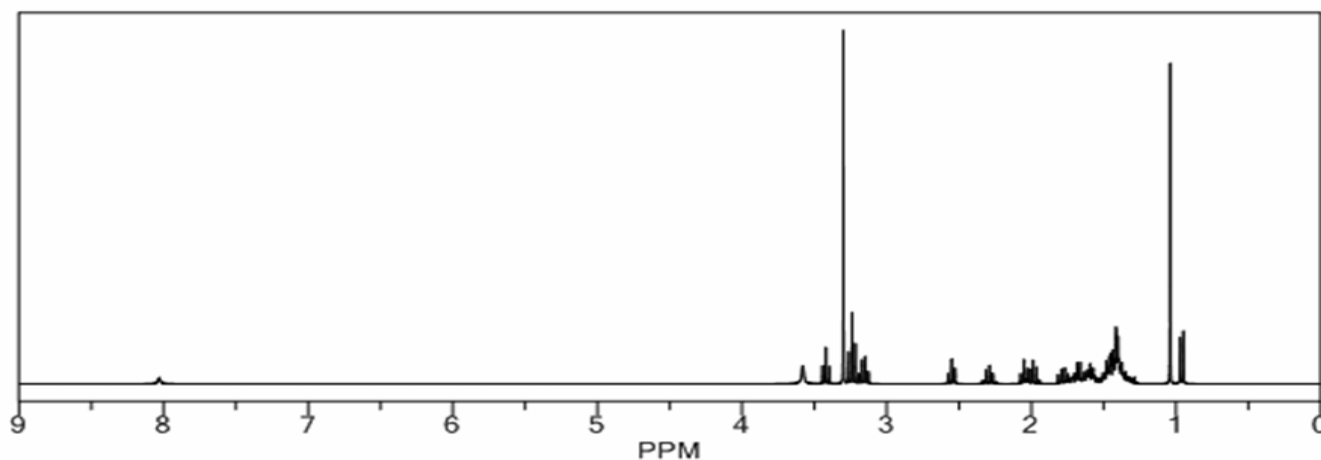


(3-[(3-choamidopropyl)dimethylammonio]-1-propanesulfonate).

¹H 1-Dimensional Proton NMR as predicted by ChemBioDraw

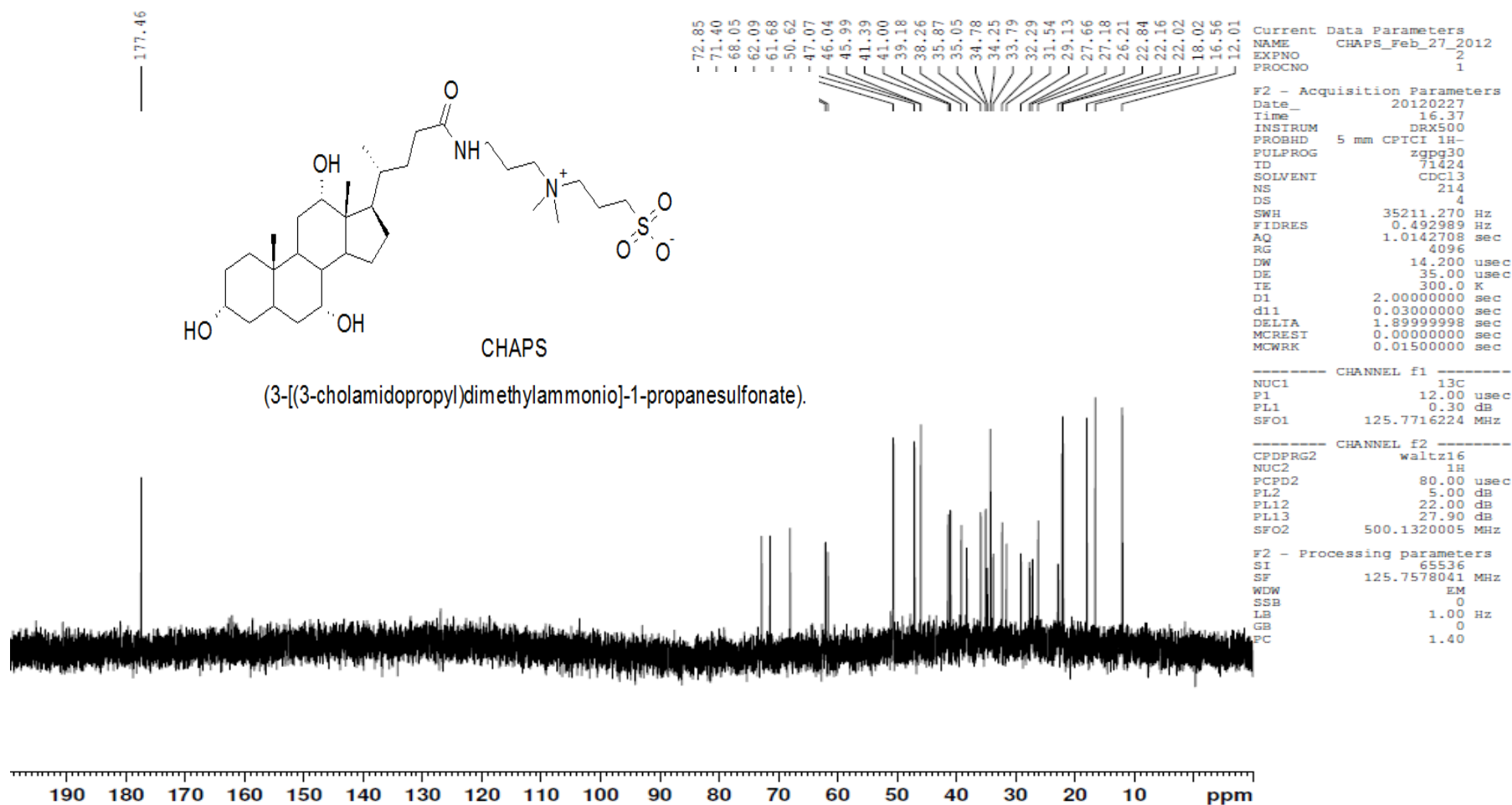
ChemNMR ¹H Estimation

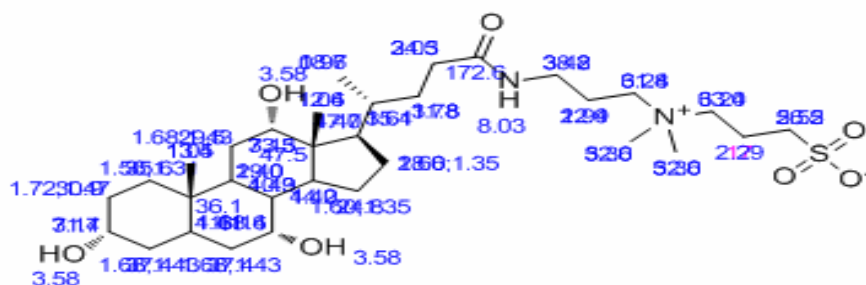
Estimation quality is indicated by color: **good**, **medium**, **rough**



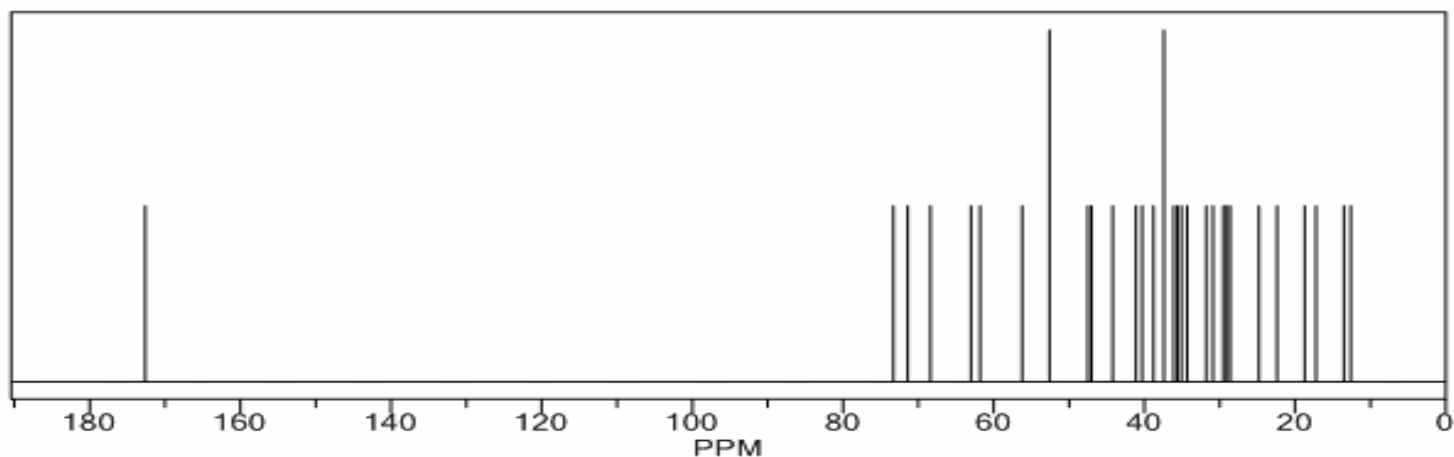
^{13}C 1-Dimensional NMR

^{13}C NMR

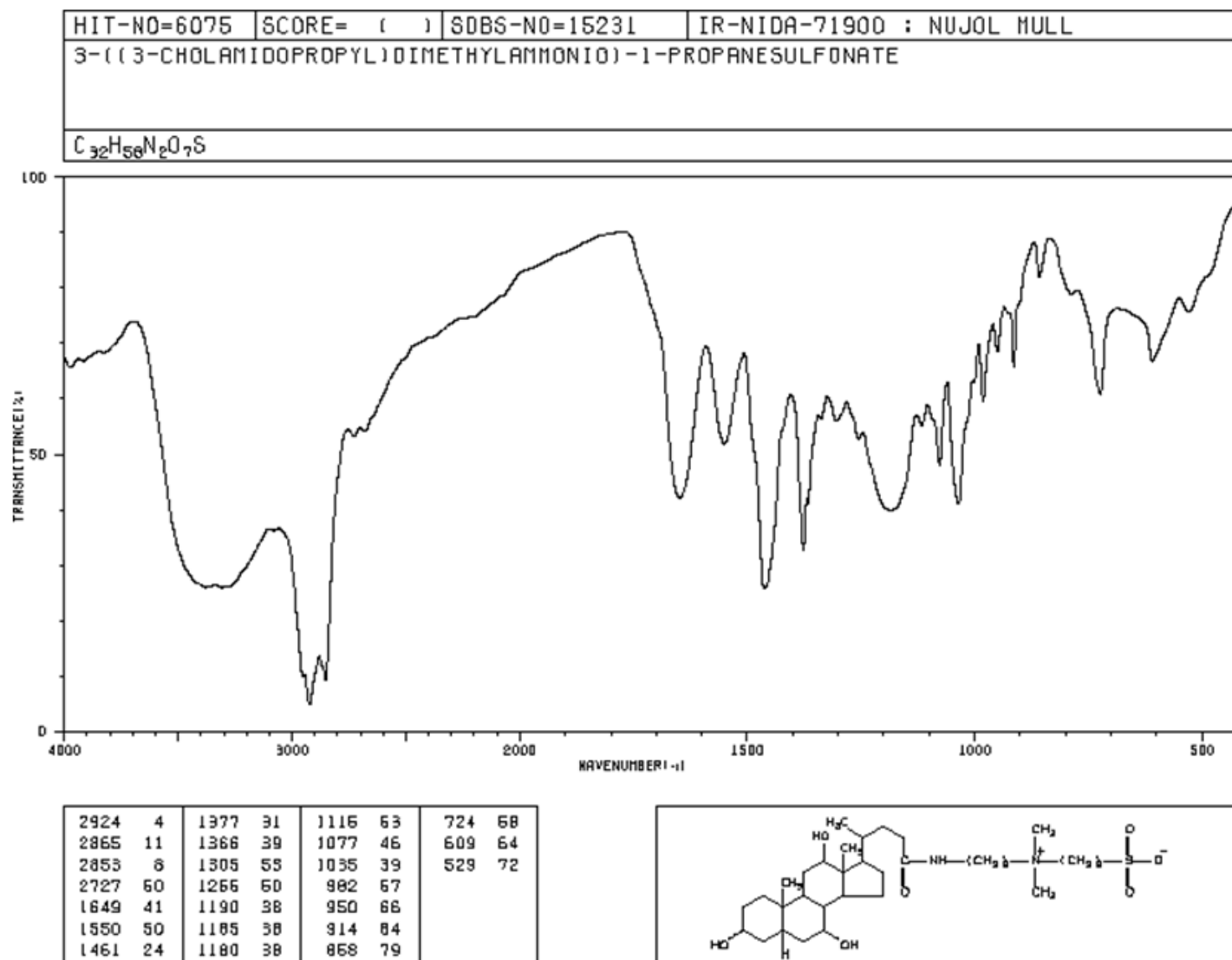


ChemNMR ¹³C Estimation

Estimation quality is indicated by color: **good**, **medium**, **rough**



IR



Spectral Data base for Organic Compounds http://riodb01.ibase.aist.go.jp/sdbs/cgi-bin/cre_index.cgi?lang=eng



Thanks Everyone
Amanda Vo and Martial Fotso