>>**Entry Name: 1-Fluoro-2-propanone,** 9CI, 8CI
Synonym(s): Fluoroacetone

CAS Registry Number: 430-51-3

H_3CCOCH_2F

Molecular Formula: C_3H_5FO
Molecular Weight: 76.07
Accurate Mass: 76.032443
Percentage Composition: C 47.37%; H 6.62%; F 24.97%; O 21.03%
Biological Source: Tentatively identified as a metab. of *Acacia georginae*, prob. as an aberrant prod. of fatty acid biosynth. A metab. prod. of fluoroacetate in rat liver
Physical Description: Yellow or colourless liq.
Solubility: Mod. sol. H_2O, forming a const. boiling mixt.

Aldrich: 11546-0
Sigma: F4509

>>**Derivative:** 2,4-Dinitrophenylhydrazone

Physical Description: Orange cryst. (EtOH or isopropanol)
Melting Point: Mp 118-119°. Mp 134-135°

>>**Derivative:** Semicarbazone

Physical Description: Cryst.
Melting Point: Mp 142-143° (140°, 131-135°)

References:

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