>>**Entry Name: 4-Quinolinecarboxaldehyde,** 9CI
Synonym(s): 4-Formylquinoline. Cinchoninaldehyde

CAS Registry Number: 4363-93-3

Molecular Formula: $C_{10}H_7NO$
Molecular Weight: 157.171
Accurate Mass: 157.052764
Percentage Composition: C 76.42%; H 4.49%; N 8.91%; O 10.18%
Biological Source: Alkaloid prod. by *Archangium gephyra*
Physical Description: Orange needles, plates + $1H_2O$
Melting Point: Mp 51-53° (monohydrate 84-84.5°)
Boiling Point: Bp 122-123°
Solubility: Sol. H_2O, EtOH, Et_2O, toluene

Aldrich: 17696-6
Fluka: 22665

>>**Derivative:** Picrate

Physical Description: Cryst. (EtOH)
Melting Point: Mp 179° (sinters at 170°)

>>**Derivative:** Oxime

CAS Registry Number: 39977-74-7

Molecular Formula: $C_{10}H_8N_2O$
Molecular Weight: 172.186
Accurate Mass: 172.063663
Percentage Composition: C 69.76%; H 4.68%; N 16.27%; O 9.29%
Biological Source: Alkaloid prod. by *Archangium gephyra*
Physical Description: Needles (MeOH)
Melting Point: Mp 181-182°

>>**Derivative:** Phenylhydrazone

Physical Description: Yellow cryst.
Melting Point: Mp 175-176°

>>**Derivative:** Semicarbazone

Melting Point: Mp 244-245 dec.°

References:

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Forbis, R.M. *et al.*, *J.A.C.S.*, 1973, **95**, 5003
Rodriguez, J.G. *et al.*, *J. Het. Chem.*, 1988, **25**, 819, *(synth)*
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