



>>Entry Name: 3,5-Dihydroxybenzaldehyde, 9CI

Synonym(s): α -Resorcyaldehyde, 8CI. α -Resorcylic aldehyde

CAS Registry Number: 26153-38-8

Related CAS Registry Numbers(s): 34967-25-4 57179-37-0

Structure by analogy with 2,3-Dihydroxybenzaldehyde, [BJR22-I](#)

Molecular Formula: $C_7H_6O_3$

Molecular Weight: 138.123

Accurate Mass: 138.031695

Percentage Composition: C 60.87%; H 4.38%; O 34.75%

Physical Description: Needles (petrol)

Melting Point: Mp 161-162°

Aldrich: 36811-3

Fluka: 37523

>>Derivative: Semicarbazone

Physical Description: Needles

Melting Point: Mp 223-224 dec.°

>>Derivative: Mono Me ether

Synonym(s): 3-Hydroxy-5-methoxybenzaldehyde

CAS Registry Number: 57179-35-8

Molecular Formula: $C_8H_8O_3$

Molecular Weight: 152.149

Accurate Mass: 152.047345

Percentage Composition: C 63.15%; H 5.30%; O 31.55%

Biological Source: Constit. of *Ranunculus ternatus*

Melting Point: Mp 130-131°

Other Data: Sublimes at 110-20°

>>Derivative: Di-Me ether

Synonym(s): 3,5-Dimethoxybenzaldehyde, 9CI, 8CI

CAS Registry Number: 7311-34-4

Molecular Formula: $C_9H_{10}O_3$

Molecular Weight: 166.176

Accurate Mass: 166.062995

Percentage Composition: C 65.05%; H 6.07%; O 28.88%

Physical Description: Cryst. (EtOH aq. or petrol)

Melting Point: Mp 45°

Boiling Point: Bp₁₆ 151°

Aldrich: 12629-2

Fluka: 38640

Sigma: D1651

>>Derivative: Di-Me ether, oxime

CAS Registry Number: 34967-25-4

Molecular Formula: $C_9H_{11}NO_3$

Molecular Weight: 181.191

Accurate Mass: 181.073894

Percentage Composition: C 59.66%; H 6.12%; N 7.73%; O 26.49%

Physical Description: Needles (petrol)

Melting Point: Mp 119-120°

>>Derivative: Di-Me ether, semicarbazone

Physical Description: Needles (H₂O)

Melting Point: Mp 177-178°