



>>Entry Name: 2-(3-Methoxyphenyl)ethylamine

Synonym(s): 3-Methoxybenzeneethanamine, 9Cl. *m*-Methoxyphenylethylamine, 8Cl

CAS Registry Number: 2039-67-0

Molecular Formula: C₉H₁₃NO

Molecular Weight: 151.208

Accurate Mass: 151.099714

Percentage Composition: C 71.49%; H 8.67%; N 9.26%; O 10.58%

Boiling Point: Bp 122-123°

Refractive Index: n_D²⁰ 1.5380

Aldrich: 27022-9

Fluka: 65164

>>Derivative: Hydrochloride

CAS Registry Number: 2039-54-5

Melting Point: Mp 145-146°

Hazard & Toxicity: LD₅₀ (mus, ipr) 159 mg/kg

Rare Chemicals Library: S32225-3

>>Derivative: *N*-Me

CAS Registry Number: 33543-62-3

Molecular Formula: C₁₀H₁₅NO

Molecular Weight: 165.235

Accurate Mass: 165.115364

Percentage Composition: C 72.69%; H 9.15%; N 8.48%; O 9.68%

Boiling Point: Bp₇ 118°

>>Derivative: *N,N*-Di-Me

CAS Registry Number: 61134-88-1

Molecular Formula: C₁₁H₁₇NO

Molecular Weight: 179.261

Accurate Mass: 179.131014

Percentage Composition: C 73.70%; H 9.56%; N 7.81%; O 8.93%

Boiling Point: Bp₁₃ 121°. Bp₂ 69-71°

>>Derivative: *N,N*-Di-Me; hydrochloride

CAS Registry Number: 60189-31-3

Melting Point: Mp 135°

References:

Aldrich Library of 13C and 1H FT NMR Spectra, 1992, **2**, 603C, (*nmr*)

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Buck, J.S., *J.A.C.S.*, 1932, **54**, 3661, (*synth*)

Kametani, T., *J.C.S.(C)*, 1971, 2632, (*synth*)

Smith, J.R.L., *J.C.S. Perkin 1*, 1972, 228, (*synth*)

Bailey, K., *Can. J. Pharm. Sci.*, 1975, **10**, 31, (*pmr*)