>>**Entry Name: 3-Fluoro-1,2-benzenediamine,** 9CI
Synonym(s): 3-Fluoro-*o*-phenylenediamine, 8CI. 1,2-Diamino-3-fluorobenzene

CAS Registry Number: 18645-88-0
Structure by analogy with 2-Fluoro-1,3-benzenediamine, HTR26-S

Molecular Formula: $C_6H_7FN_2$
Molecular Weight: 126.133
Accurate Mass: 126.059326
Percentage Composition: C 57.13%; H 5.59%; F 15.06%; N 22.21%
Physical Description: Cryst. by subl.
Melting Point: Mp 40-41°

References:
Kirk, K.L. *et al.*, *J.O.C.*, 1969, **34**, 384, *(synth)*