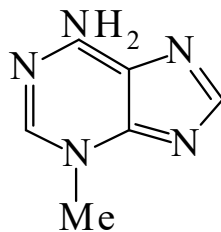




>>Entry Name: 6-Amino-3-methylpurine

Synonym(s): 3-Methyl-3*H*-purin-6-amine, 9Cl. 3-Methyladenine, 8Cl



CAS Registry Number: 5142-23-4

Molecular Formula: C₆H₇N₅

Molecular Weight: 149.155

Accurate Mass: 149.070145

Percentage Composition: C 48.32%; H 4.73%; N 46.95%

Biological Source: Metab. from the marine sponge *Topsentia genitrix*

Physical Description: Cryst. (H₂O)

Melting Point: Mp 300° (291-292° dec.)

pKa Value: pK_a 6.1

UV: [neutral] λ_{max} 273 (ε 14000) (MeOH) (Derep) [neutral] λ_{max} 273 (ε 14010) (MeOH) (Berdy)

Aldrich: 28087-9

Fluka: 8592

Sigma: M9281

>>Derivative: 9-β-D-Ribofuranosyl

Synonym(s): 3-Methyladenosine

CAS Registry Number: 72055-63-1

Molecular Formula: C₁₁H₁₅N₅O₄

Molecular Weight: 281.271

Accurate Mass: 281.112405

Percentage Composition: C 46.97%; H 5.38%; N 24.90%; O 22.75%

Physical Description: Plates (MeOH) (as tosylate salt)

Melting Point: Mp 150 dec.° (tosylate salt)

Optical Rotation: [α]_D²⁰ -28.2 (c, 1.00 in H₂O) (tosylate salt)

Other Data: Imino at C-6

>>Derivative: 9-(2-Deoxy-β-D-ribofuranosyl)

Synonym(s): 2'-Deoxy-3-methyladenosine

CAS Registry Number: 76227-26-4

Molecular Formula: C₁₁H₁₅N₅O₃

Molecular Weight: 265.271

Accurate Mass: 265.11749

Percentage Composition: C 49.81%; H 5.70%; N 26.40%; O 18.09%

Physical Description: Powder (as tosylate salt)

Melting Point: Mp 120 dec.° (tosylate salt)

Other Data: Imino at C-6

>>Derivative: 7-Oxide

Molecular Formula: C₆H₇N₅O

Molecular Weight: 165.154

Accurate Mass: 165.06506

Percentage Composition: C 43.64%; H 4.27%; N 42.40%; O 9.69%

Physical Description: Sl. dark prisms

Melting Point: Mp 250-260° (dec.)

References: