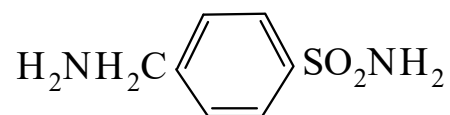




>>Entry Name: **Mafenide**, BAN, INN, USAN

**Synonym(s):** 4-(Aminomethyl)benzenesulfonamide, 9CI.  $\alpha$ -Amino-*p*-toluenesulfonamide, 8CI. 4-Homosulfanilamide. Homosul. Sulfabenzamine. Sulfamylon. NSC 34632.  
Many other names



**CAS Registry Number:** 138-39-6

**Related CAS Registry Numbers(s):** 49783-80-4

**Molecular Formula:** C<sub>7</sub>H<sub>10</sub>N<sub>2</sub>O<sub>2</sub>S

**Molecular Weight:** 186.234

**Accurate Mass:** 186.046298

**Percentage Composition:** C 45.15%; H 5.41%; N 15.04%; O 17.18%; S 17.22%

**Biological Use/Importance:** Used as an antibacterial agent in treatment of second- and third-degree burns

**Physical Description:** Cryst. (EtOH)

**Melting Point:** Mp 151-152°

**Solubility:** Sol. dil. acids, alkalis

**Calculated Log P Data:** Log P -0.74 (calc)

**Hazard & Toxicity:** Low acute toxicity

>>Derivative: Hydrochloride

**Synonym(s):** Homosulfamine

**CAS Registry Number:** 138-37-4

**Melting Point:** Mp 256°

**Sigma:** A2134

>>Derivative: Hydrochloride, hydrate

**Synonym(s):** Homosulfanilamide

**CAS Registry Number:** 65195-43-9

**Biological Use/Importance:** Potent inhibitor of carbonic anhydrase

**Melting Point:** Mp 261-263°

**Aldrich:** A6180-2

>>Derivative: Acetate salt

**Synonym(s):** **Mafenide acetate**, JAN, USAN. Mafatate. Mefamide. Napaltan. Winthrosin. TO 105

**CAS Registry Number:** 13009-99-9

**Melting Point:** Mp 169-172°

>>Derivative: Propanoate salt

**Synonym(s):** Sulfomyl

**Melting Point:** Mp 158°

>>Derivative: Compd. with 4-Amino-*N*-(aminothioxomethyl)benzenesulfonamide (1:1)

**Synonym(s):** **Sulfatolamide**, INN. Sulphatolamide, BAN. Aseptorid. Marbadal. Marbaletten. Mirbedal. Sulfathiamid M

**CAS Registry Number:** 1161-88-2

**Biological Use/Importance:** Antibacterial agent

**Physical Description:** Cryst.

**Melting Point:** Mp 179-181°

**References:**