

Sodium 1-  
octanesulfonate  
 $C_8H_{17}NaO_3S$  structure



## Contents

- Safety:

Molecular      C<sub>8</sub> H<sub>17</sub> NaO<sub>3</sub>

Formula        S

Average mass   216. 273 Da

Density

Boiling Point

Flash Point

Molar

Refractivity

Polarizability

Surface Tension

Molar Volume

- Experimental data
- Predicted - ACD/Labs
- Predicted - ChemAxon
- Experimental Physico-chemical Properties

- **Experimental Melting Point:**

300 °COakwood[009637]

300 °CAifa AesarA14292

300 °COakwood[009637]

300

°CLabNetworkLN00189149

- Miscellaneous

- **Appearance:**

White PowderNovochemistry[NC-44330]

- **Safety:**

20/21/22Novochemistry[NC-44330]

20/21/36/37/39Novochemistry[NC-44330]

GHS07; GHS09Novochemistry[NC-44330]

H332; H403Novochemistry[NC-44330]

IRRITANTMatrix Scientific091211

P102; P210; P262; P270; P302+P352;

P308+P313Novochemistry[NC-44330]

R52/347Novochemistry[NC-44330]

WarningNovochemistry[NC-44330]

Predicted data is generated using the ACD/Labs Percepta Platform -  
PhysChem Module

No predicted properties have been calculated for this compound.

Density:

Boiling Point:

Vapour Pressure:

Enthalpy of Vaporization:

Flash Point:

Index of Refraction:

Molar Refractivity:

#H bond acceptors:

#H bond donors:

#Freely Rotating Bonds:

#Rule of 5 Violations:

ACD/LogP:

ACD/LogD (pH 5. 5):

ACD/BCF (pH 5. 5):

ACD/KOC (pH 5. 5):

ACD/LogD (pH 7. 4):

ACD/BCF (pH 7. 4):

ACD/KOC (pH 7. 4):

Polar Surface Area:

Polarizability:

Surface Tension:

Molar Volume:

Click to predict properties on the Chemicalize site