

# Methylsulphonamide, N-ethyl-N-undecyl-

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C14H31NO2S/c1-4-6-7-8-9-10-11-12-13-14-15(5-2)18(3,16)17/h4-14H2,1-3H3 |
| <b>InchiKey:</b>            | GYCQCFNPIBSHOF-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C14H31NO2S  |
| <b>SMILES:</b>              | CCCCCCCCCCCN(CC)S(C)(=O)=O  |
| <b>Mol. weight [g/mol]:</b> | 277.47  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -290.76 | kJ/mol               | Joback Method  |
| hf            | -718.11 | kJ/mol               | Joback Method  |
| hfus          | 46.41   | kJ/mol               | Joback Method  |
| hvap          | 67.44   | kJ/mol               | Joback Method  |
| log10ws       | -4.08   |                      | Crippen Method |
| logp          | 3.799   |                      | Crippen Method |
| mvol          | 246.190 | ml/mol               | McGowan Method |
| pc            | 1723.16 | kPa                  | Joback Method  |
| rinpol        | 2377.00 |                      | NIST Webbook   |
| rinpol        | 2377.00 |                      | NIST Webbook   |
| tb            | 579.94  | K                    | Joback Method  |
| tc            | 737.63  | K                    | Joback Method  |
| tf            | 318.57  | K                    | Joback Method  |
| vc            | 0.964   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 638.93 | J/mol×K | 579.94          | Joback Method |
| cpg           | 657.26 | J/mol×K | 606.22          | Joback Method |
| cpg           | 674.82 | J/mol×K | 632.50          | Joback Method |
| cpg           | 691.64 | J/mol×K | 658.79          | Joback Method |
| cpg           | 707.74 | J/mol×K | 685.07          | Joback Method |
| cpg           | 723.11 | J/mol×K | 711.35          | Joback Method |
| cpg           | 737.78 | J/mol×K | 737.63          | Joback Method |

# Sources

|                        |   |
|------------------------|---|
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U415443&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U415443&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |

# Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpola:</b> | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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