

# Isobutyl angelate

|                             |                                                                 |
|-----------------------------|-----------------------------------------------------------------|
| <b>Other names:</b>         | 2-Methylpropyl angelate<br>isobutyl 2-methylisocrotonate        |
| <b>Inchi:</b>               | InChI=1S/C9H16O2/c1-5-8(4)9(10)11-6-7(2)3/h5,7H,6H2,1-4H3/b8-5- |
| <b>InchiKey:</b>            | XDEGQMQRKHFPEW-YVMONPNESA-N                                     |
| <b>Formula:</b>             | C9H16O2                                                         |
| <b>SMILES:</b>              | CC=C(C)C(=O)OCC(C)C                                             |
| <b>Mol. weight [g/mol]:</b> | 156.22                                                          |
| <b>CAS:</b>                 | 7779-81-9                                                       |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -139.79 | kJ/mol  | Joback Method  |
| hf            | -371.74 | kJ/mol  | Joback Method  |
| hfus          | 17.22   | kJ/mol  | Joback Method  |
| hvap          | 44.43   | kJ/mol  | Joback Method  |
| log10ws       | -2.06   |         | Crippen Method |
| logp          | 2.152   |         | Crippen Method |
| mcvol         | 140.810 | ml/mol  | McGowan Method |
| pc            | 2574.11 | kPa     | Joback Method  |
| rinpol        | 1030.00 |         | NIST Webbook   |
| rinpol        | 1035.00 |         | NIST Webbook   |
| rinpol        | 1060.00 |         | NIST Webbook   |
| rinpol        | 1055.40 |         | NIST Webbook   |
| rinpol        | 1030.00 |         | NIST Webbook   |
| rinpol        | 1055.40 |         | NIST Webbook   |
| rinpol        | 1035.00 |         | NIST Webbook   |
| rinpol        | 1033.00 |         | NIST Webbook   |
| rinpol        | 1035.00 |         | NIST Webbook   |
| rinpol        | 1030.00 |         | NIST Webbook   |
| ripol         | 1293.00 |         | NIST Webbook   |
| ripol         | 1287.00 |         | NIST Webbook   |
| ripol         | 1293.00 |         | NIST Webbook   |
| ripol         | 1287.00 |         | NIST Webbook   |
| tb            | 485.21  | K       | Joback Method  |
| tc            | 674.06  | K       | Joback Method  |
| tf            | 229.31  | K       | Joback Method  |
| vc            | 0.538   | m3/kmol | Joback Method  |

# Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 306.09 | J/mol×K | 485.21          | Joback Method |
| cpg           | 319.56 | J/mol×K | 516.68          | Joback Method |
| cpg           | 332.44 | J/mol×K | 548.16          | Joback Method |
| cpg           | 344.74 | J/mol×K | 579.63          | Joback Method |
| cpg           | 356.48 | J/mol×K | 611.11          | Joback Method |
| cpg           | 367.66 | J/mol×K | 642.58          | Joback Method |
| cpg           | 378.31 | J/mol×K | 674.06          | Joback Method |

## Sources

|                        |                                                                                                                                             |
|------------------------|---------------------------------------------------------------------------------------------------------------------------------------------|
| <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=C7779819&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C7779819&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>                                   |
| <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>                                       |

## 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>rinpol:</b>  | Non-polar retention indices                     |
| <b>ripol:</b>   | Polar retention indices                         |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |

**tf:** Normal melting (fusion) point

**vc:** Critical Volume

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