

# Butyl decyl ether

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C14H30O/c1-3-5-7-8-9-10-11-12-14-15-13-6-4-2/h3-14H2,1-2H3 |
| <b>InchiKey:</b>            | CGTNCTXIQDWSOL-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C14H30O   |
| <b>SMILES:</b>              | CCCCCCCCCOCCCC  |
| <b>Mol. weight [g/mol]:</b> | 214.39  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -38.00  | kJ/mol               | Joback Method  |
| hf            | -464.51 | kJ/mol               | Joback Method  |
| hfus          | 33.20   | kJ/mol               | Joback Method  |
| hvap          | 49.17   | kJ/mol               | Joback Method  |
| log10ws       | -4.77   |                      | Crippen Method |
| logp          | 4.944   |                      | Crippen Method |
| mcvol         | 213.990 | ml/mol               | McGowan Method |
| pc            | 1503.48 | kPa                  | Joback Method  |
| rinpol        | 1475.00 |                      | NIST Webbook   |
| tb            | 542.14  | K                    | Joback Method  |
| tc            | 701.88  | K                    | Joback Method  |
| tf            | 269.77  | K                    | Joback Method  |
| vc            | 0.838   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 536.33    | J/mol×K | 542.14          | Joback Method |
| cpg           | 554.10    | J/mol×K | 568.76          | Joback Method |
| cpg           | 571.22    | J/mol×K | 595.39          | Joback Method |
| cpg           | 587.72    | J/mol×K | 622.01          | Joback Method |
| cpg           | 603.58    | J/mol×K | 648.63          | Joback Method |
| cpg           | 618.84    | J/mol×K | 675.26          | Joback Method |
| cpg           | 633.50    | J/mol×K | 701.88          | Joback Method |
| dvisc         | 0.0038629 | Paxs    | 269.77          | Joback Method |
| dvisc         | 0.0015169 | Paxs    | 315.16          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0007537 | Paxs | 360.56 | Joback Method |
| dvisc | 0.0004379 | Paxs | 405.95 | Joback Method |
| dvisc | 0.0002838 | Paxs | 451.35 | Joback Method |
| dvisc | 0.0001991 | Paxs | 496.75 | Joback Method |
| dvisc | 0.0001482 | Paxs | 542.14 | 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=U406402&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U406402&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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>dvisc:</b>   | Dynamic viscosity                               |
| <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>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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