

# Hexaethylene glycol monoethyl ether

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C14H30O7/c1-2-16-5-6-18-9-10-20-13-14-21-12-11-19-8-7-17-4-3-15/h15H,2-1

UKXKPKBTMYNOFS-UHFFFAOYSA-N

C14H30O7

CCOCCOCCOCCOCCOCCOCCO

310.38

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -699.82  | kJ/mol  | Joback Method  |
| hf            | -1277.84 | kJ/mol  | Joback Method  |
| hfus          | 43.23    | kJ/mol  | Joback Method  |
| hvap          | 77.90    | kJ/mol  | Joback Method  |
| log10ws       | 0.53     |         | Crippen Method |
| logp          | 0.098    |         | Crippen Method |
| mcvol         | 249.210  | ml/mol  | McGowan Method |
| pc            | 1511.67  | kPa     | Joback Method  |
| rinpol        | 2118.00  |         | NIST Webbook   |
| rinpol        | 2118.00  |         | NIST Webbook   |
| tb            | 746.42   | K       | Joback Method  |
| tc            | 916.39   | K       | Joback Method  |
| tf            | 441.74   | K       | Joback Method  |
| vc            | 0.947    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 761.40    | J/mol×K | 746.42          | Joback Method |
| cpg           | 831.23    | J/mol×K | 888.06          | Joback Method |
| cpg           | 818.90    | J/mol×K | 859.73          | Joback Method |
| cpg           | 805.73    | J/mol×K | 831.41          | Joback Method |
| cpg           | 791.75    | J/mol×K | 803.08          | Joback Method |
| cpg           | 776.96    | J/mol×K | 774.75          | Joback Method |
| cpg           | 842.69    | J/mol×K | 916.39          | Joback Method |
| dvisc         | 0.0000085 | Pa×s    | 746.42          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000127 | Pa×s | 695.64 | Joback Method |
| dvisc | 0.0000201 | Pa×s | 644.86 | Joback Method |
| dvisc | 0.0000345 | Pa×s | 594.08 | Joback Method |
| dvisc | 0.0000656 | Pa×s | 543.30 | Joback Method |
| dvisc | 0.0001423 | Pa×s | 492.52 | Joback Method |
| dvisc | 0.0003688 | Pa×s | 441.74 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=R178842&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R178842&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>                         |

## 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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