

# 8-Oxatricyclo[3,2,1,0(1,5)]octane

**Inchi:** InChI=1S/C7H12O/c1-2-6-4-5-7(3-1)8-6/h6-7H,1-5H2  
**InchiKey:** AWEGTPPFDJUPRI-UHFFFAOYSA-N  
**Formula:** C7H10O  
**SMILES:** C1CC2CCC(C1)O2  
**Mol. weight [g/mol]:** 110.15  
**CAS:** 22341-17-9

## Physical Properties

| Property code | Value           | Unit                 | Source         |
|---------------|-----------------|----------------------|----------------|
| chl           | -4174.30 ± 0.96 | kJ/mol               | NIST Webbook   |
| gf            | 19.24           | kJ/mol               | Joback Method  |
| hf            | 27.00           | kJ/mol               | NIST Webbook   |
| hfl           | -9.54           | kJ/mol               | NIST Webbook   |
| hfus          | 13.93           | kJ/mol               | Joback Method  |
| hvap          | 36.00 ± 1.30    | kJ/mol               | NIST Webbook   |
| hvap          | 36.50           | kJ/mol               | NIST Webbook   |
| log10ws       | -1.85           |                      | Crippen Method |
| logp          | 1.718           |                      | Crippen Method |
| mcvol         | 93.640          | ml/mol               | McGowan Method |
| pc            | 3960.52         | kPa                  | Joback Method  |
| tb            | 408.53          | K                    | Joback Method  |
| tc            | 620.71          | K                    | Joback Method  |
| tf            | 224.06          | K                    | Joback Method  |
| vc            | 0.346           | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 189.08 | J/mol×K | 408.53          | Joback Method |
| cpg           | 205.70 | J/mol×K | 443.89          | Joback Method |
| cpg           | 221.26 | J/mol×K | 479.26          | Joback Method |
| cpg           | 235.82 | J/mol×K | 514.62          | Joback Method |
| cpg           | 249.44 | J/mol×K | 549.98          | Joback Method |
| cpg           | 262.16 | J/mol×K | 585.35          | Joback Method |

|       |           |         |        |               |
|-------|-----------|---------|--------|---------------|
| cpg   | 274.04    | J/molxK | 620.71 | Joback Method |
| dvisc | 0.0018676 | Paxs    | 224.06 | Joback Method |
| dvisc | 0.0013582 | Paxs    | 254.81 | Joback Method |
| dvisc | 0.0010578 | Paxs    | 285.55 | Joback Method |
| dvisc | 0.0008649 | Paxs    | 316.29 | Joback Method |
| dvisc | 0.0007328 | Paxs    | 347.04 | Joback Method |
| dvisc | 0.0006379 | Paxs    | 377.78 | Joback Method |
| dvisc | 0.0005670 | Paxs    | 408.53 | 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=C22341179&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C22341179&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>                                     |

## Legend

|                 |   |
|-----------------|---|
| <b>chl:</b>     | Standard liquid enthalpy of combustion                    |
| <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>hfl:</b>     | Liquid phase 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>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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