

# Cyclotetracosane

**Inchi:** InChI=1S/C24H48/c1-2-4-6-8-10-12-14-16-18-20-22-24-23-21-19-17-15-13-11-9-7-5-3-1  
**InchiKey:** OOTMSKDHDITFHF-UHFFFAOYSA-N  
**Formula:** C24H48  
**SMILES:** C1CCCCCCCCCCCCCCCCCCCCCCC1  
**Mol. weight [g/mol]:** 336.64  
**CAS:** 297-03-0

## Physical Properties

| Property code | Value         | Unit    | Source         |
|---------------|---------------|---------|----------------|
| gf            | -34.44        | kJ/mol  | Joback Method  |
| hf            | -574.91       | kJ/mol  | Joback Method  |
| hfus          | 10.88         | kJ/mol  | Joback Method  |
| hvap          | 72.85         | kJ/mol  | Joback Method  |
| log10ws       | -9.76         |         | Crippen Method |
| logp          | 9.362         |         | Crippen Method |
| mcvol         | 338.160       | ml/mol  | McGowan Method |
| pc            | 1245.09       | kPa     | Joback Method  |
| rinpol        | 2589.00       |         | NIST Webbook   |
| tb            | 849.60        | K       | Joback Method  |
| tc            | 1120.84       | K       | Joback Method  |
| tf            | 322.00 ± 0.30 | K       | NIST Webbook   |
| tf            | 319.90 ± 2.00 | K       | NIST Webbook   |
| tf            | 320.00 ± 3.00 | K       | NIST Webbook   |
| tf            | 320.00 ± 3.00 | K       | NIST Webbook   |
| vc            | 1.169         | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value   | Unit    | Temperature [K] | Source        |
|---------------|---------|---------|-----------------|---------------|
| cpg           | 1299.36 | J/mol×K | 1120.84         | Joback Method |
| cpg           | 1292.07 | J/mol×K | 1075.64         | Joback Method |
| cpg           | 1279.42 | J/mol×K | 1030.43         | Joback Method |
| cpg           | 1261.53 | J/mol×K | 985.22          | Joback Method |
| cpg           | 1238.53 | J/mol×K | 940.01          | Joback Method |

|       |               |         |        |               |
|-------|---------------|---------|--------|---------------|
| cpg   | 1210.54       | J/mol×K | 894.81 | Joback Method |
| cpg   | 1177.69       | J/mol×K | 849.60 | Joback Method |
| dvisc | 0.0782183     | Pa×s    | 308.50 | Joback Method |
| dvisc | 6.4841321e-08 | Pa×s    | 849.60 | Joback Method |
| dvisc | 0.0000002     | Pa×s    | 759.42 | Joback Method |
| dvisc | 0.0000006     | Pa×s    | 669.23 | Joback Method |
| dvisc | 0.0000027     | Pa×s    | 579.05 | Joback Method |
| dvisc | 0.0000235     | Pa×s    | 488.87 | Joback Method |
| dvisc | 0.0005412     | Pa×s    | 398.68 | Joback Method |
| hfust | 10.80         | kJ/mol  | 322.00 | NIST Webbook  |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C297030&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C297030&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>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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>hfust:</b>   | Enthalpy of fusion at a given temperature       |
| <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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