

# 1-Propene, 1,2,3-trichloro-, (Z)-

|                      |                                                                                                                             |
|----------------------|-----------------------------------------------------------------------------------------------------------------------------|
| Other names:         | cis-1,2,3-Trichloro-1-propene<br>1,2,3-Trichloropropene, Z-<br>(1Z)-1,2,3-Trichloro-1-propene<br>cis-1,2,3-Trichloropropene |
| Inchi:               | InChI=1S/C3H3Cl3/c4-1-3(6)2-5/h1H,2H2/b3-1-                                                                                 |
| InchiKey:            | HIILBTHBHCLUER-IWQZZHSRSA-N                                                                                                 |
| Formula:             | C3H3Cl3                                                                                                                     |
| SMILES:              | ClC=C(Cl)CCl                                                                                                                |
| Mol. weight [g/mol]: | 145.41                                                                                                                      |
| CAS:                 | 13116-57-9                                                                                                                  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 10.26   | kJ/mol  | Joback Method  |
| hf            | -45.04  | kJ/mol  | Joback Method  |
| hfus          | 15.01   | kJ/mol  | Joback Method  |
| hvap          | 35.47   | kJ/mol  | Joback Method  |
| log10ws       | -2.38   |         | Crippen Method |
| logp          | 2.544   |         | Crippen Method |
| mcvol         | 85.550  | ml/mol  | McGowan Method |
| pc            | 4109.14 | kPa     | Joback Method  |
| tb            | 384.37  | K       | Joback Method  |
| tc            | 590.66  | K       | Joback Method  |
| tf            | 194.29  | K       | Joback Method  |
| vc            | 0.332   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 114.82 | J/mol×K | 384.37          | Joback Method |
| cpg           | 119.89 | J/mol×K | 418.75          | Joback Method |
| cpg           | 124.59 | J/mol×K | 453.13          | Joback Method |
| cpg           | 128.93 | J/mol×K | 487.51          | Joback Method |
| cpg           | 132.94 | J/mol×K | 521.89          | Joback Method |

|     |        |         |        |               |
|-----|--------|---------|--------|---------------|
| cpg | 136.65 | J/mol×K | 556.28 | Joback Method |
| cpg | 140.07 | J/mol×K | 590.66 | Joback Method |

## Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C13116579&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C13116579&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>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>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

Latest version available from:

<https://www.chemeo.com/cid/56-829-0/1-Propene-1-2-3-trichloro-Z.pdf>

Generated by Cheméo on 2025-12-22 19:30:50.50355507 +0000 UTC m=+6180048.033595724.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.