

# neo-Thujan-3-ol

|                      |                                                                                    |
|----------------------|------------------------------------------------------------------------------------|
| Other names:         | neo-3-thujanol<br>Neothujol<br>3-neo-Thujanol                                      |
| Inchi:               | InChI=1S/C10H18O/c1-6(2)10-4-8(10)7(3)9(11)5-10/h6-9,11H,4-5H2,1-3H3/t7-,8-,9-,10+ |
| InchiKey:            | DZVXRFMREAADPP-KYXWUPHJSA-N                                                        |
| Formula:             | C10H18O                                                                            |
| SMILES:              | CC1C(O)CC2(C(C)C)CC12                                                              |
| Mol. weight [g/mol]: | 154.25                                                                             |
| CAS:                 | 21653-18-9                                                                         |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -5.35   | kJ/mol  | Joback Method  |
| hf            | -287.08 | kJ/mol  | Joback Method  |
| hfus          | 14.34   | kJ/mol  | Joback Method  |
| hvap          | 52.20   | kJ/mol  | Joback Method  |
| log10ws       | -2.21   |         | Crippen Method |
| logp          | 2.049   |         | Crippen Method |
| mcvol         | 135.910 | ml/mol  | McGowan Method |
| pc            | 2956.90 | kPa     | Joback Method  |
| rinpol        | 1130.00 |         | NIST Webbook   |
| rinpol        | 1138.00 |         | NIST Webbook   |
| rinpol        | 1134.00 |         | NIST Webbook   |
| rinpol        | 1151.00 |         | NIST Webbook   |
| rinpol        | 1149.00 |         | NIST Webbook   |
| rinpol        | 1149.00 |         | NIST Webbook   |
| rinpol        | 1148.00 |         | NIST Webbook   |
| rinpol        | 1134.00 |         | NIST Webbook   |
| rinpol        | 1149.00 |         | NIST Webbook   |
| ripol         | 1573.00 |         | NIST Webbook   |
| ripol         | 1620.00 |         | NIST Webbook   |
| tb            | 524.32  | K       | Joback Method  |
| tc            | 714.97  | K       | Joback Method  |
| tf            | 299.58  | K       | Joback Method  |
| vc            | 0.518   | m3/kmol | Joback Method  |

# Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 354.41 | J/mol×K | 524.32          | Joback Method |
| cpg           | 370.07 | J/mol×K | 556.10          | Joback Method |
| cpg           | 384.74 | J/mol×K | 587.87          | Joback Method |
| cpg           | 398.54 | J/mol×K | 619.65          | Joback Method |
| cpg           | 411.57 | J/mol×K | 651.42          | Joback Method |
| cpg           | 423.94 | J/mol×K | 683.20          | Joback Method |
| cpg           | 435.75 | J/mol×K | 714.97          | Joback Method |

## Sources

|                 |                                                                                                                                               |
|-----------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| Joback Method:  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                         |
| McGowan Method: | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                         |
| NIST Webbook:   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C21653189&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C21653189&amp;Units=SI</a> |
| Crippen Method: | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                     |
| Crippen Method: | <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>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>ripol:</b>   | 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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