

Ledene oxide-(II)

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| Other names: | Ledene oxide Ledenoxide |
| Inchi: | InChI=1S/C15H24O/c1-9-5-8-15-11(9)12-10(13(12,2)3)6-7-14(15,4)16-15/h9-12H,5-8H2 |
| InchiKey: | OZNHATCGPKOFBH-UHFFFAOYSA-N |
| Formula: | C15H24O |
| SMILES: | CC1CCC23OC2(C)CCC2C(C13)C2(C)C |
| Mol. weight [g/mol]: | 220.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 192.70 | kJ/mol | Joback Method |
| hf | -209.03 | kJ/mol | Joback Method |
| hfus | 19.44 | kJ/mol | Joback Method |
| hvap | 48.77 | kJ/mol | Joback Method |
| log10ws | -3.78 | | Crippen Method |
| logp | 3.626 | | Crippen Method |
| mcvol | 184.640 | ml/mol | McGowan Method |
| pc | 2244.00 | kPa | Joback Method |
| rinpol | 1631.00 | | NIST Webbook |
| rinpol | 1646.00 | | NIST Webbook |
| rinpol | 1646.00 | | NIST Webbook |
| rinpol | 1682.00 | | NIST Webbook |
| ripol | 2269.00 | | NIST Webbook |
| tb | 583.22 | K | Joback Method |
| tc | 812.69 | K | Joback Method |
| tf | 416.12 | K | Joback Method |
| vc | 0.716 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 540.98 | J/molxK | 583.22 | Joback Method |
| cpg | 563.50 | J/molxK | 621.46 | Joback Method |
| cpg | 584.43 | J/molxK | 659.71 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 604.21 | J/mol×K | 697.95 | Joback Method |
| cpg | 623.32 | J/mol×K | 736.20 | Joback Method |
| cpg | 642.20 | J/mol×K | 774.44 | Joback Method |
| cpg | 661.33 | J/mol×K | 812.69 | Joback Method |

Sources

| | |
|------------------------|---|
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U159367&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

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|-----------------|---|
| cpg: | Ideal gas heat capacity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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