

# Cyclopentane, 1-ethyl-2-methyl-, cis-

<b>Other names:</b>	1-Ethyl-2-methylcyclopentane, (Z)- 1-Ethyl-2-methylcyclopentane, cis 1-Methyl-cis-2-ethylcyclopentane CIS-1-ETHYL-2-METHYLCYCLOPENTANE CIS-1-METHYL-2-ETHYLCYCLOPENTANE c-1-Ethyl-2-methylcyclopentane
<b>Inchi:</b>	InChI=1S/C8H16/c1-3-8-6-4-5-7(8)2/h7-8H,3-6H2,1-2H3/t7-,8+/m0/s1
<b>InchiKey:</b>	BSKOLJVTLRLTHE-JGVFFNPUSA-N
<b>Formula:</b>	C8H16
<b>SMILES:</b>	CCC1CCCC1C
<b>Mol. weight [g/mol]:</b>	112.21
<b>CAS:</b>	930-89-2

## Physical Properties

Property code	Value	Unit	Source
chl	-5243.89 ± 0.92	kJ/mol	NIST Webbook
gf	45.32	kJ/mol	Joback Method
hf	-168.31	kJ/mol	Joback Method
hfl	-190.80 ± 1.00	kJ/mol	NIST Webbook
hfus	11.48	kJ/mol	Joback Method
hvap	40.20	kJ/mol	NIST Webbook
log10ws	-2.58		Crippen Method
logp	2.833		Crippen Method
mcvol	112.720	ml/mol	McGowan Method
pc	2956.90	kPa	Joback Method
rinpol	825.30		NIST Webbook
rinpol	820.80		NIST Webbook
rinpol	825.30		NIST Webbook
rinpol	825.40		NIST Webbook
rinpol	833.00		NIST Webbook
rinpol	823.00		NIST Webbook
rinpol	828.00		NIST Webbook
rinpol	821.00		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	820.00		NIST Webbook
rinpol	823.00		NIST Webbook
rinpol	824.00		NIST Webbook

rinpol	826.00		NIST Webbook
rinpol	829.00		NIST Webbook
rinpol	821.10		NIST Webbook
rinpol	820.00		NIST Webbook
rinpol	846.00		NIST Webbook
rinpol	816.00		NIST Webbook
rinpol	821.00		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	830.00		NIST Webbook
rinpol	837.00		NIST Webbook
rinpol	820.60		NIST Webbook
rinpol	815.95		NIST Webbook
rinpol	820.30		NIST Webbook
rinpol	818.90		NIST Webbook
rinpol	817.00		NIST Webbook
rinpol	832.00		NIST Webbook
rinpol	817.00		NIST Webbook
rinpol	825.00		NIST Webbook
rinpol	816.00		NIST Webbook
rinpol	822.00		NIST Webbook
rinpol	790.00		NIST Webbook
rinpol	822.00		NIST Webbook
rinpol	821.30		NIST Webbook
rinpol	823.20		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	828.10		NIST Webbook
rinpol	829.80		NIST Webbook
rinpol	831.30		NIST Webbook
rinpol	820.00		NIST Webbook
rinpol	820.70		NIST Webbook
rinpol	825.00		NIST Webbook
rinpol	827.00		NIST Webbook
rinpol	829.00		NIST Webbook
rinpol	831.00		NIST Webbook
rinpol	820.10		NIST Webbook
rinpol	825.30		NIST Webbook
rinpol	821.00		NIST Webbook
rinpol	826.00		NIST Webbook
rinpol	816.00		NIST Webbook
rinpol	820.10		NIST Webbook
rinpol	827.00		NIST Webbook
rinpol	822.00		NIST Webbook
tb	401.20	K	KDB
tc	585.20	K	Joback Method

tf	186.58	K	Joback Method
vc	0.423	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	290.01	J/mol×K	553.18	Joback Method
cpg	246.41	J/mol×K	457.10	Joback Method
cpg	230.45	J/mol×K	425.08	Joback Method
cpg	213.75	J/mol×K	393.05	Joback Method
cpg	261.64	J/mol×K	489.13	Joback Method
cpg	303.19	J/mol×K	585.20	Joback Method
cpg	276.17	J/mol×K	521.15	Joback Method
dvisc	0.0022106	Paxs	186.58	Joback Method
dvisc	0.0002855	Paxs	393.05	Joback Method
dvisc	0.0003410	Paxs	358.64	Joback Method
dvisc	0.0004228	Paxs	324.23	Joback Method
dvisc	0.0005518	Paxs	289.81	Joback Method
dvisc	0.0007737	Paxs	255.40	Joback Method
dvisc	0.0012051	Paxs	220.99	Joback Method
hvapt	38.30	kJ/mol	362.00	NIST Webbook
hvapt	39.30	kJ/mol	353.00	NIST Webbook
hvapt	42.50	kJ/mol	271.00	NIST Webbook
hvapt	41.60	kJ/mol	263.00	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.39419e+01
Coeff. B	-3.21987e+03
Coeff. C	-5.57510e+01
Temperature range (K), min.	291.57
Temperature range (K), max.	428.84

Information

Value

Property code	pvap
Equation	$\ln(P_{vp}) = A + B/T + C \cdot \ln(T) + D \cdot T^2$
Coeff. A	7.66373e+01
Coeff. B	-7.06332e+03
Coeff. C	-9.36311e+00
Coeff. D	1.22005e-05
Temperature range (K), min.	238.15
Temperature range (K), max.	288.15

## Sources

<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>
<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>KDB:</b>	<a href="https://www.thermopedia.com/research/kdb/hcprop/showprop.php?cmpid=483">https://www.thermopedia.com/research/kdb/hcprop/showprop.php?cmpid=483</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=C930892&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C930892&amp;Units=SI</a>
<b>The Yaws Handbook of Vapor Pressure:</b>	<a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</a>
<b>KDB Vapor Pressure Data:</b>	<a href="https://www.thermopedia.com/research/kdb/hcprop/showprop.php?cmpid=483">https://www.thermopedia.com/research/kdb/hcprop/showprop.php?cmpid=483</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>hvapt:</b>	Enthalpy of vaporization at a given temperature
<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>pvap:</b>	Vapor pressure
<b>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature

**tc:** Critical Temperature  
**tf:** Normal melting (fusion) point  
**vc:** Critical Volume

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