

Cyclopentane, 1,2,4-trimethyl-, (1«alpha»,2«beta»,4«alpha»)-

Other names:	Cyclopentane, 1,2,4-trimethyl-, trans,cis-1«alpha»,2«beta»,4«alpha»-Trimethylcyclopentane 1-trans-2,cis-4-Trimethylcyclopentane 1-trans-2-trans-4-Trimethylcyclopentane 1,2,4-Trimethylcyclopentane, trans, trans (1«alpha»,2«beta»,4«alpha»)-1,2,4-trimethylcyclopentane
Inchi:	InChI=1S/C8H16/c1-6-4-7(2)8(3)5-6/h6-8H,4-5H2,1-3H3/t7-,8-/m1/s1
InchiKey:	PNUFYSGVPVMNRN-HTQZYQBOSA-N
Formula:	C8H16
SMILES:	CC1CC(C)C(C)C1
Mol. weight [g/mol]:	112.21
CAS:	16883-48-0

Physical Properties

Property code	Value	Unit	Source
gf	37.61	kJ/mol	Joback Method
hf	-188.65	kJ/mol	Joback Method
hfus	12.55	kJ/mol	Joback Method
hvap	33.04	kJ/mol	Joback Method
log10ws	-2.34		Crippen Method
logp	2.688		Crippen Method
mcvol	112.720	ml/mol	McGowan Method
pc	2853.57	kPa	Joback Method
rinpol	741.00		NIST Webbook
rinpol	746.00		NIST Webbook
rinpol	745.00		NIST Webbook
rinpol	744.00		NIST Webbook
rinpol	732.00		NIST Webbook
rinpol	744.10		NIST Webbook
rinpol	738.00		NIST Webbook
rinpol	741.00		NIST Webbook
rinpol	744.00		NIST Webbook
rinpol	740.50		NIST Webbook
rinpol	743.90		NIST Webbook
rinpol	738.40		NIST Webbook
rinpol	740.10		NIST Webbook
rinpol	741.80		NIST Webbook

rinpol	743.30	NIST Webbook
rinpol	741.00	NIST Webbook
rinpol	736.00	NIST Webbook
rinpol	738.00	NIST Webbook
rinpol	739.00	NIST Webbook
rinpol	741.00	NIST Webbook
rinpol	743.00	NIST Webbook
rinpol	744.00	NIST Webbook
rinpol	744.20	NIST Webbook
rinpol	745.10	NIST Webbook
rinpol	745.20	NIST Webbook
rinpol	745.20	NIST Webbook
rinpol	741.00	NIST Webbook
rinpol	744.00	NIST Webbook
rinpol	739.00	NIST Webbook
rinpol	736.00	NIST Webbook
rinpol	742.00	NIST Webbook
rinpol	744.00	NIST Webbook
rinpol	746.00	NIST Webbook
rinpol	741.30	NIST Webbook
rinpol	739.00	NIST Webbook
rinpol	759.00	NIST Webbook
rinpol	738.00	NIST Webbook
rinpol	741.00	NIST Webbook
rinpol	744.00	NIST Webbook
rinpol	748.00	NIST Webbook
rinpol	740.40	NIST Webbook
rinpol	725.90	NIST Webbook
rinpol	737.80	NIST Webbook
rinpol	735.05	NIST Webbook
rinpol	735.35	NIST Webbook
rinpol	735.37	NIST Webbook
rinpol	735.00	NIST Webbook
rinpol	749.00	NIST Webbook
rinpol	739.00	NIST Webbook
rinpol	744.00	NIST Webbook
rinpol	736.00	NIST Webbook
rinpol	741.00	NIST Webbook
rinpol	740.40	NIST Webbook
rinpol	741.30	NIST Webbook
rinpol	742.70	NIST Webbook
rinpol	743.90	NIST Webbook
rinpol	745.00	NIST Webbook
rinpol	732.00	NIST Webbook

rinpol	730.00		NIST Webbook
rinpol	741.00		NIST Webbook
rinpol	743.00		NIST Webbook
rinpol	745.90		NIST Webbook
tb	388.38	K	Joback Method
tc	579.97	K	Joback Method
tf	142.37 ± 0.07	K	NIST Webbook
tf	142.17 ± 0.30	K	NIST Webbook
vc	0.422	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	213.62	J/mol×K	388.38	Joback Method
cpg	230.52	J/mol×K	420.31	Joback Method
cpg	246.72	J/mol×K	452.24	Joback Method
cpg	262.24	J/mol×K	484.17	Joback Method
cpg	277.08	J/mol×K	516.11	Joback Method
cpg	291.26	J/mol×K	548.04	Joback Method
cpg	304.79	J/mol×K	579.97	Joback Method
dvisc	0.0010560	Pa×s	182.34	Joback Method
dvisc	0.0006962	Pa×s	216.68	Joback Method
dvisc	0.0005145	Pa×s	251.02	Joback Method
dvisc	0.0004089	Pa×s	285.36	Joback Method
dvisc	0.0003414	Pa×s	319.70	Joback Method
dvisc	0.0002952	Pa×s	354.04	Joback Method
dvisc	0.0002619	Pa×s	388.38	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C16883480&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.cheméo.com/cid/63-722-0/Cyclopentane-1-2-4-trimethyl-1-alpha-2-beta-4-alpha.pdf>

Generated by Cheméo on 2025-12-05 07:21:23.231138731 +0000 UTC m=+4667480.761179395.

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