

# cis-1(7),5-p-Menthadien-2-ol

<b>Inchi:</b>	InChI=1S/C10H16O/c1-7(2)9-5-4-8(3)10(11)6-9/h4-5,7,9-11H,3,6H2,1-2H3/t9-,10-/m0/s1
<b>InchiKey:</b>	VKAGFXRPRUAOHV-UWVGGGRQHSA-N
<b>Formula:</b>	C10H16O
<b>SMILES:</b>	C=C1C=CC(C(C)C)CC1O
<b>Mol. weight [g/mol]:</b>	152.23

## Physical Properties

Property code	Value	Unit	Source
gf	-6.16	kJ/mol	Joback Method
hf	-231.24	kJ/mol	Joback Method
hfus	15.19	kJ/mol	Joback Method
hvap	54.72	kJ/mol	Joback Method
log10ws	-2.50		Crippen Method
logp	2.136		Crippen Method
mcvol	138.170	ml/mol	McGowan Method
pc	2918.68	kPa	Joback Method
rinpol	1224.00		NIST Webbook
tb	533.14	K	Joback Method
tc	726.14	K	Joback Method
tf	265.86	K	Joback Method
vc	0.510	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	332.43	J/molxK	533.14	Joback Method
cpg	399.57	J/molxK	693.98	Joback Method
cpg	387.52	J/molxK	661.81	Joback Method
cpg	374.81	J/molxK	629.64	Joback Method
cpg	361.39	J/molxK	597.47	Joback Method
cpg	347.27	J/molxK	565.31	Joback Method
cpg	410.94	J/molxK	726.14	Joback Method
dvisc	0.0001414	Paxs	533.14	Joback Method
dvisc	0.0002220	Paxs	488.59	Joback Method

dvisc	0.0003816	Paxs	444.05	Joback Method
dvisc	0.0007401	Paxs	399.50	Joback Method
dvisc	0.0016951	Paxs	354.95	Joback Method
dvisc	0.0049246	Paxs	310.41	Joback Method
dvisc	0.0204533	Paxs	265.86	Joback Method

## Sources

<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>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R589574&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R589574&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>

## Legend

<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>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>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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