

# Cadina-4,1(10)-dien-7«beta»-ol

<b>Inchi:</b>	InChI=1S/C15H24O/c1-10(2)15(16)8-7-12(4)13-6-5-11(3)9-14(13)15/h9-10,14,16H,5-8H2
<b>InchiKey:</b>	QQRRTCZOIIMRRM-LOACHALJSA-N
<b>Formula:</b>	C15H24O
<b>SMILES:</b>	CC1=CC2C(=C(C)CCC2(O)C(C)C)CC1
<b>Mol. weight [g/mol]:</b>	220.35

## Physical Properties

Property code	Value	Unit	Source
gf	34.80	kJ/mol	Joback Method
hf	-293.09	kJ/mol	Joback Method
hfus	18.02	kJ/mol	Joback Method
hvap	67.21	kJ/mol	Joback Method
log10ws	-4.49		Crippen Method
logp	3.840		Crippen Method
mvol	197.760	ml/mol	McGowan Method
pc	2212.45	kPa	Joback Method
rinpol	1663.00		NIST Webbook
rinpol	1663.00		NIST Webbook
tb	678.40	K	Joback Method
tc	887.06	K	Joback Method
tf	389.41	K	Joback Method
vc	0.741	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	570.40	J/mol×K	678.40	Joback Method
cpg	588.20	J/mol×K	713.18	Joback Method
cpg	605.16	J/mol×K	747.95	Joback Method
cpg	621.42	J/mol×K	782.73	Joback Method
cpg	637.10	J/mol×K	817.50	Joback Method
cpg	652.30	J/mol×K	852.28	Joback Method
cpg	667.17	J/mol×K	887.06	Joback Method

# Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R407284&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R407284&amp;Units=SI</a>
<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>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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>hvp:</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>rinp:</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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