

# 3,4-dimethyl-isochroman, 3e, 4e'

<b>Inchi:</b>	InChI=1S/C11H14O/c1-8-9(2)12-7-10-5-3-4-6-11(8)10/h3-6,8-9H,7H2,1-2H3/t8-,9+/m0/s
<b>InchiKey:</b>	FPFYZFBSRBZHR-DTWKUNHWSA-N
<b>Formula:</b>	C11H14O
<b>SMILES:</b>	CC1OCc2ccccc2C1C
<b>Mol. weight [g/mol]:</b>	162.23

## Physical Properties

Property code	Value	Unit	Source
gf	99.34	kJ/mol	Joback Method
hf	-131.01	kJ/mol	Joback Method
hfus	22.98	kJ/mol	Joback Method
hvap	47.30	kJ/mol	Joback Method
log10ws	-3.05		Crippen Method
logp	2.709		Crippen Method
mcvol	137.100	ml/mol	McGowan Method
pc	2937.70	kPa	Joback Method
rinpol	1287.90		NIST Webbook
ripol	1789.90		NIST Webbook
ripol	1789.90		NIST Webbook
tb	516.03	K	Joback Method
tc	742.26	K	Joback Method
tf	289.42	K	Joback Method
vc	0.512	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	318.39	J/molxK	516.03	Joback Method
cpg	335.94	J/molxK	553.74	Joback Method
cpg	352.37	J/molxK	591.44	Joback Method
cpg	367.75	J/molxK	629.15	Joback Method
cpg	382.12	J/molxK	666.85	Joback Method
cpg	395.53	J/molxK	704.56	Joback Method
cpg	408.03	J/molxK	742.26	Joback Method

dvisc	0.0018157	Paxs	289.42	Joback Method
dvisc	0.0012161	Paxs	327.19	Joback Method
dvisc	0.0008849	Paxs	364.96	Joback Method
dvisc	0.0006835	Paxs	402.73	Joback Method
dvisc	0.0005518	Paxs	440.49	Joback Method
dvisc	0.0004608	Paxs	478.26	Joback Method
dvisc	0.0003951	Paxs	516.03	Joback Method

## 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>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=R256749&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R256749&amp;Units=SI</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>mvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	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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