

# 2-(3-methyl-2-cyclopenten-1-yl)-2-methylpropionamide

<b>Inchi:</b>	InChI=1S/C10H16O/c1-8-4-5-9(6-8)10(2,3)7-11/h6-7,9H,4-5H2,1-3H3
<b>InchiKey:</b>	SVDLNNVLBMXPTF-UHFFFAOYSA-N
<b>Formula:</b>	C10H16O
<b>SMILES:</b>	CC1=CC(C(C)(C)C=O)CC1
<b>Mol. weight [g/mol]:</b>	152.23

## Physical Properties

Property code	Value	Unit	Source
gf	-6.48	kJ/mol	Joback Method
hf	-237.27	kJ/mol	Joback Method
hfus	11.30	kJ/mol	Joback Method
hvap	44.49	kJ/mol	Joback Method
log10ws	-2.55		Crippen Method
logp	2.568		Crippen Method
mcvol	138.170	ml/mol	McGowan Method
pc	2817.33	kPa	Joback Method
ripol	1442.00		NIST Webbook
tb	493.05	K	Joback Method
tc	703.76	K	Joback Method
tf	271.06	K	Joback Method
vc	0.528	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	315.50	J/molxK	493.05	Joback Method
cpg	332.21	J/molxK	528.17	Joback Method
cpg	347.88	J/molxK	563.29	Joback Method
cpg	362.56	J/molxK	598.41	Joback Method
cpg	376.31	J/molxK	633.52	Joback Method
cpg	389.17	J/molxK	668.64	Joback Method
cpg	401.19	J/molxK	703.76	Joback Method
dvisc	0.0039717	Paxs	271.06	Joback Method
dvisc	0.0020900	Paxs	308.06	Joback Method

dvisc	0.0012621	Paxs	345.06	Joback Method
dvisc	0.0008404	Paxs	382.06	Joback Method
dvisc	0.0006012	Paxs	419.05	Joback Method
dvisc	0.0004541	Paxs	456.05	Joback Method
dvisc	0.0003578	Paxs	493.05	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=U365940&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U365940&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>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<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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