

# (1S,2R,4R,7R)-4-Isopropyl-7-methyl-3,8-dioxatricyclo

<b>Inchi:</b>	InChI=1S/C10H16O2/c1-6(2)10-5-4-9(3)7(11-9)8(10)12-10/h6-8H,4-5H2,1-3H3
<b>InchiKey:</b>	LEZWCCRTFNBOBU-UHFFFAOYSA-N
<b>Formula:</b>	C10H16O2
<b>SMILES:</b>	CC(C)C12CCC3(C)OC3C1O2
<b>Mol. weight [g/mol]:</b>	168.23
<b>CAS:</b>	1619-26-7

## Physical Properties

Property code	Value	Unit	Source
gf	34.30	kJ/mol	Joback Method
hf	-284.31	kJ/mol	Joback Method
hfus	19.07	kJ/mol	Joback Method
hvap	43.44	kJ/mol	Joback Method
log10ws	-2.07		Crippen Method
logp	1.731		Crippen Method
mcvol	130.920	ml/mol	McGowan Method
pc	3206.41	kPa	Joback Method
rinpol	1306.00		NIST Webbook
rinpol	1309.90		NIST Webbook
rinpol	1302.00		NIST Webbook
rinpol	1308.00		NIST Webbook
rinpol	1303.00		NIST Webbook
rinpol	1306.00		NIST Webbook
tb	493.42	K	Joback Method
tc	711.67	K	Joback Method
tf	341.50	K	Joback Method
vc	0.505	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	340.55	J/molxK	493.42	Joback Method
cpg	358.15	J/molxK	529.80	Joback Method
cpg	374.04	J/molxK	566.17	Joback Method

cpg	388.50	J/mol×K	602.55	Joback Method
cpg	401.82	J/mol×K	638.92	Joback Method
cpg	414.29	J/mol×K	675.30	Joback Method
cpg	426.18	J/mol×K	711.67	Joback Method

## Sources

<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=C1619267&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1619267&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>

## 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>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

Latest version available from:

<https://www.chemeo.com/cid/71-104-7/1S-2R-4R-7R-4-Isopropyl-7-methyl-3-8-dioxatricyclo-5-1-0-02-4-octane.pdf>

Generated by Cheméo on 2024-04-24 17:16:07.987649442 +0000 UTC m=+16268216.908226764.

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