

2,7-Dimethyl-4-octene

Inchi:	InChI=1S/C10H20/c1-9(2)7-5-6-8-10(3)4/h5-6,9-10H,7-8H2,1-4H3/b6-5+
InchiKey:	LMDKCGFCRGFSDU-AATRIKPKSA-N
Formula:	C10H20
SMILES:	CC(C)CC=CCC(C)C
Mol. weight [g/mol]:	140.27

Physical Properties

Property code	Value	Unit	Source
gf	108.66	kJ/mol	Joback Method
hf	-143.07	kJ/mol	Joback Method
hfus	14.81	kJ/mol	Joback Method
hvap	37.04	kJ/mol	Joback Method
log10ws	-3.38		Crippen Method
logp	3.635		Crippen Method
mcvol	147.460	ml/mol	McGowan Method
pc	2246.13	kPa	Joback Method
rinpol	900.00		NIST Webbook
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tb	431.48	K	Joback Method
tc	609.35	K	Joback Method
tf	167.38	K	Joback Method
vc	0.564	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	300.30	J/mol×K	431.48	Joback Method
cpg	316.10	J/mol×K	461.12	Joback Method
cpg	331.20	J/mol×K	490.77	Joback Method
cpg	345.63	J/mol×K	520.41	Joback Method
cpg	359.41	J/mol×K	550.06	Joback Method
cpg	372.56	J/mol×K	579.70	Joback Method
cpg	385.12	J/mol×K	609.35	Joback Method
dvisc	0.0212319	Pa×s	167.38	Joback Method

dvisc	0.0041931	Pa×s	211.40	Joback Method
dvisc	0.0014484	Pa×s	255.41	Joback Method
dvisc	0.0006839	Pa×s	299.43	Joback Method
dvisc	0.0003914	Pa×s	343.45	Joback Method
dvisc	0.0002543	Pa×s	387.46	Joback Method
dvisc	0.0001804	Pa×s	431.48	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R141241&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log_{10ws}:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
rin_{pol}:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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