

1H-Indene, 2,3-dimethyl-

Inchi:	InChI=1S/C11H12/c1-8-7-10-5-3-4-6-11(10)9(8)2/h3-6H,7H2,1-2H3
InchiKey:	CXLNYETYUQMZKI-UHFFFAOYSA-N
Formula:	C11H12
SMILES:	CC1=C(C)c2ccccc2C1
Mol. weight [g/mol]:	144.21
CAS:	4773-82-4

Physical Properties

Property code	Value	Unit	Source
gf	223.68	kJ/mol	Joback Method
hf	82.67	kJ/mol	Joback Method
hfus	15.41	kJ/mol	Joback Method
hvap	44.86	kJ/mol	Joback Method
log10ws	-3.41		Crippen Method
logp	3.036		Crippen Method
mcvol	126.930	ml/mol	McGowan Method
pc	3124.49	kPa	Joback Method
tb	503.27	K	Joback Method
tc	728.03	K	Joback Method
tf	300.65	K	Joback Method
vc	0.487	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	269.58	J/molxK	503.27	Joback Method
cpg	332.10	J/molxK	690.57	Joback Method
cpg	321.25	J/molxK	653.11	Joback Method
cpg	309.63	J/molxK	615.65	Joback Method
cpg	297.19	J/molxK	578.19	Joback Method
cpg	283.86	J/molxK	540.73	Joback Method
cpg	342.25	J/molxK	728.03	Joback Method
dvisc	0.0003656	Paxs	503.27	Joback Method
dvisc	0.0004135	Paxs	469.50	Joback Method

dvisc	0.0004767	Paxs	435.73	Joback Method
dvisc	0.0005629	Paxs	401.96	Joback Method
dvisc	0.0006853	Paxs	368.19	Joback Method
dvisc	0.0008680	Paxs	334.42	Joback Method
dvisc	0.0011595	Paxs	300.65	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=C4773824&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mccvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/18-546-6/1H-Indene-2-3-dimethyl.pdf>

Generated by Cheméo on 2024-04-24 19:14:29.554584389 +0000 UTC m=+16275318.475161710.

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