

1H-Indene, 3-methyl-

Other names:	1-Methyl-3H-indene 3-Methyl-1H-indene 3-Methylindene Indene, 3-methyl-
Inchi:	InChI=1S/C10H10/c1-8-6-7-9-4-2-3-5-10(8)9/h2-6H,7H2,1H3
InchiKey:	COOKKJGOGWACMY-UHFFFAOYSA-N
Formula:	C10H10
SMILES:	CC1=CCc2ccccc21
Mol. weight [g/mol]:	130.19
CAS:	767-60-2

Physical Properties

Property code	Value	Unit	Source
gf	224.89	kJ/mol	Joback Method
hf	114.78	kJ/mol	Joback Method
hfus	13.21	kJ/mol	Joback Method
hvap	41.97	kJ/mol	Joback Method
ie	7.97	eV	NIST Webbook
ie	8.05	eV	NIST Webbook
log10ws	-2.99		Crippen Method
logp	2.646		Crippen Method
mcvol	112.840	ml/mol	McGowan Method
pc	3551.53	kPa	Joback Method
rinpol	1155.00		NIST Webbook
rinpol	183.80		NIST Webbook
rinpol	184.10		NIST Webbook
rinpol	193.40		NIST Webbook
tb	475.41	K	Joback Method
tc	702.69	K	Joback Method
tf	276.86	K	Joback Method
vc	0.431	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	226.56	J/mol×K	475.41	Joback Method
cpg	240.30	J/mol×K	513.29	Joback Method
cpg	253.04	J/mol×K	551.17	Joback Method
cpg	264.84	J/mol×K	589.05	Joback Method
cpg	275.79	J/mol×K	626.93	Joback Method
cpg	285.94	J/mol×K	664.81	Joback Method
cpg	295.38	J/mol×K	702.69	Joback Method
dvisc	0.0012738	Pa×s	276.86	Joback Method
dvisc	0.0009289	Pa×s	309.95	Joback Method
dvisc	0.0007199	Pa×s	343.04	Joback Method
dvisc	0.0005835	Pa×s	376.13	Joback Method
dvisc	0.0004893	Pa×s	409.23	Joback Method
dvisc	0.0004212	Pa×s	442.32	Joback Method
dvisc	0.0003703	Pa×s	475.41	Joback Method

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.40784e+01
Coeff. B	-3.80216e+03
Coeff. C	-7.21300e+01
Temperature range (K), min.	347.83
Temperature range (K), max.	505.82

Sources

The Yaws Handbook of Vapor
Pressure:
Crippen Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>
<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

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=C767602&Units=SI>

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
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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