

TRICYCLO[3.3.1.1(3,7)]DECANE, 2-METHOXY-

Other names:	Adamantane, 2-methoxy- 2-Methoxyadamantane 2-Adamantyl methyl ether Adamantylmethylether
Inchi:	InChI=1S/C11H18O/c1-12-11-9-3-7-2-8(5-9)6-10(11)4-7/h7-11H,2-6H2,1H3
InchiKey:	FYDHICUFQYUQPE-UHFFFAOYSA-N
Formula:	C11H18O
SMILES:	COC1C2CC3CC(C2)CC1C3
Mol. weight [g/mol]:	166.26

Physical Properties

Property code	Value	Unit	Source
affp	860.20	kJ/mol	NIST Webbook
basg	831.00	kJ/mol	NIST Webbook
hf	-272.79	kJ/mol	Cheméo Relay (1.0)
hfus	19.88	kJ/mol	Joback Method
ie	8.90	eV	Cheméo Relay (1.0)
logp	2.458		Crippen Method
mcvol	139.140	ml/mol	McGowan Method
tf	448.51	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	348.03	J/mol×K	488.65	Joback Method
cpg	369.22	J/mol×K	523.42	Joback Method
cpg	389.13	J/mol×K	558.19	Joback Method
cpg	407.82	J/mol×K	592.96	Joback Method
cpg	425.35	J/mol×K	627.73	Joback Method
cpg	441.81	J/mol×K	662.51	Joback Method
cpg	457.26	J/mol×K	697.28	Joback Method
dvisc	0.0007253	Pa×s	277.78	Joback Method
dvisc	0.0008394	Pa×s	312.92	Joback Method
dvisc	0.0009432	Pa×s	348.07	Joback Method

dvisc	0.0010374	Pa×s	383.21	Joback Method
dvisc	0.0011228	Pa×s	418.36	Joback Method
dvisc	0.0012006	Pa×s	453.50	Joback Method
dvisc	0.0012713	Pa×s	488.65	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C19066230&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

affp:	Proton affinity
basg:	Gas basicity
cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
ie:	Ionization energy
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
tf:	Normal melting (fusion) point

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