

Allyl bromoacetate

Inchi:	InChI=1S/C5H7BrO2/c1-2-3-8-5(7)4-6/h2H,1,3-4H2
InchiKey:	MUWVIMJPXKIXJK-UHFFFAOYSA-N
Formula:	C5H7BrO2
SMILES:	C=CCOC(=O)CBr
Mol. weight [g/mol]:	179.01

Physical Properties

Property code	Value	Unit	Source
gf	-140.54	kJ/mol	Joback Method
hf	-239.57	kJ/mol	Joback Method
hfus	15.50	kJ/mol	Joback Method
hvap	41.64	kJ/mol	Joback Method
log10ws	-1.06		Crippen Method
logp	1.110		Crippen Method
mcvol	101.950	ml/mol	McGowan Method
pc	4211.09	kPa	Joback Method
rinpol	943.00		NIST Webbook
rinpol	943.00		NIST Webbook
rinpol	943.00		NIST Webbook
ripol	1519.00		NIST Webbook
tb	452.93	K	Joback Method
tc	653.94	K	Joback Method
tf	276.31	K	Joback Method
vc	0.383	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	179.89	J/mol×K	452.93	Joback Method
cpg	214.85	J/mol×K	620.44	Joback Method
cpg	208.54	J/mol×K	586.93	Joback Method
cpg	201.90	J/mol×K	553.43	Joback Method
cpg	194.92	J/mol×K	519.93	Joback Method
cpg	187.58	J/mol×K	486.43	Joback Method

cpg	220.83	J/mol×K	653.94	Joback Method
dvisc	0.0003545	Pa×s	452.93	Joback Method
dvisc	0.0004374	Pa×s	423.49	Joback Method
dvisc	0.0005569	Pa×s	394.06	Joback Method
dvisc	0.0007373	Pa×s	364.62	Joback Method
dvisc	0.0010253	Pa×s	335.18	Joback Method
dvisc	0.0015193	Pa×s	305.75	Joback Method
dvisc	0.0024481	Pa×s	276.31	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=R26603&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	Polar retention indices
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/52-509-9/Allyl-bromoacetate.pdf>

Generated by Cheméo on 2025-12-24 09:00:05.935325901 +0000 UTC m=+6315003.465366564.

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