

BICYCLO[3.2.0]HEPT-2-EN-6-ONE, 7,7-DICHLORO-

Other names:	Bicyclo[3.2.0]hept-2-en-6-one, 7,7-dichloro-7,7-dichlorobicyclo[3.2.0]hept-2-en-6-one
Inchi:	InChI=1S/C7H6Cl2O/c8-7(9)5-3-1-2-4(5)6(7)10/h1,3-5H,2H2
InchiKey:	JBPBARAOHIDZPU-UHFFFAOYSA-N
Formula:	C7H6Cl2O
SMILES:	O=C1C2CC=CC2C1(Cl)Cl
Mol. weight [g/mol]:	177.03

Physical Properties

Property code	Value	Unit	Source
hfus	11.95	kJ/mol	Joback Method
logp	1.935		Crippen Method
mcvol	109.520	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	228.72	J/mol×K	514.72	Joback Method
cpg	240.88	J/mol×K	556.24	Joback Method
cpg	251.97	J/mol×K	597.76	Joback Method
cpg	262.16	J/mol×K	639.29	Joback Method
cpg	271.64	J/mol×K	680.81	Joback Method
cpg	280.58	J/mol×K	722.33	Joback Method
cpg	289.17	J/mol×K	763.85	Joback Method

Sources

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=C5307993&Units=SI

Joback Method: https://en.wikipedia.org/wiki/Joback_method
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Legend

cpg: Ideal gas heat capacity
hfus: Enthalpy of fusion at standard conditions
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume

Latest version available from:

<https://www.cheméo.com/cid/68-780-1/BICYCLO-3-2-0-HEPT-2-EN-6-ONE-7-7-DICHLORO.pdf>

Generated by Cheméo on 2026-07-21 22:25:46.933783292 +0000 UTC m=+183842.775668689.

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