

Methylenediformamide

Other names:	Methylenediformamide Formamide, N,N'-methylenebis- Methylenebisformamide
Inchi:	InChI=1S/C3H6N2O2/c6-2-4-1-5-3-7/h2-3H,1H2,(H,4,6)(H,5,7)
InchiKey:	QPJQPYQZFKFTHG-UHFFFAOYSA-N
Formula:	C3H6N2O2
SMILES:	O=CNCNC=O
Mol. weight [g/mol]:	102.09

Physical Properties

Property code	Value	Unit	Source
hfus	18.30	kJ/mol	Joback Method
hvap	77.57	kJ/mol	Cheméo Relay (1.0)
logp	-1.564		Crippen Method
mcvol	76.230	ml/mol	McGowan Method
tb	529.50	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	158.15	J/mol×K	465.70	Joback Method
cpg	164.68	J/mol×K	497.60	Joback Method
cpg	170.87	J/mol×K	529.51	Joback Method
cpg	176.75	J/mol×K	561.41	Joback Method
cpg	182.32	J/mol×K	593.31	Joback Method
cpg	187.58	J/mol×K	625.22	Joback Method
cpg	192.55	J/mol×K	657.12	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=C6921988&Units=SI>
Joback Method: https://en.wikipedia.org/wiki/Joback_method

Legend

cpg: Ideal gas heat capacity
hfus: Enthalpy of fusion at standard conditions
hvap: Enthalpy of vaporization at standard conditions
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
tb: Normal Boiling Point Temperature

Latest version available from:

<https://www.chemeo.com/cid/68-123-0/Methylenediformamide.pdf>

Generated by Cheméo on 2026-07-21 23:52:39.162604202 +0000 UTC m=+189055.004489589.

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