

Butylone

Inchi:	InChI=1S/C12H15NO3/c1-3-9(13-2)12(14)8-4-5-10-11(6-8)16-7-15-10/h4-6,9,13H,3,7H2
InchiKey:	CGKQZIULZRXRRJ-UHFFFAOYSA-N
Formula:	C12H15NO3
SMILES:	CCC(NC)C(=O)c1ccc2c(c1)OCO2
Mol. weight [g/mol]:	221.25
CAS:	802575-11-7

Physical Properties

Property code	Value	Unit	Source
gf	-2.44	kJ/mol	Joback Method
hf	-312.67	kJ/mol	Joback Method
hfus	36.30	kJ/mol	Joback Method
hvap	67.94	kJ/mol	Joback Method
log10ws	-2.90		Crippen Method
logp	1.596		Crippen Method
mcvol	168.610	ml/mol	McGowan Method
pc	2953.69	kPa	Joback Method
rinpol	1803.80		NIST Webbook
tb	679.51	K	Joback Method
tc	903.50	K	Joback Method
tf	439.37	K	Joback Method
vc	0.634	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	461.49	J/mol×K	679.51	Joback Method
cpg	475.08	J/mol×K	716.84	Joback Method
cpg	487.72	J/mol×K	754.17	Joback Method
cpg	499.50	J/mol×K	791.50	Joback Method
cpg	510.47	J/mol×K	828.84	Joback Method
cpg	520.70	J/mol×K	866.17	Joback Method
cpg	530.27	J/mol×K	903.50	Joback Method

Sources

Crippen Method:	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=C802575117&Units=SI

Legend

cpg:	Ideal gas heat capacity
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
rlnpol:	Non-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/94-093-5/Butylone.pdf>

Generated by Cheméo on 2025-12-23 06:45:10.270756065 +0000 UTC m=+6220507.800796716.

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