

1-Sulfanyl-3-butyl acetate

Inchi:	InChI=1S/C6H12O2S/c1-5(3-4-9)8-6(2)7/h5,9H,3-4H2,1-2H3
InchiKey:	PPGZGGPLMMKFKW-UHFFFAOYSA-N
Formula:	C6H12O2S
SMILES:	CC(=O)OC(C)CCS
Mol. weight [g/mol]:	148.22

Physical Properties

Property code	Value	Unit	Source
gf	-207.33	kJ/mol	Joback Method
hf	-378.77	kJ/mol	Joback Method
hfus	14.60	kJ/mol	Joback Method
hvap	44.45	kJ/mol	Joback Method
log10ws	-1.38		Crippen Method
logp	1.258		Crippen Method
mcvol	119.190	ml/mol	McGowan Method
pc	3560.02	kPa	Joback Method
rinpol	1024.00		NIST Webbook
rinpol	1024.00		NIST Webbook
ripol	1542.00		NIST Webbook
ripol	1542.00		NIST Webbook
tb	475.39	K	Joback Method
tc	679.41	K	Joback Method
tf	251.00	K	Joback Method
vc	0.444	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	244.02	J/molxK	475.39	Joback Method
cpg	254.95	J/molxK	509.39	Joback Method
cpg	265.43	J/molxK	543.40	Joback Method
cpg	275.45	J/molxK	577.40	Joback Method
cpg	285.02	J/molxK	611.41	Joback Method
cpg	294.14	J/molxK	645.41	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=R621188&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
mccvol:	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

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