

Succinic acid, but-3-yn-2-yl 2-heptyl ester

Inchi:	InChI=1S/C15H24O4/c1-5-7-8-9-13(4)19-15(17)11-10-14(16)18-12(3)6-2/h2,12-13H,5,7-
InchiKey:	XKJGVYJDZALHGR-UHFFFAOYSA-N
Formula:	C15H24O4
SMILES:	<chem>C#CC(C)OC(=O)CCC(=O)OC(C)CCCC</chem>
Mol. weight [g/mol]:	268.35

Physical Properties

Property code	Value	Unit	Source
gf	-174.23	kJ/mol	Joback Method
hf	-561.19	kJ/mol	Joback Method
hfus	36.11	kJ/mol	Joback Method
hvap	66.38	kJ/mol	Joback Method
log10ws	-3.84		Crippen Method
logp	2.843		Crippen Method
mvol	228.490	ml/mol	McGowan Method
pc	1733.22	kPa	Joback Method
rinpol	1680.00		NIST Webbook
tb	684.42	K	Joback Method
tc	872.41	K	Joback Method
tf	420.10	K	Joback Method
vc	0.874	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	629.90	J/mol×K	684.42	Joback Method
cpg	645.54	J/mol×K	715.75	Joback Method
cpg	660.36	J/mol×K	747.08	Joback Method
cpg	674.38	J/mol×K	778.42	Joback Method
cpg	687.61	J/mol×K	809.75	Joback Method
cpg	700.07	J/mol×K	841.08	Joback Method
cpg	711.75	J/mol×K	872.41	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U390567&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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

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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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