

Succinic acid, hexyl pent-4-enyl ester

Inchi:	InChI=1S/C15H26O4/c1-3-5-7-9-13-19-15(17)11-10-14(16)18-12-8-6-4-2/h4H,2-3,5-13H
InchiKey:	RXWMKNMUSIQHRD-UHFFFAOYSA-N
Formula:	C15H26O4
SMILES:	<chem>C=CCCCOC(=O)CCC(=O)OCCCCC</chem>
Mol. weight [g/mol]:	270.36

Physical Properties

Property code	Value	Unit	Source
gf	-304.58	kJ/mol	Joback Method
hf	-717.10	kJ/mol	Joback Method
hfus	38.90	kJ/mol	Joback Method
hvap	66.63	kJ/mol	Joback Method
log10ws	-3.68		Crippen Method
logp	3.399		Crippen Method
mvol	232.790	ml/mol	McGowan Method
pc	1569.72	kPa	Joback Method
rinpol	1839.00		NIST Webbook
rinpol	1839.00		NIST Webbook
tb	691.86	K	Joback Method
tc	869.60	K	Joback Method
tf	401.37	K	Joback Method
vc	0.904	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	653.38	J/molxK	691.86	Joback Method
cpg	668.97	J/molxK	721.48	Joback Method
cpg	683.81	J/molxK	751.11	Joback Method
cpg	697.90	J/molxK	780.73	Joback Method
cpg	711.24	J/molxK	810.35	Joback Method
cpg	723.86	J/molxK	839.98	Joback Method
cpg	735.75	J/molxK	869.60	Joback Method
dvisc	0.0013617	Paxs	401.37	Joback Method

dvisc	0.0007165	Paxs	449.78	Joback Method
dvisc	0.0004271	Paxs	498.20	Joback Method
dvisc	0.0002791	Paxs	546.62	Joback Method
dvisc	0.0001954	Paxs	595.03	Joback Method
dvisc	0.0001444	Paxs	643.44	Joback Method
dvisc	0.0001113	Paxs	691.86	Joback Method

Sources

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=U353372&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
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
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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