

Succinic acid, 2-methylhex-3-yl 2-nitrophenethyl ester

Inchi: InChI=1S/C19H27NO6/c1-4-7-17(14(2)3)26-19(22)11-10-18(21)25-13-12-15-8-5-6-9-16(
InchiKey: VTTHLSPAYIESTD-UHFFFAOYSA-N
Formula: C19H27NO6
SMILES: CCCC(OC(=O)CCC(=O)OCCc1ccccc1[N+](=O)[O-])C(C)C
Mol. weight [g/mol]: 365.42

Physical Properties

Property code	Value	Unit	Source
gf	-225.29	kJ/mol	Joback Method
hf	-721.35	kJ/mol	Joback Method
hfus	48.51	kJ/mol	Joback Method
hvap	94.95	kJ/mol	Joback Method
log10ws	-5.12		Crippen Method
logp	3.829		Crippen Method
mcvol	287.110	ml/mol	McGowan Method
pc	1472.49	kPa	Joback Method
rinpol	2569.00		NIST Webbook
rinpol	2569.00		NIST Webbook
tb	969.32	K	Joback Method
tc	1195.93	K	Joback Method
tf	600.76	K	Joback Method
vc	1.109	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	930.05	J/molxK	969.32	Joback Method
cpg	942.58	J/molxK	1007.09	Joback Method
cpg	953.75	J/molxK	1044.86	Joback Method
cpg	963.59	J/molxK	1082.62	Joback Method
cpg	972.14	J/molxK	1120.39	Joback Method
cpg	979.43	J/molxK	1158.16	Joback Method
cpg	985.50	J/molxK	1195.93	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.cheméo.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=U381272&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
hvp:	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
rinp:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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