

Succinic acid, tridecyl 2,4,4-trimethylpentyl ester

Inchi:	InChI=1S/C25H48O4/c1-6-7-8-9-10-11-12-13-14-15-16-19-28-23(26)17-18-24(27)29-21-
InchiKey:	NVLRWCNBFlyTJS-UHFFFAOYSA-N
Formula:	C25H48O4
SMILES:	CCCCCCCCCCCCOC(=O)CCC(=O)OCC(C)CC(C)(C)C
Mol. weight [g/mol]:	412.65

Physical Properties

Property code	Value	Unit	Source
gf	-307.82	kJ/mol	Joback Method
hf	-1062.96	kJ/mol	Joback Method
hfus	55.14	kJ/mol	Joback Method
hvap	87.87	kJ/mol	Joback Method
log10ws	-7.53		Crippen Method
logp	7.236		Crippen Method
mvol	377.990	ml/mol	McGowan Method
pc	816.79	kPa	Joback Method
rinpol	2707.00		NIST Webbook
rinpol	2707.00		NIST Webbook
tb	920.31	K	Joback Method
tc	1127.45	K	Joback Method
tf	503.25	K	Joback Method
vc	1.466	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1281.45	J/molxK	920.31	Joback Method
cpg	1301.73	J/molxK	954.83	Joback Method
cpg	1320.61	J/molxK	989.36	Joback Method
cpg	1338.12	J/molxK	1023.88	Joback Method
cpg	1354.33	J/molxK	1058.41	Joback Method
cpg	1369.31	J/molxK	1092.93	Joback Method
cpg	1383.09	J/molxK	1127.45	Joback Method
dvisc	0.0004880	Paxs	503.25	Joback Method

dvisc	0.0002009	Paxs	572.76	Joback Method
dvisc	0.0001003	Paxs	642.27	Joback Method
dvisc	0.0000573	Paxs	711.78	Joback Method
dvisc	0.0000362	Paxs	781.29	Joback Method
dvisc	0.0000246	Paxs	850.80	Joback Method
dvisc	0.0000178	Paxs	920.31	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U381329&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

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_{cvol}:	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

Latest version available from:

<https://www.chemeo.com/cid/86-707-2/Succinic-acid-tridecyl-2-4-4-trimethylpentyl-ester.pdf>

Generated by Cheméo on 2024-05-06 14:39:28.358609036 +0000 UTC m=+17295617.279186353.

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