

Sebacic acid, hexyl non-5-yn-3-yl ester

Inchi:	InChI=1S/C25H44O4/c1-4-7-9-15-19-23(6-3)29-25(27)21-17-14-12-11-13-16-20-24(26)2
InchiKey:	NBMHTGWWIVEXQD-UHFFFAOYSA-N
Formula:	C25H44O4
SMILES:	CCCC#CCC(CC)OC(=O)CCCCCCCCC(=O)OCCCCC
Mol. weight [g/mol]:	408.61

Physical Properties

Property code	Value	Unit	Source
gf	-107.86	kJ/mol	Joback Method
hf	-781.91	kJ/mol	Joback Method
hfus	65.68	kJ/mol	Joback Method
hvap	91.32	kJ/mol	Joback Method
log10ws	-7.92		Crippen Method
logp	6.746		Crippen Method
mcvol	369.390	ml/mol	McGowan Method
pc	886.83	kPa	Joback Method
rinsol	2778.00		NIST Webbook
tb	932.54	K	Joback Method
tc	1141.86	K	Joback Method
tf	606.93	K	Joback Method
vc	1.440	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1227.96	J/mol×K	932.54	Joback Method
cpg	1247.04	J/mol×K	967.43	Joback Method
cpg	1264.68	J/mol×K	1002.31	Joback Method
cpg	1280.93	J/mol×K	1037.20	Joback Method
cpg	1295.80	J/mol×K	1072.09	Joback Method
cpg	1309.35	J/mol×K	1106.98	Joback Method
cpg	1321.59	J/mol×K	1141.86	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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=U355802&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
h vap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 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

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