

Sebacic acid, ethyl 2-methylbenzyl ester

Inchi:	InChI=1S/C20H30O4/c1-3-23-19(21)14-8-6-4-5-7-9-15-20(22)24-16-18-13-11-10-12-17(
InchiKey:	XVCNNQWXKOYVEW-UHFFFAOYSA-N
Formula:	C20H30O4
SMILES:	CCOC(=O)CCCCCCCC(=O)OCc1ccccc1C
Mol. weight [g/mol]:	334.45

Physical Properties

Property code	Value	Unit	Source
gf	-247.54	kJ/mol	Joback Method
hf	-720.67	kJ/mol	Joback Method
hfus	46.78	kJ/mol	Joback Method
hvap	81.36	kJ/mol	Joback Method
log10ws	-5.57		Crippen Method
logp	4.722		Crippen Method
mvol	283.780	ml/mol	McGowan Method
pc	1341.76	kPa	Joback Method
rinpol	2471.00		NIST Webbook
rinpol	2471.00		NIST Webbook
tb	841.24	K	Joback Method
tc	1041.25	K	Joback Method
tf	498.42	K	Joback Method
vc	1.095	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	877.72	J/molxK	841.24	Joback Method
cpg	893.85	J/molxK	874.57	Joback Method
cpg	908.85	J/molxK	907.91	Joback Method
cpg	922.75	J/molxK	941.24	Joback Method
cpg	935.56	J/molxK	974.58	Joback Method
cpg	947.32	J/molxK	1007.91	Joback Method
cpg	958.03	J/molxK	1041.25	Joback Method
dvisc	0.0005930	Paxs	498.42	Joback Method

dvisc	0.0003251	Paxs	555.56	Joback Method
dvisc	0.0001994	Paxs	612.69	Joback Method
dvisc	0.0001329	Paxs	669.83	Joback Method
dvisc	0.0000945	Paxs	726.97	Joback Method
dvisc	0.0000705	Paxs	784.10	Joback Method
dvisc	0.0000548	Paxs	841.24	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U380712&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
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

Latest version available from:

<https://www.chemeo.com/cid/87-665-8/Sebacic-acid-ethyl-2-methylbenzyl-ester.pdf>

Generated by Cheméo on 2024-04-28 23:19:40.165669288 +0000 UTC m=+16635629.086246649.

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