

L-Methionine, N-heptafluorobutyryl-, undecyl ester

Inchi:	InChI=1S/C20H32F7NO3S/c1-3-4-5-6-7-8-9-10-11-13-31-16(29)15(12-14-32-2)28-17(30)
InchiKey:	NFPQZRCTTDPBW-UHFFFAOYSA-N
Formula:	C20H32F7NO3S
SMILES:	CCCCCCCCCCCCOC(=O)C(CCSC)NC(=O)C(F)(F)C(F)(F)C(F)(F)F
Mol. weight [g/mol]:	499.53

Physical Properties

Property code	Value	Unit	Source
hfus	56.97	kJ/mol	Joback Method
logp	6.131		Crippen Method
mcvol	340.390	ml/mol	McGowan Method
rinpol	2279.00		NIST Webbook
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tb	618.89	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1119.10	J/mol×K	890.87	Joback Method
cpg	1134.48	J/mol×K	924.27	Joback Method
cpg	1148.81	J/mol×K	957.67	Joback Method
cpg	1162.18	J/mol×K	991.06	Joback Method
cpg	1174.66	J/mol×K	1024.46	Joback Method
cpg	1186.36	J/mol×K	1057.86	Joback Method
cpg	1197.34	J/mol×K	1091.26	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	

Cheméo Relay (1.0, beta):

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=U320858&Units=SI>

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
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature

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