



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:35 AM GMT

PDB ID : 1K86
Title : Crystal structure of caspase-7
Authors : Chai, J.; Wu, Q.; Shiozaki, E.; Srinivasa, S.M.; Alnemri, E.S.; Shi, Y.
Deposited on : 2001-10-23
Resolution : 2.60 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

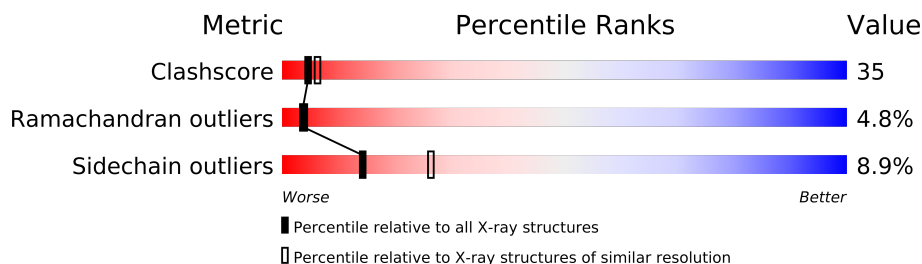
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	253	
1	B	253	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3755 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called caspase-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1853	1177	316	345	15			
1	B	232	Total	C	N	O	S	0	0	0
			1853	1177	316	345	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	ALA	ASP	ENGINEERED	UNP P55210
B	169	ALA	ASP	ENGINEERED	UNP P55210

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	14	Total	O	0	0
			14	14		
2	B	35	Total	O	0	0
			35	35		

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.90Å 88.90Å 186.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.60)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3755	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.55	2/1892 (0.1%)	1.18	18/2546 (0.7%)
1	B	0.39	0/1892	0.69	1/2546 (0.0%)
All	All	0.48	2/3784 (0.1%)	0.97	19/5092 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	272	HIS	N-CA	-7.24	1.31	1.46
1	A	213	ILE	N-CA	-6.50	1.33	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	GLY	CA-C-O	-22.76	79.63	120.60
1	A	188	GLY	CA-C-N	20.61	162.53	117.20
1	A	213	ILE	C-N-CD	-17.56	81.96	120.60
1	A	188	GLY	O-C-N	-16.76	95.88	122.70
1	A	215	VAL	N-CA-C	-12.40	77.52	111.00
1	A	213	ILE	N-CA-C	-11.76	79.25	111.00
1	A	213	ILE	C-N-CA	8.33	156.99	122.00
1	A	214	PRO	N-CA-C	-8.22	90.74	112.10
1	B	189	THR	N-CA-C	6.83	129.45	111.00
1	A	271	ARG	CA-C-N	-6.75	102.34	117.20
1	A	214	PRO	CA-CB-CG	-6.71	91.25	104.00
1	A	272	HIS	CB-CA-C	6.65	123.70	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	GLU	N-CA-C	-6.58	93.23	111.00
1	A	214	PRO	CA-N-CD	-6.48	102.42	111.50
1	A	212	LYS	CA-C-N	-6.32	103.30	117.20
1	A	271	ARG	N-CA-C	-6.02	94.75	111.00
1	A	213	ILE	N-CA-CB	-5.77	97.52	110.80
1	A	152	GLY	N-CA-C	-5.54	99.25	113.10
1	A	58	TYR	N-CA-C	5.20	125.04	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	GLY	Mainchain

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1853	0	1818	156	0
1	B	1853	0	1818	117	0
2	A	14	0	0	0	0
2	B	35	0	0	1	0
All	All	3755	0	3636	257	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 35.

All (257) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:213:ILE:CG1	1:A:214:PRO:HD3	1.80	1.11
1:A:195:ILE:HG13	1:A:196:GLN:H	1.17	1.06
1:A:213:ILE:HG13	1:A:214:PRO:HD3	1.06	1.02
1:A:214:PRO:HD2	1:A:221:PHE:HZ	1.23	0.99
1:A:214:PRO:HD2	1:A:221:PHE:CZ	1.98	0.98
1:A:160:LYS:HA	1:A:213:ILE:HG12	1.45	0.97

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:215:VAL:HG12	1:B:226:VAL:HB	1.48	0.95
1:A:213:ILE:HG13	1:A:214:PRO:CD	1.96	0.95
1:B:258:ILE:N	1:B:258:ILE:HD13	1.83	0.94
1:B:272:HIS:ND1	1:B:273:PHE:N	2.20	0.89
1:B:133:ASN:HD22	1:B:133:ASN:H	1.25	0.83
1:B:258:ILE:CD1	1:B:299:LEU:HB3	2.11	0.81
1:B:258:ILE:HD13	1:B:258:ILE:H	1.42	0.80
1:A:215:VAL:HG21	1:B:229:TYR:CE1	2.15	0.80
1:B:237:ARG:HD2	1:B:237:ARG:H	1.46	0.80
1:A:192:ASP:HB2	1:A:287:GLN:HE22	1.44	0.79
1:B:131:HIS:H	1:B:173:THR:HG21	1.45	0.79
1:B:226:VAL:HG23	1:B:288:ILE:O	1.82	0.79
1:A:191:LEU:H	1:A:191:LEU:HD12	1.48	0.78
1:B:131:HIS:N	1:B:173:THR:HG21	1.98	0.78
1:A:225:THR:HG22	1:A:226:VAL:N	1.99	0.78
1:B:258:ILE:CD1	1:B:258:ILE:H	1.96	0.78
1:B:234:SER:HB3	1:B:237:ARG:CD	2.15	0.77
1:A:195:ILE:HG13	1:A:196:GLN:N	1.98	0.77
1:A:191:LEU:HD22	1:A:285:LYS:HG3	1.67	0.77
1:B:296:THR:O	1:B:297:LYS:HG2	1.85	0.76
1:A:275:SER:HB3	1:A:285:LYS:O	1.87	0.75
1:B:258:ILE:HD12	1:B:299:LEU:HB3	1.69	0.74
1:B:272:HIS:CG	1:B:273:PHE:H	2.04	0.74
1:A:120:GLN:NE2	1:A:162:LEU:HD23	2.04	0.73
1:B:258:ILE:N	1:B:258:ILE:CD1	2.51	0.73
1:A:274:GLU:HA	1:A:286:LYS:NZ	2.05	0.71
1:B:237:ARG:HD2	1:B:237:ARG:N	2.05	0.71
1:A:191:LEU:CD2	1:A:285:LYS:HG3	2.21	0.70
1:B:133:ASN:H	1:B:133:ASN:ND2	1.90	0.69
1:A:192:ASP:H	1:A:285:LYS:HB3	1.57	0.69
1:A:149:VAL:HG22	1:A:156:VAL:HG22	1.73	0.68
1:B:233:ARG:HA	1:B:239:SER:HA	1.75	0.68
1:B:279:ASP:OD1	1:B:280:PRO:HD2	1.94	0.67
1:B:226:VAL:HG23	1:B:288:ILE:HG23	1.76	0.66
1:A:257:GLU:OE2	1:A:258:ILE:HG22	1.95	0.65
1:A:271:ARG:HG3	1:A:272:HIS:HB2	1.77	0.65
1:B:269:VAL:HG21	1:B:289:PRO:CD	2.28	0.64
1:A:118:LYS:O	1:A:122:LEU:HB2	1.98	0.64
1:A:191:LEU:N	1:A:191:LEU:HD12	2.13	0.64
1:A:184:GLN:HG3	1:A:184:GLN:O	1.97	0.64
1:B:136:CYS:HB3	1:B:178:PRO:HG2	1.80	0.64
1:A:189:THR:O	1:A:190:GLU:HB2	1.98	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:269:VAL:HG21	1:B:289:PRO:HD3	1.80	0.62
1:A:272:HIS:O	1:A:273:PHE:C	2.33	0.62
1:B:258:ILE:HD11	1:B:299:LEU:HB3	1.81	0.62
1:A:274:GLU:HA	1:A:286:LYS:CD	2.30	0.61
1:B:133:ASN:HD22	1:B:133:ASN:N	1.90	0.61
1:A:163:THR:HB	1:A:214:PRO:HG2	1.81	0.61
1:B:167:ARG:HB3	1:B:216:GLU:OE2	2.00	0.61
1:A:275:SER:O	1:A:284:GLU:HA	2.02	0.60
1:A:160:LYS:HA	1:A:213:ILE:CG1	2.25	0.60
1:A:225:THR:O	1:A:226:VAL:HB	2.01	0.59
1:B:234:SER:HB3	1:B:237:ARG:HD2	1.84	0.59
1:B:228:GLY:C	1:B:229:TYR:HD1	2.04	0.59
1:A:213:ILE:CB	1:A:214:PRO:HD3	2.26	0.59
1:A:62:MET:HB2	1:A:303:GLN:HA	1.84	0.59
1:A:233:ARG:HG3	1:A:234:SER:N	2.18	0.58
1:A:231:SER:HA	1:A:287:GLN:HG3	1.86	0.58
1:A:276:GLN:HG2	1:A:277:SER:H	1.67	0.58
1:A:120:GLN:HE21	1:A:157:THR:HG21	1.68	0.58
1:B:244:ALA:O	1:B:248:ILE:HG12	2.04	0.58
1:B:186:CYS:SG	1:B:231:SER:HB3	2.43	0.58
1:B:226:VAL:CG2	1:B:288:ILE:HG23	2.33	0.58
1:A:192:ASP:CB	1:A:287:GLN:HE22	2.15	0.58
1:A:62:MET:HB3	1:A:303:GLN:HG2	1.85	0.58
1:B:163:THR:HG22	1:B:181:PHE:CE2	2.39	0.57
1:B:269:VAL:HG21	1:B:289:PRO:CG	2.34	0.57
1:A:215:VAL:O	1:A:216:GLU:HG2	2.04	0.57
1:A:160:LYS:H	1:A:160:LYS:HZ3	1.51	0.57
1:A:233:ARG:CB	1:A:239:SER:HA	2.35	0.57
1:A:167:ARG:HG3	1:A:216:GLU:CG	2.35	0.57
1:A:274:GLU:HA	1:A:286:LYS:HD3	1.86	0.57
1:A:64:PHE:HB3	1:A:133:ASN:O	2.04	0.57
1:A:229:TYR:HE1	1:B:215:VAL:O	1.86	0.57
1:B:131:HIS:H	1:B:173:THR:CG2	2.16	0.57
1:A:97:LEU:HD13	1:A:140:ILE:HG21	1.86	0.57
1:A:239:SER:O	1:A:243:GLN:HG3	2.05	0.56
1:A:150:ILE:HD11	1:A:162:LEU:HD11	1.86	0.56
1:A:160:LYS:CA	1:A:213:ILE:HG12	2.28	0.56
1:A:60:TYR:CD1	1:A:178:PRO:HD3	2.40	0.56
1:A:233:ARG:HB2	1:A:239:SER:HA	1.87	0.56
1:A:227:PRO:O	1:B:212:LYS:HD2	2.06	0.55
1:A:145:GLY:HA3	1:A:185:ALA:HB1	1.88	0.55
1:A:225:THR:HG22	1:A:226:VAL:H	1.70	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:62:MET:CB	1:A:303:GLN:HA	2.37	0.54
1:A:215:VAL:HG21	1:B:229:TYR:HE1	1.70	0.54
1:A:271:ARG:HG3	1:A:272:HIS:N	2.23	0.54
1:A:229:TYR:CE1	1:B:215:VAL:O	2.61	0.54
1:A:167:ARG:HG3	1:A:216:GLU:HG3	1.90	0.54
1:A:213:ILE:CG1	1:A:214:PRO:CD	2.70	0.54
1:A:163:THR:HG22	1:A:181:PHE:CE2	2.43	0.54
1:A:226:VAL:HG11	1:B:215:VAL:HB	1.88	0.54
1:A:212:LYS:NZ	1:B:226:VAL:HG13	2.23	0.54
1:B:234:SER:HB2	1:B:238:GLY:O	2.09	0.53
1:A:163:THR:CB	1:A:214:PRO:HG2	2.39	0.53
1:A:294:MET:SD	1:B:290:CYS:HB2	2.48	0.53
1:B:191:LEU:HG	1:B:281:HIS:HB3	1.90	0.53
1:A:226:VAL:HG12	1:A:229:TYR:HB2	1.89	0.53
1:A:192:ASP:HB2	1:A:287:GLN:NE2	2.17	0.53
1:A:275:SER:H	1:A:286:LYS:HZ2	1.57	0.52
1:A:229:TYR:CZ	1:B:215:VAL:HG23	2.44	0.52
1:B:184:GLN:HE22	1:B:231:SER:CB	2.22	0.52
1:B:146:GLU:O	1:B:147:GLU:C	2.47	0.52
1:A:226:VAL:HG23	1:A:290:CYS:HB2	1.90	0.52
1:A:167:ARG:HD3	1:A:214:PRO:HB3	1.92	0.52
1:B:113:ASP:HB3	1:B:153:LYS:HE2	1.90	0.52
1:B:124:LYS:O	1:B:128:GLU:HG3	2.10	0.52
1:A:229:TYR:CD1	1:A:287:GLN:OE1	2.63	0.52
1:A:87:ARG:O	1:A:90:THR:HG22	2.11	0.51
1:B:277:SER:HB3	1:B:282:PHE:O	2.10	0.51
1:B:174:LEU:HD23	1:B:177:LYS:HD2	1.92	0.51
1:B:160:LYS:HA	1:B:214:PRO:HD3	1.93	0.51
1:A:111:TYR:CG	1:A:122:LEU:HD21	2.46	0.51
1:A:167:ARG:NH2	1:A:213:ILE:HD11	2.26	0.50
1:A:271:ARG:CG	1:A:272:HIS:N	2.73	0.50
1:B:163:THR:OG1	1:B:214:PRO:HG2	2.12	0.50
1:A:270:ALA:HB2	1:A:288:ILE:CD1	2.41	0.50
1:A:276:GLN:HA	1:A:284:GLU:H	1.76	0.50
1:B:93:ASP:HB3	1:B:242:VAL:HG11	1.94	0.50
1:B:237:ARG:CD	1:B:237:ARG:N	2.72	0.49
1:B:275:SER:O	1:B:277:SER:N	2.45	0.49
1:A:191:LEU:HB2	1:A:285:LYS:CD	2.43	0.49
1:B:132:THR:N	1:B:173:THR:CG2	2.75	0.49
1:A:225:THR:HA	1:A:231:SER:HB3	1.94	0.49
1:B:184:GLN:HE22	1:B:231:SER:HB3	1.78	0.49
1:A:191:LEU:HB2	1:A:285:LYS:HD2	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:277:SER:HB3	1:B:283:HIS:HA	1.95	0.49
1:A:60:TYR:CD2	1:A:297:LYS:HB2	2.48	0.49
1:B:274:GLU:HB3	1:B:284:GLU:OE2	2.13	0.49
1:B:136:CYS:CB	1:B:178:PRO:HG2	2.41	0.49
1:A:276:GLN:HG2	1:A:277:SER:N	2.27	0.48
1:A:274:GLU:HA	1:A:286:LYS:HZ2	1.77	0.48
1:A:70:CYS:HA	1:A:138:ALA:O	2.12	0.48
1:B:112:ASN:HB3	2:B:6:HOH:O	2.12	0.48
1:A:191:LEU:HD11	1:A:282:PHE:HD1	1.78	0.48
1:B:131:HIS:N	1:B:173:THR:CG2	2.73	0.48
1:B:274:GLU:HB3	1:B:284:GLU:HG2	1.94	0.48
1:B:269:VAL:CG2	1:B:289:PRO:HD3	2.44	0.48
1:A:270:ALA:C	1:A:271:ARG:O	2.47	0.47
1:B:234:SER:CB	1:B:237:ARG:HD3	2.44	0.47
1:A:240:TRP:CZ3	1:A:273:PHE:HB3	2.49	0.47
1:A:191:LEU:HD22	1:A:285:LYS:HE3	1.95	0.47
1:A:116:CYS:HB3	1:A:157:THR:OG1	2.14	0.47
1:A:272:HIS:O	1:A:273:PHE:O	2.32	0.47
1:A:234:SER:C	1:A:236:GLY:H	2.18	0.47
1:A:160:LYS:H	1:A:160:LYS:NZ	2.13	0.47
1:B:219:PHE:O	1:B:294:MET:HB2	2.15	0.47
1:A:168:GLY:N	1:A:216:GLU:HG3	2.30	0.47
1:A:274:GLU:O	1:A:274:GLU:HG2	2.13	0.47
1:A:175:LEU:O	1:A:176:GLU:HB2	2.14	0.46
1:B:191:LEU:HG	1:B:281:HIS:CB	2.45	0.46
1:A:168:GLY:H	1:A:216:GLU:HG3	1.81	0.46
1:B:95:GLU:O	1:B:99:LYS:HG3	2.14	0.46
1:B:260:GLN:O	1:B:264:ARG:HG3	2.15	0.46
1:B:59:GLN:CB	1:B:298:GLU:HB2	2.46	0.46
1:A:84:MET:HB3	1:A:144:HIS:CD2	2.51	0.46
1:B:213:ILE:HB	1:B:214:PRO:HD2	1.98	0.46
1:A:226:VAL:HG13	1:B:215:VAL:HG11	1.96	0.46
1:A:115:SER:HA	1:A:153:LYS:HG2	1.97	0.46
1:A:121:ASP:OD2	1:A:125:LYS:HE3	2.15	0.46
1:A:167:ARG:HG3	1:A:216:GLU:HG2	1.98	0.46
1:A:240:TRP:CE3	1:A:273:PHE:HB3	2.51	0.45
1:A:184:GLN:HA	1:A:224:SER:HB3	1.99	0.45
1:B:267:ASP:O	1:B:270:ALA:HB3	2.16	0.45
1:A:74:ASN:ND2	1:A:74:ASN:C	2.70	0.45
1:B:163:THR:OG1	1:B:214:PRO:CD	2.63	0.45
1:A:270:ALA:HB2	1:A:288:ILE:HD13	1.97	0.45
1:B:232:TRP:HE3	1:B:233:ARG:N	2.14	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:251:GLU:HG2	1:A:252:HIS:CD2	2.52	0.45
1:A:193:ASP:N	1:A:285:LYS:HA	2.32	0.45
1:B:228:GLY:C	1:B:229:TYR:CD1	2.89	0.45
1:B:245:LEU:O	1:B:249:LEU:HB2	2.17	0.45
1:B:59:GLN:HB3	1:B:298:GLU:HB2	1.97	0.45
1:B:59:GLN:C	1:B:297:LYS:HE3	2.37	0.45
1:B:277:SER:CB	1:B:282:PHE:O	2.65	0.45
1:A:225:THR:CG2	1:A:226:VAL:N	2.68	0.45
1:B:212:LYS:O	1:B:213:ILE:HG23	2.16	0.45
1:A:74:ASN:C	1:A:74:ASN:HD22	2.18	0.45
1:B:272:HIS:CG	1:B:273:PHE:N	2.77	0.45
1:A:118:LYS:HB2	1:A:118:LYS:NZ	2.32	0.45
1:A:278:ASP:OD1	1:A:278:ASP:O	2.34	0.44
1:A:215:VAL:HG21	1:B:229:TYR:CZ	2.51	0.44
1:B:226:VAL:HG22	1:B:290:CYS:HB3	2.00	0.44
1:A:226:VAL:HG22	1:A:227:PRO:HD2	2.00	0.44
1:A:149:VAL:HG23	1:A:157:THR:C	2.38	0.44
1:B:230:TYR:C	1:B:230:TYR:CD1	2.90	0.44
1:A:267:ASP:HB2	1:B:296:THR:O	2.17	0.44
1:B:192:ASP:HB3	1:B:193:ASP:H	1.66	0.44
1:A:136:CYS:CB	1:A:178:PRO:HG2	2.48	0.44
1:B:149:VAL:HG13	1:B:156:VAL:HG22	2.00	0.44
1:B:230:TYR:HE1	1:B:232:TRP:HA	1.82	0.44
1:A:230:TYR:C	1:A:230:TYR:CD1	2.92	0.44
1:A:160:LYS:N	1:A:160:LYS:HZ3	2.13	0.43
1:A:249:LEU:C	1:A:251:GLU:H	2.21	0.43
1:B:143:SER:OG	1:B:144:HIS:N	2.51	0.43
1:A:177:LYS:O	1:A:179:LYS:HE2	2.18	0.43
1:A:259:MET:HE2	1:B:259:MET:HG3	1.99	0.43
1:A:158:PRO:HG2	1:A:161:ASP:OD2	2.18	0.43
1:A:225:THR:O	1:A:226:VAL:CB	2.66	0.43
1:A:191:LEU:HD21	1:A:282:PHE:HB3	1.99	0.43
1:B:273:PHE:O	1:B:274:GLU:HB2	2.19	0.43
1:A:226:VAL:HG11	1:B:215:VAL:CB	2.48	0.43
1:B:132:THR:N	1:B:173:THR:HG22	2.34	0.43
1:A:233:ARG:HB3	1:A:239:SER:HA	2.01	0.43
1:B:75:ASN:HB2	1:B:143:SER:HB2	2.01	0.43
1:A:213:ILE:HG23	1:A:214:PRO:HD3	2.00	0.43
1:B:272:HIS:O	1:B:273:PHE:HB2	2.19	0.43
1:A:287:GLN:OE1	1:A:287:GLN:HA	2.18	0.42
1:B:296:THR:C	1:B:297:LYS:HG2	2.39	0.42
1:B:84:MET:CE	1:B:144:HIS:HB3	2.49	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:149:VAL:HG21	1:A:156:VAL:HG13	2.01	0.42
1:A:257:GLU:OE1	1:A:259:MET:HB2	2.19	0.42
1:A:263:THR:O	1:A:266:ASN:HB2	2.19	0.42
1:A:252:HIS:HA	1:A:255:ASP:OD1	2.19	0.42
1:B:158:PRO:HG2	1:B:161:ASP:HB2	2.00	0.42
1:B:84:MET:HE2	1:B:144:HIS:HB3	2.01	0.42
1:B:229:TYR:N	1:B:229:TYR:CD1	2.87	0.42
1:A:178:PRO:HG3	1:A:299:LEU:HD13	2.02	0.42
1:B:121:ASP:OD1	1:B:125:LYS:HE3	2.20	0.42
1:A:226:VAL:HG11	1:A:229:TYR:CD1	2.55	0.42
1:A:136:CYS:HB3	1:A:178:PRO:HG2	2.01	0.42
1:B:246:CYS:O	1:B:250:GLU:HG2	2.19	0.42
1:A:191:LEU:H	1:A:191:LEU:CD1	2.25	0.41
1:A:191:LEU:HD22	1:A:282:PHE:HA	2.02	0.41
1:B:167:ARG:O	1:B:168:GLY:C	2.57	0.41
1:A:187:ARG:HG3	1:A:230:TYR:CE2	2.55	0.41
1:B:116:CYS:HA	1:B:119:MET:HE2	2.02	0.41
1:A:136:CYS:HB2	1:A:178:PRO:O	2.21	0.41
1:B:229:TYR:N	1:B:229:TYR:HD1	2.17	0.41
1:B:213:ILE:O	1:B:213:ILE:HD12	2.20	0.41
1:B:254:LYS:HD2	1:B:254:LYS:N	2.36	0.41
1:A:163:THR:HG22	1:A:181:PHE:CZ	2.54	0.41
1:A:215:VAL:HG11	1:B:229:TYR:CD1	2.55	0.41
1:A:192:ASP:CB	1:A:287:GLN:NE2	2.80	0.41
1:B:100:CYS:SG	1:B:249:LEU:HD23	2.61	0.41
1:A:121:ASP:OD1	1:A:125:LYS:HE3	2.21	0.41
1:A:246:CYS:O	1:A:250:GLU:HB2	2.20	0.41
1:A:213:ILE:CG2	1:A:214:PRO:HD3	2.51	0.40
1:A:283:HIS:CE1	1:A:284:GLU:HG2	2.55	0.40
1:A:266:ASN:OD1	1:A:289:PRO:HB2	2.22	0.40
1:A:195:ILE:HG23	1:A:196:GLN:N	2.36	0.40
1:A:184:GLN:HA	1:A:224:SER:CB	2.51	0.40
1:A:233:ARG:CG	1:A:234:SER:N	2.82	0.40
1:B:150:ILE:HG12	1:B:151:TYR:N	2.36	0.40
1:B:60:TYR:CD1	1:B:178:PRO:HD3	2.56	0.40
1:B:84:MET:HB3	1:B:144:HIS:CD2	2.56	0.40
1:B:163:THR:OG1	1:B:214:PRO:CG	2.70	0.40
1:A:84:MET:HE2	1:A:144:HIS:HB3	2.02	0.40
1:A:73:ILE:HD12	1:A:141:LEU:CD2	2.52	0.40
1:B:123:LEU:HD12	1:B:162:LEU:HD22	2.02	0.40
1:A:180:LEU:HD13	1:A:180:LEU:N	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	228/253 (90%)	193 (85%)	25 (11%)	10 (4%)	4	5
1	B	228/253 (90%)	193 (85%)	23 (10%)	12 (5%)	3	3
All	All	456/506 (90%)	386 (85%)	48 (10%)	22 (5%)	4	4

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	TYR
1	A	190	GLU
1	B	190	GLU
1	B	233	ARG
1	B	276	GLN
1	A	193	ASP
1	A	272	HIS
1	A	273	PHE
1	A	284	GLU
1	B	147	GLU
1	B	192	ASP
1	B	272	HIS
1	B	227	PRO
1	B	273	PHE
1	A	226	VAL
1	A	227	PRO
1	A	235	PRO
1	B	274	GLU
1	B	168	GLY
1	B	235	PRO
1	B	234	SER
1	A	195	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	203/221 (92%)	186 (92%)	17 (8%)	16	29
1	B	203/221 (92%)	184 (91%)	19 (9%)	13	23
All	All	406/442 (92%)	370 (91%)	36 (9%)	14	26

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	LEU
1	A	74	ASN
1	A	148	ASN
1	A	150	ILE
1	A	156	VAL
1	A	160	LYS
1	A	173	THR
1	A	175	LEU
1	A	179	LYS
1	A	180	LEU
1	A	191	LEU
1	A	212	LYS
1	A	213	ILE
1	A	214	PRO
1	A	215	VAL
1	A	226	VAL
1	A	230	TYR
1	B	57	THR
1	B	67	LEU
1	B	98	PHE
1	B	133	ASN
1	B	139	CYS
1	B	148	ASN
1	B	149	VAL
1	B	170	ARG
1	B	180	LEU
1	B	191	LEU
1	B	215	VAL
1	B	232	TRP

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Mol	Chain	Res	Type
1	B	237	ARG
1	B	248	ILE
1	B	258	ILE
1	B	271	ARG
1	B	272	HIS
1	B	279	ASP
1	B	287	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	120	GLN
1	A	276	GLN
1	A	287	GLN
1	B	74	ASN
1	B	112	ASN
1	B	133	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.