



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:29 AM GMT

PDB ID : 3M7M
Title : Crystal structure of monomeric hsp33
Authors : Chi, S.W.; Jeong, D.G.; Woo, J.R.; Park, B.C.; Ryu, S.E.; Kim, S.J.
Deposited on : 2010-03-16
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure 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/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

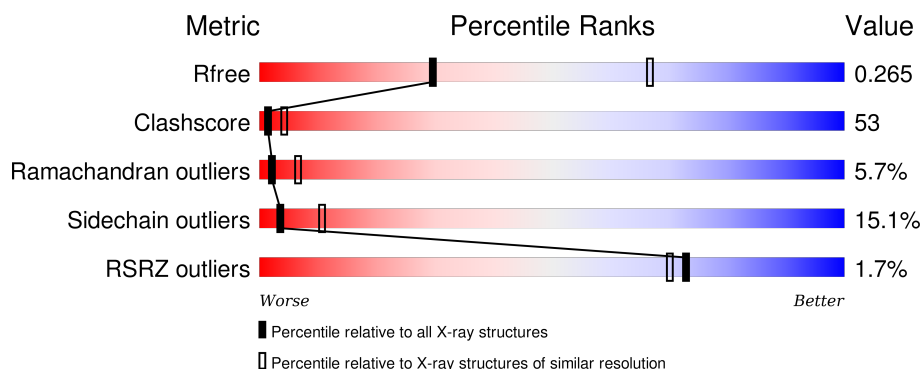
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
R_{free}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	X	235	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1810 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 33 kDa chaperonin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	231	Total	C	N	O	S	0	0	0
			1810	1138	303	362	7			

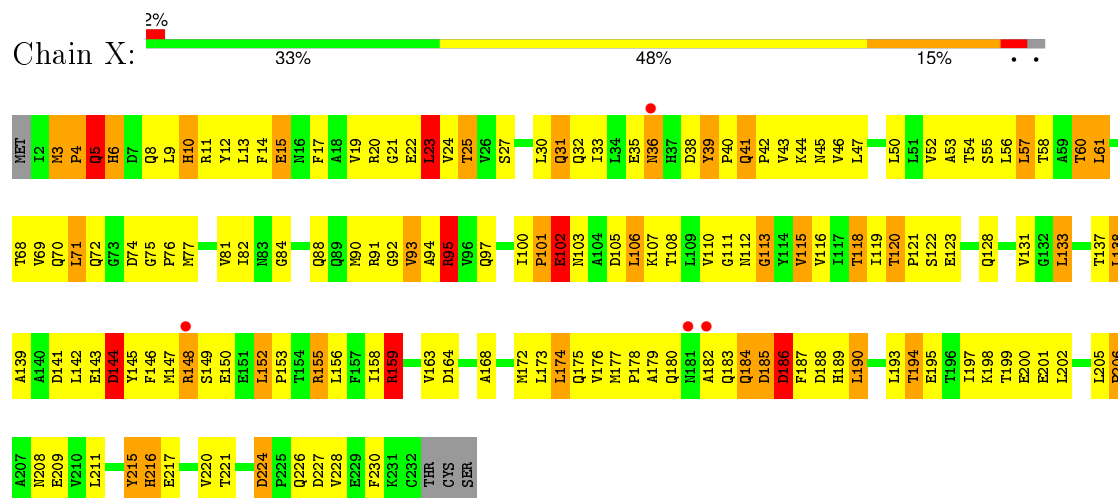
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	1	MET	-	SEE REMARK 999	UNP P0A6Y5
X	2	ILE	-	SEE REMARK 999	UNP P0A6Y5
X	141	ASP	CYS	ENGINEERED MUTATION	UNP P0A6Y5
X	151	GLU	GLN	ENGINEERED MUTATION	UNP P0A6Y5

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 33 kDa chaperonin



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	65.98 Å 65.98 Å 145.95 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.06 – 2.90 30.06 – 2.89	Depositor EDS
% Data completeness (in resolution range)	92.3 (30.06-2.90) 92.0 (30.06-2.89)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.33 (at 2.90 Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.224 , 0.281 0.200 , 0.265	Depositor DCC
R_{free} test set	326 reflections (4.83%)	DCC
Wilson B-factor (Å ²)	61.2	Xtriage
Anisotropy	0.718	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 68.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 7116 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1810	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	X	0.85	0/1841	1.54	24/2507 (1.0%)

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	X	0	3

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	159	ARG	NE-CZ-NH1	-9.10	115.75	120.30
1	X	141	ASP	CB-CG-OD2	8.34	125.80	118.30
1	X	138	LEU	CB-CG-CD1	-8.21	97.05	111.00
1	X	155	ARG	NE-CZ-NH2	-7.87	116.37	120.30
1	X	144	ASP	CB-CG-OD2	7.42	124.98	118.30
1	X	144	ASP	CB-CG-OD1	-7.23	111.79	118.30
1	X	159	ARG	NE-CZ-NH2	7.22	123.91	120.30
1	X	227	ASP	CB-CG-OD1	-6.70	112.27	118.30
1	X	148	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	X	156	LEU	CB-CG-CD2	6.56	122.14	111.00
1	X	195	GLU	CA-CB-CG	6.32	127.31	113.40
1	X	61	LEU	CA-CB-CG	-6.00	101.51	115.30
1	X	38	ASP	CB-CG-OD2	5.99	123.69	118.30
1	X	190	LEU	CA-CB-CG	-5.75	102.08	115.30
1	X	156	LEU	CB-CG-CD1	-5.68	101.34	111.00
1	X	119	ILE	CG1-CB-CG2	-5.59	99.10	111.40
1	X	23	LEU	N-CA-C	-5.52	96.09	111.00
1	X	95	ARG	NE-CZ-NH2	-5.46	117.57	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	206	PRO	N-CD-CG	-5.46	95.02	103.20
1	X	211	LEU	CB-CG-CD1	-5.37	101.87	111.00
1	X	38	ASP	N-CA-C	-5.21	96.94	111.00
1	X	195	GLU	OE1-CD-OE2	-5.18	117.09	123.30
1	X	174	LEU	CB-CG-CD1	5.16	119.77	111.00
1	X	60	THR	CA-CB-CG2	-5.04	105.35	112.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	215	TYR	Sidechain
1	X	224	ASP	Sidechain
1	X	39	TYR	Sidechain

5.2 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	1810	0	1770	188	0
All	All	1810	0	1770	188	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 53.

All (188) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:71:LEU:HD23	1:X:115:VAL:HG13	1.35	1.09
1:X:74:ASP:H	1:X:111:GLY:HA3	1.14	1.08
1:X:74:ASP:N	1:X:111:GLY:HA3	1.70	1.06
1:X:115:VAL:CG2	1:X:131:VAL:HG13	1.96	0.96
1:X:159:ARG:HB3	1:X:199:THR:HG22	1.48	0.95
1:X:71:LEU:CD2	1:X:115:VAL:HG13	1.96	0.95
1:X:115:VAL:HG23	1:X:131:VAL:HG13	1.48	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:174:LEU:HD11	1:X:194:THR:HG21	1.47	0.92
1:X:198:LYS:HB2	1:X:201:GLU:HG3	1.56	0.87
1:X:53:ALA:O	1:X:57:LEU:HD22	1.75	0.85
1:X:35:GLU:O	1:X:36:ASN:HB2	1.76	0.85
1:X:137:THR:HG22	1:X:139:ALA:H	1.42	0.84
1:X:116:VAL:HG12	1:X:118:THR:HG22	1.59	0.83
1:X:184:GLN:HE21	1:X:184:GLN:HA	1.42	0.83
1:X:100:ILE:HG22	1:X:100:ILE:O	1.77	0.81
1:X:110:VAL:HG12	1:X:111:GLY:N	1.97	0.79
1:X:137:THR:HG22	1:X:139:ALA:N	1.99	0.78
1:X:41:GLN:HA	1:X:41:GLN:NE2	1.97	0.78
1:X:159:ARG:HD3	1:X:199:THR:HG22	1.67	0.77
1:X:174:LEU:HD11	1:X:194:THR:CG2	2.15	0.76
1:X:159:ARG:NH1	1:X:159:ARG:HG2	2.02	0.74
1:X:190:LEU:O	1:X:194:THR:CG2	2.36	0.73
1:X:159:ARG:HG2	1:X:159:ARG:HH11	1.51	0.72
1:X:190:LEU:O	1:X:194:THR:HG22	1.89	0.72
1:X:5:GLN:O	1:X:6:HIS:HB3	1.88	0.72
1:X:118:THR:HB	1:X:128:GLN:OE1	1.89	0.71
1:X:27:SER:HB3	1:X:168:ALA:HA	1.73	0.70
1:X:159:ARG:CG	1:X:159:ARG:HH11	2.01	0.69
1:X:184:GLN:NE2	1:X:184:GLN:HA	2.08	0.69
1:X:152:LEU:HD21	1:X:175:GLN:HG3	1.74	0.68
1:X:13:LEU:HD12	1:X:14:PHE:H	1.60	0.67
1:X:106:LEU:O	1:X:106:LEU:HD22	1.94	0.66
1:X:23:LEU:HD23	1:X:24:VAL:N	2.10	0.66
1:X:110:VAL:CG1	1:X:111:GLY:N	2.59	0.66
1:X:10:HIS:N	1:X:10:HIS:ND1	2.44	0.65
1:X:100:ILE:CG2	1:X:100:ILE:O	2.45	0.65
1:X:17:PHE:CE2	1:X:186:ASP:HB3	2.31	0.65
1:X:88:GLN:HG2	1:X:228:VAL:HG22	1.79	0.64
1:X:173:LEU:HD23	1:X:174:LEU:N	2.14	0.63
1:X:41:GLN:HB2	1:X:42:PRO:HD3	1.80	0.63
1:X:121:PRO:O	1:X:123:GLU:N	2.32	0.63
1:X:13:LEU:HD12	1:X:14:PHE:N	2.13	0.63
1:X:41:GLN:CB	1:X:42:PRO:HD3	2.28	0.62
1:X:30:LEU:O	1:X:31:GLN:C	2.36	0.62
1:X:159:ARG:HB3	1:X:199:THR:CG2	2.27	0.61
1:X:23:LEU:C	1:X:23:LEU:HD23	2.20	0.61
1:X:20:ARG:HH22	1:X:60:THR:HA	1.66	0.60
1:X:91:ARG:HG2	1:X:92:GLY:N	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:68:THR:OG1	1:X:118:THR:HG23	2.02	0.60
1:X:186:ASP:N	1:X:186:ASP:OD2	2.32	0.59
1:X:41:GLN:NE2	1:X:41:GLN:CA	2.65	0.59
1:X:198:LYS:O	1:X:199:THR:C	2.39	0.59
1:X:11:ARG:HH22	1:X:226:GLN:HE22	1.50	0.59
1:X:107:LYS:HE2	1:X:133:LEU:O	2.03	0.58
1:X:56:LEU:O	1:X:57:LEU:C	2.38	0.58
1:X:198:LYS:HB2	1:X:201:GLU:CG	2.31	0.58
1:X:41:GLN:CB	1:X:42:PRO:CD	2.81	0.58
1:X:184:GLN:O	1:X:187:PHE:HB3	2.04	0.57
1:X:139:ALA:O	1:X:143:GLU:HG3	2.04	0.57
1:X:41:GLN:HE21	1:X:41:GLN:CA	2.18	0.57
1:X:40:PRO:O	1:X:41:GLN:C	2.43	0.57
1:X:115:VAL:HB	1:X:131:VAL:CG1	2.35	0.56
1:X:173:LEU:HD23	1:X:173:LEU:C	2.26	0.56
1:X:184:GLN:HE21	1:X:184:GLN:CA	2.07	0.56
1:X:13:LEU:HD12	1:X:19:VAL:O	2.05	0.56
1:X:3:MET:O	1:X:3:MET:CE	2.54	0.56
1:X:143:GLU:OE1	1:X:155:ARG:HD2	2.06	0.56
1:X:187:PHE:O	1:X:188:ASP:C	2.42	0.55
1:X:144:ASP:O	1:X:145:TYR:C	2.43	0.55
1:X:163:VAL:O	1:X:164:ASP:HB2	2.05	0.55
1:X:12:TYR:CE1	1:X:21:GLY:HA3	2.41	0.55
1:X:172:MET:CE	1:X:197:ILE:HG12	2.37	0.55
1:X:194:THR:O	1:X:197:ILE:HG22	2.07	0.55
1:X:177:MET:HB3	1:X:180:GLN:NE2	2.22	0.55
1:X:81:VAL:O	1:X:81:VAL:HG13	2.07	0.55
1:X:101:PRO:O	1:X:103:ASN:N	2.40	0.54
1:X:30:LEU:C	1:X:32:GLN:N	2.56	0.54
1:X:183:GLN:O	1:X:184:GLN:C	2.46	0.54
1:X:159:ARG:HD3	1:X:199:THR:CG2	2.35	0.54
1:X:190:LEU:O	1:X:194:THR:HG23	2.07	0.54
1:X:146:PHE:HA	1:X:150:GLU:HB2	1.90	0.54
1:X:40:PRO:HG2	1:X:43:VAL:HG23	1.90	0.53
1:X:185:ASP:O	1:X:188:ASP:N	2.42	0.53
1:X:155:ARG:HH11	1:X:155:ARG:HB3	1.72	0.53
1:X:69:VAL:O	1:X:81:VAL:HG23	2.09	0.52
1:X:198:LYS:O	1:X:201:GLU:N	2.41	0.52
1:X:116:VAL:CG1	1:X:118:THR:HG22	2.37	0.52
1:X:216:HIS:ND1	1:X:217:GLU:N	2.57	0.52
1:X:159:ARG:HG3	1:X:202:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:185:ASP:O	1:X:186:ASP:C	2.49	0.52
1:X:12:TYR:HE1	1:X:21:GLY:HA3	1.75	0.51
1:X:50:LEU:HB3	1:X:82:ILE:HG12	1.93	0.51
1:X:45:ASN:O	1:X:46:VAL:C	2.47	0.50
1:X:27:SER:O	1:X:31:GLN:HG2	2.12	0.50
1:X:200:GLU:HA	1:X:200:GLU:OE2	2.11	0.50
1:X:22:GLU:HG3	1:X:173:LEU:HB3	1.94	0.50
1:X:9:LEU:HD12	1:X:24:VAL:HB	1.92	0.50
1:X:11:ARG:NH1	1:X:11:ARG:HG3	2.27	0.50
1:X:105:ASP:HB3	1:X:108:THR:OG1	2.12	0.49
1:X:206:PRO:HB2	1:X:208:ASN:HD21	1.77	0.49
1:X:14:PHE:O	1:X:15:GLU:C	2.48	0.49
1:X:112:ASN:O	1:X:113:GLY:O	2.29	0.49
1:X:186:ASP:O	1:X:189:HIS:HB2	2.12	0.49
1:X:57:LEU:HD13	1:X:57:LEU:N	2.22	0.49
1:X:201:GLU:O	1:X:205:LEU:HB2	2.12	0.49
1:X:115:VAL:CB	1:X:131:VAL:HG13	2.42	0.48
1:X:27:SER:CB	1:X:168:ALA:HA	2.42	0.48
1:X:177:MET:HE3	1:X:178:PRO:HD2	1.95	0.48
1:X:30:LEU:O	1:X:32:GLN:N	2.46	0.48
1:X:110:VAL:CG1	1:X:111:GLY:H	2.27	0.48
1:X:172:MET:HE3	1:X:197:ILE:HG21	1.95	0.48
1:X:3:MET:O	1:X:3:MET:HE3	2.13	0.48
1:X:155:ARG:NH1	1:X:155:ARG:HB3	2.29	0.47
1:X:101:PRO:O	1:X:102:GLU:C	2.53	0.47
1:X:52:VAL:HG11	1:X:158:ILE:HG23	1.96	0.47
1:X:115:VAL:CG2	1:X:131:VAL:CG1	2.83	0.47
1:X:93:VAL:HG23	1:X:94:ALA:N	2.29	0.47
1:X:41:GLN:HA	1:X:44:LYS:HZ3	1.80	0.47
1:X:91:ARG:HG2	1:X:92:GLY:H	1.77	0.46
1:X:33:ILE:HG12	1:X:230:PHE:HE1	1.81	0.46
1:X:46:VAL:O	1:X:50:LEU:N	2.45	0.46
1:X:33:ILE:HG12	1:X:230:PHE:CE1	2.51	0.46
1:X:206:PRO:HB2	1:X:208:ASN:ND2	2.31	0.46
1:X:57:LEU:O	1:X:60:THR:OG1	2.29	0.46
1:X:197:ILE:HG23	1:X:197:ILE:O	2.16	0.46
1:X:142:LEU:O	1:X:143:GLU:C	2.54	0.46
1:X:70:GLN:CG	1:X:116:VAL:HB	2.46	0.45
1:X:58:THR:C	1:X:60:THR:H	2.20	0.45
1:X:115:VAL:HG23	1:X:131:VAL:O	2.17	0.45
1:X:60:THR:O	1:X:61:LEU:HD23	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:120:THR:HA	1:X:121:PRO:HD2	1.84	0.45
1:X:95:ARG:HD3	1:X:95:ARG:HA	1.36	0.45
1:X:20:ARG:HD2	1:X:177:MET:SD	2.56	0.45
1:X:177:MET:CE	1:X:180:GLN:HE22	2.31	0.44
1:X:100:ILE:HA	1:X:101:PRO:HD2	1.74	0.44
1:X:22:GLU:OE1	1:X:173:LEU:HD13	2.17	0.44
1:X:198:LYS:HB2	1:X:201:GLU:OE1	2.18	0.44
1:X:205:LEU:HA	1:X:206:PRO:HD2	1.76	0.44
1:X:23:LEU:HD23	1:X:24:VAL:CA	2.48	0.44
1:X:50:LEU:HD23	1:X:138:LEU:CD2	2.48	0.44
1:X:142:LEU:HD23	1:X:142:LEU:HA	1.84	0.44
1:X:72:GLN:O	1:X:110:VAL:CG1	2.66	0.43
1:X:11:ARG:HH22	1:X:226:GLN:NE2	2.14	0.43
1:X:54:THR:O	1:X:55:SER:C	2.54	0.43
1:X:20:ARG:HG2	1:X:177:MET:SD	2.57	0.43
1:X:41:GLN:HA	1:X:44:LYS:NZ	2.33	0.43
1:X:178:PRO:O	1:X:178:PRO:HG2	2.19	0.43
1:X:198:LYS:O	1:X:200:GLU:N	2.51	0.43
1:X:105:ASP:OD1	1:X:106:LEU:N	2.51	0.43
1:X:61:LEU:HA	1:X:61:LEU:HD23	1.84	0.43
1:X:41:GLN:HB2	1:X:42:PRO:CD	2.44	0.43
1:X:43:VAL:HG13	1:X:77:MET:SD	2.59	0.43
1:X:175:GLN:O	1:X:176:VAL:C	2.57	0.43
1:X:200:GLU:OE2	1:X:200:GLU:CA	2.64	0.42
1:X:3:MET:O	1:X:3:MET:HE2	2.18	0.42
1:X:75:GLY:H	1:X:111:GLY:N	2.17	0.42
1:X:176:VAL:HG13	1:X:182:ALA:HB3	2.01	0.42
1:X:177:MET:HE2	1:X:180:GLN:HE22	1.83	0.42
1:X:184:GLN:O	1:X:185:ASP:C	2.55	0.42
1:X:39:TYR:CE2	1:X:94:ALA:HB3	2.55	0.42
1:X:190:LEU:HD23	1:X:190:LEU:HA	1.85	0.42
1:X:4:PRO:HB2	1:X:5:GLN:H	1.64	0.42
1:X:75:GLY:N	1:X:111:GLY:N	2.68	0.42
1:X:17:PHE:CE2	1:X:186:ASP:CB	3.02	0.42
1:X:24:VAL:HG22	1:X:25:THR:N	2.35	0.42
1:X:84:GLY:HA3	1:X:90:MET:CE	2.49	0.42
1:X:88:GLN:CG	1:X:228:VAL:HG22	2.48	0.41
1:X:76:PRO:O	1:X:97:GLN:HG2	2.20	0.41
1:X:172:MET:HE3	1:X:197:ILE:HG12	2.02	0.41
1:X:27:SER:HB3	1:X:168:ALA:CA	2.47	0.41
1:X:194:THR:HB	1:X:215:TYR:OH	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:31:GLN:H	1:X:31:GLN:HG2	1.59	0.41
1:X:106:LEU:C	1:X:106:LEU:HD22	2.40	0.41
1:X:46:VAL:HG23	1:X:47:LEU:N	2.34	0.41
1:X:21:GLY:HA2	1:X:173:LEU:O	2.21	0.41
1:X:102:GLU:CG	1:X:103:ASN:N	2.83	0.41
1:X:60:THR:HG21	1:X:173:LEU:CD1	2.51	0.41
1:X:185:ASP:O	1:X:187:PHE:N	2.54	0.41
1:X:12:TYR:CE1	1:X:21:GLY:CA	3.04	0.41
1:X:115:VAL:HB	1:X:131:VAL:HG13	2.01	0.41
1:X:72:GLN:O	1:X:110:VAL:HG11	2.21	0.41
1:X:159:ARG:HG3	1:X:202:LEU:CD2	2.51	0.41
1:X:57:LEU:HD12	1:X:57:LEU:HA	1.65	0.41
1:X:153:PRO:HD2	1:X:176:VAL:HB	2.01	0.41
1:X:220:VAL:CG1	1:X:221:THR:N	2.84	0.41
1:X:58:THR:C	1:X:60:THR:N	2.74	0.40
1:X:173:LEU:HD23	1:X:174:LEU:CA	2.51	0.40
1:X:50:LEU:HB3	1:X:82:ILE:CG1	2.51	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	X	229/235 (97%)	187 (82%)	29 (13%)	13 (6%)	2 6

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	4	PRO
1	X	102	GLU
1	X	122	SER
1	X	185	ASP

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Mol	Chain	Res	Type
1	X	5	GLN
1	X	113	GLY
1	X	179	ALA
1	X	186	ASP
1	X	15	GLU
1	X	148	ARG
1	X	6	HIS
1	X	36	ASN
1	X	101	PRO

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	X	199/203 (98%)	169 (85%)	30 (15%)	3 11

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	3	MET
1	X	5	GLN
1	X	8	GLN
1	X	10	HIS
1	X	23	LEU
1	X	25	THR
1	X	31	GLN
1	X	41	GLN
1	X	57	LEU
1	X	71	LEU
1	X	93	VAL
1	X	95	ARG
1	X	102	GLU
1	X	106	LEU
1	X	115	VAL
1	X	118	THR
1	X	120	THR

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Mol	Chain	Res	Type
1	X	133	LEU
1	X	144	ASP
1	X	147	MET
1	X	149	SER
1	X	152	LEU
1	X	159	ARG
1	X	184	GLN
1	X	186	ASP
1	X	193	LEU
1	X	194	THR
1	X	209	GLU
1	X	216	HIS
1	X	224	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	8	GLN
1	X	41	GLN
1	X	45	ASN
1	X	70	GLN
1	X	88	GLN
1	X	97	GLN
1	X	103	ASN
1	X	180	GLN
1	X	184	GLN
1	X	208	ASN
1	X	226	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	X	231/235 (98%)	-0.37	4 (1%) 73 70	32, 66, 109, 133	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	36	ASN	2.8
1	X	148	ARG	2.3
1	X	182	ALA	2.1
1	X	181	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.