



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2023 – 02:31 PM EDT

PDB ID : 3KKL  
Title : Crystal structure of functionally unknown HSP33 from *Saccharomyces cerevisiae*  
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Deposited on : 2009-11-05  
Resolution : 2.03 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

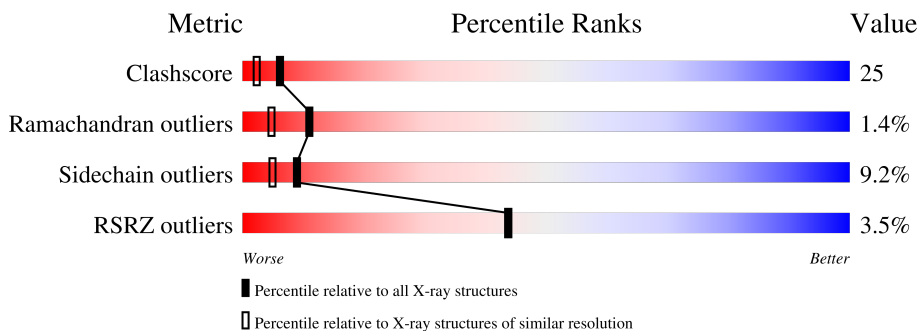
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.03 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	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	A	244	 4% 52% 28% 6% • 12%
1	B	244	 2% 59% 24% • • 11%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 3387 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 Probable chaperone protein HSP33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	215	Total	C	N	O	S	0	0	0
			1665	1073	280	311	1			
1	B	216	Total	C	N	O	S	0	0	0
			1672	1077	281	313	1			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	238	VAL	-	expression tag	UNP Q08914
A	239	GLU	-	expression tag	UNP Q08914
A	240	HIS	-	expression tag	UNP Q08914
A	241	HIS	-	expression tag	UNP Q08914
A	242	HIS	-	expression tag	UNP Q08914
A	243	HIS	-	expression tag	UNP Q08914
A	244	HIS	-	expression tag	UNP Q08914
B	238	VAL	-	expression tag	UNP Q08914
B	239	GLU	-	expression tag	UNP Q08914
B	240	HIS	-	expression tag	UNP Q08914
B	241	HIS	-	expression tag	UNP Q08914
B	242	HIS	-	expression tag	UNP Q08914
B	243	HIS	-	expression tag	UNP Q08914
B	244	HIS	-	expression tag	UNP Q08914

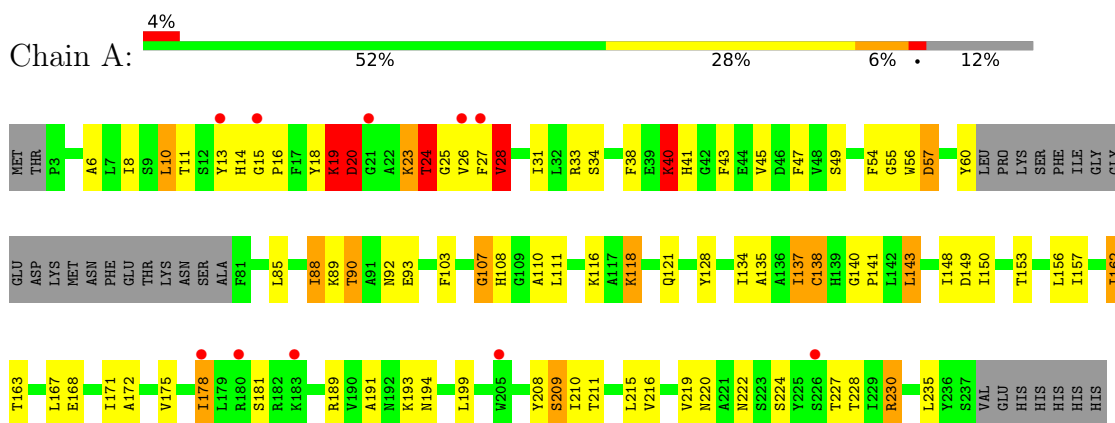
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	24	Total	O	0	0
			24	24		
2	B	26	Total	O	0	0
			26	26		

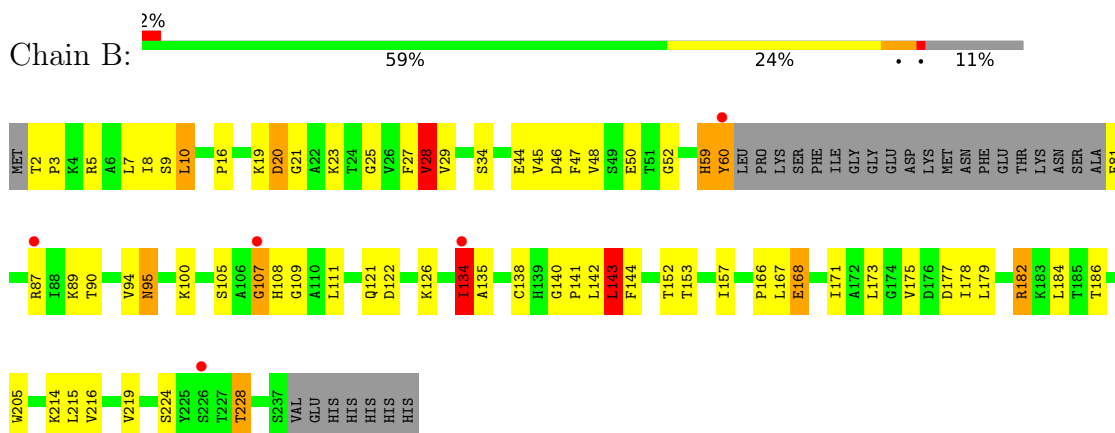
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Probable chaperone protein HSP33



- Molecule 1: Probable chaperone protein HSP33



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.45Å 68.74Å 66.31Å 90.00° 100.08° 90.00°	Depositor
Resolution (Å)	19.76 – 2.03 36.32 – 2.03	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.76-2.03) 92.4 (36.32-2.03)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.25 (at 2.03Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.234 , 0.247 0.236 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.2	Xtrriage
Anisotropy	0.608	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 55.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3387	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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	A	0.89	1/1704 (0.1%)	0.90	7/2307 (0.3%)
1	B	0.97	0/1711	0.91	4/2318 (0.2%)
All	All	0.93	1/3415 (0.0%)	0.91	11/4625 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	110	ALA	CA-CB	5.68	1.64	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	VAL	CB-CA-C	-7.93	96.33	111.40
1	A	230	ARG	NE-CZ-NH2	-6.93	116.83	120.30
1	A	137	ILE	CG1-CB-CG2	-6.60	96.89	111.40
1	B	107	GLY	N-CA-C	-5.73	98.77	113.10
1	B	143	LEU	CB-CG-CD2	5.53	120.41	111.00
1	B	134	ILE	CG1-CB-CG2	-5.52	99.25	111.40
1	A	10	LEU	CA-CB-CG	5.46	127.86	115.30
1	A	28	VAL	CB-CA-C	-5.25	101.43	111.40
1	A	107	GLY	N-CA-C	-5.12	100.29	113.10
1	A	19	LYS	N-CA-C	-5.09	97.24	111.00
1	A	10	LEU	CB-CG-CD1	-5.02	102.47	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	A	1665	0	1649	92	0
1	B	1672	0	1655	73	0
2	A	24	0	0	4	0
2	B	26	0	0	2	0
All	All	3387	0	3304	163	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 25.

All (163) 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:A:24:THR:HG21	1:A:108:HIS:HB2	1.24	1.19
1:A:11:THR:HG22	1:A:26:VAL:CG1	1.73	1.18
1:A:11:THR:HG22	1:A:26:VAL:HG12	1.17	1.15
1:B:152:THR:O	1:B:153:THR:HG22	1.58	1.02
1:B:7:LEU:HD11	1:B:48:VAL:HG23	1.40	1.00
1:A:224:SER:O	1:A:228:THR:HG23	1.65	0.96
1:A:209:SER:HB3	1:A:230:ARG:HE	1.32	0.94
1:A:175:VAL:HA	1:A:178:ILE:HD13	1.51	0.92
1:B:134:ILE:H	1:B:134:ILE:HD12	1.35	0.91
1:A:19:LYS:O	1:A:20:ASP:HB3	1.68	0.90
1:B:7:LEU:HD11	1:B:48:VAL:CG2	2.03	0.88
1:B:175:VAL:HA	1:B:178:ILE:HD13	1.56	0.84
1:A:11:THR:CG2	1:A:26:VAL:CG1	2.55	0.83
1:A:24:THR:CG2	1:A:108:HIS:HB2	2.06	0.83
1:A:178:ILE:H	1:A:178:ILE:HD12	1.43	0.82
1:A:28:VAL:HG13	1:A:56:TRP:CZ3	2.15	0.82
1:A:137:ILE:HD11	1:A:227:THR:HB	1.64	0.80
1:A:19:LYS:O	1:A:20:ASP:CB	2.30	0.79
1:A:178:ILE:H	1:A:178:ILE:CD1	1.97	0.78
1:A:11:THR:HG22	1:A:26:VAL:HG13	1.66	0.78
1:B:52:GLY:HA3	1:B:90:THR:HG22	1.67	0.76
1:B:134:ILE:HD13	1:B:215:LEU:HA	1.66	0.76
1:A:103:PHE:HZ	1:A:137:ILE:HD12	1.48	0.75
1:A:11:THR:CG2	1:A:26:VAL:HG13	2.16	0.75
1:B:111:LEU:HD23	1:B:142:LEU:HD11	1.68	0.74
1:B:2:THR:HB	1:B:3:PRO:HD3	1.71	0.72
1:B:28:VAL:CG2	1:B:60:TYR:HE2	2.03	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:LYS:HA	1:B:20:ASP:O	1.89	0.72
1:A:103:PHE:CE1	1:A:228:THR:HG22	2.24	0.71
1:B:167:LEU:O	1:B:171:ILE:HD12	1.90	0.71
1:B:34:SER:HA	1:B:228:THR:HG21	1.72	0.71
1:A:137:ILE:HD13	1:A:224:SER:HA	1.73	0.70
1:A:137:ILE:CD1	1:A:224:SER:HA	2.22	0.70
1:A:24:THR:HG21	1:A:108:HIS:CB	2.13	0.70
1:A:175:VAL:CA	1:A:178:ILE:HD13	2.21	0.69
1:A:178:ILE:HD12	1:A:178:ILE:N	2.07	0.69
1:A:162:ILE:HD11	1:A:191:ALA:HB2	1.75	0.68
1:B:28:VAL:CG2	1:B:60:TYR:CE2	2.77	0.68
1:A:103:PHE:CZ	1:A:137:ILE:HD12	2.28	0.68
1:B:27:PHE:CE1	1:B:29:VAL:HG22	2.29	0.67
1:B:48:VAL:HG22	1:B:89:LYS:HB2	1.75	0.67
1:A:16:PRO:HA	1:A:23:LYS:HG3	1.77	0.67
1:A:209:SER:HB3	1:A:230:ARG:NE	2.06	0.66
1:B:134:ILE:HD13	1:B:215:LEU:HD12	1.78	0.66
1:B:152:THR:O	1:B:153:THR:CG2	2.41	0.66
1:B:28:VAL:HG21	1:B:60:TYR:CE2	2.30	0.66
1:A:14:HIS:CG	1:A:55:GLY:HA3	2.30	0.65
1:B:134:ILE:HD12	1:B:214:LYS:O	1.96	0.65
1:B:166:PRO:HB2	1:B:168:GLU:HG3	1.79	0.64
1:A:175:VAL:C	1:A:178:ILE:HD13	2.18	0.64
1:A:108:HIS:ND1	2:A:269:HOH:O	2.30	0.63
1:A:175:VAL:HA	1:A:178:ILE:CD1	2.27	0.63
1:B:167:LEU:HG	1:B:171:ILE:HD11	1.81	0.63
1:B:224:SER:O	1:B:228:THR:HG22	1.99	0.62
1:A:137:ILE:HD11	1:A:227:THR:CB	2.28	0.62
1:B:8:ILE:HD12	1:B:45:VAL:HG11	1.81	0.62
1:A:163:THR:HG22	1:A:199:LEU:O	2.00	0.62
1:A:121:GLN:HG3	1:A:143:LEU:HD22	1.82	0.62
1:B:20:ASP:OD2	1:B:21:GLY:N	2.33	0.62
1:A:16:PRO:CA	1:A:23:LYS:HG3	2.30	0.61
1:B:134:ILE:CD1	1:B:215:LEU:HA	2.30	0.61
1:A:54:PHE:HZ	1:A:85:LEU:HD11	1.66	0.61
1:B:167:LEU:HG	1:B:171:ILE:CD1	2.32	0.59
1:A:10:LEU:CD2	1:A:31:ILE:HG21	2.32	0.59
1:B:121:GLN:HG3	1:B:143:LEU:HD22	1.83	0.59
1:B:107:GLY:O	1:B:109:GLY:N	2.36	0.58
1:B:173:LEU:HD21	1:B:205:TRP:HH2	1.68	0.58
1:B:48:VAL:HG21	1:B:94:VAL:HG21	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:LEU:HD13	1:B:186:THR:HG21	1.85	0.57
1:B:134:ILE:HD12	1:B:134:ILE:N	2.14	0.57
1:A:6:ALA:HB3	1:A:45:VAL:HG22	1.85	0.57
1:A:18:TYR:HB2	1:A:20:ASP:OD2	2.03	0.57
1:A:34:SER:HA	1:A:228:THR:HG21	1.87	0.56
1:B:224:SER:O	1:B:228:THR:CG2	2.53	0.56
1:B:177:ASP:OD2	1:B:178:ILE:HD12	2.05	0.56
1:A:88:ILE:HD13	1:A:89:LYS:O	2.07	0.55
1:B:173:LEU:HD21	1:B:205:TRP:CH2	2.40	0.55
1:B:16:PRO:HB2	1:B:182:ARG:HH22	1.70	0.55
1:A:108:HIS:N	2:A:269:HOH:O	2.00	0.55
1:A:14:HIS:CD2	1:A:55:GLY:HA3	2.42	0.55
1:A:128:TYR:HB2	1:A:134:ILE:HD11	1.90	0.54
1:A:162:ILE:HD11	1:A:191:ALA:CB	2.37	0.54
1:A:148:ILE:HG22	1:A:150:ILE:HD13	1.89	0.54
1:B:27:PHE:CZ	1:B:29:VAL:HG22	2.42	0.54
1:B:175:VAL:CA	1:B:178:ILE:HD13	2.32	0.54
1:A:18:TYR:CB	1:A:20:ASP:OD2	2.57	0.53
1:A:171:ILE:HD12	1:A:172:ALA:N	2.23	0.53
1:A:220:ASN:OD1	1:A:222:ASN:HB2	2.08	0.53
1:B:178:ILE:HD12	1:B:178:ILE:H	1.75	0.52
1:B:28:VAL:HG21	1:B:60:TYR:HE2	1.67	0.52
1:A:135:ALA:HA	1:A:216:VAL:O	2.10	0.51
1:A:27:PHE:HD1	1:A:107:GLY:HA2	1.76	0.51
1:A:54:PHE:O	1:B:23:LYS:NZ	2.41	0.51
1:A:116:LYS:O	1:A:118:LYS:HD2	2.11	0.51
1:A:11:THR:HG23	1:A:26:VAL:HG13	1.94	0.50
1:B:19:LYS:HA	1:B:20:ASP:C	2.32	0.50
1:A:189:ARG:O	1:A:193:LYS:HB2	2.11	0.50
1:B:20:ASP:CG	1:B:21:GLY:H	2.16	0.49
1:B:134:ILE:HD13	1:B:215:LEU:CD1	2.41	0.49
1:A:13:TYR:CZ	1:A:15:GLY:O	2.65	0.49
1:A:19:LYS:O	1:A:20:ASP:OD2	2.30	0.49
1:B:25:GLY:O	1:B:107:GLY:HA3	2.12	0.49
1:B:144:PHE:CD2	1:B:157:ILE:HG21	2.49	0.48
1:A:168:GLU:O	1:A:171:ILE:HG13	2.14	0.48
1:A:211:THR:OG1	1:A:230:ARG:HD3	2.13	0.48
1:B:140:GLY:N	1:B:141:PRO:CD	2.77	0.48
1:B:3:PRO:HA	2:B:262:HOH:O	2.13	0.47
1:B:138:CYS:HA	1:B:219:VAL:O	2.15	0.47
1:B:177:ASP:CG	1:B:178:ILE:HD12	2.34	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LYS:HD2	1:A:118:LYS:H	1.80	0.47
1:B:134:ILE:H	1:B:134:ILE:CD1	2.15	0.47
1:A:16:PRO:HD3	1:B:52:GLY:O	2.15	0.47
1:A:90:THR:HG23	1:A:93:GLU:OE1	2.12	0.47
1:A:8:ILE:HD12	1:A:45:VAL:HG11	1.96	0.46
1:B:19:LYS:O	1:B:19:LYS:HG2	2.14	0.46
1:B:28:VAL:HG22	1:B:60:TYR:HE2	1.77	0.46
1:B:107:GLY:C	1:B:109:GLY:N	2.67	0.46
1:A:47:PHE:O	1:A:89:LYS:HG3	2.15	0.46
1:B:81:PHE:N	2:B:255:HOH:O	2.48	0.46
1:B:122:ASP:O	1:B:126:LYS:HG3	2.15	0.46
1:A:175:VAL:O	1:A:178:ILE:HD13	2.16	0.46
1:A:138:CYS:HA	1:A:219:VAL:O	2.15	0.45
1:A:162:ILE:CG2	1:A:215:LEU:HD21	2.46	0.45
1:A:149:ASP:HB2	1:A:156:LEU:HA	1.98	0.45
1:B:175:VAL:HA	1:B:178:ILE:CD1	2.38	0.45
1:B:27:PHE:CZ	1:B:29:VAL:CG2	3.00	0.45
1:A:128:TYR:HB2	1:A:134:ILE:CD1	2.47	0.45
1:A:10:LEU:HB3	1:A:26:VAL:HG11	1.98	0.45
1:A:162:ILE:HG22	1:A:215:LEU:HD21	1.98	0.45
1:A:108:HIS:O	1:A:111:LEU:HB2	2.17	0.44
1:A:171:ILE:HD12	1:A:171:ILE:C	2.38	0.44
1:B:135:ALA:HA	1:B:216:VAL:O	2.17	0.44
1:B:7:LEU:HD12	1:B:46:ASP:O	2.17	0.44
1:B:107:GLY:O	1:B:108:HIS:C	2.57	0.43
1:A:163:THR:HG21	1:A:208:TYR:CE1	2.53	0.43
1:A:40:LYS:NZ	1:A:41:HIS:HE1	2.16	0.43
1:A:57:ASP:OD2	1:A:57:ASP:C	2.56	0.43
1:A:40:LYS:NZ	1:A:41:HIS:CE1	2.87	0.43
1:A:140:GLY:N	1:A:141:PRO:CD	2.81	0.43
1:B:178:ILE:HD12	1:B:178:ILE:N	2.34	0.43
1:A:38:PHE:O	1:A:43:PHE:HB2	2.19	0.43
1:B:5:ARG:NE	1:B:44:GLU:OE1	2.46	0.42
1:A:24:THR:HB	1:A:25:GLY:H	1.55	0.42
1:B:8:ILE:HD12	1:B:45:VAL:CG1	2.47	0.42
1:A:137:ILE:HG13	1:A:227:THR:HG21	2.01	0.42
1:A:219:VAL:HG12	1:A:220:ASN:ND2	2.34	0.42
1:A:163:THR:HG23	1:A:210:ILE:CG1	2.49	0.42
1:A:148:ILE:CG2	1:A:150:ILE:HD13	2.49	0.41
1:B:47:PHE:O	1:B:48:VAL:HG23	2.20	0.41
1:B:182:ARG:HB3	1:B:184:LEU:HG	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ASN:HB2	2:A:257:HOH:O	2.18	0.41
1:A:24:THR:HG22	1:A:57:ASP:OD1	2.21	0.41
1:A:40:LYS:HZ3	1:A:41:HIS:CE1	2.39	0.41
1:A:28:VAL:HG21	1:A:60:TYR:CE1	2.56	0.41
1:B:50:GLU:HA	1:B:50:GLU:OE1	2.20	0.41
1:B:95:ASN:HD22	1:B:95:ASN:C	2.24	0.40
1:A:49:SER:N	1:A:88:ILE:HD11	2.36	0.40
1:A:54:PHE:CZ	1:A:85:LEU:HD11	2.52	0.40
1:B:10:LEU:HD11	1:B:47:PHE:CD1	2.56	0.40
1:A:157:ILE:HD11	1:A:194:ASN:HB2	2.04	0.40
1:A:209:SER:HB2	2:A:263:HOH:O	2.20	0.40
1:B:9:SER:HA	1:B:48:VAL:O	2.22	0.40
1:B:179:LEU:HD22	1:B:184:LEU:HB2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	211/244 (86%)	193 (92%)	14 (7%)	4 (2%)	<b>8</b> <b>3</b>
1	B	212/244 (87%)	198 (93%)	12 (6%)	2 (1%)	<b>17</b> <b>10</b>
All	All	423/488 (87%)	391 (92%)	26 (6%)	6 (1%)	<b>11</b> <b>5</b>

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	20	ASP
1	B	59	HIS
1	A	138	CYS
1	A	40	LYS
1	A	20	ASP

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Mol	Chain	Res	Type
1	A	24	THR

### 5.3.2 Protein sidechains [i](#)

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	173/199 (87%)	154 (89%)	19 (11%)	<b>6</b> <b>3</b>
1	B	174/199 (87%)	161 (92%)	13 (8%)	<b>13</b> <b>8</b>
All	All	347/398 (87%)	315 (91%)	32 (9%)	<b>9</b> <b>5</b>

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LYS
1	A	20	ASP
1	A	23	LYS
1	A	24	THR
1	A	28	VAL
1	A	33	ARG
1	A	40	LYS
1	A	57	ASP
1	A	88	ILE
1	A	90	THR
1	A	118	LYS
1	A	143	LEU
1	A	153	THR
1	A	162	ILE
1	A	167	LEU
1	A	178	ILE
1	A	181	SER
1	A	209	SER
1	A	235	LEU
1	B	10	LEU
1	B	28	VAL
1	B	59	HIS

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Mol	Chain	Res	Type
1	B	60	TYR
1	B	87	ARG
1	B	95	ASN
1	B	100	LYS
1	B	105	SER
1	B	134	ILE
1	B	143	LEU
1	B	168	GLU
1	B	182	ARG
1	B	228	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	41	HIS
1	A	130	ASN
1	B	92	ASN
1	B	95	ASN
1	B	130	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	215/244 (88%)	0.34	10 (4%) 31 31	23, 38, 57, 68	1 (0%)
1	B	216/244 (88%)	0.10	5 (2%) 60 59	19, 36, 52, 62	1 (0%)
All	All	431/488 (88%)	0.22	15 (3%) 44 44	19, 37, 55, 68	2 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	GLY	4.6
1	A	15	GLY	3.8
1	A	205	TRP	3.4
1	A	226	SER	3.1
1	A	26	VAL	3.1
1	A	180	ARG	2.8
1	A	13	TYR	2.8
1	A	27	PHE	2.8
1	B	107	GLY	2.6
1	B	226	SER	2.4
1	B	134	ILE	2.2
1	A	183	LYS	2.1
1	B	87	ARG	2.0
1	B	60	TYR	2.0
1	A	178	ILE	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 monosaccharides in this entry.

## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.