



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:00 PM UTC

PDB ID : 1STZ / pdb_00001stz
Title : Crystal structure of a hypothetical protein at 2.2 Å resolution
Authors : Liu, J.; Adams, P.D.; Shin, D.-H.; Huang, C.; Yokota, H.; Jancarik, J.; Kim, R.; Kim, S.-H.; Berkeley Structural Genomics Center (BSGC)
Deposited on : 2004-03-25
Resolution : 2.20 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

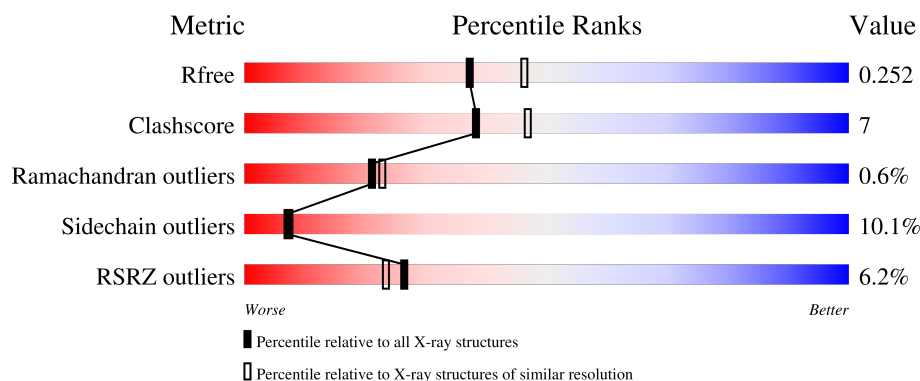
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 ≥ 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	338	9% 74% 18% • •
1	B	338	3% 74% 16% • 8%
1	C	338	6% 72% 17% • 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8166 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 Heat-inducible transcription repressor hrcA homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2634	1690	444	494	6			
1	B	311	Total	C	N	O	S	0	0	0
			2541	1634	431	470	6			
1	C	311	Total	C	N	O	S	0	0	0
			2541	1634	431	470	6			

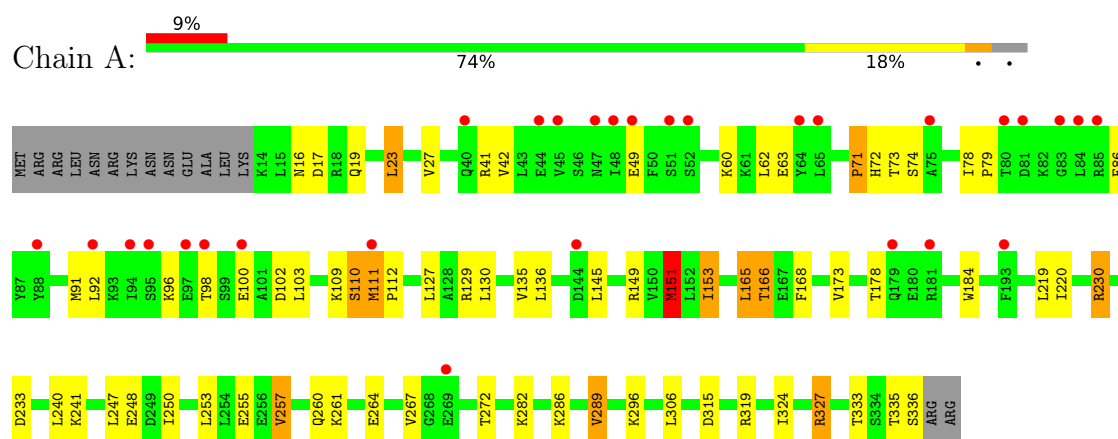
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	184	Total	O	0	0
			184	184		
2	B	131	Total	O	0	0
			131	131		
2	C	135	Total	O	0	0
			135	135		

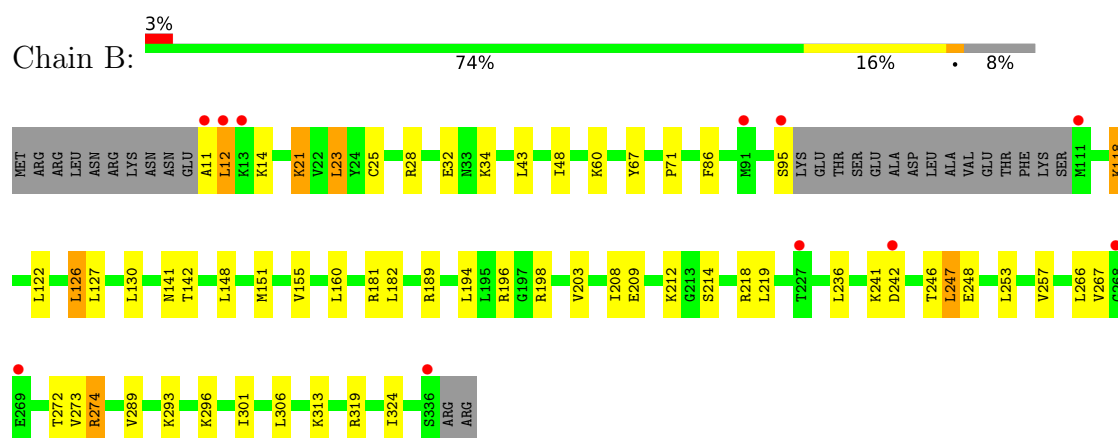
3 Residue-property plots [i](#)

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.

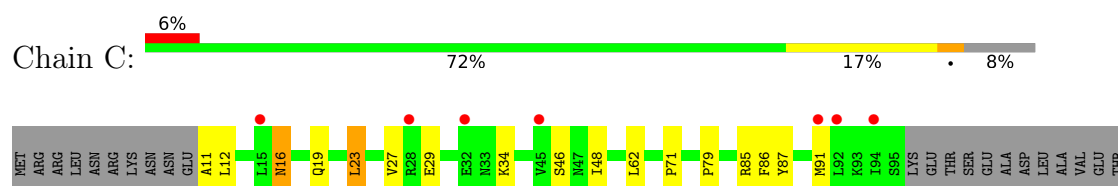
- Molecule 1: Heat-inducible transcription repressor hrcA homolog

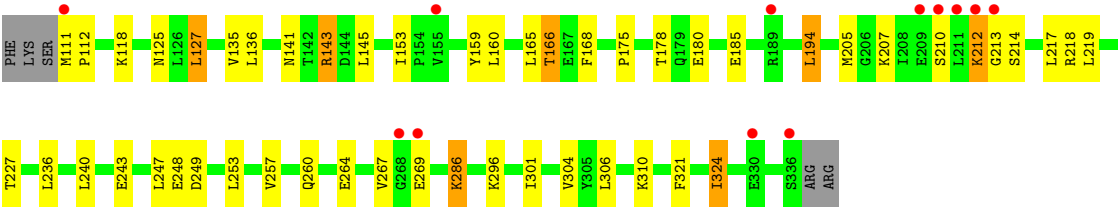


- Molecule 1: Heat-inducible transcription repressor hrcA homolog



- Molecule 1: Heat-inducible transcription repressor hrcA homolog





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	115.30Å 115.30Å 185.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	91.7 (20.00-2.20) 91.5 (20.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.11Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.215 , 0.258 0.216 , 0.252	Depositor DCC
R_{free} test set	3405 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8166	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% 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, 0.333 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	0.69	1/2676 (0.0%)	0.87	3/3599 (0.1%)
1	B	0.66	0/2581	0.85	0/3469
1	C	0.59	0/2581	0.85	0/3469
All	All	0.65	1/7838 (0.0%)	0.86	3/10537 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	151	MET	SD-CE	-7.82	1.59	1.79

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	ILE	N-CA-C	6.45	113.94	107.55
1	A	289	VAL	CB-CA-C	-6.37	101.20	110.63
1	A	110	SER	N-CA-C	-6.06	105.89	113.28

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	2634	0	2717	48	0
1	B	2541	0	2634	26	0
1	C	2541	0	2634	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	184	0	0	15	0
2	B	131	0	0	2	0
2	C	135	0	0	12	0
All	All	8166	0	7985	108	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 7.

All (108) 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:C:136:LEU:O	2:C:431:HOH:O	1.75	1.02
1:A:178:THR:HB	2:A:509:HOH:O	1.76	0.84
1:A:335:THR:C	2:A:460:HOH:O	2.21	0.83
1:B:23:LEU:HD13	1:B:86:PHE:CD2	2.14	0.81
1:C:136:LEU:C	2:C:431:HOH:O	2.21	0.81
1:A:260:GLN:NE2	2:A:502:HOH:O	2.02	0.81
1:A:111:MET:HB2	1:A:112:PRO:HD3	1.62	0.80
1:B:242:ASP:HB3	2:B:422:HOH:O	1.82	0.77
1:A:178:THR:HG21	1:A:184:TRP:HZ2	1.50	0.75
1:C:166:THR:HG23	1:C:168:PHE:H	1.52	0.75
1:A:23:LEU:HD13	1:A:86:PHE:CD2	2.24	0.72
1:A:282:LYS:NZ	2:A:364:HOH:O	2.22	0.71
1:C:321:PHE:HA	2:C:428:HOH:O	1.91	0.70
1:A:71:PRO:HG3	1:A:78:ILE:HD12	1.73	0.69
1:C:23:LEU:HD13	1:C:86:PHE:CD2	2.28	0.68
1:A:166:THR:HG22	1:A:168:PHE:H	1.59	0.67
1:A:173:VAL:HG21	1:C:175:PRO:HG3	1.77	0.66
1:C:212:LYS:NZ	2:C:453:HOH:O	2.29	0.65
1:C:141:ASN:HD21	1:C:143:ARG:HB3	1.61	0.65
1:C:260:GLN:HE22	1:C:264:GLU:CD	2.05	0.65
1:B:32:GLU:OE2	1:B:313:LYS:NZ	2.30	0.64
1:B:25:CYS:HA	1:B:28:ARG:HG2	1.80	0.62
1:B:296:LYS:CB	1:B:301:ILE:HD11	2.30	0.61
1:A:336:SER:N	2:A:460:HOH:O	2.35	0.60
1:C:125:ASN:ND2	2:C:456:HOH:O	2.35	0.60
1:A:333:THR:O	1:A:336:SER:HB3	2.02	0.59
1:A:327:ARG:NH1	2:A:421:HOH:O	2.37	0.57
1:C:111:MET:N	1:C:112:PRO:HD2	2.20	0.57
1:A:178:THR:HG21	1:A:184:TRP:CZ2	2.37	0.57
1:A:63:GLU:OE1	2:A:353:HOH:O	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:VAL:HG22	1:B:208:ILE:HD12	1.86	0.56
1:A:149:ARG:HG2	1:A:151:MET:HE2	1.89	0.55
1:B:296:LYS:HB2	1:B:301:ILE:HD11	1.88	0.54
1:A:102:ASP:HB2	1:A:319:ARG:HD3	1.90	0.54
1:C:34:LYS:NZ	2:C:441:HOH:O	2.39	0.54
1:A:111:MET:HB2	1:A:112:PRO:CD	2.33	0.54
1:A:296:LYS:HD2	2:A:341:HOH:O	2.07	0.53
1:A:151:MET:HB3	1:A:153:ILE:HD11	1.90	0.53
1:A:109:LYS:CD	1:A:111:MET:HB3	2.40	0.52
1:A:111:MET:CB	1:A:112:PRO:CD	2.87	0.52
1:B:296:LYS:HB3	1:B:301:ILE:HD11	1.91	0.51
1:A:109:LYS:HD2	1:A:111:MET:HB3	1.92	0.51
1:C:16:ASN:HD22	1:C:16:ASN:C	2.19	0.51
1:C:324:ILE:HB	2:C:428:HOH:O	2.10	0.51
1:A:240:LEU:HD21	1:C:240:LEU:HD21	1.91	0.51
1:B:189:ARG:NH2	1:C:227:THR:OG1	2.44	0.51
1:A:109:LYS:HB3	1:B:214:SER:HB3	1.92	0.50
1:C:178:THR:HG23	1:C:178:THR:O	2.11	0.50
1:B:11:ALA:N	1:B:14:LYS:HB2	2.26	0.50
1:B:236:LEU:C	1:B:236:LEU:HD13	2.37	0.50
1:B:126:LEU:HD22	1:B:130:LEU:HD11	1.95	0.49
1:C:249:ASP:OD2	1:C:286:LYS:NZ	2.45	0.48
1:C:210:SER:O	1:C:212:LYS:NZ	2.45	0.48
1:C:296:LYS:CB	1:C:301:ILE:HD11	2.44	0.48
1:C:304:VAL:HG13	2:C:431:HOH:O	2.13	0.47
1:A:165:LEU:HD11	1:C:159:TYR:CE1	2.50	0.47
1:A:336:SER:HA	2:A:341:HOH:O	2.14	0.47
1:A:151:MET:N	1:A:151:MET:HE3	2.29	0.47
1:B:267:VAL:HG11	1:B:293:LYS:HD2	1.96	0.47
1:A:96:LYS:HD2	2:A:491:HOH:O	2.14	0.46
1:C:296:LYS:HB2	1:C:301:ILE:HD11	1.96	0.46
1:A:129:ARG:NH2	2:A:343:HOH:O	2.49	0.46
1:C:127:LEU:HG	2:C:428:HOH:O	2.14	0.46
1:C:205:MET:HB2	1:C:207:LYS:HG3	1.97	0.46
1:B:118:LYS:HE3	1:B:122:LEU:HD11	1.98	0.45
1:A:315:ASP:HB3	2:A:425:HOH:O	2.16	0.45
1:C:87:TYR:CZ	1:C:91:MET:SD	3.09	0.45
1:A:230:ARG:HD3	2:A:502:HOH:O	2.17	0.45
1:A:112:PRO:HG3	1:C:219:LEU:HD11	1.97	0.45
1:B:241:LYS:HA	1:B:241:LYS:HE2	1.99	0.45
1:B:274:ARG:CZ	2:B:368:HOH:O	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:MET:HG3	1:A:112:PRO:HD2	1.98	0.45
1:C:213:GLY:O	1:C:214:SER:OG	2.29	0.45
1:C:23:LEU:HG	1:C:62:LEU:HD22	1.99	0.45
1:C:194:LEU:HD21	1:C:210:SER:HB2	1.99	0.44
1:A:16:ASN:HD22	1:A:19:GLN:CD	2.25	0.44
1:A:41:ARG:NH2	2:A:470:HOH:O	2.49	0.44
1:A:264:GLU:O	1:A:267:VAL:HG12	2.18	0.44
1:C:27:VAL:HG22	1:C:79:PRO:HB3	1.99	0.44
1:C:46:SER:OG	1:C:48:ILE:HG12	2.18	0.44
1:A:153:ILE:HG21	1:C:153:ILE:HG21	2.00	0.44
1:C:11:ALA:N	2:C:464:HOH:O	2.51	0.44
1:C:85:ARG:NH2	2:C:465:HOH:O	2.50	0.44
1:A:272:THR:HG23	2:A:462:HOH:O	2.18	0.43
1:C:87:TYR:OH	1:C:91:MET:SD	2.69	0.43
1:A:327:ARG:N	1:A:327:ARG:HD2	2.33	0.42
1:C:11:ALA:N	2:C:445:HOH:O	2.52	0.42
1:A:23:LEU:HG	1:A:62:LEU:HD22	2.01	0.42
1:A:240:LEU:HD23	1:A:250:ILE:HD13	2.01	0.42
1:B:141:ASN:HD22	1:B:142:THR:N	2.18	0.42
1:C:16:ASN:ND2	1:C:19:GLN:H	2.17	0.42
1:C:127:LEU:HD13	1:C:135:VAL:HG22	2.01	0.42
1:C:127:LEU:HB3	1:C:135:VAL:HG21	2.02	0.42
1:B:12:LEU:N	1:B:12:LEU:HD22	2.34	0.41
1:A:136:LEU:HD11	1:A:257:VAL:HG13	2.02	0.41
1:B:246:THR:HG22	1:B:248:GLU:H	1.84	0.41
1:A:27:VAL:HG22	1:A:79:PRO:HB3	2.03	0.41
1:B:11:ALA:HA	1:B:67:TYR:OH	2.20	0.41
1:B:196:ARG:O	1:C:218:ARG:NH1	2.54	0.41
1:A:72:HIS:HB2	1:A:73:THR:H	1.72	0.41
1:A:135:VAL:O	1:A:233:ASP:HA	2.21	0.41
1:C:111:MET:N	1:C:112:PRO:CD	2.82	0.41
1:A:109:LYS:HD2	1:A:111:MET:SD	2.61	0.40
1:B:21:LYS:HD2	1:B:48:ILE:HD11	2.03	0.40
1:B:246:THR:HG22	1:B:247:LEU:N	2.37	0.40
1:C:160:LEU:C	1:C:160:LEU:HD12	2.46	0.40
1:A:130:LEU:HD21	1:B:182:LEU:HD13	2.04	0.40
1:B:212:LYS:O	1:B:218:ARG:NH1	2.55	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	321/338 (95%)	312 (97%)	6 (2%)	3 (1%)	14	14
1	B	307/338 (91%)	299 (97%)	7 (2%)	1 (0%)	36	42
1	C	307/338 (91%)	294 (96%)	11 (4%)	2 (1%)	18	19
All	All	935/1014 (92%)	905 (97%)	24 (3%)	6 (1%)	21	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	111	MET
1	C	71	PRO
1	B	71	PRO
1	C	269	GLU
1	A	100	GLU
1	A	71	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	A	294/308 (96%)	263 (90%)	31 (10%)	6	6
1	B	283/308 (92%)	253 (89%)	30 (11%)	6	6
1	C	283/308 (92%)	257 (91%)	26 (9%)	8	9
All	All	860/924 (93%)	773 (90%)	87 (10%)	7	7

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASP
1	A	23	LEU
1	A	42	VAL
1	A	49	GLU
1	A	60	LYS
1	A	74	SER
1	A	91	MET
1	A	92	LEU
1	A	98	THR
1	A	103	LEU
1	A	110	SER
1	A	127	LEU
1	A	145	LEU
1	A	151	MET
1	A	165	LEU
1	A	166	THR
1	A	219	LEU
1	A	220	ILE
1	A	230	ARG
1	A	241	LYS
1	A	247	LEU
1	A	248	GLU
1	A	253	LEU
1	A	255	GLU
1	A	257	VAL
1	A	261	LYS
1	A	286	LYS
1	A	289	VAL
1	A	306	LEU
1	A	324	ILE
1	A	327	ARG
1	B	12	LEU
1	B	21	LYS
1	B	23	LEU
1	B	34	LYS
1	B	43	LEU
1	B	60	LYS
1	B	95	SER
1	B	118	LYS
1	B	126	LEU
1	B	127	LEU
1	B	148	LEU

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Mol	Chain	Res	Type
1	B	151	MET
1	B	155	VAL
1	B	160	LEU
1	B	181	ARG
1	B	194	LEU
1	B	198	ARG
1	B	209	GLU
1	B	219	LEU
1	B	247	LEU
1	B	253	LEU
1	B	257	VAL
1	B	266	LEU
1	B	272	THR
1	B	273	VAL
1	B	274	ARG
1	B	289	VAL
1	B	306	LEU
1	B	319	ARG
1	B	324	ILE
1	C	12	LEU
1	C	16	ASN
1	C	23	LEU
1	C	29	GLU
1	C	118	LYS
1	C	127	LEU
1	C	143	ARG
1	C	145	LEU
1	C	165	LEU
1	C	166	THR
1	C	180	GLU
1	C	185	GLU
1	C	194	LEU
1	C	212	LYS
1	C	217	LEU
1	C	236	LEU
1	C	243	GLU
1	C	247	LEU
1	C	248	GLU
1	C	253	LEU
1	C	257	VAL
1	C	267	VAL
1	C	286	LYS

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Mol	Chain	Res	Type
1	C	306	LEU
1	C	310	LYS
1	C	324	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	57	ASN
1	A	70	GLN
1	B	125	ASN
1	B	141	ASN
1	B	179	GLN
1	B	183	ASN
1	B	318	HIS
1	C	16	ASN
1	C	125	ASN
1	C	141	ASN
1	C	179	GLN
1	C	260	GLN

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 oligosaccharides 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

6.1 Protein, DNA and RNA chains

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	A	323/338 (95%)	0.61	29 (8%)	15 12	15, 22, 34, 57	0
1	B	311/338 (92%)	-0.01	11 (3%)	47 44	13, 21, 31, 56	0
1	C	311/338 (92%)	0.44	19 (6%)	27 24	10, 22, 30, 48	0
All	All	945/1014 (93%)	0.35	59 (6%)	26 23	10, 22, 32, 57	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	211	LEU	4.6
1	A	144	ASP	4.2
1	C	268	GLY	3.6
1	B	269	GLU	3.6
1	A	92	LEU	3.5
1	B	11	ALA	3.4
1	A	85	ARG	3.3
1	A	40	GLN	3.3
1	A	97	GLU	3.2
1	C	210	SER	3.2
1	C	336	SER	3.2
1	C	92	LEU	3.2
1	C	209	GLU	3.1
1	A	88	TYR	3.0
1	A	98	THR	3.0
1	A	75	ALA	3.0
1	A	45	VAL	3.0
1	B	13	LYS	2.9
1	A	47	ASN	2.9
1	B	242	ASP	2.8
1	B	12	LEU	2.8
1	A	111	MET	2.8
1	B	111	MET	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	111	MET	2.8
1	A	179	GLN	2.7
1	B	95	SER	2.7
1	A	83	GLY	2.7
1	C	91	MET	2.6
1	A	81	ASP	2.6
1	A	80	THR	2.6
1	C	213	GLY	2.6
1	B	268	GLY	2.6
1	C	155	VAL	2.6
1	A	193	PHE	2.6
1	C	330	GLU	2.5
1	C	32	GLU	2.4
1	B	227	THR	2.4
1	C	45	VAL	2.4
1	A	65	LEU	2.4
1	A	49	GLU	2.4
1	B	91	MET	2.3
1	A	94	ILE	2.3
1	A	84	LEU	2.3
1	A	44	GLU	2.3
1	A	64	TYR	2.3
1	C	269	GLU	2.3
1	A	52	SER	2.2
1	C	15	LEU	2.2
1	B	336	SER	2.2
1	C	28	ARG	2.1
1	A	48	ILE	2.1
1	C	94	ILE	2.1
1	A	269	GLU	2.1
1	A	51	SER	2.1
1	C	189	ARG	2.1
1	A	100	GLU	2.0
1	A	181	ARG	2.0
1	A	95	SER	2.0
1	C	212	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.