



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 10:42 PM UTC

PDB ID : 3VZ8 / pdb_00003vz8
Title : Crystal Structure Analysis of the Mini-chaperonin variant with Leu 185, Val 186, Pro 187, Arg 188 and Ser 190 replaced with all Gly
Authors : Saijo, S.; Sato, T.
Deposited on : 2012-10-09
Resolution : 1.90 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

<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

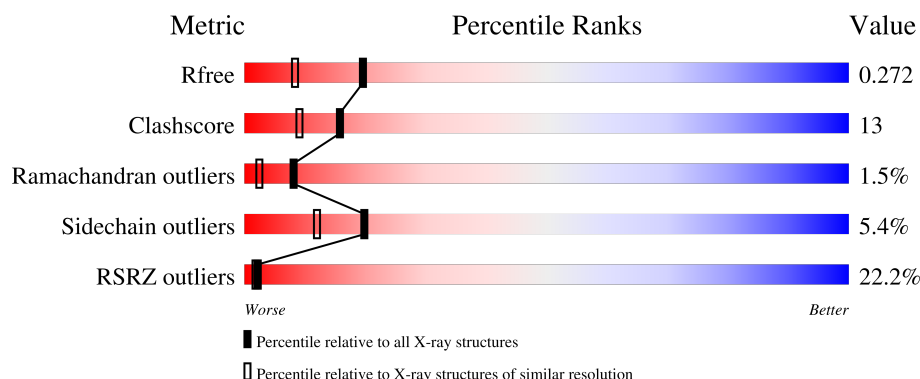
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	199	22% 66% 23% 7%
1	B	199	23% 62% 27% 7%
1	C	199	17% 72% 18% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4396 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 60 kDa chaperonin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1401	881	240	275	5			
1	B	185	Total	C	N	O	S	0	0	0
			1401	881	240	275	5			
1	C	185	Total	C	N	O	S	0	0	0
			1401	881	240	275	5			

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	178	HIS	-	expression tag	UNP P0A6F5
A	179	HIS	-	expression tag	UNP P0A6F5
A	180	HIS	-	expression tag	UNP P0A6F5
A	181	HIS	-	expression tag	UNP P0A6F5
A	182	HIS	-	expression tag	UNP P0A6F5
A	183	HIS	-	expression tag	UNP P0A6F5
A	184	GLY	-	SEE REMARK 999	UNP P0A6F5
A	185	GLY	-	SEE REMARK 999	UNP P0A6F5
A	186	GLY	-	SEE REMARK 999	UNP P0A6F5
A	187	GLY	-	SEE REMARK 999	UNP P0A6F5
A	188	GLY	-	SEE REMARK 999	UNP P0A6F5
A	189	GLY	-	SEE REMARK 999	UNP P0A6F5
A	190	GLY	-	SEE REMARK 999	UNP P0A6F5
B	178	HIS	-	expression tag	UNP P0A6F5
B	179	HIS	-	expression tag	UNP P0A6F5
B	180	HIS	-	expression tag	UNP P0A6F5
B	181	HIS	-	expression tag	UNP P0A6F5
B	182	HIS	-	expression tag	UNP P0A6F5
B	183	HIS	-	expression tag	UNP P0A6F5
B	184	GLY	-	SEE REMARK 999	UNP P0A6F5
B	185	GLY	-	SEE REMARK 999	UNP P0A6F5
B	186	GLY	-	SEE REMARK 999	UNP P0A6F5
B	187	GLY	-	SEE REMARK 999	UNP P0A6F5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	188	GLY	-	SEE REMARK 999	UNP P0A6F5
B	189	GLY	-	SEE REMARK 999	UNP P0A6F5
B	190	GLY	-	SEE REMARK 999	UNP P0A6F5
C	178	HIS	-	expression tag	UNP P0A6F5
C	179	HIS	-	expression tag	UNP P0A6F5
C	180	HIS	-	expression tag	UNP P0A6F5
C	181	HIS	-	expression tag	UNP P0A6F5
C	182	HIS	-	expression tag	UNP P0A6F5
C	183	HIS	-	expression tag	UNP P0A6F5
C	184	GLY	-	SEE REMARK 999	UNP P0A6F5
C	185	GLY	-	SEE REMARK 999	UNP P0A6F5
C	186	GLY	-	SEE REMARK 999	UNP P0A6F5
C	187	GLY	-	SEE REMARK 999	UNP P0A6F5
C	188	GLY	-	SEE REMARK 999	UNP P0A6F5
C	189	GLY	-	SEE REMARK 999	UNP P0A6F5
C	190	GLY	-	SEE REMARK 999	UNP P0A6F5

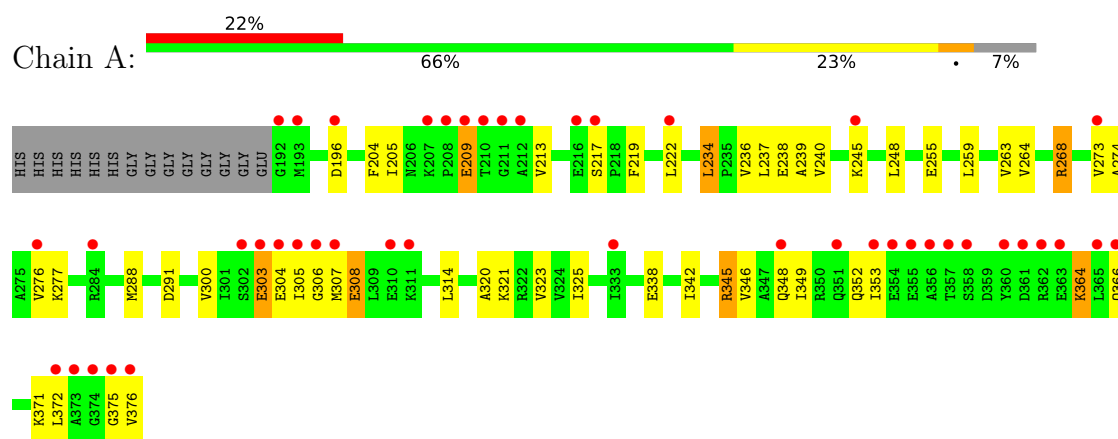
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	62	Total O 62 62	0	0
2	B	64	Total O 64 64	0	0
2	C	67	Total O 67 67	0	0

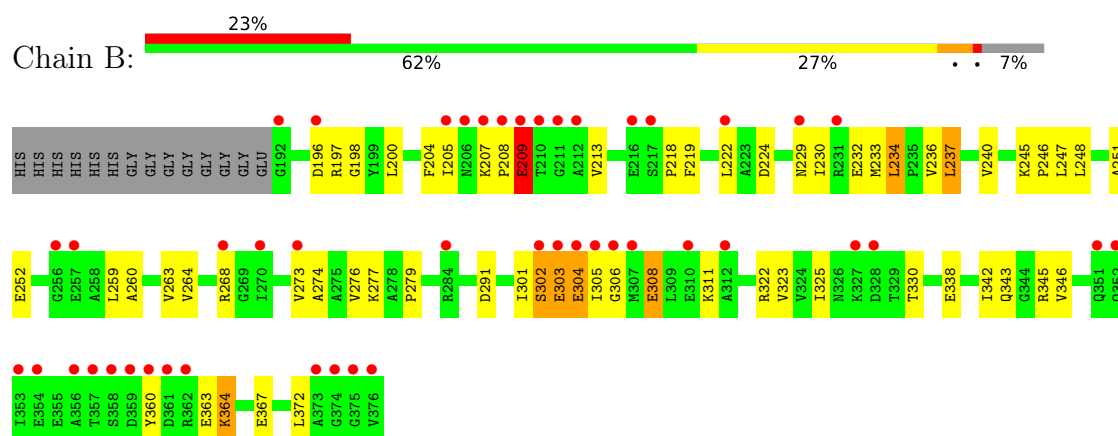
3 Residue-property plots [i](#)

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.

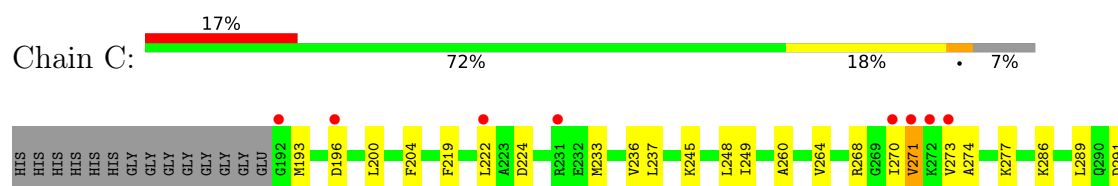
- Molecule 1: 60 kDa chaperonin

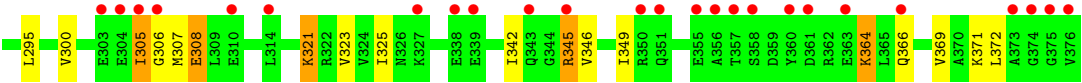


- Molecule 1: 60 kDa chaperonin



- Molecule 1: 60 kDa chaperonin





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.06Å 38.01Å 102.20Å 90.00° 116.23° 90.00°	Depositor
Resolution (Å)	47.51 – 1.90 47.51 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.4 (47.51-1.90) 99.5 (47.51-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 1.90Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.263 , 0.272 0.263 , 0.272	Depositor DCC
R_{free} test set	2605 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4396	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.54% 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 ⓘ

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	0.41	0/1413	0.93	3/1903 (0.2%)
1	B	0.40	0/1413	0.94	5/1903 (0.3%)
1	C	0.40	0/1413	0.92	4/1903 (0.2%)
All	All	0.40	0/4239	0.93	12/5709 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	219	PHE	N-CA-C	-8.12	98.50	110.52
1	B	204	PHE	N-CA-C	-7.38	104.42	113.50
1	A	204	PHE	N-CA-C	-7.36	104.45	113.50
1	C	204	PHE	N-CA-C	-7.27	104.21	113.23
1	A	219	PHE	N-CA-C	-7.24	99.81	110.52
1	C	219	PHE	N-CA-C	-6.81	100.44	110.52
1	A	196	ASP	N-CA-C	6.21	119.31	110.68
1	B	196	ASP	N-CA-C	5.92	119.88	110.70
1	C	196	ASP	N-CA-C	5.90	119.84	110.70
1	C	289	LEU	N-CA-C	-5.26	105.62	111.36
1	B	207	LYS	CA-C-N	5.13	126.25	119.84
1	B	207	LYS	C-N-CA	5.13	126.25	119.84

There are no chirality outliers.

There are no planarity outliers.

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	A	1401	0	1455	37	0
1	B	1401	0	1455	45	0
1	C	1401	0	1455	32	0
2	A	62	0	0	2	0
2	B	64	0	0	1	0
2	C	67	0	0	1	0
All	All	4396	0	4365	114	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 13.

All (114) 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:364:LYS:HZ2	1:A:364:LYS:HA	1.22	1.02
1:C:364:LYS:NZ	1:C:364:LYS:HA	1.94	0.82
1:C:233:MET:HE3	1:C:237:LEU:HD11	1.61	0.81
1:B:200:LEU:HD21	1:B:277:LYS:HG3	1.63	0.80
1:B:198:GLY:O	1:B:276:VAL:HG23	1.82	0.79
1:C:224:ASP:OD1	1:C:286:LYS:HD3	1.84	0.78
1:A:364:LYS:HA	1:A:364:LYS:NZ	2.01	0.76
1:A:248:LEU:HD22	1:A:323:VAL:HG11	1.70	0.74
1:C:364:LYS:HA	1:C:364:LYS:HZ3	1.53	0.71
1:B:364:LYS:HZ3	1:B:364:LYS:HA	1.57	0.69
1:B:364:LYS:HA	1:B:364:LYS:NZ	2.08	0.69
1:C:233:MET:O	1:C:237:LEU:HD13	1.93	0.68
1:C:248:LEU:HD22	1:C:323:VAL:HG11	1.74	0.68
1:A:345:ARG:HG2	1:A:372:LEU:CD1	2.24	0.67
1:A:345:ARG:HG2	1:A:372:LEU:HD11	1.76	0.65
1:B:224:ASP:HB3	1:B:302:SER:HB3	1.79	0.64
1:A:305:ILE:HG13	1:A:306:GLY:H	1.62	0.64
1:C:273:VAL:HG22	1:C:274:ALA:N	2.13	0.64
1:C:233:MET:HE3	1:C:237:LEU:CD1	2.29	0.62
1:A:240:VAL:CG1	1:A:245:LYS:O	2.50	0.59
1:A:300:VAL:O	1:A:307:MET:HE3	2.03	0.59
1:A:273:VAL:HG22	1:A:274:ALA:N	2.18	0.58
1:C:308:GLU:H	1:C:308:GLU:CD	2.10	0.58
1:B:291:ASP:OD1	1:B:345:ARG:HD3	2.04	0.57
1:A:291:ASP:OD1	1:A:345:ARG:HD3	2.04	0.57
1:A:234:LEU:HD22	1:A:238:GLU:HG2	1.86	0.57
1:B:248:LEU:HD13	1:B:325:ILE:HD11	1.85	0.57
1:B:209:GLU:H	1:B:209:GLU:CD	2.12	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:LEU:HD22	1:B:323:VAL:HG11	1.86	0.55
1:B:342:ILE:O	1:B:346:VAL:HG23	2.06	0.55
1:B:240:VAL:HG13	1:B:245:LYS:O	2.05	0.55
1:B:306:GLY:HA2	1:B:308:GLU:OE2	2.06	0.55
1:C:273:VAL:HG22	1:C:274:ALA:H	1.72	0.55
1:B:233:MET:O	1:B:236:VAL:HG22	2.07	0.54
1:A:217:SER:O	1:A:245:LYS:HD3	2.06	0.54
1:C:236:VAL:HG23	1:C:237:LEU:HD12	1.90	0.54
1:A:276:VAL:HG12	1:A:277:LYS:O	2.07	0.54
1:C:248:LEU:HD13	1:C:325:ILE:HD11	1.90	0.53
1:A:205:ILE:HA	1:A:213:VAL:HG22	1.91	0.53
1:C:364:LYS:HA	1:C:364:LYS:HZ2	1.72	0.52
1:A:305:ILE:O	1:A:308:GLU:OE1	2.28	0.52
1:C:200:LEU:HD21	1:C:277:LYS:HG3	1.90	0.52
1:A:222:LEU:HD22	1:A:222:LEU:N	2.25	0.51
1:C:291:ASP:OD1	1:C:345:ARG:HD3	2.10	0.51
1:A:375:GLY:O	1:A:376:VAL:C	2.53	0.51
1:A:314:LEU:HD11	2:A:446:HOH:O	2.11	0.51
1:B:208:PRO:HG2	1:B:209:GLU:OE2	2.11	0.50
1:C:193:MET:HG3	1:C:371:LYS:HB3	1.93	0.50
1:B:205:ILE:HA	1:B:213:VAL:HG22	1.93	0.50
1:B:240:VAL:CG1	1:B:245:LYS:O	2.59	0.50
1:B:247:LEU:O	1:B:273:VAL:HG23	2.11	0.50
1:C:300:VAL:O	1:C:307:MET:HE3	2.11	0.50
1:C:321:LYS:HB2	1:C:321:LYS:NZ	2.26	0.50
1:B:197:ARG:HE	1:B:279:PRO:HA	1.78	0.49
1:B:308:GLU:HG2	1:B:311:LYS:HB2	1.94	0.49
1:C:222:LEU:HD22	1:C:222:LEU:N	2.27	0.49
1:A:288:MET:CE	1:A:371:LYS:HZ1	2.26	0.49
1:B:308:GLU:H	1:B:308:GLU:CD	2.19	0.49
1:A:248:LEU:HD13	1:A:325:ILE:HD11	1.94	0.49
1:B:301:ILE:O	1:B:302:SER:CB	2.60	0.48
1:B:345:ARG:HG2	1:B:372:LEU:HD11	1.96	0.48
1:C:270:ILE:HG22	1:C:271:VAL:HG13	1.95	0.48
1:A:303:GLU:O	1:A:304:GLU:C	2.56	0.48
1:A:308:GLU:H	1:A:308:GLU:CD	2.22	0.48
1:B:360:TYR:CZ	1:B:364:LYS:HD2	2.50	0.47
1:C:236:VAL:HG23	1:C:237:LEU:CD1	2.44	0.47
1:B:343:GLN:NE2	1:B:343:GLN:HA	2.30	0.47
1:A:240:VAL:HG12	1:A:245:LYS:O	2.14	0.47
1:A:273:VAL:CG2	1:A:274:ALA:N	2.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:LEU:O	1:B:263:VAL:HG23	2.16	0.46
1:A:239:ALA:HB1	1:A:314:LEU:HG	1.97	0.45
1:C:345:ARG:HG2	1:C:372:LEU:HD11	1.98	0.45
1:C:342:ILE:O	1:C:346:VAL:HG23	2.17	0.45
1:C:193:MET:HE3	1:C:295:LEU:HD22	1.98	0.45
1:A:342:ILE:O	1:A:346:VAL:HG23	2.17	0.45
1:A:306:GLY:HA2	1:A:308:GLU:OE2	2.17	0.44
1:B:273:VAL:HG22	1:B:274:ALA:N	2.33	0.44
1:A:348:GLN:O	1:A:352:GLN:HG3	2.18	0.44
1:C:305:ILE:HG13	2:C:456:HOH:O	2.16	0.44
1:B:260:ALA:O	1:B:264:VAL:HG23	2.18	0.44
1:C:273:VAL:CG2	1:C:274:ALA:N	2.80	0.44
1:C:236:VAL:HG23	1:C:237:LEU:N	2.33	0.43
1:C:260:ALA:O	1:C:264:VAL:HG23	2.19	0.43
1:A:217:SER:HA	1:A:320:ALA:O	2.19	0.43
1:B:325:ILE:HG12	1:B:330:THR:HG23	2.00	0.43
1:C:245:LYS:HA	1:C:245:LYS:HD3	1.84	0.43
1:C:349:ILE:HG21	1:C:369:VAL:HG23	2.01	0.43
1:B:364:LYS:HZ3	1:B:367:GLU:CD	2.26	0.43
1:B:234:LEU:HD23	1:B:234:LEU:HA	1.88	0.43
1:B:342:ILE:HG23	1:B:372:LEU:HD22	2.00	0.43
1:B:229:ASN:O	1:B:232:GLU:HG2	2.19	0.42
1:A:264:VAL:HG12	1:A:268:ARG:CZ	2.49	0.42
1:B:322:ARG:HD2	2:B:462:HOH:O	2.19	0.42
1:C:233:MET:HE2	1:C:249:ILE:HD13	2.01	0.42
1:A:305:ILE:HG13	1:A:306:GLY:N	2.33	0.42
1:B:363:GLU:O	1:B:367:GLU:HG3	2.20	0.42
1:A:209:GLU:H	1:A:209:GLU:CD	2.27	0.42
1:B:264:VAL:O	1:B:268:ARG:HG3	2.19	0.41
1:B:209:GLU:CD	1:B:209:GLU:N	2.78	0.41
1:B:251:ALA:O	1:B:252:GLU:C	2.64	0.41
1:B:301:ILE:O	1:B:302:SER:HB3	2.20	0.41
1:A:236:VAL:HG22	2:A:404:HOH:O	2.19	0.41
1:B:208:PRO:O	1:B:209:GLU:C	2.62	0.41
1:B:308:GLU:CG	1:B:311:LYS:HD2	2.51	0.41
1:C:345:ARG:HG2	1:C:372:LEU:CD1	2.51	0.41
1:A:349:ILE:O	1:A:353:ILE:HG13	2.20	0.41
1:B:218:PRO:HB3	1:B:246:PRO:HG2	2.02	0.41
1:A:259:LEU:O	1:A:263:VAL:HG23	2.20	0.41
1:B:345:ARG:HG2	1:B:372:LEU:CD1	2.50	0.41
1:A:264:VAL:CG1	1:A:268:ARG:NH1	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:VAL:HG22	1:B:277:LYS:N	2.36	0.41
1:B:303:GLU:HB2	1:B:304:GLU:H	1.63	0.40
1:A:321:LYS:HB2	1:A:321:LYS:HE3	1.81	0.40
1:B:230:ILE:HD11	1:B:237:LEU:HD21	2.04	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	170/199 (85%)	165 (97%)	4 (2%)	1 (1%)	21	13
1	B	183/199 (92%)	170 (93%)	8 (4%)	5 (3%)	4	1
1	C	183/199 (92%)	174 (95%)	7 (4%)	2 (1%)	11	4
All	All	536/597 (90%)	509 (95%)	19 (4%)	8 (2%)	8	2

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	302	SER
1	C	305	ILE
1	B	305	ILE
1	B	209	GLU
1	C	306	GLY
1	A	303	GLU
1	B	303	GLU
1	B	304	GLU

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	147/154 (96%)	137 (93%)	10 (7%)	14	7
1	B	147/154 (96%)	140 (95%)	7 (5%)	23	15
1	C	147/154 (96%)	140 (95%)	7 (5%)	23	15
All	All	441/462 (96%)	417 (95%)	24 (5%)	20	12

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	209	GLU
1	A	234	LEU
1	A	237	LEU
1	A	255	GLU
1	A	268	ARG
1	A	308	GLU
1	A	338	GLU
1	A	345	ARG
1	A	364	LYS
1	A	366	GLN
1	B	209	GLU
1	B	222	LEU
1	B	234	LEU
1	B	237	LEU
1	B	308	GLU
1	B	338	GLU
1	B	364	LYS
1	C	268	ARG
1	C	271	VAL
1	C	308	GLU
1	C	321	LYS
1	C	345	ARG
1	C	364	LYS
1	C	366	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	GLN
1	A	319	GLN
1	A	343	GLN
1	B	265	ASN
1	B	343	GLN
1	B	366	GLN
1	C	194	GLN
1	C	265	ASN
1	C	352	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	185/199 (92%)	1.29	44 (23%) 2 1	9, 22, 47, 54	4 (2%)
1	B	185/199 (92%)	1.55	46 (24%) 2 1	10, 24, 52, 61	4 (2%)
1	C	185/199 (92%)	1.15	33 (17%) 4 3	8, 22, 40, 62	4 (2%)
All	All	555/597 (92%)	1.33	123 (22%) 2 2	8, 23, 47, 62	12 (2%)

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	208	PRO	10.1
1	C	376	VAL	8.6
1	B	376	VAL	8.2
1	C	305	ILE	7.9
1	B	305	ILE	7.8
1	C	375	GLY	7.5
1	B	361	ASP	7.4
1	A	305	ILE	7.0
1	A	376	VAL	7.0
1	B	210	THR	6.5
1	B	284	ARG	5.9
1	A	303	GLU	5.9
1	A	356	ALA	5.4
1	B	304	GLU	5.4
1	A	304	GLU	5.2
1	B	209	GLU	5.1
1	C	356	ALA	5.0
1	A	361	ASP	5.0
1	B	211	GLY	5.0
1	B	303	GLU	4.9
1	C	374	GLY	4.9
1	B	375	GLY	4.7
1	B	192	GLY	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	273	VAL	4.6
1	C	351	GLN	4.4
1	C	361	ASP	4.4
1	A	354	GLU	4.3
1	B	207	LYS	4.2
1	B	358	SER	4.2
1	A	375	GLY	4.1
1	A	210	THR	4.1
1	A	357	THR	4.1
1	A	358	SER	4.0
1	A	353	ILE	4.0
1	B	374	GLY	4.0
1	A	192	GLY	3.9
1	B	360	TYR	3.9
1	A	351	GLN	3.9
1	B	357	THR	3.8
1	B	373	ALA	3.8
1	B	353	ILE	3.8
1	A	208	PRO	3.7
1	B	268	ARG	3.7
1	B	196	ASP	3.7
1	C	360	TYR	3.7
1	B	205	ILE	3.7
1	B	354	GLU	3.7
1	C	310	GLU	3.7
1	B	306	GLY	3.7
1	B	222	LEU	3.6
1	C	366	GLN	3.6
1	A	374	GLY	3.5
1	A	211	GLY	3.5
1	B	212	ALA	3.5
1	A	302	SER	3.5
1	B	356	ALA	3.5
1	C	373	ALA	3.5
1	B	229	ASN	3.4
1	B	310	GLU	3.4
1	C	339	GLU	3.4
1	A	362	ARG	3.4
1	A	207	LYS	3.3
1	C	306	GLY	3.3
1	A	363	GLU	3.2
1	A	310	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	373	ALA	3.1
1	C	222	LEU	3.1
1	A	209	GLU	3.1
1	A	196	ASP	3.0
1	A	360	TYR	3.0
1	C	304	GLU	3.0
1	C	271	VAL	3.0
1	B	256	GLY	3.0
1	C	196	ASP	2.9
1	C	270	ILE	2.9
1	A	348	GLN	2.9
1	B	302	SER	2.9
1	A	273	VAL	2.8
1	C	338	GLU	2.8
1	B	206	ASN	2.8
1	B	231	ARG	2.8
1	A	307	MET	2.8
1	A	222	LEU	2.7
1	B	307	MET	2.7
1	B	273	VAL	2.7
1	C	231	ARG	2.7
1	C	357	THR	2.6
1	C	363	GLU	2.6
1	A	212	ALA	2.6
1	A	276	VAL	2.6
1	B	216	GLU	2.6
1	C	343	GLN	2.6
1	A	355	GLU	2.5
1	A	284	ARG	2.5
1	B	351	GLN	2.5
1	B	217	SER	2.5
1	C	355	GLU	2.4
1	C	358	SER	2.4
1	B	352	GLN	2.3
1	B	312	ALA	2.3
1	A	193	MET	2.3
1	A	372	LEU	2.3
1	C	192	GLY	2.3
1	C	303	GLU	2.3
1	A	333	ILE	2.3
1	B	328	ASP	2.3
1	B	270	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	327	LYS	2.2
1	A	366	GLN	2.2
1	C	345	ARG	2.2
1	C	350	ARG	2.2
1	B	362	ARG	2.2
1	B	359	ASP	2.1
1	B	327	LYS	2.1
1	B	257	GLU	2.1
1	A	245	LYS	2.1
1	A	365	LEU	2.1
1	A	311	LYS	2.1
1	C	272	LYS	2.1
1	A	306	GLY	2.1
1	A	217	SER	2.0
1	C	314	LEU	2.0
1	A	216	GLU	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 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.