



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:03 PM GMT

PDB ID : 3VZ8
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
<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

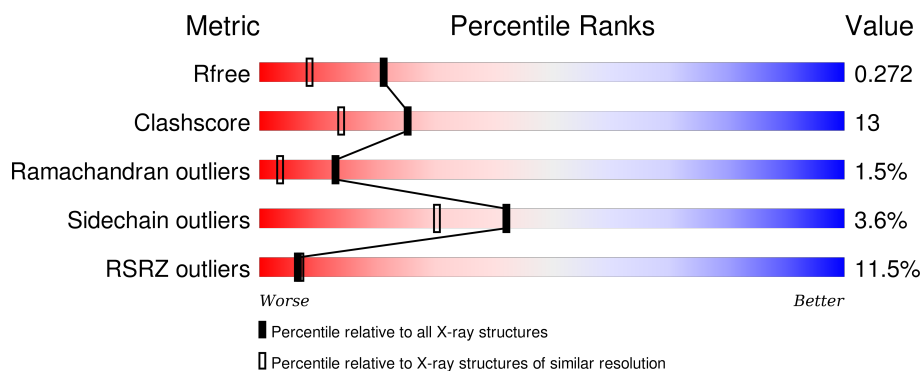
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (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 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	A	199	 11% 70% 21% • 7%
1	B	199	 13% 65% 24% • • 7%
1	C	199	 9% 76% 15% • 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

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%)	DCC
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.1	EDS
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	4 of 52211 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.91	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, 0.375 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	A	0.34	0/1413	0.64	0/1903
1	B	0.33	0/1413	0.64	0/1903
1	C	0.32	0/1413	0.63	0/1903
All	All	0.33	0/4239	0.64	0/5709

There are no bond length outliers.

There are no bond angle outliers.

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	1401	0	1455	37	0
1	B	1401	0	1455	42	0
1	C	1401	0	1455	29	0
2	A	62	0	0	3	0
2	B	64	0	0	1	0
2	C	67	0	0	1	0
All	All	4396	0	4365	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 13.

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:364:LYS:HA	1:C:364:LYS:NZ	1.94	0.82
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:NZ	1:A:364:LYS:HA	2.01	0.76
1:A:248:LEU:HD22	1:A:323:VAL:HG11	1.70	0.74
1:B:364:LYS:NZ	1:B:364:LYS:HA	2.08	0.69
1:C:248:LEU:HD22	1:C:323:VAL:HG11	1.74	0.68
1:C:233:MET:O	1:C:237:LEU:HD13	1.93	0.68
1:A:364:LYS:HZ2	1:A:364:LYS:HA	1.59	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:364:LYS:HA	1:C:364:LYS:HZ2	1.60	0.64
1:C:273:VAL:HG22	1:C:274:ALA:N	2.13	0.64
1:C:364:LYS:HZ3	1:C:364:LYS:HA	1.65	0.60
1:A:240:VAL:CG1	1:A:245:LYS:O	2.50	0.59
1:A:273:VAL:HG22	1:A:274:ALA:N	2.18	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: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:B:364:LYS:HZ3	1:B:364:LYS:HA	1.71	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:308:GLU:H	1:C:308:GLU:CD	2.10	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: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:N	1:A:222:LEU:HD22	2.25	0.51
1:B:209:GLU:H	1:B:209:GLU:CD	2.12	0.51
1:C:291:ASP:OD1	1:C:345:ARG:HD3	2.10	0.51
1:A:314:LEU:HD11	2:A:446:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:240:VAL:CG1	1:B:245:LYS:O	2.59	0.50
1:B:205:ILE:HA	1:B:213:VAL:HG22	1.93	0.50
1:B:247:LEU:O	1:B:273:VAL:HG23	2.11	0.50
1:C:321:LYS:NZ	1:C:321:LYS:HB2	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:N	1:C:222:LEU:HD22	2.27	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:307:MET:HE3	2:A:440:HOH:O	2.13	0.47
1:B:360:TYR:CZ	1:B:364:LYS:HD2	2.50	0.47
1:B:364:LYS:HZ3	1:B:367:GLU:CD	2.17	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
1:A:375:GLY:O	1:A:376:VAL:C	2.53	0.46
1:B:259:LEU:O	1:B:263:VAL:HG23	2.16	0.46
1:A:364:LYS:HZ3	1:A:364:LYS:HA	1.78	0.45
1:C:345:ARG:HG2	1:C:372:LEU:HD11	1.98	0.45
1:A:239:ALA:HB1	1:A:314:LEU:HG	1.97	0.45
1:C:342:ILE:O	1:C:346:VAL:HG23	2.17	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:A:303:GLU:O	1:A:304:GLU:C	2.56	0.43
1:B:308:GLU:CD	1:B:308:GLU:H	2.19	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:349:ILE:HG21	1:C:369:VAL:HG23	2.01	0.43
1:C:245:LYS:HA	1:C:245:LYS:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:342:ILE:HG23	1:B:372:LEU:HD22	2.00	0.43
1:B:234:LEU:HD23	1:B:234:LEU:HA	1.88	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:C:233:MET:HE3	1:C:237:LEU:HD11	2.02	0.42
1:A:308:GLU:H	1:A:308:GLU:CD	2.22	0.42
1:B:322:ARG:HD2	2:B:462:HOH:O	2.19	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:B:264:VAL:O	1:B:268:ARG:HG3	2.19	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: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:364:LYS:CA	1:A:364:LYS:HZ3	2.33	0.41
1:C:233:MET:HE3	1:C:237:LEU:CD1	2.51	0.41
1:B:218:PRO:HB3	1:B:246:PRO:HG2	2.02	0.41
1:A:349:ILE:O	1:A:353:ILE:HG13	2.20	0.41
1:B:345:ARG:HG2	1:B:372:LEU:CD1	2.50	0.41
1:A:259:LEU:O	1:A:263:VAL:HG23	2.20	0.41
1:B:276:VAL:HG22	1:B:277:LYS:N	2.36	0.41
1:A:264:VAL:CG1	1:A:268:ARG:NH1	2.84	0.41
1:B:303:GLU:HB2	1:B:304:GLU:H	1.63	0.40
1:B:230:ILE:HD11	1:B:237:LEU:HD21	2.04	0.40
1:A:321:LYS:HE3	1:A:321:LYS:HB2	1.81	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	183/199 (92%)	178 (97%)	4 (2%)	1 (0%)	34 21

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	183/199 (92%)	170 (93%)	8 (4%)	5 (3%)	6	1
1	C	183/199 (92%)	174 (95%)	7 (4%)	2 (1%)	17	6
All	All	549/597 (92%)	522 (95%)	19 (4%)	8 (2%)	13	3

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%)	140 (95%)	7 (5%)	31	19
1	B	147/154 (96%)	141 (96%)	6 (4%)	37	25
1	C	147/154 (96%)	144 (98%)	3 (2%)	63	57
All	All	441/462 (96%)	425 (96%)	16 (4%)	42	30

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	234	LEU
1	A	237	LEU
1	A	255	GLU
1	A	338	GLU
1	A	345	ARG
1	A	364	LYS

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Mol	Chain	Res	Type
1	A	366	GLN
1	B	199	TYR
1	B	209	GLU
1	B	234	LEU
1	B	237	LEU
1	B	338	GLU
1	B	364	LYS
1	C	345	ARG
1	C	364	LYS
1	C	366	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	265	ASN
1	A	290	GLN
1	A	319	GLN
1	A	343	GLN
1	B	265	ASN
1	B	319	GLN
1	B	343	GLN
1	B	366	GLN
1	C	265	ASN
1	C	319	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

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%)	0.83	21 (11%) 7 7	12, 22, 47, 54	4 (2%)
1	B	185/199 (92%)	0.91	26 (14%) 4 4	13, 24, 52, 61	4 (2%)
1	C	185/199 (92%)	0.76	17 (9%) 11 12	11, 22, 40, 62	4 (2%)
All	All	555/597 (92%)	0.83	64 (11%) 6 7	11, 23, 47, 62	12 (2%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	376	VAL	7.6
1	C	305	ILE	7.3
1	C	375	GLY	6.9
1	C	374	GLY	6.8
1	A	376	VAL	6.7
1	B	208	PRO	6.3
1	B	211	GLY	6.1
1	A	305	ILE	6.1
1	C	376	VAL	5.6
1	B	305	ILE	5.5
1	B	375	GLY	4.8
1	C	360	TYR	4.8
1	B	192	GLY	4.7
1	B	210	THR	4.6
1	B	304	GLU	4.5
1	A	304	GLU	4.2
1	B	361	ASP	4.1
1	A	192	GLY	4.0
1	A	303	GLU	3.9
1	A	357	THR	3.9
1	A	351	GLN	3.8
1	A	375	GLY	3.7
1	B	209	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	373	ALA	3.6
1	A	354	GLU	3.6
1	B	360	TYR	3.5
1	B	310	GLU	3.5
1	A	356	ALA	3.4
1	A	273	VAL	3.2
1	C	273	VAL	3.2
1	C	351	GLN	3.0
1	C	361	ASP	2.9
1	A	360	TYR	2.9
1	A	222	LEU	2.9
1	A	361	ASP	2.8
1	C	356	ALA	2.7
1	C	339	GLU	2.7
1	C	358	SER	2.6
1	B	222	LEU	2.6
1	B	373	ALA	2.6
1	A	363	GLU	2.6
1	B	284	ARG	2.6
1	C	310	GLU	2.6
1	B	357	THR	2.6
1	A	210	THR	2.5
1	A	373	ALA	2.4
1	A	333	ILE	2.4
1	B	351	GLN	2.3
1	B	273	VAL	2.2
1	B	196	ASP	2.2
1	B	354	GLU	2.2
1	B	268	ARG	2.2
1	B	374	GLY	2.2
1	A	310	GLU	2.2
1	A	233	MET	2.2
1	B	353	ILE	2.1
1	B	231	ARG	2.1
1	B	233	MET	2.1
1	C	233	MET	2.1
1	C	196	ASP	2.1
1	A	358	SER	2.1
1	B	303	GLU	2.0
1	C	357	THR	2.0
1	C	363	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 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.