



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 04:59 PM UTC

PDB ID : 8VBT / pdb_00008vbt
Title : Structure of the monofunctional Staphylococcus aureus PBP1 in its apo form
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Deposited on : 2023-12-12
Resolution : 2.00 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

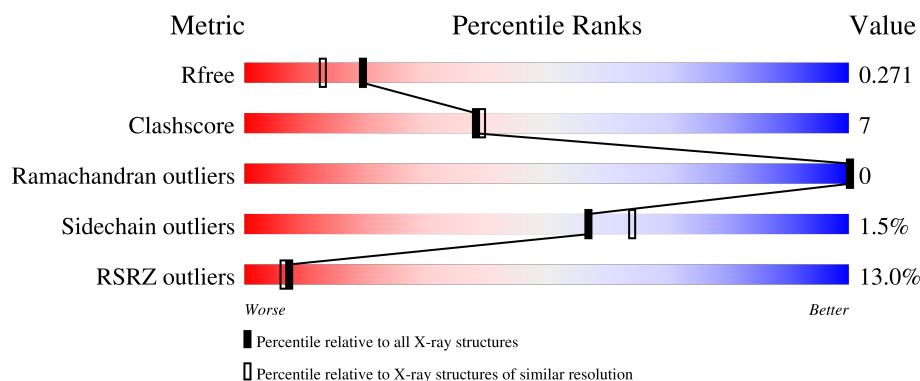
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 ≥ 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	595	10% 72% 13% • 15%
1	B	595	12% 71% 13% 15%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8064 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 Penicillin-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	504	Total	C	N	O	S	0	0	0
			3832	2427	647	743	15			
1	B	505	Total	C	N	O	S	0	1	0
			3846	2436	650	745	15			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	MET	-	initiating methionine	UNP Q2FZ94
A	15	GLY	-	expression tag	UNP Q2FZ94
A	16	SER	-	expression tag	UNP Q2FZ94
A	17	SER	-	expression tag	UNP Q2FZ94
A	18	HIS	-	expression tag	UNP Q2FZ94
A	19	HIS	-	expression tag	UNP Q2FZ94
A	20	HIS	-	expression tag	UNP Q2FZ94
A	21	HIS	-	expression tag	UNP Q2FZ94
A	22	HIS	-	expression tag	UNP Q2FZ94
A	23	HIS	-	expression tag	UNP Q2FZ94
A	24	HIS	-	expression tag	UNP Q2FZ94
A	25	HIS	-	expression tag	UNP Q2FZ94
A	26	HIS	-	expression tag	UNP Q2FZ94
A	27	HIS	-	expression tag	UNP Q2FZ94
A	28	SER	-	expression tag	UNP Q2FZ94
A	29	SER	-	expression tag	UNP Q2FZ94
A	30	GLY	-	expression tag	UNP Q2FZ94
A	31	LEU	-	expression tag	UNP Q2FZ94
A	32	VAL	-	expression tag	UNP Q2FZ94
A	33	PRO	-	expression tag	UNP Q2FZ94
A	34	ARG	-	expression tag	UNP Q2FZ94
A	35	GLY	-	expression tag	UNP Q2FZ94
A	36	SER	-	expression tag	UNP Q2FZ94
A	37	HIS	-	expression tag	UNP Q2FZ94
A	38	MET	-	expression tag	UNP Q2FZ94

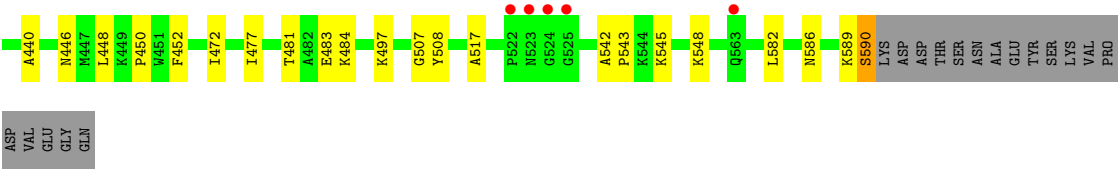
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Chain	Residue	Modelled	Actual	Comment	Reference
A	118	ASP	ASN	conflict	UNP Q2FZ94
B	14	MET	-	initiating methionine	UNP Q2FZ94
B	15	GLY	-	expression tag	UNP Q2FZ94
B	16	SER	-	expression tag	UNP Q2FZ94
B	17	SER	-	expression tag	UNP Q2FZ94
B	18	HIS	-	expression tag	UNP Q2FZ94
B	19	HIS	-	expression tag	UNP Q2FZ94
B	20	HIS	-	expression tag	UNP Q2FZ94
B	21	HIS	-	expression tag	UNP Q2FZ94
B	22	HIS	-	expression tag	UNP Q2FZ94
B	23	HIS	-	expression tag	UNP Q2FZ94
B	24	HIS	-	expression tag	UNP Q2FZ94
B	25	HIS	-	expression tag	UNP Q2FZ94
B	26	HIS	-	expression tag	UNP Q2FZ94
B	27	HIS	-	expression tag	UNP Q2FZ94
B	28	SER	-	expression tag	UNP Q2FZ94
B	29	SER	-	expression tag	UNP Q2FZ94
B	30	GLY	-	expression tag	UNP Q2FZ94
B	31	LEU	-	expression tag	UNP Q2FZ94
B	32	VAL	-	expression tag	UNP Q2FZ94
B	33	PRO	-	expression tag	UNP Q2FZ94
B	34	ARG	-	expression tag	UNP Q2FZ94
B	35	GLY	-	expression tag	UNP Q2FZ94
B	36	SER	-	expression tag	UNP Q2FZ94
B	37	HIS	-	expression tag	UNP Q2FZ94
B	38	MET	-	expression tag	UNP Q2FZ94
B	118	ASP	ASN	conflict	UNP Q2FZ94

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	171	Total O 171 171	0	0
2	B	215	Total O 215 215	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.73Å 72.18Å 86.91Å 90.00° 103.58° 90.00°	Depositor
Resolution (Å)	86.38 – 2.00 86.38 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (86.38-2.00) 99.6 (86.38-2.00)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.235 , 0.273 0.234 , 0.271	Depositor DCC
R_{free} test set	3558 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	1.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8064	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.03% 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	0.33	0/3915	0.51	0/5299
1	B	0.35	0/3930	0.52	0/5320
All	All	0.34	0/7845	0.52	0/10619

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

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	3832	0	3636	45	0
1	B	3846	0	3652	58	0
2	A	171	0	0	6	0
2	B	215	0	0	7	0
All	All	8064	0	7288	103	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 (103) 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:359:ILE:HB	1:A:363:LEU:HD23	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:MET:CE	1:B:376:LEU:HD21	2.23	0.68
1:B:545:LYS:NZ	2:B:705:HOH:O	2.27	0.67
1:A:360:PRO:HD2	1:A:363:LEU:HD23	1.75	0.66
1:B:330:PHE:CD2	1:B:361:MET:HE2	2.31	0.66
1:B:164:LEU:HD13	1:B:165:PRO:HD2	1.77	0.65
1:B:330:PHE:HA	1:B:361:MET:HE1	1.78	0.65
1:B:401:MET:HG3	2:B:876:HOH:O	1.97	0.64
1:A:412:TRP:CE2	1:A:418:GLN:HG2	2.34	0.63
1:A:414:ASN:OD1	1:A:416:LEU:N	2.33	0.60
1:B:361:MET:HE1	1:B:376:LEU:HD21	1.83	0.58
1:B:167:THR:O	1:B:186:LYS:NZ	2.35	0.58
1:B:412:TRP:CD2	1:B:418:GLN:HG2	2.38	0.58
1:B:348:ILE:HB	1:B:374:MET:HE1	1.85	0.57
1:B:508:TYR:CE2	1:B:589:LYS:HG3	2.39	0.57
1:B:420:THR:HA	1:B:423:PHE:CE1	2.40	0.57
1:A:490:ASP:OD1	1:A:494:ASN:ND2	2.31	0.55
1:A:84:LYS:HG3	1:A:143:THR:HG22	1.90	0.54
1:A:340:HIS:HB3	1:A:349:SER:HB3	1.90	0.53
1:B:110:ALA:HA	1:B:124:ILE:HG21	1.90	0.53
1:B:85:LEU:HD12	1:B:153:ILE:HD13	1.90	0.53
1:A:412:TRP:CD2	1:A:418:GLN:HG2	2.43	0.53
1:A:363:LEU:HD11	1:A:367:TYR:CZ	2.44	0.53
1:B:378:ASP:OD1	1:B:419:LYS:NZ	2.43	0.51
1:B:414:ASN:N	2:B:722:HOH:O	2.43	0.51
1:B:477:ILE:HB	1:B:481:THR:HB	1.93	0.51
1:B:483:GLU:HG3	2:B:768:HOH:O	2.11	0.51
1:A:145:LEU:HD11	1:A:163:LEU:HD12	1.93	0.51
1:A:256:LEU:HD22	1:A:268:LEU:HD21	1.93	0.51
1:A:356:TRP:HB2	1:A:359:ILE:HD13	1.94	0.50
1:B:412:TRP:CE2	1:B:418:GLN:HG2	2.46	0.50
1:B:283:TYR:OH	1:B:304:ASP:HB2	2.11	0.49
1:A:341:ARG:HG3	1:A:374:MET:HE3	1.95	0.49
1:A:307:GLN:OE1	2:A:701:HOH:O	2.20	0.49
1:B:325:ILE:HD13	1:B:484:LYS:HB3	1.94	0.49
1:A:484:LYS:NZ	2:A:725:HOH:O	2.46	0.48
1:A:283:TYR:OH	1:A:304:ASP:HB2	2.12	0.48
1:B:330:PHE:HD2	1:B:361:MET:HE2	1.77	0.48
1:B:313:GLY:HA3	1:B:517:ALA:HB2	1.95	0.48
1:A:386:LYS:HD2	1:A:412:TRP:CD1	2.49	0.48
1:A:429:VAL:HG21	1:A:434:MET:SD	2.54	0.48
1:B:324:ALA:CB	1:B:376:LEU:HD22	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:542:ALA:HA	1:B:543:PRO:C	2.38	0.48
1:B:374:MET:HE2	1:B:423:PHE:HE1	1.80	0.47
1:B:321:LEU:HB2	1:B:365:PHE:CE2	2.50	0.47
1:B:243:THR:OG1	1:B:452:PHE:HA	2.15	0.47
1:A:184:ALA:HA	1:A:195:GLY:HA2	1.97	0.47
1:B:317:LYS:HG2	1:B:373:MET:HG3	1.96	0.47
1:B:63:PRO:HB2	1:B:209:GLY:HA3	1.97	0.46
1:B:127:ARG:HB3	1:B:136:ILE:HD13	1.97	0.46
1:B:446:ASN:HB3	1:B:472:ILE:HD12	1.97	0.46
1:A:491:LEU:O	1:A:495:SER:HB2	2.15	0.46
1:A:83:TYR:HB2	1:A:145:LEU:HD12	1.97	0.46
1:A:267:ASP:HB3	1:A:556:SER:HB3	1.96	0.46
1:A:140:ARG:HG2	1:A:141:LYS:HD3	1.98	0.46
1:A:246:SER:O	1:A:250:VAL:HG23	2.15	0.46
1:B:85:LEU:HD11	1:B:113:LEU:HD22	1.97	0.46
1:B:361:MET:HE3	1:B:376:LEU:HD21	1.96	0.46
1:A:315:THR:HA	1:A:513:LYS:HG2	1.99	0.45
1:A:502:ASN:O	1:A:575:LYS:NZ	2.35	0.45
1:B:313:GLY:CA	1:B:517:ALA:HB2	2.46	0.45
1:B:376:LEU:HA	1:B:376:LEU:HD23	1.68	0.45
1:B:330:PHE:HA	1:B:361:MET:CE	2.44	0.45
1:B:184:ALA:HA	1:B:195:GLY:HA2	1.99	0.45
1:B:85:LEU:HA	1:B:163:LEU:HD13	1.99	0.44
1:A:177:ALA:HB3	1:A:181:ILE:HD12	1.97	0.44
1:B:448:LEU:O	1:B:450:PRO:HD3	2.17	0.44
1:A:197:LEU:HD23	1:A:198:GLY:N	2.32	0.44
1:A:532:ASN:HB2	1:A:562:ASP:HB2	2.00	0.44
1:B:497:LYS:HD2	1:B:497:LYS:N	2.32	0.44
1:A:147:TYR:CE2	1:A:151:LEU:HD11	2.52	0.44
1:A:310:TYR:O	1:A:428:THR:HA	2.17	0.44
1:B:103:VAL:HG23	1:B:103:VAL:O	2.18	0.44
1:B:413:SER:HB2	2:B:722:HOH:O	2.18	0.44
1:A:423:PHE:HE1	1:A:425:GLN:HG3	1.83	0.43
1:A:338:SER:HG	1:A:356:TRP:H	1.66	0.43
1:B:586:ASN:ND2	2:B:720:HOH:O	2.42	0.43
1:B:113:LEU:HD21	1:B:161:ILE:HD13	2.00	0.43
1:B:331:ASP:HB3	1:B:334:LYS:HB2	2.01	0.42
1:B:344:MET:HE2	1:B:416:LEU:HB2	2.01	0.42
1:A:572:LYS:HB2	2:A:782:HOH:O	2.20	0.42
1:A:576:PRO:HB2	2:A:837:HOH:O	2.19	0.42
1:B:543:PRO:HB3	2:B:791:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:PHE:CE1	1:A:361:MET:HB3	2.55	0.41
1:B:548:LYS:HA	1:B:548:LYS:HD2	1.63	0.41
1:A:471:GLN:NE2	2:A:728:HOH:O	2.47	0.41
1:B:340:HIS:ND1	1:B:340:HIS:C	2.78	0.41
1:B:582:LEU:HA	1:B:582:LEU:HD23	1.80	0.41
1:A:80:VAL:CG2	1:A:170:PHE:HB2	2.50	0.41
1:A:248:ILE:HA	1:A:251:PHE:CD2	2.56	0.41
1:B:250:VAL:O	1:B:254:GLU:HG3	2.21	0.41
1:A:201:LYS:NZ	1:A:403:ASP:OD2	2.49	0.41
1:A:492:VAL:O	1:A:498:SER:OG	2.34	0.41
1:B:66:GLY:HA3	1:B:238:ASP:O	2.20	0.41
1:A:70:ASP:OD1	1:A:70:ASP:C	2.64	0.40
1:B:542:ALA:HB3	1:B:582:LEU:HD11	2.03	0.40
1:B:440:ALA:HA	1:B:446:ASN:O	2.20	0.40
1:B:507:GLY:HA2	1:B:590:SER:C	2.46	0.40
1:A:186:LYS:NZ	2:A:741:HOH:O	2.53	0.40
1:B:146:THR:HB	1:B:148:GLN:OE1	2.20	0.40
1:B:361:MET:HG2	1:B:372:LEU:CD1	2.51	0.40
1:A:295:LYS:HE2	1:A:295:LYS:HB3	1.67	0.40
1:A:508:TYR:CE2	1:A:589:LYS:HD3	2.56	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	500/595 (84%)	486 (97%)	14 (3%)	0	100	100
1	B	502/595 (84%)	488 (97%)	14 (3%)	0	100	100
All	All	1002/1190 (84%)	974 (97%)	28 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	390/503 (78%)	383 (98%)	7 (2%)	51	58
1	B	392/503 (78%)	387 (99%)	5 (1%)	61	68
All	All	782/1006 (78%)	770 (98%)	12 (2%)	57	64

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ASP
1	A	94	SER
1	A	136	ILE
1	A	179	HIS
1	A	295	LYS
1	A	387	SER
1	A	495	SER
1	B	70	ASP
1	B	145	LEU
1	B	179	HIS
1	B	327	GLU
1	B	590	SER

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	264	GLN
1	A	443	ASN
1	A	546	ASN
1	A	561	ASN
1	B	102	HIS
1	B	247	ASN
1	B	375	HIS
1	B	409	GLN
1	B	446	ASN
1	B	494	ASN

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Mol	Chain	Res	Type
1	B	499	HIS
1	B	563	GLN
1	B	580	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 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	504/595 (84%)	0.84	60 (11%) 9 8	18, 39, 68, 106	0
1	B	505/595 (84%)	0.77	71 (14%) 6 5	18, 33, 80, 117	1 (0%)
All	All	1009/1190 (84%)	0.81	131 (12%) 7 6	18, 36, 76, 117	1 (0%)

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	525	GLY	4.6
1	B	96	ASN	4.4
1	B	116	VAL	4.2
1	B	110	ALA	4.1
1	A	96	ASN	4.0
1	B	104	VAL	3.9
1	B	117	ILE	3.9
1	A	354	VAL	3.9
1	A	95	ALA	3.8
1	B	125	GLU	3.8
1	B	142	GLY	3.8
1	A	110	ALA	3.8
1	B	106	LYS	3.7
1	A	140	ARG	3.6
1	B	143	THR	3.6
1	B	525	GLY	3.5
1	B	124	ILE	3.5
1	A	359	ILE	3.5
1	A	358	GLU	3.4
1	A	356	TRP	3.4
1	A	338	SER	3.4
1	B	523	ASN	3.3
1	A	339	GLY	3.3
1	B	153	ILE	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	111	LYS	3.2
1	A	61	GLN	3.2
1	A	93	ALA	3.1
1	B	95	ALA	3.1
1	A	380	VAL	3.1
1	B	119	MET	3.1
1	A	519	VAL	3.0
1	A	147	TYR	3.0
1	B	128	LEU	2.9
1	A	523	ASN	2.9
1	B	105	ASP	2.9
1	A	348	ILE	2.9
1	B	161	ILE	2.9
1	A	145	LEU	2.8
1	B	108	GLU	2.8
1	B	344	MET	2.8
1	A	363	LEU	2.8
1	B	100	PRO	2.8
1	A	117	ILE	2.8
1	A	413	SER	2.8
1	A	114	SER	2.7
1	A	113	LEU	2.7
1	A	83	TYR	2.7
1	A	104	VAL	2.7
1	B	94	SER	2.7
1	B	93	ALA	2.7
1	A	524	GLY	2.7
1	B	134	PHE	2.6
1	B	330	PHE	2.6
1	B	158	LEU	2.6
1	B	87	ALA	2.6
1	B	138	PHE	2.6
1	B	365	PHE	2.6
1	A	153	ILE	2.6
1	B	123	GLU	2.6
1	B	113	LEU	2.6
1	B	524	GLY	2.6
1	B	348	ILE	2.5
1	A	345	GLY	2.5
1	A	355	GLY	2.5
1	B	98	LYS	2.5
1	B	351	TRP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	419	LYS	2.5
1	A	85	LEU	2.5
1	B	86	VAL	2.5
1	B	118	ASP	2.5
1	B	343	ILE	2.4
1	A	367	TYR	2.4
1	A	501	ALA	2.4
1	B	133	ALA	2.4
1	B	140	ARG	2.4
1	B	563	GLN	2.4
1	B	89	ILE	2.4
1	A	235	LYS	2.4
1	B	346	SER	2.4
1	A	336	TYR	2.3
1	B	147	TYR	2.3
1	A	139	GLY	2.3
1	A	142	GLY	2.3
1	B	331	ASP	2.3
1	B	107	LYS	2.3
1	B	165	PRO	2.3
1	B	522	PRO	2.3
1	A	520	ALA	2.3
1	A	372	LEU	2.3
1	B	209	GLY	2.3
1	A	161	ILE	2.3
1	B	345	GLY	2.3
1	A	498	SER	2.2
1	A	134	PHE	2.2
1	B	121	PRO	2.2
1	B	91	LYS	2.2
1	B	112	LYS	2.2
1	B	115	THR	2.2
1	A	98	LYS	2.2
1	B	335	LYS	2.2
1	A	100	PRO	2.1
1	B	109	THR	2.2
1	A	151	LEU	2.1
1	B	164	LEU	2.1
1	B	235	LYS	2.1
1	A	351	TRP	2.1
1	A	116	VAL	2.1
1	A	499	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	102	HIS	2.1
1	B	380	VAL	2.1
1	A	158	LEU	2.1
1	B	92	LYS	2.1
1	A	414	ASN	2.1
1	A	561	ASN	2.1
1	A	340	HIS	2.1
1	A	347	ARG	2.1
1	B	354	VAL	2.1
1	B	97	SER	2.1
1	B	114	SER	2.1
1	B	413	SER	2.1
1	A	344	MET	2.1
1	A	101	ARG	2.1
1	A	343	ILE	2.1
1	B	356	TRP	2.1
1	B	367	TYR	2.0
1	B	122	GLU	2.0
1	A	97	SER	2.0
1	A	368	SER	2.0
1	A	157	ASN	2.0
1	B	84	LYS	2.0
1	B	339	GLY	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.