



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:13 AM UTC

PDB ID : 3LWT / pdb_00003lwt
Title : Crystal structure of the Yeast Sac1: Implications for its phosphoinositide phosphatase function
Authors : Mao, Y.; Manford, A.; Xia, T.; Saxena, A.K.; Stefan, C.; Hu, F.; Emr, S.D.
Deposited on : 2010-02-24
Resolution : 1.96 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

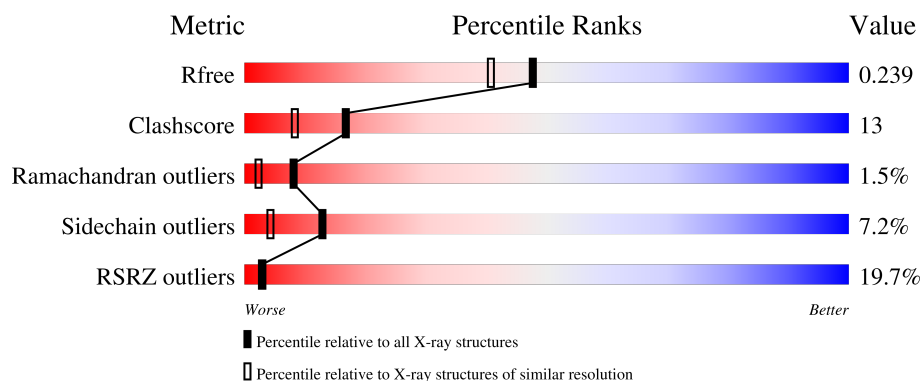
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.96 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	X	505	18% 65% 20% 5% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3851 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 Phosphoinositide phosphatase SAC1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	457	Total	C	N	O	S	0	0	0
			3683	2344	637	695	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	0	SER	-	expression tag	UNP P32368

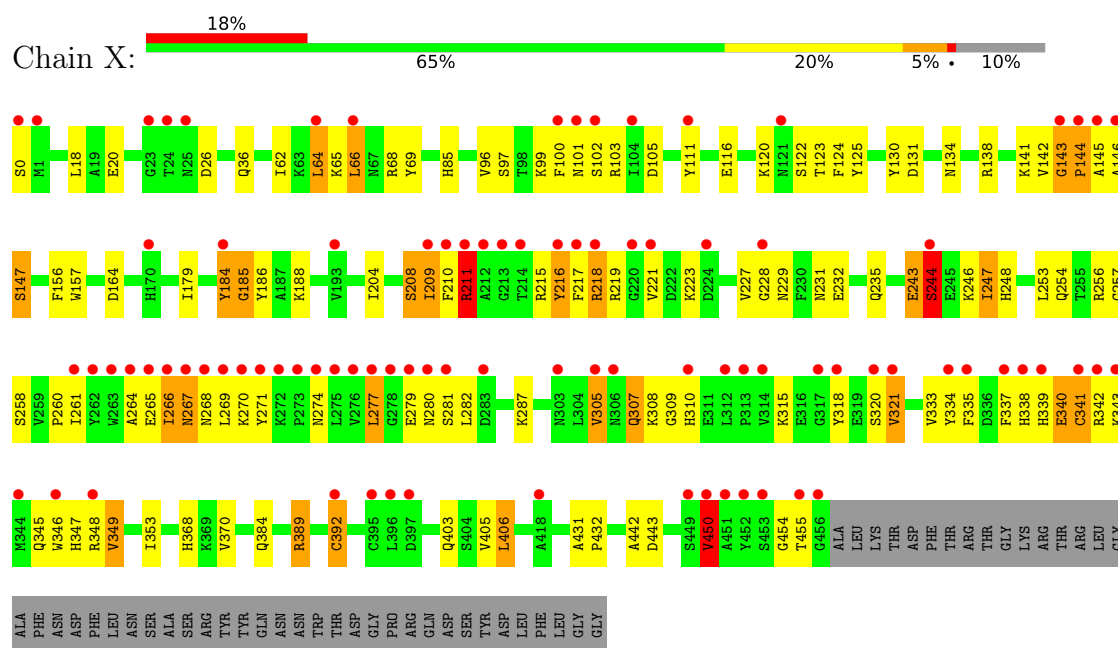
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	X	168	Total	O	0	0
			168	168		

3 Residue-property plots

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: Phosphoinositide phosphatase SAC1



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	86.16Å 94.73Å 155.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.18 – 1.96 33.18 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.2 (33.18-1.96) 99.1 (33.18-1.96)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.12 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.5.0093	Depositor
R, R_{free}	0.199 , 0.243 0.196 , 0.239	Depositor DCC
R_{free} test set	2327 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3851	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% 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	X	1.47	17/3768 (0.5%)	1.27	17/5113 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	X	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	156	PHE	N-CA	9.74	1.59	1.46
1	X	96	VAL	CA-CB	7.78	1.64	1.54
1	X	184	TYR	C-N	-7.22	1.23	1.33
1	X	184	TYR	CA-C	7.21	1.62	1.52
1	X	389	ARG	CG-CD	-7.15	1.31	1.52
1	X	450	VAL	CA-C	6.60	1.61	1.52
1	X	442	ALA	CA-CB	6.46	1.63	1.53
1	X	392	CYS	CB-SG	-6.28	1.60	1.81
1	X	450	VAL	CA-CB	5.66	1.62	1.54
1	X	204	ILE	N-CA	-5.53	1.39	1.46
1	X	384	GLN	C-O	5.39	1.30	1.24
1	X	164	ASP	N-CA	5.38	1.52	1.46
1	X	321	VAL	CA-CB	-5.34	1.48	1.54
1	X	131	ASP	CA-C	5.24	1.59	1.52
1	X	257	GLY	N-CA	5.21	1.51	1.45
1	X	209	ILE	CG1-CD1	5.02	1.71	1.51
1	X	179	ILE	CA-CB	5.01	1.60	1.53

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	184	TYR	N-CA-C	7.94	121.09	111.40
1	X	208	SER	CA-C-N	7.85	133.79	123.11
1	X	208	SER	C-N-CA	7.85	133.79	123.11
1	X	156	PHE	N-CA-CB	-6.77	100.84	110.46
1	X	147	SER	N-CA-CB	-6.62	99.25	110.50
1	X	185	GLY	N-CA-C	6.56	128.73	113.18
1	X	184	TYR	O-C-N	-6.56	114.24	122.17
1	X	143	GLY	CA-C-N	6.50	127.96	119.84
1	X	143	GLY	C-N-CA	6.50	127.96	119.84
1	X	66	LEU	N-CA-C	6.33	119.16	111.82
1	X	184	TYR	CB-CA-C	6.25	121.29	110.79
1	X	320	SER	N-CA-C	-6.12	104.52	112.23
1	X	349	VAL	CB-CA-C	-5.86	104.14	112.22
1	X	216	TYR	N-CA-C	5.57	116.71	108.86
1	X	333	VAL	CB-CA-C	-5.38	102.48	110.33
1	X	265	GLU	N-CA-C	5.05	116.63	108.34
1	X	341	CYS	N-CA-C	5.03	119.06	113.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	389	ARG	Sidechain

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	X	3683	0	3582	93	0
2	X	168	0	0	9	0
All	All	3851	0	3582	93	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 (93) 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:X:185:GLY:N	2:X:660:HOH:O	1.62	1.22
1:X:123:THR:O	1:X:184:TYR:O	1.67	1.10
1:X:146:ALA:HB1	1:X:147:SER:HA	1.39	1.02
1:X:209:ILE:HG23	2:X:642:HOH:O	1.65	0.97
1:X:338:HIS:HB2	1:X:348:ARG:NH1	1.80	0.95
1:X:64:LEU:HD11	1:X:111:TYR:CE2	2.04	0.92
1:X:62:ILE:HD11	1:X:111:TYR:OH	1.70	0.91
1:X:208:SER:O	2:X:661:HOH:O	1.90	0.88
1:X:36:GLN:HE21	1:X:68:ARG:H	1.23	0.86
1:X:267:ASN:O	1:X:267:ASN:ND2	2.16	0.78
1:X:210:PHE:HB2	1:X:211:ARG:HB2	1.63	0.78
1:X:266:ILE:HD12	1:X:266:ILE:H	1.49	0.77
1:X:134:ASN:HD22	1:X:138:ARG:HE	1.32	0.77
1:X:184:TYR:CA	2:X:660:HOH:O	2.33	0.77
1:X:184:TYR:C	2:X:660:HOH:O	2.10	0.77
1:X:431:ALA:HB3	1:X:432:PRO:HD3	1.70	0.74
1:X:146:ALA:HB1	1:X:147:SER:CA	2.18	0.73
1:X:353:ILE:HG12	1:X:406:LEU:HD13	1.72	0.72
1:X:0:SER:HB2	2:X:599:HOH:O	1.90	0.71
1:X:64:LEU:CD1	1:X:111:TYR:CE2	2.76	0.69
1:X:116:GLU:OE2	1:X:120:LYS:HE3	1.93	0.69
1:X:143:GLY:HA2	2:X:651:HOH:O	1.93	0.68
1:X:338:HIS:HB2	1:X:348:ARG:HH12	1.56	0.67
1:X:20:GLU:HG3	2:X:624:HOH:O	1.96	0.65
1:X:146:ALA:CB	1:X:147:SER:HA	2.19	0.64
1:X:64:LEU:HG	1:X:111:TYR:CD2	2.32	0.64
1:X:337:PHE:CE1	1:X:349:VAL:HG23	2.33	0.63
1:X:270:LYS:HG3	1:X:271:TYR:HD1	1.62	0.63
1:X:62:ILE:HD11	1:X:111:TYR:HH	1.64	0.63
1:X:247:ILE:HD12	1:X:247:ILE:N	2.14	0.62
1:X:246:LYS:HZ2	1:X:370:VAL:HG11	1.64	0.61
1:X:210:PHE:HB3	1:X:228:GLY:HA2	1.81	0.61
1:X:337:PHE:CE1	1:X:345:GLN:HG2	2.34	0.61
1:X:338:HIS:CB	1:X:348:ARG:NH1	2.60	0.60
1:X:243:GLU:O	1:X:244:SER:HB3	2.02	0.59
1:X:341:CYS:HA	1:X:345:GLN:HB2	1.85	0.59
1:X:157:TRP:HA	1:X:209:ILE:HG13	1.84	0.59
1:X:99:LYS:CB	1:X:102:SER:HB2	2.32	0.59
1:X:282:LEU:HD23	1:X:321:VAL:HG13	1.85	0.58
1:X:143:GLY:O	1:X:145:ALA:N	2.33	0.57
1:X:218:ARG:HG3	1:X:219:ARG:N	2.20	0.57
1:X:248:HIS:HD2	1:X:368:HIS:NE2	2.01	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:307:GLN:HG2	1:X:309:GLY:H	1.70	0.56
1:X:142:VAL:O	1:X:144:PRO:HD2	2.07	0.55
1:X:99:LYS:HB3	1:X:102:SER:HB2	1.89	0.54
1:X:36:GLN:NE2	1:X:68:ARG:H	2.00	0.53
1:X:270:LYS:HG3	1:X:271:TYR:CD1	2.44	0.53
1:X:269:LEU:HD23	1:X:270:LYS:HE3	1.91	0.53
1:X:338:HIS:CB	1:X:348:ARG:HH12	2.21	0.53
1:X:210:PHE:CB	1:X:211:ARG:HB2	2.37	0.52
1:X:337:PHE:CE1	1:X:349:VAL:CG2	2.92	0.52
1:X:346:TRP:O	1:X:347:HIS:C	2.51	0.52
1:X:211:ARG:O	1:X:216:TYR:HE1	1.91	0.52
1:X:122:SER:HB2	1:X:124:PHE:CZ	2.45	0.52
1:X:65:LYS:HE2	1:X:105:ASP:OD2	2.11	0.51
1:X:339:HIS:HB2	1:X:340:GLU:HA	1.92	0.51
1:X:254:GLN:HE22	1:X:403:GLN:HA	1.75	0.51
1:X:229:ASN:ND2	1:X:392:CYS:SG	2.85	0.50
1:X:146:ALA:CB	1:X:147:SER:CA	2.86	0.50
1:X:343:LYS:N	2:X:659:HOH:O	2.44	0.50
1:X:143:GLY:C	1:X:145:ALA:H	2.19	0.49
1:X:335:PHE:CZ	1:X:348:ARG:O	2.66	0.48
1:X:221:VAL:HB	1:X:227:VAL:HG12	1.96	0.48
1:X:100:PHE:CG	1:X:101:ASN:N	2.81	0.48
1:X:218:ARG:NH1	1:X:258:SER:OG	2.47	0.48
1:X:341:CYS:HB2	1:X:345:GLN:CD	2.40	0.47
1:X:260:PRO:HD2	1:X:318:TYR:CE1	2.50	0.47
1:X:337:PHE:HE1	1:X:349:VAL:HG23	1.77	0.46
1:X:307:GLN:HG2	1:X:308:LYS:N	2.30	0.46
1:X:243:GLU:O	1:X:244:SER:CB	2.65	0.45
1:X:307:GLN:HG3	1:X:310:HIS:O	2.16	0.45
1:X:279:GLU:O	1:X:279:GLU:HG2	2.16	0.45
1:X:111:TYR:CD2	1:X:111:TYR:C	2.95	0.44
1:X:232:GLU:HB3	1:X:253:LEU:HD11	1.99	0.44
1:X:85:HIS:CG	1:X:130:TYR:HB2	2.53	0.44
1:X:266:ILE:H	1:X:266:ILE:CD1	2.23	0.44
1:X:122:SER:HB2	1:X:124:PHE:CE2	2.53	0.43
1:X:62:ILE:CD1	1:X:111:TYR:OH	2.56	0.43
1:X:231:ASN:HD22	1:X:256:ARG:HH21	1.65	0.43
1:X:246:LYS:NZ	1:X:370:VAL:HG11	2.31	0.43
1:X:341:CYS:HA	1:X:345:GLN:CB	2.48	0.43
1:X:215:ARG:HH22	1:X:223:LYS:HD2	1.85	0.42
1:X:188:LYS:HE3	1:X:188:LYS:HB2	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:335:PHE:HZ	1:X:348:ARG:O	2.02	0.42
1:X:277:LEU:HB3	1:X:281:SER:HB2	2.02	0.41
1:X:211:ARG:O	1:X:216:TYR:CE1	2.72	0.41
1:X:243:GLU:H	1:X:243:GLU:HG3	1.62	0.41
1:X:337:PHE:CZ	1:X:345:GLN:HG2	2.55	0.41
1:X:125:TYR:CE2	1:X:186:TYR:HA	2.56	0.41
1:X:349:VAL:HG11	1:X:405:VAL:HB	2.02	0.40
1:X:454:GLY:HA3	1:X:455:THR:HA	1.94	0.40
1:X:64:LEU:CG	1:X:111:TYR:CD2	3.02	0.40
1:X:69:TYR:CE1	1:X:97:SER:HB3	2.57	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	X	455/505 (90%)	425 (93%)	23 (5%)	7 (2%)	8 2

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	144	PRO
1	X	244	SER
1	X	450	VAL
1	X	217	PHE
1	X	264	ALA
1	X	305	VAL
1	X	211	ARG

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	X	400/441 (91%)	371 (93%)	29 (7%)	13 4

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	18	LEU
1	X	26	ASP
1	X	64	LEU
1	X	66	LEU
1	X	103	ARG
1	X	141	LYS
1	X	211	ARG
1	X	218	ARG
1	X	235	GLN
1	X	243	GLU
1	X	244	SER
1	X	247	ILE
1	X	261	ILE
1	X	266	ILE
1	X	267	ASN
1	X	268	ASN
1	X	274	ASN
1	X	277	LEU
1	X	280	ASN
1	X	287	LYS
1	X	305	VAL
1	X	307	GLN
1	X	315	LYS
1	X	334	TYR
1	X	340	GLU
1	X	342	ARG
1	X	406	LEU
1	X	443	ASP
1	X	450	VAL

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

Mol	Chain	Res	Type
1	X	9	GLN
1	X	36	GLN
1	X	83	ASN
1	X	134	ASN
1	X	229	ASN
1	X	231	ASN
1	X	248	HIS
1	X	254	GLN
1	X	280	ASN
1	X	391	ASN
1	X	403	GLN
1	X	408	GLN
1	X	444	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	X	457/505 (90%)	1.02	90 (19%) 3 3	23, 45, 105, 145	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	266	ILE	9.9
1	X	312	LEU	7.9
1	X	273	PRO	7.5
1	X	341	CYS	6.9
1	X	275	LEU	6.6
1	X	337	PHE	6.6
1	X	210	PHE	6.4
1	X	455	THR	6.3
1	X	452	TYR	6.0
1	X	277	LEU	5.4
1	X	24	THR	5.4
1	X	450	VAL	5.3
1	X	111	TYR	5.3
1	X	456	GLY	5.3
1	X	144	PRO	5.2
1	X	342	ARG	5.1
1	X	276	VAL	5.0
1	X	271	TYR	4.9
1	X	264	ALA	4.9
1	X	451	ALA	4.9
1	X	278	GLY	4.8
1	X	263	TRP	4.7
1	X	212	ALA	4.6
1	X	100	PHE	4.5
1	X	217	PHE	4.5
1	X	216	TYR	4.4
1	X	267	ASN	4.4

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Mol	Chain	Res	Type	RSRZ
1	X	261	ILE	4.4
1	X	228	GLY	4.2
1	X	25	ASN	4.2
1	X	214	THR	4.1
1	X	395	CYS	4.1
1	X	146	ALA	4.0
1	X	209	ILE	4.0
1	X	221	VAL	3.9
1	X	334	TYR	3.8
1	X	262	TYR	3.7
1	X	101	ASN	3.7
1	X	453	SER	3.7
1	X	211	ARG	3.6
1	X	310	HIS	3.6
1	X	269	LEU	3.5
1	X	244	SER	3.5
1	X	339	HIS	3.4
1	X	272	LYS	3.4
1	X	281	SER	3.4
1	X	305	VAL	3.3
1	X	145	ALA	3.3
1	X	338	HIS	3.3
1	X	265	GLU	3.2
1	X	313	PRO	3.2
1	X	121	ASN	3.1
1	X	344	MET	3.0
1	X	0	SER	3.0
1	X	268	ASN	3.0
1	X	318	TYR	3.0
1	X	274	ASN	3.0
1	X	102	SER	2.9
1	X	306	ASN	2.8
1	X	314	VAL	2.8
1	X	66	LEU	2.8
1	X	418	ALA	2.8
1	X	346	TRP	2.8
1	X	220	GLY	2.7
1	X	343	LYS	2.7
1	X	64	LEU	2.6
1	X	280	ASN	2.6
1	X	335	PHE	2.5
1	X	283	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	X	270	LYS	2.5
1	X	449	SER	2.5
1	X	104	ILE	2.4
1	X	184	TYR	2.4
1	X	143	GLY	2.4
1	X	317	GLY	2.4
1	X	396	LEU	2.3
1	X	303	ASN	2.3
1	X	320	SER	2.3
1	X	193	VAL	2.3
1	X	218	ARG	2.2
1	X	392	CYS	2.2
1	X	397	ASP	2.2
1	X	213	GLY	2.2
1	X	170	HIS	2.1
1	X	23	GLY	2.1
1	X	321	VAL	2.1
1	X	224	ASP	2.1
1	X	348	ARG	2.0
1	X	279	GLU	2.0
1	X	1	MET	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.