



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 10:01 AM UTC

PDB ID : 7QPK / pdb_00007qpk
Title : Structure of the A-region of Awp14 from Candida glabrata
Authors : Essen, L.-O.; Reithofer, V.; de Groot, P.
Deposited on : 2022-01-05
Resolution : 2.75 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

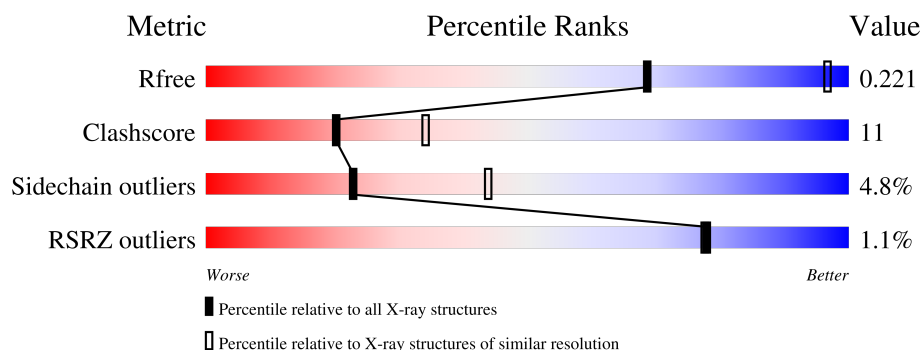
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	413	65% 24% • 9%
1	B	413	72% 17% • 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6126 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 A-region of Awp14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2967	1908	475	578	6			
1	B	378	Total	C	N	O	S	0	0	0
			2981	1916	478	581	6			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	initiating methionine	UNP A0A0W0CDE7
A	-11	GLY	-	expression tag	UNP A0A0W0CDE7
A	-10	SER	-	expression tag	UNP A0A0W0CDE7
A	-9	SER	-	expression tag	UNP A0A0W0CDE7
A	-8	HIS	-	expression tag	UNP A0A0W0CDE7
A	-7	HIS	-	expression tag	UNP A0A0W0CDE7
A	-6	HIS	-	expression tag	UNP A0A0W0CDE7
A	-5	HIS	-	expression tag	UNP A0A0W0CDE7
A	-4	HIS	-	expression tag	UNP A0A0W0CDE7
A	-3	HIS	-	expression tag	UNP A0A0W0CDE7
A	-2	SER	-	expression tag	UNP A0A0W0CDE7
A	-1	SER	-	expression tag	UNP A0A0W0CDE7
A	0	GLY	-	expression tag	UNP A0A0W0CDE7
A	1	LEU	-	expression tag	UNP A0A0W0CDE7
A	2	VAL	-	expression tag	UNP A0A0W0CDE7
A	3	PRO	-	expression tag	UNP A0A0W0CDE7
A	4	ARG	-	expression tag	UNP A0A0W0CDE7
A	5	GLY	-	expression tag	UNP A0A0W0CDE7
A	6	SER	-	expression tag	UNP A0A0W0CDE7
A	7	HIS	-	expression tag	UNP A0A0W0CDE7
A	8	MET	-	expression tag	UNP A0A0W0CDE7
A	9	ALA	-	expression tag	UNP A0A0W0CDE7
A	10	SER	-	expression tag	UNP A0A0W0CDE7
A	11	MET	-	expression tag	UNP A0A0W0CDE7
A	12	THR	-	expression tag	UNP A0A0W0CDE7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	13	GLY	-	expression tag	UNP A0A0W0CDE7
A	14	GLY	-	expression tag	UNP A0A0W0CDE7
A	15	GLN	-	expression tag	UNP A0A0W0CDE7
A	16	GLN	-	expression tag	UNP A0A0W0CDE7
A	17	MET	-	expression tag	UNP A0A0W0CDE7
A	18	GLY	-	expression tag	UNP A0A0W0CDE7
A	19	ARG	-	expression tag	UNP A0A0W0CDE7
A	20	GLY	-	expression tag	UNP A0A0W0CDE7
A	21	SER	-	expression tag	UNP A0A0W0CDE7
B	-12	MET	-	initiating methionine	UNP A0A0W0CDE7
B	-11	GLY	-	expression tag	UNP A0A0W0CDE7
B	-10	SER	-	expression tag	UNP A0A0W0CDE7
B	-9	SER	-	expression tag	UNP A0A0W0CDE7
B	-8	HIS	-	expression tag	UNP A0A0W0CDE7
B	-7	HIS	-	expression tag	UNP A0A0W0CDE7
B	-6	HIS	-	expression tag	UNP A0A0W0CDE7
B	-5	HIS	-	expression tag	UNP A0A0W0CDE7
B	-4	HIS	-	expression tag	UNP A0A0W0CDE7
B	-3	HIS	-	expression tag	UNP A0A0W0CDE7
B	-2	SER	-	expression tag	UNP A0A0W0CDE7
B	-1	SER	-	expression tag	UNP A0A0W0CDE7
B	0	GLY	-	expression tag	UNP A0A0W0CDE7
B	1	LEU	-	expression tag	UNP A0A0W0CDE7
B	2	VAL	-	expression tag	UNP A0A0W0CDE7
B	3	PRO	-	expression tag	UNP A0A0W0CDE7
B	4	ARG	-	expression tag	UNP A0A0W0CDE7
B	5	GLY	-	expression tag	UNP A0A0W0CDE7
B	6	SER	-	expression tag	UNP A0A0W0CDE7
B	7	HIS	-	expression tag	UNP A0A0W0CDE7
B	8	MET	-	expression tag	UNP A0A0W0CDE7
B	9	ALA	-	expression tag	UNP A0A0W0CDE7
B	10	SER	-	expression tag	UNP A0A0W0CDE7
B	11	MET	-	expression tag	UNP A0A0W0CDE7
B	12	THR	-	expression tag	UNP A0A0W0CDE7
B	13	GLY	-	expression tag	UNP A0A0W0CDE7
B	14	GLY	-	expression tag	UNP A0A0W0CDE7
B	15	GLN	-	expression tag	UNP A0A0W0CDE7
B	16	GLN	-	expression tag	UNP A0A0W0CDE7
B	17	MET	-	expression tag	UNP A0A0W0CDE7
B	18	GLY	-	expression tag	UNP A0A0W0CDE7
B	19	ARG	-	expression tag	UNP A0A0W0CDE7
B	20	GLY	-	expression tag	UNP A0A0W0CDE7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	21	SER	-	expression tag	UNP A0A0W0CDE7

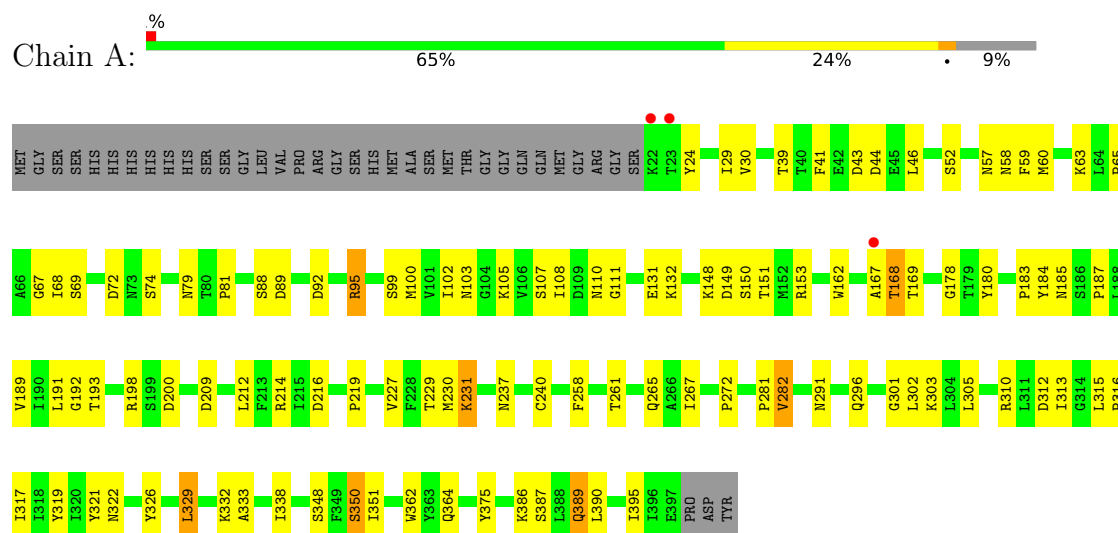
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	84	Total O 84 84	0	0
2	B	94	Total O 94 94	0	0

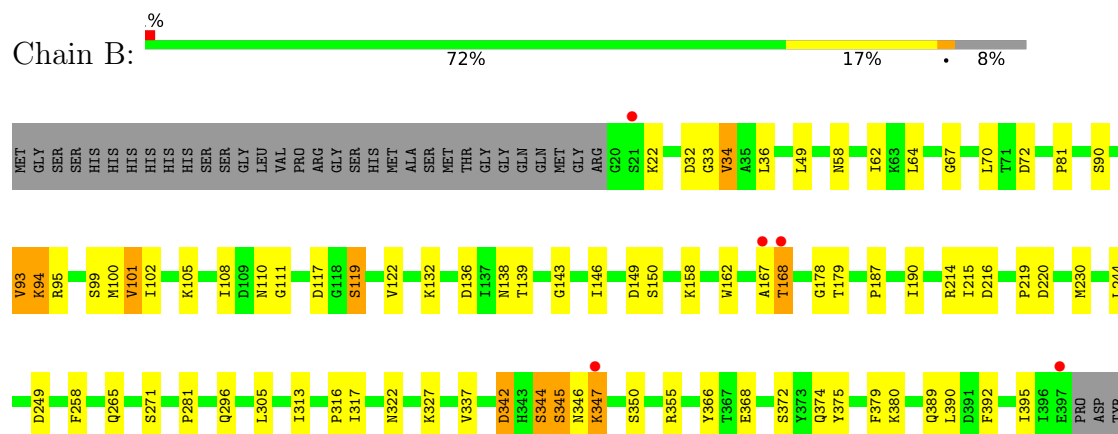
3 Residue-property plots

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: A-region of Awp14



• Molecule 1: A-region of Awp14



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	78.17Å 172.81Å 140.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.98 – 2.75 28.98 – 2.75	Depositor EDS
% Data completeness (in resolution range)	95.4 (28.98-2.75) 95.3 (28.98-2.75)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.76Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.178 , 0.243 (Not available) , 0.221	Depositor DCC
R_{free} test set	780 reflections (3.09%)	wwPDB-VP
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.9	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	0.95	EDS
Total number of atoms	6126	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% 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	0.47	4/3036 (0.1%)	0.70	4/4126 (0.1%)
1	B	0.49	0/3050	0.83	17/4143 (0.4%)
All	All	0.48	4/6086 (0.1%)	0.77	21/8269 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	231	LYS	CE-NZ	7.94	1.73	1.49
1	A	315	LEU	C-O	6.93	1.27	1.23
1	A	231	LYS	CD-CE	5.77	1.69	1.52
1	A	231	LYS	CG-CD	5.28	1.68	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	347	LYS	CA-CB-CG	8.73	131.55	114.10
1	B	346	ASN	CA-C-N	8.72	133.15	120.95
1	B	346	ASN	C-N-CA	8.72	133.15	120.95
1	B	347	LYS	CB-CG-CD	-8.46	91.85	111.30
1	A	43	ASP	CA-C-N	8.38	132.34	120.28
1	A	43	ASP	C-N-CA	8.38	132.34	120.28
1	B	345	SER	CA-C-N	8.31	134.46	122.40
1	B	345	SER	C-N-CA	8.31	134.46	122.40
1	B	347	LYS	CD-CE-NZ	8.28	138.38	111.90
1	B	94	LYS	CB-CG-CD	-7.65	93.70	111.30
1	A	44	ASP	N-CA-CB	-7.61	98.50	110.22
1	B	347	LYS	CA-C-N	-6.46	110.58	121.94
1	B	347	LYS	C-N-CA	-6.46	110.58	121.94
1	B	94	LYS	CA-CB-CG	6.36	126.82	114.10
1	B	347	LYS	N-CA-C	-5.85	101.19	109.96
1	B	347	LYS	CG-CD-CE	-5.66	98.29	111.30
1	B	168	THR	CA-C-N	5.50	132.05	121.54
1	B	168	THR	C-N-CA	5.50	132.05	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ASP	N-CA-C	5.31	117.81	111.71
1	B	342	ASP	CB-CA-C	5.03	118.43	110.19
1	B	94	LYS	CD-CE-NZ	-5.02	95.83	111.90

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	2967	0	2906	64	0
1	B	2981	0	2925	69	0
2	A	84	0	0	3	0
2	B	94	0	0	2	0
All	All	6126	0	5831	133	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 11.

All (133) 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:231:LYS:CE	1:A:231:LYS:NZ	1.73	1.51
1:B:345:SER:HB2	1:B:347:LYS:HD2	1.19	1.13
1:B:345:SER:HB2	1:B:347:LYS:CD	1.84	1.06
1:B:345:SER:H	1:B:347:LYS:NZ	1.58	1.00
1:A:58:ASN:HA	1:A:81:PRO:HB3	1.57	0.86
1:B:345:SER:H	1:B:347:LYS:HZ2	1.18	0.86
1:B:327:LYS:NZ	2:B:501:HOH:O	2.09	0.86
1:A:303:LYS:NZ	2:A:501:HOH:O	2.15	0.80
1:B:265:GLN:NE2	1:B:375:TYR:H	1.81	0.79
1:A:60:MET:HE2	1:A:81:PRO:HG2	1.69	0.74
1:B:342:ASP:HB3	1:B:347:LYS:HE3	1.70	0.74
1:A:162:TRP:HE1	1:A:167:ALA:HB3	1.53	0.73
1:B:345:SER:CB	1:B:347:LYS:HD2	2.10	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:ASN:HD21	1:A:79:ASN:HD22	1.39	0.70
1:A:168:THR:HB	1:A:169:THR:HG23	1.72	0.70
1:A:111:GLY:H	1:A:149:ASP:HB3	1.57	0.69
1:B:368:GLU:CD	1:B:368:GLU:H	1.99	0.69
1:B:22:LYS:NZ	2:B:502:HOH:O	2.25	0.69
1:B:342:ASP:HB3	1:B:347:LYS:CE	2.21	0.69
1:A:265:GLN:HE22	1:A:375:TYR:HB2	1.60	0.67
1:B:265:GLN:HE22	1:B:375:TYR:H	1.40	0.67
1:B:58:ASN:HA	1:B:81:PRO:HB3	1.77	0.66
1:B:162:TRP:HE1	1:B:167:ALA:HB3	1.61	0.66
1:B:345:SER:N	1:B:347:LYS:HZ2	1.92	0.66
1:A:39:THR:HB	1:A:60:MET:HB2	1.77	0.65
1:B:317:ILE:HB	1:B:390:LEU:HD22	1.77	0.65
1:A:95:ARG:O	1:A:132:LYS:HE3	1.96	0.65
1:A:212:LEU:HD11	1:A:282:VAL:HG21	1.78	0.65
1:B:146:ILE:HD12	1:B:190:ILE:HD12	1.80	0.63
1:B:342:ASP:CG	1:B:347:LYS:HE3	2.25	0.62
1:B:342:ASP:CB	1:B:347:LYS:HE3	2.30	0.61
1:A:231:LYS:NZ	1:A:231:LYS:CD	2.63	0.61
1:B:214:ARG:HD2	1:B:216:ASP:OD1	2.00	0.60
1:B:111:GLY:N	1:B:149:ASP:HB3	2.17	0.60
1:A:65:PRO:HA	1:A:103:ASN:HB2	1.83	0.59
1:A:214:ARG:NH1	1:A:216:ASP:OD2	2.35	0.59
1:A:387:SER:OG	1:A:389:GLN:OE1	2.19	0.59
1:A:272:PRO:HG3	1:A:301:GLY:HA3	1.85	0.59
1:B:138:ASN:OD1	1:B:179:THR:HG23	2.03	0.59
1:A:183:PRO:HD2	1:A:184:TYR:CD1	2.39	0.58
1:A:265:GLN:HE22	1:A:375:TYR:H	1.51	0.57
1:A:265:GLN:NE2	1:A:375:TYR:HB2	2.19	0.57
1:B:100:MET:HE3	1:B:102:ILE:HG13	1.87	0.57
1:B:101:VAL:HG13	1:B:136:ASP:HB3	1.87	0.56
1:B:95:ARG:NH2	1:B:122:VAL:O	2.37	0.56
1:A:178:GLY:O	1:A:219:PRO:HA	2.07	0.55
1:A:281:PRO:HG3	1:A:313:ILE:HD13	1.87	0.55
1:B:342:ASP:HB3	1:B:347:LYS:HB2	1.87	0.55
1:B:108:ILE:HG22	1:B:150:SER:HB3	1.89	0.55
1:B:345:SER:N	1:B:347:LYS:NZ	2.41	0.54
1:A:237:ASN:HD22	1:A:237:ASN:C	2.16	0.54
1:A:100:MET:HE3	1:A:102:ILE:HG13	1.89	0.53
1:B:111:GLY:H	1:B:149:ASP:HB3	1.73	0.53
1:B:249:ASP:OD2	1:B:271:SER:OG	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:TYR:CE1	1:A:329:LEU:HD22	2.44	0.53
1:A:189:VAL:HG12	1:A:191:LEU:HD13	1.90	0.53
1:A:296:GLN:NE2	1:A:322:ASN:H	2.06	0.53
1:B:90:SER:HA	1:B:93:VAL:HG23	1.91	0.53
1:A:72:ASP:H	1:A:110:ASN:HB2	1.73	0.52
1:A:362:TRP:HZ3	1:A:364:GLN:HE21	1.56	0.52
1:A:108:ILE:HG22	1:A:150:SER:HB3	1.91	0.51
1:B:342:ASP:OD1	1:B:347:LYS:NZ	2.37	0.51
1:A:265:GLN:NE2	1:A:375:TYR:H	2.08	0.51
1:B:95:ARG:O	1:B:132:LYS:NZ	2.38	0.50
1:A:258:PHE:HE1	1:A:305:LEU:HD12	1.77	0.50
1:B:345:SER:HB2	1:B:347:LYS:HD3	1.87	0.50
1:A:321:TYR:CE2	1:A:329:LEU:HD13	2.47	0.50
1:A:24:TYR:HD2	1:A:29:ILE:HD13	1.77	0.49
1:A:57:ASN:ND2	1:A:79:ASN:HD22	2.08	0.49
1:B:111:GLY:CA	1:B:149:ASP:HB3	2.43	0.49
1:B:347:LYS:HE2	1:B:347:LYS:H	1.78	0.48
1:A:92:ASP:O	1:A:132:LYS:HE2	2.13	0.48
1:A:148:LYS:HA	1:A:192:GLY:HA2	1.94	0.48
1:A:265:GLN:HG2	1:A:267:ILE:HD11	1.95	0.48
1:A:111:GLY:N	1:A:149:ASP:HB3	2.26	0.48
1:B:342:ASP:CB	1:B:347:LYS:CE	2.90	0.48
1:B:178:GLY:O	1:B:219:PRO:HA	2.13	0.48
1:B:296:GLN:NE2	1:B:322:ASN:H	2.11	0.48
1:A:168:THR:HG22	1:A:169:THR:H	1.79	0.47
1:A:151:THR:HG23	1:A:193:THR:HB	1.96	0.47
1:B:187:PRO:HB2	1:B:230:MET:HE3	1.95	0.47
1:B:72:ASP:H	1:B:110:ASN:HB2	1.79	0.47
1:A:24:TYR:CD2	1:A:29:ILE:HD13	2.49	0.47
1:A:180:TYR:HD1	1:A:185:ASN:HA	1.80	0.47
1:B:342:ASP:CG	1:B:347:LYS:HZ1	2.21	0.47
1:A:41:PHE:HB2	1:A:59:PHE:CD1	2.50	0.47
1:A:302:LEU:HD23	1:A:351:ILE:HG12	1.97	0.46
1:B:316:PRO:HG3	1:B:392:PHE:CE2	2.50	0.46
1:B:342:ASP:C	1:B:347:LYS:NZ	2.74	0.46
1:B:62:ILE:HB	1:B:100:MET:HG3	1.97	0.45
1:B:117:ASP:OD1	1:B:119:SER:OG	2.33	0.45
1:B:220:ASP:OD1	1:B:220:ASP:N	2.46	0.45
1:B:122:VAL:HG22	1:B:158:LYS:HG2	1.99	0.45
1:A:338:ILE:O	1:A:350:SER:HA	2.17	0.44
1:B:67:GLY:HA3	1:B:105:LYS:O	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:LYS:HD3	1:B:347:LYS:N	2.32	0.44
1:B:32:ASP:OD1	1:B:34:VAL:HG22	2.17	0.44
1:A:316:PRO:HD2	1:A:390:LEU:O	2.18	0.44
1:B:345:SER:H	1:B:347:LYS:HZ3	1.55	0.44
1:B:344:SER:N	1:B:347:LYS:NZ	2.65	0.44
1:A:395:ILE:HD12	1:A:395:ILE:O	2.18	0.43
1:B:33:GLY:HA3	1:B:58:ASN:O	2.18	0.43
1:B:101:VAL:HA	1:B:136:ASP:O	2.18	0.43
1:B:179:THR:HG22	1:B:379:PHE:HE1	1.83	0.43
1:B:342:ASP:HB3	1:B:347:LYS:HE2	1.96	0.43
1:B:258:PHE:HE1	1:B:305:LEU:HD12	1.84	0.43
1:B:366:TYR:CE1	1:B:372:SER:HB2	2.54	0.43
1:A:240:CYS:O	1:A:261:THR:HA	2.19	0.43
1:B:139:THR:OG1	1:B:143:GLY:HA2	2.19	0.43
1:A:229:THR:HG23	2:A:549:HOH:O	2.18	0.42
1:A:189:VAL:HG21	1:A:230:MET:SD	2.59	0.42
1:A:310:ARG:HB3	1:A:312:ASP:OD1	2.19	0.42
1:A:63:LYS:C	1:A:65:PRO:HD3	2.45	0.42
1:A:282:VAL:HB	1:A:291:ASN:HD21	1.85	0.42
1:B:215:ILE:HB	1:B:244:LEU:HD23	2.02	0.42
1:B:265:GLN:NE2	1:B:375:TYR:HB2	2.34	0.42
1:B:265:GLN:NE2	1:B:374:GLN:HB2	2.34	0.42
1:A:41:PHE:HB2	1:A:59:PHE:CE1	2.55	0.41
1:A:187:PRO:HB3	1:A:230:MET:O	2.20	0.41
1:A:326:TYR:HB3	1:A:386:LYS:HE3	2.02	0.41
1:A:198:ARG:HB3	1:A:200:ASP:OD1	2.20	0.41
1:A:332:LYS:HE3	1:A:333:ALA:O	2.20	0.41
1:B:316:PRO:HD2	1:B:390:LEU:O	2.20	0.41
1:A:316:PRO:O	1:A:317:ILE:HG13	2.21	0.41
1:B:281:PRO:HG3	1:B:313:ILE:CD1	2.50	0.41
1:A:153:ARG:HG2	2:A:512:HOH:O	2.19	0.41
1:A:67:GLY:HA3	1:A:105:LYS:O	2.21	0.41
1:B:49:LEU:HD22	1:B:70:LEU:HD22	2.03	0.40
1:A:209:ASP:OD1	1:A:209:ASP:C	2.64	0.40
1:B:162:TRP:NE1	1:B:167:ALA:HB3	2.31	0.40
1:A:46:LEU:HD11	1:A:69:SER:HB2	2.03	0.40
1:B:64:LEU:HA	1:B:64:LEU:HD23	1.71	0.40
1:B:347:LYS:HB2	1:B:347:LYS:HE2	1.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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	333/363 (92%)	316 (95%)	17 (5%)	21	40
1	B	335/363 (92%)	320 (96%)	15 (4%)	24	46
All	All	668/726 (92%)	636 (95%)	32 (5%)	23	44

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	52	SER
1	A	68	ILE
1	A	74	SER
1	A	88	SER
1	A	89	ASP
1	A	95	ARG
1	A	99	SER
1	A	107	SER
1	A	131	GLU
1	A	168	THR
1	A	227	VAL
1	A	282	VAL
1	A	329	LEU
1	A	348	SER
1	A	350	SER
1	A	389	GLN
1	B	34	VAL
1	B	36	LEU
1	B	93	VAL
1	B	94	LYS

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Mol	Chain	Res	Type
1	B	99	SER
1	B	101	VAL
1	B	119	SER
1	B	168	THR
1	B	337	VAL
1	B	344	SER
1	B	350	SER
1	B	355	ARG
1	B	380	LYS
1	B	389	GLN
1	B	395	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	77	ASN
1	A	97	GLN
1	A	103	ASN
1	A	138	ASN
1	A	237	ASN
1	A	265	GLN
1	A	291	ASN
1	A	296	GLN
1	A	336	ASN
1	A	364	GLN
1	B	58	ASN
1	B	237	ASN
1	B	265	GLN
1	B	283	GLN
1	B	291	ASN
1	B	296	GLN
1	B	325	ASN
1	B	374	GLN

5.3.3 RNA ⓘ

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	376/413 (91%)	-0.37	3 (0%)	82 82	35, 50, 75, 120	0
1	B	378/413 (91%)	-0.48	5 (1%)	75 75	35, 44, 71, 101	0
All	All	754/826 (91%)	-0.43	8 (1%)	78 77	35, 46, 74, 120	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	347	LYS	2.7
1	A	22	LYS	2.7
1	B	167	ALA	2.4
1	B	168	THR	2.3
1	B	397	GLU	2.2
1	A	23	THR	2.2
1	B	21	SER	2.2
1	A	167	ALA	2.1

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.