



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2026 – 09:50 PM UTC

PDB ID : 2HSG / pdb_00002hsg
Title : Structure of transcription regulator CcpA in its DNA-free state
Authors : Loll, B.; Alings, C.; Saenger, W.; Biesiadka, J.
Deposited on : 2006-07-21
Resolution : 2.50 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

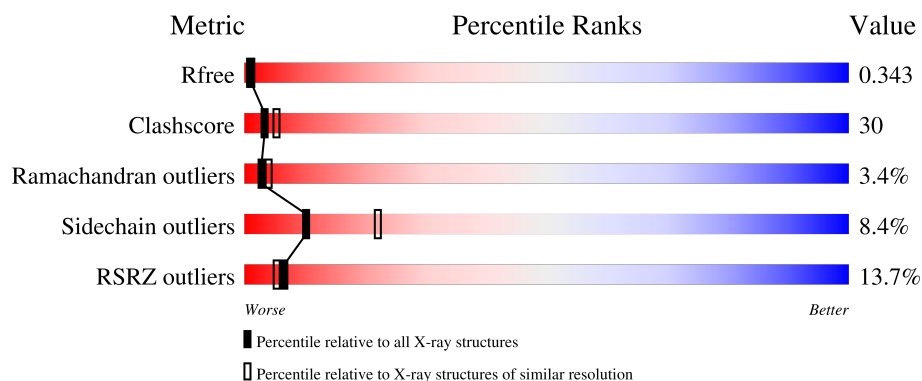
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	332	14% 48% 41% 8% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2506 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.

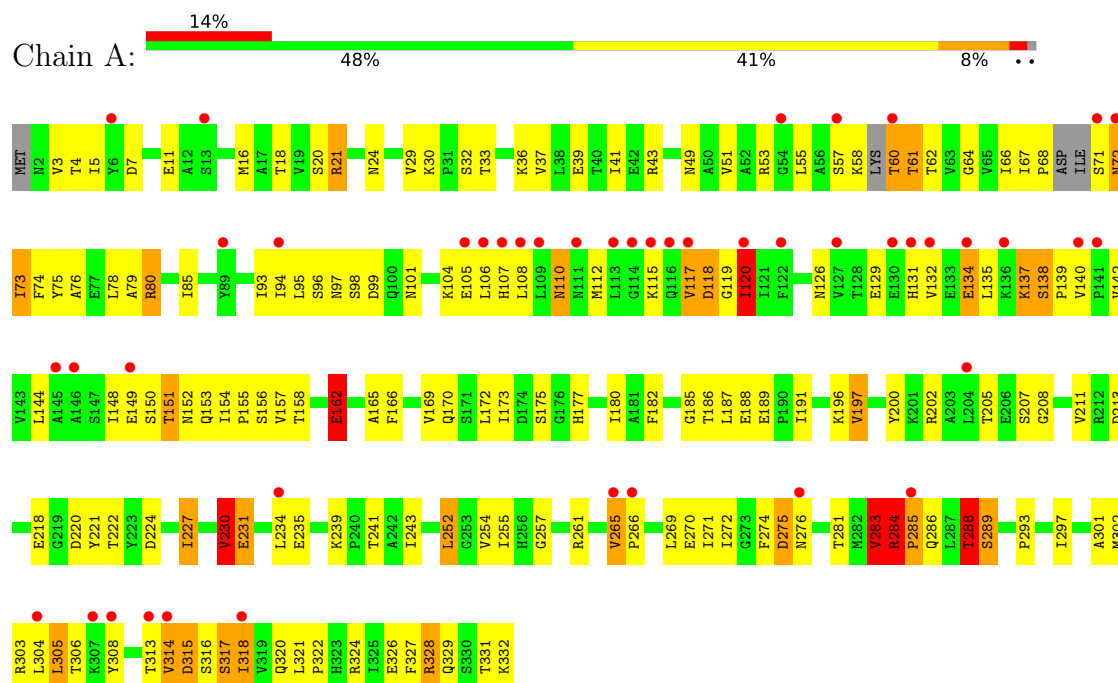
- Molecule 1 is a protein called Glucose-resistance amylase regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	0
			2506	1573	430	495	8			

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: Glucose-resistance amylase regulator



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	74.46Å 74.46Å 238.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.50 30.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.3 (30.00-2.50) 98.2 (30.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 2.51Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.271 , 0.337 0.271 , 0.343	Depositor DCC
R_{free} test set	711 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	56.1	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2506	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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	0.85	1/2541 (0.0%)	1.15	12/3447 (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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	288	THR	CA-C	-5.24	1.46	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	283	VAL	CB-CA-C	-8.71	95.58	110.30
1	A	328	ARG	NE-CZ-NH1	-8.04	113.46	121.50
1	A	288	THR	N-CA-C	-6.95	99.22	110.20
1	A	285	PRO	N-CA-C	-6.77	98.52	112.47
1	A	230	VAL	CB-CA-C	-6.16	103.72	112.22
1	A	284	ARG	N-CA-C	5.97	123.00	109.81
1	A	328	ARG	NE-CZ-NH2	5.96	124.56	119.20
1	A	231	GLU	N-CA-C	-5.88	103.84	111.02
1	A	120	ILE	N-CA-C	5.79	117.02	108.45
1	A	21	ARG	N-CA-C	-5.56	105.31	111.71
1	A	162	GLU	N-CA-C	-5.47	104.96	111.69
1	A	252	LEU	N-CA-C	-5.16	105.66	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	170	GLN	Sidechain

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	2506	0	2499	152	0
All	All	2506	0	2499	152	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 30.

All (152) 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:265:VAL:O	1:A:269:LEU:O	1.72	1.08
1:A:132:VAL:HG21	1:A:151:THR:HG21	1.46	0.96
1:A:149:GLU:CB	1:A:151:THR:OG1	2.17	0.93
1:A:49:ASN:OD1	1:A:51:VAL:HG12	1.76	0.85
1:A:120:ILE:HG23	1:A:142:VAL:HG22	1.59	0.84
1:A:288:THR:HG21	1:A:331:THR:OG1	1.77	0.83
1:A:108:LEU:O	1:A:112:MET:HB2	1.84	0.78
1:A:150:SER:C	1:A:152:ASN:HD22	1.93	0.76
1:A:73:ILE:HG23	1:A:74:PHE:N	2.01	0.75
1:A:150:SER:O	1:A:152:ASN:ND2	2.20	0.75
1:A:318:ILE:HG22	1:A:318:ILE:O	1.86	0.73
1:A:288:THR:HB	1:A:327:PHE:HA	1.69	0.73
1:A:281:THR:HG22	1:A:286:GLN:HE21	1.53	0.73
1:A:68:PRO:O	1:A:71:SER:N	2.22	0.72
1:A:286:GLN:HB2	1:A:329:GLN:NE2	2.05	0.72
1:A:188:GLU:N	1:A:188:GLU:OE1	2.21	0.72
1:A:162:GLU:OE2	1:A:202:ARG:NH1	2.23	0.71
1:A:132:VAL:CG2	1:A:151:THR:HG21	2.19	0.71
1:A:316:SER:O	1:A:318:ILE:N	2.24	0.70
1:A:32:SER:O	1:A:36:LYS:HG2	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:HIS:O	1:A:135:LEU:HG	1.91	0.69
1:A:243:ILE:HG13	1:A:269:LEU:HD11	1.75	0.69
1:A:5:ILE:HG22	1:A:16:MET:HG2	1.75	0.69
1:A:150:SER:C	1:A:152:ASN:ND2	2.51	0.68
1:A:154:ILE:O	1:A:154:ILE:HG23	1.94	0.68
1:A:68:PRO:HA	1:A:98:SER:O	1.94	0.68
1:A:101:ASN:O	1:A:105:GLU:HG2	1.94	0.67
1:A:58:LYS:CB	1:A:60:THR:HG23	2.25	0.67
1:A:220:ASP:O	1:A:221:TYR:HB2	1.94	0.67
1:A:265:VAL:O	1:A:269:LEU:C	2.39	0.66
1:A:284:ARG:O	1:A:286:GLN:N	2.28	0.66
1:A:288:THR:HG22	1:A:328:ARG:H	1.61	0.65
1:A:51:VAL:O	1:A:55:LEU:HG	1.96	0.64
1:A:67:ILE:HD11	1:A:75:TYR:HB3	1.79	0.64
1:A:106:LEU:O	1:A:110:ASN:N	2.31	0.63
1:A:107:HIS:HA	1:A:110:ASN:HB3	1.81	0.63
1:A:4:THR:N	1:A:7:ASP:OD2	2.31	0.62
1:A:61:THR:HG23	1:A:117:VAL:HG11	1.81	0.62
1:A:274:PHE:C	1:A:275:ASP:O	2.40	0.62
1:A:112:MET:CE	1:A:139:PRO:HD2	2.29	0.62
1:A:281:THR:HG22	1:A:286:GLN:NE2	2.15	0.61
1:A:308:TYR:OH	1:A:316:SER:HB3	2.00	0.61
1:A:61:THR:HA	1:A:117:VAL:HB	1.81	0.61
1:A:112:MET:O	1:A:112:MET:HE3	2.01	0.60
1:A:151:THR:O	1:A:154:ILE:HG22	2.01	0.60
1:A:205:THR:O	1:A:208:GLY:N	2.30	0.60
1:A:220:ASP:OD1	1:A:222:THR:HG23	2.00	0.60
1:A:275:ASP:O	1:A:289:SER:OG	2.19	0.60
1:A:308:TYR:OH	1:A:316:SER:CB	2.50	0.60
1:A:177:HIS:HB2	1:A:180:ILE:HD11	1.84	0.60
1:A:73:ILE:CG2	1:A:74:PHE:N	2.64	0.60
1:A:37:VAL:O	1:A:41:ILE:HG13	2.02	0.58
1:A:321:LEU:HB3	1:A:322:PRO:HD2	1.84	0.58
1:A:118:ASP:HB3	1:A:140:VAL:HG11	1.85	0.58
1:A:186:THR:HA	1:A:218:GLU:OE1	2.03	0.57
1:A:57:SER:O	1:A:58:LYS:CB	2.53	0.57
1:A:67:ILE:O	1:A:97:ASN:HA	2.05	0.56
1:A:257:GLY:O	1:A:261:ARG:HG2	2.06	0.55
1:A:288:THR:CG2	1:A:328:ARG:H	2.19	0.55
1:A:254:VAL:HG12	1:A:271:ILE:CD1	2.36	0.55
1:A:284:ARG:O	1:A:285:PRO:C	2.47	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:VAL:O	1:A:30:LYS:C	2.50	0.55
1:A:230:VAL:HG13	1:A:254:VAL:HG13	1.88	0.55
1:A:73:ILE:HG23	1:A:74:PHE:H	1.72	0.55
1:A:303:ARG:O	1:A:306:THR:HB	2.07	0.55
1:A:108:LEU:HD21	1:A:134:GLU:HB3	1.89	0.54
1:A:255:ILE:HG23	1:A:265:VAL:HG21	1.89	0.54
1:A:230:VAL:HG22	1:A:254:VAL:HA	1.89	0.54
1:A:149:GLU:C	1:A:151:THR:H	2.16	0.54
1:A:254:VAL:HG12	1:A:271:ILE:HD13	1.89	0.53
1:A:314:VAL:HG12	1:A:315:ASP:N	2.23	0.53
1:A:211:VAL:HG12	1:A:211:VAL:O	2.09	0.53
1:A:18:THR:OG1	1:A:21:ARG:NH1	2.43	0.52
1:A:144:LEU:HG	1:A:154:ILE:CG1	2.40	0.52
1:A:265:VAL:O	1:A:266:PRO:C	2.49	0.51
1:A:304:LEU:O	1:A:305:LEU:C	2.52	0.51
1:A:177:HIS:HD2	1:A:241:THR:OG1	1.94	0.51
1:A:126:ASN:ND2	1:A:189:GLU:OE2	2.43	0.51
1:A:20:SER:O	1:A:24:ASN:ND2	2.43	0.51
1:A:3:VAL:O	1:A:53:ARG:HD2	2.11	0.50
1:A:272:ILE:HD11	1:A:331:THR:HG21	1.91	0.50
1:A:108:LEU:CD2	1:A:134:GLU:HB3	2.41	0.50
1:A:73:ILE:CG2	1:A:74:PHE:H	2.23	0.50
1:A:149:GLU:C	1:A:151:THR:N	2.69	0.50
1:A:107:HIS:CA	1:A:110:ASN:HB3	2.41	0.50
1:A:230:VAL:CG2	1:A:254:VAL:HA	2.41	0.50
1:A:154:ILE:O	1:A:154:ILE:CG2	2.59	0.49
1:A:234:LEU:O	1:A:239:LYS:HE2	2.12	0.49
1:A:144:LEU:HG	1:A:154:ILE:HG13	1.94	0.49
1:A:112:MET:SD	1:A:138:SER:HB3	2.52	0.49
1:A:150:SER:HA	1:A:152:ASN:HD21	1.77	0.49
1:A:286:GLN:HB2	1:A:329:GLN:HE21	1.77	0.48
1:A:173:ILE:C	1:A:175:SER:H	2.20	0.48
1:A:71:SER:N	1:A:97:ASN:OD1	2.47	0.47
1:A:95:LEU:HD23	1:A:96:SER:N	2.29	0.47
1:A:148:ILE:HD11	1:A:191:ILE:N	2.29	0.47
1:A:157:VAL:HG21	1:A:301:ALA:HA	1.96	0.47
1:A:265:VAL:C	1:A:269:LEU:O	2.52	0.47
1:A:137:LYS:O	1:A:138:SER:C	2.57	0.47
1:A:324:ARG:NH2	1:A:326:GLU:OE1	2.42	0.47
1:A:39:GLU:O	1:A:43:ARG:HB2	2.14	0.47
1:A:293:PRO:O	1:A:297:ILE:HG13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLY:HA2	1:A:140:VAL:HB	1.97	0.46
1:A:252:LEU:HG	1:A:283:VAL:HG22	1.96	0.46
1:A:60:THR:O	1:A:61:THR:HB	2.15	0.46
1:A:156:SER:H	1:A:317:SER:HA	1.79	0.46
1:A:165:ALA:HB2	1:A:196:LYS:HA	1.98	0.46
1:A:11:GLU:OE1	1:A:43:ARG:NH2	2.48	0.46
1:A:85:ILE:HB	1:A:302:MET:HG2	1.98	0.46
1:A:169:VAL:HG21	1:A:200:TYR:HD2	1.80	0.46
1:A:108:LEU:O	1:A:112:MET:CB	2.59	0.46
1:A:60:THR:O	1:A:61:THR:CB	2.63	0.45
1:A:172:LEU:HD23	1:A:172:LEU:HA	1.78	0.45
1:A:182:PHE:HE1	1:A:197:VAL:HG23	1.82	0.45
1:A:286:GLN:HB2	1:A:329:GLN:HE22	1.80	0.45
1:A:265:VAL:HG12	1:A:266:PRO:N	2.31	0.45
1:A:231:GLU:HG2	1:A:235:GLU:OE2	2.16	0.45
1:A:78:LEU:C	1:A:78:LEU:HD23	2.42	0.45
1:A:99:ASP:O	1:A:104:LYS:HE2	2.17	0.45
1:A:117:VAL:HG12	1:A:118:ASP:N	2.31	0.45
1:A:284:ARG:HD3	1:A:284:ARG:HA	1.49	0.44
1:A:173:ILE:C	1:A:175:SER:N	2.76	0.44
1:A:220:ASP:O	1:A:221:TYR:CB	2.65	0.44
1:A:224:ASP:O	1:A:227:ILE:N	2.51	0.44
1:A:112:MET:HE1	1:A:139:PRO:HD2	1.97	0.44
1:A:154:ILE:HD12	1:A:155:PRO:HD2	1.99	0.44
1:A:187:LEU:N	1:A:218:GLU:OE1	2.45	0.44
1:A:110:ASN:C	1:A:112:MET:N	2.74	0.44
1:A:144:LEU:CD1	1:A:154:ILE:HG21	2.48	0.44
1:A:144:LEU:HD11	1:A:154:ILE:HG21	1.99	0.43
1:A:76:ALA:O	1:A:79:ALA:HB3	2.19	0.43
1:A:270:GLU:OE2	1:A:332:LYS:HB2	2.18	0.42
1:A:110:ASN:C	1:A:112:MET:H	2.28	0.42
1:A:173:ILE:O	1:A:175:SER:N	2.52	0.42
1:A:144:LEU:HG	1:A:154:ILE:HG12	2.03	0.41
1:A:67:ILE:HG13	1:A:68:PRO:HD2	2.02	0.41
1:A:66:ILE:HD12	1:A:107:HIS:HB3	2.03	0.41
1:A:30:LYS:O	1:A:33:THR:HB	2.20	0.41
1:A:62:THR:O	1:A:118:ASP:HA	2.20	0.41
1:A:187:LEU:O	1:A:188:GLU:C	2.63	0.41
1:A:252:LEU:CG	1:A:283:VAL:HG22	2.49	0.41
1:A:72:ASN:ND2	1:A:73:ILE:HG22	2.36	0.41
1:A:276:ASN:OD1	1:A:328:ARG:NH2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ARG:HA	1:A:80:ARG:HD3	1.91	0.40
1:A:115:LYS:O	1:A:118:ASP:HB2	2.21	0.40
1:A:158:THR:O	1:A:320:GLN:HA	2.21	0.40
1:A:165:ALA:O	1:A:166:PHE:C	2.64	0.40
1:A:101:ASN:HA	1:A:104:LYS:HB2	2.03	0.40
1:A:230:VAL:O	1:A:234:LEU:HD12	2.22	0.40
1:A:49:ASN:OD1	1:A:51:VAL:CG1	2.60	0.40
1:A:64:GLY:HA2	1:A:94:ILE:HB	2.04	0.40
1:A:185:GLY:HA3	1:A:221:TYR:CZ	2.56	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	322/332 (97%)	275 (85%)	36 (11%)	11 (3%)	3 4

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	275	ASP
1	A	314	VAL
1	A	110	ASN
1	A	117	VAL
1	A	151	THR
1	A	317	SER
1	A	61	THR
1	A	305	LEU
1	A	313	THR
1	A	318	ILE

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	275/289 (95%)	252 (92%)	23 (8%)	10	22

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	THR
1	A	73	ILE
1	A	80	ARG
1	A	93	ILE
1	A	118	ASP
1	A	120	ILE
1	A	129	GLU
1	A	134	GLU
1	A	137	LYS
1	A	138	SER
1	A	153	GLN
1	A	162	GLU
1	A	197	VAL
1	A	207	SER
1	A	213	ASP
1	A	227	ILE
1	A	230	VAL
1	A	265	VAL
1	A	283	VAL
1	A	284	ARG
1	A	288	THR
1	A	289	SER
1	A	315	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	72	ASN

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Mol	Chain	Res	Type
1	A	126	ASN
1	A	152	ASN
1	A	177	HIS
1	A	286	GLN
1	A	320	GLN
1	A	329	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	328/332 (98%)	0.86	45 (13%) 6 5	33, 54, 101, 110	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	54	GLY	4.5
1	A	131	HIS	4.3
1	A	318	ILE	3.7
1	A	71	SER	3.7
1	A	116	GLN	3.7
1	A	57	SER	3.6
1	A	130	GLU	3.4
1	A	132	VAL	3.4
1	A	105	GLU	3.3
1	A	13	SER	3.2
1	A	120	ILE	3.2
1	A	111	ASN	3.1
1	A	72	ASN	3.1
1	A	106	LEU	3.0
1	A	145	ALA	3.0
1	A	114	GLY	3.0
1	A	6	TYR	2.9
1	A	127	VAL	2.8
1	A	141	PRO	2.7
1	A	117	VAL	2.7
1	A	136	LYS	2.7
1	A	122	PHE	2.7
1	A	314	VAL	2.6
1	A	89	TYR	2.6
1	A	140	VAL	2.5
1	A	265	VAL	2.5
1	A	60	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	266	PRO	2.5
1	A	108	LEU	2.5
1	A	304	LEU	2.5
1	A	113	LEU	2.4
1	A	276	ASN	2.4
1	A	115	LYS	2.4
1	A	109	LEU	2.3
1	A	107	HIS	2.3
1	A	313	THR	2.3
1	A	134	GLU	2.3
1	A	94	ILE	2.2
1	A	234	LEU	2.2
1	A	285	PRO	2.1
1	A	204	LEU	2.1
1	A	308	TYR	2.0
1	A	146	ALA	2.0
1	A	149	GLU	2.0
1	A	307	LYS	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.