



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:33 AM UTC

PDB ID : 1ZC6 / pdb_00001zc6
Title : Crystal Structure of Putative N-acetylglucosamine Kinase from *Chromobacterium violaceum*. Northeast Structural Genomics Target Cvr23.
Authors : Vorobiev, S.M.; Kuzin, A.; Forouhar, F.; Abashidze, M.; Acton, T.B.; Xiao, R.; Ma, L.-C.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2005-04-11
Resolution : 2.20 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

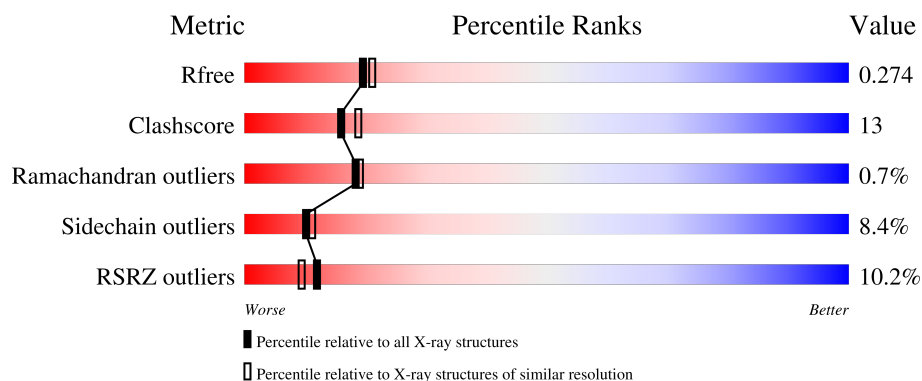
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	305	9% 70% 19% 7%
1	B	305	10% 71% 20% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4266 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 probable N-acetylglucosamine kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	Se	0	0	0
			2066	1293	378	389	2	4			
1	B	293	Total	C	N	O	S	Se	0	0	0
			2116	1323	388	398	2	5			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	MSE	MET	modified residue	UNP Q7NU07
A	37	MSE	MET	modified residue	UNP Q7NU07
A	173	MSE	MET	modified residue	UNP Q7NU07
A	199	MSE	MET	modified residue	UNP Q7NU07
A	199	MSE	MET	modified residue	UNP Q7NU07
A	298	LEU	-	cloning artifact	UNP Q7NU07
A	299	GLU	-	cloning artifact	UNP Q7NU07
A	300	HIS	-	cloning artifact	UNP Q7NU07
A	301	HIS	-	cloning artifact	UNP Q7NU07
A	302	HIS	-	cloning artifact	UNP Q7NU07
A	303	HIS	-	cloning artifact	UNP Q7NU07
A	304	HIS	-	cloning artifact	UNP Q7NU07
A	305	HIS	-	cloning artifact	UNP Q7NU07
B	6	MSE	MET	modified residue	UNP Q7NU07
B	37	MSE	MET	modified residue	UNP Q7NU07
B	173	MSE	MET	modified residue	UNP Q7NU07
B	199	MSE	MET	modified residue	UNP Q7NU07
B	199	MSE	MET	modified residue	UNP Q7NU07
B	298	LEU	-	cloning artifact	UNP Q7NU07
B	299	GLU	-	cloning artifact	UNP Q7NU07
B	300	HIS	-	cloning artifact	UNP Q7NU07
B	301	HIS	-	cloning artifact	UNP Q7NU07
B	302	HIS	-	cloning artifact	UNP Q7NU07
B	303	HIS	-	cloning artifact	UNP Q7NU07
B	304	HIS	-	cloning artifact	UNP Q7NU07

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Chain	Residue	Modelled	Actual	Comment	Reference
B	305	HIS	-	cloning artifact	UNP Q7NU07

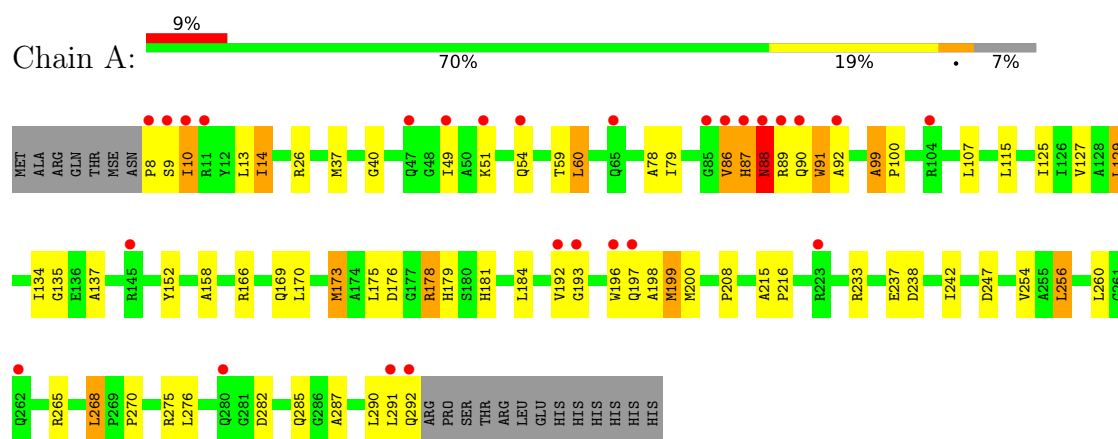
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	29	Total	O	0	0
			29	29		
2	B	55	Total	O	0	0
			55	55		

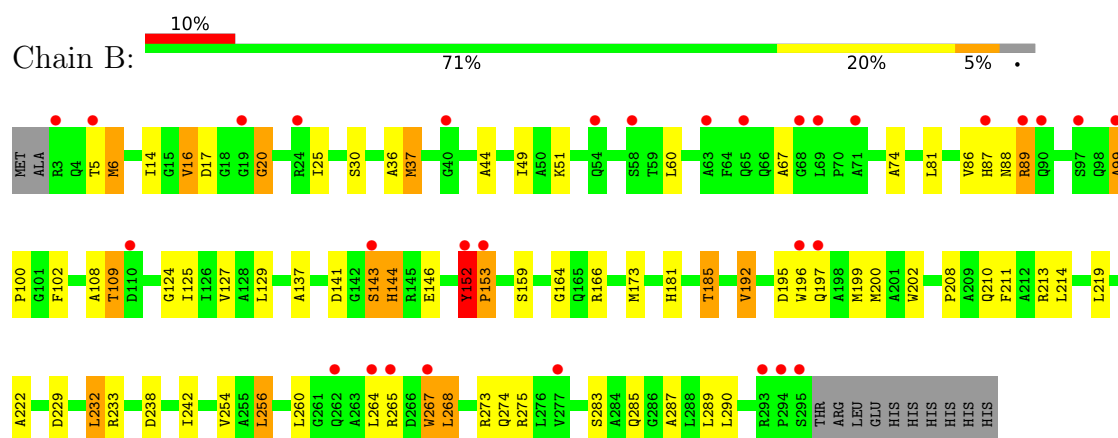
3 Residue-property plots [i](#)

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: probable N-acetylglucosamine kinase



- Molecule 1: probable N-acetylglucosamine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	178.25Å 178.25Å 117.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.42 – 2.20 29.42 – 2.20	Depositor EDS
% Data completeness (in resolution range)	92.6 (29.42-2.20) 92.6 (29.42-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.241 , 0.267 0.252 , 0.274	Depositor DCC
R_{free} test set	3569 reflections (3.93%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 28.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4266	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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	0.53	4/2110 (0.2%)	0.94	8/2867 (0.3%)
1	B	0.59	6/2160 (0.3%)	0.96	9/2936 (0.3%)
All	All	0.56	10/4270 (0.2%)	0.95	17/5803 (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	B	0	1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	173	MSE	SE-CE	-8.40	1.70	1.95
1	B	200	MSE	SE-CE	-8.12	1.71	1.95
1	B	6	MSE	SE-CE	-6.91	1.74	1.95
1	A	200	MSE	SE-CE	-6.36	1.76	1.95
1	B	6	MSE	CG-SE	-6.17	1.76	1.95
1	B	173	MSE	SE-CE	-5.94	1.77	1.95
1	A	199	MSE	SE-CE	-5.65	1.78	1.95
1	A	37	MSE	SE-CE	-5.49	1.78	1.95
1	B	37	MSE	CG-SE	-5.31	1.79	1.95
1	B	199	MSE	SE-CE	-5.14	1.80	1.95

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	152	TYR	N-CA-C	7.60	126.60	109.81
1	B	159	SER	N-CA-C	-6.85	100.09	109.95
1	A	87	HIS	N-CA-C	-6.29	105.88	112.93
1	A	79	ILE	N-CA-C	6.29	116.91	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	ALA	N-CA-C	5.57	116.91	109.83
1	A	14	ILE	N-CA-C	5.49	115.80	108.11
1	B	16	VAL	N-CA-C	5.43	115.71	108.11
1	A	91	TRP	N-CA-C	-5.30	105.59	111.36
1	B	99	ALA	N-CA-C	5.17	116.34	109.93
1	B	109	THR	CA-C-N	5.16	127.51	120.54
1	B	109	THR	C-N-CA	5.16	127.51	120.54
1	B	14	ILE	N-CA-C	5.08	115.22	108.11
1	A	127	VAL	N-CA-C	-5.06	100.29	107.77
1	A	247	ASP	CA-C-N	5.03	125.06	119.32
1	A	247	ASP	C-N-CA	5.03	125.06	119.32
1	B	195	ASP	N-CA-C	5.02	115.67	108.74
1	B	222	ALA	N-CA-C	5.01	117.47	111.71

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	152	TYR	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	A	2066	0	2015	54	0
1	B	2116	0	2053	62	0
2	A	29	0	0	2	0
2	B	55	0	0	2	0
All	All	4266	0	4068	109	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 (109) 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:169:GLN:HG2	1:A:173:MSE:HE2	1.39	1.05
1:A:87:HIS:HB2	1:A:91:TRP:HB2	1.53	0.88
1:A:54:GLN:HB3	2:A:585:HOH:O	1.74	0.87
1:B:36:ALA:HB2	1:B:67:ALA:HB2	1.66	0.78
1:A:282:ASP:H	1:A:285:GLN:HE21	1.35	0.74
1:B:20:GLY:H	1:B:44:ALA:HB2	1.53	0.74
1:A:88:ASN:HA	1:A:92:ALA:HB3	1.68	0.73
1:A:125:ILE:HG13	1:A:254:VAL:HG13	1.72	0.71
1:B:254:VAL:HG21	2:B:552:HOH:O	1.91	0.71
1:B:256:LEU:HD12	1:B:268:LEU:HD21	1.73	0.70
1:B:229:ASP:OD1	1:B:267:TRP:HZ3	1.76	0.68
1:A:233:ARG:HG2	1:B:233:ARG:NH1	2.07	0.68
1:B:181:HIS:HA	1:B:185:THR:HG21	1.76	0.67
1:B:265:ARG:NH2	1:B:273:ARG:HH11	1.93	0.67
1:B:265:ARG:NH1	1:B:265:ARG:HB2	2.09	0.66
1:A:192:VAL:HG11	1:A:199:MSE:HA	1.78	0.66
1:A:233:ARG:CG	1:B:233:ARG:HD2	2.26	0.66
1:A:193:GLY:HA3	1:A:198:ALA:HB1	1.77	0.65
1:A:170:LEU:HA	1:A:173:MSE:HE3	1.78	0.64
1:A:87:HIS:HB3	1:A:90:GLN:HB3	1.80	0.64
1:B:210:GLN:HG3	1:B:213:ARG:NH2	2.14	0.63
1:B:5:THR:HG21	1:B:30:SER:HB2	1.81	0.63
1:B:125:ILE:HG13	1:B:137:ALA:HB2	1.80	0.62
1:A:282:ASP:H	1:A:285:GLN:NE2	1.98	0.62
1:B:197:GLN:HB3	2:B:502:HOH:O	2.00	0.61
1:B:124:GLY:C	1:B:125:ILE:HD12	2.26	0.60
1:B:181:HIS:CD2	1:B:185:THR:CG2	2.85	0.60
1:A:197:GLN:H	1:A:197:GLN:CD	2.10	0.59
1:A:54:GLN:CB	2:A:585:HOH:O	2.39	0.59
1:A:291:LEU:HD12	1:A:292:GLN:HG3	1.83	0.58
1:B:49:ILE:HD12	1:B:49:ILE:N	2.17	0.58
1:A:233:ARG:HG2	1:B:233:ARG:HD2	1.86	0.57
1:A:40:GLY:HA3	1:A:59:THR:OG1	2.04	0.57
1:A:233:ARG:HG2	1:B:233:ARG:HH11	1.71	0.56
1:B:267:TRP:CD1	1:B:267:TRP:N	2.73	0.56
1:B:192:VAL:HG22	1:B:202:TRP:CG	2.41	0.55
1:A:176:ASP:OD1	1:A:178:ARG:HG2	2.07	0.54
1:A:265:ARG:HA	1:A:268:LEU:HD22	1.90	0.54
1:B:256:LEU:CD1	1:B:268:LEU:HD21	2.38	0.53
1:B:219:LEU:HD23	1:B:232:LEU:HD21	1.91	0.53
1:A:256:LEU:HD12	1:A:268:LEU:HD21	1.91	0.53
1:A:51:LYS:HA	1:A:54:GLN:HE21	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:ILE:HD12	1:B:49:ILE:H	1.75	0.51
1:B:265:ARG:HB2	1:B:265:ARG:CZ	2.40	0.51
1:A:193:GLY:HA3	1:A:198:ALA:CB	2.41	0.51
1:A:87:HIS:O	1:A:88:ASN:C	2.54	0.51
1:B:287:ALA:O	1:B:290:LEU:HB2	2.11	0.51
1:B:129:LEU:HB3	1:B:260:LEU:HD12	1.93	0.51
1:A:8:PRO:HG2	1:A:9:SER:H	1.76	0.50
1:A:129:LEU:HD22	1:A:256:LEU:HD21	1.94	0.50
1:B:89:ARG:HB2	1:B:89:ARG:HH11	1.76	0.50
1:B:285:GLN:O	1:B:289:LEU:HD13	2.12	0.49
1:B:229:ASP:OD1	1:B:267:TRP:CZ3	2.63	0.49
1:A:125:ILE:HG22	1:A:137:ALA:CB	2.43	0.48
1:A:254:VAL:HG21	1:A:275:ARG:NH2	2.27	0.48
1:A:233:ARG:NE	1:B:233:ARG:HD2	2.28	0.48
1:B:152:TYR:CG	1:B:208:PRO:HG3	2.48	0.48
1:B:88:ASN:C	1:B:88:ASN:HD22	2.19	0.48
1:A:49:ILE:HD12	1:A:49:ILE:N	2.28	0.48
1:A:51:LYS:HD2	1:A:54:GLN:NE2	2.28	0.48
1:A:10:ILE:HG13	1:A:78:ALA:HB2	1.95	0.48
1:A:10:ILE:HD13	1:A:10:ILE:O	2.14	0.48
1:B:264:LEU:O	1:B:267:TRP:HD1	1.97	0.47
1:A:270:PRO:HD2	1:B:267:TRP:CZ3	2.49	0.47
1:A:129:LEU:HG	1:A:158:ALA:O	2.15	0.47
1:B:74:ALA:O	1:B:102:PHE:HA	2.15	0.47
1:A:173:MSE:HB3	1:A:179:HIS:HB2	1.96	0.46
1:A:10:ILE:HD11	1:A:13:LEU:HD21	1.98	0.45
1:A:287:ALA:O	1:A:290:LEU:HB3	2.16	0.45
1:B:89:ARG:HH11	1:B:89:ARG:CB	2.29	0.45
1:A:134:ILE:HG12	1:A:135:GLY:N	2.30	0.45
1:A:125:ILE:CG1	1:A:254:VAL:HG13	2.44	0.45
1:A:26:ARG:CD	1:A:285:GLN:HG3	2.47	0.45
1:A:14:ILE:HD13	1:A:60:LEU:HD11	1.98	0.44
1:B:233:ARG:HH11	1:B:233:ARG:HG3	1.83	0.44
1:B:232:LEU:HD12	1:B:232:LEU:HA	1.79	0.44
1:B:274:GLN:HG3	1:B:275:ARG:HG3	2.00	0.43
1:A:87:HIS:O	1:A:90:GLN:HB3	2.17	0.43
1:A:152:TYR:CG	1:A:208:PRO:HG3	2.52	0.43
1:B:152:TYR:CG	1:B:153:PRO:N	2.86	0.43
1:B:124:GLY:O	1:B:125:ILE:HD12	2.19	0.43
1:B:238:ASP:O	1:B:242:ILE:HG12	2.18	0.43
1:B:197:GLN:CD	1:B:197:GLN:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:HIS:HD2	1:B:185:THR:CG2	2.30	0.42
1:A:197:GLN:CD	1:A:197:GLN:N	2.77	0.42
1:B:17:ASP:CG	1:B:283:SER:HB3	2.45	0.42
1:B:127:VAL:HG13	1:B:127:VAL:O	2.19	0.42
1:B:141:ASP:OD1	1:B:143:SER:HB2	2.20	0.42
1:B:265:ARG:NH2	1:B:273:ARG:HD3	2.35	0.42
1:B:267:TRP:CD1	1:B:267:TRP:H	2.37	0.42
1:A:233:ARG:HE	1:B:233:ARG:HD2	1.84	0.42
1:B:86:VAL:HB	1:B:109:THR:HG22	2.02	0.42
1:A:99:ALA:HA	1:A:100:PRO:HD3	1.94	0.41
1:A:88:ASN:O	1:A:89:ARG:HB2	2.20	0.41
1:A:233:ARG:O	1:A:237:GLU:HG3	2.19	0.41
1:B:16:VAL:HG22	1:B:25:ILE:HG12	2.01	0.41
1:B:86:VAL:HG21	1:B:108:ALA:O	2.21	0.41
1:B:232:LEU:HG	1:B:264:LEU:HD23	2.01	0.41
1:B:87:HIS:HE2	1:B:146:GLU:CD	2.29	0.41
1:B:6:MSE:HE2	1:B:6:MSE:HB3	1.99	0.41
1:A:238:ASP:O	1:A:242:ILE:HG12	2.21	0.41
1:B:99:ALA:HA	1:B:100:PRO:HD3	1.89	0.41
1:B:260:LEU:HD13	1:B:264:LEU:HD12	2.03	0.41
1:A:215:ALA:N	1:A:216:PRO:HD2	2.36	0.41
1:B:137:ALA:O	1:B:144:HIS:HA	2.22	0.40
1:B:88:ASN:C	1:B:88:ASN:ND2	2.79	0.40
1:B:164:GLY:HA3	1:B:211:PHE:O	2.22	0.40
1:A:86:VAL:C	1:A:88:ASN:H	2.29	0.40
1:A:178:ARG:HG2	1:A:178:ARG:H	1.78	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	283/305 (93%)	270 (95%)	12 (4%)	1 (0%)	30	34
1	B	291/305 (95%)	281 (97%)	7 (2%)	3 (1%)	12	11
All	All	574/610 (94%)	551 (96%)	19 (3%)	4 (1%)	18	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	B	152	TYR
1	B	153	PRO
1	B	20	GLY

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	195/209 (93%)	178 (91%)	17 (9%)	9	10
1	B	199/209 (95%)	183 (92%)	16 (8%)	11	13
All	All	394/418 (94%)	361 (92%)	33 (8%)	10	11

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE
1	A	60	LEU
1	A	86	VAL
1	A	88	ASN
1	A	107	LEU
1	A	115	LEU
1	A	129	LEU
1	A	166	ARG
1	A	175	LEU
1	A	178	ARG
1	A	181	HIS
1	A	184	LEU

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Mol	Chain	Res	Type
1	A	196	TRP
1	A	256	LEU
1	A	260	LEU
1	A	268	LEU
1	A	276	LEU
1	B	37	MSE
1	B	51	LYS
1	B	60	LEU
1	B	81	LEU
1	B	89	ARG
1	B	143	SER
1	B	144	HIS
1	B	166	ARG
1	B	185	THR
1	B	192	VAL
1	B	196	TRP
1	B	214	LEU
1	B	232	LEU
1	B	256	LEU
1	B	267	TRP
1	B	268	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	90	GLN
1	A	172	GLN
1	A	197	GLN
1	A	210	GLN
1	A	285	GLN
1	B	28	HIS
1	B	66	GLN
1	B	88	ASN
1	B	169	GLN
1	B	179	HIS
1	B	181	HIS
1	B	210	GLN
1	B	280	GLN
1	B	285	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	281/305 (92%)	0.82	27 (9%) 13 11	26, 40, 60, 88	0
1	B	288/305 (94%)	0.68	31 (10%) 11 8	22, 37, 58, 67	0
All	All	569/610 (93%)	0.75	58 (10%) 12 9	22, 39, 60, 88	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	88	ASN	8.2
1	B	295	SER	7.7
1	A	86	VAL	7.2
1	A	89	ARG	6.9
1	A	292	GLN	6.4
1	B	3	ARG	5.4
1	A	85	GLY	5.3
1	A	87	HIS	5.0
1	A	196	TRP	4.9
1	A	8	PRO	4.7
1	A	90	GLN	4.4
1	B	197	GLN	3.8
1	A	291	LEU	3.6
1	B	65	GLN	3.5
1	B	196	TRP	3.5
1	A	49	ILE	3.4
1	B	267	TRP	3.4
1	A	192	VAL	3.3
1	B	97	SER	3.3
1	B	58	SER	3.2
1	B	87	HIS	3.2
1	B	19	GLY	3.2
1	B	110	ASP	3.1
1	A	262	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	197	GLN	3.1
1	B	294	PRO	3.1
1	A	9	SER	3.0
1	A	47	GLN	2.9
1	A	280	GLN	2.9
1	B	69	LEU	2.9
1	B	143	SER	2.7
1	A	54	GLN	2.7
1	B	24	ARG	2.6
1	B	63	ALA	2.6
1	B	68	GLY	2.5
1	A	10	ILE	2.5
1	B	90	GLN	2.5
1	B	152	TYR	2.5
1	A	223	ARG	2.5
1	A	65	GLN	2.5
1	B	262	GLN	2.4
1	A	145	ARG	2.4
1	B	71	ALA	2.4
1	B	265	ARG	2.3
1	B	5	THR	2.3
1	A	11	ARG	2.3
1	B	89	ARG	2.3
1	B	153	PRO	2.3
1	A	51	LYS	2.3
1	B	293	ARG	2.2
1	B	54	GLN	2.2
1	B	277	VAL	2.2
1	B	264	LEU	2.2
1	B	40	GLY	2.1
1	A	104	ARG	2.1
1	B	99	ALA	2.1
1	A	92	ALA	2.0
1	A	193	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.