



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:18 AM UTC

PDB ID : 1PCZ / pdb_00001pcz
Title : STRUCTURE OF TATA-BINDING PROTEIN
Authors : Dedecker, B.S.; Sigler, P.B.
Deposited on : 1996-10-04
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
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

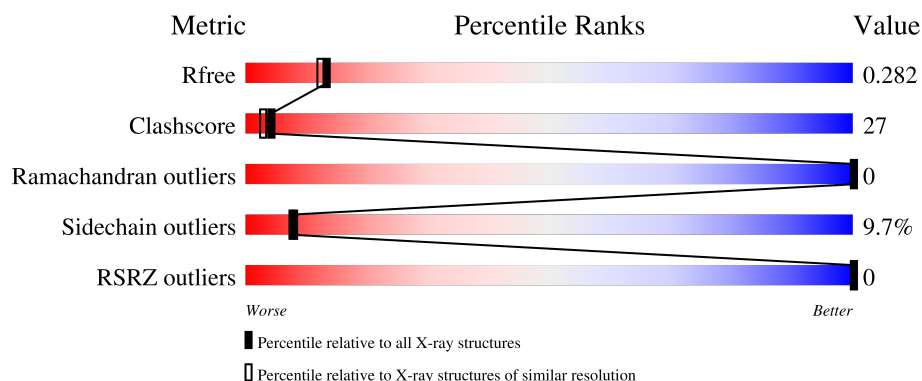
1 Overall quality at a glance

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	191	
1	B	191	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3120 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 TATA-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1426	919	237	264	6			
1	B	183	Total	C	N	O	S	0	0	0
			1426	919	237	264	6			

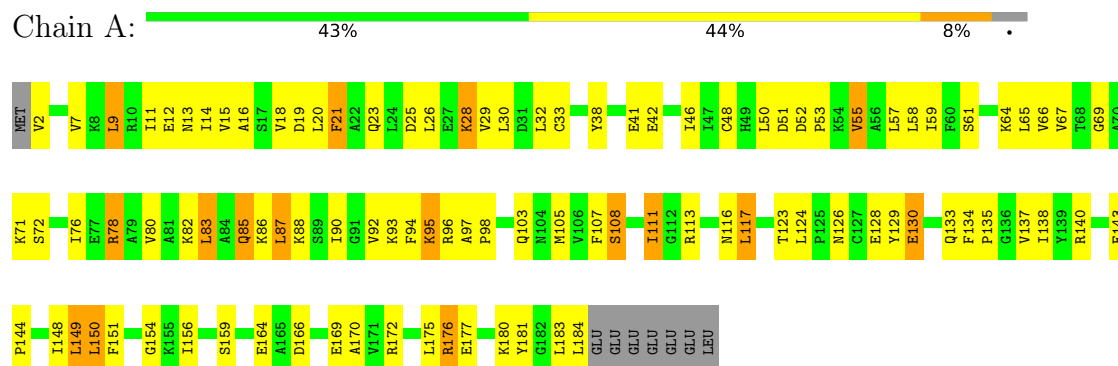
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	136	Total	O	0	0
			136	136		
2	B	132	Total	O	0	0
			132	132		

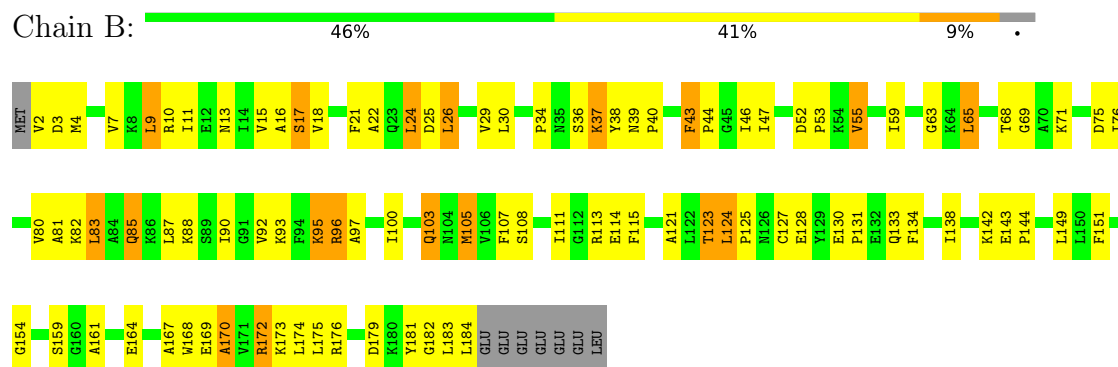
3 Residue-property plots

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: TATA-BINDING PROTEIN



• Molecule 1: TATA-BINDING PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	89.62Å 89.62Å 141.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.20 6.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (6.00-2.20) 94.2 (6.00-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.10Å)	Xtriage
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.218 , 0.300 0.206 , 0.282	Depositor DCC
R_{free} test set	3594 reflections (10.17%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 77.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.058 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3120	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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	0.84	0/1449	1.20	15/1956 (0.8%)
1	B	0.88	1/1449 (0.1%)	1.24	17/1956 (0.9%)
All	All	0.86	1/2898 (0.0%)	1.22	32/3912 (0.8%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	105	MET	SD-CE	-6.49	1.63	1.79

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	37	LYS	N-CA-C	-7.43	96.29	108.41
1	A	130	GLU	CA-C-N	7.31	127.65	119.32
1	A	130	GLU	C-N-CA	7.31	127.65	119.32
1	A	108	SER	N-CA-C	-7.17	98.63	110.17
1	B	108	SER	N-CA-C	-7.14	99.36	110.42
1	A	97	ALA	N-CA-C	-7.12	100.47	110.29
1	A	116	ASN	N-CA-C	-7.12	96.03	108.20
1	B	39	ASN	CA-C-N	7.01	128.61	119.84
1	B	39	ASN	C-N-CA	7.01	128.61	119.84
1	B	97	ALA	N-CA-C	-6.60	101.75	110.07
1	B	43	PHE	CA-C-N	6.46	126.04	119.19
1	B	43	PHE	C-N-CA	6.46	126.04	119.19
1	A	33	CYS	CA-C-N	6.42	127.43	119.98
1	A	33	CYS	C-N-CA	6.42	127.43	119.98
1	B	170	ALA	N-CA-C	-6.19	104.45	111.14
1	B	34	PRO	N-CA-C	-6.09	104.76	113.47
1	B	154	GLY	N-CA-C	-5.72	107.48	114.92
1	A	154	GLY	N-CA-C	-5.61	107.63	114.92
1	B	114	GLU	N-CA-C	-5.56	103.05	110.55
1	A	58	LEU	N-CA-C	-5.52	98.28	107.99
1	A	85	GLN	N-CA-C	-5.43	105.00	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	124	LEU	CA-C-N	5.43	125.26	119.28
1	B	124	LEU	C-N-CA	5.43	125.26	119.28
1	A	64	LYS	N-CA-C	-5.41	101.65	110.20
1	B	17	SER	N-CA-C	-5.38	100.92	109.96
1	A	21	PHE	N-CA-C	-5.26	104.76	110.91
1	A	42	GLU	N-CA-C	-5.23	105.72	112.68
1	B	123	THR	N-CA-C	5.22	117.77	111.40
1	B	93	LYS	N-CA-C	-5.20	101.68	109.95
1	A	149	LEU	N-CA-C	-5.11	100.20	108.52
1	B	176	ARG	N-CA-C	-5.06	105.77	111.28
1	A	51	ASP	N-CA-C	-5.02	107.32	113.50

There are no chirality outliers.

There are no planarity outliers.

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	1426	0	1486	86	0
1	B	1426	0	1486	80	0
2	A	136	0	0	4	0
2	B	132	0	0	7	0
All	All	3120	0	2972	158	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 27.

All (158) 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:117:LEU:HG	1:A:137:VAL:HG23	1.50	0.93
1:B:113:ARG:NH1	1:B:113:ARG:HB3	1.97	0.79
1:A:149:LEU:HD21	1:B:149:LEU:HD21	1.67	0.76
1:A:11:ILE:HD13	1:A:105:MET:HE2	1.68	0.76
1:A:16:ALA:HB2	1:A:76:ILE:HD13	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ARG:NH1	1:A:184:LEU:HA	2.03	0.74
1:A:180:LYS:HD2	1:A:181:TYR:CE1	2.24	0.73
1:B:169:GLU:HG2	2:B:208:HOH:O	1.87	0.73
1:A:18:VAL:HB	1:A:65:LEU:HG	1.69	0.73
1:B:103:GLN:HE21	1:B:103:GLN:HA	1.57	0.70
1:A:57:LEU:HD13	1:A:83:LEU:HG	1.74	0.68
1:B:7:VAL:HG21	1:B:175:LEU:HD11	1.76	0.68
1:B:95:LYS:HD2	1:B:95:LYS:H	1.61	0.66
1:A:177:GLU:O	1:A:181:TYR:HD1	1.78	0.65
1:A:144:PRO:HG3	1:A:166:ASP:HB3	1.79	0.64
1:B:30:LEU:HD21	1:B:38:TYR:HB2	1.80	0.64
1:B:113:ARG:HB3	1:B:113:ARG:HH11	1.61	0.64
1:A:123:THR:HG22	1:A:181:TYR:HB3	1.78	0.64
1:A:12:GLU:HG3	1:A:108:SER:HB3	1.79	0.64
1:B:30:LEU:HD11	1:B:38:TYR:CD2	2.33	0.64
1:B:83:LEU:HD13	1:B:87:LEU:HD11	1.78	0.63
1:A:88:LYS:CG	1:A:94:PHE:HE1	2.11	0.63
1:A:88:LYS:HG2	1:A:94:PHE:HE1	1.63	0.63
1:A:13:ASN:HD21	1:A:15:VAL:HG23	1.64	0.62
1:A:113:ARG:HH11	1:A:184:LEU:HA	1.64	0.62
1:A:123:THR:HB	1:A:183:LEU:HD11	1.82	0.62
1:B:24:LEU:HD22	1:B:63:GLY:HA2	1.81	0.62
1:A:111:ILE:HD11	1:A:150:LEU:HD21	1.83	0.61
1:A:130:GLU:HG3	1:A:133:GLN:HB3	1.83	0.61
1:B:38:TYR:CZ	1:B:40:PRO:HG3	2.36	0.61
1:A:176:ARG:HD2	2:A:269:HOH:O	2.01	0.60
1:A:19:ASP:O	1:A:98:PRO:HA	2.01	0.60
1:A:7:VAL:HG21	1:A:175:LEU:HD11	1.83	0.60
1:B:36:SER:HA	1:B:47:ILE:O	2.01	0.60
1:B:111:ILE:HD11	1:B:115:PHE:CE2	2.36	0.60
1:A:88:LYS:HG2	1:A:94:PHE:CE1	2.37	0.59
1:A:149:LEU:HD21	1:B:149:LEU:CD2	2.31	0.59
1:A:30:LEU:HD11	1:A:38:TYR:CB	2.33	0.59
1:B:25:ASP:O	1:B:29:VAL:HG23	2.03	0.59
1:B:30:LEU:HD11	1:B:38:TYR:HD2	1.66	0.58
1:B:71:LYS:HG3	2:B:199:HOH:O	2.04	0.58
1:A:25:ASP:OD2	1:A:28:LYS:HG2	2.03	0.58
1:B:111:ILE:HB	1:B:175:LEU:HD21	1.85	0.58
1:A:30:LEU:HD11	1:A:38:TYR:HB3	1.85	0.57
1:B:128:GLU:HB2	1:B:138:ILE:HB	1.86	0.57
1:A:78:ARG:HG3	1:A:78:ARG:HH11	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:CD2	1:B:149:LEU:HD21	2.34	0.56
1:B:182:GLY:HA2	2:B:212:HOH:O	2.05	0.56
1:A:130:GLU:HG2	2:A:267:HOH:O	2.04	0.56
1:A:21:PHE:CD2	1:A:96:ARG:HB2	2.41	0.56
1:B:179:ASP:HB2	1:B:184:LEU:HD12	1.88	0.55
1:B:24:LEU:HD23	1:B:59:ILE:CG2	2.37	0.55
1:B:131:PRO:HA	1:B:134:PHE:O	2.06	0.55
1:A:78:ARG:HH11	1:A:78:ARG:CG	2.20	0.54
1:A:111:ILE:HG23	1:A:156:ILE:HD13	1.89	0.54
1:B:81:ALA:O	1:B:85:GLN:HG2	2.07	0.53
1:B:130:GLU:HB3	1:B:133:GLN:HB3	1.90	0.53
1:A:9:LEU:HD22	1:A:107:PHE:HB3	1.90	0.53
1:B:167:ALA:O	1:B:170:ALA:HB3	2.07	0.53
1:A:69:GLY:HA2	2:A:287:HOH:O	2.07	0.53
1:A:123:THR:CG2	1:A:181:TYR:HB3	2.38	0.53
1:A:87:LEU:HD23	1:A:94:PHE:HE2	1.74	0.52
1:A:149:LEU:HG	1:A:151:PHE:CZ	2.43	0.52
1:A:57:LEU:HD23	1:A:67:VAL:HA	1.90	0.52
1:A:76:ILE:O	1:A:80:VAL:HG23	2.09	0.52
1:B:144:PRO:HD3	2:B:293:HOH:O	2.10	0.52
1:A:48:CYS:HB2	1:A:59:ILE:HD12	1.91	0.51
1:A:95:LYS:HA	1:A:95:LYS:HE3	1.91	0.51
1:B:7:VAL:HG21	1:B:175:LEU:CD1	2.40	0.51
1:A:88:LYS:NZ	1:A:93:LYS:HA	2.26	0.50
1:B:9:LEU:HB2	1:B:168:TRP:CZ2	2.47	0.50
1:B:18:VAL:HG22	1:B:100:ILE:HG23	1.94	0.50
1:A:15:VAL:HG12	1:A:103:GLN:HB3	1.92	0.50
1:A:25:ASP:O	1:A:29:VAL:HG23	2.12	0.50
1:B:22:ALA:O	1:B:24:LEU:HD13	2.12	0.50
1:B:2:VAL:HG23	1:B:3:ASP:H	1.76	0.50
1:A:117:LEU:HB3	1:A:129:TYR:CE1	2.47	0.49
1:B:107:PHE:CZ	1:B:164:GLU:HA	2.47	0.49
1:B:43:PHE:CD1	1:B:44:PRO:HD2	2.48	0.49
1:B:16:ALA:HB2	1:B:76:ILE:HD13	1.94	0.49
1:B:90:ILE:HG13	1:B:92:VAL:HG23	1.94	0.49
1:B:17:SER:HB3	1:B:103:GLN:HG3	1.93	0.49
1:A:130:GLU:HG3	1:A:130:GLU:O	2.12	0.48
1:B:169:GLU:HB3	2:B:293:HOH:O	2.13	0.48
1:A:148:ILE:HD11	1:A:170:ALA:C	2.38	0.48
1:B:125:PRO:HD3	1:B:181:TYR:CZ	2.49	0.48
1:B:24:LEU:HD23	1:B:59:ILE:HG22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:PHE:CD2	1:B:138:ILE:HD11	2.48	0.48
1:B:26:LEU:HG	1:B:46:ILE:HG13	1.95	0.48
1:A:88:LYS:HZ1	1:A:93:LYS:HA	1.79	0.48
1:A:135:PRO:O	1:A:149:LEU:HD11	2.15	0.47
1:B:21:PHE:O	1:B:95:LYS:HD3	2.14	0.47
1:A:2:VAL:HG11	1:A:175:LEU:HD13	1.96	0.47
1:A:32:LEU:HG	1:A:86:LYS:HB3	1.96	0.47
1:B:142:LYS:N	1:B:142:LYS:HD2	2.29	0.47
1:A:95:LYS:HA	1:A:95:LYS:CE	2.45	0.47
1:B:143:GLU:O	1:B:173:LYS:HE2	2.16	0.46
1:B:13:ASN:OD1	1:B:15:VAL:HG23	2.15	0.46
1:B:113:ARG:HB3	1:B:113:ARG:CZ	2.45	0.45
1:A:82:LYS:O	1:A:86:LYS:HG3	2.16	0.45
1:A:172:ARG:O	1:A:176:ARG:HB2	2.16	0.45
1:A:32:LEU:HD21	1:A:90:ILE:HD13	1.97	0.45
1:A:134:PHE:CE2	1:B:138:ILE:HD11	2.52	0.45
1:A:149:LEU:O	1:A:156:ILE:HA	2.16	0.45
1:A:29:VAL:HG12	1:A:46:ILE:HD11	1.98	0.45
1:B:103:GLN:HA	1:B:103:GLN:NE2	2.27	0.45
1:A:9:LEU:HD23	1:A:9:LEU:HA	1.72	0.45
1:A:72:SER:HB2	2:A:266:HOH:O	2.16	0.45
1:B:4:MET:SD	1:B:172:ARG:HG2	2.57	0.45
1:B:95:LYS:HG2	1:B:96:ARG:H	1.82	0.45
1:B:95:LYS:H	1:B:95:LYS:CD	2.26	0.45
1:A:11:ILE:HD11	1:A:14:ILE:HG12	1.99	0.44
1:B:113:ARG:NH2	1:B:183:LEU:O	2.50	0.44
1:A:128:GLU:HB2	1:A:138:ILE:HB	1.99	0.44
1:B:9:LEU:HD22	1:B:107:PHE:HB3	1.99	0.44
1:A:55:VAL:HG11	1:A:67:VAL:HG13	1.99	0.44
1:A:107:PHE:CZ	1:A:164:GLU:HA	2.53	0.44
1:B:123:THR:HG22	1:B:181:TYR:HB3	2.00	0.44
1:B:7:VAL:CG2	1:B:175:LEU:HD11	2.44	0.43
1:B:59:ILE:HG23	1:B:65:LEU:CD2	2.47	0.43
1:A:135:PRO:HB3	1:B:159:SER:HB2	2.00	0.43
1:A:183:LEU:HD21	1:B:37:LYS:HZ2	1.84	0.43
1:B:169:GLU:OE1	1:B:172:ARG:NH2	2.50	0.43
1:B:55:VAL:HG22	1:B:68:THR:O	2.18	0.43
1:B:131:PRO:HB3	2:B:316:HOH:O	2.18	0.43
1:A:12:GLU:CG	1:A:108:SER:HB3	2.46	0.43
1:B:52:ASP:HA	1:B:53:PRO:C	2.43	0.43
1:A:134:PHE:HE1	1:A:149:LEU:HD13	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:ILE:HD13	1:B:105:MET:HE2	2.01	0.43
1:B:37:LYS:HB2	1:B:47:ILE:HB	2.01	0.43
1:A:140:ARG:HG2	1:A:140:ARG:HH11	1.84	0.42
1:A:144:PRO:HG3	1:A:166:ASP:CB	2.49	0.42
1:A:52:ASP:HA	1:A:53:PRO:HA	1.88	0.42
1:B:18:VAL:HG21	1:B:80:VAL:HG21	2.02	0.42
1:A:71:LYS:HA	1:A:71:LYS:HD3	1.76	0.42
1:A:57:LEU:HA	1:A:66:VAL:O	2.20	0.42
1:A:30:LEU:HD11	1:A:38:TYR:HB2	2.02	0.41
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.79	0.41
1:B:161:ALA:HB3	2:B:204:HOH:O	2.20	0.41
1:A:23:GLN:OE1	1:A:61:SER:HB2	2.20	0.41
1:A:87:LEU:HG	1:A:92:VAL:HG21	2.03	0.41
1:B:83:LEU:HD13	1:B:87:LEU:CD1	2.48	0.41
1:A:48:CYS:HB2	1:A:59:ILE:CD1	2.51	0.41
1:A:159:SER:OG	1:B:131:PRO:O	2.38	0.41
1:B:149:LEU:HG	1:B:151:PHE:CZ	2.56	0.41
1:B:143:GLU:HA	1:B:144:PRO:HA	1.84	0.41
1:A:124:LEU:HD12	1:A:124:LEU:HA	1.89	0.41
1:A:111:ILE:HB	1:A:175:LEU:HD21	2.01	0.41
1:A:15:VAL:HA	1:A:67:VAL:O	2.20	0.41
1:B:9:LEU:HD23	1:B:9:LEU:HA	1.74	0.41
1:B:13:ASN:HD21	1:B:69:GLY:H	1.67	0.41
1:B:121:ALA:HA	1:B:127:CYS:SG	2.61	0.41
1:A:144:PRO:HD3	1:A:169:GLU:OE1	2.21	0.40
1:B:53:PRO:HG3	1:B:75:ASP:O	2.21	0.40
1:B:88:LYS:HA	1:B:92:VAL:O	2.21	0.40
1:A:20:LEU:HD11	1:A:83:LEU:CD1	2.51	0.40
1:A:30:LEU:HA	1:A:30:LEU:HD23	1.91	0.40
1:B:18:VAL:HG21	1:B:80:VAL:CG2	2.51	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	181/191 (95%)	175 (97%)	6 (3%)	0	100	100
1	B	181/191 (95%)	178 (98%)	3 (2%)	0	100	100
All	All	362/382 (95%)	353 (98%)	9 (2%)	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	160/168 (95%)	143 (89%)	17 (11%)	6	6
1	B	160/168 (95%)	146 (91%)	14 (9%)	9	10
All	All	320/336 (95%)	289 (90%)	31 (10%)	8	8

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	26	LEU
1	A	28	LYS
1	A	41	GLU
1	A	50	LEU
1	A	55	VAL
1	A	78	ARG
1	A	83	LEU
1	A	85	GLN
1	A	87	LEU
1	A	95	LYS
1	A	111	ILE
1	A	117	LEU
1	A	126	ASN
1	A	143	GLU

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Mol	Chain	Res	Type
1	A	150	LEU
1	A	176	ARG
1	B	9	LEU
1	B	10	ARG
1	B	24	LEU
1	B	26	LEU
1	B	55	VAL
1	B	65	LEU
1	B	82	LYS
1	B	83	LEU
1	B	85	GLN
1	B	95	LYS
1	B	96	ARG
1	B	103	GLN
1	B	172	ARG
1	B	174	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	103	GLN
1	B	13	ASN
1	B	49	HIS
1	B	85	GLN
1	B	103	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	183/191 (95%)	-0.58	0 100 100	20, 41, 66, 84	0
1	B	183/191 (95%)	-0.68	0 100 100	21, 38, 65, 81	0
All	All	366/382 (95%)	-0.63	0 100 100	20, 39, 65, 84	0

There are no RSRZ outliers to report.

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.