



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:42 PM UTC

PDB ID : 6A2V / pdb_00006a2v
Title : Crystal structure of Hcp protein
Authors : Jobichen, C.; Sivaraman, J.
Deposited on : 2018-06-13
Resolution : 2.59 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

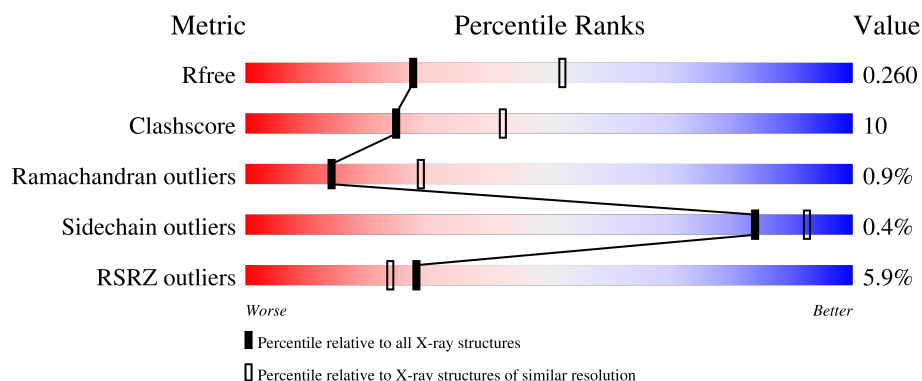
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	171	5% 80% 18% ••
1	B	171	4% 76% 17% • 6%
1	C	171	5% 83% 12% 5%
1	D	171	9% 70% 23% • 5%
1	E	171	5% 70% 16% • 13%

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Mol	Chain	Length	Quality of chain
1	F	171	6% 69% 17% 14%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7183 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 Type VI secretion system tube protein Hcp.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	Se	0	0	0
			1264	792	216	252	1	3			
1	B	160	Total	C	N	O	S	Se	0	0	0
			1218	765	209	240	1	3			
1	C	162	Total	C	N	O	S	Se	0	0	0
			1221	767	211	239	1	3			
1	D	162	Total	C	N	O	S	Se	0	0	0
			1181	743	203	231	1	3			
1	E	148	Total	C	N	O	S	Se	0	0	0
			1099	697	181	217	1	3			
1	F	147	Total	C	N	O	S	Se	0	0	0
			1102	698	187	213	1	3			

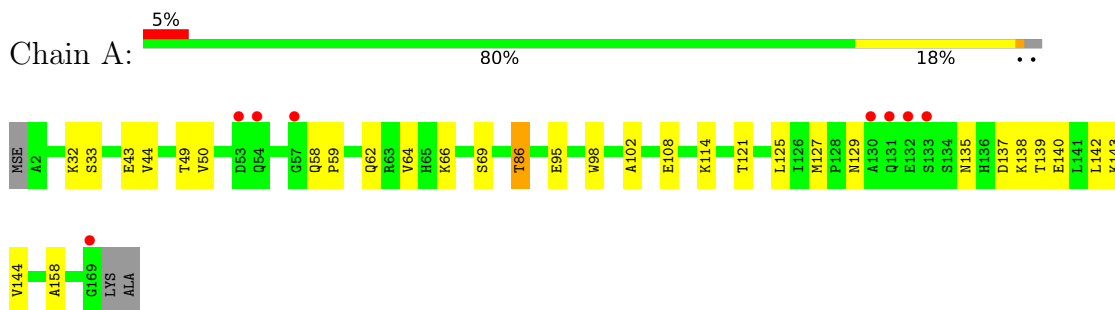
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	33	Total	O	0	0
			33	33		
2	C	18	Total	O	0	0
			18	18		
2	D	11	Total	O	0	0
			11	11		
2	E	6	Total	O	0	0
			6	6		
2	F	11	Total	O	0	0
			11	11		

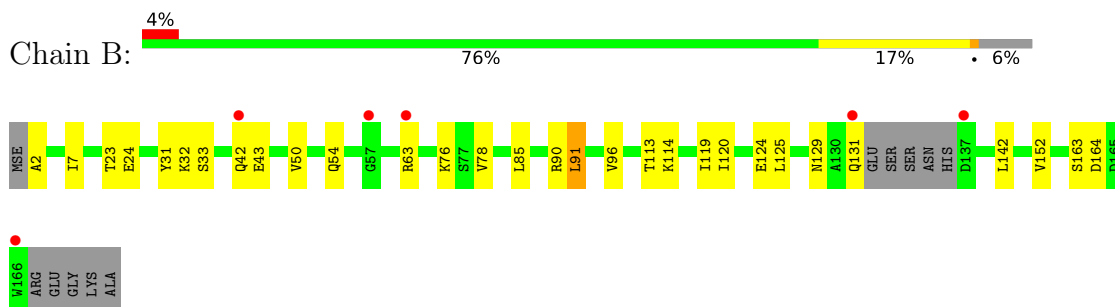
3 Residue-property plots [i](#)

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.

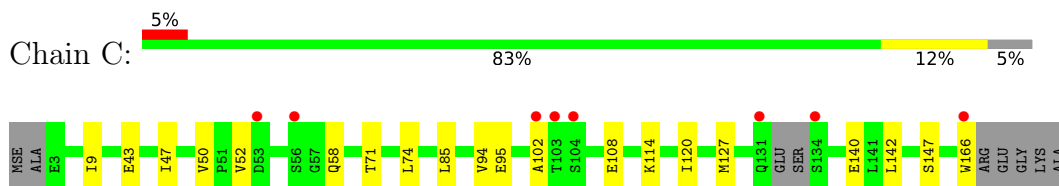
- Molecule 1: Type VI secretion system tube protein Hcp



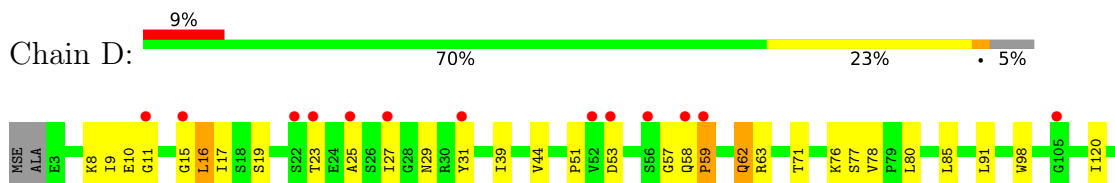
- Molecule 1: Type VI secretion system tube protein Hcp



- Molecule 1: Type VI secretion system tube protein Hcp

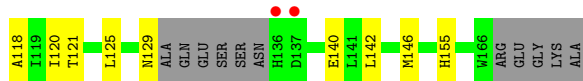
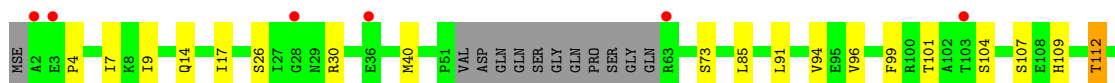


- Molecule 1: Type VI secretion system tube protein Hcp

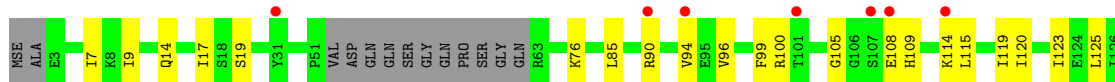




- Molecule 1: Type VI secretion system tube protein Hcp



- Molecule 1: Type VI secretion system tube protein Hcp



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.37Å 91.92Å 165.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.97 – 2.59 19.97 – 2.59	Depositor EDS
% Data completeness (in resolution range)	86.9 (19.97-2.59) 83.1 (19.97-2.59)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.58Å)	Xtriage
Refinement program	PHENIX (dev_2733: ???)	Depositor
R, R_{free}	0.215 , 0.255 0.223 , 0.260	Depositor DCC
R_{free} test set	2074 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	56.1	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.031 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7183	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.40	0/1289	0.66	1/1746 (0.1%)
1	B	0.37	0/1241	0.58	0/1679
1	C	0.36	0/1244	0.67	0/1683
1	D	0.52	0/1201	0.85	3/1628 (0.2%)
1	E	0.34	0/1120	0.67	3/1521 (0.2%)
1	F	0.38	0/1122	0.72	0/1519
All	All	0.40	0/7217	0.70	7/9776 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	51	PRO	N-CA-CB	7.03	109.57	103.31
1	D	59	PRO	N-CA-CB	6.36	109.92	103.25
1	E	107	SER	CA-C-N	-5.63	114.49	122.94
1	E	107	SER	C-N-CA	-5.63	114.49	122.94
1	D	10	GLU	N-CA-C	5.47	117.82	108.90
1	E	112	THR	OG1-CB-CG2	5.23	119.75	109.30
1	A	86	THR	OG1-CB-CG2	5.07	119.43	109.30

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	1264	0	1203	36	0
1	B	1218	0	1177	36	0
1	C	1221	0	1170	12	0
1	D	1181	0	1100	33	0
1	E	1099	0	1029	16	0
1	F	1102	0	1060	27	0
2	A	19	0	0	1	0
2	B	33	0	0	3	0
2	C	18	0	0	0	0
2	D	11	0	0	0	0
2	E	6	0	0	0	0
2	F	11	0	0	0	0
All	All	7183	0	6739	136	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 10.

All (136) 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:125:LEU:HD13	1:A:142:LEU:HD21	1.43	0.96
1:D:62:GLN:HE21	1:D:62:GLN:HA	1.26	0.96
1:F:7:ILE:HD13	1:F:96:VAL:HG22	1.59	0.85
1:A:125:LEU:HD13	1:A:142:LEU:CD2	2.07	0.83
1:E:9:ILE:HG12	1:E:94:VAL:HG13	1.62	0.81
1:F:123:ILE:HG23	1:F:142:LEU:HD11	1.68	0.74
1:D:11:GLY:HA2	1:D:91:LEU:HA	1.70	0.71
1:B:96:VAL:HG22	1:B:113:THR:HB	1.74	0.70
1:F:7:ILE:CD1	1:F:96:VAL:HG22	2.22	0.69
1:D:62:GLN:HA	1:D:62:GLN:NE2	2.04	0.69
1:D:53:ASP:O	1:D:57:GLY:HA2	1.93	0.68
1:B:91:LEU:N	1:B:91:LEU:HD13	2.08	0.68
1:B:90:ARG:C	1:B:91:LEU:HD13	2.18	0.68
1:D:44:VAL:HG11	1:D:98:TRP:CH2	2.31	0.65
1:D:8:LYS:C	1:D:9:ILE:HD12	2.22	0.64
1:B:76:LYS:NZ	2:B:202:HOH:O	2.31	0.64
1:B:85:LEU:HD13	1:B:120:ILE:HG22	1.80	0.63
1:A:86:THR:HA	1:B:63:ARG:HD2	1.81	0.62
1:E:125:LEU:HB2	1:E:142:LEU:HD12	1.82	0.62
1:A:32:LYS:NZ	1:A:33:SER:O	2.25	0.61
1:B:7:ILE:HG12	1:B:96:VAL:HG12	1.83	0.61
1:A:86:THR:HG21	1:B:164:ASP:CG	2.26	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:VAL:HG22	1:D:142:LEU:HD12	1.84	0.60
1:A:125:LEU:HB2	1:A:142:LEU:HD11	1.85	0.58
1:C:102:ALA:HB2	1:C:108:GLU:HB3	1.85	0.58
1:A:158:ALA:HB1	1:F:76:LYS:HD3	1.85	0.58
1:B:114:LYS:HB2	1:B:152:VAL:HG22	1.85	0.58
1:A:86:THR:O	1:B:63:ARG:HD3	2.04	0.57
1:D:78:VAL:CG2	1:D:142:LEU:HD12	2.34	0.57
1:B:119:ILE:HD12	2:B:204:HOH:O	2.04	0.57
1:C:85:LEU:HD13	1:C:120:ILE:HG22	1.86	0.57
1:A:125:LEU:HB2	1:A:142:LEU:CD1	2.34	0.57
1:A:135:ASN:HA	1:A:138:LYS:HG3	1.85	0.57
1:E:91:LEU:HD12	1:E:118:ALA:O	2.06	0.56
1:A:58:GLN:HB3	1:F:90:ARG:HH12	1.70	0.56
1:F:14:GLN:HB2	1:F:17:ILE:HD11	1.88	0.56
1:F:94:VAL:HG22	1:F:115:LEU:HB2	1.87	0.56
1:E:26:SER:C	1:F:100:ARG:HH22	2.13	0.55
1:A:125:LEU:CD1	1:A:142:LEU:CD2	2.83	0.55
1:C:74:LEU:HG	1:C:140:GLU:HG3	1.87	0.55
1:A:135:ASN:HD22	1:A:138:LYS:HE2	1.72	0.55
1:B:78:VAL:HG22	1:B:142:LEU:HD22	1.89	0.54
1:B:129:ASN:ND2	1:C:43:GLU:OE1	2.41	0.54
1:D:62:GLN:HE21	1:D:62:GLN:CA	2.09	0.53
1:E:4:PRO:HG3	1:E:40:MSE:HE3	1.90	0.53
1:A:121:THR:HA	1:B:50:VAL:HG22	1.91	0.53
1:D:9:ILE:HD12	1:D:9:ILE:N	2.25	0.52
1:F:7:ILE:HG22	1:F:9:ILE:HG13	1.91	0.52
1:B:125:LEU:HB2	1:B:142:LEU:HD12	1.92	0.52
1:A:49:THR:HG22	1:A:64:VAL:HG22	1.92	0.51
1:A:139:THR:HG23	1:A:140:GLU:O	2.09	0.51
1:B:152:VAL:HG12	1:B:163:SER:HB2	1.92	0.51
1:D:17:ILE:HG23	1:D:80:LEU:HB3	1.92	0.50
1:D:9:ILE:HD13	1:D:39:ILE:HD12	1.93	0.50
1:F:85:LEU:HD23	1:F:120:ILE:HG22	1.94	0.50
1:A:135:ASN:ND2	1:A:138:LYS:HE2	2.27	0.50
1:B:42:GLN:HA	1:B:42:GLN:OE1	2.12	0.50
1:D:23:THR:O	1:D:27:ILE:HG12	2.11	0.50
1:D:147:SER:HA	1:D:166:TRP:CZ3	2.48	0.49
1:F:114:LYS:CG	1:F:152:VAL:HG22	2.42	0.49
1:E:85:LEU:HD23	1:E:120:ILE:HG22	1.95	0.48
1:F:123:ILE:CG2	1:F:142:LEU:HD11	2.43	0.48
1:A:102:ALA:HB2	1:A:108:GLU:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:62:GLN:NE2	1:D:62:GLN:CA	2.73	0.48
1:E:129:ASN:OD1	1:E:129:ASN:N	2.46	0.48
1:F:114:LYS:HG3	1:F:152:VAL:HG22	1.96	0.48
1:B:114:LYS:HE3	1:B:152:VAL:HG21	1.96	0.48
1:B:32:LYS:HD2	1:B:33:SER:N	2.29	0.48
1:A:137:ASP:OD2	1:A:138:LYS:HG2	2.14	0.47
1:D:19:SER:O	1:D:76:LYS:HE2	2.14	0.47
1:A:95:GLU:HG3	1:A:114:LYS:HG2	1.96	0.47
1:C:52:VAL:HA	1:C:58:GLN:O	2.15	0.47
1:D:29:ASN:HD22	1:E:101:THR:HB	1.80	0.47
1:F:123:ILE:HG23	1:F:142:LEU:CD1	2.39	0.47
1:B:124:GLU:HG3	1:C:47:ILE:HG12	1.96	0.47
1:D:62:GLN:O	1:D:63:ARG:C	2.58	0.47
1:D:15:GLY:O	1:D:17:ILE:N	2.48	0.47
1:F:90:ARG:HG2	1:F:119:ILE:HD13	1.97	0.46
1:B:91:LEU:N	1:B:91:LEU:CD1	2.78	0.46
1:B:24:GLU:N	1:B:31:TYR:HD2	2.13	0.46
1:F:125:LEU:HB2	1:F:142:LEU:HD23	1.98	0.46
1:A:44:VAL:H	1:F:127:MSE:HB3	1.81	0.46
1:A:127:MSE:HE2	1:B:42:GLN:O	2.16	0.46
1:A:86:THR:HA	1:B:63:ARG:CD	2.46	0.45
1:F:125:LEU:HB2	1:F:142:LEU:CD2	2.47	0.45
1:D:71:THR:HA	1:D:142:LEU:O	2.16	0.45
1:E:7:ILE:HG13	1:E:96:VAL:HG13	1.97	0.45
1:F:19:SER:O	1:F:76:LYS:HE3	2.16	0.45
1:A:129:ASN:HB3	1:B:43:GLU:HG3	1.98	0.45
1:F:99:PHE:CE1	1:F:109:HIS:HB2	2.51	0.45
1:F:142:LEU:C	1:F:142:LEU:HD13	2.42	0.45
1:C:95:GLU:HG3	1:C:114:LYS:HG2	1.99	0.45
1:E:99:PHE:CE1	1:E:109:HIS:HB2	2.53	0.44
1:C:147:SER:HB2	1:C:166:TRP:CH2	2.52	0.44
1:D:125:LEU:HB2	1:D:142:LEU:CD2	2.47	0.44
1:D:143:LYS:NZ	1:D:143:LYS:HB3	2.33	0.44
1:A:129:ASN:HA	1:B:42:GLN:O	2.17	0.44
1:B:23:THR:HA	1:B:31:TYR:CE2	2.52	0.44
1:D:124:GLU:OE1	1:D:126:ILE:HD11	2.18	0.44
1:B:96:VAL:CG2	1:B:113:THR:HB	2.44	0.44
1:D:17:ILE:HG22	1:D:77:SER:OG	2.18	0.44
1:C:9:ILE:HG12	1:C:94:VAL:HG13	1.99	0.43
1:E:14:GLN:HB2	1:E:17:ILE:CD1	2.48	0.43
1:A:43:GLU:OE1	1:F:130:ALA:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:THR:HA	1:C:142:LEU:O	2.18	0.43
1:D:63:ARG:HD3	1:D:164:ASP:OD2	2.18	0.43
1:A:129:ASN:HB3	1:B:43:GLU:CG	2.48	0.43
1:D:27:ILE:HD11	1:D:31:TYR:HA	1.99	0.43
1:F:154:GLU:O	1:F:156:THR:N	2.51	0.43
1:A:69:SER:HA	1:A:144:VAL:O	2.19	0.43
1:D:146:MSE:HE3	1:D:146:MSE:HB2	1.88	0.43
1:E:26:SER:O	1:F:100:ARG:NH2	2.51	0.43
1:C:127:MSE:HB3	1:D:44:VAL:HG12	2.00	0.43
1:A:135:ASN:O	1:B:2:ALA:N	2.51	0.42
1:D:16:LEU:HB3	1:D:19:SER:HB2	2.01	0.42
1:A:50:VAL:HG11	1:A:59:PRO:HB3	2.02	0.42
1:F:152:VAL:HG12	1:F:163:SER:HB2	2.01	0.42
1:E:121:THR:HG22	1:E:146:MSE:HA	2.01	0.42
1:A:127:MSE:HE2	1:B:42:GLN:C	2.44	0.42
1:E:30:ARG:NH2	1:E:73:SER:OG	2.53	0.42
1:A:62:GLN:HG2	2:A:210:HOH:O	2.19	0.41
1:F:94:VAL:CG2	1:F:115:LEU:HD12	2.51	0.41
1:A:44:VAL:HG11	1:A:98:TRP:CH2	2.56	0.41
1:D:25:ALA:O	1:E:155:HIS:HE1	2.04	0.41
1:A:125:LEU:CD1	1:A:142:LEU:HD22	2.51	0.41
1:D:85:LEU:HD13	1:D:120:ILE:HG22	2.03	0.41
1:E:96:VAL:O	1:E:112:THR:HA	2.20	0.41
1:F:100:ARG:HD3	1:F:108:GLU:OE2	2.21	0.41
1:B:85:LEU:HD11	1:C:50:VAL:HG23	2.02	0.41
1:B:90:ARG:HD2	1:B:119:ILE:HD13	2.02	0.41
1:A:66:LYS:NZ	1:B:54:GLN:OE1	2.48	0.40
1:A:125:LEU:CD1	1:A:142:LEU:HD21	2.31	0.40
1:B:114:LYS:HE3	1:B:152:VAL:CG2	2.51	0.40
1:D:71:THR:OG1	1:D:143:LYS:HG2	2.21	0.40
1:D:155:HIS:O	1:D:159:GLY:O	2.39	0.40
1:B:131:GLN:O	2:B:201:HOH:O	2.22	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	166/171 (97%)	155 (93%)	11 (7%)	0	100	100
1	B	156/171 (91%)	152 (97%)	4 (3%)	0	100	100
1	C	158/171 (92%)	148 (94%)	10 (6%)	0	100	100
1	D	158/171 (92%)	147 (93%)	7 (4%)	4 (2%)	4	7
1	E	142/171 (83%)	133 (94%)	7 (5%)	2 (1%)	9	18
1	F	141/171 (82%)	131 (93%)	8 (6%)	2 (1%)	9	18
All	All	921/1026 (90%)	866 (94%)	47 (5%)	8 (1%)	14	29

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	59	PRO
1	F	155	HIS
1	E	140	GLU
1	D	16	LEU
1	D	160	THR
1	E	104	SER
1	D	58	GLN
1	F	105	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	136/142 (96%)	135 (99%)	1 (1%)	76	88
1	B	133/142 (94%)	132 (99%)	1 (1%)	73	87
1	C	131/142 (92%)	131 (100%)	0	100	100
1	D	120/142 (84%)	119 (99%)	1 (1%)	73	87
1	E	116/142 (82%)	116 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	118/142 (83%)	118 (100%)	0	100	100
All	All	754/852 (88%)	751 (100%)	3 (0%)	84	92

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	LYS
1	B	91	LEU
1	D	62	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	D	42	GLN
1	D	62	GLN
1	E	42	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 ⓘ

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	165/171 (96%)	0.29	8 (4%)	35	31	43, 55, 99, 113	0
1	B	157/171 (91%)	0.20	6 (3%)	44	39	40, 50, 91, 103	0
1	C	159/171 (92%)	0.33	8 (5%)	34	29	43, 57, 104, 123	1 (0%)
1	D	159/171 (92%)	0.65	15 (9%)	14	11	49, 70, 102, 110	0
1	E	145/171 (84%)	0.72	8 (5%)	30	26	56, 76, 100, 127	0
1	F	144/171 (84%)	0.47	10 (6%)	23	19	44, 61, 92, 121	0
All	All	929/1026 (90%)	0.44	55 (5%)	28	24	40, 62, 100, 127	1 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	131	GLN	4.3
1	A	53	ASP	4.2
1	D	130	ALA	4.2
1	A	133	SER	4.0
1	D	15	GLY	4.0
1	D	56	SER	3.9
1	A	132	GLU	3.8
1	E	137	ASP	3.5
1	A	169	GLY	3.4
1	F	166	TRP	3.4
1	C	134	SER	3.3
1	D	166	TRP	3.3
1	B	57	GLY	3.3
1	D	23	THR	3.2
1	B	166	TRP	3.2
1	A	130	ALA	3.0
1	C	166	TRP	3.0
1	D	25	ALA	2.8
1	A	57	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	137	ASP	2.7
1	F	101	THR	2.6
1	F	94	VAL	2.6
1	A	54	GLN	2.6
1	B	137	ASP	2.6
1	F	90	ARG	2.5
1	B	63	ARG	2.5
1	E	28	GLY	2.5
1	B	42	GLN	2.5
1	A	131	GLN	2.5
1	D	22	SER	2.5
1	E	3	GLU	2.5
1	C	102	ALA	2.5
1	E	136	HIS	2.4
1	E	63	ARG	2.4
1	D	31	TYR	2.4
1	B	131	GLN	2.3
1	D	53	ASP	2.3
1	E	103	THR	2.3
1	D	58	GLN	2.3
1	E	2	ALA	2.3
1	F	114	LYS	2.3
1	D	27	ILE	2.2
1	C	56	SER	2.2
1	E	36	GLU	2.2
1	F	129	ASN	2.2
1	D	59	PRO	2.2
1	D	105	GLY	2.1
1	F	108	GLU	2.1
1	F	31	TYR	2.1
1	C	104	SER	2.1
1	D	52	VAL	2.1
1	C	103	THR	2.1
1	F	107	SER	2.1
1	C	53	ASP	2.0
1	D	11	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.