



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 19, 2022 – 05:33 AM JST

PDB ID : 7FCF
Title : Crystal structure of T6SS Hcp protein
Authors : Jobichen, C.; Sivaraman, J.
Deposited on : 2021-07-14
Resolution : 3.30 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.29
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

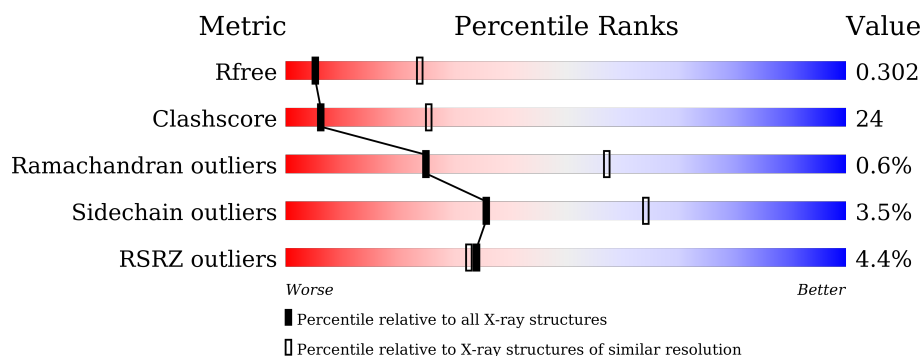
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	154	<div> <div>43%</div> <div>37%</div> <div>19%</div> </div>
2	B	163	<div>7%</div> <div>56%</div> <div>31%</div> <div>12%</div>
2	C	163	<div>2%</div> <div>55%</div> <div>35%</div> <div>10%</div>
2	D	163	<div>6%</div> <div>61%</div> <div>28%</div> <div>11%</div>
2	E	163	<div>4%</div> <div>60%</div> <div>28%</div> <div>11%</div>
2	F	163	<div>3%</div> <div>53%</div> <div>36%</div> <div>10%</div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6448 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.

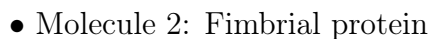
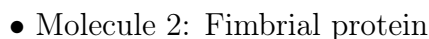
- Molecule 1 is a protein called Fimbrial protein.

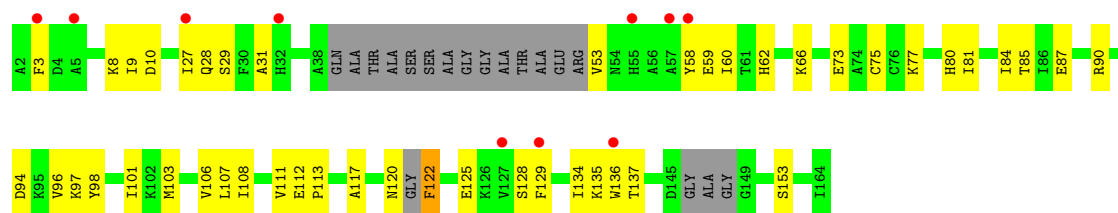
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	124	Total	C	N	O	S	0	0	0
			907	571	154	178	4			

- Molecule 2 is a protein called Fimbrial protein.

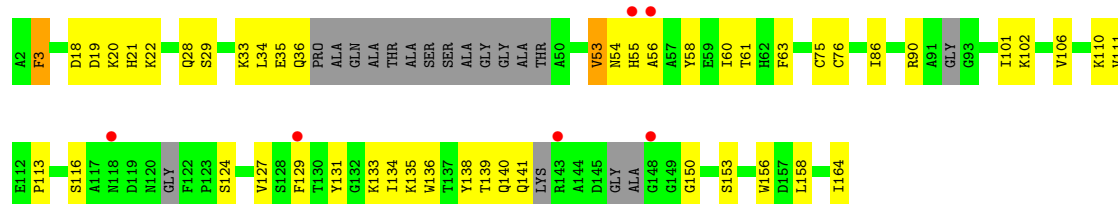
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	143	Total	C	N	O	S	0	0	0
			1109	702	188	214	5			
2	C	147	Total	C	N	O	S	0	0	0
			1117	708	188	216	5			
2	D	145	Total	C	N	O	S	0	0	0
			1075	681	179	210	5			
2	E	145	Total	C	N	O	S	0	0	0
			1121	712	188	216	5			
2	F	146	Total	C	N	O	S	0	0	0
			1119	712	187	215	5			

- Molecule 1: Fimbrial protein

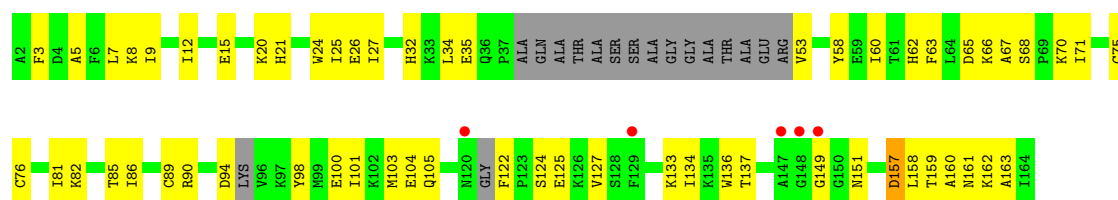




● Molecule 2: Fimbrial protein



● Molecule 2: Fimbrial protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.20Å 83.15Å 156.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.41 – 3.30 46.41 – 3.30	Depositor EDS
% Data completeness (in resolution range)	97.2 (46.41-3.30) 97.2 (46.41-3.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 3.32Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.271 , 0.306 0.270 , 0.302	Depositor DCC
R_{free} test set	759 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	153.6	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 112.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.039 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6448	wwPDB-VP
Average B, all atoms (Å ²)	165.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% 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	0/921	0.74	0/1236
2	B	0.56	0/1131	0.72	0/1520
2	C	0.59	0/1140	0.76	0/1536
2	D	0.56	0/1096	0.76	0/1483
2	E	0.57	0/1142	0.72	0/1534
2	F	0.62	0/1143	0.75	0/1540
All	All	0.57	0/6573	0.74	0/8849

There are no bond length outliers.

There are no bond angle outliers.

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	907	0	846	62	0
2	B	1109	0	1065	64	0
2	C	1117	0	1058	60	0
2	D	1075	0	977	44	0
2	E	1121	0	1058	35	0
2	F	1119	0	1062	65	0
All	All	6448	0	6066	304	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 24.

All (304) 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)
2:B:90:ARG:NH1	2:B:90:ARG:HB2	1.58	1.17
2:F:157:ASP:O	2:F:157:ASP:OD1	1.65	1.14
2:D:9:ILE:HG23	2:D:84:ILE:HG22	1.18	1.11
2:B:134:ILE:HD12	2:B:136:TRP:CH2	1.88	1.08
1:A:111:VAL:CG1	2:B:34:LEU:HD21	1.84	1.06
1:A:89:CYS:HG	1:A:97:LYS:N	1.53	1.04
2:B:101:ILE:HD13	2:B:136:TRP:CD1	1.94	1.02
2:D:9:ILE:HG23	2:D:84:ILE:CG2	1.93	0.97
2:B:101:ILE:CD1	2:B:136:TRP:CD1	2.48	0.97
2:D:87:GLU:HG2	2:D:97:LYS:HE2	1.43	0.96
2:D:90:ARG:HB2	2:D:96:VAL:HG22	1.47	0.94
2:D:90:ARG:HD2	2:D:96:VAL:HG23	1.49	0.92
2:B:134:ILE:HD12	2:B:136:TRP:HH2	1.25	0.91
2:B:35:GLU:HB2	2:B:53:VAL:HG23	1.52	0.91
1:A:111:VAL:HG12	2:B:34:LEU:HD21	1.53	0.89
2:B:101:ILE:CD1	2:B:136:TRP:HD1	1.85	0.89
2:B:7:LEU:HD22	2:B:60:ILE:CD1	2.01	0.89
2:B:74:ALA:HB3	2:B:108:ILE:HD12	1.53	0.88
2:F:157:ASP:CG	2:F:160:ALA:HB3	1.94	0.87
2:B:55:HIS:ND1	2:B:158:LEU:HD13	1.89	0.87
2:B:101:ILE:HD13	2:B:136:TRP:HD1	1.37	0.87
1:A:111:VAL:HG13	2:B:34:LEU:HD21	1.57	0.86
2:F:157:ASP:OD2	2:F:160:ALA:HB3	1.74	0.86
2:B:90:ARG:HB2	2:B:90:ARG:HH11	1.39	0.86
2:B:7:LEU:HD22	2:B:60:ILE:HD13	1.58	0.86
2:F:157:ASP:OD1	2:F:160:ALA:HB3	1.76	0.84
2:C:55:HIS:HE2	2:F:76:CYS:HG	1.21	0.83
2:D:90:ARG:HD2	2:D:96:VAL:CG2	2.07	0.83
2:B:76:CYS:HA	2:E:53:VAL:HG11	1.61	0.82
2:B:90:ARG:HH22	2:B:98:TYR:CB	1.92	0.82
2:B:90:ARG:NH1	2:B:90:ARG:CB	2.43	0.81
2:B:90:ARG:HB2	2:B:90:ARG:CZ	2.07	0.81
2:D:96:VAL:HG23	2:D:96:VAL:O	1.82	0.79
2:C:133:LYS:HZ3	2:C:159:THR:HG21	1.48	0.79
2:E:75:CYS:SG	2:F:34:LEU:HD12	2.23	0.78
2:B:90:ARG:NH2	2:B:98:TYR:CG	2.52	0.78
2:B:9:ILE:HG22	2:B:12:ILE:H	1.49	0.77
2:B:90:ARG:HH11	2:B:90:ARG:CB	1.96	0.77
1:A:111:VAL:HG13	2:B:34:LEU:CD2	2.15	0.76
2:D:90:ARG:CD	2:D:96:VAL:CG2	2.64	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:106:VAL:HG12	2:C:107:LEU:H	1.51	0.74
2:C:30:PHE:HB2	2:C:58:TYR:HE1	1.53	0.74
2:F:157:ASP:OD2	2:F:160:ALA:CB	2.36	0.73
2:D:9:ILE:CG2	2:D:84:ILE:HG22	2.10	0.73
2:F:27:ILE:HD12	2:F:60:ILE:HB	1.70	0.73
2:F:162:LYS:HG3	2:F:163:ALA:H	1.53	0.73
2:C:9:ILE:HA	2:C:84:ILE:HG22	1.71	0.72
2:F:104:GLU:HB2	2:F:133:LYS:HB2	1.72	0.72
2:F:12:ILE:HD13	2:F:70:LYS:HB2	1.70	0.72
2:D:87:GLU:HG2	2:D:97:LYS:CE	2.16	0.71
2:C:8:LYS:HB2	2:C:24:TRP:CH2	2.27	0.70
1:A:89:CYS:SG	1:A:97:LYS:N	2.62	0.70
1:A:111:VAL:CG1	2:B:34:LEU:CD2	2.66	0.70
2:C:76:CYS:HA	2:D:53:VAL:HG21	1.72	0.69
2:E:60:ILE:HD13	2:E:129:PHE:CE2	2.27	0.69
2:C:4:ASP:HB2	2:C:89:CYS:HB2	1.72	0.69
2:D:96:VAL:CG2	2:D:96:VAL:O	2.42	0.68
2:F:9:ILE:HD11	2:F:12:ILE:HB	1.75	0.68
2:B:134:ILE:CD1	2:B:136:TRP:CH2	2.73	0.68
1:A:31:ALA:HB3	1:A:59:GLU:HG2	1.75	0.67
2:C:106:VAL:HG12	2:C:107:LEU:N	2.09	0.67
2:C:133:LYS:HZ3	2:C:159:THR:CG2	2.07	0.67
2:B:110:LYS:HD3	2:E:35:GLU:HG3	1.77	0.67
2:B:134:ILE:CD1	2:B:136:TRP:HH2	2.05	0.67
1:A:29:SER:HB3	1:A:61:THR:HB	1.75	0.66
2:C:55:HIS:HE1	2:C:156:TRP:CD1	2.13	0.66
2:E:86:ILE:HB	2:E:101:ILE:HB	1.78	0.66
1:A:89:CYS:SG	1:A:97:LYS:CA	2.84	0.66
2:C:133:LYS:NZ	2:C:159:THR:HG21	2.10	0.66
1:A:103:MET:HG2	1:A:134:ILE:HG13	1.76	0.65
2:F:8:LYS:NZ	2:F:24:TRP:HA	2.11	0.65
2:B:90:ARG:CG	2:B:91:ALA:H	2.10	0.65
2:C:12:ILE:HG23	2:C:70:LYS:HB2	1.77	0.65
1:A:89:CYS:SG	1:A:97:LYS:HA	2.37	0.65
1:A:114:HIS:NE2	1:A:124:SER:OG	2.25	0.65
1:A:152:VAL:HG11	2:D:66:LYS:HA	1.77	0.65
2:B:90:ARG:HH22	2:B:98:TYR:HB3	1.61	0.64
2:C:10:ASP:HB2	2:C:81:ILE:HD11	1.78	0.64
2:B:90:ARG:HG2	2:B:91:ALA:H	1.61	0.64
1:A:76:CYS:HG	2:B:55:HIS:HE2	1.41	0.64
2:D:81:ILE:HG22	2:D:106:VAL:HB	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:133:LYS:NZ	2:C:159:THR:CG2	2.60	0.64
2:C:55:HIS:NE2	2:F:76:CYS:SG	2.59	0.64
2:C:104:GLU:HB2	2:C:133:LYS:HB2	1.79	0.63
2:C:109:ALA:HB2	2:C:130:THR:HG23	1.79	0.63
2:C:142:LYS:HD2	2:C:144:ALA:HB3	1.79	0.63
2:F:21:HIS:HB3	2:F:24:TRP:CG	2.33	0.63
1:A:21:HIS:NE2	1:A:87:GLU:OE1	2.30	0.63
2:D:103:MET:HG2	2:D:134:ILE:HD12	1.80	0.62
2:B:90:ARG:NH2	2:B:98:TYR:CD1	2.67	0.62
1:A:85:THR:HG22	1:A:102:LYS:HG2	1.80	0.61
1:A:135:LYS:HE3	1:A:153:SER:HB3	1.81	0.61
2:D:75:CYS:SG	2:D:111:VAL:HG12	2.40	0.61
2:F:101:ILE:HD12	2:F:136:TRP:HD1	1.65	0.61
2:C:94:ASP:N	2:C:94:ASP:OD1	2.33	0.61
2:F:157:ASP:OD1	2:F:160:ALA:N	2.33	0.61
2:F:157:ASP:OD1	2:F:160:ALA:CB	2.48	0.61
2:C:8:LYS:HB2	2:C:24:TRP:CZ3	2.37	0.60
2:F:67:ALA:O	2:F:71:ILE:HG13	2.01	0.60
2:C:30:PHE:HB2	2:C:58:TYR:CE1	2.35	0.60
2:C:36:GLN:NE2	2:F:75:CYS:O	2.34	0.60
2:B:15:GLU:HB3	2:B:66:LYS:HG2	1.84	0.60
2:C:36:GLN:OE1	2:C:36:GLN:HA	2.02	0.60
2:F:12:ILE:HD11	2:F:70:LYS:O	2.02	0.60
2:E:76:CYS:HA	2:F:53:VAL:HG21	1.83	0.59
1:A:66:LYS:HG3	2:B:152:VAL:HG21	1.84	0.59
2:F:12:ILE:CD1	2:F:70:LYS:HB2	2.33	0.59
2:E:60:ILE:CD1	2:E:129:PHE:HD2	2.16	0.59
2:F:62:HIS:NE2	2:F:125:GLU:OE1	2.36	0.58
2:F:90:ARG:HG3	2:F:98:TYR:CD1	2.37	0.58
2:E:60:ILE:HD13	2:E:129:PHE:CD2	2.38	0.58
1:A:108:ILE:HG12	1:A:129:PHE:CE1	2.39	0.58
1:A:21:HIS:HB3	1:A:24:TRP:CG	2.39	0.58
1:A:7:LEU:HD11	1:A:84:ILE:HG23	1.85	0.57
2:E:75:CYS:SG	2:F:34:LEU:CD1	2.92	0.57
2:E:60:ILE:CD1	2:E:129:PHE:CD2	2.87	0.57
2:F:68:SER:HA	2:F:71:ILE:HD12	1.85	0.56
2:D:73:GLU:O	2:D:77:LYS:N	2.28	0.56
1:A:84:ILE:HD12	1:A:106:VAL:HG11	1.88	0.56
2:B:84:ILE:HG12	2:B:106:VAL:HG21	1.87	0.56
2:F:103:MET:HG2	2:F:134:ILE:HG13	1.87	0.56
2:C:89:CYS:HB3	2:C:94:ASP:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:60:ILE:HD13	2:E:129:PHE:HE2	1.71	0.55
1:A:82:LYS:HE2	1:A:82:LYS:H	1.71	0.55
1:A:9:ILE:HG12	1:A:84:ILE:HG12	1.89	0.55
2:C:107:LEU:HD23	2:C:108:ILE:N	2.22	0.54
2:F:157:ASP:CG	2:F:160:ALA:CB	2.73	0.54
2:E:19:ASP:OD1	2:E:20:LYS:N	2.38	0.54
2:E:113:PRO:HG2	2:F:32:HIS:HB3	1.87	0.54
2:F:62:HIS:CD2	2:F:125:GLU:HG3	2.42	0.54
2:C:30:PHE:CD1	2:C:58:TYR:OH	2.59	0.54
2:D:58:TYR:O	2:D:128:SER:HA	2.08	0.54
2:C:9:ILE:HG22	2:C:12:ILE:H	1.73	0.53
2:C:27:ILE:HG12	2:C:60:ILE:HD13	1.91	0.53
1:A:62:HIS:ND1	1:A:63:PHE:O	2.42	0.53
2:B:156:TRP:NE1	2:B:161:ASN:HA	2.24	0.52
2:F:8:LYS:HB2	2:F:24:TRP:CD2	2.44	0.52
2:E:101:ILE:HD12	2:E:136:TRP:HD1	1.74	0.52
2:F:27:ILE:CD1	2:F:60:ILE:HB	2.39	0.52
2:F:8:LYS:HZ3	2:F:24:TRP:HA	1.73	0.52
1:A:22:LYS:O	1:A:24:TRP:HD1	1.92	0.52
2:D:101:ILE:HD12	2:D:136:TRP:HD1	1.73	0.52
2:E:33:LYS:NZ	2:E:56:ALA:HB2	2.24	0.52
2:C:107:LEU:HD23	2:C:107:LEU:C	2.30	0.52
2:E:34:LEU:HD22	2:E:55:HIS:ND1	2.25	0.52
2:D:134:ILE:HG21	2:D:136:TRP:CH2	2.45	0.52
2:C:36:GLN:HA	2:C:53:VAL:HA	1.91	0.52
2:C:53:VAL:HG21	2:F:76:CYS:SG	2.50	0.52
2:D:31:ALA:O	2:D:59:GLU:OE1	2.28	0.52
2:F:12:ILE:CD1	2:F:70:LYS:CB	2.88	0.52
2:B:90:ARG:NH1	2:B:96:VAL:HG12	2.24	0.52
2:F:12:ILE:HD11	2:F:70:LYS:C	2.30	0.52
2:B:106:VAL:HG12	2:B:107:LEU:N	2.25	0.51
2:C:106:VAL:CG1	2:C:107:LEU:H	2.21	0.51
2:F:70:LYS:HZ3	2:F:70:LYS:HB3	1.74	0.51
2:C:16:SER:HB3	2:C:24:TRP:H	1.75	0.51
2:D:90:ARG:HD3	2:D:96:VAL:HG21	1.92	0.51
2:B:89:CYS:SG	2:B:97:LYS:HG3	2.50	0.51
2:E:35:GLU:HB3	2:E:54:ASN:HB2	1.92	0.51
2:F:15:GLU:HB3	2:F:66:LYS:HB2	1.93	0.51
2:B:63:PHE:HB3	2:B:122:PHE:HD2	1.76	0.51
2:E:111:VAL:HG23	2:E:127:VAL:HG22	1.92	0.51
2:B:8:LYS:HB3	2:B:85:THR:HG22	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:160:ALA:O	2:B:161:ASN:C	2.50	0.50
2:C:21:HIS:HB3	2:C:24:TRP:CG	2.46	0.50
2:F:7:LEU:HD13	2:F:86:ILE:HG22	1.94	0.50
2:B:24:TRP:CH2	2:B:85:THR:HG23	2.46	0.50
2:B:34:LEU:HB2	2:B:55:HIS:HD2	1.75	0.50
2:B:107:LEU:HB3	2:B:130:THR:HG23	1.93	0.50
2:F:90:ARG:HG3	2:F:98:TYR:CE1	2.46	0.50
1:A:27:ILE:CD1	1:A:86:ILE:HD11	2.42	0.50
2:D:90:ARG:HD3	2:D:96:VAL:CG2	2.41	0.50
2:B:74:ALA:HB3	2:B:108:ILE:CD1	2.34	0.49
2:F:94:ASP:N	2:F:94:ASP:OD1	2.45	0.49
2:E:138:TYR:HE2	2:E:140:GLN:HB2	1.76	0.49
2:B:104:GLU:HB2	2:B:133:LYS:HB2	1.94	0.49
2:F:12:ILE:CD1	2:F:70:LYS:C	2.81	0.49
2:C:12:ILE:HD12	2:C:70:LYS:HB3	1.94	0.49
2:E:156:TRP:HD1	2:E:158:LEU:HB3	1.78	0.49
2:D:134:ILE:HG12	2:D:136:TRP:CZ2	2.47	0.49
2:F:85:THR:HA	2:F:101:ILE:O	2.13	0.49
2:D:8:LYS:HE3	2:D:85:THR:HG22	1.95	0.49
1:A:71:ILE:HG23	1:A:108:ILE:HD13	1.95	0.49
2:D:120:ASN:O	2:D:122:PHE:N	2.46	0.49
2:C:7:LEU:HD12	2:C:8:LYS:N	2.28	0.48
1:A:76:CYS:SG	2:B:55:HIS:NE2	2.75	0.48
2:B:34:LEU:H	2:B:34:LEU:HD23	1.78	0.48
2:C:62:HIS:ND1	2:C:63:PHE:O	2.46	0.48
2:D:90:ARG:HB2	2:D:96:VAL:CG2	2.30	0.48
2:F:162:LYS:HG3	2:F:163:ALA:N	2.26	0.48
1:A:27:ILE:HD13	1:A:86:ILE:HD11	1.94	0.48
2:B:60:ILE:HG12	2:B:127:VAL:HG13	1.95	0.48
2:E:63:PHE:HD1	2:E:124:SER:HB2	1.79	0.48
2:F:149:GLY:O	2:F:151:ASN:N	2.47	0.48
2:B:21:HIS:HB3	2:B:24:TRP:CG	2.48	0.48
2:F:157:ASP:OD1	2:F:160:ALA:CA	2.62	0.48
2:C:22:LYS:O	2:C:24:TRP:HD1	1.97	0.48
2:D:28:GLN:OE1	2:D:62:HIS:HA	2.14	0.47
2:B:90:ARG:CG	2:B:91:ALA:N	2.75	0.47
2:C:100:GLU:HB3	2:C:137:THR:HB	1.96	0.47
2:D:134:ILE:HG13	2:D:135:LYS:N	2.29	0.47
2:C:142:LYS:HB2	2:C:148:GLY:H	1.80	0.47
2:E:28:GLN:NE2	2:E:63:PHE:CD2	2.82	0.47
1:A:134:ILE:HG12	1:A:135:LYS:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:90:ARG:HH22	2:C:140:GLN:HE22	1.63	0.47
2:F:100:GLU:HB3	2:F:137:THR:HB	1.96	0.47
1:A:82:LYS:H	1:A:82:LYS:CE	2.27	0.47
2:C:133:LYS:HG3	2:C:157:ASP:HB2	1.96	0.47
2:E:110:LYS:HE2	2:F:35:GLU:HG3	1.97	0.47
1:A:106:VAL:HA	1:A:130:THR:O	2.15	0.47
2:C:34:LEU:HD12	2:C:34:LEU:H	1.80	0.46
1:A:113:PRO:HA	1:A:125:GLU:HB3	1.98	0.46
2:C:107:LEU:HB3	2:C:130:THR:OG1	2.16	0.46
2:F:159:THR:HG23	2:F:160:ALA:N	2.31	0.46
2:D:80:HIS:HA	2:D:107:LEU:HD23	1.96	0.46
2:F:82:LYS:O	2:F:104:GLU:HA	2.15	0.46
1:A:115:GLY:HA2	1:A:119:ASP:OD2	2.16	0.46
2:D:10:ASP:OD1	2:D:81:ILE:HG13	2.16	0.46
2:D:137:THR:HA	2:D:153:SER:OG	2.16	0.46
2:F:5:ALA:HB3	2:F:27:ILE:HG23	1.98	0.46
1:A:60:ILE:O	1:A:126:LYS:HA	2.15	0.46
1:A:76:CYS:SG	2:B:34:LEU:HD12	2.56	0.46
2:F:8:LYS:HB2	2:F:24:TRP:CE2	2.51	0.45
1:A:68:SER:HB3	1:A:72:TYR:OH	2.15	0.45
2:B:96:VAL:HG12	2:B:96:VAL:O	2.17	0.45
2:C:7:LEU:HD13	2:C:86:ILE:HG22	1.97	0.45
2:D:134:ILE:HG21	2:D:136:TRP:CZ2	2.51	0.45
2:E:116:SER:H	2:F:3:PHE:HE2	1.64	0.45
2:C:81:ILE:HD12	2:C:82:LYS:H	1.81	0.45
2:D:90:ARG:HG3	2:D:98:TYR:CD1	2.51	0.45
2:E:131:TYR:CE2	2:E:134:ILE:HB	2.52	0.45
2:B:160:ALA:O	2:B:162:LYS:N	2.50	0.45
2:F:63:PHE:HA	2:F:124:SER:HA	1.97	0.45
1:A:27:ILE:CD1	1:A:60:ILE:HB	2.46	0.45
1:A:108:ILE:HG12	1:A:129:PHE:CD1	2.51	0.45
2:D:29:SER:O	2:D:60:ILE:HA	2.16	0.45
2:D:85:THR:HA	2:D:101:ILE:O	2.16	0.45
2:B:161:ASN:OD1	2:B:161:ASN:O	2.35	0.44
1:A:135:LYS:HD2	1:A:154:SER:O	2.18	0.44
2:E:29:SER:HB2	2:E:61:THR:OG1	2.16	0.44
2:F:7:LEU:O	2:F:25:ILE:HB	2.18	0.44
1:A:110:LYS:HG2	1:A:112:GLU:OE1	2.17	0.44
2:E:35:GLU:HG2	2:E:36:GLN:H	1.82	0.44
1:A:114:HIS:CD2	1:A:124:SER:OG	2.71	0.44
2:B:90:ARG:HB3	2:B:96:VAL:HB	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:106:VAL:HA	2:E:131:TYR:HA	1.99	0.44
2:B:82:LYS:HA	2:B:82:LYS:HD2	1.87	0.44
2:E:19:ASP:OD1	2:E:20:LYS:HG3	2.18	0.44
1:A:9:ILE:HG22	1:A:12:ILE:H	1.82	0.44
2:C:14:GLY:HA3	2:C:24:TRP:O	2.18	0.44
1:A:135:LYS:HA	1:A:154:SER:O	2.18	0.43
2:F:20:LYS:HA	2:F:20:LYS:HD3	1.78	0.43
2:F:60:ILE:HG12	2:F:127:VAL:HG13	2.00	0.43
1:A:93:GLY:O	1:A:94:ASP:C	2.57	0.43
2:D:27:ILE:HG13	2:D:60:ILE:HB	2.01	0.43
2:C:82:LYS:HD3	2:C:82:LYS:N	2.33	0.43
2:E:102:LYS:HB3	2:E:102:LYS:HE2	1.75	0.43
2:F:157:ASP:OD1	2:F:157:ASP:C	2.45	0.43
1:A:87:GLU:HG3	1:A:100:GLU:HG3	2.01	0.43
2:E:139:THR:HA	2:E:150:GLY:O	2.19	0.43
2:F:81:ILE:HD13	2:F:81:ILE:HA	1.76	0.43
2:B:5:ALA:HB2	2:B:88:LEU:HD22	2.00	0.43
2:C:110:LYS:HZ3	2:C:112:GLU:HG3	1.84	0.43
2:D:101:ILE:HA	2:D:135:LYS:O	2.19	0.43
2:C:156:TRP:HA	2:C:163:ALA:HA	2.00	0.42
2:F:82:LYS:O	2:F:105:GLN:N	2.50	0.42
1:A:29:SER:HA	2:D:117:ALA:CB	2.49	0.42
2:E:110:LYS:HE2	2:E:110:LYS:HB2	1.71	0.42
2:C:8:LYS:HB3	2:C:85:THR:HG23	2.01	0.42
2:C:82:LYS:HB2	2:C:82:LYS:HE2	1.64	0.42
2:C:142:LYS:HE3	2:C:148:GLY:H	1.84	0.42
2:F:158:LEU:HD12	2:F:158:LEU:HA	1.85	0.42
1:A:59:GLU:HA	1:A:128:SER:HA	2.01	0.42
1:A:112:GLU:O	1:A:125:GLU:HA	2.19	0.42
1:A:7:LEU:HD23	1:A:25:ILE:HG21	2.01	0.42
2:B:156:TRP:HE1	2:B:161:ASN:HA	1.85	0.42
2:C:69:PRO:HB3	2:D:136:TRP:HB2	2.02	0.42
2:E:133:LYS:HG2	2:E:164:ILE:HD11	2.00	0.42
2:C:67:ALA:O	2:C:71:ILE:HG13	2.20	0.42
2:B:99:MET:HA	2:B:137:THR:O	2.20	0.42
1:A:9:ILE:HG21	1:A:12:ILE:HD12	2.01	0.41
2:C:86:ILE:HD11	2:C:101:ILE:HD12	2.01	0.41
2:E:18:ASP:HB3	2:E:21:HIS:HB2	2.02	0.41
2:D:27:ILE:HG21	2:D:27:ILE:HD13	1.86	0.41
2:B:90:ARG:HH12	2:B:96:VAL:HG12	1.84	0.41
2:C:133:LYS:HD2	2:C:157:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:LEU:HD12	1:A:85:THR:O	2.21	0.41
2:C:35:GLU:O	2:C:53:VAL:HG23	2.20	0.41
2:F:15:GLU:HG2	2:F:65:ASP:HB2	2.01	0.41
1:A:7:LEU:O	1:A:25:ILE:HG22	2.20	0.41
1:A:7:LEU:CD1	1:A:84:ILE:HG23	2.51	0.41
1:A:133:LYS:HE3	1:A:133:LYS:HB2	1.79	0.41
1:A:108:ILE:HG12	1:A:129:PHE:HE1	1.85	0.41
1:A:111:VAL:HG12	2:B:34:LEU:CD2	2.37	0.41
2:D:66:LYS:HE3	2:D:66:LYS:HB2	1.93	0.41
2:D:108:ILE:HG13	2:D:129:PHE:CE1	2.56	0.41
2:C:162:LYS:HD2	2:C:162:LYS:HA	1.37	0.41
2:E:3:PHE:C	2:E:3:PHE:CD1	2.94	0.41
2:F:12:ILE:HG12	2:F:70:LYS:NZ	2.36	0.41
1:A:21:HIS:HB3	1:A:24:TRP:CD2	2.56	0.40
2:F:89:CYS:HB3	2:F:94:ASP:HA	2.02	0.40
1:A:86:ILE:HG22	1:A:101:ILE:HB	2.04	0.40
2:D:113:PRO:HA	2:D:125:GLU:HB3	2.03	0.40
1:A:64:LEU:HD13	1:A:123:PRO:HB2	2.03	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	114/154 (74%)	104 (91%)	8 (7%)	2 (2%)	8	35
2	B	135/163 (83%)	112 (83%)	21 (16%)	2 (2%)	10	38
2	C	141/163 (86%)	123 (87%)	18 (13%)	0	100	100
2	D	137/163 (84%)	122 (89%)	15 (11%)	0	100	100
2	E	133/163 (82%)	116 (87%)	16 (12%)	1 (1%)	19	51
2	F	138/163 (85%)	128 (93%)	10 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	798/969 (82%)	705 (88%)	88 (11%)	5 (1%)	25 57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	161	ASN
2	E	22	LYS
1	A	56	ALA
1	A	118	ASN
2	B	125	GLU

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	89/121 (74%)	88 (99%)	1 (1%)	73 85
2	B	115/128 (90%)	112 (97%)	3 (3%)	46 71
2	C	112/128 (88%)	109 (97%)	3 (3%)	44 71
2	D	104/128 (81%)	100 (96%)	4 (4%)	33 62
2	E	114/128 (89%)	107 (94%)	7 (6%)	18 48
2	F	114/128 (89%)	109 (96%)	5 (4%)	28 59
All	All	648/761 (85%)	625 (96%)	23 (4%)	36 64

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

Mol	Chain	Res	Type
1	A	102	LYS
2	B	23	ASP
2	B	54	ASN
2	B	58	TYR
2	C	3	PHE
2	C	58	TYR
2	C	80	HIS

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Mol	Chain	Res	Type
2	D	3	PHE
2	D	94	ASP
2	D	112	GLU
2	D	122	PHE
2	E	3	PHE
2	E	53	VAL
2	E	58	TYR
2	E	90	ARG
2	E	135	LYS
2	E	141	GLN
2	E	153	SER
2	F	26	GLU
2	F	58	TYR
2	F	122	PHE
2	F	157	ASP
2	F	161	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides 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	124/154 (80%)	-0.15	1 (0%) 86 86	136, 173, 214, 224	0
2	B	143/163 (87%)	0.44	11 (7%) 13 12	123, 160, 205, 243	0
2	C	147/163 (90%)	0.12	4 (2%) 54 52	124, 167, 218, 226	0
2	D	145/163 (88%)	0.08	10 (6%) 16 16	134, 171, 214, 257	0
2	E	145/163 (88%)	0.08	6 (4%) 37 35	114, 156, 201, 226	0
2	F	146/163 (89%)	0.11	5 (3%) 45 43	116, 163, 211, 250	0
All	All	850/969 (87%)	0.12	37 (4%) 34 33	114, 165, 212, 257	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	164	ILE	6.1
2	D	129	PHE	3.8
2	E	129	PHE	3.4
2	D	136	TRP	3.3
2	B	163	ALA	3.2
2	F	147	ALA	3.2
2	C	55	HIS	3.1
2	B	148	GLY	3.1
2	F	148	GLY	3.1
2	C	119	ASP	3.1
1	A	2	ALA	3.0
2	B	140	GLN	3.0
2	D	127	VAL	2.9
2	D	3	PHE	2.7
2	B	55	HIS	2.6
2	E	118	ASN	2.6
2	B	99	MET	2.5
2	E	148	GLY	2.4
2	F	120	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
2	F	149	GLY	2.4
2	D	27	ILE	2.4
2	F	129	PHE	2.3
2	B	154	SER	2.3
2	B	58	TYR	2.2
2	D	55	HIS	2.2
2	B	103	MET	2.2
2	D	5	ALA	2.2
2	C	12	ILE	2.2
2	B	149	GLY	2.1
2	E	143	ARG	2.1
2	D	32	HIS	2.1
2	B	145	ASP	2.1
2	E	56	ALA	2.1
2	C	129	PHE	2.1
2	E	55	HIS	2.1
2	D	57	ALA	2.0
2	D	58	TYR	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 monosaccharides 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.