



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:07 PM GMT

PDB ID : 3SQB
Title : Structure of the major type 1 pilus subunit FimA bound to the FimC chaperone
Authors : Scharer, M.A.; Eidam, O.; Grutter, M.G.; Glockshuber, R.; Capitani, G.
Deposited on : 2011-07-05
Resolution : 3.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

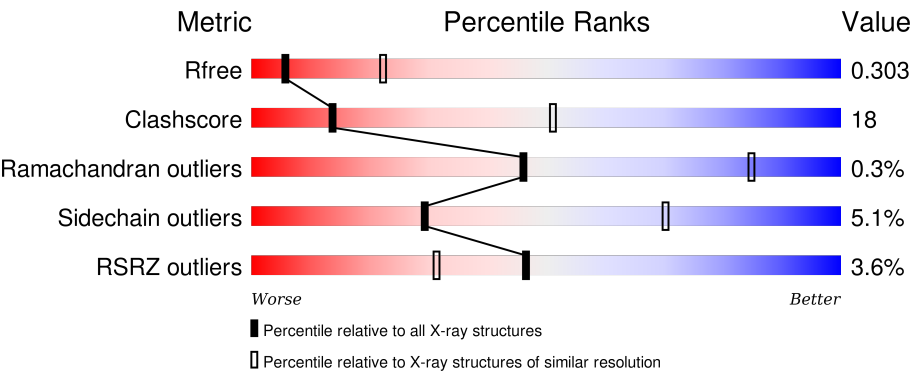
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	211	<div><div></div><div><div></div><div>63%</div><div>27%</div><div>• 9%</div></div></div>
1	C	211	<div><div></div><div><div></div><div>61%</div><div>32%</div><div>• 5%</div></div></div>
1	E	211	<div><div>5%</div><div></div><div><div></div><div>55%</div><div>28%</div><div>• 17%</div></div></div>
1	G	211	<div><div>4%</div><div></div><div><div></div><div>30%</div><div>6%</div><div>64%</div></div></div>
2	B	152	<div><div></div><div><div></div><div>69%</div><div>24%</div><div>• 5%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	D	152	
2	F	152	
2	H	152	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PEG	A	212	-	-	-	X
5	TRS	C	213	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7840 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 Chaperone protein fimC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1447	918	246	278	5			
1	C	200	Total	C	N	O	S	0	0	0
			1543	976	268	294	5			
1	E	176	Total	C	N	O	S	0	0	0
			1198	756	201	236	5			
1	G	76	Total	C	N	O		0	0	0
			485	302	84	99				

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	206	HIS	-	EXPRESSION TAG	UNP P31697
A	207	HIS	-	EXPRESSION TAG	UNP P31697
A	208	HIS	-	EXPRESSION TAG	UNP P31697
A	209	HIS	-	EXPRESSION TAG	UNP P31697
A	210	HIS	-	EXPRESSION TAG	UNP P31697
A	211	HIS	-	EXPRESSION TAG	UNP P31697
C	206	HIS	-	EXPRESSION TAG	UNP P31697
C	207	HIS	-	EXPRESSION TAG	UNP P31697
C	208	HIS	-	EXPRESSION TAG	UNP P31697
C	209	HIS	-	EXPRESSION TAG	UNP P31697
C	210	HIS	-	EXPRESSION TAG	UNP P31697
C	211	HIS	-	EXPRESSION TAG	UNP P31697
E	206	HIS	-	EXPRESSION TAG	UNP P31697
E	207	HIS	-	EXPRESSION TAG	UNP P31697
E	208	HIS	-	EXPRESSION TAG	UNP P31697
E	209	HIS	-	EXPRESSION TAG	UNP P31697
E	210	HIS	-	EXPRESSION TAG	UNP P31697
E	211	HIS	-	EXPRESSION TAG	UNP P31697
G	206	HIS	-	EXPRESSION TAG	UNP P31697
G	207	HIS	-	EXPRESSION TAG	UNP P31697
G	208	HIS	-	EXPRESSION TAG	UNP P31697

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Chain	Residue	Modelled	Actual	Comment	Reference
G	209	HIS	-	EXPRESSION TAG	UNP P31697
G	210	HIS	-	EXPRESSION TAG	UNP P31697
G	211	HIS	-	EXPRESSION TAG	UNP P31697

- Molecule 2 is a protein called Type-1 fimbrial protein, A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	145	Total	C	N	O	S	0	0	0
			1014	619	178	215	2			
2	D	145	Total	C	N	O	S	0	0	0
			1002	615	171	214	2			
2	F	89	Total	C	N	O	S	0	0	0
			592	365	100	125	2			
2	H	78	Total	C	N	O	S	0	0	0
			466	278	84	102	2			

There are 24 discrepancies between the modelled and reference sequences:

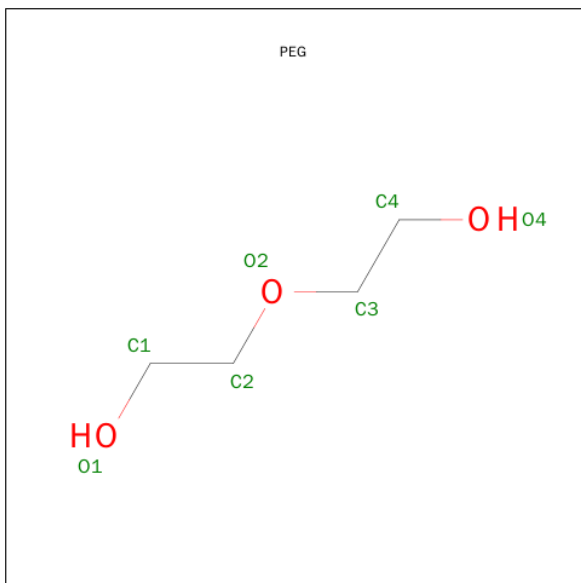
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	HIS	-	EXPRESSION TAG	UNP P04128
B	2	HIS	-	EXPRESSION TAG	UNP P04128
B	3	HIS	-	EXPRESSION TAG	UNP P04128
B	4	HIS	-	EXPRESSION TAG	UNP P04128
B	5	HIS	-	EXPRESSION TAG	UNP P04128
B	6	HIS	-	EXPRESSION TAG	UNP P04128
D	1	HIS	-	EXPRESSION TAG	UNP P04128
D	2	HIS	-	EXPRESSION TAG	UNP P04128
D	3	HIS	-	EXPRESSION TAG	UNP P04128
D	4	HIS	-	EXPRESSION TAG	UNP P04128
D	5	HIS	-	EXPRESSION TAG	UNP P04128
D	6	HIS	-	EXPRESSION TAG	UNP P04128
F	1	HIS	-	EXPRESSION TAG	UNP P04128
F	2	HIS	-	EXPRESSION TAG	UNP P04128
F	3	HIS	-	EXPRESSION TAG	UNP P04128
F	4	HIS	-	EXPRESSION TAG	UNP P04128
F	5	HIS	-	EXPRESSION TAG	UNP P04128
F	6	HIS	-	EXPRESSION TAG	UNP P04128
H	1	HIS	-	EXPRESSION TAG	UNP P04128
H	2	HIS	-	EXPRESSION TAG	UNP P04128
H	3	HIS	-	EXPRESSION TAG	UNP P04128
H	4	HIS	-	EXPRESSION TAG	UNP P04128
H	5	HIS	-	EXPRESSION TAG	UNP P04128

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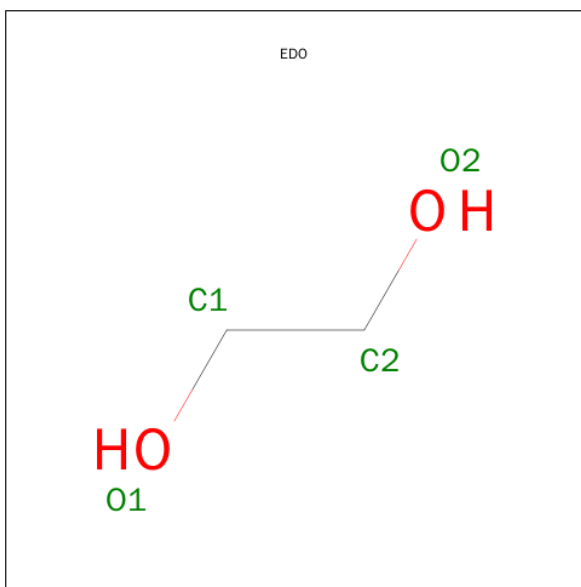
Chain	Residue	Modelled	Actual	Comment	Reference
H	6	HIS	-	EXPRESSION TAG	UNP P04128

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



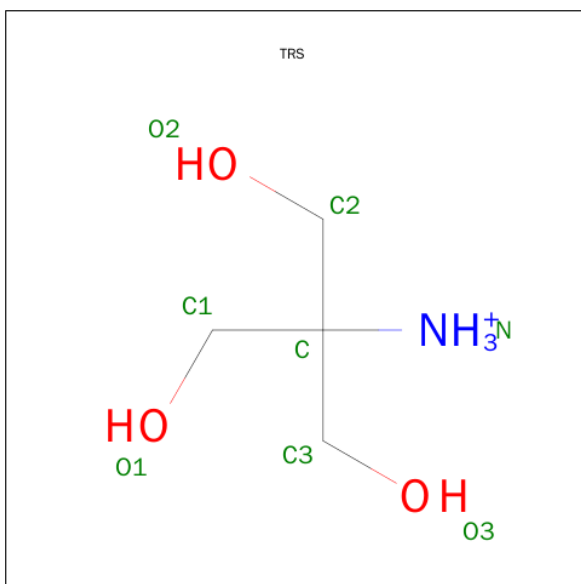
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	N	O	0	0
			8	4	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	D	1	Total	C	N	O	0	0
			8	4	1	3		

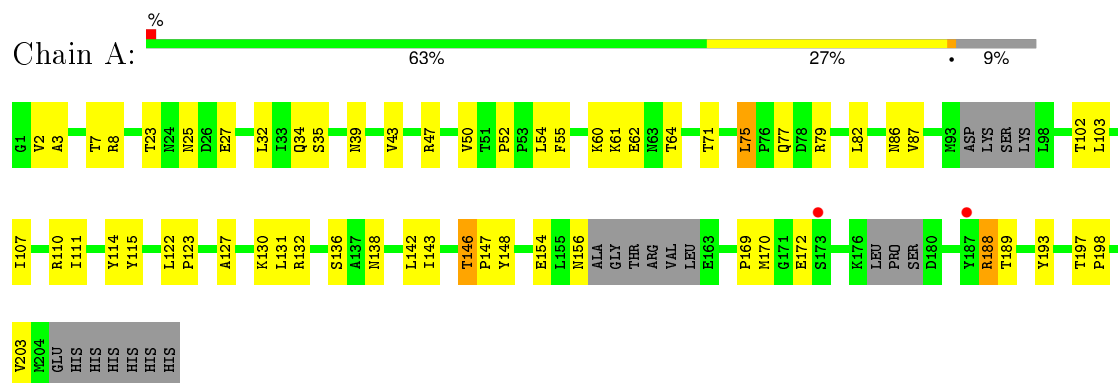
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	8	Total	O	0	0
			8	8		
6	B	13	Total	O	0	0
			13	13		
6	C	13	Total	O	0	0
			13	13		
6	D	14	Total	O	0	0
			14	14		

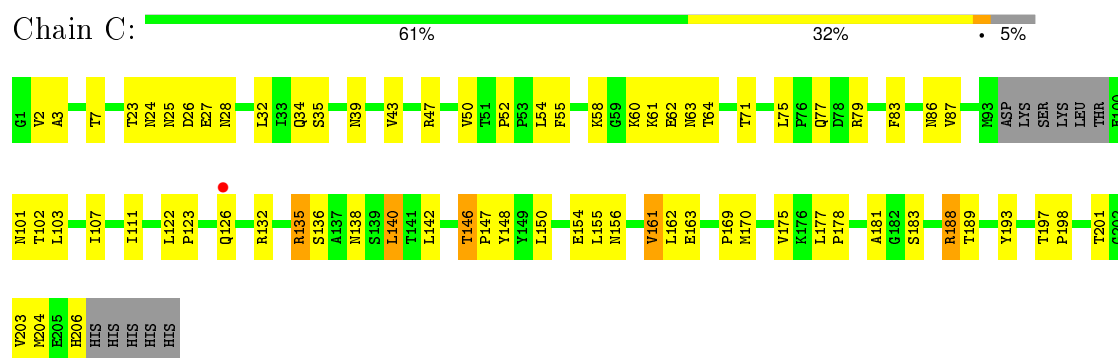
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

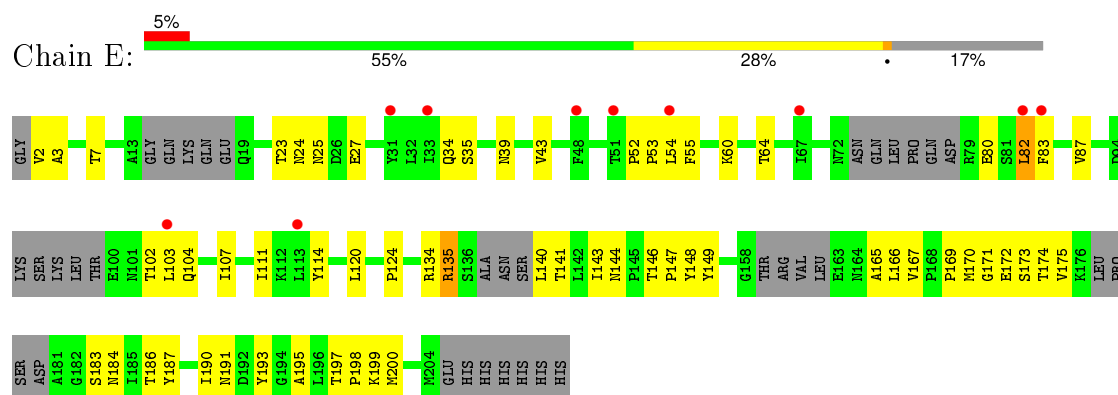
• Molecule 1: Chaperone protein fimC



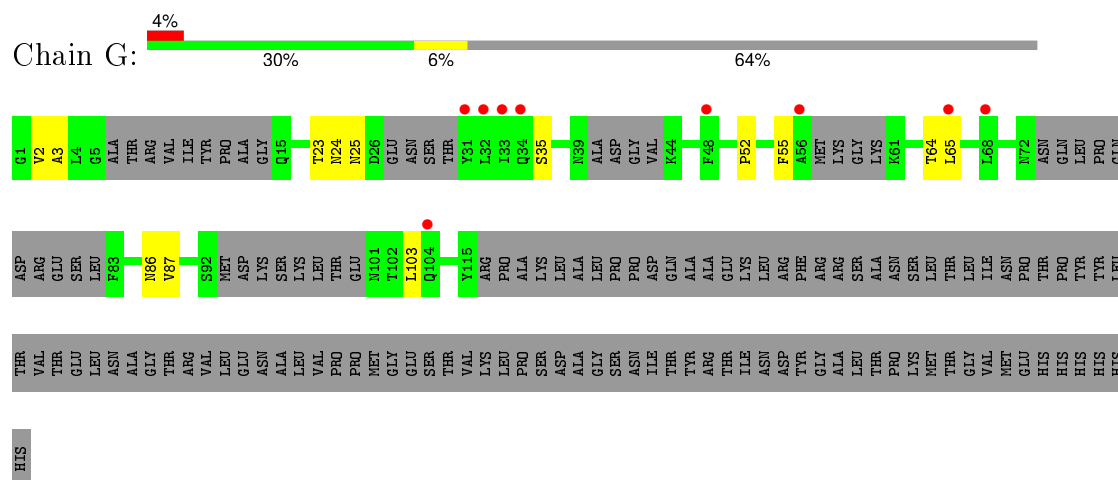
• Molecule 1: Chaperone protein fimC



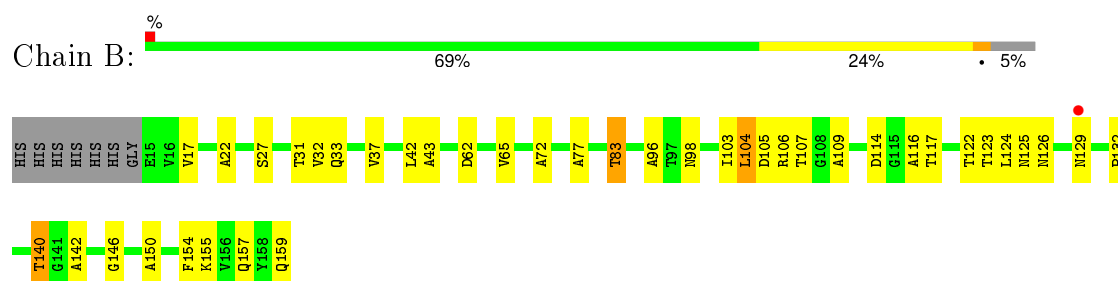
• Molecule 1: Chaperone protein fimC



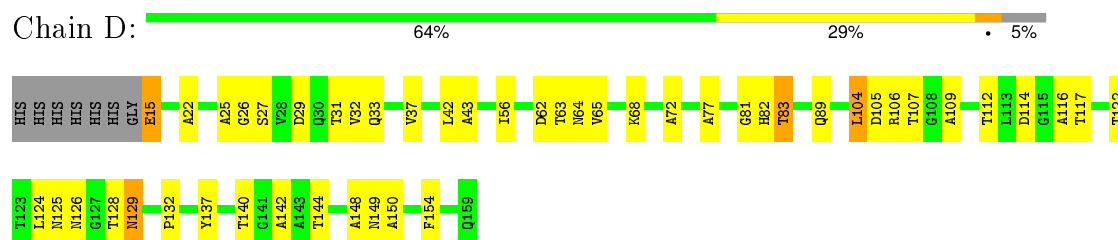
- Molecule 1: Chaperone protein fimC



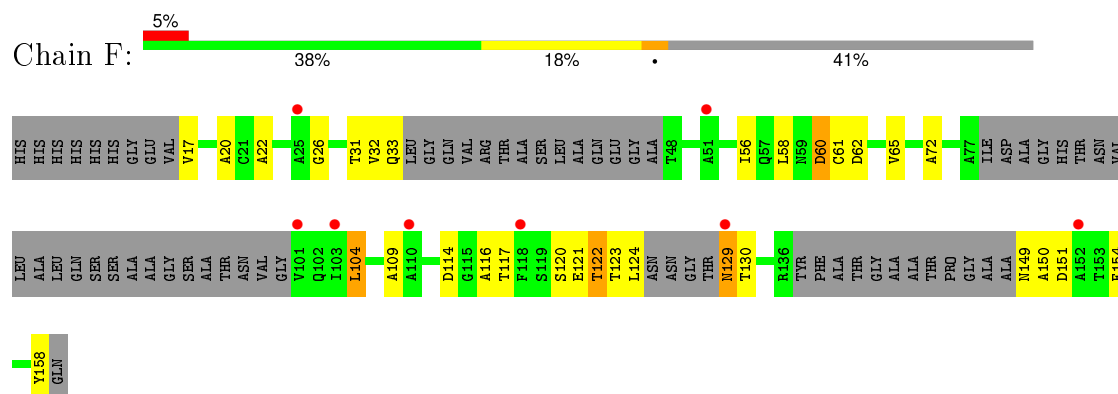
- Molecule 2: Type-1 fimbrial protein, A chain



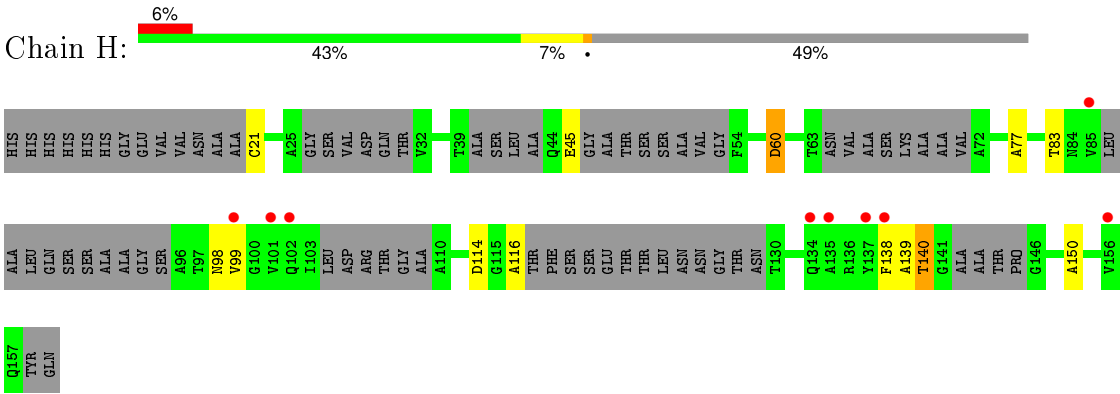
- Molecule 2: Type-1 fimbrial protein, A chain



- Molecule 2: Type-1 fimbrial protein, A chain



- Molecule 2: Type-1 fimbrial protein, A chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	79.62Å 142.18Å 171.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.39 – 3.20 58.39 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (58.39-3.20) 99.8 (58.39-3.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.261 , 0.300 0.260 , 0.303	Depositor DCC
R_{free} test set	1021 reflections (3.20%)	DCC
Wilson B-factor (Å ²)	84.2	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 89.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 32912 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7840	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, PEG, EDO

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	0.30	0/1472	0.52	1/2007 (0.0%)
1	C	0.32	0/1572	0.55	2/2143 (0.1%)
1	E	0.27	0/1216	0.52	2/1669 (0.1%)
1	G	0.23	0/487	0.43	0/667
2	B	0.33	0/1024	0.52	0/1399
2	D	0.32	0/1012	0.50	0/1384
2	F	0.27	0/593	0.52	0/810
2	H	0.22	0/460	0.42	0/623
All	All	0.30	0/7836	0.51	5/10702 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	135	ARG	NE-CZ-NH1	-7.26	116.67	120.30
1	C	135	ARG	NE-CZ-NH2	7.13	123.87	120.30
1	E	135	ARG	NE-CZ-NH2	-6.86	116.87	120.30
1	E	135	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	50	VAL	CB-CA-C	-5.27	101.39	111.40

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	1447	0	1412	58	0
1	C	1543	0	1549	59	0
1	E	1198	0	1038	59	0
1	G	485	0	360	8	0
2	B	1014	0	975	33	0
2	D	1002	0	955	53	0
2	F	592	0	526	32	0
2	H	466	0	353	13	0
3	A	14	0	20	1	0
3	B	7	0	10	1	0
4	C	4	0	6	0	0
4	D	4	0	6	3	0
5	C	8	0	12	1	0
5	D	8	0	12	4	0
6	A	8	0	0	0	0
6	B	13	0	0	0	0
6	C	13	0	0	1	0
6	D	14	0	0	1	0
All	All	7840	0	7234	278	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 18.

All (278) 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:D:83:THR:HB	4:D:160:EDO:H11	1.50	0.92
2:F:72:ALA:HB2	2:F:117:THR:HA	1.54	0.88
2:D:72:ALA:HB2	2:D:117:THR:HA	1.55	0.87
2:D:105:ASP:OD2	2:D:107:THR:HG22	1.77	0.85
1:E:120:LEU:HD13	1:E:148:TYR:HE2	1.40	0.85
1:C:122:LEU:HD23	1:C:146:THR:HG23	1.59	0.84
2:B:72:ALA:HB2	2:B:117:THR:HA	1.59	0.82
1:A:122:LEU:HD23	1:A:146:THR:HG23	1.62	0.81
2:B:105:ASP:OD2	2:B:107:THR:HG22	1.79	0.81
1:A:193:TYR:HB3	2:B:17:VAL:HA	1.61	0.80
1:C:177:LEU:HD12	1:C:178:PRO:HD2	1.65	0.78
1:C:156:ASN:HB3	1:C:161:VAL:HA	1.66	0.76
2:F:129:ASN:N	2:F:129:ASN:HD22	1.85	0.75
2:D:124:LEU:HG	2:D:129:ASN:ND2	2.00	0.75
2:D:124:LEU:HG	2:D:129:ASN:HD21	1.50	0.74
2:D:83:THR:H	5:D:161:TRS:H12	1.54	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:122:THR:HB	2:F:124:LEU:H	1.54	0.73
1:E:149:TYR:HA	1:E:169:PRO:HD3	1.71	0.73
1:A:110:ARG:HD3	2:B:157:GLN:OE1	1.90	0.70
2:D:149:ASN:HB2	6:D:170:HOH:O	1.91	0.70
2:D:82:HIS:HA	5:D:161:TRS:H32	1.74	0.69
1:E:140:LEU:HD22	1:E:175:VAL:HG23	1.76	0.67
1:E:120:LEU:HD13	1:E:148:TYR:CE2	2.29	0.67
2:D:129:ASN:HD22	2:D:129:ASN:N	1.91	0.67
2:B:62:ASP:HB3	2:B:65:VAL:HG13	1.77	0.67
1:A:146:THR:HG22	1:A:148:TYR:HD1	1.60	0.66
1:A:27:GLU:HB2	1:A:60:LYS:HG3	1.77	0.66
1:C:102:THR:HG22	2:D:33:GLN:NE2	2.09	0.66
1:C:102:THR:HG22	2:D:33:GLN:HE21	1.61	0.66
1:E:191:ASN:OD1	1:E:195:ALA:N	2.29	0.66
2:D:62:ASP:HB3	2:D:65:VAL:HG13	1.78	0.64
2:F:62:ASP:HB3	2:F:65:VAL:HG13	1.79	0.64
1:C:193:TYR:CE1	2:D:15:GLU:HA	2.33	0.64
1:C:146:THR:HG22	1:C:148:TYR:HD1	1.63	0.64
2:F:123:THR:HG22	2:F:130:THR:HG23	1.80	0.63
1:C:178:PRO:HG2	1:C:181:ALA:HB2	1.80	0.63
1:E:149:TYR:HD1	1:E:166:LEU:HD21	1.64	0.63
1:A:154:GLU:OE1	1:A:188:ARG:HD3	1.99	0.63
1:E:103:LEU:HA	2:F:150:ALA:O	1.98	0.63
1:C:146:THR:HG22	1:C:147:PRO:HD2	1.80	0.62
1:C:27:GLU:HG2	1:C:60:LYS:HG3	1.80	0.62
1:A:47:ARG:NH1	2:D:109:ALA:HB2	2.14	0.62
1:E:82:LEU:HD23	1:E:114:TYR:CE1	2.34	0.62
1:C:122:LEU:HD11	1:C:126:GLN:HB2	1.82	0.62
2:H:45:GLU:HB2	2:H:140:THR:HA	1.82	0.61
1:E:149:TYR:CD1	1:E:166:LEU:HD21	2.35	0.61
2:D:140:THR:O	2:D:140:THR:HG22	2.00	0.61
1:A:35:SER:HB3	1:A:87:VAL:HG22	1.82	0.61
1:C:154:GLU:OE1	1:C:188:ARG:HD3	2.00	0.61
1:E:34:GLN:HB3	1:E:54:LEU:HD13	1.83	0.60
1:C:122:LEU:HD12	1:C:123:PRO:HD2	1.82	0.60
1:C:52:PRO:HG2	1:C:55:PHE:CD2	2.36	0.60
1:A:52:PRO:HG2	1:A:55:PHE:CD2	2.37	0.60
2:F:62:ASP:O	2:F:65:VAL:HG22	2.02	0.60
1:A:102:THR:HB	2:B:31:THR:CG2	2.32	0.59
1:A:71:THR:HG21	1:A:75:LEU:HD22	1.84	0.59
1:A:122:LEU:HD12	1:A:123:PRO:HD2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:102:THR:HG22	2:F:33:GLN:NE2	2.16	0.59
1:A:146:THR:HG22	1:A:147:PRO:HD2	1.84	0.58
1:A:136:SER:O	1:A:138:ASN:N	2.34	0.58
2:D:122:THR:HB	2:D:129:ASN:HB3	1.84	0.58
1:E:140:LEU:HD23	1:E:141:THR:N	2.18	0.58
1:E:35:SER:HB3	1:E:87:VAL:HG22	1.84	0.58
2:D:81:GLY:O	5:D:161:TRS:H11	2.03	0.58
1:E:34:GLN:HG3	1:E:34:GLN:O	2.04	0.58
1:E:197:THR:HB	1:E:198:PRO:HD2	1.84	0.58
2:D:62:ASP:O	2:D:65:VAL:HG22	2.04	0.57
1:A:107:ILE:HD13	2:B:154:PHE:HB3	1.85	0.57
2:D:43:ALA:O	2:D:142:ALA:HA	2.05	0.57
1:E:34:GLN:CB	1:E:54:LEU:HD13	2.35	0.57
1:E:7:THR:O	1:E:111:ILE:HB	2.03	0.57
1:E:143:ILE:N	1:E:143:ILE:HD12	2.20	0.56
1:C:34:GLN:CB	1:C:54:LEU:HD13	2.35	0.56
1:A:34:GLN:OE1	3:A:212:PEG:H12	2.06	0.56
1:G:35:SER:HB3	1:G:87:VAL:HG22	1.88	0.56
2:D:106:ARG:HB3	2:D:132:PRO:HG2	1.88	0.56
1:E:135:ARG:HA	1:E:140:LEU:HA	1.89	0.55
1:E:52:PRO:HG2	1:E:55:PHE:CD2	2.42	0.55
1:A:39:ASN:HD21	1:A:43:VAL:HB	1.71	0.55
1:C:156:ASN:CB	1:C:161:VAL:HA	2.35	0.55
2:D:77:ALA:CB	2:D:83:THR:HA	2.37	0.55
2:B:98:ASN:HA	2:B:140:THR:HG22	1.89	0.55
1:G:52:PRO:HG2	1:G:55:PHE:CD2	2.41	0.55
1:E:104:GLN:HB2	2:F:151:ASP:HB3	1.89	0.54
1:A:102:THR:HB	2:B:31:THR:HG22	1.90	0.54
1:E:103:LEU:HD12	2:F:151:ASP:HA	1.90	0.54
1:C:63:ASN:HB3	6:C:214:HOH:O	2.07	0.54
1:C:34:GLN:HG3	1:C:34:GLN:O	2.07	0.54
2:B:62:ASP:O	2:B:65:VAL:HG22	2.08	0.54
2:D:125:ASN:OD1	1:E:200:MET:HA	2.08	0.53
1:C:25:ASN:O	1:C:60:LYS:HG2	2.08	0.53
1:E:143:ILE:HG13	1:E:172:GLU:HG2	1.90	0.53
1:E:141:THR:HG23	1:E:174:THR:HA	1.91	0.53
1:A:34:GLN:O	1:A:34:GLN:HG3	2.08	0.53
1:E:193:TYR:O	2:F:17:VAL:HA	2.09	0.53
1:A:8:ARG:NH2	2:B:159:GLN:O	2.42	0.53
1:E:82:LEU:HD22	1:E:83:PHE:N	2.24	0.53
2:B:106:ARG:HB3	2:B:132:PRO:HG2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:SER:HB3	1:C:87:VAL:HG22	1.90	0.52
1:C:7:THR:O	1:C:111:ILE:HB	2.09	0.52
2:D:25:ALA:O	2:D:26:GLY:C	2.47	0.52
1:E:147:PRO:C	1:E:169:PRO:HB3	2.30	0.52
2:D:105:ASP:CG	2:D:107:THR:HG22	2.29	0.52
2:B:105:ASP:CG	2:B:107:THR:HG22	2.29	0.52
1:C:102:THR:HB	2:D:31:THR:CG2	2.39	0.52
1:A:146:THR:CG2	1:A:147:PRO:HD2	2.40	0.52
2:B:77:ALA:CB	2:B:83:THR:HA	2.40	0.52
2:H:77:ALA:CB	2:H:83:THR:HA	2.40	0.52
1:E:3:ALA:HB2	2:F:22:ALA:HA	1.91	0.51
1:C:146:THR:CG2	1:C:147:PRO:HD2	2.40	0.51
2:D:89:GLN:HB3	2:D:149:ASN:O	2.11	0.51
1:C:136:SER:O	1:C:138:ASN:N	2.34	0.51
1:C:58:LYS:O	1:C:61:LYS:HE2	2.11	0.51
1:E:82:LEU:HD23	1:E:114:TYR:HE1	1.76	0.50
1:C:156:ASN:OD1	1:C:156:ASN:N	2.44	0.50
1:E:102:THR:N	2:F:149:ASN:OD1	2.31	0.50
2:B:125:ASN:O	2:B:126:ASN:HB2	2.11	0.50
1:A:27:GLU:HA	1:A:60:LYS:H	1.77	0.50
1:A:34:GLN:CB	1:A:54:LEU:HD13	2.42	0.50
1:A:127:ALA:O	1:A:130:LYS:HG2	2.11	0.50
1:E:183:SER:HB3	1:E:184:ASN:HA	1.94	0.50
1:G:2:VAL:HA	1:G:23:THR:O	2.11	0.50
1:A:102:THR:HG22	2:B:33:GLN:HA	1.92	0.50
1:A:3:ALA:HB2	2:B:22:ALA:HA	1.93	0.50
2:D:125:ASN:O	2:D:126:ASN:HB2	2.11	0.50
1:C:39:ASN:HD21	1:C:43:VAL:HB	1.76	0.50
2:D:112:THR:HB	4:D:160:EDO:H22	1.94	0.50
1:A:75:LEU:O	1:A:77:GLN:NE2	2.44	0.50
2:F:60:ASP:N	2:F:60:ASP:OD2	2.45	0.49
1:C:75:LEU:O	1:C:77:GLN:NE2	2.44	0.49
2:D:124:LEU:HD23	2:D:125:ASN:N	2.27	0.49
1:C:189:THR:O	1:C:197:THR:HG23	2.13	0.49
1:A:3:ALA:HB2	2:B:22:ALA:CA	2.43	0.49
2:D:63:THR:HB	2:D:124:LEU:HD12	1.94	0.49
2:H:140:THR:OG1	2:H:140:THR:O	2.23	0.49
1:E:167:VAL:HA	1:E:173:SER:HB3	1.95	0.49
2:F:114:ASP:N	2:F:114:ASP:OD2	2.46	0.49
2:D:114:ASP:C	2:D:116:ALA:H	2.17	0.48
1:A:27:GLU:HB2	1:A:60:LYS:CG	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:GLU:OE1	5:C:213:TRS:H12	2.12	0.48
1:C:34:GLN:HB3	1:C:54:LEU:HD13	1.94	0.48
1:E:80:GLU:OE2	1:E:149:TYR:N	2.47	0.48
2:D:129:ASN:ND2	2:D:129:ASN:N	2.61	0.48
1:G:103:LEU:HA	2:H:150:ALA:O	2.13	0.48
2:D:112:THR:HB	4:D:160:EDO:C2	2.44	0.48
1:A:156:ASN:OD1	1:A:156:ASN:N	2.47	0.48
2:B:27:SER:OG	2:B:31:THR:O	2.28	0.48
2:D:140:THR:O	2:D:140:THR:CG2	2.62	0.47
1:A:142:LEU:HD12	1:A:142:LEU:N	2.29	0.47
1:C:47:ARG:HB3	1:C:71:THR:HG22	1.96	0.47
2:F:123:THR:O	2:F:124:LEU:HB2	2.14	0.47
1:E:149:TYR:CA	1:E:169:PRO:HD3	2.42	0.47
1:C:32:LEU:HG	1:C:54:LEU:HD11	1.97	0.47
2:B:123:THR:C	2:B:125:ASN:H	2.18	0.47
2:F:114:ASP:C	2:F:116:ALA:H	2.17	0.47
2:B:96:ALA:N	2:B:146:GLY:HA3	2.29	0.47
2:D:56:ILE:HD13	2:D:154:PHE:CD1	2.50	0.47
1:A:32:LEU:HG	1:A:54:LEU:HD11	1.97	0.46
1:A:189:THR:O	1:A:197:THR:HG23	2.16	0.46
2:B:114:ASP:C	2:B:116:ALA:H	2.17	0.46
1:E:124:PRO:HA	1:E:148:TYR:CE1	2.50	0.46
2:F:129:ASN:ND2	2:F:129:ASN:N	2.58	0.46
1:E:167:VAL:HG12	1:E:171:GLY:O	2.15	0.46
1:A:7:THR:O	1:A:111:ILE:HB	2.16	0.46
2:D:27:SER:OG	2:D:31:THR:HB	2.17	0.45
2:D:114:ASP:OD2	2:D:114:ASP:N	2.44	0.45
1:A:146:THR:O	1:A:169:PRO:HA	2.16	0.45
1:E:187:TYR:O	1:E:200:MET:N	2.43	0.45
2:D:77:ALA:HB1	2:D:83:THR:HA	1.98	0.45
1:E:82:LEU:HD22	1:E:83:PHE:H	1.82	0.45
2:H:114:ASP:C	2:H:116:ALA:H	2.19	0.45
2:B:122:THR:HB	2:B:129:ASN:HB3	1.98	0.45
2:B:104:LEU:HD23	2:B:104:LEU:N	2.31	0.45
2:F:56:ILE:HD13	2:F:154:PHE:CD1	2.50	0.45
1:A:23:THR:HG22	1:A:62:GLU:HB2	1.98	0.45
2:F:58:LEU:N	2:F:58:LEU:HD23	2.31	0.45
2:F:31:THR:HG22	2:F:32:VAL:N	2.32	0.45
1:A:197:THR:HB	1:A:198:PRO:HD2	1.99	0.45
1:E:107:ILE:CD1	2:F:154:PHE:HB3	2.47	0.44
1:E:2:VAL:HA	1:E:23:THR:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:122:THR:HB	2:F:124:LEU:N	2.27	0.44
1:C:26:ASP:OD1	1:C:28:ASN:N	2.50	0.44
1:C:197:THR:HB	1:C:198:PRO:HD2	1.98	0.44
1:C:146:THR:O	1:C:169:PRO:HA	2.17	0.44
1:E:140:LEU:HD23	1:E:140:LEU:C	2.38	0.44
2:D:56:ILE:HD13	2:D:154:PHE:CE1	2.52	0.44
1:C:101:ASN:ND2	2:D:37:VAL:HG23	2.32	0.44
2:H:114:ASP:N	2:H:114:ASP:OD2	2.46	0.44
2:F:123:THR:CG2	2:F:130:THR:HG23	2.48	0.44
1:E:167:VAL:HG22	1:E:173:SER:OG	2.18	0.44
2:B:31:THR:HG22	2:B:32:VAL:N	2.33	0.44
2:H:60:ASP:OD1	2:H:60:ASP:N	2.51	0.44
1:E:144:ASN:HB2	1:E:167:VAL:HB	1.98	0.43
2:H:45:GLU:HG3	2:H:139:ALA:O	2.17	0.43
1:A:34:GLN:HB3	1:A:54:LEU:HD13	2.00	0.43
1:E:39:ASN:HD21	1:E:43:VAL:HB	1.83	0.43
1:C:79:ARG:HB3	1:C:170:MET:HE2	2.00	0.43
1:C:142:LEU:N	1:C:142:LEU:HD12	2.32	0.43
1:C:103:LEU:HB3	2:D:32:VAL:HB	1.99	0.43
1:E:165:ALA:HB2	1:E:175:VAL:HG11	2.00	0.43
2:H:77:ALA:HB1	2:H:83:THR:HA	2.00	0.43
2:F:56:ILE:HD13	2:F:154:PHE:CE1	2.53	0.43
1:C:23:THR:HG22	1:C:62:GLU:HB2	2.00	0.43
1:C:163:GLU:HG3	1:C:175:VAL:HB	2.00	0.43
2:D:82:HIS:HA	5:D:161:TRS:H12	2.01	0.43
1:A:25:ASN:O	1:A:60:LYS:HE2	2.19	0.43
2:D:31:THR:HG22	2:D:32:VAL:N	2.33	0.43
1:A:75:LEU:HD23	1:A:115:TYR:CD2	2.53	0.43
1:G:24:ASN:OD1	1:G:25:ASN:N	2.51	0.43
1:A:47:ARG:HH12	2:D:109:ALA:HB2	1.83	0.43
2:D:126:ASN:O	2:D:128:THR:HG23	2.19	0.43
1:A:147:PRO:C	1:A:169:PRO:HB3	2.39	0.43
2:D:137:TYR:CE2	2:D:148:ALA:HB1	2.54	0.43
1:C:132:ARG:HH11	1:C:203:VAL:HB	1.84	0.43
1:A:60:LYS:O	1:A:61:LYS:HG3	2.18	0.42
1:E:141:THR:HG21	1:E:172:GLU:OE1	2.19	0.42
1:C:147:PRO:C	1:C:169:PRO:HB3	2.39	0.42
1:E:134:ARG:O	1:E:141:THR:N	2.51	0.42
1:C:79:ARG:HB3	1:C:170:MET:CE	2.50	0.42
1:A:103:LEU:HD12	2:B:150:ALA:O	2.20	0.42
1:E:183:SER:CB	1:E:184:ASN:HA	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:98:ASN:CA	2:H:140:THR:HG23	2.48	0.42
1:C:107:ILE:CD1	2:D:154:PHE:HB3	2.49	0.42
2:H:98:ASN:HA	2:H:140:THR:HG23	2.01	0.42
1:E:24:ASN:OD1	1:E:25:ASN:N	2.53	0.42
1:A:79:ARG:HB3	1:A:170:MET:CE	2.50	0.42
1:E:169:PRO:O	1:E:170:MET:HB2	2.19	0.42
1:G:35:SER:HA	1:G:86:ASN:O	2.20	0.42
1:C:150:LEU:HA	1:C:150:LEU:HD23	1.92	0.42
2:H:99:VAL:HA	2:H:138:PHE:O	2.20	0.42
1:A:2:VAL:HA	1:A:23:THR:O	2.20	0.41
1:E:143:ILE:HG22	1:E:144:ASN:N	2.35	0.41
1:C:24:ASN:OD1	1:C:25:ASN:N	2.53	0.41
1:E:52:PRO:HA	1:E:53:PRO:HD3	1.89	0.41
1:A:82:LEU:HD13	1:A:114:TYR:HE1	1.85	0.41
1:A:35:SER:HA	1:A:86:ASN:O	2.19	0.41
1:E:3:ALA:HB2	2:F:22:ALA:CA	2.49	0.41
2:F:56:ILE:HG21	2:F:154:PHE:CE1	2.55	0.41
1:C:2:VAL:HA	1:C:23:THR:O	2.20	0.41
2:B:43:ALA:O	2:B:142:ALA:HA	2.19	0.41
1:A:102:THR:CB	2:B:31:THR:CG2	2.97	0.41
1:A:79:ARG:HB3	1:A:170:MET:HE2	2.02	0.41
1:A:131:LEU:O	1:A:132:ARG:HD3	2.20	0.41
1:C:60:LYS:O	1:C:61:LYS:HG3	2.20	0.41
1:G:55:PHE:HE2	1:G:65:LEU:HD21	1.85	0.41
2:B:77:ALA:HB1	2:B:83:THR:HA	2.01	0.41
1:C:47:ARG:HD2	1:C:83:PHE:HZ	1.85	0.41
2:B:104:LEU:HA	2:B:109:ALA:O	2.20	0.41
2:D:37:VAL:HG22	2:D:137:TYR:CE1	2.55	0.41
1:E:146:THR:O	1:E:169:PRO:HA	2.20	0.41
1:E:166:LEU:HD23	1:E:166:LEU:O	2.20	0.41
1:A:27:GLU:HA	1:A:60:LYS:N	2.35	0.41
1:C:25:ASN:O	1:C:26:ASP:C	2.58	0.41
2:B:103:ILE:C	2:B:104:LEU:HD23	2.41	0.41
1:A:47:ARG:NH2	2:D:107:THR:O	2.54	0.41
1:A:193:TYR:HB3	2:B:17:VAL:CA	2.43	0.41
1:E:102:THR:HG22	2:F:33:GLN:HE21	1.85	0.41
1:C:34:GLN:O	1:C:34:GLN:CG	2.69	0.41
1:A:34:GLN:O	1:A:34:GLN:CG	2.69	0.41
1:A:132:ARG:HH11	1:A:203:VAL:HB	1.86	0.41
2:F:104:LEU:HA	2:F:109:ALA:O	2.20	0.41
1:E:186:THR:OG1	1:E:199:LYS:HE3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ARG:HG2	1:C:136:SER:N	2.36	0.41
2:D:104:LEU:HA	2:D:109:ALA:O	2.21	0.40
2:F:61:CYS:HB2	2:F:65:VAL:HG21	2.02	0.40
1:C:135:ARG:HA	1:C:140:LEU:HD23	2.01	0.40
2:F:104:LEU:N	2:F:104:LEU:HD23	2.36	0.40
1:C:103:LEU:HA	2:D:150:ALA:O	2.21	0.40
2:F:20:ALA:HB1	2:F:158:TYR:CZ	2.56	0.40
1:C:35:SER:HA	1:C:86:ASN:O	2.21	0.40
2:B:142:ALA:HB1	3:B:160:PEG:H41	2.03	0.40
1:G:3:ALA:HB2	2:H:21:CYS:O	2.21	0.40
2:D:64:ASN:O	2:D:68:LYS:HG2	2.21	0.40
1:A:136:SER:C	1:A:138:ASN:H	2.23	0.40
1:C:3:ALA:HB2	2:D:22:ALA:HA	2.02	0.40
1:E:27:GLU:HG2	1:E:60:LYS:CG	2.52	0.40
1:A:143:ILE:HG23	1:A:172:GLU:HG2	2.04	0.40
1:C:155:LEU:HD23	1:C:162:LEU:HB2	2.03	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	183/211 (87%)	172 (94%)	11 (6%)	0	100	100
1	C	196/211 (93%)	183 (93%)	13 (7%)	0	100	100
1	E	162/211 (77%)	141 (87%)	21 (13%)	0	100	100
1	G	62/211 (29%)	61 (98%)	1 (2%)	0	100	100
2	B	143/152 (94%)	131 (92%)	11 (8%)	1 (1%)	26	72
2	D	143/152 (94%)	131 (92%)	12 (8%)	0	100	100
2	F	79/152 (52%)	73 (92%)	4 (5%)	2 (2%)	7	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	60/152 (40%)	55 (92%)	5 (8%)	0	100	100
All	All	1028/1452 (71%)	947 (92%)	78 (8%)	3 (0%)	46	85

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	120	SER
2	B	124	LEU
2	F	26	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	149/182 (82%)	145 (97%)	4 (3%)	52	85
1	C	166/182 (91%)	156 (94%)	10 (6%)	24	65
1	E	104/182 (57%)	101 (97%)	3 (3%)	50	83
1	G	34/182 (19%)	33 (97%)	1 (3%)	50	83
2	B	102/109 (94%)	96 (94%)	6 (6%)	24	65
2	D	99/109 (91%)	92 (93%)	7 (7%)	18	57
2	F	54/109 (50%)	49 (91%)	5 (9%)	11	41
2	H	31/109 (28%)	29 (94%)	2 (6%)	21	61
All	All	739/1164 (64%)	701 (95%)	38 (5%)	29	70

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	THR
1	A	75	LEU
1	A	146	THR
1	A	188	ARG
2	B	37	VAL

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Mol	Chain	Res	Type
2	B	42	LEU
2	B	83	THR
2	B	104	LEU
2	B	140	THR
2	B	155	LYS
1	C	50	VAL
1	C	64	THR
1	C	140	LEU
1	C	146	THR
1	C	161	VAL
1	C	183	SER
1	C	188	ARG
1	C	201	THR
1	C	204	MET
1	C	206	HIS
2	D	15	GLU
2	D	29	ASP
2	D	42	LEU
2	D	83	THR
2	D	104	LEU
2	D	129	ASN
2	D	144	THR
1	E	64	THR
1	E	82	LEU
1	E	190	ILE
2	F	60	ASP
2	F	104	LEU
2	F	121	GLU
2	F	122	THR
2	F	129	ASN
1	G	64	THR
2	H	60	ASP
2	H	140	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	33	GLN
2	D	129	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PEG	A	212	-	6,6,6	0.51	0	5,5,5	1.61	1 (20%)
3	PEG	A	213	-	6,6,6	0.63	0	5,5,5	1.31	0
3	PEG	B	160	-	6,6,6	0.62	0	5,5,5	1.34	0
4	EDO	C	212	-	3,3,3	0.51	0	2,2,2	0.36	0
5	TRS	C	213	-	7,7,7	0.93	1 (14%)	9,9,9	0.44	0
4	EDO	D	160	-	3,3,3	0.56	0	2,2,2	0.13	0
5	TRS	D	161	-	7,7,7	0.96	1 (14%)	9,9,9	0.44	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PEG	A	212	-	-	0/4/4/4	0/0/0/0
3	PEG	A	213	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PEG	B	160	-	-	0/4/4/4	0/0/0/0
4	EDO	C	212	-	-	0/1/1/1	0/0/0/0
5	TRS	C	213	-	-	0/9/9/9	0/0/0/0
4	EDO	D	160	-	-	0/1/1/1	0/0/0/0
5	TRS	D	161	-	-	0/9/9/9	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	161	TRS	C-N	-2.49	1.47	1.50
5	C	213	TRS	C-N	-2.43	1.47	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	212	PEG	O2-C2-C1	2.24	120.75	110.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	212	PEG	1	0
3	B	160	PEG	1	0
5	C	213	TRS	1	0
4	D	160	EDO	3	0
5	D	161	TRS	4	0

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	191/211 (90%)	-0.05	2 (1%) 84 75	46, 84, 154, 191	0
1	C	200/211 (94%)	-0.14	1 (0%) 91 87	41, 71, 120, 151	0
1	E	176/211 (83%)	0.21	10 (5%) 27 15	86, 138, 191, 234	0
1	G	76/211 (36%)	0.44	9 (11%) 6 4	116, 158, 198, 257	0
2	B	145/152 (95%)	-0.10	1 (0%) 89 83	41, 69, 127, 206	0
2	D	145/152 (95%)	-0.13	0 100 100	46, 77, 126, 196	0
2	F	89/152 (58%)	0.51	8 (8%) 12 6	90, 136, 188, 223	0
2	H	78/152 (51%)	0.63	9 (11%) 6 4	95, 153, 206, 220	0
All	All	1100/1452 (75%)	0.09	40 (3%) 46 31	41, 99, 184, 257	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	135	ALA	7.6
2	H	138	PHE	6.2
2	H	137	TYR	6.1
1	A	187	TYR	4.7
1	G	56	ALA	4.6
2	H	85	VAL	4.0
1	G	65	LEU	4.0
2	F	101	VAL	3.7
2	H	134	GLN	3.6
1	E	82	LEU	3.5
1	E	31	TYR	3.1
1	E	113	LEU	2.9
1	G	32	LEU	2.9
2	F	129	ASN	2.9
2	H	101	VAL	2.8
2	F	110	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
2	F	25	ALA	2.6
2	F	118	PHE	2.6
2	H	156	VAL	2.5
1	E	103	LEU	2.5
2	F	103	ILE	2.5
1	E	51	THR	2.5
1	G	48	PHE	2.4
1	G	31	TYR	2.4
1	E	67	ILE	2.4
2	H	99	VAL	2.3
1	E	33	ILE	2.3
2	B	129	ASN	2.3
1	G	34	GLN	2.2
1	A	173	SER	2.2
1	G	33	ILE	2.2
2	F	152	ALA	2.2
1	E	48	PHE	2.2
2	F	51	ALA	2.2
1	E	83	PHE	2.1
1	G	68	LEU	2.1
2	H	102	GLN	2.1
1	C	126	GLN	2.1
1	G	104	GLN	2.0
1	E	54	LEU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	TRS	C	213	8/8	0.73	0.35	8.56	138,138,138,138	0
3	PEG	A	212	7/7	0.77	0.39	7.82	67,67,67,67	0
5	TRS	D	161	8/8	0.74	0.39	-	140,140,140,140	0
4	EDO	C	212	4/4	0.54	0.31	-	109,109,109,109	0
4	EDO	D	160	4/4	0.67	0.35	-	77,77,77,77	0
3	PEG	A	213	7/7	0.82	0.35	-	111,111,111,111	0
3	PEG	B	160	7/7	0.40	1.06	-	148,148,148,148	0

6.5 Other polymers [i](#)

There are no such residues in this entry.