



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:00 PM UTC

PDB ID : 3SQB / pdb_00003sqb
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

<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:

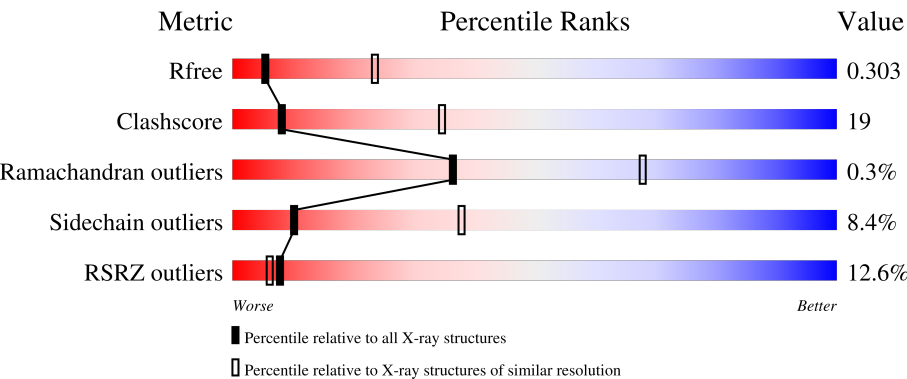
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	211	6% 58%29%•9%
1	C	211	% 60%32%•5%
1	E	211	16% 52%30%•17%
1	G	211	9% 29%6%64%
2	B	152	8% 67%26%••5%

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Mol	Chain	Length	Quality of chain
2	D	152	6% 61% 31% 5%
2	F	152	15% 37% 17% 5% 41%
2	H	152	17% 43% 8% 49%

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:

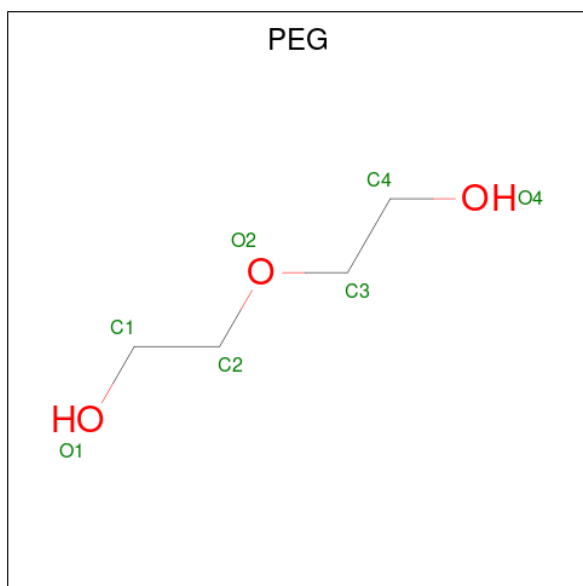
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	HIS	-	expression tag	UNP P04128
B	2	HIS	-	expression tag	UNP P04128
B	3	HIS	-	expression tag	UNP P04128
B	4	HIS	-	expression tag	UNP P04128
B	5	HIS	-	expression tag	UNP P04128
B	6	HIS	-	expression tag	UNP P04128
D	1	HIS	-	expression tag	UNP P04128
D	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 (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



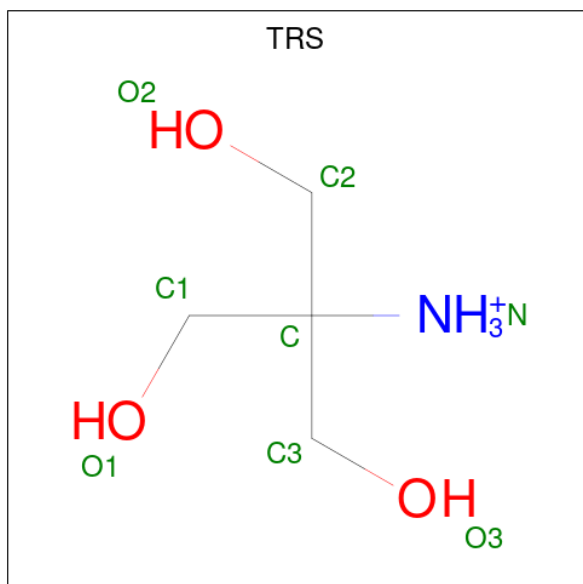
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: 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 (CCD ID: 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		

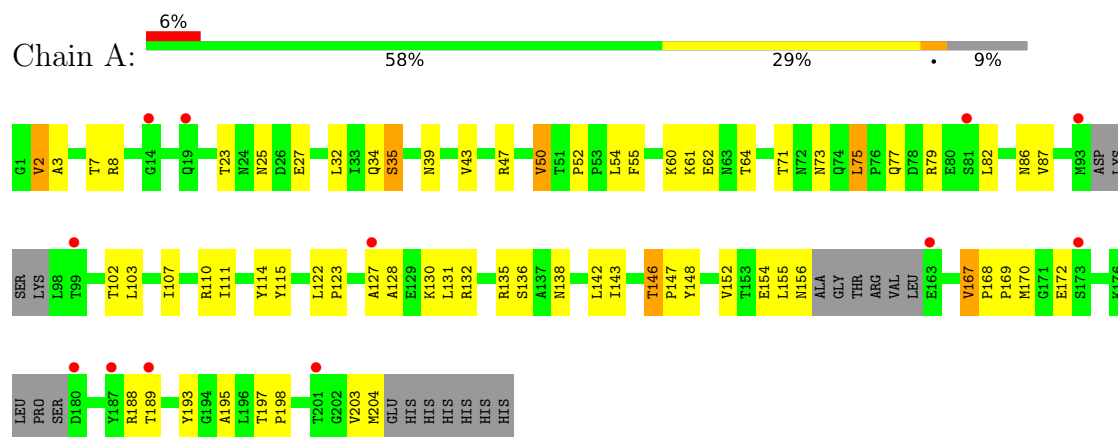
- Molecule 6 is water.

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		

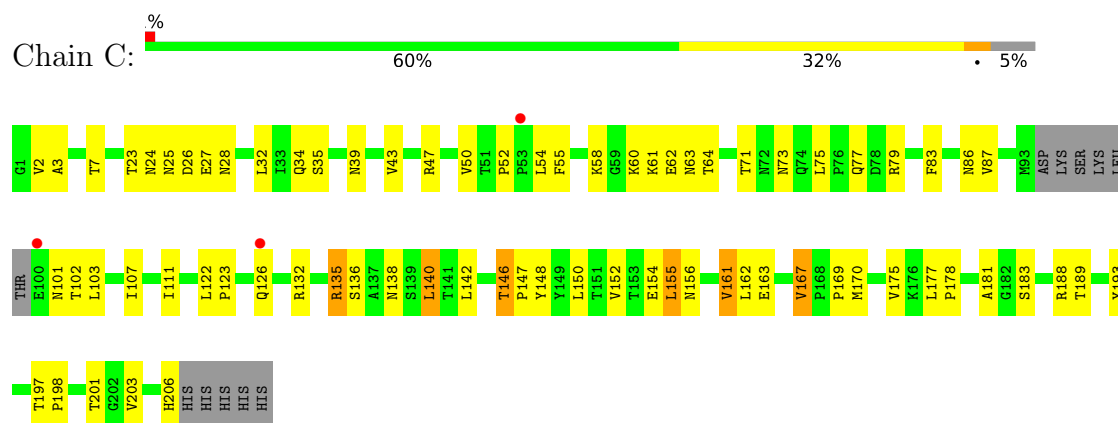
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, 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.

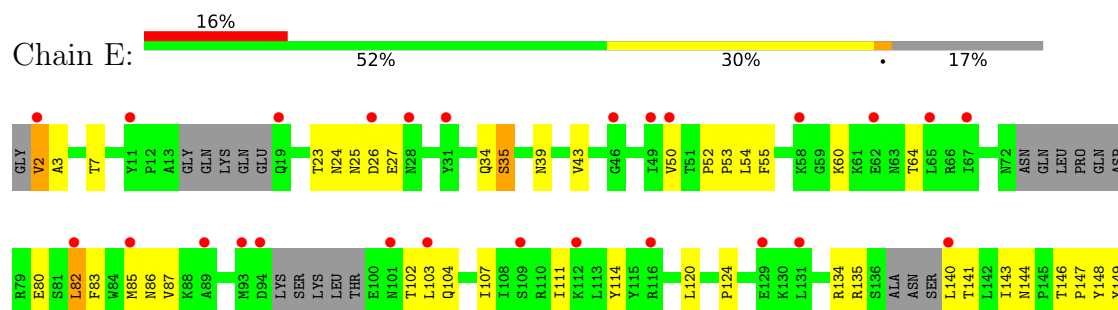
• Molecule 1: Chaperone protein fimC

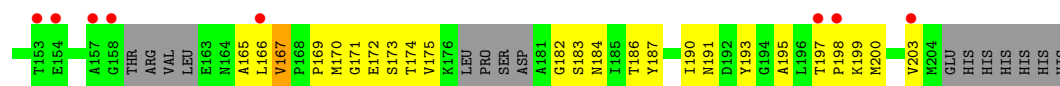


• Molecule 1: Chaperone protein fimC

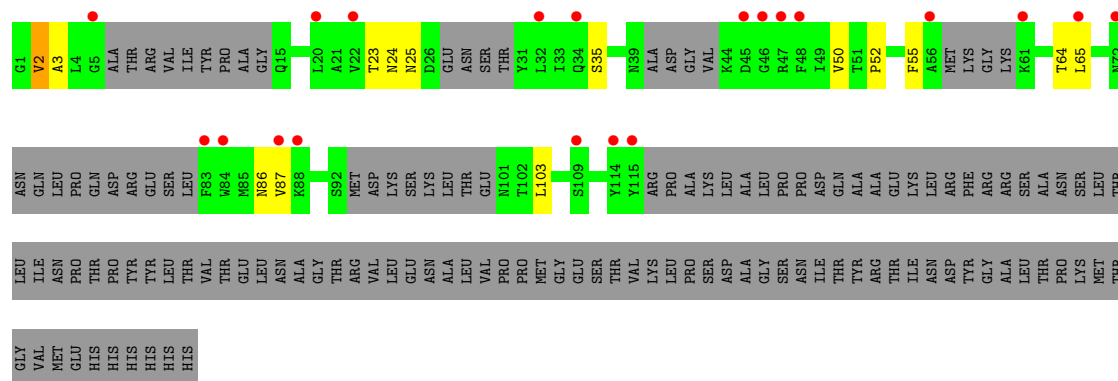


• Molecule 1: Chaperone protein fimC

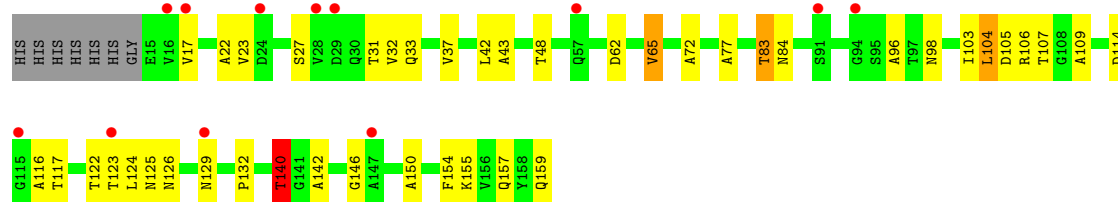




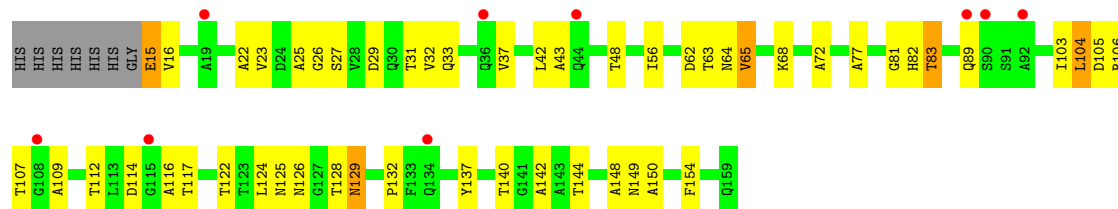
- Molecule 1: Chaperone protein fimC



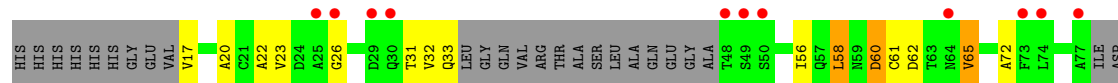
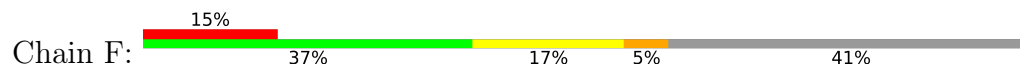
- Molecule 2: Type-1 fimbrial protein, A chain

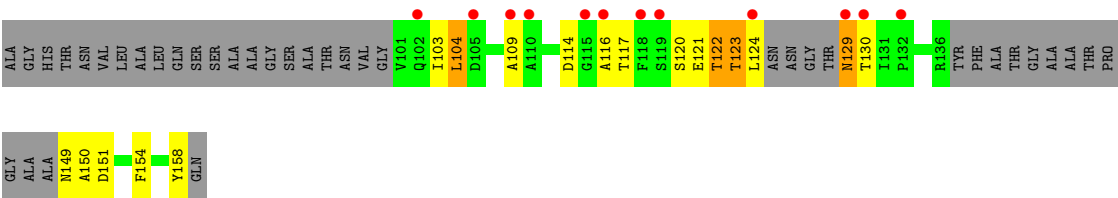


- Molecule 2: Type-1 fimbrial protein, A chain

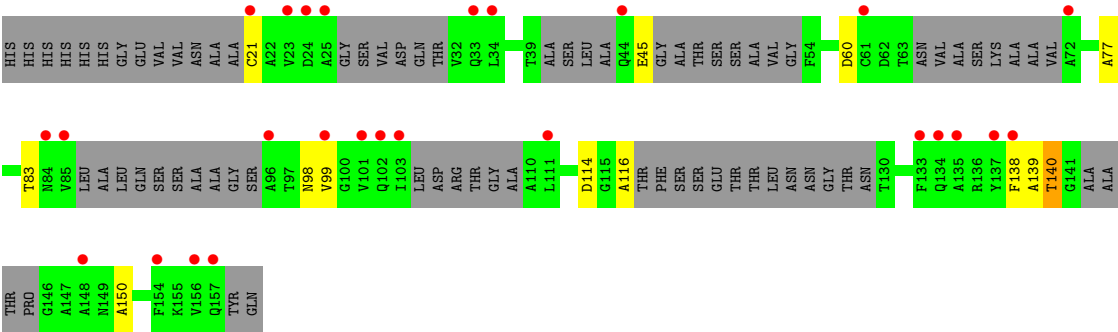
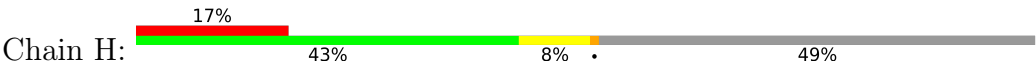


- Molecule 2: Type-1 fimbrial protein, A chain





● 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 1.7.1 _743	Depositor
R, R_{free}	0.261 , 0.300 0.260 , 0.303	Depositor DCC
R_{free} test set	1021 reflections (3.10%)	wwPDB-VP
Wilson B-factor (Å ²)	84.2	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 112.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.39	0/1472	0.82	4/2007 (0.2%)
1	C	0.41	0/1572	0.82	3/2143 (0.1%)
1	E	0.34	0/1216	0.79	0/1669
1	G	0.29	0/487	0.74	0/667
2	B	0.44	0/1024	0.81	1/1399 (0.1%)
2	D	0.40	0/1012	0.76	0/1384
2	F	0.33	0/593	0.73	1/810 (0.1%)
2	H	0.29	0/460	0.63	0/623
All	All	0.38	0/7836	0.78	9/10702 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	123	THR	N-CA-C	-6.44	105.43	113.28
1	A	50	VAL	CB-CA-C	-6.34	101.39	110.83
2	B	140	THR	N-CA-C	-5.79	106.08	114.12
1	C	167	VAL	CA-C-N	5.61	123.70	119.66
1	C	167	VAL	C-N-CA	5.61	123.70	119.66
1	A	167	VAL	CA-C-N	5.29	123.47	119.66
1	A	167	VAL	C-N-CA	5.29	123.47	119.66
1	C	135	ARG	NE-CZ-NH2	5.18	123.87	119.20
1	A	168	PRO	O-C-N	5.04	123.52	121.15

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	62	0
1	C	1543	0	1549	60	0
1	E	1198	0	1038	62	0
1	G	485	0	360	8	0
2	B	1014	0	975	37	0
2	D	1002	0	955	54	0
2	F	592	0	526	33	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	291	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 19.

All (291) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:B:62:ASP:HB3	2:B:65:VAL:HG13	1.77	0.67
2:D:129:ASN:N	2:D:129:ASN:HD22	1.91	0.67
1:A:27:GLU:HB2	1:A:60:LYS:HG3	1.77	0.66
1:A:146:THR:HG22	1:A:148:TYR:HD1	1.60	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:27:GLU:HG2	1:C:60:LYS:HG3	1.80	0.62
1:C:146:THR:HG22	1:C:147:PRO:HD2	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:A:35:SER:HB3	1:A:87:VAL:HG22	1.82	0.61
2:D:140:THR:HG22	2:D:140:THR:O	2.00	0.61
1:E:149:TYR:CD1	1:E:166:LEU:HD21	2.35	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:O	1:E:34:GLN:HG3	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
2:D:25:ALA:O	2:D:26:GLY:C	2.47	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
2:D:105:ASP:CG	2:D:107:THR:HG22	2.29	0.57
2:H:140:THR:O	2:H:140:THR:OG1	2.23	0.57
2:B:105:ASP:CG	2:B:107:THR:HG22	2.29	0.57
1:E:147:PRO:C	1:E:169:PRO:HB3	2.30	0.56
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:A:39:ASN:HD21	1:A:43:VAL:HB	1.71	0.55
1:E:52:PRO:HG2	1:E:55:PHE:CD2	2.42	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:ASN:HB3	6:C:214:HOH:O	2.07	0.54
1:C:34:GLN:O	1:C:34:GLN:HG3	2.07	0.54
2:B:62:ASP:O	2:B:65:VAL:HG22	2.08	0.54
1:C:25:ASN:O	1:C:60:LYS:HG2	2.08	0.53
2:D:125:ASN:OD1	1:E:200:MET:HA	2.08	0.53
2:D:114:ASP:C	2:D:116:ALA:H	2.17	0.53
1:E:141:THR:HG23	1:E:174:THR:HA	1.91	0.53
1:E:143:ILE:HG13	1:E:172:GLU:HG2	1.90	0.53
1:A:34:GLN:O	1:A:34:GLN:HG3	2.08	0.53
1:A:8:ARG:NH2	2:B:159:GLN:O	2.42	0.53
1:E:193:TYR:O	2:F:17:VAL:HA	2.09	0.53
2:B:106:ARG:HB3	2:B:132:PRO:HG2	1.90	0.53
1:E:82:LEU:HD22	1:E:83:PHE:N	2.24	0.53
1:C:7:THR:O	1:C:111:ILE:HB	2.09	0.52
1:C:35:SER:HB3	1:C:87:VAL:HG22	1.90	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:B:114:ASP:C	2:B:116:ALA:H	2.17	0.52
2:H:77:ALA:CB	2:H:83:THR:HA	2.40	0.52
2:B:123:THR:C	2:B:125:ASN:H	2.18	0.52
1:E:3:ALA:HB2	2:F:22:ALA:HA	1.91	0.51
2:F:114:ASP:C	2:F:116:ALA:H	2.17	0.51
1:C:146:THR:CG2	1:C:147:PRO:HD2	2.40	0.51
1:C:136:SER:O	1:C:138:ASN:N	2.34	0.51
2:D:89:GLN:HB3	2:D:149:ASN:O	2.11	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
2:B:125:ASN:O	2:B:126:ASN:HB2	2.11	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
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:3:ALA:HB2	2:B:22:ALA:HA	1.93	0.50
1:A:102:THR:HG22	2:B:33:GLN:HA	1.92	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:H:114:ASP:C	2:H:116:ALA:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:LEU:O	1:A:77:GLN:NE2	2.44	0.50
1:A:135:ARG:H	1:A:204:MET:HE1	1.77	0.50
2:D:112:THR:HB	4:D:160:EDO:H22	1.94	0.50
1:C:75:LEU:O	1:C:77:GLN:NE2	2.44	0.49
2:F:60:ASP:OD2	2:F:60:ASP:N	2.45	0.49
1:C:189:THR:O	1:C:197:THR:HG23	2.13	0.49
2:D:124:LEU:HD23	2:D:125:ASN:N	2.27	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
1:E:167:VAL:HA	1:E:173:SER:HB3	1.95	0.49
2:F:114:ASP:OD2	2:F:114:ASP:N	2.46	0.49
1:E:140:LEU:HD23	1:E:140:LEU:C	2.38	0.48
1:A:27:GLU:HB2	1:A:60:LYS:CG	2.41	0.48
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
1:A:156:ASN:OD1	1:A:156:ASN:N	2.47	0.48
2:D:112:THR:HB	4:D:160:EDO:C2	2.44	0.48
2:B:27:SER:OG	2:B:31:THR:O	2.28	0.48
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:D:140:THR:O	2:D:140:THR:CG2	2.62	0.47
1:A:147:PRO:C	1:A:169:PRO:HB3	2.39	0.47
1:C:32:LEU:HG	1:C:54:LEU:HD11	1.97	0.47
1:E:149:TYR:CA	1:E:169:PRO:HD3	2.42	0.47
2:F:123:THR:O	2:F:124:LEU:HB2	2.14	0.47
1:C:147:PRO:C	1:C:169:PRO:HB3	2.39	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
1:E:124:PRO:HA	1:E:148:TYR:CE1	2.50	0.46
2:F:129:ASN:N	2:F:129:ASN:ND2	2.58	0.46
1:E:167:VAL:HG12	1:E:171:GLY:O	2.15	0.46
2:B:103:ILE:C	2:B:104:LEU:HD23	2.41	0.46
1:C:25:ASN:O	1:C:26:ASP:C	2.58	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:104:LEU:HD23	2:B:104:LEU:N	2.31	0.45
2:B:122:THR:HB	2:B:129:ASN:HB3	1.98	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
1:A:136:SER:C	1:A:138:ASN:H	2.23	0.45
2:F:31:THR:HG22	2:F:32:VAL:N	2.32	0.45
2:F:58:LEU:N	2:F:58:LEU:HD23	2.31	0.45
1:A:197:THR:HB	1:A:198:PRO:HD2	1.99	0.45
1:E:2:VAL:HA	1:E:23:THR:O	2.17	0.44
1:E:107:ILE:CD1	2:F:154:PHE:HB3	2.47	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
2:F:122:THR:HB	2:F:124:LEU:N	2.27	0.44
1:C:101:ASN:ND2	2:D:37:VAL:HG23	2.32	0.44
1:C:136:SER:C	1:C:138:ASN:H	2.22	0.44
1:C:146:THR:O	1:C:169:PRO:HA	2.17	0.44
2:D:56:ILE:HD13	2:D:154:PHE:CE1	2.52	0.44
2:H:114:ASP:OD2	2:H:114:ASP:N	2.46	0.44
2:B:31:THR:HG22	2:B:32:VAL:N	2.33	0.44
1:E:167:VAL:HG22	1:E:173:SER:OG	2.18	0.44
2:F:123:THR:CG2	2:F:130:THR:HG23	2.48	0.44
2:H:60:ASP:OD1	2:H:60:ASP:N	2.51	0.44
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: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:C:142:LEU:HD12	1:C:142:LEU:N	2.32	0.43
1:E:85:MET:HE3	1:E:86:ASN:H	1.83	0.43
1:C:103:LEU:HB3	2:D:32:VAL:HB	1.99	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
1:E:165:ALA:HB2	1:E:175:VAL:HG11	2.00	0.43
2:F:56:ILE:HD13	2:F:154:PHE:CE1	2.53	0.43
2:H:77:ALA:HB1	2:H:83:THR:HA	2.00	0.43
1:A:25:ASN:O	1:A:60:LYS:HE2	2.19	0.43
2:D:82:HIS:HA	5:D:161:TRS:H12	2.01	0.43
1:A:75:LEU:HD23	1:A:115:TYR:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:31:THR:HG22	2:D:32:VAL:N	2.33	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:E:135:ARG:NH2	1:E:182:GLY:C	2.77	0.43
2:B:105:ASP:OD2	2:B:105:ASP:C	2.62	0.43
1:C:132:ARG:HH11	1:C:203:VAL:HB	1.84	0.43
2:D:137:TYR:CE2	2:D:148:ALA:HB1	2.54	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:A:128:ALA:C	1:A:130:LYS:H	2.26	0.42
1:A:103:LEU:HD12	2:B:150:ALA:O	2.20	0.42
1:C:79:ARG:HB3	1:C:170:MET:CE	2.50	0.42
1:E:134:ARG:O	1:E:141:THR:N	2.51	0.42
1:A:8:ARG:HH22	2:B:159:GLN:C	2.27	0.42
1:E:183:SER:CB	1:E:184:ASN:HA	2.49	0.42
1:C:107:ILE:CD1	2:D:154:PHE:HB3	2.49	0.42
2:D:103:ILE:C	2:D:104:LEU:HD23	2.45	0.42
2:H:98:ASN:CA	2:H:140:THR:HG23	2.48	0.42
1:A:79:ARG:HB3	1:A:170:MET:CE	2.50	0.42
1:E:24:ASN:OD1	1:E:25:ASN:N	2.53	0.42
2:H:98:ASN:HA	2:H:140:THR:HG23	2.01	0.42
1:E:169:PRO:O	1:E:170:MET:HB2	2.19	0.42
1:C:150:LEU:HD23	1:C:150:LEU:HA	1.92	0.42
1:G:35:SER:HA	1:G:86:ASN:O	2.20	0.42
2:H:99:VAL:HA	2:H:138:PHE:O	2.20	0.42
2:F:103:ILE:C	2:F:104:LEU:HD23	2.45	0.42
1:A:2:VAL:HA	1:A:23:THR:O	2.20	0.41
1:A:82:LEU:HD13	1:A:114:TYR:HE1	1.85	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:E:143:ILE:HG22	1:E:144:ASN:N	2.35	0.41
1:A:35:SER:HA	1:A:86:ASN:O	2.19	0.41
2:B:43:ALA:O	2:B:142:ALA:HA	2.19	0.41
1:C:2:VAL:HA	1:C:23:THR:O	2.20	0.41
1:C:79:ARG:HB3	1:C:170:MET:HE1	2.03	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:A:102:THR:CB	2:B:31:THR:CG2	2.97	0.41
1:A:131:LEU:O	1:A:132:ARG:HD3	2.20	0.41
2:B:77:ALA:HB1	2:B:83:THR:HA	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:104:LEU:HA	2:B:109:ALA:O	2.20	0.41
2:B:105:ASP:C	2:B:107:THR:H	2.29	0.41
1:C:47:ARG:HD2	1:C:83:PHE:HZ	1.85	0.41
1:C:60:LYS:O	1:C:61:LYS:HG3	2.20	0.41
2:D:37:VAL:HG22	2:D:137:TYR:CE1	2.55	0.41
1:G:55:PHE:HE2	1:G:65:LEU:HD21	1.85	0.41
1:A:27:GLU:HA	1:A:60:LYS:N	2.35	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:34:GLN:O	1:A:34:GLN:CG	2.69	0.41
1:A:47:ARG:NH2	2:D:107:THR:O	2.54	0.41
1:A:132:ARG:HH11	1:A:203:VAL:HB	1.86	0.41
1:A:193:TYR:HB3	2:B:17:VAL:CA	2.43	0.41
2:B:84:ASN:OD1	2:B:84:ASN:C	2.64	0.41
1:C:34:GLN:O	1:C:34:GLN:CG	2.69	0.41
1:E:102:THR:HG22	2:F:33:GLN:HE21	1.85	0.41
1:E:186:THR:OG1	1:E:199:LYS:HE3	2.21	0.41
2:F:104:LEU:HA	2:F:109:ALA:O	2.20	0.41
1:C:135:ARG:HG2	1:C:136:SER:N	2.36	0.41
1:E:25:ASN:O	1:E:26:ASP:C	2.64	0.41
1:A:79:ARG:HB3	1:A:170:MET:HE1	2.03	0.40
1:C:135:ARG:HA	1:C:140:LEU:HD23	2.01	0.40
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
2:F:104:LEU:HD23	2:F:104:LEU:N	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
2:B:142:ALA:HB1	3:B:160:PEG:H41	2.03	0.40
1:C:35:SER:HA	1:C:86:ASN:O	2.21	0.40
1:G:3:ALA:HB2	2:H:21:CYS:O	2.21	0.40
1:A:193:TYR:C	1:A:195:ALA:N	2.79	0.40
2:D:64:ASN:O	2:D:68:LYS:HG2	2.21	0.40
1:A:143:ILE:HG23	1:A:172:GLU:HG2	2.04	0.40
1:C:3:ALA:HB2	2:D:22:ALA:HA	2.02	0.40
1:C:155:LEU:HD23	1:C:162:LEU:HB2	2.03	0.40
1:E:27:GLU:HG2	1:E:60:LYS:CG	2.52	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%)	18	52
2	D	143/152 (94%)	131 (92%)	12 (8%)	0	100	100
2	F	79/152 (52%)	73 (92%)	4 (5%)	2 (2%)	4	27
2	H	60/152 (40%)	55 (92%)	5 (8%)	0	100	100
All	All	1028/1452 (71%)	947 (92%)	78 (8%)	3 (0%)	36	68

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%)	139 (93%)	10 (7%)	15	47
1	C	166/182 (91%)	154 (93%)	12 (7%)	13	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	104/182 (57%)	96 (92%)	8 (8%)	12	41
1	G	34/182 (19%)	31 (91%)	3 (9%)	9	35
2	B	102/109 (94%)	93 (91%)	9 (9%)	9	35
2	D	99/109 (91%)	88 (89%)	11 (11%)	6	25
2	F	54/109 (50%)	46 (85%)	8 (15%)	3	15
2	H	31/109 (28%)	30 (97%)	1 (3%)	34	65
All	All	739/1164 (64%)	677 (92%)	62 (8%)	10	38

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	35	SER
1	A	50	VAL
1	A	64	THR
1	A	73	ASN
1	A	75	LEU
1	A	146	THR
1	A	152	VAL
1	A	155	LEU
1	A	167	VAL
2	B	23	VAL
2	B	37	VAL
2	B	42	LEU
2	B	48	THR
2	B	65	VAL
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	73	ASN
1	C	140	LEU
1	C	146	THR
1	C	152	VAL
1	C	155	LEU
1	C	161	VAL
1	C	167	VAL
1	C	183	SER

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Mol	Chain	Res	Type
1	C	201	THR
1	C	206	HIS
2	D	15	GLU
2	D	16	VAL
2	D	23	VAL
2	D	29	ASP
2	D	42	LEU
2	D	48	THR
2	D	65	VAL
2	D	83	THR
2	D	104	LEU
2	D	129	ASN
2	D	144	THR
1	E	2	VAL
1	E	35	SER
1	E	50	VAL
1	E	64	THR
1	E	82	LEU
1	E	167	VAL
1	E	190	ILE
1	E	203	VAL
2	F	23	VAL
2	F	58	LEU
2	F	60	ASP
2	F	65	VAL
2	F	104	LEU
2	F	121	GLU
2	F	122	THR
2	F	129	ASN
1	G	2	VAL
1	G	50	VAL
1	G	64	THR
2	H	140	THR

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	164	ASN
2	B	125	ASN
2	B	149	ASN
2	D	33	GLN
2	D	129	ASN

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Mol	Chain	Res	Type
2	H	98	ASN

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](#)

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	$\# Z > 2$
5	TRS	D	161	-	7,7,7	0.32	0	9,9,9	0.32	0
5	TRS	C	213	-	7,7,7	0.31	0	9,9,9	0.27	0
3	PEG	B	160	-	6,6,6	0.63	0	5,5,5	1.43	0
3	PEG	A	213	-	6,6,6	0.65	0	5,5,5	1.39	0
4	EDO	D	160	-	3,3,3	0.52	0	2,2,2	0.11	0
3	PEG	A	212	-	6,6,6	0.53	0	5,5,5	1.71	1 (20%)
4	EDO	C	212	-	3,3,3	0.47	0	2,2,2	0.32	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
5	TRS	D	161	-	-	3/9/9/9	-
5	TRS	C	213	-	-	0/9/9/9	-
3	PEG	B	160	-	-	0/4/4/4	-
3	PEG	A	213	-	-	1/4/4/4	-
4	EDO	D	160	-	-	1/1/1/1	-
3	PEG	A	212	-	-	3/4/4/4	-
4	EDO	C	212	-	-	0/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	212	PEG	O2-C2-C1	2.41	120.75	110.11

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	212	PEG	O2-C3-C4-O4
4	D	160	EDO	O1-C1-C2-O2
3	A	212	PEG	O1-C1-C2-O2
3	A	212	PEG	C4-C3-O2-C2
5	D	161	TRS	C1-C-C2-O2
5	D	161	TRS	C3-C-C2-O2
5	D	161	TRS	N-C-C2-O2
3	A	213	PEG	C1-C2-O2-C3

There are no ring outliers.

5 monomers are involved in 10 short contacts:

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

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

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%)	0.44	12 (6%) 26 17	46, 84, 154, 191	0
1	C	200/211 (94%)	0.20	3 (1%) 72 52	41, 71, 120, 151	0
1	E	176/211 (83%)	1.09	34 (19%) 3 2	86, 138, 191, 234	0
1	G	76/211 (36%)	1.60	20 (26%) 1 2	116, 158, 198, 257	0
2	B	145/152 (95%)	0.31	12 (8%) 17 11	41, 69, 127, 206	0
2	D	145/152 (95%)	0.42	9 (6%) 26 17	46, 77, 126, 196	0
2	F	89/152 (58%)	1.34	23 (25%) 1 2	90, 136, 188, 223	0
2	H	78/152 (51%)	1.88	26 (33%) 1 1	95, 153, 206, 220	0
All	All	1100/1452 (75%)	0.74	139 (12%) 8 6	41, 99, 184, 257	0

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	135	ALA	9.6
2	H	137	TYR	7.5
2	H	156	VAL	7.3
2	H	138	PHE	6.2
2	H	134	GLN	5.9
2	H	85	VAL	5.6
2	F	29	ASP	4.9
2	H	101	VAL	4.5
1	G	114	TYR	4.5
1	G	56	ALA	4.4
1	G	45	ASP	4.4
1	G	5	GLY	4.4
2	F	129	ASN	4.4
1	E	94	ASP	4.3
1	A	187	TYR	3.9
1	G	34	GLN	3.9

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Mol	Chain	Res	Type	RSRZ
1	E	140	LEU	3.7
2	F	124	LEU	3.6
2	F	50	SER	3.4
2	B	17	VAL	3.4
1	G	72	ASN	3.3
2	H	72	ALA	3.3
1	C	53	PRO	3.3
2	D	108	GLY	3.3
2	H	21	CYS	3.3
1	E	197	THR	3.3
1	E	67	ILE	3.3
2	H	25	ALA	3.2
2	B	28	VAL	3.2
1	G	84	TRP	3.2
2	D	115	GLY	3.2
2	H	34	LEU	3.2
1	E	203	VAL	3.2
1	E	129	GLU	3.1
2	F	30	GLN	3.1
1	E	85	MET	3.1
2	F	48	THR	3.1
2	F	119	SER	3.0
1	G	47	ARG	3.0
2	H	103	ILE	3.0
1	E	31	TYR	3.0
2	D	89	GLN	3.0
2	H	154	PHE	3.0
1	E	153	THR	3.0
1	A	127	ALA	3.0
2	H	33	GLN	3.0
2	B	147	ALA	2.9
2	F	25	ALA	2.9
1	A	93	MET	2.9
2	H	102	GLN	2.9
2	F	64	ASN	2.9
2	B	94	GLY	2.9
2	F	118	PHE	2.9
2	F	110	ALA	2.9
1	E	109	SER	2.8
1	E	198	PRO	2.8
1	E	154	GLU	2.8
1	A	201	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	G	20	LEU	2.8
1	G	32	LEU	2.8
1	A	14	GLY	2.7
1	G	88	LYS	2.7
2	D	134	GLN	2.7
1	A	99	THR	2.7
1	G	61	LYS	2.7
1	E	131	LEU	2.7
1	C	100	GLU	2.7
1	G	65	LEU	2.7
1	E	19	GLN	2.6
2	H	133	PHE	2.6
1	G	115	TYR	2.6
2	H	84	ASN	2.6
1	A	180	ASP	2.6
1	A	81	SER	2.6
1	A	19	GLN	2.6
1	E	101	ASN	2.5
1	G	83	PHE	2.5
1	E	2	VAL	2.5
1	E	89	ALA	2.5
1	E	28	ASN	2.5
2	B	29	ASP	2.5
2	F	74	LEU	2.5
2	B	16	VAL	2.5
2	D	44	GLN	2.5
1	E	82	LEU	2.5
1	C	126	GLN	2.4
2	H	157	GLN	2.4
1	E	62	GLU	2.4
1	G	48	PHE	2.4
1	A	173	SER	2.4
2	D	90	SER	2.4
1	E	58	LYS	2.4
2	B	57	GLN	2.4
2	F	102	GLN	2.4
2	B	129	ASN	2.3
2	D	36	GLN	2.3
2	H	23	VAL	2.3
2	B	115	GLY	2.3
2	F	73	PHE	2.3
1	A	163	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	158	GLY	2.3
1	G	46	GLY	2.3
1	G	87	VAL	2.3
1	G	109	SER	2.3
2	F	109	ALA	2.3
1	E	103	LEU	2.3
2	B	123	THR	2.3
2	F	49	SER	2.3
2	B	91	SER	2.2
2	H	99	VAL	2.2
2	B	24	ASP	2.2
2	F	130	THR	2.2
1	E	65	LEU	2.2
2	F	105	ASP	2.2
1	E	157	ALA	2.2
2	D	19	ALA	2.2
2	H	96	ALA	2.2
2	H	24	ASP	2.2
1	G	22	VAL	2.2
2	H	111	LEU	2.2
1	E	46	GLY	2.2
1	A	189	THR	2.1
1	E	26	ASP	2.1
2	F	77	ALA	2.1
2	F	115	GLY	2.1
1	E	49	ILE	2.1
2	F	132	PRO	2.1
1	E	50	VAL	2.1
2	H	148	ALA	2.1
2	H	44	GLN	2.1
1	E	116	ARG	2.1
1	E	93	MET	2.0
1	E	11	TYR	2.0
2	F	116	ALA	2.0
2	H	61	CYS	2.0
2	F	26	GLY	2.0
2	D	92	ALA	2.0
1	E	112	LYS	2.0
1	E	166	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
5	TRS	D	161	8/8	0.50	0.23	140,140,140,140	0
4	EDO	C	212	4/4	0.59	0.18	109,109,109,109	0
3	PEG	B	160	7/7	0.61	0.26	148,148,148,148	0
5	TRS	C	213	8/8	0.63	0.14	138,138,138,138	0
4	EDO	D	160	4/4	0.79	0.16	77,77,77,77	0
3	PEG	A	212	7/7	0.89	0.18	67,67,67,67	0
3	PEG	A	213	7/7	0.90	0.17	111,111,111,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.