



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:04 PM UTC

PDB ID : 5XW7 / pdb_00005xw7
Title : Crystal structure of the flexible tandem repeat domain of bacterial cellulose synthase subunit C
Authors : Nojima, S.; Kato, K.; Yao, M.
Deposited on : 2017-06-29
Resolution : 3.27 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

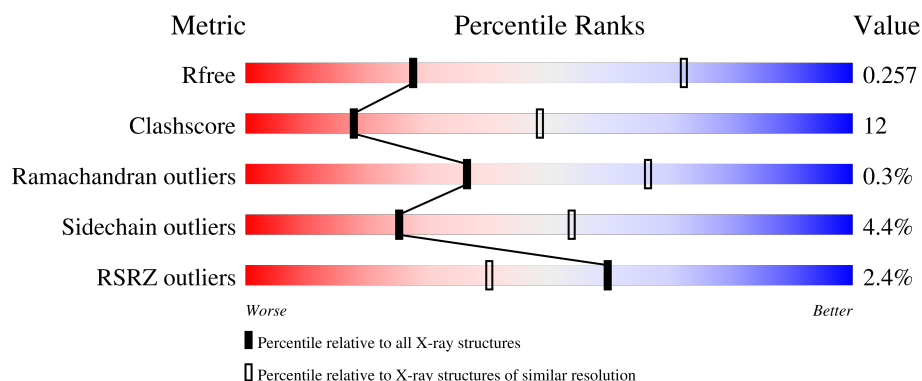
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.27 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	1303 (3.30-3.26)
Clashscore	190562	1354 (3.30-3.26)
Ramachandran outliers	187476	1334 (3.30-3.26)
Sidechain outliers	187428	1333 (3.30-3.26)
RSRZ outliers	180081	1303 (3.30-3.26)

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	260	0% 69% 23% 7%
1	B	260	3% 58% 24% 16%
1	C	260	2% 66% 25% 6%
1	D	260	2% 67% 23% 7%
1	E	260	3% 55% 18% 23%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8740 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 Cellulose synthase subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1858	1151	343	360	4			
1	B	218	Total	C	N	O	S	0	0	0
			1645	1018	305	318	4			
1	C	245	Total	C	N	O	S	0	0	0
			1870	1159	345	362	4			
1	D	241	Total	C	N	O	S	0	0	0
			1843	1142	341	356	4			
1	E	199	Total	C	N	O	S	0	0	0
			1524	945	282	294	3			

There are 55 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	GLY	-	expression tag	UNP K0J1W8
A	22	HIS	-	expression tag	UNP K0J1W8
A	23	MET	-	expression tag	UNP K0J1W8
A	273	LEU	-	expression tag	UNP K0J1W8
A	274	GLU	-	expression tag	UNP K0J1W8
A	275	HIS	-	expression tag	UNP K0J1W8
A	276	HIS	-	expression tag	UNP K0J1W8
A	277	HIS	-	expression tag	UNP K0J1W8
A	278	HIS	-	expression tag	UNP K0J1W8
A	279	HIS	-	expression tag	UNP K0J1W8
A	280	HIS	-	expression tag	UNP K0J1W8
B	21	GLY	-	expression tag	UNP K0J1W8
B	22	HIS	-	expression tag	UNP K0J1W8
B	23	MET	-	expression tag	UNP K0J1W8
B	273	LEU	-	expression tag	UNP K0J1W8
B	274	GLU	-	expression tag	UNP K0J1W8
B	275	HIS	-	expression tag	UNP K0J1W8
B	276	HIS	-	expression tag	UNP K0J1W8
B	277	HIS	-	expression tag	UNP K0J1W8

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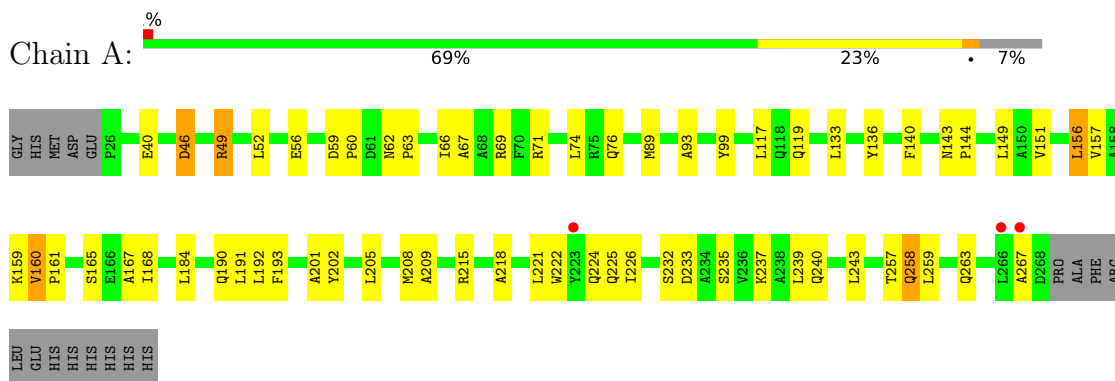
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Chain	Residue	Modelled	Actual	Comment	Reference
B	278	HIS	-	expression tag	UNP K0J1W8
B	279	HIS	-	expression tag	UNP K0J1W8
B	280	HIS	-	expression tag	UNP K0J1W8
C	21	GLY	-	expression tag	UNP K0J1W8
C	22	HIS	-	expression tag	UNP K0J1W8
C	23	MET	-	expression tag	UNP K0J1W8
C	273	LEU	-	expression tag	UNP K0J1W8
C	274	GLU	-	expression tag	UNP K0J1W8
C	275	HIS	-	expression tag	UNP K0J1W8
C	276	HIS	-	expression tag	UNP K0J1W8
C	277	HIS	-	expression tag	UNP K0J1W8
C	278	HIS	-	expression tag	UNP K0J1W8
C	279	HIS	-	expression tag	UNP K0J1W8
C	280	HIS	-	expression tag	UNP K0J1W8
D	21	GLY	-	expression tag	UNP K0J1W8
D	22	HIS	-	expression tag	UNP K0J1W8
D	23	MET	-	expression tag	UNP K0J1W8
D	273	LEU	-	expression tag	UNP K0J1W8
D	274	GLU	-	expression tag	UNP K0J1W8
D	275	HIS	-	expression tag	UNP K0J1W8
D	276	HIS	-	expression tag	UNP K0J1W8
D	277	HIS	-	expression tag	UNP K0J1W8
D	278	HIS	-	expression tag	UNP K0J1W8
D	279	HIS	-	expression tag	UNP K0J1W8
D	280	HIS	-	expression tag	UNP K0J1W8
E	21	GLY	-	expression tag	UNP K0J1W8
E	22	HIS	-	expression tag	UNP K0J1W8
E	23	MET	-	expression tag	UNP K0J1W8
E	273	LEU	-	expression tag	UNP K0J1W8
E	274	GLU	-	expression tag	UNP K0J1W8
E	275	HIS	-	expression tag	UNP K0J1W8
E	276	HIS	-	expression tag	UNP K0J1W8
E	277	HIS	-	expression tag	UNP K0J1W8
E	278	HIS	-	expression tag	UNP K0J1W8
E	279	HIS	-	expression tag	UNP K0J1W8
E	280	HIS	-	expression tag	UNP K0J1W8

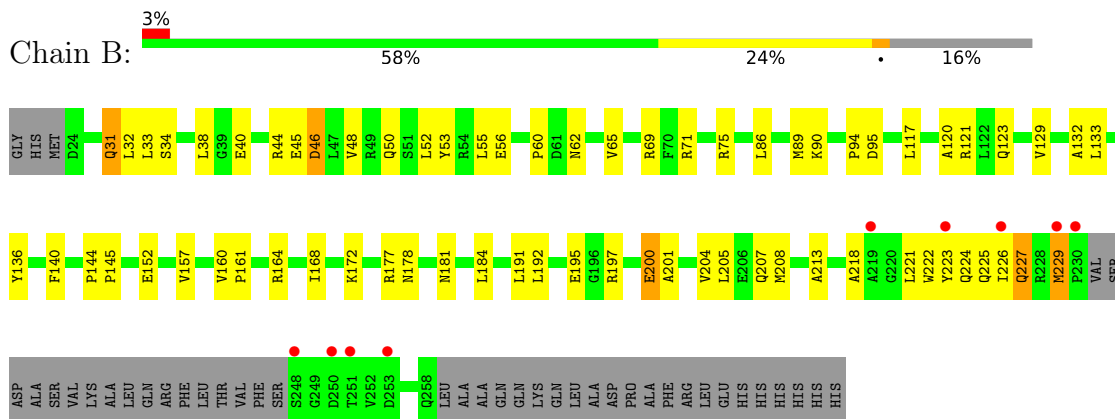
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

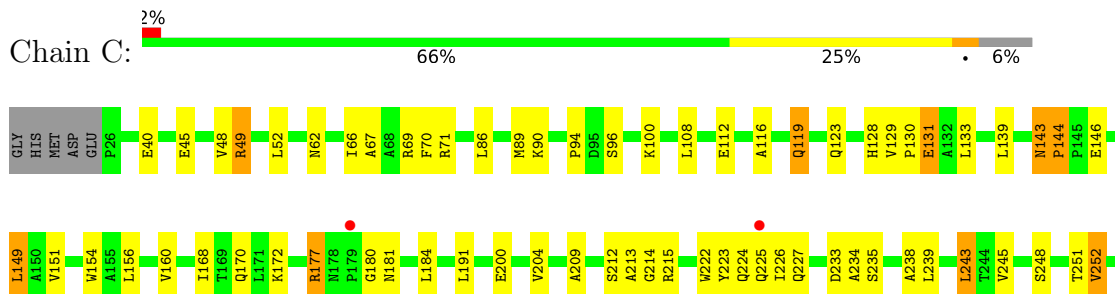
• Molecule 1: Cellulose synthase subunit C

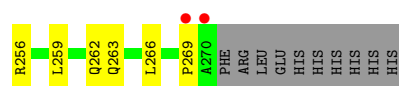


• Molecule 1: Cellulose synthase subunit C

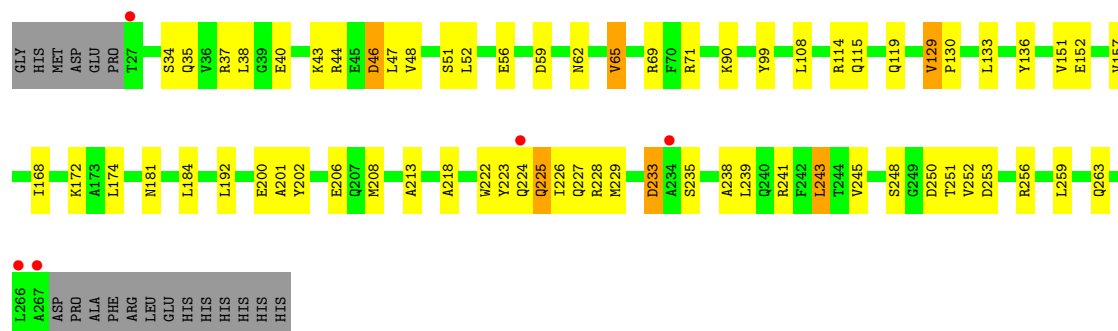


• Molecule 1: Cellulose synthase subunit C

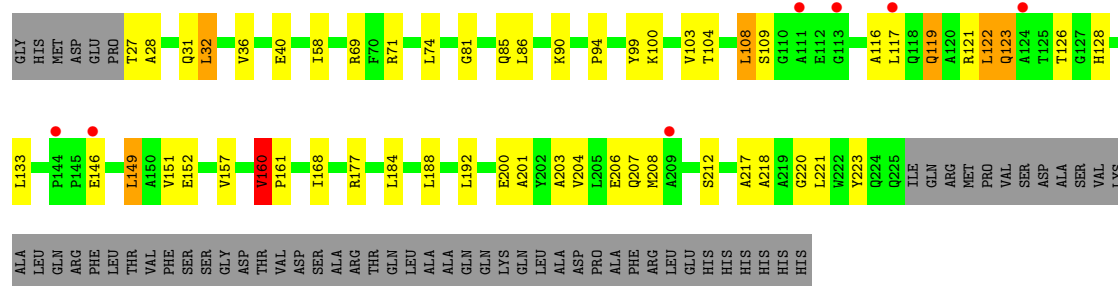




• Molecule 1: Cellulose synthase subunit C



• Molecule 1: Cellulose synthase subunit C



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	115.81Å 55.62Å 165.02Å 90.00° 95.56° 90.00°	Depositor
Resolution (Å)	47.70 – 3.27 47.70 – 3.27	Depositor EDS
% Data completeness (in resolution range)	98.6 (47.70-3.27) 98.6 (47.70-3.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.212 , 0.257 0.216 , 0.257	Depositor DCC
R_{free} test set	2222 reflections (6.73%)	wwPDB-VP
Wilson B-factor (Å ²)	81.7	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8740	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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	0.65	0/1884	1.07	11/2549 (0.4%)
1	B	0.65	0/1668	1.07	5/2258 (0.2%)
1	C	0.67	0/1897	1.04	9/2568 (0.4%)
1	D	0.66	0/1868	1.08	6/2527 (0.2%)
1	E	0.65	0/1546	1.13	6/2092 (0.3%)
All	All	0.66	0/8863	1.08	37/11994 (0.3%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	225	GLN	N-CA-C	-9.32	102.04	113.97
1	D	224	GLN	N-CA-C	-9.03	98.92	113.19
1	D	251	THR	N-CA-C	-8.98	101.57	111.36
1	C	224	GLN	N-CA-C	-8.57	100.81	114.09
1	A	143	ASN	CA-C-N	8.07	125.47	119.66
1	A	143	ASN	C-N-CA	8.07	125.47	119.66
1	A	59	ASP	CA-C-N	7.62	127.27	119.19
1	A	59	ASP	C-N-CA	7.62	127.27	119.19
1	A	224	GLN	N-CA-C	-7.02	102.09	113.19
1	E	32	LEU	N-CA-C	6.99	120.28	111.69
1	A	160	VAL	CA-C-N	6.98	126.79	119.05
1	A	160	VAL	C-N-CA	6.98	126.79	119.05
1	E	108	LEU	N-CA-C	6.88	121.71	113.12
1	C	144	PRO	O-C-N	6.80	124.35	121.15
1	C	143	ASN	CA-C-N	6.78	124.54	119.66
1	C	143	ASN	C-N-CA	6.78	124.54	119.66
1	A	144	PRO	O-C-N	6.61	124.25	121.15
1	E	212	SER	N-CA-C	6.09	117.92	111.28
1	A	225	GLN	N-CA-C	-6.07	106.20	113.97
1	D	225	GLN	N-CA-C	-6.07	106.49	114.31
1	B	160	VAL	CA-C-N	5.96	125.66	119.05
1	B	160	VAL	C-N-CA	5.96	125.66	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	129	VAL	CB-CA-C	-5.80	108.20	114.35
1	E	58	ILE	N-CA-C	5.78	116.42	110.82
1	B	178	ASN	CA-C-N	-5.77	114.31	120.03
1	B	178	ASN	C-N-CA	-5.77	114.31	120.03
1	C	251	THR	N-CA-C	-5.66	105.19	111.36
1	D	129	VAL	CB-CA-C	-5.44	108.58	114.35
1	C	180	GLY	N-CA-C	5.40	125.98	113.18
1	A	93	ALA	CA-C-N	5.33	124.95	119.56
1	A	93	ALA	C-N-CA	5.33	124.95	119.56
1	D	59	ASP	CA-C-N	5.28	124.90	119.56
1	D	59	ASP	C-N-CA	5.28	124.90	119.56
1	C	160	VAL	CA-C-N	5.20	125.25	119.32
1	C	160	VAL	C-N-CA	5.20	125.25	119.32
1	E	160	VAL	CA-C-N	5.02	124.62	119.05
1	E	160	VAL	C-N-CA	5.02	124.62	119.05

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1858	0	1863	37	0
1	B	1645	0	1612	52	0
1	C	1870	0	1875	48	0
1	D	1843	0	1851	51	0
1	E	1524	0	1522	40	0
All	All	8740	0	8723	214	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 12.

All (214) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:VAL:HA	1:B:207:GLN:HG2	1.50	0.92
1:A:226:ILE:HG21	1:A:239:LEU:HD13	1.58	0.86
1:B:62:ASN:HB3	1:B:65:VAL:HG12	1.66	0.77
1:E:151:VAL:HG22	1:E:184:LEU:HD12	1.68	0.76
1:E:121:ARG:NH1	1:E:152:GLU:OE2	2.20	0.74
1:E:117:LEU:HD11	1:E:149:LEU:HD11	1.68	0.74
1:B:34:SER:OG	1:D:44:ARG:NH2	2.21	0.73
1:C:49:ARG:NH2	1:D:40:GLU:OE1	2.20	0.71
1:B:140:PHE:CD1	1:B:145:PRO:HG3	2.26	0.70
1:D:243:LEU:HD21	1:D:256:ARG:HG2	1.74	0.68
1:B:31:GLN:HA	1:B:31:GLN:HE21	1.58	0.68
1:D:250:ASP:OD1	1:D:253:ASP:HB2	1.94	0.68
1:C:119:GLN:HG3	1:D:213:ALA:HA	1.75	0.67
1:A:209:ALA:O	1:A:215:ARG:NH1	2.27	0.66
1:B:227:GLN:O	1:B:227:GLN:NE2	2.28	0.66
1:B:56:GLU:O	1:B:60:PRO:HG3	1.96	0.65
1:A:151:VAL:HG13	1:A:184:LEU:HD12	1.79	0.65
1:D:208:MET:HE2	1:D:218:ALA:HB2	1.79	0.64
1:D:226:ILE:HD12	1:D:238:ALA:HB1	1.81	0.62
1:E:123:GLN:HG2	1:E:128:HIS:O	2.00	0.62
1:E:123:GLN:HG3	1:E:128:HIS:HB2	1.82	0.62
1:C:233:ASP:OD1	1:C:234:ALA:N	2.33	0.61
1:C:226:ILE:HD12	1:C:238:ALA:HB1	1.83	0.61
1:B:32:LEU:HB3	1:B:55:LEU:HG	1.84	0.60
1:C:213:ALA:HA	1:D:119:GLN:NE2	2.17	0.60
1:C:214:GLY:H	1:D:119:GLN:HE22	1.49	0.60
1:C:181:ASN:HB3	1:C:184:LEU:HB3	1.82	0.59
1:B:205:LEU:HD22	1:B:222:TRP:CE3	2.36	0.59
1:D:90:LYS:HB2	1:D:99:TYR:CE1	2.37	0.59
1:A:52:LEU:HB3	1:A:69:ARG:HH12	1.68	0.59
1:C:151:VAL:HG13	1:C:184:LEU:HD13	1.85	0.59
1:C:116:ALA:HB1	1:C:139:LEU:HD21	1.85	0.59
1:B:197:ARG:HD3	1:B:200:GLU:OE2	2.02	0.59
1:B:133:LEU:HD11	1:B:157:VAL:HA	1.84	0.58
1:B:31:GLN:OE1	1:D:46:ASP:HB3	2.03	0.58
1:A:232:SER:H	1:A:235:SER:HB3	1.69	0.58
1:D:56:GLU:OE2	1:D:69:ARG:HD2	2.03	0.57
1:B:136:TYR:OH	1:B:152:GLU:OE2	2.20	0.57
1:D:202:TYR:OH	1:D:225:GLN:HG2	2.04	0.57
1:C:90:LYS:O	1:C:94:PRO:HG3	2.04	0.57
1:C:146:GLU:OE2	1:C:177:ARG:NH2	2.31	0.57
1:D:62:ASN:HB3	1:D:65:VAL:HG13	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:GLU:OE2	1:C:71:ARG:NH1	2.38	0.56
1:C:259:LEU:O	1:C:263:GLN:HG2	2.06	0.56
1:D:192:LEU:HD13	1:D:200:GLU:HB2	1.88	0.56
1:C:154:TRP:CE2	1:C:170:GLN:HG2	2.41	0.55
1:B:181:ASN:HB3	1:B:184:LEU:HB3	1.88	0.55
1:C:62:ASN:O	1:C:66:ILE:HG13	2.05	0.55
1:C:209:ALA:O	1:C:215:ARG:NH1	2.40	0.55
1:A:136:TYR:HD2	1:A:140:PHE:CE1	2.25	0.54
1:B:45:GLU:OE2	1:B:75:ARG:NH1	2.41	0.54
1:E:122:LEU:HD13	1:E:123:GLN:N	2.22	0.54
1:D:40:GLU:OE2	1:D:71:ARG:NH1	2.40	0.54
1:B:223:TYR:C	1:B:225:GLN:H	2.16	0.54
1:D:192:LEU:HB2	1:D:201:ALA:HB2	1.91	0.53
1:E:146:GLU:OE2	1:E:177:ARG:NH1	2.41	0.53
1:A:259:LEU:O	1:A:263:GLN:HG2	2.08	0.53
1:A:190:GLN:HG3	1:A:221:LEU:HD21	1.90	0.53
1:B:44:ARG:HD3	1:D:38:LEU:HD13	1.91	0.53
1:C:235:SER:O	1:C:239:LEU:HD12	2.09	0.52
1:C:262:GLN:O	1:C:266:LEU:HD12	2.09	0.52
1:C:52:LEU:HD21	1:C:69:ARG:HA	1.91	0.52
1:C:168:ILE:O	1:C:172:LYS:HG3	2.10	0.52
1:A:40:GLU:OE2	1:A:71:ARG:NH1	2.43	0.51
1:A:117:LEU:HD11	1:A:149:LEU:HD21	1.91	0.51
1:C:222:TRP:CH2	1:C:245:VAL:HG21	2.45	0.51
1:C:212:SER:HA	1:C:215:ARG:HH21	1.75	0.51
1:A:46:ASP:OD1	1:B:71:ARG:HD2	2.11	0.51
1:A:208:MET:HE2	1:A:218:ALA:HB2	1.91	0.51
1:B:226:ILE:O	1:B:229:MET:HB2	2.10	0.51
1:C:96:SER:O	1:C:100:LYS:HG3	2.10	0.51
1:D:222:TRP:HD1	1:D:222:TRP:O	1.92	0.51
1:B:223:TYR:O	1:B:225:GLN:N	2.44	0.51
1:B:38:LEU:HD22	1:D:44:ARG:HD3	1.93	0.51
1:B:223:TYR:O	1:B:223:TYR:HD1	1.94	0.51
1:D:229:MET:SD	1:D:235:SER:HA	2.51	0.51
1:B:208:MET:HE2	1:B:218:ALA:HB2	1.93	0.51
1:D:222:TRP:O	1:D:222:TRP:CD1	2.64	0.50
1:B:227:GLN:HE21	1:B:227:GLN:C	2.19	0.50
1:D:248:SER:HA	1:D:252:VAL:HG11	1.92	0.50
1:D:62:ASN:O	1:D:65:VAL:HG22	2.11	0.50
1:E:119:GLN:C	1:E:122:LEU:HD12	2.36	0.50
1:B:90:LYS:O	1:B:94:PRO:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:GLU:O	1:C:204:VAL:HG23	2.12	0.50
1:A:237:LYS:HA	1:A:240:GLN:HG2	1.93	0.49
1:B:95:ASP:OD1	1:B:177:ARG:NH1	2.42	0.49
1:E:40:GLU:OE2	1:E:71:ARG:NH1	2.45	0.49
1:D:206:GLU:OE2	1:D:241:ARG:NH2	2.46	0.49
1:E:99:TYR:O	1:E:103:VAL:HG23	2.13	0.49
1:B:161:PRO:HA	1:B:164:ARG:HB2	1.94	0.49
1:C:222:TRP:O	1:C:222:TRP:CD1	2.65	0.49
1:C:128:HIS:HB3	1:C:131:GLU:CG	2.43	0.48
1:D:151:VAL:HG22	1:D:174:LEU:HD13	1.94	0.48
1:B:201:ALA:HA	1:B:204:VAL:HG22	1.95	0.48
1:E:133:LEU:HD21	1:E:157:VAL:HA	1.95	0.48
1:C:222:TRP:O	1:C:222:TRP:HD1	1.95	0.48
1:E:36:VAL:O	1:E:40:GLU:HG3	2.14	0.48
1:E:208:MET:HG2	1:E:208:MET:O	2.14	0.47
1:B:223:TYR:C	1:B:223:TYR:HD1	2.21	0.47
1:B:223:TYR:C	1:B:223:TYR:CD1	2.92	0.47
1:B:44:ARG:NE	1:B:46:ASP:OD2	2.43	0.47
1:A:56:GLU:OE2	1:A:69:ARG:HD2	2.15	0.47
1:A:239:LEU:HD21	1:A:258:GLN:HG2	1.96	0.47
1:B:120:ALA:HB1	1:B:136:TYR:CE2	2.50	0.47
1:C:108:LEU:HD21	1:C:139:LEU:HD12	1.97	0.47
1:A:222:TRP:HD1	1:A:222:TRP:O	1.97	0.47
1:B:52:LEU:HD21	1:B:69:ARG:HA	1.97	0.47
1:B:192:LEU:HB2	1:B:201:ALA:HB2	1.96	0.47
1:D:181:ASN:HB3	1:D:184:LEU:HB3	1.95	0.47
1:D:259:LEU:O	1:D:263:GLN:HG2	2.15	0.47
1:E:192:LEU:CB	1:E:201:ALA:HB2	2.45	0.46
1:B:50:GLN:O	1:B:53:TYR:HB3	2.16	0.46
1:E:28:ALA:O	1:E:32:LEU:HD12	2.15	0.46
1:E:217:ALA:O	1:E:221:LEU:HD13	2.16	0.46
1:E:168:ILE:HD13	1:E:192:LEU:HD12	1.97	0.46
1:A:133:LEU:HB2	1:A:156:LEU:HD12	1.98	0.46
1:D:151:VAL:HG13	1:D:184:LEU:HD12	1.98	0.46
1:E:123:GLN:HE21	1:E:128:HIS:HB2	1.81	0.46
1:E:203:ALA:O	1:E:207:GLN:HG3	2.17	0.45
1:D:223:TYR:O	1:D:227:GLN:HG2	2.17	0.45
1:D:34:SER:HA	1:D:37:ARG:NH2	2.31	0.45
1:E:188:LEU:O	1:E:192:LEU:HD13	2.16	0.45
1:D:233:ASP:OD1	1:D:233:ASP:N	2.50	0.45
1:E:81:GLY:O	1:E:85:GLN:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:263:GLN:HA	1:C:266:LEU:HD12	1.98	0.45
1:A:67:ALA:HB2	1:A:89:MET:HE1	1.98	0.45
1:B:168:ILE:HD11	1:B:191:LEU:HB3	1.99	0.45
1:C:131:GLU:CD	1:C:131:GLU:H	2.24	0.45
1:C:133:LEU:HD11	1:C:156:LEU:HG	1.99	0.45
1:D:47:LEU:O	1:D:51:SER:OG	2.26	0.45
1:C:149:LEU:HA	1:C:149:LEU:HD13	1.77	0.45
1:E:200:GLU:O	1:E:204:VAL:HG23	2.17	0.45
1:E:208:MET:HE3	1:E:218:ALA:HB2	1.98	0.44
1:E:28:ALA:C	1:E:32:LEU:HD12	2.42	0.44
1:A:160:VAL:HA	1:A:161:PRO:HD3	1.82	0.44
1:C:129:VAL:HB	1:C:130:PRO:HD3	1.97	0.44
1:A:133:LEU:HD13	1:A:156:LEU:HD12	1.99	0.44
1:E:100:LYS:O	1:E:104:THR:HG23	2.18	0.44
1:A:62:ASN:O	1:A:66:ILE:HG13	2.17	0.44
1:B:86:LEU:HA	1:B:89:MET:HE3	1.99	0.44
1:E:86:LEU:HD23	1:E:86:LEU:HA	1.88	0.44
1:A:193:PHE:CE1	1:A:201:ALA:HB3	2.53	0.44
1:C:214:GLY:H	1:D:119:GLN:NE2	2.15	0.44
1:A:89:MET:HE2	1:A:99:TYR:HA	2.00	0.44
1:D:168:ILE:O	1:D:172:LYS:HG3	2.18	0.44
1:C:154:TRP:NE1	1:C:170:GLN:HG2	2.33	0.44
1:C:248:SER:C	1:C:252:VAL:HG21	2.42	0.44
1:B:223:TYR:C	1:B:225:GLN:N	2.76	0.43
1:E:126:THR:HG22	1:E:128:HIS:CE1	2.52	0.43
1:B:32:LEU:HA	1:B:32:LEU:HD23	1.73	0.43
1:D:172:LYS:HE2	1:D:172:LYS:HB3	1.77	0.43
1:A:156:LEU:O	1:A:159:LYS:HB2	2.19	0.43
1:B:120:ALA:HB1	1:B:136:TYR:CD2	2.53	0.43
1:C:52:LEU:HD12	1:C:52:LEU:HA	1.83	0.43
1:E:74:LEU:HD23	1:E:74:LEU:HA	1.80	0.43
1:A:74:LEU:HA	1:A:74:LEU:HD23	1.77	0.43
1:A:233:ASP:OD1	1:A:233:ASP:N	2.52	0.43
1:B:34:SER:HG	1:D:44:ARG:HH22	1.63	0.43
1:C:133:LEU:CD1	1:C:156:LEU:HG	2.48	0.43
1:D:108:LEU:HD12	1:D:114:ARG:HA	2.01	0.43
1:D:241:ARG:O	1:D:245:VAL:HG23	2.19	0.43
1:B:227:GLN:HE21	1:B:227:GLN:CA	2.31	0.43
1:D:48:VAL:O	1:D:52:LEU:HB2	2.19	0.43
1:D:222:TRP:CH2	1:D:245:VAL:HG21	2.54	0.43
1:A:56:GLU:O	1:A:60:PRO:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:LEU:HD23	1:B:33:LEU:HA	1.85	0.42
1:A:156:LEU:HD22	1:A:156:LEU:HA	1.83	0.42
1:D:208:MET:HB3	1:D:218:ALA:HB2	2.00	0.42
1:A:205:LEU:HD23	1:A:205:LEU:HA	1.78	0.42
1:E:119:GLN:O	1:E:122:LEU:HD12	2.19	0.42
1:A:62:ASN:HA	1:A:63:PRO:HD2	1.88	0.42
1:B:168:ILE:O	1:B:172:LYS:HG3	2.20	0.42
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.71	0.42
1:C:45:GLU:O	1:C:49:ARG:HB2	2.19	0.42
1:E:116:ALA:HA	1:E:119:GLN:OE1	2.20	0.42
1:A:49:ARG:NH2	1:B:40:GLU:OE1	2.35	0.42
1:E:69:ARG:HG2	1:E:85:GLN:HG3	2.02	0.42
1:B:48:VAL:O	1:B:52:LEU:HB2	2.20	0.42
1:D:226:ILE:O	1:D:229:MET:HG2	2.19	0.42
1:B:123:GLN:HB2	1:B:132:ALA:HB2	2.02	0.42
1:B:191:LEU:HD23	1:B:191:LEU:HA	1.89	0.42
1:C:191:LEU:HD23	1:C:191:LEU:HA	1.82	0.42
1:D:252:VAL:O	1:D:256:ARG:HG3	2.20	0.42
1:B:40:GLU:OE2	1:B:71:ARG:NH1	2.53	0.42
1:C:139:LEU:HD13	1:C:139:LEU:HA	1.91	0.42
1:B:144:PRO:HA	1:B:145:PRO:HD3	1.92	0.41
1:E:192:LEU:HB2	1:E:201:ALA:HB2	2.00	0.41
1:C:243:LEU:HD21	1:C:256:ARG:HG2	2.01	0.41
1:C:67:ALA:O	1:C:70:PHE:HB3	2.20	0.41
1:A:192:LEU:HB2	1:A:201:ALA:HB2	2.03	0.41
1:C:223:TYR:O	1:C:227:GLN:HG2	2.20	0.41
1:D:43:LYS:HB3	1:D:43:LYS:HE2	1.80	0.41
1:E:220:GLY:O	1:E:223:TYR:HD2	2.03	0.41
1:C:212:SER:HB3	1:D:115:GLN:OE1	2.21	0.41
1:D:90:LYS:HB2	1:D:99:TYR:CZ	2.55	0.41
1:E:122:LEU:HD13	1:E:123:GLN:H	1.85	0.41
1:E:146:GLU:OE2	1:E:177:ARG:CZ	2.68	0.41
1:E:160:VAL:HA	1:E:161:PRO:HD3	1.92	0.41
1:A:157:VAL:HG12	1:A:167:ALA:HB2	2.02	0.41
1:E:90:LYS:O	1:E:94:PRO:HG3	2.20	0.41
1:B:117:LEU:O	1:B:121:ARG:HG2	2.21	0.41
1:C:86:LEU:HA	1:C:89:MET:HE3	2.03	0.41
1:C:143:ASN:HA	1:C:144:PRO:HD2	1.95	0.41
1:E:90:LYS:HB2	1:E:99:TYR:CZ	2.56	0.41
1:D:133:LEU:HD21	1:D:157:VAL:HA	2.02	0.40
1:D:136:TYR:OH	1:D:152:GLU:OE1	2.32	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ILE:HG12	1:A:191:LEU:HD13	2.02	0.40
1:D:129:VAL:N	1:D:130:PRO:HD2	2.36	0.40
1:A:202:TYR:HD2	1:A:202:TYR:HA	1.74	0.40
1:C:123:GLN:OE1	1:C:131:GLU:HB2	2.22	0.40
1:E:100:LYS:HE2	1:E:100:LYS:HB3	1.82	0.40
1:E:108:LEU:H	1:E:109:SER:HA	1.87	0.40
1:A:119:GLN:OE1	1:B:213:ALA:HA	2.22	0.40
1:D:239:LEU:HD22	1:D:259:LEU:HA	2.04	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	241/260 (93%)	235 (98%)	5 (2%)	1 (0%)	30	60
1	B	214/260 (82%)	208 (97%)	5 (2%)	1 (0%)	24	55
1	C	243/260 (94%)	236 (97%)	6 (2%)	1 (0%)	30	60
1	D	239/260 (92%)	232 (97%)	7 (3%)	0	100	100
1	E	197/260 (76%)	194 (98%)	3 (2%)	0	100	100
All	All	1134/1300 (87%)	1105 (97%)	26 (2%)	3 (0%)	36	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	224	GLN
1	A	267	ALA
1	C	269	PRO

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	188/203 (93%)	180 (96%)	8 (4%)	26	54
1	B	159/203 (78%)	152 (96%)	7 (4%)	25	54
1	C	189/203 (93%)	180 (95%)	9 (5%)	23	52
1	D	186/203 (92%)	180 (97%)	6 (3%)	34	60
1	E	151/203 (74%)	143 (95%)	8 (5%)	20	49
All	All	873/1015 (86%)	835 (96%)	38 (4%)	25	54

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

Mol	Chain	Res	Type
1	A	46	ASP
1	A	49	ARG
1	A	76	GLN
1	A	156	LEU
1	A	165	SER
1	A	243	LEU
1	A	257	THR
1	A	258	GLN
1	B	31	GLN
1	B	46	ASP
1	B	195	GLU
1	B	200	GLU
1	B	221	LEU
1	B	227	GLN
1	B	229	MET
1	C	48	VAL
1	C	49	ARG
1	C	112	GLU
1	C	119	GLN
1	C	131	GLU
1	C	149	LEU
1	C	177	ARG
1	C	243	LEU

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Mol	Chain	Res	Type
1	C	252	VAL
1	D	35	GLN
1	D	46	ASP
1	D	65	VAL
1	D	228	ARG
1	D	233	ASP
1	D	243	LEU
1	E	27	THR
1	E	31	GLN
1	E	119	GLN
1	E	122	LEU
1	E	123	GLN
1	E	149	LEU
1	E	160	VAL
1	E	206	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	115	GLN
1	A	123	GLN
1	B	31	GLN
1	B	79	ASN
1	B	91	GLN
1	B	227	GLN
1	C	265	GLN
1	D	119	GLN
1	D	190	GLN
1	E	30	GLN
1	E	85	GLN
1	E	123	GLN
1	E	128	HIS
1	E	181	ASN
1	E	186	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	243/260 (93%)	-0.06	3 (1%) 76 58	50, 73, 98, 126	0
1	B	218/260 (83%)	0.20	9 (4%) 41 27	46, 83, 139, 172	0
1	C	245/260 (94%)	-0.06	4 (1%) 70 52	53, 76, 99, 116	0
1	D	241/260 (92%)	-0.00	5 (2%) 63 44	46, 77, 116, 165	0
1	E	199/260 (76%)	0.35	7 (3%) 47 31	54, 89, 134, 146	0
All	All	1146/1300 (88%)	0.07	28 (2%) 59 40	46, 78, 127, 172	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	267	ALA	4.7
1	C	269	PRO	3.8
1	B	219	ALA	3.7
1	C	270	ALA	3.4
1	D	266	LEU	3.1
1	E	113	GLY	2.8
1	A	266	LEU	2.7
1	E	209	ALA	2.7
1	E	124	ALA	2.6
1	A	223	TYR	2.6
1	B	251	THR	2.6
1	B	230	PRO	2.6
1	D	27	THR	2.6
1	E	144	PRO	2.5
1	E	117	LEU	2.5
1	B	223	TYR	2.4
1	B	226	ILE	2.4
1	D	234	ALA	2.3
1	D	224	GLN	2.2
1	B	250	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	253	ASP	2.1
1	E	146	GLU	2.1
1	C	225	GLN	2.1
1	B	248	SER	2.1
1	B	229	MET	2.1
1	C	179	PRO	2.1
1	A	267	ALA	2.0
1	E	111	ALA	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.