



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 12:18 AM UTC

PDB ID : 7YKQ / pdb_00007ykq
Title : Crystal structure of a novel alpha/beta hydrolase mutant from thermomonospora curvata in apo form
Authors : Han, X.; Jian, G.; Bornscheuer, U.T.; Wei, R.; Liu, W.
Deposited on : 2022-07-23
Resolution : 2.36 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

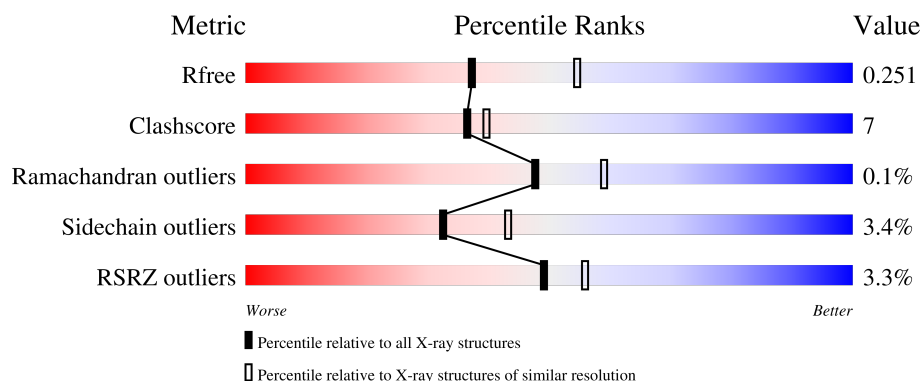
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	269	
1	B	269	
1	C	269	
1	D	269	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8231 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 Triacylglycerol lipase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	0	0
			1989	1252	352	379	6			
1	B	259	Total	C	N	O	S	0	0	0
			1989	1252	352	379	6			
1	C	259	Total	C	N	O	S	0	0	0
			1989	1252	352	379	6			
1	D	259	Total	C	N	O	S	0	0	0
			1989	1252	352	379	6			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	ASN	-	expression tag	UNP D1A9G5
A	-7	SER	-	expression tag	UNP D1A9G5
A	-6	MET	-	expression tag	UNP D1A9G5
A	-5	HIS	-	expression tag	UNP D1A9G5
A	-4	HIS	-	expression tag	UNP D1A9G5
A	-3	HIS	-	expression tag	UNP D1A9G5
A	-2	HIS	-	expression tag	UNP D1A9G5
A	-1	HIS	-	expression tag	UNP D1A9G5
A	0	HIS	-	expression tag	UNP D1A9G5
A	13	ALA	SER	engineered mutation	UNP D1A9G5
A	14	SER	LEU	engineered mutation	UNP D1A9G5
A	18	PRO	ALA	engineered mutation	UNP D1A9G5
A	41	THR	ARG	engineered mutation	UNP D1A9G5
A	47	ASP	THR	engineered mutation	UNP D1A9G5
A	92	TYR	LEU	engineered mutation	UNP D1A9G5
A	117	ASP	ASN	engineered mutation	UNP D1A9G5
A	141	ARG	LYS	engineered mutation	UNP D1A9G5
A	142	ARG	SER	engineered mutation	UNP D1A9G5
A	144	PRO	THR	engineered mutation	UNP D1A9G5
A	165	THR	ARG	engineered mutation	UNP D1A9G5
A	175	ASN	LEU	engineered mutation	UNP D1A9G5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	185	ALA	SER	engineered mutation	UNP D1A9G5
A	196	THR	ARG	engineered mutation	UNP D1A9G5
A	205	GLY	ASN	engineered mutation	UNP D1A9G5
A	213	THR	PHE	engineered mutation	UNP D1A9G5
A	214	PRO	SER	engineered mutation	UNP D1A9G5
A	224	ALA	SER	engineered mutation	UNP D1A9G5
A	246	PRO	ALA	engineered mutation	UNP D1A9G5
A	247	SER	ILE	engineered mutation	UNP D1A9G5
A	252	GLU	ASP	engineered mutation	UNP D1A9G5
B	-8	ASN	-	expression tag	UNP D1A9G5
B	-7	SER	-	expression tag	UNP D1A9G5
B	-6	MET	-	expression tag	UNP D1A9G5
B	-5	HIS	-	expression tag	UNP D1A9G5
B	-4	HIS	-	expression tag	UNP D1A9G5
B	-3	HIS	-	expression tag	UNP D1A9G5
B	-2	HIS	-	expression tag	UNP D1A9G5
B	-1	HIS	-	expression tag	UNP D1A9G5
B	0	HIS	-	expression tag	UNP D1A9G5
B	13	ALA	SER	engineered mutation	UNP D1A9G5
B	14	SER	LEU	engineered mutation	UNP D1A9G5
B	18	PRO	ALA	engineered mutation	UNP D1A9G5
B	41	THR	ARG	engineered mutation	UNP D1A9G5
B	47	ASP	THR	engineered mutation	UNP D1A9G5
B	92	TYR	LEU	engineered mutation	UNP D1A9G5
B	117	ASP	ASN	engineered mutation	UNP D1A9G5
B	141	ARG	LYS	engineered mutation	UNP D1A9G5
B	142	ARG	SER	engineered mutation	UNP D1A9G5
B	144	PRO	THR	engineered mutation	UNP D1A9G5
B	165	THR	ARG	engineered mutation	UNP D1A9G5
B	175	ASN	LEU	engineered mutation	UNP D1A9G5
B	185	ALA	SER	engineered mutation	UNP D1A9G5
B	196	THR	ARG	engineered mutation	UNP D1A9G5
B	205	GLY	ASN	engineered mutation	UNP D1A9G5
B	213	THR	PHE	engineered mutation	UNP D1A9G5
B	214	PRO	SER	engineered mutation	UNP D1A9G5
B	224	ALA	SER	engineered mutation	UNP D1A9G5
B	246	PRO	ALA	engineered mutation	UNP D1A9G5
B	247	SER	ILE	engineered mutation	UNP D1A9G5
B	252	GLU	ASP	engineered mutation	UNP D1A9G5
C	-8	ASN	-	expression tag	UNP D1A9G5
C	-7	SER	-	expression tag	UNP D1A9G5
C	-6	MET	-	expression tag	UNP D1A9G5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	HIS	-	expression tag	UNP D1A9G5
C	-4	HIS	-	expression tag	UNP D1A9G5
C	-3	HIS	-	expression tag	UNP D1A9G5
C	-2	HIS	-	expression tag	UNP D1A9G5
C	-1	HIS	-	expression tag	UNP D1A9G5
C	0	HIS	-	expression tag	UNP D1A9G5
C	13	ALA	SER	engineered mutation	UNP D1A9G5
C	14	SER	LEU	engineered mutation	UNP D1A9G5
C	18	PRO	ALA	engineered mutation	UNP D1A9G5
C	41	THR	ARG	engineered mutation	UNP D1A9G5
C	47	ASP	THR	engineered mutation	UNP D1A9G5
C	92	TYR	LEU	engineered mutation	UNP D1A9G5
C	117	ASP	ASN	engineered mutation	UNP D1A9G5
C	141	ARG	LYS	engineered mutation	UNP D1A9G5
C	142	ARG	SER	engineered mutation	UNP D1A9G5
C	144	PRO	THR	engineered mutation	UNP D1A9G5
C	165	THR	ARG	engineered mutation	UNP D1A9G5
C	175	ASN	LEU	engineered mutation	UNP D1A9G5
C	185	ALA	SER	engineered mutation	UNP D1A9G5
C	196	THR	ARG	engineered mutation	UNP D1A9G5
C	205	GLY	ASN	engineered mutation	UNP D1A9G5
C	213	THR	PHE	engineered mutation	UNP D1A9G5
C	214	PRO	SER	engineered mutation	UNP D1A9G5
C	224	ALA	SER	engineered mutation	UNP D1A9G5
C	246	PRO	ALA	engineered mutation	UNP D1A9G5
C	247	SER	ILE	engineered mutation	UNP D1A9G5
C	252	GLU	ASP	engineered mutation	UNP D1A9G5
D	-8	ASN	-	expression tag	UNP D1A9G5
D	-7	SER	-	expression tag	UNP D1A9G5
D	-6	MET	-	expression tag	UNP D1A9G5
D	-5	HIS	-	expression tag	UNP D1A9G5
D	-4	HIS	-	expression tag	UNP D1A9G5
D	-3	HIS	-	expression tag	UNP D1A9G5
D	-2	HIS	-	expression tag	UNP D1A9G5
D	-1	HIS	-	expression tag	UNP D1A9G5
D	0	HIS	-	expression tag	UNP D1A9G5
D	13	ALA	SER	engineered mutation	UNP D1A9G5
D	14	SER	LEU	engineered mutation	UNP D1A9G5
D	18	PRO	ALA	engineered mutation	UNP D1A9G5
D	41	THR	ARG	engineered mutation	UNP D1A9G5
D	47	ASP	THR	engineered mutation	UNP D1A9G5
D	92	TYR	LEU	engineered mutation	UNP D1A9G5

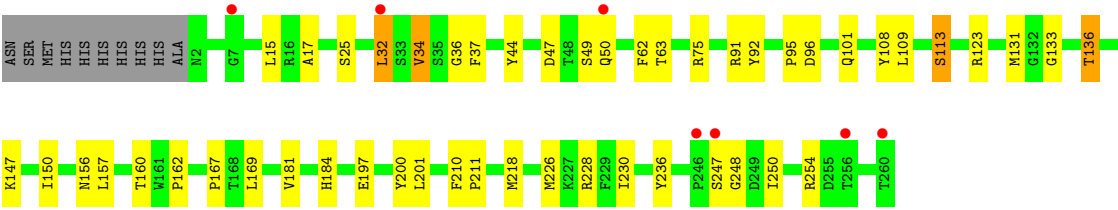
Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	117	ASP	ASN	engineered mutation	UNP D1A9G5
D	141	ARG	LYS	engineered mutation	UNP D1A9G5
D	142	ARG	SER	engineered mutation	UNP D1A9G5
D	144	PRO	THR	engineered mutation	UNP D1A9G5
D	165	THR	ARG	engineered mutation	UNP D1A9G5
D	175	ASN	LEU	engineered mutation	UNP D1A9G5
D	185	ALA	SER	engineered mutation	UNP D1A9G5
D	196	THR	ARG	engineered mutation	UNP D1A9G5
D	205	GLY	ASN	engineered mutation	UNP D1A9G5
D	213	THR	PHE	engineered mutation	UNP D1A9G5
D	214	PRO	SER	engineered mutation	UNP D1A9G5
D	224	ALA	SER	engineered mutation	UNP D1A9G5
D	246	PRO	ALA	engineered mutation	UNP D1A9G5
D	247	SER	ILE	engineered mutation	UNP D1A9G5
D	252	GLU	ASP	engineered mutation	UNP D1A9G5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	79	Total O 79 79	0	0
2	B	64	Total O 64 64	0	0
2	C	62	Total O 62 62	0	0
2	D	70	Total O 70 70	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.25Å 63.59Å 91.54Å 90.00° 97.80° 90.00°	Depositor
Resolution (Å)	45.34 – 2.36 45.34 – 2.36	Depositor EDS
% Data completeness (in resolution range)	97.7 (45.34-2.36) 93.2 (45.34-2.36)	Depositor EDS
R_{merge}	0.31	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.37Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.208 , 0.246 0.214 , 0.251	Depositor DCC
R_{free} test set	2013 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.658	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.033 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8231	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.39	0/2046	0.56	0/2794
1	B	0.33	0/2046	0.55	1/2794 (0.0%)
1	C	0.29	0/2046	0.48	0/2794
1	D	0.35	0/2046	0.54	0/2794
All	All	0.34	0/8184	0.53	1/11176 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	111	GLN	N-CA-C	5.01	116.74	111.28

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	1989	0	1918	28	0
1	B	1989	0	1918	35	0
1	C	1989	0	1918	17	0
1	D	1989	0	1918	33	0
2	A	79	0	0	1	0
2	B	64	0	0	2	0
2	C	62	0	0	2	0
2	D	70	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8231	0	7672	112	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 7.

All (112) 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:D:95:PRO:HG3	1:D:131:MET:HG2	1.59	0.84
1:C:160:THR:HG22	1:C:162:PRO:HD3	1.60	0.81
1:B:160:THR:HG22	1:B:162:PRO:HD3	1.67	0.76
1:B:95:PRO:HG3	1:B:131:MET:HG2	1.69	0.74
1:B:63:THR:HA	1:B:91:ARG:HB2	1.67	0.74
1:A:215:ASN:HB3	1:A:218:MET:HE3	1.70	0.71
1:D:17:ALA:O	1:D:75:ARG:NH1	2.23	0.71
1:D:49:SER:OG	1:D:50:GLN:OE1	2.04	0.71
1:A:42:ILE:HG21	1:A:115:VAL:HG21	1.72	0.71
1:B:19:ARG:NH1	2:B:301:HOH:O	2.24	0.70
1:D:123:ARG:HB3	1:D:147:LYS:HD2	1.74	0.70
1:C:142:ARG:NH2	2:C:302:HOH:O	2.26	0.69
1:C:245:ARG:NH1	2:C:303:HOH:O	2.26	0.68
1:C:109:LEU:HA	1:C:113:SER:HB3	1.77	0.67
1:C:62:PHE:O	1:C:63:THR:HG22	1.95	0.66
1:B:109:LEU:HA	1:B:113:SER:HB3	1.78	0.65
1:D:160:THR:HG22	1:D:162:PRO:HD3	1.80	0.63
1:A:95:PRO:HD3	1:A:131:MET:HE3	1.79	0.63
1:D:109:LEU:HA	1:D:113:SER:HB3	1.82	0.61
1:D:200:TYR:HB3	1:D:254:ARG:HB2	1.81	0.61
1:A:160:THR:HG22	1:A:162:PRO:HD3	1.82	0.61
1:A:25:SER:HB2	1:A:44:TYR:CZ	2.37	0.59
1:D:34:VAL:HG23	1:D:108:TYR:CD2	2.39	0.58
1:B:62:PHE:CE1	1:B:92:TYR:HA	2.39	0.58
1:B:42:ILE:HG21	1:B:115:VAL:HG21	1.86	0.58
1:A:211:PRO:HA	1:A:218:MET:HE1	1.85	0.57
1:B:165:THR:HA	1:B:194:ASN:O	2.06	0.55
1:D:47:ASP:OD1	1:D:50:GLN:NE2	2.35	0.55
1:D:34:VAL:HG21	1:D:37:PHE:CZ	2.41	0.55
1:B:200:TYR:HB3	1:B:254:ARG:HB2	1.87	0.55
1:B:201:LEU:HD11	1:B:250:ILE:HG12	1.87	0.55
1:A:241:CYS:HA	1:A:242:PRO:C	2.32	0.54
1:B:117:ASP:OD2	1:B:117:ASP:N	2.41	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:PHE:O	1:A:63:THR:HG22	2.08	0.53
1:C:167:PRO:HA	1:C:197:GLU:O	2.09	0.53
1:C:200:TYR:HD2	1:C:254:ARG:HG3	1.74	0.53
1:A:82:VAL:HG21	1:A:119:VAL:HG22	1.91	0.53
1:D:150:ILE:HD13	1:D:169:LEU:HB3	1.91	0.53
1:B:72:LEU:HD21	1:B:218:MET:HE2	1.89	0.53
1:B:126:VAL:HG23	1:B:136:THR:HG23	1.91	0.52
1:B:244:PRO:HD2	1:B:253:TYR:CD2	2.45	0.51
1:D:226:MET:O	1:D:230:ILE:HB	2.11	0.51
1:D:201:LEU:HD11	1:D:250:ILE:HG12	1.93	0.51
1:B:195:ALA:O	1:B:196:THR:OG1	2.25	0.50
1:B:123:ARG:HG2	1:B:147:LYS:HE2	1.95	0.49
1:D:147:LYS:O	1:D:167:PRO:HD2	2.13	0.49
1:A:109:LEU:HA	1:A:113:SER:HB3	1.94	0.48
1:B:100:ARG:NE	2:B:309:HOH:O	2.46	0.48
1:D:167:PRO:HA	1:D:197:GLU:O	2.12	0.48
1:A:22:PHE:HE2	1:A:50:GLN:HG3	1.78	0.47
1:D:15:LEU:HD21	1:D:236:TYR:HE2	1.79	0.47
1:D:34:VAL:HG21	1:D:37:PHE:CE1	2.48	0.47
1:A:204:ASN:HD22	1:A:251:SER:HB3	1.78	0.47
1:B:244:PRO:HD2	1:B:253:TYR:CE2	2.49	0.47
1:A:60:PRO:O	1:A:98:ARG:NH1	2.47	0.47
1:D:247:SER:OG	1:D:248:GLY:N	2.46	0.47
1:A:189:TYR:CE1	1:A:198:LYS:HD2	2.50	0.46
1:D:62:PHE:CD2	1:D:63:THR:HG23	2.50	0.46
1:D:211:PRO:HB3	1:D:218:MET:HE1	1.97	0.46
1:B:241:CYS:HA	1:B:242:PRO:C	2.40	0.46
1:A:67:SER:HA	1:A:70:ALA:HB2	1.98	0.46
1:B:62:PHE:CD2	1:B:63:THR:HG23	2.51	0.46
1:D:147:LYS:HD3	1:D:230:ILE:HA	1.98	0.46
1:A:5:GLN:NE2	2:A:312:HOH:O	2.48	0.46
1:A:96:ASP:OD2	1:A:97:SER:N	2.48	0.46
1:C:162:PRO:O	1:C:194:ASN:ND2	2.47	0.45
1:A:13:ALA:H	1:B:118:ARG:HH12	1.64	0.45
1:B:123:ARG:HA	1:B:147:LYS:HG3	1.98	0.45
1:B:34:VAL:HG21	1:B:37:PHE:CZ	2.52	0.45
1:B:228:ARG:O	1:B:228:ARG:NH1	2.48	0.45
1:B:167:PRO:HA	1:B:197:GLU:O	2.17	0.45
1:A:129:HIS:CD2	1:A:211:PRO:HG2	2.52	0.45
1:B:210:PHE:CD1	1:B:211:PRO:HD3	2.52	0.45
1:A:160:THR:CG2	1:A:162:PRO:HD3	2.45	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:PHE:CE2	1:A:92:TYR:HA	2.52	0.44
1:D:123:ARG:CB	1:D:147:LYS:HD2	2.46	0.44
1:B:58:ILE:HA	1:B:127:ALA:O	2.17	0.44
1:C:91:ARG:H	1:C:91:ARG:HG2	1.53	0.44
1:C:42:ILE:HG21	1:C:115:VAL:HG21	1.99	0.44
1:C:58:ILE:HA	1:C:127:ALA:O	2.18	0.44
1:B:175:ASN:ND2	1:B:205:GLY:HA2	2.33	0.43
1:A:62:PHE:HB2	1:A:131:MET:SD	2.58	0.43
1:D:63:THR:HA	1:D:91:ARG:HB3	2.00	0.43
1:C:31:ARG:NH2	1:C:89:ASN:O	2.45	0.43
1:D:34:VAL:HG11	1:D:37:PHE:CZ	2.53	0.43
1:B:34:VAL:HG21	1:B:37:PHE:CE1	2.54	0.43
1:B:233:ASP:OD1	1:B:235:ARG:HG3	2.19	0.43
1:C:37:PHE:CE1	1:C:105:ALA:HA	2.53	0.43
1:B:43:TYR:CE2	1:B:74:PRO:HD3	2.54	0.43
1:D:32:LEU:H	1:D:32:LEU:HD22	1.84	0.43
1:A:87:GLU:HB3	1:A:91:ARG:HH22	1.84	0.42
1:D:133:GLY:HA2	1:D:136:THR:HG23	2.01	0.42
1:A:244:PRO:HG2	1:A:253:TYR:CG	2.55	0.42
1:A:207:SER:O	1:A:210:PHE:HD2	2.02	0.42
1:D:181:VAL:HG22	2:D:329:HOH:O	2.19	0.42
1:D:25:SER:HB2	1:D:44:TYR:CZ	2.55	0.42
1:A:26:GLU:HA	1:A:42:ILE:O	2.20	0.42
1:B:100:ARG:HE	1:B:100:ARG:HB3	1.54	0.41
1:C:212:GLN:H	1:C:212:GLN:HG2	1.72	0.41
1:B:110:THR:O	1:B:116:ARG:HB2	2.20	0.41
1:B:221:PHE:CE2	1:B:244:PRO:HG2	2.55	0.41
1:D:15:LEU:HD21	1:D:236:TYR:CE2	2.55	0.41
1:C:243:PRO:HA	1:C:244:PRO:HD3	1.96	0.41
1:C:249:ASP:OD1	1:C:249:ASP:N	2.50	0.41
1:A:22:PHE:CE2	1:A:50:GLN:HG3	2.55	0.41
1:C:147:LYS:HE2	1:C:147:LYS:HB3	1.79	0.41
1:D:36:GLY:O	1:D:101:GLN:HG2	2.21	0.41
1:D:156:ASN:H	1:D:184:HIS:HD2	1.69	0.41
1:B:252:GLU:HG3	1:B:253:TYR:N	2.35	0.41
1:D:62:PHE:CE1	1:D:92:TYR:HA	2.56	0.41
1:A:76:LEU:HD13	1:A:226:MET:HE1	2.03	0.41
1:D:210:PHE:CG	1:D:211:PRO:HD3	2.56	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	257/269 (96%)	253 (98%)	4 (2%)	0	100	100
1	B	257/269 (96%)	252 (98%)	5 (2%)	0	100	100
1	C	257/269 (96%)	253 (98%)	4 (2%)	0	100	100
1	D	257/269 (96%)	251 (98%)	5 (2%)	1 (0%)	30	34
All	All	1028/1076 (96%)	1009 (98%)	18 (2%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	228	ARG

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	211/220 (96%)	207 (98%)	4 (2%)	50	65
1	B	211/220 (96%)	199 (94%)	12 (6%)	18	23
1	C	211/220 (96%)	204 (97%)	7 (3%)	33	44
1	D	211/220 (96%)	205 (97%)	6 (3%)	38	51
All	All	844/880 (96%)	815 (97%)	29 (3%)	32	43

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	SER
1	A	136	THR
1	A	249	ASP
1	A	260	THR
1	B	25	SER
1	B	34	VAL
1	B	75	ARG
1	B	90	THR
1	B	91	ARG
1	B	97	SER
1	B	112	ARG
1	B	113	SER
1	B	245	ARG
1	B	251	SER
1	B	252	GLU
1	B	255	ASP
1	C	27	GLN
1	C	46	THR
1	C	113	SER
1	C	136	THR
1	C	196	THR
1	C	203	LEU
1	C	249	ASP
1	D	32	LEU
1	D	34	VAL
1	D	96	ASP
1	D	113	SER
1	D	136	THR
1	D	157	LEU

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	5	GLN
1	A	50	GLN
1	A	204	ASN
1	B	50	GLN
1	C	204	ASN
1	D	184	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	259/269 (96%)	0.45	10 (3%) 43 49	24, 34, 44, 60	0
1	B	259/269 (96%)	0.56	15 (5%) 29 33	24, 32, 44, 60	0
1	C	259/269 (96%)	0.37	2 (0%) 82 86	22, 30, 39, 60	0
1	D	259/269 (96%)	0.45	7 (2%) 56 63	24, 32, 43, 54	0
All	All	1036/1076 (96%)	0.46	34 (3%) 49 56	22, 32, 43, 60	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	247	SER	5.2
1	A	51	GLY	4.3
1	B	248	GLY	3.5
1	D	50	GLN	3.3
1	C	260	THR	3.3
1	B	92	TYR	3.1
1	C	50	GLN	3.0
1	B	112	ARG	3.0
1	A	92	TYR	2.9
1	A	260	THR	2.9
1	B	246	PRO	2.8
1	A	48	THR	2.7
1	B	260	THR	2.7
1	A	50	GLN	2.5
1	B	91	ARG	2.4
1	B	210	PHE	2.4
1	D	260	THR	2.4
1	B	255	ASP	2.4
1	A	63	THR	2.4
1	A	241	CYS	2.4
1	D	247	SER	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	32	LEU	2.3
1	A	20	GLY	2.3
1	B	207	SER	2.3
1	B	204	ASN	2.2
1	D	256	THR	2.2
1	D	246	PRO	2.1
1	B	113	SER	2.1
1	B	195	ALA	2.1
1	B	249	ASP	2.1
1	D	7	GLY	2.1
1	A	35	SER	2.0
1	B	193	THR	2.0
1	A	117	ASP	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.