



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 6, 2023 – 10:35 AM EDT

PDB ID : 4EP0
Title : Structure of the bacteriophage C1 tail knob protein, gp12
Authors : Aksyuk, A.A.; Rossmann, M.G.
Deposited on : 2012-04-16
Resolution : 4.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

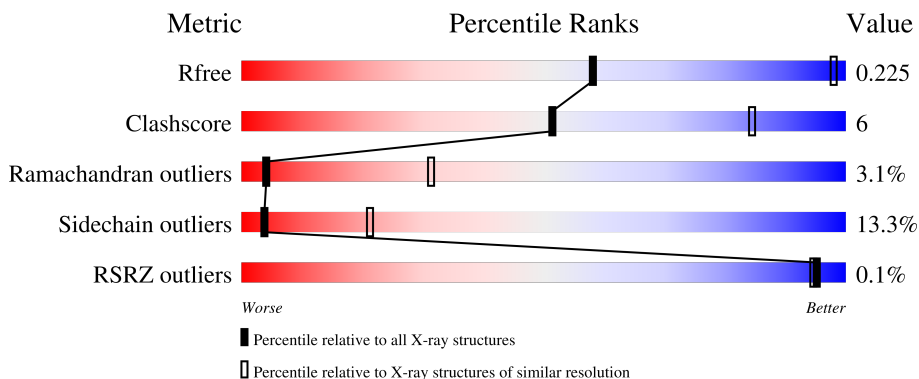
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1087 (4.30-3.70)
Clashscore	141614	1148 (4.30-3.70)
Ramachandran outliers	138981	1108 (4.30-3.70)
Sidechain outliers	138945	1099 (4.30-3.70)
RSRZ outliers	127900	1028 (4.34-3.66)

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	583	 64% 17% 16%
1	B	583	 63% 17% 16%
1	C	583	 62% 19% 16%
1	D	583	 62% 18% 16%
1	E	583	 62% 18% 16%

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Mol	Chain	Length	Quality of chain
1	F	583	 63% 17% • 16%
1	G	583	 63% 17% • 16%
1	H	583	 63% 17% • 16%
1	I	583	 62% 19% • 16%
1	J	583	 64% 16% • 16%
1	K	583	 63% 17% • 16%
1	L	583	 63% 17% • 16%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 47376 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 Major tail protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	488	3948	2504	665	761	18	0	0	0
1	B	488	3948	2504	665	761	18	0	0	0
1	C	488	3948	2504	665	761	18	0	0	0
1	D	488	3948	2504	665	761	18	0	0	0
1	E	488	3948	2504	665	761	18	0	0	0
1	F	488	3948	2504	665	761	18	0	0	0
1	G	488	3948	2504	665	761	18	0	0	0
1	H	488	3948	2504	665	761	18	0	0	0
1	I	488	3948	2504	665	761	18	0	0	0
1	J	488	3948	2504	665	761	18	0	0	0
1	K	488	3948	2504	665	761	18	0	0	0
1	L	488	3948	2504	665	761	18	0	0	0

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	575	SER	-	expression tag	UNP Q7Y3F0
A	576	LEU	-	expression tag	UNP Q7Y3F0
A	577	GLU	-	expression tag	UNP Q7Y3F0
A	578	HIS	-	expression tag	UNP Q7Y3F0
A	579	HIS	-	expression tag	UNP Q7Y3F0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	580	HIS	-	expression tag	UNP Q7Y3F0
A	581	HIS	-	expression tag	UNP Q7Y3F0
A	582	HIS	-	expression tag	UNP Q7Y3F0
A	583	HIS	-	expression tag	UNP Q7Y3F0
B	575	SER	-	expression tag	UNP Q7Y3F0
B	576	LEU	-	expression tag	UNP Q7Y3F0
B	577	GLU	-	expression tag	UNP Q7Y3F0
B	578	HIS	-	expression tag	UNP Q7Y3F0
B	579	HIS	-	expression tag	UNP Q7Y3F0
B	580	HIS	-	expression tag	UNP Q7Y3F0
B	581	HIS	-	expression tag	UNP Q7Y3F0
B	582	HIS	-	expression tag	UNP Q7Y3F0
B	583	HIS	-	expression tag	UNP Q7Y3F0
C	575	SER	-	expression tag	UNP Q7Y3F0
C	576	LEU	-	expression tag	UNP Q7Y3F0
C	577	GLU	-	expression tag	UNP Q7Y3F0
C	578	HIS	-	expression tag	UNP Q7Y3F0
C	579	HIS	-	expression tag	UNP Q7Y3F0
C	580	HIS	-	expression tag	UNP Q7Y3F0
C	581	HIS	-	expression tag	UNP Q7Y3F0
C	582	HIS	-	expression tag	UNP Q7Y3F0
C	583	HIS	-	expression tag	UNP Q7Y3F0
D	575	SER	-	expression tag	UNP Q7Y3F0
D	576	LEU	-	expression tag	UNP Q7Y3F0
D	577	GLU	-	expression tag	UNP Q7Y3F0
D	578	HIS	-	expression tag	UNP Q7Y3F0
D	579	HIS	-	expression tag	UNP Q7Y3F0
D	580	HIS	-	expression tag	UNP Q7Y3F0
D	581	HIS	-	expression tag	UNP Q7Y3F0
D	582	HIS	-	expression tag	UNP Q7Y3F0
D	583	HIS	-	expression tag	UNP Q7Y3F0
E	575	SER	-	expression tag	UNP Q7Y3F0
E	576	LEU	-	expression tag	UNP Q7Y3F0
E	577	GLU	-	expression tag	UNP Q7Y3F0
E	578	HIS	-	expression tag	UNP Q7Y3F0
E	579	HIS	-	expression tag	UNP Q7Y3F0
E	580	HIS	-	expression tag	UNP Q7Y3F0
E	581	HIS	-	expression tag	UNP Q7Y3F0
E	582	HIS	-	expression tag	UNP Q7Y3F0
E	583	HIS	-	expression tag	UNP Q7Y3F0
F	575	SER	-	expression tag	UNP Q7Y3F0
F	576	LEU	-	expression tag	UNP Q7Y3F0

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Chain	Residue	Modelled	Actual	Comment	Reference
F	577	GLU	-	expression tag	UNP Q7Y3F0
F	578	HIS	-	expression tag	UNP Q7Y3F0
F	579	HIS	-	expression tag	UNP Q7Y3F0
F	580	HIS	-	expression tag	UNP Q7Y3F0
F	581	HIS	-	expression tag	UNP Q7Y3F0
F	582	HIS	-	expression tag	UNP Q7Y3F0
F	583	HIS	-	expression tag	UNP Q7Y3F0
G	575	SER	-	expression tag	UNP Q7Y3F0
G	576	LEU	-	expression tag	UNP Q7Y3F0
G	577	GLU	-	expression tag	UNP Q7Y3F0
G	578	HIS	-	expression tag	UNP Q7Y3F0
G	579	HIS	-	expression tag	UNP Q7Y3F0
G	580	HIS	-	expression tag	UNP Q7Y3F0
G	581	HIS	-	expression tag	UNP Q7Y3F0
G	582	HIS	-	expression tag	UNP Q7Y3F0
G	583	HIS	-	expression tag	UNP Q7Y3F0
H	575	SER	-	expression tag	UNP Q7Y3F0
H	576	LEU	-	expression tag	UNP Q7Y3F0
H	577	GLU	-	expression tag	UNP Q7Y3F0
H	578	HIS	-	expression tag	UNP Q7Y3F0
H	579	HIS	-	expression tag	UNP Q7Y3F0
H	580	HIS	-	expression tag	UNP Q7Y3F0
H	581	HIS	-	expression tag	UNP Q7Y3F0
H	582	HIS	-	expression tag	UNP Q7Y3F0
H	583	HIS	-	expression tag	UNP Q7Y3F0
I	575	SER	-	expression tag	UNP Q7Y3F0
I	576	LEU	-	expression tag	UNP Q7Y3F0
I	577	GLU	-	expression tag	UNP Q7Y3F0
I	578	HIS	-	expression tag	UNP Q7Y3F0
I	579	HIS	-	expression tag	UNP Q7Y3F0
I	580	HIS	-	expression tag	UNP Q7Y3F0
I	581	HIS	-	expression tag	UNP Q7Y3F0
I	582	HIS	-	expression tag	UNP Q7Y3F0
I	583	HIS	-	expression tag	UNP Q7Y3F0
J	575	SER	-	expression tag	UNP Q7Y3F0
J	576	LEU	-	expression tag	UNP Q7Y3F0
J	577	GLU	-	expression tag	UNP Q7Y3F0
J	578	HIS	-	expression tag	UNP Q7Y3F0
J	579	HIS	-	expression tag	UNP Q7Y3F0
J	580	HIS	-	expression tag	UNP Q7Y3F0
J	581	HIS	-	expression tag	UNP Q7Y3F0
J	582	HIS	-	expression tag	UNP Q7Y3F0

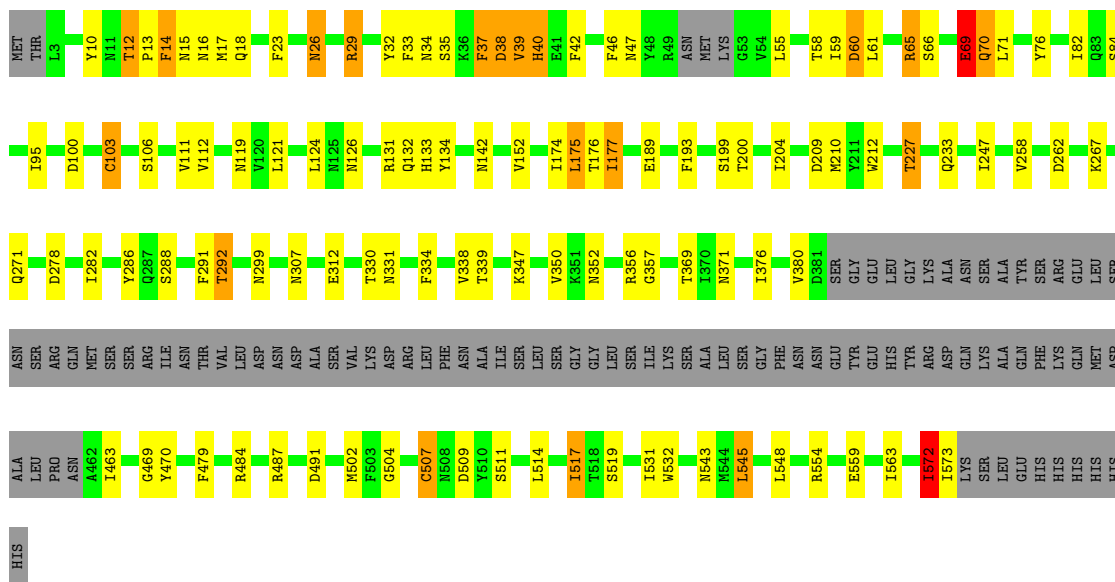
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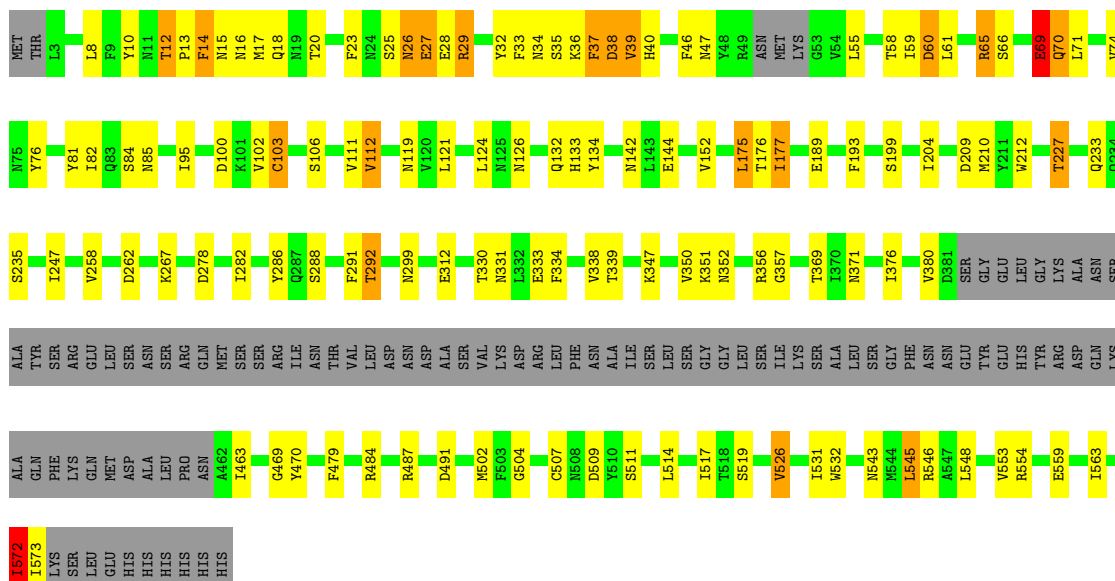
Chain	Residue	Modelled	Actual	Comment	Reference
J	583	HIS	-	expression tag	UNP Q7Y3F0
K	575	SER	-	expression tag	UNP Q7Y3F0
K	576	LEU	-	expression tag	UNP Q7Y3F0
K	577	GLU	-	expression tag	UNP Q7Y3F0
K	578	HIS	-	expression tag	UNP Q7Y3F0
K	579	HIS	-	expression tag	UNP Q7Y3F0
K	580	HIS	-	expression tag	UNP Q7Y3F0
K	581	HIS	-	expression tag	UNP Q7Y3F0
K	582	HIS	-	expression tag	UNP Q7Y3F0
K	583	HIS	-	expression tag	UNP Q7Y3F0
L	575	SER	-	expression tag	UNP Q7Y3F0
L	576	LEU	-	expression tag	UNP Q7Y3F0
L	577	GLU	-	expression tag	UNP Q7Y3F0
L	578	HIS	-	expression tag	UNP Q7Y3F0
L	579	HIS	-	expression tag	UNP Q7Y3F0
L	580	HIS	-	expression tag	UNP Q7Y3F0
L	581	HIS	-	expression tag	UNP Q7Y3F0
L	582	HIS	-	expression tag	UNP Q7Y3F0
L	583	HIS	-	expression tag	UNP Q7Y3F0



• Molecule 1: Major tail protein

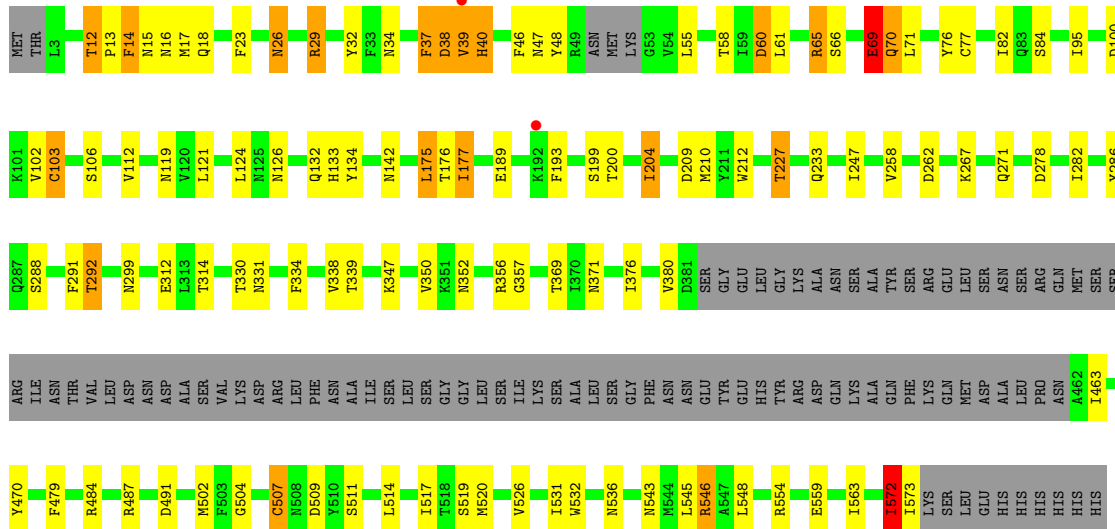


• Molecule 1: Major tail protein



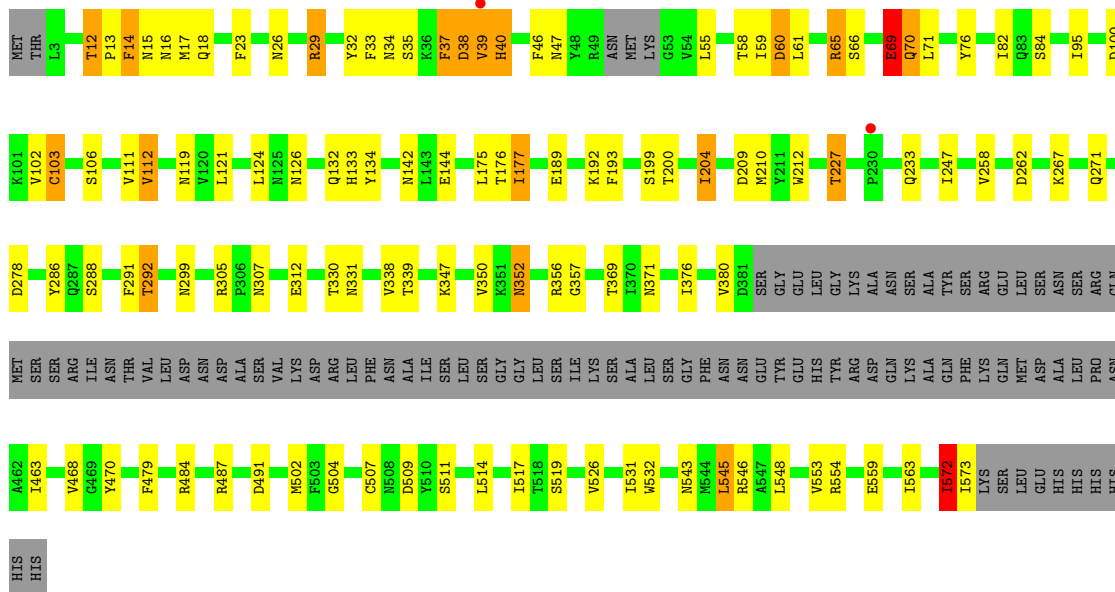
• Molecule 1: Major tail protein

Chain J:  64% 16% 16%



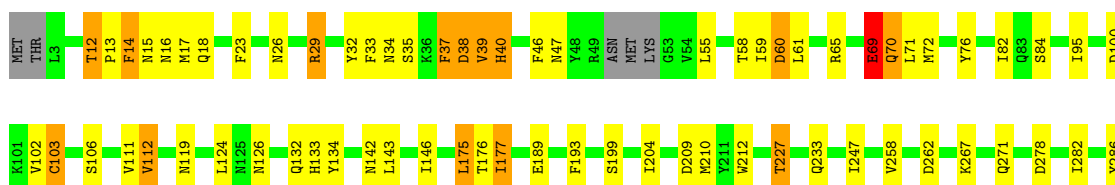
• Molecule 1: Major tail protein

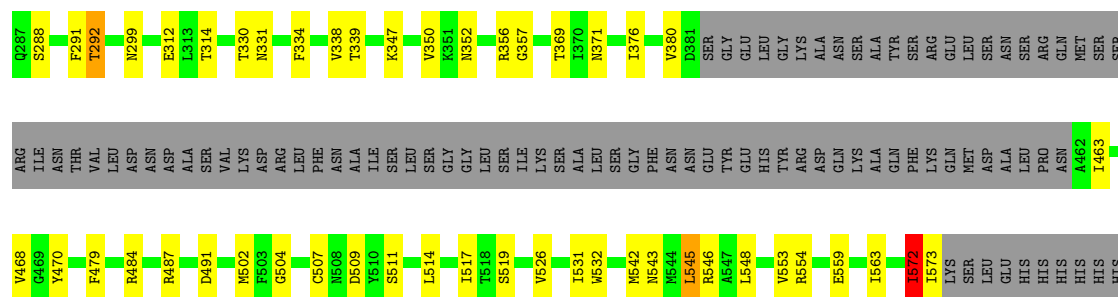
Chain K:  63% 17% 16%



• Molecule 1: Major tail protein

Chain L:  63% 17% 16%





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	316.87Å 211.38Å 178.60Å 90.00° 116.53° 90.00°	Depositor
Resolution (Å)	49.73 – 4.00 49.73 – 4.00	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.73-4.00) 99.4 (49.73-4.00)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	0.25	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.13 (at 4.00Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.195 , 0.233 0.184 , 0.225	Depositor DCC
R_{free} test set	4429 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	58.5	Xtrriage
Anisotropy	0.593	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 97.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	47376	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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	0.50	0/4030	0.71	0/5464
1	B	0.50	0/4030	0.71	0/5464
1	C	0.55	2/4030 (0.0%)	0.73	4/5464 (0.1%)
1	D	0.52	0/4030	0.73	3/5464 (0.1%)
1	E	0.49	0/4030	0.70	0/5464
1	F	0.51	0/4030	0.72	1/5464 (0.0%)
1	G	0.47	0/4030	0.70	2/5464 (0.0%)
1	H	0.48	0/4030	0.70	1/5464 (0.0%)
1	I	0.51	0/4030	0.72	1/5464 (0.0%)
1	J	0.49	0/4030	0.71	1/5464 (0.0%)
1	K	0.48	0/4030	0.70	1/5464 (0.0%)
1	L	0.48	0/4030	0.70	1/5464 (0.0%)
All	All	0.50	2/48360 (0.0%)	0.71	15/65568 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	291	PHE	CE1-CZ	-8.24	1.21	1.37
1	C	291	PHE	CD1-CE1	-5.88	1.27	1.39

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	291	PHE	CB-CG-CD1	-6.34	116.36	120.80
1	G	484	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	J	546	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	C	291	PHE	CB-CG-CD2	5.71	124.79	120.80
1	L	546	ARG	NE-CZ-NH1	5.58	123.09	120.30

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	3948	0	3832	43	0
1	B	3948	0	3832	52	0
1	C	3948	0	3832	63	2
1	D	3948	0	3832	55	1
1	E	3948	0	3832	48	0
1	F	3948	0	3832	45	0
1	G	3948	0	3832	44	0
1	H	3948	0	3832	44	0
1	I	3948	0	3832	68	0
1	J	3948	0	3832	42	0
1	K	3948	0	3832	45	1
1	L	3948	0	3832	47	0
All	All	47376	0	45984	524	2

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 6.

The worst 5 of 524 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:C:291:PHE:HZ	1:I:26:ASN:N	1.74	0.85
1:C:291:PHE:CZ	1:I:25:SER:HB2	2.22	0.74
1:L:312:GLU:OE2	1:L:484:ARG:NH1	2.23	0.72
1:F:286:TYR:N	1:F:331:ASN:OD1	2.24	0.71
1:F:32:TYR:OH	1:F:559:GLU:OE2	2.09	0.70

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:262:ASP:O	1:K:192:LYS:NZ[4_547]	2.08	0.12
1:C:260:GLY:O	1:D:187:THR:OG1[2_556]	2.10	0.10

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	482/583 (83%)	408 (85%)	58 (12%)	16 (3%)	4	30
1	B	482/583 (83%)	408 (85%)	60 (12%)	14 (3%)	4	32
1	C	482/583 (83%)	407 (84%)	59 (12%)	16 (3%)	4	30
1	D	482/583 (83%)	410 (85%)	57 (12%)	15 (3%)	4	31
1	E	482/583 (83%)	412 (86%)	56 (12%)	14 (3%)	4	32
1	F	482/583 (83%)	408 (85%)	58 (12%)	16 (3%)	4	30
1	G	482/583 (83%)	409 (85%)	59 (12%)	14 (3%)	4	32
1	H	482/583 (83%)	410 (85%)	56 (12%)	16 (3%)	4	30
1	I	482/583 (83%)	406 (84%)	60 (12%)	16 (3%)	4	30
1	J	482/583 (83%)	410 (85%)	56 (12%)	16 (3%)	4	30
1	K	482/583 (83%)	411 (85%)	57 (12%)	14 (3%)	4	32
1	L	482/583 (83%)	410 (85%)	58 (12%)	14 (3%)	4	32
All	All	5784/6996 (83%)	4909 (85%)	694 (12%)	181 (3%)	4	31

5 of 181 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	70	GLN
1	A	84	SER
1	A	258	VAL
1	B	18	GLN

5.3.2 Protein sidechains [i](#)

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	446/529 (84%)	386 (86%)	60 (14%)	4	21
1	B	446/529 (84%)	386 (86%)	60 (14%)	4	21
1	C	446/529 (84%)	387 (87%)	59 (13%)	4	21
1	D	446/529 (84%)	389 (87%)	57 (13%)	4	22
1	E	446/529 (84%)	385 (86%)	61 (14%)	3	20
1	F	446/529 (84%)	388 (87%)	58 (13%)	4	22
1	G	446/529 (84%)	384 (86%)	62 (14%)	3	20
1	H	446/529 (84%)	387 (87%)	59 (13%)	4	21
1	I	446/529 (84%)	386 (86%)	60 (14%)	4	21
1	J	446/529 (84%)	385 (86%)	61 (14%)	3	20
1	K	446/529 (84%)	387 (87%)	59 (13%)	4	21
1	L	446/529 (84%)	389 (87%)	57 (13%)	4	22
All	All	5352/6348 (84%)	4639 (87%)	713 (13%)	4	21

5 of 713 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	376	ILE
1	J	291	PHE
1	I	10	TYR
1	H	369	THR
1	I	350	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	233	GLN
1	I	233	GLN
1	L	233	GLN
1	J	233	GLN
1	D	233	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	488/583 (83%)	-0.51	0 100 100	25, 83, 153, 207	0
1	B	488/583 (83%)	-0.45	0 100 100	28, 81, 148, 232	0
1	C	488/583 (83%)	-0.54	0 100 100	16, 72, 151, 222	0
1	D	488/583 (83%)	-0.60	1 (0%) 95 93	18, 72, 140, 212	0
1	E	488/583 (83%)	-0.47	1 (0%) 95 93	27, 82, 146, 215	0
1	F	488/583 (83%)	-0.52	0 100 100	17, 79, 158, 247	0
1	G	488/583 (83%)	-0.35	1 (0%) 95 93	41, 97, 171, 227	0
1	H	488/583 (83%)	-0.45	0 100 100	35, 91, 170, 258	0
1	I	488/583 (83%)	-0.41	0 100 100	20, 84, 160, 222	0
1	J	488/583 (83%)	-0.34	2 (0%) 92 87	35, 93, 162, 211	0
1	K	488/583 (83%)	-0.39	2 (0%) 92 87	36, 94, 159, 226	0
1	L	488/583 (83%)	-0.38	0 100 100	37, 90, 174, 247	0
All	All	5856/6996 (83%)	-0.45	7 (0%) 95 94	16, 85, 161, 258	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	39	VAL	3.6
1	J	39	VAL	3.0
1	E	39	VAL	3.0
1	K	230	PRO	2.9
1	D	39	VAL	2.4

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.