



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 08:03 PM UTC

PDB ID : 1NLZ / pdb_00001nlz
Title : Crystal structure of unliganded traffic ATPase of the type IV secretion system of helicobacter pylori
Authors : Savvides, S.N.; Yeo, H.J.; Beck, M.R.; Blaesing, F.; Lurz, R.; Lanka, E.; Buhrdorf, R.; Fischer, W.; Haas, R.; Waksman, G.
Deposited on : 2003-01-08
Resolution : 3.00 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 : 4-5-2 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

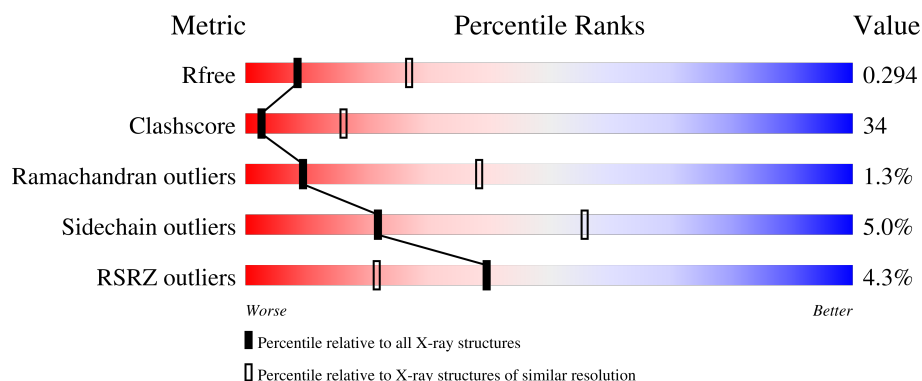
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	330	52% 42% ••
1	B	330	2% 42% 46% 5% 7%
1	C	330	14% 38% 40% 6% 15%
1	D	330	2% 41% 47% 8% •
1	E	330	4% 42% 45% 5% 8%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	330	% 49% 40% • 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14503 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 virB11 homolog.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	Se	0	0	0
			2550	1617	436	483	8	6			
1	B	307	Total	C	N	O	S	Se	0	0	0
			2419	1533	414	458	8	6			
1	C	279	Total	C	N	O	S	Se	0	0	0
			2194	1386	379	418	6	5			
1	D	320	Total	C	N	O	S	Se	0	0	0
			2516	1597	432	473	8	6			
1	E	305	Total	C	N	O	S	Se	0	0	0
			2394	1518	409	453	8	6			
1	F	305	Total	C	N	O	S	Se	0	0	0
			2409	1528	413	454	8	6			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	MSE	MET	modified residue	UNP Q7BK04
A	82	MSE	MET	modified residue	UNP Q7BK04
A	192	MSE	MET	modified residue	UNP Q7BK04
A	239	MSE	MET	modified residue	UNP Q7BK04
A	287	MSE	MET	modified residue	UNP Q7BK04
A	312	MSE	MET	modified residue	UNP Q7BK04
B	42	MSE	MET	modified residue	UNP Q7BK04
B	82	MSE	MET	modified residue	UNP Q7BK04
B	192	MSE	MET	modified residue	UNP Q7BK04
B	239	MSE	MET	modified residue	UNP Q7BK04
B	287	MSE	MET	modified residue	UNP Q7BK04
B	312	MSE	MET	modified residue	UNP Q7BK04
C	42	MSE	MET	modified residue	UNP Q7BK04
C	82	MSE	MET	modified residue	UNP Q7BK04
C	192	MSE	MET	modified residue	UNP Q7BK04
C	239	MSE	MET	modified residue	UNP Q7BK04
C	287	MSE	MET	modified residue	UNP Q7BK04

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	312	MSE	MET	modified residue	UNP Q7BK04
D	42	MSE	MET	modified residue	UNP Q7BK04
D	82	MSE	MET	modified residue	UNP Q7BK04
D	192	MSE	MET	modified residue	UNP Q7BK04
D	239	MSE	MET	modified residue	UNP Q7BK04
D	287	MSE	MET	modified residue	UNP Q7BK04
D	312	MSE	MET	modified residue	UNP Q7BK04
E	42	MSE	MET	modified residue	UNP Q7BK04
E	82	MSE	MET	modified residue	UNP Q7BK04
E	192	MSE	MET	modified residue	UNP Q7BK04
E	239	MSE	MET	modified residue	UNP Q7BK04
E	287	MSE	MET	modified residue	UNP Q7BK04
E	312	MSE	MET	modified residue	UNP Q7BK04
F	42	MSE	MET	modified residue	UNP Q7BK04
F	82	MSE	MET	modified residue	UNP Q7BK04
F	192	MSE	MET	modified residue	UNP Q7BK04
F	239	MSE	MET	modified residue	UNP Q7BK04
F	287	MSE	MET	modified residue	UNP Q7BK04
F	312	MSE	MET	modified residue	UNP Q7BK04

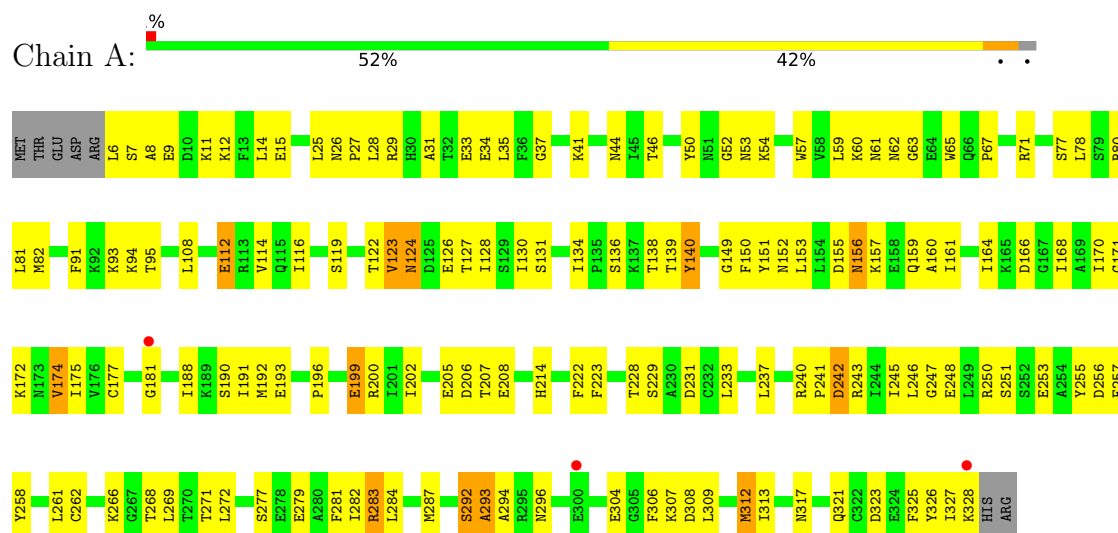
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total O 4 4	0	0
2	B	7	Total O 7 7	0	0
2	D	4	Total O 4 4	0	0
2	E	3	Total O 3 3	0	0
2	F	3	Total O 3 3	0	0

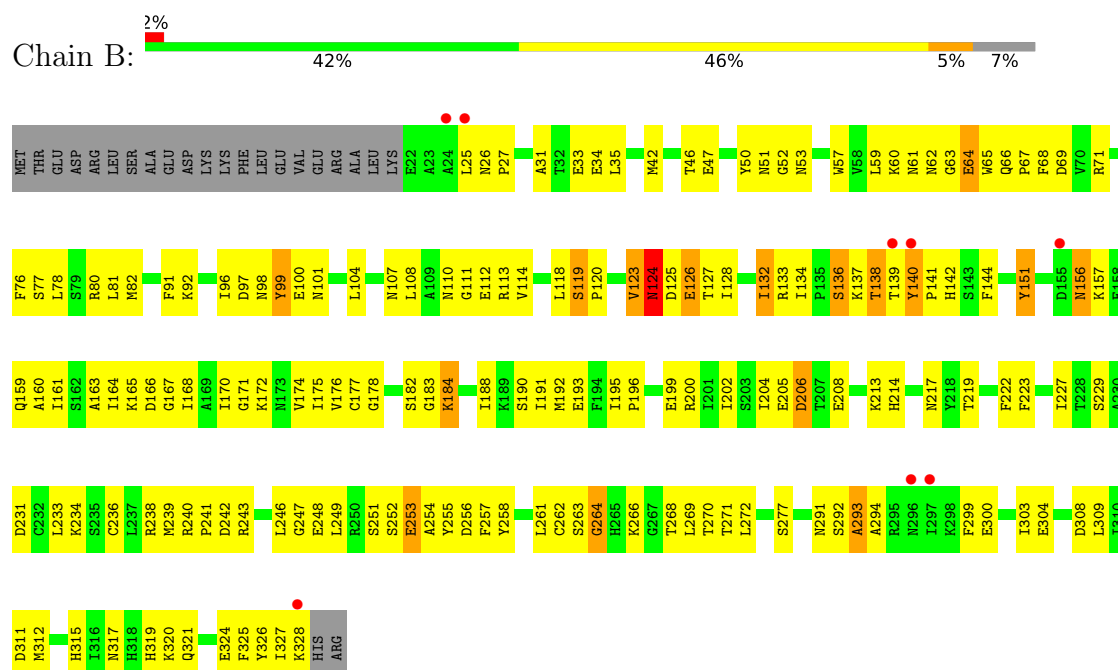
3 Residue-property plots [i](#)

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.

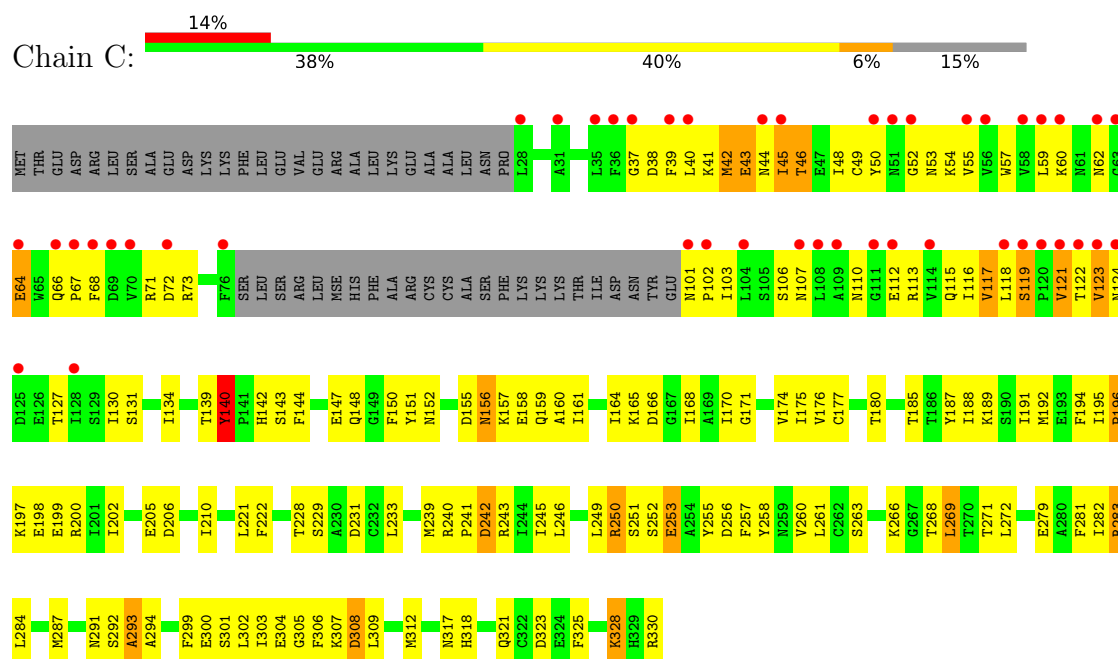
• Molecule 1: virB11 homolog



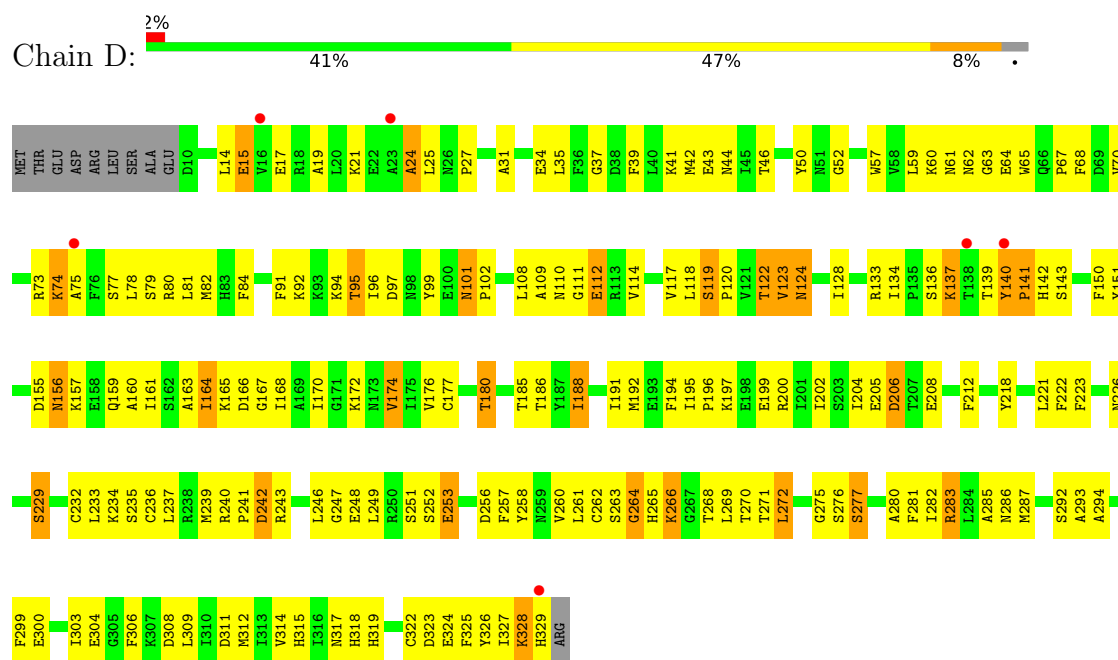
• Molecule 1: virB11 homolog



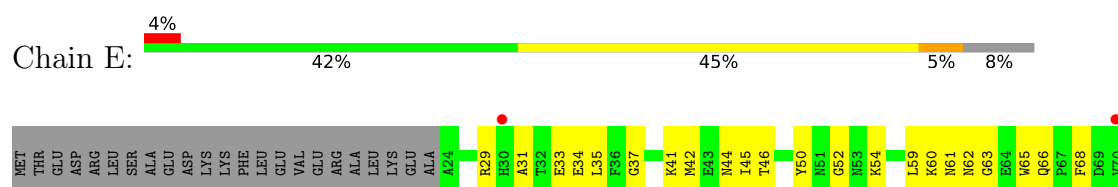
- Molecule 1: virB11 homolog

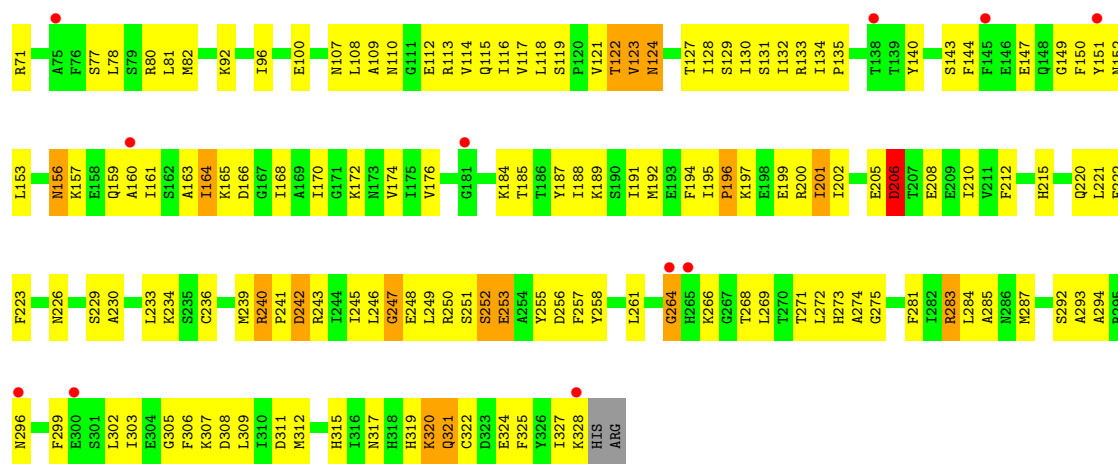


- Molecule 1: virB11 homolog

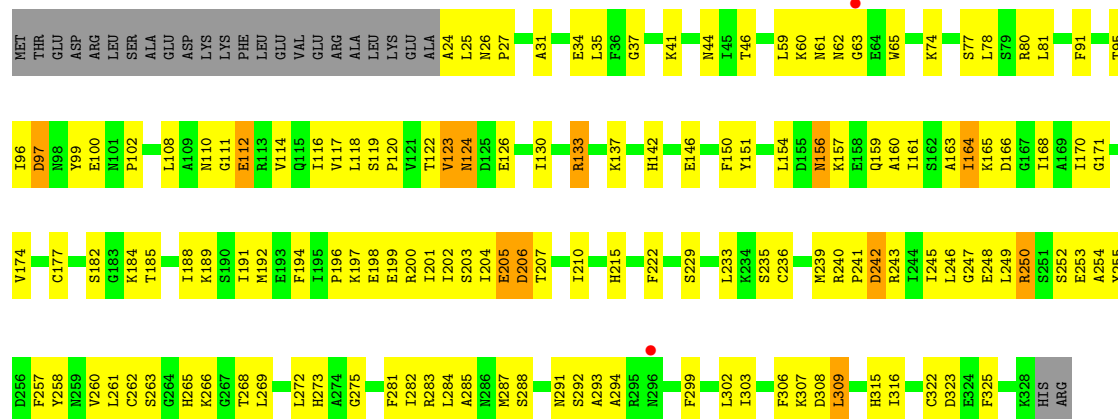


- Molecule 1: virB11 homolog





• Molecule 1: virB11 homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	104.53Å 104.53Å 168.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.70 – 3.00 29.70 – 3.00	Depositor EDS
% Data completeness (in resolution range)	88.9 (29.70-3.00) 88.9 (29.70-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.53 (at 3.00Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.239 , 0.298 0.241 , 0.294	Depositor DCC
R_{free} test set	1858 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l 0.047 for h,-h-k,-l 0.033 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	14503	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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	0.81	0/2593	1.09	9/3486 (0.3%)
1	B	0.79	1/2461 (0.0%)	1.14	14/3311 (0.4%)
1	C	0.83	0/2230	1.13	10/2999 (0.3%)
1	D	0.75	0/2560	1.08	17/3443 (0.5%)
1	E	0.78	3/2435 (0.1%)	1.05	9/3276 (0.3%)
1	F	0.73	1/2451 (0.0%)	1.00	4/3296 (0.1%)
All	All	0.78	5/14730 (0.0%)	1.08	63/19811 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	127	THR	CA-C	6.17	1.60	1.52
1	E	309	LEU	CA-CB	5.38	1.61	1.53
1	E	309	LEU	CA-C	-5.30	1.45	1.52
1	F	309	LEU	CA-C	-5.23	1.45	1.52
1	B	141	PRO	C-O	-5.06	1.18	1.23

The worst 5 of 63 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	309	LEU	N-CA-C	10.27	126.36	112.90
1	E	242	ASP	N-CA-C	-8.37	103.59	113.88
1	E	252	SER	N-CA-C	8.18	120.20	111.28
1	B	140	TYR	CB-CA-C	-7.75	94.39	110.50
1	D	123	VAL	N-CA-C	7.74	118.49	110.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	99	TYR	Sidechain

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	2550	0	2509	144	0
1	B	2419	0	2369	199	0
1	C	2194	0	2146	179	0
1	D	2516	0	2463	196	0
1	E	2394	0	2340	158	0
1	F	2409	0	2369	146	0
2	A	4	0	0	0	0
2	B	7	0	0	0	0
2	D	4	0	0	0	0
2	E	3	0	0	0	0
2	F	3	0	0	0	0
All	All	14503	0	14196	973	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 34.

The worst 5 of 973 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:54:LYS:HD2	1:C:71:ARG:HA	1.29	1.14
1:F:60:LYS:HE3	1:F:62:ASN:HD21	1.17	1.07
1:C:45:ILE:HG13	1:C:134:ILE:HG21	1.38	1.06
1:A:54:LYS:HD2	1:A:71:ARG:HA	1.41	1.02
1:D:74:LYS:HD3	1:D:74:LYS:N	1.74	1.02

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	321/330 (97%)	286 (89%)	31 (10%)	4 (1%)	10	40
1	B	305/330 (92%)	270 (88%)	31 (10%)	4 (1%)	9	38
1	C	275/330 (83%)	230 (84%)	43 (16%)	2 (1%)	18	53
1	D	318/330 (96%)	280 (88%)	31 (10%)	7 (2%)	5	26
1	E	303/330 (92%)	272 (90%)	27 (9%)	4 (1%)	9	38
1	F	303/330 (92%)	271 (89%)	29 (10%)	3 (1%)	12	45
All	All	1825/1980 (92%)	1609 (88%)	192 (10%)	24 (1%)	9	38

5 of 24 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	GLU
1	D	137	LYS
1	A	136	SER
1	A	293	ALA
1	B	136	SER

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	279/287 (97%)	266 (95%)	13 (5%)	23	58
1	B	265/287 (92%)	255 (96%)	10 (4%)	29	63
1	C	240/287 (84%)	221 (92%)	19 (8%)	11	39

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	273/287 (95%)	257 (94%)	16 (6%)	18	50
1	E	261/287 (91%)	250 (96%)	11 (4%)	26	61
1	F	265/287 (92%)	255 (96%)	10 (4%)	29	63
All	All	1583/1722 (92%)	1504 (95%)	79 (5%)	22	56

5 of 79 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	283	ARG
1	F	112	GLU
1	E	156	ASN
1	E	253	GLU
1	F	164	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 46 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	286	ASN
1	E	220	GLN
1	D	318	HIS
1	E	66	GLN
1	E	273	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no 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 [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	317/330 (96%)	-0.01	3 (0%) 81 61	12, 40, 63, 74	0
1	B	301/330 (91%)	0.17	8 (2%) 56 33	10, 46, 71, 89	0
1	C	274/330 (83%)	0.57	45 (16%) 4 3	11, 46, 90, 99	0
1	D	314/330 (95%)	0.32	6 (1%) 66 43	29, 52, 80, 98	0
1	E	299/330 (90%)	0.35	13 (4%) 40 21	28, 54, 84, 95	0
1	F	299/330 (90%)	0.14	2 (0%) 84 66	25, 49, 71, 80	0
All	All	1804/1980 (91%)	0.25	77 (4%) 40 21	10, 48, 80, 99	0

The worst 5 of 77 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	328	LYS	4.7
1	C	40	LEU	4.5
1	C	119	SER	4.4
1	C	56	VAL	4.3
1	C	125	ASP	4.3

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.