



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 31, 2022 – 04:11 PM JST

PDB ID : 7W64  
Title : Crystal structure of minor pilin TcpB from *Vibrio cholerae* complexed with N-terminal peptide fragment of TcpF  
Authors : Oki, H.; Kawahara, K.; Iimori, M.; Imoto, Y.; Maruno, T.; Uchiyama, S.; Muroga, Y.; Yoshida, A.; Yoshida, T.; Ohkubo, T.; Matsuda, S.; Iida, T.; Nakamura, S.  
Deposited on : 2021-12-01  
Resolution : 2.30 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
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)	:	2.31.2

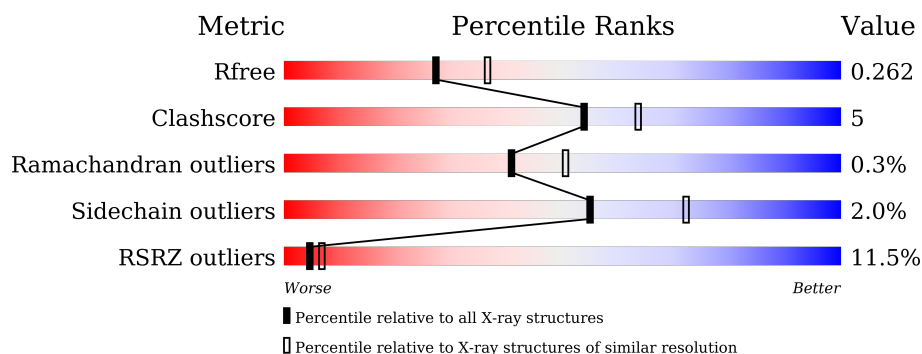
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	397	<div> <div>20%</div> <div>86%</div> <div>12%</div> <div>..</div> </div>
1	B	397	<div> <div>8%</div> <div>90%</div> <div>9%</div> <div>.</div> </div>
1	C	397	<div> <div>6%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>
1	D	397	<div> <div>15%</div> <div>84%</div> <div>13%</div> <div>..</div> </div>
1	E	397	<div> <div>8%</div> <div>89%</div> <div>10%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	397	<div><div></div><div>12%</div><div>84%</div><div>15%</div><div></div></div>
2	G	15	<div><div></div><div>67%</div><div>7%</div><div>27%</div><div></div></div>
2	H	15	<div><div></div><div>60%</div><div>13%</div><div>27%</div><div></div></div>
2	I	15	<div><div></div><div>67%</div><div>7%</div><div>27%</div><div></div></div>
2	J	15	<div><div></div><div>53%</div><div>13%</div><div>7%</div><div>27%</div><div></div></div>
2	K	15	<div><div></div><div>33%</div><div>67%</div><div>7%</div><div>27%</div><div></div></div>
2	L	15	<div><div></div><div>7%</div><div>73%</div><div>27%</div><div></div></div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 19163 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 Toxin-coregulated pilus biosynthesis protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	0	0
			3020	1889	513	610	8			
1	B	395	Total	C	N	O	S	0	0	0
			3053	1911	517	616	9			
1	C	395	Total	C	N	O	S	0	0	0
			3053	1911	517	616	9			
1	D	391	Total	C	N	O	S	0	0	0
			3020	1889	513	610	8			
1	E	395	Total	C	N	O	S	0	0	0
			3053	1911	517	616	9			
1	F	395	Total	C	N	O	S	0	0	0
			3053	1911	517	616	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	GLY	-	expression tag	UNP Q9AGX1
A	28	GLY	-	expression tag	UNP Q9AGX1
B	27	GLY	-	expression tag	UNP Q9AGX1
B	28	GLY	-	expression tag	UNP Q9AGX1
C	27	GLY	-	expression tag	UNP Q9AGX1
C	28	GLY	-	expression tag	UNP Q9AGX1
D	27	GLY	-	expression tag	UNP Q9AGX1
D	28	GLY	-	expression tag	UNP Q9AGX1
E	27	GLY	-	expression tag	UNP Q9AGX1
E	28	GLY	-	expression tag	UNP Q9AGX1
F	27	GLY	-	expression tag	UNP Q9AGX1
F	28	GLY	-	expression tag	UNP Q9AGX1

- Molecule 2 is a protein called TcpF.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	11	Total	C	N	O	0	0	0
			86	52	13	21			
2	H	11	Total	C	N	O	0	0	0
			86	52	13	21			
2	I	11	Total	C	N	O	0	0	0
			86	52	13	21			
2	J	11	Total	C	N	O	0	0	0
			86	52	13	21			
2	K	11	Total	C	N	O	0	0	0
			86	52	13	21			
2	L	11	Total	C	N	O	0	0	0
			86	52	13	21			

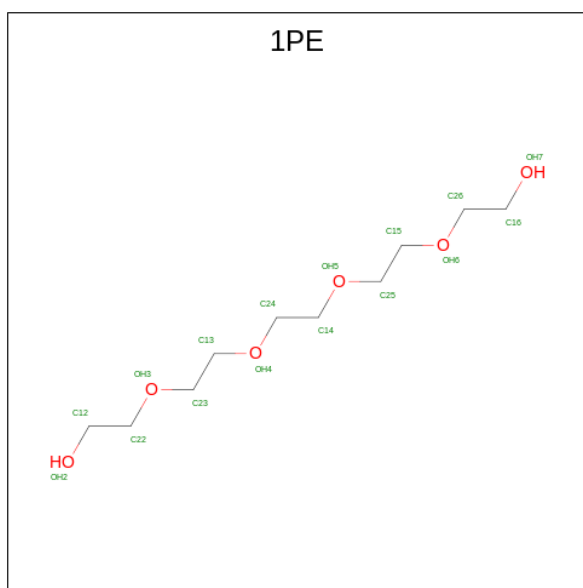
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	3	Total	Ca	0	0
			3	3		
3	C	2	Total	Ca	0	0
			2	2		
3	D	1	Total	Ca	0	0
			1	1		
3	E	3	Total	Ca	0	0
			3	3		
3	F	1	Total	Ca	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	E	1	Total	Cl	0	0
			1	1		
4	G	1	Total	Cl	0	0
			1	1		
4	I	1	Total	Cl	0	0
			1	1		
4	L	1	Total	Cl	0	0
			1	1		

- Molecule 5 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula:  $C_{10}H_{22}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			16	10	6		

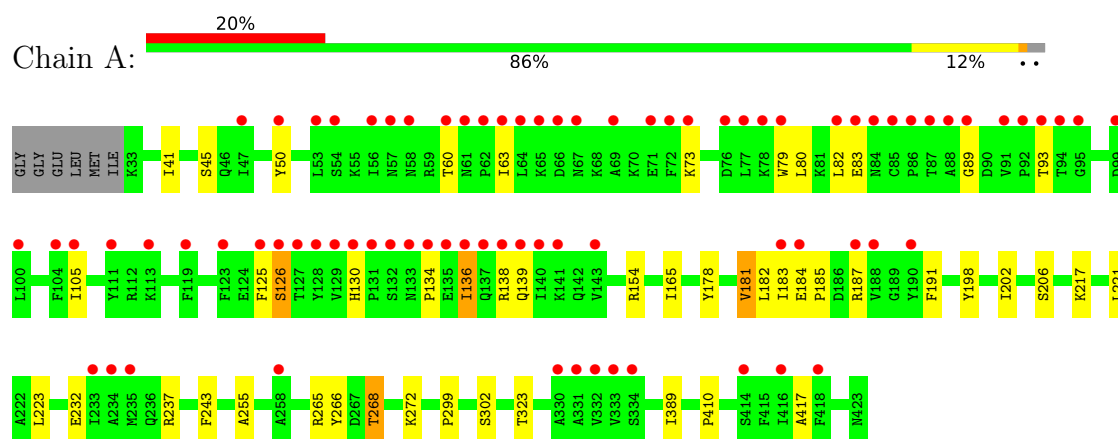
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	76	Total	O	0	0
			76	76		
6	B	103	Total	O	0	0
			103	103		
6	C	66	Total	O	0	0
			66	66		
6	D	20	Total	O	0	0
			20	20		
6	E	55	Total	O	0	0
			55	55		
6	F	35	Total	O	0	0
			35	35		
6	G	2	Total	O	0	0
			2	2		
6	H	5	Total	O	0	0
			5	5		

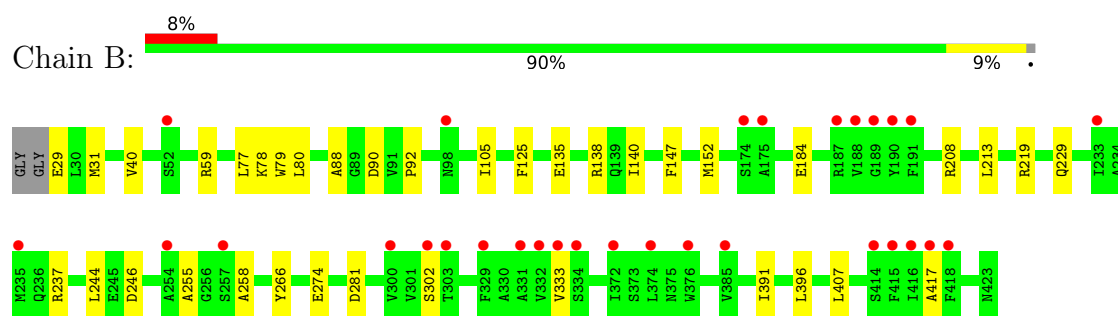
### 3 Residue-property plots [i](#)

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.

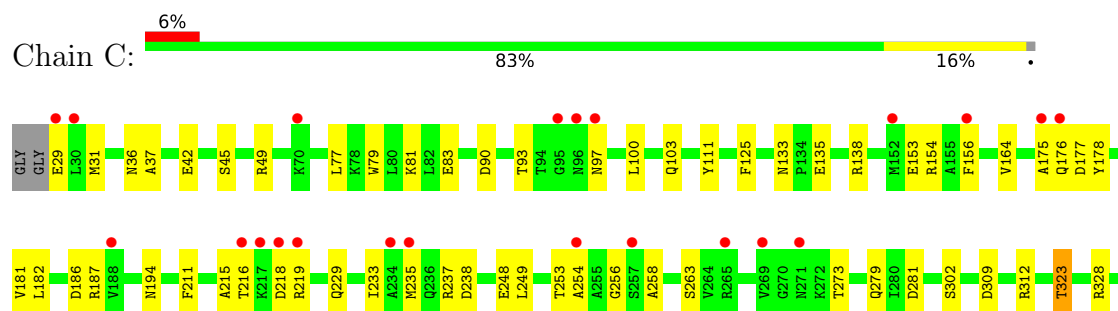
- Molecule 1: Toxin-coregulated pilus biosynthesis protein B



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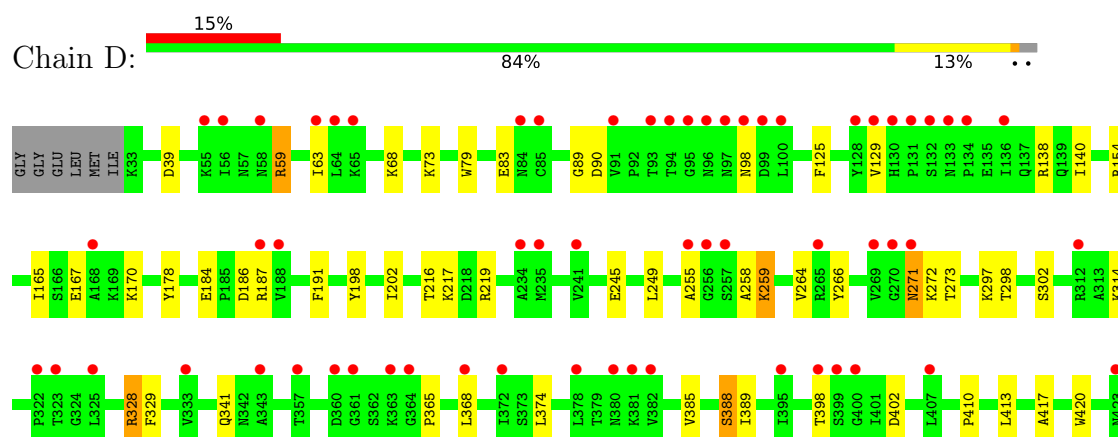


- Molecule 1: Toxin-coregulated pilus biosynthesis protein B

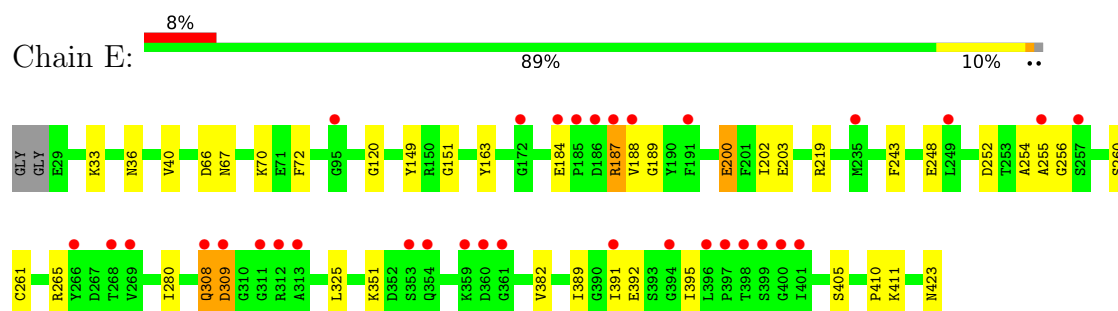




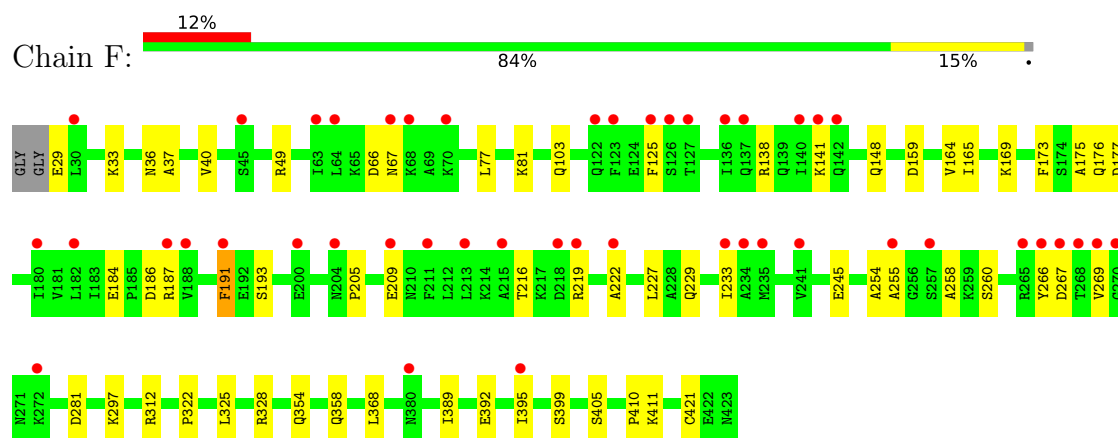
- Molecule 1: Toxin-coregulated pilus biosynthesis protein B



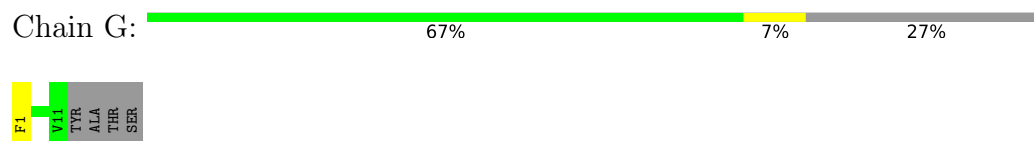
- Molecule 1: Toxin-coregulated pilus biosynthesis protein B



- Molecule 1: Toxin-coregulated pilus biosynthesis protein B



- Molecule 2: TcpF





## ● Molecule 2: TcpF

Chain H:  60% 13% 27%



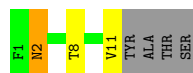
## ● Molecule 2: TcpF

Chain I:  67% 7% 27%



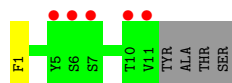
## ● Molecule 2: TcpF

Chain J:  53% 13% 7% 27%




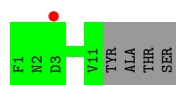
## ● Molecule 2: TcpF

Chain K:  33% 67% 7% 27%



## ● Molecule 2: TcpF

Chain L:  7% 73% 27%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.88Å 128.56Å 327.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.20 – 2.30 48.77 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.6 (37.20-2.30) 98.8 (48.77-2.30)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.19-4092	Depositor
R, $R_{free}$	0.222 , 0.264 0.220 , 0.262	Depositor DCC
$R_{free}$ test set	13805 reflections (9.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.6	Xtriage
Anisotropy	0.669	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 46.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	19163	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, CL, 1PE

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	0.49	0/3075	0.67	0/4157
1	B	0.54	0/3108	0.66	0/4201
1	C	0.50	0/3108	0.66	0/4201
1	D	0.42	0/3075	0.63	0/4157
1	E	0.49	0/3108	0.64	0/4201
1	F	0.45	0/3108	0.62	0/4201
2	G	0.55	0/87	0.57	0/118
2	H	0.48	0/87	0.63	0/118
2	I	0.45	0/87	0.54	0/118
2	J	0.36	0/87	0.51	0/118
2	K	0.34	0/87	0.55	0/118
2	L	0.41	0/87	0.54	0/118
All	All	0.48	0/19104	0.64	0/25826

There are no bond length outliers.

There are no bond angle outliers.

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	3020	0	2933	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3053	0	2970	22	0
1	C	3053	0	2970	45	0
1	D	3020	0	2933	38	0
1	E	3053	0	2972	27	0
1	F	3053	0	2970	40	0
2	G	86	0	74	1	0
2	H	86	0	74	2	0
2	I	86	0	74	1	0
2	J	86	0	74	3	0
2	K	86	0	74	1	0
2	L	86	0	74	0	0
3	A	1	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
3	E	3	0	0	0	0
3	F	1	0	0	0	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	1	0
4	I	1	0	0	0	0
4	L	1	0	0	0	0
5	B	16	0	22	1	0
6	A	76	0	0	2	0
6	B	103	0	0	3	0
6	C	66	0	0	6	0
6	D	20	0	0	0	0
6	E	55	0	0	0	0
6	F	35	0	0	4	0
6	G	2	0	0	0	0
6	H	5	0	0	1	0
All	All	19163	0	18214	196	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 5.

The worst 5 of 196 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:255:ALA:HB1	1:D:259:LYS:HE2	1.61	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ARG:NH2	1:F:354:GLN:HG3	2.01	0.76
1:A:237:ARG:HA	1:C:235:MET:HE3	1.71	0.73
1:F:229:GLN:HB3	1:F:233:ILE:HG13	1.71	0.72
1:C:29:GLU:HG3	1:C:31:MET:HG3	1.72	0.70

There are no symmetry-related clashes.

## 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	389/397 (98%)	374 (96%)	14 (4%)	1 (0%)	41	50
1	B	393/397 (99%)	381 (97%)	12 (3%)	0	100	100
1	C	393/397 (99%)	380 (97%)	13 (3%)	0	100	100
1	D	389/397 (98%)	373 (96%)	15 (4%)	1 (0%)	41	50
1	E	393/397 (99%)	375 (95%)	14 (4%)	4 (1%)	15	17
1	F	393/397 (99%)	381 (97%)	11 (3%)	1 (0%)	41	50
2	G	9/15 (60%)	9 (100%)	0	0	100	100
2	H	9/15 (60%)	9 (100%)	0	0	100	100
2	I	9/15 (60%)	8 (89%)	1 (11%)	0	100	100
2	J	9/15 (60%)	9 (100%)	0	0	100	100
2	K	9/15 (60%)	9 (100%)	0	0	100	100
2	L	9/15 (60%)	9 (100%)	0	0	100	100
All	All	2404/2472 (97%)	2317 (96%)	80 (3%)	7 (0%)	41	50

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	258	ALA

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Mol	Chain	Res	Type
1	E	308	GLN
1	E	309	ASP
1	F	176	GLN
1	E	255	ALA

### 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	330/334 (99%)	323 (98%)	7 (2%)	53	70
1	B	334/334 (100%)	331 (99%)	3 (1%)	78	89
1	C	334/334 (100%)	328 (98%)	6 (2%)	59	75
1	D	330/334 (99%)	318 (96%)	12 (4%)	35	49
1	E	334/334 (100%)	330 (99%)	4 (1%)	71	84
1	F	334/334 (100%)	327 (98%)	7 (2%)	53	70
2	G	11/14 (79%)	11 (100%)	0	100	100
2	H	11/14 (79%)	11 (100%)	0	100	100
2	I	11/14 (79%)	11 (100%)	0	100	100
2	J	11/14 (79%)	9 (82%)	2 (18%)	1	1
2	K	11/14 (79%)	11 (100%)	0	100	100
2	L	11/14 (79%)	11 (100%)	0	100	100
All	All	2062/2088 (99%)	2021 (98%)	41 (2%)	55	72

5 of 41 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	187	ARG
1	F	216	THR
1	E	200	GLU
1	F	159	ASP
1	F	399	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	B	148	GLN
1	C	176	GLN
1	D	271	ASN
2	J	2	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 17 are monoatomic - leaving 1 for Mogul analysis.

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	1PE	B	504	-	15,15,15	0.53	0	14,14,14	0.50	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	1PE	B	504	-	-	9/13/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	504	1PE	OH6-C15-C25-OH5
5	B	504	1PE	OH4-C13-C23-OH3
5	B	504	1PE	OH5-C14-C24-OH4
5	B	504	1PE	C15-C25-OH5-C14
5	B	504	1PE	C16-C26-OH6-C15

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	504	1PE	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	391/397 (98%)	1.11	79 (20%) 1 1	34, 61, 112, 130	0
1	B	395/397 (99%)	0.57	30 (7%) 13 18	35, 54, 82, 111	0
1	C	395/397 (99%)	0.49	25 (6%) 20 25	34, 59, 91, 105	0
1	D	391/397 (98%)	0.92	61 (15%) 2 2	49, 77, 105, 117	0
1	E	395/397 (99%)	0.49	33 (8%) 11 15	46, 62, 93, 104	0
1	F	395/397 (99%)	0.69	46 (11%) 4 6	47, 71, 97, 115	0
2	G	11/15 (73%)	-0.27	0 100 100	46, 51, 66, 69	0
2	H	11/15 (73%)	-0.01	0 100 100	43, 53, 61, 67	0
2	I	11/15 (73%)	-0.20	0 100 100	45, 51, 68, 71	0
2	J	11/15 (73%)	0.90	0 100 100	72, 82, 87, 90	0
2	K	11/15 (73%)	2.07	5 (45%) 0 0	91, 98, 105, 106	0
2	L	11/15 (73%)	0.88	1 (9%) 9 12	63, 75, 93, 94	0
All	All	2428/2472 (98%)	0.71	280 (11%) 4 7	34, 63, 100, 130	0

The worst 5 of 280 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	GLN	8.8
1	A	129	VAL	8.6
1	A	63	ILE	7.9
1	A	128	TYR	7.8
1	A	188	VAL	7.3

### 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 monosaccharides 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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
3	CA	B	503	1/1	0.78	0.33	88,88,88,88	0
3	CA	D	501	1/1	0.80	0.11	85,85,85,85	0
3	CA	B	502	1/1	0.84	0.12	59,59,59,59	0
4	CL	D	502	1/1	0.84	0.13	86,86,86,86	0
3	CA	E	502	1/1	0.86	0.19	69,69,69,69	0
5	1PE	B	504	16/16	0.89	0.17	54,64,71,74	0
3	CA	C	502	1/1	0.90	0.24	85,85,85,85	0
4	CL	E	504	1/1	0.90	0.07	95,95,95,95	0
3	CA	E	503	1/1	0.90	0.22	102,102,102,102	0
4	CL	L	101	1/1	0.92	0.13	66,66,66,66	0
3	CA	B	501	1/1	0.93	0.28	67,67,67,67	0
4	CL	I	101	1/1	0.94	0.17	53,53,53,53	0
3	CA	F	501	1/1	0.95	0.21	82,82,82,82	0
3	CA	E	501	1/1	0.95	0.25	73,73,73,73	0
4	CL	A	502	1/1	0.97	0.10	54,54,54,54	0
3	CA	A	501	1/1	0.97	0.07	66,66,66,66	0
4	CL	G	101	1/1	0.98	0.07	54,54,54,54	0
3	CA	C	501	1/1	0.98	0.27	65,65,65,65	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.