



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 11:44 AM UTC

PDB ID : 2PMB / pdb\_00002pmb  
Title : Crystal structure of predicted nucleotide-binding protein from *Vibrio cholerae*  
Authors : Patskovsky, Y.; Zhan, C.; Shi, W.; Toro, R.; Sauder, J.M.; Gilmore, J.; Iizuka, M.; Maletic, M.; Gheyi, T.; Wasserman, S.R.; Smith, D.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2007-04-20  
Resolution : 1.99 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 : 4-5-2 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

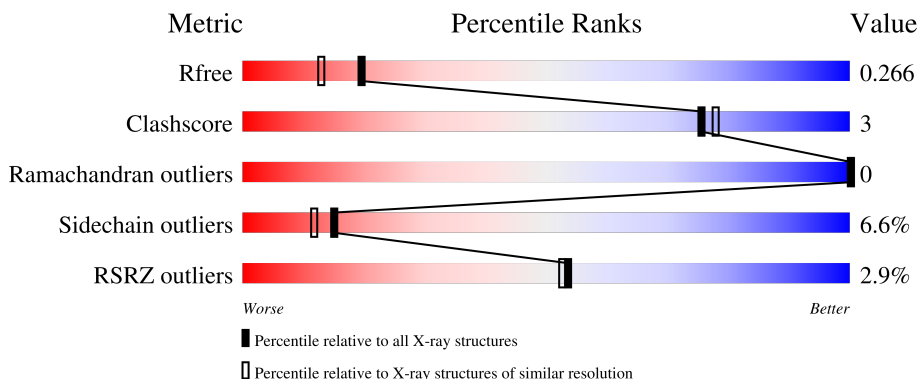
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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  $\geq 3$ , 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	462	
1	B	462	
1	C	462	
1	D	462	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14539 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	437	3483	2214	606	640	5	18	0	9	0
1	B	437	3479	2210	605	641	5	18	0	8	0
1	C	434	3484	2217	609	634	5	19	0	12	0
1	D	437	3478	2210	607	638	5	18	0	8	0

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	MSE	-	cloning artifact	UNP Q9KTK3
A	5	SER	-	cloning artifact	UNP Q9KTK3
A	6	LEU	-	cloning artifact	UNP Q9KTK3
A	16	MSE	MET	modified residue	UNP Q9KTK3
A	71	MSE	MET	modified residue	UNP Q9KTK3
A	114	MSE	MET	modified residue	UNP Q9KTK3
A	192	MSE	MET	modified residue	UNP Q9KTK3
A	196	MSE	MET	modified residue	UNP Q9KTK3
A	237	MSE	MET	modified residue	UNP Q9KTK3
A	250	MSE	MET	modified residue	UNP Q9KTK3
A	274	MSE	MET	modified residue	UNP Q9KTK3
A	275	MSE	MET	modified residue	UNP Q9KTK3
A	284	MSE	MET	modified residue	UNP Q9KTK3
A	331	MSE	MET	modified residue	UNP Q9KTK3
A	335	MSE	MET	modified residue	UNP Q9KTK3
A	372	MSE	MET	modified residue	UNP Q9KTK3
A	418	MSE	MET	modified residue	UNP Q9KTK3
A	425	MSE	MET	modified residue	UNP Q9KTK3
A	428	MSE	MET	modified residue	UNP Q9KTK3
A	441	MSE	MET	modified residue	UNP Q9KTK3
A	458	GLU	-	cloning artifact	UNP Q9KTK3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	459	GLY	-	cloning artifact	UNP Q9KTK3
A	460	HIS	-	cloning artifact	UNP Q9KTK3
A	461	HIS	-	cloning artifact	UNP Q9KTK3
A	462	HIS	-	cloning artifact	UNP Q9KTK3
A	463	HIS	-	cloning artifact	UNP Q9KTK3
A	464	HIS	-	cloning artifact	UNP Q9KTK3
A	465	HIS	-	cloning artifact	UNP Q9KTK3
B	4	MSE	-	cloning artifact	UNP Q9KTK3
B	5	SER	-	cloning artifact	UNP Q9KTK3
B	6	LEU	-	cloning artifact	UNP Q9KTK3
B	16	MSE	MET	modified residue	UNP Q9KTK3
B	71	MSE	MET	modified residue	UNP Q9KTK3
B	114	MSE	MET	modified residue	UNP Q9KTK3
B	192	MSE	MET	modified residue	UNP Q9KTK3
B	196	MSE	MET	modified residue	UNP Q9KTK3
B	237	MSE	MET	modified residue	UNP Q9KTK3
B	250	MSE	MET	modified residue	UNP Q9KTK3
B	274	MSE	MET	modified residue	UNP Q9KTK3
B	275	MSE	MET	modified residue	UNP Q9KTK3
B	284	MSE	MET	modified residue	UNP Q9KTK3
B	331	MSE	MET	modified residue	UNP Q9KTK3
B	335	MSE	MET	modified residue	UNP Q9KTK3
B	372	MSE	MET	modified residue	UNP Q9KTK3
B	418	MSE	MET	modified residue	UNP Q9KTK3
B	425	MSE	MET	modified residue	UNP Q9KTK3
B	428	MSE	MET	modified residue	UNP Q9KTK3
B	441	MSE	MET	modified residue	UNP Q9KTK3
B	458	GLU	-	cloning artifact	UNP Q9KTK3
B	459	GLY	-	cloning artifact	UNP Q9KTK3
B	460	HIS	-	cloning artifact	UNP Q9KTK3
B	461	HIS	-	cloning artifact	UNP Q9KTK3
B	462	HIS	-	cloning artifact	UNP Q9KTK3
B	463	HIS	-	cloning artifact	UNP Q9KTK3
B	464	HIS	-	cloning artifact	UNP Q9KTK3
B	465	HIS	-	cloning artifact	UNP Q9KTK3
C	4	MSE	-	cloning artifact	UNP Q9KTK3
C	5	SER	-	cloning artifact	UNP Q9KTK3
C	6	LEU	-	cloning artifact	UNP Q9KTK3
C	16	MSE	MET	modified residue	UNP Q9KTK3
C	71	MSE	MET	modified residue	UNP Q9KTK3
C	114	MSE	MET	modified residue	UNP Q9KTK3
C	192	MSE	MET	modified residue	UNP Q9KTK3

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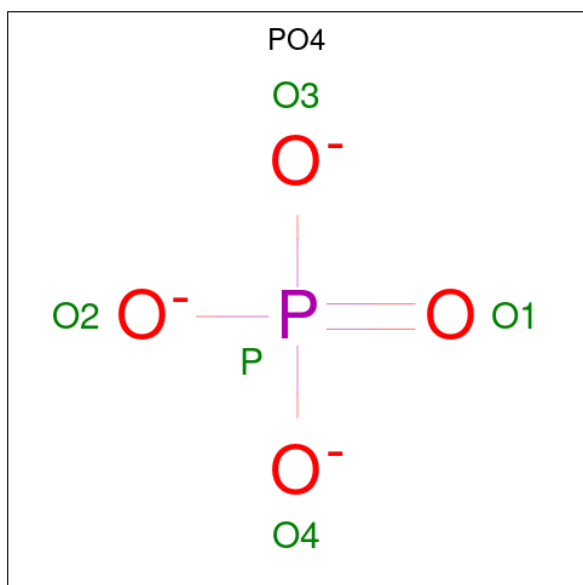
Chain	Residue	Modelled	Actual	Comment	Reference
C	196	MSE	MET	modified residue	UNP Q9KTK3
C	237	MSE	MET	modified residue	UNP Q9KTK3
C	250	MSE	MET	modified residue	UNP Q9KTK3
C	274	MSE	MET	modified residue	UNP Q9KTK3
C	275	MSE	MET	modified residue	UNP Q9KTK3
C	284	MSE	MET	modified residue	UNP Q9KTK3
C	331	MSE	MET	modified residue	UNP Q9KTK3
C	335	MSE	MET	modified residue	UNP Q9KTK3
C	372	MSE	MET	modified residue	UNP Q9KTK3
C	418	MSE	MET	modified residue	UNP Q9KTK3
C	425	MSE	MET	modified residue	UNP Q9KTK3
C	428	MSE	MET	modified residue	UNP Q9KTK3
C	441	MSE	MET	modified residue	UNP Q9KTK3
C	458	GLU	-	cloning artifact	UNP Q9KTK3
C	459	GLY	-	cloning artifact	UNP Q9KTK3
C	460	HIS	-	cloning artifact	UNP Q9KTK3
C	461	HIS	-	cloning artifact	UNP Q9KTK3
C	462	HIS	-	cloning artifact	UNP Q9KTK3
C	463	HIS	-	cloning artifact	UNP Q9KTK3
C	464	HIS	-	cloning artifact	UNP Q9KTK3
C	465	HIS	-	cloning artifact	UNP Q9KTK3
D	4	MSE	-	cloning artifact	UNP Q9KTK3
D	5	SER	-	cloning artifact	UNP Q9KTK3
D	6	LEU	-	cloning artifact	UNP Q9KTK3
D	16	MSE	MET	modified residue	UNP Q9KTK3
D	71	MSE	MET	modified residue	UNP Q9KTK3
D	114	MSE	MET	modified residue	UNP Q9KTK3
D	192	MSE	MET	modified residue	UNP Q9KTK3
D	196	MSE	MET	modified residue	UNP Q9KTK3
D	237	MSE	MET	modified residue	UNP Q9KTK3
D	250	MSE	MET	modified residue	UNP Q9KTK3
D	274	MSE	MET	modified residue	UNP Q9KTK3
D	275	MSE	MET	modified residue	UNP Q9KTK3
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D	418	MSE	MET	modified residue	UNP Q9KTK3
D	425	MSE	MET	modified residue	UNP Q9KTK3
D	428	MSE	MET	modified residue	UNP Q9KTK3
D	441	MSE	MET	modified residue	UNP Q9KTK3
D	458	GLU	-	cloning artifact	UNP Q9KTK3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	459	GLY	-	cloning artifact	UNP Q9KTK3
D	460	HIS	-	cloning artifact	UNP Q9KTK3
D	461	HIS	-	cloning artifact	UNP Q9KTK3
D	462	HIS	-	cloning artifact	UNP Q9KTK3
D	463	HIS	-	cloning artifact	UNP Q9KTK3
D	464	HIS	-	cloning artifact	UNP Q9KTK3
D	465	HIS	-	cloning artifact	UNP Q9KTK3

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O<sub>4</sub>P).



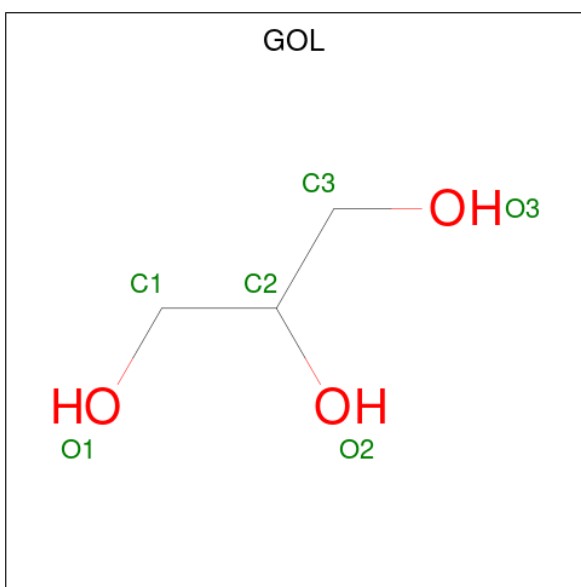
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	138	Total	O	0	0
			138	138		
4	B	149	Total	O	0	0
			149	149		
4	C	124	Total	O	0	0
			124	124		

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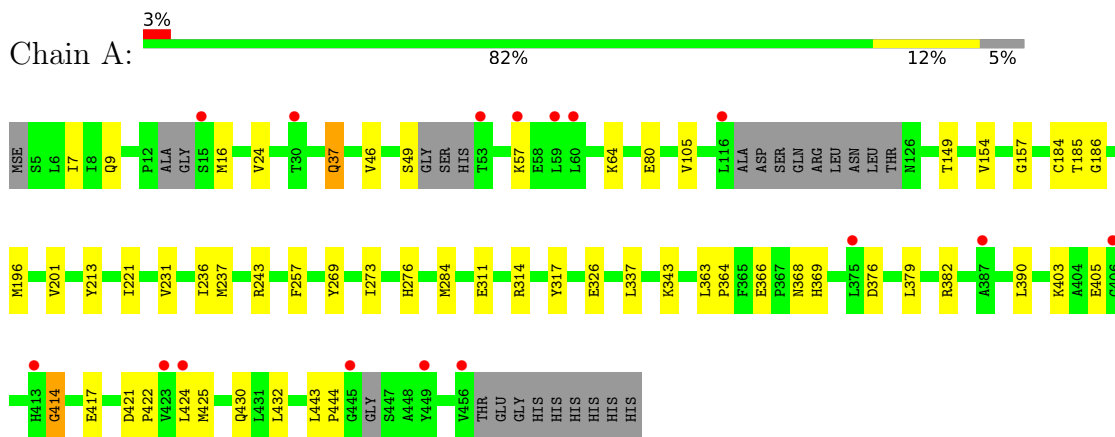
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	D	132	Total 132	O 132	0	0

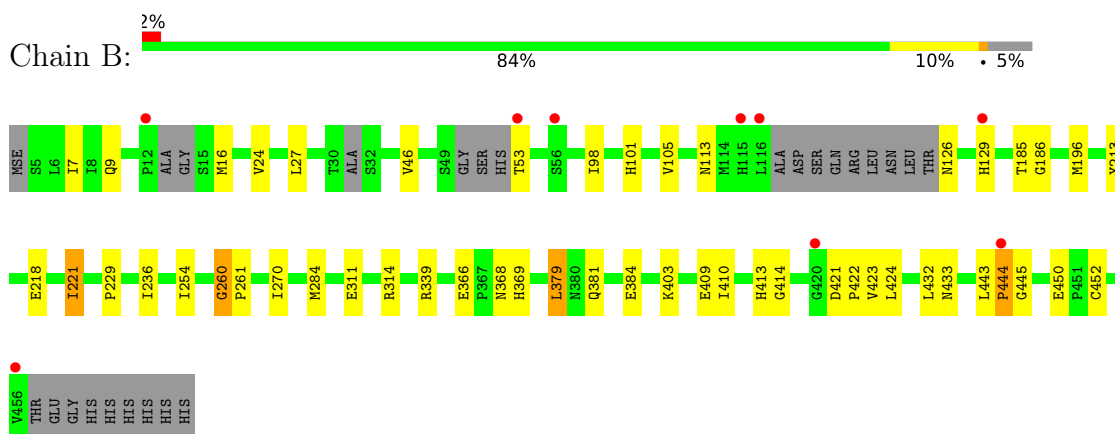
### 3 Residue-property plots [i](#)

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.

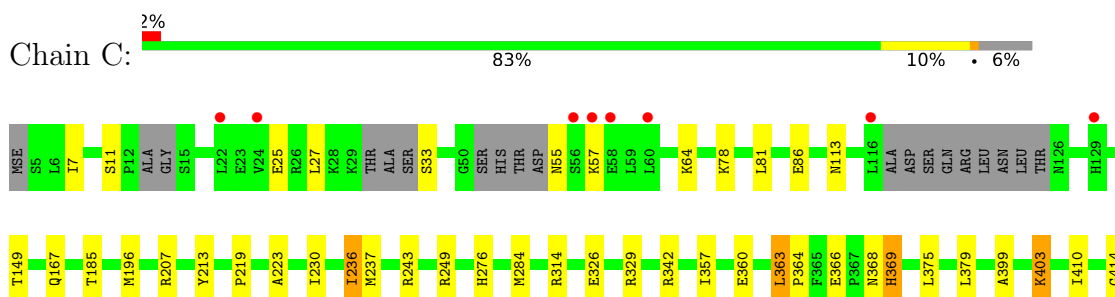
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein

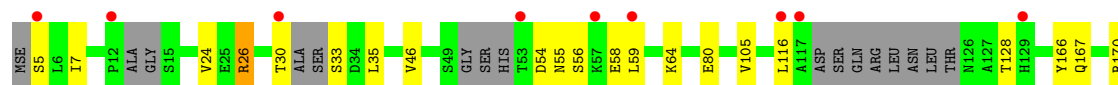
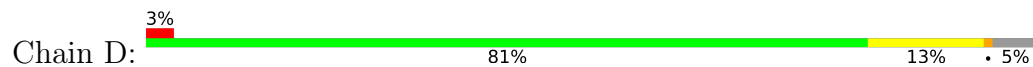


- Molecule 1: Uncharacterized protein





• Molecule 1: Uncharacterized protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.80Å 178.22Å 89.51Å 90.00° 103.78° 90.00°	Depositor
Resolution (Å)	20.00 – 1.99 20.00 – 1.99	Depositor EDS
% Data completeness (in resolution range)	90.5 (20.00-1.99) 90.4 (20.00-1.99)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.01 (at 1.99Å)	Xtrriage
Refinement program	REFMAC 5.3.0034	Depositor
R, $R_{free}$	0.208 , 0.266 0.208 , 0.266	Depositor DCC
$R_{free}$ test set	3331 reflections (2.70%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.3	Xtrriage
Anisotropy	0.026	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 46.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14539	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4

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	0.58	0/3562	0.95	4/4787 (0.1%)
1	B	0.55	0/3556	0.94	5/4777 (0.1%)
1	C	0.58	1/3573 (0.0%)	0.96	2/4797 (0.0%)
1	D	0.56	0/3555	0.94	3/4775 (0.1%)
All	All	0.57	1/14246 (0.0%)	0.95	14/19136 (0.1%)

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	2
1	D	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	369	HIS	C-O	-5.43	1.17	1.24

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	260	GLY	CA-C-N	6.87	126.84	119.28
1	B	260	GLY	C-N-CA	6.87	126.84	119.28
1	A	276	HIS	CA-C-N	5.85	125.99	119.32
1	A	276	HIS	C-N-CA	5.85	125.99	119.32
1	D	166	TYR	N-CA-C	5.55	117.01	111.07

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	444	PRO	Peptide
1	B	445	GLY	Peptide
1	D	58	GLU	Peptide

## 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	3483	0	3514	22	0
1	B	3479	0	3501	16	0
1	C	3484	0	3520	21	0
1	D	3478	0	3507	27	0
2	A	10	0	0	0	0
2	B	20	0	0	0	0
2	C	15	0	0	0	0
2	D	15	0	0	0	0
3	B	6	0	8	0	0
3	C	6	0	8	3	0
4	A	138	0	0	1	0
4	B	149	0	0	0	0
4	C	124	0	0	1	0
4	D	132	0	0	1	0
All	All	14539	0	14058	84	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 3.

The worst 5 of 84 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:167[A]:GLN:HG3	1:D:170[A]:ARG:HH22	1.38	0.86
1:C:249:ARG:HH22	3:C:701:GOL:H11	1.49	0.78
1:D:167[A]:GLN:HG3	1:D:170[A]:ARG:NH2	2.07	0.70
1:C:357:ILE:H	3:C:701:GOL:H12	1.57	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:185:THR:HG22	1:C:196:MSE:HE3	1.75	0.68

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	436/462 (94%)	423 (97%)	13 (3%)	0	100	100
1	B	435/462 (94%)	419 (96%)	16 (4%)	0	100	100
1	C	436/462 (94%)	424 (97%)	12 (3%)	0	100	100
1	D	434/462 (94%)	426 (98%)	8 (2%)	0	100	100
All	All	1741/1848 (94%)	1692 (97%)	49 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	381/373 (102%)	355 (93%)	26 (7%)	14	11
1	B	380/373 (102%)	355 (93%)	25 (7%)	15	12
1	C	380/373 (102%)	353 (93%)	27 (7%)	13	10
1	D	379/373 (102%)	351 (93%)	28 (7%)	13	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1520/1492 (102%)	1414 (93%)	106 (7%)	15 10

5 of 106 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	64	LYS
1	C	379	LEU
1	D	410	ILE
1	C	81	LEU
1	C	360	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	433	ASN
1	C	381	GLN
1	D	378	HIS
1	C	389	ASN
1	C	368	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

14 ligands are modelled in this entry.

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
2	PO4	B	607	-	4,4,4	0.77	0	6,6,6	0.68	0
2	PO4	B	609	-	4,4,4	0.94	0	6,6,6	0.55	0
2	PO4	B	612	-	4,4,4	0.83	0	6,6,6	0.60	0
2	PO4	D	601	-	4,4,4	1.09	0	6,6,6	0.32	0
2	PO4	A	604	-	4,4,4	0.92	0	6,6,6	0.57	0
2	PO4	B	603	-	4,4,4	1.20	0	6,6,6	0.61	0
2	PO4	D	606	-	4,4,4	1.02	0	6,6,6	0.48	0
3	GOL	C	701	-	5,5,5	0.34	0	5,5,5	0.37	0
2	PO4	A	610	-	4,4,4	1.00	0	6,6,6	0.54	0
2	PO4	D	602	-	4,4,4	0.87	0	6,6,6	0.51	0
2	PO4	C	611	-	4,4,4	0.93	0	6,6,6	0.66	0
2	PO4	C	608	-	4,4,4	0.83	0	6,6,6	0.57	0
3	GOL	B	702	-	5,5,5	0.35	0	5,5,5	0.31	0
2	PO4	C	605	-	4,4,4	0.94	0	6,6,6	0.62	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
3	GOL	C	701	-	-	2/4/4/4	-
3	GOL	B	702	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	702	GOL	C1-C2-C3-O3
3	C	701	GOL	O2-C2-C3-O3
3	B	702	GOL	O1-C1-C2-C3
3	C	701	GOL	C1-C2-C3-O3
3	B	702	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	701	GOL	3	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	420/462 (90%)	0.08	16 (3%) 44 43	12, 34, 75, 92	8 (1%)
1	B	420/462 (90%)	0.07	9 (2%) 63 63	16, 34, 73, 95	7 (1%)
1	C	417/462 (90%)	-0.01	10 (2%) 59 59	16, 33, 68, 97	10 (2%)
1	D	420/462 (90%)	0.13	14 (3%) 49 48	17, 35, 69, 88	7 (1%)
All	All	1677/1848 (90%)	0.07	49 (2%) 53 52	12, 34, 71, 97	32 (1%)

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	117	ALA	4.5
1	D	379	LEU	4.2
1	D	59	LEU	3.8
1	A	424	LEU	3.8
1	C	116	LEU	3.7

### 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](#)

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( $\text{\AA}^2$ )	Q<0.9
2	PO4	B	612	5/5	0.77	0.09	55,62,71,72	0
3	GOL	C	701	6/6	0.81	0.14	32,44,52,55	0
3	GOL	B	702	6/6	0.85	0.12	57,67,69,78	0
2	PO4	B	607	5/5	0.89	0.14	32,60,65,71	0
2	PO4	C	608	5/5	0.89	0.10	47,52,55,60	0
2	PO4	C	611	5/5	0.90	0.08	42,57,60,65	0
2	PO4	D	606	5/5	0.90	0.14	55,63,66,68	0
2	PO4	A	610	5/5	0.91	0.12	55,55,65,68	0
2	PO4	B	609	5/5	0.92	0.06	57,58,69,69	0
2	PO4	D	602	5/5	0.93	0.11	35,44,55,57	0
2	PO4	A	604	5/5	0.98	0.08	22,22,28,28	0
2	PO4	C	605	5/5	0.98	0.05	23,25,30,35	0
2	PO4	B	603	5/5	0.99	0.02	16,20,25,28	0
2	PO4	D	601	5/5	0.99	0.04	28,28,29,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.