



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:54 PM UTC

PDB ID : 5ROB / pdb\_00005rob  
Title : PanDDA analysis group deposition of ground-state model of SARS-CoV-2 helicase  
Authors : Newman, J.A.; Yosaatmadja, Y.; Douangamath, A.; Aimon, A.; Powell, A.J.; Dias, A.; Fearon, D.; Dunnett, L.; Brandao-Neto, J.; Krojer, T.; Skyner, R.; Gorrie-Stone, T.; Thompson, W.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.; Bountra, C.; Gileadi, O.  
Deposited on : 2020-09-22  
Resolution : 1.87 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](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)

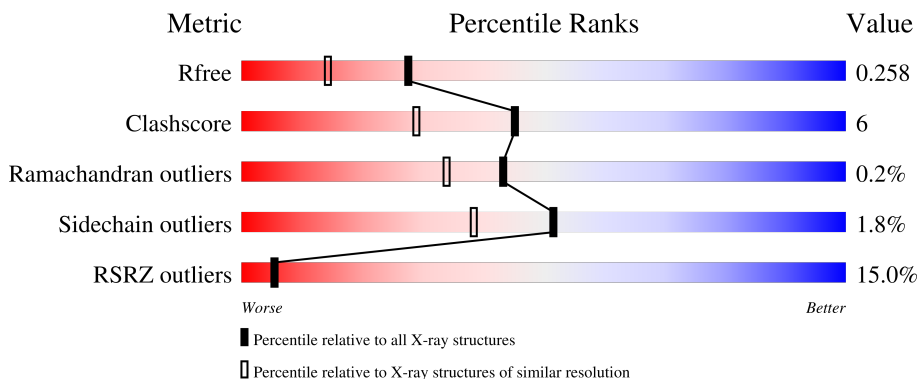
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.87 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	601	
1	B	601	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9407 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 Helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	585	Total 4508	C 2875	N 750	O 848	S 35	0	1	0
1	A	572	Total 4417	C 2816	N 737	O 832	S 32	0	0	0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total 3	Zn 3	0	0
2	A	3	Total 3	Zn 3	0	0

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O<sub>4</sub>P).



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

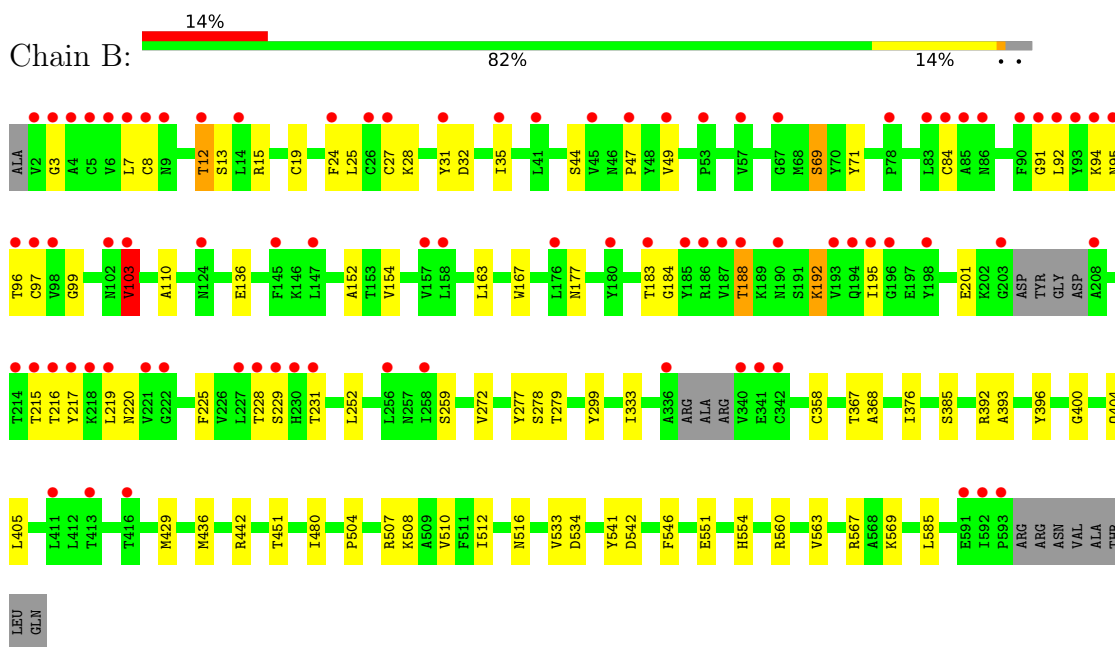
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	250	Total	O	0	0
			250	250		
4	A	206	Total	O	0	0
			206	206		

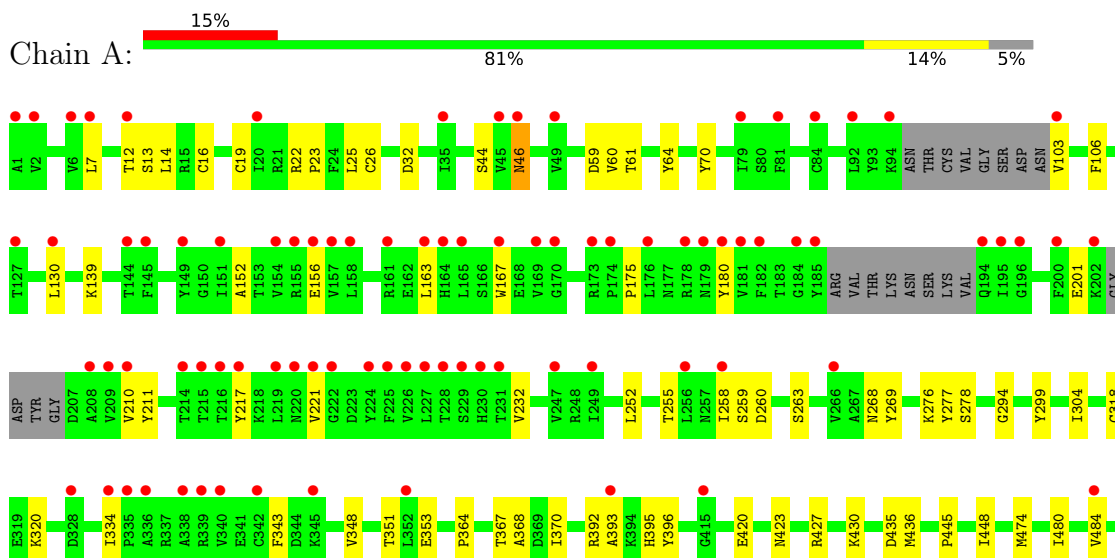
### 3 Residue-property plots [i](#)

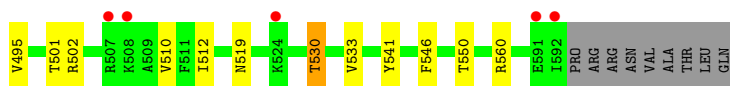
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Helicase



- Molecule 1: Helicase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.24Å 70.28Å 84.62Å 102.40° 95.71° 112.84°	Depositor
Resolution (Å)	80.92 – 1.87 80.92 – 1.87	Depositor EDS
% Data completeness (in resolution range)	94.7 (80.92-1.87) 94.9 (80.92-1.87)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.03 (at 1.87Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.219 , 0.254 0.225 , 0.258	Depositor DCC
$R_{free}$ test set	5109 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.0	Xtrriage
Anisotropy	0.169	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 22.8	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	9407	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.86% 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: PO4, ZN

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	1.06	0/4517	1.36	1/6156 (0.0%)
1	B	1.04	1/4610 (0.0%)	1.37	6/6283 (0.1%)
All	All	1.05	1/9127 (0.0%)	1.37	7/12439 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	272	VAL	C-O	5.02	1.30	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	103	VAL	N-CA-C	-6.04	106.43	112.17
1	A	423	ASN	CA-C-O	-5.98	114.73	121.25
1	B	534	ASP	CB-CA-C	5.29	119.57	110.79
1	B	136	GLU	CB-CA-C	-5.16	102.78	110.88
1	B	94	LYS	CA-C-N	5.04	127.29	120.44
1	B	94	LYS	C-N-CA	5.04	127.29	120.44
1	B	3	GLY	CA-C-O	-5.03	117.07	121.04

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	4417	0	4321	51	1
1	B	4508	0	4423	55	1
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	1	0
4	A	206	0	0	3	0
4	B	250	0	0	11	0
All	All	9407	0	8744	105	1

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.

All (105) 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:A:12:THR:HG21	1:A:25:LEU:O	1.76	0.85
1:B:201:GLU:OE2	4:B:801:HOH:O	1.94	0.84
1:B:12:THR:HG21	1:B:25:LEU:O	1.80	0.81
1:B:279:THR:HB	1:B:429:MET:HE2	1.67	0.75
1:B:183:THR:HG1	1:B:228:THR:HG1	1.31	0.75
1:B:183:THR:OG1	1:B:228:THR:OG1	2.03	0.74
1:B:7:LEU:HD12	1:B:103:VAL:HG22	1.73	0.70
1:B:27:CYS:SG	4:B:816:HOH:O	2.52	0.68
1:B:7:LEU:CD1	1:B:103:VAL:HG22	2.23	0.67
1:B:510:VAL:HG21	1:B:541:TYR:CD1	2.30	0.67
1:B:8:CYS:SG	1:B:99:GLY:N	2.69	0.66
1:A:12:THR:HG22	1:A:14:LEU:H	1.60	0.65
1:A:510:VAL:HG21	1:A:541:TYR:CD1	2.34	0.63
1:A:13:SER:O	1:A:44:SER:HB3	1.99	0.62
1:A:46:ASN:OD1	1:A:46:ASN:N	2.33	0.60
1:B:13:SER:O	1:B:44:SER:HA	2.01	0.60
1:B:279:THR:HB	1:B:429:MET:CE	2.31	0.60
1:B:84:CYS:HB3	4:B:951:HOH:O	2.02	0.59
1:B:47:PRO:O	4:B:803:HOH:O	2.17	0.59
1:A:7:LEU:HD21	1:A:106:PHE:HB2	1.84	0.59
1:B:13:SER:HB2	1:B:92:LEU:HB2	1.86	0.57
1:B:278:SER:HB2	1:B:436:MET:HE2	1.87	0.56
1:A:175:PRO:HG2	1:A:180:TYR:CE1	2.41	0.55
1:A:163:LEU:HD23	1:A:211:TYR:CD2	2.40	0.55
1:A:368:ALA:O	1:A:393:ALA:HA	2.07	0.55
1:A:519:ASN:HB3	1:A:530:THR:HG23	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:ILE:O	1:A:546:PHE:HA	2.08	0.54
1:B:252:LEU:HB3	1:B:299:TYR:CD1	2.43	0.54
1:A:430:LYS:NZ	4:A:806:HOH:O	2.38	0.53
1:B:91:GLY:N	4:B:816:HOH:O	2.40	0.53
1:A:7:LEU:HD21	1:A:130:LEU:HD21	1.92	0.52
1:A:156:GLU:HA	1:A:221:VAL:HG22	1.92	0.52
1:B:512:ILE:O	1:B:546:PHE:HA	2.10	0.52
1:B:69:SER:HB2	1:B:71:TYR:CE2	2.46	0.51
1:B:277:TYR:HA	1:B:396:TYR:O	2.10	0.51
1:A:445:PRO:HD2	1:A:448:ILE:HD12	1.91	0.51
1:A:276:LYS:O	1:A:395:HIS:HA	2.10	0.51
1:A:13:SER:O	1:A:44:SER:CB	2.58	0.51
1:A:320:LYS:NZ	4:A:815:HOH:O	2.43	0.51
1:A:420:GLU:OE1	1:A:427:ARG:NH1	2.44	0.51
1:A:480:ILE:HG12	1:A:550:THR:HG22	1.93	0.50
1:B:177:ASN:HB3	1:B:516:ASN:ND2	2.28	0.49
1:A:318:CYS:HB3	1:A:343:PHE:CD2	2.47	0.49
1:B:385:SER:OG	4:B:805:HOH:O	2.20	0.48
1:B:219:LEU:O	1:B:219:LEU:HD12	2.14	0.48
1:B:404:GLN:NE2	3:B:705:PO4:O2	2.35	0.48
1:A:480:ILE:HG21	1:A:550:THR:HG22	1.95	0.48
1:B:152:ALA:HB2	1:B:167:TRP:CZ3	2.50	0.47
1:A:32:ASP:HB2	1:A:103:VAL:HG11	1.94	0.47
1:B:508:LYS:HD3	4:B:998:HOH:O	2.15	0.47
1:A:64:TYR:O	1:A:70:TYR:HA	2.15	0.46
1:A:351:THR:HG23	1:A:364:PRO:HG3	1.98	0.46
1:B:376:ILE:HG22	1:B:400:GLY:HA3	1.98	0.46
1:A:519:ASN:HB3	1:A:530:THR:CG2	2.45	0.46
1:A:278:SER:HA	1:A:435:ASP:OD1	2.16	0.46
1:B:103:VAL:CG1	1:B:103:VAL:O	2.64	0.45
1:B:480:ILE:HD11	4:B:875:HOH:O	2.15	0.45
1:B:19:CYS:SG	1:B:110:ALA:HB1	2.56	0.45
1:B:404:GLN:NE2	4:B:837:HOH:O	2.50	0.45
1:A:60:VAL:HB	4:A:882:HOH:O	2.15	0.45
1:B:217:TYR:HD1	1:A:217:TYR:CE1	2.35	0.45
1:B:551:GLU:HG3	4:B:907:HOH:O	2.17	0.45
1:B:154:VAL:HG13	1:B:163:LEU:HD22	1.98	0.45
1:A:152:ALA:HB2	1:A:167:TRP:CZ3	2.52	0.45
1:A:269:TYR:OH	1:A:294:GLY:HA3	2.16	0.45
1:B:451:THR:HG21	1:B:585:LEU:HD23	2.00	0.44
1:A:304:ILE:HA	1:A:370:ILE:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:THR:O	1:B:217:TYR:CD2	2.71	0.44
1:A:353:GLU:HA	1:A:353:GLU:OE2	2.18	0.44
1:B:533:VAL:HG11	1:B:560:ARG:O	2.17	0.44
1:B:367:THR:HA	1:B:392:ARG:O	2.18	0.44
1:B:184:GLY:C	1:B:195:ILE:HG22	2.44	0.43
1:A:277:TYR:HA	1:A:396:TYR:O	2.18	0.43
1:B:333:ILE:HB	1:B:358:CYS:SG	2.58	0.43
1:A:12:THR:HG22	1:A:14:LEU:N	2.31	0.43
1:A:318:CYS:HB3	1:A:343:PHE:CE2	2.53	0.43
1:B:31:TYR:CZ	1:B:35:ILE:HG21	2.53	0.43
1:A:16:CYS:O	1:A:22:ARG:HA	2.18	0.43
1:B:405:LEU:HD21	1:B:560:ARG:HA	2.00	0.43
1:B:47:PRO:O	1:B:49:VAL:HG12	2.19	0.42
1:B:554:HIS:ND1	4:B:811:HOH:O	2.37	0.42
1:A:139:LYS:HG2	1:A:232:VAL:HG22	2.01	0.42
1:B:28:LYS:O	1:B:32:ASP:OD1	2.37	0.42
1:A:260:ASP:HA	1:A:263:SER:OG	2.19	0.42
1:A:304:ILE:HG12	1:A:370:ILE:HB	2.01	0.42
1:A:334:ILE:HD12	1:A:348:VAL:HG13	2.01	0.42
1:A:19:CYS:HB2	1:A:23:PRO:HD2	2.02	0.42
1:A:59:ASP:CG	1:A:61:THR:HG1	2.23	0.42
1:B:368:ALA:O	1:B:393:ALA:HA	2.19	0.42
1:B:563:VAL:O	1:B:567:ARG:NE	2.51	0.42
1:B:376:ILE:HD12	1:B:376:ILE:HA	1.95	0.42
1:A:533:VAL:HG11	1:A:560:ARG:O	2.20	0.42
1:A:268:ASN:HB3	1:A:436:MET:SD	2.60	0.41
1:A:474:MET:SD	1:A:495:VAL:HG11	2.60	0.41
1:B:183:THR:O	1:B:225:PHE:HA	2.21	0.41
1:A:12:THR:OG1	1:A:26:CYS:HA	2.19	0.41
1:A:252:LEU:HB3	1:A:299:TYR:CD2	2.55	0.41
1:A:367:THR:HA	1:A:392:ARG:O	2.20	0.41
1:B:15:ARG:HG3	1:B:24:PHE:CD2	2.56	0.41
1:B:504:PRO:HB3	1:B:507:ARG:NH2	2.36	0.41
1:B:542:ASP:OD1	1:B:569:LYS:HE3	2.20	0.41
1:A:13:SER:OG	1:A:44:SER:HB2	2.20	0.40
1:B:195:ILE:HG23	1:B:195:ILE:O	2.21	0.40
1:A:201:GLU:O	1:A:210:VAL:N	2.43	0.40
1:B:188:THR:HG23	1:B:192:LYS:HA	2.03	0.40

All (1) 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:B:442:ARG:NH1	1:A:501:THR:O[1_666]	2.11	0.09

### 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	564/601 (94%)	543 (96%)	20 (4%)	1 (0%)	43	34
1	B	580/601 (96%)	545 (94%)	34 (6%)	1 (0%)	43	34
All	All	1144/1202 (95%)	1088 (95%)	54 (5%)	2 (0%)	43	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	97	CYS
1	A	484	VAL

#### 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	485/523 (93%)	479 (99%)	6 (1%)	63	53
1	B	498/523 (95%)	486 (98%)	12 (2%)	43	27
All	All	983/1046 (94%)	965 (98%)	18 (2%)	51	38

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	12	THR
1	B	69	SER
1	B	95	ASN
1	B	96	THR
1	B	103	VAL
1	B	188	THR
1	B	192	LYS
1	B	215	THR
1	B	220	ASN
1	B	229	SER
1	B	231	THR
1	B	259	SER
1	A	46	ASN
1	A	255	THR
1	A	258	ILE
1	A	259	SER
1	A	502	ARG
1	A	530	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	46	ASN
1	B	88	GLN
1	B	179	ASN
1	B	268	ASN
1	B	404	GLN
1	B	459	ASN
1	B	470	GLN
1	B	482	HIS
1	A	51	ASN
1	A	257	ASN
1	A	404	GLN
1	A	459	ASN
1	A	531	GLN
1	A	586	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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	$\# Z  > 2$
3	PO4	A	705	-	4,4,4	0.76	0	6,6,6	0.53	0
3	PO4	B	704	-	4,4,4	1.09	0	6,6,6	0.45	0
3	PO4	B	705	-	4,4,4	1.46	0	6,6,6	0.54	0
3	PO4	A	704	-	4,4,4	0.85	0	6,6,6	0.71	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	705	PO4	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 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	572/601 (95%)	0.87	91 (15%) <b>5</b> <b>5</b>	26, 44, 82, 114	0
1	B	585/601 (97%)	0.78	83 (14%) <b>6</b> <b>6</b>	20, 40, 78, 114	1 (0%)
All	All	1157/1202 (96%)	0.82	174 (15%) <b>5</b> <b>5</b>	20, 42, 81, 114	1 (0%)

All (174) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	219	LEU	8.3
1	A	592	ILE	7.9
1	B	214	THR	7.0
1	B	187	VAL	6.7
1	A	94	LYS	6.0
1	B	217	TYR	5.5
1	A	228	THR	5.3
1	A	169	VAL	5.3
1	A	157	VAL	5.2
1	B	91	GLY	5.2
1	B	340	VAL	5.0
1	B	193	VAL	4.9
1	A	81	PHE	4.9
1	A	336	ALA	4.9
1	B	219	LEU	4.7
1	B	2	VAL	4.5
1	B	7	LEU	4.5
1	B	592	ILE	4.5
1	B	593	PRO	4.5
1	A	247	VAL	4.4
1	B	195	ILE	4.3
1	B	336	ALA	4.3
1	A	215	THR	4.2
1	A	225	PHE	4.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	338	ALA	4.2
1	A	195	ILE	4.0
1	B	157	VAL	4.0
1	B	98	VAL	4.0
1	B	93	TYR	3.8
1	A	216	THR	3.7
1	B	92	LEU	3.7
1	A	149	TYR	3.6
1	B	230	HIS	3.6
1	A	1	ALA	3.6
1	A	229	SER	3.6
1	B	203	GLY	3.6
1	A	200	PHE	3.6
1	B	14	LEU	3.5
1	A	151	ILE	3.5
1	B	53	PRO	3.4
1	A	185	TYR	3.4
1	A	170	GLY	3.4
1	B	218	LYS	3.3
1	A	256	LEU	3.3
1	B	96	THR	3.2
1	A	92	LEU	3.2
1	A	227	LEU	3.2
1	A	156	GLU	3.2
1	B	186	ARG	3.2
1	B	158	LEU	3.2
1	B	4	ALA	3.2
1	B	26	CYS	3.2
1	B	41	LEU	3.1
1	B	97	CYS	3.1
1	B	85	ALA	3.1
1	B	78	PRO	3.1
1	B	258	ILE	3.1
1	A	182	PHE	3.1
1	B	6	VAL	3.1
1	A	221	VAL	3.1
1	B	228	THR	3.1
1	A	127	THR	3.1
1	A	224	TYR	3.0
1	B	208	ALA	3.0
1	A	167	TRP	3.0
1	B	231	THR	3.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	94	LYS	3.0
1	A	154	VAL	3.0
1	A	217	TYR	3.0
1	A	226	VAL	3.0
1	B	67	GLY	3.0
1	A	176	LEU	2.9
1	A	6	VAL	2.9
1	B	102	ASN	2.9
1	A	155	ARG	2.9
1	A	230	HIS	2.8
1	B	215	THR	2.8
1	B	45	VAL	2.8
1	B	124	ASN	2.8
1	B	227	LEU	2.8
1	A	194	GLN	2.8
1	A	181	VAL	2.8
1	B	90	PHE	2.7
1	A	202	LYS	2.7
1	A	222	GLY	2.7
1	B	194	GLN	2.7
1	A	352	LEU	2.7
1	A	45	VAL	2.7
1	B	183	THR	2.7
1	B	12	THR	2.6
1	B	196	GLY	2.6
1	A	46	ASN	2.6
1	A	328	ASP	2.6
1	B	3	GLY	2.6
1	A	345	LYS	2.6
1	A	209	VAL	2.6
1	B	8	CYS	2.5
1	B	185	TYR	2.5
1	B	147	LEU	2.5
1	A	20	ILE	2.5
1	A	249	ILE	2.5
1	A	334	ILE	2.5
1	B	416	THR	2.5
1	A	214	THR	2.5
1	A	145	PHE	2.5
1	A	79	ILE	2.5
1	A	340	VAL	2.5
1	B	221	VAL	2.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	178	ARG	2.4
1	A	180	TYR	2.4
1	A	339	ARG	2.4
1	A	7	LEU	2.4
1	B	413	THR	2.4
1	A	220	ASN	2.4
1	B	84	CYS	2.4
1	A	163	LEU	2.4
1	A	524	LYS	2.4
1	B	198	TYR	2.4
1	B	188	THR	2.4
1	A	144	THR	2.4
1	B	49	VAL	2.4
1	A	196	GLY	2.4
1	A	342	CYS	2.4
1	A	173	ARG	2.3
1	B	145	PHE	2.3
1	B	86	ASN	2.3
1	B	229	SER	2.3
1	B	341	GLU	2.3
1	B	176	LEU	2.3
1	B	9	ASN	2.3
1	A	231	THR	2.3
1	A	210	VAL	2.3
1	A	130	LEU	2.3
1	B	222	GLY	2.2
1	A	415	GLY	2.2
1	A	179	ASN	2.2
1	B	216	THR	2.2
1	A	161	ARG	2.2
1	A	2	VAL	2.2
1	A	184	GLY	2.2
1	B	47	PRO	2.2
1	A	208	ALA	2.2
1	A	35	ILE	2.2
1	A	507	ARG	2.2
1	A	174	PRO	2.2
1	B	190	ASN	2.1
1	B	35	ILE	2.1
1	A	49	VAL	2.1
1	A	266	VAL	2.1
1	B	342	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	158	LEU	2.1
1	B	95	ASN	2.1
1	B	31	TYR	2.1
1	A	258	ILE	2.1
1	B	57	VAL	2.1
1	B	103	VAL	2.1
1	A	103	VAL	2.1
1	A	484	VAL	2.1
1	A	164	HIS	2.1
1	B	83	LEU	2.1
1	B	256	LEU	2.1
1	B	591	GLU	2.1
1	A	591	GLU	2.1
1	A	335	PRO	2.1
1	B	5	CYS	2.1
1	B	27	CYS	2.1
1	A	165	LEU	2.1
1	B	24	PHE	2.0
1	B	411	LEU	2.0
1	A	84	CYS	2.0
1	A	393	ALA	2.0
1	A	508	LYS	2.0
1	A	12	THR	2.0
1	B	180	TYR	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](#)

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
3	PO4	B	705	5/5	0.95	0.08	41,42,48,59	0
3	PO4	B	704	5/5	0.96	0.07	32,35,40,40	0
2	ZN	B	702	1/1	0.98	0.07	52,52,52,52	0
3	PO4	A	704	5/5	0.98	0.04	32,35,37,39	0
3	PO4	A	705	5/5	0.98	0.05	30,34,37,38	0
2	ZN	B	701	1/1	0.99	0.03	37,37,37,37	0
2	ZN	B	703	1/1	0.99	0.03	64,64,64,64	0
2	ZN	A	702	1/1	0.99	0.02	39,39,39,39	0
2	ZN	A	703	1/1	0.99	0.04	57,57,57,57	0
2	ZN	A	701	1/1	1.00	0.03	48,48,48,48	0

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

There are no such residues in this entry.