



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 8, 2018 – 08:14 PM EST

PDB ID : 5TIQ  
Title : The Structure of the Major Capsid protein of PBCV-1  
Authors : Klose, T.; De Castro, C.; Speciale, I.; Molinaro, A.; Van Etten, J.L.; Rossmann, M.G.  
Deposited on : 2016-10-03  
Resolution : 2.54 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030736

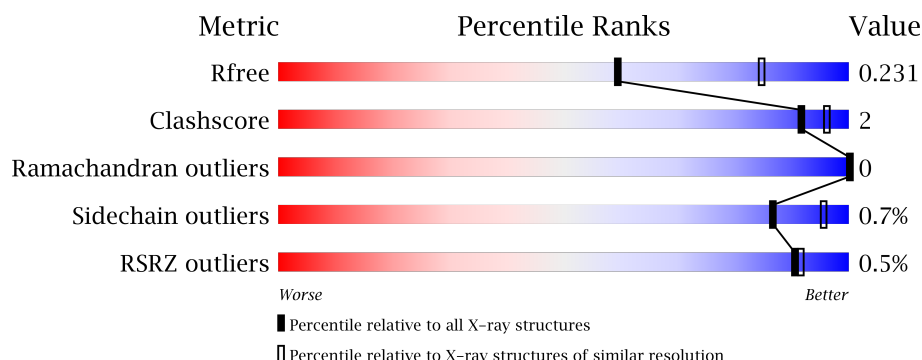
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

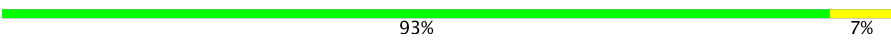
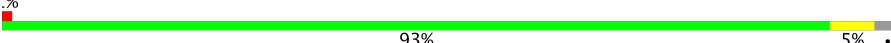
The reported resolution of this entry is 2.54 Å.

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))
$R_{free}$	100719	4993 (2.58-2.50)
Clashscore	112137	5755 (2.58-2.50)
Ramachandran outliers	110173	5652 (2.58-2.50)
Sidechain outliers	110143	5654 (2.58-2.50)
RSRZ outliers	101464	5026 (2.58-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	436	 93% 7%
1	B	436	 93% 5% 2%

## 2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 7633 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.

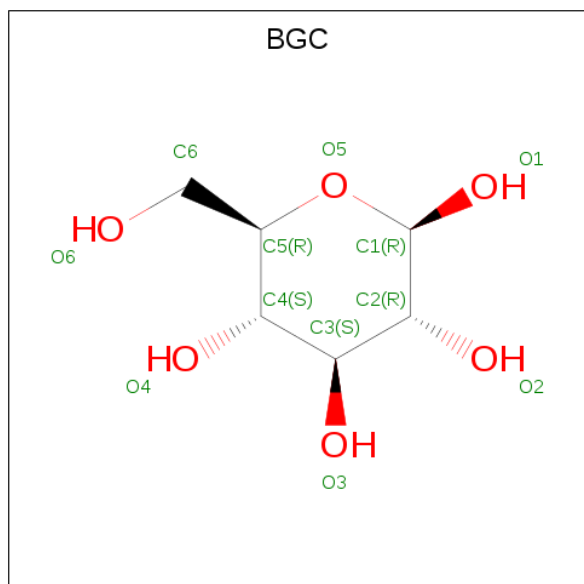
- Molecule 1 is a protein called Major capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	436	Total	C	N	O	S	0	0	0
			3396	2156	576	656	8			
1	B	425	Total	C	N	O	S	0	0	0
			3324	2110	564	642	8			

- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

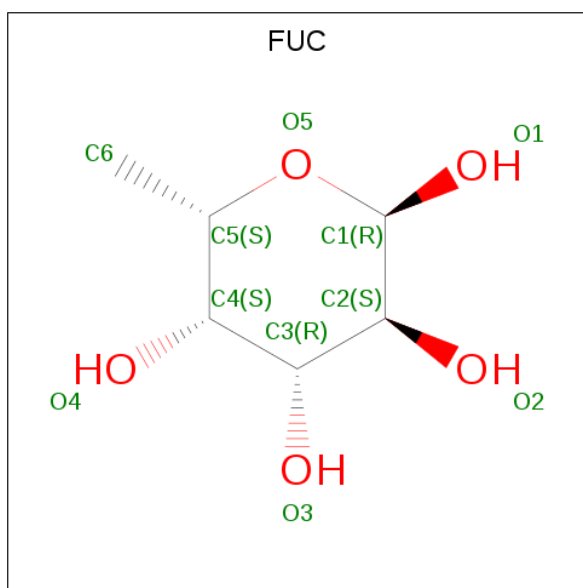
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Hg	0	0
			2	2		
2	A	2	Total	Hg	0	0
			2	2		

- Molecule 3 is BETA-D-GLUCOSE (three-letter code: BGC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	6	5		
3	A	1	Total	C	O	0	0
			11	6	5		
3	A	1	Total	C	O	0	0
			11	6	5		
3	A	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>5</sub>).



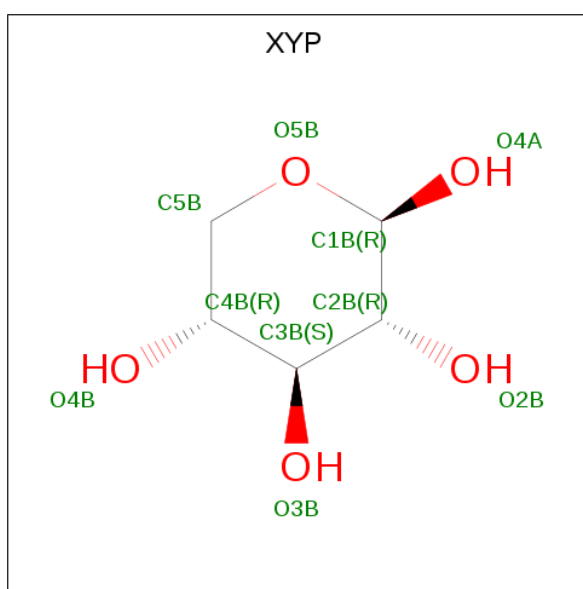
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is BETA-D-XYLOPYRANOSE (three-letter code: XYP) (formula: C<sub>5</sub>H<sub>10</sub>O<sub>5</sub>).



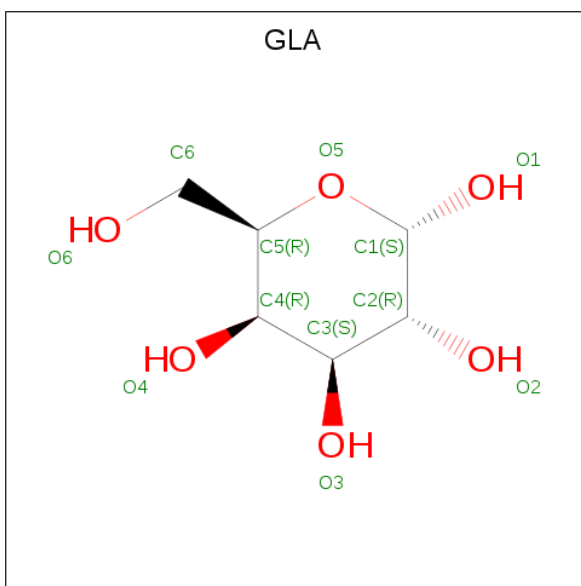
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			9	5	4		
5	A	1	Total	C	O	0	0
			9	5	4		
5	A	1	Total	C	O	0	0
			9	5	4		
5	A	1	Total	C	O	0	0
			9	5	4		
5	A	1	Total	C	O	0	0
			9	5	4		
5	A	1	Total	C	O	0	0
			9	5	4		
5	B	1	Total	C	O	0	0
			9	5	4		

*Continued on next page...*

Continued from previous page...

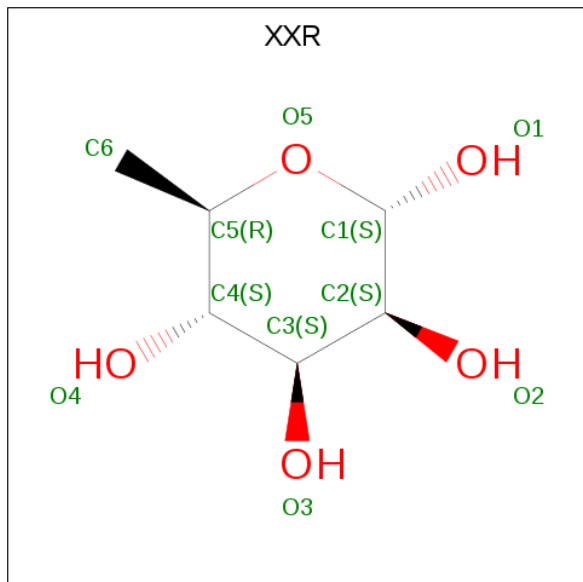
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			9	5	4		
5	B	1	Total	C	O	0	0
			9	5	4		
5	B	1	Total	C	O	0	0
			9	5	4		
5	B	1	Total	C	O	0	0
			9	5	4		
5	B	1	Total	C	O	0	0
			9	5	4		

- Molecule 6 is ALPHA D-GALACTOSE (three-letter code: GLA) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 7 is 6-deoxy-alpha-D-mannopyranose (three-letter code: XXR) (formula:  $C_6H_{12}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			10	6	4		
7	A	1	Total	C	O	0	0
			10	6	4		
7	A	1	Total	C	O	0	0
			10	6	4		
7	B	1	Total	C	O	0	0
			10	6	4		
7	B	1	Total	C	O	0	0
			10	6	4		
7	B	1	Total	C	O	0	0
			10	6	4		

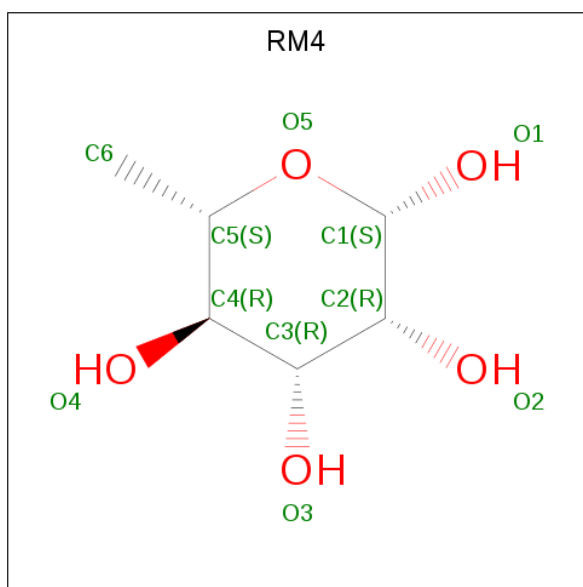
- Molecule 8 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula:  $C_6H_{12}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			11	6	5		
8	A	1	Total	C	O	0	0
			11	6	5		
8	A	1	Total	C	O	0	0
			11	6	5		
8	B	1	Total	C	O	0	0
			11	6	5		
8	B	1	Total	C	O	0	0
			11	6	5		
8	B	1	Total	C	O	0	0
			11	6	5		

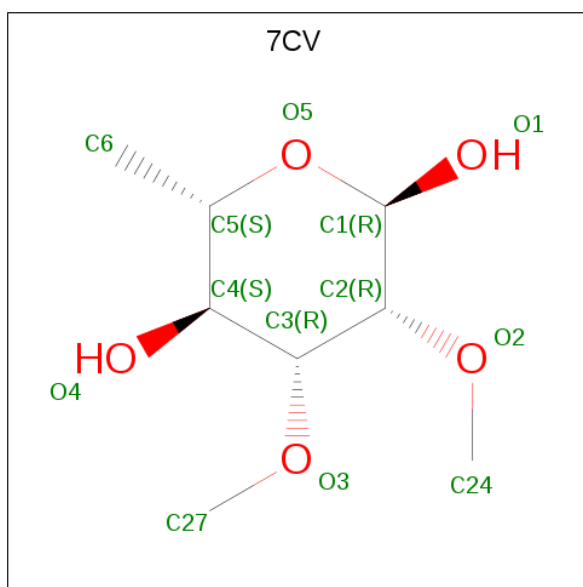
- Molecule 9 is 6-deoxy-beta-L-mannopyranose (three-letter code: RM4) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>5</sub>).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			10	6	4		
9	A	1	Total	C	O	0	0
			10	6	4		
9	B	1	Total	C	O	0	0
			10	6	4		
9	B	1	Total	C	O	0	0
			10	6	4		
9	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 10 is 6-deoxy-2,3-di-O-methyl-alpha-L-mannopyranose (three-letter code: 7CV) (formula: C<sub>8</sub>H<sub>16</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			12	8	4		
10	A	1	Total	C	O	0	0
			12	8	4		
10	B	1	Total	C	O	0	0
			12	8	4		
10	B	1	Total	C	O	0	0
			12	8	4		
10	B	1	Total	C	O	0	0
			12	8	4		

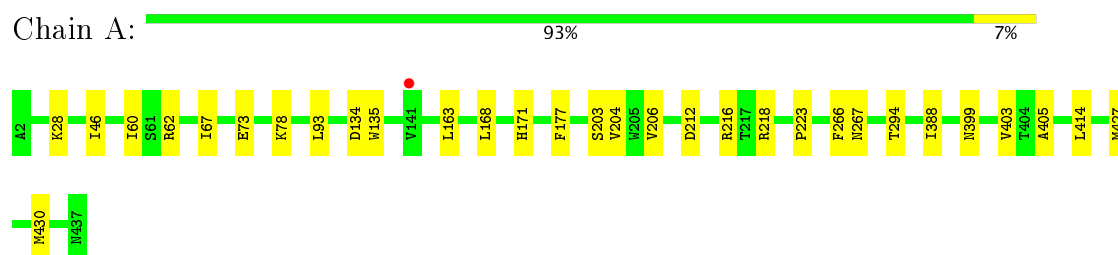
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	166	Total	O	0	0
			166	166		
11	B	165	Total	O	0	0
			165	165		

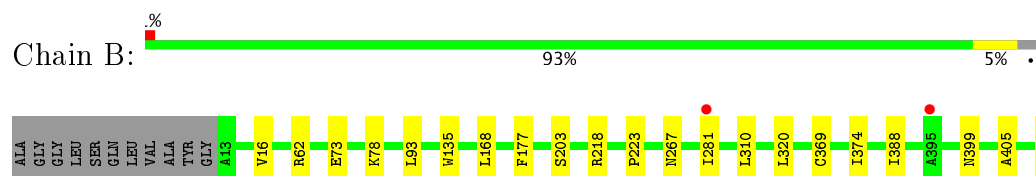
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Major capsid protein



- Molecule 1: Major capsid protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	188.76 Å   188.76 Å   188.76 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	43.30 – 2.54 84.42 – 2.54	Depositor EDS
% Data completeness (in resolution range)	99.3 (43.30-2.54) 99.6 (84.42-2.54)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	9.87 (at 2.55 Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, $R_{free}$	0.179   ,   0.223 0.184   ,   0.231	Depositor DCC
$R_{free}$ test set	1516 reflections (3.95%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.8	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	7633	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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: XYP, XXR, BGC, RM4, 7CV, GLA, FUC, HG, MAN

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	0.26	0/3474	0.45	0/4735
1	B	0.26	0/3401	0.45	0/4636
All	All	0.26	0/6875	0.45	0/9371

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

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	3396	0	3283	16	0
1	B	3324	0	3211	12	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	44	0	33	0	0
3	B	44	0	33	0	0
4	A	40	0	31	0	0
4	B	40	0	31	0	0
5	A	54	0	46	0	0
5	B	54	0	45	0	0
6	A	33	0	30	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	33	0	30	0	0
7	A	30	0	27	0	0
7	B	30	0	27	0	0
8	A	33	0	30	0	0
8	B	33	0	30	0	0
9	A	20	0	18	0	0
9	B	30	0	27	0	0
10	A	24	0	0	0	0
10	B	36	0	0	0	0
11	A	166	0	0	0	0
11	B	165	0	0	2	0
All	All	7633	0	6932	27	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 2.

All (27) 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:B:369:CYS:SG	11:B:756:HOH:O	2.33	0.85
1:B:62:ARG:NH2	1:B:168:LEU:O	2.24	0.70
1:A:218:ARG:HG2	1:B:16:VAL:HG13	1.76	0.68
1:A:62:ARG:NH2	1:A:168:LEU:O	2.30	0.64
1:B:281:ILE:HD11	1:B:310:LEU:HD11	1.84	0.60
1:B:320:LEU:HD22	1:B:374:ILE:HD13	1.85	0.59
1:B:218:ARG:NE	11:B:611:HOH:O	2.43	0.52
1:B:73:GLU:HB2	1:B:203:SER:HB2	1.97	0.47
1:A:60:ILE:HD12	1:A:206:VAL:HG21	1.96	0.47
1:A:93:LEU:HD11	1:A:177:PHE:HD1	1.81	0.45
1:B:78:LYS:HG3	1:B:135:TRP:CE2	2.50	0.45
1:A:388:ILE:HD11	1:A:403:VAL:HG23	1.99	0.45
1:B:93:LEU:HD11	1:B:177:PHE:HD1	1.80	0.45
1:B:369:CYS:SG	1:B:374:ILE:HD11	2.56	0.44
1:B:388:ILE:HB	1:B:405:ALA:HA	1.98	0.44
1:B:223:PRO:HB3	1:B:427:MET:HG3	1.99	0.44
1:A:134:ASP:OD1	1:A:135:TRP:N	2.51	0.43
1:A:212:ASP:O	1:A:216:ARG:HB2	2.19	0.43
1:A:171:HIS:CE1	1:A:430:MET:HB3	2.53	0.42
1:A:78:LYS:HG3	1:A:135:TRP:CE2	2.56	0.41
1:A:266:PHE:CE2	1:A:414:LEU:HD13	2.56	0.41
1:A:28:LYS:HA	1:A:28:LYS:HD2	1.91	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ILE:HD11	1:A:163:LEU:HD23	2.03	0.41
1:A:46:ILE:HD11	1:A:204:VAL:HG23	2.03	0.41
1:A:223:PRO:HB3	1:A:427:MET:HG3	2.03	0.41
1:A:73:GLU:HB2	1:A:203:SER:HB3	2.02	0.41
1:A:388:ILE:HB	1:A:405:ALA:HA	2.04	0.40

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	434/436 (100%)	417 (96%)	17 (4%)	0	100	100
1	B	423/436 (97%)	408 (96%)	15 (4%)	0	100	100
All	All	857/872 (98%)	825 (96%)	32 (4%)	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	357/357 (100%)	354 (99%)	3 (1%)	85	94
1	B	351/357 (98%)	349 (99%)	2 (1%)	89	96
All	All	708/714 (99%)	703 (99%)	5 (1%)	87	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	A	294	THR
1	A	399	ASN
1	B	267	ASN
1	B	399	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 60 ligands modelled in this entry, 4 are monoatomic - leaving 56 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	BGC	A	503	1,5,4	11,11,12	0.34	0	13,15,17	0.88	0
4	FUC	A	504	3,5,7,6	9,10,11	0.38	0	13,14,16	1.10	0
5	XYP	A	505	9,4	9,9,10	0.22	0	10,12,14	0.54	0
5	XYP	A	506	3	9,9,10	0.20	0	10,12,14	0.54	0
6	GLA	A	507	4	11,11,12	0.24	0	13,15,17	0.47	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	XXR	A	508	8,4	9,10,11	0.18	0	13,14,16	0.41	0
8	MAN	A	509	7	11,11,12	0.27	0	13,15,17	0.76	0
9	RM4	A	510	10,5	9,10,11	0.28	0	13,14,16	0.52	0
10	7CV	A	511	9	11,12,13	0.69	0	15,16,18	0.35	0
3	BGC	A	512	1,5,4	11,11,12	0.40	0	13,15,17	0.82	0
4	FUC	A	513	3,5,7,6	9,10,11	0.28	0	13,14,16	0.76	0
5	XYP	A	514	9,4	9,9,10	0.20	0	10,12,14	0.41	0
5	XYP	A	515	3	9,9,10	0.22	0	10,12,14	0.47	0
6	GLA	A	516	4	11,11,12	0.24	0	13,15,17	0.47	0
7	XXR	A	517	8,4	9,10,11	0.18	0	13,14,16	0.35	0
8	MAN	A	518	7	11,11,12	0.24	0	13,15,17	0.65	0
9	RM4	A	519	10,5	9,10,11	0.29	0	13,14,16	0.52	0
10	7CV	A	520	9	11,12,13	0.93	0	15,16,18	0.67	0
3	BGC	A	521	1,5,4	11,11,12	0.52	0	13,15,17	0.71	0
4	FUC	A	522	3,5,7,6	9,10,11	0.67	0	13,14,16	0.75	0
5	XYP	A	523	4	9,9,10	0.36	0	10,12,14	0.71	0
5	XYP	A	524	3	9,9,10	0.21	0	10,12,14	0.52	0
6	GLA	A	525	4	11,11,12	0.24	0	13,15,17	0.47	0
7	XXR	A	526	8,4	9,10,11	0.19	0	13,14,16	0.30	0
8	MAN	A	527	7	11,11,12	0.32	0	13,15,17	0.69	0
3	BGC	A	528	1,4	11,11,12	0.39	0	13,15,17	0.93	0
4	FUC	A	529	3	9,10,11	0.29	0	13,14,16	0.70	0
3	BGC	B	503	1,5,4	11,11,12	0.76	0	13,15,17	0.88	0
4	FUC	B	504	3,5,7,6	9,10,11	0.79	0	13,14,16	0.75	0
5	XYP	B	505	9,4	9,9,10	0.20	0	10,12,14	0.38	0
5	XYP	B	506	3	9,9,10	0.27	0	10,12,14	0.57	0
6	GLA	B	507	4	11,11,12	0.22	0	13,15,17	0.60	0
7	XXR	B	508	8,4	9,10,11	0.19	0	13,14,16	0.45	0
8	MAN	B	509	7	11,11,12	0.24	0	13,15,17	0.71	0
9	RM4	B	510	10,5	9,10,11	0.29	0	13,14,16	0.52	0
10	7CV	B	511	9	11,12,13	0.70	0	15,16,18	0.33	0
3	BGC	B	512	1,5,4	11,11,12	0.35	0	13,15,17	0.78	0
4	FUC	B	513	3,5,7,6	9,10,11	0.28	0	13,14,16	0.77	0
5	XYP	B	514	9,4	9,9,10	0.21	0	10,12,14	0.40	0
5	XYP	B	515	3	9,9,10	0.21	0	10,12,14	0.52	0
6	GLA	B	516	4	11,11,12	0.25	0	13,15,17	0.48	0
7	XXR	B	517	8,4	9,10,11	0.20	0	13,14,16	0.29	0
8	MAN	B	518	7	11,11,12	0.26	0	13,15,17	0.54	0
9	RM4	B	519	10,5	9,10,11	0.29	0	13,14,16	0.52	0
10	7CV	B	520	9	11,12,13	0.69	0	15,16,18	0.26	0
3	BGC	B	521	1,5,4	11,11,12	0.60	0	13,15,17	0.76	0
4	FUC	B	522	3,5,7,6	9,10,11	0.48	0	13,14,16	0.82	0
5	XYP	B	523	9,4	9,9,10	0.20	0	10,12,14	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	XYP	B	524	3	9,9,10	0.21	0	10,12,14	0.44	0
6	GLA	B	525	4	11,11,12	0.24	0	13,15,17	0.47	0
7	XXR	B	526	8,4	9,10,11	0.20	0	13,14,16	0.30	0
8	MAN	B	527	7	11,11,12	0.40	0	13,15,17	0.71	0
9	RM4	B	528	10,5	9,10,11	0.29	0	13,14,16	0.52	0
10	7CV	B	529	9	11,12,13	0.69	0	15,16,18	0.26	0
3	BGC	B	530	1,4	11,11,12	0.39	0	13,15,17	0.93	0
4	FUC	B	531	3	9,10,11	0.28	0	13,14,16	0.65	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	BGC	A	503	1,5,4	-	0/2/19/22	0/1/1/1
4	FUC	A	504	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	A	505	9,4	-	0/0/14/17	0/1/1/1
5	XYP	A	506	3	-	0/0/14/17	0/1/1/1
6	GLA	A	507	4	-	0/2/19/22	0/1/1/1
7	XXR	A	508	8,4	-	0/0/17/20	0/1/1/1
8	MAN	A	509	7	-	0/2/19/22	0/1/1/1
9	RM4	A	510	10,5	-	0/0/17/20	0/1/1/1
10	7CV	A	511	9	-	0/4/21/24	0/1/1/1
3	BGC	A	512	1,5,4	-	0/2/19/22	0/1/1/1
4	FUC	A	513	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	A	514	9,4	-	0/0/14/17	0/1/1/1
5	XYP	A	515	3	-	0/0/14/17	0/1/1/1
6	GLA	A	516	4	-	0/2/19/22	0/1/1/1
7	XXR	A	517	8,4	-	0/0/17/20	0/1/1/1
8	MAN	A	518	7	-	0/2/19/22	0/1/1/1
9	RM4	A	519	10,5	-	0/0/17/20	0/1/1/1
10	7CV	A	520	9	-	0/4/21/24	0/1/1/1
3	BGC	A	521	1,5,4	-	0/2/19/22	0/1/1/1
4	FUC	A	522	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	A	523	4	-	0/0/14/17	0/1/1/1
5	XYP	A	524	3	-	0/0/14/17	0/1/1/1
6	GLA	A	525	4	-	0/2/19/22	0/1/1/1
7	XXR	A	526	8,4	-	0/0/17/20	0/1/1/1
8	MAN	A	527	7	-	0/2/19/22	0/1/1/1
3	BGC	A	528	1,4	-	0/2/19/22	0/1/1/1
4	FUC	A	529	3	-	0/0/17/20	0/1/1/1
3	BGC	B	503	1,5,4	-	0/2/19/22	0/1/1/1

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FUC	B	504	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	B	505	9,4	-	0/0/14/17	0/1/1/1
5	XYP	B	506	3	-	0/0/14/17	0/1/1/1
6	GLA	B	507	4	-	0/2/19/22	0/1/1/1
7	XXR	B	508	8,4	-	0/0/17/20	0/1/1/1
8	MAN	B	509	7	-	0/2/19/22	0/1/1/1
9	RM4	B	510	10,5	-	0/0/17/20	0/1/1/1
10	7CV	B	511	9	-	0/4/21/24	0/1/1/1
3	BGC	B	512	1,5,4	-	0/2/19/22	0/1/1/1
4	FUC	B	513	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	B	514	9,4	-	0/0/14/17	0/1/1/1
5	XYP	B	515	3	-	0/0/14/17	0/1/1/1
6	GLA	B	516	4	-	0/2/19/22	0/1/1/1
7	XXR	B	517	8,4	-	0/0/17/20	0/1/1/1
8	MAN	B	518	7	-	0/2/19/22	0/1/1/1
9	RM4	B	519	10,5	-	0/0/17/20	0/1/1/1
10	7CV	B	520	9	-	0/4/21/24	0/1/1/1
3	BGC	B	521	1,5,4	-	0/2/19/22	0/1/1/1
4	FUC	B	522	3,5,7,6	-	0/0/17/20	0/1/1/1
5	XYP	B	523	9,4	-	0/0/14/17	0/1/1/1
5	XYP	B	524	3	-	0/0/14/17	0/1/1/1
6	GLA	B	525	4	-	0/2/19/22	0/1/1/1
7	XXR	B	526	8,4	-	0/0/17/20	0/1/1/1
8	MAN	B	527	7	-	0/2/19/22	0/1/1/1
9	RM4	B	528	10,5	-	0/0/17/20	0/1/1/1
10	7CV	B	529	9	-	0/4/21/24	0/1/1/1
3	BGC	B	530	1,4	-	0/2/19/22	0/1/1/1
4	FUC	B	531	3	-	0/0/17/20	0/1/1/1

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.

No monomer is involved in short contacts.

## 5.7 Other polymers

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 [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	436/436 (100%)	-0.22	1 (0%) 94 95	23, 35, 51, 64	0
1	B	425/436 (97%)	-0.17	3 (0%) 87 89	20, 36, 54, 85	0
All	All	861/872 (98%)	-0.19	4 (0%) 90 91	20, 35, 52, 85	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	437	ASN	3.9
1	A	141	VAL	2.6
1	B	281	ILE	2.2
1	B	395	ALA	2.1

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	BGC	B	530	11/12	0.83	0.26	1.98	76,79,82,87	0
3	BGC	A	528	11/12	0.87	0.23	1.81	41,59,65,66	0
6	GLA	A	516	11/12	0.94	0.21	1.75	32,34,40,43	0
3	BGC	B	512	11/12	0.97	0.16	0.84	28,35,41,41	0
3	BGC	A	521	11/12	0.93	0.17	0.12	34,42,45,46	0
6	GLA	B	516	11/12	0.92	0.14	-0.03	37,41,47,49	0
6	GLA	A	525	11/12	0.95	0.18	-0.17	34,36,43,45	0
3	BGC	A	512	11/12	0.94	0.13	-0.26	30,40,46,46	0
2	HG	A	501	1/1	0.99	0.12	-0.57	42,42,42,42	1
3	BGC	B	503	11/12	0.95	0.15	-0.72	37,45,50,50	0
6	GLA	B	525	11/12	0.94	0.16	-0.73	37,42,46,49	0
3	BGC	B	521	11/12	0.91	0.15	-1.15	34,38,44,49	0
2	HG	B	502	1/1	0.97	0.10	-1.17	72,72,72,72	1
2	HG	A	502	1/1	1.00	0.09	-2.42	61,61,61,61	1
2	HG	B	501	1/1	0.99	0.09	-2.57	60,60,60,60	1
3	BGC	A	503	11/12	0.94	0.14	-4.08	33,41,46,48	0
5	XYP	B	523	9/10	0.90	0.21	-	59,66,74,80	0
5	XYP	B	515	9/10	0.89	0.21	-	40,47,49,52	0
4	FUC	A	529	10/11	0.61	0.24	-	70,74,81,82	0
9	RM4	B	528	10/11	0.66	0.28	-	80,83,84,85	0
5	XYP	A	505	9/10	0.81	0.24	-	64,67,76,80	0
8	MAN	A	527	11/12	0.91	0.27	-	63,65,69,71	0
10	7CV	B	520	12/13	0.87	0.27	-	44,47,51,51	0
4	FUC	B	513	10/11	0.96	0.12	-	29,43,48,53	0
4	FUC	A	522	10/11	0.94	0.17	-	36,48,54,58	0
4	FUC	B	531	10/11	0.65	0.26	-	77,82,83,84	0
5	XYP	A	524	9/10	0.90	0.19	-	36,42,52,56	0
9	RM4	A	510	10/11	0.85	0.46	-	81,85,85,86	0
7	XXR	B	517	10/11	0.96	0.12	-	42,54,56,59	0
7	XXR	B	526	10/11	0.91	0.21	-	57,62,70,71	0
4	FUC	A	504	10/11	0.97	0.17	-	39,49,54,62	0
10	7CV	A	511	12/13	0.89	0.34	-	69,77,84,84	0
5	XYP	B	505	9/10	0.75	0.26	-	63,70,84,90	0
7	XXR	A	517	10/11	0.93	0.10	-	45,52,53,59	0
5	XYP	B	514	9/10	0.96	0.15	-	39,43,47,49	0
10	7CV	B	529	12/13	0.72	0.33	-	70,84,90,91	0
8	MAN	B	509	11/12	0.90	0.31	-	63,69,74,75	0
8	MAN	A	509	11/12	0.91	0.23	-	65,68,71,74	0
5	XYP	A	506	9/10	0.89	0.23	-	49,52,55,56	0
8	MAN	B	527	11/12	0.81	0.27	-	69,73,74,77	0
9	RM4	B	519	10/11	0.93	0.17	-	47,50,53,57	0
6	GLA	A	507	11/12	0.92	0.24	-	48,50,53,55	0
4	FUC	A	513	10/11	0.96	0.13	-	34,40,45,50	0

Continued on next page...

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
8	MAN	A	518	11/12	0.89	0.19	-	46,59,60,62	0
7	XXR	A	526	10/11	0.95	0.17	-	52,55,58,61	0
10	7CV	A	520	12/13	0.92	0.20	-	37,45,48,48	0
5	XYP	B	506	9/10	0.94	0.25	-	51,52,54,56	0
5	XYP	A	515	9/10	0.94	0.18	-	41,45,48,52	0
7	XXR	A	508	10/11	0.93	0.18	-	37,50,55,57	0
7	XXR	B	508	10/11	0.96	0.19	-	44,58,62,67	0
5	XYP	A	523	9/10	0.85	0.21	-	58,59,60,61	0
6	GLA	B	507	11/12	0.89	0.25	-	45,49,52,53	0
9	RM4	A	519	10/11	0.94	0.19	-	38,48,50,52	0
4	FUC	B	504	10/11	0.93	0.16	-	43,52,57,64	0
9	RM4	B	510	10/11	0.81	0.36	-	92,97,98,99	0
4	FUC	B	522	10/11	0.94	0.11	-	41,50,54,55	0
8	MAN	B	518	11/12	0.93	0.20	-	46,59,63,63	0
10	7CV	B	511	12/13	0.78	0.36	-	86,90,96,96	0
5	XYP	A	514	9/10	0.96	0.15	-	43,45,46,47	0
5	XYP	B	524	9/10	0.93	0.23	-	43,55,58,60	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.