



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 24, 2023 – 08:17 PM EDT

PDB ID : 5V7S  
Title : Crystal structure of Influenza A virus matrix protein M1 (NLS-88E, pH 6.2)  
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Deposited on : 2017-03-20  
Resolution : 2.50 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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

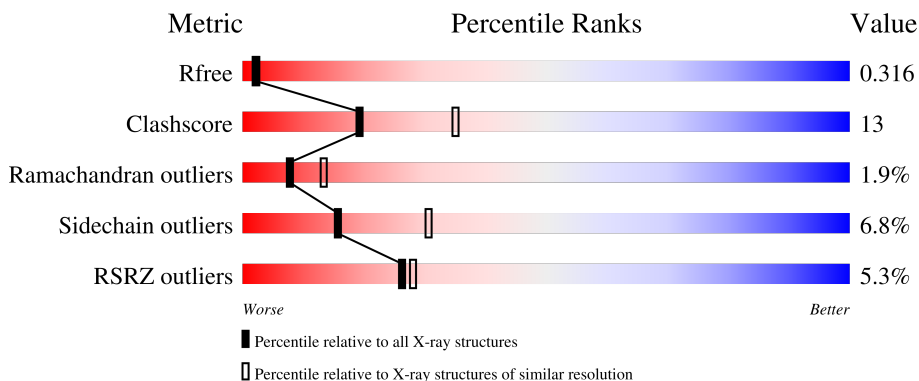
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	171	
1	B	171	
1	C	171	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 3658 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 Matrix protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	164	1267	801	220	238	8	0	0	0
1	B	154	1180	751	197	225	7	0	0	0
1	C	155	1188	755	201	225	7	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

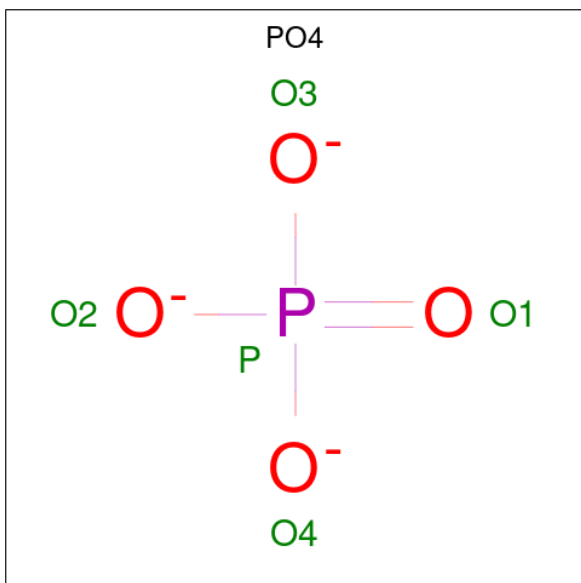
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	expression tag	UNP P05777
A	-4	HIS	-	expression tag	UNP P05777
A	-3	HIS	-	expression tag	UNP P05777
A	-2	HIS	-	expression tag	UNP P05777
A	-1	HIS	-	expression tag	UNP P05777
A	0	HIS	-	expression tag	UNP P05777
A	1	HIS	-	expression tag	UNP P05777
A	88	GLU	GLY	engineered mutation	UNP P05777
A	101	SER	ARG	engineered mutation	UNP P05777
A	105	SER	ARG	engineered mutation	UNP P05777
B	-5	MET	-	expression tag	UNP P05777
B	-4	HIS	-	expression tag	UNP P05777
B	-3	HIS	-	expression tag	UNP P05777
B	-2	HIS	-	expression tag	UNP P05777
B	-1	HIS	-	expression tag	UNP P05777
B	0	HIS	-	expression tag	UNP P05777
B	1	HIS	-	expression tag	UNP P05777
B	88	GLU	GLY	engineered mutation	UNP P05777
B	101	SER	ARG	engineered mutation	UNP P05777
B	105	SER	ARG	engineered mutation	UNP P05777
C	-5	MET	-	expression tag	UNP P05777
C	-4	HIS	-	expression tag	UNP P05777
C	-3	HIS	-	expression tag	UNP P05777

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP P05777
C	-1	HIS	-	expression tag	UNP P05777
C	0	HIS	-	expression tag	UNP P05777
C	1	HIS	-	expression tag	UNP P05777
C	88	GLU	GLY	engineered mutation	UNP P05777
C	101	SER	ARG	engineered mutation	UNP P05777
C	105	SER	ARG	engineered mutation	UNP P05777

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

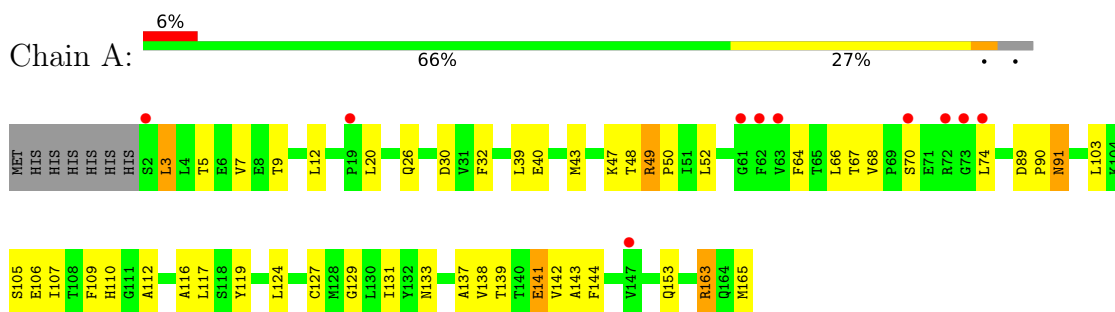
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
3	A	4	Total	O	0	0
			4	4		
3	B	3	Total	O	0	0
			3	3		
3	C	6	Total	O	0	0
			6	6		

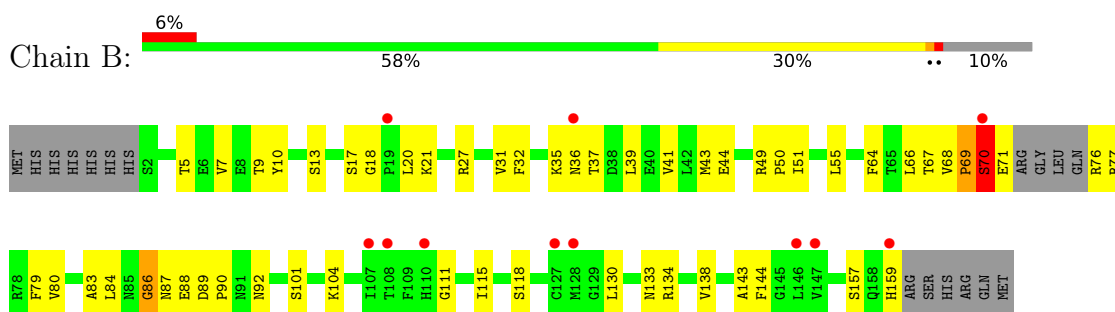
### 3 Residue-property plots [i](#)

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.

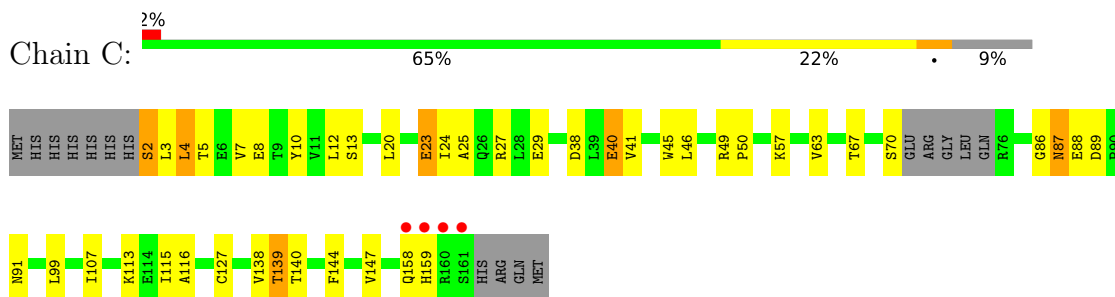
- Molecule 1: Matrix protein 1



- Molecule 1: Matrix protein 1



- Molecule 1: Matrix protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.61Å 133.26Å 39.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.23 – 2.50 26.23 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.6 (26.23-2.50) 90.6 (26.23-2.50)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.07 (at 2.50Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.220 , 0.311 0.231 , 0.316	Depositor DCC
$R_{free}$ test set	767 reflections (5.21%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.4	Xtrriage
Anisotropy	0.099	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3658	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.95% 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

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	0.40	0/1287	0.56	0/1739
1	B	0.43	0/1198	0.59	0/1621
1	C	0.45	0/1206	0.59	0/1631
All	All	0.43	0/3691	0.58	0/4991

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	1267	0	1294	37	0
1	B	1180	0	1203	36	0
1	C	1188	0	1215	27	0
2	A	5	0	0	0	0
2	B	5	0	0	1	0
3	A	4	0	0	1	0
3	B	3	0	0	1	0
3	C	6	0	0	0	0
All	All	3658	0	3712	95	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 13.

All (95) 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:119:TYR:O	1:A:153:GLN:NE2	2.21	0.72
1:B:49:ARG:HD3	1:B:50:PRO:HD2	1.76	0.68
1:B:49:ARG:HH11	1:B:49:ARG:HA	1.60	0.67
1:A:91:ASN:N	1:A:91:ASN:OD1	2.27	0.66
1:A:32:PHE:HE1	1:A:67:THR:HG21	1.64	0.63
1:A:137:ALA:O	1:B:87:ASN:ND2	2.32	0.63
1:B:35:LYS:HG2	1:B:36:ASN:H	1.64	0.62
1:C:2:SER:OG	1:C:2:SER:O	2.13	0.62
1:B:20:LEU:HD11	1:B:49:ARG:HD2	1.81	0.62
1:B:18:GLY:HA3	1:B:21:LYS:H	1.66	0.60
1:B:77:ARG:HG2	1:B:79:PHE:CE2	2.37	0.60
1:B:10:TYR:O	1:B:13:SER:HB3	2.02	0.60
1:C:49:ARG:HA	1:C:49:ARG:HH11	1.67	0.59
1:C:20:LEU:O	1:C:24:ILE:HG22	2.01	0.59
1:B:5:THR:O	1:B:9:THR:HG23	2.05	0.57
1:A:138:VAL:HA	1:B:87:ASN:HD21	1.70	0.56
1:B:92:ASN:ND2	3:B:302:HOH:O	2.39	0.56
1:B:39:LEU:HD13	1:B:68:VAL:HB	1.88	0.56
1:B:31:VAL:HG12	1:B:37:THR:HG21	1.88	0.56
1:A:112:ALA:HB2	1:A:142:VAL:HG22	1.88	0.55
1:C:2:SER:O	1:C:4:LEU:N	2.40	0.55
1:C:99:LEU:HD23	1:C:127:CYS:SG	2.48	0.54
1:C:144:PHE:HA	1:C:147:VAL:HG13	1.90	0.54
1:C:38:ASP:HB3	1:C:41:VAL:HB	1.89	0.53
1:A:40:GLU:HB2	1:A:74:LEU:HD22	1.91	0.53
1:B:76:ARG:NH2	1:B:133:ASN:O	2.41	0.53
1:C:63:VAL:O	1:C:67:THR:HG23	2.09	0.52
1:B:69:PRO:O	1:B:71:GLU:HG3	2.09	0.52
1:B:55:LEU:HA	1:B:83:ALA:HB1	1.91	0.52
1:A:5:THR:O	1:A:9:THR:HG23	2.11	0.51
1:B:49:ARG:HH11	1:B:49:ARG:CA	2.22	0.51
1:A:129:GLY:O	1:A:133:ASN:HB2	2.11	0.50
1:B:27:ARG:O	1:B:31:VAL:HG13	2.12	0.50
1:C:107:ILE:HD13	1:C:139:THR:HG21	1.93	0.49
1:A:47:LYS:HE2	3:A:302:HOH:O	2.12	0.49
1:C:87:ASN:HB3	1:C:89:ASP:H	1.76	0.49
1:A:70:SER:O	1:A:70:SER:OG	2.28	0.49
1:A:103:LEU:O	1:A:106:GLU:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:VAL:HG22	1:A:139:THR:H	1.79	0.48
1:A:138:VAL:HG21	1:A:142:VAL:HG11	1.94	0.48
1:C:12:LEU:HD11	1:C:25:ALA:HB2	1.96	0.48
1:B:49:ARG:HD3	1:B:49:ARG:HA	1.70	0.47
1:C:4:LEU:HG	1:C:5:THR:N	2.28	0.47
1:C:67:THR:HG22	1:C:144:PHE:HZ	1.78	0.47
1:B:43:MET:HE3	1:B:64:PHE:HB2	1.97	0.47
1:C:25:ALA:O	1:C:29:GLU:HG2	2.15	0.46
1:C:10:TYR:CZ	1:C:113:LYS:HE3	2.50	0.46
1:B:134:ARG:NH1	1:C:88:GLU:OE2	2.48	0.46
1:C:8:GLU:O	1:C:12:LEU:HB2	2.14	0.46
1:A:3:LEU:HD21	1:A:144:PHE:CD2	2.51	0.45
1:A:20:LEU:HD11	1:A:49:ARG:HG3	1.99	0.45
1:B:89:ASP:HA	1:B:90:PRO:HD3	1.71	0.45
1:C:24:ILE:HG23	1:C:25:ALA:H	1.81	0.45
1:C:107:ILE:HA	1:C:139:THR:HG21	1.99	0.45
1:B:101:SER:HA	1:B:104:LYS:HD2	1.98	0.45
1:C:49:ARG:HD3	1:C:50:PRO:HD2	1.99	0.45
1:C:38:ASP:OD2	1:C:40:GLU:HG2	2.17	0.45
1:A:50:PRO:O	1:A:52:LEU:N	2.46	0.44
1:B:111:GLY:O	1:B:115:ILE:HG13	2.17	0.44
1:A:67:THR:HG22	1:A:144:PHE:HZ	1.82	0.44
1:C:45:TRP:O	1:C:49:ARG:HG2	2.17	0.44
1:B:130:LEU:HD23	1:B:130:LEU:HA	1.86	0.43
1:A:116:ALA:HB1	1:A:124:LEU:HD22	1.99	0.43
1:B:35:LYS:HG2	1:B:36:ASN:N	2.31	0.43
1:A:26:GLN:NE2	1:A:30:ASP:OD2	2.52	0.43
1:B:79:PHE:N	2:B:201:PO4:O4	2.47	0.42
1:C:115:ILE:HG13	1:C:116:ALA:N	2.34	0.42
1:B:32:PHE:HE1	1:B:67:THR:HG21	1.85	0.42
1:C:10:TYR:O	1:C:13:SER:HB3	2.20	0.42
1:C:12:LEU:HD23	1:C:12:LEU:HA	1.80	0.42
1:A:68:VAL:HG22	1:A:70:SER:H	1.84	0.42
1:A:127:CYS:O	1:A:131:ILE:HG12	2.20	0.42
1:A:12:LEU:HD23	1:A:12:LEU:HA	1.85	0.42
1:A:139:THR:HG23	1:A:142:VAL:H	1.85	0.42
1:C:23:GLU:H	1:C:23:GLU:HG2	1.53	0.42
1:A:39:LEU:HB3	1:A:74:LEU:HD13	2.01	0.42
1:B:69:PRO:HB2	1:B:70:SER:H	1.67	0.42
1:A:110:HIS:CE1	1:B:159:HIS:H	2.37	0.41
1:A:117:LEU:HD23	1:A:117:LEU:HA	1.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:MET:HE1	1:A:64:PHE:HB2	2.01	0.41
1:A:20:LEU:HD11	1:A:49:ARG:HD2	2.01	0.41
1:B:80:VAL:O	1:B:84:LEU:HG	2.20	0.41
1:B:86:GLY:C	1:B:88:GLU:H	2.24	0.41
1:B:7:VAL:HG21	1:B:144:PHE:CG	2.55	0.41
1:B:66:LEU:HD21	1:B:143:ALA:CB	2.50	0.41
1:A:138:VAL:HA	1:B:87:ASN:ND2	2.35	0.41
1:A:66:LEU:HD11	1:A:143:ALA:CB	2.51	0.41
1:A:49:ARG:HA	1:A:49:ARG:HD3	1.54	0.41
1:A:39:LEU:HB3	1:A:74:LEU:CD1	2.51	0.40
1:A:89:ASP:HA	1:A:90:PRO:HD2	1.72	0.40
1:A:109:PHE:HA	1:A:141:GLU:O	2.21	0.40
1:B:50:PRO:HG2	1:B:51:ILE:HD12	2.03	0.40
1:C:46:LEU:HD21	1:C:57:LYS:O	2.20	0.40
1:A:7:VAL:HG21	1:A:144:PHE:CD1	2.56	0.40
1:A:66:LEU:HD11	1:A:143:ALA:HB1	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	162/171 (95%)	147 (91%)	13 (8%)	2 (1%)	13	24
1	B	150/171 (88%)	141 (94%)	6 (4%)	3 (2%)	7	12
1	C	151/171 (88%)	139 (92%)	8 (5%)	4 (3%)	5	8
All	All	463/513 (90%)	427 (92%)	27 (6%)	9 (2%)	8	13

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	69	PRO

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Mol	Chain	Res	Type
1	B	70	SER
1	C	87	ASN
1	B	86	GLY
1	C	3	LEU
1	C	86	GLY
1	C	158	GLN
1	A	163	ARG
1	A	107	ILE

### 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	139/146 (95%)	131 (94%)	8 (6%)	20	38
1	B	130/146 (89%)	123 (95%)	7 (5%)	22	42
1	C	131/146 (90%)	119 (91%)	12 (9%)	9	18
All	All	400/438 (91%)	373 (93%)	27 (7%)	16	30

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	48	THR
1	A	49	ARG
1	A	91	ASN
1	A	105	SER
1	A	141	GLU
1	A	163	ARG
1	A	165	MET
1	B	17	SER
1	B	41	VAL
1	B	44	GLU
1	B	70	SER
1	B	118	SER
1	B	138	VAL

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Mol	Chain	Res	Type
1	B	157	SER
1	C	2	SER
1	C	4	LEU
1	C	7	VAL
1	C	23	GLU
1	C	27	ARG
1	C	40	GLU
1	C	70	SER
1	C	91	ASN
1	C	138	VAL
1	C	139	THR
1	C	140	THR
1	C	159	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	87	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](#)

2 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	A	201	-	4,4,4	0.96	0	6,6,6	0.44	0
2	PO4	B	201	-	4,4,4	0.64	0	6,6,6	0.55	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
2	B	201	PO4	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

### 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	164/171 (95%)	0.23	10 (6%) 21 22	41, 58, 83, 97	0
1	B	154/171 (90%)	0.26	11 (7%) 16 16	41, 56, 72, 99	0
1	C	155/171 (90%)	-0.01	4 (2%) 56 59	37, 50, 71, 103	0
All	All	473/513 (92%)	0.16	25 (5%) 26 28	37, 55, 75, 103	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	SER	5.1
1	A	72	ARG	4.2
1	C	161	SER	3.8
1	A	74	LEU	3.4
1	A	73	GLY	3.3
1	B	107	ILE	3.2
1	B	36	ASN	3.0
1	A	19	PRO	3.0
1	C	160	ARG	3.0
1	B	127	CYS	2.9
1	B	147	VAL	2.8
1	C	159	HIS	2.7
1	B	70	SER	2.7
1	C	158	GLN	2.7
1	A	2	SER	2.5
1	B	110	HIS	2.5
1	A	147	VAL	2.4
1	B	128	MET	2.4
1	B	159	HIS	2.4
1	B	19	PRO	2.4
1	A	62	PHE	2.4
1	A	61	GLY	2.3
1	A	63	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	108	THR	2.2
1	B	146	LEU	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 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
2	PO4	A	201	5/5	0.90	0.52	70,80,88,91	0
2	PO4	B	201	5/5	0.92	0.40	81,81,92,96	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.