



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2017 – 08:55 pm GMT

PDB ID : 5MZ4  
Title : Crystal Structure of full-length CSFV NS3/4A  
Authors : Tortorici, M.A.; Rey, F.A.  
Deposited on : 2017-01-30  
Resolution : 3.05 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.

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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  
Xtriage (Phenix) : 1.9-1692  
EDS : recalc29047  
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) : recalc29047

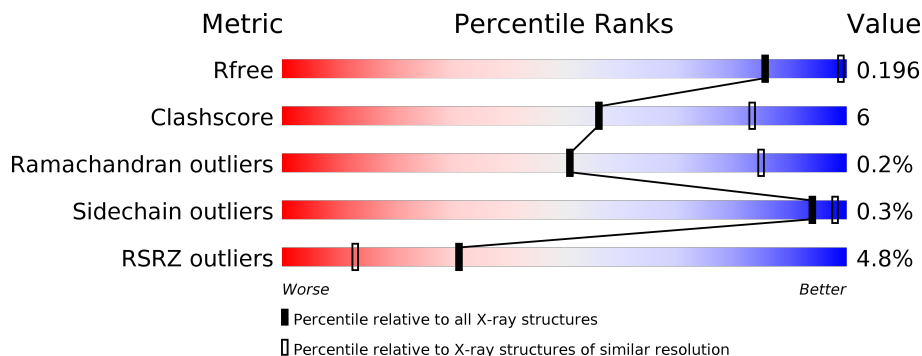
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.05 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	2176 (3.08-3.00)
Clashscore	112137	2542 (3.08-3.00)
Ramachandran outliers	110173	2458 (3.08-3.00)
Sidechain outliers	110143	2461 (3.08-3.00)
RSRZ outliers	101464	2202 (3.08-3.00)

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 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	752	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>14%</div> <div>6%</div> </div> </div>
1	B	752	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>12%</div> <div>10%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10798 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 Genome polyprotein,Genome polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	704	Total	C	N	O	S	0	0	0
			5480	3453	947	1045	35			
1	B	679	Total	C	N	O	S	0	0	0
			5308	3349	915	1008	36			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P19712
A	-18	ALA	-	expression tag	UNP P19712
A	-17	SER	-	expression tag	UNP P19712
A	-16	HIS	-	expression tag	UNP P19712
A	-15	HIS	-	expression tag	UNP P19712
A	-14	HIS	-	expression tag	UNP P19712
A	-13	HIS	-	expression tag	UNP P19712
A	-12	HIS	-	expression tag	UNP P19712
A	-11	HIS	-	expression tag	UNP P19712
A	-10	HIS	-	expression tag	UNP P19712
A	-9	HIS	-	expression tag	UNP P19712
A	-8	HIS	-	expression tag	UNP P19712
A	-7	HIS	-	expression tag	UNP P19712
A	-6	GLU	-	expression tag	UNP P19712
A	-5	ASN	-	expression tag	UNP P19712
A	-4	LEU	-	expression tag	UNP P19712
A	-3	TYR	-	expression tag	UNP P19712
A	-2	PHE	-	expression tag	UNP P19712
A	-1	GLN	-	expression tag	UNP P19712
A	0	GLY	-	expression tag	UNP P19712
A	38	GLY	-	linker	UNP P19712
A	39	SER	-	linker	UNP P19712
A	40	GLY	-	linker	UNP P19712
A	41	SER	-	linker	UNP P19712
A	204	ALA	SER	engineered mutation	UNP P19712

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP P19712
B	-18	ALA	-	expression tag	UNP P19712
B	-17	SER	-	expression tag	UNP P19712
B	-16	HIS	-	expression tag	UNP P19712
B	-15	HIS	-	expression tag	UNP P19712
B	-14	HIS	-	expression tag	UNP P19712
B	-13	HIS	-	expression tag	UNP P19712
B	-12	HIS	-	expression tag	UNP P19712
B	-11	HIS	-	expression tag	UNP P19712
B	-10	HIS	-	expression tag	UNP P19712
B	-9	HIS	-	expression tag	UNP P19712
B	-8	HIS	-	expression tag	UNP P19712
B	-7	HIS	-	expression tag	UNP P19712
B	-6	GLU	-	expression tag	UNP P19712
B	-5	ASN	-	expression tag	UNP P19712
B	-4	LEU	-	expression tag	UNP P19712
B	-3	TYR	-	expression tag	UNP P19712
B	-2	PHE	-	expression tag	UNP P19712
B	-1	GLN	-	expression tag	UNP P19712
B	0	GLY	-	expression tag	UNP P19712
B	38	GLY	-	linker	UNP P19712
B	39	SER	-	linker	UNP P19712
B	40	GLY	-	linker	UNP P19712
B	41	SER	-	linker	UNP P19712
B	204	ALA	SER	engineered mutation	UNP P19712

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	7	Total O 7 7	0	0
2	B	3	Total O 3 3	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.21Å 168.84Å 98.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.95 – 3.05 38.95 – 3.05	Depositor EDS
% Data completeness (in resolution range)	97.5 (38.95-3.05) 97.6 (38.95-3.05)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 3.06Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, $R_{free}$	0.193 , 0.233 0.195 , 0.196	Depositor DCC
$R_{free}$ test set	2041 reflections (5.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	69.8	Xtriage
Anisotropy	0.653	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 58.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10798	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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

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	0.30	0/5586	0.50	0/7562
1	B	0.29	0/5406	0.49	0/7300
All	All	0.30	0/10992	0.50	0/14862

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	5480	0	5481	74	0
1	B	5308	0	5305	58	0
2	A	7	0	0	0	0
2	B	3	0	0	0	0
All	All	10798	0	10786	125	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 6.

All (125) 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:496:GLU:O	1:A:547:ARG:NH2	2.12	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:ASN:HB2	1:A:544:ARG:HH21	1.47	0.80
1:B:132:SER:HB3	1:B:141:GLU:HB2	1.66	0.76
1:A:471:SER:HA	1:A:495:ILE:HD11	1.69	0.75
1:B:101:HIS:NE2	1:B:216:VAL:HG21	2.04	0.72
1:A:141:GLU:OE2	1:A:219:ARG:NH2	2.23	0.70
1:A:132:SER:HB2	1:A:141:GLU:HB2	1.73	0.69
1:B:221:LYS:HD2	1:B:235:SER:HA	1.74	0.69
1:A:106:SER:HB2	1:A:219:ARG:HH21	1.58	0.69
1:B:100:THR:O	1:B:101:HIS:ND1	2.27	0.67
1:B:-2:PHE:HB3	1:B:721:VAL:HA	1.78	0.65
1:B:74:THR:HG22	1:B:125:ARG:HB3	1.78	0.65
1:B:138:ASP:HB2	1:B:221:LYS:HD3	1.80	0.64
1:A:410:PHE:HB2	1:A:556:TYR:O	1.98	0.64
1:B:720:VAL:O	1:B:721:VAL:HG22	1.98	0.62
1:B:493:ASN:HB2	1:B:543:ARG:HH21	1.66	0.60
1:A:469:TYR:HE2	1:A:495:ILE:HG21	1.66	0.60
1:B:99:TYR:CE1	1:B:106:SER:HB3	2.37	0.60
1:A:449:ARG:NH2	1:A:471:SER:OG	2.36	0.58
1:A:74:THR:HG22	1:A:125:ARG:HB3	1.85	0.58
1:A:507:VAL:HB	1:A:556:TYR:CD1	2.39	0.58
1:A:437:GLU:OE1	1:A:506:VAL:HG21	2.05	0.57
1:A:492:THR:HG22	1:A:494:ALA:H	1.70	0.57
1:A:484:GLN:HG2	1:A:486:PRO:HD2	1.86	0.56
1:A:121:ASP:OD1	1:A:122:THR:N	2.36	0.56
1:B:222:VAL:HB	1:B:234:MET:HB2	1.88	0.55
1:B:277:LEU:HB3	1:B:278:PRO:HD3	1.87	0.55
1:A:485:SER:HB2	1:A:486:PRO:HD3	1.89	0.54
1:B:101:HIS:CE1	1:B:216:VAL:HG21	2.44	0.53
1:B:290:ARG:HD2	1:B:330:ALA:O	2.09	0.53
1:A:277:LEU:HB3	1:A:278:PRO:HD3	1.91	0.53
1:B:139:GLU:HG2	1:B:219:ARG:NH2	2.25	0.52
1:B:628:ASN:HB3	1:B:632:ARG:NH1	2.25	0.52
1:A:383:GLU:HG2	1:B:585:ASN:CB	2.40	0.52
1:A:517:ARG:NH2	1:A:534:MET:SD	2.83	0.52
1:B:75:THR:O	1:B:126:THR:HA	2.10	0.51
1:B:547:VAL:HG23	1:B:555:TYR:CZ	2.45	0.51
1:A:250:MET:O	1:A:254:ILE:HG12	2.11	0.51
1:A:655:PHE:CG	1:A:692:GLU:HG2	2.46	0.51
1:B:152:GLU:OE1	1:B:180:GLY:N	2.43	0.50
1:B:338:TYR:CZ	1:B:371:GLN:HG3	2.47	0.50
1:B:580:ILE:HD11	1:B:617:LEU:HD21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:THR:OG1	1:A:371:GLN:HB2	2.12	0.49
1:B:275:THR:OG1	1:B:276:GLU:N	2.44	0.49
1:A:222:VAL:HG22	1:B:730:LEU:HD23	1.93	0.49
1:B:267:LEU:O	1:B:392:ALA:HA	2.13	0.49
1:A:588:THR:H	1:A:615:ASN:HD22	1.61	0.49
1:B:684:TYR:HB2	1:B:702:TRP:CZ3	2.46	0.49
1:A:383:GLU:HG2	1:B:585:ASN:HB3	1.95	0.49
1:A:64:THR:HG22	1:A:69:VAL:O	2.14	0.48
1:A:545:ARG:HG2	1:A:556:TYR:CD2	2.48	0.48
1:A:588:THR:H	1:A:615:ASN:ND2	2.11	0.48
1:A:394:PRO:HD2	1:A:397:THR:HG21	1.95	0.47
1:A:182:GLU:HG3	1:A:232:LYS:HD3	1.96	0.47
1:B:182:GLU:HG2	1:B:234:MET:HA	1.94	0.47
1:B:516:ARG:NH2	1:B:533:MET:SD	2.88	0.47
1:B:175:HIS:HB2	1:B:186:VAL:O	2.14	0.47
1:A:713:GLU:OE1	1:A:716:ARG:HD3	2.16	0.46
1:A:434:PRO:HD2	1:A:557:TYR:CE2	2.50	0.46
1:A:300:ALA:O	1:A:304:VAL:HG23	2.15	0.46
1:A:398:VAL:HG13	1:A:539:GLY:HA2	1.96	0.46
1:A:361:LEU:HD13	1:A:375:MET:HG2	1.97	0.46
1:A:667:TYR:CZ	1:A:677:LYS:HD3	2.51	0.46
1:B:250:MET:O	1:B:254:ILE:HG12	2.15	0.46
1:B:407:ILE:HB	1:B:555:TYR:CE2	2.51	0.46
1:A:537:THR:HA	1:A:564:VAL:O	2.16	0.45
1:A:636:HIS:ND1	1:A:637:PRO:HD2	2.31	0.45
1:B:654:PHE:CG	1:B:691:GLU:HG2	2.51	0.45
1:A:713:GLU:OE2	1:A:716:ARG:NH1	2.47	0.45
1:A:586:ASN:OD1	1:A:615:ASN:ND2	2.46	0.45
1:B:77:ARG:HB2	1:B:126:THR:HB	1.98	0.45
1:A:713:GLU:HA	1:A:716:ARG:HG2	1.98	0.45
1:B:139:GLU:HG2	1:B:219:ARG:HH21	1.81	0.45
1:A:380:ARG:O	1:B:380:ARG:NH2	2.50	0.45
1:B:425:TYR:HD1	1:B:434:PRO:HA	1.81	0.45
1:B:108:VAL:HG23	1:B:205:GLY:HA2	1.99	0.44
1:A:507:VAL:HB	1:A:556:TYR:HD1	1.80	0.44
1:A:548:VAL:HG11	1:A:554:GLY:HA3	1.97	0.44
1:A:545:ARG:O	1:A:556:TYR:OH	2.25	0.44
1:A:548:VAL:CG1	1:A:554:GLY:HA3	2.48	0.44
1:B:527:VAL:HA	1:B:687:ALA:O	2.18	0.44
1:B:324:MET:SD	1:B:344:MET:HG2	2.58	0.44
1:B:657:ILE:HD11	1:B:694:ALA:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:ASN:HD22	1:B:602:GLU:HG3	1.82	0.43
1:A:433:ILE:HG13	1:A:557:TYR:CD2	2.54	0.43
1:A:717:ALA:O	1:A:721:VAL:HG23	2.18	0.43
1:A:441:ASN:ND2	1:A:503:ASP:O	2.41	0.43
1:A:151:PRO:HG2	1:A:154:ALA:HB2	1.99	0.43
1:A:449:ARG:HB2	1:A:470:TYR:CD1	2.53	0.43
1:A:487:TYR:CE1	1:A:501:LEU:HD12	2.53	0.43
1:B:338:TYR:CE2	1:B:371:GLN:HG3	2.53	0.43
1:B:444:VAL:HB	1:B:490:VAL:HG22	2.00	0.43
1:B:535:VAL:HG22	1:B:539:GLU:HB2	2.01	0.43
1:A:275:THR:OG1	1:A:276:GLU:N	2.52	0.43
1:A:642:LEU:O	1:A:646:SER:HB3	2.19	0.43
1:B:316:ALA:O	1:B:332:GLY:HA3	2.19	0.43
1:A:224:LYS:HA	1:B:727:GLU:O	2.19	0.42
1:A:77:ARG:HH12	1:A:121:ASP:CB	2.32	0.42
1:A:478:LEU:O	1:A:482:THR:HG23	2.19	0.42
1:B:618:ILE:HG22	1:B:618:ILE:O	2.19	0.42
1:A:158:VAL:HG21	1:A:220:VAL:HG21	2.00	0.42
1:A:324:MET:HA	1:A:340:TYR:CE1	2.54	0.42
1:A:42:GLY:N	1:A:43:PRO:HD3	2.34	0.42
1:A:426:LEU:HD23	1:A:433:ILE:O	2.20	0.42
1:B:369:PRO:HA	1:B:575:ALA:HB2	2.02	0.42
1:B:544:ARG:HG3	1:B:555:TYR:CD1	2.56	0.41
1:A:518:ILE:HG22	1:A:529:THR:HG22	2.03	0.41
1:A:685:TYR:HB2	1:A:703:TRP:CZ3	2.55	0.41
1:B:276:GLU:OE1	1:B:279:ARG:NH2	2.50	0.41
1:A:200:LEU:HB3	1:B:731:LEU:HD13	2.02	0.41
1:A:528:VAL:HA	1:A:688:ALA:O	2.20	0.41
1:B:42:GLY:N	1:B:43:PRO:HD3	2.35	0.41
1:A:262:PHE:HA	1:A:387:VAL:O	2.20	0.41
1:B:364:TYR:HA	1:B:371:GLN:NE2	2.35	0.41
1:B:517:ILE:HG22	1:B:528:THR:HG22	2.01	0.41
1:A:383:GLU:HG2	1:B:585:ASN:HB2	2.02	0.41
1:B:648:THR:HG22	1:B:650:VAL:H	1.85	0.41
1:A:364:TYR:CZ	1:A:600:LEU:HD22	2.55	0.41
1:A:364:TYR:CE1	1:A:600:LEU:HD22	2.56	0.41
1:A:501:LEU:HA	1:A:501:LEU:HD13	1.91	0.41
1:A:539:GLY:O	1:A:543:GLN:HG3	2.21	0.41
1:A:391:THR:HG21	1:A:394:PRO:HA	2.03	0.40
1:A:595:ASN:ND2	1:A:603:GLU:HG3	2.36	0.40
1:B:99:TYR:HE2	1:B:219:ARG:HG2	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:628:LYS:HE3	1:A:628:LYS:HB2	1.82	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	694/752 (92%)	676 (97%)	16 (2%)	2 (0%)	44	79
1	B	663/752 (88%)	649 (98%)	13 (2%)	1 (0%)	51	85
All	All	1357/1504 (90%)	1325 (98%)	29 (2%)	3 (0%)	51	85

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	500	THR
1	B	721	VAL
1	A	486	PRO

### 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	599/640 (94%)	597 (100%)	2 (0%)	94	98
1	B	577/640 (90%)	576 (100%)	1 (0%)	94	98

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1176/1280 (92%)	1173 (100%)	3 (0%)	94 98

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	PRO
1	A	547	ARG
1	B	29	PRO

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

Mol	Chain	Res	Type
1	A	615	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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	704/752 (93%)	0.03	19 (2%) 55 26	41, 75, 140, 221	0
1	B	679/752 (90%)	0.28	48 (7%) 17 6	38, 92, 164, 220	0
All	All	1383/1504 (91%)	0.15	67 (4%) 31 13	38, 82, 154, 221	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	555	TYR	7.3
1	B	478	LEU	6.7
1	A	499	VAL	5.0
1	A	556	TYR	4.7
1	A	432	LYS	4.7
1	B	467	GLY	4.7
1	A	433	ILE	4.4
1	B	408	GLU	3.8
1	A	488	VAL	3.8
1	B	18	ARG	3.5
1	B	547	VAL	3.4
1	B	480	VAL	3.2
1	B	71	PRO	3.1
1	B	468	TYR	3.1
1	A	724	LEU	3.1
1	B	101	HIS	3.0
1	A	565	GLY	3.0
1	A	425	TYR	3.0
1	B	192	PRO	3.0
1	A	555	ARG	2.9
1	B	442	MET	2.9
1	B	490	VAL	2.9
1	A	429	ALA	2.9
1	B	193	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	177	GLN	2.8
1	B	39	SER	2.8
1	B	507	VAL	2.7
1	B	38	GLY	2.7
1	B	545	GLY	2.7
1	A	662	GLU	2.7
1	B	310	GLN	2.7
1	B	477	ASN	2.7
1	B	424	GLU	2.7
1	B	474	ASP	2.6
1	B	444	VAL	2.6
1	B	269	THR	2.6
1	B	556	TYR	2.6
1	B	176	LEU	2.5
1	B	40	GLY	2.4
1	B	189	SER	2.4
1	B	45	VAL	2.4
1	B	188	ALA	2.4
1	A	557	TYR	2.3
1	B	61	ASP	2.3
1	B	422	GLY	2.3
1	B	559	GLN	2.3
1	B	479	ARG	2.2
1	B	476	SER	2.2
1	B	138	ASP	2.2
1	A	548	VAL	2.2
1	B	473	GLU	2.2
1	A	428	ILE	2.2
1	A	326	GLU	2.1
1	B	459	LEU	2.1
1	B	72	ARG	2.1
1	A	426	LEU	2.1
1	B	118	LEU	2.1
1	B	546	ARG	2.1
1	B	69	VAL	2.1
1	B	179	THR	2.1
1	B	472	GLY	2.1
1	B	182	GLU	2.1
1	B	441	ASN	2.1
1	A	507	VAL	2.0
1	A	399	THR	2.0
1	B	70	MET	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	415	VAL	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.