



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 11:30 pm BST

PDB ID : 3PUM
Title : Crystal structure of P domain dimer of Norovirus VA207
Authors : Chen, Y.; Tan, M.; Xia, M.; Hao, N.; Zhang, X.C.; Huang, P.; Jiang, X.; Li, X.; Rao, Z.
Deposited on : 2010-12-06
Resolution : 2.25 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

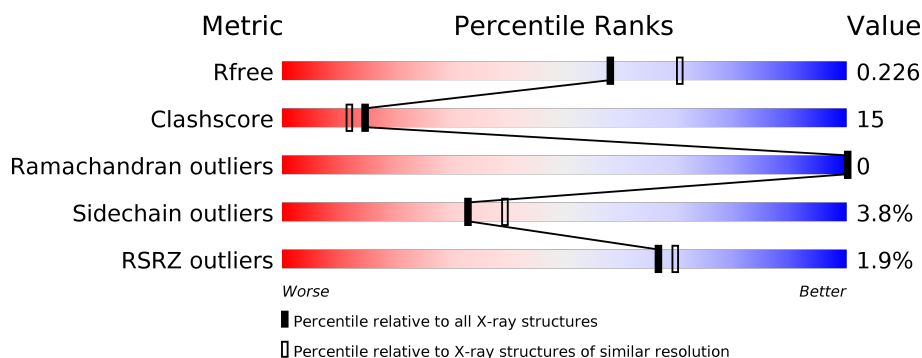
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 ≥ 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	316	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%; text-align: center;">%</div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 22%, green 70%, grey 6%);"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; text-align: center;"> • </div> </div> </div>
1	B	316	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%; text-align: center;">%</div> <div style="position: absolute; top: 0; left: 0; width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 3%, yellow 24%, green 68%, grey 6%);"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; text-align: center;"> • </div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4978 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 Capsid.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2302	1464	392	438	8			
1	B	297	Total	C	N	O	S	0	0	0
			2311	1470	394	439	8			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	289	ASN	THR	ENGINEERED MUTATION	UNP Q91H09
A	374	ASP	ASN	ENGINEERED MUTATION	UNP Q91H09
A	425	GLY	ARG	ENGINEERED MUTATION	UNP Q91H09
A	466	ARG	GLN	ENGINEERED MUTATION	UNP Q91H09
A	482	ALA	VAL	ENGINEERED MUTATION	UNP Q91H09
B	289	ASN	THR	ENGINEERED MUTATION	UNP Q91H09
B	374	ASP	ASN	ENGINEERED MUTATION	UNP Q91H09
B	425	GLY	ARG	ENGINEERED MUTATION	UNP Q91H09
B	466	ARG	GLN	ENGINEERED MUTATION	UNP Q91H09
B	482	ALA	VAL	ENGINEERED MUTATION	UNP Q91H09

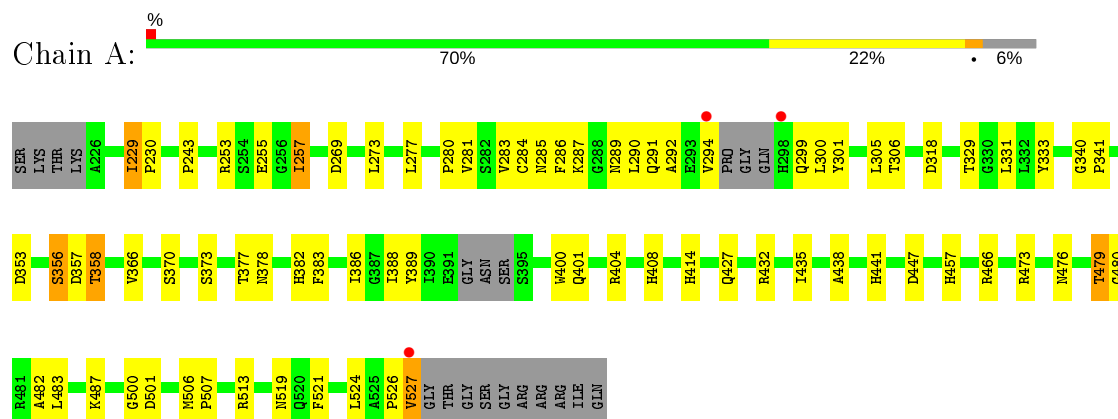
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	193	Total	O	0	0
			193	193		
2	B	172	Total	O	0	0
			172	172		

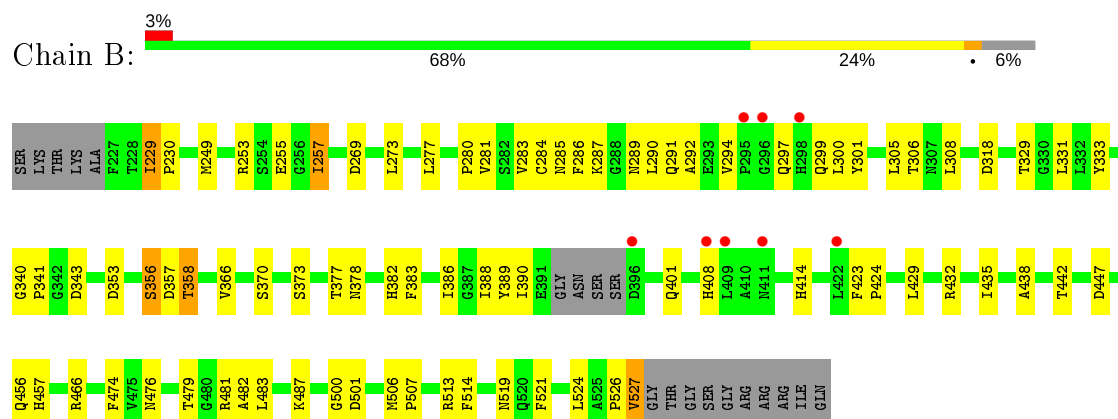
3 Residue-property plots [i](#)

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: Capsid



• Molecule 1: Capsid



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	93.89Å 96.24Å 66.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.10 – 2.25 47.29 – 2.25	Depositor EDS
% Data completeness (in resolution range)	90.1 (23.10-2.25) 93.5 (47.29-2.25)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.15 (at 2.24Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.225 , 0.232 0.222 , 0.226	Depositor DCC
R_{free} test set	1386 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4978	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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](#)

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	0.48	0/2366	0.65	1/3224 (0.0%)
1	B	0.45	0/2377	0.66	1/3241 (0.0%)
All	All	0.46	0/4743	0.65	2/6465 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	THR	N-CA-C	-5.41	96.39	111.00
1	B	306	THR	N-CA-C	-5.27	96.78	111.00

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	2302	0	2200	71	0
1	B	2311	0	2209	70	0
2	A	193	0	0	0	0
2	B	172	0	0	1	0
All	All	4978	0	4409	138	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 15.

All (138) 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:386:ILE:HD11	1:B:386:ILE:HD11	1.21	1.12
1:A:473:ARG:HD3	1:A:482:ALA:HB2	1.37	1.05
1:A:294:VAL:HG21	1:A:300:LEU:HD23	1.49	0.93
1:B:289:ASN:HD21	1:B:378:ASN:HA	1.32	0.92
1:A:289:ASN:HD21	1:A:378:ASN:HA	1.35	0.91
1:B:285:ASN:HD22	1:B:382:HIS:HE1	1.22	0.87
1:B:479:THR:HB	1:B:481:ARG:HG2	1.55	0.85
1:B:294:VAL:HG21	1:B:300:LEU:HD23	1.58	0.83
1:A:229:ILE:HD13	1:A:229:ILE:H	1.49	0.77
1:A:285:ASN:HD22	1:A:382:HIS:HE1	1.35	0.74
1:B:229:ILE:H	1:B:229:ILE:HD13	1.51	0.74
1:A:476:ASN:ND2	1:A:479:THR:CG2	2.51	0.73
1:B:285:ASN:HD22	1:B:382:HIS:CE1	2.04	0.72
1:B:229:ILE:HD11	1:B:513:ARG:HA	1.70	0.72
1:A:229:ILE:HD11	1:A:513:ARG:HA	1.72	0.71
1:A:329:THR:HG23	1:A:401:GLN:O	1.92	0.70
1:A:473:ARG:HD3	1:A:482:ALA:CB	2.21	0.69
1:B:301:TYR:OH	1:B:373:SER:HA	1.93	0.68
1:A:301:TYR:OH	1:A:373:SER:HA	1.94	0.67
1:B:388:ILE:HD13	1:B:435:ILE:CD1	2.25	0.66
1:A:526:PRO:O	1:A:527:VAL:HG12	1.96	0.66
1:A:285:ASN:HD22	1:A:382:HIS:CE1	2.14	0.64
1:A:476:ASN:HB3	1:A:479:THR:HG23	1.79	0.64
1:A:253:ARG:NH1	1:A:500:GLY:HA2	2.12	0.64
1:A:466:ARG:HB3	1:A:466:ARG:HH11	1.62	0.63
1:A:487:LYS:HG3	1:A:524:LEU:HD11	1.81	0.62
1:B:289:ASN:ND2	1:B:378:ASN:HA	2.10	0.62
1:B:526:PRO:O	1:B:527:VAL:HG12	1.99	0.62
1:B:318:ASP:HA	1:B:414:HIS:CE1	2.35	0.61
1:B:466:ARG:HB3	1:B:466:ARG:HH11	1.63	0.61
1:A:329:THR:HG21	1:A:401:GLN:HG2	1.82	0.60
1:B:487:LYS:HG3	1:B:524:LEU:HD11	1.83	0.60
1:A:253:ARG:CZ	1:A:500:GLY:HA2	2.30	0.60
1:B:294:VAL:CG2	1:B:300:LEU:HD23	2.31	0.60
1:A:229:ILE:N	1:A:229:ILE:HD13	2.18	0.59
1:A:289:ASN:ND2	1:A:378:ASN:HA	2.14	0.59
1:B:329:THR:HG23	1:B:401:GLN:O	2.02	0.58
1:B:388:ILE:HD13	1:B:435:ILE:HD12	1.84	0.58
1:B:432:ARG:HB2	1:B:447:ASP:OD1	2.03	0.58
1:A:290:LEU:HD12	1:A:377:THR:HA	1.84	0.58
1:B:229:ILE:N	1:B:229:ILE:HD13	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:VAL:O	1:A:527:VAL:HG13	2.04	0.58
1:A:357:ASP:OD1	1:A:408:HIS:HB3	2.04	0.57
1:A:388:ILE:HD13	1:A:435:ILE:CD1	2.35	0.57
1:B:479:THR:CB	1:B:481:ARG:HG2	2.30	0.57
1:A:294:VAL:CG2	1:A:300:LEU:HD23	2.29	0.57
1:B:527:VAL:O	1:B:527:VAL:HG13	2.04	0.56
1:A:331:LEU:HD23	1:A:389:TYR:HB3	1.88	0.55
1:A:432:ARG:HB2	1:A:447:ASP:OD1	2.06	0.55
1:B:294:VAL:HG21	1:B:297:GLN:OE1	2.07	0.54
1:A:476:ASN:HB2	1:A:483:LEU:HD11	1.89	0.54
1:A:333:TYR:CD2	1:A:438:ALA:HB2	2.42	0.54
1:A:281:VAL:HG12	1:B:281:VAL:HG12	1.88	0.54
1:A:331:LEU:HD23	1:A:389:TYR:CD2	2.43	0.54
1:A:257:ILE:HD13	1:A:257:ILE:C	2.29	0.53
1:B:483:LEU:HD21	1:B:507:PRO:HD2	1.90	0.53
1:B:257:ILE:HD13	1:B:257:ILE:C	2.29	0.53
1:B:290:LEU:HD12	1:B:377:THR:HA	1.90	0.53
1:A:299:GLN:HB3	1:A:370:SER:O	2.09	0.52
1:B:299:GLN:HB3	1:B:370:SER:O	2.09	0.52
1:A:301:TYR:HH	1:A:373:SER:HA	1.74	0.52
1:A:466:ARG:HB3	1:A:466:ARG:NH1	2.25	0.52
1:A:483:LEU:HD21	1:A:507:PRO:HD2	1.92	0.51
1:A:476:ASN:ND2	1:A:479:THR:HG22	2.25	0.51
1:B:331:LEU:HD23	1:B:389:TYR:CD2	2.45	0.51
1:B:331:LEU:HD23	1:B:389:TYR:HB3	1.92	0.51
1:A:476:ASN:CG	1:A:479:THR:HG23	2.31	0.51
1:B:476:ASN:HB2	1:B:483:LEU:HD11	1.93	0.50
1:B:283:VAL:O	1:B:284:CYS:HB2	2.12	0.50
1:B:483:LEU:CD2	1:B:507:PRO:HD2	2.41	0.50
1:B:388:ILE:HD13	1:B:435:ILE:HD11	1.94	0.50
1:A:476:ASN:CB	1:A:479:THR:HG23	2.40	0.50
1:B:329:THR:HG21	1:B:401:GLN:HG2	1.94	0.50
1:A:340:GLY:HA2	1:A:341:PRO:C	2.32	0.49
1:B:340:GLY:HA2	1:B:341:PRO:C	2.32	0.49
1:B:466:ARG:HB3	1:B:466:ARG:NH1	2.26	0.49
1:A:329:THR:CG2	1:A:401:GLN:HG2	2.42	0.49
1:B:229:ILE:HD11	1:B:514:PHE:N	2.27	0.49
1:B:474:PHE:HD2	1:B:483:LEU:HD22	1.78	0.49
1:A:473:ARG:CD	1:A:482:ALA:HB2	2.26	0.48
1:B:353:ASP:HB3	1:B:356:SER:HB3	1.95	0.48
1:A:283:VAL:O	1:A:284:CYS:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:LEU:CD2	1:A:507:PRO:HD2	2.43	0.47
1:A:253:ARG:HG2	1:A:427:GLN:OE1	2.14	0.47
1:B:287:LYS:NZ	1:B:382:HIS:HD2	2.12	0.47
1:B:483:LEU:HD23	1:B:506:MET:SD	2.55	0.47
1:A:388:ILE:HD13	1:A:435:ILE:HD12	1.97	0.46
1:A:388:ILE:HD13	1:A:435:ILE:HD11	1.97	0.46
1:A:253:ARG:HB2	1:A:255:GLU:HG3	1.97	0.46
1:B:390:ILE:HG13	1:B:442:THR:HB	1.98	0.46
1:A:318:ASP:HA	1:A:414:HIS:CE1	2.51	0.46
1:A:291:GLN:O	1:A:292:ALA:HB2	2.16	0.46
1:B:357:ASP:OD1	1:B:408:HIS:HB3	2.16	0.45
1:B:318:ASP:HA	1:B:414:HIS:ND1	2.30	0.45
1:A:476:ASN:ND2	1:A:479:THR:HG23	2.29	0.45
1:A:519:ASN:OD1	1:A:521:PHE:N	2.46	0.45
1:A:353:ASP:HB3	1:A:356:SER:HB3	1.98	0.45
1:A:230:PRO:HD3	1:A:457:HIS:CG	2.51	0.45
1:A:476:ASN:HD22	1:A:479:THR:CG2	2.28	0.44
1:A:483:LEU:HD23	1:A:506:MET:SD	2.58	0.44
1:B:253:ARG:NH1	1:B:500:GLY:HA2	2.33	0.44
1:B:294:VAL:CG2	1:B:297:GLN:HB2	2.47	0.44
1:B:474:PHE:CD2	1:B:483:LEU:HD22	2.51	0.44
1:B:305:LEU:HD11	1:B:366:VAL:HG21	1.98	0.44
1:B:286:PHE:HB2	1:B:383:PHE:HB3	1.99	0.43
1:B:333:TYR:CD2	1:B:438:ALA:HB2	2.53	0.43
1:B:483:LEU:CD1	1:B:483:LEU:N	2.81	0.43
1:B:229:ILE:HD11	1:B:513:ARG:CA	2.42	0.43
1:B:483:LEU:HD12	1:B:483:LEU:N	2.34	0.43
1:B:308:LEU:HD23	1:B:308:LEU:HA	1.89	0.43
1:A:286:PHE:HB2	1:A:383:PHE:HB3	2.00	0.43
1:A:479:THR:OG1	1:A:480:GLY:N	2.50	0.43
1:B:253:ARG:HB2	1:B:255:GLU:HG3	2.01	0.43
1:A:243:PRO:HD3	1:A:283:VAL:HG13	2.01	0.42
1:A:287:LYS:NZ	1:A:382:HIS:HD2	2.16	0.42
1:B:329:THR:CG2	1:B:401:GLN:HG2	2.50	0.42
1:B:356:SER:OG	1:B:358:THR:HG23	2.19	0.42
1:B:285:ASN:ND2	1:B:382:HIS:HE1	2.04	0.42
1:A:483:LEU:N	1:A:483:LEU:CD1	2.83	0.42
1:B:253:ARG:CZ	1:B:500:GLY:HA2	2.50	0.42
1:A:357:ASP:OD1	1:A:404:ARG:NH2	2.51	0.42
1:B:230:PRO:HD3	1:B:457:HIS:CG	2.55	0.42
1:B:291:GLN:O	1:B:292:ALA:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:482:ALA:C	1:B:483:LEU:HD12	2.40	0.42
1:A:257:ILE:HG13	1:A:400:TRP:CD1	2.55	0.41
1:A:356:SER:OG	1:A:358:THR:HG23	2.20	0.41
1:A:280:PRO:HG3	2:B:11:HOH:O	2.20	0.41
1:B:423:PHE:HA	1:B:424:PRO:HD3	1.90	0.41
1:A:305:LEU:HD11	1:A:366:VAL:HG21	2.02	0.41
1:A:483:LEU:N	1:A:483:LEU:HD12	2.36	0.41
1:A:441:HIS:HD2	1:B:343:ASP:OD1	2.04	0.41
1:B:229:ILE:HD11	1:B:514:PHE:H	1.85	0.41
1:B:280:PRO:HD3	1:B:456:GLN:NE2	2.34	0.41
1:B:519:ASN:OD1	1:B:521:PHE:N	2.48	0.40
1:B:479:THR:HB	1:B:481:ARG:H	1.86	0.40
1:A:229:ILE:CD1	1:A:229:ILE:N	2.85	0.40
1:A:331:LEU:HD12	1:A:331:LEU:HA	1.87	0.40
1:B:249:MET:HB3	1:B:429:LEU:HD11	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	290/316 (92%)	279 (96%)	11 (4%)	0	100	100
1	B	293/316 (93%)	285 (97%)	8 (3%)	0	100	100
All	All	583/632 (92%)	564 (97%)	19 (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	251/266 (94%)	241 (96%)	10 (4%)	31	37
1	B	252/266 (95%)	243 (96%)	9 (4%)	35	42
All	All	503/532 (94%)	484 (96%)	19 (4%)	33	39

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	ILE
1	A	257	ILE
1	A	269	ASP
1	A	273	LEU
1	A	277	LEU
1	A	356	SER
1	A	358	THR
1	A	479	THR
1	A	501	ASP
1	A	527	VAL
1	B	229	ILE
1	B	257	ILE
1	B	269	ASP
1	B	273	LEU
1	B	277	LEU
1	B	356	SER
1	B	358	THR
1	B	501	ASP
1	B	527	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	285	ASN
1	A	289	ASN
1	A	291	GLN
1	A	307	ASN
1	A	378	ASN
1	A	382	HIS
1	A	441	HIS

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Mol	Chain	Res	Type
1	A	509	ASN
1	B	248	GLN
1	B	285	ASN
1	B	289	ASN
1	B	291	GLN
1	B	307	ASN
1	B	378	ASN
1	B	382	HIS
1	B	401	GLN
1	B	441	HIS
1	B	509	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 [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, 95th 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 ²)	Q<0.9
1	A	296/316 (93%)	0.09	3 (1%) 82 84	19, 30, 45, 60	0
1	B	297/316 (93%)	0.05	8 (2%) 54 57	18, 31, 47, 68	0
All	All	593/632 (93%)	0.07	11 (1%) 66 69	18, 30, 47, 68	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	295	PRO	4.4
1	B	298	HIS	4.1
1	B	409	LEU	4.1
1	A	527	VAL	3.8
1	A	294	VAL	3.6
1	B	296	GLY	2.9
1	B	422	LEU	2.9
1	B	408	HIS	2.8
1	A	298	HIS	2.3
1	B	396	ASP	2.3
1	B	411	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.