



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2017 – 09:48 AM EDT

PDB ID : 2OBR  
Title : Crystal Structures of P Domain of Norovirus VA387  
Authors : Cao, S.; Lou, Z.; Jiang, X.; Zhang, X.C.; Li, X.; Rao, Z.  
Deposited on : unknown  
Resolution : 2.20 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 : rb-20030345  
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-20030345

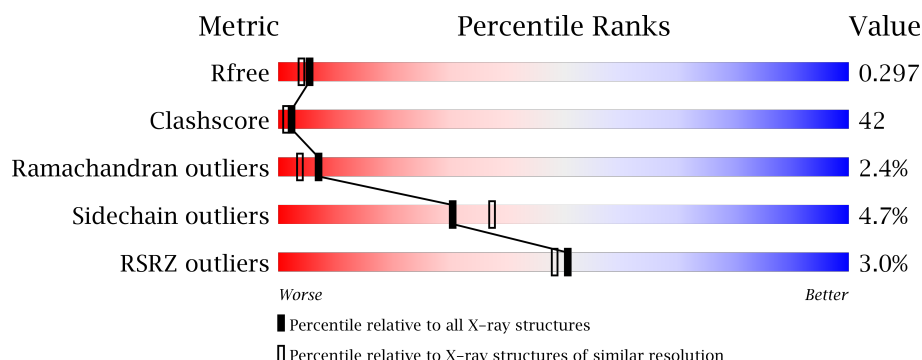
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	327	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2535 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 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2309	1468	394	436	11			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	213	GLY	-	CLONING ARTIFACT	UNP Q913Z3
A	355	SER	THR	ENGINEERED	UNP Q913Z3
A	375	LEU	PHE	ENGINEERED	UNP Q913Z3

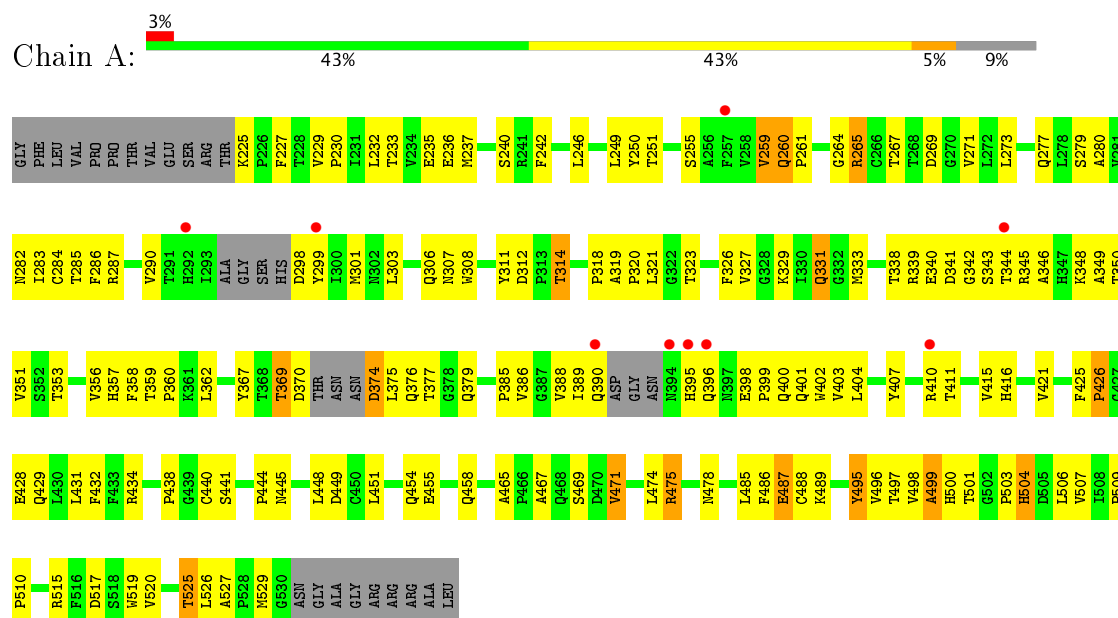
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	226	Total	O	0	0
			226	226		

### 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 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.33Å 97.44Å 118.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 37.68 – 2.21	Depositor EDS
% Data completeness (in resolution range)	95.0 (50.00-2.20) 90.4 (37.68-2.21)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.44 (at 2.22Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.229 , 0.297 0.235 , 0.297	Depositor DCC
$R_{free}$ test set	727 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.1	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 71.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.039 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2535	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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](#)

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	0.61	0/2375	0.87	5/3247 (0.2%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	GLU	N-CA-C	7.32	130.76	111.00
1	A	340	GLU	N-CA-CB	-7.24	97.57	110.60
1	A	339	ARG	CB-CA-C	7.15	124.69	110.40
1	A	374	ASP	N-CA-C	-6.86	92.47	111.00
1	A	342	GLY	N-CA-C	-6.06	97.95	113.10

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	2309	0	2220	190	0
2	A	226	0	0	16	1
All	All	2535	0	2220	190	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 42.

All (190) 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:286:PHE:HB3	1:A:301:MET:HE1	1.34	1.06
1:A:390:GLN:HB3	1:A:398:GLU:HB2	1.39	1.04
1:A:327:VAL:HG23	1:A:403:VAL:HG23	1.39	1.02
1:A:233:THR:HG21	2:A:671:HOH:O	1.78	0.82
1:A:469:SER:HB3	1:A:519:TRP:HB3	1.62	0.82
1:A:327:VAL:CG2	1:A:403:VAL:HG23	2.08	0.81
1:A:265:ARG:HB3	1:A:265:ARG:HH11	1.46	0.80
1:A:312:ASP:OD2	1:A:314:THR:HB	1.80	0.80
1:A:499:ALA:HB2	1:A:526:LEU:HB2	1.65	0.79
1:A:400:GLN:HB2	1:A:403:VAL:CG2	2.13	0.78
1:A:400:GLN:HB2	1:A:403:VAL:HG22	1.67	0.77
1:A:267:THR:CG2	1:A:269:ASP:HB3	2.15	0.76
1:A:331:GLN:NE2	1:A:389:ILE:HG21	2.02	0.74
1:A:390:GLN:HB2	1:A:444:PRO:HB3	1.70	0.73
1:A:389:ILE:HG23	1:A:440:CYS:HB2	1.68	0.73
1:A:267:THR:HG22	1:A:269:ASP:H	1.53	0.73
1:A:280:ALA:HB2	1:A:454:GLN:HG2	1.71	0.72
1:A:290:VAL:O	1:A:377:THR:HG22	1.89	0.71
1:A:399:PRO:HD2	1:A:445:ASN:HB3	1.73	0.71
1:A:395:HIS:O	1:A:396:GLN:HG3	1.90	0.70
1:A:360:PRO:HD2	1:A:407:TYR:O	1.92	0.70
1:A:301:MET:HE3	1:A:303:LEU:HD23	1.73	0.70
1:A:390:GLN:NE2	1:A:398:GLU:HB3	2.07	0.69
1:A:431:LEU:HD23	1:A:496:VAL:HG11	1.74	0.68
1:A:428:GLU:OE2	1:A:489:LYS:HD2	1.93	0.68
1:A:400:GLN:O	1:A:403:VAL:HG22	1.94	0.68
1:A:299:TYR:OH	1:A:375:LEU:HB2	1.94	0.68
1:A:349:ALA:HA	1:A:369:THR:HA	1.75	0.67
1:A:390:GLN:CB	1:A:398:GLU:HB2	2.22	0.67
1:A:338:THR:HA	1:A:379:GLN:NE2	2.09	0.67
1:A:475:ARG:HD3	1:A:517:ASP:OD2	1.95	0.67
1:A:390:GLN:HE22	1:A:399:PRO:HA	1.60	0.66
1:A:267:THR:HG22	1:A:269:ASP:HB3	1.76	0.66
1:A:348:LYS:HB3	1:A:370:ASP:O	1.94	0.66
1:A:415:VAL:HG22	2:A:690:HOH:O	1.94	0.66
1:A:515:ARG:HD3	2:A:761:HOH:O	1.96	0.66
1:A:286:PHE:CB	1:A:301:MET:HE1	2.18	0.65
1:A:344:THR:HG22	1:A:345:ARG:N	2.11	0.65
1:A:489:LYS:HG3	1:A:526:LEU:HD11	1.79	0.65
1:A:338:THR:HG21	1:A:376:GLN:OE1	1.97	0.64
1:A:389:ILE:CG2	1:A:440:CYS:HB2	2.27	0.63
1:A:233:THR:HG22	1:A:236:GLU:CG	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:THR:HG21	1:A:389:ILE:HD11	1.79	0.63
1:A:475:ARG:HB2	1:A:517:ASP:OD2	1.98	0.63
1:A:478:ASN:HD22	1:A:509:PRO:HG3	1.62	0.63
1:A:329:LYS:NZ	2:A:604:HOH:O	2.31	0.63
1:A:301:MET:HE2	1:A:367:TYR:OH	1.99	0.63
1:A:338:THR:OG1	1:A:343:SER:HB3	2.00	0.62
1:A:351:VAL:CG2	1:A:367:TYR:CD2	2.83	0.62
1:A:390:GLN:HE22	1:A:399:PRO:CA	2.12	0.61
1:A:390:GLN:NE2	1:A:399:PRO:HA	2.15	0.61
1:A:504:HIS:HD2	1:A:529:MET:CE	2.12	0.60
1:A:338:THR:CG2	1:A:376:GLN:OE1	2.49	0.60
1:A:233:THR:HG22	1:A:236:GLU:CB	2.32	0.60
1:A:265:ARG:HH11	1:A:265:ARG:CB	2.12	0.60
1:A:415:VAL:HB	1:A:416:HIS:ND1	2.17	0.59
1:A:475:ARG:HE	1:A:487:GLU:HB3	1.66	0.59
1:A:277:GLN:HB3	1:A:321:LEU:HB3	1.83	0.59
1:A:233:THR:HG22	1:A:236:GLU:HB2	1.83	0.59
1:A:265:ARG:NH1	2:A:565:HOH:O	2.35	0.58
1:A:227:PHE:HD1	1:A:465:ALA:HB3	1.69	0.58
1:A:489:LYS:HB2	1:A:497:THR:OG1	2.03	0.58
1:A:469:SER:CB	1:A:519:TRP:HB3	2.34	0.58
1:A:265:ARG:NH1	1:A:265:ARG:HB3	2.17	0.58
1:A:260:GLN:HE21	1:A:260:GLN:HA	1.68	0.57
1:A:400:GLN:HB2	1:A:403:VAL:HG21	1.86	0.57
1:A:260:GLN:CA	1:A:260:GLN:HE21	2.18	0.57
1:A:267:THR:HB	1:A:271:VAL:HB	1.85	0.57
1:A:425:PHE:CZ	1:A:525:THR:HG23	2.39	0.57
1:A:265:ARG:NE	2:A:554:HOH:O	2.37	0.57
1:A:331:GLN:HE21	1:A:389:ILE:HG21	1.70	0.56
1:A:237:MET:HB2	1:A:246:LEU:HD12	1.87	0.56
1:A:504:HIS:HD2	1:A:529:MET:HE3	1.69	0.56
1:A:242:PHE:HB2	1:A:448:LEU:CD2	2.36	0.56
1:A:283:ILE:O	1:A:284:CYS:HB2	2.06	0.56
1:A:467:ALA:C	1:A:469:SER:H	2.09	0.55
1:A:478:ASN:ND2	1:A:509:PRO:HG3	2.21	0.55
1:A:299:TYR:CZ	1:A:375:LEU:HB2	2.41	0.55
1:A:280:ALA:HB1	2:A:622:HOH:O	2.05	0.55
1:A:467:ALA:HA	1:A:519:TRP:CH2	2.40	0.55
1:A:434:ARG:HB2	1:A:449:ASP:OD1	2.07	0.54
1:A:290:VAL:HG22	1:A:299:TYR:HD2	1.71	0.54
1:A:259:VAL:CG1	1:A:421:VAL:HB	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ALA:HB1	1:A:320:PRO:HD2	1.89	0.53
1:A:390:GLN:HE22	1:A:399:PRO:CB	2.21	0.53
1:A:235:GLU:HG3	2:A:640:HOH:O	2.08	0.53
1:A:327:VAL:HG23	1:A:403:VAL:O	2.07	0.53
1:A:403:VAL:O	1:A:403:VAL:HG23	2.09	0.53
1:A:265:ARG:NH1	1:A:273:LEU:HD12	2.24	0.52
1:A:421:VAL:HG13	1:A:495:TYR:CD1	2.44	0.52
1:A:267:THR:HG22	1:A:269:ASP:N	2.22	0.52
1:A:285:THR:HG22	1:A:385:PRO:HD2	1.91	0.52
1:A:471:VAL:HG13	1:A:520:VAL:O	2.10	0.52
1:A:259:VAL:HG12	1:A:421:VAL:HB	1.91	0.51
1:A:311:TYR:CE1	1:A:320:PRO:HG3	2.45	0.51
1:A:351:VAL:HG22	1:A:367:TYR:CD2	2.46	0.51
1:A:390:GLN:NE2	1:A:399:PRO:CA	2.73	0.51
1:A:390:GLN:OE1	1:A:399:PRO:HD3	2.10	0.51
1:A:432:PHE:HD2	1:A:449:ASP:HB3	1.75	0.51
1:A:233:THR:HG22	1:A:236:GLU:HG3	1.92	0.50
1:A:389:ILE:O	1:A:389:ILE:HD12	2.12	0.50
1:A:390:GLN:HE22	1:A:399:PRO:HB3	1.76	0.50
1:A:350:THR:HG22	1:A:351:VAL:N	2.27	0.50
1:A:489:LYS:O	1:A:496:VAL:HA	2.10	0.50
1:A:225:LYS:NZ	1:A:225:LYS:HB2	2.27	0.49
1:A:341:ASP:HB2	1:A:343:SER:HB2	1.94	0.49
1:A:504:HIS:CD2	1:A:529:MET:HE3	2.47	0.49
1:A:331:GLN:OE1	1:A:348:LYS:HD3	2.13	0.49
1:A:301:MET:HE3	1:A:303:LEU:CD2	2.42	0.49
1:A:432:PHE:HB3	1:A:449:ASP:HB3	1.95	0.49
1:A:301:MET:CE	1:A:367:TYR:OH	2.61	0.48
1:A:429:GLN:HB2	1:A:500:HIS:O	2.13	0.48
1:A:411:THR:HG22	2:A:687:HOH:O	2.14	0.48
1:A:425:PHE:HZ	1:A:525:THR:HG23	1.78	0.48
1:A:301:MET:CE	1:A:303:LEU:HD23	2.39	0.48
1:A:267:THR:HG21	1:A:269:ASP:HB3	1.94	0.48
1:A:432:PHE:CD2	1:A:449:ASP:HB3	2.48	0.48
1:A:282:ASN:HB3	1:A:306:GLN:HE22	1.79	0.48
1:A:386:VAL:O	1:A:438:PRO:HG2	2.13	0.48
1:A:344:THR:CG2	1:A:345:ARG:N	2.76	0.47
1:A:510:PRO:HG3	2:A:651:HOH:O	2.13	0.47
1:A:307:ASN:O	1:A:308:TRP:HB2	2.14	0.47
1:A:265:ARG:HH12	1:A:273:LEU:HD12	1.79	0.47
1:A:487:GLU:OE1	1:A:527:ALA:N	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:PRO:O	1:A:504:HIS:HB2	2.15	0.46
1:A:249:LEU:O	1:A:506:LEU:HB2	2.16	0.46
1:A:411:THR:HG21	2:A:665:HOH:O	2.16	0.46
1:A:265:ARG:HA	1:A:265:ARG:HD2	1.52	0.46
1:A:454:GLN:O	1:A:458:GLN:HG3	2.16	0.46
1:A:425:PHE:CG	1:A:426:PRO:HD2	2.50	0.46
1:A:475:ARG:HE	1:A:487:GLU:CG	2.28	0.46
1:A:515:ARG:NH1	1:A:517:ASP:OD1	2.49	0.46
1:A:240:SER:OG	1:A:264:GLY:HA3	2.15	0.46
1:A:298:ASP:HB3	1:A:369:THR:HG22	1.98	0.45
1:A:402:TRP:CZ2	1:A:432:PHE:HE2	2.35	0.45
1:A:232:LEU:HD13	1:A:455:GLU:HG3	1.99	0.45
1:A:376:GLN:HB3	2:A:594:HOH:O	2.15	0.45
1:A:390:GLN:HB2	1:A:444:PRO:CB	2.43	0.45
1:A:356:VAL:HG23	1:A:357:HIS:CD2	2.52	0.45
1:A:485:LEU:HG	1:A:486:PHE:HD1	1.81	0.45
1:A:279:SER:HB2	2:A:558:HOH:O	2.16	0.45
1:A:348:LYS:HB2	1:A:348:LYS:HE3	1.81	0.45
1:A:259:VAL:HA	1:A:402:TRP:CE3	2.52	0.45
1:A:395:HIS:O	1:A:396:GLN:CG	2.60	0.45
1:A:415:VAL:HB	1:A:416:HIS:CE1	2.51	0.45
1:A:280:ALA:CB	1:A:454:GLN:HG2	2.44	0.45
1:A:260:GLN:N	1:A:261:PRO:CD	2.80	0.44
1:A:267:THR:HG22	1:A:269:ASP:CB	2.43	0.44
1:A:374:ASP:O	1:A:375:LEU:HG	2.18	0.44
1:A:475:ARG:HE	1:A:487:GLU:CB	2.31	0.44
1:A:475:ARG:NE	1:A:487:GLU:HB3	2.32	0.44
1:A:475:ARG:HE	1:A:487:GLU:HG2	1.83	0.44
1:A:232:LEU:CD1	1:A:455:GLU:HG3	2.48	0.43
1:A:333:MET:CE	1:A:346:ALA:HB1	2.47	0.43
1:A:499:ALA:CB	1:A:526:LEU:HB2	2.43	0.43
1:A:345:ARG:HG3	2:A:747:HOH:O	2.18	0.43
1:A:287:ARG:HB3	1:A:308:TRP:CZ3	2.54	0.43
1:A:399:PRO:HD2	1:A:445:ASN:O	2.19	0.43
1:A:428:GLU:OE1	1:A:489:LYS:NZ	2.51	0.43
1:A:250:TYR:CD1	1:A:251:THR:N	2.86	0.43
1:A:260:GLN:NE2	1:A:260:GLN:HA	2.33	0.43
1:A:358:PHE:C	1:A:360:PRO:HD3	2.39	0.43
1:A:250:TYR:HE1	1:A:504:HIS:H	1.67	0.42
1:A:421:VAL:HG21	1:A:451:LEU:CD1	2.49	0.42
1:A:428:GLU:OE1	1:A:489:LYS:CE	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:388:VAL:O	1:A:388:VAL:HG23	2.19	0.42
1:A:389:ILE:HG22	1:A:440:CYS:C	2.40	0.42
1:A:301:MET:CE	1:A:303:LEU:CD2	2.97	0.42
1:A:360:PRO:HD2	1:A:407:TYR:HA	2.01	0.42
1:A:359:THR:HB	1:A:362:LEU:HD12	2.00	0.42
1:A:474:LEU:HB2	1:A:488:CYS:SG	2.60	0.42
1:A:504:HIS:HD2	1:A:529:MET:HE1	1.82	0.42
1:A:267:THR:CG2	1:A:269:ASP:H	2.27	0.42
1:A:318:PRO:HG3	1:A:416:HIS:O	2.19	0.41
1:A:509:PRO:HA	1:A:510:PRO:HD3	1.86	0.41
1:A:323:THR:HA	2:A:634:HOH:O	2.20	0.41
1:A:390:GLN:OE1	1:A:399:PRO:CD	2.68	0.41
1:A:421:VAL:HG21	1:A:451:LEU:HD13	2.02	0.41
1:A:299:TYR:CE2	1:A:375:LEU:HB2	2.55	0.41
1:A:353:THR:HA	1:A:358:PHE:CG	2.55	0.41
1:A:233:THR:CG2	1:A:236:GLU:HG3	2.51	0.41
1:A:467:ALA:C	1:A:469:SER:N	2.74	0.41
1:A:326:PHE:HB2	1:A:401:GLN:HA	2.03	0.41
1:A:501:THR:HG23	2:A:656:HOH:O	2.20	0.41
1:A:488:CYS:HB3	1:A:498:VAL:HG12	2.02	0.41
1:A:259:VAL:O	1:A:259:VAL:HG12	2.20	0.40
1:A:319:ALA:HB1	1:A:320:PRO:CD	2.51	0.40
1:A:319:ALA:O	1:A:320:PRO:C	2.58	0.40
1:A:229:VAL:HG22	1:A:230:PRO:HD2	2.03	0.40
1:A:290:VAL:CG2	1:A:299:TYR:HD2	2.34	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)
2:A:613:HOH:O	2:A:613:HOH:O[3_555]	1.86	0.34

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	288/327 (88%)	249 (86%)	32 (11%)	7 (2%)	7	4

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	255	SER
1	A	441	SER
1	A	499	ALA
1	A	504	HIS
1	A	369	THR
1	A	259	VAL
1	A	426	PRO

### 5.3.2 Protein sidechains ⓘ

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	258/281 (92%)	246 (95%)	12 (5%)	30	37

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	260	GLN
1	A	265	ARG
1	A	314	THR
1	A	331	GLN
1	A	404	LEU
1	A	410	ARG
1	A	471	VAL
1	A	475	ARG
1	A	487	GLU
1	A	495	TYR
1	A	507	VAL
1	A	525	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	GLN
1	A	282	ASN
1	A	292	HIS
1	A	306	GLN
1	A	331	GLN
1	A	336	GLN
1	A	366	GLN
1	A	379	GLN
1	A	447	ASN
1	A	500	HIS
1	A	504	HIS
1	A	511	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, 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	296/327 (90%)	0.08	9 (3%)	51 48	16, 40, 64, 92	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	HIS	3.2
1	A	396	GLN	3.1
1	A	299	TYR	2.8
1	A	394	ASN	2.8
1	A	257	PHE	2.6
1	A	390	GLN	2.4
1	A	395	HIS	2.3
1	A	344	THR	2.2
1	A	410	ARG	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.