



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 12, 2017 – 06:05 AM EST

PDB ID : 4NWV  
Title : Crystal structure of Orsay virus-like particle  
Authors : Tao, Y.J.; Guo, Y.R.  
Deposited on : unknown  
Resolution : 3.25 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](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

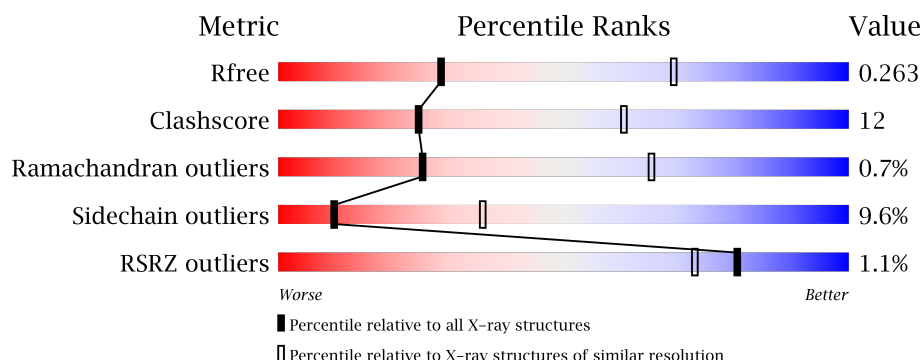
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.25 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))
$R_{free}$	100719	1852 (3.32-3.20)
Clashscore	112137	2036 (3.32-3.20)
Ramachandran outliers	110173	2000 (3.32-3.20)
Sidechain outliers	110143	1998 (3.32-3.20)
RSRZ outliers	101464	1861 (3.32-3.20)

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	391	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 62%, yellow 26%, orange 3%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>100%</span> <span>62%</span> <span>26%</span> <span>•</span> <span>9%</span> </div> </div>
1	B	391	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 63%, yellow 21%, orange 12%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>100%</span> <span>63%</span> <span>21%</span> <span>•</span> <span>12%</span> </div> </div>
1	C	391	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 60%, yellow 25%, orange 12%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>100%</span> <span>60%</span> <span>25%</span> <span>•</span> <span>12%</span> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	A	901	-	-	-	X

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8122 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	355	Total	C	N	O	S	0	0	0
			2747	1763	462	516	6			
1	C	346	Total	C	N	O	S	0	0	0
			2689	1729	448	506	6			
1	B	345	Total	C	N	O	S	0	0	0
			2683	1726	447	504	6			

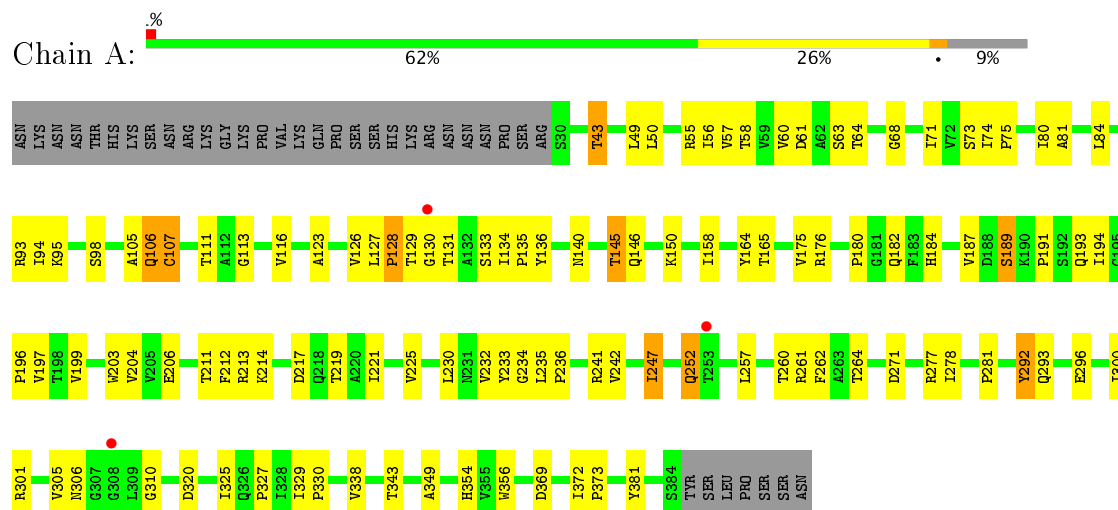
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		
2	C	1	Total	Ca	0	0
			1	1		

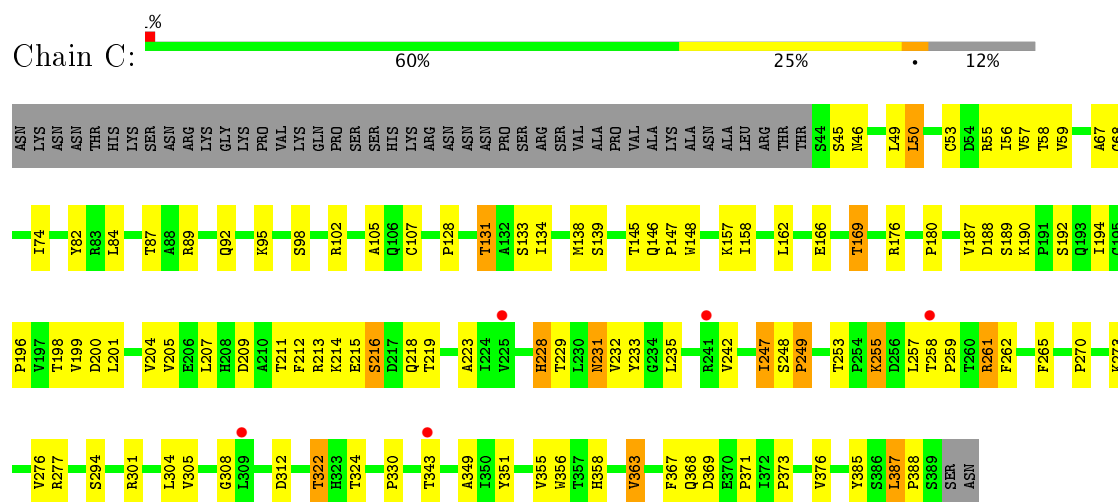
### 3 Residue-property plots [i](#)

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

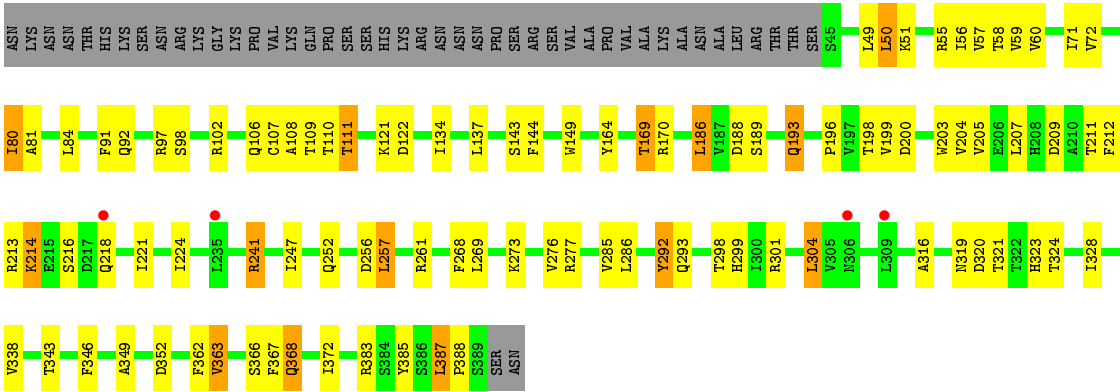


#### • Molecule 1: Capsid protein



#### • Molecule 1: Capsid protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	402.20Å 369.86Å 410.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.25 49.68 – 3.26	Depositor EDS
% Data completeness (in resolution range)	86.4 (50.00-3.25) 87.0 (49.68-3.26)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 3.25Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.278 , 0.279 0.263 , 0.263	Depositor DCC
$R_{free}$ test set	20302 reflections (5.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.0	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 55.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -l,-k,-h	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	8122	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.68% 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: CA

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	0.43	0/2823	0.65	0/3873
1	B	0.42	0/2760	0.66	0/3787
1	C	0.41	0/2766	0.64	0/3795
All	All	0.42	0/8349	0.65	0/11455

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	144	PHE	Peptide

### 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	2747	0	2709	72	0
1	B	2683	0	2634	56	0
1	C	2689	0	2641	80	0
2	A	2	0	0	0	0
2	C	1	0	0	0	0
All	All	8122	0	7984	197	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 12.

All (197) 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:C:228:HIS:HB3	1:C:249:PRO:HA	1.42	1.00
1:C:228:HIS:CB	1:C:249:PRO:HA	1.99	0.91
1:C:107:CYS:SG	1:C:148:TRP:HB3	2.12	0.89
1:A:176:ARG:HG2	1:B:164:TYR:CE2	2.22	0.73
1:A:247:ILE:HD11	1:A:257:LEU:HA	1.71	0.72
1:C:247:ILE:HD11	1:C:257:LEU:HG	1.72	0.71
1:C:82:TYR:HH	1:C:385:TYR:HH	1.36	0.70
1:A:130:GLY:HA2	1:A:133:SER:OG	1.91	0.70
1:A:164:TYR:CE2	1:C:176:ARG:HG2	2.28	0.69
1:C:211:THR:HG22	1:C:212:PHE:N	2.08	0.69
1:A:43:THR:HG21	1:C:138:MET:O	1.94	0.68
1:A:158:ILE:HD13	1:A:180:PRO:HB2	1.76	0.67
1:C:247:ILE:HD13	1:C:261:ARG:HG2	1.76	0.66
1:B:49:LEU:HD12	1:B:205:VAL:O	1.95	0.66
1:C:218:GLN:HG2	1:C:371:PRO:HB3	1.77	0.66
1:B:57:VAL:HA	1:B:388:PRO:HD3	1.77	0.65
1:B:81:ALA:HB3	1:B:84:LEU:HB3	1.78	0.64
1:C:259:PRO:HB2	1:C:270:PRO:HG2	1.79	0.64
1:C:218:GLN:HG2	1:C:371:PRO:CB	2.29	0.63
1:B:109:THR:HG22	1:B:149:TRP:CD2	2.34	0.62
1:B:211:THR:HG22	1:B:212:PHE:N	2.14	0.62
1:C:67:ALA:HA	1:C:187:VAL:HG23	1.81	0.61
1:A:176:ARG:HG2	1:B:164:TYR:CZ	2.35	0.61
1:A:123:ALA:HB1	1:A:175:VAL:HG22	1.81	0.61
1:C:262:PHE:CD2	1:C:270:PRO:HG3	2.36	0.61
1:B:387:LEU:O	1:B:388:PRO:C	2.36	0.61
1:C:57:VAL:CG1	1:C:199:VAL:HB	2.32	0.60
1:C:211:THR:HG22	1:C:212:PHE:H	1.67	0.60
1:C:223:ALA:HB2	1:C:265:PHE:CZ	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:GLY:HA2	1:A:242:VAL:HG22	1.83	0.60
1:B:277:ARG:HB3	1:B:363:VAL:HG13	1.83	0.59
1:C:134:ILE:HD13	1:C:188:ASP:HA	1.83	0.59
1:C:262:PHE:CG	1:C:270:PRO:HG3	2.38	0.59
1:A:211:THR:HG22	1:A:212:PHE:N	2.18	0.59
1:C:55:ARG:HD2	1:C:200:ASP:OD2	2.03	0.57
1:C:105:ALA:HB1	1:C:148:TRP:O	2.04	0.57
1:C:57:VAL:HG12	1:C:199:VAL:HB	1.86	0.57
1:B:211:THR:HG22	1:B:212:PHE:H	1.70	0.57
1:C:257:LEU:HD12	1:C:257:LEU:N	2.19	0.56
1:C:128:PRO:HG2	1:C:133:SER:HA	1.86	0.55
1:C:95:LYS:HB2	1:C:162:LEU:HD23	1.89	0.54
1:C:228:HIS:HB2	1:C:249:PRO:HA	1.88	0.54
1:B:107:CYS:HB2	1:B:111:THR:HG21	1.89	0.54
1:A:57:VAL:HG12	1:A:199:VAL:HB	1.90	0.54
1:A:75:PRO:O	1:A:80:ILE:HD11	2.07	0.54
1:C:55:ARG:CD	1:C:200:ASP:OD2	2.55	0.54
1:C:387:LEU:HD23	1:C:388:PRO:HA	1.90	0.54
1:A:64:THR:HG21	1:A:225:VAL:HG23	1.90	0.53
1:B:58:THR:HG23	1:B:387:LEU:HB2	1.89	0.53
1:B:286:LEU:HB3	1:B:346:PHE:CD1	2.44	0.53
1:A:277:ARG:NH1	1:A:369:ASP:OD1	2.42	0.53
1:B:213:ARG:HE	1:B:214:LYS:H	1.55	0.53
1:C:305:VAL:HG21	1:C:312:ASP:HB2	1.90	0.53
1:C:49:LEU:HD12	1:C:205:VAL:O	2.09	0.52
1:C:277:ARG:HB3	1:C:363:VAL:HG13	1.92	0.52
1:C:231:ASN:HD21	1:C:248:SER:HB2	1.73	0.52
1:A:194:ILE:HG22	1:A:196:PRO:HD3	1.92	0.52
1:A:330:PRO:HG3	1:A:349:ALA:HB2	1.92	0.52
1:A:165:THR:OG1	1:A:213:ARG:NH2	2.43	0.51
1:A:305:VAL:HG12	1:A:310:GLY:HA3	1.92	0.51
1:A:123:ALA:O	1:B:170:ARG:NH1	2.44	0.51
1:C:56:ILE:CG1	1:C:201:LEU:HB2	2.40	0.51
1:B:80:ILE:O	1:B:80:ILE:HD12	2.11	0.51
1:C:187:VAL:O	1:C:187:VAL:HG23	2.10	0.51
1:C:50:LEU:HD11	1:C:87:THR:HG21	1.91	0.51
1:B:276:VAL:HG11	1:B:362:PHE:HB3	1.92	0.51
1:C:211:THR:CG2	1:C:212:PHE:N	2.73	0.51
1:C:294:SER:HB3	1:B:352:ASP:OD2	2.11	0.51
1:B:304:LEU:H	1:B:304:LEU:HD23	1.75	0.50
1:B:109:THR:HG22	1:B:149:TRP:CE3	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:ASN:ND2	1:B:321:THR:HB	2.27	0.50
1:A:74:ILE:HD12	1:A:80:ILE:HD13	1.94	0.50
1:B:111:THR:HB	1:B:193:GLN:HG2	1.91	0.50
1:C:92:GLN:HE21	1:C:211:THR:HB	1.77	0.50
1:B:55:ARG:HD3	1:B:200:ASP:OD1	2.11	0.50
1:B:276:VAL:CG1	1:B:362:PHE:HB3	2.42	0.49
1:A:232:VAL:HG23	1:A:356:TRP:HZ3	1.76	0.49
1:B:56:ILE:HG23	1:B:80:ILE:HG12	1.94	0.49
1:C:134:ILE:CD1	1:C:188:ASP:HA	2.42	0.49
1:A:211:THR:HG22	1:A:212:PHE:H	1.78	0.49
1:A:217:ASP:HA	1:A:373:PRO:HA	1.95	0.49
1:C:247:ILE:HD12	1:C:255:LYS:O	2.13	0.49
1:A:252:GLN:HE21	1:A:252:GLN:N	2.10	0.49
1:C:146:GLN:HG3	1:C:147:PRO:HD2	1.95	0.49
1:C:242:VAL:HG12	1:C:351:TYR:CD2	2.47	0.48
1:C:89:ARG:NE	1:C:215:GLU:HG3	2.28	0.48
1:A:140:ASN:HD21	1:B:211:THR:HG23	1.79	0.48
1:C:258:THR:N	1:C:259:PRO:HA	2.29	0.48
1:B:134:ILE:HG12	1:B:186:LEU:HD23	1.95	0.48
1:A:128:PRO:HG2	1:A:136:TYR:CD2	2.49	0.47
1:C:166:GLU:OE1	1:C:213:ARG:HG2	2.13	0.47
1:C:223:ALA:HB3	1:C:358:HIS:HA	1.96	0.47
1:A:235:LEU:HB3	1:A:236:PRO:HD3	1.96	0.47
1:A:260:THR:O	1:A:264:THR:HG23	2.14	0.47
1:B:92:GLN:NE2	1:B:211:THR:OG1	2.47	0.47
1:C:258:THR:OG1	1:C:261:ARG:NH1	2.46	0.47
1:A:73:SER:HB3	1:A:184:HIS:CD2	2.50	0.47
1:A:63:SER:HA	1:A:191:PRO:HG2	1.96	0.47
1:C:128:PRO:O	1:C:133:SER:HB3	2.14	0.47
1:B:134:ILE:CD1	1:B:188:ASP:HA	2.45	0.47
1:C:214:LYS:HB3	1:C:216:SER:N	2.30	0.47
1:A:106:GLN:CG	1:A:196:PRO:HG2	2.45	0.47
1:C:194:ILE:HG22	1:C:196:PRO:HD3	1.97	0.47
1:A:105:ALA:HB1	1:A:107:CYS:SG	2.55	0.46
1:A:116:VAL:HG21	1:A:134:ILE:HG23	1.96	0.46
1:A:232:VAL:HG23	1:A:356:TRP:CZ3	2.50	0.46
1:C:53:CYS:HA	1:C:201:LEU:O	2.14	0.46
1:A:49:LEU:C	1:A:49:LEU:HD23	2.34	0.46
1:B:108:ALA:O	1:B:111:THR:CG2	2.64	0.46
1:B:292:TYR:CD2	1:B:292:TYR:C	2.88	0.46
1:C:232:VAL:HG23	1:C:356:TRP:CZ3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:TYR:HB2	1:B:338:VAL:HG12	1.96	0.46
1:B:366:SER:C	1:B:367:PHE:O	2.48	0.46
1:C:211:THR:CG2	1:C:212:PHE:H	2.29	0.46
1:B:247:ILE:HD12	1:B:261:ARG:HB3	1.96	0.46
1:C:74:ILE:HG23	1:C:74:ILE:O	2.16	0.46
1:A:111:THR:HG22	1:A:193:GLN:HB2	1.97	0.46
1:A:126:VAL:HG12	1:B:169:THR:HB	1.97	0.46
1:C:330:PRO:HG3	1:C:349:ALA:HB2	1.98	0.45
1:B:91:PHE:CG	1:B:207:LEU:HB3	2.51	0.45
1:A:128:PRO:HG2	1:A:136:TYR:CG	2.51	0.45
1:B:108:ALA:O	1:B:111:THR:HG23	2.16	0.45
1:C:59:VAL:HA	1:C:198:THR:HG22	1.98	0.45
1:A:271:ASP:OD1	1:A:381:TYR:CE1	2.69	0.45
1:B:268:PHE:HB3	1:B:383:ARG:HG3	1.99	0.45
1:C:242:VAL:HG12	1:C:351:TYR:HD2	1.82	0.45
1:A:194:ILE:H	1:A:194:ILE:HD12	1.81	0.45
1:B:50:LEU:HB3	1:B:203:TRP:HZ3	1.82	0.44
1:B:51:LYS:HG2	1:B:204:VAL:HG22	1.99	0.44
1:A:50:LEU:HB3	1:A:203:TRP:HZ3	1.82	0.44
1:B:323:HIS:O	1:B:368:GLN:NE2	2.50	0.44
1:A:94:ILE:HB	1:A:165:THR:HG22	1.99	0.44
1:B:137:LEU:CD2	1:B:186:LEU:HD22	2.48	0.44
1:A:71:ILE:CD1	1:A:197:VAL:HG21	2.48	0.44
1:A:93:ARG:HG2	1:A:93:ARG:HH11	1.83	0.44
1:C:169:THR:HG21	1:C:213:ARG:CZ	2.47	0.44
1:C:301:ARG:CZ	1:C:367:PHE:HE2	2.30	0.44
1:C:45:SER:OG	1:C:46:ASN:N	2.51	0.44
1:A:292:TYR:CD2	1:A:292:TYR:C	2.91	0.43
1:A:301:ARG:NH1	1:A:325:ILE:O	2.50	0.43
1:A:330:PRO:HG3	1:A:349:ALA:CB	2.48	0.43
1:A:56:ILE:HD12	1:A:56:ILE:HA	1.90	0.43
1:B:97:ARG:HH11	1:B:97:ARG:HG2	1.82	0.43
1:A:146:GLN:HG3	1:A:150:LYS:HB2	2.00	0.43
1:A:257:LEU:HD22	1:A:262:PHE:HB2	1.99	0.43
1:B:247:ILE:CG1	1:B:257:LEU:HG	2.48	0.43
1:C:322:THR:HB	1:C:324:THR:HG22	1.99	0.43
1:C:213:ARG:HG3	1:C:215:GLU:HA	2.01	0.43
1:B:57:VAL:HG22	1:B:199:VAL:HG23	2.01	0.43
1:C:373:PRO:HB2	1:C:376:VAL:HG23	2.01	0.43
1:C:58:THR:HG23	1:C:387:LEU:HB3	2.01	0.43
1:A:305:VAL:HG13	1:A:306:ASN:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:HIS:HD2	1:A:356:TRP:CE2	2.37	0.43
1:C:92:GLN:OE1	1:B:122:ASP:HB2	2.19	0.43
1:C:102:ARG:NH1	1:C:102:ARG:HB2	2.34	0.43
1:A:281:PRO:HB3	1:A:296:GLU:HG3	2.00	0.43
1:B:316:ALA:HA	1:B:328:ILE:HG13	2.00	0.43
1:C:158:ILE:HD12	1:C:180:PRO:HB2	2.00	0.43
1:C:229:THR:HB	1:C:356:TRP:O	2.19	0.43
1:A:130:GLY:O	1:A:131:THR:C	2.58	0.42
1:A:95:LYS:HB3	1:A:206:GLU:HB2	2.01	0.42
1:A:338:VAL:HG23	1:C:235:LEU:O	2.18	0.42
1:C:68:GLY:HA3	1:C:131:THR:HA	2.00	0.42
1:C:304:LEU:HB3	1:C:308:GLY:HA2	2.01	0.42
1:B:59:VAL:HG22	1:B:198:THR:HG22	2.01	0.42
1:A:127:LEU:C	1:A:128:PRO:O	2.57	0.42
1:A:278:ILE:HD12	1:A:278:ILE:HA	1.91	0.42
1:B:299:HIS:O	1:B:316:ALA:N	2.50	0.42
1:B:385:TYR:OH	1:B:387:LEU:HD13	2.20	0.42
1:C:262:PHE:CD2	1:C:262:PHE:C	2.92	0.42
1:A:123:ALA:CB	1:A:175:VAL:HG22	2.46	0.42
1:A:56:ILE:HG13	1:A:80:ILE:HG21	2.01	0.42
1:A:116:VAL:HG22	1:A:145:THR:HB	2.02	0.42
1:A:278:ILE:HD13	1:A:300:ILE:HG13	2.02	0.42
1:B:57:VAL:HG11	1:B:72:VAL:HG22	2.02	0.42
1:C:56:ILE:CG2	1:C:74:ILE:HD13	2.50	0.42
1:C:261:ARG:HB2	1:C:261:ARG:HH11	1.85	0.41
1:B:241:ARG:NH1	1:B:256:ASP:HB3	2.34	0.41
1:C:231:ASN:ND2	1:C:233:TYR:OH	2.53	0.41
1:A:134:ILE:N	1:A:135:PRO:HD2	2.35	0.41
1:C:187:VAL:O	1:C:187:VAL:CG2	2.68	0.41
1:A:327:PRO:HB2	1:A:329:ILE:HD13	2.01	0.41
1:C:231:ASN:HB3	1:C:355:VAL:HA	2.02	0.41
1:B:71:ILE:HG13	1:B:72:VAL:HG23	2.02	0.41
1:C:367:PHE:HD1	1:C:367:PHE:H	1.68	0.41
1:A:113:GLY:HA2	1:A:189:SER:HB3	2.03	0.41
1:A:278:ILE:HD11	1:A:356:TRP:CE2	2.56	0.41
1:A:56:ILE:HG23	1:A:74:ILE:HD11	2.03	0.41
1:A:68:GLY:O	1:A:130:GLY:HA3	2.20	0.41
1:B:285:VAL:O	1:B:349:ALA:HA	2.21	0.41
1:A:305:VAL:HG12	1:A:310:GLY:CA	2.51	0.41
1:C:218:GLN:HG2	1:C:371:PRO:HB2	2.03	0.40
1:B:60:VAL:O	1:B:196:PRO:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:ARG:HH11	1:C:102:ARG:HB2	1.86	0.40
1:B:269:LEU:HA	1:B:269:LEU:HD12	1.92	0.40
1:A:247:ILE:HD13	1:A:261:ARG:HD2	2.02	0.40
1:A:81:ALA:HB3	1:A:84:LEU:HB3	2.03	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	353/391 (90%)	334 (95%)	17 (5%)	2 (1%)	28	68
1	B	343/391 (88%)	324 (94%)	17 (5%)	2 (1%)	28	68
1	C	344/391 (88%)	322 (94%)	19 (6%)	3 (1%)	20	60
All	All	1040/1173 (89%)	980 (94%)	53 (5%)	7 (1%)	25	65

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	PRO
1	C	369	ASP
1	C	209	ASP
1	A	61	ASP
1	B	209	ASP
1	C	249	PRO
1	B	387	LEU

### 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	303/338 (90%)	278 (92%)	25 (8%)	13	45
1	B	297/338 (88%)	264 (89%)	33 (11%)	7	29
1	C	298/338 (88%)	270 (91%)	28 (9%)	10	37
All	All	898/1014 (89%)	812 (90%)	86 (10%)	10	36

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	THR
1	A	55	ARG
1	A	58	THR
1	A	60	VAL
1	A	98	SER
1	A	106	GLN
1	A	107	CYS
1	A	129	THR
1	A	145	THR
1	A	182	GLN
1	A	187	VAL
1	A	189	SER
1	A	204	VAL
1	A	214	LYS
1	A	219	THR
1	A	221	ILE
1	A	230	LEU
1	A	241	ARG
1	A	247	ILE
1	A	252	GLN
1	A	292	TYR
1	A	293	GLN
1	A	320	ASP
1	A	343	THR
1	A	372	ILE
1	C	50	LEU
1	C	84	LEU
1	C	98	SER
1	C	131	THR
1	C	139	SER

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Mol	Chain	Res	Type
1	C	145	THR
1	C	157	LYS
1	C	169	THR
1	C	189	SER
1	C	190	LYS
1	C	192	SER
1	C	204	VAL
1	C	207	LEU
1	C	216	SER
1	C	219	THR
1	C	228	HIS
1	C	231	ASN
1	C	247	ILE
1	C	253	THR
1	C	255	LYS
1	C	261	ARG
1	C	273	LYS
1	C	276	VAL
1	C	322	THR
1	C	343	THR
1	C	363	VAL
1	C	368	GLN
1	C	387	LEU
1	B	50	LEU
1	B	80	ILE
1	B	98	SER
1	B	102	ARG
1	B	106	GLN
1	B	110	THR
1	B	111	THR
1	B	121	LYS
1	B	143	SER
1	B	169	THR
1	B	186	LEU
1	B	189	SER
1	B	193	GLN
1	B	214	LYS
1	B	216	SER
1	B	218	GLN
1	B	221	ILE
1	B	224	ILE
1	B	241	ARG

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Mol	Chain	Res	Type
1	B	252	GLN
1	B	257	LEU
1	B	273	LYS
1	B	292	TYR
1	B	293	GLN
1	B	298	THR
1	B	301	ARG
1	B	304	LEU
1	B	320	ASP
1	B	324	THR
1	B	343	THR
1	B	363	VAL
1	B	368	GLN
1	B	372	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	140	ASN
1	A	184	HIS
1	A	252	GLN
1	A	319	ASN
1	A	354	HIS
1	C	182	GLN
1	C	231	ASN
1	C	326	GLN
1	B	140	ASN
1	B	228	HIS
1	B	244	HIS
1	B	319	ASN
1	B	368	GLN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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 [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	355/391 (90%)	0.11	3 (0%) 86 80	15, 39, 82, 185	0
1	B	345/391 (88%)	0.15	4 (1%) 79 71	14, 43, 84, 127	0
1	C	346/391 (88%)	0.26	5 (1%) 75 67	15, 47, 95, 162	0
All	All	1046/1173 (89%)	0.17	12 (1%) 80 73	14, 43, 91, 185	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	309	LEU	3.8
1	C	343	THR	2.7
1	C	258	THR	2.6
1	B	306	ASN	2.5
1	A	130	GLY	2.4
1	A	308	GLY	2.4
1	B	235	LEU	2.4
1	C	225	VAL	2.3
1	C	309	LEU	2.2
1	B	218	GLN	2.2
1	A	253	THR	2.2
1	C	241	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](#)

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(Å <sup>2</sup> )	Q<0.9
2	CA	A	901	1/1	0.83	0.29	3.75	94,94,94,94	0
2	CA	A	902	1/1	0.89	0.27	1.32	80,80,80,80	0
2	CA	C	901	1/1	0.90	0.14	-1.48	73,73,73,73	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.