



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:55 AM UTC

PDB ID : 2CKW / pdb_00002ckw
Title : The 2.3 Å resolution structure of the Sapporo virus RNA dependant RNA polymerase.
Authors : Fullerton, S.W.B.; Tucker, P.A.; Rohayem, J.; Coutard, B.; Gebhardt, J.; Gorbalenya, A.; Canard, B.
Deposited on : 2006-04-24
Resolution : 2.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

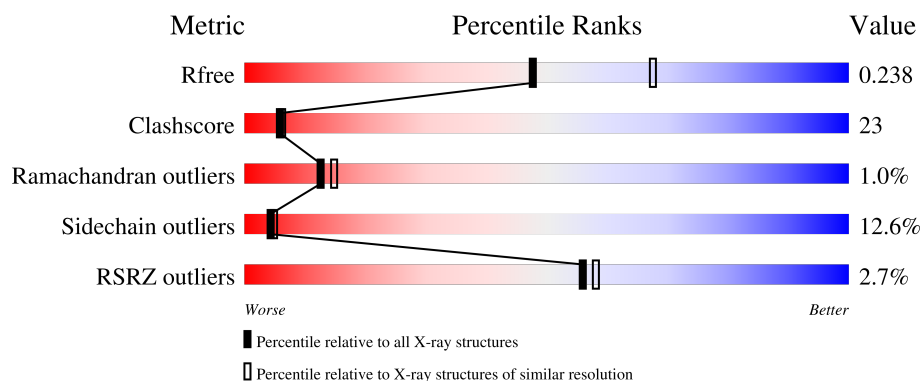
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 ≥ 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	515	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4037 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 RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	0	0
			3766	2392	649	708	17			

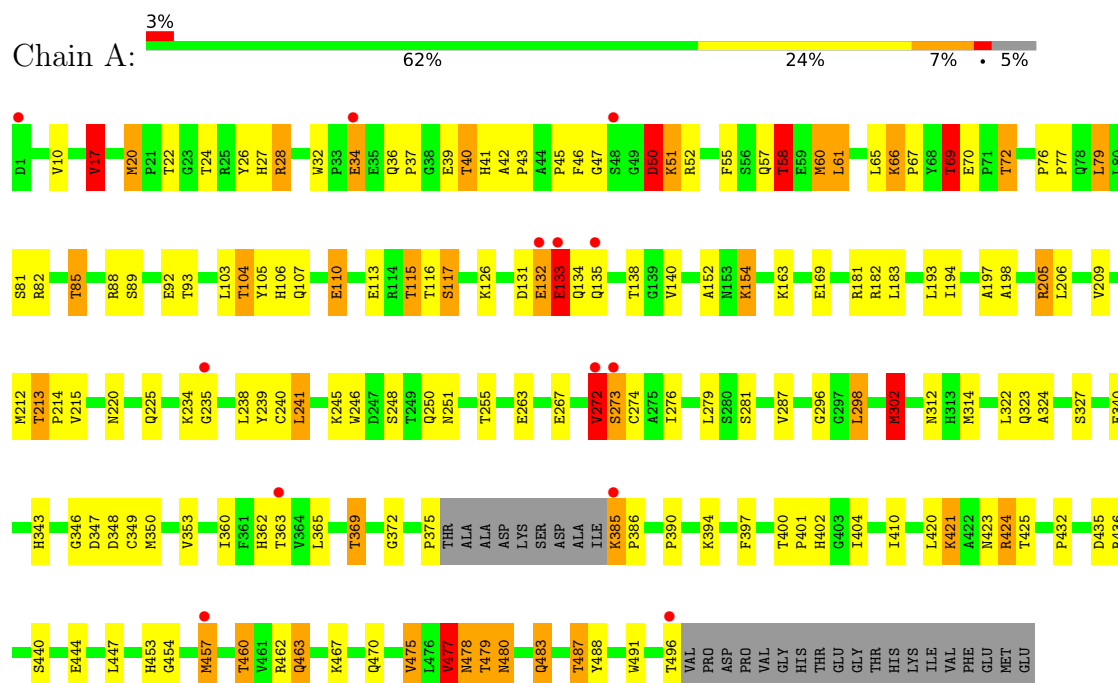
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	271	Total	O	0	0
			271	271		

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: RNA-DIRECTED RNA POLYMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.43Å 93.88Å 94.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.67 – 2.30 66.67 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (66.67-2.30) 99.0 (66.67-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.171 , 0.229 0.190 , 0.238	Depositor DCC
R_{free} test set	1209 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.023 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4037	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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

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	1.21	1/3863 (0.0%)	1.22	20/5272 (0.4%)

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	A	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	205	ARG	CD-NE	-5.36	1.38	1.46

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	ARG	NE-CZ-NH2	-7.58	112.37	119.20
1	A	50	ASP	N-CA-C	6.80	119.81	108.73
1	A	20	MET	CA-C-N	6.78	126.81	119.90
1	A	20	MET	C-N-CA	6.78	126.81	119.90
1	A	17	VAL	N-CA-C	-6.76	106.41	112.90
1	A	69	THR	N-CA-C	-6.61	105.37	113.50
1	A	477	VAL	N-CA-C	6.33	117.26	111.81
1	A	447	LEU	N-CA-C	6.29	118.13	111.28
1	A	460	THR	N-CA-C	5.78	117.26	111.07
1	A	17	VAL	N-CA-CB	-5.63	105.00	112.26
1	A	60	MET	CG-SD-CE	-5.61	88.55	100.90
1	A	475	VAL	N-CA-C	5.61	115.96	108.11
1	A	267	GLU	CA-C-N	-5.27	114.14	119.83
1	A	267	GLU	C-N-CA	-5.27	114.14	119.83
1	A	58	THR	N-CA-CB	5.13	117.67	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	302	MET	N-CA-CB	5.12	117.75	110.03
1	A	372	GLY	N-CA-C	-5.06	108.08	115.63
1	A	480	ASN	CA-C-N	5.05	127.35	120.54
1	A	480	ASN	C-N-CA	5.05	127.35	120.54
1	A	193	LEU	N-CA-C	5.02	117.13	111.11

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	GLU	Peptide
1	A	133	GLU	Peptide
1	A	198	ALA	Peptide
1	A	272	VAL	Peptide
1	A	273	SER	Peptide
1	A	454	GLY	Peptide

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	3766	0	3710	172	0
2	A	271	0	0	32	0
All	All	4037	0	3710	172	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 23.

All (172) 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:46:PHE:CZ	1:A:50:ASP:HB3	1.69	1.25
1:A:45:PRO:CG	1:A:60:MET:HE1	1.67	1.25
1:A:360:ILE:O	1:A:363:THR:HG22	1.38	1.21
1:A:154:LYS:HE3	2:A:2112:HOH:O	1.52	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:MET:HE1	1:A:390:PRO:HB3	1.32	1.07
1:A:404:ILE:HD12	1:A:404:ILE:H	1.21	1.01
1:A:45:PRO:HG2	1:A:60:MET:HE1	1.01	1.01
1:A:46:PHE:CE2	1:A:50:ASP:HB3	1.96	1.01
1:A:45:PRO:HG2	1:A:60:MET:CE	1.92	1.00
1:A:46:PHE:CZ	1:A:50:ASP:CB	2.45	0.99
1:A:34:GLU:HG2	2:A:2254:HOH:O	1.62	0.97
1:A:250:GLN:HE21	1:A:312:ASN:HD22	0.97	0.95
1:A:404:ILE:HD12	1:A:404:ILE:N	1.83	0.92
1:A:34:GLU:HG3	2:A:2025:HOH:O	1.69	0.90
1:A:213:THR:HG22	1:A:215:VAL:H	1.39	0.87
1:A:40:THR:HG22	1:A:41:HIS:ND1	1.91	0.86
1:A:410:ILE:CD1	1:A:457:MET:SD	2.64	0.85
1:A:61:LEU:HD23	1:A:183:LEU:CD2	2.07	0.84
1:A:477:VAL:C	1:A:479:THR:H	1.85	0.83
1:A:205:ARG:CD	2:A:2072:HOH:O	2.30	0.80
1:A:245:LYS:HG2	2:A:2062:HOH:O	1.81	0.79
1:A:51:LYS:HB3	2:A:2110:HOH:O	1.84	0.77
1:A:272:VAL:O	1:A:272:VAL:CG1	2.31	0.77
1:A:350:MET:HE1	1:A:390:PRO:CB	2.15	0.76
1:A:324:ALA:HB1	1:A:363:THR:CG2	2.17	0.74
1:A:42:ALA:HB1	1:A:43:PRO:HD2	1.70	0.73
1:A:238:LEU:HD11	1:A:350:MET:HE2	1.71	0.73
1:A:263:GLU:HB2	1:A:276:ILE:HD11	1.69	0.73
1:A:246:TRP:CG	1:A:347:ASP:HB3	2.24	0.72
1:A:72:THR:HG22	1:A:251:ASN:HD22	1.53	0.72
1:A:410:ILE:HD13	1:A:457:MET:SD	2.29	0.72
1:A:131:ASP:OD1	1:A:134:GLN:HB2	1.89	0.72
1:A:404:ILE:H	1:A:404:ILE:CD1	2.00	0.72
1:A:47:GLY:HA2	1:A:55:PHE:O	1.90	0.71
1:A:324:ALA:CB	1:A:363:THR:HG23	2.21	0.71
1:A:402:HIS:HD2	2:A:2102:HOH:O	1.74	0.71
1:A:61:LEU:CD2	1:A:183:LEU:CD2	2.69	0.71
1:A:205:ARG:HD3	2:A:2072:HOH:O	1.90	0.70
1:A:79:LEU:HD11	1:A:323:GLN:HB2	1.72	0.70
1:A:250:GLN:NE2	1:A:312:ASN:HD22	1.80	0.70
1:A:113:GLU:OE1	1:A:115:THR:HB	1.90	0.70
1:A:40:THR:CG2	1:A:41:HIS:ND1	2.55	0.70
1:A:205:ARG:HD2	2:A:2072:HOH:O	1.92	0.70
1:A:457:MET:HE3	1:A:457:MET:C	2.17	0.69
1:A:46:PHE:CE1	1:A:50:ASP:HB3	2.26	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:VAL:O	1:A:272:VAL:HG12	1.92	0.69
1:A:213:THR:HG22	1:A:215:VAL:N	2.07	0.68
1:A:104:THR:HG23	2:A:2085:HOH:O	1.94	0.67
1:A:169:GLU:OE2	1:A:182:ARG:HD3	1.95	0.67
1:A:135:GLN:O	1:A:135:GLN:HG3	1.94	0.67
1:A:477:VAL:C	1:A:479:THR:N	2.52	0.66
1:A:22:THR:O	1:A:424:ARG:NH2	2.28	0.66
1:A:82:ARG:HD2	2:A:2059:HOH:O	1.95	0.66
1:A:410:ILE:HG12	1:A:457:MET:SD	2.37	0.65
1:A:61:LEU:HD23	1:A:183:LEU:HD21	1.76	0.64
1:A:42:ALA:HB1	1:A:43:PRO:CD	2.26	0.64
1:A:220:ASN:H	1:A:225:GLN:NE2	1.95	0.64
1:A:483:GLN:O	1:A:487:THR:HG22	1.98	0.64
1:A:81:SER:O	1:A:85:THR:CG2	2.46	0.64
1:A:324:ALA:HB1	1:A:363:THR:HG23	1.76	0.64
1:A:365:LEU:O	1:A:369:THR:HG23	1.98	0.63
1:A:302:MET:HE3	2:A:2181:HOH:O	1.98	0.63
1:A:212:MET:CG	2:A:2149:HOH:O	2.46	0.63
1:A:213:THR:CG2	1:A:215:VAL:H	2.10	0.62
1:A:483:GLN:O	1:A:487:THR:CG2	2.47	0.62
1:A:66:LYS:O	1:A:69:THR:HB	2.00	0.62
1:A:26:TYR:CE1	1:A:421:LYS:NZ	2.69	0.61
1:A:479:THR:HG21	2:A:2259:HOH:O	2.01	0.60
1:A:477:VAL:O	1:A:479:THR:N	2.34	0.60
1:A:103:LEU:HB2	1:A:197:ALA:HA	1.84	0.59
1:A:213:THR:HB	2:A:2150:HOH:O	2.02	0.58
1:A:45:PRO:HG3	1:A:60:MET:HE1	1.77	0.58
1:A:61:LEU:CD2	1:A:183:LEU:HD21	2.33	0.57
1:A:212:MET:HG2	2:A:2149:HOH:O	2.04	0.57
1:A:89:SER:O	1:A:93:THR:HG23	2.05	0.57
1:A:360:ILE:HA	1:A:362:HIS:CE1	2.40	0.57
1:A:88:ARG:O	1:A:92:GLU:HG3	2.05	0.57
1:A:133:GLU:HG3	1:A:133:GLU:O	2.05	0.57
1:A:46:PHE:CE1	1:A:50:ASP:CB	2.86	0.56
1:A:22:THR:HG22	2:A:2017:HOH:O	2.04	0.56
1:A:61:LEU:CD2	1:A:183:LEU:HD22	2.37	0.55
1:A:85:THR:HG21	2:A:2057:HOH:O	2.05	0.55
1:A:462:ARG:HH11	1:A:463:GLN:HE22	1.53	0.55
1:A:302:MET:CE	2:A:2181:HOH:O	2.53	0.55
1:A:410:ILE:HD11	1:A:457:MET:SD	2.45	0.55
1:A:410:ILE:CG1	1:A:457:MET:SD	2.94	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:HIS:HE1	1:A:423:ASN:O	1.90	0.55
1:A:324:ALA:HB1	1:A:363:THR:HG21	1.87	0.55
1:A:234:LYS:HD2	1:A:235:GLY:H	1.72	0.54
1:A:453:HIS:HB2	1:A:457:MET:HE2	1.88	0.54
1:A:279:LEU:HD22	1:A:302:MET:HE2	1.90	0.54
1:A:152:ALA:HA	1:A:274:CYS:SG	2.48	0.54
1:A:239:TYR:HB3	1:A:241:LEU:HD21	1.88	0.54
1:A:246:TRP:CD1	1:A:347:ASP:HB3	2.43	0.54
1:A:281:SER:HB2	2:A:2078:HOH:O	2.07	0.53
1:A:324:ALA:HB2	1:A:363:THR:HG23	1.90	0.53
1:A:400:THR:HB	1:A:401:PRO:CD	2.39	0.53
1:A:250:GLN:HE21	1:A:312:ASN:ND2	1.83	0.53
1:A:45:PRO:CG	1:A:60:MET:CE	2.62	0.52
1:A:104:THR:HG23	1:A:105:TYR:N	2.26	0.51
1:A:385:LYS:HB2	1:A:386:PRO:HD2	1.92	0.51
1:A:478:ASN:ND2	1:A:487:THR:HG21	2.26	0.51
1:A:81:SER:O	1:A:85:THR:HG23	2.10	0.51
1:A:245:LYS:CG	2:A:2062:HOH:O	2.51	0.51
1:A:81:SER:O	1:A:85:THR:HG22	2.11	0.51
1:A:206:LEU:HD22	1:A:213:THR:HG21	1.93	0.50
1:A:420:LEU:HD21	1:A:432:PRO:HG3	1.93	0.50
1:A:420:LEU:HD21	1:A:432:PRO:CG	2.42	0.50
1:A:66:LYS:HE3	2:A:2046:HOH:O	2.11	0.50
1:A:272:VAL:O	1:A:272:VAL:HG13	2.11	0.50
1:A:369:THR:HG22	1:A:375:PRO:CG	2.42	0.50
1:A:214:PRO:HG2	1:A:314:MET:HE3	1.94	0.50
1:A:104:THR:CG2	1:A:106:HIS:H	2.24	0.49
1:A:444:GLU:OE2	1:A:477:VAL:N	2.31	0.49
1:A:480:ASN:HB3	2:A:2261:HOH:O	2.11	0.49
1:A:116:THR:HG23	1:A:126:LYS:NZ	2.28	0.49
1:A:131:ASP:CG	1:A:134:GLN:HB2	2.38	0.48
1:A:212:MET:HG3	2:A:2149:HOH:O	2.11	0.48
1:A:82:ARG:HG2	1:A:322:LEU:HD13	1.94	0.48
1:A:20:MET:HB3	2:A:2124:HOH:O	2.14	0.48
1:A:104:THR:HG22	1:A:107:GLN:H	1.77	0.48
1:A:340:GLU:HG3	1:A:353:VAL:HA	1.96	0.48
1:A:296:GLY:HA2	2:A:2189:HOH:O	2.14	0.48
1:A:26:TYR:CZ	1:A:421:LYS:NZ	2.82	0.48
1:A:46:PHE:CZ	1:A:50:ASP:HB2	2.43	0.48
1:A:478:ASN:HD21	1:A:487:THR:HG21	1.79	0.48
1:A:36:GLN:HB3	1:A:37:PRO:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:THR:CG2	1:A:251:ASN:HB2	2.44	0.47
1:A:69:THR:HG22	1:A:70:GLU:HG2	1.97	0.47
1:A:240:CYS:SG	1:A:350:MET:HE3	2.54	0.47
1:A:51:LYS:O	1:A:51:LYS:HG3	2.13	0.47
1:A:17:VAL:HG13	1:A:163:LYS:HE3	1.97	0.47
1:A:234:LYS:HD2	1:A:235:GLY:N	2.30	0.47
1:A:239:TYR:HB3	1:A:241:LEU:CD2	2.44	0.46
1:A:57:GLN:HE21	1:A:181:ARG:HH12	1.63	0.46
1:A:72:THR:HB	2:A:2164:HOH:O	2.15	0.46
1:A:241:LEU:HD23	1:A:241:LEU:N	2.30	0.46
1:A:116:THR:OG1	1:A:117:SER:N	2.49	0.46
1:A:453:HIS:CB	1:A:457:MET:HE2	2.45	0.46
1:A:444:GLU:HG3	2:A:2244:HOH:O	2.15	0.45
1:A:134:GLN:O	1:A:135:GLN:HG2	2.17	0.45
1:A:20:MET:HE1	1:A:58:THR:HA	1.99	0.45
1:A:61:LEU:HD22	1:A:65:LEU:HG	1.98	0.45
1:A:106:HIS:O	1:A:110:GLU:HG2	2.17	0.44
1:A:163:LYS:HE2	2:A:2123:HOH:O	2.17	0.44
1:A:302:MET:HE2	1:A:302:MET:HB2	1.56	0.44
1:A:385:LYS:CB	1:A:386:PRO:HD2	2.47	0.44
1:A:462:ARG:NH1	1:A:463:GLN:HE22	2.16	0.44
1:A:28:ARG:HD3	1:A:32:TRP:O	2.18	0.44
1:A:220:ASN:H	1:A:225:GLN:HE21	1.64	0.43
1:A:298:LEU:HD13	2:A:2167:HOH:O	2.18	0.43
1:A:24:THR:O	1:A:425:THR:HG22	2.19	0.43
1:A:246:TRP:CG	1:A:347:ASP:CB	3.00	0.43
1:A:255:THR:HG21	1:A:298:LEU:HD11	1.99	0.43
1:A:76:PRO:HA	1:A:77:PRO:HD3	1.81	0.43
1:A:390:PRO:HB2	1:A:397:PHE:CG	2.53	0.43
1:A:72:THR:CG2	1:A:251:ASN:HD22	2.27	0.43
1:A:104:THR:HG23	1:A:106:HIS:H	1.83	0.43
1:A:400:THR:HB	1:A:401:PRO:HD2	2.00	0.42
1:A:343:HIS:O	1:A:349:CYS:HA	2.20	0.42
1:A:402:HIS:CD2	2:A:2102:HOH:O	2.59	0.42
1:A:272:VAL:O	1:A:276:ILE:HD12	2.20	0.41
1:A:113:GLU:HB3	1:A:116:THR:HG22	2.02	0.41
1:A:67:PRO:HB2	1:A:248:SER:HB2	2.01	0.41
1:A:79:LEU:HD23	1:A:79:LEU:HA	1.86	0.41
1:A:340:GLU:HG3	1:A:353:VAL:HG12	2.02	0.41
1:A:477:VAL:HG12	1:A:478:ASN:N	2.35	0.41
1:A:483:GLN:HE21	1:A:483:GLN:HA	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD21	1:A:183:LEU:HD22	2.03	0.41
1:A:28:ARG:NH2	1:A:39:GLU:OE1	2.54	0.40
1:A:478:ASN:ND2	2:A:2258:HOH:O	2.51	0.40
1:A:488:TYR:O	1:A:491:TRP:HB2	2.21	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	483/515 (94%)	461 (95%)	17 (4%)	5 (1%)	12	15

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	133	GLU
1	A	478	ASN
1	A	346	GLY
1	A	424	ARG
1	A	273	SER

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	404/427 (95%)	353 (87%)	51 (13%)	4 5

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	17	VAL
1	A	28	ARG
1	A	34	GLU
1	A	40	THR
1	A	50	ASP
1	A	51	LYS
1	A	52	ARG
1	A	58	THR
1	A	61	LEU
1	A	66	LYS
1	A	69	THR
1	A	72	THR
1	A	79	LEU
1	A	85	THR
1	A	104	THR
1	A	110	GLU
1	A	115	THR
1	A	117	SER
1	A	132	GLU
1	A	138	THR
1	A	140	VAL
1	A	154	LYS
1	A	194	ILE
1	A	209	VAL
1	A	213	THR
1	A	241	LEU
1	A	272	VAL
1	A	287	VAL
1	A	298	LEU
1	A	302	MET
1	A	327	SER
1	A	348	ASP
1	A	369	THR
1	A	385	LYS
1	A	394	LYS
1	A	421	LYS
1	A	435	ASP

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Mol	Chain	Res	Type
1	A	436	ARG
1	A	440	SER
1	A	457	MET
1	A	460	THR
1	A	463	GLN
1	A	467	LYS
1	A	470	GLN
1	A	475	VAL
1	A	477	VAL
1	A	479	THR
1	A	483	GLN
1	A	487	THR
1	A	496	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	HIS
1	A	57	GLN
1	A	86	HIS
1	A	106	HIS
1	A	134	GLN
1	A	147	GLN
1	A	225	GLN
1	A	250	GLN
1	A	251	ASN
1	A	288	ASN
1	A	323	GLN
1	A	328	HIS
1	A	362	HIS
1	A	402	HIS
1	A	452	GLN
1	A	453	HIS
1	A	463	GLN
1	A	478	ASN
1	A	483	GLN

5.3.3 RNA ⓘ

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 oligosaccharides 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	487/515 (94%)	-0.09	13 (2%) 56 58	6, 17, 35, 55	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	496	THR	4.0
1	A	273	SER	3.7
1	A	457	MET	3.1
1	A	48	SER	2.9
1	A	132	GLU	2.9
1	A	1	ASP	2.9
1	A	385	LYS	2.6
1	A	272	VAL	2.6
1	A	363	THR	2.5
1	A	235	GLY	2.4
1	A	133	GLU	2.3
1	A	135	GLN	2.3
1	A	34	GLU	2.2

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 oligosaccharides 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.