Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:54 AM GMT

PDB ID	:	3CM8
Title	:	A RNA polymerase subunit structure from virus
Authors	:	He, X.; Zhou, J.; Zeng, Z.; Ma, J.; Zhang, R.; Rao, Z.; Liu, Y.
Deposited on	:	2008-03-21
Resolution	:	2.90 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

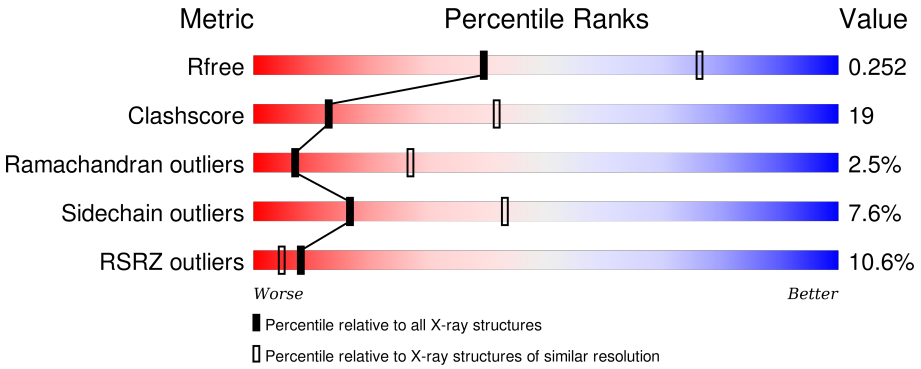
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. 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 <=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	471	10%, 59%, 28%, 6%, 8%
2	B	30	43%, 10%, 47%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3674 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 Polymerase acidic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3502	2234	590	650	28			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	246	GLY	-	EXPRESSION TAG	UNP Q9EA60
A	247	PRO	-	EXPRESSION TAG	UNP Q9EA60
A	248	LEU	-	EXPRESSION TAG	UNP Q9EA60
A	249	GLY	-	EXPRESSION TAG	UNP Q9EA60
A	250	SER	-	EXPRESSION TAG	UNP Q9EA60
A	251	PRO	-	EXPRESSION TAG	UNP Q9EA60
A	252	GLU	-	EXPRESSION TAG	UNP Q9EA60
A	253	PHE	-	EXPRESSION TAG	UNP Q9EA60
A	254	PRO	-	EXPRESSION TAG	UNP Q9EA60
A	255	GLY	-	EXPRESSION TAG	UNP Q9EA60

- Molecule 2 is a protein called peptide from RNA-directed RNA polymerase catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	16	Total	C	N	O	S	0	0	0
			123	81	19	22	1			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	EXPRESSION TAG	UNP Q9WLS3
B	-2	PRO	-	EXPRESSION TAG	UNP Q9WLS3
B	-1	LEU	-	EXPRESSION TAG	UNP Q9WLS3
B	0	GLY	-	EXPRESSION TAG	UNP Q9WLS3
B	1	SER	-	EXPRESSION TAG	UNP Q9WLS3

- Molecule 3 is water.

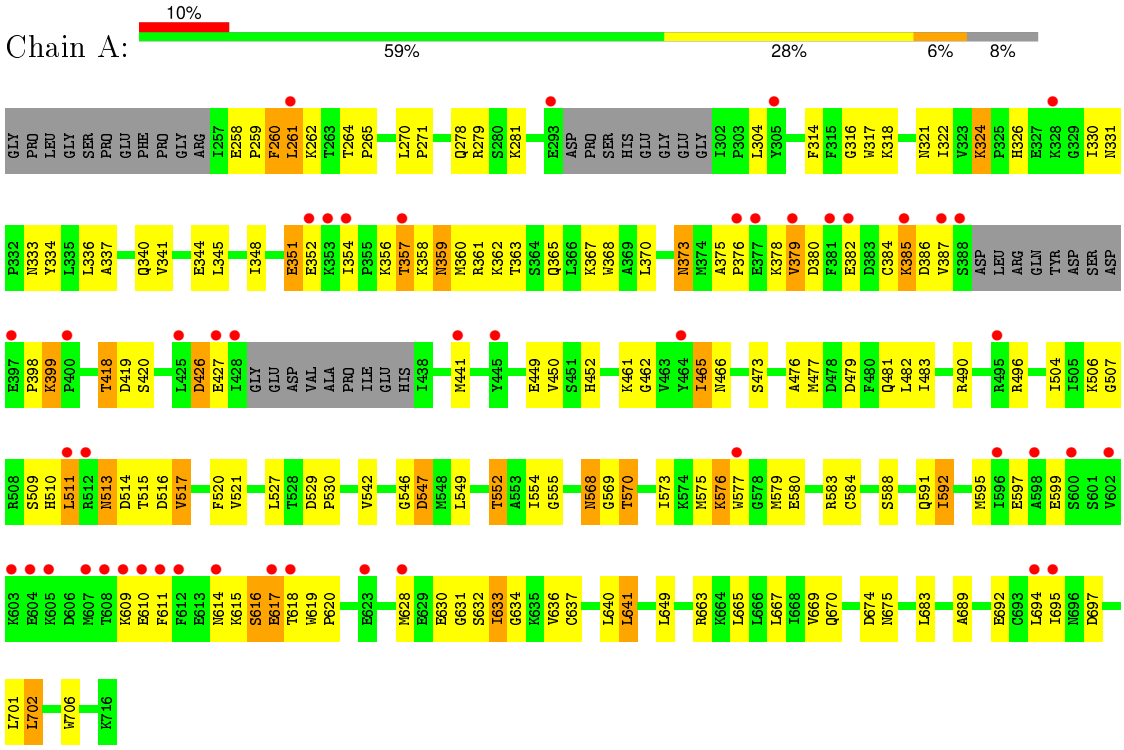
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	49	Total O 49 49	0	0

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 errors 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.

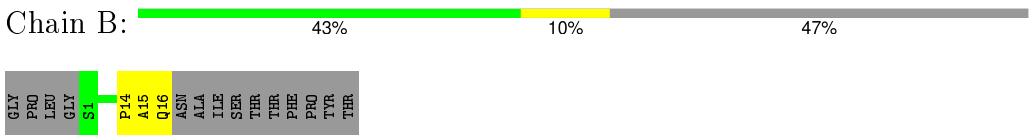
- Molecule 1: Polymerase acidic protein

Chain A:



- Molecule 2: peptide from RNA-directed RNA polymerase catalytic subunit

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	121.94Å 121.94Å 134.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.18 – 2.90 45.18 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (45.18-2.90) 99.3 (45.18-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.62 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.224 , 0.261 0.210 , 0.252	Depositor DCC
R_{free} test set	1174 reflections (5.12%)	DCC
Wilson B-factor (Å^2)	75.5	Xtriage
Anisotropy	0.520	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 79.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Outliers	0 of 22964 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3674	wwPDB-VP
Average B, all atoms (Å^2)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.34	0/3575	0.51	0/4816
2	B	0.31	0/125	0.48	0/170
All	All	0.34	0/3700	0.51	0/4986

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	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	A	261	LEU	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	3502	0	3530	138	0
2	B	123	0	133	6	0
3	A	49	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3674	0	3663	140	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 19.

All (140) 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:611:PHE:HB3	1:A:633:ILE:HG12	1.39	1.05
1:A:552:THR:HG22	1:A:554:ILE:H	1.30	0.96
1:A:278:GLN:HE21	1:A:279:ARG:H	1.08	0.95
1:A:385:LYS:H	1:A:385:LYS:HD3	1.33	0.92
1:A:633:ILE:H	1:A:633:ILE:HD12	1.37	0.88
1:A:461:LYS:O	1:A:465:ILE:HG23	1.74	0.85
1:A:592:ILE:HD11	1:A:636:VAL:HG12	1.56	0.85
1:A:568:ASN:HD22	1:A:569:GLY:H	1.31	0.79
1:A:278:GLN:HE21	1:A:279:ARG:N	1.81	0.77
1:A:530:PRO:HG3	1:A:542:VAL:HG11	1.66	0.76
1:A:378:LYS:O	1:A:379:VAL:HB	1.85	0.75
1:A:281:LYS:HE3	1:A:465:ILE:HD11	1.69	0.75
1:A:552:THR:HB	1:A:555:GLY:O	1.85	0.75
1:A:449:GLU:HG2	1:A:634:GLY:HA2	1.68	0.75
1:A:633:ILE:N	1:A:633:ILE:HD12	2.01	0.74
1:A:348:ILE:HG22	1:A:354:ILE:HG13	1.70	0.72
1:A:633:ILE:CD1	1:A:633:ILE:H	1.98	0.71
1:A:281:LYS:CE	1:A:465:ILE:HD11	2.21	0.71
1:A:385:LYS:HD3	1:A:385:LYS:N	2.04	0.70
1:A:568:ASN:ND2	1:A:569:GLY:H	1.90	0.70
1:A:418:THR:HG22	1:A:420:SER:H	1.57	0.69
1:A:318:LYS:HG2	1:A:546:GLY:HA2	1.75	0.68
1:A:370:LEU:HD13	1:A:507:GLY:HA2	1.76	0.68
1:A:375:ALA:HB3	1:A:378:LYS:NZ	2.10	0.67
1:A:357:THR:OG1	1:A:481:GLN:HG3	1.95	0.67
1:A:477:MET:HB3	1:A:510:HIS:CD2	2.29	0.66
1:A:418:THR:HG22	1:A:420:SER:N	2.09	0.66
1:A:552:THR:CG2	1:A:554:ILE:H	2.08	0.65
1:A:379:VAL:HG22	1:A:380:ASP:H	1.61	0.65
1:A:476:ALA:HB1	1:A:479:ASP:HB2	1.78	0.65
1:A:611:PHE:HB3	1:A:633:ILE:CG1	2.22	0.65
1:A:552:THR:HG22	1:A:554:ILE:N	2.08	0.64
1:A:591:GLN:HE21	1:A:640:LEU:HD13	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:618:THR:O	1:A:618:THR:HG22	1.98	0.63
1:A:398:PRO:HG2	1:A:695:ILE:HG21	1.81	0.63
1:A:568:ASN:HD22	1:A:569:GLY:N	1.96	0.62
1:A:330:ILE:HD12	1:A:368:TRP:NE1	2.14	0.62
1:A:336:LEU:CB	1:A:365:GLN:HG2	2.29	0.62
2:B:15:ALA:O	2:B:16:GLN:HB2	1.99	0.61
1:A:510:HIS:O	1:A:511:LEU:HB2	2.00	0.61
1:A:375:ALA:HB3	1:A:378:LYS:HZ1	1.66	0.61
1:A:357:THR:O	1:A:358:LYS:HB3	2.01	0.61
1:A:617:GLU:HG3	1:A:631:GLY:HA2	1.83	0.61
1:A:336:LEU:HB2	1:A:365:GLN:HG2	1.81	0.60
1:A:670:GLN:HE22	2:B:14:PRO:HA	1.67	0.59
1:A:618:THR:HA	1:A:630:GLU:OE1	2.02	0.59
1:A:465:ILE:HD13	1:A:504:ILE:HD12	1.83	0.59
1:A:689:ALA:O	1:A:692:GLU:HG2	2.02	0.59
1:A:337:ALA:O	1:A:341:VAL:HG23	2.03	0.58
1:A:330:ILE:O	1:A:333:ASN:HB2	2.04	0.58
1:A:373:ASN:H	1:A:373:ASN:ND2	2.01	0.57
1:A:490:ARG:HE	1:A:496:ARG:NH1	2.03	0.57
1:A:592:ILE:CD1	1:A:636:VAL:HG12	2.34	0.56
1:A:575:MET:O	1:A:579:MET:HG3	2.04	0.56
1:A:513:ASN:HB2	1:A:516:ASP:HB2	1.86	0.56
1:A:616:SER:C	1:A:618:THR:N	2.59	0.56
1:A:326:HIS:HD2	1:A:334:TYR:OH	1.89	0.55
1:A:584:CYS:O	1:A:588:SER:HB2	2.06	0.55
1:A:331:ASN:HA	1:A:334:TYR:HD2	1.72	0.55
1:A:632:SER:HB2	1:A:633:ILE:HD12	1.87	0.55
1:A:576:LYS:HA	1:A:576:LYS:HE3	1.89	0.55
1:A:357:THR:HG23	1:A:481:GLN:HE21	1.71	0.54
1:A:616:SER:C	1:A:618:THR:H	2.11	0.54
1:A:322:ILE:HD13	1:A:331:ASN:HB3	1.87	0.54
1:A:490:ARG:HH21	1:A:496:ARG:NH2	2.06	0.54
1:A:580:GLU:O	1:A:583:ARG:HG2	2.08	0.53
1:A:258:GLU:HB3	1:A:260:PHE:CE1	2.43	0.53
1:A:513:ASN:N	1:A:513:ASN:ND2	2.55	0.53
1:A:617:GLU:HG3	1:A:631:GLY:CA	2.38	0.53
1:A:359:ASN:H	1:A:359:ASN:ND2	2.07	0.53
1:A:663:ARG:HH12	2:B:16:GLN:HA	1.74	0.52
1:A:510:HIS:O	1:A:511:LEU:CB	2.57	0.52
1:A:667:LEU:HD23	2:B:15:ALA:HB3	1.92	0.52
1:A:344:GLU:O	1:A:348:ILE:HG13	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:THR:HB	1:A:452:HIS:O	2.09	0.51
1:A:591:GLN:O	1:A:595:MET:HB2	2.11	0.51
1:A:615:LYS:O	1:A:616:SER:HB2	2.10	0.51
2:B:15:ALA:O	2:B:16:GLN:CB	2.58	0.51
1:A:259:PRO:O	1:A:261:LEU:N	2.38	0.50
1:A:281:LYS:HE2	1:A:465:ILE:HD11	1.94	0.50
1:A:697:ASP:O	1:A:701:LEU:HG	2.12	0.49
1:A:620:PRO:HB3	1:A:628:MET:SD	2.53	0.49
1:A:477:MET:HA	1:A:509:SER:OG	2.13	0.49
1:A:264:THR:HB	1:A:265:PRO:HD3	1.95	0.48
1:A:385:LYS:O	1:A:385:LYS:HG2	2.13	0.48
1:A:616:SER:O	1:A:618:THR:N	2.40	0.48
1:A:340:GLN:HG2	1:A:360:MET:CE	2.43	0.47
1:A:331:ASN:HA	1:A:334:TYR:CD2	2.49	0.47
1:A:314:PHE:HD2	1:A:317:TRP:CE2	2.34	0.46
1:A:520:PHE:CZ	1:A:568:ASN:HB3	2.50	0.46
1:A:465:ILE:O	1:A:465:ILE:HD12	2.15	0.46
1:A:592:ILE:HD11	1:A:636:VAL:CG1	2.39	0.46
1:A:362:LYS:O	1:A:367:LYS:HE3	2.16	0.46
1:A:609:LYS:HE3	1:A:609:LYS:HB2	1.80	0.45
1:A:517:VAL:HA	1:A:570:THR:O	2.16	0.45
1:A:418:THR:CG2	1:A:419:ASP:N	2.80	0.45
1:A:271:PRO:HG3	1:A:694:LEU:O	2.16	0.45
1:A:592:ILE:HG12	1:A:637:CYS:SG	2.56	0.45
1:A:351:GLU:HB3	1:A:352:GLU:H	1.59	0.45
1:A:324:LYS:C	1:A:324:LYS:HD2	2.36	0.45
1:A:278:GLN:NE2	1:A:278:GLN:HA	2.32	0.45
1:A:577:TRP:HA	1:A:577:TRP:CE3	2.52	0.45
1:A:426:ASP:HB2	1:A:427:GLU:H	1.59	0.44
1:A:702:LEU:HD13	1:A:706:TRP:CZ2	2.53	0.44
1:A:513:ASN:N	1:A:513:ASN:HD22	2.15	0.44
1:A:316:GLY:HA3	1:A:547:ASP:O	2.18	0.44
1:A:385:LYS:H	1:A:385:LYS:CD	2.18	0.44
1:A:259:PRO:HB3	1:A:675:ASN:C	2.37	0.44
1:A:665:LEU:O	1:A:669:VAL:HG23	2.18	0.44
1:A:359:ASN:ND2	1:A:359:ASN:N	2.66	0.43
1:A:473:SER:HA	1:A:506:LYS:HD2	2.00	0.43
1:A:378:LYS:O	1:A:379:VAL:CB	2.59	0.43
1:A:462:GLY:O	1:A:466:ASN:HB2	2.18	0.43
1:A:674:ASP:O	1:A:675:ASN:HB2	2.19	0.43
1:A:670:GLN:NE2	2:B:14:PRO:HA	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:ARG:O	1:A:363:THR:HG23	2.19	0.43
1:A:615:LYS:O	1:A:616:SER:CB	2.66	0.42
1:A:357:THR:OG1	1:A:481:GLN:CG	2.65	0.42
1:A:258:GLU:HA	1:A:259:PRO:HD3	1.63	0.42
1:A:481:GLN:HB3	1:A:483:ILE:HD11	2.02	0.42
1:A:336:LEU:HB3	1:A:365:GLN:HG2	1.99	0.42
1:A:382:GLU:C	1:A:384:CYS:H	2.23	0.42
1:A:304:LEU:HD11	1:A:358:LYS:HB3	2.01	0.42
1:A:345:LEU:HA	1:A:348:ILE:HD12	2.02	0.42
1:A:348:ILE:H	1:A:348:ILE:HG13	1.64	0.42
1:A:641:LEU:HA	1:A:641:LEU:HD23	1.90	0.41
1:A:375:ALA:HA	1:A:376:PRO:HD3	1.82	0.41
1:A:357:THR:O	1:A:482:LEU:O	2.39	0.41
1:A:356:LYS:HB3	1:A:356:LYS:HE3	1.93	0.41
1:A:270:LEU:HA	1:A:271:PRO:HD3	1.96	0.41
1:A:490:ARG:HE	1:A:496:ARG:CZ	2.33	0.41
1:A:573:ILE:HD13	1:A:573:ILE:HA	1.93	0.41
1:A:398:PRO:O	1:A:399:LYS:HG3	2.21	0.41
1:A:483:ILE:HD12	1:A:483:ILE:N	2.36	0.40
1:A:330:ILE:HD12	1:A:368:TRP:CD1	2.56	0.40
1:A:259:PRO:HB3	1:A:675:ASN:O	2.20	0.40
1:A:597:GLU:C	1:A:599:GLU:H	2.22	0.40
1:A:529:ASP:OD1	1:A:530:PRO:HD2	2.22	0.40
1:A:373:ASN:HA	1:A:378:LYS:HE3	2.03	0.40
1:A:577:TRP:HE3	1:A:577:TRP:HA	1.85	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	427/471 (91%)	375 (88%)	41 (10%)	11 (3%)	7 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	14/30 (47%)	12 (86%)	2 (14%)	0	100 100
All	All	441/501 (88%)	387 (88%)	43 (10%)	11 (2%)	7 27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	PHE
1	A	379	VAL
1	A	387	VAL
1	A	511	LEU
1	A	616	SER
1	A	357	THR
1	A	617	GLU
1	A	262	LYS
1	A	399	LYS
1	A	514	ASP
1	A	614	ASN

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	391/420 (93%)	360 (92%)	31 (8%)	15 41
2	B	15/26 (58%)	15 (100%)	0	100 100
All	All	406/446 (91%)	375 (92%)	31 (8%)	16 43

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	321	ASN
1	A	324	LYS
1	A	351	GLU
1	A	359	ASN
1	A	373	ASN

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Mol	Chain	Res	Type
1	A	385	LYS
1	A	386	ASP
1	A	418	THR
1	A	426	ASP
1	A	441	MET
1	A	450	VAL
1	A	465	ILE
1	A	513	ASN
1	A	515	THR
1	A	517	VAL
1	A	521	VAL
1	A	527	LEU
1	A	547	ASP
1	A	549	LEU
1	A	552	THR
1	A	568	ASN
1	A	570	THR
1	A	576	LYS
1	A	592	ILE
1	A	610	GLU
1	A	619	TRP
1	A	633	ILE
1	A	641	LEU
1	A	649	LEU
1	A	683	LEU
1	A	702	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	GLN
1	A	326	HIS
1	A	373	ASN
1	A	510	HIS
1	A	519	ASN
1	A	568	ASN
1	A	591	GLN
2	B	16	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	435/471 (92%)	0.70	48 (11%) 7 4	53, 77, 136, 160	0
2	B	16/30 (53%)	0.43	0 100 100	65, 74, 103, 109	0
All	All	451/501 (90%)	0.69	48 (10%) 8 5	53, 77, 136, 160	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	428	ILE	4.9
1	A	445	TYR	4.8
1	A	512	ARG	4.5
1	A	617	GLU	4.2
1	A	603	LYS	4.1
1	A	382	GLU	4.1
1	A	293	GLU	3.8
1	A	376	PRO	3.7
1	A	357	THR	3.5
1	A	379	VAL	3.4
1	A	618	THR	3.3
1	A	611	PHE	3.3
1	A	598	ALA	3.3
1	A	596	ILE	3.3
1	A	607	MET	3.2
1	A	628	MET	3.2
1	A	623	GLU	3.2
1	A	397	GLU	3.2
1	A	387	VAL	3.1
1	A	695	ILE	3.1
1	A	608	THR	3.0
1	A	425	LEU	3.0
1	A	605	LYS	2.8
1	A	354	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	602	VAL	2.7
1	A	577	TRP	2.6
1	A	600	SER	2.6
1	A	381	PHE	2.5
1	A	511	LEU	2.5
1	A	261	LEU	2.5
1	A	612	PHE	2.5
1	A	427	GLU	2.4
1	A	495	ARG	2.4
1	A	441	MET	2.3
1	A	328	LYS	2.3
1	A	388	SER	2.3
1	A	305	TYR	2.2
1	A	385	LYS	2.2
1	A	694	LEU	2.2
1	A	377	GLU	2.2
1	A	604	GLU	2.2
1	A	609	LYS	2.1
1	A	614	ASN	2.1
1	A	353	LYS	2.0
1	A	400	PRO	2.0
1	A	610	GLU	2.0
1	A	464	TYR	2.0
1	A	352	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.