Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:07 AM GMT

PDB ID	:	2ZNL
Title	:	Crystal structure of PA-PB1 complex form influenza virus RNA polymerase
Authors	:	Obayashi, E.; Yoshida, H.; Kawai, F.; Shibayama, N.; Kawaguchi, A.; Nagata, K.; Tame, J.R.H.; Park, S.-Y.
Deposited on	:	2008-04-28
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
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:

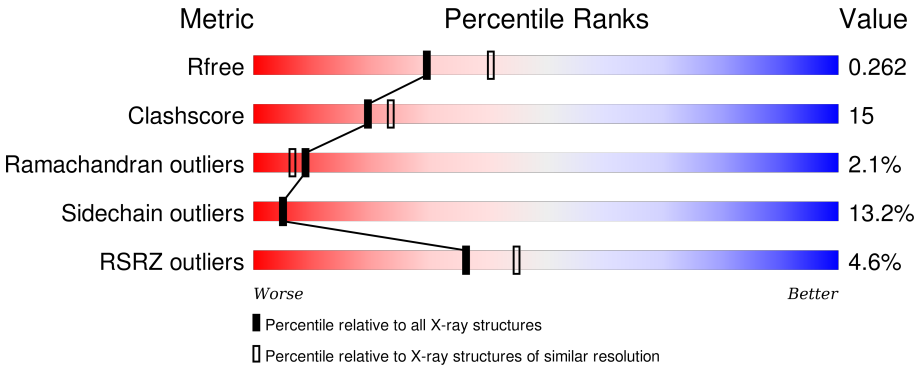
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 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	478	4% 52% 28% 7% • 12%
2	B	81	12% 5% • 81%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3569 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	421	Total	C	N	O	S	0	0	0
			3382	2157	571	629	25			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	15	Total	C	N	O	S	0	0	0
			117	78	18	20	1			

- Molecule 3 is water.

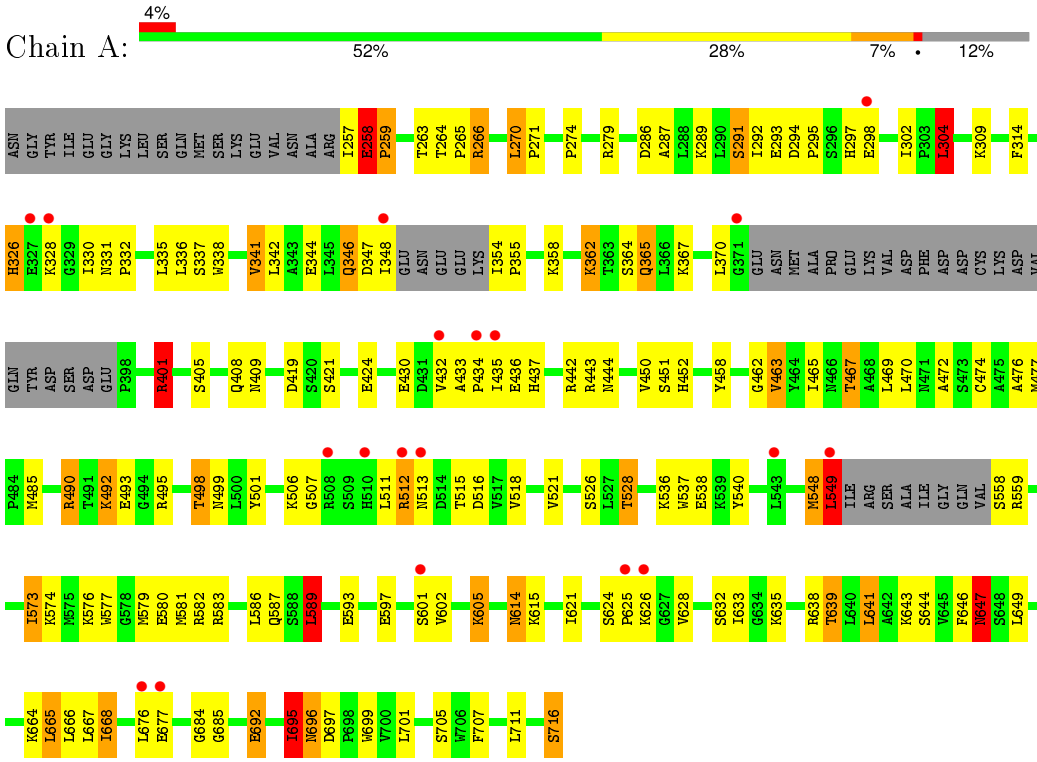
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total	O	0	0
			70	70		

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.

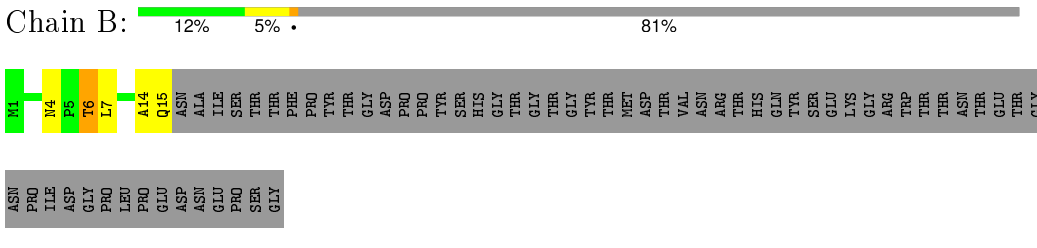
- Molecule 1: Polymerase acidic protein

Chain A:



- Molecule 2: RNA-directed RNA polymerase catalytic subunit

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	101.96Å 101.96Å 115.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.95 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.30) 98.5 (19.95-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.62 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.207 , 0.262 0.206 , 0.262	Depositor DCC
R_{free} test set	1545 reflections (5.30%)	DCC
Wilson B-factor (Å^2)	44.8	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 50.5	EDS
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 30669 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3569	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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	1.46	25/3456 (0.7%)	1.32	34/4661 (0.7%)
2	B	1.45	0/119	1.59	1/162 (0.6%)
All	All	1.45	25/3575 (0.7%)	1.33	35/4823 (0.7%)

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	2
2	B	0	1
All	All	0	3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	705	SER	CB-OG	-9.07	1.30	1.42
1	A	647	ASN	CB-CG	-8.53	1.31	1.51
1	A	450	VAL	CB-CG2	7.88	1.69	1.52
1	A	632	SER	CB-OG	7.87	1.52	1.42
1	A	615	LYS	N-CA	7.23	1.60	1.46
1	A	696	ASN	CB-CG	-7.00	1.34	1.51
1	A	458	TYR	CD1-CE1	6.92	1.49	1.39
1	A	638	ARG	CG-CD	6.34	1.67	1.51
1	A	474	CYS	CB-SG	-6.27	1.71	1.82
1	A	647	ASN	N-CA	-6.23	1.33	1.46
1	A	298	GLU	CB-CG	6.18	1.63	1.52
1	A	298	GLU	CG-CD	5.91	1.60	1.51
1	A	424	GLU	CG-CD	5.89	1.60	1.51
1	A	654	GLN	CG-CD	5.89	1.64	1.51
1	A	583	ARG	CG-CD	5.88	1.66	1.51
1	A	501	TYR	CE2-CZ	-5.75	1.31	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	492	LYS	CD-CE	5.69	1.65	1.51
1	A	309	LYS	CD-CE	5.68	1.65	1.51
1	A	408	GLN	CD-OE1	5.40	1.35	1.24
1	A	287	ALA	CA-CB	-5.34	1.41	1.52
1	A	451	SER	CB-OG	-5.32	1.35	1.42
1	A	424	GLU	CB-CG	5.18	1.61	1.52
1	A	540	TYR	CE2-CZ	5.17	1.45	1.38
1	A	463	VAL	CB-CG2	-5.16	1.42	1.52
1	A	430	GLU	CG-CD	5.07	1.59	1.51

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	266	ARG	NE-CZ-NH1	-9.92	115.34	120.30
1	A	614	ASN	C-N-CA	-9.38	98.24	121.70
1	A	647	ASN	CB-CA-C	9.10	128.60	110.40
2	B	7	LEU	CB-CG-CD1	-8.64	96.32	111.00
1	A	419	ASP	CB-CG-OD2	-7.79	111.29	118.30
1	A	685	GLY	N-CA-C	-7.02	95.54	113.10
1	A	638	ARG	NE-CZ-NH2	-6.94	116.83	120.30
1	A	566	ARG	NE-CZ-NH2	-6.92	116.84	120.30
1	A	419	ASP	CB-CG-OD1	6.87	124.48	118.30
1	A	646	PHE	C-N-CA	-6.71	104.91	121.70
1	A	292	ILE	CB-CA-C	-6.55	98.51	111.60
1	A	638	ARG	NE-CZ-NH1	6.43	123.52	120.30
1	A	304	LEU	CB-CG-CD1	-6.42	100.08	111.00
1	A	583	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	583	ARG	CG-CD-NE	6.05	124.50	111.80
1	A	667	LEU	CB-CG-CD1	-5.95	100.88	111.00
1	A	614	ASN	O-C-N	-5.83	113.37	122.70
1	A	270	LEU	CA-CB-CG	5.78	128.59	115.30
1	A	458	TYR	CB-CG-CD2	-5.64	117.62	121.00
1	A	614	ASN	N-CA-C	-5.63	95.80	111.00
1	A	701	LEU	CA-CB-CG	5.46	127.85	115.30
1	A	401	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	A	490	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	A	643	LYS	CD-CE-NZ	-5.24	99.65	111.70
1	A	676	LEU	CB-CA-C	5.21	120.09	110.20
1	A	605	LYS	CD-CE-NZ	5.17	123.58	111.70
1	A	582	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	665	LEU	CB-CG-CD2	5.15	119.76	111.00
1	A	430	GLU	CA-CB-CG	5.14	124.72	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	492	LYS	CD-CE-NZ	5.14	123.53	111.70
1	A	549	LEU	CA-CB-CG	-5.12	103.52	115.30
1	A	270	LEU	CB-CG-CD2	5.08	119.63	111.00
1	A	589	LEU	CB-CG-CD2	5.07	119.62	111.00
1	A	716	SER	CB-CA-C	5.07	119.73	110.10
1	A	641	LEU	CB-CG-CD2	5.02	119.54	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	258	GLU	Peptide
1	A	695	ILE	Peptide
2	B	14	ALA	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	3382	0	3392	102	0
2	B	117	0	128	3	0
3	A	70	0	0	5	0
All	All	3569	0	3520	105	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 (105) 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:476:ALA:HB1	1:A:479:ASP:HB2	1.27	1.13
1:A:462:GLY:HA3	1:A:581:MET:CE	1.79	1.12
1:A:463:VAL:O	1:A:467:THR:HG23	1.65	0.96
1:A:467:THR:HG22	1:A:579:MET:HG2	1.51	0.92
1:A:462:GLY:HA3	1:A:581:MET:HE3	1.53	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:HIS:H	1:A:331:ASN:HD21	1.20	0.90
1:A:647:ASN:HD22	1:A:647:ASN:C	1.74	0.89
1:A:576:LYS:HE2	1:A:580:GLU:OE1	1.74	0.87
1:A:647:ASN:HB2	1:A:655:LEU:HD11	1.55	0.87
1:A:506:LYS:HD2	1:A:507:GLY:O	1.77	0.83
1:A:512:ARG:HH11	1:A:512:ARG:HB2	1.42	0.83
1:A:289:LYS:HG2	1:A:499:ASN:HB3	1.64	0.80
1:A:442:ARG:HH12	1:A:597:GLU:HG2	1.47	0.80
1:A:472:ALA:CB	1:A:483:ILE:HD11	2.11	0.79
1:A:476:ALA:CB	1:A:479:ASP:HB2	2.11	0.77
1:A:365:GLN:NE2	1:A:365:GLN:H	1.83	0.77
1:A:344:GLU:HB3	1:A:358:LYS:HE3	1.65	0.76
1:A:326:HIS:N	1:A:331:ASN:HD21	1.83	0.76
1:A:336:LEU:CB	1:A:365:GLN:HG3	2.15	0.76
1:A:286:ASP:OD1	1:A:528:THR:HG21	1.89	0.72
1:A:326:HIS:H	1:A:331:ASN:ND2	1.90	0.70
3:A:4:HOH:O	2:B:6:THR:HG21	1.91	0.70
1:A:324:LYS:HG3	1:A:324:LYS:O	1.91	0.69
1:A:635:LYS:O	1:A:639:THR:HG23	1.93	0.68
1:A:279:ARG:NH1	3:A:64:HOH:O	2.26	0.68
1:A:647:ASN:CB	1:A:699:TRP:HE1	2.06	0.67
2:B:15:GLN:HA	2:B:15:GLN:OE1	1.93	0.67
1:A:479:ASP:O	1:A:506:LYS:HE2	1.95	0.67
1:A:472:ALA:HB3	1:A:483:ILE:HD11	1.77	0.66
1:A:336:LEU:HB3	1:A:365:GLN:HG3	1.75	0.66
1:A:647:ASN:HB3	1:A:699:TRP:HE1	1.60	0.65
1:A:443:ARG:HH21	1:A:444:ASN:HD21	1.42	0.65
1:A:442:ARG:NH1	1:A:597:GLU:HG2	2.11	0.64
1:A:336:LEU:HB2	1:A:365:GLN:HG3	1.79	0.64
1:A:434:PRO:O	1:A:435:ILE:HG13	1.98	0.64
1:A:295:PRO:HG3	1:A:498:THR:HG23	1.83	0.60
1:A:528:THR:HG22	3:A:49:HOH:O	2.01	0.59
1:A:330:ILE:HG22	1:A:332:PRO:HD2	1.83	0.59
1:A:695:ILE:HD13	1:A:695:ILE:N	2.18	0.58
1:A:344:GLU:HB3	1:A:358:LYS:CE	2.35	0.57
1:A:635:LYS:O	1:A:639:THR:CG2	2.53	0.57
1:A:574:LYS:O	1:A:577:TRP:HB3	2.05	0.57
1:A:462:GLY:HA3	1:A:581:MET:HE1	1.77	0.57
1:A:258:GLU:OE2	1:A:258:GLU:N	2.38	0.56
1:A:346:GLN:C	1:A:347:ASP:OD1	2.44	0.56
1:A:348:ILE:O	1:A:348:ILE:HG22	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LYS:CG	1:A:499:ASN:HB3	2.36	0.55
1:A:342:LEU:O	1:A:346:GLN:OE1	2.25	0.55
1:A:436:GLU:HG2	1:A:437:HIS:N	2.22	0.55
1:A:512:ARG:NH1	1:A:512:ARG:HB2	2.18	0.54
1:A:589:LEU:HD22	1:A:593:GLU:HG3	1.89	0.54
1:A:647:ASN:HB3	1:A:699:TRP:NE1	2.23	0.54
1:A:291:SER:HB3	1:A:498:THR:O	2.07	0.54
1:A:337:SER:O	1:A:341:VAL:HG12	2.09	0.53
1:A:587:GLN:NE2	1:A:644:SER:OG	2.42	0.53
1:A:257:ILE:C	1:A:259:PRO:HD2	2.29	0.53
1:A:264:THR:HB	1:A:265:PRO:CD	2.39	0.52
1:A:668:ILE:HD13	1:A:668:ILE:C	2.30	0.52
1:A:576:LYS:HE2	1:A:580:GLU:CD	2.30	0.51
1:A:536:LYS:HD3	1:A:537:TRP:CZ2	2.45	0.51
1:A:692:GLU:O	1:A:695:ILE:HD11	2.11	0.51
1:A:511:LEU:HD21	1:A:518:VAL:HG22	1.93	0.50
1:A:320:PRO:HB2	1:A:335:LEU:HD22	1.93	0.50
1:A:436:GLU:HG2	1:A:437:HIS:H	1.77	0.50
1:A:624:SER:HB2	1:A:625:PRO:HD2	1.93	0.50
1:A:338:TRP:O	1:A:341:VAL:HG13	2.12	0.49
1:A:370:LEU:CD2	1:A:521:VAL:HG23	2.42	0.49
1:A:421:SER:HB3	1:A:452:HIS:CE1	2.47	0.49
1:A:365:GLN:N	1:A:365:GLN:NE2	2.56	0.48
1:A:472:ALA:CB	1:A:483:ILE:CD1	2.90	0.48
1:A:304:LEU:HB2	3:A:36:HOH:O	2.14	0.47
1:A:325:PRO:HB2	1:A:331:ASN:OD1	2.15	0.47
1:A:624:SER:HB2	1:A:625:PRO:CD	2.45	0.47
1:A:370:LEU:HD21	1:A:521:VAL:HG23	1.96	0.47
1:A:654:GLN:HE22	1:A:697:ASP:H	1.63	0.47
1:A:270:LEU:HB3	1:A:271:PRO:HD2	1.97	0.46
1:A:563:LEU:HD13	1:A:565:VAL:HG12	1.98	0.45
1:A:490:ARG:HA	1:A:495:ARG:O	2.16	0.45
1:A:472:ALA:HB3	1:A:483:ILE:CD1	2.46	0.45
1:A:257:ILE:C	1:A:258:GLU:OE2	2.56	0.45
1:A:549:LEU:CD2	1:A:558:SER:HB2	2.47	0.45
1:A:409:ASN:ND2	3:A:58:HOH:O	2.50	0.44
1:A:293:GLU:O	1:A:294:ASP:C	2.55	0.44
1:A:465:ILE:HD13	1:A:485:MET:HE3	1.99	0.44
1:A:654:GLN:HE22	1:A:696:ASN:H	1.65	0.44
1:A:549:LEU:HA	1:A:549:LEU:HD23	1.46	0.44
1:A:362:LYS:O	1:A:367:LYS:HE3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:614:ASN:N	1:A:614:ASN:HD22	2.16	0.43
1:A:462:GLY:CA	1:A:581:MET:HE3	2.36	0.43
1:A:274:PRO:O	1:A:401:ARG:CG	2.67	0.43
1:A:526:SER:OG	1:A:528:THR:HG23	2.19	0.43
1:A:462:GLY:HA3	1:A:581:MET:HE2	1.88	0.42
1:A:647:ASN:CB	1:A:699:TRP:NE1	2.79	0.42
1:A:707:PHE:CE2	1:A:711:LEU:HD21	2.55	0.42
1:A:354:ILE:N	1:A:355:PRO:HD3	2.34	0.42
1:A:257:ILE:HG22	1:A:257:ILE:O	2.19	0.42
1:A:354:ILE:N	1:A:355:PRO:CD	2.83	0.42
1:A:647:ASN:HB2	1:A:699:TRP:HE1	1.82	0.41
1:A:314:PHE:CE2	1:A:548:MET:HG2	2.54	0.41
1:A:513:ASN:HB2	1:A:516:ASP:OD1	2.21	0.41
1:A:476:ALA:O	1:A:477:MET:HG2	2.20	0.41
2:B:4:ASN:C	2:B:4:ASN:OD1	2.59	0.41
1:A:264:THR:HB	1:A:265:PRO:HD2	2.01	0.41
1:A:664:LYS:O	1:A:668:ILE:HG22	2.21	0.40
1:A:573:ILE:HD13	1:A:573:ILE:HA	1.72	0.40

There are no symmetry-related clashes.

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	413/478 (86%)	379 (92%)	25 (6%)	9 (2%)	8 6
2	B	13/81 (16%)	11 (85%)	2 (15%)	0	100 100
All	All	426/559 (76%)	390 (92%)	27 (6%)	9 (2%)	9 7

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	GLU
1	A	297	HIS
1	A	433	ALA
1	A	346	GLN
1	A	684	GLY
1	A	362	LYS
1	A	259	PRO
1	A	621	ILE
1	A	602	VAL

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	379/429 (88%)	328 (86%)	51 (14%)	5 4
2	B	14/70 (20%)	13 (93%)	1 (7%)	18 23
All	All	393/499 (79%)	341 (87%)	52 (13%)	5 5

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	258	GLU
1	A	263	THR
1	A	266	ARG
1	A	291	SER
1	A	302	ILE
1	A	304	LEU
1	A	324	LYS
1	A	326	HIS
1	A	328	LYS
1	A	341	VAL
1	A	364	SER
1	A	365	GLN
1	A	401	ARG
1	A	405	SER
1	A	432	VAL

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Mol	Chain	Res	Type
1	A	467	THR
1	A	469	LEU
1	A	470	LEU
1	A	492	LYS
1	A	493	GLU
1	A	498	THR
1	A	512	ARG
1	A	515	THR
1	A	528	THR
1	A	538	GLU
1	A	548	MET
1	A	549	LEU
1	A	559	ARG
1	A	563	LEU
1	A	565	VAL
1	A	573	ILE
1	A	586	LEU
1	A	589	LEU
1	A	601	SER
1	A	605	LYS
1	A	626	LYS
1	A	628	VAL
1	A	633	ILE
1	A	639	THR
1	A	641	LEU
1	A	647	ASN
1	A	649	LEU
1	A	655	LEU
1	A	659	SER
1	A	665	LEU
1	A	666	LEU
1	A	668	ILE
1	A	677	GLU
1	A	692	GLU
1	A	695	ILE
1	A	716	SER
2	B	6	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	331	ASN

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Mol	Chain	Res	Type
1	A	333	ASN
1	A	340	GLN
1	A	365	GLN
1	A	409	ASN
1	A	444	ASN
1	A	452	HIS
1	A	481	GLN
1	A	587	GLN
1	A	614	ASN
1	A	647	ASN
1	A	703	ASN

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	421/478 (88%)	0.09	20 (4%) 34 43	29, 48, 77, 93	0
2	B	15/81 (18%)	-0.38	0 100 100	32, 39, 64, 73	0
All	All	436/559 (77%)	0.07	20 (4%) 36 45	29, 48, 77, 93	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	434	PRO	8.3
1	A	676	LEU	6.1
1	A	432	VAL	5.0
1	A	512	ARG	4.2
1	A	563	LEU	3.7
1	A	328	LYS	3.6
1	A	348	ILE	3.3
1	A	435	ILE	2.9
1	A	298	GLU	2.7
1	A	513	ASN	2.6
1	A	508	ARG	2.6
1	A	371	GLY	2.5
1	A	327	GLU	2.3
1	A	510	HIS	2.3
1	A	677	GLU	2.2
1	A	549	LEU	2.1
1	A	626	LYS	2.1
1	A	543	LEU	2.1
1	A	601	SER	2.1
1	A	625	PRO	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.