Full wwPDB X-ray Structure Validation Report (i)

Aug 28, 2023 – 05:08 AM EDT

PDB ID	:	3KN6
Title	:	Crystal structure of the C-terminal kinase domain of MSK1
Authors	:	D'Angelo, I.; Malakhova, M.; Dong, Z.
Deposited on	:	2009-11-12
Resolution	:	2.00 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 (i) 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 (1)) were used in the production of this report:

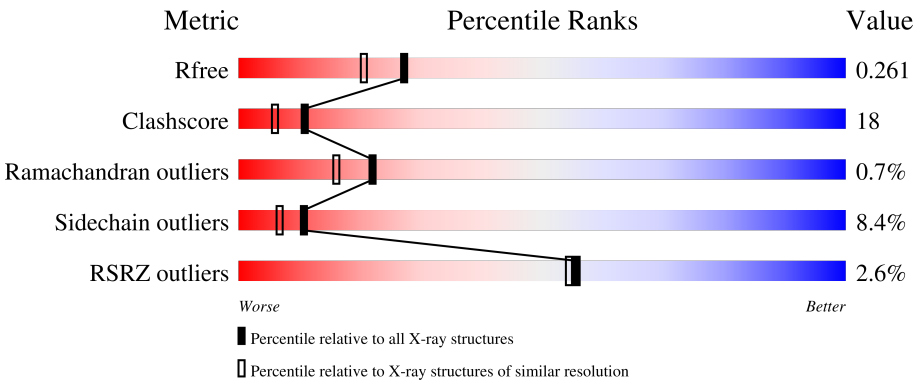
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	325	2% 63% 21% • 13%
1	B	325	3% 58% 24% 6% • 11%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4744 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 Ribosomal protein S6 kinase alpha-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	282	2251	1438	387	412	14	0	0	0
1	B	288	2284	1457	393	420	14	0	0	0

- Molecule 2 is water.

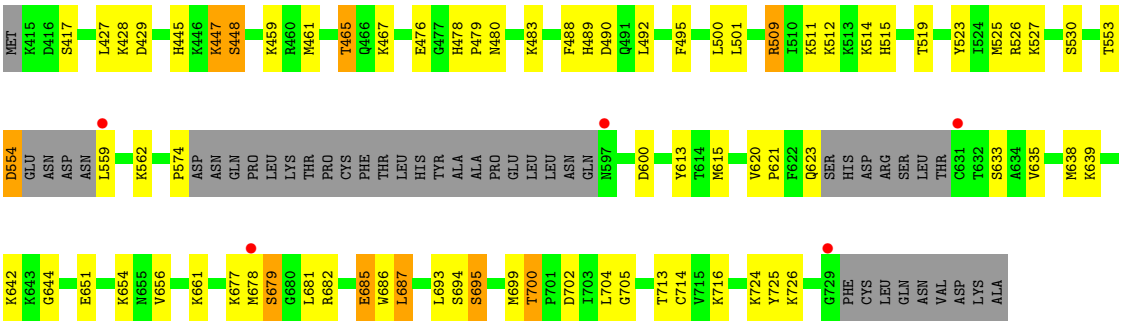
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	102	102	102	0	0
2	B	107	107	107	0	0

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.

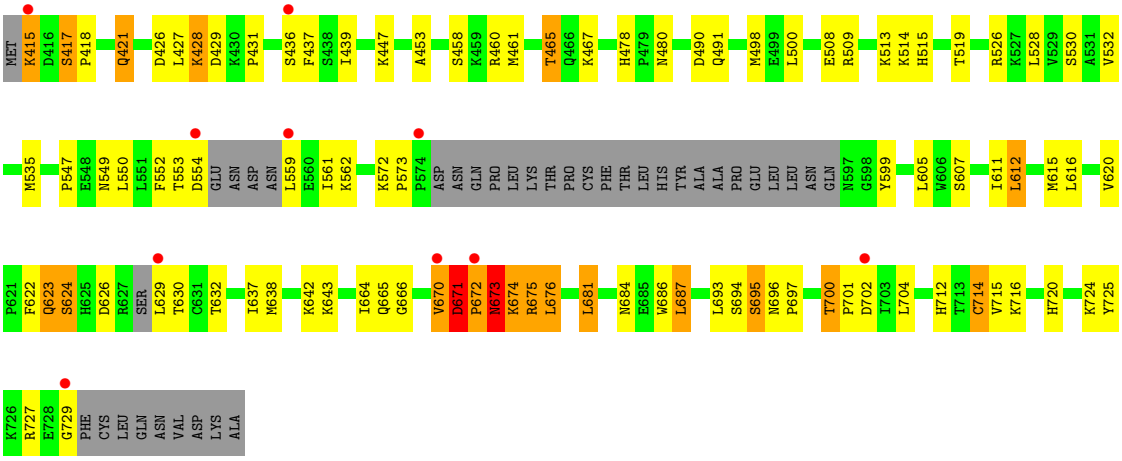
• Molecule 1: Ribosomal protein S6 kinase alpha-5

Chain A:



• Molecule 1: Ribosomal protein S6 kinase alpha-5

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.39Å 91.74Å 134.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.87 – 2.00 45.87 – 2.00	Depositor EDS
% Data completeness (in resolution range)	52.4 (45.87-2.00) 99.2 (45.87-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	3.64 (at 2.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.216 , 0.263 0.211 , 0.261	Depositor DCC
R_{free} test set	2183 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	27.8	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 45.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4744	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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, 0.333 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	0.31	0/2297	0.50	0/3083
1	B	0.37	0/2330	0.53	0/3129
All	All	0.34	0/4627	0.51	0/6212

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	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	671	ASP	Peptide
1	B	673	ASN	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	2251	0	2264	67	0
1	B	2284	0	2283	101	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	102	0	0	17	0
2	B	107	0	0	19	0
All	All	4744	0	4547	165	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 18.

All (165) 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:B:611:ILE:HG12	2:B:163:HOH:O	1.51	1.10
1:B:674:LYS:O	1:B:674:LYS:HG2	1.52	1.03
1:A:501:LEU:HD22	1:A:553:THR:HG22	1.45	0.98
1:B:426:ASP:OD2	1:B:428:LYS:HG3	1.68	0.93
1:A:445:HIS:ND1	1:A:448:SER:HB3	1.87	0.90
1:B:671:ASP:HB3	1:B:672:PRO:C	1.95	0.86
1:B:671:ASP:OD2	1:B:673:ASN:HA	1.75	0.85
1:B:671:ASP:C	1:B:673:ASN:N	2.30	0.84
1:B:671:ASP:HB3	1:B:672:PRO:O	1.77	0.84
1:A:478:HIS:HD2	1:A:480:ASN:H	1.23	0.83
1:B:491:GLN:HB3	2:B:151:HOH:O	1.79	0.82
1:A:519:THR:HG23	1:A:694:SER:O	1.79	0.82
1:B:670:VAL:O	1:B:670:VAL:CG2	2.29	0.81
1:B:671:ASP:OD2	1:B:673:ASN:CA	2.30	0.80
1:B:670:VAL:O	1:B:670:VAL:HG22	1.79	0.79
1:A:615:MET:O	1:A:700:THR:HG21	1.82	0.79
1:B:615:MET:O	1:B:700:THR:HG21	1.81	0.78
1:B:672:PRO:HD2	1:B:674:LYS:HD2	1.67	0.77
1:B:417:SER:HB2	1:B:490:ASP:HB3	1.68	0.74
1:A:519:THR:HG22	1:A:693:LEU:HB2	1.70	0.73
1:B:478:HIS:HD2	1:B:480:ASN:H	1.33	0.73
1:A:679:SER:HB3	1:A:682:ARG:HG2	1.71	0.73
1:A:478:HIS:CD2	1:A:480:ASN:H	2.07	0.72
1:B:673:ASN:OD1	1:B:674:LYS:CB	2.37	0.72
1:B:458:SER:OG	1:B:460:ARG:HG2	1.89	0.72
1:B:671:ASP:C	1:B:673:ASN:H	1.92	0.71
1:A:526:ARG:HD2	2:A:48:HOH:O	1.92	0.70
1:B:672:PRO:HG2	1:B:674:LYS:HE2	1.73	0.70
1:A:553:THR:HG21	1:A:562:LYS:NZ	2.07	0.69
1:B:526:ARG:HD3	1:B:687:LEU:O	1.93	0.69
1:B:623:GLN:HG3	1:B:632:THR:HG22	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:465:THR:HG21	2:B:10:HOH:O	1.93	0.68
1:B:638:MET:HE1	1:B:724:LYS:HD3	1.74	0.68
1:A:678:MET:HG3	1:A:679:SER:HB2	1.76	0.68
1:A:526:ARG:HD3	1:A:687:LEU:O	1.94	0.66
1:B:673:ASN:OD1	1:B:673:ASN:C	2.30	0.66
1:B:673:ASN:OD1	1:B:674:LYS:N	2.30	0.65
1:A:553:THR:HG21	1:A:562:LYS:HZ3	1.62	0.65
1:A:554:ASP:OD2	1:A:554:ASP:N	2.30	0.64
1:B:674:LYS:HB3	1:B:674:LYS:HZ3	1.63	0.64
1:B:478:HIS:CD2	1:B:480:ASN:H	2.17	0.63
1:A:465:THR:HG22	2:A:52:HOH:O	1.97	0.62
1:B:562:LYS:HE2	2:B:178:HOH:O	1.99	0.62
1:B:426:ASP:OD2	1:B:428:LYS:CG	2.44	0.61
1:B:528:LEU:HD21	1:B:561:ILE:HD13	1.81	0.61
1:B:460:ARG:HG3	2:B:172:HOH:O	2.00	0.61
1:A:562:LYS:HE2	2:A:42:HOH:O	2.00	0.61
1:B:674:LYS:CB	1:B:674:LYS:NZ	2.61	0.61
1:B:702:ASP:HB2	2:B:96:HOH:O	2.00	0.61
1:B:415:LYS:HA	2:B:107:HOH:O	1.99	0.60
1:A:702:ASP:HB3	2:A:197:HOH:O	2.00	0.60
1:B:624:SER:HG	1:B:629:LEU:N	2.00	0.60
1:B:638:MET:HG3	1:B:725:TYR:HB2	1.84	0.60
1:B:673:ASN:OD1	1:B:674:LYS:CA	2.50	0.60
1:A:427:LEU:HA	2:A:59:HOH:O	2.02	0.59
1:B:727:ARG:O	1:B:729:GLY:HA3	2.02	0.59
1:A:644:GLY:HA2	2:A:87:HOH:O	2.01	0.59
1:B:626:ASP:O	1:B:629:LEU:N	2.35	0.59
1:B:720:HIS:HB2	2:B:198:HOH:O	2.02	0.59
1:B:532:VAL:HA	1:B:535:MET:HE3	1.85	0.59
1:B:519:THR:HG22	1:B:693:LEU:HB2	1.84	0.58
1:B:671:ASP:O	1:B:673:ASN:N	2.30	0.58
1:A:479:PRO:O	1:A:562:LYS:HE3	2.04	0.57
1:A:638:MET:O	1:A:642:LYS:HG2	2.04	0.56
1:A:511:LYS:HD2	1:A:705:GLY:HA2	1.88	0.56
1:A:623:GLN:HB2	2:A:196:HOH:O	2.05	0.55
2:A:105:HOH:O	1:B:695:SER:HB2	2.06	0.55
1:A:685:GLU:HG2	1:A:686:TRP:N	2.21	0.55
1:A:638:MET:HE1	1:A:724:LYS:NZ	2.22	0.55
1:A:489:HIS:HE1	2:A:17:HOH:O	1.90	0.54
1:A:417:SER:HB2	1:A:490:ASP:HB3	1.89	0.54
1:A:429:ASP:HB2	2:A:53:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:VAL:HG21	1:A:714:CYS:HB3	1.89	0.54
1:B:674:LYS:O	1:B:674:LYS:CG	2.41	0.54
1:A:677:LYS:O	1:A:679:SER:HA	2.07	0.53
1:A:465:THR:HG21	2:A:6:HOH:O	2.08	0.53
1:A:639:LYS:HA	1:A:642:LYS:HE3	1.90	0.53
1:A:461:MET:HE2	2:A:184:HOH:O	2.08	0.53
1:B:418:PRO:HA	1:B:421:GLN:HG3	1.90	0.53
1:B:675:ARG:HD2	2:B:190:HOH:O	2.09	0.53
1:A:478:HIS:CD2	1:A:479:PRO:HD2	2.44	0.53
1:B:632:THR:HG21	2:B:55:HOH:O	2.08	0.53
1:B:508:GLU:HG3	2:B:23:HOH:O	2.09	0.52
1:B:553:THR:CG2	1:B:562:LYS:NZ	2.72	0.52
1:B:622:PHE:C	1:B:623:GLN:HE21	2.12	0.52
1:B:612:LEU:HD13	1:B:664:ILE:HG13	1.91	0.52
1:B:712:HIS:NE2	1:B:716:LYS:HE3	2.25	0.52
1:A:651:GLU:HG3	1:A:699:MET:SD	2.50	0.51
1:B:643:LYS:O	1:B:643:LYS:HG2	2.08	0.51
1:A:514:LYS:HG3	1:B:696:ASN:HD21	1.76	0.51
1:B:431:PRO:HG3	1:B:439:ILE:HD11	1.93	0.51
1:B:478:HIS:HE1	1:B:530:SER:OG	1.94	0.51
1:A:613:TYR:CD1	1:A:621:PRO:HD3	2.46	0.51
1:A:678:MET:HA	1:A:679:SER:CB	2.38	0.51
1:A:574:PRO:HD2	2:A:153:HOH:O	2.10	0.51
1:B:498:MET:HB2	2:B:169:HOH:O	2.10	0.50
1:B:671:ASP:OD2	1:B:673:ASN:C	2.49	0.50
1:A:515:HIS:HE1	1:A:695:SER:O	1.95	0.50
1:B:514:LYS:HB2	2:B:168:HOH:O	2.12	0.50
1:B:519:THR:HG23	1:B:694:SER:O	2.12	0.50
1:B:461:MET:O	1:B:465:THR:HG23	2.12	0.50
1:B:465:THR:HG22	2:B:31:HOH:O	2.12	0.50
1:B:623:GLN:OE1	1:B:714:CYS:SG	2.70	0.49
1:B:436:SER:HA	2:B:83:HOH:O	2.13	0.49
1:B:572:LYS:HB2	1:B:599:TYR:CZ	2.48	0.49
1:A:514:LYS:HA	1:B:696:ASN:HD21	1.76	0.49
1:B:509:ARG:HG3	1:B:552:PHE:CD2	2.48	0.49
1:B:553:THR:CG2	1:B:562:LYS:HZ2	2.25	0.49
1:B:632:THR:HG23	1:B:637:ILE:HG13	1.95	0.49
1:A:447:LYS:HE3	1:A:447:LYS:HA	1.95	0.48
1:A:559:LEU:HB3	2:A:181:HOH:O	2.12	0.48
1:B:453:ALA:HB3	2:B:169:HOH:O	2.12	0.48
1:A:478:HIS:HE1	1:A:530:SER:OG	1.97	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:547:PRO:HG2	1:B:715:VAL:HG11	1.94	0.48
1:B:727:ARG:C	1:B:729:GLY:HA3	2.34	0.47
1:B:467:LYS:HD2	1:B:467:LYS:C	2.35	0.47
1:A:716:LYS:HG2	2:A:138:HOH:O	2.15	0.47
1:B:461:MET:HE2	2:B:100:HOH:O	2.14	0.47
1:B:550:LEU:HD11	1:B:611:ILE:HD13	1.96	0.47
1:A:525:MET:HE3	1:A:615:MET:CE	2.45	0.47
1:B:553:THR:HG22	1:B:562:LYS:NZ	2.30	0.46
1:A:509:ARG:CZ	1:A:559:LEU:HD11	2.46	0.46
1:B:418:PRO:O	1:B:421:GLN:HG3	2.15	0.46
1:B:559:LEU:HD13	1:B:559:LEU:O	2.15	0.46
1:A:427:LEU:HD12	2:A:59:HOH:O	2.14	0.46
1:A:600:ASP:OD2	1:A:726:LYS:HE3	2.15	0.46
1:A:523:TYR:CE2	1:A:527:LYS:HE2	2.51	0.46
1:B:607:SER:O	1:B:611:ILE:HG13	2.15	0.46
1:B:431:PRO:CG	1:B:439:ILE:HD11	2.45	0.45
1:B:673:ASN:OD1	1:B:674:LYS:HB3	2.16	0.45
1:B:515:HIS:ND1	1:B:697:PRO:HG3	2.32	0.45
1:A:656:VAL:O	1:A:661:LYS:HE3	2.17	0.45
1:A:686:TRP:HE3	1:A:687:LEU:HD13	1.82	0.45
1:B:612:LEU:HD13	1:B:664:ILE:CG1	2.47	0.45
1:B:612:LEU:HD22	1:B:616:LEU:HG	1.97	0.45
1:A:480:ASN:HD21	1:A:527:LYS:HD3	1.82	0.44
1:B:620:VAL:HG21	1:B:714:CYS:HB3	1.99	0.44
1:A:459:LYS:NZ	1:A:489:HIS:HD2	2.15	0.44
1:B:671:ASP:CB	1:B:672:PRO:C	2.79	0.44
1:A:467:LYS:HB3	1:A:467:LYS:HE2	1.82	0.43
1:B:674:LYS:HB3	1:B:674:LYS:NZ	2.24	0.43
1:B:712:HIS:CD2	1:B:716:LYS:HE3	2.54	0.43
1:A:679:SER:HB3	1:A:682:ARG:CG	2.44	0.43
1:B:526:ARG:NE	2:B:173:HOH:O	2.52	0.43
1:A:488:PHE:HB2	1:A:495:PHE:HB2	2.00	0.43
1:A:686:TRP:CE3	1:A:687:LEU:HD13	2.54	0.42
1:B:700:THR:HG23	1:B:701:PRO:HD3	2.00	0.42
1:B:437:PHE:HB3	1:B:458:SER:HB3	2.00	0.42
1:B:513:LYS:HB3	1:B:513:LYS:HE2	1.86	0.42
1:A:654:LYS:HE2	2:A:164:HOH:O	2.20	0.42
1:B:666:GLY:HA3	1:B:676:LEU:HD22	2.01	0.42
1:A:638:MET:HE1	1:A:724:LYS:CB	2.50	0.41
1:B:572:LYS:HA	1:B:573:PRO:HD3	1.82	0.41
1:B:605:LEU:CD2	1:B:681:LEU:HD12	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:684:ASN:ND2	1:B:686:TRP:H	2.19	0.41
1:A:638:MET:HE2	1:A:638:MET:HB2	1.79	0.41
1:A:467:LYS:C	1:A:467:LYS:HD2	2.41	0.41
1:A:428:LYS:HB3	1:A:428:LYS:HE2	1.60	0.40
1:A:553:THR:HG21	1:A:562:LYS:HZ2	1.81	0.40
1:B:429:ASP:HB2	2:B:150:HOH:O	2.21	0.40
1:B:674:LYS:CB	1:B:674:LYS:HZ3	2.24	0.40
1:A:685:GLU:HG2	1:A:686:TRP:H	1.84	0.40
1:B:549:ASN:HD22	1:B:549:ASN:HA	1.72	0.40
1:A:476:GLU:HG3	1:A:483:LYS:HZ3	1.85	0.40
1:A:512:LYS:HG2	1:B:697:PRO:HG2	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	274/325 (84%)	265 (97%)	9 (3%)	0	100 100
1	B	280/325 (86%)	267 (95%)	9 (3%)	4 (1%)	11 5
All	All	554/650 (85%)	532 (96%)	18 (3%)	4 (1%)	22 16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	672	PRO
1	B	673	ASN
1	B	670	VAL
1	B	671	ASP

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	250/290 (86%)	232 (93%)	18 (7%)	14 9
1	B	252/290 (87%)	228 (90%)	24 (10%)	8 5
All	All	502/580 (87%)	460 (92%)	42 (8%)	11 7

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	447	LYS
1	A	448	SER
1	A	465	THR
1	A	492	LEU
1	A	500	LEU
1	A	509	ARG
1	A	554	ASP
1	A	633	SER
1	A	635	VAL
1	A	679	SER
1	A	681	LEU
1	A	685	GLU
1	A	687	LEU
1	A	695	SER
1	A	700	THR
1	A	704	LEU
1	A	713	THR
1	A	725	TYR
1	B	415	LYS
1	B	417	SER
1	B	421	GLN
1	B	427	LEU
1	B	428	LYS
1	B	447	LYS
1	B	465	THR
1	B	500	LEU
1	B	554	ASP

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Mol	Chain	Res	Type
1	B	612	LEU
1	B	623	GLN
1	B	624	SER
1	B	630	THR
1	B	642	LYS
1	B	665	GLN
1	B	674	LYS
1	B	675	ARG
1	B	676	LEU
1	B	681	LEU
1	B	687	LEU
1	B	695	SER
1	B	700	THR
1	B	704	LEU
1	B	714	CYS

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	478	HIS
1	A	485	HIS
1	A	489	HIS
1	A	515	HIS
1	A	549	ASN
1	A	655	ASN
1	A	684	ASN
1	A	692	GLN
1	A	720	HIS
1	A	723	ASN
1	B	478	HIS
1	B	489	HIS
1	B	493	HIS
1	B	549	ASN
1	B	623	GLN
1	B	655	ASN
1	B	684	ASN
1	B	696	ASN
1	B	723	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 monosaccharides 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	282/325 (86%)	0.20	5 (1%)	68	66	17, 30, 52, 68	0
1	B	288/325 (88%)	0.17	10 (3%)	44	43	18, 30, 56, 68	0
All	All	570/650 (87%)	0.19	15 (2%)	56	54	17, 30, 55, 68	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	436	SER	4.2
1	B	729	GLY	3.8
1	A	597	ASN	3.7
1	A	631	CYS	3.7
1	B	559	LEU	3.0
1	B	670	VAL	2.9
1	B	415	LYS	2.8
1	B	702	ASP	2.6
1	B	574	PRO	2.5
1	B	629	LEU	2.4
1	A	559	LEU	2.4
1	A	729	GLY	2.3
1	B	554	ASP	2.3
1	B	672	PRO	2.3
1	A	678	MET	2.0

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 monosaccharides 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.