



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 10:26 am BST

PDB ID : 6Q83  
Title : Crystal structure of the biportin Pdr6 in complex with UBC9  
Authors : Aksu, M.; Trakhanov, S.; Vera-Rodriguez, A.; Gorlich, D.  
Deposited on : 2018-12-14  
Resolution : 4.53 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

<https://www.wwpdb.org/validation/2017/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) : 1.13  
EDS : 2.11  
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) : 2.11

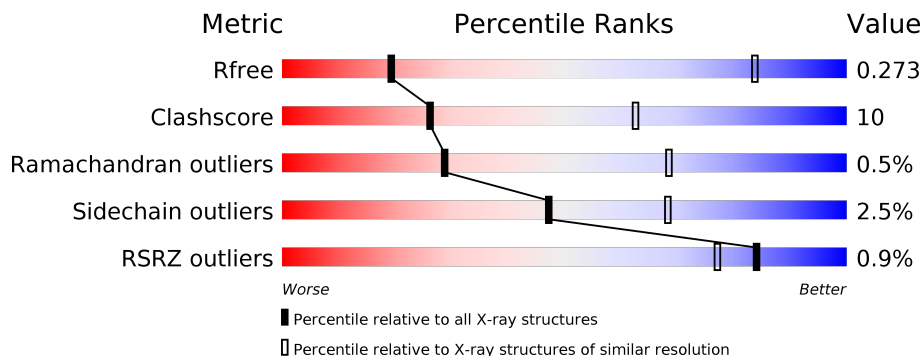
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.53 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	1058 (5.28-3.80)
Clashscore	141614	1126 (5.28-3.80)
Ramachandran outliers	138981	1071 (5.26-3.80)
Sidechain outliers	138945	1052 (5.26-3.80)
RSRZ outliers	127900	1101 (5.30-3.70)

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 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	1080	
2	B	157	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9380 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 Importin beta-like protein KAP122.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1012	8122	5238	1303	1536	45	0	0	0

- Molecule 2 is a protein called UBC9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	156	1258	810	216	229	3	0	0	0

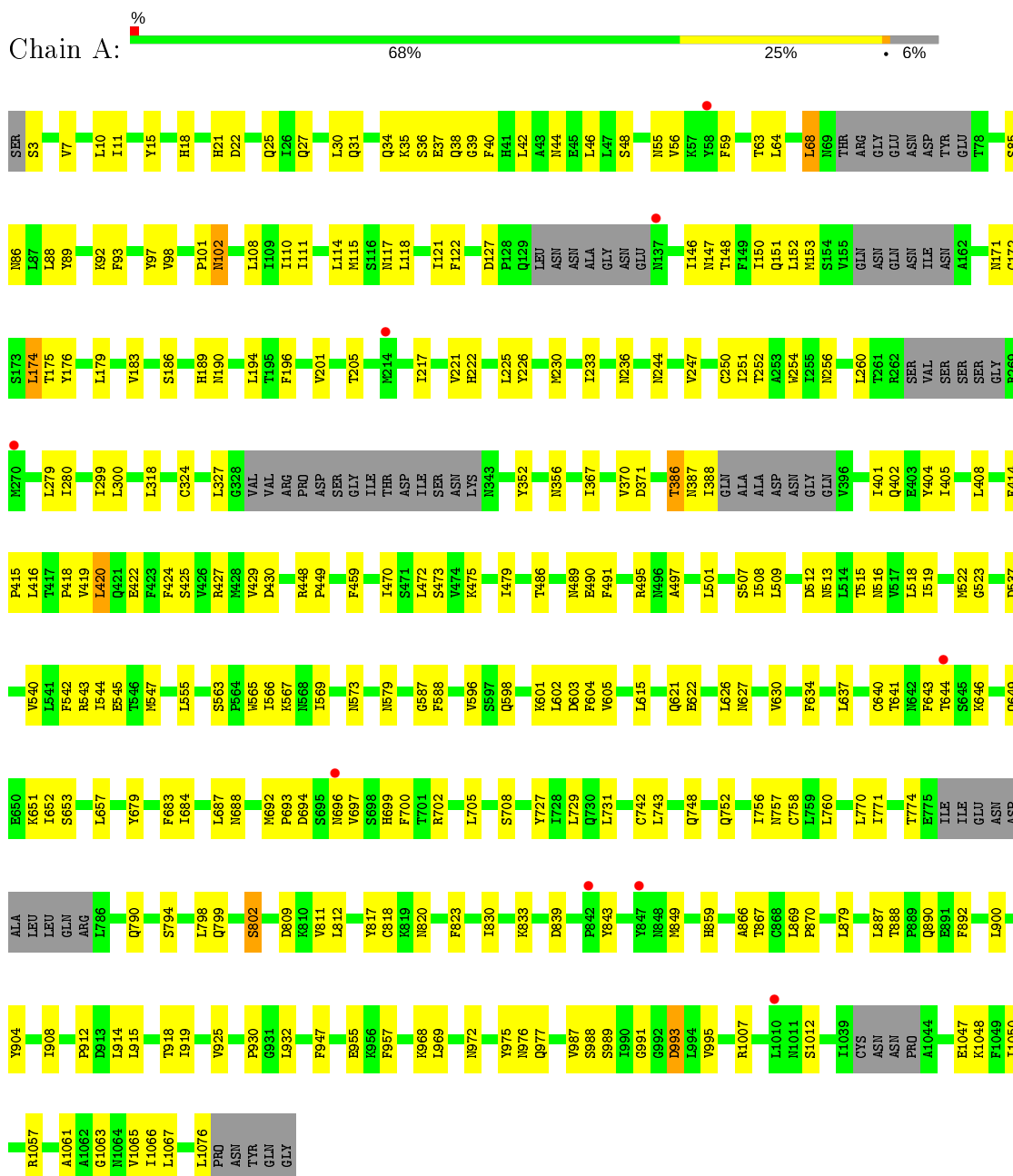
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLY	-	expression tag	UNP P50623


### 3 Residue-property plots

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: Importin beta-like protein KAP122



## ● Molecule 2: UBC9

Chain B:  % 85% 14%



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	198.26Å 198.26Å 289.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.56 – 4.53 49.57 – 4.53	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.56-4.53) 99.5 (49.57-4.53)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.94 (at 4.45Å)	Xtrriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.259 , 0.271 0.263 , 0.273	Depositor DCC
$R_{free}$ test set	388 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	293.0	Xtrriage
Anisotropy	0.208	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 248.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	9380	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	311.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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](#)

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	0/8280	0.53	1/11239 (0.0%)
2	B	0.31	0/1298	0.50	0/1762
All	All	0.31	0/9578	0.53	1/13001 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	LEU	CA-CB-CG	5.29	127.47	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	8122	0	8102	185	0
2	B	1258	0	1243	24	0
All	All	9380	0	9345	195	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 10.

All (195) 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:602:LEU:HD11	1:A:653:SER:HB3	1.49	0.93
1:A:588:PHE:HZ	1:A:640:CYS:HB3	1.38	0.89
1:A:867:THR:O	2:B:135:ARG:NH2	2.14	0.80
1:A:151:GLN:HB2	1:A:174:LEU:HD11	1.69	0.74
1:A:418:PRO:O	1:A:420:LEU:N	2.21	0.73
1:A:700:PHE:HB2	2:B:150:LEU:HD22	1.71	0.71
1:A:430:ASP:HA	2:B:17:ARG:HH21	1.55	0.71
1:A:1063:GLY:H	1:A:1066:ILE:HD13	1.55	0.70
1:A:64:LEU:O	1:A:89:TYR:OH	2.09	0.69
2:B:87:TYR:HD2	2:B:91:THR:HB	1.56	0.69
1:A:1063:GLY:HA2	1:A:1066:ILE:HB	1.76	0.68
1:A:870:PRO:HB2	2:B:135:ARG:HH12	1.59	0.68
1:A:108:LEU:HA	1:A:111:ILE:HB	1.75	0.67
1:A:605:VAL:HG11	1:A:657:LEU:HG	1.77	0.65
1:A:7:VAL:HG11	1:A:42:LEU:HD21	1.79	0.64
1:A:900:LEU:HD13	1:A:918:THR:HG23	1.79	0.64
1:A:147:ASN:HB3	1:A:174:LEU:CB	2.28	0.64
1:A:799:GLN:HB3	1:A:802:SER:HB3	1.78	0.64
1:A:429:VAL:HG11	1:A:497:ALA:HB1	1.80	0.63
1:A:489:ASN:HB3	2:B:22:PHE:CG	2.34	0.63
1:A:280:ILE:HD12	1:A:318:LEU:HD12	1.80	0.62
1:A:148:THR:OG1	1:A:171:ASN:O	2.15	0.62
1:A:870:PRO:HB2	2:B:135:ARG:NH1	2.15	0.62
1:A:36:SER:O	1:A:38:GLN:N	2.33	0.61
1:A:415:PRO:O	1:A:416:LEU:HB2	2.00	0.61
1:A:422:GLU:OE1	1:A:424:PHE:N	2.33	0.61
1:A:679:TYR:HB3	1:A:683:PHE:CZ	2.35	0.61
1:A:729:LEU:HD22	1:A:798:LEU:HD22	1.83	0.61
2:B:146:LYS:O	2:B:150:LEU:HG	2.01	0.61
1:A:976:ASN:OD1	1:A:977:GLN:N	2.34	0.61
1:A:569:ILE:O	1:A:573:ASN:ND2	2.32	0.60
1:A:489:ASN:HB3	2:B:22:PHE:CD1	2.38	0.59
1:A:371:ASP:OD1	1:A:427:ARG:NH2	2.36	0.58
1:A:151:GLN:HB2	1:A:174:LEU:HD21	1.85	0.58
1:A:888:THR:HG22	1:A:890:GLN:H	1.67	0.58
1:A:352:TYR:O	1:A:356:ASN:ND2	2.26	0.58
1:A:544:ILE:HD11	1:A:579:ASN:HB3	1.85	0.58
1:A:1050:ILE:HG13	1:A:1065:VAL:HG12	1.86	0.57
1:A:849:MET:HG3	1:A:879:LEU:HD21	1.85	0.57
1:A:887:LEU:HB3	1:A:932:LEU:HD21	1.85	0.57
1:A:370:VAL:HG11	1:A:427:ARG:HB3	1.85	0.57
1:A:40:PHE:CZ	1:A:88:LEU:HD12	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:TYR:CZ	1:A:111:ILE:HG23	2.39	0.56
1:A:48:SER:HB3	1:A:92:LYS:HE3	1.88	0.56
1:A:563:SER:O	1:A:621:GLN:NE2	2.31	0.55
1:A:179:LEU:O	1:A:183:VAL:HG23	2.06	0.55
1:A:221:VAL:HG13	1:A:225:LEU:HB3	1.89	0.55
1:A:11:ILE:HG21	1:A:56:VAL:HG13	1.88	0.55
1:A:175:THR:OG1	1:A:176:TYR:N	2.38	0.55
1:A:147:ASN:HB3	1:A:174:LEU:HB2	1.89	0.54
2:B:87:TYR:CD2	2:B:91:THR:HB	2.41	0.54
1:A:919:ILE:HG12	1:A:947:PHE:CE1	2.42	0.54
1:A:233:ILE:HD13	1:A:251:ILE:HG12	1.88	0.54
1:A:102:ASN:N	1:A:102:ASN:OD1	2.31	0.54
1:A:148:THR:O	1:A:152:LEU:HG	2.08	0.54
2:B:29:VAL:HG12	2:B:42:GLU:HG3	1.90	0.53
1:A:386:THR:HG23	1:A:388:ILE:H	1.73	0.53
1:A:475:LYS:HB2	1:A:491:PHE:CE2	2.44	0.53
1:A:174:LEU:HD13	1:A:175:THR:H	1.74	0.53
1:A:687:LEU:HD13	1:A:705:LEU:HD23	1.92	0.52
1:A:637:LEU:O	1:A:641:THR:OG1	2.14	0.52
1:A:27:GLN:NE2	1:A:31:GLN:OE1	2.39	0.52
1:A:692:MET:C	1:A:694:ASP:H	2.11	0.52
1:A:915:LEU:O	1:A:919:ILE:HG13	2.08	0.52
1:A:279:LEU:HD22	1:A:299:ILE:HG12	1.91	0.52
2:B:95:SER:HA	2:B:101:GLN:OE1	2.10	0.52
1:A:771:ILE:HG12	1:A:843:TYR:CE1	2.45	0.52
1:A:97:TYR:OH	1:A:111:ILE:O	2.27	0.52
1:A:537:ASP:HB3	1:A:540:VAL:HG12	1.91	0.52
1:A:1057:ARG:NE	1:A:1061:ALA:HB3	2.25	0.52
1:A:470:ILE:O	1:A:495:ARG:NH1	2.42	0.51
1:A:201:VAL:HG11	1:A:254:TRP:HA	1.91	0.51
1:A:429:VAL:HG13	1:A:501:LEU:HD12	1.91	0.51
1:A:699:HIS:CG	1:A:758:CYS:HB2	2.46	0.51
1:A:869:LEU:HB3	1:A:870:PRO:HD3	1.93	0.51
1:A:89:TYR:O	1:A:93:PHE:HB2	2.11	0.51
2:B:94:LEU:HG	2:B:96:ILE:HG12	1.92	0.51
1:A:448:ARG:HG2	1:A:449:PRO:HD2	1.94	0.50
1:A:968:LYS:O	1:A:972:ASN:HB2	2.11	0.50
1:A:89:TYR:CE1	1:A:117:ASN:HB3	2.46	0.50
1:A:86:ASN:HA	1:A:89:TYR:HD2	1.77	0.49
1:A:110:ILE:HD12	1:A:110:ILE:H	1.77	0.49
1:A:190:ASN:ND2	1:A:236:ASN:OD1	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:ASP:CG	2:B:17:ARG:HE	2.15	0.49
1:A:371:ASP:OD2	2:B:18:LYS:NZ	2.46	0.49
1:A:523:GLY:HA3	1:A:573:ASN:HB3	1.93	0.49
1:A:870:PRO:HD2	2:B:135:ARG:NH2	2.27	0.49
1:A:790:GLN:O	1:A:794:SER:OG	2.20	0.49
1:A:118:LEU:HD23	1:A:121:ILE:HD12	1.94	0.49
1:A:486:THR:O	1:A:490:GLU:HG3	2.13	0.49
1:A:692:MET:O	1:A:694:ASP:N	2.43	0.49
1:A:68:LEU:HD23	1:A:68:LEU:H	1.79	0.48
1:A:186:SER:HB3	1:A:189:HIS:ND1	2.28	0.48
1:A:22:ASP:HB3	1:A:25:GLN:HB2	1.95	0.48
1:A:151:GLN:CB	1:A:174:LEU:HD21	2.43	0.48
1:A:15:TYR:CE2	1:A:55:ASN:HB3	2.48	0.48
1:A:598:GLN:O	1:A:602:LEU:HD13	2.14	0.48
1:A:401:ILE:O	1:A:405:ILE:HG12	2.13	0.48
1:A:812:LEU:O	1:A:818:CYS:HB3	2.14	0.47
1:A:40:PHE:O	1:A:44:ASN:ND2	2.47	0.47
1:A:752:GLN:O	1:A:756:ILE:HG12	2.14	0.47
1:A:993:ASP:N	1:A:993:ASP:OD1	2.45	0.47
1:A:1063:GLY:CA	1:A:1066:ILE:HB	2.45	0.47
1:A:174:LEU:HD22	1:A:174:LEU:HA	1.59	0.47
1:A:148:THR:HG23	1:A:171:ASN:HA	1.95	0.47
1:A:475:LYS:NZ	1:A:603:ASP:OD2	2.46	0.47
1:A:684:ILE:O	1:A:688:ASN:ND2	2.40	0.47
1:A:42:LEU:O	1:A:46:LEU:HG	2.15	0.47
1:A:473:SER:HB3	1:A:542:PHE:CD2	2.49	0.47
1:A:122:PHE:CE2	1:A:217:ILE:HD11	2.50	0.47
1:A:743:LEU:HD11	1:A:811:VAL:HG22	1.96	0.47
1:A:757:ASN:OD1	2:B:146:LYS:NZ	2.39	0.47
1:A:401:ILE:HA	1:A:404:TYR:HB3	1.97	0.46
1:A:699:HIS:CD2	1:A:758:CYS:HB2	2.50	0.46
1:A:748:GLN:HB3	1:A:817:TYR:OH	2.15	0.46
1:A:729:LEU:HB2	1:A:798:LEU:HD13	1.98	0.46
1:A:40:PHE:HE1	1:A:85:SER:HA	1.81	0.46
1:A:912:PRO:HB3	1:A:955:GLU:HG2	1.97	0.46
1:A:97:TYR:OH	1:A:115:MET:HG3	2.16	0.46
1:A:1007:ARG:HA	1:A:1007:ARG:HD2	1.79	0.46
1:A:567:LYS:HG3	1:A:622:GLU:OE1	2.16	0.46
1:A:40:PHE:CE1	1:A:85:SER:HA	2.51	0.45
2:B:72:PRO:HD3	2:B:99:GLU:HB3	1.98	0.45
1:A:544:ILE:CD1	1:A:579:ASN:HB3	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:830:ILE:HA	1:A:833:LYS:HG2	1.98	0.45
1:A:866:ALA:O	1:A:914:LEU:HD11	2.16	0.45
1:A:930:PRO:HG3	1:A:972:ASN:HD21	1.81	0.45
1:A:820:ASN:OD1	2:B:140:ASN:ND2	2.50	0.45
1:A:30:LEU:HD22	1:A:59:PHE:CE1	2.52	0.45
1:A:646:LYS:O	1:A:646:LYS:HG3	2.17	0.45
1:A:1047:GLU:O	1:A:1050:ILE:HG22	2.17	0.45
1:A:425:SER:O	1:A:429:VAL:HG23	2.17	0.45
1:A:201:VAL:HG21	1:A:254:TRP:CE2	2.53	0.44
1:A:327:LEU:HD13	1:A:408:LEU:HD11	1.98	0.44
1:A:697:VAL:HG21	1:A:702:ARG:HG3	2.00	0.44
1:A:587:GLY:O	1:A:601:LYS:HG3	2.17	0.44
1:A:627:ASN:OD1	1:A:679:TYR:OH	2.33	0.44
1:A:957:PHE:CE1	2:B:128:PRO:HG2	2.52	0.44
1:A:644:THR:HG22	1:A:696:ASN:HD21	1.81	0.44
1:A:300:LEU:HA	1:A:300:LEU:HD23	1.78	0.44
1:A:97:TYR:CE2	1:A:111:ILE:HG12	2.53	0.44
1:A:742:CYS:SG	1:A:752:GLN:HG3	2.57	0.44
1:A:833:LYS:HA	1:A:833:LYS:HD3	1.65	0.44
1:A:175:THR:HG21	1:A:179:LEU:HD23	2.00	0.44
1:A:892:PHE:CE1	1:A:925:VAL:HG21	2.53	0.44
1:A:147:ASN:HB3	1:A:174:LEU:HB3	2.00	0.43
1:A:459:PHE:CE2	1:A:508:ILE:HD13	2.53	0.43
1:A:770:LEU:HD23	1:A:770:LEU:HA	1.89	0.43
1:A:64:LEU:HB3	1:A:89:TYR:CE1	2.53	0.43
1:A:904:TYR:HA	1:A:908:ILE:HB	2.00	0.43
1:A:98:VAL:O	1:A:101:PRO:HD2	2.18	0.43
1:A:588:PHE:CZ	1:A:640:CYS:HB3	2.30	0.43
1:A:727:TYR:O	1:A:731:LEU:HG	2.18	0.43
1:A:509:LEU:O	1:A:513:ASN:HB2	2.18	0.43
1:A:430:ASP:OD1	2:B:17:ARG:NE	2.45	0.43
1:A:566:ILE:HA	1:A:569:ILE:HG22	2.01	0.43
1:A:97:TYR:HB3	1:A:153:MET:HE1	2.00	0.43
1:A:820:ASN:HB3	1:A:823:PHE:HD2	1.83	0.42
2:B:9:LEU:HD13	2:B:38:LEU:O	2.18	0.42
1:A:743:LEU:HD21	1:A:756:ILE:HG13	2.02	0.42
1:A:809:ASP:HA	1:A:859:HIS:CE1	2.54	0.42
1:A:115:MET:HG2	1:A:196:PHE:CD1	2.55	0.42
1:A:543:ARG:HG3	1:A:547:MET:HE2	2.01	0.42
1:A:479:ILE:HG22	1:A:596:VAL:HG22	2.00	0.42
1:A:649:GLN:O	1:A:652:ILE:HG12	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:756:ILE:O	1:A:760:LEU:HG	2.18	0.42
1:A:252:THR:HG22	1:A:256:ASN:HD21	1.83	0.42
1:A:34:GLN:HB3	1:A:63:THR:HG23	2.01	0.42
1:A:35:LYS:HG2	1:A:35:LYS:O	2.20	0.42
1:A:849:MET:HA	1:A:879:LEU:HD11	2.01	0.42
1:A:991:GLY:O	1:A:995:VAL:HG23	2.20	0.42
1:A:36:SER:O	1:A:39:GLY:N	2.27	0.42
1:A:367:ILE:HG23	1:A:427:ARG:HH12	1.84	0.42
1:A:93:PHE:CD1	1:A:114:LEU:HB3	2.54	0.42
1:A:1057:ARG:HD2	1:A:1057:ARG:HA	1.75	0.41
1:A:518:LEU:O	1:A:522:MET:HG2	2.20	0.41
1:A:226:TYR:O	1:A:230:MET:HG2	2.20	0.41
1:A:146:ILE:O	1:A:150:ILE:HG13	2.20	0.41
1:A:205:THR:HG23	1:A:260:LEU:HD12	2.03	0.41
1:A:545:GLU:HG3	1:A:604:PHE:CE1	2.55	0.41
1:A:555:LEU:HD23	1:A:615:LEU:HD21	2.01	0.41
1:A:644:THR:HG22	1:A:696:ASN:OD1	2.19	0.41
1:A:472:LEU:HD12	1:A:472:LEU:HA	1.83	0.41
1:A:515:THR:O	1:A:519:ILE:HG13	2.21	0.41
2:B:72:PRO:CD	2:B:99:GLU:HB3	2.50	0.41
1:A:1048:LYS:HB3	1:A:1048:LYS:HE3	1.81	0.41
1:A:108:LEU:CA	1:A:111:ILE:HB	2.47	0.41
1:A:3:SER:O	1:A:7:VAL:HG23	2.21	0.41
1:A:244:ASN:O	1:A:247:VAL:HG12	2.21	0.41
1:A:512:ASP:HA	1:A:516:ASN:HB2	2.02	0.41
1:A:634:PHE:HA	1:A:637:LEU:HB3	2.03	0.41
1:A:969:LEU:HD12	1:A:975:TYR:OH	2.21	0.40
2:B:144:TYR:O	2:B:148:VAL:HG23	2.21	0.40
1:A:987:VAL:C	1:A:989:SER:H	2.25	0.40
1:A:630:VAL:O	1:A:634:PHE:HD2	2.04	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	994/1080 (92%)	925 (93%)	63 (6%)	6 (1%)	25	65
2	B	154/157 (98%)	150 (97%)	4 (3%)	0	100	100
All	All	1148/1237 (93%)	1075 (94%)	67 (6%)	6 (0%)	29	68

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	419	VAL
1	A	37	GLU
1	A	988	SER
1	A	414	PHE
1	A	693	PRO

### 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	931/1000 (93%)	904 (97%)	27 (3%)	42	64
2	B	136/136 (100%)	136 (100%)	0	100	100
All	All	1067/1136 (94%)	1040 (98%)	27 (2%)	47	68

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	68	LEU
1	A	102	ASN
1	A	127	ASP
1	A	172	CYS
1	A	174	LEU
1	A	194	LEU

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Mol	Chain	Res	Type
1	A	222	HIS
1	A	250	CYS
1	A	324	CYS
1	A	386	THR
1	A	387	ASN
1	A	402	GLN
1	A	420	LEU
1	A	507	SER
1	A	565	TRP
1	A	626	LEU
1	A	643	PHE
1	A	651	LYS
1	A	708	SER
1	A	774	THR
1	A	802	SER
1	A	839	ASP
1	A	993	ASP
1	A	1012	SER
1	A	1067	LEU
1	A	1076	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	44	ASN
1	A	256	ASN
1	A	402	GLN
1	A	577	ASN
1	A	696	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 carbohydrates 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, 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	1012/1080 (93%)	-0.06	9 (0%) 84 77	289, 312, 329, 351	0
2	B	156/157 (99%)	-0.06	1 (0%) 89 84	299, 308, 317, 322	0
All	All	1168/1237 (94%)	-0.06	10 (0%) 84 77	289, 312, 327, 351	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	644	THR	3.1
1	A	58	TYR	2.7
2	B	82	TYR	2.5
1	A	847	TYR	2.3
1	A	137	ASN	2.2
1	A	1010	LEU	2.2
1	A	842	PRO	2.0
1	A	214	MET	2.0
1	A	696	ASN	2.0
1	A	270	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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.



## 6.5 Other polymers

There are no such residues in this entry.