Full wwPDB X-ray Structure Validation Report (i)

Feb 15, 2024 – 05:57 PM EST

PDB ID	:	3QYB
Title	:	X-ray Crystal Structure of Human TBC1D4 (AS160) RabGAP domain
Authors	:	Park, S.Y.; Shoelson, S.E.
Deposited on	:	2011-03-03
Resolution	:	3.50 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:

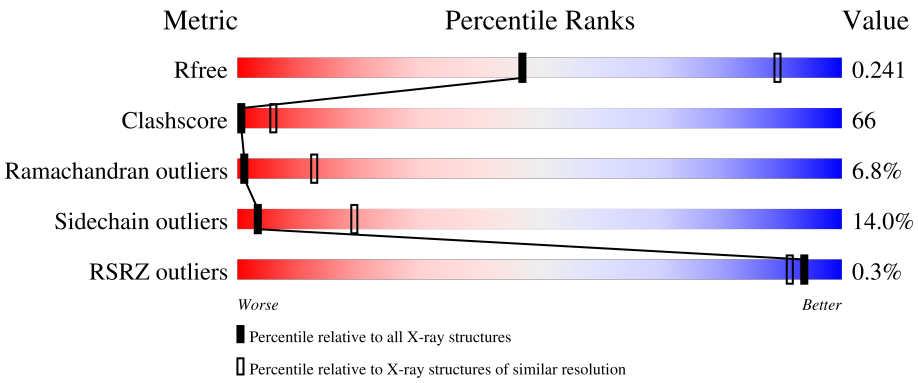
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	301	31% 53% 13% ••

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2453 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 TBC1 domain family member 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	297	2444	1584	403	440	17	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	871	GLY	-	expression tag	UNP O60343
A	872	SER	-	expression tag	UNP O60343
A	873	HIS	-	expression tag	UNP O60343
A	874	MET	-	expression tag	UNP O60343

- Molecule 2 is water.

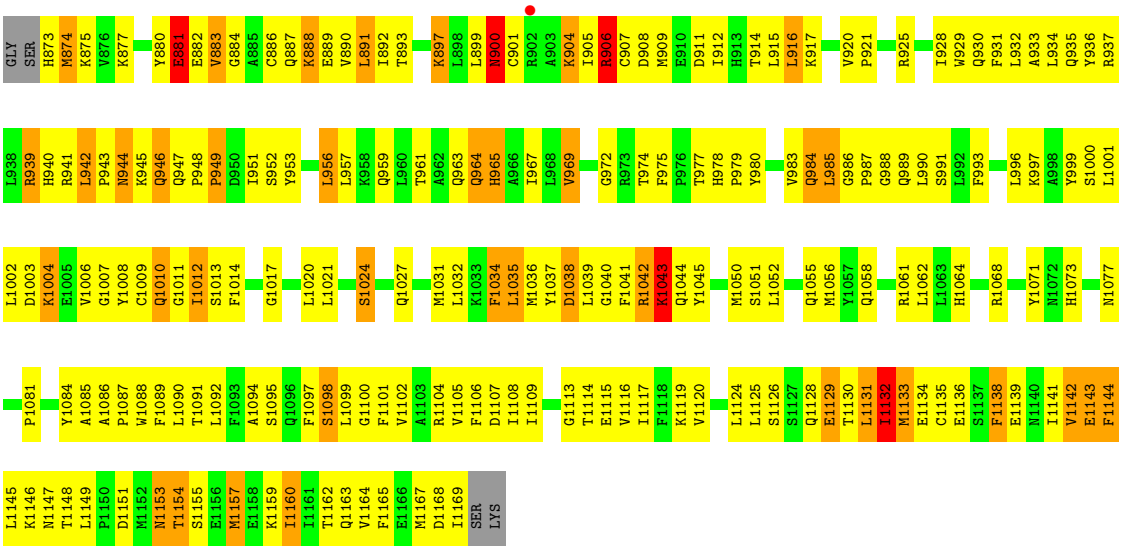
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	9	9	9	0	0

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: TBC1 domain family member 4

Chain A: 31% 53% 13%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	151.21Å 151.21Å 53.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.50 49.25 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.50) 99.3 (49.25-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.06 (at 3.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.300 0.230 , 0.241	Depositor DCC
R_{free} test set	424 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å^2)	123.4	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 106.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2453	wwPDB-VP
Average B, all atoms (Å^2)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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	0.54	2/2499 (0.1%)	0.75	2/3372 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	969	VAL	CB-CG2	-6.72	1.38	1.52
1	A	1157	MET	SD-CE	-5.23	1.48	1.77

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	969	VAL	CG1-CB-CG2	-5.53	102.05	110.90
1	A	1042	ARG	NE-CZ-NH1	-5.51	117.54	120.30

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	2444	0	2462	326	0
2	A	9	0	0	0	0
All	All	2453	0	2462	326	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 66.

All (326) 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:1034:PHE:CE1	1:A:1039:LEU:HG	1.56	1.39
1:A:1043:LYS:CD	1:A:1044:GLN:H	1.41	1.33
1:A:1043:LYS:NZ	1:A:1055:GLN:HE22	1.36	1.23
1:A:1037:TYR:CD2	1:A:1042:ARG:HD2	1.82	1.15
1:A:1043:LYS:HD2	1:A:1044:GLN:H	0.96	1.13
1:A:906:ARG:HD2	1:A:939:ARG:HG3	1.32	1.11
1:A:909:MET:HE2	1:A:912:ILE:HD11	1.25	1.09
1:A:948:PRO:HB3	1:A:1037:TYR:CE1	1.88	1.08
1:A:1043:LYS:HD2	1:A:1044:GLN:N	1.70	1.07
1:A:1037:TYR:HD2	1:A:1042:ARG:HD2	1.24	1.03
1:A:1034:PHE:HE1	1:A:1039:LEU:CG	1.72	1.02
1:A:944:ASN:HA	1:A:947:GLN:HG3	1.37	1.02
1:A:897:LYS:HE3	1:A:907:CYS:SG	2.01	1.00
1:A:909:MET:CE	1:A:912:ILE:HD11	1.91	1.00
1:A:948:PRO:HB3	1:A:1037:TYR:CD1	1.97	0.99
1:A:909:MET:HE1	1:A:939:ARG:NH1	1.78	0.99
1:A:1084:TYR:HA	1:A:1146:LYS:NZ	1.78	0.97
1:A:1034:PHE:CE1	1:A:1039:LEU:CG	2.47	0.97
1:A:1043:LYS:CD	1:A:1044:GLN:N	2.25	0.97
1:A:906:ARG:HB3	1:A:939:ARG:HD2	1.46	0.96
1:A:1144:PHE:CE2	1:A:1145:LEU:HD23	2.01	0.95
1:A:901:CYS:HB3	1:A:904:LYS:HD3	1.51	0.93
1:A:1128:GLN:HG3	1:A:1131:LEU:CG	1.97	0.93
1:A:1043:LYS:HD3	1:A:1044:GLN:H	1.32	0.93
1:A:1105:VAL:O	1:A:1109:ILE:HG12	1.66	0.93
1:A:1084:TYR:HA	1:A:1146:LYS:HZ1	1.27	0.93
1:A:909:MET:HE1	1:A:939:ARG:HH12	1.33	0.90
1:A:1043:LYS:NZ	1:A:1055:GLN:NE2	2.20	0.90
1:A:909:MET:HE3	1:A:939:ARG:HH22	1.39	0.88
1:A:887:GLN:O	1:A:890:VAL:HG12	1.72	0.88
1:A:906:ARG:HD2	1:A:939:ARG:CG	2.02	0.88
1:A:933:ALA:HA	1:A:1034:PHE:HE2	1.38	0.88
1:A:1043:LYS:HZ3	1:A:1055:GLN:HE22	1.15	0.87
1:A:909:MET:CE	1:A:939:ARG:HH22	1.86	0.87
1:A:1126:SER:HA	1:A:1129:GLU:HB3	1.56	0.86
1:A:901:CYS:HB3	1:A:904:LYS:CD	2.05	0.85
1:A:1113:GLY:O	1:A:1115:GLU:N	2.09	0.84
1:A:909:MET:HE2	1:A:909:MET:HA	1.59	0.84
1:A:1034:PHE:HE1	1:A:1039:LEU:HG	0.78	0.83
1:A:1043:LYS:HZ2	1:A:1055:GLN:HE22	1.22	0.83

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:942:LEU:HD23	1:A:943:PRO:HD2	1.57	0.83
1:A:1128:GLN:HE21	1:A:1131:LEU:HD21	1.44	0.83
1:A:1044:GLN:HG3	1:A:1045:TYR:CD2	2.15	0.82
1:A:1124:LEU:HD13	1:A:1160:ILE:HG13	1.62	0.82
1:A:943:PRO:HG2	1:A:946:GLN:NE2	1.95	0.81
1:A:1037:TYR:HE2	1:A:1042:ARG:HE	1.27	0.81
1:A:1128:GLN:NE2	1:A:1131:LEU:HD21	1.98	0.79
1:A:1043:LYS:HZ3	1:A:1055:GLN:NE2	1.79	0.79
1:A:948:PRO:HB3	1:A:1037:TYR:HE1	1.48	0.78
1:A:1009:CYS:O	1:A:1012:ILE:HG22	1.83	0.78
1:A:1037:TYR:CE2	1:A:1042:ARG:HD2	2.18	0.78
1:A:905:ILE:HG22	1:A:906:ARG:H	1.48	0.77
1:A:1043:LYS:HE3	1:A:1044:GLN:HG2	1.65	0.77
1:A:906:ARG:CB	1:A:939:ARG:HD2	2.13	0.77
1:A:1128:GLN:CG	1:A:1131:LEU:HD11	2.15	0.77
1:A:1146:LYS:C	1:A:1147:ASN:HD22	1.87	0.76
1:A:909:MET:CE	1:A:939:ARG:NH2	2.48	0.76
1:A:1128:GLN:HG3	1:A:1131:LEU:HG	1.66	0.76
1:A:1094:ALA:HA	1:A:1102:VAL:HG21	1.68	0.76
1:A:1086:ALA:HB3	1:A:1087:PRO:HD3	1.66	0.76
1:A:909:MET:HE3	1:A:939:ARG:NH2	2.01	0.75
1:A:1086:ALA:O	1:A:1090:LEU:HD23	1.87	0.74
1:A:1128:GLN:HG3	1:A:1131:LEU:CD1	2.16	0.74
1:A:1128:GLN:HG3	1:A:1131:LEU:HD11	1.70	0.74
1:A:1034:PHE:C	1:A:1034:PHE:HD1	1.92	0.73
1:A:881:GLU:OE1	1:A:1099:LEU:N	2.22	0.73
1:A:1037:TYR:HE2	1:A:1042:ARG:NE	1.86	0.73
1:A:909:MET:CE	1:A:939:ARG:HH12	2.01	0.72
1:A:873:HIS:O	1:A:875:LYS:N	2.23	0.72
1:A:1159:LYS:O	1:A:1162:THR:HG22	1.89	0.72
1:A:991:SER:HB2	1:A:1020:LEU:HD21	1.71	0.72
1:A:984:GLN:HA	1:A:984:GLN:OE1	1.89	0.72
1:A:1128:GLN:HE21	1:A:1131:LEU:CD2	2.02	0.71
1:A:1037:TYR:CD2	1:A:1042:ARG:CD	2.70	0.71
1:A:933:ALA:CA	1:A:1034:PHE:HE2	2.01	0.71
1:A:996:LEU:HD11	1:A:1013:SER:CB	2.21	0.71
1:A:1101:PHE:HB2	1:A:1157:MET:CE	2.20	0.71
1:A:1034:PHE:C	1:A:1034:PHE:CD1	2.64	0.71
1:A:1084:TYR:CA	1:A:1146:LYS:NZ	2.54	0.70
1:A:964:GLN:O	1:A:965:HIS:C	2.30	0.70
1:A:948:PRO:CB	1:A:1037:TYR:CE1	2.70	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:899:LEU:O	1:A:901:CYS:N	2.24	0.69
1:A:916:LEU:CD1	1:A:1107:ASP:HB3	2.22	0.69
1:A:1101:PHE:HB2	1:A:1157:MET:HE3	1.72	0.69
1:A:1088:TRP:HH2	1:A:1125:LEU:HD21	1.57	0.69
1:A:963:GLN:O	1:A:967:ILE:HG13	1.93	0.68
1:A:1163:GLN:O	1:A:1167:MET:HB2	1.94	0.68
1:A:1034:PHE:CD1	1:A:1038:ASP:HB2	2.28	0.68
1:A:1044:GLN:HG3	1:A:1045:TYR:HD2	1.59	0.68
1:A:1144:PHE:CE2	1:A:1145:LEU:CD2	2.76	0.68
1:A:933:ALA:HB1	1:A:1034:PHE:CD2	2.29	0.67
1:A:943:PRO:O	1:A:945:LYS:N	2.28	0.67
1:A:909:MET:HE1	1:A:939:ARG:CZ	2.23	0.67
1:A:1037:TYR:CE2	1:A:1042:ARG:CD	2.78	0.67
1:A:1037:TYR:CE2	1:A:1042:ARG:NE	2.61	0.67
1:A:1128:GLN:HG3	1:A:1131:LEU:CD2	2.24	0.67
1:A:1128:GLN:O	1:A:1131:LEU:HG	1.94	0.67
1:A:920:VAL:HG13	1:A:928:ILE:HD12	1.76	0.66
1:A:1135:CYS:HB2	1:A:1141:ILE:HG12	1.77	0.66
1:A:946:GLN:NE2	1:A:946:GLN:H	1.93	0.66
1:A:1041:PHE:HA	1:A:1043:LYS:HE2	1.76	0.66
1:A:906:ARG:CD	1:A:939:ARG:HG3	2.19	0.65
1:A:1128:GLN:HG3	1:A:1131:LEU:HD21	1.79	0.65
1:A:1043:LYS:HD3	1:A:1044:GLN:N	2.02	0.65
1:A:882:GLU:O	1:A:883:VAL:HG22	1.97	0.65
1:A:933:ALA:HA	1:A:1034:PHE:CE2	2.29	0.65
1:A:961:THR:OG1	1:A:997:LYS:HG3	1.97	0.65
1:A:1132:ILE:HG12	1:A:1133:MET:N	2.12	0.64
1:A:944:ASN:HA	1:A:947:GLN:CG	2.19	0.64
1:A:983:VAL:CG2	1:A:986:GLY:HA3	2.28	0.64
1:A:909:MET:CE	1:A:909:MET:HA	2.25	0.63
1:A:996:LEU:HD22	1:A:1008:TYR:CE1	2.33	0.63
1:A:975:PHE:CE1	1:A:1099:LEU:HD11	2.34	0.63
1:A:1002:LEU:HD23	1:A:1002:LEU:C	2.19	0.63
1:A:996:LEU:HD11	1:A:1013:SER:HB3	1.80	0.63
1:A:1168:ASP:OD1	1:A:1169:ILE:O	2.17	0.63
1:A:948:PRO:CB	1:A:1037:TYR:CD1	2.79	0.62
1:A:891:LEU:HD13	1:A:921:PRO:HG3	1.80	0.62
1:A:1153:ASN:O	1:A:1155:SER:N	2.33	0.62
1:A:1035:LEU:HD12	1:A:1041:PHE:CD2	2.35	0.61
1:A:984:GLN:O	1:A:986:GLY:N	2.34	0.61
1:A:905:ILE:HG22	1:A:906:ARG:N	2.15	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1160:ILE:HD13	1:A:1160:ILE:O	2.01	0.61
1:A:946:GLN:H	1:A:946:GLN:HE21	1.47	0.60
1:A:1011:GLY:O	1:A:1013:SER:N	2.34	0.60
1:A:997:LYS:O	1:A:1000:SER:HB3	2.01	0.60
1:A:1041:PHE:HE1	1:A:1117:ILE:HD11	1.67	0.60
1:A:972:GLY:HA2	1:A:989:GLN:HE22	1.67	0.59
1:A:1089:PHE:O	1:A:1092:LEU:HD23	2.03	0.59
1:A:1090:LEU:HD22	1:A:1090:LEU:N	2.16	0.59
1:A:1124:LEU:CD1	1:A:1160:ILE:HG13	2.31	0.59
1:A:909:MET:CE	1:A:939:ARG:NH1	2.59	0.59
1:A:1052:LEU:HD11	1:A:1090:LEU:HD21	1.85	0.59
1:A:940:HIS:O	1:A:941:ARG:HB2	2.01	0.58
1:A:948:PRO:CB	1:A:1037:TYR:HE1	2.10	0.58
1:A:1051:SER:O	1:A:1055:GLN:HG3	2.03	0.58
1:A:882:GLU:C	1:A:883:VAL:HG13	2.23	0.58
1:A:942:LEU:HD23	1:A:943:PRO:CD	2.31	0.58
1:A:964:GLN:O	1:A:967:ILE:N	2.35	0.58
1:A:1149:LEU:HD23	1:A:1149:LEU:O	2.04	0.58
1:A:906:ARG:HA	1:A:935:GLN:NE2	2.19	0.57
1:A:1157:MET:HA	1:A:1160:ILE:HG22	1.85	0.57
1:A:1119:LYS:HG2	1:A:1167:MET:HE2	1.87	0.57
1:A:1043:LYS:HE3	1:A:1044:GLN:HE21	1.69	0.57
1:A:1160:ILE:HD13	1:A:1160:ILE:C	2.25	0.57
1:A:983:VAL:HG23	1:A:986:GLY:HA3	1.86	0.57
1:A:1089:PHE:CD2	1:A:1117:ILE:HG21	2.39	0.57
1:A:1084:TYR:CB	1:A:1146:LYS:HZ3	2.17	0.56
1:A:901:CYS:HB3	1:A:904:LYS:HD2	1.84	0.56
1:A:930:GLN:O	1:A:933:ALA:HB3	2.05	0.56
1:A:987:PRO:HG2	1:A:988:GLY:H	1.70	0.56
1:A:1035:LEU:HD13	1:A:1039:LEU:HD12	1.87	0.56
1:A:978:HIS:HD2	1:A:980:TYR:HB3	1.70	0.56
1:A:1131:LEU:O	1:A:1132:ILE:O	2.24	0.55
1:A:1119:LYS:HA	1:A:1167:MET:HE1	1.88	0.55
1:A:1131:LEU:HD23	1:A:1131:LEU:N	2.20	0.55
1:A:1149:LEU:HD23	1:A:1149:LEU:C	2.26	0.55
1:A:933:ALA:CB	1:A:1034:PHE:CE2	2.89	0.55
1:A:1132:ILE:O	1:A:1134:GLU:N	2.39	0.55
1:A:996:LEU:HD22	1:A:1008:TYR:HE1	1.71	0.55
1:A:1000:SER:HA	1:A:1007:GLY:O	2.06	0.55
1:A:933:ALA:HB1	1:A:1034:PHE:CE2	2.41	0.55
1:A:1144:PHE:CD2	1:A:1145:LEU:HD23	2.40	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:904:LYS:O	1:A:905:ILE:HD13	2.07	0.54
1:A:1041:PHE:O	1:A:1043:LYS:HD2	2.06	0.54
1:A:925:ARG:HD2	1:A:929:TRP:CZ2	2.43	0.54
1:A:925:ARG:O	1:A:925:ARG:HG3	2.07	0.54
1:A:933:ALA:HB1	1:A:1034:PHE:HD2	1.72	0.54
1:A:909:MET:CE	1:A:939:ARG:CZ	2.86	0.54
1:A:956:LEU:HD11	1:A:1002:LEU:HB2	1.89	0.54
1:A:946:GLN:NE2	1:A:946:GLN:N	2.56	0.54
1:A:1084:TYR:CA	1:A:1146:LYS:HZ3	2.19	0.54
1:A:1036:MET:O	1:A:1042:ARG:HG2	2.07	0.53
1:A:1084:TYR:HB2	1:A:1146:LYS:HZ3	1.74	0.53
1:A:1088:TRP:CH2	1:A:1125:LEU:HD21	2.41	0.53
1:A:1116:VAL:O	1:A:1120:VAL:HG23	2.08	0.53
1:A:956:LEU:O	1:A:1001:LEU:HD12	2.09	0.53
1:A:883:VAL:CG2	1:A:1104:ARG:HG2	2.38	0.53
1:A:911:ASP:O	1:A:914:THR:HG22	2.08	0.53
1:A:987:PRO:O	1:A:990:LEU:HB2	2.10	0.52
1:A:906:ARG:HA	1:A:935:GLN:HE22	1.74	0.52
1:A:1052:LEU:O	1:A:1056:MET:HG2	2.10	0.52
1:A:1043:LYS:CE	1:A:1044:GLN:HG2	2.38	0.52
1:A:883:VAL:CG2	1:A:884:GLY:N	2.72	0.52
1:A:1044:GLN:HG3	1:A:1045:TYR:CE2	2.45	0.52
1:A:917:LYS:HA	1:A:1165:PHE:CE2	2.45	0.52
1:A:969:VAL:O	1:A:969:VAL:HG22	2.09	0.51
1:A:1038:ASP:C	1:A:1040:GLY:H	2.12	0.51
1:A:1147:ASN:N	1:A:1147:ASN:ND2	2.58	0.51
1:A:963:GLN:O	1:A:964:GLN:O	2.28	0.51
1:A:1147:ASN:HD22	1:A:1147:ASN:N	2.03	0.51
1:A:1144:PHE:O	1:A:1148:THR:HB	2.11	0.51
1:A:916:LEU:HD21	1:A:932:LEU:HD11	1.92	0.51
1:A:957:LEU:CD2	1:A:997:LYS:HD3	2.40	0.51
1:A:972:GLY:HA2	1:A:989:GLN:NE2	2.26	0.51
1:A:893:THR:HG21	1:A:915:LEU:HD21	1.92	0.51
1:A:911:ASP:HA	1:A:914:THR:HG22	1.92	0.51
1:A:978:HIS:NE2	1:A:1021:LEU:HD21	2.26	0.50
1:A:986:GLY:O	1:A:990:LEU:HG	2.11	0.50
1:A:983:VAL:HG21	1:A:986:GLY:HA3	1.92	0.50
1:A:1013:SER:O	1:A:1017:GLY:N	2.44	0.50
1:A:1132:ILE:O	1:A:1133:MET:C	2.50	0.50
1:A:906:ARG:O	1:A:907:CYS:C	2.47	0.50
1:A:943:PRO:HG2	1:A:946:GLN:HE21	1.71	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:948:PRO:HB3	1:A:1037:TYR:HD1	1.67	0.50
1:A:1052:LEU:CD1	1:A:1090:LEU:HD21	2.42	0.50
1:A:1128:GLN:CG	1:A:1131:LEU:HD21	2.39	0.50
1:A:1089:PHE:CD2	1:A:1117:ILE:CG2	2.95	0.50
1:A:880:TYR:O	1:A:881:GLU:O	2.29	0.50
1:A:920:VAL:HG13	1:A:928:ILE:CD1	2.40	0.50
1:A:959:GLN:HB2	1:A:1001:LEU:HD13	1.93	0.50
1:A:891:LEU:CD1	1:A:921:PRO:HG3	2.42	0.49
1:A:1081:PRO:HA	1:A:1084:TYR:CE2	2.46	0.49
1:A:943:PRO:C	1:A:945:LYS:H	2.15	0.49
1:A:886:CYS:SG	1:A:921:PRO:N	2.85	0.49
1:A:935:GLN:HB3	1:A:939:ARG:NE	2.26	0.49
1:A:1104:ARG:O	1:A:1107:ASP:N	2.46	0.49
1:A:936:TYR:CD2	1:A:936:TYR:O	2.66	0.49
1:A:933:ALA:CB	1:A:1034:PHE:HE2	2.25	0.49
1:A:1146:LYS:C	1:A:1147:ASN:ND2	2.64	0.49
1:A:1101:PHE:CB	1:A:1157:MET:HE3	2.40	0.49
1:A:937:ARG:NH2	1:A:1038:ASP:OD1	2.39	0.49
1:A:1035:LEU:HD12	1:A:1041:PHE:CE2	2.47	0.49
1:A:1091:THR:HG21	1:A:1095:SER:HB3	1.95	0.49
1:A:1130:THR:HG22	1:A:1130:THR:O	2.13	0.48
1:A:905:ILE:O	1:A:907:CYS:N	2.46	0.48
1:A:1011:GLY:O	1:A:1014:PHE:N	2.41	0.48
1:A:1104:ARG:HD2	1:A:1107:ASP:OD2	2.13	0.48
1:A:935:GLN:HB3	1:A:939:ARG:HE	1.78	0.48
1:A:1101:PHE:CG	1:A:1157:MET:HE3	2.48	0.48
1:A:1115:GLU:C	1:A:1117:ILE:N	2.62	0.48
1:A:1153:ASN:O	1:A:1153:ASN:ND2	2.43	0.48
1:A:886:CYS:SG	1:A:920:VAL:C	2.93	0.48
1:A:1011:GLY:C	1:A:1013:SER:N	2.66	0.48
1:A:1101:PHE:HB2	1:A:1157:MET:HE1	1.93	0.47
1:A:978:HIS:CD2	1:A:980:TYR:HB3	2.49	0.47
1:A:948:PRO:O	1:A:949:PRO:O	2.32	0.47
1:A:996:LEU:CD1	1:A:1013:SER:HB3	2.44	0.47
1:A:1144:PHE:HE2	1:A:1145:LEU:CD2	2.26	0.47
1:A:914:THR:O	1:A:917:LYS:HB3	2.15	0.47
1:A:1034:PHE:CE1	1:A:1039:LEU:CD1	2.98	0.47
1:A:1089:PHE:CE2	1:A:1117:ILE:HG22	2.49	0.47
1:A:1106:PHE:HA	1:A:1109:ILE:CG1	2.45	0.47
1:A:1119:LYS:HA	1:A:1167:MET:CE	2.45	0.47
1:A:1131:LEU:O	1:A:1132:ILE:C	2.52	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:929:TRP:CZ2	1:A:1106:PHE:HB3	2.50	0.47
1:A:1153:ASN:O	1:A:1154:THR:C	2.53	0.47
1:A:906:ARG:HB3	1:A:939:ARG:CD	2.32	0.47
1:A:1002:LEU:HD23	1:A:1003:ASP:N	2.30	0.46
1:A:1077:ASN:ND2	1:A:1136:GLU:O	2.48	0.46
1:A:1120:VAL:HA	1:A:1164:VAL:HG21	1.96	0.46
1:A:874:MET:O	1:A:877:LYS:HB2	2.15	0.46
1:A:936:TYR:CD2	1:A:936:TYR:C	2.88	0.46
1:A:990:LEU:O	1:A:993:PHE:HB3	2.15	0.46
1:A:1024:SER:OG	1:A:1027:GLN:HG3	2.14	0.46
1:A:985:LEU:HD12	1:A:985:LEU:HA	1.73	0.46
1:A:948:PRO:CA	1:A:1037:TYR:HE1	2.28	0.46
1:A:900:ASN:O	1:A:904:LYS:HD3	2.16	0.46
1:A:974:THR:HG23	1:A:1014:PHE:CD1	2.51	0.46
1:A:959:GLN:HB2	1:A:1001:LEU:CD1	2.45	0.46
1:A:1132:ILE:CG1	1:A:1133:MET:N	2.79	0.46
1:A:984:GLN:O	1:A:985:LEU:C	2.54	0.45
1:A:999:TYR:CE2	1:A:1003:ASP:HB3	2.51	0.45
1:A:1129:GLU:C	1:A:1131:LEU:N	2.65	0.45
1:A:1085:ALA:O	1:A:1086:ALA:C	2.53	0.45
1:A:1090:LEU:N	1:A:1090:LEU:CD2	2.80	0.45
1:A:916:LEU:HD11	1:A:1107:ASP:HB3	1.96	0.45
1:A:1097:PHE:N	1:A:1097:PHE:CD2	2.84	0.45
1:A:883:VAL:CG2	1:A:1104:ARG:CG	2.94	0.45
1:A:920:VAL:CG1	1:A:928:ILE:HD12	2.46	0.45
1:A:1011:GLY:O	1:A:1012:ILE:C	2.53	0.45
1:A:1043:LYS:HZ2	1:A:1055:GLN:NE2	1.97	0.45
1:A:1115:GLU:O	1:A:1116:VAL:C	2.55	0.45
1:A:1043:LYS:HE3	1:A:1044:GLN:NE2	2.32	0.44
1:A:948:PRO:CA	1:A:1037:TYR:CE1	3.01	0.44
1:A:1124:LEU:HG	1:A:1149:LEU:HD21	1.99	0.44
1:A:1084:TYR:CA	1:A:1146:LYS:HZ1	2.12	0.44
1:A:1045:TYR:CD2	1:A:1045:TYR:N	2.85	0.44
1:A:1098:SER:O	1:A:1101:PHE:HB3	2.18	0.44
1:A:1153:ASN:ND2	1:A:1153:ASN:H	2.14	0.44
1:A:1043:LYS:HE3	1:A:1044:GLN:CG	2.41	0.44
1:A:1159:LYS:HA	1:A:1162:THR:HG22	1.99	0.43
1:A:1056:MET:HE2	1:A:1089:PHE:CD1	2.53	0.43
1:A:1151:ASP:OD1	1:A:1151:ASP:O	2.36	0.43
1:A:920:VAL:HG23	1:A:1104:ARG:HH12	1.84	0.43
1:A:983:VAL:HG21	1:A:987:PRO:HD2	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1089:PHE:CE2	1:A:1117:ILE:CG2	3.01	0.43
1:A:1105:VAL:HG12	1:A:1109:ILE:HD11	2.01	0.43
1:A:1109:ILE:HD13	1:A:1116:VAL:HB	2.01	0.43
1:A:1090:LEU:HD22	1:A:1090:LEU:H	1.84	0.43
1:A:1058:GLN:O	1:A:1062:LEU:HB2	2.18	0.43
1:A:1119:LYS:HB3	1:A:1164:VAL:HG13	2.01	0.43
1:A:1040:GLY:O	1:A:1043:LYS:HG3	2.18	0.43
1:A:888:LYS:O	1:A:892:ILE:HG12	2.19	0.42
1:A:1090:LEU:CD2	1:A:1090:LEU:H	2.32	0.42
1:A:1159:LYS:O	1:A:1162:THR:CG2	2.64	0.42
1:A:891:LEU:HD13	1:A:891:LEU:HA	1.77	0.42
1:A:937:ARG:HH21	1:A:1038:ASP:CG	2.20	0.42
1:A:967:ILE:HG12	1:A:1008:TYR:CD1	2.55	0.42
1:A:1031:MET:O	1:A:1032:LEU:C	2.58	0.42
1:A:1099:LEU:N	1:A:1099:LEU:HD22	2.34	0.42
1:A:1064:HIS:HD2	1:A:1071:TYR:CE2	2.37	0.42
1:A:1034:PHE:O	1:A:1035:LEU:C	2.57	0.42
1:A:1101:PHE:CD1	1:A:1157:MET:CE	3.03	0.42
1:A:880:TYR:CE1	1:A:1100:GLY:HA2	2.55	0.41
1:A:935:GLN:O	1:A:936:TYR:C	2.59	0.41
1:A:1106:PHE:HA	1:A:1109:ILE:HG12	2.01	0.41
1:A:880:TYR:C	1:A:881:GLU:O	2.59	0.41
1:A:920:VAL:HG23	1:A:1104:ARG:NH1	2.35	0.41
1:A:935:GLN:HB3	1:A:939:ARG:NH1	2.35	0.41
1:A:1138:PHE:O	1:A:1142:VAL:N	2.54	0.41
1:A:1003:ASP:OD2	1:A:1006:VAL:HB	2.21	0.41
1:A:1139:GLU:O	1:A:1143:GLU:HB2	2.20	0.41
1:A:875:LYS:HE2	1:A:875:LYS:HA	2.02	0.41
1:A:909:MET:CE	1:A:909:MET:CA	2.97	0.41
1:A:952:SER:O	1:A:953:TYR:C	2.58	0.41
1:A:987:PRO:HG2	1:A:988:GLY:N	2.35	0.41
1:A:1146:LYS:HE3	1:A:1146:LYS:HB3	1.85	0.41
1:A:928:ILE:O	1:A:931:PHE:HB3	2.21	0.41
1:A:906:ARG:C	1:A:907:CYS:O	2.57	0.40
1:A:1041:PHE:C	1:A:1043:LYS:N	2.74	0.40
1:A:948:PRO:O	1:A:949:PRO:C	2.60	0.40
1:A:957:LEU:HD23	1:A:997:LYS:HD3	2.03	0.40
1:A:1056:MET:CE	1:A:1089:PHE:CD1	3.04	0.40
1:A:996:LEU:HD11	1:A:1013:SER:HB2	2.00	0.40
1:A:1073:HIS:CD2	1:A:1132:ILE:HG12	2.56	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	295/301 (98%)	213 (72%)	62 (21%)	20 (7%)	1 13

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	874	MET
1	A	881	GLU
1	A	906	ARG
1	A	944	ASN
1	A	964	GLN
1	A	965	HIS
1	A	985	LEU
1	A	1114	THR
1	A	1133	MET
1	A	1154	THR
1	A	900	ASN
1	A	1004	LYS
1	A	1012	ILE
1	A	949	PRO
1	A	1132	ILE
1	A	1010	GLN
1	A	1043	LYS
1	A	1143	GLU
1	A	979	PRO
1	A	1142	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	272/275 (99%)	234 (86%)	38 (14%)	3 19

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	881	GLU
1	A	883	VAL
1	A	888	LYS
1	A	889	GLU
1	A	891	LEU
1	A	897	LYS
1	A	900	ASN
1	A	904	LYS
1	A	906	ARG
1	A	908	ASP
1	A	916	LEU
1	A	934	LEU
1	A	939	ARG
1	A	942	LEU
1	A	946	GLN
1	A	951	ILE
1	A	956	LEU
1	A	977	THR
1	A	984	GLN
1	A	1004	LYS
1	A	1010	GLN
1	A	1024	SER
1	A	1034	PHE
1	A	1035	LEU
1	A	1038	ASP
1	A	1043	LYS
1	A	1050	MET
1	A	1061	ARG
1	A	1068	ARG
1	A	1098	SER
1	A	1108	ILE
1	A	1129	GLU
1	A	1131	LEU
1	A	1132	ILE
1	A	1138	PHE
1	A	1144	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	1153	ASN
1	A	1160	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	930	GLN
1	A	935	GLN
1	A	946	GLN
1	A	1044	GLN
1	A	1053	GLN
1	A	1055	GLN
1	A	1073	HIS
1	A	1128	GLN
1	A	1140	ASN
1	A	1147	ASN
1	A	1153	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	297/301 (98%)	-0.19	1 (0%) 94 91	33, 91, 170, 199	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	902	ARG	3.5

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.