Full wwPDB X-ray Structure Validation Report (i)

Sep 6, 2023 – 04:36 PM EDT

PDB ID	:	4GZU
Title	:	Crystal structure of the DH-PH-PH domain of FARP2
Authors	:	He, X.; Zhang, X.
Deposited on	:	2012-09-06
Resolution	:	3.20 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:

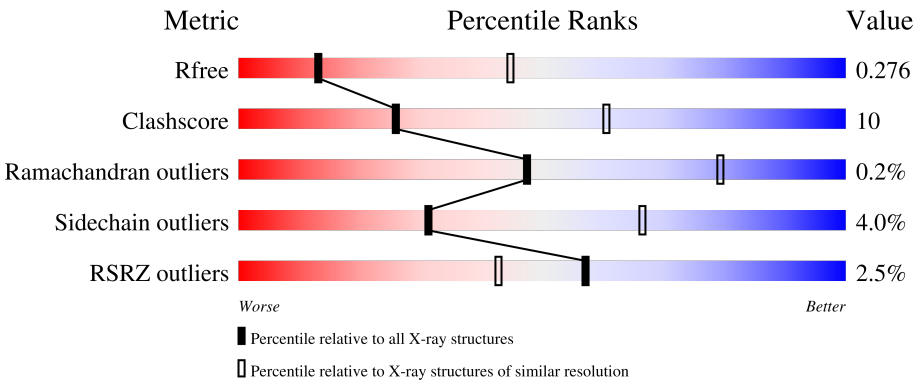
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation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

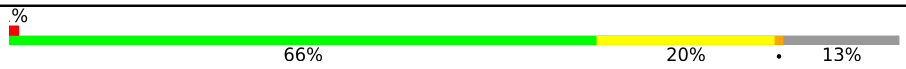
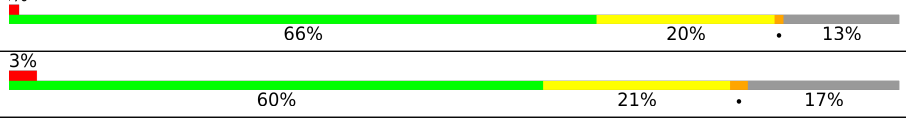
The reported resolution of this entry is 3.20 Å.

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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	501	
1	B	501	

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6724 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 FERM, RhoGEF and pleckstrin domain-containing protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	436	3463	2245	592	609	17	0	0	0
1	B	418	3242	2111	550	566	15	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	532	GLY	-	expression tag	UNP Q91VS8
A	533	PRO	-	expression tag	UNP Q91VS8
A	534	HIS	-	expression tag	UNP Q91VS8
A	535	MET	-	expression tag	UNP Q91VS8
A	821	LEU	PRO	SEE REMARK 999	UNP Q91VS8
B	532	GLY	-	expression tag	UNP Q91VS8
B	533	PRO	-	expression tag	UNP Q91VS8
B	534	HIS	-	expression tag	UNP Q91VS8
B	535	MET	-	expression tag	UNP Q91VS8
B	821	LEU	PRO	SEE REMARK 999	UNP Q91VS8

- Molecule 2 is water.

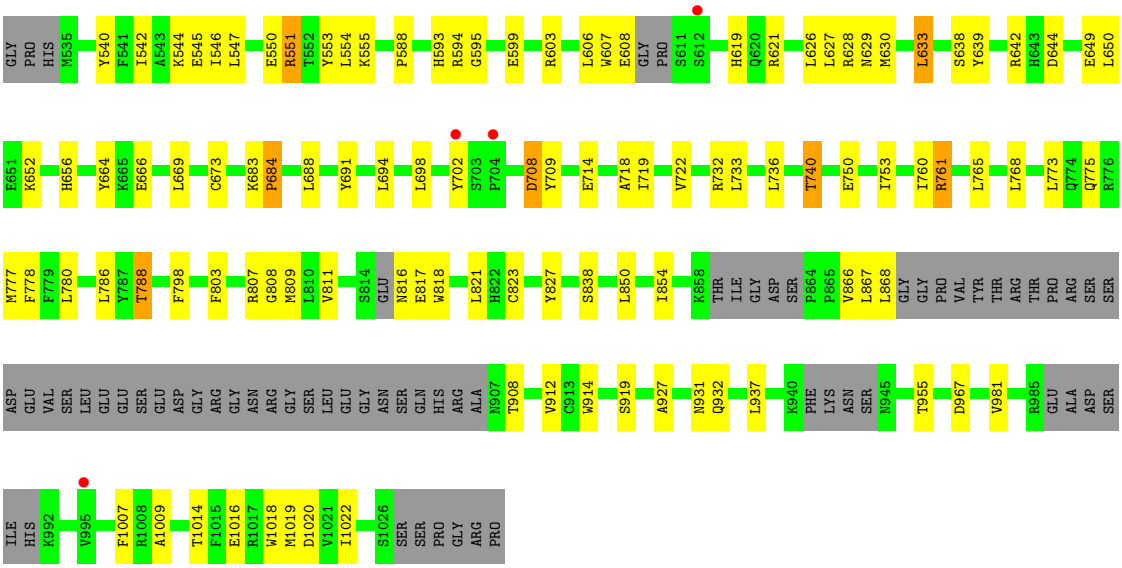
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	12	12	12	0	0
2	B	7	7	7	0	0

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.

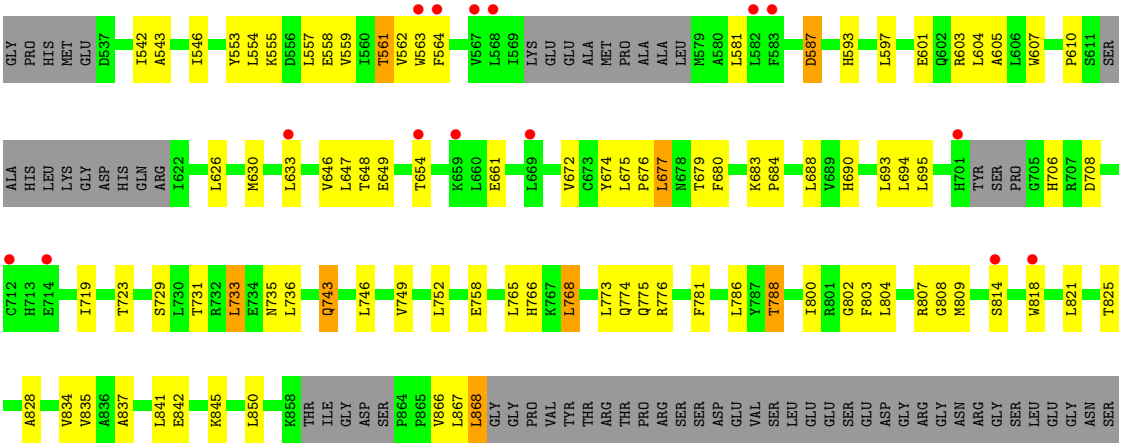
• Molecule 1: FERM, RhoGEF and pleckstrin domain-containing protein 2

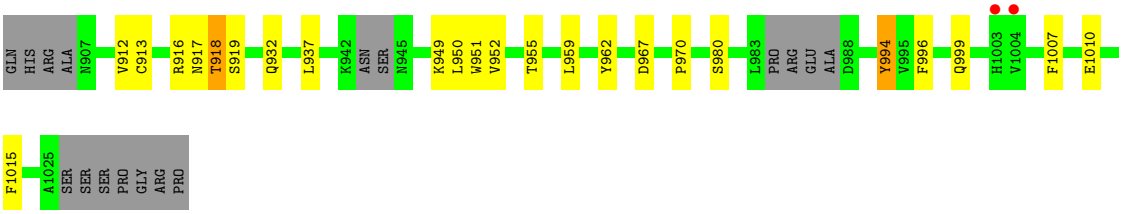
Chain A:



• Molecule 1: FERM, RhoGEF and pleckstrin domain-containing protein 2

Chain B:




GLN
HIS
ARG
ALA
N907
V912
C913
R916
N917
T918
S919
Q932
L937
K942
ASN
SER
N945
K949
L950
W951
V952
T955
L959
Y962
D967
P970
S980
L983
PRO
ARG
GLU
ALA
D988
Y994
V995
F996
Q999
H1003
V1004
F1007
E1010
F1015
A1025
SER
SER
SER
PRO
GLY
ARG
PRO

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	184.39Å 85.25Å 103.45Å 90.00° 118.76° 90.00°	Depositor
Resolution (Å)	45.34 – 3.20 45.34 – 3.17	Depositor EDS
% Data completeness (in resolution range)	92.4 (45.34-3.20) 92.1 (45.34-3.17)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.26 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.7.2_869	Depositor
R, R_{free}	0.232 , 0.278 0.225 , 0.276	Depositor DCC
R_{free} test set	1167 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	70.5	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 87.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6724	wwPDB-VP
Average B, all atoms (Å^2)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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	0.39	0/3543	0.55	1/4806 (0.0%)
1	B	0.34	0/3315	0.53	2/4505 (0.0%)
All	All	0.37	0/6858	0.54	3/9311 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	610	PRO	N-CA-CB	5.74	110.19	103.30
1	A	551	ARG	NE-CZ-NH1	-5.23	117.69	120.30
1	B	868	LEU	CA-CB-CG	5.20	127.25	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	3463	0	3339	64	0
1	B	3242	0	3062	74	0
2	A	12	0	0	0	0
2	B	7	0	0	0	0
All	All	6724	0	6401	137	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 (137) 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:554:LEU:HD11	1:A:593:HIS:HB3	1.61	0.80
1:B:546:ILE:HG12	1:B:694:LEU:HD13	1.65	0.79
1:A:816:ASN:HB3	1:A:818:TRP:HD1	1.53	0.73
1:A:937:LEU:HD12	1:A:1007:PHE:HB3	1.70	0.73
1:B:937:LEU:HD11	1:B:952:VAL:HG21	1.70	0.72
1:B:746:LEU:HD22	1:B:800:ILE:HD11	1.76	0.68
1:A:554:LEU:HD21	1:A:594:ARG:HA	1.77	0.67
1:A:542:ILE:HG21	1:A:698:LEU:HG	1.77	0.64
1:B:937:LEU:HD13	1:B:1007:PHE:HD2	1.63	0.64
1:A:555:LYS:NZ	1:A:967:ASP:O	2.20	0.64
1:B:786:LEU:HD22	1:B:800:ILE:HG21	1.79	0.64
1:B:994:TYR:HB3	1:B:1015:PHE:HB2	1.81	0.63
1:A:777:MET:H	1:A:788:THR:HB	1.64	0.62
1:B:647:LEU:HD23	1:B:733:LEU:HD22	1.85	0.58
1:B:932:GLN:HG2	1:B:955:THR:HG22	1.85	0.58
1:B:601:GLU:O	1:B:605:ALA:N	2.37	0.58
1:A:603:ARG:NH1	1:A:621:ARG:O	2.32	0.58
1:B:951:TRP:HB3	1:B:962:TYR:HB2	1.87	0.57
1:B:802:GLY:HA2	1:B:868:LEU:HD11	1.87	0.57
1:B:675:LEU:HB2	1:B:680:PHE:CE1	2.40	0.56
1:B:735:ASN:HD21	1:B:758:GLU:HA	1.71	0.56
1:A:550:GLU:OE1	1:A:691:TYR:OH	2.22	0.56
1:A:867:LEU:O	1:A:868:LEU:HD23	2.06	0.55
1:B:776:ARG:HD2	1:B:788:THR:HG22	1.88	0.55
1:A:760:ILE:HB	1:A:780:LEU:HG	1.88	0.55
1:B:867:LEU:O	1:B:868:LEU:HD23	2.07	0.55
1:B:654:THR:HG23	1:B:661:GLU:HB2	1.89	0.55
1:A:547:LEU:HG	1:A:551:ARG:HD2	1.88	0.54
1:A:932:GLN:HG2	1:A:955:THR:HG22	1.89	0.54
1:B:706:HIS:CD2	1:B:708:ASP:HB2	2.43	0.54
1:A:666:GLU:HA	1:A:669:LEU:HD12	1.89	0.54
1:A:818:TRP:HZ2	1:B:773:LEU:HG	1.73	0.53
1:B:676:PRO:HB2	1:B:679:THR:HG23	1.91	0.53
1:B:804:LEU:HD13	1:B:828:ALA:HB2	1.91	0.53
1:A:599:GLU:OE1	1:A:628:ARG:NH1	2.41	0.53
1:A:765:LEU:HB2	1:A:778:PHE:CE2	2.43	0.52
1:B:962:TYR:HD2	1:B:967:ASP:HB2	1.74	0.52
1:A:626:LEU:O	1:A:630:MET:HG2	2.10	0.52
1:B:912:VAL:HG13	1:B:916:ARG:HG3	1.91	0.52
1:B:749:VAL:HG13	1:B:800:ILE:HG13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:683:LYS:N	1:B:684:PRO:HD2	2.26	0.51
1:B:564:PHE:CE1	1:B:677:LEU:HD11	2.46	0.51
1:B:554:LEU:HD21	1:B:593:HIS:HB3	1.93	0.51
1:A:740:THR:HA	1:A:753:ILE:HD11	1.94	0.50
1:B:543:ALA:HB1	1:B:604:LEU:HD21	1.94	0.50
1:B:587:ASP:N	1:B:587:ASP:OD1	2.45	0.50
1:B:809:MET:HE1	1:B:850:LEU:HD22	1.95	0.49
1:A:588:PRO:HG2	1:A:639:TYR:CE1	2.47	0.49
1:A:719:ILE:HD12	1:A:722:VAL:HB	1.95	0.49
1:A:540:TYR:CZ	1:A:544:LYS:HD2	2.47	0.49
1:A:761:ARG:HH11	1:A:919:SER:HB2	1.78	0.49
1:A:638:SER:O	1:A:642:ARG:NH2	2.45	0.49
1:A:786:LEU:HD23	1:A:803:PHE:HB3	1.95	0.49
1:A:629:ASN:O	1:A:633:LEU:N	2.45	0.48
1:A:702:TYR:HB3	1:A:709:TYR:HD1	1.78	0.48
1:B:674:TYR:CD2	1:B:675:LEU:HG	2.48	0.48
1:B:825:THR:HG23	1:B:834:VAL:HG22	1.94	0.48
1:B:603:ARG:O	1:B:607:TRP:HB2	2.13	0.48
1:A:777:MET:HG3	1:A:798:PHE:CE1	2.49	0.48
1:A:553:TYR:HE2	1:A:593:HIS:CE1	2.31	0.48
1:B:918:THR:OG1	1:B:919:SER:N	2.47	0.48
1:B:786:LEU:HD23	1:B:803:PHE:HB3	1.95	0.48
1:A:818:TRP:HB3	1:A:823:CYS:SG	2.54	0.47
1:B:554:LEU:HD12	1:B:597:LEU:HD22	1.95	0.47
1:A:683:LYS:N	1:A:684:PRO:HD2	2.30	0.47
1:B:733:LEU:HD23	1:B:733:LEU:HA	1.71	0.47
1:B:980:SER:HB3	1:B:999:GLN:HG3	1.95	0.47
1:B:996:PHE:CE1	1:B:1007:PHE:HB2	2.50	0.47
1:A:649:GLU:OE1	1:A:652:LYS:NZ	2.39	0.47
1:B:646:VAL:HA	1:B:649:GLU:HG2	1.96	0.46
1:B:633:LEU:HD22	1:B:688:LEU:HD11	1.97	0.46
1:A:666:GLU:O	1:A:669:LEU:HB2	2.15	0.46
1:A:688:LEU:HD12	1:A:719:ILE:HD11	1.96	0.46
1:B:809:MET:CE	1:B:850:LEU:HD22	2.45	0.46
1:A:736:LEU:O	1:A:740:THR:HB	2.16	0.45
1:B:558:GLU:O	1:B:562:VAL:HB	2.17	0.45
1:B:581:LEU:HD11	1:B:649:GLU:HG3	1.98	0.45
1:B:814:SER:HB3	1:B:818:TRP:HB2	1.99	0.45
1:B:626:LEU:HD13	1:B:695:LEU:HD21	1.99	0.45
1:A:908:THR:O	1:A:912:VAL:HG23	2.17	0.44
1:A:927:ALA:O	1:A:931:ASN:ND2	2.43	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:626:LEU:O	1:B:630:MET:HG2	2.17	0.44
1:A:850:LEU:O	1:A:854:ILE:HG13	2.16	0.44
1:A:718:ALA:O	1:A:722:VAL:HG23	2.18	0.44
1:B:768:LEU:HD12	1:B:834:VAL:HG11	1.99	0.43
1:A:1018:TRP:O	1:A:1022:ILE:HG13	2.18	0.43
1:A:644:ASP:HB2	1:A:732:ARG:HH21	1.83	0.43
1:B:719:ILE:O	1:B:723:THR:HG23	2.18	0.43
1:B:781:PHE:HE2	1:B:786:LEU:HD12	1.84	0.43
1:A:775:GLN:NE2	1:A:914:TRP:O	2.48	0.43
1:B:765:LEU:HB3	1:B:835:VAL:CG2	2.48	0.43
1:A:809:MET:HB2	1:A:827:TYR:O	2.18	0.43
1:A:540:TYR:CZ	1:A:608:GLU:HG3	2.54	0.43
1:B:555:LYS:O	1:B:559:VAL:HG23	2.19	0.43
1:B:837:ALA:HB1	1:B:842:GLU:HB3	1.99	0.43
1:A:850:LEU:HD23	1:A:850:LEU:HA	1.88	0.42
1:A:1016:GLU:O	1:A:1020:ASP:HB2	2.19	0.42
1:A:816:ASN:CG	1:A:817:GLU:H	2.23	0.42
1:A:595:GLY:O	1:A:599:GLU:HG3	2.20	0.42
1:A:866:VAL:O	1:A:867:LEU:HD23	2.19	0.42
1:B:563:TRP:CZ2	1:B:672:VAL:HG21	2.54	0.42
1:A:777:MET:HG3	1:A:798:PHE:HE1	1.84	0.42
1:B:542:ILE:O	1:B:546:ILE:HG13	2.19	0.42
1:B:913:CYS:HB3	1:B:918:THR:O	2.19	0.42
1:B:729:SER:O	1:B:733:LEU:HB2	2.20	0.42
1:B:866:VAL:O	1:B:867:LEU:HD23	2.20	0.42
1:B:557:LEU:O	1:B:561:THR:HG23	2.20	0.42
1:B:559:VAL:HG22	1:B:672:VAL:HG12	2.01	0.42
1:A:627:LEU:HD21	1:A:714:GLU:HB3	2.01	0.42
1:A:768:LEU:HA	1:A:768:LEU:HD12	1.86	0.42
1:A:773:LEU:HA	1:A:773:LEU:HD23	1.73	0.42
1:A:807:ARG:HA	1:A:808:GLY:HA2	1.77	0.42
1:A:702:TYR:CE2	1:A:708:ASP:HB3	2.55	0.41
1:B:807:ARG:HA	1:B:808:GLY:HA2	1.51	0.41
1:A:638:SER:O	1:A:642:ARG:HD2	2.20	0.41
1:A:811:VAL:HG11	1:A:850:LEU:HD13	2.02	0.41
1:B:690:HIS:HA	1:B:693:LEU:HD12	2.02	0.41
1:B:743:GLN:HG2	1:B:752:LEU:HB3	2.02	0.41
1:B:841:LEU:O	1:B:845:LYS:HG3	2.19	0.41
1:B:821:LEU:HD23	1:B:821:LEU:HA	1.83	0.41
1:A:786:LEU:HD23	1:A:786:LEU:HA	1.80	0.41
1:B:554:LEU:HD23	1:B:557:LEU:HD12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:949:LYS:O	1:B:949:LYS:HG3	2.18	0.41
1:A:546:ILE:HG12	1:A:694:LEU:HD13	2.02	0.41
1:B:554:LEU:HD23	1:B:554:LEU:HA	1.71	0.41
1:B:773:LEU:HD23	1:B:773:LEU:HA	1.92	0.41
1:B:766:HIS:CD2	1:B:775:GLN:HA	2.56	0.41
1:B:774:GLN:OE1	1:B:776:ARG:NH1	2.41	0.41
1:B:959:LEU:HD23	1:B:959:LEU:HA	1.90	0.41
1:B:553:TYR:HE2	1:B:593:HIS:CE1	2.38	0.41
1:B:563:TRP:NE1	1:B:672:VAL:HG11	2.36	0.41
1:B:962:TYR:OH	1:B:970:PRO:HG3	2.21	0.41
1:A:644:ASP:HA	1:A:733:LEU:HD13	2.02	0.40
1:A:606:LEU:HB3	1:A:619:HIS:ND1	2.36	0.40
1:A:981:VAL:HG11	1:A:1019:MET:HB2	2.02	0.40
1:A:652:LYS:O	1:A:656:HIS:ND1	2.48	0.40
1:A:1009:ALA:HB1	1:A:1014:THR:HG23	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	422/501 (84%)	392 (93%)	29 (7%)	1 (0%)	47	79
1	B	402/501 (80%)	363 (90%)	38 (10%)	1 (0%)	47	79
All	All	824/1002 (82%)	755 (92%)	67 (8%)	2 (0%)	47	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	607	TRP
1	B	1010	GLU

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	352/444 (79%)	339 (96%)	13 (4%)	34	68
1	B	317/444 (71%)	303 (96%)	14 (4%)	28	64
All	All	669/888 (75%)	642 (96%)	27 (4%)	31	66

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

Mol	Chain	Res	Type
1	A	545	GLU
1	A	633	LEU
1	A	650	LEU
1	A	664	TYR
1	A	673	CYS
1	A	684	PRO
1	A	708	ASP
1	A	740	THR
1	A	750	GLU
1	A	761	ARG
1	A	788	THR
1	A	821	LEU
1	A	838	SER
1	B	561	THR
1	B	587	ASP
1	B	648	THR
1	B	677	LEU
1	B	731	THR
1	B	733	LEU
1	B	736	LEU
1	B	743	GLN
1	B	768	LEU
1	B	788	THR
1	B	917	ASN
1	B	918	THR
1	B	950	LEU
1	B	994	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	643	HIS
1	A	925	HIS
1	B	766	HIS

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	436/501 (87%)	-0.27	4 (0%)	84	75	54, 90, 142, 181	0
1	B	418/501 (83%)	0.02	17 (4%)	37	24	58, 111, 176, 203	0
All	All	854/1002 (85%)	-0.13	21 (2%)	57	43	54, 98, 164, 203	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	712	CYS	4.8
1	B	564	PHE	4.6
1	B	714	GLU	3.6
1	B	582	LEU	3.1
1	B	814	SER	3.1
1	B	669	LEU	3.1
1	A	702	TYR	3.0
1	B	633	LEU	2.9
1	B	659	LYS	2.7
1	B	701	HIS	2.6
1	B	1003	HIS	2.6
1	B	1004	VAL	2.5
1	B	583	PHE	2.5
1	B	818	TRP	2.4
1	B	563	TRP	2.4
1	B	568	LEU	2.3
1	A	704	PRO	2.3
1	B	654	THR	2.2
1	B	567	VAL	2.2
1	A	612	SER	2.1
1	A	995	VAL	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.